

Synthesis, Characterization, Antimicrobial, DFT and Molecular Docking Studies of novel Mn(II), Fe(III) and Cr(III) Complexes Incorporating 4-(2-hydroxyphenyl azo)-1-naphthol (Az)

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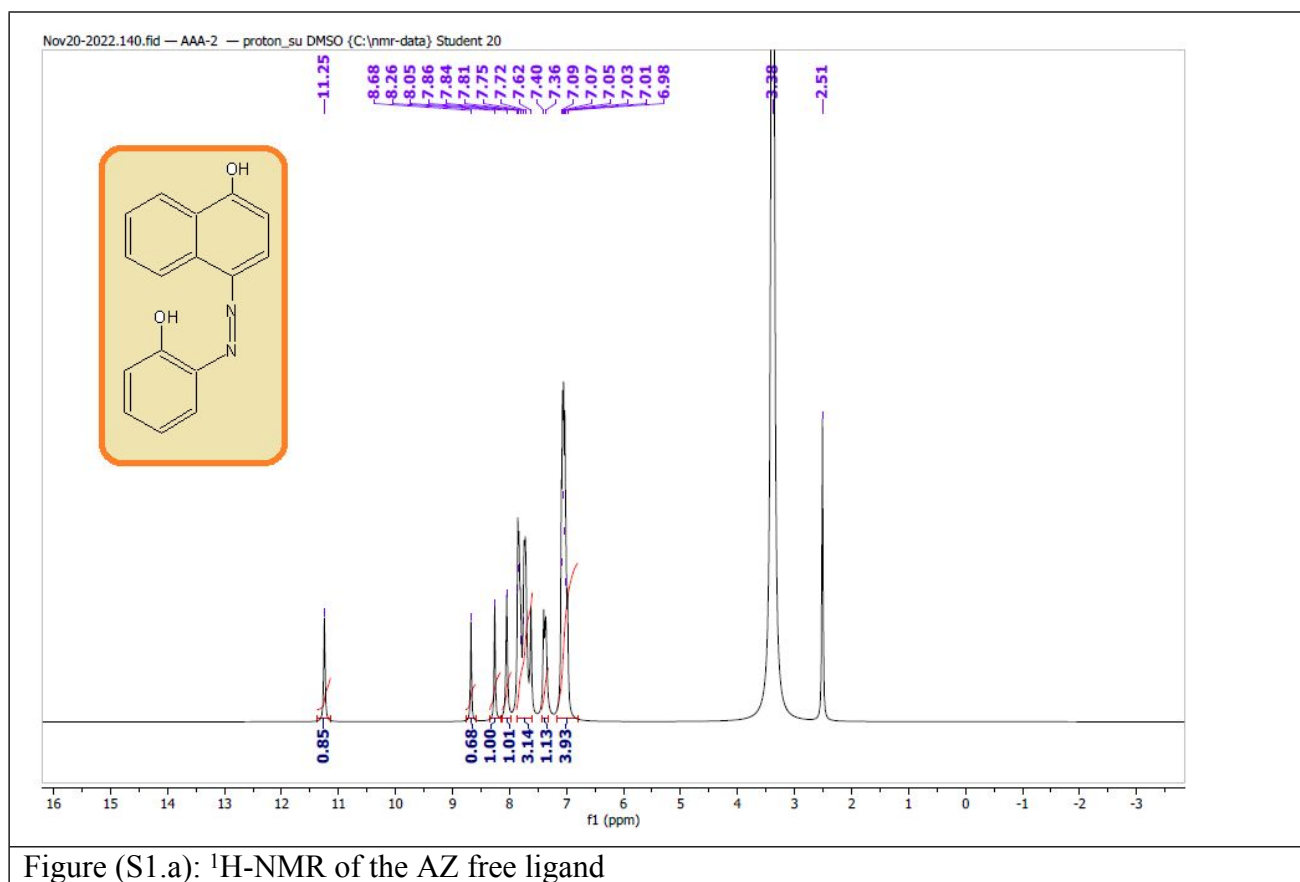
1. Instruments for Characterization

The compositions of the complexes were determined using elemental analysis (for C, H, and N content), molar conductance (10^{-3} mol/L DMF and acetonitrile solution), thermal decomposition (TGA), magnetic measurements, infrared (IR) spectroscopy, and electronic spectra (10^{-3} mol/L acetonitrile solution). Elemental analysis was performed at the University of Cairo's Central Laboratory using a Perkin-Elmer 2408 analyzer. Infrared spectra of KBr pellets were obtained using a Shimadzu DR-8001 spectrometer and an infrared spectrometer. $^1\text{H-NMR}$ spectra of DMSO- d_6 were obtained using a Bruker DRX 400 MHz instrument with TMS as a reference. Mass spectra were carried out using Direct Inlet part to the mass analyzer in Thermo Scientific GCMS model ISQ. UV-Vis spectra were obtained using a Jenway UV-Vis spectrophotometer. Thermal decomposition (TGA) analysis was performed on the required compounds using a Shimadzu type 60 H analyzer. The molar conductance of a DMF and acetonitrile solution (10^{-3} mol/L) was measured using a JENWAY model 4320 electronic conductivity meter. The molar magnetic susceptibility of powdered materials was determined using a Bartington Susceptibility machine (model 4320). The stoichiometry of the compounds was determined using a continuous-variation spectrophotometric method.

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Figure (S1.a): ^1H -NMR of the AZ free ligand

Figure (S1.b): ^{13}C -NMR of the AZ free ligand



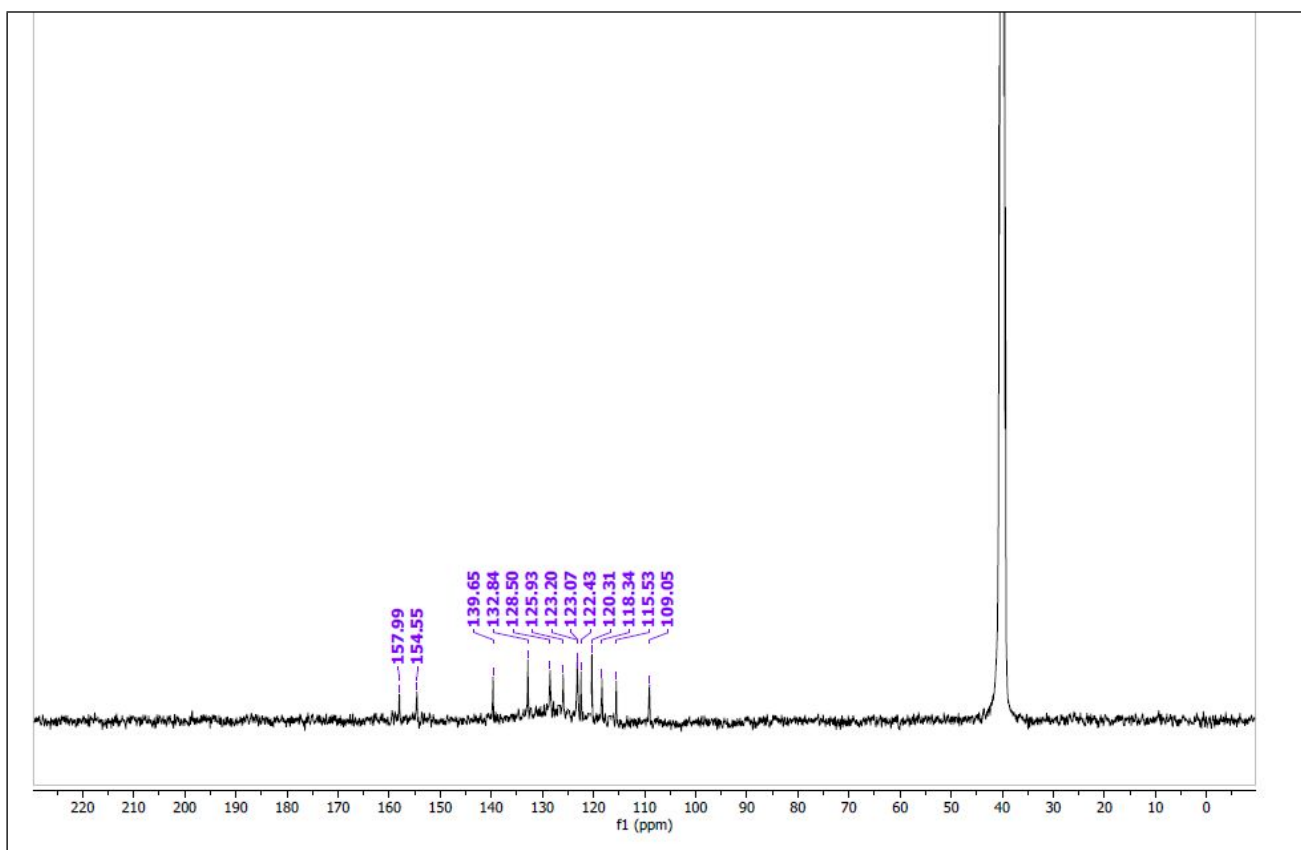


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Table (S10): Summary of Natural Population Analysis of the Fe(III) complex:

Table (S11): NBO data of the Fe(III) complex:

Table (S1): Binding energy of the studied complexes based on DFT/B3LYP with cc-PVQZ basis set in ethanol as solvent.

	E _{Total}	Binding
Az	-29028.12	
CrAz₂	-117049.33	-475.78
MnAz₂	-112342.85	-589.06
FeAz₂	-123228.14	-692.26
Cr(III)	-33569.30	
Mn(II)	-39198.19	
Fe(III)	-39531.63	
H₂O	-7249.68	
Cl(-1)	-17698.33	

Table (S2): Comparison between the minimum inhibition concentrations (MIC) of the current compounds with previously reported compounds in the literature survey against <i>E. coli</i> .			
	E. coli (G-)		
		MIC	
1	CrHPN	12.5	This work
2	MnHPN	6.25	
3	FeHPN	6.25	
4	Co:FMBT (1:1)	100.00	1
5	Co:FMBT (1:2)	125.00	
6	Ni:FMBT (1:1)	125.00	
7	Ni:FMBT (1:2)	62.50	
8	Pd:FMBT (1:1)	125.00	
9	Pd:FMBT (1:2)	100.00	
10	FeLQ	25.50	2
11	FeLB	22.50	
12	FeLI	23.50	
13	CuLQ	21.50	3
14	CuLB	22.50	
15	CuLI	22.50	
16	NiLQ	35.50	4
17	NiLB	34.50	
18	NiLI	37.50	
19	FeL	50.00	5
20	CoL	55.00	
21	NiL	60.00	
22	CuL	50.00	
23	ZnL	70.00	
24	3a	25.20	6
25	3b	78.50	
26	3c	18.20	
27	3e	15.40	
28	[Ag(3-Py-CHO)2NO3]	62.50	7
29	[Pd(HL)(Cl2)].H2O	12.50	8
30	[Pt(HL)(Cl2)].2H2O	25.00	

¹ A. Skotnicka, E. Kolehmainen, P. Czelen, A. Valkonen, R. Gawinecki, Synthesis and structural characterization of substituted 2-phenacylbenzoxazoles. *Int. J. Mol. Sci.* 14 (2013) 4444–4460. <https://doi.org/10.3390/ijms14034444>.

² M. Rahimizadeh, H. Eshghi, A. Shiri, Z. Ghadamyari, M.M. Matin, F. Oroojalian, P. Pordeli, Fe(HSO₄)₃ as an Efficient Catalyst for Diazotization and Diazo Coupling Reactions. *J. Korean Chem. Soc.* 56 (2012) 716-719. <https://doi.org/10.5012/jkcs.2012.56.6.716>.

³ A. Shomali, H. Valizadeh, S. Noorshargh, New Generation of Nitrite Functionalized Star-like Polyvinyl Imidazolium Compound: Application as a Nitrosonium Source and Three Dimensional Nanocatalyst for the Synthesis of Azo Dyes. *Lett. Org. Chem.* 14 (2017) 14, 409–418. DOI: <https://doi.org/10.2174/1570178614666170321125216>.

- ⁴ M. Akki, D.S. Reddy, K.S. Katagi, A. Kumar, H.C. Devarajegowda, V. Babagond, S. Mane, S.D. Joshi, Synthesis of coumarin-thioether conjugates as potential anti-tubercular agents: Their molecular docking and X-ray crystal studies. *J. Mol. Struct.* 1266 (2022) 133452. <https://doi.org/10.1016/j.molstruc.2022.133452>.
- ⁵ F. Perveen, R. Qureshi, F.L. Ansari, S. Kalsoom, S. Ahmed, Investigations of drug–DNA interactions using molecular docking, cyclic voltammetry and UV–Vis spectroscopy. *J. Mol. Struct.* 1004 (2011) 67–73. <https://doi.org/10.1016/j.molstruc.2011.07.027>.
- ⁶ X.-Y. Meng, H.-X. Zhang, M. Mezei, M. Cui, Molecular Docking: A Powerful Approach for Structure-Based Drug Discovery. *Curr. Comput. Aided Drug Des.* 7 (2011) 146–157. DOI: <https://doi.org/10.2174/157340911795677602>.
- ⁷ M. Akki, D.S. Reddy, K.S. Katagi, A. Kumar, H.C. Devarajegowda, V. Babagond, S. Mane, S.D. Joshi, Synthesis of coumarin-thioether conjugates as potential anti-tubercular agents: Their molecular docking and X-ray crystal studies. *J. Mol. Struct.* 1266 (2022) 133452. <https://doi.org/10.1016/j.molstruc.2022.133452>.
- ⁸ R.G. Deghadi, A.A. Abbas, G.G. Mohamed, Theoretical and experimental investigations of new bis(amino triazole) Schiff base ligand: Preparation of its UO₂(II), Er (III) and La (III) complexes, studying of their antibacterial, anticancer and molecular docking. *Appl. Organomet. Chem.* 35 (2021) e6292. <https://doi.org/10.1002/aoc.6292>.

Table (S3): The bond parameters (bond length, bond angles, and dihedral angles) of the free ligand and its metal complexes							
Ligand		Cr(III) complex		Mn(II) complex		Fe(III) complex	
Bond Length	value	Bond Length	value	Bond Length	value	Bond Length	value
O(20)-H(31)	0.9495	Cl(66)-Cr(67)	2.2794	O(68)-Mn(69)	1.9131	Cl(66)-Fe(67)	2.2282
O(19)-H(32)	0.9501	O(65)-Cr(67)	1.9235	H(67)-O(68)	0.9892	O(65)-Fe(67)	1.8708
C(18)-H(30)	1.0952	O(62)-Cr(67)	1.9218	H(66)-O(68)	0.9888	O(62)-Fe(67)	1.8652
C(17)-H(29)	1.0962	O(62)-H(64)	0.9876	O(65)-Mn(69)	1.9237	O(62)-H(64)	0.9881
C(17)-C(18)	1.3857	O(62)-H(63)	0.9896	O(62)-Mn(69)	1.9247	O(62)-H(63)	0.9907
C(16)-H(28)	1.0948	O(50)-H(60)	1.0079	O(62)-H(64)	0.9879	O(50)-H(60)	1.0073
C(16)-C(18)	1.3943	C(49)-H(59)	1.0824	O(62)-H(63)	0.9866	C(49)-H(59)	1.0823
C(15)-O(19)	1.3618	C(48)-H(58)	1.0814	O(50)-H(60)	1.008	C(48)-H(58)	1.0825
C(15)-C(17)	1.4049	C(48)-C(49)	1.3368	C(49)-H(59)	1.0824	C(48)-C(49)	1.337
C(14)-H(27)	1.0977	C(47)-H(57)	1.0824	C(48)-H(58)	1.0827	C(47)-H(57)	1.0823
C(14)-C(16)	1.3866	C(47)-C(49)	1.4731	C(48)-C(49)	1.3369	C(47)-C(49)	1.4745
C(13)-C(15)	1.4126	C(46)-O(65)	1.3936	C(47)-H(57)	1.0823	C(46)-O(65)	1.3918
C(13)-C(14)	1.4013	C(46)-C(48)	1.4813	C(47)-C(49)	1.4737	C(46)-C(48)	1.4801
N(12)-C(13)	1.4422	C(45)-H(56)	1.0829	C(46)-O(65)	1.3918	C(45)-H(56)	1.0827
N(11)-N(12)	1.234	C(45)-C(47)	1.3371	C(46)-C(48)	1.4805	C(45)-C(47)	1.3379
C(10)-H(26)	1.0952	C(44)-C(46)	1.3524	C(45)-H(56)	1.0828	C(44)-C(46)	1.3511
C(9)-H(25)	1.0972	C(44)-C(45)	1.482	C(45)-C(47)	1.3375	C(44)-C(45)	1.4814
C(9)-C(10)	1.3696	N(43)-C(44)	1.4304	C(44)-C(46)	1.352	N(43)-C(44)	1.4314
C(8)-H(24)	1.0948	N(42)-Cr(67)	2.001	C(44)-C(45)	1.4819	N(42)-Fe(67)	1.9384
C(8)-C(10)	1.4111	N(42)-N(43)	1.2712	N(43)-C(44)	1.4313	N(42)-N(43)	1.2723
C(7)-H(23)	1.0962	C(41)-H(55)	1.0824	N(42)-Mn(69)	2.0063	C(41)-H(55)	1.0825
C(7)-C(8)	1.3692	C(40)-H(54)	1.0822	N(42)-N(43)	1.2716	C(40)-H(54)	1.0816
C(6)-N(11)	1.4418	C(40)-C(41)	1.3352	C(41)-H(55)	1.0824	C(40)-C(41)	1.3349
C(5)-H(22)	1.0977	C(39)-H(53)	1.0825	C(40)-H(54)	1.0816	C(39)-H(53)	1.0826
C(5)-C(6)	1.3798	C(39)-C(41)	1.4716	C(40)-C(41)	1.3352	C(39)-C(41)	1.4709
C(4)-C(9)	1.4216	C(38)-H(52)	1.0788	C(39)-H(53)	1.0826	C(38)-H(52)	1.0783
C(4)-C(6)	1.4355	C(38)-C(39)	1.3374	C(39)-C(41)	1.4716	C(38)-C(39)	1.3372
C(3)-H(21)	1.0965	C(37)-N(42)	1.4324	C(38)-H(52)	1.0789	C(37)-N(42)	1.4328
C(3)-C(5)	1.4062	C(36)-H(61)	1.0824	C(38)-C(39)	1.3373	C(36)-H(61)	1.0796
C(2)-C(4)	1.4136	C(36)-C(37)	1.3383	C(37)-N(42)	1.4328	C(36)-C(37)	1.3371
C(2)-C(7)	1.4206	C(35)-C(40)	1.4869	C(36)-H(61)	1.0821	C(35)-C(40)	1.4874
C(1)-O(20)	1.3647	C(35)-C(37)	1.4904	C(36)-C(37)	1.3386	C(35)-C(37)	1.4918
C(1)-C(3)	1.3818	C(34)-H(51)	1.083	C(35)-C(40)	1.4864	C(34)-H(51)	1.083
C(1)-C(2)	1.4337	C(34)-C(36)	1.4713	C(35)-C(37)	1.4903	C(34)-C(36)	1.4704
Bond Angles		C(33)-C(35)	1.3604	C(34)-H(51)	1.0829	C(33)-C(35)	1.3611
H(31)-O(20)-C(1)	108.0501	C(33)-C(38)	1.4922	C(34)-C(36)	1.4714	C(33)-C(38)	1.4928
H(32)-O(19)-C(15)	108.8189	C(32)-O(50)	1.3987	C(33)-C(35)	1.3603	C(32)-O(50)	1.399
H(30)-C(18)-C(17)	119.6629	C(32)-C(34)	1.3388	C(33)-C(38)	1.4921	C(32)-C(34)	1.3384
H(30)-C(18)-C(16)	119.7381	C(32)-C(33)	1.5016	C(32)-O(50)	1.3986	C(32)-C(33)	1.5027
C(17)-C(18)-C(16)	120.599	O(20)-H(30)	1.0079	C(32)-C(34)	1.3386	O(20)-H(30)	1.0075
H(29)-C(17)-C(18)	120.4956	O(19)-Cr(67)	1.9341	C(32)-C(33)	1.5014	O(19)-Fe(67)	1.885
H(29)-C(17)-C(15)	119.9881	C(18)-H(29)	1.0823	O(20)-H(30)	1.0076	C(18)-H(29)	1.0824
C(18)-C(17)-C(15)	119.5163	C(17)-H(28)	1.0827	O(19)-Mn(69)	1.9331	C(17)-H(28)	1.0826
H(28)-C(16)-C(18)	119.7716	C(17)-C(18)	1.3368	C(18)-H(29)	1.0824	C(17)-C(18)	1.3379
H(28)-C(16)-C(14)	119.7977	C(16)-H(27)	1.0824	C(17)-H(28)	1.0827	C(16)-H(27)	1.0822
C(18)-C(16)-C(14)	120.4307	C(16)-C(18)	1.4729	C(17)-C(18)	1.3367	C(16)-C(18)	1.4754

O(19)-C(15)-C(17)	115.3294	C(15)-O(19)	1.3926	C(16)-H(27)	1.0824	C(15)-O(19)	1.3908
O(19)-C(15)-C(13)	124.5944	C(15)-C(17)	1.4811	C(16)-C(18)	1.473	C(15)-C(17)	1.4803
C(17)-C(15)-C(13)	120.0762	C(14)-H(26)	1.0829	C(15)-O(19)	1.3942	C(14)-H(26)	1.0827
H(27)-C(14)-C(16)	119.4703	C(14)-C(16)	1.3372	C(15)-C(17)	1.481	C(14)-C(16)	1.3383
H(27)-C(14)-C(13)	120.4325	C(13)-C(15)	1.3504	C(14)-H(26)	1.0828	C(13)-C(15)	1.3437
C(16)-C(14)-C(13)	120.0971	C(13)-C(14)	1.4823	C(14)-C(16)	1.3372	C(13)-C(14)	1.4809
C(15)-C(13)-C(14)	119.2806	N(12)-C(13)	1.4288	C(13)-C(15)	1.3526	N(12)-C(13)	1.4259
C(15)-C(13)-N(12)	116.4941	N(11)-Cr(67)	2.0106	C(13)-C(14)	1.4822	N(11)-Fe(67)	1.9549
C(14)-C(13)-N(12)	124.2253	N(11)-N(12)	1.2712	N(12)-C(13)	1.4299	N(11)-N(12)	1.2717
C(13)-N(12)-N(11)	119.7143	C(10)-H(25)	1.0825	N(11)-Mn(69)	2.0009	C(10)-H(25)	1.0824
N(12)-N(11)-C(6)	120.2291	C(9)-H(24)	1.0823	N(11)-N(12)	1.2706	C(9)-H(24)	1.082
H(26)-C(10)-C(9)	120.3505	C(9)-C(10)	1.3352	C(10)-H(25)	1.0825	C(9)-C(10)	1.335
H(26)-C(10)-C(8)	119.1607	C(8)-H(23)	1.0826	C(9)-H(24)	1.0785	C(8)-H(23)	1.0826
C(9)-C(10)-C(8)	120.4888	C(8)-C(10)	1.4716	C(9)-C(10)	1.3351	C(8)-C(10)	1.4715
H(25)-C(9)-C(10)	119.6605	C(7)-H(22)	1.079	C(8)-H(23)	1.0826	C(7)-H(22)	1.0787
H(25)-C(9)-C(4)	119.6154	C(7)-C(8)	1.3375	C(8)-C(10)	1.4711	C(7)-C(8)	1.3376
C(10)-C(9)-C(4)	120.7241	C(6)-N(11)	1.4315	C(7)-H(22)	1.0787	C(6)-N(11)	1.4312
H(24)-C(8)-C(10)	119.3209	C(5)-H(31)	1.0832	C(7)-C(8)	1.3372	C(5)-H(31)	1.0819
H(24)-C(8)-C(7)	120.4939	C(5)-C(6)	1.3388	C(6)-N(11)	1.4328	C(5)-C(6)	1.3369
C(10)-C(8)-C(7)	120.1852	C(4)-C(9)	1.4855	C(5)-H(31)	1.0826	C(4)-C(9)	1.4868
H(23)-C(7)-C(8)	119.9077	C(4)-C(6)	1.4884	C(5)-C(6)	1.3383	C(4)-C(6)	1.489
H(23)-C(7)-C(2)	119.6249	C(3)-H(21)	1.0829	C(4)-C(9)	1.4851	C(3)-H(21)	1.0829
C(8)-C(7)-C(2)	120.4673	C(3)-C(5)	1.4715	C(4)-C(6)	1.4904	C(3)-C(5)	1.4714
N(11)-C(6)-C(5)	123.4007	C(2)-C(4)	1.3601	C(3)-H(21)	1.0829	C(2)-C(4)	1.3599
N(11)-C(6)-C(4)	116.0376	C(2)-C(7)	1.4922	C(3)-C(5)	1.4706	C(2)-C(7)	1.4924
C(5)-C(6)-C(4)	120.5617	C(1)-O(20)	1.3986	C(2)-C(4)	1.361	C(1)-O(20)	1.3988
H(22)-C(5)-C(6)	121.4636	C(1)-C(3)	1.3392	C(2)-C(7)	1.4925	C(1)-C(3)	1.3392
H(22)-C(5)-C(3)	117.8788	C(1)-C(2)	1.5017	C(1)-O(20)	1.3988	C(1)-C(2)	1.5023
C(6)-C(5)-C(3)	120.6576	Bond Angles		C(1)-C(3)	1.3384	Bond Angles	
C(9)-C(4)-C(6)	122.2952	Cl(66)-Cr(67)-O(65)	89.345	C(1)-C(2)	1.5022	Cl(66)-Fe(67)-O(65)	74.7263
C(9)-C(4)-C(2)	118.6136	Cl(66)-Cr(67)-O(62)	88.5937	Bond Angles		Cl(66)-Fe(67)-O(62)	71.548
C(6)-C(4)-C(2)	119.0912	Cl(66)-Cr(67)-N(42)	88.9102	O(68)-Mn(69)-O(65)	89.7527	Cl(66)-Fe(67)-N(42)	90.1183
H(21)-C(3)-C(5)	118.8444	Cl(66)-Cr(67)-O(19)	178.1789	O(68)-Mn(69)-O(62)	89.1426	Cl(66)-Fe(67)-O(19)	163.77
H(21)-C(3)-C(1)	121.2686	Cl(66)-Cr(67)-N(11)	91.5894	O(68)-Mn(69)-N(42)	90.7184	Cl(66)-Fe(67)-N(11)	108.9723
C(5)-C(3)-C(1)	119.8871	O(65)-Cr(67)-O(62)	177.9297	O(68)-Mn(69)-O(19)	179.5957	O(65)-Fe(67)-O(62)	144.0427
C(4)-C(2)-C(7)	119.5209	O(65)-Cr(67)-N(42)	89.8324	O(68)-Mn(69)-N(11)	89.3339	O(65)-Fe(67)-N(42)	90.7546
C(4)-C(2)-C(1)	118.6178	O(65)-Cr(67)-O(19)	90.0151	O(65)-Mn(69)-O(62)	178.8762	O(65)-Fe(67)-O(19)	110.6881
C(7)-C(2)-C(1)	121.8613	O(65)-Cr(67)-N(11)	89.2169	O(65)-Mn(69)-N(42)	89.5823	O(65)-Fe(67)-N(11)	85.751
O(20)-C(1)-C(3)	122.3321	O(62)-Cr(67)-N(42)	90.3206	O(65)-Mn(69)-O(19)	89.94	O(62)-Fe(67)-N(42)	101.3284
O(20)-C(1)-C(2)	116.4833	O(62)-Cr(67)-O(19)	92.0511	O(65)-Mn(69)-N(11)	88.918	O(62)-Fe(67)-O(19)	105.1625
C(3)-C(1)-C(2)	121.1846	O(62)-Cr(67)-N(11)	90.6479	O(62)-Mn(69)-N(42)	90.2254	O(62)-Fe(67)-N(11)	93.5875
Dihedral Angles		N(42)-Cr(67)-O(19)	89.383	O(62)-Mn(69)-O(19)	91.1657	N(42)-Fe(67)-O(19)	74.7676
C(15)-C(17)-C(18)-C(16)	-0.0044	N(42)-Cr(67)-N(11)	178.9209	O(62)-Mn(69)-N(11)	91.2752	N(42)-Fe(67)-N(11)	158.8249
C(15)-C(17)-C(18)-H(30)	-179.9962	O(19)-Cr(67)-N(11)	90.1066	N(42)-Mn(69)-O(19)	89.5421	O(19)-Fe(67)-N(11)	86.9391
H(29)-C(17)-C(18)-C(16)	179.9948	Cr(67)-O(65)-C(46)	116.3996	N(42)-Mn(69)-N(11)	178.4992	Fe(67)-O(65)-C(46)	113.8384
H(29)-C(17)-C(18)-H(30)	0.003	Cr(67)-O(62)-H(64)	109.3842	O(19)-Mn(69)-N(11)	90.3976	Fe(67)-O(62)-H(64)	106.804
C(14)-C(16)-C(18)-C(17)	0.0052	Cr(67)-O(62)-H(63)	107.858	Mn(69)-O(68)-H(67)	107.368	Fe(67)-O(62)-H(63)	107.4385
C(14)-C(16)-C(18)-H(30)	179.997	H(64)-O(62)-H(63)	105.3912	Mn(69)-O(68)-H(66)	106.8329	H(64)-O(62)-H(63)	104.0478
H(28)-C(16)-C(18)-C(17)	179.9974	H(60)-O(50)-C(32)	114.9478	H(67)-O(68)-H(66)	105.2546	H(60)-O(50)-C(32)	115.2868
H(28)-C(16)-C(18)-H(30)	-0.0108	H(59)-C(49)-C(48)	120.2013	Mn(69)-O(65)-C(46)	116.244	H(59)-C(49)-C(48)	120.1722
C(13)-C(15)-O(19)-H(32)	-0.0291	H(59)-C(49)-C(47)	119.8797	Mn(69)-O(62)-H(64)	108.3596	H(59)-C(49)-C(47)	119.8962
C(17)-C(15)-O(19)-H(32)	179.975	C(48)-C(49)-C(47)	119.9167	Mn(69)-O(62)-H(63)	108.6036	C(48)-C(49)-C(47)	119.9272
C(13)-C(15)-C(17)-C(18)	0.0039	H(58)-C(48)-C(49)	119.5573	H(64)-O(62)-H(63)	104.6257	H(58)-C(48)-C(49)	119.773
C(13)-C(15)-C(17)-H(29)	-179.9953	H(58)-C(48)-C(46)	119.8136	H(60)-O(50)-C(32)	114.9085	H(58)-C(48)-C(46)	119.7645

O(19)-C(15)-C(17)-C(18)	180	C(49)-C(48)-C(46)	120.6291	H(59)-C(49)-C(48)	120.2001	C(49)-C(48)-C(46)	120.4623
O(19)-C(15)-C(17)-H(29)	0.0008	H(57)-C(47)-C(49)	119.8663	H(59)-C(49)-C(47)	119.8842	H(57)-C(47)-C(49)	119.8419
C(13)-C(14)-C(16)-C(18)	-0.0054	H(57)-C(47)-C(45)	120.1958	C(48)-C(49)-C(47)	119.9147	H(57)-C(47)-C(45)	120.17
C(13)-C(14)-C(16)-H(28)	-179.9976	C(49)-C(47)-C(45)	119.9334	H(58)-C(48)-C(49)	119.7808	C(49)-C(47)-C(45)	119.9814
H(27)-C(14)-C(16)-C(18)	-180	O(65)-C(46)-C(48)	117.566	H(58)-C(48)-C(46)	119.6735	O(65)-C(46)-C(48)	118.0448
H(27)-C(14)-C(16)-H(28)	0.0078	O(65)-C(46)-C(44)	122.8463	C(49)-C(48)-C(46)	120.5448	O(65)-C(46)-C(44)	122.0893
N(12)-C(13)-C(15)-C(17)	-179.9952	C(48)-C(46)-C(44)	119.5149	H(57)-C(47)-C(49)	119.8512	C(48)-C(46)-C(44)	119.7132
N(12)-C(13)-C(15)-O(19)	0.0091	H(56)-C(45)-C(47)	119.596	H(57)-C(47)-C(45)	120.1922	H(56)-C(45)-C(47)	119.705
C(14)-C(13)-C(15)-C(17)	-0.0041	H(56)-C(45)-C(44)	119.6968	C(49)-C(47)-C(45)	119.9538	H(56)-C(45)-C(44)	119.7301
C(14)-C(13)-C(15)-O(19)	-179.9998	C(47)-C(45)-C(44)	120.707	O(65)-C(46)-C(48)	117.4855	C(47)-C(45)-C(44)	120.5648
N(12)-C(13)-C(14)-C(16)	179.9952	C(46)-C(44)-C(45)	119.2737	O(65)-C(46)-C(44)	122.7585	C(46)-C(44)-C(45)	119.303
N(12)-C(13)-C(14)-H(27)	-0.0102	C(46)-C(44)-N(43)	123.2844	C(48)-C(46)-C(44)	119.6855	C(46)-C(44)-N(43)	122.5354
C(15)-C(13)-C(14)-C(16)	0.0049	C(45)-C(44)-N(43)	117.1462	H(56)-C(45)-C(47)	119.6119	C(45)-C(44)-N(43)	117.8158
C(15)-C(13)-C(14)-H(27)	179.9994	C(44)-N(43)-N(42)	125.1601	H(56)-C(45)-C(44)	119.6994	C(44)-N(43)-N(42)	123.668
N(11)-N(12)-C(13)-C(14)	0.051	Cr(67)-N(42)-N(43)	120.6673	C(47)-C(45)-C(44)	120.6884	Fe(67)-N(42)-N(43)	120.1676
N(11)-N(12)-C(13)-C(15)	-179.9583	Cr(67)-N(42)-C(37)	121.4338	C(46)-C(44)-C(45)	119.2017	Fe(67)-N(42)-C(37)	121.0363
C(6)-N(11)-N(12)-C(13)	179.9964	N(43)-N(42)-C(37)	117.8395	C(46)-C(44)-N(43)	123.254	N(43)-N(42)-C(37)	118.7947
C(4)-C(9)-C(10)-C(8)	0.0032	H(55)-C(41)-C(40)	120.1347	C(45)-C(44)-N(43)	117.2699	H(55)-C(41)-C(40)	120.142
C(4)-C(9)-C(10)-H(26)	-179.995	H(55)-C(41)-C(39)	119.7741	C(44)-N(43)-N(42)	125.1607	H(55)-C(41)-C(39)	119.7507
H(25)-C(9)-C(10)-C(8)	179.9969	C(40)-C(41)-C(39)	120.0869	Mn(69)-N(42)-N(43)	120.5497	C(40)-C(41)-C(39)	120.1039
H(25)-C(9)-C(10)-H(26)	-0.0012	H(54)-C(40)-C(41)	118.89	Mn(69)-N(42)-C(37)	121.5983	H(54)-C(40)-C(41)	118.6915
C(7)-C(8)-C(10)-C(9)	-0.0046	H(54)-C(40)-C(35)	120.6236	N(43)-N(42)-C(37)	117.7916	H(54)-C(40)-C(35)	120.791
C(7)-C(8)-C(10)-H(26)	179.9936	C(41)-C(40)-C(35)	120.4864	H(55)-C(41)-C(40)	120.134	C(41)-C(40)-C(35)	120.5152
H(24)-C(8)-C(10)-C(9)	179.9989	H(53)-C(39)-C(41)	119.5656	H(55)-C(41)-C(39)	119.7607	H(53)-C(39)-C(41)	119.5655
H(24)-C(8)-C(10)-H(26)	-0.0029	H(53)-C(39)-C(38)	120.0916	C(40)-C(41)-C(39)	120.0946	H(53)-C(39)-C(38)	120.089
C(2)-C(7)-C(8)-C(10)	0.0058	C(41)-C(39)-C(38)	120.3358	H(54)-C(40)-C(41)	118.8317	C(41)-C(39)-C(38)	120.3416
C(2)-C(7)-C(8)-H(24)	-179.9977	H(52)-C(38)-C(39)	117.5795	H(54)-C(40)-C(35)	120.669	H(52)-C(38)-C(39)	117.3864
H(23)-C(7)-C(8)-C(10)	179.9975	H(52)-C(38)-C(33)	121.8448	C(41)-C(40)-C(35)	120.4917	H(52)-C(38)-C(33)	122.0137
H(23)-C(7)-C(8)-H(24)	-0.006	C(39)-C(38)-C(33)	120.558	H(53)-C(39)-C(41)	119.5761	C(39)-C(38)-C(33)	120.591
C(4)-C(6)-N(11)-N(12)	179.8859	N(42)-C(37)-C(36)	119.598	H(53)-C(39)-C(38)	120.0925	N(42)-C(37)-C(36)	119.6227
C(5)-C(6)-N(11)-N(12)	-0.1306	N(42)-C(37)-C(35)	120.654	C(41)-C(39)-C(38)	120.319	N(42)-C(37)-C(35)	120.7777
C(3)-C(5)-C(6)-C(4)	-0.0015	C(36)-C(37)-C(35)	119.5357	H(52)-C(38)-C(39)	117.5907	C(36)-C(37)-C(35)	119.58
C(3)-C(5)-C(6)-N(11)	-179.9843	H(61)-C(36)-C(37)	120.469	H(52)-C(38)-C(33)	121.8385	H(61)-C(36)-C(37)	120.5852
H(22)-C(5)-C(6)-C(4)	179.9977	H(61)-C(36)-C(34)	118.9842	C(39)-C(38)-C(33)	120.5556	H(61)-C(36)-C(34)	118.727
H(22)-C(5)-C(6)-N(11)	0.0148	C(37)-C(36)-C(34)	120.5457	N(42)-C(37)-C(36)	119.6413	C(37)-C(36)-C(34)	120.663
C(2)-C(4)-C(9)-C(10)	-0.003	C(40)-C(35)-C(37)	120.1459	N(42)-C(37)-C(35)	120.7133	C(40)-C(35)-C(37)	120.1708
C(2)-C(4)-C(9)-H(25)	-179.9968	C(40)-C(35)-C(33)	119.7465	C(36)-C(37)-C(35)	119.4752	C(40)-C(35)-C(33)	119.7108
C(6)-C(4)-C(9)-C(10)	-179.9984	C(37)-C(35)-C(33)	120.0893	H(61)-C(36)-C(37)	120.523	C(37)-C(35)-C(33)	120.1044
C(6)-C(4)-C(9)-H(25)	0.0079	H(51)-C(34)-C(36)	118.9444	H(61)-C(36)-C(34)	118.9415	H(51)-C(34)-C(36)	118.8794
C(2)-C(4)-C(6)-C(5)	0.0059	H(51)-C(34)-C(32)	120.0586	C(37)-C(36)-C(34)	120.5286	H(51)-C(34)-C(32)	120.0469
C(2)-C(4)-C(6)-N(11)	179.99	C(36)-C(34)-C(32)	120.9693	C(40)-C(35)-C(37)	120.1688	C(36)-C(34)-C(32)	121.0526
C(9)-C(4)-C(6)-C(5)	-179.9987	C(35)-C(33)-C(38)	118.7694	C(40)-C(35)-C(33)	119.7325	C(35)-C(33)-C(38)	118.7268
C(9)-C(4)-C(6)-N(11)	-0.0147	C(35)-C(33)-C(32)	119.2159	C(37)-C(35)-C(33)	120.0669	C(35)-C(33)-C(32)	119.2047
C(1)-C(3)-C(5)-C(6)	-0.0046	C(38)-C(33)-C(32)	121.9748	H(51)-C(34)-C(36)	118.942	C(38)-C(33)-C(32)	122.0556
C(1)-C(3)-C(5)-H(22)	179.9962	O(50)-C(32)-C(34)	117.4363	H(51)-C(34)-C(32)	120.0615	O(50)-C(32)-C(34)	117.2542
H(21)-C(3)-C(5)-C(6)	-179.9996	O(50)-C(32)-C(33)	123.1914	C(36)-C(34)-C(32)	120.9524	O(50)-C(32)-C(33)	123.3718
H(21)-C(3)-C(5)-H(22)	0.0013	C(34)-C(32)-C(33)	119.371	C(35)-C(33)-C(38)	118.7919	C(34)-C(32)-C(33)	119.3701
C(1)-C(2)-C(4)-C(6)	-0.0043	H(30)-O(20)-C(1)	114.9208	C(35)-C(33)-C(32)	119.2212	H(30)-O(20)-C(1)	115.0838
C(1)-C(2)-C(4)-C(9)	-179.9998	Cr(67)-O(19)-C(15)	119.0678	C(38)-C(33)-C(32)	121.9337	Fe(67)-O(19)-C(15)	115.4443
C(7)-C(2)-C(4)-C(6)	179.9997	H(29)-C(18)-C(17)	120.2228	O(50)-C(32)-C(34)	117.4695	H(29)-C(18)-C(17)	120.1982
C(7)-C(2)-C(4)-C(9)	0.0042	H(29)-C(18)-C(16)	119.9025	O(50)-C(32)-C(33)	123.1755	H(29)-C(18)-C(16)	119.8786
C(1)-C(2)-C(7)-C(8)	179.9984	C(17)-C(18)-C(16)	119.8742	C(34)-C(32)-C(33)	119.3547	C(17)-C(18)-C(16)	119.9208
C(1)-C(2)-C(7)-H(23)	0.0068	H(28)-C(17)-C(18)	119.7222	H(30)-O(20)-C(1)	115.1596	H(28)-C(17)-C(18)	119.8695
C(4)-C(2)-C(7)-C(8)	-0.0057	H(28)-C(17)-C(15)	119.6346	Mn(69)-O(19)-C(15)	118.2087	H(28)-C(17)-C(15)	119.8077

C(4)-C(2)-C(7)-H(23)	-179.9973	C(18)-C(17)-C(15)	120.6422	H(29)-C(18)-C(17)	120.2093	C(18)-C(17)-C(15)	120.322
C(2)-C(1)-O(20)-H(31)	-179.9896	H(27)-C(16)-C(18)	119.878	H(29)-C(18)-C(16)	119.8916	H(27)-C(16)-C(18)	119.863
C(3)-C(1)-O(20)-H(31)	0.0176	H(27)-C(16)-C(14)	120.2081	C(17)-C(18)-C(16)	119.8985	H(27)-C(16)-C(14)	120.1856
C(2)-C(1)-C(3)-C(5)	0.0062	C(18)-C(16)-C(14)	119.9114	H(28)-C(17)-C(18)	119.7239	C(18)-C(16)-C(14)	119.9456
C(2)-C(1)-C(3)-H(21)	-179.999	O(19)-C(15)-C(17)	117.0507	H(28)-C(17)-C(15)	119.6187	O(19)-C(15)-C(17)	118.6547
O(20)-C(1)-C(3)-C(5)	179.9986	O(19)-C(15)-C(13)	123.3288	C(18)-C(17)-C(15)	120.6561	O(19)-C(15)-C(13)	121.4124
O(20)-C(1)-C(3)-H(21)	-0.0066	C(17)-C(15)-C(13)	119.5958	H(27)-C(16)-C(18)	119.8654	C(17)-C(15)-C(13)	119.8125
C(3)-C(1)-C(2)-C(7)	179.9942	H(26)-C(14)-C(16)	119.5755	H(27)-C(16)-C(14)	120.2008	H(26)-C(14)-C(16)	119.7884
C(3)-C(1)-C(2)-C(4)	-0.0017	H(26)-C(14)-C(13)	119.6778	C(18)-C(16)-C(14)	119.9314	H(26)-C(14)-C(13)	119.8177
O(20)-C(1)-C(2)-C(7)	0.0014	C(16)-C(14)-C(13)	120.7461	O(19)-C(15)-C(17)	117.0414	C(16)-C(14)-C(13)	120.3931
O(20)-C(1)-C(2)-C(4)	-179.9945	C(15)-C(13)-C(14)	119.2222	O(19)-C(15)-C(13)	123.383	C(15)-C(13)-C(14)	119.5752
		C(15)-C(13)-N(12)	123.4595	C(17)-C(15)-C(13)	119.5482	C(15)-C(13)-N(12)	121.6129
		C(14)-C(13)-N(12)	117.0019	H(26)-C(14)-C(16)	119.5789	C(14)-C(13)-N(12)	118.3655
		C(13)-N(12)-N(11)	125.9684	H(26)-C(14)-C(13)	119.6797	C(13)-N(12)-N(11)	123.6942
		Cr(67)-N(11)-N(12)	121.6602	C(16)-C(14)-C(13)	120.7412	Fe(67)-N(11)-N(12)	121.1678
		Cr(67)-N(11)-C(6)	121.0832	C(15)-C(13)-C(14)	119.2156	Fe(67)-N(11)-C(6)	120.6264
		N(12)-N(11)-C(6)	117.2556	C(15)-C(13)-N(12)	123.465	N(12)-N(11)-C(6)	117.8456
		H(25)-C(10)-C(9)	120.1306	C(14)-C(13)-N(12)	117.029	H(25)-C(10)-C(9)	120.1324
		H(25)-C(10)-C(8)	119.7692	C(13)-N(12)-N(11)	125.3502	H(25)-C(10)-C(8)	119.7797
		C(9)-C(10)-C(8)	120.0954	Mn(69)-N(11)-N(12)	122.0391	C(9)-C(10)-C(8)	120.0814
		H(24)-C(9)-C(10)	118.9773	Mn(69)-N(11)-C(6)	119.556	H(24)-C(9)-C(10)	118.8653
		H(24)-C(9)-C(4)	120.5611	N(12)-N(11)-C(6)	118.2929	H(24)-C(9)-C(4)	120.6881
		C(10)-C(9)-C(4)	120.4586	H(25)-C(10)-C(9)	120.1509	C(10)-C(9)-C(4)	120.4369
		H(23)-C(8)-C(10)	119.5794	H(25)-C(10)-C(8)	119.7673	H(23)-C(8)-C(10)	119.5497
		H(23)-C(8)-C(7)	120.0793	C(9)-C(10)-C(8)	120.0817	H(23)-C(8)-C(7)	120.0873
		C(10)-C(8)-C(7)	120.3325	H(24)-C(9)-C(10)	118.7855	C(10)-C(8)-C(7)	120.3584
		H(22)-C(7)-C(8)	117.5988	H(24)-C(9)-C(4)	120.6514	H(22)-C(7)-C(8)	117.5211
		H(22)-C(7)-C(2)	121.8446	C(10)-C(9)-C(4)	120.5507	H(22)-C(7)-C(2)	121.901
		C(8)-C(7)-C(2)	120.54	H(23)-C(8)-C(10)	119.5856	C(8)-C(7)-C(2)	120.5608
		N(11)-C(6)-C(5)	119.5361	H(23)-C(8)-C(7)	120.1079	N(11)-C(6)-C(5)	119.7103
		N(11)-C(6)-C(4)	120.6784	C(10)-C(8)-C(7)	120.3061	N(11)-C(6)-C(4)	120.4716
		C(5)-C(6)-C(4)	119.621	H(22)-C(7)-C(8)	117.5007	C(5)-C(6)-C(4)	119.8181
		H(31)-C(5)-C(6)	120.2959	H(22)-C(7)-C(2)	121.9347	H(31)-C(5)-C(6)	120.4872
		H(31)-C(5)-C(3)	119.1898	C(8)-C(7)-C(2)	120.5456	H(31)-C(5)-C(3)	119.031
		C(6)-C(5)-C(3)	120.468	N(11)-C(6)-C(5)	119.2035	C(6)-C(5)-C(3)	120.453
		C(9)-C(4)-C(6)	120.0532	N(11)-C(6)-C(4)	121.2036	C(9)-C(4)-C(6)	120.0352
		C(9)-C(4)-C(2)	119.8241	C(5)-C(6)-C(4)	119.5895	C(9)-C(4)-C(2)	119.8495
		C(6)-C(4)-C(2)	120.0774	H(31)-C(5)-C(6)	120.4755	C(6)-C(4)-C(2)	120.0876
		H(21)-C(3)-C(5)	118.9526	H(31)-C(5)-C(3)	118.8752	H(21)-C(3)-C(5)	118.9104
		H(21)-C(3)-C(1)	120.0669	C(6)-C(5)-C(3)	120.6474	H(21)-C(3)-C(1)	120.044
		C(5)-C(3)-C(1)	120.9372	C(9)-C(4)-C(6)	120.2694	C(5)-C(3)-C(1)	121.0233
		C(4)-C(2)-C(7)	118.7419	C(9)-C(4)-C(2)	119.6676	C(4)-C(2)-C(7)	118.6895
		C(4)-C(2)-C(1)	119.2063	C(6)-C(4)-C(2)	120.0628	C(4)-C(2)-C(1)	119.1765
		C(7)-C(2)-C(1)	121.9976	H(21)-C(3)-C(5)	118.9245	C(7)-C(2)-C(1)	122.1121
		O(20)-C(1)-C(3)	117.4555	H(21)-C(3)-C(1)	120.054	O(20)-C(1)-C(3)	117.3591
		O(20)-C(1)-C(2)	123.1734	C(5)-C(3)-C(1)	121.0129	O(20)-C(1)-C(2)	123.2395
		C(3)-C(1)-C(2)	119.3704	C(4)-C(2)-C(7)	118.7432	C(3)-C(1)-C(2)	119.3937
		Dihedral Angles		C(4)-C(2)-C(1)	119.2398	Dihedral Angles	
		C(46)-O(65)-Cr(67)-N(11)	131.88	C(7)-C(2)-C(1)	122.0105	C(46)-O(65)-Fe(67)-N(11)	106.3112
		C(46)-O(65)-Cr(67)-O(19)	41.7735	O(20)-C(1)-C(3)	117.3183	C(46)-O(65)-Fe(67)-O(19)	21.204
		C(46)-O(65)-Cr(67)-N(42)	-47.6094	O(20)-C(1)-C(2)	123.2873	C(46)-O(65)-Fe(67)-N(42)	-52.7806
		C(46)-O(65)-Cr(67)-O(62)	-141.851	C(3)-C(1)-C(2)	119.3785	C(46)-O(65)-Fe(67)-O(62)	-163.4344
		C(46)-O(65)-Cr(67)-Cl(66)	-136.5214	Dihedral Angles		C(46)-O(65)-Fe(67)-Cl(66)	-142.6971
		H(63)-O(62)-Cr(67)-N(11)	14.9145	H(66)-O(68)-Mn(69)-N(11)	120.689	H(63)-O(62)-Fe(67)-N(11)	14.1734

	H(63)-O(62)-Cr(67)-O(19)	105.0444	H(66)-O(68)-Mn(69)-O(19)	72.2968	H(63)-O(62)-Fe(67)-O(19)	101.9702
	H(63)-O(62)-Cr(67)-N(42)	-165.5615	H(66)-O(68)-Mn(69)-N(42)	-57.811	H(63)-O(62)-Fe(67)-N(42)	179.0655
	H(63)-O(62)-Cr(67)-O(65)	-71.3287	H(66)-O(68)-Mn(69)-O(62)	-148.0257	H(63)-O(62)-Fe(67)-O(65)	-73.5343
	H(63)-O(62)-Cr(67)-Cl(66)	-76.6595	H(66)-O(68)-Mn(69)-O(65)	31.7682	H(63)-O(62)-Fe(67)-Cl(66)	-94.6399
	H(64)-O(62)-Cr(67)-N(11)	-99.2249	H(67)-O(68)-Mn(69)-N(11)	-126.8135	H(64)-O(62)-Fe(67)-N(11)	-96.9659
	H(64)-O(62)-Cr(67)-O(19)	-9.095	H(67)-O(68)-Mn(69)-O(19)	-175.2057	H(64)-O(62)-Fe(67)-O(19)	-9.169
	H(64)-O(62)-Cr(67)-N(42)	80.2991	H(67)-O(68)-Mn(69)-N(42)	54.6865	H(64)-O(62)-Fe(67)-N(42)	67.9262
	H(64)-O(62)-Cr(67)-O(65)	174.5319	H(67)-O(68)-Mn(69)-O(62)	-35.5281	H(64)-O(62)-Fe(67)-O(65)	175.3264
	H(64)-O(62)-Cr(67)-Cl(66)	169.2011	H(67)-O(68)-Mn(69)-O(65)	144.2657	H(64)-O(62)-Fe(67)-Cl(66)	154.2208
	C(46)-C(48)-C(49)-C(47)	0.8486	C(46)-O(65)-Mn(69)-N(11)	131.5251	C(46)-C(48)-C(49)-C(47)	1.2477
	C(46)-C(48)-C(49)-H(59)	-179.7019	C(46)-O(65)-Mn(69)-O(19)	41.1262	C(46)-C(48)-C(49)-H(59)	-179.5164
	H(58)-C(48)-C(49)-C(47)	-179.1917	C(46)-O(65)-Mn(69)-N(42)	-48.4163	H(58)-C(48)-C(49)-C(47)	-178.899
	H(58)-C(48)-C(49)-H(59)	0.2579	C(46)-O(65)-Mn(69)-O(62)	-128.5681	H(58)-C(48)-C(49)-H(59)	0.337
	C(45)-C(47)-C(49)-C(48)	0.4586	C(46)-O(65)-Mn(69)-O(68)	-139.1365	C(45)-C(47)-C(49)-C(48)	0.5303
	C(45)-C(47)-C(49)-H(59)	-178.9927	H(63)-O(62)-Mn(69)-N(11)	3.6704	C(45)-C(47)-C(49)-H(59)	-178.7077
	H(57)-C(47)-C(49)-C(48)	179.685	H(63)-O(62)-Mn(69)-O(19)	94.0942	H(57)-C(47)-C(49)-C(48)	179.5829
	H(57)-C(47)-C(49)-H(59)	0.2337	H(63)-O(62)-Mn(69)-N(42)	-176.3592	H(57)-C(47)-C(49)-H(59)	0.3449
	C(44)-C(46)-O(65)-Cr(67)	36.9544	H(63)-O(62)-Mn(69)-O(65)	-96.2136	C(44)-C(46)-O(65)-Fe(67)	40.8314
	C(48)-C(46)-O(65)-Cr(67)	-146.1671	H(63)-O(62)-Mn(69)-O(68)	-85.6441	C(48)-C(46)-O(65)-Fe(67)	-143.6759
	C(44)-C(46)-C(48)-C(49)	-1.8439	H(64)-O(62)-Mn(69)-N(11)	-109.435	C(44)-C(46)-C(48)-C(49)	-2.5822
	C(44)-C(46)-C(48)-H(58)	178.1964	H(64)-O(62)-Mn(69)-O(19)	-19.0112	C(44)-C(46)-C(48)-H(58)	177.5644
	O(65)-C(46)-C(48)-C(49)	-178.8305	H(64)-O(62)-Mn(69)-N(42)	70.5354	O(65)-C(46)-C(48)-C(49)	-178.1856
	O(65)-C(46)-C(48)-H(58)	1.2099	H(64)-O(62)-Mn(69)-O(65)	150.681	O(65)-C(46)-C(48)-H(58)	1.961
	C(44)-C(45)-C(47)-C(49)	-0.8092	H(64)-O(62)-Mn(69)-O(68)	161.2505	C(44)-C(45)-C(47)-C(49)	-1.0377
	C(44)-C(45)-C(47)-H(57)	179.967	C(46)-C(48)-C(49)-C(47)	0.6589	C(44)-C(45)-C(47)-H(57)	179.9128
	H(56)-C(45)-C(47)-C(49)	179.0034	C(46)-C(48)-C(49)-H(59)	-179.7056	H(56)-C(45)-C(47)-C(49)	178.8246
	H(56)-C(45)-C(47)-H(57)	-0.2205	H(58)-C(48)-C(49)-C(47)	-179.6818	H(56)-C(45)-C(47)-H(57)	-0.2249
	N(43)-C(44)-C(46)-C(48)	-172.0924	H(58)-C(48)-C(49)-H(59)	-0.0462	N(43)-C(44)-C(46)-C(48)	-171.0409
	N(43)-C(44)-C(46)-O(65)	4.7278	C(45)-C(47)-C(49)-C(48)	0.2481	N(43)-C(44)-C(46)-O(65)	4.3786
	C(45)-C(44)-C(46)-C(48)	1.4785	C(45)-C(47)-C(49)-H(59)	-179.3886	C(45)-C(44)-C(46)-C(48)	2.0565
	C(45)-C(44)-C(46)-O(65)	178.2987	H(57)-C(47)-C(49)-C(48)	179.6458	C(45)-C(44)-C(46)-O(65)	177.476
	N(43)-C(44)-C(45)-C(47)	173.7701	H(57)-C(47)-C(49)-H(59)	0.0091	N(43)-C(44)-C(45)-C(47)	173.1228
	N(43)-C(44)-C(45)-H(56)	-6.0423	C(44)-C(46)-O(65)-Mn(69)	38.209	N(43)-C(44)-C(45)-H(56)	-6.7394
	C(46)-C(44)-C(45)-C(47)	-0.1916	C(48)-C(46)-O(65)-Mn(69)	-144.8569	C(46)-C(44)-C(45)-C(47)	-0.2991
	C(46)-C(44)-C(45)-H(56)	179.996	C(44)-C(46)-C(48)-C(49)	-1.2279	C(46)-C(44)-C(45)-H(56)	179.8387
	N(42)-N(43)-C(44)-C(45)	163.3135	C(44)-C(46)-C(48)-H(58)	179.1124	N(42)-N(43)-C(44)-C(45)	161.4045
	N(42)-N(43)-C(44)-C(46)	-22.9883	O(65)-C(46)-C(48)-C(49)	-178.2602	N(42)-N(43)-C(44)-C(46)	-25.4008
	C(37)-N(42)-Cr(67)-N(11)	-177.5062	O(65)-C(46)-C(48)-H(58)	2.0801	C(37)-N(42)-Fe(67)-N(11)	137.1828
	C(37)-N(42)-Cr(67)-O(19)	120.7176	C(44)-C(45)-C(47)-C(49)	-0.6114	C(37)-N(42)-Fe(67)-O(19)	106.07
	C(37)-N(42)-Cr(67)-O(62)	28.6698	C(44)-C(45)-C(47)-H(57)	179.993	C(37)-N(42)-Fe(67)-O(62)	3.2485
	C(37)-N(42)-Cr(67)-O(65)	-149.2656	H(56)-C(45)-C(47)-C(49)	179.1701	C(37)-N(42)-Fe(67)-O(65)	-142.6707
	C(37)-N(42)-Cr(67)-Cl(66)	-59.9175	H(56)-C(45)-C(47)-H(57)	-0.2255	C(37)-N(42)-Fe(67)-Cl(66)	-67.9442
	N(43)-N(42)-Cr(67)-N(11)	5.3558	N(43)-C(44)-C(46)-C(48)	-172.956	N(43)-N(42)-Fe(67)-N(11)	-43.2648
	N(43)-N(42)-Cr(67)-O(19)	-56.4204	N(43)-C(44)-C(46)-O(65)	3.9132	N(43)-N(42)-Fe(67)-O(19)	-74.3775
	N(43)-N(42)-Cr(67)-O(62)	-148.4682	C(45)-C(44)-C(46)-C(48)	0.8571	N(43)-N(42)-Fe(67)-O(62)	-177.199
	N(43)-N(42)-Cr(67)-O(65)	33.5964	C(45)-C(44)-C(46)-O(65)	177.7263	N(43)-N(42)-Fe(67)-O(65)	36.8817
	N(43)-N(42)-Cr(67)-Cl(66)	122.9445	N(43)-C(44)-C(45)-C(47)	174.2205	N(43)-N(42)-Fe(67)-Cl(66)	111.6082
	C(37)-N(42)-N(43)-C(44)	177.9821	N(43)-C(44)-C(45)-H(56)	-5.5608	C(37)-N(42)-N(43)-C(44)	175.4884
	Cr(67)-N(42)-N(43)-C(44)	-4.7794	C(46)-C(44)-C(45)-C(47)	0.0399	Fe(67)-N(42)-N(43)-C(44)	-4.074
	C(35)-C(40)-C(41)-C(39)	-0.8275	C(46)-C(44)-C(45)-H(56)	-179.7414	C(35)-C(40)-C(41)-C(39)	0.4728
	C(35)-C(40)-C(41)-H(55)	179.9295	N(42)-N(43)-C(44)-C(45)	163.1811	C(35)-C(40)-C(41)-H(55)	179.8028
	H(54)-C(40)-C(41)-C(39)	179.1703	N(42)-N(43)-C(44)-C(46)	-22.8943	H(54)-C(40)-C(41)-C(39)	-178.9698
	H(54)-C(40)-C(41)-H(55)	-0.0727	C(37)-N(42)-Mn(69)-N(11)	-151.38	H(54)-C(40)-C(41)-H(55)	0.3602
	C(38)-C(39)-C(41)-C(40)	0.4984	C(37)-N(42)-Mn(69)-O(19)	120.9147	C(38)-C(39)-C(41)-C(40)	-0.3393

	C(38)-C(39)-C(41)-H(55)	179.7442	C(37)-N(42)-Mn(69)-O(62)	29.7508	C(38)-C(39)-C(41)-H(55)	-179.6719
	H(53)-C(39)-C(41)-C(40)	-178.5407	C(37)-N(42)-Mn(69)-O(65)	-149.142	H(53)-C(39)-C(41)-C(40)	178.9517
	H(53)-C(39)-C(41)-H(55)	0.7051	C(37)-N(42)-Mn(69)-O(68)	-59.3946	H(53)-C(39)-C(41)-H(55)	-0.3809
	C(33)-C(38)-C(39)-C(41)	0.6376	N(43)-N(42)-Mn(69)-N(11)	31.5099	C(33)-C(38)-C(39)-C(41)	-0.4745
	C(33)-C(38)-C(39)-H(53)	179.6716	N(43)-N(42)-Mn(69)-O(19)	-56.1954	C(33)-C(38)-C(39)-H(53)	-179.7617
	H(52)-C(38)-C(39)-C(41)	-177.8601	N(43)-N(42)-Mn(69)-O(62)	-147.3593	H(52)-C(38)-C(39)-C(41)	178.4621
	H(52)-C(38)-C(39)-H(53)	1.1739	N(43)-N(42)-Mn(69)-O(65)	33.7479	H(52)-C(38)-C(39)-H(53)	-0.8252
	C(35)-C(37)-N(42)-N(43)	-77.4729	N(43)-N(42)-Mn(69)-O(68)	123.4953	C(35)-C(37)-N(42)-N(43)	-67.0411
	C(35)-C(37)-N(42)-Cr(67)	105.311	C(37)-N(42)-N(43)-C(44)	178.1364	C(35)-C(37)-N(42)-Fe(67)	112.5174
	C(36)-C(37)-N(42)-N(43)	97.2088	Mn(69)-N(42)-N(43)-C(44)	-4.6459	C(36)-C(37)-N(42)-N(43)	111.3431
	C(36)-C(37)-N(42)-Cr(67)	-80.0073	C(35)-C(40)-C(41)-C(39)	-1.0197	C(36)-C(37)-N(42)-Fe(67)	-69.0985
	C(34)-C(36)-C(37)-C(35)	-5.2788	C(35)-C(40)-C(41)-H(55)	-179.8269	C(34)-C(36)-C(37)-C(35)	0.1197
	C(34)-C(36)-C(37)-N(42)	179.9796	H(54)-C(40)-C(41)-C(39)	177.9846	C(34)-C(36)-C(37)-N(42)	-178.284
	H(61)-C(36)-C(37)-C(35)	175.1038	H(54)-C(40)-C(41)-H(55)	-0.8226	H(61)-C(36)-C(37)-C(35)	178.2827
	H(61)-C(36)-C(37)-N(42)	0.3621	C(38)-C(39)-C(41)-C(40)	0.0676	H(61)-C(36)-C(37)-N(42)	-0.121
	C(33)-C(35)-C(40)-C(41)	-0.006	C(38)-C(39)-C(41)-H(55)	178.8793	C(33)-C(35)-C(40)-C(41)	0.2365
	C(33)-C(35)-C(40)-H(54)	179.9963	H(53)-C(39)-C(41)-C(40)	-178.6568	C(33)-C(35)-C(40)-H(54)	179.6672
	C(37)-C(35)-C(40)-C(41)	178.436	H(53)-C(39)-C(41)-H(55)	0.1548	C(37)-C(35)-C(40)-C(41)	-178.4035
	C(37)-C(35)-C(40)-H(54)	-1.5617	C(33)-C(38)-C(39)-C(41)	0.964	C(37)-C(35)-C(40)-H(54)	1.0273
	C(33)-C(35)-C(37)-C(36)	4.1977	C(33)-C(38)-C(39)-H(53)	179.6818	C(33)-C(35)-C(37)-C(36)	0.134
	C(33)-C(35)-C(37)-N(42)	178.8826	H(52)-C(38)-C(39)-C(41)	-177.6384	C(33)-C(35)-C(37)-N(42)	178.5189
	C(40)-C(35)-C(37)-C(36)	-174.239	H(52)-C(38)-C(39)-H(53)	1.0794	C(40)-C(35)-C(37)-C(36)	178.7686
	C(40)-C(35)-C(37)-N(42)	0.446	C(35)-C(37)-N(42)-N(43)	-80.5671	C(40)-C(35)-C(37)-N(42)	-2.8466
	C(32)-C(34)-C(36)-C(37)	1.7369	C(35)-C(37)-N(42)-Mn(69)	102.2461	C(32)-C(34)-C(36)-C(37)	0.716
	C(32)-C(34)-C(36)-H(61)	-178.64	C(36)-C(37)-N(42)-N(43)	94.6662	C(32)-C(34)-C(36)-H(61)	-177.4807
	H(51)-C(34)-C(36)-C(37)	-176.3496	C(36)-C(37)-N(42)-Mn(69)	-82.5206	H(51)-C(34)-C(36)-C(37)	179.0463
	H(51)-C(34)-C(36)-H(61)	3.2735	C(34)-C(36)-C(37)-C(35)	-6.3705	H(51)-C(34)-C(36)-H(61)	0.8496
	C(32)-C(33)-C(35)-C(37)	0.4163	C(34)-C(36)-C(37)-N(42)	178.3367	C(32)-C(33)-C(35)-C(37)	-1.1012
	C(32)-C(33)-C(35)-C(40)	178.8591	H(61)-C(36)-C(37)-C(35)	174.5882	C(32)-C(33)-C(35)-C(40)	-179.7421
	C(38)-C(33)-C(35)-C(37)	-177.3302	H(61)-C(36)-C(37)-N(42)	-0.7046	C(38)-C(33)-C(35)-C(37)	177.6169
	C(38)-C(33)-C(35)-C(40)	1.1127	C(33)-C(35)-C(40)-C(41)	0.9358	C(38)-C(33)-C(35)-C(40)	-1.024
	C(32)-C(33)-C(38)-C(39)	-179.1291	C(33)-C(35)-C(40)-H(54)	-178.0501	C(32)-C(33)-C(38)-C(39)	179.8451
	C(32)-C(33)-C(38)-H(52)	-0.6967	C(37)-C(35)-C(40)-C(41)	178.8852	C(32)-C(33)-C(38)-H(52)	0.9588
	C(35)-C(33)-C(38)-C(39)	-1.4479	C(37)-C(35)-C(40)-H(54)	-0.1007	C(35)-C(33)-C(38)-C(39)	1.1653
	C(35)-C(33)-C(38)-H(52)	176.9846	C(33)-C(35)-C(37)-C(36)	5.7681	C(35)-C(33)-C(38)-H(52)	-177.721
	C(33)-C(32)-O(50)-H(60)	-10.8808	C(33)-C(35)-C(37)-N(42)	-178.9907	C(33)-C(32)-O(50)-H(60)	7.932
	C(34)-C(32)-O(50)-H(60)	168.6986	C(40)-C(35)-C(37)-C(36)	-172.1743	C(34)-C(32)-O(50)-H(60)	-171.3489
	C(33)-C(32)-C(34)-C(36)	2.9373	C(40)-C(35)-C(37)-N(42)	3.0669	C(33)-C(32)-C(34)-C(36)	-1.6881
	C(33)-C(32)-C(34)-H(51)	-178.9974	C(32)-C(34)-C(36)-C(37)	1.692	C(33)-C(32)-C(34)-H(51)	-179.999
	O(50)-C(32)-C(34)-C(36)	-176.6588	C(32)-C(34)-C(36)-H(61)	-179.2518	O(50)-C(32)-C(34)-C(36)	177.6228
	O(50)-C(32)-C(34)-H(51)	1.4066	H(51)-C(34)-C(36)-C(37)	-175.8943	O(50)-C(32)-C(34)-H(51)	-0.6882
	C(34)-C(32)-C(33)-C(38)	173.7428	H(51)-C(34)-C(36)-H(61)	3.162	C(34)-C(32)-C(33)-C(38)	-176.7668
	C(34)-C(32)-C(33)-C(35)	-3.9285	C(32)-C(33)-C(35)-C(37)	-0.4669	C(34)-C(32)-C(33)-C(35)	1.9069
	O(50)-C(32)-C(33)-C(38)	-6.6856	C(32)-C(33)-C(35)-C(40)	177.4845	O(50)-C(32)-C(33)-C(38)	3.9668
	O(50)-C(32)-C(33)-C(35)	175.6431	C(38)-C(33)-C(35)-C(37)	-177.8674	O(50)-C(32)-C(33)-C(35)	-177.3595
	C(15)-O(19)-Cr(67)-N(11)	41.4882	C(38)-C(33)-C(35)-C(40)	0.084	C(15)-O(19)-Fe(67)-N(11)	55.3659
	C(15)-O(19)-Cr(67)-N(42)	-139.4626	C(32)-C(33)-C(38)-C(39)	-178.3554	C(15)-O(19)-Fe(67)-N(42)	-135.4071
	C(15)-O(19)-Cr(67)-O(62)	-49.1639	C(32)-C(33)-C(38)-H(52)	0.1866	C(15)-O(19)-Fe(67)-O(62)	-37.5249
	C(15)-O(19)-Cr(67)-O(65)	130.7051	C(35)-C(33)-C(38)-C(39)	-1.0287	C(15)-O(19)-Fe(67)-O(65)	139.6553
	C(15)-O(19)-Cr(67)-Cl(66)	-159.8729	C(35)-C(33)-C(38)-H(52)	177.5133	C(15)-O(19)-Fe(67)-Cl(66)	-113.4997
	C(15)-C(17)-C(18)-C(16)	-0.5076	C(33)-C(32)-O(50)-H(60)	-11.6098	C(15)-C(17)-C(18)-C(16)	-0.9821
	C(15)-C(17)-C(18)-H(29)	179.745	C(34)-C(32)-O(50)-H(60)	168.2081	C(15)-C(17)-C(18)-H(29)	179.584
	H(28)-C(17)-C(18)-C(16)	179.8473	C(33)-C(32)-C(34)-C(36)	3.6728	H(28)-C(17)-C(18)-C(16)	179.3389
	H(28)-C(17)-C(18)-H(29)	0.0998	C(33)-C(32)-C(34)-H(51)	-178.7678	H(28)-C(17)-C(18)-H(29)	-0.095

		C(14)-C(16)-C(18)-C(17)	-0.2736	O(50)-C(32)-C(34)-C(36)	-176.1523	C(14)-C(16)-C(18)-C(17)	-0.4563
		C(14)-C(16)-C(18)-H(29)	179.4746	O(50)-C(32)-C(34)-H(51)	1.4071	C(14)-C(16)-C(18)-H(29)	178.9794
		H(27)-C(16)-C(18)-C(17)	-179.7015	C(34)-C(32)-C(33)-C(38)	173.1421	H(27)-C(16)-C(18)-C(17)	-179.5869
		H(27)-C(16)-C(18)-H(29)	0.0467	C(34)-C(32)-C(33)-C(35)	-4.1734	H(27)-C(16)-C(18)-H(29)	-0.1512
		C(13)-C(15)-O(19)-Cr(67)	-31.7568	O(50)-C(32)-C(33)-C(38)	-7.0433	C(13)-C(15)-O(19)-Fe(67)	-42.1835
		C(17)-C(15)-O(19)-Cr(67)	150.0641	O(50)-C(32)-C(33)-C(35)	175.6411	C(17)-C(15)-O(19)-Fe(67)	141.814
		C(13)-C(15)-C(17)-C(18)	1.0393	C(15)-O(19)-Mn(69)-N(11)	42.2092	C(13)-C(15)-C(17)-C(18)	2.0553
		C(13)-C(15)-C(17)-H(28)	-179.3153	C(15)-O(19)-Mn(69)-N(42)	-139.2905	C(13)-C(15)-C(17)-H(28)	-178.2655
		O(19)-C(15)-C(17)-C(18)	179.2896	C(15)-O(19)-Mn(69)-O(62)	-49.0744	O(19)-C(15)-C(17)-C(18)	178.1234
		O(19)-C(15)-C(17)-H(28)	-1.0649	C(15)-O(19)-Mn(69)-O(65)	131.1267	O(19)-C(15)-C(17)-H(28)	-2.1974
		C(13)-C(14)-C(16)-C(18)	0.5433	C(15)-O(19)-Mn(69)-O(68)	90.5985	C(13)-C(14)-C(16)-C(18)	0.8767
		C(13)-C(14)-C(16)-H(27)	179.9693	C(15)-C(17)-C(18)-C(16)	-0.597	C(13)-C(14)-C(16)-H(27)	-179.9955
		H(26)-C(14)-C(16)-C(18)	-179.1628	C(15)-C(17)-C(18)-H(29)	179.6849	H(26)-C(14)-C(16)-C(18)	-178.7894
		H(26)-C(14)-C(16)-H(27)	0.2633	H(28)-C(17)-C(18)-C(16)	179.821	H(26)-C(14)-C(16)-H(27)	0.3384
		N(12)-C(13)-C(15)-C(17)	172.5778	H(28)-C(17)-C(18)-H(29)	0.1029	N(12)-C(13)-C(15)-C(17)	170.5809
		N(12)-C(13)-C(15)-O(19)	-5.5572	C(14)-C(16)-C(18)-C(17)	-0.2377	N(12)-C(13)-C(15)-O(19)	-5.3761
		C(14)-C(13)-C(15)-C(17)	-0.7625	C(14)-C(16)-C(18)-H(29)	179.4812	C(14)-C(13)-C(15)-C(17)	-1.626
		C(14)-C(13)-C(15)-O(19)	-178.8975	H(27)-C(16)-C(18)-C(17)	-179.6763	C(14)-C(13)-C(15)-O(19)	-177.583
		N(12)-C(13)-C(14)-C(16)	-173.775	H(27)-C(16)-C(18)-H(29)	0.0427	N(12)-C(13)-C(14)-C(16)	-172.2618
		N(12)-C(13)-C(14)-H(26)	5.9308	C(13)-C(15)-O(19)-Mn(69)	-32.9767	N(12)-C(13)-C(14)-H(26)	7.4041
		C(15)-C(13)-C(14)-C(16)	-0.0091	C(17)-C(15)-O(19)-Mn(69)	148.9383	C(15)-C(13)-C(14)-C(16)	0.1974
		C(15)-C(13)-C(14)-H(26)	179.6967	C(13)-C(15)-C(17)-C(18)	1.1005	C(15)-C(13)-C(14)-H(26)	179.8633
		N(11)-N(12)-C(13)-C(14)	-165.2797	C(13)-C(15)-C(17)-H(28)	-179.3171	N(11)-N(12)-C(13)-C(14)	-163.781
		N(11)-N(12)-C(13)-C(15)	21.2429	O(19)-C(15)-C(17)-C(18)	179.2624	N(11)-N(12)-C(13)-C(15)	23.9211
		C(6)-N(11)-Cr(67)-O(19)	151.3239	O(19)-C(15)-C(17)-H(28)	-1.1552	C(6)-N(11)-Fe(67)-O(19)	146.1276
		C(6)-N(11)-Cr(67)-N(42)	89.5537	C(13)-C(14)-C(16)-C(18)	0.5881	C(6)-N(11)-Fe(67)-N(42)	116.1751
		C(6)-N(11)-Cr(67)-O(62)	-116.6237	C(13)-C(14)-C(16)-H(27)	-179.9752	C(6)-N(11)-Fe(67)-O(62)	-108.8562
		C(6)-N(11)-Cr(67)-O(65)	61.3103	H(26)-C(14)-C(16)-C(18)	-179.2114	C(6)-N(11)-Fe(67)-O(65)	35.1052
		C(6)-N(11)-Cr(67)-Cl(66)	-28.0126	H(26)-C(14)-C(16)-H(27)	0.2252	C(6)-N(11)-Fe(67)-Cl(66)	-37.1442
		N(12)-N(11)-Cr(67)-O(19)	-29.0558	N(12)-C(13)-C(15)-C(17)	172.8769	N(12)-N(11)-Fe(67)-O(19)	-40.9114
		N(12)-N(11)-Cr(67)-N(42)	-90.826	N(12)-C(13)-C(15)-O(19)	-5.1624	N(12)-N(11)-Fe(67)-N(42)	-70.8639
		N(12)-N(11)-Cr(67)-O(62)	62.9967	C(14)-C(13)-C(15)-C(17)	-0.7434	N(12)-N(11)-Fe(67)-O(62)	64.1048
		N(12)-N(11)-Cr(67)-O(65)	-119.0694	C(14)-C(13)-C(15)-O(19)	-178.7826	N(12)-N(11)-Fe(67)-O(65)	-151.9338
		N(12)-N(11)-Cr(67)-Cl(66)	151.6077	N(12)-C(13)-C(14)-C(16)	-174.1074	N(12)-N(11)-Fe(67)-Cl(66)	135.8167
		C(6)-N(11)-N(12)-C(13)	-176.5864	N(12)-C(13)-C(14)-H(26)	5.692	C(6)-N(11)-N(12)-C(13)	-178.0729
		Cr(67)-N(11)-N(12)-C(13)	3.7793	C(15)-C(13)-C(14)-C(16)	-0.0806	Fe(67)-N(11)-N(12)-C(13)	8.7765
		C(4)-C(9)-C(10)-C(8)	0.6683	C(15)-C(13)-C(14)-H(26)	179.7188	C(4)-C(9)-C(10)-C(8)	-0.3022
		C(4)-C(9)-C(10)-H(25)	179.8748	N(11)-N(12)-C(13)-C(14)	-163.6025	C(4)-C(9)-C(10)-H(25)	-179.375
		H(24)-C(9)-C(10)-C(8)	-178.7011	N(11)-N(12)-C(13)-C(15)	22.6479	H(24)-C(9)-C(10)-C(8)	178.574
		H(24)-C(9)-C(10)-H(25)	0.5054	C(6)-N(11)-Mn(69)-O(19)	147.4326	H(24)-C(9)-C(10)-H(25)	-0.4988
		C(7)-C(8)-C(10)-C(9)	-0.0365	C(6)-N(11)-Mn(69)-N(42)	59.7385	C(7)-C(8)-C(10)-C(9)	0.5563
		C(7)-C(8)-C(10)-H(25)	-179.2458	C(6)-N(11)-Mn(69)-O(62)	-121.3926	C(7)-C(8)-C(10)-H(25)	179.6324
		H(23)-C(8)-C(10)-C(9)	178.8861	C(6)-N(11)-Mn(69)-O(65)	57.5001	H(23)-C(8)-C(10)-C(9)	-178.6673
		H(23)-C(8)-C(10)-H(25)	-0.3232	C(6)-N(11)-Mn(69)-O(68)	-32.2651	H(23)-C(8)-C(10)-H(25)	0.4088
		C(2)-C(7)-C(8)-C(10)	-0.719	N(12)-N(11)-Mn(69)-O(19)	-28.6529	C(2)-C(7)-C(8)-C(10)	0.3543
		C(2)-C(7)-C(8)-H(23)	-179.6362	N(12)-N(11)-Mn(69)-N(42)	-116.347	C(2)-C(7)-C(8)-H(23)	179.5737
		H(22)-C(7)-C(8)-C(10)	177.8283	N(12)-N(11)-Mn(69)-O(62)	62.522	H(22)-C(7)-C(8)-C(10)	-178.167
		H(22)-C(7)-C(8)-H(23)	-1.0889	N(12)-N(11)-Mn(69)-O(65)	-118.5853	H(22)-C(7)-C(8)-H(23)	1.0524
		C(4)-C(6)-N(11)-N(12)	-103.0673	N(12)-N(11)-Mn(69)-O(68)	151.6495	C(4)-C(6)-N(11)-N(12)	74.0072
		C(4)-C(6)-N(11)-Cr(67)	76.5692	C(6)-N(11)-N(12)-C(13)	-174.027	C(4)-C(6)-N(11)-Fe(67)	-112.8034
		C(5)-C(6)-N(11)-N(12)	81.6119	Mn(69)-N(11)-N(12)-C(13)	2.1059	C(5)-C(6)-N(11)-N(12)	-106.0483
		C(5)-C(6)-N(11)-Cr(67)	-98.7516	C(4)-C(9)-C(10)-C(8)	0.1177	C(5)-C(6)-N(11)-Fe(67)	67.1411
		C(3)-C(5)-C(6)-C(4)	5.6564	C(4)-C(9)-C(10)-H(25)	-179.8231	C(3)-C(5)-C(6)-C(4)	-0.5372
		C(3)-C(5)-C(6)-N(11)	-178.9728	H(24)-C(9)-C(10)-C(8)	178.8423	C(3)-C(5)-C(6)-N(11)	179.5179

	H(31)-C(5)-C(6)-C(4)	-176.8391	H(24)-C(9)-C(10)-H(25)	-1.0984	H(31)-C(5)-C(6)-C(4)	-178.5665
	H(31)-C(5)-C(6)-N(11)	-1.4684	C(7)-C(8)-C(10)-C(9)	-1.4078	H(31)-C(5)-C(6)-N(11)	1.4887
	C(2)-C(4)-C(9)-C(10)	-0.5351	C(7)-C(8)-C(10)-H(25)	178.5332	C(2)-C(4)-C(9)-C(10)	-0.9145
	C(2)-C(4)-C(9)-H(24)	178.8242	H(23)-C(8)-C(10)-C(9)	178.3525	C(2)-C(4)-C(9)-H(24)	-179.77
	C(6)-C(4)-C(9)-C(10)	-178.0879	H(23)-C(8)-C(10)-H(25)	-1.7065	C(6)-C(4)-C(9)-C(10)	177.174
	C(6)-C(4)-C(9)-H(24)	1.2714	C(2)-C(7)-C(8)-C(10)	0.0707	C(6)-C(4)-C(9)-H(24)	-1.6815
	C(2)-C(4)-C(6)-C(5)	-4.8409	C(2)-C(7)-C(8)-H(23)	-179.6884	C(2)-C(4)-C(6)-C(5)	-0.4734
	C(2)-C(4)-C(6)-N(11)	179.8424	H(22)-C(7)-C(8)-C(10)	178.5186	C(2)-C(4)-C(6)-N(11)	179.471
	C(9)-C(4)-C(6)-C(5)	172.7057	H(22)-C(7)-C(8)-H(23)	-1.2405	C(9)-C(4)-C(6)-C(5)	-178.5573
	C(9)-C(4)-C(6)-N(11)	-2.6111	C(4)-C(6)-N(11)-N(12)	60.4752	C(9)-C(4)-C(6)-N(11)	1.3871
	C(1)-C(3)-C(5)-C(6)	-1.5966	C(4)-C(6)-N(11)-Mn(69)	-115.7565	C(1)-C(3)-C(5)-C(6)	-0.1539
	C(1)-C(3)-C(5)-H(31)	-179.1285	C(5)-C(6)-N(11)-N(12)	-118.8517	C(1)-C(3)-C(5)-H(31)	177.9038
	H(21)-C(3)-C(5)-C(6)	176.0112	C(5)-C(6)-N(11)-Mn(69)	64.9166	H(21)-C(3)-C(5)-C(6)	-178.4372
	H(21)-C(3)-C(5)-H(31)	-1.5207	C(3)-C(5)-C(6)-C(4)	-1.1406	H(21)-C(3)-C(5)-H(31)	-0.3795
	C(1)-C(2)-C(4)-C(6)	-0.033	C(3)-C(5)-C(6)-N(11)	178.1973	C(1)-C(2)-C(4)-C(6)	2.0271
	C(1)-C(2)-C(4)-C(9)	-177.5853	H(31)-C(5)-C(6)-C(4)	178.3491	C(1)-C(2)-C(4)-C(9)	-179.8854
	C(7)-C(2)-C(4)-C(6)	177.3433	H(31)-C(5)-C(6)-N(11)	-2.313	C(7)-C(2)-C(4)-C(6)	-176.3021
	C(7)-C(2)-C(4)-C(9)	-0.209	C(2)-C(4)-C(9)-C(10)	2.605	C(7)-C(2)-C(4)-C(9)	1.7854
	C(1)-C(2)-C(7)-C(8)	178.1368	C(2)-C(4)-C(9)-H(24)	-176.0957	C(1)-C(2)-C(7)-C(8)	-179.824
	C(1)-C(2)-C(7)-H(22)	-0.3478	C(6)-C(4)-C(9)-C(10)	-177.214	C(1)-C(2)-C(7)-H(22)	-1.3687
	C(4)-C(2)-C(7)-C(8)	0.8372	C(6)-C(4)-C(9)-H(24)	4.0853	C(4)-C(2)-C(7)-C(8)	-1.5463
	C(4)-C(2)-C(7)-H(22)	-177.6473	C(2)-C(4)-C(6)-C(5)	2.6685	C(4)-C(2)-C(7)-H(22)	176.909
	C(2)-C(1)-O(20)-H(30)	11.1298	C(2)-C(4)-C(6)-N(11)	-176.6559	C(2)-C(1)-O(20)-H(30)	-9.3173
	C(3)-C(1)-O(20)-H(30)	-168.5605	C(9)-C(4)-C(6)-C(5)	-177.5132	C(3)-C(1)-O(20)-H(30)	169.6659
	C(2)-C(1)-C(3)-C(5)	-3.3265	C(9)-C(4)-C(6)-N(11)	3.1624	C(2)-C(1)-C(3)-C(5)	1.7274
	C(2)-C(1)-C(3)-H(21)	179.0922	C(1)-C(3)-C(5)-C(6)	0.2695	C(2)-C(1)-C(3)-H(21)	179.9915
	O(20)-C(1)-C(3)-C(5)	176.376	C(1)-C(3)-C(5)-H(31)	-179.2282	O(20)-C(1)-C(3)-C(5)	-177.2964
	O(20)-C(1)-C(3)-H(21)	-1.2053	H(21)-C(3)-C(5)-C(6)	179.2022	O(20)-C(1)-C(3)-H(21)	0.9676
	C(3)-C(1)-C(2)-C(7)	-173.2364	H(21)-C(3)-C(5)-H(31)	-0.2956	C(3)-C(1)-C(2)-C(7)	175.5778
	C(3)-C(1)-C(2)-C(4)	4.051	C(1)-C(2)-C(4)-C(6)	-3.1387	C(3)-C(1)-C(2)-C(4)	-2.6918
	O(20)-C(1)-C(2)-C(7)	7.079	C(1)-C(2)-C(4)-C(9)	177.0419	O(20)-C(1)-C(2)-C(7)	-5.4587
	O(20)-C(1)-C(2)-C(4)	-175.6336	C(7)-C(2)-C(4)-C(6)	175.9549	O(20)-C(1)-C(2)-C(4)	176.2717
			C(7)-C(2)-C(4)-C(9)	-3.8645		
			C(1)-C(2)-C(7)-C(8)	-178.3138		
			C(1)-C(2)-C(7)-H(22)	3.3085		
			C(4)-C(2)-C(7)-C(8)	2.6189		
			C(4)-C(2)-C(7)-H(22)	-175.7588		
			C(2)-C(1)-O(20)-H(30)	8.7846		
			C(3)-C(1)-O(20)-H(30)	-169.7535		
			C(2)-C(1)-C(3)-C(5)	-0.753		
			C(2)-C(1)-C(3)-H(21)	-179.6737		
			O(20)-C(1)-C(3)-C(5)	177.8446		
			O(20)-C(1)-C(3)-H(21)	-1.0761		
			C(3)-C(1)-C(2)-C(7)	-176.8059		
			C(3)-C(1)-C(2)-C(4)	2.2569		
			O(20)-C(1)-C(2)-C(7)	4.6847		
			O(20)-C(1)-C(2)-C(4)	-176.2525		

Table (S4): Summary of Natural Population Analysis:

Atom	No	Natural Charge	Natural Population			
			Core	Valence	Rydberg	Total
C	1	0.36179	1.99860	3.62338	0.01624	5.63821
C	2	-0.09554	1.99880	4.08392	0.01283	6.09554
C	3	-0.28405	1.99899	4.27248	0.01258	6.28405
C	4	-0.03854	1.99883	4.02707	0.01264	6.03854
C	5	-0.15823	1.99899	4.14748	0.01176	6.15823
C	6	0.09752	1.99878	3.88001	0.02369	5.90248
C	7	-0.18411	1.99904	4.17088	0.01420	6.18411
C	8	-0.19923	1.99908	4.18695	0.01320	6.19923
C	9	-0.18722	1.99903	4.17407	0.01412	6.18722
C	10	-0.19372	1.99907	4.18150	0.01315	6.19372
N	11	-0.23399	1.99942	5.21141	0.02317	7.23399
N	12	-0.25605	1.99940	5.23289	0.02375	7.25605
C	13	0.03170	1.99882	3.94717	0.02231	5.96830
C	14	-0.19007	1.99903	4.17795	0.01309	6.19007
C	15	0.34550	1.99861	3.63864	0.01726	5.65450
C	16	-0.22227	1.99908	4.21016	0.01303	6.22227
C	17	-0.24276	1.99900	4.23008	0.01368	6.24276
C	18	-0.17167	1.99908	4.15979	0.01280	6.17167
O	19	-0.69995	1.99979	6.69686	0.00331	8.69995
O	20	-0.69907	1.99978	6.69608	0.00321	8.69907
H	21	0.22280	0.00000	0.77455	0.00265	0.77720
H	22	0.22881	0.00000	0.76569	0.00550	0.77119
H	23	0.22475	0.00000	0.77302	0.00224	0.77525
H	24	0.21403	0.00000	0.78473	0.00124	0.78597
H	25	0.22777	0.00000	0.76909	0.00314	0.77223
H	26	0.21378	0.00000	0.78501	0.00121	0.78622
H	27	0.22925	0.00000	0.76700	0.00376	0.77075
H	28	0.21481	0.00000	0.78333	0.00186	0.78519
H	29	0.22815	0.00000	0.76960	0.00225	0.77185
H	30	0.21879	0.00000	0.78010	0.00111	0.78121
H	31	0.49594	0.00000	0.50317	0.00089	0.50406
H	32	0.50109	0.00000	0.49513	0.00378	0.49891
=====						
* Total *		0.00000	39.98121	97.69917	0.31961	138.00000

Table (S5): NBO data of the Ligand:

Donor NBO (i)		Acceptor NBO (j)		kcal/mol	a.u.	a.u.
within unit 1						
1. BD (1) C	1 - C 2	/ 86. RY* (1) C	3	1.19	1.93	0.043
1. BD (1) C	1 - C 2	/ 94. RY* (1) C	4	1.16	2.27	0.046
1. BD (1) C	1 - C 2	/118. RY* (1) C	7	1.30	1.87	0.044
1. BD (1) C	1 - C 2	/255. BD* (1) C	1 - C 3	3.71	1.27	0.061
1. BD (1) C	1 - C 2	/258. BD* (1) C	2 - C 4	3.61	1.29	0.061
1. BD (1) C	1 - C 2	/260. BD* (1) C	2 - C 7	1.71	1.11	0.039
1. BD (1) C	1 - C 2	/262. BD* (1) C	3 - H 21	2.89	1.09	0.050
1. BD (1) C	1 - C 2	/264. BD* (1) C	4 - C 9	3.96	1.11	0.059
1. BD (1) C	1 - C 2	/269. BD* (1) C	7 - C 8	1.27	1.31	0.037
1. BD (1) C	1 - C 2	/296. BD* (1) O	20 - H 31	1.60	0.96	0.035
2. BD (1) C	1 - C 3	/ 79. RY* (2) C	2	1.83	2.23	0.057
2. BD (1) C	1 - C 3	/103. RY* (2) C	5	0.99	2.00	0.040
2. BD (1) C	1 - C 3	/254. BD* (1) C	1 - C 2	4.09	1.20	0.063
2. BD (1) C	1 - C 3	/260. BD* (1) C	2 - C 7	2.43	1.21	0.049
2. BD (1) C	1 - C 3	/261. BD* (1) C	3 - C 5	2.48	1.22	0.049
2. BD (1) C	1 - C 3	/262. BD* (1) C	3 - H 21	1.73	1.19	0.041
2. BD (1) C	1 - C 3	/267. BD* (1) C	5 - H 22	1.26	1.20	0.035
3. BD (2) C	1 - C 3	/ 81. RY* (4) C	2	0.80	0.87	0.025
3. BD (2) C	1 - C 3	/105. RY* (4) C	5	0.94	0.80	0.026
3. BD (2) C	1 - C 3	/222. RY* (1) O	20	0.74	1.13	0.028
3. BD (2) C	1 - C 3	/256. BD* (2) C	1 - C 3	0.99	0.31	0.016
3. BD (2) C	1 - C 3	/259. BD* (2) C	2 - C 4	11.24	0.32	0.056
3. BD (2) C	1 - C 3	/266. BD* (2) C	5 - C 6	15.57	0.32	0.063
4. BD (1) C	1 - O 20	/ 70. RY* (1) C	1	0.61	2.03	0.031
4. BD (1) C	1 - O 20	/ 79. RY* (2) C	2	0.67	2.37	0.036
4. BD (1) C	1 - O 20	/255. BD* (1) C	1 - C 3	1.23	1.52	0.039
4. BD (1) C	1 - O 20	/258. BD* (1) C	2 - C 4	1.27	1.53	0.039
4. BD (1) C	1 - O 20	/261. BD* (1) C	3 - C 5	2.12	1.36	0.048
5. BD (1) C	2 - C 4	/ 70. RY* (1) C	1	0.77	1.87	0.034
5. BD (1) C	2 - C 4	/111. RY* (2) C	6	0.96	2.09	0.040
5. BD (1) C	2 - C 4	/118. RY* (1) C	7	0.73	1.96	0.034
5. BD (1) C	2 - C 4	/134. RY* (1) C	9	0.80	1.96	0.036
5. BD (1) C	2 - C 4	/254. BD* (1) C	1 - C 2	3.63	1.18	0.058
5. BD (1) C	2 - C 4	/257. BD* (1) C	1 - O 20	2.71	1.01	0.047
5. BD (1) C	2 - C 4	/260. BD* (1) C	2 - C 7	4.04	1.19	0.062
5. BD (1) C	2 - C 4	/263. BD* (1) C	4 - C 6	3.78	1.19	0.060
5. BD (1) C	2 - C 4	/264. BD* (1) C	4 - C 9	3.59	1.19	0.059
5. BD (1) C	2 - C 4	/268. BD* (1) C	6 - N 11	2.67	1.12	0.049
5. BD (1) C	2 - C 4	/271. BD* (1) C	7 - H 23	1.36	1.19	0.036
5. BD (1) C	2 - C 4	/276. BD* (1) C	9 - H 25	1.37	1.19	0.036
6. BD (2) C	2 - C 4	/ 73. RY* (4) C	1	0.67	0.81	0.023
6. BD (2) C	2 - C 4	/113. RY* (4) C	6	0.83	0.83	0.025
6. BD (2) C	2 - C 4	/120. RY* (3) C	7	1.17	0.77	0.029
6. BD (2) C	2 - C 4	/136. RY* (3) C	9	1.10	0.78	0.029
6. BD (2) C	2 - C 4	/256. BD* (2) C	1 - C 3	14.68	0.29	0.059
6. BD (2) C	2 - C 4	/266. BD* (2) C	5 - C 6	11.81	0.30	0.054
6. BD (2) C	2 - C 4	/270. BD* (2) C	7 - C 8	12.06	0.31	0.057
6. BD (2) C	2 - C 4	/275. BD* (2) C	9 - C 10	11.08	0.31	0.055
7. BD (1) C	2 - C 7	/ 70. RY* (1) C	1	1.31	1.76	0.043
7. BD (1) C	2 - C 7	/ 94. RY* (1) C	4	0.61	2.25	0.033
7. BD (1) C	2 - C 7	/ 95. RY* (2) C	4	0.71	2.16	0.035
7. BD (1) C	2 - C 7	/126. RY* (1) C	8	0.70	1.82	0.032
7. BD (1) C	2 - C 7	/127. RY* (2) C	8	1.53	1.96	0.049
7. BD (1) C	2 - C 7	/254. BD* (1) C	1 - C 2	1.52	1.08	0.036
7. BD (1) C	2 - C 7	/255. BD* (1) C	1 - C 3	1.72	1.25	0.041
7. BD (1) C	2 - C 7	/258. BD* (1) C	2 - C 4	4.66	1.26	0.069
7. BD (1) C	2 - C 7	/263. BD* (1) C	4 - C 6	4.81	1.09	0.065
7. BD (1) C	2 - C 7	/269. BD* (1) C	7 - C 8	2.65	1.29	0.052
7. BD (1) C	2 - C 7	/271. BD* (1) C	7 - H 23	0.62	1.09	0.023
7. BD (1) C	2 - C 7	/273. BD* (1) C	8 - H 24	2.99	1.09	0.051
8. BD (1) C	3 - C 5	/ 71. RY* (2) C	1	1.46	2.07	0.049
8. BD (1) C	3 - C 5	/110. RY* (1) C	6	0.91	1.62	0.034
8. BD (1) C	3 - C 5	/111. RY* (2) C	6	1.01	1.99	0.040

8. BD (1) C 3 - C 5	/255. BD* (1) C 1 - C 3	2.66	1.26	0.052
8. BD (1) C 3 - C 5	/257. BD* (1) C 1 - O 20	5.41	0.92	0.063
8. BD (1) C 3 - C 5	/262. BD* (1) C 3 - H 21	0.76	1.08	0.026
8. BD (1) C 3 - C 5	/265. BD* (1) C 5 - C 6	2.76	1.27	0.053
8. BD (1) C 3 - C 5	/267. BD* (1) C 5 - H 22	0.56	1.09	0.022
8. BD (1) C 3 - C 5	/268. BD* (1) C 6 - N 11	5.26	1.03	0.066
9. BD (1) C 3 - H 21	/ 70. RY* (1) C 1	1.10	1.65	0.038
9. BD (1) C 3 - H 21	/ 74. RY* (5) C 1	0.60	9.46	0.068
9. BD (1) C 3 - H 21	/102. RY* (1) C 5	1.02	1.76	0.038
9. BD (1) C 3 - H 21	/254. BD* (1) C 1 - C 2	5.76	0.96	0.067
9. BD (1) C 3 - H 21	/255. BD* (1) C 1 - C 3	1.19	1.14	0.033
9. BD (1) C 3 - H 21	/257. BD* (1) C 1 - O 20	1.62	0.80	0.032
9. BD (1) C 3 - H 21	/265. BD* (1) C 5 - C 6	2.74	1.14	0.050
10. BD (1) C 4 - C 6	/ 78. RY* (1) C 2	1.44	2.29	0.051
10. BD (1) C 4 - C 6	/102. RY* (1) C 5	0.75	1.88	0.034
10. BD (1) C 4 - C 6	/134. RY* (1) C 9	1.21	1.87	0.043
10. BD (1) C 4 - C 6	/258. BD* (1) C 2 - C 4	3.50	1.27	0.060
10. BD (1) C 4 - C 6	/260. BD* (1) C 2 - C 7	4.08	1.10	0.060
10. BD (1) C 4 - C 6	/264. BD* (1) C 4 - C 9	1.82	1.10	0.040
10. BD (1) C 4 - C 6	/265. BD* (1) C 5 - C 6	3.98	1.27	0.064
10. BD (1) C 4 - C 6	/267. BD* (1) C 5 - H 22	2.87	1.09	0.050
10. BD (1) C 4 - C 6	/268. BD* (1) C 6 - N 11	1.17	1.02	0.031
10. BD (1) C 4 - C 6	/274. BD* (1) C 9 - C 10	1.28	1.30	0.037
10. BD (1) C 4 - C 6	/278. BD* (1) N 11 - N 12	2.49	1.17	0.048
11. BD (1) C 4 - C 9	/ 78. RY* (1) C 2	0.78	2.27	0.038
11. BD (1) C 4 - C 9	/110. RY* (1) C 6	1.00	1.61	0.036
11. BD (1) C 4 - C 9	/142. RY* (1) C 10	0.74	1.81	0.033
11. BD (1) C 4 - C 9	/143. RY* (2) C 10	1.48	1.98	0.049
11. BD (1) C 4 - C 9	/254. BD* (1) C 1 - C 2	4.87	1.07	0.065
11. BD (1) C 4 - C 9	/258. BD* (1) C 2 - C 4	4.36	1.26	0.066
11. BD (1) C 4 - C 9	/263. BD* (1) C 4 - C 6	1.68	1.09	0.038
11. BD (1) C 4 - C 9	/265. BD* (1) C 5 - C 6	1.95	1.26	0.044
11. BD (1) C 4 - C 9	/274. BD* (1) C 9 - C 10	2.71	1.29	0.053
11. BD (1) C 4 - C 9	/276. BD* (1) C 9 - H 25	0.67	1.09	0.024
11. BD (1) C 4 - C 9	/277. BD* (1) C 10 - H 26	3.00	1.09	0.051
12. BD (1) C 5 - C 6	/ 87. RY* (2) C 3	1.33	2.04	0.047
12. BD (1) C 5 - C 6	/ 95. RY* (2) C 4	1.52	2.27	0.053
12. BD (1) C 5 - C 6	/151. RY* (2) N 11	0.64	1.86	0.031
12. BD (1) C 5 - C 6	/261. BD* (1) C 3 - C 5	2.42	1.21	0.048
12. BD (1) C 5 - C 6	/262. BD* (1) C 3 - H 21	1.43	1.18	0.037
12. BD (1) C 5 - C 6	/263. BD* (1) C 4 - C 6	4.33	1.20	0.065
12. BD (1) C 5 - C 6	/264. BD* (1) C 4 - C 9	2.45	1.20	0.049
12. BD (1) C 5 - C 6	/267. BD* (1) C 5 - H 22	1.66	1.19	0.040
12. BD (1) C 5 - C 6	/268. BD* (1) C 6 - N 11	1.32	1.13	0.035
13. BD (2) C 5 - C 6	/ 88. RY* (3) C 3	0.95	0.78	0.026
13. BD (2) C 5 - C 6	/ 97. RY* (4) C 4	0.83	0.87	0.026
13. BD (2) C 5 - C 6	/152. RY* (3) N 11	0.84	0.99	0.028
13. BD (2) C 5 - C 6	/256. BD* (2) C 1 - C 3	11.78	0.30	0.053
13. BD (2) C 5 - C 6	/259. BD* (2) C 2 - C 4	14.20	0.32	0.061
13. BD (2) C 5 - C 6	/279. BD* (2) N 11 - N 12	19.46	0.23	0.060
14. BD (1) C 5 - H 22	/ 86. RY* (1) C 3	0.90	1.79	0.036
14. BD (1) C 5 - H 22	/110. RY* (1) C 6	1.05	1.49	0.035
14. BD (1) C 5 - H 22	/112. RY* (3) C 6	0.83	1.64	0.033
14. BD (1) C 5 - H 22	/255. BD* (1) C 1 - C 3	2.84	1.13	0.051
14. BD (1) C 5 - H 22	/262. BD* (1) C 3 - H 21	0.53	0.95	0.020
14. BD (1) C 5 - H 22	/263. BD* (1) C 4 - C 6	6.18	0.97	0.069
14. BD (1) C 5 - H 22	/265. BD* (1) C 5 - C 6	1.56	1.14	0.038
14. BD (1) C 5 - H 22	/268. BD* (1) C 6 - N 11	1.27	0.90	0.030
15. BD (1) C 6 - N 11	/ 95. RY* (2) C 4	1.06	2.30	0.044
15. BD (1) C 6 - N 11	/102. RY* (1) C 5	0.63	2.01	0.032
15. BD (1) C 6 - N 11	/158. RY* (1) N 12	1.04	1.71	0.038
15. BD (1) C 6 - N 11	/258. BD* (1) C 2 - C 4	1.45	1.40	0.040
15. BD (1) C 6 - N 11	/261. BD* (1) C 3 - C 5	2.31	1.23	0.048
15. BD (1) C 6 - N 11	/263. BD* (1) C 4 - C 6	0.93	1.22	0.030
15. BD (1) C 6 - N 11	/265. BD* (1) C 5 - C 6	2.21	1.40	0.050
15. BD (1) C 6 - N 11	/280. BD* (1) N 12 - C 13	4.62	1.15	0.065
16. BD (1) C 7 - C 8	/ 78. RY* (1) C 2	1.43	2.38	0.052
16. BD (1) C 7 - C 8	/ 79. RY* (2) C 2	0.63	2.21	0.033
16. BD (1) C 7 - C 8	/142. RY* (1) C 10	0.91	1.92	0.037
16. BD (1) C 7 - C 8	/143. RY* (2) C 10	0.82	2.08	0.037
16. BD (1) C 7 - C 8	/254. BD* (1) C 1 - C 2	2.54	1.18	0.049

16. BD (1) C 7 - C 8	/260. BD* (1) C 2 - C 7	2.84	1.19	0.052
16. BD (1) C 7 - C 8	/271. BD* (1) C 7 - H 23	1.64	1.19	0.040
16. BD (1) C 7 - C 8	/272. BD* (1) C 8 - C 10	2.36	1.21	0.048
16. BD (1) C 7 - C 8	/273. BD* (1) C 8 - H 24	1.61	1.19	0.039
16. BD (1) C 7 - C 8	/277. BD* (1) C 10 - H 26	1.57	1.19	0.039
17. BD (2) C 7 - C 8	/ 81. RY* (4) C 2	0.71	0.85	0.023
17. BD (2) C 7 - C 8	/144. RY* (3) C 10	0.98	0.77	0.026
17. BD (2) C 7 - C 8	/259. BD* (2) C 2 - C 4	12.56	0.31	0.058
17. BD (2) C 7 - C 8	/275. BD* (2) C 9 - C 10	13.29	0.31	0.058
18. BD (1) C 7 - H 23	/ 78. RY* (1) C 2	1.53	2.15	0.051
18. BD (1) C 7 - H 23	/126. RY* (1) C 8	1.24	1.69	0.041
18. BD (1) C 7 - H 23	/258. BD* (1) C 2 - C 4	3.47	1.14	0.056
18. BD (1) C 7 - H 23	/269. BD* (1) C 7 - C 8	1.50	1.16	0.037
18. BD (1) C 7 - H 23	/272. BD* (1) C 8 - C 10	5.50	0.98	0.065
18. BD (1) C 7 - H 23	/273. BD* (1) C 8 - H 24	1.00	0.96	0.028
19. BD (1) C 8 - C 10	/119. RY* (2) C 7	1.89	1.76	0.052
19. BD (1) C 8 - C 10	/135. RY* (2) C 9	1.90	1.73	0.051
19. BD (1) C 8 - C 10	/269. BD* (1) C 7 - C 8	2.61	1.29	0.052
19. BD (1) C 8 - C 10	/271. BD* (1) C 7 - H 23	3.20	1.09	0.053
19. BD (1) C 8 - C 10	/273. BD* (1) C 8 - H 24	0.58	1.09	0.022
19. BD (1) C 8 - C 10	/274. BD* (1) C 9 - C 10	2.59	1.29	0.052
19. BD (1) C 8 - C 10	/276. BD* (1) C 9 - H 25	3.19	1.09	0.053
19. BD (1) C 8 - C 10	/277. BD* (1) C 10 - H 26	0.57	1.09	0.022
20. BD (1) C 8 - H 24	/118. RY* (1) C 7	1.48	1.73	0.045
20. BD (1) C 8 - H 24	/142. RY* (1) C 10	1.21	1.69	0.040
20. BD (1) C 8 - H 24	/260. BD* (1) C 2 - C 7	6.03	0.96	0.068
20. BD (1) C 8 - H 24	/269. BD* (1) C 7 - C 8	1.40	1.16	0.036
20. BD (1) C 8 - H 24	/271. BD* (1) C 7 - H 23	0.99	0.96	0.028
20. BD (1) C 8 - H 24	/274. BD* (1) C 9 - C 10	2.76	1.16	0.051
21. BD (1) C 9 - C 10	/ 94. RY* (1) C 4	1.75	2.35	0.058
21. BD (1) C 9 - C 10	/126. RY* (1) C 8	0.91	1.92	0.037
21. BD (1) C 9 - C 10	/127. RY* (2) C 8	0.80	2.07	0.036
21. BD (1) C 9 - C 10	/263. BD* (1) C 4 - C 6	2.55	1.19	0.049
21. BD (1) C 9 - C 10	/264. BD* (1) C 4 - C 9	2.85	1.19	0.052
21. BD (1) C 9 - C 10	/272. BD* (1) C 8 - C 10	2.38	1.21	0.048
21. BD (1) C 9 - C 10	/273. BD* (1) C 8 - H 24	1.57	1.19	0.039
21. BD (1) C 9 - C 10	/276. BD* (1) C 9 - H 25	1.66	1.19	0.040
21. BD (1) C 9 - C 10	/277. BD* (1) C 10 - H 26	1.60	1.19	0.039
22. BD (2) C 9 - C 10	/ 97. RY* (4) C 4	0.68	0.86	0.023
22. BD (2) C 9 - C 10	/128. RY* (3) C 8	0.98	0.77	0.026
22. BD (2) C 9 - C 10	/259. BD* (2) C 2 - C 4	13.43	0.31	0.060
22. BD (2) C 9 - C 10	/270. BD* (2) C 7 - C 8	13.24	0.31	0.058
23. BD (1) C 9 - H 25	/ 94. RY* (1) C 4	1.52	2.12	0.051
23. BD (1) C 9 - H 25	/142. RY* (1) C 10	1.18	1.69	0.040
23. BD (1) C 9 - H 25	/258. BD* (1) C 2 - C 4	3.48	1.13	0.056
23. BD (1) C 9 - H 25	/272. BD* (1) C 8 - C 10	5.52	0.97	0.065
23. BD (1) C 9 - H 25	/274. BD* (1) C 9 - C 10	1.51	1.16	0.037
23. BD (1) C 9 - H 25	/277. BD* (1) C 10 - H 26	1.00	0.96	0.028
24. BD (1) C 10 - H 26	/126. RY* (1) C 8	1.21	1.69	0.041
24. BD (1) C 10 - H 26	/134. RY* (1) C 9	1.54	1.73	0.046
24. BD (1) C 10 - H 26	/264. BD* (1) C 4 - C 9	6.09	0.96	0.068
24. BD (1) C 10 - H 26	/269. BD* (1) C 7 - C 8	2.77	1.16	0.051
24. BD (1) C 10 - H 26	/274. BD* (1) C 9 - C 10	1.41	1.16	0.036
24. BD (1) C 10 - H 26	/276. BD* (1) C 9 - H 25	0.98	0.96	0.027
25. BD (1) N 11 - N 12	/110. RY* (1) C 6	0.76	1.95	0.034
25. BD (1) N 11 - N 12	/111. RY* (2) C 6	3.29	2.32	0.078
25. BD (1) N 11 - N 12	/151. RY* (2) N 11	0.58	2.08	0.031
25. BD (1) N 11 - N 12	/159. RY* (2) N 12	0.64	2.05	0.033
25. BD (1) N 11 - N 12	/166. RY* (1) C 13	1.12	1.99	0.042
25. BD (1) N 11 - N 12	/167. RY* (2) C 13	3.13	2.38	0.077
25. BD (1) N 11 - N 12	/263. BD* (1) C 4 - C 6	1.58	1.43	0.043
25. BD (1) N 11 - N 12	/282. BD* (1) C 13 - C 15	1.27	1.57	0.040
26. BD (2) N 11 - N 12	/113. RY* (4) C 6	0.87	0.93	0.026
26. BD (2) N 11 - N 12	/169. RY* (4) C 13	0.93	0.92	0.027
26. BD (2) N 11 - N 12	/266. BD* (2) C 5 - C 6	8.50	0.40	0.055
26. BD (2) N 11 - N 12	/283. BD* (2) C 13 - C 15	9.61	0.39	0.059
27. BD (1) N 12 - C 13	/150. RY* (1) N 11	1.06	1.76	0.039
27. BD (1) N 12 - C 13	/182. RY* (1) C 15	0.55	1.93	0.029
27. BD (1) N 12 - C 13	/183. RY* (2) C 15	0.69	2.10	0.034
27. BD (1) N 12 - C 13	/268. BD* (1) C 6 - N 11	4.71	1.16	0.066
27. BD (1) N 12 - C 13	/281. BD* (1) C 13 - C 14	1.05	1.22	0.032

27. BD (1) N 12 - C 13	/282. BD* (1) C 13 - C 15	1.79	1.37	0.044
27. BD (1) N 12 - C 13	/284. BD* (1) C 14 - C 16	0.88	1.43	0.032
27. BD (1) N 12 - C 13	/287. BD* (1) C 15 - C 17	2.51	1.21	0.049
28. BD (1) C 13 - C 14	/159. RY* (2) N 12	0.51	1.72	0.027
28. BD (1) C 13 - C 14	/183. RY* (2) C 15	1.34	1.97	0.046
28. BD (1) C 13 - C 14	/190. RY* (1) C 16	0.75	1.89	0.034
28. BD (1) C 13 - C 14	/191. RY* (2) C 16	1.56	1.98	0.050
28. BD (1) C 13 - C 14	/282. BD* (1) C 13 - C 15	4.03	1.24	0.063
28. BD (1) C 13 - C 14	/284. BD* (1) C 14 - C 16	2.65	1.29	0.052
28. BD (1) C 13 - C 14	/288. BD* (1) C 15 - O 19	5.19	0.92	0.062
28. BD (1) C 13 - C 14	/290. BD* (1) C 16 - H 28	3.14	1.09	0.052
29. BD (1) C 13 - C 15	/158. RY* (1) N 12	0.61	1.71	0.029
29. BD (1) C 13 - C 15	/174. RY* (1) C 14	0.73	2.04	0.035
29. BD (1) C 13 - C 15	/198. RY* (1) C 17	1.38	2.03	0.047
29. BD (1) C 13 - C 15	/278. BD* (1) N 11 - N 12	2.35	1.29	0.049
29. BD (1) C 13 - C 15	/280. BD* (1) N 12 - C 13	2.36	1.14	0.046
29. BD (1) C 13 - C 15	/281. BD* (1) C 13 - C 14	3.43	1.20	0.058
29. BD (1) C 13 - C 15	/286. BD* (1) C 14 - H 27	1.26	1.21	0.035
29. BD (1) C 13 - C 15	/287. BD* (1) C 15 - C 17	3.19	1.20	0.055
29. BD (1) C 13 - C 15	/293. BD* (1) C 17 - H 29	1.27	1.20	0.035
30. BD (2) C 13 - C 15	/160. RY* (3) N 12	0.87	0.99	0.028
30. BD (2) C 13 - C 15	/177. RY* (4) C 14	1.02	0.80	0.027
30. BD (2) C 13 - C 15	/200. RY* (3) C 17	0.98	0.79	0.027
30. BD (2) C 13 - C 15	/214. RY* (1) O 19	0.72	1.14	0.028
30. BD (2) C 13 - C 15	/279. BD* (2) N 11 - N 12	21.23	0.24	0.065
30. BD (2) C 13 - C 15	/283. BD* (2) C 13 - C 15	1.23	0.31	0.018
30. BD (2) C 13 - C 15	/285. BD* (2) C 14 - C 16	13.15	0.33	0.060
30. BD (2) C 13 - C 15	/292. BD* (2) C 17 - C 18	10.67	0.32	0.054
31. BD (1) C 14 - C 16	/166. RY* (1) C 13	1.38	1.76	0.044
31. BD (1) C 14 - C 16	/168. RY* (3) C 13	0.58	1.77	0.029
31. BD (1) C 14 - C 16	/206. RY* (1) C 18	0.99	1.94	0.039
31. BD (1) C 14 - C 16	/207. RY* (2) C 18	0.83	2.06	0.037
31. BD (1) C 14 - C 16	/280. BD* (1) N 12 - C 13	2.60	1.12	0.048
31. BD (1) C 14 - C 16	/281. BD* (1) C 13 - C 14	2.35	1.18	0.047
31. BD (1) C 14 - C 16	/286. BD* (1) C 14 - H 27	1.71	1.20	0.040
31. BD (1) C 14 - C 16	/289. BD* (1) C 16 - C 18	2.39	1.21	0.048
31. BD (1) C 14 - C 16	/290. BD* (1) C 16 - H 28	1.65	1.19	0.040
31. BD (1) C 14 - C 16	/294. BD* (1) C 18 - H 30	1.58	1.19	0.039
32. BD (2) C 14 - C 16	/169. RY* (4) C 13	0.69	0.82	0.022
32. BD (2) C 14 - C 16	/208. RY* (3) C 18	0.97	0.77	0.026
32. BD (2) C 14 - C 16	/283. BD* (2) C 13 - C 15	12.05	0.29	0.055
32. BD (2) C 14 - C 16	/292. BD* (2) C 17 - C 18	14.92	0.31	0.060
33. BD (1) C 14 - H 27	/166. RY* (1) C 13	0.81	1.53	0.031
33. BD (1) C 14 - H 27	/190. RY* (1) C 16	1.33	1.76	0.043
33. BD (1) C 14 - H 27	/280. BD* (1) N 12 - C 13	0.70	0.89	0.022
33. BD (1) C 14 - H 27	/282. BD* (1) C 13 - C 15	3.16	1.11	0.053
33. BD (1) C 14 - H 27	/284. BD* (1) C 14 - C 16	1.53	1.16	0.038
33. BD (1) C 14 - H 27	/289. BD* (1) C 16 - C 18	5.62	0.98	0.066
33. BD (1) C 14 - H 27	/290. BD* (1) C 16 - H 28	0.98	0.96	0.027
34. BD (1) C 15 - C 17	/167. RY* (2) C 13	1.48	2.07	0.050
34. BD (1) C 15 - C 17	/206. RY* (1) C 18	0.71	1.85	0.033
34. BD (1) C 15 - C 17	/207. RY* (2) C 18	1.46	1.97	0.048
34. BD (1) C 15 - C 17	/280. BD* (1) N 12 - C 13	4.97	1.04	0.064
34. BD (1) C 15 - C 17	/282. BD* (1) C 13 - C 15	3.77	1.26	0.061
34. BD (1) C 15 - C 17	/291. BD* (1) C 17 - C 18	2.38	1.30	0.050
34. BD (1) C 15 - C 17	/293. BD* (1) C 17 - H 29	0.56	1.10	0.022
34. BD (1) C 15 - C 17	/294. BD* (1) C 18 - H 30	2.87	1.11	0.050
34. BD (1) C 15 - C 17	/295. BD* (1) O 19 - H 32	1.79	0.98	0.037
35. BD (1) C 15 - O 19	/182. RY* (1) C 15	0.60	2.04	0.031
35. BD (1) C 15 - O 19	/281. BD* (1) C 13 - C 14	2.46	1.33	0.052
35. BD (1) C 15 - O 19	/282. BD* (1) C 13 - C 15	1.34	1.49	0.040
35. BD (1) C 15 - O 19	/291. BD* (1) C 17 - C 18	0.83	1.54	0.032
36. BD (1) C 16 - C 18	/175. RY* (2) C 14	1.45	1.77	0.045
36. BD (1) C 16 - C 18	/176. RY* (3) C 14	0.50	1.78	0.027
36. BD (1) C 16 - C 18	/199. RY* (2) C 17	1.91	1.90	0.054
36. BD (1) C 16 - C 18	/284. BD* (1) C 14 - C 16	2.57	1.30	0.052
36. BD (1) C 16 - C 18	/286. BD* (1) C 14 - H 27	3.14	1.10	0.052
36. BD (1) C 16 - C 18	/290. BD* (1) C 16 - H 28	0.59	1.09	0.023
36. BD (1) C 16 - C 18	/291. BD* (1) C 17 - C 18	2.44	1.29	0.050
36. BD (1) C 16 - C 18	/293. BD* (1) C 17 - H 29	3.28	1.08	0.053
36. BD (1) C 16 - C 18	/294. BD* (1) C 18 - H 30	0.59	1.09	0.023

37. BD (1) C 16 - H 28	/174. RY* (1) C 14	1.40	1.79	0.045
37. BD (1) C 16 - H 28	/206. RY* (1) C 18	1.04	1.70	0.038
37. BD (1) C 16 - H 28	/281. BD* (1) C 13 - C 14	5.82	0.95	0.067
37. BD (1) C 16 - H 28	/284. BD* (1) C 14 - C 16	1.50	1.16	0.037
37. BD (1) C 16 - H 28	/286. BD* (1) C 14 - H 27	0.94	0.96	0.027
37. BD (1) C 16 - H 28	/291. BD* (1) C 17 - C 18	2.66	1.15	0.050
38. BD (1) C 17 - C 18	/182. RY* (1) C 15	1.53	1.90	0.048
38. BD (1) C 17 - C 18	/190. RY* (1) C 16	1.06	2.00	0.041
38. BD (1) C 17 - C 18	/191. RY* (2) C 16	0.86	2.08	0.038
38. BD (1) C 17 - C 18	/287. BD* (1) C 15 - C 17	2.11	1.19	0.045
38. BD (1) C 17 - C 18	/288. BD* (1) C 15 - O 19	2.75	1.03	0.048
38. BD (1) C 17 - C 18	/289. BD* (1) C 16 - C 18	2.29	1.22	0.047
38. BD (1) C 17 - C 18	/290. BD* (1) C 16 - H 28	1.61	1.20	0.039
38. BD (1) C 17 - C 18	/293. BD* (1) C 17 - H 29	1.84	1.19	0.042
38. BD (1) C 17 - C 18	/294. BD* (1) C 18 - H 30	1.47	1.20	0.038
39. BD (2) C 17 - C 18	/185. RY* (4) C 15	0.56	0.81	0.020
39. BD (2) C 17 - C 18	/192. RY* (3) C 16	0.98	0.79	0.026
39. BD (2) C 17 - C 18	/283. BD* (2) C 13 - C 15	15.15	0.30	0.062
39. BD (2) C 17 - C 18	/285. BD* (2) C 14 - C 16	12.02	0.32	0.055
40. BD (1) C 17 - H 29	/182. RY* (1) C 15	0.57	1.67	0.028
40. BD (1) C 17 - H 29	/206. RY* (1) C 18	1.31	1.71	0.042
40. BD (1) C 17 - H 29	/282. BD* (1) C 13 - C 15	3.06	1.11	0.052
40. BD (1) C 17 - H 29	/288. BD* (1) C 15 - O 19	0.81	0.79	0.023
40. BD (1) C 17 - H 29	/289. BD* (1) C 16 - C 18	5.45	0.98	0.065
40. BD (1) C 17 - H 29	/291. BD* (1) C 17 - C 18	1.59	1.16	0.038
40. BD (1) C 17 - H 29	/294. BD* (1) C 18 - H 30	0.99	0.96	0.028
41. BD (1) C 18 - H 30	/190. RY* (1) C 16	1.30	1.76	0.043
41. BD (1) C 18 - H 30	/198. RY* (1) C 17	1.59	1.79	0.048
41. BD (1) C 18 - H 30	/284. BD* (1) C 14 - C 16	2.74	1.17	0.051
41. BD (1) C 18 - H 30	/287. BD* (1) C 15 - C 17	6.06	0.95	0.068
41. BD (1) C 18 - H 30	/291. BD* (1) C 17 - C 18	1.36	1.16	0.035
41. BD (1) C 18 - H 30	/293. BD* (1) C 17 - H 29	1.00	0.95	0.028
42. BD (1) O 19 - H 32	/182. RY* (1) C 15	2.86	1.87	0.065
42. BD (1) O 19 - H 32	/287. BD* (1) C 15 - C 17	4.00	1.15	0.061
43. BD (1) O 20 - H 31	/ 70. RY* (1) C 1	2.68	1.84	0.063
43. BD (1) O 20 - H 31	/254. BD* (1) C 1 - C 2	4.47	1.16	0.065
44. CR (1) C 1	/ 78. RY* (1) C 2	1.23	11.77	0.108
44. CR (1) C 1	/ 87. RY* (2) C 3	2.32	11.43	0.145
44. CR (1) C 1	/255. BD* (1) C 1 - C 3	1.92	10.75	0.129
44. CR (1) C 1	/257. BD* (1) C 1 - O 20	1.32	10.41	0.105
44. CR (1) C 1	/258. BD* (1) C 2 - C 4	0.69	10.76	0.078
44. CR (1) C 1	/260. BD* (1) C 2 - C 7	0.67	10.58	0.075
44. CR (1) C 1	/261. BD* (1) C 3 - C 5	0.94	10.59	0.089
45. CR (1) C 2	/ 71. RY* (2) C 1	1.59	11.48	0.121
45. CR (1) C 2	/ 94. RY* (1) C 4	1.07	11.67	0.100
45. CR (1) C 2	/ 95. RY* (2) C 4	1.79	11.58	0.128
45. CR (1) C 2	/119. RY* (2) C 7	1.09	11.18	0.099
45. CR (1) C 2	/255. BD* (1) C 1 - C 3	0.69	10.68	0.077
45. CR (1) C 2	/257. BD* (1) C 1 - O 20	0.59	10.34	0.070
45. CR (1) C 2	/258. BD* (1) C 2 - C 4	1.29	10.69	0.105
45. CR (1) C 2	/263. BD* (1) C 4 - C 6	1.21	10.51	0.101
45. CR (1) C 2	/264. BD* (1) C 4 - C 9	1.03	10.51	0.093
45. CR (1) C 2	/269. BD* (1) C 7 - C 8	0.56	10.71	0.069
46. CR (1) C 3	/ 70. RY* (1) C 1	1.99	11.17	0.133
46. CR (1) C 3	/ 72. RY* (3) C 1	1.38	11.31	0.111
46. CR (1) C 3	/104. RY* (3) C 5	0.96	11.53	0.094
46. CR (1) C 3	/231. RY* (2) H 21	0.58	12.28	0.076
46. CR (1) C 3	/254. BD* (1) C 1 - C 2	1.29	10.48	0.105
46. CR (1) C 3	/255. BD* (1) C 1 - C 3	0.64	10.66	0.074
46. CR (1) C 3	/257. BD* (1) C 1 - O 20	0.75	10.32	0.079
46. CR (1) C 3	/265. BD* (1) C 5 - C 6	0.67	10.67	0.076
47. CR (1) C 4	/ 78. RY* (1) C 2	0.56	11.70	0.072
47. CR (1) C 4	/ 79. RY* (2) C 2	2.22	11.53	0.143
47. CR (1) C 4	/111. RY* (2) C 6	0.75	11.41	0.083
47. CR (1) C 4	/112. RY* (3) C 6	0.53	11.18	0.068
47. CR (1) C 4	/135. RY* (2) C 9	1.12	11.15	0.100
47. CR (1) C 4	/254. BD* (1) C 1 - C 2	1.19	10.50	0.101
47. CR (1) C 4	/258. BD* (1) C 2 - C 4	1.20	10.69	0.102
47. CR (1) C 4	/260. BD* (1) C 2 - C 7	1.04	10.51	0.094
47. CR (1) C 4	/265. BD* (1) C 5 - C 6	0.68	10.68	0.076
47. CR (1) C 4	/268. BD* (1) C 6 - N 11	0.60	10.44	0.071

47. CR (1) C 4	/274. BD* (1) C 9 - C 10	0.56	10.71	0.069
48. CR (1) C 5	/ 89. RY* (4) C 3	0.54	11.26	0.069
48. CR (1) C 5	/111. RY* (2) C 6	3.03	11.39	0.166
48. CR (1) C 5	/233. RY* (2) H 22	0.54	12.29	0.073
48. CR (1) C 5	/255. BD* (1) C 1 - C 3	0.66	10.66	0.075
48. CR (1) C 5	/263. BD* (1) C 4 - C 6	1.30	10.50	0.105
48. CR (1) C 5	/265. BD* (1) C 5 - C 6	0.77	10.67	0.082
48. CR (1) C 5	/268. BD* (1) C 6 - N 11	0.81	10.43	0.083
49. CR (1) C 6	/ 94. RY* (1) C 4	1.15	11.70	0.104
49. CR (1) C 6	/ 96. RY* (3) C 4	0.50	11.19	0.067
49. CR (1) C 6	/103. RY* (2) C 5	2.16	11.33	0.140
49. CR (1) C 6	/258. BD* (1) C 2 - C 4	0.70	10.71	0.078
49. CR (1) C 6	/261. BD* (1) C 3 - C 5	0.95	10.55	0.090
49. CR (1) C 6	/264. BD* (1) C 4 - C 9	0.70	10.54	0.077
49. CR (1) C 6	/265. BD* (1) C 5 - C 6	1.50	10.71	0.114
49. CR (1) C 6	/278. BD* (1) N 11 - N 12	0.81	10.61	0.083
50. CR (1) C 7	/ 79. RY* (2) C 2	0.79	11.50	0.085
50. CR (1) C 7	/127. RY* (2) C 8	2.16	11.36	0.140
50. CR (1) C 7	/235. RY* (2) H 23	0.60	12.29	0.076
50. CR (1) C 7	/254. BD* (1) C 1 - C 2	0.66	10.47	0.075
50. CR (1) C 7	/258. BD* (1) C 2 - C 4	0.82	10.66	0.084
50. CR (1) C 7	/269. BD* (1) C 7 - C 8	0.84	10.69	0.085
50. CR (1) C 7	/272. BD* (1) C 8 - C 10	1.06	10.50	0.095
51. CR (1) C 8	/118. RY* (1) C 7	0.67	11.25	0.077
51. CR (1) C 8	/119. RY* (2) C 7	1.23	11.16	0.105
51. CR (1) C 8	/121. RY* (4) C 7	0.97	11.17	0.093
51. CR (1) C 8	/143. RY* (2) C 10	0.92	11.38	0.091
51. CR (1) C 8	/237. RY* (2) H 24	0.64	11.99	0.078
51. CR (1) C 8	/260. BD* (1) C 2 - C 7	1.17	10.49	0.100
51. CR (1) C 8	/269. BD* (1) C 7 - C 8	0.78	10.69	0.082
51. CR (1) C 8	/274. BD* (1) C 9 - C 10	0.62	10.69	0.073
52. CR (1) C 9	/ 95. RY* (2) C 4	1.07	11.56	0.099
52. CR (1) C 9	/143. RY* (2) C 10	2.13	11.38	0.139
52. CR (1) C 9	/239. RY* (2) H 25	0.54	12.28	0.073
52. CR (1) C 9	/258. BD* (1) C 2 - C 4	0.83	10.66	0.084
52. CR (1) C 9	/263. BD* (1) C 4 - C 6	0.68	10.49	0.076
52. CR (1) C 9	/272. BD* (1) C 8 - C 10	1.07	10.50	0.095
52. CR (1) C 9	/274. BD* (1) C 9 - C 10	0.84	10.69	0.085
53. CR (1) C 10	/127. RY* (2) C 8	0.90	11.36	0.090
53. CR (1) C 10	/134. RY* (1) C 9	0.71	11.26	0.080
53. CR (1) C 10	/135. RY* (2) C 9	1.15	11.12	0.101
53. CR (1) C 10	/137. RY* (4) C 9	1.04	11.21	0.096
53. CR (1) C 10	/241. RY* (2) H 26	0.62	11.97	0.077
53. CR (1) C 10	/264. BD* (1) C 4 - C 9	1.18	10.49	0.100
53. CR (1) C 10	/269. BD* (1) C 7 - C 8	0.62	10.69	0.073
53. CR (1) C 10	/274. BD* (1) C 9 - C 10	0.78	10.69	0.082
54. CR (1) N 11	/110. RY* (1) C 6	2.45	15.17	0.173
54. CR (1) N 11	/159. RY* (2) N 12	1.87	15.28	0.151
54. CR (1) N 11	/265. BD* (1) C 5 - C 6	0.70	14.82	0.091
54. CR (1) N 11	/280. BD* (1) N 12 - C 13	1.23	14.58	0.121
55. CR (1) N 12	/151. RY* (2) N 11	1.81	15.31	0.149
55. CR (1) N 12	/166. RY* (1) C 13	2.20	15.22	0.164
55. CR (1) N 12	/268. BD* (1) C 6 - N 11	1.26	14.58	0.122
55. CR (1) N 12	/281. BD* (1) C 13 - C 14	0.58	14.65	0.083
56. CR (1) C 13	/175. RY* (2) C 14	1.22	11.21	0.104
56. CR (1) C 13	/182. RY* (1) C 15	1.93	11.24	0.132
56. CR (1) C 13	/183. RY* (2) C 15	0.68	11.42	0.079
56. CR (1) C 13	/184. RY* (3) C 15	0.64	11.31	0.076
56. CR (1) C 13	/278. BD* (1) N 11 - N 12	0.81	10.62	0.083
56. CR (1) C 13	/282. BD* (1) C 13 - C 15	1.25	10.69	0.104
56. CR (1) C 13	/284. BD* (1) C 14 - C 16	0.57	10.74	0.070
56. CR (1) C 13	/287. BD* (1) C 15 - C 17	1.01	10.53	0.093
56. CR (1) C 13	/288. BD* (1) C 15 - O 19	0.79	10.37	0.081
57. CR (1) C 14	/167. RY* (2) C 13	1.50	11.45	0.117
57. CR (1) C 14	/191. RY* (2) C 16	2.21	11.37	0.142
57. CR (1) C 14	/243. RY* (2) H 27	0.53	12.28	0.072
57. CR (1) C 14	/282. BD* (1) C 13 - C 15	0.76	10.64	0.081
57. CR (1) C 14	/284. BD* (1) C 14 - C 16	0.89	10.69	0.087
57. CR (1) C 14	/289. BD* (1) C 16 - C 18	1.07	10.50	0.095
58. CR (1) C 15	/166. RY* (1) C 13	0.92	11.15	0.091
58. CR (1) C 15	/167. RY* (2) C 13	1.14	11.54	0.103

58. CR (1) C 15	/168. RY* (3) C 13	1.07	11.16	0.098
58. CR (1) C 15	/199. RY* (2) C 17	1.00	11.39	0.095
58. CR (1) C 15	/280. BD* (1) N 12 - C 13	1.00	10.52	0.092
58. CR (1) C 15	/281. BD* (1) C 13 - C 14	1.00	10.58	0.093
58. CR (1) C 15	/282. BD* (1) C 13 - C 15	1.73	10.73	0.123
58. CR (1) C 15	/288. BD* (1) C 15 - O 19	1.21	10.41	0.101
58. CR (1) C 15	/291. BD* (1) C 17 - C 18	0.56	10.78	0.069
59. CR (1) C 16	/174. RY* (1) C 14	0.58	11.31	0.073
59. CR (1) C 16	/175. RY* (2) C 14	0.55	11.16	0.070
59. CR (1) C 16	/176. RY* (3) C 14	1.70	11.17	0.123
59. CR (1) C 16	/207. RY* (2) C 18	0.93	11.35	0.092
59. CR (1) C 16	/245. RY* (2) H 28	0.66	12.29	0.081
59. CR (1) C 16	/281. BD* (1) C 13 - C 14	1.10	10.48	0.097
59. CR (1) C 16	/284. BD* (1) C 14 - C 16	0.78	10.69	0.082
59. CR (1) C 16	/291. BD* (1) C 17 - C 18	0.61	10.68	0.072
60. CR (1) C 17	/183. RY* (2) C 15	1.74	11.37	0.126
60. CR (1) C 17	/207. RY* (2) C 18	2.16	11.36	0.140
60. CR (1) C 17	/247. RY* (2) H 29	0.63	12.30	0.079
60. CR (1) C 17	/282. BD* (1) C 13 - C 15	0.82	10.64	0.084
60. CR (1) C 17	/288. BD* (1) C 15 - O 19	0.51	10.32	0.065
60. CR (1) C 17	/289. BD* (1) C 16 - C 18	1.04	10.51	0.094
60. CR (1) C 17	/291. BD* (1) C 17 - C 18	0.91	10.69	0.088
61. CR (1) C 18	/191. RY* (2) C 16	0.89	11.38	0.090
61. CR (1) C 18	/198. RY* (1) C 17	0.80	11.32	0.085
61. CR (1) C 18	/199. RY* (2) C 17	1.64	11.30	0.122
61. CR (1) C 18	/249. RY* (2) H 30	0.64	11.97	0.078
61. CR (1) C 18	/284. BD* (1) C 14 - C 16	0.61	10.70	0.072
61. CR (1) C 18	/287. BD* (1) C 15 - C 17	1.10	10.48	0.097
61. CR (1) C 18	/291. BD* (1) C 17 - C 18	0.72	10.69	0.079
62. CR (1) O 19	/182. RY* (1) C 15	2.24	20.10	0.190
62. CR (1) O 19	/253. RY* (2) H 32	0.58	20.97	0.099
62. CR (1) O 19	/282. BD* (1) C 13 - C 15	0.59	19.55	0.097
63. CR (1) O 20	/ 70. RY* (1) C 1	2.08	20.09	0.183
63. CR (1) O 20	/255. BD* (1) C 1 - C 3	0.55	19.57	0.093
64. LP (1) N 11	/110. RY* (1) C 6	3.30	1.34	0.060
64. LP (1) N 11	/158. RY* (1) N 12	1.62	1.31	0.042
64. LP (1) N 11	/159. RY* (2) N 12	1.04	1.45	0.035
64. LP (1) N 11	/263. BD* (1) C 4 - C 6	0.94	0.82	0.025
64. LP (1) N 11	/265. BD* (1) C 5 - C 6	6.93	0.99	0.074
65. LP (1) N 12	/150. RY* (1) N 11	2.06	1.36	0.048
65. LP (1) N 12	/151. RY* (2) N 11	1.08	1.48	0.036
65. LP (1) N 12	/166. RY* (1) C 13	2.84	1.39	0.057
65. LP (1) N 12	/281. BD* (1) C 13 - C 14	7.99	0.82	0.072
65. LP (1) N 12	/282. BD* (1) C 13 - C 15	0.80	0.97	0.025
65. LP (1) N 12	/295. BD* (1) O 19 - H 32	0.78	0.69	0.021
66. LP (1) O 19	/182. RY* (1) C 15	2.72	1.73	0.061
66. LP (1) O 19	/253. RY* (2) H 32	0.68	2.60	0.038
66. LP (1) O 19	/282. BD* (1) C 13 - C 15	6.19	1.18	0.076
67. LP (2) O 19	/283. BD* (2) C 13 - C 15	27.81	0.34	0.091
68. LP (1) O 20	/ 70. RY* (1) C 1	2.63	1.72	0.060
68. LP (1) O 20	/255. BD* (1) C 1 - C 3	5.28	1.20	0.071
69. LP (2) O 20	/256. BD* (2) C 1 - C 3	27.19	0.35	0.090
256. BD* (2) C 1 - C 3	/ 73. RY* (4) C 1	1.13	0.51	0.058
256. BD* (2) C 1 - C 3	/ 88. RY* (3) C 3	2.18	0.48	0.078
256. BD* (2) C 1 - C 3	/259. BD* (2) C 2 - C 4	132.62	0.02	0.073
259. BD* (2) C 2 - C 4	/ 81. RY* (4) C 2	2.32	0.54	0.075
259. BD* (2) C 2 - C 4	/ 97. RY* (4) C 4	2.31	0.55	0.075
266. BD* (2) C 5 - C 6	/105. RY* (4) C 5	2.35	0.48	0.082
266. BD* (2) C 5 - C 6	/113. RY* (4) C 6	1.23	0.52	0.062
270. BD* (2) C 7 - C 8	/120. RY* (3) C 7	1.43	0.45	0.075
270. BD* (2) C 7 - C 8	/128. RY* (3) C 8	1.44	0.46	0.077
275. BD* (2) C 9 - C 10	/136. RY* (3) C 9	1.40	0.47	0.076
275. BD* (2) C 9 - C 10	/144. RY* (3) C 10	1.43	0.46	0.076
279. BD* (2) N 11 - N 12	/152. RY* (3) N 11	1.04	0.75	0.070
279. BD* (2) N 11 - N 12	/160. RY* (3) N 12	0.98	0.75	0.067
279. BD* (2) N 11 - N 12	/266. BD* (2) C 5 - C 6	19.28	0.08	0.068
279. BD* (2) N 11 - N 12	/283. BD* (2) C 13 - C 15	21.60	0.07	0.061
283. BD* (2) C 13 - C 15	/169. RY* (4) C 13	1.71	0.53	0.065
283. BD* (2) C 13 - C 15	/185. RY* (4) C 15	1.72	0.51	0.064
283. BD* (2) C 13 - C 15	/285. BD* (2) C 14 - C 16	76.25	0.02	0.069
283. BD* (2) C 13 - C 15	/292. BD* (2) C 17 - C 18	116.70	0.02	0.072

285. BD* (2) C 14 - C 16	/177. RY* (4) C 14	1.47	0.47	0.076
285. BD* (2) C 14 - C 16	/192. RY* (3) C 16	1.48	0.47	0.076
292. BD* (2) C 17 - C 18	/200. RY* (3) C 17	1.58	0.46	0.076
292. BD* (2) C 17 - C 18	/208. RY* (3) C 18	1.56	0.47	0.076

Table (S6): Summary of Natural Population Analysis of the Cr(III) complex:

Atom	No	Natural Charge	Natural Population			Total
			Core	Valence	Rydberg	
C	1	0.30563	1.99867	3.67636	0.01934	5.69437
C	2	-0.08377	1.99885	4.06796	0.01696	6.08377
C	3	-0.18124	1.99903	4.16339	0.01882	6.18124
C	4	-0.03697	1.99885	4.02025	0.01787	6.03697
C	5	-0.22476	1.99894	4.20641	0.01942	6.22476
C	6	0.14716	1.99876	3.83109	0.02298	5.85284
C	7	-0.24318	1.99907	4.22811	0.01600	6.24318
C	8	-0.15576	1.99909	4.13997	0.01669	6.15576
C	9	-0.22471	1.99905	4.21058	0.01508	6.22471
C	10	-0.21236	1.99907	4.19690	0.01638	6.21236
N	11	-0.35329	1.99930	5.32830	0.02569	7.35329
N	12	-0.01950	1.99916	5.00064	0.01971	7.01950
C	13	0.06139	1.99873	3.92217	0.01771	5.93861
C	14	-0.23787	1.99905	4.21456	0.02426	6.23787
C	15	0.29681	1.99867	3.68396	0.02056	5.70319
C	16	-0.17334	1.99912	4.15446	0.01976	6.17334
C	17	-0.23487	1.99901	4.21354	0.02232	6.23487
C	18	-0.21038	1.99913	4.19655	0.01470	6.21038
O	19	-0.68886	1.99982	6.68378	0.00526	8.68886
O	20	-0.64528	1.99979	6.64168	0.00381	8.64528
H	21	0.21917	0.00000	0.77836	0.00248	0.78083
H	22	0.20463	0.00000	0.79255	0.00282	0.79537
H	23	0.20517	0.00000	0.79306	0.00177	0.79483
H	24	0.22735	0.00000	0.76972	0.00293	0.77265
H	25	0.20564	0.00000	0.79257	0.00180	0.79436
H	26	0.20955	0.00000	0.78813	0.00232	0.79045
H	27	0.20757	0.00000	0.79125	0.00118	0.79243
H	28	0.21477	0.00000	0.78263	0.00259	0.78523
H	29	0.21489	0.00000	0.78391	0.00120	0.78511
H	30	0.49914	0.00000	0.49975	0.00112	0.50086
H	31	0.22841	0.00000	0.76871	0.00287	0.77159
C	32	0.30866	1.99868	3.67359	0.01908	5.69134
C	33	-0.08055	1.99883	4.06454	0.01717	6.08055
C	34	-0.16290	1.99902	4.14537	0.01851	6.16290
C	35	-0.07761	1.99886	4.05967	0.01908	6.07761
C	36	-0.20982	1.99893	4.18983	0.02106	6.20982
C	37	0.13207	1.99875	3.84524	0.02395	5.86793
C	38	-0.23789	1.99907	4.22290	0.01593	6.23789
C	39	-0.16951	1.99909	4.15414	0.01627	6.16951
C	40	-0.25088	1.99906	4.23580	0.01602	6.25088
C	41	-0.19769	1.99907	4.18109	0.01753	6.19769
N	42	-0.39542	1.99931	5.36935	0.02676	7.39542
N	43	0.00615	1.99915	4.97485	0.01985	6.99385
C	44	0.05764	1.99873	3.92590	0.01772	5.94236
C	45	-0.22003	1.99905	4.19782	0.02315	6.22003
C	46	0.30565	1.99867	3.67526	0.02042	5.69435
C	47	-0.16798	1.99912	4.14916	0.01970	6.16798
C	48	-0.22603	1.99901	4.20597	0.02105	6.22603
C	49	-0.20794	1.99913	4.19454	0.01427	6.20794
O	50	-0.64330	1.99979	6.63974	0.00378	8.64330
H	51	0.22254	0.00000	0.77498	0.00248	0.77746
H	52	0.20665	0.00000	0.79058	0.00277	0.79335
H	53	0.20894	0.00000	0.78985	0.00121	0.79106
H	54	0.21106	0.00000	0.78553	0.00341	0.78894
H	55	0.20760	0.00000	0.79059	0.00181	0.79240
H	56	0.21213	0.00000	0.78565	0.00222	0.78787
H	57	0.20894	0.00000	0.78988	0.00118	0.79106
H	58	0.21794	0.00000	0.77991	0.00215	0.78206
H	59	0.21743	0.00000	0.78140	0.00117	0.78257
H	60	0.49998	0.00000	0.49885	0.00117	0.50002
H	61	0.22623	0.00000	0.77050	0.00327	0.77377
O	62	-0.70008	1.99981	6.69514	0.00513	8.70008
Cr	63	1.44120	17.98827	4.54653	0.02399	22.55880

H	64	0.51876	0.00000	0.47826	0.00298	0.48124
H	65	0.51891	0.00000	0.47765	0.00344	0.48109
O	66	-0.86176	1.99979	6.85673	0.00524	8.86176
Cl	67	-0.64023	9.99978	7.63787	0.00258	17.64023
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* Total *		0.00000	109.95011	214.25595	0.79394	325.00000

Table (S7): NBO data of the Cr(III) complex:

Donor NBO (i)		Acceptor NBO (j)		kcal/mol	a.u.	a.u.
within unit 1						
1.	BD (1) C	1 - C	2	/198. RY* (2) C	3	0.55 1.99 0.042
1.	BD (1) C	1 - C	2	/199. RY* (3) C	3	0.25 2.02 0.029
1.	BD (1) C	1 - C	2	/206. RY* (2) C	4	0.73 2.25 0.052
1.	BD (1) C	1 - C	2	/229. RY* (1) C	7	0.41 1.53 0.032
1.	BD (1) C	1 - C	2	/230. RY* (2) C	7	0.41 1.53 0.032
1.	BD (1) C	1 - C	2	/592. BD* (1) C	1 - C	3 1.47 1.26 0.055
1.	BD (1) C	1 - C	2	/594. BD* (1) C	2 - C	4 1.20 1.27 0.049
1.	BD (1) C	1 - C	2	/595. BD* (1) C	2 - C	7 0.61 1.10 0.033
1.	BD (1) C	1 - C	2	/597. BD* (1) C	3 - H	21 1.36 1.09 0.049
1.	BD (1) C	1 - C	2	/599. BD* (1) C	4 - C	9 2.02 1.12 0.060
1.	BD (1) C	1 - C	2	/603. BD* (1) C	7 - C	8 0.70 1.30 0.038
2.	BD (2) C	1 - C	2	/ 9. BD* (2) C	3 - C	5 5.68 0.23 0.066
2.	BD (2) C	1 - C	2	/ 13. BD* (2) C	4 - C	6 5.28 0.24 0.062
2.	BD (2) C	1 - C	2	/197. RY* (1) C	3	1.20 0.77 0.039
2.	BD (2) C	1 - C	2	/205. RY* (1) C	4	0.80 0.87 0.034
2.	BD (2) C	1 - C	2	/333. RY* (1) O	20	0.38 1.15 0.027
2.	BD (2) C	1 - C	2	/604. BD* (2) C	7 - C	8 6.60 0.30 0.064
3.	BD (1) C	1 - C	3	/191. RY* (3) C	2	0.69 2.23 0.050
3.	BD (1) C	1 - C	3	/215. RY* (3) C	5	0.50 2.01 0.040
3.	BD (1) C	1 - C	3	/590. BD* (1) C	1 - C	2 1.44 1.19 0.053
3.	BD (1) C	1 - C	3	/593. BD* (1) C	1 - O	20 0.28 1.04 0.021
3.	BD (1) C	1 - C	3	/595. BD* (1) C	2 - C	7 1.13 1.22 0.047
3.	BD (1) C	1 - C	3	/596. BD* (1) C	3 - C	5 0.73 1.23 0.038
3.	BD (1) C	1 - C	3	/597. BD* (1) C	3 - H	21 0.60 1.20 0.034
3.	BD (1) C	1 - C	3	/601. BD* (1) C	5 - H	31 0.64 1.22 0.035
3.	BD (1) C	1 - C	3	/626. BD* (1) O	20 - H	30 0.74 1.08 0.036
4.	BD (1) C	1 - O	20	/198. RY* (2) C	3	0.29 2.24 0.032
4.	BD (1) C	1 - O	20	/592. BD* (1) C	1 - C	3 0.37 1.51 0.030
4.	BD (1) C	1 - O	20	/594. BD* (1) C	2 - C	4 0.56 1.52 0.037
4.	BD (1) C	1 - O	20	/596. BD* (1) C	3 - C	5 0.99 1.36 0.046
5.	BD (1) C	2 - C	4	/183. RY* (3) C	1	0.45 2.21 0.040
5.	BD (1) C	2 - C	4	/222. RY* (2) C	6	0.48 1.95 0.039
5.	BD (1) C	2 - C	4	/229. RY* (1) C	7	0.29 1.62 0.028
5.	BD (1) C	2 - C	4	/245. RY* (1) C	9	0.39 1.98 0.035
5.	BD (1) C	2 - C	4	/590. BD* (1) C	1 - C	2 1.21 1.16 0.047
5.	BD (1) C	2 - C	4	/593. BD* (1) C	1 - O	20 1.32 1.02 0.046
5.	BD (1) C	2 - C	4	/595. BD* (1) C	2 - C	7 1.57 1.19 0.055
5.	BD (1) C	2 - C	4	/598. BD* (1) C	4 - C	6 1.50 1.20 0.054
5.	BD (1) C	2 - C	4	/599. BD* (1) C	4 - C	9 1.48 1.21 0.053
5.	BD (1) C	2 - C	4	/602. BD* (1) C	6 - N	11 1.44 1.12 0.051
5.	BD (1) C	2 - C	4	/605. BD* (1) C	7 - H	22 0.76 1.19 0.038
5.	BD (1) C	2 - C	4	/610. BD* (1) C	9 - H	24 0.69 1.22 0.037
6.	BD (1) C	2 - C	7	/183. RY* (3) C	1	0.35 2.10 0.034
6.	BD (1) C	2 - C	7	/206. RY* (2) C	4	0.48 2.24 0.042
6.	BD (1) C	2 - C	7	/238. RY* (2) C	8	0.32 1.98 0.032
6.	BD (1) C	2 - C	7	/239. RY* (3) C	8	0.70 1.99 0.047
6.	BD (1) C	2 - C	7	/590. BD* (1) C	1 - C	2 0.48 1.05 0.029
6.	BD (1) C	2 - C	7	/592. BD* (1) C	1 - C	3 0.73 1.25 0.038
6.	BD (1) C	2 - C	7	/594. BD* (1) C	2 - C	4 1.77 1.25 0.060
6.	BD (1) C	2 - C	7	/598. BD* (1) C	4 - C	6 2.41 1.09 0.065
6.	BD (1) C	2 - C	7	/603. BD* (1) C	7 - C	8 1.26 1.28 0.051
6.	BD (1) C	2 - C	7	/605. BD* (1) C	7 - H	22 0.30 1.08 0.023
6.	BD (1) C	2 - C	7	/607. BD* (1) C	8 - H	23 1.36 1.08 0.049
7.	BD (1) C	3 - C	5	/182. RY* (2) C	1	0.50 1.57 0.036
7.	BD (1) C	3 - C	5	/183. RY* (3) C	1	0.37 2.12 0.036
7.	BD (1) C	3 - C	5	/223. RY* (3) C	6	0.97 1.76 0.053
7.	BD (1) C	3 - C	5	/592. BD* (1) C	1 - C	3 0.84 1.26 0.041
7.	BD (1) C	3 - C	5	/593. BD* (1) C	1 - O	20 2.68 0.92 0.063
7.	BD (1) C	3 - C	5	/600. BD* (1) C	5 - C	6 1.08 1.29 0.047
7.	BD (1) C	3 - C	5	/602. BD* (1) C	6 - N	11 2.94 1.02 0.069
8.	BD (2) C	3 - C	5	/ 13. BD* (2) C	4 - C	6 6.01 0.23 0.067

8. BD (2) C 3 - C 5	/181. RY* (1) C 1	0.28	1.04	0.022
8. BD (2) C 3 - C 5	/221. RY* (1) C 6	0.30	1.14	0.024
8. BD (2) C 3 - C 5	/591. BD* (2) C 1 - C 2	7.08	0.23	0.069
9. BD* (2) C 3 - C 5	/197. RY* (1) C 3	1.76	0.54	0.051
9. BD* (2) C 3 - C 5	/213. RY* (1) C 5	1.68	0.59	0.052
10. BD (1) C 3 - H 21	/183. RY* (3) C 1	0.47	1.99	0.039
10. BD (1) C 3 - H 21	/186. RY* (6) C 1	0.29	8.73	0.064
10. BD (1) C 3 - H 21	/214. RY* (2) C 5	0.55	1.92	0.041
10. BD (1) C 3 - H 21	/590. BD* (1) C 1 - C 2	3.08	0.94	0.068
10. BD (1) C 3 - H 21	/592. BD* (1) C 1 - C 3	0.39	1.13	0.027
10. BD (1) C 3 - H 21	/593. BD* (1) C 1 - O 20	0.94	0.79	0.034
10. BD (1) C 3 - H 21	/600. BD* (1) C 5 - C 6	1.44	1.15	0.051
10. BD (1) C 3 - H 21	/601. BD* (1) C 5 - H 31	0.29	0.97	0.021
11. BD (1) C 4 - C 6	/190. RY* (2) C 2	0.77	2.34	0.054
11. BD (1) C 4 - C 6	/214. RY* (2) C 5	0.39	2.05	0.036
11. BD (1) C 4 - C 6	/245. RY* (1) C 9	0.62	1.89	0.044
11. BD (1) C 4 - C 6	/261. RY* (1) N 11	0.54	1.52	0.036
11. BD (1) C 4 - C 6	/594. BD* (1) C 2 - C 4	1.44	1.27	0.054
11. BD (1) C 4 - C 6	/595. BD* (1) C 2 - C 7	1.97	1.10	0.059
11. BD (1) C 4 - C 6	/599. BD* (1) C 4 - C 9	0.77	1.11	0.037
11. BD (1) C 4 - C 6	/600. BD* (1) C 5 - C 6	1.59	1.28	0.057
11. BD (1) C 4 - C 6	/601. BD* (1) C 5 - H 31	1.43	1.10	0.050
11. BD (1) C 4 - C 6	/602. BD* (1) C 6 - N 11	0.29	1.02	0.022
11. BD (1) C 4 - C 6	/608. BD* (1) C 9 - C 10	0.71	1.31	0.039
12. BD (2) C 4 - C 6	/ 9. BD* (2) C 3 - C 5	6.14	0.22	0.066
12. BD (2) C 4 - C 6	/189. RY* (1) C 2	0.90	0.86	0.036
12. BD (2) C 4 - C 6	/213. RY* (1) C 5	1.21	0.81	0.041
12. BD (2) C 4 - C 6	/246. RY* (2) C 9	0.57	0.88	0.029
12. BD (2) C 4 - C 6	/591. BD* (2) C 1 - C 2	6.68	0.22	0.065
12. BD (2) C 4 - C 6	/609. BD* (2) C 9 - C 10	7.25	0.31	0.066
12. BD (2) C 4 - C 6	/612. BD* (1) N 11 - N 12	2.11	0.76	0.052
13. BD* (2) C 4 - C 6	/189. RY* (1) C 2	0.77	0.62	0.038
13. BD* (2) C 4 - C 6	/205. RY* (1) C 4	1.65	0.63	0.055
13. BD* (2) C 4 - C 6	/213. RY* (1) C 5	1.37	0.58	0.048
13. BD* (2) C 4 - C 6	/221. RY* (1) C 6	0.72	0.90	0.043
13. BD* (2) C 4 - C 6	/222. RY* (2) C 6	0.27	1.25	0.032
13. BD* (2) C 4 - C 6	/246. RY* (2) C 9	0.44	0.65	0.029
13. BD* (2) C 4 - C 6	/604. BD* (2) C 7 - C 8	0.28	0.06	0.006
13. BD* (2) C 4 - C 6	/609. BD* (2) C 9 - C 10	38.89	0.08	0.078
13. BD* (2) C 4 - C 6	/612. BD* (1) N 11 - N 12	2.40	0.53	0.060
14. BD (1) C 4 - C 9	/190. RY* (2) C 2	0.41	2.32	0.039
14. BD (1) C 4 - C 9	/221. RY* (1) C 6	0.45	1.49	0.033
14. BD (1) C 4 - C 9	/222. RY* (2) C 6	0.80	1.83	0.049
14. BD (1) C 4 - C 9	/254. RY* (2) C 10	0.31	2.01	0.032
14. BD (1) C 4 - C 9	/255. RY* (3) C 10	0.74	2.01	0.049
14. BD (1) C 4 - C 9	/590. BD* (1) C 1 - C 2	2.40	1.05	0.063
14. BD (1) C 4 - C 9	/594. BD* (1) C 2 - C 4	1.72	1.25	0.059
14. BD (1) C 4 - C 9	/598. BD* (1) C 4 - C 6	0.75	1.08	0.036
14. BD (1) C 4 - C 9	/600. BD* (1) C 5 - C 6	1.02	1.26	0.045
14. BD (1) C 4 - C 9	/608. BD* (1) C 9 - C 10	1.28	1.29	0.052
14. BD (1) C 4 - C 9	/610. BD* (1) C 9 - H 24	0.34	1.10	0.025
14. BD (1) C 4 - C 9	/611. BD* (1) C 10 - H 25	1.45	1.08	0.050
15. BD (1) C 5 - C 6	/199. RY* (3) C 3	0.65	2.13	0.047
15. BD (1) C 5 - C 6	/207. RY* (3) C 4	0.70	2.36	0.051
15. BD (1) C 5 - C 6	/261. RY* (1) N 11	0.70	1.63	0.043
15. BD (1) C 5 - C 6	/596. BD* (1) C 3 - C 5	0.94	1.22	0.043
15. BD (1) C 5 - C 6	/597. BD* (1) C 3 - H 21	0.59	1.19	0.034
15. BD (1) C 5 - C 6	/598. BD* (1) C 4 - C 6	1.65	1.21	0.057
15. BD (1) C 5 - C 6	/599. BD* (1) C 4 - C 9	1.09	1.22	0.046
15. BD (1) C 5 - C 6	/601. BD* (1) C 5 - H 31	0.81	1.21	0.040
15. BD (1) C 5 - C 6	/602. BD* (1) C 6 - N 11	0.71	1.13	0.036
15. BD (1) C 5 - C 6	/612. BD* (1) N 11 - N 12	0.51	1.24	0.032
16. BD (1) C 5 - H 31	/198. RY* (2) C 3	0.45	1.85	0.037
16. BD (1) C 5 - H 31	/221. RY* (1) C 6	0.42	1.36	0.030
16. BD (1) C 5 - H 31	/222. RY* (2) C 6	0.79	1.71	0.047
16. BD (1) C 5 - H 31	/592. BD* (1) C 1 - C 3	1.36	1.12	0.049
16. BD (1) C 5 - H 31	/597. BD* (1) C 3 - H 21	0.29	0.94	0.021
16. BD (1) C 5 - H 31	/598. BD* (1) C 4 - C 6	3.32	0.96	0.072
16. BD (1) C 5 - H 31	/600. BD* (1) C 5 - C 6	0.71	1.14	0.036
16. BD (1) C 5 - H 31	/602. BD* (1) C 6 - N 11	0.69	0.88	0.031
17. BD (1) C 6 - N 11	/207. RY* (3) C 4	0.68	2.39	0.051

17. BD (1) C 6 - N 11	/214. RY* (2) C 5	0.60	2.19	0.046
17. BD (1) C 6 - N 11	/269. RY* (1) N 12	0.46	1.58	0.034
17. BD (1) C 6 - N 11	/270. RY* (2) N 12	0.65	1.61	0.041
17. BD (1) C 6 - N 11	/594. BD* (1) C 2 - C 4	0.72	1.41	0.040
17. BD (1) C 6 - N 11	/596. BD* (1) C 3 - C 5	1.22	1.25	0.049
17. BD (1) C 6 - N 11	/598. BD* (1) C 4 - C 6	0.46	1.24	0.030
17. BD (1) C 6 - N 11	/600. BD* (1) C 5 - C 6	0.86	1.42	0.044
17. BD (1) C 6 - N 11	/612. BD* (1) N 11 - N 12	0.60	1.27	0.035
17. BD (1) C 6 - N 11	/613. BD* (1) N 12 - C 13	2.81	1.14	0.072
18. BD (1) C 7 - C 8	/190. RY* (2) C 2	0.49	2.43	0.044
18. BD (1) C 7 - C 8	/191. RY* (3) C 2	0.45	2.20	0.040
18. BD (1) C 7 - C 8	/254. RY* (2) C 10	0.44	2.12	0.038
18. BD (1) C 7 - C 8	/255. RY* (3) C 10	0.39	2.13	0.037
18. BD (1) C 7 - C 8	/590. BD* (1) C 1 - C 2	1.26	1.16	0.049
18. BD (1) C 7 - C 8	/595. BD* (1) C 2 - C 7	1.35	1.19	0.051
18. BD (1) C 7 - C 8	/605. BD* (1) C 7 - H 22	0.80	1.19	0.039
18. BD (1) C 7 - C 8	/606. BD* (1) C 8 - C 10	0.94	1.21	0.043
18. BD (1) C 7 - C 8	/607. BD* (1) C 8 - H 23	0.65	1.19	0.035
18. BD (1) C 7 - C 8	/611. BD* (1) C 10 - H 25	0.78	1.20	0.039
19. BD (2) C 7 - C 8	/189. RY* (1) C 2	0.40	0.86	0.024
19. BD (2) C 7 - C 8	/253. RY* (1) C 10	0.73	0.78	0.031
19. BD (2) C 7 - C 8	/591. BD* (2) C 1 - C 2	3.47	0.23	0.049
19. BD (2) C 7 - C 8	/604. BD* (2) C 7 - C 8	0.36	0.30	0.015
19. BD (2) C 7 - C 8	/609. BD* (2) C 9 - C 10	4.59	0.32	0.054
20. BD (1) C 7 - H 22	/190. RY* (2) C 2	0.67	2.20	0.049
20. BD (1) C 7 - H 22	/238. RY* (2) C 8	0.72	1.86	0.046
20. BD (1) C 7 - H 22	/594. BD* (1) C 2 - C 4	1.59	1.13	0.054
20. BD (1) C 7 - H 22	/603. BD* (1) C 7 - C 8	0.65	1.16	0.035
20. BD (1) C 7 - H 22	/606. BD* (1) C 8 - C 10	2.59	0.99	0.064
20. BD (1) C 7 - H 22	/607. BD* (1) C 8 - H 23	0.46	0.96	0.027
21. BD (1) C 8 - C 10	/231. RY* (3) C 7	0.94	1.78	0.052
21. BD (1) C 8 - C 10	/247. RY* (3) C 9	0.92	1.77	0.051
21. BD (1) C 8 - C 10	/603. BD* (1) C 7 - C 8	1.05	1.29	0.047
21. BD (1) C 8 - C 10	/605. BD* (1) C 7 - H 22	1.66	1.09	0.054
21. BD (1) C 8 - C 10	/608. BD* (1) C 9 - C 10	1.11	1.30	0.048
21. BD (1) C 8 - C 10	/610. BD* (1) C 9 - H 24	1.66	1.11	0.054
22. BD (1) C 8 - H 23	/229. RY* (1) C 7	0.44	1.38	0.031
22. BD (1) C 8 - H 23	/230. RY* (2) C 7	0.45	1.38	0.032
22. BD (1) C 8 - H 23	/254. RY* (2) C 10	0.58	1.88	0.042
22. BD (1) C 8 - H 23	/595. BD* (1) C 2 - C 7	3.22	0.95	0.070
22. BD (1) C 8 - H 23	/603. BD* (1) C 7 - C 8	0.56	1.16	0.032
22. BD (1) C 8 - H 23	/605. BD* (1) C 7 - H 22	0.57	0.96	0.030
22. BD (1) C 8 - H 23	/608. BD* (1) C 9 - C 10	1.32	1.17	0.050
22. BD (1) C 8 - H 23	/611. BD* (1) C 10 - H 25	0.26	0.96	0.020
23. BD (1) C 9 - C 10	/206. RY* (2) C 4	0.63	2.34	0.048
23. BD (1) C 9 - C 10	/207. RY* (3) C 4	0.34	2.33	0.036
23. BD (1) C 9 - C 10	/238. RY* (2) C 8	0.38	2.08	0.036
23. BD (1) C 9 - C 10	/239. RY* (3) C 8	0.38	2.09	0.036
23. BD (1) C 9 - C 10	/598. BD* (1) C 4 - C 6	1.20	1.18	0.048
23. BD (1) C 9 - C 10	/599. BD* (1) C 4 - C 9	1.33	1.20	0.051
23. BD (1) C 9 - C 10	/606. BD* (1) C 8 - C 10	1.02	1.20	0.044
23. BD (1) C 9 - C 10	/607. BD* (1) C 8 - H 23	0.76	1.18	0.038
23. BD (1) C 9 - C 10	/610. BD* (1) C 9 - H 24	0.87	1.21	0.041
23. BD (1) C 9 - C 10	/611. BD* (1) C 10 - H 25	0.76	1.19	0.038
24. BD (2) C 9 - C 10	/ 13. BD* (2) C 4 - C 6	3.31	0.23	0.049
24. BD (2) C 9 - C 10	/205. RY* (1) C 4	0.35	0.86	0.022
24. BD (2) C 9 - C 10	/237. RY* (1) C 8	0.68	0.76	0.029
24. BD (2) C 9 - C 10	/604. BD* (2) C 7 - C 8	4.88	0.29	0.055
24. BD (2) C 9 - C 10	/609. BD* (2) C 9 - C 10	0.41	0.30	0.016
25. BD (1) C 9 - H 24	/206. RY* (2) C 4	0.66	2.10	0.047
25. BD (1) C 9 - H 24	/254. RY* (2) C 10	0.66	1.87	0.045
25. BD (1) C 9 - H 24	/594. BD* (1) C 2 - C 4	1.72	1.12	0.055
25. BD (1) C 9 - H 24	/606. BD* (1) C 8 - C 10	2.69	0.97	0.065
25. BD (1) C 9 - H 24	/608. BD* (1) C 9 - C 10	0.80	1.16	0.039
25. BD (1) C 9 - H 24	/611. BD* (1) C 10 - H 25	0.47	0.95	0.027
26. BD (1) C 10 - H 25	/238. RY* (2) C 8	0.54	1.85	0.040
26. BD (1) C 10 - H 25	/245. RY* (1) C 9	0.68	1.74	0.044
26. BD (1) C 10 - H 25	/599. BD* (1) C 4 - C 9	3.16	0.96	0.070
26. BD (1) C 10 - H 25	/603. BD* (1) C 7 - C 8	1.32	1.15	0.049
26. BD (1) C 10 - H 25	/607. BD* (1) C 8 - H 23	0.26	0.95	0.020
26. BD (1) C 10 - H 25	/608. BD* (1) C 9 - C 10	0.70	1.16	0.036

26. BD (1) C 10 - H 25	/610. BD* (1) C 9 - H 24	0.47	0.97	0.027
27. BD (1) N 11 - N 12	/221. RY* (1) C 6	1.02	1.85	0.055
27. BD (1) N 11 - N 12	/277. RY* (1) C 13	0.74	2.11	0.050
27. BD (1) N 11 - N 12	/278. RY* (2) C 13	0.79	2.42	0.055
27. BD (1) N 11 - N 12	/602. BD* (1) C 6 - N 11	0.61	1.36	0.037
27. BD (1) N 11 - N 12	/613. BD* (1) N 12 - C 13	0.41	1.35	0.030
27. BD (1) N 11 - N 12	/614. BD* (1) C 13 - C 14	0.56	1.43	0.036
28. BD (1) N 12 - C 13	/221. RY* (1) C 6	0.31	1.65	0.029
28. BD (1) N 12 - C 13	/262. RY* (2) N 11	0.89	2.73	0.063
28. BD (1) N 12 - C 13	/263. RY* (3) N 11	0.63	2.14	0.047
28. BD (1) N 12 - C 13	/286. RY* (2) C 14	0.36	2.11	0.035
28. BD (1) N 12 - C 13	/293. RY* (1) C 15	0.41	1.98	0.036
28. BD (1) N 12 - C 13	/602. BD* (1) C 6 - N 11	2.43	1.17	0.068
28. BD (1) N 12 - C 13	/614. BD* (1) C 13 - C 14	0.30	1.24	0.024
28. BD (1) N 12 - C 13	/615. BD* (1) C 13 - C 15	0.95	1.40	0.046
28. BD (1) N 12 - C 13	/617. BD* (1) C 14 - C 16	0.40	1.44	0.030
28. BD (1) N 12 - C 13	/619. BD* (1) C 15 - C 17	1.19	1.24	0.049
29. BD (1) C 13 - C 14	/269. RY* (1) N 12	0.34	1.45	0.028
29. BD (1) C 13 - C 14	/295. RY* (3) C 15	0.34	2.17	0.034
29. BD (1) C 13 - C 14	/296. RY* (4) C 15	0.25	1.66	0.026
29. BD (1) C 13 - C 14	/302. RY* (2) C 16	0.38	1.88	0.034
29. BD (1) C 13 - C 14	/303. RY* (3) C 16	0.69	2.03	0.048
29. BD (1) C 13 - C 14	/612. BD* (1) N 11 - N 12	1.60	1.13	0.054
29. BD (1) C 13 - C 14	/613. BD* (1) N 12 - C 13	0.38	1.01	0.025
29. BD (1) C 13 - C 14	/615. BD* (1) C 13 - C 15	1.81	1.25	0.060
29. BD (1) C 13 - C 14	/617. BD* (1) C 14 - C 16	0.89	1.29	0.043
29. BD (1) C 13 - C 14	/620. BD* (1) C 15 - O 19	2.64	0.97	0.064
29. BD (1) C 13 - C 14	/622. BD* (1) C 16 - H 27	1.41	1.10	0.050
30. BD (1) C 13 - C 15	/271. RY* (3) N 12	0.30	1.89	0.030
30. BD (1) C 13 - C 15	/286. RY* (2) C 14	0.43	2.08	0.038
30. BD (1) C 13 - C 15	/310. RY* (2) C 17	0.63	2.10	0.046
30. BD (1) C 13 - C 15	/613. BD* (1) N 12 - C 13	0.68	1.13	0.035
30. BD (1) C 13 - C 15	/614. BD* (1) C 13 - C 14	1.77	1.21	0.059
30. BD (1) C 13 - C 15	/618. BD* (1) C 14 - H 26	0.55	1.21	0.033
30. BD (1) C 13 - C 15	/619. BD* (1) C 15 - C 17	1.43	1.21	0.053
30. BD (1) C 13 - C 15	/620. BD* (1) C 15 - O 19	0.34	1.09	0.024
30. BD (1) C 13 - C 15	/624. BD* (1) C 17 - H 28	0.57	1.22	0.033
31. BD (2) C 13 - C 15	/ 36. BD* (1) C 16 - C 17	0.50	0.13	0.053
31. BD (2) C 13 - C 15	/270. RY* (2) N 12	0.26	1.12	0.022
31. BD (2) C 13 - C 15	/285. RY* (1) C 14	0.76	0.80	0.031
31. BD (2) C 13 - C 15	/309. RY* (1) C 17	1.19	0.81	0.039
31. BD (2) C 13 - C 15	/325. RY* (1) O 19	0.37	1.40	0.029
31. BD (2) C 13 - C 15	/616. BD* (2) C 13 - C 15	0.48	0.32	0.017
32. BD (1) C 14 - C 16	/278. RY* (2) C 13	0.99	2.19	0.059
32. BD (1) C 14 - C 16	/317. RY* (1) C 18	0.54	1.95	0.041
32. BD (1) C 14 - C 16	/319. RY* (3) C 18	0.43	2.06	0.038
32. BD (1) C 14 - C 16	/613. BD* (1) N 12 - C 13	1.41	1.11	0.050
32. BD (1) C 14 - C 16	/614. BD* (1) C 13 - C 14	0.86	1.20	0.041
32. BD (1) C 14 - C 16	/618. BD* (1) C 14 - H 26	0.66	1.19	0.035
32. BD (1) C 14 - C 16	/621. BD* (1) C 16 - C 18	0.85	1.22	0.041
32. BD (1) C 14 - C 16	/622. BD* (1) C 16 - H 27	0.54	1.20	0.032
32. BD (1) C 14 - C 16	/625. BD* (1) C 18 - H 29	0.84	1.21	0.040
33. BD (1) C 14 - H 26	/277. RY* (1) C 13	0.43	1.63	0.033
33. BD (1) C 14 - H 26	/302. RY* (2) C 16	0.61	1.73	0.041
33. BD (1) C 14 - H 26	/613. BD* (1) N 12 - C 13	0.37	0.86	0.023
33. BD (1) C 14 - H 26	/615. BD* (1) C 13 - C 15	1.75	1.11	0.056
33. BD (1) C 14 - H 26	/617. BD* (1) C 14 - C 16	0.50	1.15	0.030
33. BD (1) C 14 - H 26	/621. BD* (1) C 16 - C 18	2.72	0.98	0.065
33. BD (1) C 14 - H 26	/622. BD* (1) C 16 - H 27	0.54	0.96	0.029
34. BD (1) C 15 - C 17	/277. RY* (1) C 13	0.57	1.78	0.040
34. BD (1) C 15 - C 17	/317. RY* (1) C 18	0.34	1.85	0.032
34. BD (1) C 15 - C 17	/319. RY* (3) C 18	0.78	1.95	0.050
34. BD (1) C 15 - C 17	/613. BD* (1) N 12 - C 13	2.80	1.01	0.067
34. BD (1) C 15 - C 17	/615. BD* (1) C 13 - C 15	1.64	1.25	0.057
34. BD (1) C 15 - C 17	/623. BD* (1) C 17 - C 18	0.96	1.30	0.045
34. BD (1) C 15 - C 17	/625. BD* (1) C 18 - H 29	1.56	1.11	0.053
35. BD (1) C 15 - O 19	/277. RY* (1) C 13	0.27	1.97	0.029
35. BD (1) C 15 - O 19	/293. RY* (1) C 15	0.50	2.02	0.040
35. BD (1) C 15 - O 19	/614. BD* (1) C 13 - C 14	1.27	1.29	0.051
35. BD (1) C 15 - O 19	/615. BD* (1) C 13 - C 15	0.67	1.45	0.039
35. BD (1) C 15 - O 19	/623. BD* (1) C 17 - C 18	0.47	1.49	0.034

36. BD* (1) C 16 - C 17	/285. RY* (1) C 14	1.51	0.67	0.040
36. BD* (1) C 16 - C 17	/294. RY* (2) C 15	0.28	0.81	0.019
36. BD* (1) C 16 - C 17	/616. BD* (2) C 13 - C 15	5.86	0.19	0.047
37. BD (1) C 16 - C 17	/285. RY* (1) C 14	0.75	0.63	0.037
37. BD (1) C 16 - C 17	/318. RY* (2) C 18	1.55	0.65	0.055
37. BD (1) C 16 - C 17	/616. BD* (2) C 13 - C 15	10.59	0.15	0.058
38. BD (1) C 16 - C 18	/287. RY* (3) C 14	0.84	1.93	0.051
38. BD (1) C 16 - C 18	/311. RY* (3) C 17	0.84	1.90	0.051
38. BD (1) C 16 - C 18	/617. BD* (1) C 14 - C 16	0.90	1.28	0.043
38. BD (1) C 16 - C 18	/618. BD* (1) C 14 - H 26	1.59	1.08	0.053
38. BD (1) C 16 - C 18	/623. BD* (1) C 17 - C 18	1.07	1.29	0.047
38. BD (1) C 16 - C 18	/624. BD* (1) C 17 - H 28	1.54	1.09	0.052
38. BD (1) C 16 - C 18	/625. BD* (1) C 18 - H 29	0.27	1.10	0.022
39. BD (1) C 16 - H 27	/286. RY* (2) C 14	0.76	1.82	0.047
39. BD (1) C 16 - H 27	/317. RY* (1) C 18	0.65	1.71	0.042
39. BD (1) C 16 - H 27	/614. BD* (1) C 13 - C 14	3.00	0.95	0.068
39. BD (1) C 16 - H 27	/617. BD* (1) C 14 - C 16	0.43	1.15	0.028
39. BD (1) C 16 - H 27	/618. BD* (1) C 14 - H 26	0.54	0.95	0.029
39. BD (1) C 16 - H 27	/623. BD* (1) C 17 - C 18	1.45	1.16	0.052
39. BD (1) C 16 - H 27	/625. BD* (1) C 18 - H 29	0.25	0.97	0.020
40. BD (1) C 17 - C 18	/293. RY* (1) C 15	0.59	1.93	0.043
40. BD (1) C 17 - C 18	/295. RY* (3) C 15	0.47	2.26	0.041
40. BD (1) C 17 - C 18	/302. RY* (2) C 16	0.51	1.97	0.040
40. BD (1) C 17 - C 18	/303. RY* (3) C 16	0.38	2.13	0.036
40. BD (1) C 17 - C 18	/619. BD* (1) C 15 - C 17	0.77	1.19	0.038
40. BD (1) C 17 - C 18	/620. BD* (1) C 15 - O 19	1.23	1.06	0.046
40. BD (1) C 17 - C 18	/621. BD* (1) C 16 - C 18	1.00	1.22	0.044
40. BD (1) C 17 - C 18	/622. BD* (1) C 16 - H 27	0.73	1.20	0.037
40. BD (1) C 17 - C 18	/624. BD* (1) C 17 - H 28	0.75	1.19	0.038
40. BD (1) C 17 - C 18	/625. BD* (1) C 18 - H 29	0.70	1.20	0.037
41. BD (1) C 17 - H 28	/293. RY* (1) C 15	0.37	1.69	0.032
41. BD (1) C 17 - H 28	/317. RY* (1) C 18	0.67	1.71	0.043
41. BD (1) C 17 - H 28	/615. BD* (1) C 13 - C 15	1.61	1.11	0.054
41. BD (1) C 17 - H 28	/620. BD* (1) C 15 - O 19	0.53	0.82	0.027
41. BD (1) C 17 - H 28	/621. BD* (1) C 16 - C 18	2.90	0.98	0.067
41. BD (1) C 17 - H 28	/623. BD* (1) C 17 - C 18	0.61	1.16	0.034
41. BD (1) C 17 - H 28	/625. BD* (1) C 18 - H 29	0.52	0.96	0.028
42. BD (1) C 18 - H 29	/302. RY* (2) C 16	0.59	1.73	0.041
42. BD (1) C 18 - H 29	/310. RY* (2) C 17	0.75	1.83	0.047
42. BD (1) C 18 - H 29	/617. BD* (1) C 14 - C 16	1.30	1.15	0.049
42. BD (1) C 18 - H 29	/619. BD* (1) C 15 - C 17	2.84	0.95	0.066
42. BD (1) C 18 - H 29	/623. BD* (1) C 17 - C 18	0.61	1.16	0.034
42. BD (1) C 18 - H 29	/624. BD* (1) C 17 - H 28	0.48	0.95	0.027
43. BD (1) O 20 - H 30	/181. RY* (1) C 1	0.64	1.49	0.039
43. BD (1) O 20 - H 30	/182. RY* (2) C 1	0.84	1.65	0.047
43. BD (1) O 20 - H 30	/592. BD* (1) C 1 - C 3	1.59	1.34	0.058
89. CR (1) C 1	/190. RY* (2) C 2	0.56	11.81	0.102
89. CR (1) C 1	/199. RY* (3) C 3	1.08	11.49	0.140
89. CR (1) C 1	/592. BD* (1) C 1 - C 3	0.92	10.73	0.126
89. CR (1) C 1	/593. BD* (1) C 1 - O 20	0.66	10.39	0.106
89. CR (1) C 1	/594. BD* (1) C 2 - C 4	0.29	10.74	0.071
89. CR (1) C 1	/595. BD* (1) C 2 - C 7	0.31	10.57	0.073
89. CR (1) C 1	/596. BD* (1) C 3 - C 5	0.40	10.58	0.082
90. CR (1) C 2	/182. RY* (2) C 1	0.32	10.98	0.075
90. CR (1) C 2	/207. RY* (3) C 4	1.15	11.66	0.146
90. CR (1) C 2	/231. RY* (3) C 7	0.51	11.20	0.096
90. CR (1) C 2	/592. BD* (1) C 1 - C 3	0.27	10.67	0.069
90. CR (1) C 2	/594. BD* (1) C 2 - C 4	0.60	10.68	0.101
90. CR (1) C 2	/598. BD* (1) C 4 - C 6	0.53	10.51	0.096
90. CR (1) C 2	/599. BD* (1) C 4 - C 9	0.46	10.53	0.088
90. CR (1) C 2	/603. BD* (1) C 7 - C 8	0.27	10.71	0.068
91. CR (1) C 3	/183. RY* (3) C 1	1.52	11.51	0.167
91. CR (1) C 3	/216. RY* (4) C 5	0.47	11.20	0.092
91. CR (1) C 3	/342. RY* (2) H 21	0.29	12.28	0.075
91. CR (1) C 3	/590. BD* (1) C 1 - C 2	0.59	10.46	0.100
91. CR (1) C 3	/592. BD* (1) C 1 - C 3	0.29	10.66	0.070
91. CR (1) C 3	/593. BD* (1) C 1 - O 20	0.36	10.32	0.077
91. CR (1) C 3	/600. BD* (1) C 5 - C 6	0.31	10.68	0.073
92. CR (1) C 4	/191. RY* (3) C 2	1.21	11.51	0.149
92. CR (1) C 4	/223. RY* (3) C 6	0.79	11.16	0.118
92. CR (1) C 4	/247. RY* (3) C 9	0.57	11.18	0.101

92.	CR	(1)	C	4	/590.	BD*	(1)	C	1 - C	2	0.51	10.47	0.094
92.	CR	(1)	C	4	/594.	BD*	(1)	C	2 - C	4	0.52	10.67	0.095
92.	CR	(1)	C	4	/595.	BD*	(1)	C	2 - C	7	0.46	10.50	0.088
92.	CR	(1)	C	4	/600.	BD*	(1)	C	5 - C	6	0.32	10.69	0.075
92.	CR	(1)	C	4	/602.	BD*	(1)	C	6 - N	11	0.27	10.42	0.067
92.	CR	(1)	C	4	/608.	BD*	(1)	C	9 - C	10	0.28	10.72	0.069
93.	CR	(1)	C	5	/222.	RY*	(2)	C	6		0.54	11.23	0.099
93.	CR	(1)	C	5	/223.	RY*	(3)	C	6		0.71	11.14	0.112
93.	CR	(1)	C	5	/224.	RY*	(4)	C	6		0.49	11.07	0.093
93.	CR	(1)	C	5	/362.	RY*	(2)	H	31		0.26	12.34	0.071
93.	CR	(1)	C	5	/592.	BD*	(1)	C	1 - C	3	0.29	10.64	0.070
93.	CR	(1)	C	5	/598.	BD*	(1)	C	4 - C	6	0.64	10.48	0.105
93.	CR	(1)	C	5	/600.	BD*	(1)	C	5 - C	6	0.43	10.67	0.086
93.	CR	(1)	C	5	/602.	BD*	(1)	C	6 - N	11	0.44	10.40	0.086
94.	CR	(1)	C	6	/206.	RY*	(2)	C	4		0.58	11.69	0.104
94.	CR	(1)	C	6	/215.	RY*	(3)	C	5		1.11	11.33	0.142
94.	CR	(1)	C	6	/594.	BD*	(1)	C	2 - C	4	0.33	10.70	0.076
94.	CR	(1)	C	6	/596.	BD*	(1)	C	3 - C	5	0.43	10.55	0.086
94.	CR	(1)	C	6	/599.	BD*	(1)	C	4 - C	9	0.31	10.55	0.072
94.	CR	(1)	C	6	/600.	BD*	(1)	C	5 - C	6	0.71	10.72	0.111
94.	CR	(1)	C	6	/612.	BD*	(1)	N	11 - N	12	0.40	10.57	0.082
95.	CR	(1)	C	7	/190.	RY*	(2)	C	2		0.28	11.72	0.073
95.	CR	(1)	C	7	/239.	RY*	(3)	C	8		1.04	11.39	0.137
95.	CR	(1)	C	7	/344.	RY*	(2)	H	22		0.27	12.26	0.073
95.	CR	(1)	C	7	/590.	BD*	(1)	C	1 - C	2	0.33	10.45	0.075
95.	CR	(1)	C	7	/594.	BD*	(1)	C	2 - C	4	0.36	10.65	0.078
95.	CR	(1)	C	7	/603.	BD*	(1)	C	7 - C	8	0.38	10.68	0.081
95.	CR	(1)	C	7	/606.	BD*	(1)	C	8 - C	10	0.48	10.50	0.090
96.	CR	(1)	C	8	/231.	RY*	(3)	C	7		0.66	11.17	0.108
96.	CR	(1)	C	8	/232.	RY*	(4)	C	7		0.39	11.07	0.083
96.	CR	(1)	C	8	/255.	RY*	(3)	C	10		0.41	11.41	0.086
96.	CR	(1)	C	8	/346.	RY*	(2)	H	23		0.33	12.30	0.080
96.	CR	(1)	C	8	/595.	BD*	(1)	C	2 - C	7	0.60	10.48	0.101
96.	CR	(1)	C	8	/603.	BD*	(1)	C	7 - C	8	0.39	10.68	0.081
96.	CR	(1)	C	8	/608.	BD*	(1)	C	9 - C	10	0.28	10.69	0.069
97.	CR	(1)	C	9	/207.	RY*	(3)	C	4		0.36	11.62	0.082
97.	CR	(1)	C	9	/255.	RY*	(3)	C	10		1.00	11.40	0.135
97.	CR	(1)	C	9	/348.	RY*	(2)	H	24		0.27	12.36	0.072
97.	CR	(1)	C	9	/594.	BD*	(1)	C	2 - C	4	0.40	10.64	0.083
97.	CR	(1)	C	9	/598.	BD*	(1)	C	4 - C	6	0.32	10.47	0.074
97.	CR	(1)	C	9	/606.	BD*	(1)	C	8 - C	10	0.51	10.49	0.093
97.	CR	(1)	C	9	/608.	BD*	(1)	C	9 - C	10	0.40	10.68	0.083
98.	CR	(1)	C	10	/239.	RY*	(3)	C	8		0.38	11.38	0.083
98.	CR	(1)	C	10	/245.	RY*	(1)	C	9		0.37	11.26	0.082
98.	CR	(1)	C	10	/247.	RY*	(3)	C	9		0.56	11.15	0.099
98.	CR	(1)	C	10	/248.	RY*	(4)	C	9		0.54	11.13	0.098
98.	CR	(1)	C	10	/350.	RY*	(2)	H	25		0.33	12.30	0.080
98.	CR	(1)	C	10	/599.	BD*	(1)	C	4 - C	9	0.60	10.48	0.101
98.	CR	(1)	C	10	/603.	BD*	(1)	C	7 - C	8	0.29	10.67	0.070
98.	CR	(1)	C	10	/608.	BD*	(1)	C	9 - C	10	0.43	10.68	0.086
99.	CR	(1)	N	11	/221.	RY*	(1)	C	6		0.38	15.07	0.096
99.	CR	(1)	N	11	/222.	RY*	(2)	C	6		0.68	15.42	0.129
99.	CR	(1)	N	11	/271.	RY*	(3)	N	12		0.85	15.34	0.144
99.	CR	(1)	N	11	/598.	BD*	(1)	C	4 - C	6	0.27	14.67	0.080
99.	CR	(1)	N	11	/600.	BD*	(1)	C	5 - C	6	0.26	14.85	0.078
99.	CR	(1)	N	11	/613.	BD*	(1)	N	12 - C	13	0.70	14.57	0.129
100.	CR	(1)	N	12	/263.	RY*	(3)	N	11		0.94	15.60	0.153
100.	CR	(1)	N	12	/277.	RY*	(1)	C	13		0.67	15.38	0.128
100.	CR	(1)	N	12	/278.	RY*	(2)	C	13		0.33	15.69	0.091
100.	CR	(1)	N	12	/602.	BD*	(1)	C	6 - N	11	0.60	14.62	0.120
100.	CR	(1)	N	12	/615.	BD*	(1)	C	13 - C	15	0.34	14.86	0.091
101.	CR	(1)	C	13	/287.	RY*	(3)	C	14		0.42	11.39	0.088
101.	CR	(1)	C	13	/293.	RY*	(1)	C	15		0.58	11.29	0.103
101.	CR	(1)	C	13	/295.	RY*	(3)	C	15		0.97	11.62	0.134
101.	CR	(1)	C	13	/612.	BD*	(1)	N	11 - N	12	0.47	10.59	0.090
101.	CR	(1)	C	13	/615.	BD*	(1)	C	13 - C	15	0.62	10.71	0.104
101.	CR	(1)	C	13	/619.	BD*	(1)	C	15 - C	17	0.46	10.55	0.088
101.	CR	(1)	C	13	/620.	BD*	(1)	C	15 - O	19	0.41	10.42	0.083
102.	CR	(1)	C	14	/277.	RY*	(1)	C	13		0.42	11.15	0.087
102.	CR	(1)	C	14	/279.	RY*	(3)	C	13		0.31	10.95	0.074
102.	CR	(1)	C	14	/303.	RY*	(3)	C	16		0.99	11.41	0.134

102. CR (1) C 14	/352. RY* (2) H 26	0.30	12.30	0.076
102. CR (1) C 14	/613. BD* (1) N 12 - C 13	0.26	10.38	0.066
102. CR (1) C 14	/615. BD* (1) C 13 - C 15	0.39	10.63	0.082
102. CR (1) C 14	/617. BD* (1) C 14 - C 16	0.40	10.67	0.083
102. CR (1) C 14	/621. BD* (1) C 16 - C 18	0.45	10.50	0.088
103. CR (1) C 15	/278. RY* (2) C 13	1.57	11.54	0.170
103. CR (1) C 15	/311. RY* (3) C 17	0.48	11.37	0.093
103. CR (1) C 15	/613. BD* (1) N 12 - C 13	0.47	10.47	0.089
103. CR (1) C 15	/614. BD* (1) C 13 - C 14	0.48	10.55	0.091
103. CR (1) C 15	/615. BD* (1) C 13 - C 15	0.83	10.71	0.120
103. CR (1) C 15	/620. BD* (1) C 15 - O 19	0.45	10.43	0.087
103. CR (1) C 15	/623. BD* (1) C 17 - C 18	0.26	10.76	0.067
104. CR (1) C 16	/286. RY* (2) C 14	0.28	11.35	0.071
104. CR (1) C 16	/287. RY* (3) C 14	0.76	11.32	0.117
104. CR (1) C 16	/288. RY* (4) C 14	0.27	10.96	0.069
104. CR (1) C 16	/319. RY* (3) C 18	0.50	11.34	0.095
104. CR (1) C 16	/354. RY* (2) H 27	0.31	11.95	0.076
104. CR (1) C 16	/614. BD* (1) C 13 - C 14	0.49	10.48	0.091
104. CR (1) C 16	/617. BD* (1) C 14 - C 16	0.32	10.68	0.074
104. CR (1) C 16	/623. BD* (1) C 17 - C 18	0.29	10.69	0.071
105. CR (1) C 17	/295. RY* (3) C 15	0.48	11.54	0.094
105. CR (1) C 17	/296. RY* (4) C 15	0.39	11.04	0.083
105. CR (1) C 17	/319. RY* (3) C 18	1.06	11.33	0.139
105. CR (1) C 17	/356. RY* (2) H 28	0.30	12.31	0.076
105. CR (1) C 17	/615. BD* (1) C 13 - C 15	0.40	10.63	0.083
105. CR (1) C 17	/621. BD* (1) C 16 - C 18	0.51	10.50	0.092
105. CR (1) C 17	/623. BD* (1) C 17 - C 18	0.45	10.68	0.088
106. CR (1) C 18	/303. RY* (3) C 16	0.42	11.41	0.087
106. CR (1) C 18	/310. RY* (2) C 17	0.36	11.36	0.081
106. CR (1) C 18	/311. RY* (3) C 17	0.72	11.29	0.114
106. CR (1) C 18	/358. RY* (2) H 29	0.32	11.98	0.078
106. CR (1) C 18	/617. BD* (1) C 14 - C 16	0.27	10.68	0.068
106. CR (1) C 18	/619. BD* (1) C 15 - C 17	0.49	10.48	0.091
106. CR (1) C 18	/623. BD* (1) C 17 - C 18	0.33	10.69	0.075
107. CR (1) O 19	/293. RY* (1) C 15	1.37	20.10	0.210
107. CR (1) O 19	/615. BD* (1) C 13 - C 15	0.29	19.53	0.096
108. CR (1) O 20	/181. RY* (1) C 1	0.49	19.74	0.125
108. CR (1) O 20	/182. RY* (2) C 1	0.59	19.89	0.137
108. CR (1) O 20	/360. RY* (2) H 30	0.29	20.83	0.099
144. LP (1) N 11	/269. RY* (1) N 12	0.99	1.10	0.042
144. LP (1) N 11	/270. RY* (2) N 12	0.65	1.12	0.035
144. LP (1) N 11	/598. BD* (1) C 4 - C 6	4.42	0.75	0.073
144. LP (1) N 11	/600. BD* (1) C 5 - C 6	2.66	0.94	0.064
145. LP (2) N 11	/ 13. BD* (2) C 4 - C 6	2.93	0.42	0.057
145. LP (2) N 11	/221. RY* (1) C 6	0.64	1.32	0.040
145. LP (2) N 11	/262. RY* (2) N 11	0.98	2.40	0.066
145. LP (2) N 11	/263. RY* (3) N 11	0.46	1.81	0.039
145. LP (2) N 11	/270. RY* (2) N 12	0.37	1.28	0.030
145. LP (2) N 11	/598. BD* (1) C 4 - C 6	0.33	0.92	0.023
145. LP (2) N 11	/600. BD* (1) C 5 - C 6	0.36	1.10	0.027
146. LP (1) N 12	/262. RY* (2) N 11	0.34	2.38	0.036
146. LP (1) N 12	/263. RY* (3) N 11	0.34	1.79	0.031
146. LP (1) N 12	/277. RY* (1) C 13	0.82	1.57	0.046
146. LP (1) N 12	/614. BD* (1) C 13 - C 14	0.56	0.89	0.028
146. LP (1) N 12	/615. BD* (1) C 13 - C 15	3.51	1.05	0.077
147. LP (2) N 12	/261. RY* (1) N 11	1.34	1.17	0.052
147. LP (2) N 12	/280. RY* (4) C 13	0.51	0.91	0.028
147. LP (2) N 12	/616. BD* (2) C 13 - C 15	16.46	0.31	0.097
148. LP (1) C 14	/280. RY* (4) C 13	0.29	0.78	0.021
148. LP (1) C 14	/301. RY* (1) C 16	2.90	0.66	0.059
148. LP (1) C 14	/616. BD* (2) C 13 - C 15	29.13	0.19	0.097
149. LP (1) C 18	/301. RY* (1) C 16	1.02	0.60	0.043
149. LP (1) C 18	/309. RY* (1) C 17	2.02	0.62	0.062
149. LP (1) C 18	/318. RY* (2) C 18	1.34	0.62	0.051
150. LP (1) O 19	/293. RY* (1) C 15	1.10	1.79	0.056
150. LP (1) O 19	/327. RY* (3) O 19	0.29	2.47	0.034
150. LP (1) O 19	/615. BD* (1) C 13 - C 15	2.87	1.21	0.075
151. LP (2) O 19	/615. BD* (1) C 13 - C 15	0.50	0.94	0.027
151. LP (2) O 19	/616. BD* (2) C 13 - C 15	5.74	0.35	0.063
151. LP (2) O 19	/619. BD* (1) C 15 - C 17	1.41	0.78	0.042
152. LP (3) O 19	/262. RY* (2) N 11	0.30	2.37	0.036

152. LP (3) O 19	/293. RY* (1) C 15	0.33	1.61	0.031
152. LP (3) O 19	/326. RY* (2) O 19	0.30	2.09	0.034
152. LP (3) O 19	/327. RY* (3) O 19	0.35	2.29	0.039
152. LP (3) O 19	/615. BD* (1) C 13 - C 15	0.28	1.03	0.023
152. LP (3) O 19	/616. BD* (2) C 13 - C 15	2.78	0.44	0.046
152. LP (3) O 19	/619. BD* (1) C 15 - C 17	2.48	0.87	0.063
153. LP (1) O 20	/181. RY* (1) C 1	0.67	1.37	0.038
153. LP (1) O 20	/182. RY* (2) C 1	0.80	1.52	0.044
153. LP (1) O 20	/590. BD* (1) C 1 - C 2	3.22	1.02	0.073
154. LP (2) O 20	/591. BD* (2) C 1 - C 2	6.84	0.29	0.078
591. BD* (2) C 1 - C 2	/181. RY* (1) C 1	0.43	0.82	0.033
591. BD* (2) C 1 - C 2	/182. RY* (2) C 1	0.28	0.97	0.029
591. BD* (2) C 1 - C 2	/189. RY* (1) C 2	0.93	0.63	0.043
591. BD* (2) C 1 - C 2	/197. RY* (1) C 3	1.03	0.54	0.042
591. BD* (2) C 1 - C 2	/205. RY* (1) C 4	0.58	0.64	0.034
591. BD* (2) C 1 - C 2	/333. RY* (1) O 20	0.36	0.92	0.032
591. BD* (2) C 1 - C 2	/604. BD* (2) C 7 - C 8	41.96	0.07	0.078
604. BD* (2) C 7 - C 8	/229. RY* (1) C 7	0.65	0.86	0.056
604. BD* (2) C 7 - C 8	/230. RY* (2) C 7	0.64	0.86	0.055
604. BD* (2) C 7 - C 8	/237. RY* (1) C 8	1.77	0.47	0.068
604. BD* (2) C 7 - C 8	/609. BD* (2) C 9 - C 10	88.93	0.01	0.061
609. BD* (2) C 9 - C 10	/246. RY* (2) C 9	1.62	0.57	0.076
609. BD* (2) C 9 - C 10	/253. RY* (1) C 10	1.61	0.47	0.068
616. BD* (2) C 13 - C 15	/280. RY* (4) C 13	0.93	0.60	0.063
616. BD* (2) C 13 - C 15	/285. RY* (1) C 14	0.41	0.48	0.037
616. BD* (2) C 13 - C 15	/294. RY* (2) C 15	0.58	0.61	0.050
616. BD* (2) C 13 - C 15	/309. RY* (1) C 17	0.28	0.49	0.031
from unit 1 to unit 2				
9. BD* (2) C 3 - C 5	/657. BD* (1) C 46 - O 62	0.05	0.36	0.007
15. BD (1) C 5 - C 6	/539. RY* (3) O 62	0.03	2.38	0.011
35. BD (1) C 15 - O 19	/459. RY* (1) C 44	0.05	2.00	0.012
35. BD (1) C 15 - O 19	/477. RY* (3) C 46	0.03	2.37	0.010
36. BD* (1) C 16 - C 17	/535. RY* (1) H 61	0.03	0.66	0.006
144. LP (1) N 11	/657. BD* (1) C 46 - O 62	0.10	0.61	0.010
145. LP (2) N 11	/404. RY* (2) C 37	0.12	1.56	0.019
145. LP (2) N 11	/408. RY* (6) C 37	0.05	4.03	0.020
145. LP (2) N 11	/444. RY* (2) N 42	0.13	2.47	0.024
145. LP (2) N 11	/445. RY* (3) N 42	0.07	1.79	0.016
145. LP (2) N 11	/453. RY* (3) N 43	0.03	1.55	0.009
145. LP (2) N 11	/475. RY* (1) C 46	0.08	1.65	0.016
145. LP (2) N 11	/538. RY* (2) O 62	0.14	2.32	0.025
145. LP (2) N 11	/539. RY* (3) O 62	0.26	2.08	0.032
145. LP (2) N 11	/657. BD* (1) C 46 - O 62	0.07	0.78	0.010
150. LP (1) O 19	/404. RY* (2) C 37	0.06	1.71	0.012
150. LP (1) O 19	/444. RY* (2) N 42	0.21	2.61	0.030
150. LP (1) O 19	/445. RY* (3) N 42	0.14	1.93	0.021
150. LP (1) O 19	/454. RY* (4) N 43	0.03	2.28	0.010
150. LP (1) O 19	/475. RY* (1) C 46	0.04	1.80	0.011
150. LP (1) O 19	/480. RY* (6) C 46	0.04	3.29	0.015
150. LP (1) O 19	/538. RY* (2) O 62	0.04	2.47	0.013
150. LP (1) O 19	/539. RY* (3) O 62	0.07	2.23	0.016
150. LP (1) O 19	/639. BD* (1) C 37 - N 42	0.04	0.97	0.008
150. LP (1) O 19	/650. BD* (1) N 43 - C 44	0.05	0.96	0.009
150. LP (1) O 19	/657. BD* (1) C 46 - O 62	0.05	0.92	0.009
151. LP (2) O 19	/491. RY* (1) C 48	0.03	0.84	0.006
151. LP (2) O 19	/535. RY* (1) H 61	0.03	0.82	0.006
151. LP (2) O 19	/653. BD* (2) C 44 - C 46	0.13	0.34	0.009
151. LP (2) O 19	/657. BD* (1) C 46 - O 62	0.16	0.65	0.013
152. LP (3) O 19	/404. RY* (2) C 37	0.11	1.52	0.017
152. LP (3) O 19	/444. RY* (2) N 42	0.36	2.43	0.040
152. LP (3) O 19	/445. RY* (3) N 42	0.11	1.75	0.019
152. LP (3) O 19	/475. RY* (1) C 46	0.08	1.61	0.015
152. LP (3) O 19	/476. RY* (2) C 46	0.03	1.05	0.008
152. LP (3) O 19	/477. RY* (3) C 46	0.04	1.96	0.012
152. LP (3) O 19	/480. RY* (6) C 46	0.04	3.11	0.016
152. LP (3) O 19	/538. RY* (2) O 62	0.08	2.29	0.018
152. LP (3) O 19	/539. RY* (3) O 62	0.07	2.05	0.016
152. LP (3) O 19	/639. BD* (1) C 37 - N 42	0.09	0.79	0.011
152. LP (3) O 19	/657. BD* (1) C 46 - O 62	0.05	0.74	0.008

from unit			1 to unit			3				
11.	BD	(1) C	4 - C	6	/168.	LP (6)Cr	63	0.04	1.29	0.010
11.	BD	(1) C	4 - C	6	/545.	RY* (1)Cr	63	0.03	1.16	0.007
12.	BD	(2) C	4 - C	6	/166.	LP (4)Cr	63	0.10	0.13	0.006
12.	BD	(2) C	4 - C	6	/167.	LP (5)Cr	63	0.06	0.15	0.005
12.	BD	(2) C	4 - C	6	/171.	LP* (9)Cr	63	0.05	2.09	0.013
13.	BD*	(2) C	4 - C	6	/551.	RY* (7)Cr	63	0.03	1.95	0.012
15.	BD	(1) C	5 - C	6	/171.	LP* (9)Cr	63	0.03	2.57	0.012
15.	BD	(1) C	5 - C	6	/556.	RY* (12)Cr	63	0.03	8.80	0.020
17.	BD	(1) C	6 - N	11	/166.	LP (4)Cr	63	0.17	0.65	0.017
17.	BD	(1) C	6 - N	11	/167.	LP (5)Cr	63	0.06	0.66	0.009
17.	BD	(1) C	6 - N	11	/168.	LP (6)Cr	63	0.85	1.43	0.047
17.	BD	(1) C	6 - N	11	/170.	LP* (8)Cr	63	0.10	2.60	0.020
17.	BD	(1) C	6 - N	11	/546.	RY* (2)Cr	63	0.04	2.37	0.013
17.	BD	(1) C	6 - N	11	/548.	RY* (4)Cr	63	0.05	2.33	0.014
17.	BD	(1) C	6 - N	11	/549.	RY* (5)Cr	63	0.04	2.18	0.012
17.	BD	(1) C	6 - N	11	/550.	RY* (6)Cr	63	0.13	3.59	0.027
17.	BD	(1) C	6 - N	11	/552.	RY* (8)Cr	63	0.04	9.48	0.024
17.	BD	(1) C	6 - N	11	/555.	RY* (11)Cr	63	0.04	5.14	0.018
27.	BD	(1) N	11 - N	12	/166.	LP (4)Cr	63	0.10	0.85	0.015
27.	BD	(1) N	11 - N	12	/167.	LP (5)Cr	63	0.04	0.87	0.009
27.	BD	(1) N	11 - N	12	/168.	LP (6)Cr	63	0.85	1.63	0.050
27.	BD	(1) N	11 - N	12	/169.	LP (7)Cr	63	0.06	2.60	0.016
27.	BD	(1) N	11 - N	12	/170.	LP* (8)Cr	63	0.06	2.80	0.016
27.	BD	(1) N	11 - N	12	/171.	LP* (9)Cr	63	0.10	2.80	0.021
27.	BD	(1) N	11 - N	12	/545.	RY* (1)Cr	63	0.06	1.50	0.012
27.	BD	(1) N	11 - N	12	/546.	RY* (2)Cr	63	0.26	2.57	0.033
27.	BD	(1) N	11 - N	12	/547.	RY* (3)Cr	63	0.07	2.69	0.017
27.	BD	(1) N	11 - N	12	/548.	RY* (4)Cr	63	0.07	2.54	0.017
27.	BD	(1) N	11 - N	12	/550.	RY* (6)Cr	63	0.05	3.79	0.017
27.	BD	(1) N	11 - N	12	/554.	RY* (10)Cr	63	0.05	10.07	0.029
27.	BD	(1) N	11 - N	12	/561.	RY* (17)Cr	63	0.06	38.53	0.060
27.	BD	(1) N	11 - N	12	/564.	RY* (20)Cr	63	0.03	675.27	0.166
28.	BD	(1) N	12 - C	13	/168.	LP (6)Cr	63	0.10	1.44	0.016
28.	BD	(1) N	12 - C	13	/171.	LP* (9)Cr	63	0.05	2.61	0.014
28.	BD	(1) N	12 - C	13	/545.	RY* (1)Cr	63	0.03	1.31	0.008
28.	BD	(1) N	12 - C	13	/546.	RY* (2)Cr	63	0.03	2.37	0.011
28.	BD	(1) N	12 - C	13	/550.	RY* (6)Cr	63	0.07	3.60	0.020
28.	BD	(1) N	12 - C	13	/556.	RY* (12)Cr	63	0.08	8.84	0.033
28.	BD	(1) N	12 - C	13	/561.	RY* (17)Cr	63	0.12	38.34	0.087
28.	BD	(1) N	12 - C	13	/564.	RY* (20)Cr	63	0.04	675.07	0.211
28.	BD	(1) N	12 - C	13	/565.	RY* (21)Cr	63	0.04	27.02	0.039
30.	BD	(1) C	13 - C	15	/168.	LP (6)Cr	63	0.09	1.41	0.015
30.	BD	(1) C	13 - C	15	/545.	RY* (1)Cr	63	0.05	1.28	0.010
30.	BD	(1) C	13 - C	15	/549.	RY* (5)Cr	63	0.03	2.16	0.011
31.	BD	(2) C	13 - C	15	/167.	LP (5)Cr	63	0.06	0.18	0.005
34.	BD	(1) C	15 - C	17	/168.	LP (6)Cr	63	0.15	1.29	0.019
34.	BD	(1) C	15 - C	17	/170.	LP* (8)Cr	63	0.03	2.46	0.011
34.	BD	(1) C	15 - C	17	/545.	RY* (1)Cr	63	0.03	1.16	0.008
34.	BD	(1) C	15 - C	17	/550.	RY* (6)Cr	63	0.03	3.46	0.012
35.	BD	(1) C	15 - O	19	/167.	LP (5)Cr	63	0.08	0.72	0.012
35.	BD	(1) C	15 - O	19	/168.	LP (6)Cr	63	0.73	1.48	0.045
35.	BD	(1) C	15 - O	19	/169.	LP (7)Cr	63	0.05	2.46	0.013
35.	BD	(1) C	15 - O	19	/170.	LP* (8)Cr	63	0.18	2.66	0.028
35.	BD	(1) C	15 - O	19	/171.	LP* (9)Cr	63	0.03	2.65	0.011
35.	BD	(1) C	15 - O	19	/547.	RY* (3)Cr	63	0.04	2.54	0.013
35.	BD	(1) C	15 - O	19	/548.	RY* (4)Cr	63	0.04	2.39	0.012
35.	BD	(1) C	15 - O	19	/549.	RY* (5)Cr	63	0.05	2.23	0.014
35.	BD	(1) C	15 - O	19	/552.	RY* (8)Cr	63	0.04	9.54	0.024
35.	BD	(1) C	15 - O	19	/553.	RY* (9)Cr	63	0.03	5.27	0.017
35.	BD	(1) C	15 - O	19	/554.	RY* (10)Cr	63	0.03	9.93	0.023
99.	CR	(1) N	11		/166.	LP (4)Cr	63	0.19	14.07	0.084
99.	CR	(1) N	11		/167.	LP (5)Cr	63	0.06	14.09	0.045
99.	CR	(1) N	11		/168.	LP (6)Cr	63	1.13	14.86	0.176
100.	CR	(1) N	12		/168.	LP (6)Cr	63	0.03	14.89	0.028
103.	CR	(1) C	15		/168.	LP (6)Cr	63	0.04	10.75	0.029
107.	CR	(1) O	19		/167.	LP (5)Cr	63	0.17	18.80	0.088
107.	CR	(1) O	19		/168.	LP (6)Cr	63	1.60	19.56	0.241
144.	LP	(1) N	11		/166.	LP (4)Cr	63	0.50	0.16	0.014
144.	LP	(1) N	11		/167.	LP (5)Cr	63	0.08	0.18	0.006
144.	LP	(1) N	11		/168.	LP (6)Cr	63	0.37	0.94	0.025

144. LP (1) N 11	/169. LP (7)Cr 63	0.04	1.92	0.012
144. LP (1) N 11	/170. LP* (8)Cr 63	0.04	2.12	0.011
144. LP (1) N 11	/550. RY* (6)Cr 63	0.07	3.11	0.019
144. LP (1) N 11	/561. RY* (17)Cr 63	0.07	37.84	0.068
145. LP (2) N 11	/166. LP (4)Cr 63	22.12	0.32	0.125
145. LP (2) N 11	/167. LP (5)Cr 63	5.56	0.34	0.061
145. LP (2) N 11	/168. LP (6)Cr 63	18.49	1.11	0.181
145. LP (2) N 11	/169. LP (7)Cr 63	0.08	2.08	0.018
145. LP (2) N 11	/170. LP* (8)Cr 63	0.60	2.28	0.050
145. LP (2) N 11	/171. LP* (9)Cr 63	1.31	2.27	0.074
145. LP (2) N 11	/545. RY* (1)Cr 63	0.41	0.98	0.027
145. LP (2) N 11	/546. RY* (2)Cr 63	0.43	2.04	0.040
145. LP (2) N 11	/547. RY* (3)Cr 63	0.08	2.17	0.017
145. LP (2) N 11	/549. RY* (5)Cr 63	0.15	1.85	0.023
145. LP (2) N 11	/550. RY* (6)Cr 63	2.38	3.27	0.120
145. LP (2) N 11	/551. RY* (7)Cr 63	0.07	2.37	0.017
145. LP (2) N 11	/554. RY* (10)Cr 63	0.04	9.55	0.027
145. LP (2) N 11	/555. RY* (11)Cr 63	0.06	4.82	0.022
145. LP (2) N 11	/556. RY* (12)Cr 63	0.15	8.50	0.048
145. LP (2) N 11	/558. RY* (14)Cr 63	0.03	5.04	0.016
145. LP (2) N 11	/560. RY* (16)Cr 63	0.30	115.36	0.254
145. LP (2) N 11	/561. RY* (17)Cr 63	1.75	38.01	0.351
145. LP (2) N 11	/562. RY* (18)Cr 63	0.05	4.97	0.022
145. LP (2) N 11	/564. RY* (20)Cr 63	0.56	674.74	0.833
145. LP (2) N 11	/565. RY* (21)Cr 63	0.07	26.68	0.060
146. LP (1) N 12	/166. LP (4)Cr 63	0.07	0.30	0.008
146. LP (1) N 12	/168. LP (6)Cr 63	0.11	1.08	0.014
146. LP (1) N 12	/171. LP* (9)Cr 63	0.12	2.25	0.021
146. LP (1) N 12	/551. RY* (7)Cr 63	0.05	2.34	0.014
150. LP (1) O 19	/167. LP (5)Cr 63	1.24	0.49	0.037
150. LP (1) O 19	/168. LP (6)Cr 63	7.16	1.25	0.128
150. LP (1) O 19	/169. LP (7)Cr 63	0.24	2.23	0.029
150. LP (1) O 19	/170. LP* (8)Cr 63	0.12	2.42	0.022
150. LP (1) O 19	/171. LP* (9)Cr 63	0.19	2.42	0.027
150. LP (1) O 19	/545. RY* (1)Cr 63	0.05	1.12	0.009
150. LP (1) O 19	/546. RY* (2)Cr 63	0.30	2.19	0.032
150. LP (1) O 19	/548. RY* (4)Cr 63	0.48	2.16	0.041
150. LP (1) O 19	/549. RY* (5)Cr 63	0.16	2.00	0.022
150. LP (1) O 19	/550. RY* (6)Cr 63	0.54	3.42	0.055
150. LP (1) O 19	/551. RY* (7)Cr 63	0.20	2.51	0.029
150. LP (1) O 19	/558. RY* (14)Cr 63	0.03	5.18	0.016
150. LP (1) O 19	/560. RY* (16)Cr 63	0.10	115.51	0.136
150. LP (1) O 19	/561. RY* (17)Cr 63	0.65	38.15	0.200
150. LP (1) O 19	/564. RY* (20)Cr 63	0.23	674.89	0.505
151. LP (2) O 19	/167. LP (5)Cr 63	0.78	0.21	0.019
151. LP (2) O 19	/168. LP (6)Cr 63	1.62	0.98	0.053
151. LP (2) O 19	/169. LP (7)Cr 63	0.17	1.95	0.023
151. LP (2) O 19	/170. LP* (8)Cr 63	0.16	2.15	0.024
151. LP (2) O 19	/546. RY* (2)Cr 63	0.03	1.91	0.009
151. LP (2) O 19	/549. RY* (5)Cr 63	0.12	1.72	0.018
151. LP (2) O 19	/550. RY* (6)Cr 63	0.13	3.14	0.026
151. LP (2) O 19	/551. RY* (7)Cr 63	0.05	2.24	0.014
151. LP (2) O 19	/561. RY* (17)Cr 63	0.16	37.88	0.098
151. LP (2) O 19	/564. RY* (20)Cr 63	0.05	674.61	0.241
152. LP (3) O 19	/166. LP (4)Cr 63	0.55	0.29	0.019
152. LP (3) O 19	/167. LP (5)Cr 63	20.05	0.30	0.109
152. LP (3) O 19	/168. LP (6)Cr 63	20.30	1.07	0.186
152. LP (3) O 19	/169. LP (7)Cr 63	0.16	2.04	0.024
152. LP (3) O 19	/170. LP* (8)Cr 63	1.34	2.24	0.074
152. LP (3) O 19	/171. LP* (9)Cr 63	0.45	2.24	0.043
152. LP (3) O 19	/545. RY* (1)Cr 63	0.31	0.94	0.023
152. LP (3) O 19	/546. RY* (2)Cr 63	0.34	2.01	0.035
152. LP (3) O 19	/548. RY* (4)Cr 63	0.34	1.98	0.035
152. LP (3) O 19	/549. RY* (5)Cr 63	0.16	1.82	0.023
152. LP (3) O 19	/550. RY* (6)Cr 63	0.87	3.23	0.072
152. LP (3) O 19	/551. RY* (7)Cr 63	0.34	2.33	0.038
152. LP (3) O 19	/554. RY* (10)Cr 63	0.04	9.51	0.028
152. LP (3) O 19	/558. RY* (14)Cr 63	0.03	5.00	0.017
152. LP (3) O 19	/560. RY* (16)Cr 63	0.20	115.32	0.204
152. LP (3) O 19	/561. RY* (17)Cr 63	1.19	37.97	0.289
152. LP (3) O 19	/563. RY* (19)Cr 63	0.07	26.08	0.059

152. LP (3) O 19	/564. RY* (20)Cr 63	0.36	674.71	0.672
from unit 1 to unit 4				
11. BD (1) C 4 - C 6	/571. RY* (2) O 66	0.04	1.68	0.010
12. BD (2) C 4 - C 6	/571. RY* (2) O 66	0.03	1.31	0.008
12. BD (2) C 4 - C 6	/665. BD* (1) H 65 - O 66	0.04	0.60	0.006
13. BD* (2) C 4 - C 6	/665. BD* (1) H 65 - O 66	0.05	0.37	0.007
28. BD (1) N 12 - C 13	/570. RY* (1) O 66	0.04	2.92	0.014
144. LP (1) N 11	/570. RY* (1) O 66	0.03	2.43	0.010
145. LP (2) N 11	/566. RY* (1) H 64	0.12	1.09	0.015
145. LP (2) N 11	/568. RY* (1) H 65	0.13	1.20	0.017
145. LP (2) N 11	/570. RY* (1) O 66	0.60	2.59	0.054
150. LP (1) O 19	/566. RY* (1) H 64	0.05	1.23	0.010
150. LP (1) O 19	/568. RY* (1) H 65	0.05	1.34	0.011
150. LP (1) O 19	/570. RY* (1) O 66	0.18	2.74	0.029
151. LP (2) O 19	/568. RY* (1) H 65	0.03	1.07	0.007
151. LP (2) O 19	/570. RY* (1) O 66	0.05	2.46	0.014
152. LP (3) O 19	/566. RY* (1) H 64	0.05	1.05	0.009
152. LP (3) O 19	/568. RY* (1) H 65	0.05	1.16	0.010
152. LP (3) O 19	/570. RY* (1) O 66	0.23	2.55	0.033
152. LP (3) O 19	/572. RY* (3) O 66	0.07	1.73	0.015
152. LP (3) O 19	/664. BD* (1) H 64 - O 66	0.03	0.75	0.006
152. LP (3) O 19	/665. BD* (1) H 65 - O 66	0.03	0.75	0.007
from unit 1 to unit 5				
30. BD (1) C 13 - C 15	/578. RY* (1)C1 67	0.04	1.33	0.009
31. BD (2) C 13 - C 15	/578. RY* (1)C1 67	0.04	0.86	0.007
144. LP (1) N 11	/580. RY* (3)C1 67	0.03	1.19	0.007
144. LP (1) N 11	/581. RY* (4)C1 67	0.03	1.93	0.009
145. LP (2) N 11	/578. RY* (1)C1 67	0.03	1.02	0.008
145. LP (2) N 11	/579. RY* (2)C1 67	0.43	1.12	0.030
145. LP (2) N 11	/580. RY* (3)C1 67	0.09	1.36	0.015
145. LP (2) N 11	/581. RY* (4)C1 67	0.32	2.09	0.035
150. LP (1) O 19	/579. RY* (2)C1 67	0.13	1.26	0.017
150. LP (1) O 19	/580. RY* (3)C1 67	0.06	1.50	0.012
150. LP (1) O 19	/581. RY* (4)C1 67	0.16	2.24	0.024
151. LP (2) O 19	/579. RY* (2)C1 67	0.04	0.99	0.008
151. LP (2) O 19	/580. RY* (3)C1 67	0.03	1.23	0.008
151. LP (2) O 19	/581. RY* (4)C1 67	0.07	1.96	0.015
152. LP (3) O 19	/579. RY* (2)C1 67	0.26	1.08	0.023
152. LP (3) O 19	/580. RY* (3)C1 67	0.08	1.32	0.014
152. LP (3) O 19	/581. RY* (4)C1 67	0.15	2.05	0.024
from unit 2 to unit 1				
71. BD (1) N 43 - C 44	/327. RY* (3) O 19	0.03	2.66	0.012
73. BD (1) C 44 - C 46	/327. RY* (3) O 19	0.08	2.64	0.018
74. BD (2) C 44 - C 46	/327. RY* (3) O 19	0.05	2.17	0.013
74. BD (2) C 44 - C 46	/620. BD* (1) C 15 - O 19	0.11	0.62	0.011
155. LP (1) N 42	/620. BD* (1) C 15 - O 19	0.05	0.63	0.007
156. LP (2) N 42	/222. RY* (2) C 6	0.08	1.68	0.016
156. LP (2) N 42	/224. RY* (4) C 6	0.07	1.51	0.014
156. LP (2) N 42	/226. RY* (6) C 6	0.05	3.47	0.017
156. LP (2) N 42	/262. RY* (2) N 11	0.14	2.41	0.025
156. LP (2) N 42	/263. RY* (3) N 11	0.04	1.81	0.011
156. LP (2) N 42	/293. RY* (1) C 15	0.09	1.65	0.016
156. LP (2) N 42	/295. RY* (3) C 15	0.03	1.99	0.010
156. LP (2) N 42	/297. RY* (5) C 15	0.03	4.28	0.016
156. LP (2) N 42	/326. RY* (2) O 19	0.14	2.14	0.023
156. LP (2) N 42	/327. RY* (3) O 19	0.31	2.33	0.037
156. LP (2) N 42	/620. BD* (1) C 15 - O 19	0.12	0.79	0.013
163. LP (1) O 62	/ 9. BD* (2) C 3 - C 5	0.09	0.56	0.013
163. LP (1) O 62	/262. RY* (2) N 11	0.19	2.55	0.028
163. LP (1) O 62	/263. RY* (3) N 11	0.13	1.95	0.020
163. LP (1) O 62	/293. RY* (1) C 15	0.04	1.79	0.011
163. LP (1) O 62	/326. RY* (2) O 19	0.05	2.28	0.013
163. LP (1) O 62	/327. RY* (3) O 19	0.08	2.47	0.018
163. LP (1) O 62	/596. BD* (1) C 3 - C 5	0.05	1.07	0.009
163. LP (1) O 62	/612. BD* (1) N 11 - N 12	0.04	1.09	0.009
164. LP (2) O 62	/327. RY* (3) O 19	0.03	2.22	0.010
164. LP (2) O 62	/601. BD* (1) C 5 - H 31	0.04	0.81	0.008
165. LP (3) O 62	/222. RY* (2) C 6	0.07	1.63	0.015

165. LP (3) O 62	/261. RY* (1) N 11	0.04	1.30	0.010
165. LP (3) O 62	/262. RY* (2) N 11	0.28	2.36	0.035
165. LP (3) O 62	/263. RY* (3) N 11	0.20	1.77	0.026
165. LP (3) O 62	/271. RY* (3) N 12	0.03	1.54	0.009
165. LP (3) O 62	/293. RY* (1) C 15	0.06	1.60	0.014
165. LP (3) O 62	/326. RY* (2) O 19	0.10	2.09	0.019
165. LP (3) O 62	/327. RY* (3) O 19	0.16	2.28	0.026
165. LP (3) O 62	/602. BD* (1) C 6 - N 11	0.04	0.79	0.008
165. LP (3) O 62	/612. BD* (1) N 11 - N 12	0.08	0.90	0.011
165. LP (3) O 62	/620. BD* (1) C 15 - O 19	0.06	0.74	0.009

within unit 2

44. BD (1) C 32 - C 33	/380. RY* (2) C 34	0.54	2.00	0.042
44. BD (1) C 32 - C 33	/381. RY* (3) C 34	0.28	2.00	0.030
44. BD (1) C 32 - C 33	/388. RY* (2) C 35	0.70	2.21	0.050
44. BD (1) C 32 - C 33	/411. RY* (1) C 38	0.59	1.83	0.042
44. BD (1) C 32 - C 33	/629. BD* (1) C 32 - C 34	1.45	1.26	0.054
44. BD (1) C 32 - C 33	/631. BD* (1) C 33 - C 35	1.25	1.27	0.050
44. BD (1) C 32 - C 33	/632. BD* (1) C 33 - C 38	0.60	1.10	0.032
44. BD (1) C 32 - C 33	/634. BD* (1) C 34 - H 51	1.33	1.09	0.048
44. BD (1) C 32 - C 33	/636. BD* (1) C 35 - C 40	2.12	1.11	0.061
44. BD (1) C 32 - C 33	/640. BD* (1) C 38 - C 39	0.69	1.30	0.038
45. BD (2) C 32 - C 33	/ 52. BD* (2) C 34 - C 36	5.84	0.22	0.066
45. BD (2) C 32 - C 33	/ 56. BD* (2) C 35 - C 37	7.32	0.23	0.071
45. BD (2) C 32 - C 33	/379. RY* (1) C 34	1.06	0.77	0.037
45. BD (2) C 32 - C 33	/387. RY* (1) C 35	0.88	0.87	0.036
45. BD (2) C 32 - C 33	/412. RY* (2) C 38	0.62	0.83	0.029
45. BD (2) C 32 - C 33	/507. RY* (1) O 50	0.34	1.15	0.026
45. BD (2) C 32 - C 33	/628. BD* (2) C 32 - C 33	0.40	0.23	0.016
45. BD (2) C 32 - C 33	/641. BD* (2) C 38 - C 39	7.07	0.31	0.065
45. BD (2) C 32 - C 33	/646. BD* (2) C 40 - C 41	0.28	0.30	0.013
46. BD (1) C 32 - C 34	/373. RY* (3) C 33	0.64	2.23	0.048
46. BD (1) C 32 - C 34	/397. RY* (3) C 36	0.58	1.94	0.043
46. BD (1) C 32 - C 34	/627. BD* (1) C 32 - C 33	1.44	1.19	0.053
46. BD (1) C 32 - C 34	/630. BD* (1) C 32 - O 50	0.28	1.04	0.021
46. BD (1) C 32 - C 34	/632. BD* (1) C 33 - C 38	1.12	1.22	0.047
46. BD (1) C 32 - C 34	/633. BD* (1) C 34 - C 36	0.73	1.23	0.038
46. BD (1) C 32 - C 34	/634. BD* (1) C 34 - H 51	0.58	1.20	0.034
46. BD (1) C 32 - C 34	/638. BD* (1) C 36 - H 61	0.65	1.22	0.036
46. BD (1) C 32 - C 34	/663. BD* (1) O 50 - H 60	0.72	1.08	0.035
47. BD (1) C 32 - O 50	/380. RY* (2) C 34	0.28	2.24	0.032
47. BD (1) C 32 - O 50	/629. BD* (1) C 32 - C 34	0.38	1.51	0.030
47. BD (1) C 32 - O 50	/631. BD* (1) C 33 - C 35	0.56	1.52	0.037
47. BD (1) C 32 - O 50	/633. BD* (1) C 34 - C 36	0.98	1.36	0.046
48. BD (1) C 33 - C 35	/365. RY* (3) C 32	0.43	2.20	0.039
48. BD (1) C 33 - C 35	/404. RY* (2) C 37	0.37	1.84	0.033
48. BD (1) C 33 - C 35	/411. RY* (1) C 38	0.41	1.93	0.036
48. BD (1) C 33 - C 35	/627. BD* (1) C 32 - C 33	1.24	1.16	0.048
48. BD (1) C 33 - C 35	/630. BD* (1) C 32 - O 50	1.31	1.02	0.046
48. BD (1) C 33 - C 35	/632. BD* (1) C 33 - C 38	1.60	1.19	0.055
48. BD (1) C 33 - C 35	/635. BD* (1) C 35 - C 37	1.64	1.19	0.056
48. BD (1) C 33 - C 35	/636. BD* (1) C 35 - C 40	1.60	1.20	0.056
48. BD (1) C 33 - C 35	/639. BD* (1) C 37 - N 42	1.44	1.11	0.051
48. BD (1) C 33 - C 35	/642. BD* (1) C 38 - H 52	0.75	1.20	0.038
48. BD (1) C 33 - C 35	/647. BD* (1) C 40 - H 54	0.72	1.20	0.037
49. BD (1) C 33 - C 38	/363. RY* (1) C 32	0.33	1.48	0.028
49. BD (1) C 33 - C 38	/365. RY* (3) C 32	0.34	2.08	0.034
49. BD (1) C 33 - C 38	/388. RY* (2) C 35	0.45	2.19	0.040
49. BD (1) C 33 - C 38	/420. RY* (2) C 39	0.34	1.87	0.032
49. BD (1) C 33 - C 38	/421. RY* (3) C 39	0.72	1.99	0.048
49. BD (1) C 33 - C 38	/627. BD* (1) C 32 - C 33	0.48	1.05	0.028
49. BD (1) C 33 - C 38	/629. BD* (1) C 32 - C 34	0.74	1.24	0.038
49. BD (1) C 33 - C 38	/631. BD* (1) C 33 - C 35	1.82	1.25	0.060
49. BD (1) C 33 - C 38	/635. BD* (1) C 35 - C 37	2.50	1.08	0.066
49. BD (1) C 33 - C 38	/640. BD* (1) C 38 - C 39	1.26	1.29	0.051
49. BD (1) C 33 - C 38	/642. BD* (1) C 38 - H 52	0.29	1.09	0.022
49. BD (1) C 33 - C 38	/644. BD* (1) C 39 - H 53	1.36	1.09	0.049
50. BD (1) C 34 - C 36	/363. RY* (1) C 32	0.28	1.50	0.026
50. BD (1) C 34 - C 36	/364. RY* (2) C 32	0.48	1.52	0.034
50. BD (1) C 34 - C 36	/365. RY* (3) C 32	0.35	2.10	0.034
50. BD (1) C 34 - C 36	/405. RY* (3) C 37	0.89	1.72	0.050

50. BD (1) C 34 - C 36	/629. BD* (1) C 32 - C 34	0.81	1.26	0.041
50. BD (1) C 34 - C 36	/630. BD* (1) C 32 - O 50	2.67	0.93	0.063
50. BD (1) C 34 - C 36	/637. BD* (1) C 36 - C 37	1.03	1.28	0.046
50. BD (1) C 34 - C 36	/639. BD* (1) C 37 - N 42	3.00	1.02	0.070
51. BD (2) C 34 - C 36	/ 56. BD* (2) C 35 - C 37	5.77	0.23	0.064
51. BD (2) C 34 - C 36	/403. RY* (1) C 37	0.44	1.12	0.028
51. BD (2) C 34 - C 36	/628. BD* (2) C 32 - C 33	6.60	0.23	0.067
52. BD* (2) C 34 - C 36	/379. RY* (1) C 34	2.12	0.54	0.056
52. BD* (2) C 34 - C 36	/395. RY* (1) C 36	1.49	0.59	0.049
53. BD (1) C 34 - H 51	/365. RY* (3) C 32	0.48	1.97	0.039
53. BD (1) C 34 - H 51	/368. RY* (6) C 32	0.30	8.49	0.064
53. BD (1) C 34 - H 51	/396. RY* (2) C 36	0.51	1.78	0.038
53. BD (1) C 34 - H 51	/627. BD* (1) C 32 - C 33	3.08	0.94	0.068
53. BD (1) C 34 - H 51	/629. BD* (1) C 32 - C 34	0.40	1.13	0.027
53. BD (1) C 34 - H 51	/630. BD* (1) C 32 - O 50	0.94	0.79	0.034
53. BD (1) C 34 - H 51	/637. BD* (1) C 36 - C 37	1.45	1.15	0.052
53. BD (1) C 34 - H 51	/638. BD* (1) C 36 - H 61	0.30	0.97	0.021
54. BD (1) C 35 - C 37	/372. RY* (2) C 33	0.83	2.34	0.056
54. BD (1) C 35 - C 37	/396. RY* (2) C 36	0.49	1.92	0.039
54. BD (1) C 35 - C 37	/427. RY* (1) C 40	0.34	1.49	0.029
54. BD (1) C 35 - C 37	/428. RY* (2) C 40	0.47	1.62	0.035
54. BD (1) C 35 - C 37	/443. RY* (1) N 42	0.59	1.52	0.038
54. BD (1) C 35 - C 37	/631. BD* (1) C 33 - C 35	1.55	1.27	0.056
54. BD (1) C 35 - C 37	/632. BD* (1) C 33 - C 38	1.92	1.10	0.058
54. BD (1) C 35 - C 37	/636. BD* (1) C 35 - C 40	0.81	1.11	0.038
54. BD (1) C 35 - C 37	/637. BD* (1) C 36 - C 37	1.63	1.28	0.058
54. BD (1) C 35 - C 37	/638. BD* (1) C 36 - H 61	1.41	1.10	0.050
54. BD (1) C 35 - C 37	/639. BD* (1) C 37 - N 42	0.29	1.02	0.022
54. BD (1) C 35 - C 37	/645. BD* (1) C 40 - C 41	0.72	1.31	0.039
54. BD (1) C 35 - C 37	/649. BD* (1) N 42 - N 43	0.29	1.13	0.023
55. BD (2) C 35 - C 37	/ 52. BD* (2) C 34 - C 36	6.29	0.23	0.069
55. BD (2) C 35 - C 37	/371. RY* (1) C 33	0.79	0.87	0.034
55. BD (2) C 35 - C 37	/395. RY* (1) C 36	1.56	0.81	0.046
55. BD (2) C 35 - C 37	/628. BD* (2) C 32 - C 33	4.54	0.23	0.055
55. BD (2) C 35 - C 37	/646. BD* (2) C 40 - C 41	5.88	0.30	0.061
55. BD (2) C 35 - C 37	/649. BD* (1) N 42 - N 43	2.60	0.76	0.057
56. BD* (2) C 35 - C 37	/371. RY* (1) C 33	0.91	0.63	0.042
56. BD* (2) C 35 - C 37	/387. RY* (1) C 35	1.49	0.64	0.054
56. BD* (2) C 35 - C 37	/395. RY* (1) C 36	1.14	0.58	0.045
56. BD* (2) C 35 - C 37	/403. RY* (1) C 37	0.51	0.89	0.037
56. BD* (2) C 35 - C 37	/635. BD* (1) C 35 - C 37	0.38	0.49	0.023
56. BD* (2) C 35 - C 37	/646. BD* (2) C 40 - C 41	53.76	0.07	0.085
56. BD* (2) C 35 - C 37	/649. BD* (1) N 42 - N 43	2.12	0.52	0.057
57. BD (1) C 35 - C 40	/372. RY* (2) C 33	0.48	2.32	0.042
57. BD (1) C 35 - C 40	/403. RY* (1) C 37	0.50	1.48	0.034
57. BD (1) C 35 - C 40	/404. RY* (2) C 37	0.68	1.73	0.044
57. BD (1) C 35 - C 40	/436. RY* (2) C 41	0.38	1.96	0.035
57. BD (1) C 35 - C 40	/437. RY* (3) C 41	0.74	2.01	0.049
57. BD (1) C 35 - C 40	/627. BD* (1) C 32 - C 33	2.33	1.05	0.063
57. BD (1) C 35 - C 40	/631. BD* (1) C 33 - C 35	1.83	1.25	0.061
57. BD (1) C 35 - C 40	/635. BD* (1) C 35 - C 37	0.80	1.08	0.037
57. BD (1) C 35 - C 40	/637. BD* (1) C 36 - C 37	1.02	1.26	0.045
57. BD (1) C 35 - C 40	/645. BD* (1) C 40 - C 41	1.38	1.29	0.054
57. BD (1) C 35 - C 40	/647. BD* (1) C 40 - H 54	0.32	1.09	0.023
57. BD (1) C 35 - C 40	/648. BD* (1) C 41 - H 55	1.41	1.09	0.050
58. BD (1) C 36 - C 37	/381. RY* (3) C 34	0.66	2.11	0.047
58. BD (1) C 36 - C 37	/389. RY* (3) C 35	0.71	2.31	0.051
58. BD (1) C 36 - C 37	/443. RY* (1) N 42	0.70	1.63	0.043
58. BD (1) C 36 - C 37	/633. BD* (1) C 34 - C 36	0.90	1.22	0.042
58. BD (1) C 36 - C 37	/634. BD* (1) C 34 - H 51	0.58	1.20	0.033
58. BD (1) C 36 - C 37	/635. BD* (1) C 35 - C 37	1.69	1.21	0.058
58. BD (1) C 36 - C 37	/636. BD* (1) C 35 - C 40	1.15	1.22	0.047
58. BD (1) C 36 - C 37	/638. BD* (1) C 36 - H 61	0.75	1.21	0.038
58. BD (1) C 36 - C 37	/639. BD* (1) C 37 - N 42	0.69	1.13	0.035
58. BD (1) C 36 - C 37	/649. BD* (1) N 42 - N 43	0.40	1.24	0.028
59. BD (1) C 36 - H 61	/380. RY* (2) C 34	0.49	1.85	0.038
59. BD (1) C 36 - H 61	/403. RY* (1) C 37	0.38	1.35	0.029
59. BD (1) C 36 - H 61	/404. RY* (2) C 37	0.68	1.61	0.042
59. BD (1) C 36 - H 61	/406. RY* (4) C 37	0.30	1.56	0.028
59. BD (1) C 36 - H 61	/629. BD* (1) C 32 - C 34	1.35	1.12	0.049
59. BD (1) C 36 - H 61	/634. BD* (1) C 34 - H 51	0.29	0.94	0.021

59. BD (1) C 36 - H 61	/635. BD* (1) C 35 - C 37	3.38	0.95	0.072
59. BD (1) C 36 - H 61	/637. BD* (1) C 36 - C 37	0.66	1.14	0.035
59. BD (1) C 36 - H 61	/639. BD* (1) C 37 - N 42	0.72	0.87	0.032
60. BD (1) C 37 - N 42	/389. RY* (3) C 35	0.69	2.34	0.051
60. BD (1) C 37 - N 42	/396. RY* (2) C 36	0.63	2.06	0.046
60. BD (1) C 37 - N 42	/404. RY* (2) C 37	0.26	1.89	0.028
60. BD (1) C 37 - N 42	/451. RY* (1) N 43	0.72	1.62	0.043
60. BD (1) C 37 - N 42	/452. RY* (2) N 43	0.35	1.62	0.030
60. BD (1) C 37 - N 42	/631. BD* (1) C 33 - C 35	0.76	1.41	0.041
60. BD (1) C 37 - N 42	/633. BD* (1) C 34 - C 36	1.21	1.25	0.049
60. BD (1) C 37 - N 42	/635. BD* (1) C 35 - C 37	0.44	1.24	0.030
60. BD (1) C 37 - N 42	/637. BD* (1) C 36 - C 37	0.88	1.42	0.045
60. BD (1) C 37 - N 42	/649. BD* (1) N 42 - N 43	0.56	1.27	0.034
60. BD (1) C 37 - N 42	/650. BD* (1) N 43 - C 44	2.82	1.15	0.072
61. BD (1) C 38 - C 39	/372. RY* (2) C 33	0.42	2.43	0.040
61. BD (1) C 38 - C 39	/373. RY* (3) C 33	0.51	2.20	0.042
61. BD (1) C 38 - C 39	/436. RY* (2) C 41	0.43	2.06	0.038
61. BD (1) C 38 - C 39	/437. RY* (3) C 41	0.41	2.11	0.037
61. BD (1) C 38 - C 39	/627. BD* (1) C 32 - C 33	1.26	1.16	0.049
61. BD (1) C 38 - C 39	/632. BD* (1) C 33 - C 38	1.36	1.19	0.051
61. BD (1) C 38 - C 39	/642. BD* (1) C 38 - H 52	0.83	1.19	0.040
61. BD (1) C 38 - C 39	/643. BD* (1) C 39 - C 41	0.97	1.21	0.043
61. BD (1) C 38 - C 39	/644. BD* (1) C 39 - H 53	0.70	1.19	0.036
61. BD (1) C 38 - C 39	/648. BD* (1) C 41 - H 55	0.76	1.19	0.038
62. BD (2) C 38 - C 39	/371. RY* (1) C 33	0.41	0.86	0.024
62. BD (2) C 38 - C 39	/435. RY* (1) C 41	0.68	0.78	0.029
62. BD (2) C 38 - C 39	/628. BD* (2) C 32 - C 33	3.49	0.22	0.049
62. BD (2) C 38 - C 39	/646. BD* (2) C 40 - C 41	4.14	0.30	0.052
63. BD (1) C 38 - H 52	/372. RY* (2) C 33	0.63	2.20	0.047
63. BD (1) C 38 - H 52	/420. RY* (2) C 39	0.68	1.75	0.044
63. BD (1) C 38 - H 52	/631. BD* (1) C 33 - C 35	1.60	1.13	0.054
63. BD (1) C 38 - H 52	/640. BD* (1) C 38 - C 39	0.70	1.16	0.036
63. BD (1) C 38 - H 52	/643. BD* (1) C 39 - C 41	2.64	0.98	0.064
63. BD (1) C 38 - H 52	/644. BD* (1) C 39 - H 53	0.47	0.97	0.027
64. BD (1) C 39 - C 41	/413. RY* (3) C 38	0.92	1.78	0.052
64. BD (1) C 39 - C 41	/429. RY* (3) C 40	0.91	1.77	0.051
64. BD (1) C 39 - C 41	/640. BD* (1) C 38 - C 39	1.07	1.29	0.047
64. BD (1) C 39 - C 41	/642. BD* (1) C 38 - H 52	1.63	1.09	0.053
64. BD (1) C 39 - C 41	/645. BD* (1) C 40 - C 41	1.15	1.30	0.049
64. BD (1) C 39 - C 41	/647. BD* (1) C 40 - H 54	1.76	1.09	0.055
65. BD (1) C 39 - H 53	/411. RY* (1) C 38	0.76	1.69	0.045
65. BD (1) C 39 - H 53	/436. RY* (2) C 41	0.65	1.83	0.044
65. BD (1) C 39 - H 53	/632. BD* (1) C 33 - C 38	3.27	0.95	0.071
65. BD (1) C 39 - H 53	/640. BD* (1) C 38 - C 39	0.60	1.16	0.033
65. BD (1) C 39 - H 53	/642. BD* (1) C 38 - H 52	0.56	0.96	0.029
65. BD (1) C 39 - H 53	/645. BD* (1) C 40 - C 41	1.37	1.16	0.051
65. BD (1) C 39 - H 53	/648. BD* (1) C 41 - H 55	0.26	0.96	0.020
66. BD (1) C 40 - C 41	/388. RY* (2) C 35	0.62	2.30	0.048
66. BD (1) C 40 - C 41	/389. RY* (3) C 35	0.36	2.29	0.036
66. BD (1) C 40 - C 41	/420. RY* (2) C 39	0.43	1.98	0.037
66. BD (1) C 40 - C 41	/421. RY* (3) C 39	0.39	2.10	0.036
66. BD (1) C 40 - C 41	/635. BD* (1) C 35 - C 37	1.24	1.19	0.049
66. BD (1) C 40 - C 41	/636. BD* (1) C 35 - C 40	1.44	1.20	0.053
66. BD (1) C 40 - C 41	/643. BD* (1) C 39 - C 41	1.01	1.21	0.044
66. BD (1) C 40 - C 41	/644. BD* (1) C 39 - H 53	0.75	1.19	0.038
66. BD (1) C 40 - C 41	/647. BD* (1) C 40 - H 54	0.85	1.19	0.040
66. BD (1) C 40 - C 41	/648. BD* (1) C 41 - H 55	0.74	1.19	0.038
67. BD (2) C 40 - C 41	/ 56. BD* (2) C 35 - C 37	3.89	0.23	0.052
67. BD (2) C 40 - C 41	/387. RY* (1) C 35	0.36	0.87	0.023
67. BD (2) C 40 - C 41	/419. RY* (1) C 39	0.67	0.77	0.029
67. BD (2) C 40 - C 41	/641. BD* (2) C 38 - C 39	5.50	0.30	0.058
67. BD (2) C 40 - C 41	/646. BD* (2) C 40 - C 41	0.64	0.30	0.020
68. BD (1) C 40 - H 54	/388. RY* (2) C 35	0.68	2.07	0.048
68. BD (1) C 40 - H 54	/436. RY* (2) C 41	0.69	1.83	0.045
68. BD (1) C 40 - H 54	/631. BD* (1) C 33 - C 35	1.76	1.13	0.056
68. BD (1) C 40 - H 54	/643. BD* (1) C 39 - C 41	2.57	0.98	0.063
68. BD (1) C 40 - H 54	/645. BD* (1) C 40 - C 41	0.75	1.17	0.037
68. BD (1) C 40 - H 54	/648. BD* (1) C 41 - H 55	0.46	0.96	0.026
69. BD (1) C 41 - H 55	/420. RY* (2) C 39	0.51	1.74	0.038
69. BD (1) C 41 - H 55	/427. RY* (1) C 40	0.35	1.34	0.028
69. BD (1) C 41 - H 55	/428. RY* (2) C 40	0.48	1.47	0.034

69. BD (1) C 41 - H 55	/636. BD* (1) C 35 - C 40	3.21	0.96	0.070
69. BD (1) C 41 - H 55	/640. BD* (1) C 38 - C 39	1.34	1.15	0.050
69. BD (1) C 41 - H 55	/644. BD* (1) C 39 - H 53	0.26	0.96	0.020
69. BD (1) C 41 - H 55	/645. BD* (1) C 40 - C 41	0.66	1.16	0.035
69. BD (1) C 41 - H 55	/647. BD* (1) C 40 - H 54	0.51	0.95	0.028
70. BD (1) N 42 - N 43	/403. RY* (1) C 37	0.87	1.83	0.051
70. BD (1) N 42 - N 43	/459. RY* (1) C 44	0.76	2.15	0.051
70. BD (1) N 42 - N 43	/460. RY* (2) C 44	0.73	2.41	0.053
70. BD (1) N 42 - N 43	/635. BD* (1) C 35 - C 37	0.25	1.44	0.024
70. BD (1) N 42 - N 43	/639. BD* (1) C 37 - N 42	0.61	1.36	0.036
70. BD (1) N 42 - N 43	/650. BD* (1) N 43 - C 44	0.42	1.34	0.030
70. BD (1) N 42 - N 43	/651. BD* (1) C 44 - C 45	0.57	1.43	0.036
71. BD (1) N 43 - C 44	/403. RY* (1) C 37	0.30	1.64	0.028
71. BD (1) N 43 - C 44	/444. RY* (2) N 42	0.83	2.81	0.061
71. BD (1) N 43 - C 44	/445. RY* (3) N 42	0.67	2.13	0.048
71. BD (1) N 43 - C 44	/468. RY* (2) C 45	0.35	2.12	0.035
71. BD (1) N 43 - C 44	/475. RY* (1) C 46	0.41	1.99	0.036
71. BD (1) N 43 - C 44	/639. BD* (1) C 37 - N 42	2.41	1.16	0.067
71. BD (1) N 43 - C 44	/651. BD* (1) C 44 - C 45	0.30	1.24	0.025
71. BD (1) N 43 - C 44	/652. BD* (1) C 44 - C 46	0.95	1.40	0.046
71. BD (1) N 43 - C 44	/654. BD* (1) C 45 - C 47	0.38	1.44	0.030
71. BD (1) N 43 - C 44	/656. BD* (1) C 46 - C 48	1.20	1.24	0.049
72. BD (1) C 44 - C 45	/451. RY* (1) N 43	0.32	1.48	0.028
72. BD (1) C 44 - C 45	/477. RY* (3) C 46	0.30	2.19	0.033
72. BD (1) C 44 - C 45	/478. RY* (4) C 46	0.26	1.69	0.027
72. BD (1) C 44 - C 45	/484. RY* (2) C 47	0.38	1.89	0.034
72. BD (1) C 44 - C 45	/485. RY* (3) C 47	0.68	2.04	0.047
72. BD (1) C 44 - C 45	/649. BD* (1) N 42 - N 43	1.57	1.13	0.053
72. BD (1) C 44 - C 45	/650. BD* (1) N 43 - C 44	0.39	1.00	0.025
72. BD (1) C 44 - C 45	/652. BD* (1) C 44 - C 46	1.82	1.25	0.060
72. BD (1) C 44 - C 45	/654. BD* (1) C 45 - C 47	0.85	1.29	0.042
72. BD (1) C 44 - C 45	/657. BD* (1) C 46 - O 62	2.57	0.97	0.063
72. BD (1) C 44 - C 45	/659. BD* (1) C 47 - H 57	1.39	1.10	0.050
73. BD (1) C 44 - C 46	/453. RY* (3) N 43	0.32	1.86	0.031
73. BD (1) C 44 - C 46	/468. RY* (2) C 45	0.42	2.09	0.037
73. BD (1) C 44 - C 46	/492. RY* (2) C 48	0.61	2.10	0.046
73. BD (1) C 44 - C 46	/650. BD* (1) N 43 - C 44	0.68	1.13	0.035
73. BD (1) C 44 - C 46	/651. BD* (1) C 44 - C 45	1.78	1.21	0.059
73. BD (1) C 44 - C 46	/655. BD* (1) C 45 - H 56	0.55	1.21	0.033
73. BD (1) C 44 - C 46	/656. BD* (1) C 46 - C 48	1.45	1.21	0.053
73. BD (1) C 44 - C 46	/657. BD* (1) C 46 - O 62	0.34	1.09	0.024
73. BD (1) C 44 - C 46	/661. BD* (1) C 48 - H 58	0.57	1.22	0.033
74. BD (2) C 44 - C 46	/ 79. BD* (1) C 47 - C 48	0.75	0.13	0.053
74. BD (2) C 44 - C 46	/452. RY* (2) N 43	0.27	1.13	0.022
74. BD (2) C 44 - C 46	/467. RY* (1) C 45	0.70	0.80	0.030
74. BD (2) C 44 - C 46	/491. RY* (1) C 48	1.14	0.81	0.039
74. BD (2) C 44 - C 46	/537. RY* (1) O 62	0.37	1.36	0.029
74. BD (2) C 44 - C 46	/653. BD* (2) C 44 - C 46	0.57	0.32	0.019
75. BD (1) C 45 - C 47	/460. RY* (2) C 44	0.98	2.17	0.058
75. BD (1) C 45 - C 47	/499. RY* (1) C 49	0.54	1.96	0.041
75. BD (1) C 45 - C 47	/501. RY* (3) C 49	0.44	2.06	0.038
75. BD (1) C 45 - C 47	/650. BD* (1) N 43 - C 44	1.44	1.11	0.051
75. BD (1) C 45 - C 47	/651. BD* (1) C 44 - C 45	0.83	1.20	0.040
75. BD (1) C 45 - C 47	/655. BD* (1) C 45 - H 56	0.64	1.20	0.035
75. BD (1) C 45 - C 47	/658. BD* (1) C 47 - C 49	0.82	1.22	0.040
75. BD (1) C 45 - C 47	/659. BD* (1) C 47 - H 57	0.52	1.21	0.032
75. BD (1) C 45 - C 47	/662. BD* (1) C 49 - H 59	0.83	1.22	0.040
76. BD (1) C 45 - H 56	/459. RY* (1) C 44	0.41	1.67	0.033
76. BD (1) C 45 - H 56	/484. RY* (2) C 47	0.61	1.75	0.041
76. BD (1) C 45 - H 56	/650. BD* (1) N 43 - C 44	0.37	0.86	0.023
76. BD (1) C 45 - H 56	/652. BD* (1) C 44 - C 46	1.74	1.11	0.056
76. BD (1) C 45 - H 56	/654. BD* (1) C 45 - C 47	0.49	1.15	0.030
76. BD (1) C 45 - H 56	/658. BD* (1) C 47 - C 49	2.70	0.98	0.065
76. BD (1) C 45 - H 56	/659. BD* (1) C 47 - H 57	0.54	0.96	0.029
77. BD (1) C 46 - C 48	/459. RY* (1) C 44	0.56	1.82	0.041
77. BD (1) C 46 - C 48	/499. RY* (1) C 49	0.34	1.86	0.032
77. BD (1) C 46 - C 48	/501. RY* (3) C 49	0.77	1.96	0.049
77. BD (1) C 46 - C 48	/650. BD* (1) N 43 - C 44	2.78	1.01	0.067
77. BD (1) C 46 - C 48	/652. BD* (1) C 44 - C 46	1.66	1.25	0.058
77. BD (1) C 46 - C 48	/660. BD* (1) C 48 - C 49	0.94	1.30	0.044
77. BD (1) C 46 - C 48	/662. BD* (1) C 49 - H 59	1.54	1.11	0.052

78. BD (1) C 46 - O 62	/459. RY* (1) C 44	0.29	2.01	0.030
78. BD (1) C 46 - O 62	/475. RY* (1) C 46	0.48	2.04	0.039
78. BD (1) C 46 - O 62	/651. BD* (1) C 44 - C 45	1.29	1.29	0.052
78. BD (1) C 46 - O 62	/652. BD* (1) C 44 - C 46	0.66	1.45	0.039
78. BD (1) C 46 - O 62	/660. BD* (1) C 48 - C 49	0.46	1.50	0.033
79. BD* (1) C 47 - C 48	/160. LP (1) C 49	2.31	0.07	0.022
79. BD* (1) C 47 - C 48	/467. RY* (1) C 45	1.54	0.67	0.041
79. BD* (1) C 47 - C 48	/476. RY* (2) C 46	0.27	0.81	0.019
79. BD* (1) C 47 - C 48	/653. BD* (2) C 44 - C 46	5.29	0.19	0.045
80. BD (1) C 47 - C 48	/160. LP (1) C 49	655.99	0.03	0.187
80. BD (1) C 47 - C 48	/467. RY* (1) C 45	0.70	0.63	0.036
80. BD (1) C 47 - C 48	/500. RY* (2) C 49	1.55	0.66	0.055
80. BD (1) C 47 - C 48	/653. BD* (2) C 44 - C 46	10.95	0.15	0.059
81. BD (1) C 47 - C 49	/469. RY* (3) C 45	0.83	1.95	0.051
81. BD (1) C 47 - C 49	/493. RY* (3) C 48	0.82	1.93	0.050
81. BD (1) C 47 - C 49	/654. BD* (1) C 45 - C 47	0.86	1.28	0.042
81. BD (1) C 47 - C 49	/655. BD* (1) C 45 - H 56	1.58	1.08	0.052
81. BD (1) C 47 - C 49	/660. BD* (1) C 48 - C 49	1.04	1.29	0.046
81. BD (1) C 47 - C 49	/661. BD* (1) C 48 - H 58	1.54	1.09	0.052
81. BD (1) C 47 - C 49	/662. BD* (1) C 49 - H 59	0.27	1.10	0.022
82. BD (1) C 47 - H 57	/468. RY* (2) C 45	0.75	1.83	0.047
82. BD (1) C 47 - H 57	/499. RY* (1) C 49	0.64	1.71	0.042
82. BD (1) C 47 - H 57	/651. BD* (1) C 44 - C 45	3.00	0.95	0.067
82. BD (1) C 47 - H 57	/654. BD* (1) C 45 - C 47	0.41	1.15	0.028
82. BD (1) C 47 - H 57	/655. BD* (1) C 45 - H 56	0.55	0.95	0.029
82. BD (1) C 47 - H 57	/660. BD* (1) C 48 - C 49	1.45	1.16	0.052
82. BD (1) C 47 - H 57	/662. BD* (1) C 49 - H 59	0.26	0.97	0.020
83. BD (1) C 48 - C 49	/475. RY* (1) C 46	0.57	1.93	0.042
83. BD (1) C 48 - C 49	/477. RY* (3) C 46	0.47	2.28	0.042
83. BD (1) C 48 - C 49	/484. RY* (2) C 47	0.50	1.98	0.040
83. BD (1) C 48 - C 49	/485. RY* (3) C 47	0.37	2.14	0.036
83. BD (1) C 48 - C 49	/656. BD* (1) C 46 - C 48	0.76	1.19	0.038
83. BD (1) C 48 - C 49	/657. BD* (1) C 46 - O 62	1.26	1.06	0.046
83. BD (1) C 48 - C 49	/658. BD* (1) C 47 - C 49	0.98	1.21	0.043
83. BD (1) C 48 - C 49	/659. BD* (1) C 47 - H 57	0.71	1.20	0.037
83. BD (1) C 48 - C 49	/661. BD* (1) C 48 - H 58	0.76	1.19	0.038
83. BD (1) C 48 - C 49	/662. BD* (1) C 49 - H 59	0.71	1.21	0.037
84. BD (1) C 48 - H 58	/475. RY* (1) C 46	0.38	1.69	0.032
84. BD (1) C 48 - H 58	/499. RY* (1) C 49	0.66	1.71	0.043
84. BD (1) C 48 - H 58	/652. BD* (1) C 44 - C 46	1.61	1.11	0.053
84. BD (1) C 48 - H 58	/657. BD* (1) C 46 - O 62	0.52	0.82	0.026
84. BD (1) C 48 - H 58	/658. BD* (1) C 47 - C 49	2.88	0.98	0.067
84. BD (1) C 48 - H 58	/660. BD* (1) C 48 - C 49	0.63	1.16	0.034
84. BD (1) C 48 - H 58	/662. BD* (1) C 49 - H 59	0.52	0.97	0.028
85. BD (1) C 49 - H 59	/484. RY* (2) C 47	0.58	1.74	0.040
85. BD (1) C 49 - H 59	/492. RY* (2) C 48	0.75	1.83	0.047
85. BD (1) C 49 - H 59	/654. BD* (1) C 45 - C 47	1.28	1.14	0.049
85. BD (1) C 49 - H 59	/656. BD* (1) C 46 - C 48	2.85	0.95	0.066
85. BD (1) C 49 - H 59	/660. BD* (1) C 48 - C 49	0.62	1.16	0.034
85. BD (1) C 49 - H 59	/661. BD* (1) C 48 - H 58	0.48	0.95	0.027
86. BD (1) O 50 - H 60	/363. RY* (1) C 32	0.74	1.57	0.043
86. BD (1) O 50 - H 60	/364. RY* (2) C 32	0.78	1.60	0.045
86. BD (1) O 50 - H 60	/629. BD* (1) C 32 - C 34	1.60	1.34	0.058
109. CR (1) C 32	/372. RY* (2) C 33	0.54	11.81	0.101
109. CR (1) C 32	/381. RY* (3) C 34	1.06	11.47	0.139
109. CR (1) C 32	/629. BD* (1) C 32 - C 34	0.90	10.73	0.125
109. CR (1) C 32	/630. BD* (1) C 32 - O 50	0.66	10.40	0.105
109. CR (1) C 32	/631. BD* (1) C 33 - C 35	0.29	10.74	0.071
109. CR (1) C 32	/632. BD* (1) C 33 - C 38	0.31	10.57	0.072
109. CR (1) C 32	/633. BD* (1) C 34 - C 36	0.40	10.58	0.082
110. CR (1) C 33	/364. RY* (2) C 32	0.30	10.93	0.072
110. CR (1) C 33	/389. RY* (3) C 35	1.17	11.61	0.147
110. CR (1) C 33	/413. RY* (3) C 38	0.52	11.20	0.096
110. CR (1) C 33	/629. BD* (1) C 32 - C 34	0.27	10.67	0.069
110. CR (1) C 33	/631. BD* (1) C 33 - C 35	0.63	10.68	0.104
110. CR (1) C 33	/635. BD* (1) C 35 - C 37	0.57	10.51	0.099
110. CR (1) C 33	/636. BD* (1) C 35 - C 40	0.50	10.52	0.092
110. CR (1) C 33	/640. BD* (1) C 38 - C 39	0.27	10.71	0.068
111. CR (1) C 34	/365. RY* (3) C 32	1.48	11.50	0.165
111. CR (1) C 34	/398. RY* (4) C 36	0.45	11.22	0.090
111. CR (1) C 34	/516. RY* (2) H 51	0.29	12.28	0.075

111. CR (1) C 34	/627. BD* (1) C 32 - C 33	0.59	10.47	0.101
111. CR (1) C 34	/629. BD* (1) C 32 - C 34	0.29	10.66	0.070
111. CR (1) C 34	/630. BD* (1) C 32 - O 50	0.36	10.32	0.077
111. CR (1) C 34	/637. BD* (1) C 36 - C 37	0.30	10.68	0.072
112. CR (1) C 35	/373. RY* (3) C 33	1.24	11.51	0.151
112. CR (1) C 35	/405. RY* (3) C 37	0.66	11.12	0.108
112. CR (1) C 35	/429. RY* (3) C 40	0.57	11.18	0.101
112. CR (1) C 35	/627. BD* (1) C 32 - C 33	0.51	10.47	0.093
112. CR (1) C 35	/631. BD* (1) C 33 - C 35	0.50	10.67	0.093
112. CR (1) C 35	/632. BD* (1) C 33 - C 38	0.46	10.50	0.088
112. CR (1) C 35	/637. BD* (1) C 36 - C 37	0.32	10.69	0.075
112. CR (1) C 35	/639. BD* (1) C 37 - N 42	0.27	10.42	0.068
112. CR (1) C 35	/645. BD* (1) C 40 - C 41	0.28	10.71	0.070
113. CR (1) C 36	/404. RY* (2) C 37	0.36	11.13	0.080
113. CR (1) C 36	/405. RY* (3) C 37	1.07	11.10	0.137
113. CR (1) C 36	/406. RY* (4) C 37	0.36	11.09	0.080
113. CR (1) C 36	/536. RY* (2) H 61	0.29	12.37	0.076
113. CR (1) C 36	/629. BD* (1) C 32 - C 34	0.28	10.64	0.069
113. CR (1) C 36	/635. BD* (1) C 35 - C 37	0.64	10.48	0.105
113. CR (1) C 36	/637. BD* (1) C 36 - C 37	0.45	10.66	0.088
113. CR (1) C 36	/639. BD* (1) C 37 - N 42	0.44	10.40	0.086
114. CR (1) C 37	/388. RY* (2) C 35	0.61	11.65	0.106
114. CR (1) C 37	/397. RY* (3) C 36	1.16	11.26	0.144
114. CR (1) C 37	/631. BD* (1) C 33 - C 35	0.35	10.71	0.077
114. CR (1) C 37	/633. BD* (1) C 34 - C 36	0.43	10.55	0.085
114. CR (1) C 37	/636. BD* (1) C 35 - C 40	0.33	10.55	0.075
114. CR (1) C 37	/637. BD* (1) C 36 - C 37	0.70	10.72	0.110
114. CR (1) C 37	/649. BD* (1) N 42 - N 43	0.37	10.56	0.080
115. CR (1) C 38	/372. RY* (2) C 33	0.31	11.72	0.076
115. CR (1) C 38	/421. RY* (3) C 39	1.06	11.39	0.139
115. CR (1) C 38	/518. RY* (2) H 52	0.28	12.27	0.074
115. CR (1) C 38	/627. BD* (1) C 32 - C 33	0.32	10.45	0.074
115. CR (1) C 38	/631. BD* (1) C 33 - C 35	0.36	10.65	0.078
115. CR (1) C 38	/640. BD* (1) C 38 - C 39	0.40	10.68	0.082
115. CR (1) C 38	/643. BD* (1) C 39 - C 41	0.49	10.50	0.091
116. CR (1) C 39	/411. RY* (1) C 38	0.37	11.21	0.082
116. CR (1) C 39	/413. RY* (3) C 38	0.65	11.17	0.108
116. CR (1) C 39	/414. RY* (4) C 38	0.39	11.05	0.083
116. CR (1) C 39	/437. RY* (3) C 41	0.38	11.40	0.083
116. CR (1) C 39	/520. RY* (2) H 53	0.30	11.90	0.075
116. CR (1) C 39	/632. BD* (1) C 33 - C 38	0.61	10.48	0.101
116. CR (1) C 39	/640. BD* (1) C 38 - C 39	0.39	10.68	0.082
116. CR (1) C 39	/645. BD* (1) C 40 - C 41	0.29	10.69	0.070
117. CR (1) C 40	/389. RY* (3) C 35	0.37	11.58	0.083
117. CR (1) C 40	/437. RY* (3) C 41	1.06	11.40	0.139
117. CR (1) C 40	/631. BD* (1) C 33 - C 35	0.42	10.65	0.084
117. CR (1) C 40	/635. BD* (1) C 35 - C 37	0.34	10.47	0.077
117. CR (1) C 40	/643. BD* (1) C 39 - C 41	0.50	10.50	0.091
117. CR (1) C 40	/645. BD* (1) C 40 - C 41	0.39	10.69	0.081
118. CR (1) C 41	/421. RY* (3) C 39	0.39	11.38	0.084
118. CR (1) C 41	/429. RY* (3) C 40	0.54	11.15	0.098
118. CR (1) C 41	/430. RY* (4) C 40	0.55	11.10	0.099
118. CR (1) C 41	/524. RY* (2) H 55	0.33	12.30	0.081
118. CR (1) C 41	/636. BD* (1) C 35 - C 40	0.62	10.48	0.102
118. CR (1) C 41	/640. BD* (1) C 38 - C 39	0.29	10.67	0.070
118. CR (1) C 41	/645. BD* (1) C 40 - C 41	0.43	10.68	0.086
119. CR (1) N 42	/403. RY* (1) C 37	0.35	15.06	0.092
119. CR (1) N 42	/404. RY* (2) C 37	0.83	15.32	0.142
119. CR (1) N 42	/453. RY* (3) N 43	0.83	15.31	0.143
119. CR (1) N 42	/635. BD* (1) C 35 - C 37	0.26	14.66	0.079
119. CR (1) N 42	/637. BD* (1) C 36 - C 37	0.26	14.85	0.079
119. CR (1) N 42	/650. BD* (1) N 43 - C 44	0.69	14.57	0.128
120. CR (1) N 43	/445. RY* (3) N 42	0.93	15.59	0.152
120. CR (1) N 43	/459. RY* (1) C 44	0.65	15.42	0.126
120. CR (1) N 43	/460. RY* (2) C 44	0.33	15.68	0.091
120. CR (1) N 43	/639. BD* (1) C 37 - N 42	0.61	14.63	0.120
120. CR (1) N 43	/652. BD* (1) C 44 - C 46	0.34	14.86	0.090
121. CR (1) C 44	/469. RY* (3) C 45	0.42	11.42	0.088
121. CR (1) C 44	/475. RY* (1) C 46	0.56	11.30	0.101
121. CR (1) C 44	/477. RY* (3) C 46	0.99	11.64	0.136
121. CR (1) C 44	/649. BD* (1) N 42 - N 43	0.46	10.58	0.089

121. CR (1) C 44	/652. BD* (1) C 44 - C 46	0.62	10.71	0.104
121. CR (1) C 44	/656. BD* (1) C 46 - C 48	0.47	10.55	0.089
121. CR (1) C 44	/657. BD* (1) C 46 - O 62	0.39	10.42	0.081
122. CR (1) C 45	/459. RY* (1) C 44	0.42	11.19	0.087
122. CR (1) C 45	/461. RY* (3) C 44	0.28	10.86	0.069
122. CR (1) C 45	/485. RY* (3) C 47	0.99	11.42	0.134
122. CR (1) C 45	/526. RY* (2) H 56	0.30	12.30	0.077
122. CR (1) C 45	/650. BD* (1) N 43 - C 44	0.26	10.38	0.067
122. CR (1) C 45	/652. BD* (1) C 44 - C 46	0.39	10.63	0.082
122. CR (1) C 45	/654. BD* (1) C 45 - C 47	0.39	10.67	0.082
122. CR (1) C 45	/658. BD* (1) C 47 - C 49	0.45	10.50	0.087
123. CR (1) C 46	/460. RY* (2) C 44	1.55	11.53	0.169
123. CR (1) C 46	/493. RY* (3) C 48	0.48	11.40	0.093
123. CR (1) C 46	/650. BD* (1) N 43 - C 44	0.47	10.47	0.089
123. CR (1) C 46	/651. BD* (1) C 44 - C 45	0.49	10.55	0.091
123. CR (1) C 46	/652. BD* (1) C 44 - C 46	0.84	10.71	0.121
123. CR (1) C 46	/657. BD* (1) C 46 - O 62	0.43	10.43	0.085
123. CR (1) C 46	/660. BD* (1) C 48 - C 49	0.26	10.76	0.067
124. CR (1) C 47	/468. RY* (2) C 45	0.26	11.36	0.069
124. CR (1) C 47	/469. RY* (3) C 45	0.76	11.34	0.117
124. CR (1) C 47	/470. RY* (4) C 45	0.27	10.96	0.068
124. CR (1) C 47	/501. RY* (3) C 49	0.49	11.34	0.094
124. CR (1) C 47	/528. RY* (2) H 57	0.30	11.94	0.076
124. CR (1) C 47	/651. BD* (1) C 44 - C 45	0.48	10.48	0.090
124. CR (1) C 47	/654. BD* (1) C 45 - C 47	0.31	10.67	0.073
124. CR (1) C 47	/660. BD* (1) C 48 - C 49	0.29	10.69	0.071
125. CR (1) C 48	/477. RY* (3) C 46	0.44	11.56	0.091
125. CR (1) C 48	/478. RY* (4) C 46	0.41	11.06	0.085
125. CR (1) C 48	/501. RY* (3) C 49	1.07	11.33	0.139
125. CR (1) C 48	/530. RY* (2) H 58	0.29	12.31	0.075
125. CR (1) C 48	/652. BD* (1) C 44 - C 46	0.39	10.63	0.083
125. CR (1) C 48	/658. BD* (1) C 47 - C 49	0.50	10.50	0.092
125. CR (1) C 48	/660. BD* (1) C 48 - C 49	0.46	10.68	0.088
126. CR (1) C 49	/485. RY* (3) C 47	0.41	11.43	0.086
126. CR (1) C 49	/492. RY* (2) C 48	0.35	11.36	0.080
126. CR (1) C 49	/493. RY* (3) C 48	0.72	11.32	0.114
126. CR (1) C 49	/532. RY* (2) H 59	0.32	11.97	0.078
126. CR (1) C 49	/654. BD* (1) C 45 - C 47	0.27	10.67	0.067
126. CR (1) C 49	/656. BD* (1) C 46 - C 48	0.49	10.48	0.091
126. CR (1) C 49	/660. BD* (1) C 48 - C 49	0.32	10.69	0.074
127. CR (1) O 50	/363. RY* (1) C 32	0.60	19.81	0.138
127. CR (1) O 50	/364. RY* (2) C 32	0.47	19.84	0.123
127. CR (1) O 50	/534. RY* (2) H 60	0.30	20.87	0.101
128. CR (1) O 62	/475. RY* (1) C 46	1.36	20.11	0.209
128. CR (1) O 62	/652. BD* (1) C 44 - C 46	0.28	19.53	0.095
155. LP (1) N 42	/451. RY* (1) N 43	0.73	1.14	0.037
155. LP (1) N 42	/452. RY* (2) N 43	0.90	1.13	0.041
155. LP (1) N 42	/635. BD* (1) C 35 - C 37	3.47	0.75	0.065
155. LP (1) N 42	/637. BD* (1) C 36 - C 37	3.29	0.94	0.071
156. LP (2) N 42	/ 56. BD* (2) C 35 - C 37	2.21	0.42	0.050
156. LP (2) N 42	/403. RY* (1) C 37	0.62	1.31	0.038
156. LP (2) N 42	/444. RY* (2) N 42	1.08	2.47	0.070
156. LP (2) N 42	/445. RY* (3) N 42	0.63	1.79	0.045
156. LP (2) N 42	/452. RY* (2) N 43	0.31	1.29	0.027
156. LP (2) N 42	/635. BD* (1) C 35 - C 37	0.99	0.91	0.040
157. LP (1) N 43	/444. RY* (2) N 42	0.29	2.46	0.034
157. LP (1) N 43	/445. RY* (3) N 42	0.34	1.78	0.031
157. LP (1) N 43	/459. RY* (1) C 44	0.82	1.61	0.046
157. LP (1) N 43	/651. BD* (1) C 44 - C 45	0.57	0.89	0.028
157. LP (1) N 43	/652. BD* (1) C 44 - C 46	3.46	1.05	0.076
158. LP (2) N 43	/443. RY* (1) N 42	1.35	1.17	0.053
158. LP (2) N 43	/462. RY* (4) C 44	0.27	1.01	0.022
158. LP (2) N 43	/653. BD* (2) C 44 - C 46	16.07	0.32	0.097
159. LP (1) C 45	/160. LP (1) C 49	1.21	0.07	0.014
159. LP (1) C 45	/483. RY* (1) C 47	2.83	0.66	0.059
159. LP (1) C 45	/653. BD* (2) C 44 - C 46	30.50	0.19	0.098
160. LP (1) C 49	/483. RY* (1) C 47	0.98	0.60	0.043
160. LP (1) C 49	/491. RY* (1) C 48	1.94	0.62	0.061
160. LP (1) C 49	/500. RY* (2) C 49	1.43	0.62	0.053
161. LP (1) O 50	/363. RY* (1) C 32	0.78	1.45	0.043
161. LP (1) O 50	/364. RY* (2) C 32	0.68	1.47	0.040

161. LP (1) O 50	/627. BD* (1) C 32 - C 33	3.23	1.02	0.073
162. LP (2) O 50	/628. BD* (2) C 32 - C 33	7.23	0.29	0.080
163. LP (1) O 62	/475. RY* (1) C 46	1.06	1.80	0.055
163. LP (1) O 62	/652. BD* (1) C 44 - C 46	2.86	1.21	0.074
164. LP (2) O 62	/652. BD* (1) C 44 - C 46	0.54	0.96	0.029
164. LP (2) O 62	/653. BD* (2) C 44 - C 46	4.33	0.37	0.056
164. LP (2) O 62	/656. BD* (1) C 46 - C 48	1.86	0.80	0.049
165. LP (3) O 62	/444. RY* (2) N 42	0.29	2.43	0.036
165. LP (3) O 62	/475. RY* (1) C 46	0.27	1.61	0.028
165. LP (3) O 62	/538. RY* (2) O 62	0.37	2.28	0.040
165. LP (3) O 62	/539. RY* (3) O 62	0.32	2.04	0.035
165. LP (3) O 62	/653. BD* (2) C 44 - C 46	3.78	0.43	0.053
165. LP (3) O 62	/656. BD* (1) C 46 - C 48	2.17	0.86	0.058
628. BD* (2) C 32 - C 33	/363. RY* (1) C 32	0.33	0.89	0.030
628. BD* (2) C 32 - C 33	/364. RY* (2) C 32	0.40	0.92	0.034
628. BD* (2) C 32 - C 33	/371. RY* (1) C 33	0.91	0.64	0.042
628. BD* (2) C 32 - C 33	/379. RY* (1) C 34	1.05	0.54	0.042
628. BD* (2) C 32 - C 33	/387. RY* (1) C 35	0.56	0.65	0.033
628. BD* (2) C 32 - C 33	/412. RY* (2) C 38	0.34	0.60	0.025
628. BD* (2) C 32 - C 33	/507. RY* (1) O 50	0.39	0.92	0.033
628. BD* (2) C 32 - C 33	/641. BD* (2) C 38 - C 39	31.24	0.08	0.073
628. BD* (2) C 32 - C 33	/646. BD* (2) C 40 - C 41	0.27	0.08	0.006
641. BD* (2) C 38 - C 39	/412. RY* (2) C 38	1.56	0.52	0.073
641. BD* (2) C 38 - C 39	/419. RY* (1) C 39	1.54	0.47	0.069
646. BD* (2) C 40 - C 41	/427. RY* (1) C 40	0.96	0.81	0.064
646. BD* (2) C 40 - C 41	/428. RY* (2) C 40	0.52	0.94	0.051
646. BD* (2) C 40 - C 41	/435. RY* (1) C 41	2.02	0.48	0.071
653. BD* (2) C 44 - C 46	/461. RY* (3) C 44	0.38	0.82	0.047
653. BD* (2) C 44 - C 46	/462. RY* (4) C 44	0.53	0.69	0.051
653. BD* (2) C 44 - C 46	/467. RY* (1) C 45	0.40	0.48	0.036
653. BD* (2) C 44 - C 46	/476. RY* (2) C 46	0.60	0.62	0.051
653. BD* (2) C 44 - C 46	/491. RY* (1) C 48	0.26	0.50	0.030

from unit 2 to unit 3

54. BD (1) C 35 - C 37	/168. LP (6)Cr 63	0.05	1.30	0.011
54. BD (1) C 35 - C 37	/545. RY* (1)Cr 63	0.03	1.17	0.008
55. BD (2) C 35 - C 37	/166. LP (4)Cr 63	0.11	0.14	0.006
55. BD (2) C 35 - C 37	/171. LP* (9)Cr 63	0.04	2.09	0.013
58. BD (1) C 36 - C 37	/171. LP* (9)Cr 63	0.03	2.58	0.012
58. BD (1) C 36 - C 37	/545. RY* (1)Cr 63	0.03	1.28	0.008
60. BD (1) C 37 - N 42	/166. LP (4)Cr 63	0.15	0.66	0.016
60. BD (1) C 37 - N 42	/167. LP (5)Cr 63	0.05	0.67	0.008
60. BD (1) C 37 - N 42	/168. LP (6)Cr 63	0.70	1.44	0.043
60. BD (1) C 37 - N 42	/169. LP (7)Cr 63	0.13	2.41	0.023
60. BD (1) C 37 - N 42	/171. LP* (9)Cr 63	0.04	2.61	0.013
60. BD (1) C 37 - N 42	/545. RY* (1)Cr 63	0.05	1.31	0.010
60. BD (1) C 37 - N 42	/546. RY* (2)Cr 63	0.06	2.37	0.016
60. BD (1) C 37 - N 42	/547. RY* (3)Cr 63	0.06	2.50	0.015
60. BD (1) C 37 - N 42	/548. RY* (4)Cr 63	0.06	2.34	0.015
60. BD (1) C 37 - N 42	/552. RY* (8)Cr 63	0.03	9.49	0.021
60. BD (1) C 37 - N 42	/556. RY* (12)Cr 63	0.03	8.84	0.020
60. BD (1) C 37 - N 42	/557. RY* (13)Cr 63	0.03	5.22	0.015
70. BD (1) N 42 - N 43	/166. LP (4)Cr 63	0.08	0.85	0.014
70. BD (1) N 42 - N 43	/167. LP (5)Cr 63	0.04	0.87	0.009
70. BD (1) N 42 - N 43	/168. LP (6)Cr 63	0.87	1.64	0.051
70. BD (1) N 42 - N 43	/169. LP (7)Cr 63	0.06	2.61	0.015
70. BD (1) N 42 - N 43	/170. LP* (8)Cr 63	0.05	2.81	0.015
70. BD (1) N 42 - N 43	/545. RY* (1)Cr 63	0.03	1.51	0.008
70. BD (1) N 42 - N 43	/546. RY* (2)Cr 63	0.22	2.57	0.030
70. BD (1) N 42 - N 43	/548. RY* (4)Cr 63	0.03	2.54	0.012
70. BD (1) N 42 - N 43	/549. RY* (5)Cr 63	0.05	2.38	0.014
70. BD (1) N 42 - N 43	/550. RY* (6)Cr 63	0.17	3.80	0.032
70. BD (1) N 42 - N 43	/552. RY* (8)Cr 63	0.03	9.69	0.021
70. BD (1) N 42 - N 43	/554. RY* (10)Cr 63	0.05	10.08	0.028
70. BD (1) N 42 - N 43	/555. RY* (11)Cr 63	0.04	5.35	0.018
70. BD (1) N 42 - N 43	/561. RY* (17)Cr 63	0.07	38.54	0.065
70. BD (1) N 42 - N 43	/564. RY* (20)Cr 63	0.03	675.27	0.180
71. BD (1) N 43 - C 44	/168. LP (6)Cr 63	0.10	1.44	0.017
71. BD (1) N 43 - C 44	/171. LP* (9)Cr 63	0.03	2.61	0.011
71. BD (1) N 43 - C 44	/545. RY* (1)Cr 63	0.03	1.31	0.008
71. BD (1) N 43 - C 44	/546. RY* (2)Cr 63	0.03	2.38	0.011

71. BD (1) N 43 - C 44	/550. RY* (6)Cr 63	0.07	3.61	0.020
71. BD (1) N 43 - C 44	/556. RY* (12)Cr 63	0.09	8.84	0.037
71. BD (1) N 43 - C 44	/561. RY* (17)Cr 63	0.12	38.34	0.087
71. BD (1) N 43 - C 44	/564. RY* (20)Cr 63	0.04	675.08	0.213
71. BD (1) N 43 - C 44	/565. RY* (21)Cr 63	0.04	27.02	0.040
73. BD (1) C 44 - C 46	/168. LP (6)Cr 63	0.14	1.42	0.019
73. BD (1) C 44 - C 46	/545. RY* (1)Cr 63	0.03	1.29	0.008
73. BD (1) C 44 - C 46	/546. RY* (2)Cr 63	0.07	2.35	0.016
73. BD (1) C 44 - C 46	/561. RY* (17)Cr 63	0.03	38.32	0.044
74. BD (2) C 44 - C 46	/166. LP (4)Cr 63	0.03	0.17	0.003
74. BD (2) C 44 - C 46	/167. LP (5)Cr 63	0.04	0.18	0.004
74. BD (2) C 44 - C 46	/168. LP (6)Cr 63	0.06	0.95	0.010
74. BD (2) C 44 - C 46	/169. LP (7)Cr 63	0.03	1.92	0.010
77. BD (1) C 46 - C 48	/168. LP (6)Cr 63	0.15	1.30	0.019
77. BD (1) C 46 - C 48	/169. LP (7)Cr 63	0.08	2.27	0.017
77. BD (1) C 46 - C 48	/545. RY* (1)Cr 63	0.03	1.17	0.008
77. BD (1) C 46 - C 48	/561. RY* (17)Cr 63	0.03	38.20	0.040
78. BD (1) C 46 - O 62	/166. LP (4)Cr 63	0.07	0.71	0.011
78. BD (1) C 46 - O 62	/167. LP (5)Cr 63	0.18	0.73	0.017
78. BD (1) C 46 - O 62	/168. LP (6)Cr 63	0.92	1.49	0.050
78. BD (1) C 46 - O 62	/169. LP (7)Cr 63	0.13	2.47	0.023
78. BD (1) C 46 - O 62	/171. LP* (9)Cr 63	0.03	2.66	0.011
78. BD (1) C 46 - O 62	/546. RY* (2)Cr 63	0.14	2.43	0.024
78. BD (1) C 46 - O 62	/549. RY* (5)Cr 63	0.05	2.24	0.013
78. BD (1) C 46 - O 62	/550. RY* (6)Cr 63	0.08	3.66	0.022
78. BD (1) C 46 - O 62	/552. RY* (8)Cr 63	0.05	9.55	0.028
78. BD (1) C 46 - O 62	/554. RY* (10)Cr 63	0.04	9.94	0.026
119. CR (1) N 42	/166. LP (4)Cr 63	0.17	14.08	0.079
119. CR (1) N 42	/167. LP (5)Cr 63	0.05	14.10	0.042
119. CR (1) N 42	/168. LP (6)Cr 63	1.08	14.86	0.172
120. CR (1) N 43	/168. LP (6)Cr 63	0.03	14.91	0.029
123. CR (1) C 46	/168. LP (6)Cr 63	0.05	10.76	0.032
128. CR (1) O 62	/166. LP (4)Cr 63	0.11	18.79	0.076
128. CR (1) O 62	/167. LP (5)Cr 63	0.22	18.80	0.099
128. CR (1) O 62	/168. LP (6)Cr 63	1.45	19.57	0.229
155. LP (1) N 42	/166. LP (4)Cr 63	0.37	0.17	0.013
155. LP (1) N 42	/167. LP (5)Cr 63	0.12	0.19	0.007
155. LP (1) N 42	/168. LP (6)Cr 63	0.38	0.95	0.025
155. LP (1) N 42	/169. LP (7)Cr 63	0.05	1.93	0.013
155. LP (1) N 42	/170. LP* (8)Cr 63	0.06	2.13	0.015
155. LP (1) N 42	/550. RY* (6)Cr 63	0.07	3.12	0.018
155. LP (1) N 42	/561. RY* (17)Cr 63	0.04	37.85	0.049
156. LP (2) N 42	/166. LP (4)Cr 63	20.07	0.33	0.121
156. LP (2) N 42	/167. LP (5)Cr 63	4.80	0.35	0.058
156. LP (2) N 42	/168. LP (6)Cr 63	18.63	1.11	0.183
156. LP (2) N 42	/169. LP (7)Cr 63	0.43	2.09	0.040
156. LP (2) N 42	/170. LP* (8)Cr 63	0.57	2.29	0.049
156. LP (2) N 42	/171. LP* (9)Cr 63	1.08	2.28	0.067
156. LP (2) N 42	/545. RY* (1)Cr 63	0.33	0.98	0.024
156. LP (2) N 42	/546. RY* (2)Cr 63	0.32	2.05	0.034
156. LP (2) N 42	/547. RY* (3)Cr 63	0.09	2.18	0.019
156. LP (2) N 42	/548. RY* (4)Cr 63	0.04	2.02	0.012
156. LP (2) N 42	/549. RY* (5)Cr 63	0.16	1.86	0.024
156. LP (2) N 42	/550. RY* (6)Cr 63	2.36	3.28	0.119
156. LP (2) N 42	/551. RY* (7)Cr 63	0.05	2.37	0.015
156. LP (2) N 42	/554. RY* (10)Cr 63	0.06	9.56	0.032
156. LP (2) N 42	/556. RY* (12)Cr 63	0.20	8.51	0.056
156. LP (2) N 42	/560. RY* (16)Cr 63	0.31	115.37	0.256
156. LP (2) N 42	/561. RY* (17)Cr 63	1.81	38.01	0.354
156. LP (2) N 42	/564. RY* (20)Cr 63	0.58	674.75	0.846
156. LP (2) N 42	/565. RY* (21)Cr 63	0.07	26.69	0.058
157. LP (1) N 43	/166. LP (4)Cr 63	0.05	0.31	0.007
157. LP (1) N 43	/168. LP (6)Cr 63	0.15	1.09	0.017
157. LP (1) N 43	/170. LP* (8)Cr 63	0.06	2.27	0.015
157. LP (1) N 43	/171. LP* (9)Cr 63	0.03	2.26	0.011
159. LP (1) C 45	/166. LP (4)Cr 63	0.03	0.03	0.002
163. LP (1) O 62	/166. LP (4)Cr 63	0.86	0.47	0.032
163. LP (1) O 62	/167. LP (5)Cr 63	1.57	0.49	0.042
163. LP (1) O 62	/168. LP (6)Cr 63	6.38	1.25	0.121
163. LP (1) O 62	/169. LP (7)Cr 63	0.69	2.23	0.050
163. LP (1) O 62	/171. LP* (9)Cr 63	0.13	2.42	0.022

163. LP (1) O 62	/545. RY* (1)Cr 63	0.11	1.12	0.014
163. LP (1) O 62	/546. RY* (2)Cr 63	0.38	2.19	0.036
163. LP (1) O 62	/547. RY* (3)Cr 63	0.59	2.32	0.047
163. LP (1) O 62	/548. RY* (4)Cr 63	0.07	2.16	0.015
163. LP (1) O 62	/549. RY* (5)Cr 63	0.07	2.00	0.015
163. LP (1) O 62	/550. RY* (6)Cr 63	0.32	3.42	0.042
163. LP (1) O 62	/551. RY* (7)Cr 63	0.09	2.51	0.019
163. LP (1) O 62	/552. RY* (8)Cr 63	0.08	9.31	0.035
163. LP (1) O 62	/560. RY* (16)Cr 63	0.09	115.51	0.133
163. LP (1) O 62	/561. RY* (17)Cr 63	0.61	38.15	0.195
163. LP (1) O 62	/564. RY* (20)Cr 63	0.22	674.89	0.491
164. LP (2) O 62	/166. LP (4)Cr 63	0.79	0.22	0.021
164. LP (2) O 62	/167. LP (5)Cr 63	1.27	0.24	0.026
164. LP (2) O 62	/168. LP (6)Cr 63	2.36	1.00	0.065
164. LP (2) O 62	/169. LP (7)Cr 63	0.41	1.98	0.036
164. LP (2) O 62	/170. LP* (8)Cr 63	0.14	2.18	0.023
164. LP (2) O 62	/545. RY* (1)Cr 63	0.08	0.87	0.011
164. LP (2) O 62	/547. RY* (3)Cr 63	0.22	2.07	0.028
164. LP (2) O 62	/550. RY* (6)Cr 63	0.13	3.17	0.026
164. LP (2) O 62	/551. RY* (7)Cr 63	0.03	2.26	0.011
164. LP (2) O 62	/552. RY* (8)Cr 63	0.05	9.06	0.028
164. LP (2) O 62	/560. RY* (16)Cr 63	0.03	115.26	0.077
164. LP (2) O 62	/561. RY* (17)Cr 63	0.20	37.90	0.111
164. LP (2) O 62	/564. RY* (20)Cr 63	0.07	674.64	0.275
165. LP (3) O 62	/166. LP (4)Cr 63	12.93	0.28	0.089
165. LP (3) O 62	/167. LP (5)Cr 63	16.19	0.30	0.097
165. LP (3) O 62	/168. LP (6)Cr 63	14.87	1.06	0.159
165. LP (3) O 62	/169. LP (7)Cr 63	3.26	2.04	0.111
165. LP (3) O 62	/171. LP* (9)Cr 63	0.03	2.23	0.010
165. LP (3) O 62	/545. RY* (1)Cr 63	0.50	0.93	0.030
165. LP (3) O 62	/546. RY* (2)Cr 63	0.25	2.00	0.031
165. LP (3) O 62	/547. RY* (3)Cr 63	0.89	2.13	0.059
165. LP (3) O 62	/548. RY* (4)Cr 63	0.07	1.97	0.016
165. LP (3) O 62	/549. RY* (5)Cr 63	0.18	1.81	0.025
165. LP (3) O 62	/550. RY* (6)Cr 63	0.40	3.23	0.049
165. LP (3) O 62	/551. RY* (7)Cr 63	0.06	2.33	0.016
165. LP (3) O 62	/552. RY* (8)Cr 63	0.21	9.12	0.060
165. LP (3) O 62	/559. RY* (15)Cr 63	0.07	26.12	0.060
165. LP (3) O 62	/560. RY* (16)Cr 63	0.21	115.32	0.212
165. LP (3) O 62	/561. RY* (17)Cr 63	1.24	37.96	0.296
165. LP (3) O 62	/564. RY* (20)Cr 63	0.38	674.70	0.695

from unit 2 to unit 4

54. BD (1) C 35 - C 37	/568. RY* (1) H 65	0.03	1.39	0.009
55. BD (2) C 35 - C 37	/665. BD* (1) H 65 - O 66	0.23	0.61	0.015
68. BD (1) C 40 - H 54	/664. BD* (1) H 64 - O 66	0.03	0.84	0.006
71. BD (1) N 43 - C 44	/570. RY* (1) O 66	0.03	2.93	0.012
156. LP (2) N 42	/566. RY* (1) H 64	0.12	1.09	0.016
156. LP (2) N 42	/568. RY* (1) H 65	0.13	1.21	0.017
156. LP (2) N 42	/570. RY* (1) O 66	0.56	2.60	0.052
156. LP (2) N 42	/572. RY* (3) O 66	0.07	1.78	0.015
156. LP (2) N 42	/573. RY* (4) O 66	0.03	1.61	0.009
156. LP (2) N 42	/664. BD* (1) H 64 - O 66	0.03	0.79	0.006
156. LP (2) N 42	/665. BD* (1) H 65 - O 66	0.06	0.80	0.009
163. LP (1) O 62	/566. RY* (1) H 64	0.03	1.23	0.008
163. LP (1) O 62	/568. RY* (1) H 65	0.07	1.35	0.012
163. LP (1) O 62	/570. RY* (1) O 66	0.19	2.74	0.029
164. LP (2) O 62	/568. RY* (1) H 65	0.03	1.10	0.008
164. LP (2) O 62	/570. RY* (1) O 66	0.08	2.49	0.018
164. LP (2) O 62	/664. BD* (1) H 64 - O 66	0.08	0.69	0.010
164. LP (2) O 62	/665. BD* (1) H 65 - O 66	0.03	0.69	0.006
165. LP (3) O 62	/566. RY* (1) H 64	0.12	1.04	0.015
165. LP (3) O 62	/568. RY* (1) H 65	0.06	1.16	0.012
165. LP (3) O 62	/570. RY* (1) O 66	0.35	2.55	0.041
165. LP (3) O 62	/573. RY* (4) O 66	0.04	1.56	0.011
165. LP (3) O 62	/664. BD* (1) H 64 - O 66	0.05	0.75	0.009
646. BD* (2) C 40 - C 41	/664. BD* (1) H 64 - O 66	0.06	0.30	0.009

from unit 2 to unit 5

58. BD (1) C 36 - C 37	/580. RY* (3)C1 67	0.03	1.66	0.009
71. BD (1) N 43 - C 44	/581. RY* (4)C1 67	0.03	2.43	0.011

156. LP (2) N 42	/579. RY* (2)C1 67	0.41	1.12	0.029
156. LP (2) N 42	/580. RY* (3)C1 67	0.16	1.36	0.020
156. LP (2) N 42	/581. RY* (4)C1 67	0.33	2.10	0.036
163. LP (1) O 62	/579. RY* (2)C1 67	0.07	1.26	0.012
163. LP (1) O 62	/580. RY* (3)C1 67	0.04	1.50	0.010
163. LP (1) O 62	/581. RY* (4)C1 67	0.18	2.24	0.026
164. LP (2) O 62	/579. RY* (2)C1 67	0.03	1.01	0.007
164. LP (2) O 62	/581. RY* (4)C1 67	0.06	1.99	0.014
165. LP (3) O 62	/579. RY* (2)C1 67	0.13	1.08	0.016
165. LP (3) O 62	/580. RY* (3)C1 67	0.06	1.32	0.012
165. LP (3) O 62	/581. RY* (4)C1 67	0.36	2.05	0.037
from unit 3 to unit 1				
129. CR (1)Cr 63	/602. BD* (1) C 6 - N 11	0.03	215.68	0.094
129. CR (1)Cr 63	/620. BD* (1) C 15 - O 19	0.03	215.63	0.097
130. CR (2)Cr 63	/602. BD* (1) C 6 - N 11	0.15	24.51	0.077
130. CR (2)Cr 63	/612. BD* (1) N 11 - N 12	0.10	24.62	0.063
130. CR (2)Cr 63	/620. BD* (1) C 15 - O 19	0.16	24.46	0.080
131. CR (3)Cr 63	/262. RY* (2) N 11	0.03	5.76	0.015
131. CR (3)Cr 63	/263. RY* (3) N 11	0.12	5.17	0.032
131. CR (3)Cr 63	/293. RY* (1) C 15	0.03	5.00	0.014
131. CR (3)Cr 63	/326. RY* (2) O 19	0.03	5.49	0.015
131. CR (3)Cr 63	/602. BD* (1) C 6 - N 11	0.73	4.19	0.070
131. CR (3)Cr 63	/612. BD* (1) N 11 - N 12	0.48	4.30	0.058
131. CR (3)Cr 63	/619. BD* (1) C 15 - C 17	0.03	4.27	0.015
131. CR (3)Cr 63	/620. BD* (1) C 15 - O 19	0.90	4.14	0.078
133. CR (5)Cr 63	/262. RY* (2) N 11	0.03	3.86	0.014
133. CR (5)Cr 63	/263. RY* (3) N 11	0.08	3.27	0.021
133. CR (5)Cr 63	/602. BD* (1) C 6 - N 11	0.28	2.30	0.032
133. CR (5)Cr 63	/612. BD* (1) N 11 - N 12	0.24	2.41	0.031
135. CR (7)Cr 63	/612. BD* (1) N 11 - N 12	0.04	2.41	0.012
135. CR (7)Cr 63	/620. BD* (1) C 15 - O 19	0.50	2.25	0.043
137. CR (9)Cr 63	/602. BD* (1) C 6 - N 11	0.04	2.30	0.013
166. LP (4)Cr 63	/ 13. BD* (2) C 4 - C 6	0.19	0.10	0.006
166. LP (4)Cr 63	/262. RY* (2) N 11	0.20	2.08	0.041
166. LP (4)Cr 63	/264. RY* (4) N 11	0.04	1.64	0.016
166. LP (4)Cr 63	/602. BD* (1) C 6 - N 11	0.04	0.51	0.009
166. LP (4)Cr 63	/612. BD* (1) N 11 - N 12	0.03	0.62	0.009
166. LP (4)Cr 63	/613. BD* (1) N 12 - C 13	0.04	0.50	0.009
167. LP (5)Cr 63	/ 13. BD* (2) C 4 - C 6	0.07	0.08	0.003
167. LP (5)Cr 63	/222. RY* (2) C 6	0.03	1.33	0.013
167. LP (5)Cr 63	/262. RY* (2) N 11	0.07	2.06	0.027
167. LP (5)Cr 63	/263. RY* (3) N 11	0.05	1.47	0.019
167. LP (5)Cr 63	/293. RY* (1) C 15	0.05	1.31	0.018
167. LP (5)Cr 63	/326. RY* (2) O 19	0.29	1.79	0.050
167. LP (5)Cr 63	/327. RY* (3) O 19	0.34	1.99	0.058
167. LP (5)Cr 63	/620. BD* (1) C 15 - O 19	0.07	0.44	0.012
168. LP (6)Cr 63	/192. RY* (4) C 2	0.07	0.43	0.018
168. LP (6)Cr 63	/206. RY* (2) C 4	0.05	0.96	0.023
168. LP (6)Cr 63	/207. RY* (3) C 4	0.11	0.96	0.035
168. LP (6)Cr 63	/208. RY* (4) C 4	0.05	0.33	0.013
168. LP (6)Cr 63	/214. RY* (2) C 5	0.08	0.76	0.026
168. LP (6)Cr 63	/215. RY* (3) C 5	0.10	0.60	0.026
168. LP (6)Cr 63	/216. RY* (4) C 5	0.04	0.51	0.016
168. LP (6)Cr 63	/218. RY* (6) C 5	0.03	3.35	0.034
168. LP (6)Cr 63	/221. RY* (1) C 6	1.36	0.22	0.057
168. LP (6)Cr 63	/222. RY* (2) C 6	4.66	0.56	0.171
168. LP (6)Cr 63	/223. RY* (3) C 6	0.72	0.47	0.062
168. LP (6)Cr 63	/224. RY* (4) C 6	3.31	0.40	0.122
168. LP (6)Cr 63	/225. RY* (5) C 6	0.04	2.40	0.034
168. LP (6)Cr 63	/226. RY* (6) C 6	0.54	2.36	0.121
168. LP (6)Cr 63	/228. RY* (8) C 6	0.05	2.17	0.035
168. LP (6)Cr 63	/245. RY* (1) C 9	0.07	0.60	0.021
168. LP (6)Cr 63	/247. RY* (3) C 9	0.13	0.49	0.027
168. LP (6)Cr 63	/248. RY* (4) C 9	0.03	0.47	0.013
168. LP (6)Cr 63	/261. RY* (1) N 11	0.19	0.23	0.022
168. LP (6)Cr 63	/262. RY* (2) N 11	16.50	1.30	0.488
168. LP (6)Cr 63	/263. RY* (3) N 11	10.17	0.70	0.284
168. LP (6)Cr 63	/264. RY* (4) N 11	0.10	0.86	0.030
168. LP (6)Cr 63	/266. RY* (6) N 11	0.04	3.35	0.038
168. LP (6)Cr 63	/269. RY* (1) N 12	0.60	0.16	0.032

168. LP (6)Cr 63	/270. RY* (2) N 12	1.33	0.18	0.051
168. LP (6)Cr 63	/271. RY* (3) N 12	1.58	0.48	0.093
168. LP (6)Cr 63	/272. RY* (4) N 12	0.64	1.03	0.086
168. LP (6)Cr 63	/275. RY* (7) N 12	0.08	3.06	0.052
168. LP (6)Cr 63	/278. RY* (2) C 13	0.12	0.79	0.032
168. LP (6)Cr 63	/279. RY* (3) C 13	1.38	0.28	0.067
168. LP (6)Cr 63	/282. RY* (6) C 13	0.05	2.52	0.039
168. LP (6)Cr 63	/287. RY* (3) C 14	0.03	0.65	0.016
168. LP (6)Cr 63	/288. RY* (4) C 14	0.16	0.28	0.022
168. LP (6)Cr 63	/293. RY* (1) C 15	8.19	0.54	0.221
168. LP (6)Cr 63	/295. RY* (3) C 15	1.42	0.88	0.118
168. LP (6)Cr 63	/296. RY* (4) C 15	0.49	0.37	0.045
168. LP (6)Cr 63	/297. RY* (5) C 15	0.82	3.17	0.172
168. LP (6)Cr 63	/299. RY* (7) C 15	0.08	2.53	0.047
168. LP (6)Cr 63	/302. RY* (2) C 16	0.06	0.59	0.019
168. LP (6)Cr 63	/310. RY* (2) C 17	0.37	0.69	0.054
168. LP (6)Cr 63	/312. RY* (4) C 17	0.08	0.26	0.015
168. LP (6)Cr 63	/317. RY* (1) C 18	0.03	0.56	0.014
168. LP (6)Cr 63	/326. RY* (2) O 19	6.21	1.02	0.268
168. LP (6)Cr 63	/327. RY* (3) O 19	7.62	1.22	0.325
168. LP (6)Cr 63	/329. RY* (5) O 19	0.06	4.53	0.054
168. LP (6)Cr 63	/608. BD* (1) C 9 - C 10	0.16	0.02	0.006
168. LP (6)Cr 63	/623. BD* (1) C 17 - C 18	0.13	0.01	0.004
169. LP (7)Cr 63	/262. RY* (2) N 11	0.07	0.32	0.052
169. LP (7)Cr 63	/272. RY* (4) N 12	0.18	0.05	0.039
169. LP (7)Cr 63	/298. RY* (6) C 15	0.03	1.99	0.097
169. LP (7)Cr 63	/326. RY* (2) O 19	1.70	0.05	0.112
169. LP (7)Cr 63	/327. RY* (3) O 19	2.04	0.25	0.279
169. LP (7)Cr 63	/328. RY* (4) O 19	0.07	0.25	0.054
172. LP (1)Cr 63	/325. RY* (1) O 19	0.04	1.37	0.009
172. LP (1)Cr 63	/602. BD* (1) C 6 - N 11	0.04	0.65	0.007
172. LP (1)Cr 63	/620. BD* (1) C 15 - O 19	0.09	0.59	0.009
173. LP (2)Cr 63	/261. RY* (1) N 11	0.19	1.15	0.019
173. LP (2)Cr 63	/620. BD* (1) C 15 - O 19	0.14	0.59	0.012
174. LP (3)Cr 63	/ 13. BD* (2) C 4 - C 6	0.07	0.23	0.007
174. LP (3)Cr 63	/262. RY* (2) N 11	0.04	2.21	0.012
174. LP (3)Cr 63	/264. RY* (4) N 11	0.12	1.77	0.018
174. LP (3)Cr 63	/265. RY* (5) N 11	0.03	4.18	0.015
174. LP (3)Cr 63	/277. RY* (1) C 13	0.03	1.40	0.009
174. LP (3)Cr 63	/602. BD* (1) C 6 - N 11	0.30	0.64	0.018
174. LP (3)Cr 63	/612. BD* (1) N 11 - N 12	0.54	0.76	0.026
174. LP (3)Cr 63	/620. BD* (1) C 15 - O 19	0.04	0.59	0.006

from unit 3 to unit 2

129. CR (1)Cr 63	/639. BD* (1) C 37 - N 42	0.03	215.67	0.094
129. CR (1)Cr 63	/657. BD* (1) C 46 - O 62	0.03	215.62	0.099
130. CR (2)Cr 63	/639. BD* (1) C 37 - N 42	0.15	24.50	0.077
130. CR (2)Cr 63	/649. BD* (1) N 42 - N 43	0.08	24.61	0.057
130. CR (2)Cr 63	/657. BD* (1) C 46 - O 62	0.17	24.45	0.081
131. CR (3)Cr 63	/445. RY* (3) N 42	0.09	5.14	0.027
131. CR (3)Cr 63	/475. RY* (1) C 46	0.03	5.01	0.016
131. CR (3)Cr 63	/538. RY* (2) O 62	0.04	5.68	0.018
131. CR (3)Cr 63	/639. BD* (1) C 37 - N 42	0.73	4.18	0.071
131. CR (3)Cr 63	/649. BD* (1) N 42 - N 43	0.41	4.29	0.053
131. CR (3)Cr 63	/650. BD* (1) N 43 - C 44	0.03	4.17	0.013
131. CR (3)Cr 63	/656. BD* (1) C 46 - C 48	0.04	4.26	0.016
131. CR (3)Cr 63	/657. BD* (1) C 46 - O 62	0.92	4.13	0.078
133. CR (5)Cr 63	/444. RY* (2) N 42	0.03	3.93	0.013
133. CR (5)Cr 63	/445. RY* (3) N 42	0.06	3.25	0.018
133. CR (5)Cr 63	/639. BD* (1) C 37 - N 42	0.30	2.29	0.033
133. CR (5)Cr 63	/649. BD* (1) N 42 - N 43	0.21	2.40	0.029
133. CR (5)Cr 63	/657. BD* (1) C 46 - O 62	0.04	2.24	0.012
135. CR (7)Cr 63	/649. BD* (1) N 42 - N 43	0.03	2.40	0.011
137. CR (9)Cr 63	/639. BD* (1) C 37 - N 42	0.05	2.29	0.013
137. CR (9)Cr 63	/657. BD* (1) C 46 - O 62	0.47	2.24	0.041
166. LP (4)Cr 63	/ 56. BD* (2) C 35 - C 37	0.20	0.09	0.006
166. LP (4)Cr 63	/444. RY* (2) N 42	0.24	2.15	0.044
166. LP (4)Cr 63	/445. RY* (3) N 42	0.06	1.46	0.019
166. LP (4)Cr 63	/446. RY* (4) N 42	0.04	1.60	0.016
166. LP (4)Cr 63	/538. RY* (2) O 62	0.22	2.00	0.042
166. LP (4)Cr 63	/539. RY* (3) O 62	0.12	1.76	0.029

166. LP (4)Cr 63	/639. BD* (1) C 37 - N 42	0.05	0.50	0.010
166. LP (4)Cr 63	/650. BD* (1) N 43 - C 44	0.05	0.49	0.009
166. LP (4)Cr 63	/657. BD* (1) C 46 - O 62	0.05	0.45	0.009
167. LP (5)Cr 63	/404. RY* (2) C 37	0.03	1.22	0.014
167. LP (5)Cr 63	/444. RY* (2) N 42	0.04	2.13	0.020
167. LP (5)Cr 63	/445. RY* (3) N 42	0.07	1.45	0.022
167. LP (5)Cr 63	/475. RY* (1) C 46	0.03	1.31	0.014
167. LP (5)Cr 63	/657. BD* (1) C 46 - O 62	0.05	0.44	0.010
168. LP (6)Cr 63	/374. RY* (4) C 33	0.05	0.38	0.015
168. LP (6)Cr 63	/388. RY* (2) C 35	0.07	0.92	0.027
168. LP (6)Cr 63	/389. RY* (3) C 35	0.15	0.90	0.038
168. LP (6)Cr 63	/390. RY* (4) C 35	0.06	0.30	0.014
168. LP (6)Cr 63	/392. RY* (6) C 35	0.03	2.54	0.028
168. LP (6)Cr 63	/396. RY* (2) C 36	0.03	0.62	0.015
168. LP (6)Cr 63	/397. RY* (3) C 36	0.11	0.52	0.026
168. LP (6)Cr 63	/398. RY* (4) C 36	0.07	0.53	0.020
168. LP (6)Cr 63	/403. RY* (1) C 37	1.95	0.20	0.065
168. LP (6)Cr 63	/404. RY* (2) C 37	7.55	0.45	0.195
168. LP (6)Cr 63	/405. RY* (3) C 37	0.77	0.42	0.061
168. LP (6)Cr 63	/406. RY* (4) C 37	0.66	0.41	0.055
168. LP (6)Cr 63	/408. RY* (6) C 37	0.56	2.92	0.137
168. LP (6)Cr 63	/410. RY* (8) C 37	0.13	2.14	0.056
168. LP (6)Cr 63	/428. RY* (2) C 40	0.04	0.32	0.011
168. LP (6)Cr 63	/429. RY* (3) C 40	0.19	0.48	0.032
168. LP (6)Cr 63	/430. RY* (4) C 40	0.04	0.43	0.014
168. LP (6)Cr 63	/443. RY* (1) N 42	0.23	0.22	0.024
168. LP (6)Cr 63	/444. RY* (2) N 42	15.38	1.36	0.483
168. LP (6)Cr 63	/445. RY* (3) N 42	13.54	0.68	0.323
168. LP (6)Cr 63	/446. RY* (4) N 42	0.08	0.81	0.027
168. LP (6)Cr 63	/447. RY* (5) N 42	0.05	3.24	0.042
168. LP (6)Cr 63	/448. RY* (6) N 42	0.05	3.37	0.043
168. LP (6)Cr 63	/451. RY* (1) N 43	1.70	0.18	0.059
168. LP (6)Cr 63	/452. RY* (2) N 43	1.39	0.18	0.052
168. LP (6)Cr 63	/453. RY* (3) N 43	2.84	0.44	0.119
168. LP (6)Cr 63	/454. RY* (4) N 43	0.24	1.03	0.053
168. LP (6)Cr 63	/455. RY* (5) N 43	0.05	3.54	0.043
168. LP (6)Cr 63	/457. RY* (7) N 43	0.05	3.11	0.042
168. LP (6)Cr 63	/460. RY* (2) C 44	0.06	0.77	0.023
168. LP (6)Cr 63	/461. RY* (3) C 44	1.45	0.19	0.055
168. LP (6)Cr 63	/462. RY* (4) C 44	0.91	0.06	0.024
168. LP (6)Cr 63	/463. RY* (5) C 44	0.03	2.10	0.029
168. LP (6)Cr 63	/464. RY* (6) C 44	0.06	2.60	0.041
168. LP (6)Cr 63	/469. RY* (3) C 45	0.04	0.66	0.016
168. LP (6)Cr 63	/470. RY* (4) C 45	0.25	0.28	0.029
168. LP (6)Cr 63	/475. RY* (1) C 46	7.19	0.54	0.208
168. LP (6)Cr 63	/477. RY* (3) C 46	1.18	0.89	0.109
168. LP (6)Cr 63	/478. RY* (4) C 46	0.35	0.39	0.040
168. LP (6)Cr 63	/479. RY* (5) C 46	0.18	3.99	0.089
168. LP (6)Cr 63	/480. RY* (6) C 46	0.43	2.04	0.100
168. LP (6)Cr 63	/484. RY* (2) C 47	0.05	0.59	0.019
168. LP (6)Cr 63	/486. RY* (4) C 47	0.03	0.42	0.012
168. LP (6)Cr 63	/492. RY* (2) C 48	0.29	0.68	0.047
168. LP (6)Cr 63	/494. RY* (4) C 48	0.04	0.25	0.011
168. LP (6)Cr 63	/499. RY* (1) C 49	0.03	0.56	0.013
168. LP (6)Cr 63	/522. RY* (2) H 54	0.13	1.70	0.050
168. LP (6)Cr 63	/537. RY* (1) O 62	0.04	0.41	0.013
168. LP (6)Cr 63	/538. RY* (2) O 62	5.66	1.22	0.280
168. LP (6)Cr 63	/539. RY* (3) O 62	6.23	0.98	0.263
168. LP (6)Cr 63	/541. RY* (5) O 62	0.20	4.38	0.100
168. LP (6)Cr 63	/542. RY* (6) O 62	0.03	4.26	0.038
168. LP (6)Cr 63	/544. RY* (8) O 62	0.06	4.23	0.052
168. LP (6)Cr 63	/645. BD* (1) C 40 - C 41	0.11	0.01	0.004
169. LP (7)Cr 63	/444. RY* (2) N 42	3.56	0.39	0.410
169. LP (7)Cr 63	/454. RY* (4) N 43	0.85	0.06	0.087
169. LP (7)Cr 63	/479. RY* (5) C 46	0.05	3.02	0.151
169. LP (7)Cr 63	/480. RY* (6) C 46	0.04	1.07	0.081
169. LP (7)Cr 63	/538. RY* (2) O 62	15.31	0.25	0.750
169. LP (7)Cr 63	/540. RY* (4) O 62	0.16	0.24	0.077
172. LP (1)Cr 63	/537. RY* (1) O 62	0.04	1.33	0.009
172. LP (1)Cr 63	/649. BD* (1) N 42 - N 43	0.06	0.75	0.008
172. LP (1)Cr 63	/657. BD* (1) C 46 - O 62	0.07	0.59	0.008

173. LP (2)Cr 63	/ 56. BD* (2) C 35 - C 37	0.03	0.22	0.005
173. LP (2)Cr 63	/408. RY* (6) C 37	0.03	3.84	0.014
173. LP (2)Cr 63	/446. RY* (4) N 42	0.09	1.73	0.016
173. LP (2)Cr 63	/447. RY* (5) N 42	0.03	4.15	0.013
173. LP (2)Cr 63	/459. RY* (1) C 44	0.03	1.43	0.008
173. LP (2)Cr 63	/539. RY* (3) O 62	0.04	1.89	0.010
173. LP (2)Cr 63	/639. BD* (1) C 37 - N 42	0.31	0.63	0.018
173. LP (2)Cr 63	/649. BD* (1) N 42 - N 43	0.53	0.75	0.025
173. LP (2)Cr 63	/656. BD* (1) C 46 - C 48	0.04	0.71	0.006
173. LP (2)Cr 63	/657. BD* (1) C 46 - O 62	0.10	0.59	0.010
174. LP (3)Cr 63	/443. RY* (1) N 42	0.13	1.14	0.016
174. LP (3)Cr 63	/657. BD* (1) C 46 - O 62	0.11	0.59	0.010

within unit 3

130. CR (2)Cr 63	/168. LP (6)Cr 63	0.80	24.78	0.191
131. CR (3)Cr 63	/168. LP (6)Cr 63	2.90	4.46	0.155
166. LP (4)Cr 63	/167. LP (5)Cr 63	0.52	0.02	0.004
166. LP (4)Cr 63	/551. RY* (7)Cr 63	0.28	2.04	0.047
167. LP (5)Cr 63	/168. LP (6)Cr 63	1.40	0.77	0.061
167. LP (5)Cr 63	/547. RY* (3)Cr 63	0.64	1.83	0.075
167. LP (5)Cr 63	/550. RY* (6)Cr 63	0.28	2.93	0.063
167. LP (5)Cr 63	/561. RY* (17)Cr 63	0.32	37.67	0.243
168. LP (6)Cr 63	/169. LP (7)Cr 63	3.40	0.97	0.188
168. LP (6)Cr 63	/546. RY* (2)Cr 63	9.96	0.94	0.325
168. LP (6)Cr 63	/547. RY* (3)Cr 63	3.93	1.06	0.217
168. LP (6)Cr 63	/548. RY* (4)Cr 63	1.15	0.91	0.108
168. LP (6)Cr 63	/549. RY* (5)Cr 63	5.71	0.75	0.220
168. LP (6)Cr 63	/550. RY* (6)Cr 63	27.44	2.16	0.821
168. LP (6)Cr 63	/551. RY* (7)Cr 63	1.43	1.26	0.143
168. LP (6)Cr 63	/554. RY* (10)Cr 63	0.98	8.44	0.307
168. LP (6)Cr 63	/555. RY* (11)Cr 63	0.42	3.71	0.134
168. LP (6)Cr 63	/558. RY* (14)Cr 63	0.33	3.93	0.122
168. LP (6)Cr 63	/560. RY* (16)Cr 63	3.83	114.25	2.234
168. LP (6)Cr 63	/561. RY* (17)Cr 63	23.31	36.90	3.131
168. LP (6)Cr 63	/564. RY* (20)Cr 63	7.38	673.64	7.526
169. LP (7)Cr 63	/547. RY* (3)Cr 63	4.07	0.09	0.230
169. LP (7)Cr 63	/551. RY* (7)Cr 63	0.36	0.29	0.127
169. LP (7)Cr 63	/552. RY* (8)Cr 63	2.98	7.08	1.853
169. LP (7)Cr 63	/559. RY* (15)Cr 63	0.81	24.08	1.782
169. LP (7)Cr 63	/561. RY* (17)Cr 63	0.32	35.93	1.366

from unit 3 to unit 4

133. CR (5)Cr 63	/665. BD* (1) H 65 - O 66	0.03	2.25	0.010
135. CR (7)Cr 63	/664. BD* (1) H 64 - O 66	0.03	2.25	0.010
137. CR (9)Cr 63	/664. BD* (1) H 64 - O 66	0.03	2.25	0.011
137. CR (9)Cr 63	/665. BD* (1) H 65 - O 66	0.03	2.26	0.011
167. LP (5)Cr 63	/566. RY* (1) H 64	0.04	0.75	0.012
167. LP (5)Cr 63	/568. RY* (1) H 65	0.03	0.86	0.012
167. LP (5)Cr 63	/570. RY* (1) O 66	0.48	2.25	0.072
168. LP (6)Cr 63	/568. RY* (1) H 65	36.37	0.09	0.195
168. LP (6)Cr 63	/569. RY* (2) H 65	0.20	1.44	0.057
168. LP (6)Cr 63	/570. RY* (1) O 66	15.67	1.48	0.514
168. LP (6)Cr 63	/571. RY* (2) O 66	0.05	0.39	0.015
168. LP (6)Cr 63	/572. RY* (3) O 66	1.62	0.66	0.110
168. LP (6)Cr 63	/573. RY* (4) O 66	0.70	0.50	0.063
168. LP (6)Cr 63	/575. RY* (6) O 66	0.05	4.24	0.047
169. LP (7)Cr 63	/570. RY* (1) O 66	0.67	0.51	0.231
172. LP (1)Cr 63	/664. BD* (1) H 64 - O 66	0.22	0.60	0.014
172. LP (1)Cr 63	/665. BD* (1) H 65 - O 66	0.11	0.60	0.010
173. LP (2)Cr 63	/664. BD* (1) H 64 - O 66	0.11	0.60	0.010
174. LP (3)Cr 63	/665. BD* (1) H 65 - O 66	0.19	0.60	0.014

from unit 3 to unit 5

166. LP (4)Cr 63	/579. RY* (2)C1 67	0.05	0.79	0.013
167. LP (5)Cr 63	/581. RY* (4)C1 67	0.15	1.75	0.036
168. LP (6)Cr 63	/579. RY* (2)C1 67	421.69	0.01	0.235
168. LP (6)Cr 63	/580. RY* (3)C1 67	9.57	0.25	0.165
168. LP (6)Cr 63	/581. RY* (4)C1 67	13.41	0.98	0.388
168. LP (6)Cr 63	/583. RY* (6)C1 67	0.34	2.13	0.091

from unit 4 to unit 1

87. BD (1) H 64 - O 66	/262. RY* (2) N 11	0.05	2.72	0.014
88. BD (1) H 65 - O 66	/221. RY* (1) C 6	0.04	1.65	0.010
88. BD (1) H 65 - O 66	/222. RY* (2) C 6	0.06	1.99	0.014
88. BD (1) H 65 - O 66	/224. RY* (4) C 6	0.05	1.83	0.012
175. LP (1) O 66	/ 13. BD* (2) C 4 - C 6	0.07	0.50	0.011
175. LP (1) O 66	/222. RY* (2) C 6	0.04	1.74	0.011
175. LP (1) O 66	/262. RY* (2) N 11	0.03	2.48	0.011
175. LP (1) O 66	/263. RY* (3) N 11	0.04	1.88	0.011
175. LP (1) O 66	/264. RY* (4) N 11	0.03	2.04	0.010
175. LP (1) O 66	/602. BD* (1) C 6 - N 11	0.10	0.91	0.012
175. LP (1) O 66	/609. BD* (2) C 9 - C 10	0.03	0.57	0.006
175. LP (1) O 66	/612. BD* (1) N 11 - N 12	0.07	1.02	0.010
176. LP (2) O 66	/ 13. BD* (2) C 4 - C 6	0.06	0.53	0.010
176. LP (2) O 66	/222. RY* (2) C 6	0.13	1.77	0.020
176. LP (2) O 66	/224. RY* (4) C 6	0.04	1.61	0.011
176. LP (2) O 66	/226. RY* (6) C 6	0.06	3.57	0.020
176. LP (2) O 66	/262. RY* (2) N 11	0.25	2.51	0.033
176. LP (2) O 66	/263. RY* (3) N 11	0.35	1.91	0.034
176. LP (2) O 66	/270. RY* (2) N 12	0.03	1.39	0.009
176. LP (2) O 66	/271. RY* (3) N 12	0.09	1.69	0.016
176. LP (2) O 66	/293. RY* (1) C 15	0.05	1.75	0.012
176. LP (2) O 66	/295. RY* (3) C 15	0.05	2.09	0.013
176. LP (2) O 66	/297. RY* (5) C 15	0.05	4.38	0.020
176. LP (2) O 66	/602. BD* (1) C 6 - N 11	0.07	0.94	0.010
176. LP (2) O 66	/612. BD* (1) N 11 - N 12	0.20	1.05	0.019

from unit 4 to unit 2

87. BD (1) H 64 - O 66	/404. RY* (2) C 37	0.04	1.88	0.011
87. BD (1) H 64 - O 66	/444. RY* (2) N 42	0.14	2.79	0.025
87. BD (1) H 64 - O 66	/445. RY* (3) N 42	0.19	2.11	0.025
87. BD (1) H 64 - O 66	/538. RY* (2) O 62	0.05	2.65	0.015
87. BD (1) H 64 - O 66	/539. RY* (3) O 62	0.05	2.40	0.014
87. BD (1) H 64 - O 66	/646. BD* (2) C 40 - C 41	0.04	0.81	0.009
87. BD (1) H 64 - O 66	/657. BD* (1) C 46 - O 62	0.03	1.10	0.007
88. BD (1) H 65 - O 66	/444. RY* (2) N 42	0.07	2.79	0.018
88. BD (1) H 65 - O 66	/445. RY* (3) N 42	0.08	2.11	0.017
88. BD (1) H 65 - O 66	/646. BD* (2) C 40 - C 41	0.03	0.81	0.007
88. BD (1) H 65 - O 66	/657. BD* (1) C 46 - O 62	0.03	1.10	0.007
175. LP (1) O 66	/404. RY* (2) C 37	0.08	1.63	0.014
175. LP (1) O 66	/427. RY* (1) C 40	0.04	1.37	0.010
175. LP (1) O 66	/445. RY* (3) N 42	0.03	1.86	0.010
176. LP (2) O 66	/404. RY* (2) C 37	0.11	1.66	0.017
176. LP (2) O 66	/408. RY* (6) C 37	0.03	4.13	0.015
176. LP (2) O 66	/443. RY* (1) N 42	0.04	1.43	0.010
176. LP (2) O 66	/444. RY* (2) N 42	0.37	2.57	0.040
176. LP (2) O 66	/445. RY* (3) N 42	0.35	1.89	0.034
176. LP (2) O 66	/451. RY* (1) N 43	0.05	1.39	0.011
176. LP (2) O 66	/453. RY* (3) N 43	0.07	1.65	0.014
176. LP (2) O 66	/475. RY* (1) C 46	0.09	1.75	0.017
176. LP (2) O 66	/476. RY* (2) C 46	0.03	1.19	0.008
176. LP (2) O 66	/537. RY* (1) O 62	0.03	1.62	0.008
176. LP (2) O 66	/538. RY* (2) O 62	0.18	2.43	0.028
176. LP (2) O 66	/539. RY* (3) O 62	0.25	2.19	0.031
176. LP (2) O 66	/540. RY* (4) O 62	0.03	2.42	0.011
176. LP (2) O 66	/639. BD* (1) C 37 - N 42	0.04	0.93	0.008
176. LP (2) O 66	/649. BD* (1) N 42 - N 43	0.11	1.04	0.014
176. LP (2) O 66	/657. BD* (1) C 46 - O 62	0.07	0.88	0.010

from unit 4 to unit 3

87. BD (1) H 64 - O 66	/167. LP (5)Cr 63	0.18	0.66	0.017
87. BD (1) H 64 - O 66	/168. LP (6)Cr 63	3.19	1.43	0.092
87. BD (1) H 64 - O 66	/169. LP (7)Cr 63	0.05	2.40	0.015
87. BD (1) H 64 - O 66	/170. LP* (8)Cr 63	0.07	2.60	0.017
87. BD (1) H 64 - O 66	/545. RY* (1)Cr 63	0.05	1.30	0.010
87. BD (1) H 64 - O 66	/546. RY* (2)Cr 63	0.12	2.36	0.021
87. BD (1) H 64 - O 66	/547. RY* (3)Cr 63	0.14	2.49	0.023
87. BD (1) H 64 - O 66	/548. RY* (4)Cr 63	0.08	2.33	0.017
87. BD (1) H 64 - O 66	/549. RY* (5)Cr 63	0.22	2.17	0.028
87. BD (1) H 64 - O 66	/550. RY* (6)Cr 63	0.50	3.59	0.053
87. BD (1) H 64 - O 66	/554. RY* (10)Cr 63	0.10	9.87	0.040
87. BD (1) H 64 - O 66	/560. RY* (16)Cr 63	0.06	115.68	0.108

87. BD (1) H 64 - O 66	/561. RY* (17)Cr 63	0.39	38.33	0.155
87. BD (1) H 64 - O 66	/564. RY* (20)Cr 63	0.14	675.06	0.394
88. BD (1) H 65 - O 66	/167. LP (5)Cr 63	0.15	0.67	0.015
88. BD (1) H 65 - O 66	/168. LP (6)Cr 63	2.64	1.43	0.084
88. BD (1) H 65 - O 66	/169. LP (7)Cr 63	0.03	2.40	0.011
88. BD (1) H 65 - O 66	/170. LP* (8)Cr 63	0.05	2.60	0.015
88. BD (1) H 65 - O 66	/545. RY* (1)Cr 63	0.04	1.30	0.010
88. BD (1) H 65 - O 66	/546. RY* (2)Cr 63	0.20	2.37	0.027
88. BD (1) H 65 - O 66	/547. RY* (3)Cr 63	0.11	2.49	0.021
88. BD (1) H 65 - O 66	/548. RY* (4)Cr 63	0.04	2.34	0.012
88. BD (1) H 65 - O 66	/550. RY* (6)Cr 63	0.36	3.59	0.045
88. BD (1) H 65 - O 66	/553. RY* (9)Cr 63	0.03	5.22	0.016
88. BD (1) H 65 - O 66	/554. RY* (10)Cr 63	0.08	9.87	0.035
88. BD (1) H 65 - O 66	/560. RY* (16)Cr 63	0.03	115.68	0.077
88. BD (1) H 65 - O 66	/561. RY* (17)Cr 63	0.20	38.33	0.110
88. BD (1) H 65 - O 66	/564. RY* (20)Cr 63	0.08	675.07	0.287
138. CR (1) O 66	/167. LP (5)Cr 63	0.21	18.86	0.096
138. CR (1) O 66	/168. LP (6)Cr 63	1.80	19.62	0.256
175. LP (1) O 66	/167. LP (5)Cr 63	0.40	0.42	0.020
175. LP (1) O 66	/168. LP (6)Cr 63	2.87	1.18	0.079
175. LP (1) O 66	/170. LP* (8)Cr 63	0.03	2.35	0.011
175. LP (1) O 66	/171. LP* (9)Cr 63	0.29	2.35	0.033
175. LP (1) O 66	/545. RY* (1)Cr 63	0.09	1.05	0.012
175. LP (1) O 66	/546. RY* (2)Cr 63	0.15	2.12	0.022
175. LP (1) O 66	/548. RY* (4)Cr 63	0.26	2.09	0.030
175. LP (1) O 66	/550. RY* (6)Cr 63	0.19	3.35	0.032
175. LP (1) O 66	/551. RY* (7)Cr 63	0.06	2.44	0.015
175. LP (1) O 66	/554. RY* (10)Cr 63	0.07	9.63	0.033
175. LP (1) O 66	/560. RY* (16)Cr 63	0.03	115.43	0.074
175. LP (1) O 66	/561. RY* (17)Cr 63	0.19	38.08	0.107
175. LP (1) O 66	/564. RY* (20)Cr 63	0.07	674.82	0.284
176. LP (2) O 66	/166. LP (4)Cr 63	0.32	0.43	0.018
176. LP (2) O 66	/167. LP (5)Cr 63	14.43	0.45	0.116
176. LP (2) O 66	/168. LP (6)Cr 63	23.02	1.21	0.216
176. LP (2) O 66	/169. LP (7)Cr 63	0.20	2.18	0.028
176. LP (2) O 66	/170. LP* (8)Cr 63	2.56	2.38	0.103
176. LP (2) O 66	/171. LP* (9)Cr 63	1.29	2.38	0.073
176. LP (2) O 66	/545. RY* (1)Cr 63	0.30	1.08	0.024
176. LP (2) O 66	/546. RY* (2)Cr 63	0.53	2.15	0.045
176. LP (2) O 66	/548. RY* (4)Cr 63	0.28	2.12	0.032
176. LP (2) O 66	/549. RY* (5)Cr 63	0.06	1.96	0.015
176. LP (2) O 66	/550. RY* (6)Cr 63	1.87	3.37	0.105
176. LP (2) O 66	/551. RY* (7)Cr 63	0.28	2.47	0.035
176. LP (2) O 66	/552. RY* (8)Cr 63	0.06	9.26	0.030
176. LP (2) O 66	/553. RY* (9)Cr 63	0.03	5.00	0.016
176. LP (2) O 66	/554. RY* (10)Cr 63	1.41	9.65	0.154
176. LP (2) O 66	/557. RY* (13)Cr 63	0.03	4.99	0.016
176. LP (2) O 66	/558. RY* (14)Cr 63	0.08	5.14	0.027
176. LP (2) O 66	/560. RY* (16)Cr 63	0.34	115.46	0.262
176. LP (2) O 66	/561. RY* (17)Cr 63	1.88	38.11	0.354
176. LP (2) O 66	/563. RY* (19)Cr 63	0.25	26.22	0.106
176. LP (2) O 66	/564. RY* (20)Cr 63	0.65	674.85	0.877
within unit 4				
87. BD (1) H 64 - O 66	/569. RY* (2) H 65	0.27	2.87	0.035
138. CR (1) O 66	/567. RY* (2) H 64	0.37	21.15	0.112
138. CR (1) O 66	/569. RY* (2) H 65	0.29	21.06	0.099
175. LP (1) O 66	/567. RY* (2) H 64	0.31	2.71	0.037
175. LP (1) O 66	/569. RY* (2) H 65	0.38	2.62	0.040
176. LP (2) O 66	/566. RY* (1) H 64	0.57	1.19	0.035
176. LP (2) O 66	/568. RY* (1) H 65	0.47	1.30	0.033
176. LP (2) O 66	/569. RY* (2) H 65	0.33	2.65	0.039
176. LP (2) O 66	/570. RY* (1) O 66	1.57	2.69	0.086
from unit 4 to unit 5				
87. BD (1) H 64 - O 66	/579. RY* (2)C1 67	0.03	1.44	0.008
87. BD (1) H 64 - O 66	/581. RY* (4)C1 67	0.08	2.41	0.017
88. BD (1) H 65 - O 66	/579. RY* (2)C1 67	0.05	1.44	0.011
175. LP (1) O 66	/579. RY* (2)C1 67	0.03	1.19	0.008
175. LP (1) O 66	/581. RY* (4)C1 67	0.03	2.16	0.010
176. LP (2) O 66	/579. RY* (2)C1 67	0.46	1.22	0.031

176. LP (2) O 66	/580. RY* (3)C1 67	0.10	1.46	0.016
176. LP (2) O 66	/581. RY* (4)C1 67	0.36	2.19	0.037
from unit 5 to unit 1				
177. LP (1)C1 67	/347. RY* (1) H 24	0.05	0.82	0.008
177. LP (1)C1 67	/609. BD* (2) C 9 - C 10	0.05	0.31	0.006
177. LP (1)C1 67	/610. BD* (1) C 9 - H 24	0.22	0.75	0.016
177. LP (1)C1 67	/616. BD* (2) C 13 - C 15	0.15	0.29	0.009
177. LP (1)C1 67	/620. BD* (1) C 15 - O 19	0.11	0.60	0.010
178. LP (2)C1 67	/602. BD* (1) C 6 - N 11	0.05	0.65	0.007
178. LP (2)C1 67	/610. BD* (1) C 9 - H 24	0.12	0.75	0.012
178. LP (2)C1 67	/613. BD* (1) N 12 - C 13	0.07	0.64	0.008
178. LP (2)C1 67	/616. BD* (2) C 13 - C 15	0.03	0.29	0.004
178. LP (2)C1 67	/620. BD* (1) C 15 - O 19	0.07	0.60	0.008
179. LP (3)C1 67	/222. RY* (2) C 6	0.04	2.05	0.012
179. LP (3)C1 67	/224. RY* (4) C 6	0.03	1.89	0.009
179. LP (3)C1 67	/262. RY* (2) N 11	0.22	2.79	0.032
179. LP (3)C1 67	/263. RY* (3) N 11	0.10	2.19	0.019
179. LP (3)C1 67	/293. RY* (1) C 15	0.06	2.03	0.013
179. LP (3)C1 67	/326. RY* (2) O 19	0.06	2.51	0.016
179. LP (3)C1 67	/327. RY* (3) O 19	0.08	2.71	0.019
179. LP (3)C1 67	/610. BD* (1) C 9 - H 24	0.05	1.32	0.010
180. LP (4)C1 67	/221. RY* (1) C 6	0.03	1.20	0.009
180. LP (4)C1 67	/222. RY* (2) C 6	0.23	1.55	0.025
180. LP (4)C1 67	/223. RY* (3) C 6	0.06	1.46	0.013
180. LP (4)C1 67	/224. RY* (4) C 6	0.14	1.38	0.019
180. LP (4)C1 67	/226. RY* (6) C 6	0.06	3.35	0.019
180. LP (4)C1 67	/262. RY* (2) N 11	1.30	2.28	0.074
180. LP (4)C1 67	/263. RY* (3) N 11	0.41	1.69	0.036
180. LP (4)C1 67	/270. RY* (2) N 12	0.03	1.17	0.008
180. LP (4)C1 67	/271. RY* (3) N 12	0.03	1.47	0.009
180. LP (4)C1 67	/272. RY* (4) N 12	0.04	2.01	0.013
180. LP (4)C1 67	/293. RY* (1) C 15	0.32	1.53	0.030
180. LP (4)C1 67	/294. RY* (2) C 15	0.07	0.97	0.011
180. LP (4)C1 67	/295. RY* (3) C 15	0.11	1.86	0.019
180. LP (4)C1 67	/297. RY* (5) C 15	0.10	4.16	0.028
180. LP (4)C1 67	/310. RY* (2) C 17	0.03	1.67	0.010
180. LP (4)C1 67	/325. RY* (1) O 19	0.04	1.44	0.010
180. LP (4)C1 67	/326. RY* (2) O 19	0.33	2.01	0.035
180. LP (4)C1 67	/327. RY* (3) O 19	0.44	2.21	0.042
180. LP (4)C1 67	/347. RY* (1) H 24	0.03	0.89	0.007
180. LP (4)C1 67	/602. BD* (1) C 6 - N 11	0.15	0.71	0.014
180. LP (4)C1 67	/620. BD* (1) C 15 - O 19	0.13	0.66	0.013
from unit 5 to unit 2				
177. LP (1)C1 67	/ 52. BD* (2) C 34 - C 36	0.19	0.22	0.013
177. LP (1)C1 67	/633. BD* (1) C 34 - C 36	0.03	0.73	0.006
178. LP (2)C1 67	/ 52. BD* (2) C 34 - C 36	0.11	0.22	0.010
178. LP (2)C1 67	/ 56. BD* (2) C 35 - C 37	0.16	0.23	0.011
178. LP (2)C1 67	/639. BD* (1) C 37 - N 42	0.04	0.64	0.006
178. LP (2)C1 67	/649. BD* (1) N 42 - N 43	0.05	0.75	0.007
179. LP (3)C1 67	/ 52. BD* (2) C 34 - C 36	0.06	0.79	0.013
179. LP (3)C1 67	/ 56. BD* (2) C 35 - C 37	0.04	0.80	0.010
179. LP (3)C1 67	/403. RY* (1) C 37	0.03	1.69	0.009
179. LP (3)C1 67	/404. RY* (2) C 37	0.04	1.94	0.011
179. LP (3)C1 67	/444. RY* (2) N 42	0.20	2.85	0.030
179. LP (3)C1 67	/445. RY* (3) N 42	0.13	2.17	0.022
179. LP (3)C1 67	/453. RY* (3) N 43	0.03	1.93	0.009
179. LP (3)C1 67	/475. RY* (1) C 46	0.06	2.03	0.014
179. LP (3)C1 67	/538. RY* (2) O 62	0.06	2.71	0.016
179. LP (3)C1 67	/539. RY* (3) O 62	0.06	2.47	0.016
180. LP (4)C1 67	/ 56. BD* (2) C 35 - C 37	0.12	0.30	0.010
180. LP (4)C1 67	/403. RY* (1) C 37	0.04	1.18	0.010
180. LP (4)C1 67	/404. RY* (2) C 37	0.32	1.44	0.029
180. LP (4)C1 67	/408. RY* (6) C 37	0.07	3.91	0.023
180. LP (4)C1 67	/444. RY* (2) N 42	0.92	2.35	0.063
180. LP (4)C1 67	/445. RY* (3) N 42	0.75	1.67	0.048
180. LP (4)C1 67	/451. RY* (1) N 43	0.04	1.17	0.009
180. LP (4)C1 67	/452. RY* (2) N 43	0.05	1.16	0.010
180. LP (4)C1 67	/453. RY* (3) N 43	0.16	1.43	0.021
180. LP (4)C1 67	/475. RY* (1) C 46	0.23	1.53	0.026

180. LP (4)C1 67	/476. RY* (2) C 46	0.06	0.97	0.010
180. LP (4)C1 67	/477. RY* (3) C 46	0.09	1.88	0.018
180. LP (4)C1 67	/480. RY* (6) C 46	0.05	3.03	0.017
180. LP (4)C1 67	/538. RY* (2) O 62	0.10	2.21	0.020
180. LP (4)C1 67	/539. RY* (3) O 62	0.22	1.96	0.028
180. LP (4)C1 67	/639. BD* (1) C 37 - N 42	0.07	0.71	0.010
180. LP (4)C1 67	/649. BD* (1) N 42 - N 43	0.21	0.82	0.017

from unit 5 to unit 3

139. CR (1)C1 67	/168. LP (6)Cr 63	0.25	100.55	0.217
140. CR (2)C1 67	/167. LP (5)Cr 63	0.10	10.62	0.049
140. CR (2)C1 67	/168. LP (6)Cr 63	2.04	11.39	0.208
143. CR (5)C1 67	/167. LP (5)Cr 63	0.04	7.02	0.024
143. CR (5)C1 67	/168. LP (6)Cr 63	0.22	7.79	0.056
177. LP (1)C1 67	/168. LP (6)Cr 63	0.03	0.92	0.007
177. LP (1)C1 67	/171. LP* (9)Cr 63	0.10	2.09	0.018
177. LP (1)C1 67	/546. RY* (2)Cr 63	0.12	1.86	0.019
177. LP (1)C1 67	/549. RY* (5)Cr 63	0.10	1.67	0.016
178. LP (2)C1 67	/168. LP (6)Cr 63	0.03	0.92	0.008
178. LP (2)C1 67	/170. LP* (8)Cr 63	0.12	2.10	0.020
178. LP (2)C1 67	/546. RY* (2)Cr 63	0.07	1.86	0.015
178. LP (2)C1 67	/549. RY* (5)Cr 63	0.09	1.67	0.016
179. LP (3)C1 67	/166. LP (4)Cr 63	0.26	0.71	0.022
179. LP (3)C1 67	/167. LP (5)Cr 63	0.79	0.73	0.037
179. LP (3)C1 67	/168. LP (6)Cr 63	7.19	1.49	0.140
179. LP (3)C1 67	/545. RY* (1)Cr 63	0.14	1.36	0.018
179. LP (3)C1 67	/546. RY* (2)Cr 63	0.25	2.43	0.031
179. LP (3)C1 67	/547. RY* (3)Cr 63	0.22	2.55	0.030
179. LP (3)C1 67	/549. RY* (5)Cr 63	0.19	2.24	0.026
179. LP (3)C1 67	/550. RY* (6)Cr 63	0.89	3.65	0.072
179. LP (3)C1 67	/551. RY* (7)Cr 63	0.07	2.75	0.018
179. LP (3)C1 67	/554. RY* (10)Cr 63	0.05	9.93	0.029
179. LP (3)C1 67	/560. RY* (16)Cr 63	0.13	115.74	0.157
179. LP (3)C1 67	/561. RY* (17)Cr 63	0.86	38.39	0.230
179. LP (3)C1 67	/564. RY* (20)Cr 63	0.32	675.13	0.586
180. LP (4)C1 67	/166. LP (4)Cr 63	10.73	0.20	0.069
180. LP (4)C1 67	/167. LP (5)Cr 63	16.41	0.22	0.085
180. LP (4)C1 67	/168. LP (6)Cr 63	29.99	0.99	0.218
180. LP (4)C1 67	/169. LP (7)Cr 63	1.99	1.96	0.084
180. LP (4)C1 67	/545. RY* (1)Cr 63	0.93	0.86	0.038
180. LP (4)C1 67	/546. RY* (2)Cr 63	0.82	1.92	0.054
180. LP (4)C1 67	/547. RY* (3)Cr 63	1.32	2.05	0.070
180. LP (4)C1 67	/549. RY* (5)Cr 63	0.90	1.73	0.053
180. LP (4)C1 67	/550. RY* (6)Cr 63	3.21	3.15	0.136
180. LP (4)C1 67	/551. RY* (7)Cr 63	0.42	2.25	0.042
180. LP (4)C1 67	/552. RY* (8)Cr 63	0.48	9.04	0.089
180. LP (4)C1 67	/554. RY* (10)Cr 63	0.31	9.43	0.073
180. LP (4)C1 67	/555. RY* (11)Cr 63	0.05	4.70	0.021
180. LP (4)C1 67	/556. RY* (12)Cr 63	0.03	8.38	0.021
180. LP (4)C1 67	/559. RY* (15)Cr 63	0.14	26.04	0.081
180. LP (4)C1 67	/560. RY* (16)Cr 63	0.56	115.24	0.343
180. LP (4)C1 67	/561. RY* (17)Cr 63	3.29	37.89	0.477
180. LP (4)C1 67	/562. RY* (18)Cr 63	0.03	4.85	0.017
180. LP (4)C1 67	/564. RY* (20)Cr 63	1.06	674.62	1.144

from unit 5 to unit 4

177. LP (1)C1 67	/664. BD* (1) H 64 - O 66	0.10	0.60	0.010
177. LP (1)C1 67	/665. BD* (1) H 65 - O 66	0.05	0.60	0.007
178. LP (2)C1 67	/664. BD* (1) H 64 - O 66	0.03	0.60	0.006
178. LP (2)C1 67	/665. BD* (1) H 65 - O 66	0.14	0.61	0.012
179. LP (3)C1 67	/566. RY* (1) H 64	0.05	1.47	0.011
179. LP (3)C1 67	/568. RY* (1) H 65	0.05	1.58	0.012
179. LP (3)C1 67	/570. RY* (1) O 66	0.23	2.97	0.033
180. LP (4)C1 67	/566. RY* (1) H 64	0.32	0.97	0.024
180. LP (4)C1 67	/568. RY* (1) H 65	0.42	1.08	0.029
180. LP (4)C1 67	/570. RY* (1) O 66	1.40	2.47	0.079
180. LP (4)C1 67	/572. RY* (3) O 66	0.11	1.65	0.018

within unit 5

180. LP (4)C1 67	/579. RY* (2)C1 67	0.63	1.00	0.034
180. LP (4)C1 67	/580. RY* (3)C1 67	0.43	1.24	0.031

180. LP (4)C1 67

/581. RY* (4)C1 67

1.43 1.97 0.072

Table (S8): Summary of Natural Population Analysis of the Mn(II) complex:

Atom	No	Natural Charge	Core	Valence	Rydberg	Total
C	1	0.31891	1.99861	3.66363	0.01885	5.68109
C	2	-0.09395	1.99882	4.08293	0.01219	6.09395
C	3	-0.21623	1.99901	4.20116	0.01606	6.21623
C	4	-0.06428	1.99883	4.05142	0.01403	6.06428
C	5	-0.21755	1.99893	4.20188	0.01674	6.21755
C	6	0.17364	1.99870	3.80431	0.02335	5.82636
C	7	-0.23549	1.99903	4.21965	0.01681	6.23549
C	8	-0.20422	1.99909	4.18991	0.01523	6.20422
C	9	-0.20385	1.99902	4.18858	0.01625	6.20385
C	10	-0.22308	1.99906	4.20948	0.01454	6.22308
N	11	-0.41509	1.99936	5.38613	0.02961	7.41509
N	12	-0.04994	1.99918	5.03041	0.02035	7.04994
C	13	0.02482	1.99873	3.95709	0.01936	5.97518
C	14	-0.28139	1.99905	4.25695	0.02539	6.28139
C	15	0.23846	1.99869	3.73920	0.02366	5.76154
C	16	-0.18251	1.99910	4.16343	0.01999	6.18251
C	17	-0.26125	1.99901	4.23864	0.02360	6.26125
C	18	-0.23377	1.99912	4.21917	0.01548	6.23377
O	19	-0.72266	1.99984	6.71658	0.00624	8.72266
O	20	-0.65330	1.99979	6.64968	0.00383	8.65330
H	21	0.22757	0.00000	0.76998	0.00245	0.77243
H	22	0.20033	0.00000	0.79672	0.00295	0.79967
H	23	0.20956	0.00000	0.78923	0.00121	0.79044
H	24	0.22107	0.00000	0.77528	0.00366	0.77893
H	25	0.20962	0.00000	0.78860	0.00178	0.79038
H	26	0.20443	0.00000	0.79310	0.00246	0.79557
H	27	0.20298	0.00000	0.79579	0.00123	0.79702
H	28	0.20880	0.00000	0.78850	0.00269	0.79120
H	29	0.20865	0.00000	0.79003	0.00132	0.79135
H	30	0.49754	0.00000	0.50129	0.00117	0.50246
H	31	0.22334	0.00000	0.77375	0.00291	0.77666
C	32	0.31251	1.99867	3.66979	0.01904	5.68749
C	33	-0.08378	1.99884	4.06779	0.01715	6.08378
C	34	-0.18629	1.99902	4.16831	0.01896	6.18629
C	35	-0.06154	1.99886	4.04373	0.01895	6.06154
C	36	-0.21584	1.99893	4.19601	0.02090	6.21584
C	37	0.14231	1.99878	3.83488	0.02402	5.85769
C	38	-0.25969	1.99907	4.24363	0.01699	6.25969
C	39	-0.17053	1.99909	4.15446	0.01697	6.17053
C	40	-0.27017	1.99906	4.25280	0.01831	6.27017
C	41	-0.19760	1.99907	4.18131	0.01722	6.19760
N	42	-0.43173	1.99936	5.40230	0.03008	7.43173
N	43	-0.14896	1.99924	5.12664	0.02308	7.14896
C	44	0.08992	1.99871	3.89183	0.01954	5.91008
C	45	-0.20943	1.99902	4.19640	0.01401	6.20943
C	46	0.31195	1.99867	3.66863	0.02076	5.68805
C	47	-0.18440	1.99908	4.17204	0.01328	6.18440
C	48	-0.22409	1.99899	4.21072	0.01438	6.22409
C	49	-0.20839	1.99908	4.19596	0.01334	6.20839
O	50	-0.64897	1.99979	6.64533	0.00384	8.64897
H	51	0.21973	0.00000	0.77772	0.00255	0.78027
H	52	0.20273	0.00000	0.79443	0.00284	0.79727
H	53	0.20437	0.00000	0.79387	0.00176	0.79563
H	54	0.21216	0.00000	0.78386	0.00399	0.78784
H	55	0.20450	0.00000	0.79370	0.00180	0.79550
H	56	0.22292	0.00000	0.77440	0.00269	0.77708
H	57	0.21616	0.00000	0.78266	0.00119	0.78384
H	58	0.22426	0.00000	0.77286	0.00288	0.77574
H	59	0.21350	0.00000	0.78517	0.00133	0.78650
H	60	0.49887	0.00000	0.49985	0.00127	0.50113
H	61	0.21166	0.00000	0.78425	0.00408	0.78834
O	62	-0.94009	1.99981	6.93369	0.00659	8.94009

H	63	0.50216	0.00000	0.49389	0.00395	0.49784
H	64	0.50612	0.00000	0.48934	0.00454	0.49388
O	65	-0.76144	1.99984	6.75617	0.00543	8.76144
H	66	0.50523	0.00000	0.49067	0.00410	0.49477
H	67	0.50059	0.00000	0.49554	0.00388	0.49941
O	68	-0.92844	1.99981	6.92157	0.00706	8.92844
Mn	69	1.51852	17.99029	5.47740	0.01379	23.48148

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* Total *      0.00000    101.95204    216.25606    0.79190    319.00000
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Table (S9): NBO data of the Mn(II) complex:

Donor NBO (i)		Acceptor NBO (j)		kcal/mol	a.u.	a.u.
within unit 1						
1. BD (1) C	1 - C 2	/190. RY* (2) C	3	0.61	1.94	0.043
1. BD (1) C	1 - C 2	/191. RY* (3) C	3	0.29	1.96	0.031
1. BD (1) C	1 - C 2	/197. RY* (1) C	4	0.68	2.20	0.049
1. BD (1) C	1 - C 2	/222. RY* (2) C	7	0.52	1.77	0.038
1. BD (1) C	1 - C 2	/584. BD* (1) C	1 - C 3	1.65	1.26	0.058
1. BD (1) C	1 - C 2	/586. BD* (1) C	2 - C 4	1.55	1.28	0.056
1. BD (1) C	1 - C 2	/587. BD* (1) C	2 - C 7	0.70	1.10	0.035
1. BD (1) C	1 - C 2	/590. BD* (1) C	3 - H 21	1.48	1.08	0.051
1. BD (1) C	1 - C 2	/593. BD* (1) C	4 - C 9	2.05	1.11	0.060
1. BD (1) C	1 - C 2	/597. BD* (1) C	7 - C 8	0.64	1.30	0.037
2. BD (2) C	1 - C 2	/189. RY* (1) C	3	1.71	0.77	0.049
2. BD (2) C	1 - C 2	/325. RY* (1) O	20	0.55	1.13	0.034
2. BD (2) C	1 - C 2	/583. BD* (2) C	1 - C 2	1.03	0.24	0.023
2. BD (2) C	1 - C 2	/589. BD* (2) C	3 - C 5	18.21	0.25	0.090
2. BD (2) C	1 - C 2	/592. BD* (2) C	4 - C 6	7.16	0.25	0.061
2. BD (2) C	1 - C 2	/598. BD* (2) C	7 - C 8	2.30	0.29	0.035
3. BD (1) C	1 - C 3	/182. RY* (2) C	2	0.81	2.21	0.054
3. BD (1) C	1 - C 3	/207. RY* (3) C	5	0.69	1.96	0.047
3. BD (1) C	1 - C 3	/582. BD* (1) C	1 - C 2	1.60	1.19	0.055
3. BD (1) C	1 - C 3	/585. BD* (1) C	1 - O 20	0.32	1.04	0.023
3. BD (1) C	1 - C 3	/587. BD* (1) C	2 - C 7	1.20	1.22	0.048
3. BD (1) C	1 - C 3	/588. BD* (1) C	3 - C 5	1.01	1.23	0.045
3. BD (1) C	1 - C 3	/590. BD* (1) C	3 - H 21	0.71	1.20	0.037
3. BD (1) C	1 - C 3	/595. BD* (1) C	5 - H 31	0.69	1.21	0.037
3. BD (1) C	1 - C 3	/620. BD* (1) O	20 - H 30	0.75	1.08	0.036
4. BD (1) C	1 - O 20	/173. RY* (1) C	1	0.25	1.87	0.028
4. BD (1) C	1 - O 20	/182. RY* (2) C	2	0.27	2.34	0.032
4. BD (1) C	1 - O 20	/190. RY* (2) C	3	0.30	2.19	0.032
4. BD (1) C	1 - O 20	/584. BD* (1) C	1 - C 3	0.41	1.51	0.032
4. BD (1) C	1 - O 20	/586. BD* (1) C	2 - C 4	0.57	1.53	0.037
4. BD (1) C	1 - O 20	/588. BD* (1) C	3 - C 5	1.08	1.37	0.049
5. BD (1) C	2 - C 4	/175. RY* (3) C	1	0.40	2.18	0.037
5. BD (1) C	2 - C 4	/214. RY* (2) C	6	0.40	1.89	0.035
5. BD (1) C	2 - C 4	/222. RY* (2) C	7	0.30	1.85	0.030
5. BD (1) C	2 - C 4	/238. RY* (2) C	9	0.30	1.82	0.030
5. BD (1) C	2 - C 4	/582. BD* (1) C	1 - C 2	1.52	1.15	0.053
5. BD (1) C	2 - C 4	/585. BD* (1) C	1 - O 20	1.32	1.00	0.046
5. BD (1) C	2 - C 4	/587. BD* (1) C	2 - C 7	1.92	1.18	0.060
5. BD (1) C	2 - C 4	/591. BD* (1) C	4 - C 6	1.88	1.18	0.060
5. BD (1) C	2 - C 4	/593. BD* (1) C	4 - C 9	1.88	1.20	0.060
5. BD (1) C	2 - C 4	/596. BD* (1) C	6 - N 11	1.33	1.10	0.049
5. BD (1) C	2 - C 4	/599. BD* (1) C	7 - H 22	0.75	1.18	0.038
5. BD (1) C	2 - C 4	/604. BD* (1) C	9 - H 24	0.67	1.19	0.036
6. BD (1) C	2 - C 7	/173. RY* (1) C	1	0.28	1.60	0.027
6. BD (1) C	2 - C 7	/175. RY* (3) C	1	0.34	2.08	0.034
6. BD (1) C	2 - C 7	/197. RY* (1) C	4	0.44	2.18	0.039
6. BD (1) C	2 - C 7	/229. RY* (1) C	8	0.34	1.82	0.032
6. BD (1) C	2 - C 7	/231. RY* (3) C	8	0.78	1.93	0.049
6. BD (1) C	2 - C 7	/582. BD* (1) C	1 - C 2	0.57	1.05	0.031
6. BD (1) C	2 - C 7	/584. BD* (1) C	1 - C 3	0.70	1.24	0.037
6. BD (1) C	2 - C 7	/586. BD* (1) C	2 - C 4	2.19	1.26	0.066
6. BD (1) C	2 - C 7	/591. BD* (1) C	4 - C 6	2.42	1.09	0.065
6. BD (1) C	2 - C 7	/597. BD* (1) C	7 - C 8	1.38	1.28	0.053
6. BD (1) C	2 - C 7	/599. BD* (1) C	7 - H 22	0.30	1.08	0.023
6. BD (1) C	2 - C 7	/601. BD* (1) C	8 - H 23	1.48	1.09	0.051
7. BD (1) C	3 - C 5	/173. RY* (1) C	1	0.48	1.62	0.035
7. BD (1) C	3 - C 5	/174. RY* (2) C	1	0.34	1.38	0.027
7. BD (1) C	3 - C 5	/175. RY* (3) C	1	0.31	2.09	0.032
7. BD (1) C	3 - C 5	/215. RY* (3) C	6	0.92	1.76	0.051
7. BD (1) C	3 - C 5	/584. BD* (1) C	1 - C 3	1.09	1.26	0.047
7. BD (1) C	3 - C 5	/585. BD* (1) C	1 - O 20	2.57	0.92	0.061
7. BD (1) C	3 - C 5	/590. BD* (1) C	3 - H 21	0.33	1.08	0.024

7. BD (1) C 3 - C 5	/594. BD* (1) C 5 - C 6	1.34	1.28	0.052
7. BD (1) C 3 - C 5	/595. BD* (1) C 5 - H 31	0.28	1.09	0.022
7. BD (1) C 3 - C 5	/596. BD* (1) C 6 - N 11	2.85	1.02	0.068
8. BD (2) C 3 - C 5	/174. RY* (2) C 1	0.36	0.99	0.025
8. BD (2) C 3 - C 5	/213. RY* (1) C 6	0.34	1.15	0.026
8. BD (2) C 3 - C 5	/583. BD* (2) C 1 - C 2	4.82	0.24	0.052
8. BD (2) C 3 - C 5	/592. BD* (2) C 4 - C 6	3.29	0.25	0.044
9. BD (1) C 3 - H 21	/175. RY* (3) C 1	0.46	1.96	0.038
9. BD (1) C 3 - H 21	/178. RY* (6) C 1	0.28	8.67	0.063
9. BD (1) C 3 - H 21	/206. RY* (2) C 5	0.52	1.77	0.039
9. BD (1) C 3 - H 21	/582. BD* (1) C 1 - C 2	2.95	0.94	0.067
9. BD (1) C 3 - H 21	/584. BD* (1) C 1 - C 3	0.47	1.13	0.029
9. BD (1) C 3 - H 21	/585. BD* (1) C 1 - O 20	0.89	0.79	0.033
9. BD (1) C 3 - H 21	/594. BD* (1) C 5 - C 6	1.47	1.15	0.052
9. BD (1) C 3 - H 21	/595. BD* (1) C 5 - H 31	0.27	0.96	0.020
10. BD (1) C 4 - C 6	/181. RY* (1) C 2	0.71	2.31	0.051
10. BD (1) C 4 - C 6	/206. RY* (2) C 5	0.53	1.90	0.040
10. BD (1) C 4 - C 6	/238. RY* (2) C 9	0.51	1.74	0.038
10. BD (1) C 4 - C 6	/254. RY* (2) N 11	0.25	2.05	0.029
10. BD (1) C 4 - C 6	/586. BD* (1) C 2 - C 4	1.85	1.28	0.061
10. BD (1) C 4 - C 6	/587. BD* (1) C 2 - C 7	1.96	1.10	0.058
10. BD (1) C 4 - C 6	/593. BD* (1) C 4 - C 9	0.90	1.11	0.040
10. BD (1) C 4 - C 6	/594. BD* (1) C 5 - C 6	1.76	1.28	0.060
10. BD (1) C 4 - C 6	/595. BD* (1) C 5 - H 31	1.59	1.09	0.053
10. BD (1) C 4 - C 6	/596. BD* (1) C 6 - N 11	0.30	1.02	0.022
10. BD (1) C 4 - C 6	/602. BD* (1) C 9 - C 10	0.66	1.31	0.037
11. BD (2) C 4 - C 6	/205. RY* (1) C 5	1.88	0.80	0.053
11. BD (2) C 4 - C 6	/583. BD* (2) C 1 - C 2	6.25	0.23	0.054
11. BD (2) C 4 - C 6	/589. BD* (2) C 3 - C 5	23.68	0.24	0.099
11. BD (2) C 4 - C 6	/592. BD* (2) C 4 - C 6	2.19	0.24	0.033
11. BD (2) C 4 - C 6	/603. BD* (2) C 9 - C 10	2.15	0.29	0.033
11. BD (2) C 4 - C 6	/606. BD* (1) N 11 - N 12	3.34	0.74	0.068
12. BD (1) C 4 - C 9	/181. RY* (1) C 2	0.35	2.29	0.036
12. BD (1) C 4 - C 9	/213. RY* (1) C 6	0.46	1.51	0.034
12. BD (1) C 4 - C 9	/214. RY* (2) C 6	0.65	1.79	0.043
12. BD (1) C 4 - C 9	/245. RY* (1) C 10	0.31	1.91	0.031
12. BD (1) C 4 - C 9	/247. RY* (3) C 10	0.79	1.94	0.050
12. BD (1) C 4 - C 9	/582. BD* (1) C 1 - C 2	2.35	1.05	0.063
12. BD (1) C 4 - C 9	/586. BD* (1) C 2 - C 4	2.18	1.26	0.066
12. BD (1) C 4 - C 9	/591. BD* (1) C 4 - C 6	0.90	1.08	0.040
12. BD (1) C 4 - C 9	/594. BD* (1) C 5 - C 6	0.96	1.26	0.044
12. BD (1) C 4 - C 9	/602. BD* (1) C 9 - C 10	1.39	1.29	0.054
12. BD (1) C 4 - C 9	/604. BD* (1) C 9 - H 24	0.34	1.09	0.024
12. BD (1) C 4 - C 9	/605. BD* (1) C 10 - H 25	1.58	1.08	0.052
13. BD (1) C 5 - C 6	/191. RY* (3) C 3	0.71	2.08	0.049
13. BD (1) C 5 - C 6	/198. RY* (2) C 4	0.79	2.30	0.054
13. BD (1) C 5 - C 6	/253. RY* (1) N 11	0.43	2.12	0.038
13. BD (1) C 5 - C 6	/588. BD* (1) C 3 - C 5	1.22	1.23	0.049
13. BD (1) C 5 - C 6	/590. BD* (1) C 3 - H 21	0.62	1.20	0.035
13. BD (1) C 5 - C 6	/591. BD* (1) C 4 - C 6	1.80	1.22	0.060
13. BD (1) C 5 - C 6	/593. BD* (1) C 4 - C 9	1.17	1.23	0.048
13. BD (1) C 5 - C 6	/595. BD* (1) C 5 - H 31	0.85	1.21	0.040
13. BD (1) C 5 - C 6	/596. BD* (1) C 6 - N 11	0.76	1.14	0.037
13. BD (1) C 5 - C 6	/606. BD* (1) N 11 - N 12	0.51	1.25	0.032
14. BD (1) C 5 - H 31	/190. RY* (2) C 3	0.50	1.80	0.038
14. BD (1) C 5 - H 31	/213. RY* (1) C 6	0.40	1.40	0.030
14. BD (1) C 5 - H 31	/214. RY* (2) C 6	0.73	1.68	0.044
14. BD (1) C 5 - H 31	/584. BD* (1) C 1 - C 3	1.37	1.13	0.050
14. BD (1) C 5 - H 31	/590. BD* (1) C 3 - H 21	0.26	0.95	0.020
14. BD (1) C 5 - H 31	/591. BD* (1) C 4 - C 6	3.18	0.97	0.070
14. BD (1) C 5 - H 31	/594. BD* (1) C 5 - C 6	0.72	1.15	0.036
14. BD (1) C 5 - H 31	/596. BD* (1) C 6 - N 11	0.66	0.89	0.031
15. BD (1) C 6 - N 11	/198. RY* (2) C 4	0.69	2.33	0.051
15. BD (1) C 6 - N 11	/206. RY* (2) C 5	0.61	2.04	0.045
15. BD (1) C 6 - N 11	/261. RY* (1) N 12	0.83	1.62	0.047
15. BD (1) C 6 - N 11	/262. RY* (2) N 12	0.25	1.58	0.025
15. BD (1) C 6 - N 11	/586. BD* (1) C 2 - C 4	0.77	1.42	0.042
15. BD (1) C 6 - N 11	/588. BD* (1) C 3 - C 5	1.32	1.25	0.051
15. BD (1) C 6 - N 11	/591. BD* (1) C 4 - C 6	0.51	1.24	0.032
15. BD (1) C 6 - N 11	/594. BD* (1) C 5 - C 6	0.93	1.42	0.046
15. BD (1) C 6 - N 11	/606. BD* (1) N 11 - N 12	0.46	1.28	0.031

15. BD (1) C 6 - N 11	/607. BD* (1) N 12 - C 13	2.78	1.15	0.072
16. BD (1) C 7 - C 8	/181. RY* (1) C 2	0.55	2.40	0.046
16. BD (1) C 7 - C 8	/182. RY* (2) C 2	0.42	2.18	0.038
16. BD (1) C 7 - C 8	/245. RY* (1) C 10	0.48	2.02	0.039
16. BD (1) C 7 - C 8	/247. RY* (3) C 10	0.39	2.06	0.036
16. BD (1) C 7 - C 8	/582. BD* (1) C 1 - C 2	1.37	1.16	0.051
16. BD (1) C 7 - C 8	/587. BD* (1) C 2 - C 7	1.47	1.19	0.053
16. BD (1) C 7 - C 8	/599. BD* (1) C 7 - H 22	0.86	1.18	0.040
16. BD (1) C 7 - C 8	/600. BD* (1) C 8 - C 10	1.20	1.22	0.048
16. BD (1) C 7 - C 8	/601. BD* (1) C 8 - H 23	0.76	1.19	0.038
16. BD (1) C 7 - C 8	/605. BD* (1) C 10 - H 25	0.79	1.19	0.039
17. BD (2) C 7 - C 8	/184. RY* (4) C 2	0.27	1.00	0.022
17. BD (2) C 7 - C 8	/246. RY* (2) C 10	0.44	0.81	0.025
17. BD (2) C 7 - C 8	/583. BD* (2) C 1 - C 2	10.43	0.25	0.076
17. BD (2) C 7 - C 8	/603. BD* (2) C 9 - C 10	5.03	0.31	0.055
18. BD (1) C 7 - H 22	/181. RY* (1) C 2	0.73	2.18	0.051
18. BD (1) C 7 - H 22	/229. RY* (1) C 8	0.70	1.71	0.044
18. BD (1) C 7 - H 22	/586. BD* (1) C 2 - C 4	1.62	1.15	0.054
18. BD (1) C 7 - H 22	/597. BD* (1) C 7 - C 8	0.69	1.17	0.036
18. BD (1) C 7 - H 22	/600. BD* (1) C 8 - C 10	2.75	1.00	0.066
18. BD (1) C 7 - H 22	/601. BD* (1) C 8 - H 23	0.47	0.97	0.027
19. BD (1) C 8 - C 10	/223. RY* (3) C 7	0.87	1.77	0.050
19. BD (1) C 8 - C 10	/239. RY* (3) C 9	0.91	1.71	0.050
19. BD (1) C 8 - C 10	/597. BD* (1) C 7 - C 8	1.32	1.28	0.052
19. BD (1) C 8 - C 10	/599. BD* (1) C 7 - H 22	1.61	1.08	0.053
19. BD (1) C 8 - C 10	/601. BD* (1) C 8 - H 23	0.30	1.09	0.023
19. BD (1) C 8 - C 10	/602. BD* (1) C 9 - C 10	1.36	1.29	0.053
19. BD (1) C 8 - C 10	/604. BD* (1) C 9 - H 24	1.65	1.09	0.054
19. BD (1) C 8 - C 10	/605. BD* (1) C 10 - H 25	0.30	1.09	0.023
20. BD (1) C 8 - H 23	/222. RY* (2) C 7	0.70	1.63	0.043
20. BD (1) C 8 - H 23	/245. RY* (1) C 10	0.68	1.79	0.044
20. BD (1) C 8 - H 23	/587. BD* (1) C 2 - C 7	3.15	0.95	0.069
20. BD (1) C 8 - H 23	/597. BD* (1) C 7 - C 8	0.64	1.16	0.034
20. BD (1) C 8 - H 23	/599. BD* (1) C 7 - H 22	0.53	0.95	0.028
20. BD (1) C 8 - H 23	/602. BD* (1) C 9 - C 10	1.41	1.17	0.051
21. BD (1) C 9 - C 10	/197. RY* (1) C 4	0.74	2.29	0.052
21. BD (1) C 9 - C 10	/198. RY* (2) C 4	0.35	2.28	0.036
21. BD (1) C 9 - C 10	/229. RY* (1) C 8	0.43	1.93	0.036
21. BD (1) C 9 - C 10	/231. RY* (3) C 8	0.42	2.03	0.037
21. BD (1) C 9 - C 10	/591. BD* (1) C 4 - C 6	1.29	1.19	0.050
21. BD (1) C 9 - C 10	/593. BD* (1) C 4 - C 9	1.43	1.20	0.052
21. BD (1) C 9 - C 10	/600. BD* (1) C 8 - C 10	1.25	1.22	0.049
21. BD (1) C 9 - C 10	/601. BD* (1) C 8 - H 23	0.78	1.19	0.039
21. BD (1) C 9 - C 10	/604. BD* (1) C 9 - H 24	0.83	1.20	0.040
21. BD (1) C 9 - C 10	/605. BD* (1) C 10 - H 25	0.80	1.19	0.039
22. BD (2) C 9 - C 10	/199. RY* (3) C 4	0.29	0.89	0.021
22. BD (2) C 9 - C 10	/230. RY* (2) C 8	0.43	0.77	0.024
22. BD (2) C 9 - C 10	/592. BD* (2) C 4 - C 6	10.77	0.25	0.078
22. BD (2) C 9 - C 10	/598. BD* (2) C 7 - C 8	6.07	0.29	0.059
23. BD (1) C 9 - H 24	/197. RY* (1) C 4	0.72	2.05	0.049
23. BD (1) C 9 - H 24	/245. RY* (1) C 10	0.63	1.78	0.043
23. BD (1) C 9 - H 24	/586. BD* (1) C 2 - C 4	1.76	1.13	0.056
23. BD (1) C 9 - H 24	/600. BD* (1) C 8 - C 10	2.80	0.98	0.066
23. BD (1) C 9 - H 24	/602. BD* (1) C 9 - C 10	0.73	1.16	0.037
23. BD (1) C 9 - H 24	/605. BD* (1) C 10 - H 25	0.49	0.95	0.027
24. BD (1) C 10 - H 25	/229. RY* (1) C 8	0.53	1.69	0.038
24. BD (1) C 10 - H 25	/238. RY* (2) C 9	0.59	1.59	0.039
24. BD (1) C 10 - H 25	/593. BD* (1) C 4 - C 9	3.05	0.97	0.069
24. BD (1) C 10 - H 25	/597. BD* (1) C 7 - C 8	1.36	1.15	0.050
24. BD (1) C 10 - H 25	/602. BD* (1) C 9 - C 10	0.71	1.16	0.036
24. BD (1) C 10 - H 25	/604. BD* (1) C 9 - H 24	0.48	0.96	0.027
25. BD (1) N 11 - N 12	/213. RY* (1) C 6	0.96	1.87	0.054
25. BD (1) N 11 - N 12	/269. RY* (1) C 13	0.85	2.09	0.053
25. BD (1) N 11 - N 12	/270. RY* (2) C 13	0.63	2.42	0.050
25. BD (1) N 11 - N 12	/594. BD* (1) C 5 - C 6	0.28	1.62	0.027
25. BD (1) N 11 - N 12	/596. BD* (1) C 6 - N 11	0.49	1.36	0.033
25. BD (1) N 11 - N 12	/607. BD* (1) N 12 - C 13	0.40	1.35	0.029
25. BD (1) N 11 - N 12	/608. BD* (1) C 13 - C 14	0.55	1.44	0.036
26. BD (1) N 12 - C 13	/213. RY* (1) C 6	0.31	1.68	0.029
26. BD (1) N 12 - C 13	/253. RY* (1) N 11	0.47	2.15	0.040
26. BD (1) N 12 - C 13	/254. RY* (2) N 11	0.53	2.20	0.043

26. BD (1) N 12 - C 13	/255. RY* (3) N 11	0.54	2.06	0.042
26. BD (1) N 12 - C 13	/256. RY* (4) N 11	0.27	2.33	0.032
26. BD (1) N 12 - C 13	/278. RY* (2) C 14	0.39	2.10	0.036
26. BD (1) N 12 - C 13	/285. RY* (1) C 15	0.46	1.96	0.038
26. BD (1) N 12 - C 13	/596. BD* (1) C 6 - N 11	2.41	1.17	0.067
26. BD (1) N 12 - C 13	/608. BD* (1) C 13 - C 14	0.32	1.24	0.025
26. BD (1) N 12 - C 13	/609. BD* (1) C 13 - C 15	0.99	1.39	0.047
26. BD (1) N 12 - C 13	/611. BD* (1) C 14 - C 16	0.42	1.44	0.031
26. BD (1) N 12 - C 13	/613. BD* (1) C 15 - C 17	1.19	1.24	0.049
27. BD (1) C 13 - C 14	/261. RY* (1) N 12	0.38	1.47	0.030
27. BD (1) C 13 - C 14	/287. RY* (3) C 15	0.29	1.96	0.030
27. BD (1) C 13 - C 14	/288. RY* (4) C 15	0.33	1.68	0.030
27. BD (1) C 13 - C 14	/294. RY* (2) C 16	0.38	1.85	0.034
27. BD (1) C 13 - C 14	/295. RY* (3) C 16	0.70	1.99	0.048
27. BD (1) C 13 - C 14	/606. BD* (1) N 11 - N 12	1.68	1.13	0.055
27. BD (1) C 13 - C 14	/607. BD* (1) N 12 - C 13	0.42	1.00	0.026
27. BD (1) C 13 - C 14	/609. BD* (1) C 13 - C 15	1.88	1.24	0.061
27. BD (1) C 13 - C 14	/611. BD* (1) C 14 - C 16	0.99	1.29	0.045
27. BD (1) C 13 - C 14	/612. BD* (1) C 14 - H 26	0.25	1.09	0.021
27. BD (1) C 13 - C 14	/614. BD* (1) C 15 - O 19	2.73	0.96	0.065
27. BD (1) C 13 - C 14	/616. BD* (1) C 16 - H 27	1.45	1.09	0.051
28. BD (1) C 13 - C 15	/263. RY* (3) N 12	0.34	1.87	0.032
28. BD (1) C 13 - C 15	/278. RY* (2) C 14	0.44	2.07	0.038
28. BD (1) C 13 - C 15	/302. RY* (2) C 17	0.68	2.09	0.048
28. BD (1) C 13 - C 15	/607. BD* (1) N 12 - C 13	0.70	1.12	0.035
28. BD (1) C 13 - C 15	/608. BD* (1) C 13 - C 14	1.85	1.21	0.060
28. BD (1) C 13 - C 15	/612. BD* (1) C 14 - H 26	0.57	1.20	0.033
28. BD (1) C 13 - C 15	/613. BD* (1) C 15 - C 17	1.42	1.21	0.053
28. BD (1) C 13 - C 15	/614. BD* (1) C 15 - O 19	0.37	1.08	0.025
28. BD (1) C 13 - C 15	/618. BD* (1) C 17 - H 28	0.57	1.21	0.033
29. BD (2) C 13 - C 15	/ 34. BD* (1) C 16 - C 17	0.40	0.13	0.053
29. BD (2) C 13 - C 15	/262. RY* (2) N 12	0.41	1.09	0.027
29. BD (2) C 13 - C 15	/277. RY (1) C 14	0.79	0.80	0.032
29. BD (2) C 13 - C 15	/301. RY (1) C 17	1.12	0.81	0.038
29. BD (2) C 13 - C 15	/318. RY* (2) O 19	0.39	1.40	0.030
29. BD (2) C 13 - C 15	/610. BD* (2) C 13 - C 15	0.30	0.33	0.014
30. BD (1) C 14 - C 16	/270. RY* (2) C 13	1.01	2.18	0.059
30. BD (1) C 14 - C 16	/309. RY* (1) C 18	0.55	1.94	0.041
30. BD (1) C 14 - C 16	/311. RY* (3) C 18	0.43	2.05	0.038
30. BD (1) C 14 - C 16	/607. BD* (1) N 12 - C 13	1.38	1.11	0.050
30. BD (1) C 14 - C 16	/608. BD* (1) C 13 - C 14	0.95	1.20	0.043
30. BD (1) C 14 - C 16	/612. BD* (1) C 14 - H 26	0.71	1.19	0.037
30. BD (1) C 14 - C 16	/615. BD* (1) C 16 - C 18	0.93	1.22	0.043
30. BD (1) C 14 - C 16	/616. BD* (1) C 16 - H 27	0.58	1.20	0.033
30. BD (1) C 14 - C 16	/619. BD* (1) C 18 - H 29	0.86	1.21	0.041
31. BD (1) C 14 - H 26	/269. RY* (1) C 13	0.35	1.61	0.030
31. BD (1) C 14 - H 26	/294. RY* (2) C 16	0.62	1.71	0.041
31. BD (1) C 14 - H 26	/607. BD* (1) N 12 - C 13	0.36	0.86	0.022
31. BD (1) C 14 - H 26	/609. BD* (1) C 13 - C 15	1.71	1.10	0.055
31. BD (1) C 14 - H 26	/611. BD* (1) C 14 - C 16	0.54	1.15	0.032
31. BD (1) C 14 - H 26	/615. BD* (1) C 16 - C 18	2.73	0.97	0.065
31. BD (1) C 14 - H 26	/616. BD* (1) C 16 - H 27	0.51	0.95	0.028
32. BD (1) C 15 - C 17	/269. RY* (1) C 13	0.59	1.75	0.041
32. BD (1) C 15 - C 17	/271. RY* (3) C 13	0.30	1.46	0.027
32. BD (1) C 15 - C 17	/309. RY* (1) C 18	0.34	1.84	0.032
32. BD (1) C 15 - C 17	/311. RY* (3) C 18	0.79	1.94	0.050
32. BD (1) C 15 - C 17	/607. BD* (1) N 12 - C 13	2.85	1.01	0.068
32. BD (1) C 15 - C 17	/609. BD* (1) C 13 - C 15	1.68	1.24	0.058
32. BD (1) C 15 - C 17	/617. BD* (1) C 17 - C 18	1.08	1.30	0.047
32. BD (1) C 15 - C 17	/619. BD* (1) C 18 - H 29	1.62	1.10	0.054
33. BD (1) C 15 - O 19	/285. RY* (1) C 15	0.55	2.01	0.042
33. BD (1) C 15 - O 19	/608. BD* (1) C 13 - C 14	1.26	1.29	0.051
33. BD (1) C 15 - O 19	/609. BD* (1) C 13 - C 15	0.63	1.44	0.038
33. BD (1) C 15 - O 19	/617. BD* (1) C 17 - C 18	0.50	1.50	0.035
34. BD* (1) C 16 - C 17	/277. RY (1) C 14	1.47	0.67	0.040
34. BD* (1) C 16 - C 17	/286. RY* (2) C 15	0.28	0.93	0.020
34. BD* (1) C 16 - C 17	/610. BD* (2) C 13 - C 15	6.55	0.20	0.051
34. BD* (1) C 16 - C 17	/615. BD* (1) C 16 - C 18	0.42	0.65	0.021
35. BD (1) C 16 - C 17	/277. RY (1) C 14	0.83	0.63	0.039
35. BD (1) C 16 - C 17	/310. RY* (2) C 18	1.61	0.64	0.056
35. BD (1) C 16 - C 17	/610. BD* (2) C 13 - C 15	9.86	0.16	0.058

36. BD (1) C 16 - C 18	/279. RY* (3) C 14	0.88	1.89	0.052
36. BD (1) C 16 - C 18	/303. RY* (3) C 17	0.88	1.86	0.051
36. BD (1) C 16 - C 18	/611. BD* (1) C 14 - C 16	1.00	1.28	0.045
36. BD (1) C 16 - C 18	/612. BD* (1) C 14 - H 26	1.65	1.08	0.053
36. BD (1) C 16 - C 18	/617. BD* (1) C 17 - C 18	1.17	1.29	0.049
36. BD (1) C 16 - C 18	/618. BD* (1) C 17 - H 28	1.56	1.08	0.052
36. BD (1) C 16 - C 18	/619. BD* (1) C 18 - H 29	0.29	1.09	0.023
37. BD (1) C 16 - H 27	/278. RY* (2) C 14	0.78	1.82	0.048
37. BD (1) C 16 - H 27	/309. RY* (1) C 18	0.66	1.70	0.043
37. BD (1) C 16 - H 27	/608. BD* (1) C 13 - C 14	3.03	0.96	0.068
37. BD (1) C 16 - H 27	/611. BD* (1) C 14 - C 16	0.46	1.16	0.029
37. BD (1) C 16 - H 27	/612. BD* (1) C 14 - H 26	0.53	0.95	0.028
37. BD (1) C 16 - H 27	/617. BD* (1) C 17 - C 18	1.45	1.17	0.052
38. BD (1) C 17 - C 18	/285. RY* (1) C 15	0.51	1.91	0.040
38. BD (1) C 17 - C 18	/287. RY* (3) C 15	0.57	2.06	0.043
38. BD (1) C 17 - C 18	/294. RY* (2) C 16	0.51	1.95	0.040
38. BD (1) C 17 - C 18	/295. RY* (3) C 16	0.38	2.09	0.036
38. BD (1) C 17 - C 18	/613. BD* (1) C 15 - C 17	0.80	1.19	0.039
38. BD (1) C 17 - C 18	/614. BD* (1) C 15 - O 19	1.19	1.06	0.045
38. BD (1) C 17 - C 18	/615. BD* (1) C 16 - C 18	1.08	1.21	0.046
38. BD (1) C 17 - C 18	/616. BD* (1) C 16 - H 27	0.76	1.19	0.038
38. BD (1) C 17 - C 18	/618. BD* (1) C 17 - H 28	0.77	1.19	0.038
38. BD (1) C 17 - C 18	/619. BD* (1) C 18 - H 29	0.75	1.20	0.038
39. BD (1) C 17 - H 28	/285. RY* (1) C 15	0.39	1.67	0.032
39. BD (1) C 17 - H 28	/309. RY* (1) C 18	0.67	1.70	0.043
39. BD (1) C 17 - H 28	/609. BD* (1) C 13 - C 15	1.64	1.10	0.054
39. BD (1) C 17 - H 28	/614. BD* (1) C 15 - O 19	0.54	0.83	0.027
39. BD (1) C 17 - H 28	/615. BD* (1) C 16 - C 18	2.90	0.97	0.067
39. BD (1) C 17 - H 28	/617. BD* (1) C 17 - C 18	0.63	1.16	0.034
39. BD (1) C 17 - H 28	/619. BD* (1) C 18 - H 29	0.51	0.96	0.028
40. BD (1) C 18 - H 29	/294. RY* (2) C 16	0.61	1.72	0.041
40. BD (1) C 18 - H 29	/302. RY* (2) C 17	0.77	1.84	0.048
40. BD (1) C 18 - H 29	/611. BD* (1) C 14 - C 16	1.31	1.15	0.049
40. BD (1) C 18 - H 29	/613. BD* (1) C 15 - C 17	2.80	0.95	0.066
40. BD (1) C 18 - H 29	/617. BD* (1) C 17 - C 18	0.64	1.16	0.034
40. BD (1) C 18 - H 29	/618. BD* (1) C 17 - H 28	0.46	0.95	0.026
41. BD (1) O 20 - H 30	/173. RY* (1) C 1	1.02	1.70	0.053
41. BD (1) O 20 - H 30	/174. RY* (2) C 1	0.45	1.46	0.032
41. BD (1) O 20 - H 30	/584. BD* (1) C 1 - C 3	1.60	1.34	0.059
89. CR (1) C 1	/181. RY* (1) C 2	0.62	11.78	0.108
89. CR (1) C 1	/191. RY* (3) C 3	1.17	11.44	0.146
89. CR (1) C 1	/584. BD* (1) C 1 - C 3	1.00	10.73	0.131
89. CR (1) C 1	/585. BD* (1) C 1 - O 20	0.66	10.39	0.105
89. CR (1) C 1	/586. BD* (1) C 2 - C 4	0.30	10.75	0.072
89. CR (1) C 1	/587. BD* (1) C 2 - C 7	0.33	10.57	0.075
89. CR (1) C 1	/588. BD* (1) C 3 - C 5	0.45	10.59	0.088
90. CR (1) C 2	/173. RY* (1) C 1	0.31	11.03	0.074
90. CR (1) C 2	/197. RY* (1) C 4	0.28	11.61	0.072
90. CR (1) C 2	/198. RY* (2) C 4	1.12	11.60	0.144
90. CR (1) C 2	/223. RY* (3) C 7	0.50	11.20	0.094
90. CR (1) C 2	/584. BD* (1) C 1 - C 3	0.29	10.67	0.070
90. CR (1) C 2	/586. BD* (1) C 2 - C 4	0.66	10.69	0.106
90. CR (1) C 2	/591. BD* (1) C 4 - C 6	0.59	10.52	0.101
90. CR (1) C 2	/593. BD* (1) C 4 - C 9	0.53	10.53	0.095
90. CR (1) C 2	/597. BD* (1) C 7 - C 8	0.27	10.71	0.068
91. CR (1) C 3	/175. RY* (3) C 1	1.45	11.49	0.163
91. CR (1) C 3	/208. RY* (4) C 5	0.41	11.10	0.085
91. CR (1) C 3	/334. RY* (2) H 21	0.30	12.29	0.076
91. CR (1) C 3	/582. BD* (1) C 1 - C 2	0.61	10.47	0.102
91. CR (1) C 3	/584. BD* (1) C 1 - C 3	0.29	10.66	0.070
91. CR (1) C 3	/585. BD* (1) C 1 - O 20	0.37	10.32	0.078
91. CR (1) C 3	/594. BD* (1) C 5 - C 6	0.34	10.68	0.077
92. CR (1) C 4	/182. RY* (2) C 2	1.18	11.49	0.147
92. CR (1) C 4	/215. RY* (3) C 6	0.77	11.17	0.117
92. CR (1) C 4	/239. RY* (3) C 9	0.56	11.13	0.100
92. CR (1) C 4	/582. BD* (1) C 1 - C 2	0.57	10.47	0.099
92. CR (1) C 4	/586. BD* (1) C 2 - C 4	0.57	10.68	0.099
92. CR (1) C 4	/587. BD* (1) C 2 - C 7	0.51	10.50	0.093
92. CR (1) C 4	/594. BD* (1) C 5 - C 6	0.34	10.69	0.077
92. CR (1) C 4	/596. BD* (1) C 6 - N 11	0.26	10.42	0.066
92. CR (1) C 4	/602. BD* (1) C 9 - C 10	0.29	10.71	0.070

93.	CR	(1)	C	5	/214.	RY*	(2)	C	6	0.46	11.20	0.090	
93.	CR	(1)	C	5	/215.	RY*	(3)	C	6	0.76	11.16	0.116	
93.	CR	(1)	C	5	/216.	RY*	(4)	C	6	0.54	11.09	0.098	
93.	CR	(1)	C	5	/584.	BD*	(1)	C	1 - C	3	0.33	10.65	0.075
93.	CR	(1)	C	5	/591.	BD*	(1)	C	4 - C	6	0.64	10.49	0.105
93.	CR	(1)	C	5	/594.	BD*	(1)	C	5 - C	6	0.42	10.67	0.085
93.	CR	(1)	C	5	/596.	BD*	(1)	C	6 - N	11	0.44	10.41	0.086
94.	CR	(1)	C	6	/197.	RY*	(1)	C	4	0.66	11.64	0.111	
94.	CR	(1)	C	6	/207.	RY*	(3)	C	5	1.25	11.28	0.150	
94.	CR	(1)	C	6	/586.	BD*	(1)	C	2 - C	4	0.36	10.72	0.079
94.	CR	(1)	C	6	/588.	BD*	(1)	C	3 - C	5	0.48	10.55	0.090
94.	CR	(1)	C	6	/593.	BD*	(1)	C	4 - C	9	0.33	10.55	0.075
94.	CR	(1)	C	6	/594.	BD*	(1)	C	5 - C	6	0.77	10.72	0.116
94.	CR	(1)	C	6	/606.	BD*	(1)	N	11 - N	12	0.39	10.58	0.082
95.	CR	(1)	C	7	/181.	RY*	(1)	C	2	0.25	11.69	0.068	
95.	CR	(1)	C	7	/182.	RY*	(2)	C	2	0.34	11.47	0.079	
95.	CR	(1)	C	7	/231.	RY*	(3)	C	8	1.14	11.33	0.144	
95.	CR	(1)	C	7	/336.	RY*	(2)	H	22	0.27	12.26	0.073	
95.	CR	(1)	C	7	/582.	BD*	(1)	C	1 - C	2	0.35	10.45	0.077
95.	CR	(1)	C	7	/586.	BD*	(1)	C	2 - C	4	0.37	10.66	0.080
95.	CR	(1)	C	7	/597.	BD*	(1)	C	7 - C	8	0.45	10.68	0.088
95.	CR	(1)	C	7	/600.	BD*	(1)	C	8 - C	10	0.52	10.51	0.094
96.	CR	(1)	C	8	/222.	RY*	(2)	C	7	0.27	11.15	0.070	
96.	CR	(1)	C	8	/223.	RY*	(3)	C	7	0.64	11.16	0.106	
96.	CR	(1)	C	8	/224.	RY*	(4)	C	7	0.42	11.07	0.086	
96.	CR	(1)	C	8	/247.	RY*	(3)	C	10	0.45	11.34	0.090	
96.	CR	(1)	C	8	/338.	RY*	(2)	H	23	0.31	11.94	0.076	
96.	CR	(1)	C	8	/587.	BD*	(1)	C	2 - C	7	0.61	10.48	0.102
96.	CR	(1)	C	8	/597.	BD*	(1)	C	7 - C	8	0.38	10.68	0.080
96.	CR	(1)	C	8	/602.	BD*	(1)	C	9 - C	10	0.32	10.69	0.074
97.	CR	(1)	C	9	/198.	RY*	(2)	C	4	0.45	11.56	0.091	
97.	CR	(1)	C	9	/247.	RY*	(3)	C	10	1.07	11.34	0.139	
97.	CR	(1)	C	9	/586.	BD*	(1)	C	2 - C	4	0.43	10.65	0.086
97.	CR	(1)	C	9	/591.	BD*	(1)	C	4 - C	6	0.35	10.48	0.077
97.	CR	(1)	C	9	/600.	BD*	(1)	C	8 - C	10	0.54	10.50	0.095
97.	CR	(1)	C	9	/602.	BD*	(1)	C	9 - C	10	0.45	10.68	0.088
98.	CR	(1)	C	10	/231.	RY*	(3)	C	8	0.42	11.32	0.088	
98.	CR	(1)	C	10	/238.	RY*	(2)	C	9	0.25	11.11	0.067	
98.	CR	(1)	C	10	/239.	RY*	(3)	C	9	0.56	11.10	0.100	
98.	CR	(1)	C	10	/240.	RY*	(4)	C	9	0.58	11.14	0.102	
98.	CR	(1)	C	10	/342.	RY*	(2)	H	25	0.33	12.33	0.081	
98.	CR	(1)	C	10	/593.	BD*	(1)	C	4 - C	9	0.60	10.49	0.100
98.	CR	(1)	C	10	/597.	BD*	(1)	C	7 - C	8	0.32	10.67	0.074
98.	CR	(1)	C	10	/602.	BD*	(1)	C	9 - C	10	0.41	10.69	0.083
99.	CR	(1)	N	11	/213.	RY*	(1)	C	6	0.43	15.09	0.103	
99.	CR	(1)	N	11	/214.	RY*	(2)	C	6	0.68	15.36	0.130	
99.	CR	(1)	N	11	/263.	RY*	(3)	N	12	0.88	15.31	0.146	
99.	CR	(1)	N	11	/591.	BD*	(1)	C	4 - C	6	0.27	14.66	0.080
99.	CR	(1)	N	11	/594.	BD*	(1)	C	5 - C	6	0.27	14.84	0.080
99.	CR	(1)	N	11	/607.	BD*	(1)	N	12 - C	13	0.71	14.56	0.130
100.	CR	(1)	N	12	/255.	RY*	(3)	N	11	0.99	15.51	0.157	
100.	CR	(1)	N	12	/269.	RY*	(1)	C	13	0.60	15.35	0.121	
100.	CR	(1)	N	12	/270.	RY*	(2)	C	13	0.43	15.68	0.104	
100.	CR	(1)	N	12	/596.	BD*	(1)	C	6 - N	11	0.62	14.62	0.122
100.	CR	(1)	N	12	/609.	BD*	(1)	C	13 - C	15	0.36	14.84	0.094
101.	CR	(1)	C	13	/279.	RY*	(3)	C	14	0.44	11.36	0.090	
101.	CR	(1)	C	13	/285.	RY*	(1)	C	15	0.52	11.26	0.097	
101.	CR	(1)	C	13	/287.	RY*	(3)	C	15	0.99	11.41	0.134	
101.	CR	(1)	C	13	/606.	BD*	(1)	N	11 - N	12	0.49	10.59	0.092
101.	CR	(1)	C	13	/609.	BD*	(1)	C	13 - C	15	0.58	10.70	0.101
101.	CR	(1)	C	13	/611.	BD*	(1)	C	14 - C	16	0.25	10.75	0.066
101.	CR	(1)	C	13	/613.	BD*	(1)	C	15 - C	17	0.47	10.55	0.090
101.	CR	(1)	C	13	/614.	BD*	(1)	C	15 - O	19	0.44	10.42	0.086
102.	CR	(1)	C	14	/269.	RY*	(1)	C	13	0.39	11.12	0.083	
102.	CR	(1)	C	14	/271.	RY*	(3)	C	13	0.34	10.83	0.076	
102.	CR	(1)	C	14	/295.	RY*	(3)	C	16	1.02	11.37	0.136	
102.	CR	(1)	C	14	/344.	RY*	(2)	H	26	0.29	12.29	0.076	
102.	CR	(1)	C	14	/607.	BD*	(1)	N	12 - C	13	0.26	10.38	0.066
102.	CR	(1)	C	14	/609.	BD*	(1)	C	13 - C	15	0.40	10.62	0.083
102.	CR	(1)	C	14	/611.	BD*	(1)	C	14 - C	16	0.42	10.67	0.085
102.	CR	(1)	C	14	/615.	BD*	(1)	C	16 - C	18	0.47	10.49	0.089

103. CR (1) C 15	/270. RY* (2) C 13	1.47	11.54	0.164
103. CR (1) C 15	/303. RY* (3) C 17	0.50	11.33	0.095
103. CR (1) C 15	/607. BD* (1) N 12 - C 13	0.50	10.46	0.092
103. CR (1) C 15	/608. BD* (1) C 13 - C 14	0.50	10.55	0.092
103. CR (1) C 15	/609. BD* (1) C 13 - C 15	0.77	10.70	0.116
103. CR (1) C 15	/614. BD* (1) C 15 - O 19	0.45	10.43	0.087
103. CR (1) C 15	/617. BD* (1) C 17 - C 18	0.28	10.76	0.069
104. CR (1) C 16	/278. RY* (2) C 14	0.31	11.34	0.075
104. CR (1) C 16	/279. RY* (3) C 14	0.75	11.29	0.116
104. CR (1) C 16	/280. RY* (4) C 14	0.28	10.94	0.070
104. CR (1) C 16	/311. RY* (3) C 18	0.50	11.33	0.095
104. CR (1) C 16	/346. RY* (2) H 27	0.31	11.97	0.077
104. CR (1) C 16	/608. BD* (1) C 13 - C 14	0.51	10.48	0.093
104. CR (1) C 16	/611. BD* (1) C 14 - C 16	0.34	10.68	0.077
104. CR (1) C 16	/617. BD* (1) C 17 - C 18	0.30	10.69	0.072
105. CR (1) C 17	/287. RY* (3) C 15	0.36	11.33	0.080
105. CR (1) C 17	/288. RY* (4) C 15	0.51	11.06	0.095
105. CR (1) C 17	/311. RY* (3) C 18	1.07	11.32	0.139
105. CR (1) C 17	/348. RY* (2) H 28	0.29	12.30	0.076
105. CR (1) C 17	/609. BD* (1) C 13 - C 15	0.42	10.62	0.085
105. CR (1) C 17	/615. BD* (1) C 16 - C 18	0.51	10.49	0.093
105. CR (1) C 17	/617. BD* (1) C 17 - C 18	0.47	10.68	0.089
106. CR (1) C 18	/295. RY* (3) C 16	0.42	11.38	0.088
106. CR (1) C 18	/302. RY* (2) C 17	0.43	11.36	0.089
106. CR (1) C 18	/303. RY* (3) C 17	0.67	11.26	0.110
106. CR (1) C 18	/350. RY* (2) H 29	0.32	11.98	0.078
106. CR (1) C 18	/611. BD* (1) C 14 - C 16	0.28	10.68	0.070
106. CR (1) C 18	/613. BD* (1) C 15 - C 17	0.49	10.48	0.091
106. CR (1) C 18	/617. BD* (1) C 17 - C 18	0.35	10.69	0.077
107. CR (1) O 19	/285. RY* (1) C 15	1.41	20.07	0.213
107. CR (1) O 19	/609. BD* (1) C 13 - C 15	0.30	19.50	0.097
108. CR (1) O 20	/173. RY* (1) C 1	0.85	19.94	0.165
108. CR (1) O 20	/352. RY* (2) H 30	0.30	20.86	0.100
140. LP (1) N 11	/261. RY* (1) N 12	0.75	1.15	0.038
140. LP (1) N 11	/262. RY* (2) N 12	0.68	1.11	0.035
140. LP (1) N 11	/591. BD* (1) C 4 - C 6	2.21	0.77	0.053
140. LP (1) N 11	/592. BD* (2) C 4 - C 6	0.68	0.31	0.022
140. LP (1) N 11	/594. BD* (1) C 5 - C 6	3.01	0.95	0.069
141. LP (2) N 11	/213. RY* (1) C 6	0.72	1.33	0.040
141. LP (2) N 11	/253. RY* (1) N 11	1.17	1.80	0.059
141. LP (2) N 11	/254. RY* (2) N 11	0.87	1.85	0.052
141. LP (2) N 11	/255. RY* (3) N 11	0.47	1.72	0.037
141. LP (2) N 11	/262. RY* (2) N 12	0.52	1.23	0.033
141. LP (2) N 11	/317. RY* (1) O 19	0.52	2.82	0.049
141. LP (2) N 11	/591. BD* (1) C 4 - C 6	2.22	0.90	0.057
141. LP (2) N 11	/592. BD* (2) C 4 - C 6	0.93	0.44	0.031
142. LP (1) N 12	/254. RY* (2) N 11	0.46	1.85	0.037
142. LP (1) N 12	/255. RY* (3) N 11	0.31	1.71	0.029
142. LP (1) N 12	/269. RY* (1) C 13	0.88	1.54	0.047
142. LP (1) N 12	/608. BD* (1) C 13 - C 14	0.56	0.89	0.028
142. LP (1) N 12	/609. BD* (1) C 13 - C 15	3.64	1.04	0.078
143. LP (2) N 12	/253. RY* (1) N 11	0.76	1.65	0.047
143. LP (2) N 12	/271. RY* (3) C 13	0.30	1.11	0.024
143. LP (2) N 12	/610. BD* (2) C 13 - C 15	15.99	0.32	0.097
144. LP (1) C 14	/293. RY* (1) C 16	2.99	0.66	0.060
144. LP (1) C 14	/610. BD* (2) C 13 - C 15	26.11	0.19	0.093
145. LP (1) C 18	/293. RY* (1) C 16	1.05	0.60	0.043
145. LP (1) C 18	/301. RY (1) C 17	2.06	0.61	0.062
145. LP (1) C 18	/310. RY* (2) C 18	1.28	0.62	0.049
145. LP (1) C 18	/610. BD* (2) C 13 - C 15	0.25	0.13	0.009
145. LP (1) C 18	/615. BD* (1) C 16 - C 18	0.73	0.58	0.036
146. LP (1) O 19	/285. RY* (1) C 15	1.13	1.72	0.056
146. LP (1) O 19	/609. BD* (1) C 13 - C 15	3.53	1.15	0.081
147. LP (2) O 19	/285. RY* (1) C 15	0.50	1.61	0.036
147. LP (2) O 19	/317. RY* (1) O 19	0.55	2.82	0.051
147. LP (2) O 19	/609. BD* (1) C 13 - C 15	0.27	1.04	0.021
147. LP (2) O 19	/610. BD* (2) C 13 - C 15	1.48	0.47	0.037
147. LP (2) O 19	/613. BD* (1) C 15 - C 17	3.17	0.89	0.068
148. LP (3) O 19	/285. RY* (1) C 15	0.27	1.54	0.026
148. LP (3) O 19	/317. RY* (1) O 19	0.50	2.75	0.048
148. LP (3) O 19	/610. BD* (2) C 13 - C 15	6.86	0.40	0.073

148. LP (3) O 19	/613. BD* (1) C 15 - C 17	0.81	0.82	0.033
149. LP (1) O 20	/173. RY* (1) C 1	1.04	1.58	0.051
149. LP (1) O 20	/174. RY* (2) C 1	0.33	1.34	0.027
149. LP (1) O 20	/582. BD* (1) C 1 - C 2	3.15	1.03	0.072
150. LP (2) O 20	/583. BD* (2) C 1 - C 2	4.52	0.32	0.060
583. BD* (2) C 1 - C 2	/174. RY* (2) C 1	0.39	0.75	0.036
583. BD* (2) C 1 - C 2	/183. RY* (3) C 2	0.30	0.95	0.035
583. BD* (2) C 1 - C 2	/184. RY* (4) C 2	1.26	0.75	0.064
583. BD* (2) C 1 - C 2	/189. RY* (1) C 3	0.44	0.53	0.031
583. BD* (2) C 1 - C 2	/199. RY* (3) C 4	0.41	0.64	0.034
583. BD* (2) C 1 - C 2	/221. RY* (1) C 7	0.39	0.61	0.032
583. BD* (2) C 1 - C 2	/592. BD* (2) C 4 - C 6	720.14	0.01	0.129
583. BD* (2) C 1 - C 2	/598. BD* (2) C 7 - C 8	47.20	0.05	0.077
583. BD* (2) C 1 - C 2	/603. BD* (2) C 9 - C 10	0.26	0.06	0.006
589. BD* (2) C 3 - C 5	/189. RY* (1) C 3	0.81	0.52	0.052
589. BD* (2) C 3 - C 5	/205. RY* (1) C 5	0.78	0.56	0.053
592. BD* (2) C 4 - C 6	/184. RY* (4) C 2	0.46	0.74	0.038
592. BD* (2) C 4 - C 6	/199. RY* (3) C 4	3.31	0.63	0.094
592. BD* (2) C 4 - C 6	/205. RY* (1) C 5	0.45	0.56	0.032
592. BD* (2) C 4 - C 6	/213. RY* (1) C 6	0.37	0.90	0.037
592. BD* (2) C 4 - C 6	/237. RY* (1) C 9	0.44	0.74	0.037
592. BD* (2) C 4 - C 6	/598. BD* (2) C 7 - C 8	0.32	0.04	0.006
592. BD* (2) C 4 - C 6	/603. BD* (2) C 9 - C 10	58.19	0.05	0.083
592. BD* (2) C 4 - C 6	/606. BD* (1) N 11 - N 12	0.86	0.50	0.042
598. BD* (2) C 7 - C 8	/221. RY* (1) C 7	1.39	0.56	0.067
598. BD* (2) C 7 - C 8	/230. RY* (2) C 8	2.38	0.48	0.081
603. BD* (2) C 9 - C 10	/237. RY* (1) C 9	1.14	0.70	0.068
603. BD* (2) C 9 - C 10	/246. RY* (2) C 10	2.18	0.51	0.080
610. BD* (2) C 13 - C 15	/272. RY* (4) C 13	0.65	0.68	0.058
610. BD* (2) C 13 - C 15	/277. RY (1) C 14	0.37	0.47	0.035
610. BD* (2) C 13 - C 15	/287. RY* (3) C 15	0.38	1.29	0.060
610. BD* (2) C 13 - C 15	/301. RY (1) C 17	0.27	0.48	0.031
610. BD* (2) C 13 - C 15	/609. BD* (1) C 13 - C 15	1.58	0.57	0.078

from unit 1 to unit 2

21. BD (1) C 9 - C 10	/521. RY* (1) H 58	0.06	1.27	0.011
22. BD (2) C 9 - C 10	/657. BD* (1) C 48 - H 58	0.09	0.72	0.011
33. BD (1) C 15 - O 19	/435. RY* (1) N 42	0.03	2.73	0.012
33. BD (1) C 15 - O 19	/451. RY* (1) C 44	0.05	1.96	0.012
33. BD (1) C 15 - O 19	/468. RY* (2) C 46	0.03	2.29	0.010
34. BD* (1) C 16 - C 17	/527. RY* (1) H 61	0.03	0.66	0.005
140. LP (1) N 11	/435. RY* (1) N 42	0.04	2.21	0.013
140. LP (1) N 11	/437. RY* (3) N 42	0.04	1.54	0.011
140. LP (1) N 11	/541. RY* (1) O 65	0.07	2.71	0.018
140. LP (1) N 11	/652. BD* (1) C 46 - O 65	0.12	0.65	0.011
141. LP (2) N 11	/396. RY* (2) C 37	0.13	1.54	0.019
141. LP (2) N 11	/398. RY* (4) C 37	0.03	1.49	0.009
141. LP (2) N 11	/400. RY* (6) C 37	0.05	3.58	0.018
141. LP (2) N 11	/435. RY* (1) N 42	0.30	2.34	0.034
141. LP (2) N 11	/436. RY* (2) N 42	0.07	1.39	0.013
141. LP (2) N 11	/437. RY* (3) N 42	0.16	1.67	0.021
141. LP (2) N 11	/438. RY* (4) N 42	0.06	1.94	0.014
141. LP (2) N 11	/445. RY* (3) N 43	0.04	1.54	0.010
141. LP (2) N 11	/467. RY* (1) C 46	0.15	1.61	0.020
141. LP (2) N 11	/468. RY* (2) C 46	0.03	1.90	0.010
141. LP (2) N 11	/470. RY* (4) C 46	0.03	1.21	0.007
141. LP (2) N 11	/471. RY* (5) C 46	0.03	5.29	0.017
141. LP (2) N 11	/541. RY* (1) O 65	0.45	2.83	0.046
141. LP (2) N 11	/543. RY* (3) O 65	0.09	1.85	0.017
141. LP (2) N 11	/544. RY* (4) O 65	0.07	2.13	0.015
141. LP (2) N 11	/652. BD* (1) C 46 - O 65	0.03	0.77	0.006
146. LP (1) O 19	/396. RY* (2) C 37	0.03	1.65	0.009
146. LP (1) O 19	/435. RY* (1) N 42	0.14	2.45	0.023
146. LP (1) O 19	/436. RY* (2) N 42	0.03	1.50	0.008
146. LP (1) O 19	/437. RY* (3) N 42	0.10	1.78	0.017
146. LP (1) O 19	/446. RY* (4) N 43	0.03	2.16	0.010
146. LP (1) O 19	/467. RY* (1) C 46	0.03	1.72	0.009
146. LP (1) O 19	/541. RY* (1) O 65	0.08	2.94	0.020
146. LP (1) O 19	/633. BD* (1) C 37 - N 42	0.03	0.93	0.006
146. LP (1) O 19	/644. BD* (1) N 43 - C 44	0.04	0.93	0.008
146. LP (1) O 19	/652. BD* (1) C 46 - O 65	0.03	0.88	0.007

147. LP (2) O 19	/396. RY* (2) C 37	0.09	1.53	0.015
147. LP (2) O 19	/400. RY* (6) C 37	0.03	3.58	0.012
147. LP (2) O 19	/435. RY* (1) N 42	0.24	2.34	0.030
147. LP (2) O 19	/437. RY* (3) N 42	0.10	1.67	0.017
147. LP (2) O 19	/467. RY* (1) C 46	0.09	1.60	0.016
147. LP (2) O 19	/468. RY* (2) C 46	0.04	1.89	0.011
147. LP (2) O 19	/541. RY* (1) O 65	0.16	2.83	0.027
147. LP (2) O 19	/544. RY* (4) O 65	0.05	2.13	0.013
147. LP (2) O 19	/647. BD* (2) C 44 - C 46	0.12	0.46	0.010
147. LP (2) O 19	/652. BD* (1) C 46 - O 65	0.19	0.77	0.015
148. LP (3) O 19	/396. RY* (2) C 37	0.06	1.47	0.012
148. LP (3) O 19	/435. RY* (1) N 42	0.27	2.27	0.032
148. LP (3) O 19	/436. RY* (2) N 42	0.03	1.32	0.007
148. LP (3) O 19	/437. RY* (3) N 42	0.09	1.60	0.015
148. LP (3) O 19	/438. RY* (4) N 42	0.08	1.87	0.016
148. LP (3) O 19	/467. RY* (1) C 46	0.06	1.53	0.012
148. LP (3) O 19	/541. RY* (1) O 65	0.14	2.76	0.025
148. LP (3) O 19	/633. BD* (1) C 37 - N 42	0.08	0.75	0.010

from unit 1 to unit 3

26. BD (1) N 12 - C 13	/529. RY* (1) O 62	0.04	2.82	0.013
26. BD (1) N 12 - C 13	/537. RY* (1) H 63	0.04	1.35	0.009
28. BD (1) C 13 - C 15	/539. RY* (1) H 64	0.04	1.32	0.009
29. BD (2) C 13 - C 15	/661. BD* (1) O 62 - H 64	0.19	0.64	0.014
32. BD (1) C 15 - C 17	/529. RY* (1) O 62	0.04	2.67	0.013
32. BD (1) C 15 - C 17	/539. RY* (1) H 64	0.05	1.21	0.010
140. LP (1) N 11	/529. RY* (1) O 62	0.12	2.35	0.022
140. LP (1) N 11	/532. RY* (4) O 62	0.04	1.68	0.011
140. LP (1) N 11	/537. RY* (1) H 63	0.06	0.87	0.009
140. LP (1) N 11	/539. RY* (1) H 64	0.03	0.88	0.007
140. LP (1) N 11	/660. BD* (1) O 62 - H 63	0.15	0.65	0.013
141. LP (2) N 11	/529. RY* (1) O 62	0.85	2.47	0.059
141. LP (2) N 11	/531. RY* (3) O 62	0.17	1.80	0.022
141. LP (2) N 11	/532. RY* (4) O 62	0.27	1.81	0.029
141. LP (2) N 11	/537. RY* (1) H 63	0.26	1.00	0.021
141. LP (2) N 11	/539. RY* (1) H 64	0.23	1.01	0.020
146. LP (1) O 19	/529. RY* (1) O 62	0.19	2.58	0.028
146. LP (1) O 19	/531. RY* (3) O 62	0.03	1.91	0.010
146. LP (1) O 19	/532. RY* (4) O 62	0.04	1.92	0.012
146. LP (1) O 19	/537. RY* (1) H 63	0.07	1.11	0.011
146. LP (1) O 19	/539. RY* (1) H 64	0.05	1.12	0.009
146. LP (1) O 19	/661. BD* (1) O 62 - H 64	0.03	0.89	0.006
147. LP (2) O 19	/529. RY* (1) O 62	0.47	2.47	0.043
147. LP (2) O 19	/531. RY* (3) O 62	0.08	1.80	0.015
147. LP (2) O 19	/532. RY* (4) O 62	0.13	1.80	0.020
147. LP (2) O 19	/537. RY* (1) H 63	0.11	1.00	0.013
147. LP (2) O 19	/539. RY* (1) H 64	0.16	1.01	0.016
147. LP (2) O 19	/660. BD* (1) O 62 - H 63	0.03	0.77	0.006
147. LP (2) O 19	/661. BD* (1) O 62 - H 64	0.09	0.78	0.011
148. LP (3) O 19	/529. RY* (1) O 62	0.28	2.40	0.033
148. LP (3) O 19	/531. RY* (3) O 62	0.07	1.73	0.014
148. LP (3) O 19	/532. RY* (4) O 62	0.11	1.73	0.018
148. LP (3) O 19	/537. RY* (1) H 63	0.07	0.93	0.010
148. LP (3) O 19	/539. RY* (1) H 64	0.06	0.94	0.010
301. RY (1) C 17	/539. RY* (1) H 64	0.29	0.06	0.050
610. BD* (2) C 13 - C 15	/661. BD* (1) O 62 - H 64	0.08	0.30	0.013

from unit 1 to unit 4

11. BD (2) C 4 - C 6	/663. BD* (1) H 67 - O 68	0.04	0.59	0.006
13. BD (1) C 5 - C 6	/555. RY* (3) O 68	0.06	1.82	0.014
25. BD (1) N 11 - N 12	/553. RY* (1) O 68	0.04	2.91	0.013
26. BD (1) N 12 - C 13	/553. RY* (1) O 68	0.04	2.72	0.013
140. LP (1) N 11	/549. RY* (1) H 66	0.04	1.01	0.008
140. LP (1) N 11	/551. RY* (1) H 67	0.06	0.95	0.009
140. LP (1) N 11	/553. RY* (1) O 68	0.11	2.25	0.021
140. LP (1) N 11	/556. RY* (4) O 68	0.05	2.12	0.013
141. LP (2) N 11	/549. RY* (1) H 66	0.36	1.14	0.026
141. LP (2) N 11	/551. RY* (1) H 67	0.36	1.07	0.025
141. LP (2) N 11	/553. RY* (1) O 68	0.84	2.37	0.058
141. LP (2) N 11	/555. RY* (3) O 68	0.04	1.51	0.009
141. LP (2) N 11	/556. RY* (4) O 68	0.28	2.25	0.033

146. LP (1) O 19	/549. RY* (1) H 66	0.07	1.25	0.012
146. LP (1) O 19	/551. RY* (1) H 67	0.07	1.18	0.012
146. LP (1) O 19	/553. RY* (1) O 68	0.12	2.48	0.022
146. LP (1) O 19	/556. RY* (4) O 68	0.05	2.36	0.014
147. LP (2) O 19	/549. RY* (1) H 66	0.12	1.14	0.015
147. LP (2) O 19	/551. RY* (1) H 67	0.16	1.07	0.017
147. LP (2) O 19	/553. RY* (1) O 68	0.31	2.37	0.035
147. LP (2) O 19	/556. RY* (4) O 68	0.17	2.25	0.025
148. LP (3) O 19	/549. RY* (1) H 66	0.08	1.07	0.012
148. LP (3) O 19	/551. RY* (1) H 67	0.09	1.00	0.012
148. LP (3) O 19	/553. RY* (1) O 68	0.20	2.30	0.028
148. LP (3) O 19	/556. RY* (4) O 68	0.13	2.18	0.021
148. LP (3) O 19	/663. BD* (1) H 67 - O 68	0.03	0.71	0.006
from unit 1 to unit 5				
10. BD (1) C 4 - C 6	/166. LP (6)Mn 69	0.05	1.41	0.011
10. BD (1) C 4 - C 6	/167. LP* (7)Mn 69	0.05	1.19	0.010
11. BD (2) C 4 - C 6	/561. RY* (1)Mn 69	0.04	2.05	0.013
13. BD (1) C 5 - C 6	/167. LP* (7)Mn 69	0.05	1.30	0.010
15. BD (1) C 6 - N 11	/166. LP (6)Mn 69	0.73	1.56	0.045
15. BD (1) C 6 - N 11	/167. LP* (7)Mn 69	0.04	1.33	0.009
15. BD (1) C 6 - N 11	/168. LP* (8)Mn 69	0.09	2.62	0.019
15. BD (1) C 6 - N 11	/169. LP* (9)Mn 69	0.03	2.54	0.011
15. BD (1) C 6 - N 11	/564. RY* (4)Mn 69	0.05	2.98	0.016
15. BD (1) C 6 - N 11	/565. RY* (5)Mn 69	0.05	2.60	0.014
15. BD (1) C 6 - N 11	/567. RY* (7)Mn 69	0.03	2.80	0.012
15. BD (1) C 6 - N 11	/569. RY* (9)Mn 69	0.04	8.82	0.023
15. BD (1) C 6 - N 11	/572. RY* (12)Mn 69	0.03	5.97	0.016
25. BD (1) N 11 - N 12	/166. LP (6)Mn 69	0.79	1.75	0.050
25. BD (1) N 11 - N 12	/167. LP* (7)Mn 69	0.05	1.53	0.011
25. BD (1) N 11 - N 12	/168. LP* (8)Mn 69	0.03	2.82	0.011
25. BD (1) N 11 - N 12	/169. LP* (9)Mn 69	0.09	2.73	0.020
25. BD (1) N 11 - N 12	/561. RY* (1)Mn 69	0.06	2.78	0.016
25. BD (1) N 11 - N 12	/562. RY* (2)Mn 69	0.19	3.98	0.034
25. BD (1) N 11 - N 12	/569. RY* (9)Mn 69	0.06	9.02	0.030
25. BD (1) N 11 - N 12	/580. RY* (20)Mn 69	0.03	149.96	0.090
25. BD (1) N 11 - N 12	/581. RY* (21)Mn 69	0.06	39.36	0.059
26. BD (1) N 12 - C 13	/166. LP (6)Mn 69	0.11	1.56	0.018
26. BD (1) N 12 - C 13	/167. LP* (7)Mn 69	0.04	1.34	0.010
26. BD (1) N 12 - C 13	/561. RY* (1)Mn 69	0.04	2.59	0.012
26. BD (1) N 12 - C 13	/562. RY* (2)Mn 69	0.08	3.78	0.022
26. BD (1) N 12 - C 13	/568. RY* (8)Mn 69	0.04	7.48	0.023
26. BD (1) N 12 - C 13	/570. RY* (10)Mn 69	0.05	9.56	0.029
26. BD (1) N 12 - C 13	/577. RY* (17)Mn 69	0.03	28.06	0.034
26. BD (1) N 12 - C 13	/578. RY* (18)Mn 69	0.03	679.52	0.195
26. BD (1) N 12 - C 13	/580. RY* (20)Mn 69	0.07	149.76	0.126
26. BD (1) N 12 - C 13	/581. RY* (21)Mn 69	0.12	39.17	0.088
27. BD (1) C 13 - C 14	/166. LP (6)Mn 69	0.03	1.41	0.009
28. BD (1) C 13 - C 15	/166. LP (6)Mn 69	0.14	1.53	0.019
28. BD (1) C 13 - C 15	/167. LP* (7)Mn 69	0.07	1.30	0.012
28. BD (1) C 13 - C 15	/562. RY* (2)Mn 69	0.03	3.75	0.014
28. BD (1) C 13 - C 15	/581. RY* (21)Mn 69	0.04	39.14	0.047
29. BD (2) C 13 - C 15	/166. LP (6)Mn 69	0.03	1.07	0.008
32. BD (1) C 15 - C 17	/166. LP (6)Mn 69	0.20	1.41	0.023
32. BD (1) C 15 - C 17	/167. LP* (7)Mn 69	0.05	1.18	0.009
32. BD (1) C 15 - C 17	/562. RY* (2)Mn 69	0.05	3.63	0.016
32. BD (1) C 15 - C 17	/581. RY* (21)Mn 69	0.04	39.02	0.050
33. BD (1) C 15 - O 19	/166. LP (6)Mn 69	0.55	1.61	0.040
33. BD (1) C 15 - O 19	/168. LP* (8)Mn 69	0.11	2.67	0.021
33. BD (1) C 15 - O 19	/169. LP* (9)Mn 69	0.08	2.59	0.018
33. BD (1) C 15 - O 19	/569. RY* (9)Mn 69	0.03	8.87	0.021
33. BD (1) C 15 - O 19	/573. RY* (13)Mn 69	0.03	8.54	0.020
94. CR (1) C 6	/166. LP (6)Mn 69	0.03	10.85	0.024
99. CR (1) N 11	/166. LP (6)Mn 69	1.01	14.97	0.166
100. CR (1) N 12	/166. LP (6)Mn 69	0.03	15.01	0.030
103. CR (1) C 15	/166. LP (6)Mn 69	0.04	10.87	0.029
107. CR (1) O 19	/166. LP (6)Mn 69	1.20	19.67	0.208
140. LP (1) N 11	/166. LP (6)Mn 69	2.54	1.09	0.070
140. LP (1) N 11	/167. LP* (7)Mn 69	0.08	0.86	0.010
140. LP (1) N 11	/561. RY* (1)Mn 69	0.06	2.11	0.015
140. LP (1) N 11	/562. RY* (2)Mn 69	0.35	3.31	0.044

140. LP (1) N 11	/565. RY* (5) Mn 69	0.04	2.13	0.012
140. LP (1) N 11	/567. RY* (7) Mn 69	0.04	2.33	0.012
140. LP (1) N 11	/575. RY* (15) Mn 69	0.05	5.88	0.022
140. LP (1) N 11	/578. RY* (18) Mn 69	0.08	679.05	0.294
140. LP (1) N 11	/580. RY* (20) Mn 69	0.15	149.29	0.195
140. LP (1) N 11	/581. RY* (21) Mn 69	0.28	38.70	0.134
141. LP (2) N 11	/166. LP (6) Mn 69	16.68	1.21	0.188
141. LP (2) N 11	/167. LP* (7) Mn 69	0.23	0.99	0.020
141. LP (2) N 11	/168. LP* (8) Mn 69	0.25	2.28	0.031
141. LP (2) N 11	/169. LP* (9) Mn 69	0.04	2.19	0.011
141. LP (2) N 11	/561. RY* (1) Mn 69	0.54	2.24	0.045
141. LP (2) N 11	/562. RY* (2) Mn 69	2.43	3.44	0.118
141. LP (2) N 11	/564. RY* (4) Mn 69	0.11	2.64	0.022
141. LP (2) N 11	/565. RY* (5) Mn 69	0.17	2.26	0.025
141. LP (2) N 11	/567. RY* (7) Mn 69	0.28	2.46	0.034
141. LP (2) N 11	/568. RY* (8) Mn 69	0.07	7.13	0.028
141. LP (2) N 11	/574. RY* (14) Mn 69	0.05	6.37	0.022
141. LP (2) N 11	/575. RY* (15) Mn 69	0.20	6.00	0.045
141. LP (2) N 11	/578. RY* (18) Mn 69	0.56	679.18	0.796
141. LP (2) N 11	/580. RY* (20) Mn 69	1.11	149.42	0.528
141. LP (2) N 11	/581. RY* (21) Mn 69	2.07	38.82	0.367
142. LP (1) N 12	/166. LP (6) Mn 69	0.16	1.20	0.019
142. LP (1) N 12	/561. RY* (1) Mn 69	0.08	2.23	0.017
146. LP (1) O 19	/166. LP (6) Mn 69	4.01	1.32	0.098
146. LP (1) O 19	/167. LP* (7) Mn 69	0.08	1.10	0.012
146. LP (1) O 19	/169. LP* (9) Mn 69	0.06	2.30	0.015
146. LP (1) O 19	/561. RY* (1) Mn 69	0.05	2.35	0.013
146. LP (1) O 19	/562. RY* (2) Mn 69	0.54	3.55	0.056
146. LP (1) O 19	/563. RY* (3) Mn 69	0.05	2.32	0.014
146. LP (1) O 19	/564. RY* (4) Mn 69	0.04	2.75	0.013
146. LP (1) O 19	/567. RY* (7) Mn 69	0.07	2.57	0.017
146. LP (1) O 19	/569. RY* (9) Mn 69	0.03	8.59	0.020
146. LP (1) O 19	/575. RY* (15) Mn 69	0.06	6.11	0.024
146. LP (1) O 19	/578. RY* (18) Mn 69	0.12	679.28	0.368
146. LP (1) O 19	/580. RY* (20) Mn 69	0.21	149.52	0.225
146. LP (1) O 19	/581. RY* (21) Mn 69	0.38	38.93	0.154
147. LP (2) O 19	/166. LP (6) Mn 69	9.25	1.21	0.141
147. LP (2) O 19	/167. LP* (7) Mn 69	0.66	0.99	0.033
147. LP (2) O 19	/168. LP* (8) Mn 69	0.08	2.27	0.017
147. LP (2) O 19	/169. LP* (9) Mn 69	0.32	2.19	0.034
147. LP (2) O 19	/562. RY* (2) Mn 69	1.08	3.43	0.078
147. LP (2) O 19	/563. RY* (3) Mn 69	0.05	2.21	0.014
147. LP (2) O 19	/564. RY* (4) Mn 69	0.05	2.64	0.014
147. LP (2) O 19	/567. RY* (7) Mn 69	0.22	2.46	0.030
147. LP (2) O 19	/575. RY* (15) Mn 69	0.12	6.00	0.035
147. LP (2) O 19	/578. RY* (18) Mn 69	0.28	679.17	0.560
147. LP (2) O 19	/580. RY* (20) Mn 69	0.54	149.41	0.365
147. LP (2) O 19	/581. RY* (21) Mn 69	1.02	38.82	0.255
148. LP (3) O 19	/166. LP (6) Mn 69	8.27	1.14	0.128
148. LP (3) O 19	/167. LP* (7) Mn 69	0.61	0.92	0.030
148. LP (3) O 19	/168. LP* (8) Mn 69	0.15	2.20	0.023
148. LP (3) O 19	/169. LP* (9) Mn 69	0.24	2.12	0.029
148. LP (3) O 19	/561. RY* (1) Mn 69	0.13	2.17	0.022
148. LP (3) O 19	/562. RY* (2) Mn 69	0.81	3.36	0.068
148. LP (3) O 19	/563. RY* (3) Mn 69	0.03	2.14	0.010
148. LP (3) O 19	/564. RY* (4) Mn 69	0.03	2.57	0.012
148. LP (3) O 19	/567. RY* (7) Mn 69	0.16	2.39	0.025
148. LP (3) O 19	/575. RY* (15) Mn 69	0.09	5.93	0.030
148. LP (3) O 19	/578. RY* (18) Mn 69	0.20	679.10	0.480
148. LP (3) O 19	/580. RY* (20) Mn 69	0.41	149.34	0.319
148. LP (3) O 19	/581. RY* (21) Mn 69	0.77	38.75	0.224
277. RY (1) C 14	/166. LP (6) Mn 69	0.05	0.27	0.012
277. RY (1) C 14	/167. LP* (7) Mn 69	0.03	0.05	0.014
301. RY (1) C 17	/166. LP (6) Mn 69	0.27	0.26	0.029
301. RY (1) C 17	/167. LP* (7) Mn 69	0.10	0.04	0.022

from unit 2 to unit 1

71. BD (1) C 44 - C 46	/319. RY* (3) O 19	0.09	2.10	0.017
72. BD (2) C 44 - C 46	/319. RY* (3) O 19	0.05	1.62	0.012
72. BD (2) C 44 - C 46	/614. BD* (1) C 15 - O 19	0.08	0.61	0.009
76. BD (1) C 46 - C 48	/237. RY* (1) C 9	0.04	1.38	0.009

77. BD (1) C 46 - O 65	/214. RY* (2) C 6	0.06	2.00	0.014
151. LP (1) N 42	/614. BD* (1) C 15 - O 19	0.04	0.62	0.006
152. LP (2) N 42	/214. RY* (2) C 6	0.13	1.61	0.019
152. LP (2) N 42	/216. RY* (4) C 6	0.08	1.49	0.015
152. LP (2) N 42	/218. RY* (6) C 6	0.05	3.66	0.017
152. LP (2) N 42	/253. RY* (1) N 11	0.24	1.80	0.027
152. LP (2) N 42	/254. RY* (2) N 11	0.23	1.85	0.027
152. LP (2) N 42	/255. RY* (3) N 11	0.15	1.71	0.021
152. LP (2) N 42	/256. RY* (4) N 11	0.07	1.98	0.015
152. LP (2) N 42	/285. RY* (1) C 15	0.17	1.61	0.021
152. LP (2) N 42	/286. RY* (2) C 15	0.03	1.21	0.008
152. LP (2) N 42	/289. RY* (5) C 15	0.08	3.89	0.022
152. LP (2) N 42	/317. RY* (1) O 19	0.53	2.82	0.050
152. LP (2) N 42	/319. RY* (3) O 19	0.10	1.78	0.018
152. LP (2) N 42	/320. RY* (4) O 19	0.04	2.25	0.012
152. LP (2) N 42	/614. BD* (1) C 15 - O 19	0.15	0.77	0.014
159. LP (1) O 65	/253. RY* (1) N 11	0.12	1.93	0.019
159. LP (1) O 65	/254. RY* (2) N 11	0.10	1.98	0.018
159. LP (1) O 65	/255. RY* (3) N 11	0.10	1.84	0.017
159. LP (1) O 65	/256. RY* (4) N 11	0.05	2.10	0.013
159. LP (1) O 65	/285. RY* (1) C 15	0.05	1.74	0.011
159. LP (1) O 65	/317. RY* (1) O 19	0.13	2.94	0.025
159. LP (1) O 65	/603. BD* (2) C 9 - C 10	0.10	0.61	0.012
159. LP (1) O 65	/606. BD* (1) N 11 - N 12	0.03	1.06	0.007
160. LP (2) O 65	/214. RY* (2) C 6	0.08	1.62	0.014
160. LP (2) O 65	/253. RY* (1) N 11	0.17	1.82	0.022
160. LP (2) O 65	/254. RY* (2) N 11	0.24	1.87	0.027
160. LP (2) O 65	/255. RY* (3) N 11	0.13	1.73	0.019
160. LP (2) O 65	/256. RY* (4) N 11	0.03	1.99	0.010
160. LP (2) O 65	/263. RY* (3) N 12	0.03	1.57	0.009
160. LP (2) O 65	/285. RY* (1) C 15	0.13	1.63	0.019
160. LP (2) O 65	/286. RY* (2) C 15	0.03	1.22	0.008
160. LP (2) O 65	/289. RY* (5) C 15	0.05	3.90	0.018
160. LP (2) O 65	/317. RY* (1) O 19	0.26	2.83	0.034
160. LP (2) O 65	/319. RY* (3) O 19	0.05	1.79	0.012
160. LP (2) O 65	/320. RY* (4) O 19	0.03	2.27	0.010
160. LP (2) O 65	/606. BD* (1) N 11 - N 12	0.03	0.95	0.007
161. LP (3) O 65	/214. RY* (2) C 6	0.03	1.47	0.008
161. LP (3) O 65	/253. RY* (1) N 11	0.06	1.66	0.013
161. LP (3) O 65	/254. RY* (2) N 11	0.08	1.72	0.016
161. LP (3) O 65	/255. RY* (3) N 11	0.09	1.58	0.015
161. LP (3) O 65	/256. RY* (4) N 11	0.03	1.84	0.009
161. LP (3) O 65	/317. RY* (1) O 19	0.07	2.68	0.018
161. LP (3) O 65	/596. BD* (1) C 6 - N 11	0.03	0.68	0.006
161. LP (3) O 65	/606. BD* (1) N 11 - N 12	0.05	0.80	0.008
161. LP (3) O 65	/614. BD* (1) C 15 - O 19	0.08	0.63	0.009

within unit 2

42. BD (1) C 32 - C 33	/372. RY* (2) C 34	0.55	2.00	0.042
42. BD (1) C 32 - C 33	/373. RY* (3) C 34	0.27	2.01	0.030
42. BD (1) C 32 - C 33	/380. RY* (2) C 35	0.72	2.25	0.051
42. BD (1) C 32 - C 33	/404. RY* (2) C 38	0.65	1.88	0.044
42. BD (1) C 32 - C 33	/623. BD* (1) C 32 - C 34	1.49	1.26	0.055
42. BD (1) C 32 - C 33	/625. BD* (1) C 33 - C 35	1.24	1.27	0.050
42. BD (1) C 32 - C 33	/626. BD* (1) C 33 - C 38	0.61	1.10	0.033
42. BD (1) C 32 - C 33	/628. BD* (1) C 34 - H 51	1.36	1.08	0.049
42. BD (1) C 32 - C 33	/630. BD* (1) C 35 - C 40	2.08	1.11	0.061
42. BD (1) C 32 - C 33	/634. BD* (1) C 38 - C 39	0.69	1.30	0.038
43. BD (2) C 32 - C 33	/ 50. BD* (2) C 34 - C 36	5.37	0.22	0.064
43. BD (2) C 32 - C 33	/ 54. BD* (2) C 35 - C 37	6.26	0.24	0.065
43. BD (2) C 32 - C 33	/371. RY* (1) C 34	1.12	0.77	0.038
43. BD (2) C 32 - C 33	/379. RY* (1) C 35	0.89	0.87	0.036
43. BD (2) C 32 - C 33	/403. RY* (1) C 38	0.59	0.81	0.028
43. BD (2) C 32 - C 33	/499. RY* (1) O 50	0.35	1.15	0.026
43. BD (2) C 32 - C 33	/622. BD* (2) C 32 - C 33	0.37	0.23	0.015
43. BD (2) C 32 - C 33	/635. BD* (2) C 38 - C 39	6.49	0.30	0.065
43. BD (2) C 32 - C 33	/640. BD* (2) C 40 - C 41	0.26	0.30	0.013
44. BD (1) C 32 - C 34	/365. RY* (3) C 33	0.69	2.22	0.049
44. BD (1) C 32 - C 34	/389. RY* (3) C 36	0.57	1.99	0.043
44. BD (1) C 32 - C 34	/621. BD* (1) C 32 - C 33	1.47	1.19	0.053
44. BD (1) C 32 - C 34	/624. BD* (1) C 32 - O 50	0.28	1.04	0.022

44. BD (1) C 32 - C 34	/626. BD* (1) C 33 - C 38	1.12	1.22	0.047
44. BD (1) C 32 - C 34	/627. BD* (1) C 34 - C 36	0.77	1.23	0.039
44. BD (1) C 32 - C 34	/628. BD* (1) C 34 - H 51	0.61	1.20	0.034
44. BD (1) C 32 - C 34	/632. BD* (1) C 36 - H 61	0.67	1.21	0.036
44. BD (1) C 32 - C 34	/659. BD* (1) O 50 - H 60	0.73	1.08	0.036
45. BD (1) C 32 - O 50	/372. RY* (2) C 34	0.28	2.24	0.032
45. BD (1) C 32 - O 50	/623. BD* (1) C 32 - C 34	0.38	1.51	0.030
45. BD (1) C 32 - O 50	/625. BD* (1) C 33 - C 35	0.56	1.52	0.037
45. BD (1) C 32 - O 50	/627. BD* (1) C 34 - C 36	1.00	1.36	0.047
46. BD (1) C 33 - C 35	/357. RY* (3) C 32	0.44	2.18	0.039
46. BD (1) C 33 - C 35	/395. RY* (1) C 37	0.27	1.65	0.027
46. BD (1) C 33 - C 35	/396. RY* (2) C 37	0.41	1.84	0.035
46. BD (1) C 33 - C 35	/404. RY* (2) C 38	0.39	1.98	0.035
46. BD (1) C 33 - C 35	/420. RY* (2) C 40	0.36	1.91	0.033
46. BD (1) C 33 - C 35	/621. BD* (1) C 32 - C 33	1.23	1.16	0.048
46. BD (1) C 33 - C 35	/624. BD* (1) C 32 - O 50	1.32	1.02	0.046
46. BD (1) C 33 - C 35	/626. BD* (1) C 33 - C 38	1.62	1.19	0.056
46. BD (1) C 33 - C 35	/629. BD* (1) C 35 - C 37	1.54	1.20	0.054
46. BD (1) C 33 - C 35	/630. BD* (1) C 35 - C 40	1.54	1.21	0.054
46. BD (1) C 33 - C 35	/633. BD* (1) C 37 - N 42	1.35	1.13	0.049
46. BD (1) C 33 - C 35	/636. BD* (1) C 38 - H 52	0.76	1.20	0.038
46. BD (1) C 33 - C 35	/641. BD* (1) C 40 - H 54	0.72	1.20	0.037
47. BD (1) C 33 - C 38	/355. RY* (1) C 32	0.34	1.54	0.029
47. BD (1) C 33 - C 38	/357. RY* (3) C 32	0.34	2.07	0.034
47. BD (1) C 33 - C 38	/380. RY* (2) C 35	0.45	2.23	0.040
47. BD (1) C 33 - C 38	/412. RY* (2) C 39	0.31	1.99	0.031
47. BD (1) C 33 - C 38	/413. RY* (3) C 39	0.71	2.00	0.048
47. BD (1) C 33 - C 38	/621. BD* (1) C 32 - C 33	0.49	1.05	0.029
47. BD (1) C 33 - C 38	/623. BD* (1) C 32 - C 34	0.75	1.24	0.039
47. BD (1) C 33 - C 38	/625. BD* (1) C 33 - C 35	1.80	1.25	0.060
47. BD (1) C 33 - C 38	/629. BD* (1) C 35 - C 37	2.42	1.08	0.065
47. BD (1) C 33 - C 38	/634. BD* (1) C 38 - C 39	1.26	1.28	0.051
47. BD (1) C 33 - C 38	/636. BD* (1) C 38 - H 52	0.30	1.08	0.023
47. BD (1) C 33 - C 38	/638. BD* (1) C 39 - H 53	1.37	1.08	0.049
48. BD (1) C 34 - C 36	/355. RY* (1) C 32	0.36	1.57	0.030
48. BD (1) C 34 - C 36	/356. RY* (2) C 32	0.39	1.43	0.030
48. BD (1) C 34 - C 36	/357. RY* (3) C 32	0.36	2.09	0.035
48. BD (1) C 34 - C 36	/397. RY* (3) C 37	0.95	1.79	0.052
48. BD (1) C 34 - C 36	/623. BD* (1) C 32 - C 34	0.84	1.26	0.041
48. BD (1) C 34 - C 36	/624. BD* (1) C 32 - O 50	2.66	0.92	0.063
48. BD (1) C 34 - C 36	/631. BD* (1) C 36 - C 37	1.08	1.29	0.047
48. BD (1) C 34 - C 36	/633. BD* (1) C 37 - N 42	2.85	1.04	0.069
49. BD (2) C 34 - C 36	/ 54. BD* (2) C 35 - C 37	6.75	0.24	0.068
49. BD (2) C 34 - C 36	/356. RY* (2) C 32	0.26	1.06	0.022
49. BD (2) C 34 - C 36	/395. RY* (1) C 37	0.35	1.18	0.026
49. BD (2) C 34 - C 36	/622. BD* (2) C 32 - C 33	7.76	0.23	0.071
50. BD* (2) C 34 - C 36	/ 54. BD* (2) C 35 - C 37	338.46	0.02	0.094
50. BD* (2) C 34 - C 36	/371. RY* (1) C 34	2.02	0.54	0.054
50. BD* (2) C 34 - C 36	/387. RY* (1) C 36	1.69	0.59	0.052
51. BD (1) C 34 - H 51	/357. RY* (3) C 32	0.48	1.95	0.039
51. BD (1) C 34 - H 51	/360. RY* (6) C 32	0.29	8.50	0.063
51. BD (1) C 34 - H 51	/388. RY* (2) C 36	0.49	1.78	0.037
51. BD (1) C 34 - H 51	/621. BD* (1) C 32 - C 33	3.08	0.94	0.068
51. BD (1) C 34 - H 51	/623. BD* (1) C 32 - C 34	0.41	1.13	0.027
51. BD (1) C 34 - H 51	/624. BD* (1) C 32 - O 50	0.94	0.79	0.034
51. BD (1) C 34 - H 51	/631. BD* (1) C 36 - C 37	1.42	1.15	0.051
51. BD (1) C 34 - H 51	/632. BD* (1) C 36 - H 61	0.29	0.96	0.021
52. BD (1) C 35 - C 37	/364. RY* (2) C 33	0.84	2.34	0.056
52. BD (1) C 35 - C 37	/388. RY* (2) C 36	0.50	1.91	0.039
52. BD (1) C 35 - C 37	/420. RY* (2) C 40	0.62	1.81	0.043
52. BD (1) C 35 - C 37	/436. RY* (2) N 42	0.49	1.60	0.035
52. BD (1) C 35 - C 37	/625. BD* (1) C 33 - C 35	1.49	1.27	0.055
52. BD (1) C 35 - C 37	/626. BD* (1) C 33 - C 38	2.00	1.10	0.059
52. BD (1) C 35 - C 37	/630. BD* (1) C 35 - C 40	0.78	1.11	0.037
52. BD (1) C 35 - C 37	/631. BD* (1) C 36 - C 37	1.59	1.28	0.057
52. BD (1) C 35 - C 37	/632. BD* (1) C 36 - H 61	1.47	1.09	0.051
52. BD (1) C 35 - C 37	/633. BD* (1) C 37 - N 42	0.29	1.03	0.022
52. BD (1) C 35 - C 37	/639. BD* (1) C 40 - C 41	0.74	1.31	0.039
53. BD (2) C 35 - C 37	/ 50. BD* (2) C 34 - C 36	5.84	0.22	0.066
53. BD (2) C 35 - C 37	/363. RY* (1) C 33	0.93	0.86	0.037
53. BD (2) C 35 - C 37	/387. RY* (1) C 36	1.25	0.81	0.041

53. BD (2) C 35 - C 37	/419. RY* (1) C 40	0.48	0.91	0.027
53. BD (2) C 35 - C 37	/622. BD* (2) C 32 - C 33	6.25	0.22	0.062
53. BD (2) C 35 - C 37	/640. BD* (2) C 40 - C 41	6.59	0.30	0.065
53. BD (2) C 35 - C 37	/643. BD* (1) N 42 - N 43	2.04	0.79	0.052
54. BD* (2) C 35 - C 37	/363. RY* (1) C 33	0.71	0.62	0.038
54. BD* (2) C 35 - C 37	/379. RY* (1) C 35	1.18	0.63	0.049
54. BD* (2) C 35 - C 37	/387. RY* (1) C 36	1.19	0.58	0.047
54. BD* (2) C 35 - C 37	/395. RY* (1) C 37	0.45	0.94	0.037
54. BD* (2) C 35 - C 37	/396. RY* (2) C 37	0.27	1.14	0.031
54. BD* (2) C 35 - C 37	/419. RY* (1) C 40	0.33	0.67	0.027
54. BD* (2) C 35 - C 37	/635. BD* (2) C 38 - C 39	0.27	0.06	0.006
54. BD* (2) C 35 - C 37	/640. BD* (2) C 40 - C 41	49.04	0.06	0.078
54. BD* (2) C 35 - C 37	/643. BD* (1) N 42 - N 43	2.12	0.55	0.061
55. BD (1) C 35 - C 40	/364. RY* (2) C 33	0.45	2.32	0.041
55. BD (1) C 35 - C 40	/395. RY* (1) C 37	0.55	1.53	0.037
55. BD (1) C 35 - C 40	/396. RY* (2) C 37	0.68	1.73	0.044
55. BD (1) C 35 - C 40	/428. RY* (2) C 41	0.34	2.00	0.033
55. BD (1) C 35 - C 40	/429. RY* (3) C 41	0.75	2.00	0.049
55. BD (1) C 35 - C 40	/621. BD* (1) C 32 - C 33	2.38	1.05	0.063
55. BD (1) C 35 - C 40	/625. BD* (1) C 33 - C 35	1.76	1.25	0.059
55. BD (1) C 35 - C 40	/629. BD* (1) C 35 - C 37	0.79	1.08	0.037
55. BD (1) C 35 - C 40	/631. BD* (1) C 36 - C 37	1.03	1.27	0.046
55. BD (1) C 35 - C 40	/639. BD* (1) C 40 - C 41	1.34	1.29	0.053
55. BD (1) C 35 - C 40	/641. BD* (1) C 40 - H 54	0.32	1.09	0.024
55. BD (1) C 35 - C 40	/642. BD* (1) C 41 - H 55	1.43	1.08	0.050
56. BD (1) C 36 - C 37	/373. RY* (3) C 34	0.65	2.11	0.047
56. BD (1) C 36 - C 37	/381. RY* (3) C 35	0.71	2.33	0.052
56. BD (1) C 36 - C 37	/436. RY* (2) N 42	0.41	1.71	0.034
56. BD (1) C 36 - C 37	/627. BD* (1) C 34 - C 36	0.97	1.22	0.044
56. BD (1) C 36 - C 37	/628. BD* (1) C 34 - H 51	0.60	1.19	0.034
56. BD (1) C 36 - C 37	/629. BD* (1) C 35 - C 37	1.64	1.21	0.057
56. BD (1) C 36 - C 37	/630. BD* (1) C 35 - C 40	1.11	1.22	0.047
56. BD (1) C 36 - C 37	/632. BD* (1) C 36 - H 61	0.77	1.20	0.038
56. BD (1) C 36 - C 37	/633. BD* (1) C 37 - N 42	0.75	1.15	0.037
56. BD (1) C 36 - C 37	/643. BD* (1) N 42 - N 43	0.48	1.27	0.031
57. BD (1) C 36 - H 61	/372. RY* (2) C 34	0.47	1.86	0.038
57. BD (1) C 36 - H 61	/395. RY* (1) C 37	0.45	1.42	0.032
57. BD (1) C 36 - H 61	/396. RY* (2) C 37	0.70	1.61	0.043
57. BD (1) C 36 - H 61	/398. RY* (4) C 37	0.28	1.56	0.027
57. BD (1) C 36 - H 61	/623. BD* (1) C 32 - C 34	1.33	1.12	0.049
57. BD (1) C 36 - H 61	/628. BD* (1) C 34 - H 51	0.28	0.95	0.020
57. BD (1) C 36 - H 61	/629. BD* (1) C 35 - C 37	3.32	0.97	0.072
57. BD (1) C 36 - H 61	/631. BD* (1) C 36 - C 37	0.68	1.15	0.035
57. BD (1) C 36 - H 61	/633. BD* (1) C 37 - N 42	0.67	0.90	0.031
58. BD (1) C 37 - N 42	/381. RY* (3) C 35	0.62	2.34	0.048
58. BD (1) C 37 - N 42	/388. RY* (2) C 36	0.59	2.03	0.044
58. BD (1) C 37 - N 42	/443. RY* (1) N 43	0.90	1.65	0.049
58. BD (1) C 37 - N 42	/625. BD* (1) C 33 - C 35	0.78	1.39	0.042
58. BD (1) C 37 - N 42	/627. BD* (1) C 34 - C 36	1.27	1.23	0.050
58. BD (1) C 37 - N 42	/629. BD* (1) C 35 - C 37	0.51	1.22	0.032
58. BD (1) C 37 - N 42	/631. BD* (1) C 36 - C 37	0.93	1.41	0.046
58. BD (1) C 37 - N 42	/643. BD* (1) N 42 - N 43	0.48	1.29	0.032
58. BD (1) C 37 - N 42	/644. BD* (1) N 43 - C 44	2.84	1.15	0.072
59. BD (1) C 38 - C 39	/364. RY* (2) C 33	0.47	2.43	0.043
59. BD (1) C 38 - C 39	/365. RY* (3) C 33	0.46	2.19	0.040
59. BD (1) C 38 - C 39	/428. RY* (2) C 41	0.41	2.11	0.037
59. BD (1) C 38 - C 39	/429. RY* (3) C 41	0.40	2.11	0.037
59. BD (1) C 38 - C 39	/621. BD* (1) C 32 - C 33	1.28	1.16	0.049
59. BD (1) C 38 - C 39	/626. BD* (1) C 33 - C 38	1.36	1.19	0.051
59. BD (1) C 38 - C 39	/636. BD* (1) C 38 - H 52	0.81	1.19	0.039
59. BD (1) C 38 - C 39	/637. BD* (1) C 39 - C 41	0.95	1.21	0.043
59. BD (1) C 38 - C 39	/638. BD* (1) C 39 - H 53	0.67	1.19	0.036
59. BD (1) C 38 - C 39	/642. BD* (1) C 41 - H 55	0.77	1.19	0.038
60. BD (2) C 38 - C 39	/363. RY* (1) C 33	0.41	0.86	0.024
60. BD (2) C 38 - C 39	/427. RY* (1) C 41	0.69	0.78	0.030
60. BD (2) C 38 - C 39	/622. BD* (2) C 32 - C 33	3.64	0.23	0.049
60. BD (2) C 38 - C 39	/640. BD* (2) C 40 - C 41	4.34	0.30	0.054
61. BD (1) C 38 - H 52	/364. RY* (2) C 33	0.67	2.20	0.049
61. BD (1) C 38 - H 52	/412. RY* (2) C 39	0.72	1.87	0.047
61. BD (1) C 38 - H 52	/625. BD* (1) C 33 - C 35	1.60	1.13	0.054
61. BD (1) C 38 - H 52	/634. BD* (1) C 38 - C 39	0.67	1.16	0.035

61. BD (1) C 38 - H 52	/637. BD* (1) C 39 - C 41	2.62	0.98	0.064
61. BD (1) C 38 - H 52	/638. BD* (1) C 39 - H 53	0.46	0.96	0.027
62. BD (1) C 39 - C 41	/405. RY* (3) C 38	0.93	1.78	0.052
62. BD (1) C 39 - C 41	/421. RY* (3) C 40	0.92	1.76	0.051
62. BD (1) C 39 - C 41	/634. BD* (1) C 38 - C 39	1.06	1.29	0.047
62. BD (1) C 39 - C 41	/636. BD* (1) C 38 - H 52	1.65	1.09	0.054
62. BD (1) C 39 - C 41	/639. BD* (1) C 40 - C 41	1.13	1.30	0.049
62. BD (1) C 39 - C 41	/641. BD* (1) C 40 - H 54	1.74	1.10	0.055
63. BD (1) C 39 - H 53	/404. RY* (2) C 38	0.68	1.74	0.044
63. BD (1) C 39 - H 53	/428. RY* (2) C 41	0.58	1.88	0.042
63. BD (1) C 39 - H 53	/626. BD* (1) C 33 - C 38	3.21	0.95	0.070
63. BD (1) C 39 - H 53	/634. BD* (1) C 38 - C 39	0.58	1.16	0.033
63. BD (1) C 39 - H 53	/636. BD* (1) C 38 - H 52	0.56	0.96	0.029
63. BD (1) C 39 - H 53	/639. BD* (1) C 40 - C 41	1.32	1.17	0.050
63. BD (1) C 39 - H 53	/642. BD* (1) C 41 - H 55	0.26	0.96	0.020
64. BD (1) C 40 - C 41	/380. RY* (2) C 35	0.64	2.33	0.049
64. BD (1) C 40 - C 41	/381. RY* (3) C 35	0.35	2.31	0.036
64. BD (1) C 40 - C 41	/412. RY* (2) C 39	0.40	2.09	0.037
64. BD (1) C 40 - C 41	/413. RY* (3) C 39	0.38	2.10	0.036
64. BD (1) C 40 - C 41	/629. BD* (1) C 35 - C 37	1.21	1.19	0.048
64. BD (1) C 40 - C 41	/630. BD* (1) C 35 - C 40	1.40	1.20	0.052
64. BD (1) C 40 - C 41	/637. BD* (1) C 39 - C 41	1.01	1.21	0.044
64. BD (1) C 40 - C 41	/638. BD* (1) C 39 - H 53	0.75	1.19	0.038
64. BD (1) C 40 - C 41	/641. BD* (1) C 40 - H 54	0.85	1.20	0.040
64. BD (1) C 40 - C 41	/642. BD* (1) C 41 - H 55	0.74	1.19	0.038
65. BD (2) C 40 - C 41	/ 54. BD* (2) C 35 - C 37	3.68	0.24	0.051
65. BD (2) C 40 - C 41	/379. RY* (1) C 35	0.38	0.87	0.023
65. BD (2) C 40 - C 41	/411. RY* (1) C 39	0.66	0.77	0.029
65. BD (2) C 40 - C 41	/635. BD* (2) C 38 - C 39	4.63	0.30	0.055
65. BD (2) C 40 - C 41	/640. BD* (2) C 40 - C 41	0.30	0.30	0.014
66. BD (1) C 40 - H 54	/380. RY* (2) C 35	0.70	2.10	0.048
66. BD (1) C 40 - H 54	/428. RY* (2) C 41	0.70	1.88	0.046
66. BD (1) C 40 - H 54	/625. BD* (1) C 33 - C 35	1.74	1.13	0.056
66. BD (1) C 40 - H 54	/637. BD* (1) C 39 - C 41	2.62	0.98	0.064
66. BD (1) C 40 - H 54	/639. BD* (1) C 40 - C 41	0.75	1.17	0.037
66. BD (1) C 40 - H 54	/642. BD* (1) C 41 - H 55	0.46	0.96	0.027
67. BD (1) C 41 - H 55	/412. RY* (2) C 39	0.55	1.86	0.041
67. BD (1) C 41 - H 55	/420. RY* (2) C 40	0.63	1.66	0.041
67. BD (1) C 41 - H 55	/630. BD* (1) C 35 - C 40	3.21	0.96	0.070
67. BD (1) C 41 - H 55	/634. BD* (1) C 38 - C 39	1.32	1.15	0.049
67. BD (1) C 41 - H 55	/638. BD* (1) C 39 - H 53	0.26	0.95	0.020
67. BD (1) C 41 - H 55	/639. BD* (1) C 40 - C 41	0.66	1.16	0.035
67. BD (1) C 41 - H 55	/641. BD* (1) C 40 - H 54	0.50	0.96	0.028
68. BD (1) N 42 - N 43	/395. RY* (1) C 37	0.79	1.87	0.049
68. BD (1) N 42 - N 43	/451. RY* (1) C 44	0.73	2.10	0.049
68. BD (1) N 42 - N 43	/452. RY* (2) C 44	0.86	2.37	0.057
68. BD (1) N 42 - N 43	/631. BD* (1) C 36 - C 37	0.28	1.60	0.027
68. BD (1) N 42 - N 43	/633. BD* (1) C 37 - N 42	0.50	1.35	0.033
68. BD (1) N 42 - N 43	/644. BD* (1) N 43 - C 44	0.34	1.34	0.027
68. BD (1) N 42 - N 43	/645. BD* (1) C 44 - C 45	0.56	1.42	0.036
69. BD (1) N 43 - C 44	/395. RY* (1) C 37	0.28	1.67	0.028
69. BD (1) N 43 - C 44	/435. RY* (1) N 42	0.77	2.67	0.057
69. BD (1) N 43 - C 44	/437. RY* (3) N 42	0.50	2.00	0.040
69. BD (1) N 43 - C 44	/438. RY* (4) N 42	0.28	2.27	0.032
69. BD (1) N 43 - C 44	/459. RY* (1) C 45	0.40	2.05	0.036
69. BD (1) N 43 - C 44	/467. RY* (1) C 46	0.45	1.94	0.037
69. BD (1) N 43 - C 44	/633. BD* (1) C 37 - N 42	2.40	1.16	0.067
69. BD (1) N 43 - C 44	/645. BD* (1) C 44 - C 45	0.33	1.22	0.025
69. BD (1) N 43 - C 44	/646. BD* (1) C 44 - C 46	1.06	1.38	0.049
69. BD (1) N 43 - C 44	/648. BD* (1) C 45 - C 47	0.49	1.42	0.033
69. BD (1) N 43 - C 44	/651. BD* (1) C 46 - C 48	1.29	1.23	0.050
70. BD (1) C 44 - C 45	/443. RY* (1) N 43	0.34	1.52	0.029
70. BD (1) C 44 - C 45	/468. RY* (2) C 46	0.40	2.10	0.037
70. BD (1) C 44 - C 45	/475. RY* (1) C 47	0.41	1.85	0.035
70. BD (1) C 44 - C 45	/476. RY* (2) C 47	0.74	1.98	0.049
70. BD (1) C 44 - C 45	/643. BD* (1) N 42 - N 43	1.58	1.16	0.054
70. BD (1) C 44 - C 45	/644. BD* (1) N 43 - C 44	0.46	1.02	0.027
70. BD (1) C 44 - C 45	/646. BD* (1) C 44 - C 46	1.77	1.25	0.060
70. BD (1) C 44 - C 45	/648. BD* (1) C 45 - C 47	1.23	1.30	0.051
70. BD (1) C 44 - C 45	/650. BD* (1) C 45 - H 56	0.30	1.09	0.023
70. BD (1) C 44 - C 45	/652. BD* (1) C 46 - O 65	2.61	0.98	0.064

70. BD (1) C 44 - C 45	/654. BD* (1) C 47 - H 57	1.43	1.10	0.050
71. BD (1) C 44 - C 46	/445. RY* (3) N 43	0.31	1.86	0.030
71. BD (1) C 44 - C 46	/459. RY* (1) C 45	0.49	2.04	0.040
71. BD (1) C 44 - C 46	/483. RY* (1) C 48	0.70	2.10	0.048
71. BD (1) C 44 - C 46	/644. BD* (1) N 43 - C 44	0.74	1.14	0.037
71. BD (1) C 44 - C 46	/645. BD* (1) C 44 - C 45	1.79	1.21	0.059
71. BD (1) C 44 - C 46	/650. BD* (1) C 45 - H 56	0.65	1.21	0.035
71. BD (1) C 44 - C 46	/651. BD* (1) C 46 - C 48	1.53	1.22	0.055
71. BD (1) C 44 - C 46	/652. BD* (1) C 46 - O 65	0.38	1.09	0.026
71. BD (1) C 44 - C 46	/657. BD* (1) C 48 - H 58	0.62	1.22	0.035
72. BD (2) C 44 - C 46	/444. RY* (2) N 43	0.44	1.10	0.030
72. BD (2) C 44 - C 46	/461. RY* (3) C 45	0.52	0.83	0.028
72. BD (2) C 44 - C 46	/485. RY* (3) C 48	0.53	0.84	0.028
72. BD (2) C 44 - C 46	/542. RY* (2) O 65	0.28	1.37	0.026
72. BD (2) C 44 - C 46	/649. BD* (2) C 45 - C 47	6.40	0.33	0.058
72. BD (2) C 44 - C 46	/656. BD* (2) C 48 - C 49	7.03	0.33	0.061
73. BD (1) C 45 - C 47	/452. RY* (2) C 44	0.95	2.14	0.057
73. BD (1) C 45 - C 47	/491. RY* (1) C 49	0.51	1.94	0.040
73. BD (1) C 45 - C 47	/492. RY* (2) C 49	0.42	2.09	0.037
73. BD (1) C 45 - C 47	/644. BD* (1) N 43 - C 44	1.31	1.12	0.048
73. BD (1) C 45 - C 47	/645. BD* (1) C 44 - C 45	1.16	1.19	0.047
73. BD (1) C 45 - C 47	/650. BD* (1) C 45 - H 56	0.86	1.19	0.041
73. BD (1) C 45 - C 47	/653. BD* (1) C 47 - C 49	1.14	1.21	0.047
73. BD (1) C 45 - C 47	/654. BD* (1) C 47 - H 57	0.76	1.20	0.038
73. BD (1) C 45 - C 47	/658. BD* (1) C 49 - H 59	0.82	1.20	0.040
74. BD (2) C 45 - C 47	/493. RY* (3) C 49	0.53	0.78	0.027
74. BD (2) C 45 - C 47	/647. BD* (2) C 44 - C 46	6.01	0.29	0.057
74. BD (2) C 45 - C 47	/656. BD* (2) C 48 - C 49	6.03	0.31	0.056
75. BD (1) C 45 - H 56	/451. RY* (1) C 44	0.40	1.63	0.032
75. BD (1) C 45 - H 56	/475. RY* (1) C 47	0.62	1.71	0.041
75. BD (1) C 45 - H 56	/644. BD* (1) N 43 - C 44	0.33	0.88	0.021
75. BD (1) C 45 - H 56	/646. BD* (1) C 44 - C 46	1.63	1.11	0.054
75. BD (1) C 45 - H 56	/648. BD* (1) C 45 - C 47	0.76	1.16	0.037
75. BD (1) C 45 - H 56	/653. BD* (1) C 47 - C 49	2.72	0.98	0.065
75. BD (1) C 45 - H 56	/654. BD* (1) C 47 - H 57	0.50	0.96	0.028
76. BD (1) C 46 - C 48	/451. RY* (1) C 44	0.60	1.77	0.041
76. BD (1) C 46 - C 48	/453. RY* (3) C 44	0.27	1.53	0.026
76. BD (1) C 46 - C 48	/491. RY* (1) C 49	0.36	1.84	0.033
76. BD (1) C 46 - C 48	/492. RY* (2) C 49	0.72	1.99	0.048
76. BD (1) C 46 - C 48	/644. BD* (1) N 43 - C 44	2.67	1.02	0.066
76. BD (1) C 46 - C 48	/646. BD* (1) C 44 - C 46	1.71	1.25	0.059
76. BD (1) C 46 - C 48	/655. BD* (1) C 48 - C 49	1.29	1.30	0.052
76. BD (1) C 46 - C 48	/657. BD* (1) C 48 - H 58	0.31	1.10	0.024
76. BD (1) C 46 - C 48	/658. BD* (1) C 49 - H 59	1.53	1.10	0.052
77. BD (1) C 46 - O 65	/451. RY* (1) C 44	0.30	1.96	0.031
77. BD (1) C 46 - O 65	/467. RY* (1) C 46	0.55	2.00	0.042
77. BD (1) C 46 - O 65	/645. BD* (1) C 44 - C 45	1.30	1.28	0.052
77. BD (1) C 46 - O 65	/646. BD* (1) C 44 - C 46	0.72	1.44	0.041
77. BD (1) C 46 - O 65	/651. BD* (1) C 46 - C 48	0.29	1.29	0.025
77. BD (1) C 46 - O 65	/655. BD* (1) C 48 - C 49	0.54	1.49	0.036
78. BD (1) C 47 - C 49	/460. RY* (2) C 45	0.98	1.84	0.054
78. BD (1) C 47 - C 49	/484. RY* (2) C 48	0.98	1.83	0.054
78. BD (1) C 47 - C 49	/648. BD* (1) C 45 - C 47	1.22	1.29	0.050
78. BD (1) C 47 - C 49	/650. BD* (1) C 45 - H 56	1.64	1.09	0.053
78. BD (1) C 47 - C 49	/654. BD* (1) C 47 - H 57	0.29	1.09	0.022
78. BD (1) C 47 - C 49	/655. BD* (1) C 48 - C 49	1.23	1.30	0.051
78. BD (1) C 47 - C 49	/657. BD* (1) C 48 - H 58	1.64	1.10	0.054
78. BD (1) C 47 - C 49	/658. BD* (1) C 49 - H 59	0.29	1.10	0.022
79. BD (1) C 47 - H 57	/459. RY* (1) C 45	0.80	1.79	0.048
79. BD (1) C 47 - H 57	/491. RY* (1) C 49	0.60	1.71	0.041
79. BD (1) C 47 - H 57	/645. BD* (1) C 44 - C 45	3.06	0.96	0.068
79. BD (1) C 47 - H 57	/648. BD* (1) C 45 - C 47	0.68	1.16	0.036
79. BD (1) C 47 - H 57	/650. BD* (1) C 45 - H 56	0.50	0.96	0.028
79. BD (1) C 47 - H 57	/655. BD* (1) C 48 - C 49	1.36	1.17	0.050
80. BD (1) C 48 - C 49	/467. RY* (1) C 46	0.51	1.90	0.039
80. BD (1) C 48 - C 49	/468. RY* (2) C 46	0.54	2.19	0.044
80. BD (1) C 48 - C 49	/475. RY* (1) C 47	0.50	1.94	0.039
80. BD (1) C 48 - C 49	/476. RY* (2) C 47	0.41	2.07	0.037
80. BD (1) C 48 - C 49	/651. BD* (1) C 46 - C 48	0.98	1.19	0.043
80. BD (1) C 48 - C 49	/652. BD* (1) C 46 - O 65	1.16	1.07	0.045
80. BD (1) C 48 - C 49	/653. BD* (1) C 47 - C 49	1.15	1.21	0.047

80. BD (1) C 48 - C 49	/654. BD* (1) C 47 - H 57	0.82	1.19	0.039
80. BD (1) C 48 - C 49	/657. BD* (1) C 48 - H 58	0.91	1.20	0.042
80. BD (1) C 48 - C 49	/658. BD* (1) C 49 - H 59	0.81	1.19	0.039
81. BD (2) C 48 - C 49	/477. RY* (3) C 47	0.52	0.78	0.027
81. BD (2) C 48 - C 49	/647. BD* (2) C 44 - C 46	5.72	0.29	0.055
81. BD (2) C 48 - C 49	/649. BD* (2) C 45 - C 47	7.14	0.31	0.060
82. BD (1) C 48 - H 58	/467. RY* (1) C 46	0.38	1.67	0.032
82. BD (1) C 48 - H 58	/491. RY* (1) C 49	0.57	1.70	0.039
82. BD (1) C 48 - H 58	/646. BD* (1) C 44 - C 46	1.58	1.11	0.053
82. BD (1) C 48 - H 58	/652. BD* (1) C 46 - O 65	0.45	0.83	0.025
82. BD (1) C 48 - H 58	/653. BD* (1) C 47 - C 49	2.75	0.97	0.065
82. BD (1) C 48 - H 58	/655. BD* (1) C 48 - C 49	0.78	1.16	0.038
82. BD (1) C 48 - H 58	/658. BD* (1) C 49 - H 59	0.48	0.96	0.027
83. BD (1) C 49 - H 59	/475. RY* (1) C 47	0.59	1.70	0.040
83. BD (1) C 49 - H 59	/483. RY* (1) C 48	0.83	1.85	0.050
83. BD (1) C 49 - H 59	/648. BD* (1) C 45 - C 47	1.37	1.16	0.050
83. BD (1) C 49 - H 59	/651. BD* (1) C 46 - C 48	2.91	0.96	0.067
83. BD (1) C 49 - H 59	/655. BD* (1) C 48 - C 49	0.71	1.16	0.036
83. BD (1) C 49 - H 59	/657. BD* (1) C 48 - H 58	0.47	0.96	0.027
84. BD (1) O 50 - H 60	/355. RY* (1) C 32	0.90	1.64	0.049
84. BD (1) O 50 - H 60	/356. RY* (2) C 32	0.60	1.51	0.038
84. BD (1) O 50 - H 60	/623. BD* (1) C 32 - C 34	1.61	1.34	0.059
109. CR (1) C 32	/364. RY* (2) C 33	0.56	11.81	0.102
109. CR (1) C 32	/373. RY* (3) C 34	1.07	11.48	0.140
109. CR (1) C 32	/623. BD* (1) C 32 - C 34	0.91	10.74	0.125
109. CR (1) C 32	/624. BD* (1) C 32 - O 50	0.67	10.39	0.106
109. CR (1) C 32	/625. BD* (1) C 33 - C 35	0.29	10.75	0.071
109. CR (1) C 32	/626. BD* (1) C 33 - C 38	0.31	10.57	0.073
109. CR (1) C 32	/627. BD* (1) C 34 - C 36	0.41	10.58	0.083
110. CR (1) C 33	/355. RY* (1) C 32	0.25	10.97	0.067
110. CR (1) C 33	/381. RY* (3) C 35	1.14	11.63	0.145
110. CR (1) C 33	/405. RY* (3) C 38	0.51	11.20	0.096
110. CR (1) C 33	/623. BD* (1) C 32 - C 34	0.28	10.67	0.069
110. CR (1) C 33	/625. BD* (1) C 33 - C 35	0.61	10.68	0.103
110. CR (1) C 33	/629. BD* (1) C 35 - C 37	0.54	10.51	0.097
110. CR (1) C 33	/630. BD* (1) C 35 - C 40	0.49	10.52	0.091
110. CR (1) C 33	/634. BD* (1) C 38 - C 39	0.27	10.71	0.068
111. CR (1) C 34	/357. RY* (3) C 32	1.49	11.48	0.165
111. CR (1) C 34	/390. RY* (4) C 36	0.49	11.17	0.093
111. CR (1) C 34	/508. RY* (2) H 51	0.29	12.28	0.075
111. CR (1) C 34	/621. BD* (1) C 32 - C 33	0.60	10.47	0.101
111. CR (1) C 34	/623. BD* (1) C 32 - C 34	0.29	10.66	0.071
111. CR (1) C 34	/624. BD* (1) C 32 - O 50	0.36	10.32	0.078
111. CR (1) C 34	/631. BD* (1) C 36 - C 37	0.30	10.68	0.072
112. CR (1) C 35	/365. RY* (3) C 33	1.22	11.50	0.150
112. CR (1) C 35	/397. RY* (3) C 37	0.75	11.19	0.116
112. CR (1) C 35	/421. RY* (3) C 40	0.59	11.17	0.103
112. CR (1) C 35	/621. BD* (1) C 32 - C 33	0.51	10.47	0.094
112. CR (1) C 35	/625. BD* (1) C 33 - C 35	0.51	10.67	0.093
112. CR (1) C 35	/626. BD* (1) C 33 - C 38	0.46	10.50	0.089
112. CR (1) C 35	/631. BD* (1) C 36 - C 37	0.32	10.69	0.075
112. CR (1) C 35	/639. BD* (1) C 40 - C 41	0.28	10.71	0.070
113. CR (1) C 36	/396. RY* (2) C 37	0.40	11.13	0.084
113. CR (1) C 36	/397. RY* (3) C 37	0.87	11.17	0.125
113. CR (1) C 36	/398. RY* (4) C 37	0.50	11.09	0.094
113. CR (1) C 36	/528. RY* (2) H 61	0.26	12.34	0.072
113. CR (1) C 36	/623. BD* (1) C 32 - C 34	0.29	10.64	0.070
113. CR (1) C 36	/629. BD* (1) C 35 - C 37	0.64	10.49	0.105
113. CR (1) C 36	/631. BD* (1) C 36 - C 37	0.46	10.67	0.089
113. CR (1) C 36	/633. BD* (1) C 37 - N 42	0.43	10.42	0.085
114. CR (1) C 37	/380. RY* (2) C 35	0.61	11.68	0.107
114. CR (1) C 37	/389. RY* (3) C 36	1.13	11.30	0.142
114. CR (1) C 37	/625. BD* (1) C 33 - C 35	0.35	10.70	0.078
114. CR (1) C 37	/627. BD* (1) C 34 - C 36	0.44	10.54	0.087
114. CR (1) C 37	/630. BD* (1) C 35 - C 40	0.32	10.54	0.074
114. CR (1) C 37	/631. BD* (1) C 36 - C 37	0.67	10.71	0.108
114. CR (1) C 37	/643. BD* (1) N 42 - N 43	0.38	10.59	0.081
115. CR (1) C 38	/364. RY* (2) C 33	0.29	11.72	0.074
115. CR (1) C 38	/413. RY* (3) C 39	1.04	11.39	0.137
115. CR (1) C 38	/510. RY* (2) H 52	0.27	12.26	0.073
115. CR (1) C 38	/621. BD* (1) C 32 - C 33	0.33	10.45	0.075

115. CR (1) C 38	/625. BD* (1) C 33 - C 35	0.36	10.65	0.078
115. CR (1) C 38	/634. BD* (1) C 38 - C 39	0.39	10.68	0.082
115. CR (1) C 38	/637. BD* (1) C 39 - C 41	0.49	10.50	0.091
116. CR (1) C 39	/404. RY* (2) C 38	0.34	11.26	0.078
116. CR (1) C 39	/405. RY* (3) C 38	0.65	11.17	0.107
116. CR (1) C 39	/406. RY* (4) C 38	0.40	11.07	0.084
116. CR (1) C 39	/429. RY* (3) C 41	0.38	11.40	0.083
116. CR (1) C 39	/512. RY* (2) H 53	0.33	12.30	0.080
116. CR (1) C 39	/626. BD* (1) C 33 - C 38	0.60	10.48	0.101
116. CR (1) C 39	/634. BD* (1) C 38 - C 39	0.39	10.68	0.082
116. CR (1) C 39	/639. BD* (1) C 40 - C 41	0.29	10.69	0.070
117. CR (1) C 40	/381. RY* (3) C 35	0.38	11.59	0.084
117. CR (1) C 40	/429. RY* (3) C 41	1.05	11.39	0.138
117. CR (1) C 40	/625. BD* (1) C 33 - C 35	0.41	10.64	0.084
117. CR (1) C 40	/629. BD* (1) C 35 - C 37	0.34	10.47	0.076
117. CR (1) C 40	/637. BD* (1) C 39 - C 41	0.50	10.49	0.092
117. CR (1) C 40	/639. BD* (1) C 40 - C 41	0.39	10.68	0.082
118. CR (1) C 41	/413. RY* (3) C 39	0.38	11.38	0.083
118. CR (1) C 41	/420. RY* (2) C 40	0.33	11.18	0.077
118. CR (1) C 41	/421. RY* (3) C 40	0.45	11.14	0.090
118. CR (1) C 41	/422. RY* (4) C 40	0.65	11.11	0.107
118. CR (1) C 41	/516. RY* (2) H 55	0.32	12.30	0.080
118. CR (1) C 41	/630. BD* (1) C 35 - C 40	0.61	10.48	0.102
118. CR (1) C 41	/634. BD* (1) C 38 - C 39	0.29	10.67	0.070
118. CR (1) C 41	/639. BD* (1) C 40 - C 41	0.43	10.68	0.086
119. CR (1) N 42	/395. RY* (1) C 37	0.43	15.09	0.102
119. CR (1) N 42	/396. RY* (2) C 37	0.71	15.28	0.132
119. CR (1) N 42	/445. RY* (3) N 43	0.87	15.28	0.145
119. CR (1) N 42	/629. BD* (1) C 35 - C 37	0.29	14.64	0.083
119. CR (1) N 42	/631. BD* (1) C 36 - C 37	0.27	14.82	0.081
119. CR (1) N 42	/644. BD* (1) N 43 - C 44	0.70	14.56	0.129
120. CR (1) N 43	/437. RY* (3) N 42	1.01	15.44	0.158
120. CR (1) N 43	/451. RY* (1) C 44	0.73	15.34	0.134
120. CR (1) N 43	/452. RY* (2) C 44	0.34	15.61	0.092
120. CR (1) N 43	/633. BD* (1) C 37 - N 42	0.62	14.60	0.121
120. CR (1) N 43	/646. BD* (1) C 44 - C 46	0.37	14.82	0.095
121. CR (1) C 44	/460. RY* (2) C 45	0.50	11.30	0.095
121. CR (1) C 44	/467. RY* (1) C 46	0.53	11.26	0.097
121. CR (1) C 44	/468. RY* (2) C 46	1.11	11.55	0.143
121. CR (1) C 44	/643. BD* (1) N 42 - N 43	0.47	10.61	0.090
121. CR (1) C 44	/646. BD* (1) C 44 - C 46	0.61	10.71	0.103
121. CR (1) C 44	/648. BD* (1) C 45 - C 47	0.29	10.75	0.070
121. CR (1) C 44	/651. BD* (1) C 46 - C 48	0.49	10.55	0.092
121. CR (1) C 44	/652. BD* (1) C 46 - O 65	0.40	10.43	0.083
122. CR (1) C 45	/451. RY* (1) C 44	0.38	11.16	0.082
122. CR (1) C 45	/453. RY* (3) C 44	0.31	10.93	0.073
122. CR (1) C 45	/476. RY* (2) C 47	1.05	11.37	0.138
122. CR (1) C 45	/518. RY* (2) H 56	0.31	12.30	0.078
122. CR (1) C 45	/644. BD* (1) N 43 - C 44	0.26	10.41	0.066
122. CR (1) C 45	/646. BD* (1) C 44 - C 46	0.39	10.64	0.083
122. CR (1) C 45	/648. BD* (1) C 45 - C 47	0.43	10.69	0.086
122. CR (1) C 45	/653. BD* (1) C 47 - C 49	0.52	10.51	0.094
123. CR (1) C 46	/452. RY* (2) C 44	1.53	11.51	0.168
123. CR (1) C 46	/484. RY* (2) C 48	0.53	11.30	0.098
123. CR (1) C 46	/644. BD* (1) N 43 - C 44	0.45	10.48	0.088
123. CR (1) C 46	/645. BD* (1) C 44 - C 45	0.50	10.56	0.092
123. CR (1) C 46	/646. BD* (1) C 44 - C 46	0.75	10.72	0.115
123. CR (1) C 46	/652. BD* (1) C 46 - O 65	0.42	10.44	0.085
123. CR (1) C 46	/655. BD* (1) C 48 - C 49	0.30	10.77	0.072
124. CR (1) C 47	/459. RY* (1) C 45	0.40	11.31	0.085
124. CR (1) C 47	/460. RY* (2) C 45	0.80	11.23	0.119
124. CR (1) C 47	/492. RY* (2) C 49	0.47	11.38	0.093
124. CR (1) C 47	/520. RY* (2) H 57	0.33	12.00	0.079
124. CR (1) C 47	/645. BD* (1) C 44 - C 45	0.57	10.48	0.098
124. CR (1) C 47	/648. BD* (1) C 45 - C 47	0.38	10.69	0.080
124. CR (1) C 47	/655. BD* (1) C 48 - C 49	0.30	10.69	0.072
125. CR (1) C 48	/468. RY* (2) C 46	0.55	11.48	0.101
125. CR (1) C 48	/469. RY* (3) C 46	0.32	10.96	0.074
125. CR (1) C 48	/492. RY* (2) C 49	1.04	11.38	0.137
125. CR (1) C 48	/522. RY* (2) H 58	0.27	12.33	0.073
125. CR (1) C 48	/646. BD* (1) C 44 - C 46	0.43	10.64	0.087

125. CR (1) C 48	/653. BD* (1) C 47 - C 49	0.53	10.50	0.095
125. CR (1) C 48	/655. BD* (1) C 48 - C 49	0.44	10.69	0.087
126. CR (1) C 49	/476. RY* (2) C 47	0.47	11.37	0.092
126. CR (1) C 49	/483. RY* (1) C 48	0.47	11.37	0.092
126. CR (1) C 49	/484. RY* (2) C 48	0.80	11.22	0.119
126. CR (1) C 49	/524. RY* (2) H 59	0.30	11.95	0.076
126. CR (1) C 49	/648. BD* (1) C 45 - C 47	0.30	10.68	0.072
126. CR (1) C 49	/651. BD* (1) C 46 - C 48	0.52	10.48	0.094
126. CR (1) C 49	/655. BD* (1) C 48 - C 49	0.39	10.69	0.082
127. CR (1) O 50	/355. RY* (1) C 32	0.76	19.88	0.156
127. CR (1) O 50	/356. RY* (2) C 32	0.33	19.75	0.101
127. CR (1) O 50	/526. RY* (2) H 60	0.31	20.89	0.101
127. CR (1) O 50	/621. BD* (1) C 32 - C 33	0.25	19.39	0.089
129. CR (1) O 65	/467. RY* (1) C 46	1.46	20.05	0.216
129. CR (1) O 65	/646. BD* (1) C 44 - C 46	0.32	19.50	0.101
151. LP (1) N 42	/ 54. BD* (2) C 35 - C 37	0.87	0.25	0.025
151. LP (1) N 42	/443. RY* (1) N 43	0.50	1.16	0.031
151. LP (1) N 42	/444. RY* (2) N 43	1.11	1.10	0.045
151. LP (1) N 42	/627. BD* (1) C 34 - C 36	0.27	0.75	0.018
151. LP (1) N 42	/629. BD* (1) C 35 - C 37	3.09	0.74	0.061
151. LP (1) N 42	/631. BD* (1) C 36 - C 37	3.31	0.92	0.071
152. LP (2) N 42	/ 54. BD* (2) C 35 - C 37	2.21	0.40	0.050
152. LP (2) N 42	/395. RY* (1) C 37	0.73	1.34	0.041
152. LP (2) N 42	/435. RY* (1) N 42	1.62	2.34	0.080
152. LP (2) N 42	/436. RY* (2) N 42	0.30	1.39	0.027
152. LP (2) N 42	/437. RY* (3) N 42	0.61	1.67	0.041
152. LP (2) N 42	/444. RY* (2) N 43	0.35	1.26	0.027
152. LP (2) N 42	/541. RY* (1) O 65	0.57	2.83	0.052
152. LP (2) N 42	/629. BD* (1) C 35 - C 37	1.73	0.89	0.050
153. LP (1) N 43	/437. RY* (3) N 42	0.42	1.63	0.033
153. LP (1) N 43	/451. RY* (1) C 44	0.96	1.53	0.049
153. LP (1) N 43	/645. BD* (1) C 44 - C 45	0.54	0.85	0.027
153. LP (1) N 43	/646. BD* (1) C 44 - C 46	3.76	1.01	0.078
154. LP (2) N 43	/436. RY* (2) N 42	0.96	1.22	0.045
154. LP (2) N 43	/454. RY* (4) C 44	0.32	0.92	0.023
154. LP (2) N 43	/647. BD* (2) C 44 - C 46	15.70	0.29	0.091
155. LP (1) O 50	/355. RY* (1) C 32	0.97	1.52	0.049
155. LP (1) O 50	/356. RY* (2) C 32	0.51	1.39	0.034
155. LP (1) O 50	/621. BD* (1) C 32 - C 33	3.24	1.02	0.073
156. LP (2) O 50	/622. BD* (2) C 32 - C 33	7.52	0.29	0.080
159. LP (1) O 65	/467. RY* (1) C 46	1.23	1.73	0.059
159. LP (1) O 65	/646. BD* (1) C 44 - C 46	3.45	1.17	0.080
160. LP (2) O 65	/435. RY* (1) N 42	0.30	2.35	0.034
160. LP (2) O 65	/467. RY* (1) C 46	0.66	1.62	0.042
160. LP (2) O 65	/541. RY* (1) O 65	0.81	2.85	0.062
160. LP (2) O 65	/646. BD* (1) C 44 - C 46	0.53	1.06	0.030
160. LP (2) O 65	/651. BD* (1) C 46 - C 48	3.94	0.91	0.076
161. LP (3) O 65	/647. BD* (2) C 44 - C 46	12.34	0.33	0.087
622. BD* (2) C 32 - C 33	/ 54. BD* (2) C 35 - C 37	400.09	0.01	0.082
622. BD* (2) C 32 - C 33	/356. RY* (2) C 32	0.50	0.83	0.037
622. BD* (2) C 32 - C 33	/363. RY* (1) C 33	0.88	0.63	0.043
622. BD* (2) C 32 - C 33	/371. RY* (1) C 34	1.02	0.54	0.042
622. BD* (2) C 32 - C 33	/379. RY* (1) C 35	0.56	0.64	0.034
622. BD* (2) C 32 - C 33	/403. RY* (1) C 38	0.35	0.58	0.026
622. BD* (2) C 32 - C 33	/499. RY* (1) O 50	0.36	0.92	0.033
622. BD* (2) C 32 - C 33	/635. BD* (2) C 38 - C 39	41.03	0.07	0.075
635. BD* (2) C 38 - C 39	/403. RY* (1) C 38	2.27	0.52	0.076
635. BD* (2) C 38 - C 39	/411. RY* (1) C 39	2.04	0.47	0.069
640. BD* (2) C 40 - C 41	/419. RY* (1) C 40	1.78	0.61	0.073
640. BD* (2) C 40 - C 41	/427. RY* (1) C 41	2.15	0.47	0.071
647. BD* (2) C 44 - C 46	/454. RY* (4) C 44	0.75	0.63	0.058
647. BD* (2) C 44 - C 46	/470. RY* (4) C 46	0.55	0.74	0.054
647. BD* (2) C 44 - C 46	/649. BD* (2) C 45 - C 47	62.76	0.02	0.071
647. BD* (2) C 44 - C 46	/656. BD* (2) C 48 - C 49	55.32	0.02	0.073
649. BD* (2) C 45 - C 47	/461. RY* (3) C 45	0.88	0.51	0.075
649. BD* (2) C 45 - C 47	/477. RY* (3) C 47	0.93	0.47	0.074
656. BD* (2) C 48 - C 49	/485. RY* (3) C 48	1.02	0.51	0.078
656. BD* (2) C 48 - C 49	/493. RY* (3) C 49	0.98	0.46	0.073
from unit 2 to unit 3				
50. BD* (2) C 34 - C 36	/660. BD* (1) O 62 - H 63	0.05	0.39	0.007

56. BD (1) C 36 - C 37	/532. RY* (4) O 62	0.03	2.13	0.010
68. BD (1) N 42 - N 43	/529. RY* (1) O 62	0.03	3.00	0.012
69. BD (1) N 43 - C 44	/529. RY* (1) O 62	0.04	2.81	0.014
151. LP (1) N 42	/529. RY* (1) O 62	0.06	2.32	0.015
152. LP (2) N 42	/529. RY* (1) O 62	0.99	2.47	0.064
152. LP (2) N 42	/531. RY* (3) O 62	0.15	1.80	0.021
152. LP (2) N 42	/532. RY* (4) O 62	0.35	1.81	0.032
152. LP (2) N 42	/537. RY* (1) H 63	0.32	1.00	0.023
152. LP (2) N 42	/539. RY* (1) H 64	0.24	1.01	0.020
159. LP (1) O 65	/529. RY* (1) O 62	0.17	2.60	0.027
159. LP (1) O 65	/531. RY* (3) O 62	0.04	1.93	0.011
159. LP (1) O 65	/532. RY* (4) O 62	0.07	1.93	0.015
159. LP (1) O 65	/537. RY* (1) H 63	0.06	1.12	0.011
159. LP (1) O 65	/539. RY* (1) H 64	0.06	1.13	0.010
160. LP (2) O 65	/529. RY* (1) O 62	0.39	2.49	0.040
160. LP (2) O 65	/531. RY* (3) O 62	0.13	1.82	0.020
160. LP (2) O 65	/532. RY* (4) O 62	0.20	1.82	0.024
160. LP (2) O 65	/537. RY* (1) H 63	0.14	1.01	0.015
160. LP (2) O 65	/539. RY* (1) H 64	0.10	1.02	0.013
160. LP (2) O 65	/660. BD* (1) O 62 - H 63	0.03	0.79	0.006
161. LP (3) O 65	/529. RY* (1) O 62	0.10	2.34	0.020
161. LP (3) O 65	/531. RY* (3) O 62	0.03	1.67	0.010
161. LP (3) O 65	/532. RY* (4) O 62	0.06	1.67	0.013

from unit 2 to unit 4

52. BD (1) C 35 - C 37	/551. RY* (1) H 67	0.03	1.28	0.008
53. BD (2) C 35 - C 37	/663. BD* (1) H 67 - O 68	0.17	0.61	0.013
66. BD (1) C 40 - H 54	/662. BD* (1) H 66 - O 68	0.03	0.85	0.006
68. BD (1) N 42 - N 43	/553. RY* (1) O 68	0.03	2.90	0.011
69. BD (1) N 43 - C 44	/553. RY* (1) O 68	0.03	2.70	0.012
151. LP (1) N 42	/553. RY* (1) O 68	0.03	2.22	0.011
152. LP (2) N 42	/549. RY* (1) H 66	0.37	1.14	0.027
152. LP (2) N 42	/551. RY* (1) H 67	0.41	1.07	0.027
152. LP (2) N 42	/553. RY* (1) O 68	0.90	2.37	0.060
152. LP (2) N 42	/556. RY* (4) O 68	0.35	2.25	0.036
152. LP (2) N 42	/662. BD* (1) H 66 - O 68	0.03	0.78	0.007
152. LP (2) N 42	/663. BD* (1) H 67 - O 68	0.08	0.78	0.010
159. LP (1) O 65	/549. RY* (1) H 66	0.07	1.26	0.012
159. LP (1) O 65	/551. RY* (1) H 67	0.12	1.20	0.015
159. LP (1) O 65	/553. RY* (1) O 68	0.20	2.50	0.029
159. LP (1) O 65	/556. RY* (4) O 68	0.07	2.37	0.016
160. LP (2) O 65	/549. RY* (1) H 66	0.15	1.15	0.017
160. LP (2) O 65	/551. RY* (1) H 67	0.23	1.09	0.020
160. LP (2) O 65	/553. RY* (1) O 68	0.49	2.39	0.044
160. LP (2) O 65	/556. RY* (4) O 68	0.21	2.26	0.028
161. LP (3) O 65	/549. RY* (1) H 66	0.12	1.00	0.015
161. LP (3) O 65	/553. RY* (1) O 68	0.12	2.24	0.021
161. LP (3) O 65	/556. RY* (4) O 68	0.04	2.11	0.011
161. LP (3) O 65	/662. BD* (1) H 66 - O 68	0.13	0.64	0.012
640. BD* (2) C 40 - C 41	/662. BD* (1) H 66 - O 68	0.07	0.32	0.011

from unit 2 to unit 5

52. BD (1) C 35 - C 37	/166. LP (6) Mn 69	0.07	1.42	0.014
52. BD (1) C 35 - C 37	/167. LP* (7) Mn 69	0.03	1.19	0.007
53. BD (2) C 35 - C 37	/561. RY* (1) Mn 69	0.03	2.07	0.010
56. BD (1) C 36 - C 37	/167. LP* (7) Mn 69	0.03	1.31	0.008
56. BD (1) C 36 - C 37	/561. RY* (1) Mn 69	0.05	2.56	0.014
56. BD (1) C 36 - C 37	/570. RY* (10) Mn 69	0.03	9.53	0.021
58. BD (1) C 37 - N 42	/166. LP (6) Mn 69	0.60	1.54	0.041
58. BD (1) C 37 - N 42	/167. LP* (7) Mn 69	0.05	1.32	0.010
58. BD (1) C 37 - N 42	/168. LP* (8) Mn 69	0.06	2.61	0.016
58. BD (1) C 37 - N 42	/169. LP* (9) Mn 69	0.04	2.53	0.013
58. BD (1) C 37 - N 42	/566. RY* (6) Mn 69	0.05	2.54	0.014
58. BD (1) C 37 - N 42	/567. RY* (7) Mn 69	0.03	2.79	0.011
58. BD (1) C 37 - N 42	/568. RY* (8) Mn 69	0.06	7.46	0.027
68. BD (1) N 42 - N 43	/166. LP (6) Mn 69	0.78	1.74	0.050
68. BD (1) N 42 - N 43	/167. LP* (7) Mn 69	0.04	1.51	0.009
68. BD (1) N 42 - N 43	/169. LP* (9) Mn 69	0.13	2.72	0.024
68. BD (1) N 42 - N 43	/561. RY* (1) Mn 69	0.03	2.77	0.011
68. BD (1) N 42 - N 43	/562. RY* (2) Mn 69	0.15	3.96	0.031
68. BD (1) N 42 - N 43	/565. RY* (5) Mn 69	0.04	2.78	0.013

68.	BD	(1)	N	42	-	N	43	/568.	RY*	(8)	Mn	69	0.03	7.66	0.019
68.	BD	(1)	N	42	-	N	43	/570.	RY*	(10)	Mn	69	0.04	9.74	0.025
68.	BD	(1)	N	42	-	N	43	/580.	RY*	(20)	Mn	69	0.03	149.94	0.089
68.	BD	(1)	N	42	-	N	43	/581.	RY*	(21)	Mn	69	0.05	39.35	0.058
69.	BD	(1)	N	43	-	C	44	/166.	LP	(6)	Mn	69	0.12	1.54	0.018
69.	BD	(1)	N	43	-	C	44	/167.	LP*	(7)	Mn	69	0.04	1.32	0.010
69.	BD	(1)	N	43	-	C	44	/561.	RY*	(1)	Mn	69	0.04	2.57	0.014
69.	BD	(1)	N	43	-	C	44	/562.	RY*	(2)	Mn	69	0.09	3.77	0.023
69.	BD	(1)	N	43	-	C	44	/570.	RY*	(10)	Mn	69	0.09	9.54	0.036
69.	BD	(1)	N	43	-	C	44	/578.	RY*	(18)	Mn	69	0.04	679.51	0.197
69.	BD	(1)	N	43	-	C	44	/580.	RY*	(20)	Mn	69	0.07	149.75	0.126
69.	BD	(1)	N	43	-	C	44	/581.	RY*	(21)	Mn	69	0.12	39.15	0.088
70.	BD	(1)	C	44	-	C	45	/166.	LP	(6)	Mn	69	0.03	1.42	0.008
71.	BD	(1)	C	44	-	C	46	/166.	LP	(6)	Mn	69	0.13	1.53	0.019
71.	BD	(1)	C	44	-	C	46	/167.	LP*	(7)	Mn	69	0.03	1.31	0.008
71.	BD	(1)	C	44	-	C	46	/562.	RY*	(2)	Mn	69	0.03	3.76	0.013
71.	BD	(1)	C	44	-	C	46	/581.	RY*	(21)	Mn	69	0.03	39.14	0.044
76.	BD	(1)	C	46	-	C	48	/166.	LP	(6)	Mn	69	0.19	1.41	0.022
76.	BD	(1)	C	46	-	C	48	/167.	LP*	(7)	Mn	69	0.06	1.19	0.010
76.	BD	(1)	C	46	-	C	48	/562.	RY*	(2)	Mn	69	0.03	3.64	0.014
76.	BD	(1)	C	46	-	C	48	/581.	RY*	(21)	Mn	69	0.03	39.02	0.045
77.	BD	(1)	C	46	-	O	65	/166.	LP	(6)	Mn	69	0.67	1.60	0.044
77.	BD	(1)	C	46	-	O	65	/168.	LP*	(8)	Mn	69	0.09	2.67	0.020
77.	BD	(1)	C	46	-	O	65	/169.	LP*	(9)	Mn	69	0.15	2.58	0.025
77.	BD	(1)	C	46	-	O	65	/562.	RY*	(2)	Mn	69	0.03	3.83	0.014
77.	BD	(1)	C	46	-	O	65	/564.	RY*	(4)	Mn	69	0.03	3.03	0.013
77.	BD	(1)	C	46	-	O	65	/568.	RY*	(8)	Mn	69	0.04	7.52	0.022
77.	BD	(1)	C	46	-	O	65	/571.	RY*	(11)	Mn	69	0.05	6.92	0.023
114.	CR	(1)	C	37				/166.	LP	(6)	Mn	69	0.03	10.85	0.022
119.	CR	(1)	N	42				/166.	LP	(6)	Mn	69	0.99	14.95	0.165
120.	CR	(1)	N	43				/166.	LP	(6)	Mn	69	0.03	14.98	0.031
123.	CR	(1)	C	46				/166.	LP	(6)	Mn	69	0.05	10.88	0.033
129.	CR	(1)	O	65				/166.	LP	(6)	Mn	69	1.23	19.66	0.211
151.	LP	(1)	N	42				/166.	LP	(6)	Mn	69	0.94	1.06	0.042
151.	LP	(1)	N	42				/168.	LP*	(8)	Mn	69	0.05	2.12	0.014
151.	LP	(1)	N	42				/562.	RY*	(2)	Mn	69	0.13	3.28	0.027
151.	LP	(1)	N	42				/575.	RY*	(15)	Mn	69	0.03	5.85	0.018
151.	LP	(1)	N	42				/578.	RY*	(18)	Mn	69	0.03	679.02	0.172
151.	LP	(1)	N	42				/580.	RY*	(20)	Mn	69	0.05	149.26	0.114
151.	LP	(1)	N	42				/581.	RY*	(21)	Mn	69	0.09	38.67	0.078
152.	LP	(2)	N	42				/166.	LP	(6)	Mn	69	18.47	1.21	0.197
152.	LP	(2)	N	42				/167.	LP*	(7)	Mn	69	0.15	0.99	0.016
152.	LP	(2)	N	42				/168.	LP*	(8)	Mn	69	0.30	2.28	0.034
152.	LP	(2)	N	42				/561.	RY*	(1)	Mn	69	0.65	2.24	0.049
152.	LP	(2)	N	42				/562.	RY*	(2)	Mn	69	2.74	3.44	0.126
152.	LP	(2)	N	42				/564.	RY*	(4)	Mn	69	0.09	2.64	0.020
152.	LP	(2)	N	42				/565.	RY*	(5)	Mn	69	0.15	2.26	0.023
152.	LP	(2)	N	42				/566.	RY*	(6)	Mn	69	0.04	2.21	0.012
152.	LP	(2)	N	42				/567.	RY*	(7)	Mn	69	0.32	2.46	0.036
152.	LP	(2)	N	42				/570.	RY*	(10)	Mn	69	0.09	9.21	0.038
152.	LP	(2)	N	42				/574.	RY*	(14)	Mn	69	0.07	6.37	0.027
152.	LP	(2)	N	42				/575.	RY*	(15)	Mn	69	0.18	6.00	0.042
152.	LP	(2)	N	42				/578.	RY*	(18)	Mn	69	0.62	679.17	0.844
152.	LP	(2)	N	42				/580.	RY*	(20)	Mn	69	1.24	149.41	0.559
152.	LP	(2)	N	42				/581.	RY*	(21)	Mn	69	2.31	38.82	0.388
153.	LP	(1)	N	43				/166.	LP	(6)	Mn	69	0.21	1.17	0.021
153.	LP	(1)	N	43				/561.	RY*	(1)	Mn	69	0.06	2.20	0.015
153.	LP	(1)	N	43				/565.	RY*	(5)	Mn	69	0.03	2.22	0.010
159.	LP	(1)	O	65				/166.	LP	(6)	Mn	69	5.04	1.34	0.110
159.	LP	(1)	O	65				/167.	LP*	(7)	Mn	69	0.13	1.11	0.015
159.	LP	(1)	O	65				/168.	LP*	(8)	Mn	69	0.08	2.40	0.017
159.	LP	(1)	O	65				/561.	RY*	(1)	Mn	69	0.14	2.36	0.023
159.	LP	(1)	O	65				/562.	RY*	(2)	Mn	69	0.64	3.56	0.061
159.	LP	(1)	O	65				/563.	RY*	(3)	Mn	69	0.03	2.34	0.010
159.	LP	(1)	O	65				/567.	RY*	(7)	Mn	69	0.14	2.58	0.024
159.	LP	(1)	O	65				/575.	RY*	(15)	Mn	69	0.03	6.13	0.018
159.	LP	(1)	O	65				/578.	RY*	(18)	Mn	69	0.15	679.30	0.412
159.	LP	(1)	O	65				/580.	RY*	(20)	Mn	69	0.27	149.54	0.254
159.	LP	(1)	O	65				/581.	RY*	(21)	Mn	69	0.48	38.95	0.175
160.	LP	(2)	O	65				/166.	LP	(6)	Mn	69	12.37	1.23	0.164
160.	LP	(2)	O	65				/167.	LP*	(7)	Mn	69	0.60	1.00	0.031

160. LP (2) O 65	/168. LP* (8) Mn 69	0.88	2.29	0.058
160. LP (2) O 65	/169. LP* (9) Mn 69	0.03	2.21	0.011
160. LP (2) O 65	/561. RY* (1) Mn 69	0.12	2.25	0.021
160. LP (2) O 65	/562. RY* (2) Mn 69	1.23	3.45	0.084
160. LP (2) O 65	/563. RY* (3) Mn 69	0.03	2.23	0.011
160. LP (2) O 65	/567. RY* (7) Mn 69	0.28	2.47	0.034
160. LP (2) O 65	/573. RY* (13) Mn 69	0.09	8.16	0.035
160. LP (2) O 65	/575. RY* (15) Mn 69	0.13	6.02	0.036
160. LP (2) O 65	/578. RY* (18) Mn 69	0.35	679.19	0.630
160. LP (2) O 65	/579. RY* (19) Mn 69	0.03	27.52	0.035
160. LP (2) O 65	/580. RY* (20) Mn 69	0.69	149.43	0.412
160. LP (2) O 65	/581. RY* (21) Mn 69	1.28	38.84	0.287
161. LP (3) O 65	/166. LP (6) Mn 69	3.67	1.07	0.082
161. LP (3) O 65	/167. LP* (7) Mn 69	0.49	0.85	0.027
161. LP (3) O 65	/168. LP* (8) Mn 69	0.11	2.14	0.020
161. LP (3) O 65	/169. LP* (9) Mn 69	0.09	2.05	0.017
161. LP (3) O 65	/561. RY* (1) Mn 69	0.12	2.10	0.021
161. LP (3) O 65	/562. RY* (2) Mn 69	0.34	3.30	0.044
161. LP (3) O 65	/567. RY* (7) Mn 69	0.05	2.32	0.015
161. LP (3) O 65	/575. RY* (15) Mn 69	0.04	5.87	0.020
161. LP (3) O 65	/578. RY* (18) Mn 69	0.09	679.04	0.316
161. LP (3) O 65	/580. RY* (20) Mn 69	0.18	149.28	0.215
161. LP (3) O 65	/581. RY* (21) Mn 69	0.35	38.68	0.151

from unit 3 to unit 1

85. BD (1) O 62 - H 63	/253. RY* (1) N 11	0.10	2.13	0.019
85. BD (1) O 62 - H 63	/254. RY* (2) N 11	0.04	2.18	0.012
85. BD (1) O 62 - H 63	/255. RY* (3) N 11	0.05	2.04	0.013
85. BD (1) O 62 - H 63	/256. RY* (4) N 11	0.05	2.30	0.013
85. BD (1) O 62 - H 63	/264. RY* (4) N 12	0.03	2.41	0.011
85. BD (1) O 62 - H 63	/289. RY* (5) C 15	0.03	4.22	0.015
85. BD (1) O 62 - H 63	/610. BD* (2) C 13 - C 15	0.03	0.80	0.007
86. BD (1) O 62 - H 64	/253. RY* (1) N 11	0.06	2.13	0.014
86. BD (1) O 62 - H 64	/254. RY* (2) N 11	0.03	2.19	0.010
86. BD (1) O 62 - H 64	/255. RY* (3) N 11	0.03	2.05	0.010
86. BD (1) O 62 - H 64	/285. RY* (1) C 15	0.04	1.94	0.011
86. BD (1) O 62 - H 64	/301. RY (1) C 17	0.03	1.28	0.007
86. BD (1) O 62 - H 64	/317. RY* (1) O 19	0.04	3.15	0.014
86. BD (1) O 62 - H 64	/610. BD* (2) C 13 - C 15	0.04	0.80	0.008
86. BD (1) O 62 - H 64	/614. BD* (1) C 15 - O 19	0.07	1.10	0.011
157. LP (1) O 62	/269. RY* (1) C 13	0.03	1.55	0.008
157. LP (1) O 62	/286. RY* (2) C 15	0.05	1.21	0.010
157. LP (1) O 62	/596. BD* (1) C 6 - N 11	0.06	0.82	0.009
157. LP (1) O 62	/614. BD* (1) C 15 - O 19	0.06	0.77	0.009
158. LP (2) O 62	/213. RY* (1) C 6	0.03	1.53	0.009
158. LP (2) O 62	/214. RY* (2) C 6	0.16	1.81	0.021
158. LP (2) O 62	/216. RY* (4) C 6	0.08	1.69	0.015
158. LP (2) O 62	/218. RY* (6) C 6	0.04	3.86	0.016
158. LP (2) O 62	/253. RY* (1) N 11	0.59	2.00	0.044
158. LP (2) O 62	/254. RY* (2) N 11	0.36	2.05	0.034
158. LP (2) O 62	/255. RY* (3) N 11	0.21	1.91	0.025
158. LP (2) O 62	/256. RY* (4) N 11	0.15	2.18	0.023
158. LP (2) O 62	/264. RY* (4) N 12	0.03	2.28	0.011
158. LP (2) O 62	/271. RY* (3) C 13	0.03	1.46	0.009
158. LP (2) O 62	/285. RY* (1) C 15	0.19	1.81	0.024
158. LP (2) O 62	/286. RY* (2) C 15	0.09	1.40	0.014
158. LP (2) O 62	/289. RY* (5) C 15	0.12	4.09	0.028
158. LP (2) O 62	/317. RY* (1) O 19	0.46	3.02	0.048
158. LP (2) O 62	/319. RY* (3) O 19	0.05	1.98	0.013
158. LP (2) O 62	/320. RY* (4) O 19	0.03	2.45	0.010
158. LP (2) O 62	/596. BD* (1) C 6 - N 11	0.06	1.02	0.010
158. LP (2) O 62	/614. BD* (1) C 15 - O 19	0.03	0.97	0.007

from unit 3 to unit 2

85. BD (1) O 62 - H 63	/387. RY* (1) C 36	0.05	1.30	0.010
85. BD (1) O 62 - H 63	/388. RY* (2) C 36	0.04	2.02	0.011
85. BD (1) O 62 - H 63	/397. RY* (3) C 37	0.05	1.90	0.013
86. BD (1) O 62 - H 64	/ 50. BD* (2) C 34 - C 36	0.04	0.72	0.011
86. BD (1) O 62 - H 64	/387. RY* (1) C 36	0.04	1.31	0.009
86. BD (1) O 62 - H 64	/388. RY* (2) C 36	0.06	2.03	0.014
86. BD (1) O 62 - H 64	/435. RY* (1) N 42	0.06	2.67	0.016

86. BD (1) O 62 - H 64	/437. RY* (3) N 42	0.10	2.00	0.018
157. LP (1) O 62	/ 50. BD* (2) C 34 - C 36	0.11	0.39	0.013
157. LP (1) O 62	/ 54. BD* (2) C 35 - C 37	0.09	0.40	0.011
158. LP (2) O 62	/ 54. BD* (2) C 35 - C 37	0.08	0.60	0.012
158. LP (2) O 62	/395. RY* (1) C 37	0.04	1.54	0.010
158. LP (2) O 62	/396. RY* (2) C 37	0.14	1.73	0.020
158. LP (2) O 62	/400. RY* (6) C 37	0.08	3.78	0.022
158. LP (2) O 62	/435. RY* (1) N 42	0.47	2.54	0.044
158. LP (2) O 62	/436. RY* (2) N 42	0.07	1.59	0.014
158. LP (2) O 62	/437. RY* (3) N 42	0.40	1.87	0.035
158. LP (2) O 62	/438. RY* (4) N 42	0.09	2.14	0.017
158. LP (2) O 62	/443. RY* (1) N 43	0.03	1.51	0.009
158. LP (2) O 62	/444. RY* (2) N 43	0.03	1.45	0.008
158. LP (2) O 62	/445. RY* (3) N 43	0.11	1.73	0.018
158. LP (2) O 62	/467. RY* (1) C 46	0.13	1.80	0.020
158. LP (2) O 62	/468. RY* (2) C 46	0.04	2.09	0.012
158. LP (2) O 62	/470. RY* (4) C 46	0.05	1.40	0.010
158. LP (2) O 62	/471. RY* (5) C 46	0.04	5.49	0.018
158. LP (2) O 62	/541. RY* (1) O 65	0.15	3.03	0.027
158. LP (2) O 62	/543. RY* (3) O 65	0.05	2.04	0.013
158. LP (2) O 62	/544. RY* (4) O 65	0.04	2.33	0.012
158. LP (2) O 62	/633. BD* (1) C 37 - N 42	0.05	1.02	0.009
158. LP (2) O 62	/643. BD* (1) N 42 - N 43	0.18	1.15	0.018

within unit 3

128. CR (1) O 62	/538. RY* (2) H 63	0.40	21.17	0.116
128. CR (1) O 62	/540. RY* (2) H 64	0.41	21.21	0.118
158. LP (2) O 62	/529. RY* (1) O 62	2.53	2.67	0.105
158. LP (2) O 62	/531. RY* (3) O 62	0.31	2.00	0.032
158. LP (2) O 62	/532. RY* (4) O 62	0.45	2.00	0.038
158. LP (2) O 62	/537. RY* (1) H 63	0.95	1.20	0.043
158. LP (2) O 62	/538. RY* (2) H 63	0.31	2.85	0.038
158. LP (2) O 62	/539. RY* (1) H 64	1.02	1.21	0.045
158. LP (2) O 62	/540. RY* (2) H 64	0.32	2.90	0.039

from unit 3 to unit 4

85. BD (1) O 62 - H 63	/549. RY* (1) H 66	0.03	1.46	0.009
85. BD (1) O 62 - H 63	/551. RY* (1) H 67	0.05	1.40	0.011
85. BD (1) O 62 - H 63	/553. RY* (1) O 68	0.06	2.70	0.016
86. BD (1) O 62 - H 64	/549. RY* (1) H 66	0.03	1.47	0.009
86. BD (1) O 62 - H 64	/553. RY* (1) O 68	0.03	2.70	0.012
158. LP (2) O 62	/549. RY* (1) H 66	0.46	1.34	0.031
158. LP (2) O 62	/551. RY* (1) H 67	0.44	1.27	0.030
158. LP (2) O 62	/553. RY* (1) O 68	0.93	2.57	0.062
158. LP (2) O 62	/554. RY* (2) O 68	0.05	1.63	0.011
158. LP (2) O 62	/556. RY* (4) O 68	0.35	2.45	0.037
158. LP (2) O 62	/663. BD* (1) H 67 - O 68	0.03	0.98	0.006

from unit 3 to unit 5

85. BD (1) O 62 - H 63	/166. LP (6) Mn 69	2.19	1.54	0.079
85. BD (1) O 62 - H 63	/167. LP* (7) Mn 69	0.04	1.31	0.009
85. BD (1) O 62 - H 63	/168. LP* (8) Mn 69	0.04	2.60	0.013
85. BD (1) O 62 - H 63	/169. LP* (9) Mn 69	0.03	2.52	0.011
85. BD (1) O 62 - H 63	/562. RY* (2) Mn 69	0.56	3.76	0.058
85. BD (1) O 62 - H 63	/563. RY* (3) Mn 69	0.09	2.54	0.019
85. BD (1) O 62 - H 63	/565. RY* (5) Mn 69	0.04	2.58	0.013
85. BD (1) O 62 - H 63	/568. RY* (8) Mn 69	0.05	7.45	0.023
85. BD (1) O 62 - H 63	/571. RY* (11) Mn 69	0.03	6.86	0.019
85. BD (1) O 62 - H 63	/578. RY* (18) Mn 69	0.07	679.50	0.272
85. BD (1) O 62 - H 63	/580. RY* (20) Mn 69	0.12	149.74	0.168
85. BD (1) O 62 - H 63	/581. RY* (21) Mn 69	0.21	39.15	0.115
86. BD (1) O 62 - H 64	/166. LP (6) Mn 69	2.30	1.54	0.081
86. BD (1) O 62 - H 64	/562. RY* (2) Mn 69	0.55	3.77	0.057
86. BD (1) O 62 - H 64	/563. RY* (3) Mn 69	0.03	2.54	0.012
86. BD (1) O 62 - H 64	/564. RY* (4) Mn 69	0.16	2.97	0.027
86. BD (1) O 62 - H 64	/568. RY* (8) Mn 69	0.03	7.46	0.017
86. BD (1) O 62 - H 64	/575. RY* (15) Mn 69	0.04	6.34	0.020
86. BD (1) O 62 - H 64	/578. RY* (18) Mn 69	0.08	679.51	0.301
86. BD (1) O 62 - H 64	/580. RY* (20) Mn 69	0.15	149.75	0.189
86. BD (1) O 62 - H 64	/581. RY* (21) Mn 69	0.27	39.15	0.129
128. CR (1) O 62	/166. LP (6) Mn 69	1.34	19.72	0.221

157. LP (1) O 62	/166. LP (6) Mn 69	0.27	1.22	0.024
157. LP (1) O 62	/168. LP* (8) Mn 69	0.04	2.28	0.011
157. LP (1) O 62	/169. LP* (9) Mn 69	0.03	2.20	0.010
158. LP (2) O 62	/166. LP (6) Mn 69	19.06	1.41	0.219
158. LP (2) O 62	/167. LP* (7) Mn 69	0.25	1.18	0.022
158. LP (2) O 62	/168. LP* (8) Mn 69	0.92	2.47	0.061
158. LP (2) O 62	/169. LP* (9) Mn 69	0.83	2.39	0.057
158. LP (2) O 62	/561. RY* (1) Mn 69	0.29	2.44	0.034
158. LP (2) O 62	/562. RY* (2) Mn 69	2.94	3.63	0.132
158. LP (2) O 62	/564. RY* (4) Mn 69	0.19	2.84	0.030
158. LP (2) O 62	/565. RY* (5) Mn 69	0.18	2.46	0.027
158. LP (2) O 62	/566. RY* (6) Mn 69	0.03	2.41	0.011
158. LP (2) O 62	/567. RY* (7) Mn 69	0.22	2.65	0.031
158. LP (2) O 62	/568. RY* (8) Mn 69	0.22	7.33	0.052
158. LP (2) O 62	/569. RY* (9) Mn 69	0.04	8.67	0.023
158. LP (2) O 62	/570. RY* (10) Mn 69	0.13	9.41	0.045
158. LP (2) O 62	/571. RY* (11) Mn 69	0.21	6.73	0.048
158. LP (2) O 62	/572. RY* (12) Mn 69	0.08	5.82	0.027
158. LP (2) O 62	/573. RY* (13) Mn 69	0.18	8.34	0.050
158. LP (2) O 62	/574. RY* (14) Mn 69	0.10	6.57	0.033
158. LP (2) O 62	/575. RY* (15) Mn 69	0.30	6.20	0.055
158. LP (2) O 62	/577. RY* (17) Mn 69	0.03	27.91	0.034
158. LP (2) O 62	/578. RY* (18) Mn 69	0.79	679.37	0.932
158. LP (2) O 62	/579. RY* (19) Mn 69	0.11	27.70	0.071
158. LP (2) O 62	/580. RY* (20) Mn 69	1.52	149.61	0.608
158. LP (2) O 62	/581. RY* (21) Mn 69	2.76	39.02	0.419

from unit 4 to unit 1

87. BD (1) H 66 - O 68	/206. RY* (2) C 5	0.05	2.02	0.013
88. BD (1) H 67 - O 68	/213. RY* (1) C 6	0.03	1.66	0.009
88. BD (1) H 67 - O 68	/214. RY* (2) C 6	0.03	1.93	0.009
88. BD (1) H 67 - O 68	/215. RY* (3) C 6	0.05	1.88	0.012
88. BD (1) H 67 - O 68	/216. RY* (4) C 6	0.03	1.82	0.010
88. BD (1) H 67 - O 68	/253. RY* (1) N 11	0.03	2.13	0.010
88. BD (1) H 67 - O 68	/255. RY* (3) N 11	0.03	2.04	0.010
162. LP (1) O 68	/353. RY* (1) H 31	0.04	1.01	0.008
162. LP (1) O 68	/354. RY* (2) H 31	0.03	2.67	0.012
162. LP (1) O 68	/589. BD* (2) C 3 - C 5	0.37	0.43	0.018
162. LP (1) O 68	/592. BD* (2) C 4 - C 6	0.06	0.43	0.008
163. LP (2) O 68	/213. RY* (1) C 6	0.03	1.53	0.009
163. LP (2) O 68	/214. RY* (2) C 6	0.14	1.80	0.020
163. LP (2) O 68	/216. RY* (4) C 6	0.03	1.69	0.009
163. LP (2) O 68	/218. RY* (6) C 6	0.08	3.85	0.022
163. LP (2) O 68	/253. RY* (1) N 11	0.30	2.00	0.031
163. LP (2) O 68	/254. RY* (2) N 11	0.27	2.05	0.030
163. LP (2) O 68	/255. RY* (3) N 11	0.34	1.91	0.032
163. LP (2) O 68	/256. RY* (4) N 11	0.11	2.17	0.020
163. LP (2) O 68	/262. RY* (2) N 12	0.03	1.43	0.008
163. LP (2) O 68	/263. RY* (3) N 12	0.09	1.75	0.016
163. LP (2) O 68	/285. RY* (1) C 15	0.13	1.80	0.019
163. LP (2) O 68	/286. RY* (2) C 15	0.06	1.40	0.011
163. LP (2) O 68	/287. RY* (3) C 15	0.03	1.95	0.010
163. LP (2) O 68	/289. RY* (5) C 15	0.10	4.08	0.026
163. LP (2) O 68	/317. RY* (1) O 19	0.12	3.01	0.024
163. LP (2) O 68	/319. RY* (3) O 19	0.05	1.97	0.013
163. LP (2) O 68	/589. BD* (2) C 3 - C 5	0.06	0.63	0.009
163. LP (2) O 68	/592. BD* (2) C 4 - C 6	0.07	0.63	0.010
163. LP (2) O 68	/596. BD* (1) C 6 - N 11	0.07	1.01	0.011
163. LP (2) O 68	/606. BD* (1) N 11 - N 12	0.22	1.13	0.020

from unit 4 to unit 2

87. BD (1) H 66 - O 68	/396. RY* (2) C 37	0.03	1.86	0.010
87. BD (1) H 66 - O 68	/435. RY* (1) N 42	0.10	2.66	0.020
87. BD (1) H 66 - O 68	/436. RY* (2) N 42	0.04	1.72	0.010
87. BD (1) H 66 - O 68	/437. RY* (3) N 42	0.17	1.99	0.023
87. BD (1) H 66 - O 68	/438. RY* (4) N 42	0.03	2.27	0.010
87. BD (1) H 66 - O 68	/541. RY* (1) O 65	0.09	3.16	0.021
87. BD (1) H 66 - O 68	/544. RY* (4) O 65	0.03	2.46	0.011
87. BD (1) H 66 - O 68	/640. BD* (2) C 40 - C 41	0.04	0.79	0.009
88. BD (1) H 67 - O 68	/395. RY* (1) C 37	0.03	1.66	0.008
88. BD (1) H 67 - O 68	/396. RY* (2) C 37	0.03	1.86	0.010

88. BD (1) H 67 - O 68	/435. RY* (1) N 42	0.05	2.66	0.015
88. BD (1) H 67 - O 68	/436. RY* (2) N 42	0.04	1.72	0.010
88. BD (1) H 67 - O 68	/437. RY* (3) N 42	0.09	1.99	0.017
88. BD (1) H 67 - O 68	/438. RY* (4) N 42	0.03	2.27	0.010
88. BD (1) H 67 - O 68	/652. BD* (1) C 46 - O 65	0.03	1.09	0.007
162. LP (1) O 68	/396. RY* (2) C 37	0.03	1.53	0.009
162. LP (1) O 68	/419. RY* (1) C 40	0.03	1.07	0.007
162. LP (1) O 68	/633. BD* (1) C 37 - N 42	0.03	0.82	0.006
162. LP (1) O 68	/643. BD* (1) N 42 - N 43	0.06	0.95	0.009
163. LP (2) O 68	/396. RY* (2) C 37	0.16	1.73	0.021
163. LP (2) O 68	/398. RY* (4) C 37	0.03	1.68	0.009
163. LP (2) O 68	/400. RY* (6) C 37	0.04	3.78	0.015
163. LP (2) O 68	/419. RY* (1) C 40	0.04	1.27	0.009
163. LP (2) O 68	/435. RY* (1) N 42	0.50	2.53	0.045
163. LP (2) O 68	/436. RY* (2) N 42	0.17	1.59	0.021
163. LP (2) O 68	/437. RY* (3) N 42	0.40	1.86	0.035
163. LP (2) O 68	/438. RY* (4) N 42	0.06	2.14	0.014
163. LP (2) O 68	/443. RY* (1) N 43	0.06	1.51	0.012
163. LP (2) O 68	/445. RY* (3) N 43	0.08	1.73	0.015
163. LP (2) O 68	/453. RY* (3) C 44	0.03	1.53	0.009
163. LP (2) O 68	/467. RY* (1) C 46	0.17	1.80	0.022
163. LP (2) O 68	/470. RY* (4) C 46	0.04	1.40	0.010
163. LP (2) O 68	/471. RY* (5) C 46	0.04	5.49	0.018
163. LP (2) O 68	/541. RY* (1) O 65	0.53	3.03	0.051
163. LP (2) O 68	/543. RY* (3) O 65	0.07	2.04	0.015
163. LP (2) O 68	/544. RY* (4) O 65	0.10	2.33	0.019
163. LP (2) O 68	/643. BD* (1) N 42 - N 43	0.07	1.15	0.011
163. LP (2) O 68	/652. BD* (1) C 46 - O 65	0.06	0.97	0.010

from unit 4 to unit 3

87. BD (1) H 66 - O 68	/529. RY* (1) O 62	0.06	2.80	0.017
87. BD (1) H 66 - O 68	/537. RY* (1) H 63	0.03	1.32	0.008
87. BD (1) H 66 - O 68	/661. BD* (1) O 62 - H 64	0.03	1.10	0.007
88. BD (1) H 67 - O 68	/529. RY* (1) O 62	0.08	2.80	0.019
163. LP (2) O 68	/529. RY* (1) O 62	0.95	2.67	0.064
163. LP (2) O 68	/531. RY* (3) O 62	0.14	2.00	0.021
163. LP (2) O 68	/532. RY* (4) O 62	0.34	2.00	0.033
163. LP (2) O 68	/537. RY* (1) H 63	0.25	1.19	0.022
163. LP (2) O 68	/539. RY* (1) H 64	0.24	1.20	0.022
163. LP (2) O 68	/660. BD* (1) O 62 - H 63	0.04	0.97	0.008

within unit 4

88. BD (1) H 67 - O 68	/550. RY* (2) H 66	0.28	2.94	0.036
130. CR (1) O 68	/550. RY* (2) H 66	0.36	21.12	0.110
130. CR (1) O 68	/552. RY* (2) H 67	0.39	21.16	0.115
163. LP (2) O 68	/549. RY* (1) H 66	0.93	1.33	0.045
163. LP (2) O 68	/550. RY* (2) H 66	0.40	2.81	0.043
163. LP (2) O 68	/551. RY* (1) H 67	1.17	1.26	0.049
163. LP (2) O 68	/552. RY* (2) H 67	0.39	2.85	0.042
163. LP (2) O 68	/553. RY* (1) O 68	2.43	2.57	0.101
163. LP (2) O 68	/556. RY* (4) O 68	0.60	2.44	0.049

from unit 4 to unit 5

87. BD (1) H 66 - O 68	/166. LP (6) Mn 69	2.51	1.53	0.084
87. BD (1) H 66 - O 68	/169. LP* (9) Mn 69	0.05	2.52	0.014
87. BD (1) H 66 - O 68	/562. RY* (2) Mn 69	0.68	3.76	0.064
87. BD (1) H 66 - O 68	/563. RY* (3) Mn 69	0.03	2.54	0.011
87. BD (1) H 66 - O 68	/564. RY* (4) Mn 69	0.05	2.96	0.015
87. BD (1) H 66 - O 68	/569. RY* (9) Mn 69	0.03	8.80	0.022
87. BD (1) H 66 - O 68	/570. RY* (10) Mn 69	0.03	9.53	0.021
87. BD (1) H 66 - O 68	/575. RY* (15) Mn 69	0.04	6.33	0.020
87. BD (1) H 66 - O 68	/578. RY* (18) Mn 69	0.10	679.50	0.324
87. BD (1) H 66 - O 68	/580. RY* (20) Mn 69	0.17	149.74	0.202
87. BD (1) H 66 - O 68	/581. RY* (21) Mn 69	0.30	39.14	0.138
88. BD (1) H 67 - O 68	/166. LP (6) Mn 69	2.43	1.53	0.083
88. BD (1) H 67 - O 68	/169. LP* (9) Mn 69	0.06	2.51	0.016
88. BD (1) H 67 - O 68	/562. RY* (2) Mn 69	0.57	3.76	0.059
88. BD (1) H 67 - O 68	/563. RY* (3) Mn 69	0.03	2.53	0.011
88. BD (1) H 67 - O 68	/566. RY* (6) Mn 69	0.03	2.53	0.011
88. BD (1) H 67 - O 68	/567. RY* (7) Mn 69	0.05	2.78	0.015
88. BD (1) H 67 - O 68	/569. RY* (9) Mn 69	0.06	8.80	0.030

88. BD (1) H 67 - O 68	/578. RY* (18) Mn 69	0.08	679.50	0.291
88. BD (1) H 67 - O 68	/580. RY* (20) Mn 69	0.14	149.74	0.182
88. BD (1) H 67 - O 68	/581. RY* (21) Mn 69	0.25	39.14	0.125
130. CR (1) O 68	/166. LP (6) Mn 69	1.37	19.72	0.222
162. LP (1) O 68	/166. LP (6) Mn 69	0.17	1.21	0.019
162. LP (1) O 68	/168. LP* (8) Mn 69	0.03	2.27	0.010
162. LP (1) O 68	/561. RY* (1) Mn 69	0.03	2.24	0.011
163. LP (2) O 68	/166. LP (6) Mn 69	18.50	1.40	0.216
163. LP (2) O 68	/167. LP* (7) Mn 69	0.18	1.18	0.019
163. LP (2) O 68	/168. LP* (8) Mn 69	0.32	2.47	0.036
163. LP (2) O 68	/169. LP* (9) Mn 69	2.15	2.39	0.091
163. LP (2) O 68	/561. RY* (1) Mn 69	0.15	2.43	0.024
163. LP (2) O 68	/562. RY* (2) Mn 69	2.62	3.63	0.124
163. LP (2) O 68	/563. RY* (3) Mn 69	0.03	2.41	0.011
163. LP (2) O 68	/565. RY* (5) Mn 69	0.08	2.45	0.018
163. LP (2) O 68	/567. RY* (7) Mn 69	0.27	2.65	0.034
163. LP (2) O 68	/568. RY* (8) Mn 69	0.12	7.32	0.038
163. LP (2) O 68	/569. RY* (9) Mn 69	0.57	8.67	0.089
163. LP (2) O 68	/570. RY* (10) Mn 69	0.03	9.40	0.022
163. LP (2) O 68	/573. RY* (13) Mn 69	0.11	8.34	0.038
163. LP (2) O 68	/574. RY* (14) Mn 69	0.24	6.56	0.051
163. LP (2) O 68	/575. RY* (15) Mn 69	0.08	6.20	0.028
163. LP (2) O 68	/576. RY* (16) Mn 69	0.15	27.60	0.083
163. LP (2) O 68	/578. RY* (18) Mn 69	0.75	679.37	0.910
163. LP (2) O 68	/580. RY* (20) Mn 69	1.45	149.61	0.594
163. LP (2) O 68	/581. RY* (21) Mn 69	2.63	39.01	0.409

from unit 5 to unit 1

132. CR (2) Mn 69	/596. BD* (1) C 6 - N 11	0.13	26.79	0.074
132. CR (2) Mn 69	/606. BD* (1) N 11 - N 12	0.09	26.91	0.062
132. CR (2) Mn 69	/614. BD* (1) C 15 - O 19	0.13	26.75	0.075
133. CR (3) Mn 69	/255. RY* (3) N 11	0.08	5.38	0.026
133. CR (3) Mn 69	/596. BD* (1) C 6 - N 11	0.64	4.49	0.068
133. CR (3) Mn 69	/606. BD* (1) N 11 - N 12	0.45	4.61	0.058
133. CR (3) Mn 69	/613. BD* (1) C 15 - C 17	0.03	4.56	0.014
133. CR (3) Mn 69	/614. BD* (1) C 15 - O 19	0.75	4.44	0.074
135. CR (5) Mn 69	/254. RY* (2) N 11	0.03	3.48	0.012
135. CR (5) Mn 69	/255. RY* (3) N 11	0.06	3.34	0.017
135. CR (5) Mn 69	/596. BD* (1) C 6 - N 11	0.19	2.44	0.028
135. CR (5) Mn 69	/606. BD* (1) N 11 - N 12	0.16	2.56	0.026
137. CR (7) Mn 69	/317. RY* (1) O 19	0.03	4.44	0.014
137. CR (7) Mn 69	/606. BD* (1) N 11 - N 12	0.05	2.56	0.015
137. CR (7) Mn 69	/614. BD* (1) C 15 - O 19	0.38	2.39	0.038
139. CR (9) Mn 69	/596. BD* (1) C 6 - N 11	0.08	2.44	0.017
139. CR (9) Mn 69	/606. BD* (1) N 11 - N 12	0.03	2.56	0.011
139. CR (9) Mn 69	/614. BD* (1) C 15 - O 19	0.04	2.40	0.013
164. LP (3) Mn 69	/216. RY* (4) C 6	0.03	1.28	0.008
164. LP (3) Mn 69	/596. BD* (1) C 6 - N 11	0.13	0.60	0.011
164. LP (3) Mn 69	/606. BD* (1) N 11 - N 12	0.41	0.72	0.022
165. LP (5) Mn 69	/214. RY* (2) C 6	0.05	1.37	0.011
165. LP (5) Mn 69	/216. RY* (4) C 6	0.04	1.26	0.009
165. LP (5) Mn 69	/253. RY* (1) N 11	0.92	1.57	0.049
165. LP (5) Mn 69	/254. RY* (2) N 11	0.82	1.62	0.047
165. LP (5) Mn 69	/255. RY* (3) N 11	0.40	1.48	0.031
165. LP (5) Mn 69	/259. RY* (7) N 11	0.03	35.37	0.041
165. LP (5) Mn 69	/261. RY* (1) N 12	0.03	1.04	0.007
165. LP (5) Mn 69	/264. RY* (4) N 12	0.04	1.85	0.011
165. LP (5) Mn 69	/285. RY* (1) C 15	0.16	1.38	0.019
165. LP (5) Mn 69	/286. RY* (2) C 15	0.03	0.97	0.007
165. LP (5) Mn 69	/592. BD* (2) C 4 - C 6	0.03	0.20	0.004
165. LP (5) Mn 69	/596. BD* (1) C 6 - N 11	0.25	0.58	0.015
165. LP (5) Mn 69	/607. BD* (1) N 12 - C 13	0.07	0.57	0.008
166. LP (6) Mn 69	/183. RY* (3) C 2	0.08	0.16	0.013
166. LP (6) Mn 69	/190. RY* (2) C 3	0.03	0.52	0.013
166. LP (6) Mn 69	/197. RY* (1) C 4	0.06	0.79	0.023
166. LP (6) Mn 69	/198. RY* (2) C 4	0.17	0.77	0.040
166. LP (6) Mn 69	/200. RY* (4) C 4	0.04	0.20	0.010
166. LP (6) Mn 69	/201. RY* (5) C 4	0.04	2.35	0.035
166. LP (6) Mn 69	/207. RY* (3) C 5	0.26	0.43	0.036
166. LP (6) Mn 69	/208. RY* (4) C 5	0.48	0.29	0.041
166. LP (6) Mn 69	/213. RY* (1) C 6	3.87	0.12	0.074

166.	LP	(6)	Mn	69	/214.	RY*	(2)	C	6	7.35	0.40	0.186
166.	LP	(6)	Mn	69	/215.	RY*	(3)	C	6	0.26	0.35	0.033
166.	LP	(6)	Mn	69	/216.	RY*	(4)	C	6	4.92	0.28	0.129
166.	LP	(6)	Mn	69	/218.	RY*	(6)	C	6	0.46	2.45	0.117
166.	LP	(6)	Mn	69	/220.	RY*	(8)	C	6	0.07	2.03	0.042
166.	LP	(6)	Mn	69	/239.	RY*	(3)	C	9	0.28	0.31	0.032
166.	LP	(6)	Mn	69	/240.	RY*	(4)	C	9	0.10	0.35	0.021
166.	LP	(6)	Mn	69	/253.	RY*	(1)	N	11	20.18	0.59	0.373
166.	LP	(6)	Mn	69	/254.	RY*	(2)	N	11	18.49	0.64	0.373
166.	LP	(6)	Mn	69	/255.	RY*	(3)	N	11	15.27	0.50	0.304
166.	LP	(6)	Mn	69	/256.	RY*	(4)	N	11	3.18	0.77	0.172
166.	LP	(6)	Mn	69	/258.	RY*	(6)	N	11	0.04	3.21	0.039
166.	LP	(6)	Mn	69	/261.	RY*	(1)	N	12	5.94	0.06	0.066
166.	LP	(6)	Mn	69	/262.	RY*	(2)	N	12	11.07	0.02	0.052
166.	LP	(6)	Mn	69	/263.	RY*	(3)	N	12	2.41	0.34	0.100
166.	LP	(6)	Mn	69	/264.	RY*	(4)	N	12	0.63	0.87	0.081
166.	LP	(6)	Mn	69	/265.	RY*	(5)	N	12	0.05	3.39	0.047
166.	LP	(6)	Mn	69	/266.	RY*	(6)	N	12	0.03	3.07	0.032
166.	LP	(6)	Mn	69	/269.	RY*	(1)	C	13	0.25	0.34	0.031
166.	LP	(6)	Mn	69	/270.	RY*	(2)	C	13	0.28	0.67	0.047
166.	LP	(6)	Mn	69	/271.	RY*	(3)	C	13	7.31	0.05	0.064
166.	LP	(6)	Mn	69	/274.	RY*	(6)	C	13	0.07	2.58	0.047
166.	LP	(6)	Mn	69	/280.	RY*	(4)	C	14	0.14	0.14	0.016
166.	LP	(6)	Mn	69	/285.	RY*	(1)	C	15	14.71	0.40	0.263
166.	LP	(6)	Mn	69	/287.	RY*	(3)	C	15	0.83	0.55	0.074
166.	LP	(6)	Mn	69	/288.	RY*	(4)	C	15	0.38	0.27	0.035
166.	LP	(6)	Mn	69	/289.	RY*	(5)	C	15	1.31	2.68	0.206
166.	LP	(6)	Mn	69	/294.	RY*	(2)	C	16	0.11	0.45	0.024
166.	LP	(6)	Mn	69	/296.	RY*	(4)	C	16	0.08	0.25	0.016
166.	LP	(6)	Mn	69	/302.	RY*	(2)	C	17	0.46	0.57	0.056
166.	LP	(6)	Mn	69	/304.	RY*	(4)	C	17	0.52	0.12	0.027
166.	LP	(6)	Mn	69	/309.	RY*	(1)	C	18	0.07	0.43	0.019
166.	LP	(6)	Mn	69	/317.	RY*	(1)	O	19	13.34	1.61	0.507
166.	LP	(6)	Mn	69	/318.	RY*	(2)	O	19	0.05	0.33	0.013
166.	LP	(6)	Mn	69	/319.	RY*	(3)	O	19	3.63	0.57	0.158
166.	LP	(6)	Mn	69	/320.	RY*	(4)	O	19	0.83	1.04	0.102
170.	LP	(1)	Mn	69	/253.	RY*	(1)	N	11	0.13	1.59	0.018
170.	LP	(1)	Mn	69	/254.	RY*	(2)	N	11	0.13	1.64	0.018
170.	LP	(1)	Mn	69	/255.	RY*	(3)	N	11	0.04	1.50	0.009
170.	LP	(1)	Mn	69	/317.	RY*	(1)	O	19	0.11	2.60	0.021
170.	LP	(1)	Mn	69	/596.	BD*	(1)	C	6 - N 11	0.04	0.60	0.006
170.	LP	(1)	Mn	69	/606.	BD*	(1)	N	11 - N 12	0.10	0.72	0.011
170.	LP	(1)	Mn	69	/614.	BD*	(1)	C	15 - O 19	0.04	0.55	0.006
171.	LP	(2)	Mn	69	/253.	RY*	(1)	N	11	0.03	1.59	0.009
171.	LP	(2)	Mn	69	/269.	RY*	(1)	C	13	0.05	1.33	0.010
171.	LP	(2)	Mn	69	/289.	RY*	(5)	C	15	0.03	3.67	0.013
171.	LP	(2)	Mn	69	/317.	RY*	(1)	O	19	0.04	2.60	0.013
171.	LP	(2)	Mn	69	/596.	BD*	(1)	C	6 - N 11	0.17	0.60	0.013
171.	LP	(2)	Mn	69	/606.	BD*	(1)	N	11 - N 12	0.33	0.72	0.020
171.	LP	(2)	Mn	69	/613.	BD*	(1)	C	15 - C 17	0.04	0.68	0.007
171.	LP	(2)	Mn	69	/614.	BD*	(1)	C	15 - O 19	0.13	0.55	0.011
172.	LP	(4)	Mn	69	/253.	RY*	(1)	N	11	0.10	1.57	0.016
172.	LP	(4)	Mn	69	/254.	RY*	(2)	N	11	0.12	1.62	0.018
172.	LP	(4)	Mn	69	/255.	RY*	(3)	N	11	0.09	1.48	0.015
172.	LP	(4)	Mn	69	/317.	RY*	(1)	O	19	0.61	2.58	0.051
172.	LP	(4)	Mn	69	/318.	RY*	(2)	O	19	0.04	1.30	0.009
172.	LP	(4)	Mn	69	/319.	RY*	(3)	O	19	0.05	1.55	0.011
172.	LP	(4)	Mn	69	/324.	RY*	(8)	O	19	0.03	49.31	0.047
172.	LP	(4)	Mn	69	/609.	BD*	(1)	C	13 - C 15	0.03	0.81	0.006
172.	LP	(4)	Mn	69	/614.	BD*	(1)	C	15 - O 19	0.43	0.54	0.019
from unit 5 to unit 2														
132.	CR	(2)	Mn	69	/633.	BD*	(1)	C	37 - N 42	0.13	26.80	0.076
132.	CR	(2)	Mn	69	/643.	BD*	(1)	N	42 - N 43	0.07	26.93	0.056
132.	CR	(2)	Mn	69	/652.	BD*	(1)	C	46 - O 65	0.14	26.75	0.077
133.	CR	(3)	Mn	69	/437.	RY*	(3)	N	42	0.06	5.34	0.022
133.	CR	(3)	Mn	69	/633.	BD*	(1)	C	37 - N 42	0.67	4.49	0.070
133.	CR	(3)	Mn	69	/643.	BD*	(1)	N	42 - N 43	0.37	4.62	0.053
133.	CR	(3)	Mn	69	/651.	BD*	(1)	C	46 - C 48	0.04	4.56	0.017
133.	CR	(3)	Mn	69	/652.	BD*	(1)	C	46 - O 65	0.77	4.44	0.075
135.	CR	(5)	Mn	69	/437.	RY*	(3)	N	42	0.06	3.30	0.017

135. CR (5)Mn 69	/633. BD* (1) C 37 - N 42	0.16	2.45	0.025
135. CR (5)Mn 69	/643. BD* (1) N 42 - N 43	0.21	2.58	0.030
135. CR (5)Mn 69	/652. BD* (1) C 46 - O 65	0.11	2.40	0.020
137. CR (7)Mn 69	/652. BD* (1) C 46 - O 65	0.05	2.40	0.014
139. CR (9)Mn 69	/467. RY* (1) C 46	0.03	3.23	0.012
139. CR (9)Mn 69	/633. BD* (1) C 37 - N 42	0.14	2.45	0.023
139. CR (9)Mn 69	/652. BD* (1) C 46 - O 65	0.25	2.40	0.031
164. LP (3)Mn 69	/435. RY* (1) N 42	0.03	2.12	0.011
164. LP (3)Mn 69	/651. BD* (1) C 46 - C 48	0.03	0.68	0.006
164. LP (3)Mn 69	/652. BD* (1) C 46 - O 65	0.23	0.56	0.014
165. LP (5)Mn 69	/ 54. BD* (2) C 35 - C 37	0.08	0.17	0.006
165. LP (5)Mn 69	/396. RY* (2) C 37	0.04	1.30	0.009
165. LP (5)Mn 69	/435. RY* (1) N 42	1.41	2.11	0.069
165. LP (5)Mn 69	/436. RY* (2) N 42	0.26	1.16	0.022
165. LP (5)Mn 69	/437. RY* (3) N 42	0.46	1.44	0.033
165. LP (5)Mn 69	/441. RY* (7) N 42	0.03	35.43	0.045
165. LP (5)Mn 69	/443. RY* (1) N 43	0.05	1.08	0.009
165. LP (5)Mn 69	/445. RY* (3) N 43	0.06	1.30	0.011
165. LP (5)Mn 69	/446. RY* (4) N 43	0.03	1.82	0.009
165. LP (5)Mn 69	/467. RY* (1) C 46	0.07	1.37	0.013
165. LP (5)Mn 69	/470. RY* (4) C 46	0.05	0.97	0.009
165. LP (5)Mn 69	/541. RY* (1) O 65	0.09	2.60	0.020
165. LP (5)Mn 69	/544. RY* (4) O 65	0.05	1.90	0.013
165. LP (5)Mn 69	/633. BD* (1) C 37 - N 42	0.20	0.59	0.014
165. LP (5)Mn 69	/644. BD* (1) N 43 - C 44	0.08	0.58	0.009
165. LP (5)Mn 69	/652. BD* (1) C 46 - O 65	0.17	0.54	0.012
166. LP (6)Mn 69	/380. RY* (2) C 35	0.08	0.82	0.028
166. LP (6)Mn 69	/381. RY* (3) C 35	0.13	0.80	0.035
166. LP (6)Mn 69	/382. RY* (4) C 35	0.06	0.16	0.011
166. LP (6)Mn 69	/383. RY* (5) C 35	0.03	2.35	0.031
166. LP (6)Mn 69	/389. RY* (3) C 36	0.68	0.45	0.061
166. LP (6)Mn 69	/390. RY* (4) C 36	0.26	0.35	0.033
166. LP (6)Mn 69	/395. RY* (1) C 37	3.79	0.13	0.076
166. LP (6)Mn 69	/396. RY* (2) C 37	9.94	0.32	0.196
166. LP (6)Mn 69	/397. RY* (3) C 37	0.63	0.36	0.053
166. LP (6)Mn 69	/398. RY* (4) C 37	2.23	0.28	0.087
166. LP (6)Mn 69	/400. RY* (6) C 37	0.58	2.37	0.129
166. LP (6)Mn 69	/402. RY* (8) C 37	0.07	2.02	0.043
166. LP (6)Mn 69	/421. RY* (3) C 40	0.27	0.35	0.033
166. LP (6)Mn 69	/422. RY* (4) C 40	0.03	0.31	0.010
166. LP (6)Mn 69	/426. RY* (8) C 40	0.06	2.23	0.041
166. LP (6)Mn 69	/435. RY* (1) N 42	19.29	1.13	0.505
166. LP (6)Mn 69	/436. RY* (2) N 42	14.14	0.18	0.174
166. LP (6)Mn 69	/437. RY* (3) N 42	17.76	0.46	0.313
166. LP (6)Mn 69	/438. RY* (4) N 42	2.95	0.73	0.162
166. LP (6)Mn 69	/440. RY* (6) N 42	0.04	3.26	0.038
166. LP (6)Mn 69	/443. RY* (1) N 43	5.71	0.10	0.083
166. LP (6)Mn 69	/444. RY* (2) N 43	5.14	0.04	0.051
166. LP (6)Mn 69	/445. RY* (3) N 43	4.11	0.33	0.127
166. LP (6)Mn 69	/446. RY* (4) N 43	0.40	0.84	0.064
166. LP (6)Mn 69	/447. RY* (5) N 43	0.07	3.44	0.054
166. LP (6)Mn 69	/449. RY* (7) N 43	0.03	3.04	0.035
166. LP (6)Mn 69	/451. RY* (1) C 44	0.03	0.36	0.011
166. LP (6)Mn 69	/452. RY* (2) C 44	0.07	0.63	0.023
166. LP (6)Mn 69	/453. RY* (3) C 44	5.49	0.12	0.090
166. LP (6)Mn 69	/455. RY* (5) C 44	0.06	1.97	0.038
166. LP (6)Mn 69	/456. RY* (6) C 44	0.05	3.09	0.043
166. LP (6)Mn 69	/460. RY* (2) C 45	0.03	0.43	0.013
166. LP (6)Mn 69	/462. RY* (4) C 45	0.32	0.16	0.025
166. LP (6)Mn 69	/467. RY* (1) C 46	14.08	0.39	0.255
166. LP (6)Mn 69	/468. RY* (2) C 46	1.22	0.68	0.100
166. LP (6)Mn 69	/469. RY* (3) C 46	0.06	0.16	0.011
166. LP (6)Mn 69	/471. RY* (5) C 46	0.51	4.08	0.159
166. LP (6)Mn 69	/472. RY* (6) C 46	0.04	2.01	0.032
166. LP (6)Mn 69	/475. RY* (1) C 47	0.10	0.43	0.023
166. LP (6)Mn 69	/478. RY* (4) C 47	0.09	0.16	0.013
166. LP (6)Mn 69	/483. RY* (1) C 48	0.50	0.57	0.058
166. LP (6)Mn 69	/486. RY* (4) C 48	0.29	0.17	0.024
166. LP (6)Mn 69	/491. RY* (1) C 49	0.07	0.43	0.019
166. LP (6)Mn 69	/514. RY* (2) H 54	0.14	1.56	0.051
166. LP (6)Mn 69	/541. RY* (1) O 65	13.31	1.62	0.509

166.	LP	(6)Mn	69	/542.	RY*	(2)	O	65	0.14	0.31	0.023
166.	LP	(6)Mn	69	/543.	RY*	(3)	O	65	3.01	0.64	0.152
166.	LP	(6)Mn	69	/544.	RY*	(4)	O	65	1.70	0.92	0.138
166.	LP	(6)Mn	69	/545.	RY*	(5)	O	65	0.06	4.27	0.055
170.	LP	(1)Mn	69	/435.	RY*	(1)	N	42	0.09	2.12	0.017
170.	LP	(1)Mn	69	/541.	RY*	(1)	O	65	0.07	2.62	0.017
170.	LP	(1)Mn	69	/633.	BD*	(1)	C	37 - N 42	0.10	0.61	0.010
170.	LP	(1)Mn	69	/643.	BD*	(1)	N	42 - N 43	0.38	0.74	0.021
171.	LP	(2)Mn	69	/396.	RY*	(2)	C	37	0.03	1.32	0.008
171.	LP	(2)Mn	69	/435.	RY*	(1)	N	42	0.06	2.12	0.014
171.	LP	(2)Mn	69	/437.	RY*	(3)	N	42	0.04	1.45	0.010
171.	LP	(2)Mn	69	/451.	RY*	(1)	C	44	0.07	1.35	0.012
171.	LP	(2)Mn	69	/633.	BD*	(1)	C	37 - N 42	0.26	0.61	0.016
171.	LP	(2)Mn	69	/643.	BD*	(1)	N	42 - N 43	0.25	0.74	0.017
171.	LP	(2)Mn	69	/651.	BD*	(1)	C	46 - C 48	0.03	0.68	0.006
172.	LP	(4)Mn	69	/435.	RY*	(1)	N	42	0.18	2.11	0.025
172.	LP	(4)Mn	69	/436.	RY*	(2)	N	42	0.03	1.16	0.008
172.	LP	(4)Mn	69	/467.	RY*	(1)	C	46	0.07	1.37	0.012
172.	LP	(4)Mn	69	/541.	RY*	(1)	O	65	0.53	2.60	0.047
172.	LP	(4)Mn	69	/543.	RY*	(3)	O	65	0.07	1.61	0.013
172.	LP	(4)Mn	69	/643.	BD*	(1)	N	42 - N 43	0.15	0.72	0.013
172.	LP	(4)Mn	69	/647.	BD*	(2)	C	44 - C 46	0.03	0.23	0.004
172.	LP	(4)Mn	69	/652.	BD*	(1)	C	46 - O 65	0.15	0.54	0.011

from unit 5 to unit 3

135.	CR	(5)Mn	69	/660.	BD*	(1)	O	62 - H 63	0.05	2.40	0.013
139.	CR	(9)Mn	69	/661.	BD*	(1)	O	62 - H 64	0.03	2.40	0.011
164.	LP	(3)Mn	69	/660.	BD*	(1)	O	62 - H 63	0.17	0.56	0.012
164.	LP	(3)Mn	69	/661.	BD*	(1)	O	62 - H 64	0.28	0.56	0.016
165.	LP	(5)Mn	69	/529.	RY*	(1)	O	62	0.03	2.24	0.010
165.	LP	(5)Mn	69	/531.	RY*	(3)	O	62	0.03	1.57	0.008
165.	LP	(5)Mn	69	/532.	RY*	(4)	O	62	0.11	1.57	0.016
165.	LP	(5)Mn	69	/537.	RY*	(1)	H	63	0.20	0.77	0.016
165.	LP	(5)Mn	69	/539.	RY*	(1)	H	64	0.10	0.77	0.011
165.	LP	(5)Mn	69	/660.	BD*	(1)	O	62 - H 63	0.21	0.54	0.013
165.	LP	(5)Mn	69	/661.	BD*	(1)	O	62 - H 64	0.13	0.54	0.011
166.	LP	(6)Mn	69	/529.	RY*	(1)	O	62	24.33	1.26	0.608
166.	LP	(6)Mn	69	/530.	RY*	(2)	O	62	0.53	0.25	0.040
166.	LP	(6)Mn	69	/531.	RY*	(3)	O	62	5.88	0.59	0.205
166.	LP	(6)Mn	69	/532.	RY*	(4)	O	62	12.07	0.59	0.295
166.	LP	(6)Mn	69	/538.	RY*	(2)	H	63	0.04	1.44	0.025
170.	LP	(1)Mn	69	/531.	RY*	(3)	O	62	0.05	1.59	0.011
170.	LP	(1)Mn	69	/532.	RY*	(4)	O	62	0.03	1.59	0.009
170.	LP	(1)Mn	69	/537.	RY*	(1)	H	63	0.04	0.78	0.007
170.	LP	(1)Mn	69	/660.	BD*	(1)	O	62 - H 63	0.04	0.56	0.006
170.	LP	(1)Mn	69	/661.	BD*	(1)	O	62 - H 64	0.03	0.56	0.005
171.	LP	(2)Mn	69	/529.	RY*	(1)	O	62	0.04	2.26	0.012
171.	LP	(2)Mn	69	/661.	BD*	(1)	O	62 - H 64	0.10	0.56	0.010
172.	LP	(4)Mn	69	/529.	RY*	(1)	O	62	0.22	2.24	0.028
172.	LP	(4)Mn	69	/532.	RY*	(4)	O	62	0.03	1.57	0.009
172.	LP	(4)Mn	69	/537.	RY*	(1)	H	63	0.04	0.77	0.007
172.	LP	(4)Mn	69	/660.	BD*	(1)	O	62 - H 63	0.10	0.54	0.009

from unit 5 to unit 4

137.	CR	(7)Mn	69	/663.	BD*	(1)	H	67 - O 68	0.03	2.40	0.010
139.	CR	(9)Mn	69	/662.	BD*	(1)	H	66 - O 68	0.05	2.41	0.013
164.	LP	(3)Mn	69	/662.	BD*	(1)	H	66 - O 68	0.11	0.56	0.010
164.	LP	(3)Mn	69	/663.	BD*	(1)	H	67 - O 68	0.33	0.56	0.017
165.	LP	(5)Mn	69	/549.	RY*	(1)	H	66	0.18	0.90	0.016
165.	LP	(5)Mn	69	/551.	RY*	(1)	H	67	0.18	0.84	0.016
165.	LP	(5)Mn	69	/553.	RY*	(1)	O	68	0.17	2.14	0.024
165.	LP	(5)Mn	69	/555.	RY*	(3)	O	68	0.04	1.27	0.009
165.	LP	(5)Mn	69	/556.	RY*	(4)	O	68	0.05	2.01	0.013
165.	LP	(5)Mn	69	/662.	BD*	(1)	H	66 - O 68	0.07	0.55	0.008
165.	LP	(5)Mn	69	/663.	BD*	(1)	H	67 - O 68	0.03	0.55	0.005
166.	LP	(6)Mn	69	/550.	RY*	(2)	H	66	0.13	1.40	0.046
166.	LP	(6)Mn	69	/552.	RY*	(2)	H	67	0.08	1.44	0.037
166.	LP	(6)Mn	69	/553.	RY*	(1)	O	68	24.02	1.16	0.579
166.	LP	(6)Mn	69	/554.	RY*	(2)	O	68	1.90	0.22	0.070
166.	LP	(6)Mn	69	/555.	RY*	(3)	O	68	1.63	0.29	0.076
166.	LP	(6)Mn	69	/556.	RY*	(4)	O	68	8.27	1.04	0.322

170.	LP	(1)	Mn	69	/549.	RY*	(1)	H	66	0.03	0.92	0.006
170.	LP	(1)	Mn	69	/553.	RY*	(1)	O	68	0.03	2.16	0.011
170.	LP	(1)	Mn	69	/662.	BD*	(1)	H	66 - O 68	0.04	0.56	0.006
170.	LP	(1)	Mn	69	/663.	BD*	(1)	H	67 - O 68	0.03	0.56	0.005
171.	LP	(2)	Mn	69	/662.	BD*	(1)	H	66 - O 68	0.14	0.56	0.011
172.	LP	(4)	Mn	69	/551.	RY*	(1)	H	67	0.03	0.84	0.006
172.	LP	(4)	Mn	69	/553.	RY*	(1)	O	68	0.10	2.14	0.018
172.	LP	(4)	Mn	69	/662.	BD*	(1)	H	66 - O 68	0.13	0.55	0.011
172.	LP	(4)	Mn	69	/663.	BD*	(1)	H	67 - O 68	0.08	0.55	0.008
within unit 5														
132.	CR	(2)	Mn	69	/166.	LP	(6)	Mn	69	0.58	27.19	0.170
133.	CR	(3)	Mn	69	/166.	LP	(6)	Mn	69	2.21	4.88	0.140
165.	LP	(5)	Mn	69	/166.	LP	(6)	Mn	69	0.66	0.98	0.034
165.	LP	(5)	Mn	69	/167.	LP*	(7)	Mn	69	0.25	0.75	0.018
165.	LP	(5)	Mn	69	/562.	RY*	(2)	Mn	69	0.48	3.20	0.050
165.	LP	(5)	Mn	69	/581.	RY*	(21)	Mn	69	0.44	38.59	0.168
166.	LP	(6)	Mn	69	/169.	LP*	(9)	Mn	69	0.36	0.98	0.065
166.	LP	(6)	Mn	69	/562.	RY*	(2)	Mn	69	37.97	2.22	1.010
166.	LP	(6)	Mn	69	/564.	RY*	(4)	Mn	69	1.40	1.43	0.156
166.	LP	(6)	Mn	69	/565.	RY*	(5)	Mn	69	2.05	1.05	0.161
166.	LP	(6)	Mn	69	/567.	RY*	(7)	Mn	69	6.26	1.25	0.308
166.	LP	(6)	Mn	69	/568.	RY*	(8)	Mn	69	0.28	5.92	0.142
166.	LP	(6)	Mn	69	/574.	RY*	(14)	Mn	69	0.38	5.16	0.153
166.	LP	(6)	Mn	69	/575.	RY*	(15)	Mn	69	3.06	4.79	0.422
166.	LP	(6)	Mn	69	/578.	RY*	(18)	Mn	69	7.02	677.96	7.604
166.	LP	(6)	Mn	69	/580.	RY*	(20)	Mn	69	13.87	148.20	4.995
166.	LP	(6)	Mn	69	/581.	RY*	(21)	Mn	69	26.43	37.61	3.474

Table (S10): Summary of Natural Population Analysis of the Fe(III) complex:

Fe(III) complex

Atom	No	Natural Charge	Core	Valence	Rydberg	Total
C	1	0.34207	1.99860	3.64275	0.01658	5.65793
C	2	-0.09671	1.99880	4.08565	0.01226	6.09671
C	3	-0.25998	1.99898	4.24769	0.01331	6.25998
C	4	-0.05333	1.99882	4.04048	0.01403	6.05333
C	5	-0.17914	1.99890	4.16773	0.01251	6.17914
C	6	0.15212	1.99874	3.82795	0.02118	5.84788
C	7	-0.19673	1.99903	4.18365	0.01405	6.19673
C	8	-0.19280	1.99908	4.18014	0.01358	6.19280
C	9	-0.18513	1.99902	4.17168	0.01443	6.18513
C	10	-0.19384	1.99905	4.18165	0.01314	6.19384
N	11	-0.48047	1.99931	5.45278	0.02839	7.48047
N	12	-0.16012	1.99929	5.13818	0.02265	7.16012
C	13	0.00772	1.99869	3.97330	0.02029	5.99228
C	14	-0.18372	1.99902	4.17066	0.01404	6.18372
C	15	0.24243	1.99864	3.73645	0.02247	5.75757
C	16	-0.21395	1.99908	4.20151	0.01336	6.21395
C	17	-0.22742	1.99900	4.21412	0.01430	6.22742
C	18	-0.19654	1.99908	4.18401	0.01344	6.19654
O	19	-0.79132	1.99982	6.78416	0.00735	8.79132
O	20	-0.69894	1.99978	6.69558	0.00358	8.69894
H	21	0.22893	0.00000	0.76855	0.00251	0.77107
H	22	0.20887	0.00000	0.78825	0.00289	0.79113
H	23	0.21648	0.00000	0.78235	0.00117	0.78352
H	24	0.22458	0.00000	0.77168	0.00374	0.77542
H	25	0.21629	0.00000	0.78191	0.00180	0.78371
H	26	0.22391	0.00000	0.77356	0.00253	0.77609
H	27	0.21707	0.00000	0.78177	0.00116	0.78293
H	28	0.22621	0.00000	0.77160	0.00220	0.77379
H	29	0.21600	0.00000	0.78275	0.00125	0.78400

H	30	0.49103	0.00000	0.50812	0.00085	0.50897
H	31	0.22581	0.00000	0.77201	0.00219	0.77419
C	32	0.34781	1.99860	3.63709	0.01651	5.65219
C	33	-0.09342	1.99880	4.08238	0.01223	6.09342
C	34	-0.25873	1.99897	4.24623	0.01352	6.25873
C	35	-0.05022	1.99883	4.03808	0.01332	6.05022
C	36	-0.20384	1.99890	4.19230	0.01264	6.20384
C	37	0.14609	1.99874	3.83425	0.02092	5.85391
C	38	-0.19686	1.99903	4.18388	0.01396	6.19686
C	39	-0.18714	1.99908	4.17450	0.01357	6.18714
C	40	-0.17247	1.99902	4.15975	0.01370	6.17247
C	41	-0.19472	1.99906	4.18225	0.01341	6.19472
N	42	-0.36179	1.99928	5.33589	0.02662	7.36179
N	43	-0.21511	1.99928	5.19181	0.02402	7.21511
C	44	0.12654	1.99874	3.85565	0.01907	5.87346
C	45	-0.20723	1.99902	4.19384	0.01437	6.20723
C	46	0.27385	1.99865	3.70604	0.02147	5.72615
C	47	-0.14118	1.99908	4.12939	0.01271	6.14118
C	48	-0.20031	1.99900	4.18664	0.01466	6.20031
C	49	-0.21049	1.99907	4.19812	0.01330	6.21049
O	50	-0.69603	1.99978	6.69270	0.00355	8.69603
H	51	0.23091	0.00000	0.76663	0.00247	0.76909
H	52	0.20890	0.00000	0.78821	0.00289	0.79110
H	53	0.21696	0.00000	0.78183	0.00121	0.78304
H	54	0.22818	0.00000	0.76889	0.00294	0.77182
H	55	0.21657	0.00000	0.78219	0.00123	0.78343
H	56	0.23065	0.00000	0.76672	0.00263	0.76935
H	57	0.22253	0.00000	0.77635	0.00112	0.77747
H	58	0.23335	0.00000	0.76363	0.00302	0.76665
H	59	0.22087	0.00000	0.77785	0.00127	0.77913
H	60	0.49274	0.00000	0.50645	0.00081	0.50726
H	61	0.22603	0.00000	0.77076	0.00321	0.77397
O	62	-0.91532	1.99980	6.90886	0.00666	8.91532
H	63	0.52542	0.00000	0.47249	0.00209	0.47458
H	64	0.50767	0.00000	0.48942	0.00291	0.49233
O	65	-0.64494	1.99981	6.63860	0.00653	8.64494
Cl	66	-0.58061	9.99977	7.57826	0.00258	17.58061
Fe	67	1.74594	17.99113	6.24533	0.01760	24.25406

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* Total * 0.00000 109.95218 216.36588 0.68194 327.00000
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Table (S11): NBO data of the Fe(III) complex:

Donor NBO (i)		Acceptor NBO (j)		kcal/mol	a.u.	a.u.
within unit 1						
1.	BD (1) C	1 - C	2	/189. RY* (1) C	3	0.67 1.97 0.046
1.	BD (1) C	1 - C	2	/190. RY* (2) C	3	0.31 2.00 0.032
1.	BD (1) C	1 - C	2	/197. RY* (1) C	4	0.65 2.18 0.048
1.	BD (1) C	1 - C	2	/221. RY* (1) C	7	0.57 1.85 0.041
1.	BD (1) C	1 - C	2	/583. BD* (1) C	1 - C 3	1.87 1.27 0.062
1.	BD (1) C	1 - C	2	/586. BD* (1) C	2 - C 4	1.75 1.27 0.060
1.	BD (1) C	1 - C	2	/588. BD* (1) C	2 - C 7	0.74 1.09 0.036
1.	BD (1) C	1 - C	2	/590. BD* (1) C	3 - H 21	1.44 1.09 0.050
1.	BD (1) C	1 - C	2	/592. BD* (1) C	4 - C 9	2.09 1.10 0.061
1.	BD (1) C	1 - C	2	/597. BD* (1) C	7 - C 8	0.65 1.29 0.037
2.	BD (1) C	1 - C	3	/182. RY* (2) C	2	0.80 2.19 0.053
2.	BD (1) C	1 - C	3	/206. RY* (2) C	5	0.61 1.97 0.044
2.	BD (1) C	1 - C	3	/582. BD* (1) C	1 - C 2	1.84 1.18 0.059
2.	BD (1) C	1 - C	3	/585. BD* (1) C	1 - O 20	0.42 1.03 0.026
2.	BD (1) C	1 - C	3	/588. BD* (1) C	2 - C 7	1.22 1.20 0.048
2.	BD (1) C	1 - C	3	/589. BD* (1) C	3 - C 5	1.21 1.22 0.049
2.	BD (1) C	1 - C	3	/590. BD* (1) C	3 - H 21	0.89 1.19 0.041
2.	BD (1) C	1 - C	3	/595. BD* (1) C	5 - H 31	0.70 1.20 0.037
2.	BD (1) C	1 - C	3	/622. BD* (1) O	20 - H 30	0.80 1.07 0.037
3.	BD (2) C	1 - C	3	/184. RY* (4) C	2	0.35 0.89 0.023
3.	BD (2) C	1 - C	3	/325. RY* (1) O	20	0.33 1.19 0.026
3.	BD (2) C	1 - C	3	/584. BD* (2) C	1 - C 3	0.28 0.31 0.012
3.	BD (2) C	1 - C	3	/587. BD* (2) C	2 - C 4	5.77 0.32 0.057
3.	BD (2) C	1 - C	3	/594. BD* (2) C	5 - C 6	6.65 0.32 0.059
4.	BD (1) C	1 - O	20	/173. RY* (1) C	1	0.33 1.97 0.032
4.	BD (1) C	1 - O	20	/189. RY* (1) C	3	0.29 2.21 0.032
4.	BD (1) C	1 - O	20	/582. BD* (1) C	1 - C 2	0.27 1.32 0.024
4.	BD (1) C	1 - O	20	/583. BD* (1) C	1 - C 3	0.54 1.52 0.036
4.	BD (1) C	1 - O	20	/586. BD* (1) C	2 - C 4	0.58 1.52 0.038
4.	BD (1) C	1 - O	20	/589. BD* (1) C	3 - C 5	1.10 1.36 0.049
5.	BD (1) C	2 - C	4	/174. RY* (2) C	1	0.44 2.15 0.039
5.	BD (1) C	2 - C	4	/213. RY* (1) C	6	0.40 1.88 0.035
5.	BD (1) C	2 - C	4	/214. RY* (2) C	6	0.35 1.71 0.031
5.	BD (1) C	2 - C	4	/221. RY* (1) C	7	0.35 1.94 0.033
5.	BD (1) C	2 - C	4	/237. RY* (1) C	9	0.36 1.94 0.034
5.	BD (1) C	2 - C	4	/582. BD* (1) C	1 - C 2	1.65 1.16 0.055
5.	BD (1) C	2 - C	4	/585. BD* (1) C	1 - O 20	1.34 1.01 0.047
5.	BD (1) C	2 - C	4	/588. BD* (1) C	2 - C 7	1.91 1.18 0.060
5.	BD (1) C	2 - C	4	/591. BD* (1) C	4 - C 6	1.99 1.19 0.062
5.	BD (1) C	2 - C	4	/592. BD* (1) C	4 - C 9	1.83 1.19 0.059
5.	BD (1) C	2 - C	4	/596. BD* (1) C	6 - N 11	1.39 1.11 0.050
5.	BD (1) C	2 - C	4	/599. BD* (1) C	7 - H 22	0.76 1.18 0.038
5.	BD (1) C	2 - C	4	/604. BD* (1) C	9 - H 24	0.69 1.19 0.036
6.	BD (2) C	2 - C	4	/176. RY* (4) C	1	0.30 0.84 0.022
6.	BD (2) C	2 - C	4	/223. RY* (3) C	7	0.55 0.77 0.028
6.	BD (2) C	2 - C	4	/239. RY* (3) C	9	0.50 0.86 0.028
6.	BD (2) C	2 - C	4	/584. BD* (2) C	1 - C 3	6.40 0.30 0.056
6.	BD (2) C	2 - C	4	/594. BD* (2) C	5 - C 6	6.35 0.31 0.057
6.	BD (2) C	2 - C	4	/598. BD* (2) C	7 - C 8	6.00 0.31 0.056
6.	BD (2) C	2 - C	4	/603. BD* (2) C	9 - C 10	5.79 0.31 0.056
7.	BD (1) C	2 - C	7	/173. RY* (1) C	1	0.29 1.72 0.028
7.	BD (1) C	2 - C	7	/174. RY* (2) C	1	0.36 2.06 0.035
7.	BD (1) C	2 - C	7	/197. RY* (1) C	4	0.41 2.17 0.038
7.	BD (1) C	2 - C	7	/198. RY* (2) C	4	0.28 2.18 0.031
7.	BD (1) C	2 - C	7	/229. RY* (1) C	8	0.36 1.83 0.033
7.	BD (1) C	2 - C	7	/230. RY* (2) C	8	0.78 1.94 0.049
7.	BD (1) C	2 - C	7	/582. BD* (1) C	1 - C 2	0.62 1.06 0.032
7.	BD (1) C	2 - C	7	/583. BD* (1) C	1 - C 3	0.71 1.27 0.038
7.	BD (1) C	2 - C	7	/586. BD* (1) C	2 - C 4	2.16 1.26 0.066
7.	BD (1) C	2 - C	7	/591. BD* (1) C	4 - C 6	2.35 1.10 0.064
7.	BD (1) C	2 - C	7	/597. BD* (1) C	7 - C 8	1.40 1.28 0.054
7.	BD (1) C	2 - C	7	/599. BD* (1) C	7 - H 22	0.31 1.08 0.023
7.	BD (1) C	2 - C	7	/601. BD* (1) C	8 - H 23	1.45 1.09 0.050

8.	BD	(1)	C	3	-	C	5	/173.	RY*	(1)	C	1	0.63	1.72	0.042			
8.	BD	(1)	C	3	-	C	5	/174.	RY*	(2)	C	1	0.38	2.06	0.036			
8.	BD	(1)	C	3	-	C	5	/214.	RY*	(2)	C	6	0.29	1.62	0.028			
8.	BD	(1)	C	3	-	C	5	/215.	RY*	(3)	C	6	0.72	1.77	0.045			
8.	BD	(1)	C	3	-	C	5	/583.	BD*	(1)	C	1	-	C	3	1.29	1.27	0.051
8.	BD	(1)	C	3	-	C	5	/585.	BD*	(1)	C	1	-	O	20	2.53	0.92	0.061
8.	BD	(1)	C	3	-	C	5	/590.	BD*	(1)	C	3	-	H	21	0.37	1.08	0.025
8.	BD	(1)	C	3	-	C	5	/593.	BD*	(1)	C	5	-	C	6	1.50	1.29	0.056
8.	BD	(1)	C	3	-	C	5	/595.	BD*	(1)	C	5	-	H	31	0.30	1.08	0.023
8.	BD	(1)	C	3	-	C	5	/596.	BD*	(1)	C	6	-	N	11	3.00	1.02	0.070
9.	BD	(1)	C	3	-	H	21	/174.	RY*	(2)	C	1	0.46	1.93	0.038			
9.	BD	(1)	C	3	-	H	21	/178.	RY*	(6)	C	1	0.27	8.78	0.062			
9.	BD	(1)	C	3	-	H	21	/205.	RY*	(1)	C	5	0.50	1.82	0.038			
9.	BD	(1)	C	3	-	H	21	/582.	BD*	(1)	C	1	-	C	2	3.06	0.93	0.068
9.	BD	(1)	C	3	-	H	21	/583.	BD*	(1)	C	1	-	C	3	0.71	1.13	0.036
9.	BD	(1)	C	3	-	H	21	/585.	BD*	(1)	C	1	-	O	20	0.84	0.79	0.032
9.	BD	(1)	C	3	-	H	21	/593.	BD*	(1)	C	5	-	C	6	1.41	1.15	0.051
10.	BD	(1)	C	4	-	C	6	/181.	RY*	(1)	C	2	0.74	2.30	0.052			
10.	BD	(1)	C	4	-	C	6	/205.	RY*	(1)	C	5	0.46	1.95	0.038			
10.	BD	(1)	C	4	-	C	6	/237.	RY*	(1)	C	9	0.59	1.85	0.042			
10.	BD	(1)	C	4	-	C	6	/253.	RY*	(1)	N	11	0.32	1.74	0.030			
10.	BD	(1)	C	4	-	C	6	/586.	BD*	(1)	C	2	-	C	4	2.02	1.27	0.064
10.	BD	(1)	C	4	-	C	6	/588.	BD*	(1)	C	2	-	C	7	2.02	1.08	0.059
10.	BD	(1)	C	4	-	C	6	/592.	BD*	(1)	C	4	-	C	9	0.91	1.10	0.040
10.	BD	(1)	C	4	-	C	6	/593.	BD*	(1)	C	5	-	C	6	2.05	1.29	0.065
10.	BD	(1)	C	4	-	C	6	/595.	BD*	(1)	C	5	-	H	31	1.51	1.09	0.051
10.	BD	(1)	C	4	-	C	6	/596.	BD*	(1)	C	6	-	N	11	0.43	1.02	0.027
10.	BD	(1)	C	4	-	C	6	/602.	BD*	(1)	C	9	-	C	10	0.70	1.30	0.038
11.	BD	(1)	C	4	-	C	9	/181.	RY*	(1)	C	2	0.34	2.29	0.036			
11.	BD	(1)	C	4	-	C	9	/182.	RY*	(2)	C	2	0.25	2.07	0.029			
11.	BD	(1)	C	4	-	C	9	/213.	RY*	(1)	C	6	0.76	1.78	0.047			
11.	BD	(1)	C	4	-	C	9	/214.	RY*	(2)	C	6	0.29	1.61	0.027			
11.	BD	(1)	C	4	-	C	9	/245.	RY*	(1)	C	10	0.32	1.90	0.031			
11.	BD	(1)	C	4	-	C	9	/246.	RY*	(2)	C	10	0.81	1.95	0.051			
11.	BD	(1)	C	4	-	C	9	/582.	BD*	(1)	C	1	-	C	2	2.31	1.06	0.063
11.	BD	(1)	C	4	-	C	9	/586.	BD*	(1)	C	2	-	C	4	2.10	1.26	0.065
11.	BD	(1)	C	4	-	C	9	/591.	BD*	(1)	C	4	-	C	6	0.95	1.09	0.041
11.	BD	(1)	C	4	-	C	9	/593.	BD*	(1)	C	5	-	C	6	0.99	1.28	0.045
11.	BD	(1)	C	4	-	C	9	/602.	BD*	(1)	C	9	-	C	10	1.43	1.29	0.054
11.	BD	(1)	C	4	-	C	9	/604.	BD*	(1)	C	9	-	H	24	0.33	1.09	0.024
11.	BD	(1)	C	4	-	C	9	/605.	BD*	(1)	C	10	-	H	25	1.52	1.08	0.051
12.	BD	(1)	C	5	-	C	6	/190.	RY*	(2)	C	3	0.73	2.11	0.050			
12.	BD	(1)	C	5	-	C	6	/198.	RY*	(2)	C	4	0.73	2.29	0.052			
12.	BD	(1)	C	5	-	C	6	/253.	RY*	(1)	N	11	0.59	1.85	0.042			
12.	BD	(1)	C	5	-	C	6	/589.	BD*	(1)	C	3	-	C	5	1.37	1.22	0.052
12.	BD	(1)	C	5	-	C	6	/590.	BD*	(1)	C	3	-	H	21	0.65	1.19	0.035
12.	BD	(1)	C	5	-	C	6	/591.	BD*	(1)	C	4	-	C	6	2.11	1.21	0.064
12.	BD	(1)	C	5	-	C	6	/592.	BD*	(1)	C	4	-	C	9	1.15	1.21	0.047
12.	BD	(1)	C	5	-	C	6	/595.	BD*	(1)	C	5	-	H	31	0.96	1.19	0.043
12.	BD	(1)	C	5	-	C	6	/596.	BD*	(1)	C	6	-	N	11	0.98	1.13	0.042
12.	BD	(1)	C	5	-	C	6	/606.	BD*	(1)	N	11	-	N	12	0.62	1.25	0.035
13.	BD	(2)	C	5	-	C	6	/191.	RY*	(3)	C	3	0.56	0.78	0.028			
13.	BD	(2)	C	5	-	C	6	/200.	RY*	(4)	C	4	0.34	0.95	0.024			
13.	BD	(2)	C	5	-	C	6	/584.	BD*	(2)	C	1	-	C	3	6.54	0.31	0.058
13.	BD	(2)	C	5	-	C	6	/587.	BD*	(2)	C	2	-	C	4	6.25	0.32	0.059
13.	BD	(2)	C	5	-	C	6	/606.	BD*	(1)	N	11	-	N	12	1.88	0.77	0.050
14.	BD	(1)	C	5	-	H	31	/189.	RY*	(1)	C	3	0.49	1.83	0.038			
14.	BD	(1)	C	5	-	H	31	/213.	RY*	(1)	C	6	0.67	1.66	0.042			
14.	BD	(1)	C	5	-	H	31	/214.	RY*	(2)	C	6	0.28	1.49	0.026			
14.	BD	(1)	C	5	-	H	31	/583.	BD*	(1)	C	1	-	C	3	1.34	1.14	0.049
14.	BD	(1)	C	5	-	H	31	/590.	BD*	(1)	C	3	-	H	21	0.25	0.95	0.020
14.	BD	(1)	C	5	-	H	31	/591.	BD*	(1)	C	4	-	C	6	3.28	0.97	0.072
14.	BD	(1)	C	5	-	H	31	/593.	BD*	(1)	C	5	-	C	6	0.87	1.16	0.040
14.	BD	(1)	C	5	-	H	31	/596.	BD*	(1)	C	6	-	N	11	0.65	0.89	0.031
15.	BD	(1)	C	6	-	N	11	/198.	RY*	(2)	C	4	0.61	2.33	0.048			
15.	BD	(1)	C	6	-	N	11	/205.	RY*	(1)	C	5	0.57	2.09	0.044			
15.	BD	(1)	C	6	-	N	11	/261.	RY*	(1)	N	12	0.74	1.61	0.044			
15.	BD	(1)	C	6	-	N	11	/262.	RY*	(2)	N	12	0.31	1.65	0.029			
15.	BD	(1)	C	6	-	N	11	/586.	BD*	(1)	C	2	-	C	4	0.73	1.41	0.040
15.	BD	(1)	C	6	-	N	11	/589.	BD*	(1)	C	3	-	C	5	1.23	1.25	0.050
15.	BD	(1)	C	6	-	N	11	/591.	BD*	(1)	C	4	-	C	6	0.63	1.24	0.035

15. BD (1) C 6 - N 11	/593. BD* (1) C 5 - C 6	1.15	1.43	0.051
15. BD (1) C 6 - N 11	/606. BD* (1) N 11 - N 12	0.74	1.28	0.039
15. BD (1) C 6 - N 11	/607. BD* (1) N 12 - C 13	2.74	1.16	0.071
16. BD (1) C 7 - C 8	/181. RY* (1) C 2	0.59	2.40	0.048
16. BD (1) C 7 - C 8	/182. RY* (2) C 2	0.36	2.18	0.036
16. BD (1) C 7 - C 8	/245. RY* (1) C 10	0.46	2.01	0.039
16. BD (1) C 7 - C 8	/246. RY* (2) C 10	0.41	2.06	0.037
16. BD (1) C 7 - C 8	/582. BD* (1) C 1 - C 2	1.35	1.17	0.050
16. BD (1) C 7 - C 8	/588. BD* (1) C 2 - C 7	1.49	1.19	0.053
16. BD (1) C 7 - C 8	/599. BD* (1) C 7 - H 22	0.88	1.19	0.041
16. BD (1) C 7 - C 8	/600. BD* (1) C 8 - C 10	1.21	1.21	0.048
16. BD (1) C 7 - C 8	/601. BD* (1) C 8 - H 23	0.80	1.19	0.039
16. BD (1) C 7 - C 8	/605. BD* (1) C 10 - H 25	0.78	1.19	0.039
17. BD (2) C 7 - C 8	/184. RY* (4) C 2	0.35	0.89	0.024
17. BD (2) C 7 - C 8	/247. RY* (3) C 10	0.46	0.82	0.026
17. BD (2) C 7 - C 8	/587. BD* (2) C 2 - C 4	6.09	0.31	0.058
17. BD (2) C 7 - C 8	/603. BD* (2) C 9 - C 10	6.43	0.32	0.057
18. BD (1) C 7 - H 22	/181. RY* (1) C 2	0.72	2.18	0.050
18. BD (1) C 7 - H 22	/229. RY* (1) C 8	0.69	1.72	0.044
18. BD (1) C 7 - H 22	/586. BD* (1) C 2 - C 4	1.61	1.15	0.054
18. BD (1) C 7 - H 22	/597. BD* (1) C 7 - C 8	0.75	1.17	0.037
18. BD (1) C 7 - H 22	/600. BD* (1) C 8 - C 10	2.74	0.99	0.066
18. BD (1) C 7 - H 22	/601. BD* (1) C 8 - H 23	0.48	0.97	0.027
19. BD (1) C 8 - C 10	/222. RY* (2) C 7	0.92	1.76	0.051
19. BD (1) C 8 - C 10	/238. RY* (2) C 9	0.96	1.70	0.051
19. BD (1) C 8 - C 10	/597. BD* (1) C 7 - C 8	1.33	1.29	0.052
19. BD (1) C 8 - C 10	/599. BD* (1) C 7 - H 22	1.58	1.08	0.052
19. BD (1) C 8 - C 10	/601. BD* (1) C 8 - H 23	0.30	1.09	0.023
19. BD (1) C 8 - C 10	/602. BD* (1) C 9 - C 10	1.37	1.30	0.053
19. BD (1) C 8 - C 10	/604. BD* (1) C 9 - H 24	1.63	1.10	0.054
19. BD (1) C 8 - C 10	/605. BD* (1) C 10 - H 25	0.29	1.09	0.022
20. BD (1) C 8 - H 23	/221. RY* (1) C 7	0.76	1.72	0.046
20. BD (1) C 8 - H 23	/245. RY* (1) C 10	0.66	1.77	0.043
20. BD (1) C 8 - H 23	/588. BD* (1) C 2 - C 7	3.14	0.95	0.069
20. BD (1) C 8 - H 23	/597. BD* (1) C 7 - C 8	0.71	1.16	0.036
20. BD (1) C 8 - H 23	/599. BD* (1) C 7 - H 22	0.54	0.95	0.029
20. BD (1) C 8 - H 23	/602. BD* (1) C 9 - C 10	1.40	1.17	0.051
21. BD (1) C 9 - C 10	/197. RY* (1) C 4	0.75	2.27	0.052
21. BD (1) C 9 - C 10	/198. RY* (2) C 4	0.30	2.29	0.033
21. BD (1) C 9 - C 10	/229. RY* (1) C 8	0.42	1.94	0.036
21. BD (1) C 9 - C 10	/230. RY* (2) C 8	0.43	2.04	0.038
21. BD (1) C 9 - C 10	/591. BD* (1) C 4 - C 6	1.26	1.20	0.050
21. BD (1) C 9 - C 10	/592. BD* (1) C 4 - C 9	1.47	1.20	0.053
21. BD (1) C 9 - C 10	/600. BD* (1) C 8 - C 10	1.24	1.21	0.049
21. BD (1) C 9 - C 10	/601. BD* (1) C 8 - H 23	0.78	1.19	0.038
21. BD (1) C 9 - C 10	/604. BD* (1) C 9 - H 24	0.85	1.20	0.040
21. BD (1) C 9 - C 10	/605. BD* (1) C 10 - H 25	0.83	1.19	0.040
22. BD (2) C 9 - C 10	/200. RY* (4) C 4	0.26	0.94	0.021
22. BD (2) C 9 - C 10	/231. RY* (3) C 8	0.48	0.78	0.026
22. BD (2) C 9 - C 10	/587. BD* (2) C 2 - C 4	6.46	0.31	0.059
22. BD (2) C 9 - C 10	/598. BD* (2) C 7 - C 8	7.04	0.31	0.059
23. BD (1) C 9 - H 24	/197. RY* (1) C 4	0.71	2.04	0.048
23. BD (1) C 9 - H 24	/245. RY* (1) C 10	0.65	1.77	0.043
23. BD (1) C 9 - H 24	/586. BD* (1) C 2 - C 4	1.75	1.13	0.056
23. BD (1) C 9 - H 24	/600. BD* (1) C 8 - C 10	2.73	0.98	0.065
23. BD (1) C 9 - H 24	/602. BD* (1) C 9 - C 10	0.78	1.16	0.038
23. BD (1) C 9 - H 24	/605. BD* (1) C 10 - H 25	0.49	0.96	0.027
24. BD (1) C 10 - H 25	/229. RY* (1) C 8	0.52	1.70	0.038
24. BD (1) C 10 - H 25	/237. RY* (1) C 9	0.64	1.71	0.042
24. BD (1) C 10 - H 25	/592. BD* (1) C 4 - C 9	3.04	0.96	0.068
24. BD (1) C 10 - H 25	/597. BD* (1) C 7 - C 8	1.36	1.16	0.050
24. BD (1) C 10 - H 25	/602. BD* (1) C 9 - C 10	0.77	1.16	0.038
24. BD (1) C 10 - H 25	/604. BD* (1) C 9 - H 24	0.50	0.96	0.028
25. BD (1) N 11 - N 12	/213. RY* (1) C 6	0.58	2.12	0.044
25. BD (1) N 11 - N 12	/269. RY* (1) C 13	0.56	2.09	0.043
25. BD (1) N 11 - N 12	/270. RY* (2) C 13	0.98	2.41	0.062
25. BD (1) N 11 - N 12	/593. BD* (1) C 5 - C 6	0.37	1.62	0.031
25. BD (1) N 11 - N 12	/594. BD* (2) C 5 - C 6	0.41	1.02	0.028
25. BD (1) N 11 - N 12	/596. BD* (1) C 6 - N 11	0.70	1.36	0.039
25. BD (1) N 11 - N 12	/607. BD* (1) N 12 - C 13	0.27	1.36	0.024
25. BD (1) N 11 - N 12	/608. BD* (1) C 13 - C 14	0.66	1.43	0.039

26. BD (1) N 12 - C 13	/213. RY* (1) C 6	0.31	1.93	0.031
26. BD (1) N 12 - C 13	/254. RY* (2) N 11	0.73	2.53	0.054
26. BD (1) N 12 - C 13	/255. RY* (3) N 11	0.51	2.02	0.041
26. BD (1) N 12 - C 13	/277. RY* (1) C 14	0.43	2.06	0.038
26. BD (1) N 12 - C 13	/285. RY* (1) C 15	0.49	1.91	0.039
26. BD (1) N 12 - C 13	/596. BD* (1) C 6 - N 11	2.42	1.17	0.067
26. BD (1) N 12 - C 13	/608. BD* (1) C 13 - C 14	0.42	1.23	0.029
26. BD (1) N 12 - C 13	/609. BD* (1) C 13 - C 15	1.21	1.40	0.052
26. BD (1) N 12 - C 13	/611. BD* (1) C 14 - C 16	0.44	1.43	0.032
26. BD (1) N 12 - C 13	/614. BD* (1) C 15 - C 17	1.47	1.23	0.054
27. BD (1) C 13 - C 14	/261. RY* (1) N 12	0.35	1.48	0.029
27. BD (1) C 13 - C 14	/287. RY* (3) C 15	0.44	1.77	0.035
27. BD (1) C 13 - C 14	/293. RY* (1) C 16	0.43	1.85	0.036
27. BD (1) C 13 - C 14	/294. RY* (2) C 16	0.75	1.98	0.049
27. BD (1) C 13 - C 14	/606. BD* (1) N 11 - N 12	1.49	1.15	0.052
27. BD (1) C 13 - C 14	/607. BD* (1) N 12 - C 13	0.60	1.03	0.031
27. BD (1) C 13 - C 14	/609. BD* (1) C 13 - C 15	2.20	1.27	0.067
27. BD (1) C 13 - C 14	/611. BD* (1) C 14 - C 16	1.22	1.30	0.050
27. BD (1) C 13 - C 14	/613. BD* (1) C 14 - H 26	0.30	1.09	0.023
27. BD (1) C 13 - C 14	/615. BD* (1) C 15 - O 19	2.60	0.97	0.063
27. BD (1) C 13 - C 14	/617. BD* (1) C 16 - H 27	1.47	1.10	0.051
28. BD (1) C 13 - C 15	/263. RY* (3) N 12	0.29	1.84	0.029
28. BD (1) C 13 - C 15	/277. RY* (1) C 14	0.52	2.05	0.041
28. BD (1) C 13 - C 15	/301. RY* (1) C 17	0.68	2.07	0.048
28. BD (1) C 13 - C 15	/607. BD* (1) N 12 - C 13	0.83	1.15	0.039
28. BD (1) C 13 - C 15	/608. BD* (1) C 13 - C 14	2.16	1.22	0.065
28. BD (1) C 13 - C 15	/613. BD* (1) C 14 - H 26	0.63	1.21	0.035
28. BD (1) C 13 - C 15	/614. BD* (1) C 15 - C 17	1.89	1.22	0.061
28. BD (1) C 13 - C 15	/615. BD* (1) C 15 - O 19	0.41	1.09	0.027
28. BD (1) C 13 - C 15	/620. BD* (1) C 17 - H 28	0.65	1.22	0.035
29. BD (2) C 13 - C 15	/262. RY* (2) N 12	0.29	1.16	0.025
29. BD (2) C 13 - C 15	/279. RY* (3) C 14	0.49	0.84	0.027
29. BD (2) C 13 - C 15	/303. RY* (3) C 17	0.60	0.87	0.031
29. BD (2) C 13 - C 15	/612. BD* (2) C 14 - C 16	5.37	0.34	0.054
29. BD (2) C 13 - C 15	/619. BD* (2) C 17 - C 18	7.33	0.33	0.063
30. BD (1) C 14 - C 16	/270. RY* (2) C 13	0.91	2.18	0.056
30. BD (1) C 14 - C 16	/309. RY* (1) C 18	0.50	1.94	0.039
30. BD (1) C 14 - C 16	/310. RY* (2) C 18	0.42	2.06	0.037
30. BD (1) C 14 - C 16	/607. BD* (1) N 12 - C 13	1.44	1.12	0.051
30. BD (1) C 14 - C 16	/608. BD* (1) C 13 - C 14	1.18	1.19	0.047
30. BD (1) C 14 - C 16	/613. BD* (1) C 14 - H 26	0.85	1.19	0.040
30. BD (1) C 14 - C 16	/616. BD* (1) C 16 - C 18	1.16	1.21	0.047
30. BD (1) C 14 - C 16	/617. BD* (1) C 16 - H 27	0.79	1.20	0.039
30. BD (1) C 14 - C 16	/621. BD* (1) C 18 - H 29	0.79	1.20	0.039
31. BD (2) C 14 - C 16	/272. RY* (4) C 13	0.26	0.93	0.020
31. BD (2) C 14 - C 16	/311. RY* (3) C 18	0.46	0.79	0.025
31. BD (2) C 14 - C 16	/610. BD* (2) C 13 - C 15	6.92	0.29	0.061
31. BD (2) C 14 - C 16	/619. BD* (2) C 17 - C 18	5.76	0.31	0.054
32. BD (1) C 14 - H 26	/269. RY* (1) C 13	0.40	1.63	0.032
32. BD (1) C 14 - H 26	/293. RY* (1) C 16	0.65	1.71	0.042
32. BD (1) C 14 - H 26	/607. BD* (1) N 12 - C 13	0.31	0.89	0.021
32. BD (1) C 14 - H 26	/609. BD* (1) C 13 - C 15	1.64	1.13	0.055
32. BD (1) C 14 - H 26	/611. BD* (1) C 14 - C 16	0.74	1.16	0.037
32. BD (1) C 14 - H 26	/616. BD* (1) C 16 - C 18	2.75	0.98	0.066
32. BD (1) C 14 - H 26	/617. BD* (1) C 16 - H 27	0.50	0.96	0.028
33. BD (1) C 15 - C 17	/269. RY* (1) C 13	0.53	1.77	0.039
33. BD (1) C 15 - C 17	/271. RY* (3) C 13	0.30	1.58	0.028
33. BD (1) C 15 - C 17	/309. RY* (1) C 18	0.40	1.85	0.034
33. BD (1) C 15 - C 17	/310. RY* (2) C 18	0.72	1.96	0.048
33. BD (1) C 15 - C 17	/607. BD* (1) N 12 - C 13	2.53	1.03	0.064
33. BD (1) C 15 - C 17	/609. BD* (1) C 13 - C 15	2.08	1.27	0.065
33. BD (1) C 15 - C 17	/618. BD* (1) C 17 - C 18	1.32	1.30	0.053
33. BD (1) C 15 - C 17	/620. BD* (1) C 17 - H 28	0.30	1.10	0.023
33. BD (1) C 15 - C 17	/621. BD* (1) C 18 - H 29	1.50	1.10	0.052
34. BD (1) C 15 - O 19	/269. RY* (1) C 13	0.29	1.97	0.030
34. BD (1) C 15 - O 19	/285. RY* (1) C 15	0.57	1.98	0.042
34. BD (1) C 15 - O 19	/608. BD* (1) C 13 - C 14	1.40	1.31	0.054
34. BD (1) C 15 - O 19	/609. BD* (1) C 13 - C 15	0.76	1.48	0.043
34. BD (1) C 15 - O 19	/614. BD* (1) C 15 - C 17	0.31	1.31	0.026
34. BD (1) C 15 - O 19	/618. BD* (1) C 17 - C 18	0.49	1.51	0.034
35. BD (1) C 16 - C 18	/278. RY* (2) C 14	0.97	1.83	0.053

35. BD (1) C 16 - C 18	/302. RY* (2) C 17	0.96	1.83	0.053
35. BD (1) C 16 - C 18	/611. BD* (1) C 14 - C 16	1.25	1.29	0.051
35. BD (1) C 16 - C 18	/613. BD* (1) C 14 - H 26	1.63	1.08	0.053
35. BD (1) C 16 - C 18	/617. BD* (1) C 16 - H 27	0.30	1.09	0.023
35. BD (1) C 16 - C 18	/618. BD* (1) C 17 - C 18	1.24	1.29	0.051
35. BD (1) C 16 - C 18	/620. BD* (1) C 17 - H 28	1.69	1.09	0.054
35. BD (1) C 16 - C 18	/621. BD* (1) C 18 - H 29	0.29	1.09	0.023
36. BD (1) C 16 - H 27	/277. RY* (1) C 14	0.78	1.78	0.047
36. BD (1) C 16 - H 27	/309. RY* (1) C 18	0.59	1.70	0.040
36. BD (1) C 16 - H 27	/608. BD* (1) C 13 - C 14	2.99	0.96	0.068
36. BD (1) C 16 - H 27	/611. BD* (1) C 14 - C 16	0.71	1.16	0.036
36. BD (1) C 16 - H 27	/613. BD* (1) C 14 - H 26	0.48	0.95	0.027
36. BD (1) C 16 - H 27	/618. BD* (1) C 17 - C 18	1.36	1.16	0.050
37. BD (1) C 17 - C 18	/285. RY* (1) C 15	0.49	1.87	0.038
37. BD (1) C 17 - C 18	/286. RY* (2) C 15	0.52	2.17	0.043
37. BD (1) C 17 - C 18	/293. RY* (1) C 16	0.52	1.94	0.040
37. BD (1) C 17 - C 18	/294. RY* (2) C 16	0.43	2.08	0.038
37. BD (1) C 17 - C 18	/614. BD* (1) C 15 - C 17	1.07	1.19	0.045
37. BD (1) C 17 - C 18	/615. BD* (1) C 15 - O 19	1.37	1.07	0.048
37. BD (1) C 17 - C 18	/616. BD* (1) C 16 - C 18	1.14	1.21	0.047
37. BD (1) C 17 - C 18	/617. BD* (1) C 16 - H 27	0.80	1.20	0.039
37. BD (1) C 17 - C 18	/620. BD* (1) C 17 - H 28	0.88	1.19	0.041
37. BD (1) C 17 - C 18	/621. BD* (1) C 18 - H 29	0.78	1.20	0.039
38. BD (2) C 17 - C 18	/288. RY* (4) C 15	0.27	0.94	0.021
38. BD (2) C 17 - C 18	/295. RY* (3) C 16	0.52	0.79	0.027
38. BD (2) C 17 - C 18	/610. BD* (2) C 13 - C 15	5.38	0.30	0.054
38. BD (2) C 17 - C 18	/612. BD* (2) C 14 - C 16	7.34	0.31	0.061
39. BD (1) C 17 - H 28	/285. RY* (1) C 15	0.40	1.64	0.033
39. BD (1) C 17 - H 28	/309. RY* (1) C 18	0.63	1.71	0.042
39. BD (1) C 17 - H 28	/609. BD* (1) C 13 - C 15	1.59	1.13	0.054
39. BD (1) C 17 - H 28	/615. BD* (1) C 15 - O 19	0.43	0.83	0.024
39. BD (1) C 17 - H 28	/616. BD* (1) C 16 - C 18	2.70	0.98	0.065
39. BD (1) C 17 - H 28	/618. BD* (1) C 17 - C 18	0.75	1.16	0.037
39. BD (1) C 17 - H 28	/621. BD* (1) C 18 - H 29	0.48	0.96	0.027
40. BD (1) C 18 - H 29	/293. RY* (1) C 16	0.60	1.71	0.041
40. BD (1) C 18 - H 29	/301. RY* (1) C 17	0.80	1.81	0.048
40. BD (1) C 18 - H 29	/611. BD* (1) C 14 - C 16	1.37	1.16	0.050
40. BD (1) C 18 - H 29	/614. BD* (1) C 15 - C 17	2.94	0.96	0.067
40. BD (1) C 18 - H 29	/618. BD* (1) C 17 - C 18	0.67	1.16	0.035
40. BD (1) C 18 - H 29	/620. BD* (1) C 17 - H 28	0.48	0.96	0.027
41. BD (1) O 20 - H 30	/173. RY* (1) C 1	1.35	1.79	0.062
41. BD (1) O 20 - H 30	/583. BD* (1) C 1 - C 3	1.72	1.34	0.061
85. CR (1) C 1	/181. RY* (1) C 2	0.63	11.78	0.109
85. CR (1) C 1	/190. RY* (2) C 3	1.15	11.48	0.145
85. CR (1) C 1	/583. BD* (1) C 1 - C 3	1.02	10.75	0.133
85. CR (1) C 1	/585. BD* (1) C 1 - O 20	0.66	10.40	0.105
85. CR (1) C 1	/586. BD* (1) C 2 - C 4	0.31	10.75	0.074
85. CR (1) C 1	/588. BD* (1) C 2 - C 7	0.34	10.57	0.077
85. CR (1) C 1	/589. BD* (1) C 3 - C 5	0.49	10.60	0.092
86. CR (1) C 2	/173. RY* (1) C 1	0.50	11.14	0.094
86. CR (1) C 2	/197. RY* (1) C 4	0.33	11.59	0.078
86. CR (1) C 2	/198. RY* (2) C 4	1.06	11.61	0.140
86. CR (1) C 2	/222. RY* (2) C 7	0.51	11.18	0.096
86. CR (1) C 2	/583. BD* (1) C 1 - C 3	0.30	10.69	0.072
86. CR (1) C 2	/585. BD* (1) C 1 - O 20	0.26	10.34	0.066
86. CR (1) C 2	/586. BD* (1) C 2 - C 4	0.67	10.69	0.107
86. CR (1) C 2	/591. BD* (1) C 4 - C 6	0.58	10.52	0.100
86. CR (1) C 2	/592. BD* (1) C 4 - C 9	0.53	10.52	0.095
86. CR (1) C 2	/597. BD* (1) C 7 - C 8	0.28	10.71	0.069
87. CR (1) C 3	/174. RY* (2) C 1	1.64	11.45	0.173
87. CR (1) C 3	/334. RY* (2) H 21	0.30	12.29	0.076
87. CR (1) C 3	/582. BD* (1) C 1 - C 2	0.68	10.46	0.107
87. CR (1) C 3	/583. BD* (1) C 1 - C 3	0.33	10.66	0.075
87. CR (1) C 3	/585. BD* (1) C 1 - O 20	0.40	10.31	0.082
87. CR (1) C 3	/593. BD* (1) C 5 - C 6	0.35	10.68	0.077
88. CR (1) C 4	/182. RY* (2) C 2	1.13	11.49	0.144
88. CR (1) C 4	/215. RY* (3) C 6	0.68	11.18	0.110
88. CR (1) C 4	/238. RY* (2) C 9	0.57	11.11	0.101
88. CR (1) C 4	/582. BD* (1) C 1 - C 2	0.56	10.48	0.098
88. CR (1) C 4	/586. BD* (1) C 2 - C 4	0.57	10.68	0.099
88. CR (1) C 4	/588. BD* (1) C 2 - C 7	0.52	10.50	0.094

88.	CR	(1)	C	4	/593.	BD*	(1)	C	5	-	C	6	0.35	10.70	0.078
88.	CR	(1)	C	4	/596.	BD*	(1)	C	6	-	N	11	0.28	10.44	0.069
88.	CR	(1)	C	4	/602.	BD*	(1)	C	9	-	C	10	0.29	10.71	0.070
89.	CR	(1)	C	5	/213.	RY*	(1)	C	6				0.33	11.18	0.076
89.	CR	(1)	C	5	/214.	RY*	(2)	C	6				1.21	11.02	0.146
89.	CR	(1)	C	5	/216.	RY*	(4)	C	6				0.43	11.03	0.086
89.	CR	(1)	C	5	/354.	RY*	(2)	H	31				0.26	12.32	0.071
89.	CR	(1)	C	5	/583.	BD*	(1)	C	1	-	C	3	0.33	10.66	0.075
89.	CR	(1)	C	5	/591.	BD*	(1)	C	4	-	C	6	0.68	10.50	0.108
89.	CR	(1)	C	5	/593.	BD*	(1)	C	5	-	C	6	0.48	10.68	0.091
89.	CR	(1)	C	5	/596.	BD*	(1)	C	6	-	N	11	0.49	10.42	0.091
90.	CR	(1)	C	6	/197.	RY*	(1)	C	4				0.66	11.61	0.111
90.	CR	(1)	C	6	/206.	RY*	(2)	C	5				1.18	11.30	0.146
90.	CR	(1)	C	6	/586.	BD*	(1)	C	2	-	C	4	0.37	10.71	0.080
90.	CR	(1)	C	6	/589.	BD*	(1)	C	3	-	C	5	0.49	10.55	0.091
90.	CR	(1)	C	6	/592.	BD*	(1)	C	4	-	C	9	0.34	10.54	0.076
90.	CR	(1)	C	6	/593.	BD*	(1)	C	5	-	C	6	0.75	10.73	0.114
90.	CR	(1)	C	6	/595.	BD*	(1)	C	5	-	H	31	0.25	10.52	0.066
90.	CR	(1)	C	6	/606.	BD*	(1)	N	11	-	N	12	0.39	10.58	0.082
91.	CR	(1)	C	7	/182.	RY*	(2)	C	2				0.35	11.48	0.081
91.	CR	(1)	C	7	/230.	RY*	(2)	C	8				1.15	11.34	0.144
91.	CR	(1)	C	7	/336.	RY*	(2)	H	22				0.27	12.26	0.073
91.	CR	(1)	C	7	/582.	BD*	(1)	C	1	-	C	2	0.35	10.47	0.077
91.	CR	(1)	C	7	/586.	BD*	(1)	C	2	-	C	4	0.37	10.67	0.080
91.	CR	(1)	C	7	/597.	BD*	(1)	C	7	-	C	8	0.43	10.69	0.086
91.	CR	(1)	C	7	/600.	BD*	(1)	C	8	-	C	10	0.53	10.51	0.095
92.	CR	(1)	C	8	/221.	RY*	(1)	C	7				0.37	11.24	0.082
92.	CR	(1)	C	8	/222.	RY*	(2)	C	7				0.63	11.15	0.106
92.	CR	(1)	C	8	/224.	RY*	(4)	C	7				0.40	11.08	0.085
92.	CR	(1)	C	8	/246.	RY*	(2)	C	10				0.44	11.35	0.089
92.	CR	(1)	C	8	/338.	RY*	(2)	H	23				0.31	11.93	0.076
92.	CR	(1)	C	8	/588.	BD*	(1)	C	2	-	C	7	0.62	10.48	0.102
92.	CR	(1)	C	8	/597.	BD*	(1)	C	7	-	C	8	0.38	10.68	0.081
92.	CR	(1)	C	8	/602.	BD*	(1)	C	9	-	C	10	0.32	10.69	0.074
93.	CR	(1)	C	9	/198.	RY*	(2)	C	4				0.50	11.58	0.096
93.	CR	(1)	C	9	/246.	RY*	(2)	C	10				1.09	11.35	0.141
93.	CR	(1)	C	9	/340.	RY*	(2)	H	24				0.26	12.40	0.071
93.	CR	(1)	C	9	/586.	BD*	(1)	C	2	-	C	4	0.42	10.66	0.085
93.	CR	(1)	C	9	/591.	BD*	(1)	C	4	-	C	6	0.34	10.49	0.077
93.	CR	(1)	C	9	/600.	BD*	(1)	C	8	-	C	10	0.54	10.50	0.095
93.	CR	(1)	C	9	/602.	BD*	(1)	C	9	-	C	10	0.43	10.69	0.086
94.	CR	(1)	C	10	/230.	RY*	(2)	C	8				0.42	11.33	0.087
94.	CR	(1)	C	10	/237.	RY*	(1)	C	9				0.39	11.24	0.084
94.	CR	(1)	C	10	/238.	RY*	(2)	C	9				0.51	11.09	0.095
94.	CR	(1)	C	10	/240.	RY*	(4)	C	9				0.57	11.20	0.101
94.	CR	(1)	C	10	/342.	RY*	(2)	H	25				0.34	12.31	0.082
94.	CR	(1)	C	10	/592.	BD*	(1)	C	4	-	C	9	0.61	10.49	0.101
94.	CR	(1)	C	10	/597.	BD*	(1)	C	7	-	C	8	0.32	10.68	0.074
94.	CR	(1)	C	10	/602.	BD*	(1)	C	9	-	C	10	0.41	10.69	0.084
95.	CR	(1)	N	11	/213.	RY*	(1)	C	6				0.89	15.34	0.148
95.	CR	(1)	N	11	/263.	RY*	(3)	N	12				0.85	15.27	0.144
95.	CR	(1)	N	11	/591.	BD*	(1)	C	4	-	C	6	0.28	14.66	0.082
95.	CR	(1)	N	11	/593.	BD*	(1)	C	5	-	C	6	0.28	14.84	0.082
95.	CR	(1)	N	11	/607.	BD*	(1)	N	12	-	C	13	0.68	14.58	0.127
96.	CR	(1)	N	12	/255.	RY*	(3)	N	11				0.99	15.44	0.156
96.	CR	(1)	N	12	/269.	RY*	(1)	C	13				0.78	15.32	0.138
96.	CR	(1)	N	12	/270.	RY*	(2)	C	13				0.30	15.64	0.086
96.	CR	(1)	N	12	/596.	BD*	(1)	C	6	-	N	11	0.63	14.59	0.122
96.	CR	(1)	N	12	/609.	BD*	(1)	C	13	-	C	15	0.36	14.83	0.093
97.	CR	(1)	C	13	/278.	RY*	(2)	C	14				0.51	11.28	0.096
97.	CR	(1)	C	13	/285.	RY*	(1)	C	15				0.51	11.21	0.096
97.	CR	(1)	C	13	/286.	RY*	(2)	C	15				1.17	11.51	0.146
97.	CR	(1)	C	13	/606.	BD*	(1)	N	11	-	N	12	0.46	10.59	0.088
97.	CR	(1)	C	13	/609.	BD*	(1)	C	13	-	C	15	0.74	10.71	0.113
97.	CR	(1)	C	13	/611.	BD*	(1)	C	14	-	C	16	0.28	10.73	0.070
97.	CR	(1)	C	13	/614.	BD*	(1)	C	15	-	C	17	0.56	10.53	0.098
97.	CR	(1)	C	13	/615.	BD*	(1)	C	15	-	O	19	0.40	10.41	0.082
98.	CR	(1)	C	14	/269.	RY*	(1)	C	13				0.29	11.16	0.072
98.	CR	(1)	C	14	/271.	RY*	(3)	C	13				0.40	10.97	0.084
98.	CR	(1)	C	14	/294.	RY*	(2)	C	16				1.07	11.37	0.139
98.	CR	(1)	C	14	/344.	RY*	(2)	H	26				0.31	12.30	0.078

98. CR (1) C 14	/607. BD* (1) N 12 - C 13	0.29	10.42	0.070
98. CR (1) C 14	/609. BD* (1) C 13 - C 15	0.41	10.66	0.084
98. CR (1) C 14	/611. BD* (1) C 14 - C 16	0.44	10.69	0.087
98. CR (1) C 14	/616. BD* (1) C 16 - C 18	0.52	10.51	0.094
99. CR (1) C 15	/270. RY* (2) C 13	1.56	11.54	0.169
99. CR (1) C 15	/302. RY* (2) C 17	0.53	11.30	0.097
99. CR (1) C 15	/607. BD* (1) N 12 - C 13	0.43	10.49	0.086
99. CR (1) C 15	/608. BD* (1) C 13 - C 14	0.56	10.56	0.097
99. CR (1) C 15	/609. BD* (1) C 13 - C 15	0.90	10.73	0.125
99. CR (1) C 15	/615. BD* (1) C 15 - O 19	0.42	10.43	0.085
99. CR (1) C 15	/618. BD* (1) C 17 - C 18	0.30	10.76	0.072
100. CR (1) C 16	/277. RY* (1) C 14	0.43	11.31	0.088
100. CR (1) C 16	/278. RY* (2) C 14	0.75	11.23	0.116
100. CR (1) C 16	/310. RY* (2) C 18	0.45	11.35	0.091
100. CR (1) C 16	/346. RY* (2) H 27	0.33	12.02	0.080
100. CR (1) C 16	/608. BD* (1) C 13 - C 14	0.56	10.48	0.098
100. CR (1) C 16	/611. BD* (1) C 14 - C 16	0.37	10.68	0.080
100. CR (1) C 16	/618. BD* (1) C 17 - C 18	0.30	10.68	0.072
101. CR (1) C 17	/286. RY* (2) C 15	0.26	11.46	0.069
101. CR (1) C 17	/287. RY* (3) C 15	0.65	11.15	0.107
101. CR (1) C 17	/310. RY* (2) C 18	1.06	11.35	0.139
101. CR (1) C 17	/348. RY* (2) H 28	0.30	12.34	0.077
101. CR (1) C 17	/609. BD* (1) C 13 - C 15	0.45	10.66	0.088
101. CR (1) C 17	/616. BD* (1) C 16 - C 18	0.52	10.50	0.094
101. CR (1) C 17	/618. BD* (1) C 17 - C 18	0.42	10.69	0.085
102. CR (1) C 18	/294. RY* (2) C 16	0.46	11.37	0.092
102. CR (1) C 18	/301. RY* (1) C 17	0.44	11.34	0.090
102. CR (1) C 18	/302. RY* (2) C 17	0.82	11.22	0.121
102. CR (1) C 18	/350. RY* (2) H 29	0.31	11.98	0.078
102. CR (1) C 18	/611. BD* (1) C 14 - C 16	0.30	10.68	0.072
102. CR (1) C 18	/614. BD* (1) C 15 - C 17	0.53	10.48	0.095
102. CR (1) C 18	/618. BD* (1) C 17 - C 18	0.38	10.69	0.081
103. CR (1) O 19	/285. RY* (1) C 15	1.45	20.03	0.216
103. CR (1) O 19	/609. BD* (1) C 13 - C 15	0.30	19.53	0.098
104. CR (1) O 20	/173. RY* (1) C 1	1.16	20.03	0.193
104. CR (1) O 20	/582. BD* (1) C 1 - C 2	0.28	19.38	0.094
140. LP (1) N 11	/214. RY* (2) C 6	0.27	1.28	0.024
140. LP (1) N 11	/261. RY* (1) N 12	0.63	1.13	0.034
140. LP (1) N 11	/262. RY* (2) N 12	1.02	1.16	0.044
140. LP (1) N 11	/589. BD* (1) C 3 - C 5	0.25	0.77	0.018
140. LP (1) N 11	/591. BD* (1) C 4 - C 6	4.40	0.76	0.074
140. LP (1) N 11	/593. BD* (1) C 5 - C 6	2.54	0.94	0.063
140. LP (1) N 11	/594. BD* (2) C 5 - C 6	0.62	0.34	0.019
141. LP (2) N 11	/213. RY* (1) C 6	0.82	1.62	0.047
141. LP (2) N 11	/253. RY* (1) N 11	0.88	1.58	0.048
141. LP (2) N 11	/254. RY* (2) N 11	1.67	2.22	0.079
141. LP (2) N 11	/255. RY* (3) N 11	0.68	1.71	0.044
141. LP (2) N 11	/261. RY* (1) N 12	0.37	1.31	0.028
141. LP (2) N 11	/317. RY* (1) O 19	0.45	2.65	0.045
141. LP (2) N 11	/591. BD* (1) C 4 - C 6	0.55	0.94	0.029
141. LP (2) N 11	/594. BD* (2) C 5 - C 6	3.49	0.53	0.056
142. LP (1) N 12	/254. RY* (2) N 11	0.35	2.15	0.035
142. LP (1) N 12	/255. RY* (3) N 11	0.45	1.63	0.035
142. LP (1) N 12	/269. RY* (1) C 13	0.96	1.52	0.049
142. LP (1) N 12	/608. BD* (1) C 13 - C 14	0.57	0.85	0.028
142. LP (1) N 12	/609. BD* (1) C 13 - C 15	3.64	1.02	0.077
143. LP (2) N 12	/253. RY* (1) N 11	1.08	1.38	0.051
143. LP (2) N 12	/272. RY* (4) C 13	0.35	0.94	0.024
143. LP (2) N 12	/610. BD* (2) C 13 - C 15	16.65	0.30	0.095
144. LP (1) O 19	/285. RY* (1) C 15	1.04	1.67	0.053
144. LP (1) O 19	/609. BD* (1) C 13 - C 15	3.39	1.17	0.080
145. LP (2) O 19	/254. RY* (2) N 11	0.37	2.24	0.037
145. LP (2) O 19	/285. RY* (1) C 15	0.69	1.62	0.043
145. LP (2) O 19	/317. RY* (1) O 19	0.54	2.67	0.049
145. LP (2) O 19	/609. BD* (1) C 13 - C 15	0.35	1.12	0.025
145. LP (2) O 19	/610. BD* (2) C 13 - C 15	0.45	0.53	0.021
145. LP (2) O 19	/614. BD* (1) C 15 - C 17	3.58	0.95	0.074
146. LP (3) O 19	/288. RY* (4) C 15	0.34	1.01	0.024
146. LP (3) O 19	/317. RY* (1) O 19	0.25	2.52	0.033
146. LP (3) O 19	/610. BD* (2) C 13 - C 15	10.31	0.37	0.085
146. LP (3) O 19	/614. BD* (1) C 15 - C 17	0.33	0.79	0.021

147. LP (1) O 20	/173. RY* (1) C 1	1.45	1.66	0.062
147. LP (1) O 20	/582. BD* (1) C 1 - C 2	3.52	1.01	0.076
148. LP (2) O 20	/584. BD* (2) C 1 - C 3	12.04	0.35	0.086
584. BD* (2) C 1 - C 3	/176. RY* (4) C 1	0.51	0.54	0.058
584. BD* (2) C 1 - C 3	/191. RY* (3) C 3	1.15	0.48	0.081
587. BD* (2) C 2 - C 4	/184. RY* (4) C 2	1.11	0.57	0.075
587. BD* (2) C 2 - C 4	/200. RY* (4) C 4	0.87	0.63	0.070
594. BD* (2) C 5 - C 6	/207. RY* (3) C 5	0.50	0.68	0.066
594. BD* (2) C 5 - C 6	/606. BD* (1) N 11 - N 12	0.51	0.45	0.053
598. BD* (2) C 7 - C 8	/223. RY* (3) C 7	0.69	0.46	0.073
598. BD* (2) C 7 - C 8	/231. RY* (3) C 8	0.74	0.48	0.077
603. BD* (2) C 9 - C 10	/239. RY* (3) C 9	0.62	0.55	0.077
603. BD* (2) C 9 - C 10	/247. RY* (3) C 10	0.65	0.51	0.076
610. BD* (2) C 13 - C 15	/272. RY* (4) C 13	1.03	0.64	0.070
610. BD* (2) C 13 - C 15	/288. RY* (4) C 15	0.77	0.64	0.061
610. BD* (2) C 13 - C 15	/612. BD* (2) C 14 - C 16	62.76	0.02	0.071
610. BD* (2) C 13 - C 15	/619. BD* (2) C 17 - C 18	82.91	0.01	0.071
612. BD* (2) C 14 - C 16	/279. RY* (3) C 14	0.75	0.50	0.072
612. BD* (2) C 14 - C 16	/295. RY* (3) C 16	0.90	0.48	0.077
619. BD* (2) C 17 - C 18	/303. RY* (3) C 17	0.96	0.54	0.078
619. BD* (2) C 17 - C 18	/311. RY* (3) C 18	0.95	0.48	0.073

from unit 1 to unit 2

5. BD (1) C 2 - C 4	/521. RY* (1) H 58	0.04	1.18	0.009
6. BD (2) C 2 - C 4	/661. BD* (1) C 48 - H 58	0.03	0.71	0.006
10. BD (1) C 4 - C 6	/467. RY* (1) C 46	0.05	1.80	0.011
10. BD (1) C 4 - C 6	/542. RY* (2) O 65	0.04	1.79	0.011
11. BD (1) C 4 - C 9	/483. RY* (1) C 48	0.06	1.92	0.013
11. BD (1) C 4 - C 9	/485. RY* (3) C 48	0.04	1.41	0.009
12. BD (1) C 5 - C 6	/542. RY* (2) O 65	0.03	1.90	0.009
13. BD (2) C 5 - C 6	/542. RY* (2) O 65	0.03	1.42	0.009
13. BD (2) C 5 - C 6	/656. BD* (1) C 46 - O 65	0.04	0.57	0.007
15. BD (1) C 6 - N 11	/467. RY* (1) C 46	0.04	1.94	0.012
15. BD (1) C 6 - N 11	/542. RY* (2) O 65	0.04	1.93	0.012
21. BD (1) C 9 - C 10	/483. RY* (1) C 48	0.05	2.02	0.013
21. BD (1) C 9 - C 10	/485. RY* (3) C 48	0.09	1.52	0.015
22. BD (2) C 9 - C 10	/660. BD* (2) C 48 - C 49	0.17	0.30	0.009
33. BD (1) C 15 - C 17	/527. RY* (1) H 61	0.12	1.34	0.016
34. BD (1) C 15 - O 19	/435. RY* (1) N 42	0.16	2.11	0.023
34. BD (1) C 15 - O 19	/436. RY* (2) N 42	0.05	2.42	0.014
34. BD (1) C 15 - O 19	/451. RY* (1) C 44	0.07	1.99	0.015
38. BD (2) C 17 - C 18	/636. BD* (1) C 36 - H 61	0.05	0.72	0.008
39. BD (1) C 17 - H 28	/636. BD* (1) C 36 - H 61	0.03	0.96	0.007
141. LP (2) N 11	/395. RY* (1) C 37	0.05	1.61	0.011
141. LP (2) N 11	/396. RY* (2) C 37	0.10	1.58	0.016
141. LP (2) N 11	/398. RY* (4) C 37	0.10	1.43	0.016
141. LP (2) N 11	/399. RY* (5) C 37	0.03	3.44	0.014
141. LP (2) N 11	/400. RY* (6) C 37	0.04	3.60	0.015
141. LP (2) N 11	/436. RY* (2) N 42	0.17	2.04	0.024
141. LP (2) N 11	/437. RY* (3) N 42	0.27	1.76	0.028
141. LP (2) N 11	/443. RY* (1) N 43	0.03	1.32	0.008
141. LP (2) N 11	/445. RY* (3) N 43	0.06	1.56	0.012
141. LP (2) N 11	/453. RY* (3) C 44	0.03	1.38	0.008
141. LP (2) N 11	/467. RY* (1) C 46	0.23	1.64	0.025
141. LP (2) N 11	/468. RY* (2) C 46	0.05	1.88	0.013
141. LP (2) N 11	/469. RY* (3) C 46	0.09	1.20	0.013
141. LP (2) N 11	/471. RY* (5) C 46	0.06	3.47	0.018
141. LP (2) N 11	/541. RY* (1) O 65	0.50	2.56	0.046
141. LP (2) N 11	/543. RY* (3) O 65	0.03	1.96	0.009
141. LP (2) N 11	/544. RY* (4) O 65	0.10	2.01	0.018
141. LP (2) N 11	/656. BD* (1) C 46 - O 65	0.14	0.78	0.013
144. LP (1) O 19	/435. RY* (1) N 42	0.15	1.80	0.021
144. LP (1) O 19	/436. RY* (2) N 42	0.19	2.11	0.025
144. LP (1) O 19	/437. RY* (3) N 42	0.17	1.83	0.022
144. LP (1) O 19	/444. RY* (2) N 43	0.03	1.41	0.008
144. LP (1) O 19	/446. RY* (4) N 43	0.03	2.19	0.010
144. LP (1) O 19	/467. RY* (1) C 46	0.04	1.71	0.011
144. LP (1) O 19	/471. RY* (5) C 46	0.03	3.54	0.014
144. LP (1) O 19	/541. RY* (1) O 65	0.06	2.63	0.016
144. LP (1) O 19	/637. BD* (1) C 37 - N 42	0.16	0.92	0.015
144. LP (1) O 19	/648. BD* (1) N 43 - C 44	0.06	0.91	0.010

144. LP (1) O 19	/651. BD* (2) C 44 - C 46	0.08	0.55	0.009
145. LP (2) O 19	/395. RY* (1) C 37	0.03	1.63	0.009
145. LP (2) O 19	/396. RY* (2) C 37	0.05	1.60	0.011
145. LP (2) O 19	/398. RY* (4) C 37	0.04	1.45	0.010
145. LP (2) O 19	/435. RY* (1) N 42	0.03	1.75	0.009
145. LP (2) O 19	/436. RY* (2) N 42	0.17	2.06	0.024
145. LP (2) O 19	/437. RY* (3) N 42	0.22	1.78	0.025
145. LP (2) O 19	/467. RY* (1) C 46	0.06	1.66	0.013
145. LP (2) O 19	/469. RY* (3) C 46	0.03	1.22	0.007
145. LP (2) O 19	/471. RY* (5) C 46	0.03	3.49	0.014
145. LP (2) O 19	/541. RY* (1) O 65	0.05	2.58	0.015
145. LP (2) O 19	/544. RY* (4) O 65	0.05	2.03	0.013
145. LP (2) O 19	/637. BD* (1) C 37 - N 42	0.17	0.87	0.015
145. LP (2) O 19	/648. BD* (1) N 43 - C 44	0.03	0.86	0.007
145. LP (2) O 19	/656. BD* (1) C 46 - O 65	0.03	0.80	0.006
146. LP (3) O 19	/395. RY* (1) C 37	0.03	1.47	0.009
146. LP (3) O 19	/397. RY* (3) C 37	0.03	1.41	0.008
146. LP (3) O 19	/435. RY* (1) N 42	0.49	1.60	0.036
146. LP (3) O 19	/436. RY* (2) N 42	0.31	1.91	0.032
146. LP (3) O 19	/437. RY* (3) N 42	0.14	1.63	0.020
146. LP (3) O 19	/438. RY* (4) N 42	0.03	1.82	0.010
146. LP (3) O 19	/444. RY* (2) N 43	0.03	1.21	0.007
146. LP (3) O 19	/445. RY* (3) N 43	0.04	1.42	0.009
146. LP (3) O 19	/453. RY* (3) C 44	0.03	1.25	0.007
146. LP (3) O 19	/467. RY* (1) C 46	0.05	1.51	0.012
146. LP (3) O 19	/468. RY* (2) C 46	0.05	1.75	0.012
146. LP (3) O 19	/469. RY* (3) C 46	0.06	1.06	0.011
146. LP (3) O 19	/541. RY* (1) O 65	0.03	2.43	0.012
146. LP (3) O 19	/544. RY* (4) O 65	0.03	1.88	0.011
146. LP (3) O 19	/637. BD* (1) C 37 - N 42	0.15	0.72	0.013
146. LP (3) O 19	/647. BD* (1) N 42 - N 43	0.04	0.84	0.007
619. BD* (2) C 17 - C 18	/636. BD* (1) C 36 - H 61	0.06	0.41	0.017

from unit 1 to unit 3

25. BD (1) N 11 - N 12	/529. RY* (1) O 62	0.03	2.74	0.011
26. BD (1) N 12 - C 13	/529. RY* (1) O 62	0.03	2.55	0.012
26. BD (1) N 12 - C 13	/537. RY* (1) H 63	0.03	1.31	0.007
28. BD (1) C 13 - C 15	/539. RY* (1) H 64	0.03	1.39	0.008
29. BD (2) C 13 - C 15	/665. BD* (1) O 62 - H 64	0.18	0.61	0.014
33. BD (1) C 15 - C 17	/529. RY* (1) O 62	0.03	2.41	0.010
33. BD (1) C 15 - C 17	/539. RY* (1) H 64	0.03	1.27	0.008
140. LP (1) N 11	/529. RY* (1) O 62	0.03	2.06	0.010
140. LP (1) N 11	/664. BD* (1) O 62 - H 63	0.12	0.61	0.011
141. LP (2) N 11	/529. RY* (1) O 62	0.91	2.24	0.059
141. LP (2) N 11	/530. RY* (2) O 62	0.21	1.85	0.025
141. LP (2) N 11	/531. RY* (3) O 62	0.10	1.55	0.016
141. LP (2) N 11	/532. RY* (4) O 62	0.14	1.70	0.020
141. LP (2) N 11	/537. RY* (1) H 63	0.28	1.01	0.022
141. LP (2) N 11	/539. RY* (1) H 64	0.28	1.10	0.023
141. LP (2) N 11	/664. BD* (1) O 62 - H 63	0.05	0.79	0.008
144. LP (1) O 19	/529. RY* (1) O 62	0.16	2.31	0.024
144. LP (1) O 19	/530. RY* (2) O 62	0.03	1.92	0.009
144. LP (1) O 19	/537. RY* (1) H 63	0.04	1.08	0.009
144. LP (1) O 19	/539. RY* (1) H 64	0.05	1.17	0.010
145. LP (2) O 19	/529. RY* (1) O 62	0.41	2.26	0.039
145. LP (2) O 19	/530. RY* (2) O 62	0.09	1.87	0.016
145. LP (2) O 19	/531. RY* (3) O 62	0.03	1.57	0.008
145. LP (2) O 19	/532. RY* (4) O 62	0.06	1.72	0.013
145. LP (2) O 19	/537. RY* (1) H 63	0.09	1.03	0.012
145. LP (2) O 19	/539. RY* (1) H 64	0.16	1.12	0.017
145. LP (2) O 19	/664. BD* (1) O 62 - H 63	0.03	0.81	0.006
145. LP (2) O 19	/665. BD* (1) O 62 - H 64	0.06	0.81	0.009
146. LP (3) O 19	/529. RY* (1) O 62	0.14	2.11	0.023
146. LP (3) O 19	/530. RY* (2) O 62	0.06	1.72	0.013
146. LP (3) O 19	/532. RY* (4) O 62	0.03	1.57	0.009
146. LP (3) O 19	/539. RY* (1) H 64	0.03	0.97	0.007
146. LP (3) O 19	/665. BD* (1) O 62 - H 64	0.04	0.66	0.006
610. BD* (2) C 13 - C 15	/665. BD* (1) O 62 - H 64	0.04	0.29	0.009

from unit 1 to unit 4

26. BD (1) N 12 - C 13	/551. RY* (3) C1 66	0.03	2.17	0.010
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141.	LP	(2)	N	11	/549.	RY*	(1)	Cl	66	0.16	1.06	0.017	
141.	LP	(2)	N	11	/550.	RY*	(2)	Cl	66	0.21	1.11	0.020	
141.	LP	(2)	N	11	/551.	RY*	(3)	Cl	66	0.51	1.86	0.040	
144.	LP	(1)	O	19	/550.	RY*	(2)	Cl	66	0.04	1.18	0.008	
144.	LP	(1)	O	19	/551.	RY*	(3)	Cl	66	0.11	1.93	0.018	
145.	LP	(2)	O	19	/549.	RY*	(1)	Cl	66	0.04	1.08	0.008	
145.	LP	(2)	O	19	/550.	RY*	(2)	Cl	66	0.13	1.13	0.016	
145.	LP	(2)	O	19	/551.	RY*	(3)	Cl	66	0.32	1.88	0.032	
146.	LP	(3)	O	19	/549.	RY*	(1)	Cl	66	0.05	0.93	0.009	
146.	LP	(3)	O	19	/550.	RY*	(2)	Cl	66	0.05	0.98	0.009	
146.	LP	(3)	O	19	/551.	RY*	(3)	Cl	66	0.11	1.73	0.018	
from unit 1 to unit 5															
10.	BD	(1)	C	4 - C	6	/169.	LP*	(6)	Fe	67	0.10	1.17	0.015
10.	BD	(1)	C	4 - C	6	/561.	RY*	(1)	Fe	67	0.03	1.28	0.008
12.	BD	(1)	C	5 - C	6	/172.	LP*	(9)	Fe	67	0.03	2.44	0.011
12.	BD	(1)	C	5 - C	6	/561.	RY*	(1)	Fe	67	0.06	1.39	0.012
12.	BD	(1)	C	5 - C	6	/571.	RY*	(11)	Fe	67	0.03	10.29	0.022
15.	BD	(1)	C	6 - N	11	/169.	LP*	(6)	Fe	67	1.14	1.31	0.053
15.	BD	(1)	C	6 - N	11	/170.	LP*	(7)	Fe	67	0.07	2.28	0.015
15.	BD	(1)	C	6 - N	11	/561.	RY*	(1)	Fe	67	0.03	1.42	0.009
15.	BD	(1)	C	6 - N	11	/562.	RY*	(2)	Fe	67	0.03	2.39	0.011
15.	BD	(1)	C	6 - N	11	/563.	RY*	(3)	Fe	67	0.09	4.94	0.026
15.	BD	(1)	C	6 - N	11	/564.	RY*	(4)	Fe	67	0.04	2.17	0.012
15.	BD	(1)	C	6 - N	11	/567.	RY*	(7)	Fe	67	0.07	2.85	0.017
15.	BD	(1)	C	6 - N	11	/571.	RY*	(11)	Fe	67	0.03	10.33	0.021
15.	BD	(1)	C	6 - N	11	/575.	RY*	(15)	Fe	67	0.04	166.16	0.100
15.	BD	(1)	C	6 - N	11	/580.	RY*	(20)	Fe	67	0.08	48.80	0.078
25.	BD	(1)	N	11 - N	12	/169.	LP*	(6)	Fe	67	1.10	1.51	0.056
25.	BD	(1)	N	11 - N	12	/563.	RY*	(3)	Fe	67	0.31	5.13	0.050
25.	BD	(1)	N	11 - N	12	/569.	RY*	(9)	Fe	67	0.04	10.16	0.024
25.	BD	(1)	N	11 - N	12	/570.	RY*	(10)	Fe	67	0.03	6.85	0.018
25.	BD	(1)	N	11 - N	12	/575.	RY*	(15)	Fe	67	0.06	166.36	0.125
25.	BD	(1)	N	11 - N	12	/579.	RY*	(19)	Fe	67	0.03	730.90	0.195
25.	BD	(1)	N	11 - N	12	/580.	RY*	(20)	Fe	67	0.11	49.00	0.094
26.	BD	(1)	N	12 - C	13	/169.	LP*	(6)	Fe	67	0.14	1.31	0.018
26.	BD	(1)	N	12 - C	13	/172.	LP*	(9)	Fe	67	0.03	2.47	0.010
26.	BD	(1)	N	12 - C	13	/561.	RY*	(1)	Fe	67	0.06	1.42	0.012
26.	BD	(1)	N	12 - C	13	/563.	RY*	(3)	Fe	67	0.12	4.94	0.031
26.	BD	(1)	N	12 - C	13	/567.	RY*	(7)	Fe	67	0.04	2.85	0.013
26.	BD	(1)	N	12 - C	13	/571.	RY*	(11)	Fe	67	0.08	10.33	0.037
26.	BD	(1)	N	12 - C	13	/575.	RY*	(15)	Fe	67	0.07	166.16	0.135
26.	BD	(1)	N	12 - C	13	/576.	RY*	(16)	Fe	67	0.03	30.87	0.035
26.	BD	(1)	N	12 - C	13	/579.	RY*	(19)	Fe	67	0.03	730.71	0.198
26.	BD	(1)	N	12 - C	13	/580.	RY*	(20)	Fe	67	0.14	48.80	0.104
27.	BD	(1)	C	13 - C	14	/169.	LP*	(6)	Fe	67	0.04	1.17	0.009
28.	BD	(1)	C	13 - C	15	/169.	LP*	(6)	Fe	67	0.18	1.30	0.021
28.	BD	(1)	C	13 - C	15	/561.	RY*	(1)	Fe	67	0.04	1.41	0.009
28.	BD	(1)	C	13 - C	15	/563.	RY*	(3)	Fe	67	0.05	4.92	0.019
28.	BD	(1)	C	13 - C	15	/580.	RY*	(20)	Fe	67	0.03	48.79	0.052
29.	BD	(2)	C	13 - C	15	/169.	LP*	(6)	Fe	67	0.20	0.82	0.016
29.	BD	(2)	C	13 - C	15	/563.	RY*	(3)	Fe	67	0.03	4.45	0.016
33.	BD	(1)	C	15 - C	17	/169.	LP*	(6)	Fe	67	0.21	1.18	0.022
33.	BD	(1)	C	15 - C	17	/170.	LP*	(7)	Fe	67	0.04	2.15	0.011
33.	BD	(1)	C	15 - C	17	/561.	RY*	(1)	Fe	67	0.03	1.29	0.008
33.	BD	(1)	C	15 - C	17	/563.	RY*	(3)	Fe	67	0.05	4.81	0.019
33.	BD	(1)	C	15 - C	17	/580.	RY*	(20)	Fe	67	0.04	48.67	0.054
34.	BD	(1)	C	15 - O	19	/169.	LP*	(6)	Fe	67	0.90	1.39	0.048
34.	BD	(1)	C	15 - O	19	/170.	LP*	(7)	Fe	67	0.17	2.36	0.025
34.	BD	(1)	C	15 - O	19	/171.	LP*	(8)	Fe	67	0.03	2.61	0.011
34.	BD	(1)	C	15 - O	19	/563.	RY*	(3)	Fe	67	0.03	5.01	0.016
34.	BD	(1)	C	15 - O	19	/568.	RY*	(8)	Fe	67	0.07	9.68	0.032
34.	BD	(1)	C	15 - O	19	/569.	RY*	(9)	Fe	67	0.06	10.04	0.032
34.	BD	(1)	C	15 - O	19	/581.	RY*	(21)	Fe	67	0.03	30.63	0.036
90.	CR	(1)	C	6	/169.	LP*	(6)	Fe	67	0.03	10.61	0.024	
95.	CR	(1)	N	11	/169.	LP*	(6)	Fe	67	1.18	14.73	0.181	
96.	CR	(1)	N	12	/169.	LP*	(6)	Fe	67	0.04	14.73	0.033	
99.	CR	(1)	C	15	/169.	LP*	(6)	Fe	67	0.06	10.64	0.034	
103.	CR	(1)	O	19	/169.	LP*	(6)	Fe	67	1.38	19.44	0.224	
140.	LP	(1)	N	11	/169.	LP*	(6)	Fe	67	1.34	0.83	0.044	
140.	LP	(1)	N	11	/171.	LP*	(8)	Fe	67	0.03	2.05	0.011	

140.	LP	(1)	N	11	/172.	LP*	(9)	Fe	67	0.10	1.98	0.018
140.	LP	(1)	N	11	/563.	RY*	(3)	Fe	67	0.15	4.45	0.033
140.	LP	(1)	N	11	/567.	RY*	(7)	Fe	67	0.07	2.36	0.017
140.	LP	(1)	N	11	/573.	RY*	(13)	Fe	67	0.03	6.98	0.020
140.	LP	(1)	N	11	/575.	RY*	(15)	Fe	67	0.05	165.67	0.118
140.	LP	(1)	N	11	/580.	RY*	(20)	Fe	67	0.10	48.32	0.089
141.	LP	(2)	N	11	/169.	LP*	(6)	Fe	67	24.59	1.01	0.209
141.	LP	(2)	N	11	/170.	LP*	(7)	Fe	67	0.45	1.98	0.039
141.	LP	(2)	N	11	/172.	LP*	(9)	Fe	67	0.99	2.16	0.060
141.	LP	(2)	N	11	/561.	RY*	(1)	Fe	67	0.50	1.12	0.031
141.	LP	(2)	N	11	/562.	RY*	(2)	Fe	67	0.66	2.09	0.048
141.	LP	(2)	N	11	/563.	RY*	(3)	Fe	67	3.78	4.63	0.172
141.	LP	(2)	N	11	/564.	RY*	(4)	Fe	67	0.21	1.87	0.026
141.	LP	(2)	N	11	/567.	RY*	(7)	Fe	67	0.95	2.54	0.064
141.	LP	(2)	N	11	/570.	RY*	(10)	Fe	67	0.22	6.35	0.049
141.	LP	(2)	N	11	/571.	RY*	(11)	Fe	67	0.05	10.02	0.030
141.	LP	(2)	N	11	/572.	RY*	(12)	Fe	67	0.17	7.20	0.045
141.	LP	(2)	N	11	/573.	RY*	(13)	Fe	67	0.23	7.16	0.053
141.	LP	(2)	N	11	/574.	RY*	(14)	Fe	67	0.22	7.00	0.051
141.	LP	(2)	N	11	/575.	RY*	(15)	Fe	67	1.45	165.86	0.636
141.	LP	(2)	N	11	/576.	RY*	(16)	Fe	67	0.03	30.56	0.041
141.	LP	(2)	N	11	/579.	RY*	(19)	Fe	67	0.67	730.40	0.907
141.	LP	(2)	N	11	/580.	RY*	(20)	Fe	67	2.89	48.50	0.486
142.	LP	(1)	N	12	/169.	LP*	(6)	Fe	67	0.27	0.93	0.021
142.	LP	(1)	N	12	/172.	LP*	(9)	Fe	67	0.10	2.08	0.018
142.	LP	(1)	N	12	/563.	RY*	(3)	Fe	67	0.03	4.55	0.015
144.	LP	(1)	O	19	/169.	LP*	(6)	Fe	67	4.27	1.08	0.092
144.	LP	(1)	O	19	/170.	LP*	(7)	Fe	67	0.10	2.05	0.018
144.	LP	(1)	O	19	/171.	LP*	(8)	Fe	67	0.04	2.30	0.012
144.	LP	(1)	O	19	/172.	LP*	(9)	Fe	67	0.14	2.23	0.022
144.	LP	(1)	O	19	/561.	RY*	(1)	Fe	67	0.04	1.19	0.008
144.	LP	(1)	O	19	/562.	RY*	(2)	Fe	67	0.32	2.16	0.033
144.	LP	(1)	O	19	/563.	RY*	(3)	Fe	67	0.64	4.70	0.070
144.	LP	(1)	O	19	/564.	RY*	(4)	Fe	67	0.05	1.94	0.012
144.	LP	(1)	O	19	/566.	RY*	(6)	Fe	67	0.07	2.32	0.017
144.	LP	(1)	O	19	/567.	RY*	(7)	Fe	67	0.16	2.61	0.026
144.	LP	(1)	O	19	/574.	RY*	(14)	Fe	67	0.05	7.07	0.024
144.	LP	(1)	O	19	/575.	RY*	(15)	Fe	67	0.24	165.93	0.252
144.	LP	(1)	O	19	/579.	RY*	(19)	Fe	67	0.13	730.47	0.390
144.	LP	(1)	O	19	/580.	RY*	(20)	Fe	67	0.48	48.57	0.193
145.	LP	(2)	O	19	/169.	LP*	(6)	Fe	67	13.85	1.03	0.160
145.	LP	(2)	O	19	/170.	LP*	(7)	Fe	67	1.29	2.00	0.065
145.	LP	(2)	O	19	/171.	LP*	(8)	Fe	67	0.03	2.25	0.011
145.	LP	(2)	O	19	/172.	LP*	(9)	Fe	67	0.06	2.18	0.015
145.	LP	(2)	O	19	/561.	RY*	(1)	Fe	67	0.63	1.14	0.034
145.	LP	(2)	O	19	/562.	RY*	(2)	Fe	67	1.05	2.11	0.060
145.	LP	(2)	O	19	/563.	RY*	(3)	Fe	67	1.22	4.65	0.097
145.	LP	(2)	O	19	/565.	RY*	(5)	Fe	67	0.11	1.97	0.019
145.	LP	(2)	O	19	/566.	RY*	(6)	Fe	67	0.09	2.27	0.019
145.	LP	(2)	O	19	/567.	RY*	(7)	Fe	67	0.44	2.56	0.043
145.	LP	(2)	O	19	/568.	RY*	(8)	Fe	67	0.05	9.32	0.029
145.	LP	(2)	O	19	/570.	RY*	(10)	Fe	67	0.08	6.37	0.029
145.	LP	(2)	O	19	/573.	RY*	(13)	Fe	67	0.03	7.18	0.018
145.	LP	(2)	O	19	/574.	RY*	(14)	Fe	67	0.10	7.02	0.034
145.	LP	(2)	O	19	/575.	RY*	(15)	Fe	67	0.67	165.88	0.428
145.	LP	(2)	O	19	/579.	RY*	(19)	Fe	67	0.32	730.42	0.620
145.	LP	(2)	O	19	/580.	RY*	(20)	Fe	67	1.36	48.52	0.330
146.	LP	(3)	O	19	/169.	LP*	(6)	Fe	67	8.48	0.87	0.114
146.	LP	(3)	O	19	/170.	LP*	(7)	Fe	67	0.92	1.84	0.053
146.	LP	(3)	O	19	/172.	LP*	(9)	Fe	67	0.35	2.03	0.035
146.	LP	(3)	O	19	/561.	RY*	(1)	Fe	67	0.39	0.98	0.026
146.	LP	(3)	O	19	/562.	RY*	(2)	Fe	67	0.38	1.96	0.035
146.	LP	(3)	O	19	/563.	RY*	(3)	Fe	67	0.60	4.50	0.068
146.	LP	(3)	O	19	/567.	RY*	(7)	Fe	67	0.23	2.41	0.031
146.	LP	(3)	O	19	/568.	RY*	(8)	Fe	67	0.03	9.16	0.023
146.	LP	(3)	O	19	/574.	RY*	(14)	Fe	67	0.07	6.87	0.029
146.	LP	(3)	O	19	/575.	RY*	(15)	Fe	67	0.34	165.72	0.311
146.	LP	(3)	O	19	/579.	RY*	(19)	Fe	67	0.15	730.27	0.434
146.	LP	(3)	O	19	/580.	RY*	(20)	Fe	67	0.71	48.36	0.242

from unit 2 to unit 1

55. BD (1) C 36 - H 61	/610. BD* (2) C 13 - C 15	0.03	0.55	0.006
55. BD (1) C 36 - H 61	/619. BD* (2) C 17 - C 18	0.05	0.56	0.007
66. BD (1) N 42 - N 43	/317. RY* (1) O 19	0.03	3.15	0.012
66. BD (1) N 42 - N 43	/318. RY* (2) O 19	0.05	2.25	0.014
67. BD (1) N 43 - C 44	/319. RY* (3) O 19	0.04	1.87	0.011
69. BD (1) C 44 - C 46	/319. RY* (3) O 19	0.04	1.88	0.011
70. BD (2) C 44 - C 46	/615. BD* (1) C 15 - O 19	0.04	0.63	0.007
74. BD (1) C 46 - C 48	/237. RY* (1) C 9	0.04	1.87	0.012
74. BD (1) C 46 - C 48	/239. RY* (3) C 9	0.07	1.26	0.012
75. BD (1) C 46 - O 65	/214. RY* (2) C 6	0.03	1.85	0.009
75. BD (1) C 46 - O 65	/215. RY* (3) C 6	0.03	2.00	0.009
80. BD (1) C 48 - H 58	/239. RY* (3) C 9	0.03	1.12	0.008
149. LP (1) N 42	/208. RY* (4) C 5	0.03	1.35	0.008
149. LP (1) N 42	/213. RY* (1) C 6	0.03	1.60	0.009
149. LP (1) N 42	/218. RY* (6) C 6	0.03	3.70	0.013
149. LP (1) N 42	/253. RY* (1) N 11	0.15	1.56	0.020
149. LP (1) N 42	/254. RY* (2) N 11	0.19	2.20	0.027
149. LP (1) N 42	/255. RY* (3) N 11	0.15	1.69	0.021
149. LP (1) N 42	/263. RY* (3) N 12	0.03	1.52	0.009
149. LP (1) N 42	/271. RY* (3) C 13	0.03	1.39	0.008
149. LP (1) N 42	/285. RY* (1) C 15	0.21	1.58	0.024
149. LP (1) N 42	/286. RY* (2) C 15	0.05	1.88	0.012
149. LP (1) N 42	/288. RY* (4) C 15	0.05	1.12	0.009
149. LP (1) N 42	/289. RY* (5) C 15	0.04	3.45	0.015
149. LP (1) N 42	/317. RY* (1) O 19	0.63	2.63	0.052
149. LP (1) N 42	/318. RY* (2) O 19	0.11	1.73	0.017
149. LP (1) N 42	/319. RY* (3) O 19	0.08	1.55	0.014
149. LP (1) N 42	/320. RY* (4) O 19	0.06	2.29	0.015
149. LP (1) N 42	/615. BD* (1) C 15 - O 19	0.95	0.77	0.035
150. LP (2) N 42	/318. RY* (2) O 19	0.10	1.60	0.017
157. LP (1) O 65	/253. RY* (1) N 11	0.04	1.71	0.010
157. LP (1) O 65	/254. RY* (2) N 11	0.08	2.35	0.018
157. LP (1) O 65	/255. RY* (3) N 11	0.03	1.84	0.010
157. LP (1) O 65	/256. RY* (4) N 11	0.04	2.14	0.011
157. LP (1) O 65	/285. RY* (1) C 15	0.04	1.73	0.011
157. LP (1) O 65	/317. RY* (1) O 19	0.05	2.78	0.015
157. LP (1) O 65	/594. BD* (2) C 5 - C 6	0.41	0.65	0.022
157. LP (1) O 65	/596. BD* (1) C 6 - N 11	0.09	0.99	0.012
157. LP (1) O 65	/606. BD* (1) N 11 - N 12	0.10	1.11	0.013
158. LP (2) O 65	/253. RY* (1) N 11	0.04	1.60	0.011
158. LP (2) O 65	/254. RY* (2) N 11	0.16	2.25	0.024
158. LP (2) O 65	/255. RY* (3) N 11	0.11	1.73	0.018
158. LP (2) O 65	/263. RY* (3) N 12	0.03	1.57	0.009
158. LP (2) O 65	/285. RY* (1) C 15	0.06	1.63	0.012
158. LP (2) O 65	/288. RY* (4) C 15	0.04	1.17	0.008
158. LP (2) O 65	/289. RY* (5) C 15	0.03	3.50	0.012
158. LP (2) O 65	/317. RY* (1) O 19	0.04	2.68	0.013
159. LP (3) O 65	/213. RY* (1) C 6	0.06	1.54	0.012
159. LP (3) O 65	/214. RY* (2) C 6	0.04	1.37	0.009
159. LP (3) O 65	/215. RY* (3) C 6	0.03	1.52	0.008
159. LP (3) O 65	/216. RY* (4) C 6	0.05	1.38	0.011
159. LP (3) O 65	/253. RY* (1) N 11	0.08	1.49	0.014
159. LP (3) O 65	/254. RY* (2) N 11	0.22	2.14	0.028
159. LP (3) O 65	/255. RY* (3) N 11	0.30	1.62	0.028
159. LP (3) O 65	/256. RY* (4) N 11	0.03	1.92	0.009
159. LP (3) O 65	/263. RY* (3) N 12	0.07	1.46	0.013
159. LP (3) O 65	/285. RY* (1) C 15	0.05	1.52	0.012
159. LP (3) O 65	/286. RY* (2) C 15	0.03	1.81	0.010
159. LP (3) O 65	/288. RY* (4) C 15	0.04	1.06	0.008
159. LP (3) O 65	/594. BD* (2) C 5 - C 6	0.24	0.44	0.013
159. LP (3) O 65	/596. BD* (1) C 6 - N 11	0.12	0.77	0.012
159. LP (3) O 65	/606. BD* (1) N 11 - N 12	0.47	0.89	0.026
651. BD* (2) C 44 - C 46	/603. BD* (2) C 9 - C 10	0.07	0.04	0.004
660. BD* (2) C 48 - C 49	/603. BD* (2) C 9 - C 10	0.74	0.01	0.009
within unit 2				
42. BD (1) C 32 - C 33	/371. RY* (1) C 34	0.66	1.97	0.046
42. BD (1) C 32 - C 33	/372. RY* (2) C 34	0.32	1.99	0.032
42. BD (1) C 32 - C 33	/379. RY* (1) C 35	0.65	2.20	0.048
42. BD (1) C 32 - C 33	/403. RY* (1) C 38	0.58	1.85	0.042
42. BD (1) C 32 - C 33	/624. BD* (1) C 32 - C 34	1.86	1.28	0.062

42. BD (1) C 32 - C 33	/627. BD* (1) C 33 - C 35	1.70	1.27	0.059
42. BD (1) C 32 - C 33	/629. BD* (1) C 33 - C 38	0.73	1.09	0.036
42. BD (1) C 32 - C 33	/631. BD* (1) C 34 - H 51	1.45	1.09	0.050
42. BD (1) C 32 - C 33	/633. BD* (1) C 35 - C 40	2.05	1.11	0.060
42. BD (1) C 32 - C 33	/638. BD* (1) C 38 - C 39	0.65	1.30	0.037
43. BD (1) C 32 - C 34	/364. RY* (2) C 33	0.81	2.18	0.053
43. BD (1) C 32 - C 34	/388. RY* (2) C 36	0.54	1.87	0.040
43. BD (1) C 32 - C 34	/623. BD* (1) C 32 - C 33	1.83	1.18	0.059
43. BD (1) C 32 - C 34	/626. BD* (1) C 32 - O 50	0.42	1.03	0.026
43. BD (1) C 32 - C 34	/629. BD* (1) C 33 - C 38	1.21	1.20	0.048
43. BD (1) C 32 - C 34	/630. BD* (1) C 34 - C 36	1.21	1.22	0.049
43. BD (1) C 32 - C 34	/631. BD* (1) C 34 - H 51	0.90	1.19	0.042
43. BD (1) C 32 - C 34	/636. BD* (1) C 36 - H 61	0.70	1.21	0.037
43. BD (1) C 32 - C 34	/663. BD* (1) O 50 - H 60	0.80	1.07	0.037
44. BD (2) C 32 - C 34	/366. RY* (4) C 33	0.35	0.89	0.023
44. BD (2) C 32 - C 34	/389. RY* (3) C 36	0.26	1.06	0.022
44. BD (2) C 32 - C 34	/499. RY* (1) O 50	0.35	1.18	0.027
44. BD (2) C 32 - C 34	/625. BD* (2) C 32 - C 34	0.27	0.31	0.012
44. BD (2) C 32 - C 34	/628. BD* (2) C 33 - C 35	5.84	0.32	0.057
44. BD (2) C 32 - C 34	/635. BD* (2) C 36 - C 37	6.63	0.31	0.059
45. BD (1) C 32 - O 50	/355. RY* (1) C 32	0.32	1.98	0.032
45. BD (1) C 32 - O 50	/371. RY* (1) C 34	0.29	2.22	0.032
45. BD (1) C 32 - O 50	/623. BD* (1) C 32 - C 33	0.27	1.32	0.024
45. BD (1) C 32 - O 50	/624. BD* (1) C 32 - C 34	0.54	1.52	0.036
45. BD (1) C 32 - O 50	/627. BD* (1) C 33 - C 35	0.58	1.52	0.037
45. BD (1) C 32 - O 50	/630. BD* (1) C 34 - C 36	1.12	1.36	0.049
46. BD (1) C 33 - C 35	/356. RY* (2) C 32	0.44	2.16	0.039
46. BD (1) C 33 - C 35	/395. RY* (1) C 37	0.38	1.86	0.034
46. BD (1) C 33 - C 35	/397. RY* (3) C 37	0.28	1.80	0.029
46. BD (1) C 33 - C 35	/403. RY* (1) C 38	0.36	1.93	0.033
46. BD (1) C 33 - C 35	/419. RY* (1) C 40	0.37	1.94	0.034
46. BD (1) C 33 - C 35	/623. BD* (1) C 32 - C 33	1.62	1.16	0.055
46. BD (1) C 33 - C 35	/626. BD* (1) C 32 - O 50	1.35	1.01	0.047
46. BD (1) C 33 - C 35	/629. BD* (1) C 33 - C 38	1.90	1.18	0.060
46. BD (1) C 33 - C 35	/632. BD* (1) C 35 - C 37	1.89	1.18	0.060
46. BD (1) C 33 - C 35	/633. BD* (1) C 35 - C 40	1.81	1.19	0.059
46. BD (1) C 33 - C 35	/637. BD* (1) C 37 - N 42	1.37	1.11	0.049
46. BD (1) C 33 - C 35	/640. BD* (1) C 38 - H 52	0.76	1.18	0.038
46. BD (1) C 33 - C 35	/645. BD* (1) C 40 - H 54	0.69	1.19	0.036
47. BD (2) C 33 - C 35	/358. RY* (4) C 32	0.30	0.83	0.022
47. BD (2) C 33 - C 35	/405. RY* (3) C 38	0.56	0.76	0.028
47. BD (2) C 33 - C 35	/421. RY* (3) C 40	0.51	0.84	0.028
47. BD (2) C 33 - C 35	/625. BD* (2) C 32 - C 34	6.35	0.30	0.056
47. BD (2) C 33 - C 35	/635. BD* (2) C 36 - C 37	6.71	0.30	0.057
47. BD (2) C 33 - C 35	/639. BD* (2) C 38 - C 39	5.80	0.31	0.055
47. BD (2) C 33 - C 35	/644. BD* (2) C 40 - C 41	5.78	0.32	0.056
48. BD (1) C 33 - C 38	/355. RY* (1) C 32	0.28	1.72	0.028
48. BD (1) C 33 - C 38	/356. RY* (2) C 32	0.36	2.06	0.035
48. BD (1) C 33 - C 38	/379. RY* (1) C 35	0.41	2.18	0.038
48. BD (1) C 33 - C 38	/380. RY* (2) C 35	0.27	2.17	0.031
48. BD (1) C 33 - C 38	/411. RY* (1) C 39	0.37	1.82	0.033
48. BD (1) C 33 - C 38	/412. RY* (2) C 39	0.77	1.95	0.049
48. BD (1) C 33 - C 38	/623. BD* (1) C 32 - C 33	0.61	1.06	0.032
48. BD (1) C 33 - C 38	/624. BD* (1) C 32 - C 34	0.71	1.26	0.038
48. BD (1) C 33 - C 38	/627. BD* (1) C 33 - C 35	2.15	1.26	0.066
48. BD (1) C 33 - C 38	/632. BD* (1) C 35 - C 37	2.35	1.09	0.064
48. BD (1) C 33 - C 38	/638. BD* (1) C 38 - C 39	1.40	1.29	0.054
48. BD (1) C 33 - C 38	/640. BD* (1) C 38 - H 52	0.31	1.08	0.023
48. BD (1) C 33 - C 38	/642. BD* (1) C 39 - H 53	1.44	1.09	0.050
49. BD (1) C 34 - C 36	/355. RY* (1) C 32	0.64	1.73	0.042
49. BD (1) C 34 - C 36	/356. RY* (2) C 32	0.38	2.07	0.035
49. BD (1) C 34 - C 36	/396. RY* (2) C 37	0.54	1.75	0.039
49. BD (1) C 34 - C 36	/397. RY* (3) C 37	0.60	1.71	0.041
49. BD (1) C 34 - C 36	/624. BD* (1) C 32 - C 34	1.29	1.27	0.051
49. BD (1) C 34 - C 36	/626. BD* (1) C 32 - O 50	2.50	0.92	0.061
49. BD (1) C 34 - C 36	/631. BD* (1) C 34 - H 51	0.36	1.09	0.025
49. BD (1) C 34 - C 36	/634. BD* (1) C 36 - C 37	1.53	1.28	0.056
49. BD (1) C 34 - C 36	/636. BD* (1) C 36 - H 61	0.35	1.10	0.025
49. BD (1) C 34 - C 36	/637. BD* (1) C 37 - N 42	3.03	1.02	0.070
50. BD (1) C 34 - H 51	/356. RY* (2) C 32	0.47	1.93	0.038
50. BD (1) C 34 - H 51	/360. RY* (6) C 32	0.27	8.59	0.061

50. BD (1) C 34 - H 51	/387. RY* (1) C 36	0.52	1.88	0.040
50. BD (1) C 34 - H 51	/623. BD* (1) C 32 - C 33	3.06	0.93	0.068
50. BD (1) C 34 - H 51	/624. BD* (1) C 32 - C 34	0.72	1.13	0.036
50. BD (1) C 34 - H 51	/626. BD* (1) C 32 - O 50	0.84	0.79	0.032
50. BD (1) C 34 - H 51	/634. BD* (1) C 36 - C 37	1.46	1.15	0.052
51. BD (1) C 35 - C 37	/363. RY* (1) C 33	0.73	2.29	0.052
51. BD (1) C 35 - C 37	/387. RY* (1) C 36	0.46	2.01	0.039
51. BD (1) C 35 - C 37	/419. RY* (1) C 40	0.55	1.84	0.041
51. BD (1) C 35 - C 37	/435. RY* (1) N 42	0.30	1.89	0.030
51. BD (1) C 35 - C 37	/627. BD* (1) C 33 - C 35	1.94	1.27	0.063
51. BD (1) C 35 - C 37	/629. BD* (1) C 33 - C 38	2.01	1.09	0.059
51. BD (1) C 35 - C 37	/633. BD* (1) C 35 - C 40	0.88	1.10	0.039
51. BD (1) C 35 - C 37	/634. BD* (1) C 36 - C 37	2.04	1.28	0.065
51. BD (1) C 35 - C 37	/636. BD* (1) C 36 - H 61	1.57	1.09	0.053
51. BD (1) C 35 - C 37	/637. BD* (1) C 37 - N 42	0.38	1.02	0.025
51. BD (1) C 35 - C 37	/643. BD* (1) C 40 - C 41	0.69	1.30	0.038
52. BD (1) C 35 - C 40	/363. RY* (1) C 33	0.38	2.28	0.037
52. BD (1) C 35 - C 40	/395. RY* (1) C 37	0.68	1.76	0.044
52. BD (1) C 35 - C 40	/427. RY* (1) C 41	0.38	1.83	0.034
52. BD (1) C 35 - C 40	/428. RY* (2) C 41	0.77	1.96	0.049
52. BD (1) C 35 - C 40	/623. BD* (1) C 32 - C 33	2.33	1.06	0.063
52. BD (1) C 35 - C 40	/627. BD* (1) C 33 - C 35	2.12	1.26	0.065
52. BD (1) C 35 - C 40	/632. BD* (1) C 35 - C 37	0.89	1.08	0.039
52. BD (1) C 35 - C 40	/634. BD* (1) C 36 - C 37	0.98	1.27	0.045
52. BD (1) C 35 - C 40	/643. BD* (1) C 40 - C 41	1.38	1.29	0.054
52. BD (1) C 35 - C 40	/645. BD* (1) C 40 - H 54	0.34	1.09	0.024
52. BD (1) C 35 - C 40	/646. BD* (1) C 41 - H 55	1.50	1.09	0.051
53. BD (1) C 36 - C 37	/372. RY* (2) C 34	0.76	2.10	0.051
53. BD (1) C 36 - C 37	/380. RY* (2) C 35	0.74	2.28	0.052
53. BD (1) C 36 - C 37	/435. RY* (1) N 42	0.60	2.01	0.044
53. BD (1) C 36 - C 37	/630. BD* (1) C 34 - C 36	1.38	1.23	0.052
53. BD (1) C 36 - C 37	/631. BD* (1) C 34 - H 51	0.65	1.20	0.035
53. BD (1) C 36 - C 37	/632. BD* (1) C 35 - C 37	2.05	1.21	0.063
53. BD (1) C 36 - C 37	/633. BD* (1) C 35 - C 40	1.13	1.21	0.047
53. BD (1) C 36 - C 37	/636. BD* (1) C 36 - H 61	1.01	1.21	0.044
53. BD (1) C 36 - C 37	/637. BD* (1) C 37 - N 42	1.00	1.13	0.043
53. BD (1) C 36 - C 37	/647. BD* (1) N 42 - N 43	0.74	1.25	0.038
54. BD (2) C 36 - C 37	/373. RY* (3) C 34	0.60	0.79	0.029
54. BD (2) C 36 - C 37	/382. RY* (4) C 35	0.44	0.90	0.026
54. BD (2) C 36 - C 37	/625. BD* (2) C 32 - C 34	6.50	0.32	0.058
54. BD (2) C 36 - C 37	/628. BD* (2) C 33 - C 35	5.74	0.32	0.057
54. BD (2) C 36 - C 37	/647. BD* (1) N 42 - N 43	1.72	0.78	0.048
55. BD (1) C 36 - H 61	/371. RY* (1) C 34	0.49	1.84	0.038
55. BD (1) C 36 - H 61	/395. RY* (1) C 37	0.69	1.65	0.043
55. BD (1) C 36 - H 61	/396. RY* (2) C 37	0.39	1.63	0.032
55. BD (1) C 36 - H 61	/624. BD* (1) C 32 - C 34	1.33	1.15	0.049
55. BD (1) C 36 - H 61	/632. BD* (1) C 35 - C 37	3.21	0.97	0.071
55. BD (1) C 36 - H 61	/634. BD* (1) C 36 - C 37	0.88	1.16	0.040
55. BD (1) C 36 - H 61	/637. BD* (1) C 37 - N 42	0.62	0.89	0.030
56. BD (1) C 37 - N 42	/380. RY* (2) C 35	0.57	2.31	0.046
56. BD (1) C 37 - N 42	/387. RY* (1) C 36	0.55	2.16	0.044
56. BD (1) C 37 - N 42	/443. RY* (1) N 43	0.64	1.62	0.041
56. BD (1) C 37 - N 42	/444. RY* (2) N 43	0.34	1.65	0.030
56. BD (1) C 37 - N 42	/627. BD* (1) C 33 - C 35	0.71	1.41	0.040
56. BD (1) C 37 - N 42	/630. BD* (1) C 34 - C 36	1.22	1.25	0.049
56. BD (1) C 37 - N 42	/632. BD* (1) C 35 - C 37	0.62	1.24	0.035
56. BD (1) C 37 - N 42	/634. BD* (1) C 36 - C 37	1.09	1.42	0.050
56. BD (1) C 37 - N 42	/647. BD* (1) N 42 - N 43	0.68	1.28	0.037
56. BD (1) C 37 - N 42	/648. BD* (1) N 43 - C 44	2.77	1.15	0.072
57. BD (1) C 38 - C 39	/363. RY* (1) C 33	0.57	2.39	0.047
57. BD (1) C 38 - C 39	/364. RY* (2) C 33	0.39	2.17	0.037
57. BD (1) C 38 - C 39	/427. RY* (1) C 41	0.48	1.94	0.038
57. BD (1) C 38 - C 39	/428. RY* (2) C 41	0.42	2.07	0.037
57. BD (1) C 38 - C 39	/623. BD* (1) C 32 - C 33	1.34	1.17	0.050
57. BD (1) C 38 - C 39	/629. BD* (1) C 33 - C 38	1.49	1.19	0.053
57. BD (1) C 38 - C 39	/640. BD* (1) C 38 - H 52	0.88	1.19	0.041
57. BD (1) C 38 - C 39	/641. BD* (1) C 39 - C 41	1.19	1.22	0.048
57. BD (1) C 38 - C 39	/642. BD* (1) C 39 - H 53	0.80	1.19	0.039
57. BD (1) C 38 - C 39	/646. BD* (1) C 41 - H 55	0.78	1.20	0.039
58. BD (2) C 38 - C 39	/366. RY* (4) C 33	0.35	0.88	0.023
58. BD (2) C 38 - C 39	/429. RY* (3) C 41	0.47	0.79	0.026

58. BD (2) C 38 - C 39	/628. BD* (2) C 33 - C 35	6.26	0.31	0.058
58. BD (2) C 38 - C 39	/644. BD* (2) C 40 - C 41	6.24	0.32	0.056
59. BD (1) C 38 - H 52	/363. RY* (1) C 33	0.70	2.17	0.050
59. BD (1) C 38 - H 52	/411. RY* (1) C 39	0.67	1.71	0.043
59. BD (1) C 38 - H 52	/627. BD* (1) C 33 - C 35	1.61	1.14	0.054
59. BD (1) C 38 - H 52	/638. BD* (1) C 38 - C 39	0.75	1.17	0.038
59. BD (1) C 38 - H 52	/641. BD* (1) C 39 - C 41	2.73	0.99	0.066
59. BD (1) C 38 - H 52	/642. BD* (1) C 39 - H 53	0.48	0.97	0.027
60. BD (1) C 39 - C 41	/404. RY* (2) C 38	0.92	1.77	0.051
60. BD (1) C 39 - C 41	/420. RY* (2) C 40	0.89	1.73	0.050
60. BD (1) C 39 - C 41	/638. BD* (1) C 38 - C 39	1.32	1.29	0.052
60. BD (1) C 39 - C 41	/640. BD* (1) C 38 - H 52	1.57	1.08	0.052
60. BD (1) C 39 - C 41	/642. BD* (1) C 39 - H 53	0.30	1.09	0.023
60. BD (1) C 39 - C 41	/643. BD* (1) C 40 - C 41	1.35	1.30	0.053
60. BD (1) C 39 - C 41	/645. BD* (1) C 40 - H 54	1.62	1.10	0.053
60. BD (1) C 39 - C 41	/646. BD* (1) C 41 - H 55	0.29	1.09	0.023
61. BD (1) C 39 - H 53	/403. RY* (1) C 38	0.74	1.71	0.045
61. BD (1) C 39 - H 53	/427. RY* (1) C 41	0.62	1.70	0.041
61. BD (1) C 39 - H 53	/629. BD* (1) C 33 - C 38	3.15	0.95	0.069
61. BD (1) C 39 - H 53	/638. BD* (1) C 38 - C 39	0.71	1.16	0.036
61. BD (1) C 39 - H 53	/640. BD* (1) C 38 - H 52	0.54	0.95	0.029
61. BD (1) C 39 - H 53	/643. BD* (1) C 40 - C 41	1.39	1.17	0.051
62. BD (1) C 40 - C 41	/379. RY* (1) C 35	0.76	2.29	0.053
62. BD (1) C 40 - C 41	/380. RY* (2) C 35	0.27	2.27	0.031
62. BD (1) C 40 - C 41	/411. RY* (1) C 39	0.44	1.93	0.037
62. BD (1) C 40 - C 41	/412. RY* (2) C 39	0.42	2.05	0.037
62. BD (1) C 40 - C 41	/632. BD* (1) C 35 - C 37	1.26	1.19	0.049
62. BD (1) C 40 - C 41	/633. BD* (1) C 35 - C 40	1.44	1.20	0.053
62. BD (1) C 40 - C 41	/641. BD* (1) C 39 - C 41	1.24	1.21	0.049
62. BD (1) C 40 - C 41	/642. BD* (1) C 39 - H 53	0.78	1.19	0.039
62. BD (1) C 40 - C 41	/645. BD* (1) C 40 - H 54	0.87	1.20	0.041
62. BD (1) C 40 - C 41	/646. BD* (1) C 41 - H 55	0.84	1.19	0.040
63. BD (2) C 40 - C 41	/382. RY* (4) C 35	0.34	0.88	0.023
63. BD (2) C 40 - C 41	/413. RY* (3) C 39	0.49	0.78	0.026
63. BD (2) C 40 - C 41	/628. BD* (2) C 33 - C 35	6.44	0.31	0.058
63. BD (2) C 40 - C 41	/639. BD* (2) C 38 - C 39	7.19	0.31	0.060
64. BD (1) C 40 - H 54	/379. RY* (1) C 35	0.69	2.05	0.048
64. BD (1) C 40 - H 54	/427. RY* (1) C 41	0.60	1.70	0.040
64. BD (1) C 40 - H 54	/627. BD* (1) C 33 - C 35	1.74	1.13	0.056
64. BD (1) C 40 - H 54	/641. BD* (1) C 39 - C 41	2.76	0.98	0.066
64. BD (1) C 40 - H 54	/643. BD* (1) C 40 - C 41	0.80	1.16	0.038
64. BD (1) C 40 - H 54	/646. BD* (1) C 41 - H 55	0.50	0.96	0.028
65. BD (1) C 41 - H 55	/411. RY* (1) C 39	0.60	1.69	0.040
65. BD (1) C 41 - H 55	/419. RY* (1) C 40	0.76	1.71	0.046
65. BD (1) C 41 - H 55	/633. BD* (1) C 35 - C 40	3.15	0.96	0.070
65. BD (1) C 41 - H 55	/638. BD* (1) C 38 - C 39	1.41	1.15	0.051
65. BD (1) C 41 - H 55	/643. BD* (1) C 40 - C 41	0.76	1.16	0.038
65. BD (1) C 41 - H 55	/645. BD* (1) C 40 - H 54	0.49	0.96	0.028
66. BD (1) N 42 - N 43	/395. RY* (1) C 37	0.58	2.11	0.044
66. BD (1) N 42 - N 43	/437. RY* (3) N 42	0.32	2.26	0.034
66. BD (1) N 42 - N 43	/451. RY* (1) C 44	0.51	2.12	0.042
66. BD (1) N 42 - N 43	/452. RY* (2) C 44	1.03	2.36	0.062
66. BD (1) N 42 - N 43	/634. BD* (1) C 36 - C 37	0.39	1.62	0.032
66. BD (1) N 42 - N 43	/635. BD* (2) C 36 - C 37	0.38	1.01	0.026
66. BD (1) N 42 - N 43	/637. BD* (1) C 37 - N 42	0.67	1.35	0.038
66. BD (1) N 42 - N 43	/648. BD* (1) N 43 - C 44	0.25	1.35	0.023
66. BD (1) N 42 - N 43	/649. BD* (1) C 44 - C 45	0.60	1.42	0.037
67. BD (1) N 43 - C 44	/435. RY* (1) N 42	0.28	2.03	0.030
67. BD (1) N 43 - C 44	/436. RY* (2) N 42	0.73	2.34	0.052
67. BD (1) N 43 - C 44	/437. RY* (3) N 42	0.57	2.06	0.043
67. BD (1) N 43 - C 44	/459. RY* (1) C 45	0.42	2.05	0.037
67. BD (1) N 43 - C 44	/467. RY* (1) C 46	0.46	1.94	0.038
67. BD (1) N 43 - C 44	/637. BD* (1) C 37 - N 42	2.52	1.15	0.068
67. BD (1) N 43 - C 44	/649. BD* (1) C 44 - C 45	0.37	1.22	0.027
67. BD (1) N 43 - C 44	/650. BD* (1) C 44 - C 46	0.97	1.37	0.046
67. BD (1) N 43 - C 44	/652. BD* (1) C 45 - C 47	0.48	1.41	0.033
67. BD (1) N 43 - C 44	/655. BD* (1) C 46 - C 48	1.41	1.22	0.052
68. BD (1) C 44 - C 45	/443. RY* (1) N 43	0.34	1.50	0.029
68. BD (1) C 44 - C 45	/470. RY* (4) C 46	0.36	1.68	0.031
68. BD (1) C 44 - C 45	/475. RY* (1) C 47	0.42	1.85	0.035
68. BD (1) C 44 - C 45	/476. RY* (2) C 47	0.71	1.99	0.048

68. BD (1) C 44 - C 45	/647. BD* (1) N 42 - N 43	1.50	1.16	0.053
68. BD (1) C 44 - C 45	/648. BD* (1) N 43 - C 44	0.52	1.03	0.029
68. BD (1) C 44 - C 45	/650. BD* (1) C 44 - C 46	1.77	1.25	0.060
68. BD (1) C 44 - C 45	/652. BD* (1) C 45 - C 47	1.16	1.30	0.049
68. BD (1) C 44 - C 45	/654. BD* (1) C 45 - H 56	0.31	1.10	0.023
68. BD (1) C 44 - C 45	/656. BD* (1) C 46 - O 65	2.69	0.96	0.064
68. BD (1) C 44 - C 45	/658. BD* (1) C 47 - H 57	1.37	1.11	0.049
69. BD (1) C 44 - C 46	/445. RY* (3) N 43	0.30	1.86	0.030
69. BD (1) C 44 - C 46	/459. RY* (1) C 45	0.53	2.06	0.042
69. BD (1) C 44 - C 46	/483. RY* (1) C 48	0.65	2.07	0.047
69. BD (1) C 44 - C 46	/648. BD* (1) N 43 - C 44	0.63	1.15	0.034
69. BD (1) C 44 - C 46	/649. BD* (1) C 44 - C 45	1.71	1.22	0.058
69. BD (1) C 44 - C 46	/654. BD* (1) C 45 - H 56	0.63	1.22	0.035
69. BD (1) C 44 - C 46	/655. BD* (1) C 46 - C 48	1.70	1.22	0.058
69. BD (1) C 44 - C 46	/656. BD* (1) C 46 - O 65	0.35	1.08	0.025
69. BD (1) C 44 - C 46	/661. BD* (1) C 48 - H 58	0.61	1.23	0.035
70. BD (2) C 44 - C 46	/444. RY* (2) N 43	0.36	1.17	0.027
70. BD (2) C 44 - C 46	/461. RY* (3) C 45	0.52	0.86	0.028
70. BD (2) C 44 - C 46	/485. RY* (3) C 48	0.27	1.08	0.023
70. BD (2) C 44 - C 46	/653. BD* (2) C 45 - C 47	5.90	0.34	0.056
70. BD (2) C 44 - C 46	/660. BD* (2) C 48 - C 49	6.90	0.34	0.062
71. BD (1) C 45 - C 47	/452. RY* (2) C 44	0.87	2.14	0.054
71. BD (1) C 45 - C 47	/491. RY* (1) C 49	0.52	1.96	0.041
71. BD (1) C 45 - C 47	/492. RY* (2) C 49	0.40	2.10	0.037
71. BD (1) C 45 - C 47	/648. BD* (1) N 43 - C 44	1.34	1.12	0.049
71. BD (1) C 45 - C 47	/649. BD* (1) C 44 - C 45	1.11	1.19	0.046
71. BD (1) C 45 - C 47	/654. BD* (1) C 45 - H 56	0.85	1.20	0.040
71. BD (1) C 45 - C 47	/657. BD* (1) C 47 - C 49	1.04	1.21	0.045
71. BD (1) C 45 - C 47	/658. BD* (1) C 47 - H 57	0.70	1.20	0.037
71. BD (1) C 45 - C 47	/662. BD* (1) C 49 - H 59	0.80	1.20	0.039
72. BD (2) C 45 - C 47	/493. RY* (3) C 49	0.54	0.79	0.027
72. BD (2) C 45 - C 47	/651. BD* (2) C 44 - C 46	6.18	0.29	0.057
72. BD (2) C 45 - C 47	/660. BD* (2) C 48 - C 49	6.05	0.32	0.057
73. BD (1) C 45 - H 56	/451. RY* (1) C 44	0.44	1.65	0.034
73. BD (1) C 45 - H 56	/475. RY* (1) C 47	0.61	1.71	0.041
73. BD (1) C 45 - H 56	/648. BD* (1) N 43 - C 44	0.31	0.88	0.021
73. BD (1) C 45 - H 56	/650. BD* (1) C 44 - C 46	1.67	1.11	0.055
73. BD (1) C 45 - H 56	/652. BD* (1) C 45 - C 47	0.74	1.15	0.037
73. BD (1) C 45 - H 56	/657. BD* (1) C 47 - C 49	2.68	0.97	0.064
73. BD (1) C 45 - H 56	/658. BD* (1) C 47 - H 57	0.50	0.96	0.028
74. BD (1) C 46 - C 48	/451. RY* (1) C 44	0.49	1.80	0.038
74. BD (1) C 46 - C 48	/491. RY* (1) C 49	0.35	1.87	0.032
74. BD (1) C 46 - C 48	/492. RY* (2) C 49	0.73	2.00	0.049
74. BD (1) C 46 - C 48	/648. BD* (1) N 43 - C 44	2.51	1.03	0.064
74. BD (1) C 46 - C 48	/650. BD* (1) C 44 - C 46	1.89	1.25	0.062
74. BD (1) C 46 - C 48	/659. BD* (1) C 48 - C 49	1.21	1.31	0.050
74. BD (1) C 46 - C 48	/661. BD* (1) C 48 - H 58	0.29	1.11	0.023
74. BD (1) C 46 - C 48	/662. BD* (1) C 49 - H 59	1.48	1.11	0.051
75. BD (1) C 46 - O 65	/451. RY* (1) C 44	0.38	2.01	0.035
75. BD (1) C 46 - O 65	/467. RY* (1) C 46	0.51	2.03	0.041
75. BD (1) C 46 - O 65	/483. RY* (1) C 48	0.27	2.16	0.031
75. BD (1) C 46 - O 65	/649. BD* (1) C 44 - C 45	1.26	1.31	0.051
75. BD (1) C 46 - O 65	/650. BD* (1) C 44 - C 46	0.68	1.46	0.040
75. BD (1) C 46 - O 65	/655. BD* (1) C 46 - C 48	0.30	1.31	0.025
75. BD (1) C 46 - O 65	/659. BD* (1) C 48 - C 49	0.48	1.52	0.034
76. BD (1) C 47 - C 49	/460. RY* (2) C 45	1.00	1.83	0.054
76. BD (1) C 47 - C 49	/483. RY* (1) C 48	0.26	1.94	0.028
76. BD (1) C 47 - C 49	/484. RY* (2) C 48	0.93	1.82	0.052
76. BD (1) C 47 - C 49	/652. BD* (1) C 45 - C 47	1.11	1.29	0.048
76. BD (1) C 47 - C 49	/654. BD* (1) C 45 - H 56	1.63	1.09	0.053
76. BD (1) C 47 - C 49	/658. BD* (1) C 47 - H 57	0.26	1.10	0.021
76. BD (1) C 47 - C 49	/659. BD* (1) C 48 - C 49	1.18	1.30	0.049
76. BD (1) C 47 - C 49	/661. BD* (1) C 48 - H 58	1.62	1.10	0.053
76. BD (1) C 47 - C 49	/662. BD* (1) C 49 - H 59	0.28	1.10	0.022
77. BD (1) C 47 - H 57	/459. RY* (1) C 45	0.80	1.79	0.048
77. BD (1) C 47 - H 57	/491. RY* (1) C 49	0.60	1.73	0.041
77. BD (1) C 47 - H 57	/649. BD* (1) C 44 - C 45	3.09	0.96	0.069
77. BD (1) C 47 - H 57	/652. BD* (1) C 45 - C 47	0.63	1.16	0.034
77. BD (1) C 47 - H 57	/654. BD* (1) C 45 - H 56	0.52	0.96	0.028
77. BD (1) C 47 - H 57	/659. BD* (1) C 48 - C 49	1.37	1.17	0.051
78. BD (1) C 48 - C 49	/467. RY* (1) C 46	0.46	1.91	0.038

78. BD (1) C 48 - C 49	/468. RY* (2) C 46	0.53	2.15	0.043
78. BD (1) C 48 - C 49	/475. RY* (1) C 47	0.50	1.94	0.039
78. BD (1) C 48 - C 49	/476. RY* (2) C 47	0.41	2.08	0.037
78. BD (1) C 48 - C 49	/521. RY* (1) H 58	0.26	1.20	0.022
78. BD (1) C 48 - C 49	/655. BD* (1) C 46 - C 48	0.95	1.19	0.043
78. BD (1) C 48 - C 49	/656. BD* (1) C 46 - O 65	1.28	1.05	0.046
78. BD (1) C 48 - C 49	/657. BD* (1) C 47 - C 49	1.10	1.21	0.046
78. BD (1) C 48 - C 49	/658. BD* (1) C 47 - H 57	0.78	1.19	0.038
78. BD (1) C 48 - C 49	/661. BD* (1) C 48 - H 58	0.90	1.20	0.041
78. BD (1) C 48 - C 49	/662. BD* (1) C 49 - H 59	0.80	1.20	0.039
79. BD (2) C 48 - C 49	/477. RY* (3) C 47	0.49	0.79	0.026
79. BD (2) C 48 - C 49	/651. BD* (2) C 44 - C 46	5.73	0.28	0.054
79. BD (2) C 48 - C 49	/653. BD* (2) C 45 - C 47	7.18	0.30	0.060
80. BD (1) C 48 - H 58	/467. RY* (1) C 46	0.36	1.68	0.031
80. BD (1) C 48 - H 58	/491. RY* (1) C 49	0.59	1.72	0.040
80. BD (1) C 48 - H 58	/650. BD* (1) C 44 - C 46	1.57	1.11	0.053
80. BD (1) C 48 - H 58	/656. BD* (1) C 46 - O 65	0.47	0.81	0.025
80. BD (1) C 48 - H 58	/657. BD* (1) C 47 - C 49	2.76	0.97	0.065
80. BD (1) C 48 - H 58	/659. BD* (1) C 48 - C 49	0.82	1.16	0.039
80. BD (1) C 48 - H 58	/662. BD* (1) C 49 - H 59	0.49	0.96	0.027
81. BD (1) C 49 - H 59	/475. RY* (1) C 47	0.57	1.70	0.040
81. BD (1) C 49 - H 59	/483. RY* (1) C 48	0.80	1.80	0.048
81. BD (1) C 49 - H 59	/652. BD* (1) C 45 - C 47	1.35	1.15	0.050
81. BD (1) C 49 - H 59	/655. BD* (1) C 46 - C 48	2.98	0.95	0.067
81. BD (1) C 49 - H 59	/659. BD* (1) C 48 - C 49	0.74	1.16	0.037
81. BD (1) C 49 - H 59	/661. BD* (1) C 48 - H 58	0.47	0.96	0.027
82. BD (1) O 50 - H 60	/355. RY* (1) C 32	1.34	1.80	0.062
82. BD (1) O 50 - H 60	/624. BD* (1) C 32 - C 34	1.72	1.34	0.061
105. CR (1) C 32	/363. RY* (1) C 33	0.63	11.77	0.108
105. CR (1) C 32	/372. RY* (2) C 34	1.14	11.47	0.145
105. CR (1) C 32	/624. BD* (1) C 32 - C 34	1.03	10.75	0.133
105. CR (1) C 32	/626. BD* (1) C 32 - O 50	0.66	10.40	0.105
105. CR (1) C 32	/627. BD* (1) C 33 - C 35	0.31	10.75	0.073
105. CR (1) C 32	/629. BD* (1) C 33 - C 38	0.34	10.57	0.076
105. CR (1) C 32	/630. BD* (1) C 34 - C 36	0.49	10.60	0.092
106. CR (1) C 33	/355. RY* (1) C 32	0.50	11.15	0.095
106. CR (1) C 33	/379. RY* (1) C 35	0.32	11.61	0.077
106. CR (1) C 33	/380. RY* (2) C 35	1.04	11.59	0.138
106. CR (1) C 33	/404. RY* (2) C 38	0.51	11.19	0.095
106. CR (1) C 33	/624. BD* (1) C 32 - C 34	0.30	10.69	0.072
106. CR (1) C 33	/626. BD* (1) C 32 - O 50	0.26	10.34	0.066
106. CR (1) C 33	/627. BD* (1) C 33 - C 35	0.66	10.69	0.106
106. CR (1) C 33	/632. BD* (1) C 35 - C 37	0.58	10.52	0.100
106. CR (1) C 33	/633. BD* (1) C 35 - C 40	0.52	10.52	0.094
106. CR (1) C 33	/638. BD* (1) C 38 - C 39	0.27	10.71	0.068
107. CR (1) C 34	/356. RY* (2) C 32	1.65	11.46	0.174
107. CR (1) C 34	/508. RY* (2) H 51	0.29	12.29	0.076
107. CR (1) C 34	/623. BD* (1) C 32 - C 33	0.67	10.46	0.107
107. CR (1) C 34	/624. BD* (1) C 32 - C 34	0.34	10.66	0.076
107. CR (1) C 34	/626. BD* (1) C 32 - O 50	0.40	10.31	0.081
107. CR (1) C 34	/634. BD* (1) C 36 - C 37	0.36	10.67	0.079
108. CR (1) C 35	/364. RY* (2) C 33	1.17	11.48	0.146
108. CR (1) C 35	/396. RY* (2) C 37	0.42	11.16	0.086
108. CR (1) C 35	/397. RY* (3) C 37	0.41	11.12	0.085
108. CR (1) C 35	/420. RY* (2) C 40	0.57	11.15	0.101
108. CR (1) C 35	/623. BD* (1) C 32 - C 33	0.56	10.48	0.098
108. CR (1) C 35	/627. BD* (1) C 33 - C 35	0.58	10.68	0.100
108. CR (1) C 35	/629. BD* (1) C 33 - C 38	0.51	10.50	0.093
108. CR (1) C 35	/634. BD* (1) C 36 - C 37	0.35	10.70	0.078
108. CR (1) C 35	/637. BD* (1) C 37 - N 42	0.27	10.43	0.067
108. CR (1) C 35	/643. BD* (1) C 40 - C 41	0.28	10.72	0.070
109. CR (1) C 36	/372. RY* (2) C 34	0.26	11.38	0.069
109. CR (1) C 36	/395. RY* (1) C 37	0.32	11.17	0.075
109. CR (1) C 36	/397. RY* (3) C 37	1.23	11.11	0.148
109. CR (1) C 36	/398. RY* (4) C 37	0.33	10.99	0.077
109. CR (1) C 36	/624. BD* (1) C 32 - C 34	0.33	10.67	0.075
109. CR (1) C 36	/632. BD* (1) C 35 - C 37	0.67	10.49	0.107
109. CR (1) C 36	/634. BD* (1) C 36 - C 37	0.47	10.68	0.090
109. CR (1) C 36	/637. BD* (1) C 37 - N 42	0.51	10.41	0.093
110. CR (1) C 37	/379. RY* (1) C 35	0.66	11.63	0.111
110. CR (1) C 37	/388. RY* (2) C 36	1.01	11.20	0.134

110.	CR	(1)	C	37	/627.	BD*	(1)	C	33 - C	35	0.36	10.71	0.079
110.	CR	(1)	C	37	/630.	BD*	(1)	C	34 - C	36	0.49	10.55	0.091
110.	CR	(1)	C	37	/633.	BD*	(1)	C	35 - C	40	0.33	10.54	0.075
110.	CR	(1)	C	37	/634.	BD*	(1)	C	36 - C	37	0.75	10.72	0.114
110.	CR	(1)	C	37	/636.	BD*	(1)	C	36 - H	61	0.27	10.54	0.067
110.	CR	(1)	C	37	/647.	BD*	(1)	N	42 - N	43	0.40	10.58	0.083
111.	CR	(1)	C	38	/363.	RY*	(1)	C	33		0.25	11.69	0.069
111.	CR	(1)	C	38	/364.	RY*	(2)	C	33		0.34	11.47	0.079
111.	CR	(1)	C	38	/412.	RY*	(2)	C	39		1.15	11.35	0.144
111.	CR	(1)	C	38	/510.	RY*	(2)	H	52		0.27	12.26	0.073
111.	CR	(1)	C	38	/623.	BD*	(1)	C	32 - C	33	0.34	10.46	0.077
111.	CR	(1)	C	38	/627.	BD*	(1)	C	33 - C	35	0.37	10.66	0.079
111.	CR	(1)	C	38	/638.	BD*	(1)	C	38 - C	39	0.44	10.69	0.087
111.	CR	(1)	C	38	/641.	BD*	(1)	C	39 - C	41	0.53	10.51	0.095
112.	CR	(1)	C	39	/403.	RY*	(1)	C	38		0.37	11.23	0.082
112.	CR	(1)	C	39	/404.	RY*	(2)	C	38		0.63	11.16	0.106
112.	CR	(1)	C	39	/406.	RY*	(4)	C	38		0.40	11.09	0.085
112.	CR	(1)	C	39	/428.	RY*	(2)	C	41		0.44	11.37	0.089
112.	CR	(1)	C	39	/512.	RY*	(2)	H	53		0.32	11.98	0.078
112.	CR	(1)	C	39	/629.	BD*	(1)	C	33 - C	38	0.62	10.48	0.102
112.	CR	(1)	C	39	/638.	BD*	(1)	C	38 - C	39	0.38	10.68	0.081
112.	CR	(1)	C	39	/643.	BD*	(1)	C	40 - C	41	0.31	10.69	0.073
113.	CR	(1)	C	40	/380.	RY*	(2)	C	35		0.50	11.56	0.096
113.	CR	(1)	C	40	/428.	RY*	(2)	C	41		1.11	11.36	0.142
113.	CR	(1)	C	40	/514.	RY*	(2)	H	54		0.26	12.35	0.071
113.	CR	(1)	C	40	/627.	BD*	(1)	C	33 - C	35	0.42	10.65	0.085
113.	CR	(1)	C	40	/632.	BD*	(1)	C	35 - C	37	0.34	10.48	0.077
113.	CR	(1)	C	40	/641.	BD*	(1)	C	39 - C	41	0.54	10.51	0.095
113.	CR	(1)	C	40	/643.	BD*	(1)	C	40 - C	41	0.43	10.69	0.085
114.	CR	(1)	C	41	/412.	RY*	(2)	C	39		0.42	11.34	0.087
114.	CR	(1)	C	41	/419.	RY*	(1)	C	40		0.35	11.23	0.079
114.	CR	(1)	C	41	/420.	RY*	(2)	C	40		0.57	11.12	0.101
114.	CR	(1)	C	41	/422.	RY*	(4)	C	40		0.51	11.13	0.095
114.	CR	(1)	C	41	/516.	RY*	(2)	H	55		0.32	11.98	0.078
114.	CR	(1)	C	41	/633.	BD*	(1)	C	35 - C	40	0.61	10.49	0.102
114.	CR	(1)	C	41	/638.	BD*	(1)	C	38 - C	39	0.32	10.68	0.074
114.	CR	(1)	C	41	/643.	BD*	(1)	C	40 - C	41	0.41	10.69	0.084
115.	CR	(1)	N	42	/395.	RY*	(1)	C	37		0.85	15.34	0.144
115.	CR	(1)	N	42	/396.	RY*	(2)	C	37		0.33	15.31	0.089
115.	CR	(1)	N	42	/445.	RY*	(3)	N	43		0.86	15.29	0.145
115.	CR	(1)	N	42	/632.	BD*	(1)	C	35 - C	37	0.29	14.66	0.084
115.	CR	(1)	N	42	/634.	BD*	(1)	C	36 - C	37	0.27	14.85	0.080
115.	CR	(1)	N	42	/648.	BD*	(1)	N	43 - C	44	0.67	14.57	0.126
116.	CR	(1)	N	43	/437.	RY*	(3)	N	42		1.00	15.48	0.157
116.	CR	(1)	N	43	/451.	RY*	(1)	C	44		0.80	15.34	0.140
116.	CR	(1)	N	43	/452.	RY*	(2)	C	44		0.25	15.58	0.080
116.	CR	(1)	N	43	/637.	BD*	(1)	C	37 - N	42	0.63	14.58	0.122
116.	CR	(1)	N	43	/650.	BD*	(1)	C	44 - C	46	0.36	14.80	0.093
117.	CR	(1)	C	44	/460.	RY*	(2)	C	45		0.50	11.29	0.095
117.	CR	(1)	C	44	/467.	RY*	(1)	C	46		0.44	11.28	0.089
117.	CR	(1)	C	44	/468.	RY*	(2)	C	46		1.17	11.52	0.147
117.	CR	(1)	C	44	/647.	BD*	(1)	N	42 - N	43	0.45	10.61	0.088
117.	CR	(1)	C	44	/650.	BD*	(1)	C	44 - C	46	0.61	10.71	0.103
117.	CR	(1)	C	44	/652.	BD*	(1)	C	45 - C	47	0.28	10.75	0.069
117.	CR	(1)	C	44	/655.	BD*	(1)	C	46 - C	48	0.51	10.55	0.093
117.	CR	(1)	C	44	/656.	BD*	(1)	C	46 - O	65	0.40	10.41	0.083
118.	CR	(1)	C	45	/451.	RY*	(1)	C	44		0.31	11.19	0.075
118.	CR	(1)	C	45	/453.	RY*	(3)	C	44		0.31	10.95	0.073
118.	CR	(1)	C	45	/476.	RY*	(2)	C	47		1.02	11.37	0.136
118.	CR	(1)	C	45	/518.	RY*	(2)	H	56		0.31	12.30	0.078
118.	CR	(1)	C	45	/648.	BD*	(1)	N	43 - C	44	0.27	10.41	0.068
118.	CR	(1)	C	45	/650.	BD*	(1)	C	44 - C	46	0.41	10.64	0.084
118.	CR	(1)	C	45	/652.	BD*	(1)	C	45 - C	47	0.41	10.68	0.084
118.	CR	(1)	C	45	/657.	BD*	(1)	C	47 - C	49	0.51	10.51	0.093
119.	CR	(1)	C	46	/452.	RY*	(2)	C	44		1.56	11.51	0.169
119.	CR	(1)	C	46	/484.	RY*	(2)	C	48		0.57	11.29	0.101
119.	CR	(1)	C	46	/648.	BD*	(1)	N	43 - C	44	0.41	10.49	0.084
119.	CR	(1)	C	46	/649.	BD*	(1)	C	44 - C	45	0.49	10.56	0.091
119.	CR	(1)	C	46	/650.	BD*	(1)	C	44 - C	46	0.80	10.72	0.118
119.	CR	(1)	C	46	/656.	BD*	(1)	C	46 - O	65	0.47	10.42	0.089
119.	CR	(1)	C	46	/659.	BD*	(1)	C	48 - C	49	0.29	10.77	0.071

120. CR (1) C 47	/459. RY* (1) C 45	0.41	11.33	0.086
120. CR (1) C 47	/460. RY* (2) C 45	0.80	11.23	0.119
120. CR (1) C 47	/492. RY* (2) C 49	0.47	11.40	0.093
120. CR (1) C 47	/520. RY* (2) H 57	0.33	12.00	0.079
120. CR (1) C 47	/649. BD* (1) C 44 - C 45	0.56	10.49	0.097
120. CR (1) C 47	/652. BD* (1) C 45 - C 47	0.36	10.69	0.079
120. CR (1) C 47	/659. BD* (1) C 48 - C 49	0.29	10.70	0.071
121. CR (1) C 48	/470. RY* (4) C 46	0.48	11.06	0.092
121. CR (1) C 48	/492. RY* (2) C 49	1.02	11.39	0.136
121. CR (1) C 48	/522. RY* (2) H 58	0.28	12.41	0.075
121. CR (1) C 48	/650. BD* (1) C 44 - C 46	0.43	10.64	0.086
121. CR (1) C 48	/657. BD* (1) C 47 - C 49	0.53	10.50	0.094
121. CR (1) C 48	/659. BD* (1) C 48 - C 49	0.44	10.69	0.087
122. CR (1) C 49	/476. RY* (2) C 47	0.48	11.37	0.093
122. CR (1) C 49	/483. RY* (1) C 48	0.40	11.32	0.085
122. CR (1) C 49	/484. RY* (2) C 48	0.81	11.21	0.120
122. CR (1) C 49	/524. RY* (2) H 59	0.31	11.98	0.077
122. CR (1) C 49	/652. BD* (1) C 45 - C 47	0.29	10.68	0.071
122. CR (1) C 49	/655. BD* (1) C 46 - C 48	0.53	10.48	0.095
122. CR (1) C 49	/659. BD* (1) C 48 - C 49	0.39	10.69	0.081
123. CR (1) O 50	/355. RY* (1) C 32	1.16	20.04	0.192
123. CR (1) O 50	/623. BD* (1) C 32 - C 33	0.28	19.38	0.094
125. CR (1) O 65	/467. RY* (1) C 46	1.38	20.11	0.211
125. CR (1) O 65	/650. BD* (1) C 44 - C 46	0.29	19.54	0.096
149. LP (1) N 42	/395. RY* (1) C 37	0.63	1.58	0.041
149. LP (1) N 42	/435. RY* (1) N 42	0.31	1.70	0.030
149. LP (1) N 42	/436. RY* (2) N 42	0.94	2.02	0.056
149. LP (1) N 42	/437. RY* (3) N 42	0.77	1.73	0.047
149. LP (1) N 42	/444. RY* (2) N 43	0.79	1.32	0.042
149. LP (1) N 42	/541. RY* (1) O 65	0.34	2.54	0.038
149. LP (1) N 42	/632. BD* (1) C 35 - C 37	3.16	0.90	0.068
149. LP (1) N 42	/635. BD* (2) C 36 - C 37	1.25	0.49	0.033
150. LP (2) N 42	/397. RY* (3) C 37	0.31	1.40	0.027
150. LP (2) N 42	/443. RY* (1) N 43	1.13	1.17	0.047
150. LP (2) N 42	/444. RY* (2) N 43	0.39	1.20	0.028
150. LP (2) N 42	/632. BD* (1) C 35 - C 37	1.32	0.78	0.041
150. LP (2) N 42	/634. BD* (1) C 36 - C 37	2.17	0.97	0.059
150. LP (2) N 42	/635. BD* (2) C 36 - C 37	4.66	0.37	0.054
151. LP (1) N 43	/436. RY* (2) N 42	0.37	1.96	0.034
151. LP (1) N 43	/437. RY* (3) N 42	0.47	1.68	0.036
151. LP (1) N 43	/451. RY* (1) C 44	0.93	1.53	0.048
151. LP (1) N 43	/649. BD* (1) C 44 - C 45	0.52	0.83	0.026
151. LP (1) N 43	/650. BD* (1) C 44 - C 46	3.80	0.99	0.077
152. LP (2) N 43	/435. RY* (1) N 42	0.79	1.51	0.047
152. LP (2) N 43	/649. BD* (1) C 44 - C 45	0.29	0.70	0.019
152. LP (2) N 43	/651. BD* (2) C 44 - C 46	18.06	0.27	0.093
153. LP (1) O 50	/355. RY* (1) C 32	1.44	1.67	0.062
153. LP (1) O 50	/623. BD* (1) C 32 - C 33	3.54	1.01	0.076
154. LP (2) O 50	/625. BD* (2) C 32 - C 34	12.11	0.35	0.086
157. LP (1) O 65	/467. RY* (1) C 46	1.13	1.77	0.057
157. LP (1) O 65	/650. BD* (1) C 44 - C 46	3.08	1.20	0.077
158. LP (2) O 65	/467. RY* (1) C 46	0.56	1.66	0.039
158. LP (2) O 65	/541. RY* (1) O 65	0.44	2.59	0.043
158. LP (2) O 65	/650. BD* (1) C 44 - C 46	0.50	1.09	0.030
158. LP (2) O 65	/651. BD* (2) C 44 - C 46	0.69	0.51	0.026
158. LP (2) O 65	/655. BD* (1) C 46 - C 48	3.46	0.94	0.072
159. LP (3) O 65	/469. RY* (3) C 46	0.39	1.11	0.027
159. LP (3) O 65	/541. RY* (1) O 65	0.37	2.48	0.039
159. LP (3) O 65	/651. BD* (2) C 44 - C 46	8.54	0.40	0.081
159. LP (3) O 65	/655. BD* (1) C 46 - C 48	0.45	0.83	0.025
625. BD* (2) C 32 - C 34	/358. RY* (4) C 32	0.52	0.53	0.058
625. BD* (2) C 32 - C 34	/373. RY* (3) C 34	1.16	0.48	0.082
628. BD* (2) C 33 - C 35	/366. RY* (4) C 33	1.12	0.57	0.075
628. BD* (2) C 33 - C 35	/382. RY* (4) C 35	1.11	0.57	0.075
635. BD* (2) C 36 - C 37	/389. RY* (3) C 36	0.45	0.75	0.064
635. BD* (2) C 36 - C 37	/647. BD* (1) N 42 - N 43	0.51	0.46	0.051
639. BD* (2) C 38 - C 39	/405. RY* (3) C 38	0.71	0.46	0.074
639. BD* (2) C 38 - C 39	/413. RY* (3) C 39	0.76	0.47	0.078
644. BD* (2) C 40 - C 41	/421. RY* (3) C 40	0.57	0.53	0.074
644. BD* (2) C 40 - C 41	/429. RY* (3) C 41	0.69	0.47	0.077
651. BD* (2) C 44 - C 46	/454. RY* (4) C 44	0.80	0.67	0.061

651. BD* (2) C 44 - C 46	/469. RY* (3) C 46	0.65	0.71	0.056
651. BD* (2) C 44 - C 46	/653. BD* (2) C 45 - C 47	44.13	0.03	0.072
651. BD* (2) C 44 - C 46	/660. BD* (2) C 48 - C 49	38.44	0.03	0.072
653. BD* (2) C 45 - C 47	/461. RY* (3) C 45	0.77	0.52	0.075
653. BD* (2) C 45 - C 47	/477. RY* (3) C 47	0.77	0.48	0.072
660. BD* (2) C 48 - C 49	/485. RY* (3) C 48	0.47	0.74	0.065
660. BD* (2) C 48 - C 49	/493. RY* (3) C 49	0.95	0.47	0.074

from unit 2 to unit 3

54. BD (2) C 36 - C 37	/664. BD* (1) O 62 - H 63	0.07	0.59	0.008
54. BD (2) C 36 - C 37	/665. BD* (1) O 62 - H 64	0.07	0.60	0.009
66. BD (1) N 42 - N 43	/529. RY* (1) O 62	0.10	2.74	0.021
66. BD (1) N 42 - N 43	/530. RY* (2) O 62	0.03	2.35	0.010
67. BD (1) N 43 - C 44	/529. RY* (1) O 62	0.04	2.54	0.012
75. BD (1) C 46 - O 65	/529. RY* (1) O 62	0.05	2.64	0.015
149. LP (1) N 42	/529. RY* (1) O 62	0.67	2.22	0.050
149. LP (1) N 42	/530. RY* (2) O 62	0.15	1.83	0.021
149. LP (1) N 42	/531. RY* (3) O 62	0.08	1.52	0.014
149. LP (1) N 42	/532. RY* (4) O 62	0.08	1.68	0.015
149. LP (1) N 42	/537. RY* (1) H 63	0.15	0.98	0.016
149. LP (1) N 42	/539. RY* (1) H 64	0.21	1.08	0.020
149. LP (1) N 42	/665. BD* (1) O 62 - H 64	0.03	0.77	0.007
150. LP (2) N 42	/529. RY* (1) O 62	0.11	2.10	0.019
150. LP (2) N 42	/530. RY* (2) O 62	0.03	1.71	0.009
150. LP (2) N 42	/537. RY* (1) H 63	0.03	0.86	0.006
150. LP (2) N 42	/539. RY* (1) H 64	0.03	0.95	0.007
157. LP (1) O 65	/529. RY* (1) O 62	0.10	2.37	0.019
157. LP (1) O 65	/537. RY* (1) H 63	0.03	1.14	0.008
157. LP (1) O 65	/539. RY* (1) H 64	0.04	1.23	0.008
158. LP (2) O 65	/529. RY* (1) O 62	0.15	2.26	0.024
158. LP (2) O 65	/530. RY* (2) O 62	0.07	1.87	0.014
158. LP (2) O 65	/532. RY* (4) O 62	0.05	1.73	0.012
158. LP (2) O 65	/537. RY* (1) H 63	0.04	1.03	0.008
158. LP (2) O 65	/539. RY* (1) H 64	0.05	1.12	0.009
159. LP (3) O 65	/529. RY* (1) O 62	0.19	2.15	0.026
159. LP (3) O 65	/530. RY* (2) O 62	0.09	1.76	0.016
159. LP (3) O 65	/531. RY* (3) O 62	0.03	1.46	0.008
159. LP (3) O 65	/532. RY* (4) O 62	0.05	1.62	0.012
159. LP (3) O 65	/537. RY* (1) H 63	0.04	0.92	0.008
159. LP (3) O 65	/539. RY* (1) H 64	0.06	1.01	0.010
159. LP (3) O 65	/664. BD* (1) O 62 - H 63	0.04	0.70	0.007
635. BD* (2) C 36 - C 37	/665. BD* (1) O 62 - H 64	0.05	0.28	0.013

from unit 2 to unit 4

66. BD (1) N 42 - N 43	/551. RY* (3)C1 66	0.03	2.36	0.011
67. BD (1) N 43 - C 44	/551. RY* (3)C1 66	0.03	2.16	0.010
149. LP (1) N 42	/549. RY* (1)C1 66	0.15	1.04	0.016
149. LP (1) N 42	/550. RY* (2)C1 66	0.18	1.09	0.018
149. LP (1) N 42	/551. RY* (3)C1 66	0.42	1.84	0.036
150. LP (2) N 42	/549. RY* (1)C1 66	0.05	0.92	0.009
150. LP (2) N 42	/551. RY* (3)C1 66	0.03	1.72	0.009
157. LP (1) O 65	/550. RY* (2)C1 66	0.05	1.24	0.010
157. LP (1) O 65	/551. RY* (3)C1 66	0.08	1.99	0.016
158. LP (2) O 65	/549. RY* (1)C1 66	0.10	1.09	0.014
158. LP (2) O 65	/551. RY* (3)C1 66	0.17	1.88	0.023
159. LP (3) O 65	/549. RY* (1)C1 66	0.07	0.97	0.010
159. LP (3) O 65	/550. RY* (2)C1 66	0.03	1.02	0.007
159. LP (3) O 65	/551. RY* (3)C1 66	0.17	1.77	0.022

from unit 2 to unit 5

51. BD (1) C 35 - C 37	/169. LP* (6)Fe 67	0.08	1.17	0.013
53. BD (1) C 36 - C 37	/172. LP* (9)Fe 67	0.05	2.44	0.015
53. BD (1) C 36 - C 37	/561. RY* (1)Fe 67	0.03	1.40	0.008
53. BD (1) C 36 - C 37	/571. RY* (11)Fe 67	0.03	10.30	0.022
55. BD (1) C 36 - H 61	/169. LP* (6)Fe 67	0.04	1.05	0.008
55. BD (1) C 36 - H 61	/563. RY* (3)Fe 67	0.03	4.68	0.015
56. BD (1) C 37 - N 42	/169. LP* (6)Fe 67	0.85	1.31	0.045
56. BD (1) C 37 - N 42	/171. LP* (8)Fe 67	0.09	2.54	0.019
56. BD (1) C 37 - N 42	/567. RY* (7)Fe 67	0.09	2.85	0.020
56. BD (1) C 37 - N 42	/568. RY* (8)Fe 67	0.03	9.61	0.022
66. BD (1) N 42 - N 43	/169. LP* (6)Fe 67	1.39	1.51	0.062

66. BD (1) N 42 - N 43	/171. LP* (8) Fe 67	0.06	2.73	0.016
66. BD (1) N 42 - N 43	/172. LP* (9) Fe 67	0.05	2.66	0.015
66. BD (1) N 42 - N 43	/561. RY* (1) Fe 67	0.03	1.62	0.009
66. BD (1) N 42 - N 43	/562. RY* (2) Fe 67	0.04	2.59	0.013
66. BD (1) N 42 - N 43	/563. RY* (3) Fe 67	0.43	5.13	0.059
66. BD (1) N 42 - N 43	/568. RY* (8) Fe 67	0.08	9.80	0.035
66. BD (1) N 42 - N 43	/573. RY* (13) Fe 67	0.03	7.66	0.018
66. BD (1) N 42 - N 43	/575. RY* (15) Fe 67	0.10	166.36	0.165
66. BD (1) N 42 - N 43	/579. RY* (19) Fe 67	0.06	730.90	0.255
66. BD (1) N 42 - N 43	/580. RY* (20) Fe 67	0.19	49.00	0.123
67. BD (1) N 43 - C 44	/169. LP* (6) Fe 67	0.18	1.31	0.021
67. BD (1) N 43 - C 44	/172. LP* (9) Fe 67	0.07	2.47	0.017
67. BD (1) N 43 - C 44	/563. RY* (3) Fe 67	0.17	4.94	0.036
67. BD (1) N 43 - C 44	/567. RY* (7) Fe 67	0.04	2.84	0.014
67. BD (1) N 43 - C 44	/571. RY* (11) Fe 67	0.08	10.32	0.037
67. BD (1) N 43 - C 44	/575. RY* (15) Fe 67	0.08	166.16	0.148
67. BD (1) N 43 - C 44	/579. RY* (19) Fe 67	0.04	730.71	0.217
67. BD (1) N 43 - C 44	/580. RY* (20) Fe 67	0.17	48.80	0.116
68. BD (1) C 44 - C 45	/169. LP* (6) Fe 67	0.04	1.19	0.009
69. BD (1) C 44 - C 46	/169. LP* (6) Fe 67	0.14	1.31	0.018
69. BD (1) C 44 - C 46	/170. LP* (7) Fe 67	0.03	2.28	0.010
69. BD (1) C 44 - C 46	/563. RY* (3) Fe 67	0.03	4.94	0.016
70. BD (2) C 44 - C 46	/169. LP* (6) Fe 67	0.07	0.84	0.010
70. BD (2) C 44 - C 46	/171. LP* (8) Fe 67	0.03	2.06	0.010
74. BD (1) C 46 - C 48	/169. LP* (6) Fe 67	0.19	1.19	0.020
74. BD (1) C 46 - C 48	/171. LP* (8) Fe 67	0.05	2.41	0.013
74. BD (1) C 46 - C 48	/563. RY* (3) Fe 67	0.03	4.82	0.015
74. BD (1) C 46 - C 48	/580. RY* (20) Fe 67	0.03	48.68	0.046
75. BD (1) C 46 - O 65	/169. LP* (6) Fe 67	1.17	1.40	0.055
75. BD (1) C 46 - O 65	/171. LP* (8) Fe 67	0.27	2.62	0.034
75. BD (1) C 46 - O 65	/562. RY* (2) Fe 67	0.04	2.49	0.013
75. BD (1) C 46 - O 65	/563. RY* (3) Fe 67	0.10	5.03	0.028
75. BD (1) C 46 - O 65	/568. RY* (8) Fe 67	0.07	9.69	0.033
75. BD (1) C 46 - O 65	/569. RY* (9) Fe 67	0.10	10.05	0.039
75. BD (1) C 46 - O 65	/572. RY* (12) Fe 67	0.09	7.59	0.032
75. BD (1) C 46 - O 65	/578. RY* (18) Fe 67	0.04	30.73	0.047
110. CR (1) C 37	/169. LP* (6) Fe 67	0.03	10.61	0.024
115. CR (1) N 42	/169. LP* (6) Fe 67	1.22	14.74	0.184
116. CR (1) N 43	/169. LP* (6) Fe 67	0.05	14.73	0.036
119. CR (1) C 46	/169. LP* (6) Fe 67	0.07	10.65	0.037
125. CR (1) O 65	/169. LP* (6) Fe 67	1.41	19.47	0.227
149. LP (1) N 42	/169. LP* (6) Fe 67	19.24	0.98	0.183
149. LP (1) N 42	/170. LP* (7) Fe 67	0.06	1.95	0.014
149. LP (1) N 42	/171. LP* (8) Fe 67	0.15	2.20	0.023
149. LP (1) N 42	/172. LP* (9) Fe 67	1.49	2.14	0.073
149. LP (1) N 42	/562. RY* (2) Fe 67	0.31	2.07	0.033
149. LP (1) N 42	/563. RY* (3) Fe 67	2.96	4.61	0.151
149. LP (1) N 42	/567. RY* (7) Fe 67	0.82	2.52	0.059
149. LP (1) N 42	/570. RY* (10) Fe 67	0.13	6.32	0.037
149. LP (1) N 42	/571. RY* (11) Fe 67	0.09	10.00	0.039
149. LP (1) N 42	/572. RY* (12) Fe 67	0.03	7.17	0.018
149. LP (1) N 42	/573. RY* (13) Fe 67	0.23	7.14	0.053
149. LP (1) N 42	/574. RY* (14) Fe 67	0.06	6.98	0.027
149. LP (1) N 42	/575. RY* (15) Fe 67	1.11	165.83	0.554
149. LP (1) N 42	/579. RY* (19) Fe 67	0.51	730.38	0.787
149. LP (1) N 42	/580. RY* (20) Fe 67	2.23	48.47	0.425
150. LP (2) N 42	/169. LP* (6) Fe 67	4.86	0.86	0.086
150. LP (2) N 42	/171. LP* (8) Fe 67	0.11	2.08	0.020
150. LP (2) N 42	/172. LP* (9) Fe 67	0.41	2.02	0.037
150. LP (2) N 42	/562. RY* (2) Fe 67	0.03	1.94	0.010
150. LP (2) N 42	/563. RY* (3) Fe 67	0.46	4.49	0.059
150. LP (2) N 42	/567. RY* (7) Fe 67	0.20	2.40	0.029
150. LP (2) N 42	/570. RY* (10) Fe 67	0.06	6.20	0.024
150. LP (2) N 42	/571. RY* (11) Fe 67	0.04	9.88	0.025
150. LP (2) N 42	/575. RY* (15) Fe 67	0.16	165.71	0.214
150. LP (2) N 42	/579. RY* (19) Fe 67	0.07	730.26	0.302
150. LP (2) N 42	/580. RY* (20) Fe 67	0.33	48.35	0.165
151. LP (1) N 43	/169. LP* (6) Fe 67	0.28	0.92	0.022
151. LP (1) N 43	/172. LP* (9) Fe 67	0.12	2.08	0.020
151. LP (1) N 43	/561. RY* (1) Fe 67	0.04	1.04	0.008
151. LP (1) N 43	/563. RY* (3) Fe 67	0.03	4.55	0.015

157. LP (1) O 65	/169. LP* (6) Fe 67	4.72	1.13	0.099
157. LP (1) O 65	/171. LP* (8) Fe 67	0.11	2.35	0.021
157. LP (1) O 65	/172. LP* (9) Fe 67	0.03	2.29	0.010
157. LP (1) O 65	/561. RY* (1) Fe 67	0.03	1.24	0.008
157. LP (1) O 65	/562. RY* (2) Fe 67	0.03	2.22	0.011
157. LP (1) O 65	/563. RY* (3) Fe 67	0.76	4.76	0.076
157. LP (1) O 65	/565. RY* (5) Fe 67	0.09	2.08	0.018
157. LP (1) O 65	/566. RY* (6) Fe 67	0.08	2.38	0.018
157. LP (1) O 65	/567. RY* (7) Fe 67	0.17	2.67	0.027
157. LP (1) O 65	/570. RY* (10) Fe 67	0.04	6.47	0.021
157. LP (1) O 65	/573. RY* (13) Fe 67	0.05	7.29	0.025
157. LP (1) O 65	/575. RY* (15) Fe 67	0.21	165.98	0.238
157. LP (1) O 65	/579. RY* (19) Fe 67	0.12	730.53	0.371
157. LP (1) O 65	/580. RY* (20) Fe 67	0.42	48.62	0.182
158. LP (2) O 65	/169. LP* (6) Fe 67	10.86	1.03	0.143
158. LP (2) O 65	/170. LP* (7) Fe 67	0.06	2.00	0.014
158. LP (2) O 65	/171. LP* (8) Fe 67	1.11	2.25	0.064
158. LP (2) O 65	/561. RY* (1) Fe 67	0.12	1.14	0.015
158. LP (2) O 65	/563. RY* (3) Fe 67	1.15	4.66	0.094
158. LP (2) O 65	/565. RY* (5) Fe 67	0.12	1.97	0.020
158. LP (2) O 65	/566. RY* (6) Fe 67	0.03	2.28	0.010
158. LP (2) O 65	/567. RY* (7) Fe 67	0.54	2.56	0.048
158. LP (2) O 65	/570. RY* (10) Fe 67	0.06	6.37	0.025
158. LP (2) O 65	/573. RY* (13) Fe 67	0.10	7.18	0.034
158. LP (2) O 65	/575. RY* (15) Fe 67	0.44	165.88	0.347
158. LP (2) O 65	/579. RY* (19) Fe 67	0.21	730.43	0.502
158. LP (2) O 65	/580. RY* (20) Fe 67	0.90	48.52	0.267
159. LP (3) O 65	/169. LP* (6) Fe 67	9.47	0.92	0.124
159. LP (3) O 65	/171. LP* (8) Fe 67	0.98	2.14	0.059
159. LP (3) O 65	/172. LP* (9) Fe 67	0.10	2.08	0.019
159. LP (3) O 65	/561. RY* (1) Fe 67	0.26	1.03	0.021
159. LP (3) O 65	/563. RY* (3) Fe 67	1.09	4.55	0.091
159. LP (3) O 65	/565. RY* (5) Fe 67	0.07	1.86	0.015
159. LP (3) O 65	/567. RY* (7) Fe 67	0.26	2.45	0.033
159. LP (3) O 65	/569. RY* (9) Fe 67	0.03	9.57	0.024
159. LP (3) O 65	/570. RY* (10) Fe 67	0.03	6.26	0.017
159. LP (3) O 65	/573. RY* (13) Fe 67	0.06	7.07	0.026
159. LP (3) O 65	/575. RY* (15) Fe 67	0.41	165.77	0.339
159. LP (3) O 65	/579. RY* (19) Fe 67	0.18	730.32	0.474
159. LP (3) O 65	/580. RY* (20) Fe 67	0.85	48.41	0.263

from unit 3 to unit 1

83. BD (1) O 62 - H 63	/253. RY* (1) N 11	0.13	1.90	0.020
83. BD (1) O 62 - H 63	/254. RY* (2) N 11	0.13	2.55	0.023
83. BD (1) O 62 - H 63	/255. RY* (3) N 11	0.09	2.03	0.017
83. BD (1) O 62 - H 63	/256. RY* (4) N 11	0.04	2.33	0.012
83. BD (1) O 62 - H 63	/264. RY* (4) N 12	0.03	2.43	0.011
83. BD (1) O 62 - H 63	/285. RY* (1) C 15	0.04	1.92	0.011
83. BD (1) O 62 - H 63	/288. RY* (4) C 15	0.03	1.47	0.008
83. BD (1) O 62 - H 63	/289. RY* (5) C 15	0.03	3.80	0.013
83. BD (1) O 62 - H 63	/610. BD* (2) C 13 - C 15	0.04	0.83	0.008
84. BD (1) O 62 - H 64	/253. RY* (1) N 11	0.05	1.90	0.013
84. BD (1) O 62 - H 64	/254. RY* (2) N 11	0.05	2.55	0.014
84. BD (1) O 62 - H 64	/255. RY* (3) N 11	0.04	2.03	0.011
84. BD (1) O 62 - H 64	/615. BD* (1) C 15 - O 19	0.03	1.12	0.007
155. LP (1) O 62	/253. RY* (1) N 11	0.03	1.62	0.008
155. LP (1) O 62	/269. RY* (1) C 13	0.03	1.64	0.008
155. LP (1) O 62	/271. RY* (3) C 13	0.03	1.45	0.008
155. LP (1) O 62	/286. RY* (2) C 15	0.03	1.94	0.009
155. LP (1) O 62	/288. RY* (4) C 15	0.04	1.19	0.009
155. LP (1) O 62	/596. BD* (1) C 6 - N 11	0.04	0.90	0.008
155. LP (1) O 62	/615. BD* (1) C 15 - O 19	0.04	0.84	0.007
156. LP (2) O 62	/208. RY* (4) C 5	0.03	1.54	0.008
156. LP (2) O 62	/213. RY* (1) C 6	0.06	1.79	0.014
156. LP (2) O 62	/253. RY* (1) N 11	0.38	1.75	0.033
156. LP (2) O 62	/254. RY* (2) N 11	0.41	2.39	0.040
156. LP (2) O 62	/255. RY* (3) N 11	0.20	1.88	0.024
156. LP (2) O 62	/256. RY* (4) N 11	0.09	2.17	0.018
156. LP (2) O 62	/262. RY* (2) N 12	0.03	1.51	0.009
156. LP (2) O 62	/263. RY* (3) N 12	0.03	1.71	0.008
156. LP (2) O 62	/271. RY* (3) C 13	0.03	1.58	0.008

156. LP (2) O 62	/285. RY* (1) C 15	0.13	1.77	0.019
156. LP (2) O 62	/286. RY* (2) C 15	0.09	2.07	0.017
156. LP (2) O 62	/288. RY* (4) C 15	0.09	1.31	0.014
156. LP (2) O 62	/289. RY* (5) C 15	0.09	3.64	0.023
156. LP (2) O 62	/317. RY* (1) O 19	0.09	2.82	0.020
156. LP (2) O 62	/320. RY* (4) O 19	0.05	2.48	0.014
156. LP (2) O 62	/596. BD* (1) C 6 - N 11	0.06	1.03	0.010

from unit 3 to unit 2

83. BD (1) O 62 - H 63	/396. RY* (2) C 37	0.04	1.90	0.010
83. BD (1) O 62 - H 63	/398. RY* (4) C 37	0.03	1.75	0.009
83. BD (1) O 62 - H 63	/436. RY* (2) N 42	0.03	2.36	0.011
83. BD (1) O 62 - H 63	/437. RY* (3) N 42	0.05	2.08	0.013
84. BD (1) O 62 - H 64	/396. RY* (2) C 37	0.04	1.90	0.010
84. BD (1) O 62 - H 64	/436. RY* (2) N 42	0.03	2.36	0.011
84. BD (1) O 62 - H 64	/437. RY* (3) N 42	0.05	2.08	0.013
156. LP (2) O 62	/395. RY* (1) C 37	0.04	1.77	0.011
156. LP (2) O 62	/396. RY* (2) C 37	0.09	1.75	0.016
156. LP (2) O 62	/398. RY* (4) C 37	0.04	1.59	0.010
156. LP (2) O 62	/399. RY* (5) C 37	0.04	3.61	0.014
156. LP (2) O 62	/400. RY* (6) C 37	0.03	3.77	0.015
156. LP (2) O 62	/435. RY* (1) N 42	0.04	1.89	0.011
156. LP (2) O 62	/436. RY* (2) N 42	0.15	2.21	0.023
156. LP (2) O 62	/437. RY* (3) N 42	0.39	1.92	0.035
156. LP (2) O 62	/443. RY* (1) N 43	0.03	1.48	0.009
156. LP (2) O 62	/445. RY* (3) N 43	0.11	1.72	0.018
156. LP (2) O 62	/467. RY* (1) C 46	0.08	1.80	0.016
156. LP (2) O 62	/468. RY* (2) C 46	0.06	2.05	0.014
156. LP (2) O 62	/469. RY* (3) C 46	0.08	1.36	0.013
156. LP (2) O 62	/471. RY* (5) C 46	0.06	3.64	0.019
156. LP (2) O 62	/541. RY* (1) O 65	0.03	2.73	0.011
156. LP (2) O 62	/544. RY* (4) O 65	0.03	2.18	0.011
156. LP (2) O 62	/635. BD* (2) C 36 - C 37	0.03	0.68	0.006
156. LP (2) O 62	/637. BD* (1) C 37 - N 42	0.03	1.02	0.007
156. LP (2) O 62	/647. BD* (1) N 42 - N 43	0.16	1.14	0.017

within unit 3

124. CR (1) O 62	/538. RY* (2) H 63	0.46	21.17	0.125
124. CR (1) O 62	/540. RY* (2) H 64	0.36	21.17	0.110
155. LP (1) O 62	/538. RY* (2) H 63	0.27	2.71	0.034
155. LP (1) O 62	/540. RY* (2) H 64	0.28	2.71	0.035
156. LP (2) O 62	/529. RY* (1) O 62	2.11	2.41	0.091
156. LP (2) O 62	/530. RY* (2) O 62	0.34	2.02	0.033
156. LP (2) O 62	/532. RY* (4) O 62	0.33	1.87	0.032
156. LP (2) O 62	/537. RY* (1) H 63	0.76	1.17	0.038
156. LP (2) O 62	/539. RY* (1) H 64	0.61	1.27	0.036
156. LP (2) O 62	/540. RY* (2) H 64	0.29	2.83	0.037

from unit 3 to unit 4

83. BD (1) O 62 - H 63	/550. RY* (2)C1 66	0.03	1.43	0.008
83. BD (1) O 62 - H 63	/551. RY* (3)C1 66	0.08	2.18	0.017
84. BD (1) O 62 - H 64	/551. RY* (3)C1 66	0.05	2.18	0.013
156. LP (2) O 62	/549. RY* (1)C1 66	0.27	1.23	0.023
156. LP (2) O 62	/550. RY* (2)C1 66	0.30	1.28	0.025
156. LP (2) O 62	/551. RY* (3)C1 66	0.61	2.03	0.045

from unit 3 to unit 5

83. BD (1) O 62 - H 63	/169. LP* (6)Fe 67	2.99	1.33	0.086
83. BD (1) O 62 - H 63	/171. LP* (8)Fe 67	0.03	2.55	0.011
83. BD (1) O 62 - H 63	/561. RY* (1)Fe 67	0.03	1.44	0.008
83. BD (1) O 62 - H 63	/562. RY* (2)Fe 67	0.24	2.41	0.031
83. BD (1) O 62 - H 63	/563. RY* (3)Fe 67	0.82	4.95	0.081
83. BD (1) O 62 - H 63	/564. RY* (4)Fe 67	0.03	2.19	0.010
83. BD (1) O 62 - H 63	/565. RY* (5)Fe 67	0.03	2.27	0.010
83. BD (1) O 62 - H 63	/567. RY* (7)Fe 67	0.04	2.86	0.013
83. BD (1) O 62 - H 63	/568. RY* (8)Fe 67	0.09	9.62	0.038
83. BD (1) O 62 - H 63	/570. RY* (10)Fe 67	0.04	6.67	0.021
83. BD (1) O 62 - H 63	/573. RY* (13)Fe 67	0.07	7.48	0.028
83. BD (1) O 62 - H 63	/575. RY* (15)Fe 67	0.23	166.18	0.245
83. BD (1) O 62 - H 63	/579. RY* (19)Fe 67	0.12	730.72	0.372
83. BD (1) O 62 - H 63	/580. RY* (20)Fe 67	0.44	48.82	0.185

84. BD (1) O 62 - H 64	/169. LP* (6) Fe 67	2.87	1.33	0.085
84. BD (1) O 62 - H 64	/562. RY* (2) Fe 67	0.29	2.41	0.033
84. BD (1) O 62 - H 64	/563. RY* (3) Fe 67	0.62	4.95	0.070
84. BD (1) O 62 - H 64	/566. RY* (6) Fe 67	0.03	2.57	0.011
84. BD (1) O 62 - H 64	/567. RY* (7) Fe 67	0.13	2.86	0.024
84. BD (1) O 62 - H 64	/569. RY* (9) Fe 67	0.07	9.98	0.034
84. BD (1) O 62 - H 64	/573. RY* (13) Fe 67	0.04	7.48	0.021
84. BD (1) O 62 - H 64	/574. RY* (14) Fe 67	0.07	7.32	0.028
84. BD (1) O 62 - H 64	/575. RY* (15) Fe 67	0.18	166.18	0.220
84. BD (1) O 62 - H 64	/579. RY* (19) Fe 67	0.10	730.72	0.333
84. BD (1) O 62 - H 64	/580. RY* (20) Fe 67	0.36	48.82	0.166
124. CR (1) O 62	/169. LP* (6) Fe 67	1.59	19.51	0.241
155. LP (1) O 62	/169. LP* (6) Fe 67	1.18	1.05	0.048
155. LP (1) O 62	/172. LP* (9) Fe 67	0.05	2.21	0.013
155. LP (1) O 62	/561. RY* (1) Fe 67	0.03	1.16	0.007
155. LP (1) O 62	/563. RY* (3) Fe 67	0.22	4.68	0.040
155. LP (1) O 62	/566. RY* (6) Fe 67	0.03	2.30	0.011
155. LP (1) O 62	/575. RY* (15) Fe 67	0.04	165.90	0.109
155. LP (1) O 62	/579. RY* (19) Fe 67	0.03	730.45	0.179
155. LP (1) O 62	/580. RY* (20) Fe 67	0.08	48.54	0.081
156. LP (2) O 62	/169. LP* (6) Fe 67	22.76	1.17	0.221
156. LP (2) O 62	/171. LP* (8) Fe 67	2.46	2.39	0.098
156. LP (2) O 62	/172. LP* (9) Fe 67	0.07	2.33	0.016
156. LP (2) O 62	/561. RY* (1) Fe 67	0.22	1.28	0.021
156. LP (2) O 62	/562. RY* (2) Fe 67	0.05	2.26	0.014
156. LP (2) O 62	/563. RY* (3) Fe 67	3.73	4.80	0.171
156. LP (2) O 62	/564. RY* (4) Fe 67	0.05	2.03	0.013
156. LP (2) O 62	/565. RY* (5) Fe 67	0.31	2.12	0.032
156. LP (2) O 62	/566. RY* (6) Fe 67	0.11	2.42	0.021
156. LP (2) O 62	/567. RY* (7) Fe 67	0.55	2.71	0.049
156. LP (2) O 62	/568. RY* (8) Fe 67	0.41	9.46	0.080
156. LP (2) O 62	/569. RY* (9) Fe 67	0.30	9.82	0.070
156. LP (2) O 62	/570. RY* (10) Fe 67	0.17	6.51	0.043
156. LP (2) O 62	/572. RY* (12) Fe 67	0.36	7.37	0.066
156. LP (2) O 62	/573. RY* (13) Fe 67	0.30	7.33	0.060
156. LP (2) O 62	/574. RY* (14) Fe 67	0.18	7.17	0.045
156. LP (2) O 62	/575. RY* (15) Fe 67	1.36	166.02	0.608
156. LP (2) O 62	/576. RY* (16) Fe 67	0.03	30.73	0.038
156. LP (2) O 62	/577. RY* (17) Fe 67	0.22	7.23	0.051
156. LP (2) O 62	/578. RY* (18) Fe 67	0.10	30.50	0.071
156. LP (2) O 62	/579. RY* (19) Fe 67	0.64	730.57	0.874
156. LP (2) O 62	/580. RY* (20) Fe 67	2.64	48.66	0.458
156. LP (2) O 62	/581. RY* (21) Fe 67	0.04	30.42	0.044

from unit 4 to unit 1

160. LP (1) C1 66	/254. RY* (2) N 11	0.04	2.48	0.013
160. LP (1) C1 66	/317. RY* (1) O 19	0.04	2.91	0.014
160. LP (1) C1 66	/594. BD* (2) C 5 - C 6	0.03	0.78	0.007
161. LP (2) C1 66	/589. BD* (1) C 3 - C 5	0.03	0.76	0.006
161. LP (2) C1 66	/594. BD* (2) C 5 - C 6	0.16	0.34	0.010
161. LP (2) C1 66	/595. BD* (1) C 5 - H 31	0.05	0.73	0.008
162. LP (3) C1 66	/595. BD* (1) C 5 - H 31	0.06	0.72	0.009
163. LP (4) C1 66	/206. RY* (2) C 5	0.03	1.72	0.009
163. LP (4) C1 66	/208. RY* (4) C 5	0.07	1.41	0.013
163. LP (4) C1 66	/213. RY* (1) C 6	0.11	1.65	0.018
163. LP (4) C1 66	/215. RY* (3) C 6	0.07	1.63	0.013
163. LP (4) C1 66	/216. RY* (4) C 6	0.05	1.50	0.011
163. LP (4) C1 66	/218. RY* (6) C 6	0.05	3.75	0.017
163. LP (4) C1 66	/253. RY* (1) N 11	0.65	1.61	0.042
163. LP (4) C1 66	/254. RY* (2) N 11	1.04	2.25	0.062
163. LP (4) C1 66	/255. RY* (3) N 11	0.79	1.74	0.048
163. LP (4) C1 66	/256. RY* (4) N 11	0.08	2.04	0.016
163. LP (4) C1 66	/262. RY* (2) N 12	0.09	1.37	0.014
163. LP (4) C1 66	/263. RY* (3) N 12	0.14	1.57	0.019
163. LP (4) C1 66	/271. RY* (3) C 13	0.04	1.44	0.010
163. LP (4) C1 66	/285. RY* (1) C 15	0.39	1.63	0.033
163. LP (4) C1 66	/286. RY* (2) C 15	0.16	1.93	0.023
163. LP (4) C1 66	/288. RY* (4) C 15	0.16	1.17	0.018
163. LP (4) C1 66	/289. RY* (5) C 15	0.09	3.51	0.023
163. LP (4) C1 66	/317. RY* (1) O 19	0.25	2.68	0.033
163. LP (4) C1 66	/320. RY* (4) O 19	0.05	2.34	0.014

163. LP (4)C1 66	/606. BD* (1) N 11 - N 12	0.06	1.01	0.010
from unit 4 to unit 2				
160. LP (1)C1 66	/437. RY* (3) N 42	0.03	2.01	0.010
160. LP (1)C1 66	/644. BD* (2) C 40 - C 41	0.03	0.78	0.006
161. LP (2)C1 66	/628. BD* (2) C 33 - C 35	0.03	0.33	0.005
161. LP (2)C1 66	/635. BD* (2) C 36 - C 37	0.14	0.32	0.009
161. LP (2)C1 66	/637. BD* (1) C 37 - N 42	0.05	0.66	0.008
161. LP (2)C1 66	/644. BD* (2) C 40 - C 41	0.21	0.34	0.011
161. LP (2)C1 66	/647. BD* (1) N 42 - N 43	0.05	0.79	0.008
161. LP (2)C1 66	/656. BD* (1) C 46 - O 65	0.07	0.59	0.008
162. LP (3)C1 66	/543. RY* (3) O 65	0.07	1.76	0.014
162. LP (3)C1 66	/656. BD* (1) C 46 - O 65	0.81	0.58	0.027
163. LP (4)C1 66	/395. RY* (1) C 37	0.14	1.63	0.019
163. LP (4)C1 66	/396. RY* (2) C 37	0.22	1.61	0.024
163. LP (4)C1 66	/398. RY* (4) C 37	0.18	1.46	0.021
163. LP (4)C1 66	/399. RY* (5) C 37	0.04	3.47	0.015
163. LP (4)C1 66	/400. RY* (6) C 37	0.05	3.63	0.018
163. LP (4)C1 66	/402. RY* (8) C 37	0.05	3.28	0.016
163. LP (4)C1 66	/435. RY* (1) N 42	0.24	1.76	0.026
163. LP (4)C1 66	/436. RY* (2) N 42	1.11	2.07	0.062
163. LP (4)C1 66	/437. RY* (3) N 42	1.02	1.79	0.055
163. LP (4)C1 66	/438. RY* (4) N 42	0.05	1.98	0.013
163. LP (4)C1 66	/443. RY* (1) N 43	0.12	1.34	0.016
163. LP (4)C1 66	/444. RY* (2) N 43	0.05	1.37	0.010
163. LP (4)C1 66	/445. RY* (3) N 43	0.17	1.58	0.021
163. LP (4)C1 66	/453. RY* (3) C 44	0.08	1.41	0.014
163. LP (4)C1 66	/467. RY* (1) C 46	0.56	1.67	0.039
163. LP (4)C1 66	/468. RY* (2) C 46	0.15	1.91	0.022
163. LP (4)C1 66	/469. RY* (3) C 46	0.17	1.22	0.019
163. LP (4)C1 66	/470. RY* (4) C 46	0.04	1.52	0.010
163. LP (4)C1 66	/471. RY* (5) C 46	0.07	3.50	0.020
163. LP (4)C1 66	/541. RY* (1) O 65	1.15	2.59	0.070
163. LP (4)C1 66	/543. RY* (3) O 65	0.23	1.98	0.028
163. LP (4)C1 66	/544. RY* (4) O 65	0.31	2.04	0.032
163. LP (4)C1 66	/635. BD* (2) C 36 - C 37	0.04	0.54	0.006
163. LP (4)C1 66	/637. BD* (1) C 37 - N 42	0.07	0.88	0.010
163. LP (4)C1 66	/647. BD* (1) N 42 - N 43	0.11	1.00	0.014
163. LP (4)C1 66	/656. BD* (1) C 46 - O 65	0.85	0.80	0.033
from unit 4 to unit 3				
160. LP (1)C1 66	/529. RY* (1) O 62	0.03	2.50	0.010
160. LP (1)C1 66	/665. BD* (1) O 62 - H 64	0.03	1.05	0.007
162. LP (3)C1 66	/530. RY* (2) O 62	0.06	1.65	0.013
162. LP (3)C1 66	/665. BD* (1) O 62 - H 64	0.77	0.59	0.027
163. LP (4)C1 66	/529. RY* (1) O 62	2.38	2.27	0.095
163. LP (4)C1 66	/530. RY* (2) O 62	0.72	1.88	0.048
163. LP (4)C1 66	/531. RY* (3) O 62	0.46	1.58	0.035
163. LP (4)C1 66	/532. RY* (4) O 62	0.14	1.73	0.020
163. LP (4)C1 66	/537. RY* (1) H 63	0.67	1.04	0.034
163. LP (4)C1 66	/539. RY* (1) H 64	0.88	1.13	0.041
163. LP (4)C1 66	/664. BD* (1) O 62 - H 63	0.03	0.81	0.007
163. LP (4)C1 66	/665. BD* (1) O 62 - H 64	0.61	0.82	0.029
within unit 4				
163. LP (4)C1 66	/550. RY* (2)C1 66	0.55	1.14	0.032
163. LP (4)C1 66	/551. RY* (3)C1 66	1.38	1.89	0.066
from unit 4 to unit 5				
126. CR (1)C1 66	/169. LP* (6)Fe 67	0.28	100.54	0.229
127. CR (2)C1 66	/169. LP* (6)Fe 67	2.20	11.14	0.215
129. CR (4)C1 66	/169. LP* (6)Fe 67	0.13	7.67	0.043
130. CR (5)C1 66	/169. LP* (6)Fe 67	0.11	7.68	0.039
160. LP (1)C1 66	/169. LP* (6)Fe 67	2.70	1.26	0.080
160. LP (1)C1 66	/170. LP* (7)Fe 67	0.13	2.23	0.022
160. LP (1)C1 66	/562. RY* (2)Fe 67	0.03	2.34	0.011
160. LP (1)C1 66	/563. RY* (3)Fe 67	0.30	4.89	0.049
160. LP (1)C1 66	/567. RY* (7)Fe 67	0.08	2.80	0.018
160. LP (1)C1 66	/575. RY* (15)Fe 67	0.08	166.11	0.148
160. LP (1)C1 66	/579. RY* (19)Fe 67	0.05	730.66	0.246
160. LP (1)C1 66	/580. RY* (20)Fe 67	0.16	48.75	0.113

161. LP (2)Cl 66	/169. LP* (6)Fe 67	0.37	0.82	0.024
161. LP (2)Cl 66	/172. LP* (9)Fe 67	0.11	1.97	0.019
161. LP (2)Cl 66	/563. RY* (3)Fe 67	0.05	4.44	0.020
161. LP (2)Cl 66	/564. RY* (4)Fe 67	0.17	1.68	0.021
161. LP (2)Cl 66	/580. RY* (20)Fe 67	0.03	48.31	0.052
162. LP (3)Cl 66	/169. LP* (6)Fe 67	0.09	0.81	0.012
162. LP (3)Cl 66	/566. RY* (6)Fe 67	0.03	2.05	0.010
162. LP (3)Cl 66	/567. RY* (7)Fe 67	0.04	2.34	0.012
163. LP (4)Cl 66	/169. LP* (6)Fe 67	33.84	1.03	0.251
163. LP (4)Cl 66	/170. LP* (7)Fe 67	2.77	2.00	0.096
163. LP (4)Cl 66	/172. LP* (9)Fe 67	0.08	2.19	0.016
163. LP (4)Cl 66	/561. RY* (1)Fe 67	0.29	1.14	0.023
163. LP (4)Cl 66	/562. RY* (2)Fe 67	1.84	2.12	0.080
163. LP (4)Cl 66	/563. RY* (3)Fe 67	6.54	4.66	0.225
163. LP (4)Cl 66	/564. RY* (4)Fe 67	0.03	1.89	0.009
163. LP (4)Cl 66	/565. RY* (5)Fe 67	0.38	1.98	0.035
163. LP (4)Cl 66	/567. RY* (7)Fe 67	1.55	2.57	0.081
163. LP (4)Cl 66	/569. RY* (9)Fe 67	0.61	9.68	0.099
163. LP (4)Cl 66	/570. RY* (10)Fe 67	0.35	6.37	0.061
163. LP (4)Cl 66	/571. RY* (11)Fe 67	0.11	10.05	0.044
163. LP (4)Cl 66	/572. RY* (12)Fe 67	0.19	7.23	0.048
163. LP (4)Cl 66	/573. RY* (13)Fe 67	0.90	7.19	0.104
163. LP (4)Cl 66	/574. RY* (14)Fe 67	0.04	7.03	0.022
163. LP (4)Cl 66	/575. RY* (15)Fe 67	2.57	165.88	0.840
163. LP (4)Cl 66	/578. RY* (18)Fe 67	0.05	30.36	0.051
163. LP (4)Cl 66	/579. RY* (19)Fe 67	1.22	730.43	1.215
163. LP (4)Cl 66	/580. RY* (20)Fe 67	5.07	48.52	0.639
163. LP (4)Cl 66	/581. RY* (21)Fe 67	0.07	30.28	0.059

from unit 5 to unit 1

131. CR (1)Fe 67	/596. BD* (1) C 6 - N 11	0.03	256.20	0.108
131. CR (1)Fe 67	/615. BD* (1) C 15 - O 19	0.03	256.14	0.115
132. CR (2)Fe 67	/596. BD* (1) C 6 - N 11	0.15	29.53	0.086
132. CR (2)Fe 67	/606. BD* (1) N 11 - N 12	0.10	29.65	0.068
132. CR (2)Fe 67	/615. BD* (1) C 15 - O 19	0.17	29.47	0.091
133. CR (3)Fe 67	/255. RY* (3) N 11	0.09	5.99	0.030
133. CR (3)Fe 67	/285. RY* (1) C 15	0.04	5.88	0.019
133. CR (3)Fe 67	/317. RY* (1) O 19	0.03	6.93	0.018
133. CR (3)Fe 67	/596. BD* (1) C 6 - N 11	0.75	5.14	0.079
133. CR (3)Fe 67	/606. BD* (1) N 11 - N 12	0.48	5.25	0.063
133. CR (3)Fe 67	/607. BD* (1) N 12 - C 13	0.03	5.13	0.015
133. CR (3)Fe 67	/614. BD* (1) C 15 - C 17	0.05	5.20	0.020
133. CR (3)Fe 67	/615. BD* (1) C 15 - O 19	0.93	5.07	0.087
135. CR (5)Fe 67	/255. RY* (3) N 11	0.05	3.67	0.017
135. CR (5)Fe 67	/596. BD* (1) C 6 - N 11	0.16	2.82	0.027
135. CR (5)Fe 67	/606. BD* (1) N 11 - N 12	0.20	2.94	0.031
137. CR (7)Fe 67	/596. BD* (1) C 6 - N 11	0.11	2.82	0.022
137. CR (7)Fe 67	/615. BD* (1) C 15 - O 19	0.15	2.76	0.026
139. CR (9)Fe 67	/615. BD* (1) C 15 - O 19	0.26	2.77	0.034
164. LP (1)Fe 67	/596. BD* (1) C 6 - N 11	0.04	0.79	0.007
164. LP (1)Fe 67	/606. BD* (1) N 11 - N 12	0.05	0.91	0.009
164. LP (1)Fe 67	/614. BD* (1) C 15 - C 17	0.03	0.85	0.006
164. LP (1)Fe 67	/615. BD* (1) C 15 - O 19	0.03	0.73	0.006
165. LP (2)Fe 67	/254. RY* (2) N 11	0.07	2.15	0.016
165. LP (2)Fe 67	/317. RY* (1) O 19	0.34	2.58	0.038
165. LP (2)Fe 67	/596. BD* (1) C 6 - N 11	0.07	0.79	0.010
165. LP (2)Fe 67	/606. BD* (1) N 11 - N 12	0.07	0.91	0.010
166. LP (3)Fe 67	/254. RY* (2) N 11	0.03	2.15	0.009
166. LP (3)Fe 67	/255. RY* (3) N 11	0.03	1.64	0.009
166. LP (3)Fe 67	/256. RY* (4) N 11	0.03	1.93	0.010
166. LP (3)Fe 67	/257. RY* (5) N 11	0.03	4.35	0.015
166. LP (3)Fe 67	/285. RY* (1) C 15	0.04	1.53	0.010
166. LP (3)Fe 67	/317. RY* (1) O 19	0.44	2.58	0.042
166. LP (3)Fe 67	/596. BD* (1) C 6 - N 11	0.11	0.79	0.012
166. LP (3)Fe 67	/606. BD* (1) N 11 - N 12	0.31	0.90	0.021
167. LP (4)Fe 67	/253. RY* (1) N 11	0.03	1.51	0.009
167. LP (4)Fe 67	/254. RY* (2) N 11	0.19	2.15	0.025
167. LP (4)Fe 67	/255. RY* (3) N 11	0.15	1.64	0.020
167. LP (4)Fe 67	/256. RY* (4) N 11	0.06	1.93	0.013
167. LP (4)Fe 67	/263. RY* (3) N 12	0.03	1.47	0.008
167. LP (4)Fe 67	/317. RY* (1) O 19	0.46	2.58	0.043

167. LP (4) Fe 67	/606. BD* (1) N 11 - N 12	0.09	0.90	0.012
167. LP (4) Fe 67	/615. BD* (1) C 15 - O 19	0.17	0.72	0.014
168. LP (5) Fe 67	/253. RY* (1) N 11	0.56	1.49	0.036
168. LP (5) Fe 67	/254. RY* (2) N 11	1.30	2.14	0.067
168. LP (5) Fe 67	/255. RY* (3) N 11	0.33	1.62	0.029
168. LP (5) Fe 67	/259. RY* (7) N 11	0.04	35.57	0.045
168. LP (5) Fe 67	/262. RY* (2) N 12	0.03	1.26	0.007
168. LP (5) Fe 67	/264. RY* (4) N 12	0.04	2.02	0.011
168. LP (5) Fe 67	/285. RY* (1) C 15	0.03	1.52	0.009
168. LP (5) Fe 67	/288. RY* (4) C 15	0.03	1.06	0.007
168. LP (5) Fe 67	/289. RY* (5) C 15	0.03	3.39	0.012
168. LP (5) Fe 67	/317. RY* (1) O 19	0.05	2.57	0.015
168. LP (5) Fe 67	/596. BD* (1) C 6 - N 11	0.09	0.77	0.010
168. LP (5) Fe 67	/607. BD* (1) N 12 - C 13	0.04	0.77	0.007
168. LP (5) Fe 67	/615. BD* (1) C 15 - O 19	0.04	0.71	0.007
169. LP* (6) Fe 67	/176. RY* (4) C 1	0.07	0.05	0.006
169. LP* (6) Fe 67	/182. RY* (2) C 2	0.03	0.91	0.018
169. LP* (6) Fe 67	/200. RY* (4) C 4	0.03	0.15	0.007
169. LP* (6) Fe 67	/205. RY* (1) C 5	0.08	0.78	0.026
169. LP* (6) Fe 67	/206. RY* (2) C 5	0.35	0.69	0.050
169. LP* (6) Fe 67	/207. RY* (3) C 5	0.57	0.20	0.034
169. LP* (6) Fe 67	/208. RY* (4) C 5	1.23	0.37	0.069
169. LP* (6) Fe 67	/213. RY* (1) C 6	1.46	0.62	0.095
169. LP* (6) Fe 67	/214. RY* (2) C 6	0.19	0.45	0.030
169. LP* (6) Fe 67	/215. RY* (3) C 6	0.54	0.60	0.058
169. LP* (6) Fe 67	/216. RY* (4) C 6	0.60	0.47	0.054
169. LP* (6) Fe 67	/217. RY* (5) C 6	0.05	2.45	0.034
169. LP* (6) Fe 67	/218. RY* (6) C 6	0.17	2.72	0.069
169. LP* (6) Fe 67	/220. RY* (8) C 6	0.08	2.29	0.043
169. LP* (6) Fe 67	/237. RY* (1) C 9	0.03	0.68	0.015
169. LP* (6) Fe 67	/238. RY* (2) C 9	0.05	0.53	0.017
169. LP* (6) Fe 67	/247. RY* (3) C 10	0.32	0.03	0.010
169. LP* (6) Fe 67	/248. RY* (4) C 10	0.04	0.38	0.012
169. LP* (6) Fe 67	/253. RY* (1) N 11	9.61	0.57	0.237
169. LP* (6) Fe 67	/254. RY* (2) N 11	11.33	1.22	0.375
169. LP* (6) Fe 67	/255. RY* (3) N 11	9.85	0.70	0.269
169. LP* (6) Fe 67	/256. RY* (4) N 11	1.30	1.00	0.117
169. LP* (6) Fe 67	/258. RY* (6) N 11	0.13	3.48	0.070
169. LP* (6) Fe 67	/261. RY* (1) N 12	0.70	0.30	0.046
169. LP* (6) Fe 67	/262. RY* (2) N 12	1.67	0.34	0.076
169. LP* (6) Fe 67	/263. RY* (3) N 12	1.72	0.54	0.098
169. LP* (6) Fe 67	/264. RY* (4) N 12	0.20	1.10	0.048
169. LP* (6) Fe 67	/265. RY* (5) N 12	0.04	3.63	0.038
169. LP* (6) Fe 67	/269. RY* (1) C 13	0.06	0.59	0.019
169. LP* (6) Fe 67	/270. RY* (2) C 13	0.07	0.91	0.026
169. LP* (6) Fe 67	/271. RY* (3) C 13	1.18	0.41	0.071
169. LP* (6) Fe 67	/273. RY* (5) C 13	0.05	2.18	0.033
169. LP* (6) Fe 67	/274. RY* (6) C 13	0.03	2.93	0.032
169. LP* (6) Fe 67	/279. RY* (3) C 14	2.55	0.02	0.021
169. LP* (6) Fe 67	/280. RY* (4) C 14	0.14	0.39	0.024
169. LP* (6) Fe 67	/282. RY* (6) C 14	0.03	2.05	0.024
169. LP* (6) Fe 67	/285. RY* (1) C 15	7.44	0.60	0.212
169. LP* (6) Fe 67	/286. RY* (2) C 15	1.34	0.90	0.112
169. LP* (6) Fe 67	/288. RY* (4) C 15	6.46	0.14	0.097
169. LP* (6) Fe 67	/289. RY* (5) C 15	0.72	2.47	0.137
169. LP* (6) Fe 67	/293. RY* (1) C 16	0.08	0.67	0.023
169. LP* (6) Fe 67	/296. RY* (4) C 16	0.05	0.39	0.015
169. LP* (6) Fe 67	/301. RY* (1) C 17	0.26	0.78	0.046
169. LP* (6) Fe 67	/303. RY* (3) C 17	1.77	0.05	0.030
169. LP* (6) Fe 67	/304. RY* (4) C 17	0.09	0.41	0.019
169. LP* (6) Fe 67	/309. RY* (1) C 18	0.03	0.67	0.013
169. LP* (6) Fe 67	/317. RY* (1) O 19	4.65	1.65	0.282
169. LP* (6) Fe 67	/318. RY* (2) O 19	0.23	0.74	0.042
169. LP* (6) Fe 67	/319. RY* (3) O 19	0.17	0.57	0.032
169. LP* (6) Fe 67	/320. RY* (4) O 19	1.13	1.31	0.124
169. LP* (6) Fe 67	/321. RY* (5) O 19	0.11	4.51	0.073
169. LP* (6) Fe 67	/340. RY* (2) H 24	0.07	1.84	0.036
169. LP* (6) Fe 67	/609. BD* (1) C 13 - C 15	0.18	0.09	0.012

from unit 5 to unit 2
131. CR (1) Fe 67

/637. BD* (1) C 37 - N 42 0.03 256.19 0.114

131. CR (1)Fe 67	/656. BD* (1) C 46 - O 65	0.03	256.12	0.115
132. CR (2)Fe 67	/637. BD* (1) C 37 - N 42	0.17	29.52	0.091
132. CR (2)Fe 67	/647. BD* (1) N 42 - N 43	0.10	29.64	0.067
132. CR (2)Fe 67	/656. BD* (1) C 46 - O 65	0.18	29.44	0.092
133. CR (3)Fe 67	/437. RY* (3) N 42	0.10	6.03	0.032
133. CR (3)Fe 67	/467. RY* (1) C 46	0.04	5.91	0.019
133. CR (3)Fe 67	/541. RY* (1) O 65	0.04	6.84	0.021
133. CR (3)Fe 67	/637. BD* (1) C 37 - N 42	0.82	5.12	0.082
133. CR (3)Fe 67	/647. BD* (1) N 42 - N 43	0.47	5.25	0.063
133. CR (3)Fe 67	/648. BD* (1) N 43 - C 44	0.03	5.12	0.016
133. CR (3)Fe 67	/655. BD* (1) C 46 - C 48	0.05	5.19	0.021
133. CR (3)Fe 67	/656. BD* (1) C 46 - O 65	0.93	5.05	0.087
135. CR (5)Fe 67	/437. RY* (3) N 42	0.06	3.72	0.018
135. CR (5)Fe 67	/637. BD* (1) C 37 - N 42	0.10	2.81	0.022
135. CR (5)Fe 67	/647. BD* (1) N 42 - N 43	0.22	2.94	0.033
135. CR (5)Fe 67	/656. BD* (1) C 46 - O 65	0.11	2.74	0.022
137. CR (7)Fe 67	/637. BD* (1) C 37 - N 42	0.04	2.81	0.013
137. CR (7)Fe 67	/656. BD* (1) C 46 - O 65	0.26	2.74	0.034
139. CR (9)Fe 67	/637. BD* (1) C 37 - N 42	0.16	2.82	0.027
164. LP (1)Fe 67	/436. RY* (2) N 42	0.09	1.97	0.017
164. LP (1)Fe 67	/438. RY* (4) N 42	0.04	1.88	0.010
164. LP (1)Fe 67	/467. RY* (1) C 46	0.04	1.57	0.010
164. LP (1)Fe 67	/470. RY* (4) C 46	0.03	1.42	0.008
164. LP (1)Fe 67	/541. RY* (1) O 65	0.47	2.49	0.043
164. LP (1)Fe 67	/543. RY* (3) O 65	0.05	1.88	0.012
164. LP (1)Fe 67	/647. BD* (1) N 42 - N 43	0.05	0.90	0.008
164. LP (1)Fe 67	/656. BD* (1) C 46 - O 65	0.09	0.70	0.010
165. LP (2)Fe 67	/435. RY* (1) N 42	0.14	1.66	0.019
165. LP (2)Fe 67	/436. RY* (2) N 42	0.09	1.97	0.017
165. LP (2)Fe 67	/446. RY* (4) N 43	0.03	2.05	0.009
165. LP (2)Fe 67	/541. RY* (1) O 65	0.04	2.49	0.012
165. LP (2)Fe 67	/637. BD* (1) C 37 - N 42	0.13	0.78	0.013
165. LP (2)Fe 67	/647. BD* (1) N 42 - N 43	0.08	0.90	0.011
166. LP (3)Fe 67	/435. RY* (1) N 42	0.27	1.65	0.026
166. LP (3)Fe 67	/436. RY* (2) N 42	0.14	1.97	0.021
166. LP (3)Fe 67	/437. RY* (3) N 42	0.09	1.68	0.015
166. LP (3)Fe 67	/471. RY* (5) C 46	0.03	3.40	0.014
166. LP (3)Fe 67	/541. RY* (1) O 65	0.06	2.49	0.015
166. LP (3)Fe 67	/544. RY* (4) O 65	0.03	1.94	0.009
166. LP (3)Fe 67	/647. BD* (1) N 42 - N 43	0.09	0.90	0.011
166. LP (3)Fe 67	/656. BD* (1) C 46 - O 65	0.05	0.70	0.008
167. LP (4)Fe 67	/438. RY* (4) N 42	0.07	1.88	0.015
167. LP (4)Fe 67	/439. RY* (5) N 42	0.05	4.29	0.018
167. LP (4)Fe 67	/467. RY* (1) C 46	0.04	1.56	0.010
167. LP (4)Fe 67	/541. RY* (1) O 65	0.58	2.49	0.048
167. LP (4)Fe 67	/543. RY* (3) O 65	0.09	1.88	0.016
167. LP (4)Fe 67	/637. BD* (1) C 37 - N 42	0.13	0.77	0.013
167. LP (4)Fe 67	/647. BD* (1) N 42 - N 43	0.37	0.90	0.023
167. LP (4)Fe 67	/656. BD* (1) C 46 - O 65	0.03	0.70	0.006
168. LP (5)Fe 67	/435. RY* (1) N 42	0.44	1.64	0.034
168. LP (5)Fe 67	/436. RY* (2) N 42	1.19	1.96	0.061
168. LP (5)Fe 67	/437. RY* (3) N 42	0.39	1.67	0.032
168. LP (5)Fe 67	/441. RY* (7) N 42	0.04	35.62	0.050
168. LP (5)Fe 67	/443. RY* (1) N 43	0.03	1.23	0.008
168. LP (5)Fe 67	/445. RY* (3) N 43	0.05	1.47	0.011
168. LP (5)Fe 67	/446. RY* (4) N 43	0.03	2.04	0.010
168. LP (5)Fe 67	/541. RY* (1) O 65	0.13	2.48	0.023
168. LP (5)Fe 67	/544. RY* (4) O 65	0.05	1.93	0.012
168. LP (5)Fe 67	/637. BD* (1) C 37 - N 42	0.04	0.76	0.007
168. LP (5)Fe 67	/648. BD* (1) N 43 - C 44	0.04	0.76	0.007
168. LP (5)Fe 67	/656. BD* (1) C 46 - O 65	0.11	0.69	0.011
169. LP* (6)Fe 67	/379. RY* (1) C 35	0.03	1.02	0.017
169. LP* (6)Fe 67	/380. RY* (2) C 35	0.06	1.00	0.025
169. LP* (6)Fe 67	/381. RY* (3) C 35	0.07	0.43	0.018
169. LP* (6)Fe 67	/382. RY* (4) C 35	0.03	0.09	0.005
169. LP* (6)Fe 67	/387. RY* (1) C 36	0.05	0.84	0.021
169. LP* (6)Fe 67	/388. RY* (2) C 36	0.06	0.59	0.018
169. LP* (6)Fe 67	/389. RY* (3) C 36	0.10	0.25	0.016
169. LP* (6)Fe 67	/390. RY* (4) C 36	0.15	0.48	0.027
169. LP* (6)Fe 67	/391. RY* (5) C 36	0.03	2.04	0.024
169. LP* (6)Fe 67	/395. RY* (1) C 37	1.62	0.60	0.099

169. LP* (6)Fe 67	/396. RY* (2) C 37	2.13	0.57	0.113
169. LP* (6)Fe 67	/397. RY* (3) C 37	0.06	0.54	0.019
169. LP* (6)Fe 67	/398. RY* (4) C 37	2.33	0.42	0.101
169. LP* (6)Fe 67	/399. RY* (5) C 37	0.27	2.43	0.083
169. LP* (6)Fe 67	/400. RY* (6) C 37	0.20	2.60	0.073
169. LP* (6)Fe 67	/402. RY* (8) C 37	0.20	2.25	0.068
169. LP* (6)Fe 67	/421. RY* (3) C 40	0.19	0.05	0.010
169. LP* (6)Fe 67	/422. RY* (4) C 40	0.06	0.57	0.018
169. LP* (6)Fe 67	/435. RY* (1) N 42	3.54	0.72	0.161
169. LP* (6)Fe 67	/436. RY* (2) N 42	9.00	1.04	0.309
169. LP* (6)Fe 67	/437. RY* (3) N 42	12.12	0.75	0.308
169. LP* (6)Fe 67	/438. RY* (4) N 42	0.87	0.94	0.093
169. LP* (6)Fe 67	/440. RY* (6) N 42	0.03	3.47	0.034
169. LP* (6)Fe 67	/443. RY* (1) N 43	1.87	0.31	0.076
169. LP* (6)Fe 67	/444. RY* (2) N 43	1.24	0.34	0.065
169. LP* (6)Fe 67	/445. RY* (3) N 43	2.26	0.55	0.114
169. LP* (6)Fe 67	/446. RY* (4) N 43	0.13	1.12	0.039
169. LP* (6)Fe 67	/447. RY* (5) N 43	0.03	3.45	0.035
169. LP* (6)Fe 67	/451. RY* (1) C 44	0.05	0.61	0.018
169. LP* (6)Fe 67	/452. RY* (2) C 44	0.07	0.85	0.025
169. LP* (6)Fe 67	/453. RY* (3) C 44	1.24	0.38	0.070
169. LP* (6)Fe 67	/454. RY* (4) C 44	0.05	0.15	0.009
169. LP* (6)Fe 67	/455. RY* (5) C 44	0.03	2.22	0.025
169. LP* (6)Fe 67	/461. RY* (3) C 45	1.38	0.02	0.018
169. LP* (6)Fe 67	/462. RY* (4) C 45	0.08	0.40	0.018
169. LP* (6)Fe 67	/467. RY* (1) C 46	6.27	0.63	0.201
169. LP* (6)Fe 67	/468. RY* (2) C 46	1.18	0.87	0.103
169. LP* (6)Fe 67	/469. RY* (3) C 46	5.01	0.19	0.100
169. LP* (6)Fe 67	/470. RY* (4) C 46	0.50	0.49	0.051
169. LP* (6)Fe 67	/471. RY* (5) C 46	0.59	2.46	0.124
169. LP* (6)Fe 67	/474. RY* (8) C 46	0.03	2.18	0.028
169. LP* (6)Fe 67	/475. RY* (1) C 47	0.05	0.66	0.019
169. LP* (6)Fe 67	/478. RY* (4) C 47	0.03	0.43	0.011
169. LP* (6)Fe 67	/483. RY* (1) C 48	0.17	0.76	0.036
169. LP* (6)Fe 67	/486. RY* (4) C 48	0.41	0.26	0.033
169. LP* (6)Fe 67	/491. RY* (1) C 49	0.03	0.68	0.014
169. LP* (6)Fe 67	/515. RY* (1) H 55	0.14	0.02	0.006
169. LP* (6)Fe 67	/523. RY* (1) H 59	0.06	0.02	0.004
169. LP* (6)Fe 67	/528. RY* (2) H 61	0.05	1.79	0.030
169. LP* (6)Fe 67	/541. RY* (1) O 65	6.41	1.56	0.322
169. LP* (6)Fe 67	/543. RY* (3) O 65	0.25	0.95	0.050
169. LP* (6)Fe 67	/544. RY* (4) O 65	1.99	1.01	0.145
169. LP* (6)Fe 67	/545. RY* (5) O 65	0.09	4.51	0.064
169. LP* (6)Fe 67	/634. BD* (1) C 36 - C 37	0.15	0.11	0.013
169. LP* (6)Fe 67	/650. BD* (1) C 44 - C 46	0.24	0.06	0.012

from unit 5 to unit 3

135. CR (5)Fe 67	/664. BD* (1) O 62 - H 63	0.07	2.75	0.017
139. CR (9)Fe 67	/665. BD* (1) O 62 - H 64	0.05	2.76	0.015
164. LP (1)Fe 67	/529. RY* (1) O 62	0.15	2.17	0.023
164. LP (1)Fe 67	/539. RY* (1) H 64	0.04	1.03	0.008
164. LP (1)Fe 67	/664. BD* (1) O 62 - H 63	0.06	0.71	0.008
164. LP (1)Fe 67	/665. BD* (1) O 62 - H 64	0.07	0.72	0.009
165. LP (2)Fe 67	/529. RY* (1) O 62	0.15	2.17	0.023
166. LP (3)Fe 67	/529. RY* (1) O 62	0.09	2.17	0.018
167. LP (4)Fe 67	/664. BD* (1) O 62 - H 63	0.25	0.71	0.017
167. LP (4)Fe 67	/665. BD* (1) O 62 - H 64	0.14	0.72	0.013
168. LP (5)Fe 67	/530. RY* (2) O 62	0.03	1.77	0.009
168. LP (5)Fe 67	/537. RY* (1) H 63	0.06	0.92	0.009
168. LP (5)Fe 67	/539. RY* (1) H 64	0.04	1.01	0.008
168. LP (5)Fe 67	/664. BD* (1) O 62 - H 63	0.04	0.70	0.007
169. LP* (6)Fe 67	/529. RY* (1) O 62	15.14	1.24	0.441
169. LP* (6)Fe 67	/530. RY* (2) O 62	3.21	0.85	0.168
169. LP* (6)Fe 67	/531. RY* (3) O 62	1.99	0.54	0.106
169. LP* (6)Fe 67	/532. RY* (4) O 62	2.17	0.70	0.126
169. LP* (6)Fe 67	/535. RY* (7) O 62	0.04	4.34	0.043
169. LP* (6)Fe 67	/536. RY* (8) O 62	0.15	4.32	0.084
169. LP* (6)Fe 67	/539. RY* (1) H 64	38.14	0.09	0.193

from unit 5 to unit 4

165. LP (2)Fe 67	/549. RY* (1)C1 66	0.06	0.99	0.009
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168. LP (5)Fe	67	/550. RY* (2)Cl	66	0.04	1.03	0.008
168. LP (5)Fe	67	/551. RY* (3)Cl	66	0.07	1.78	0.014
169. LP* (6)Fe	67	/549. RY* (1)Cl	66	9.20	0.06	0.074
169. LP* (6)Fe	67	/550. RY* (2)Cl	66	18.89	0.11	0.144
169. LP* (6)Fe	67	/551. RY* (3)Cl	66	8.74	0.86	0.280
169. LP* (6)Fe	67	/552. RY* (4)Cl	66	0.31	0.22	0.027
169. LP* (6)Fe	67	/553. RY* (5)Cl	66	0.07	1.78	0.036
169. LP* (6)Fe	67	/556. RY* (8)Cl	66	0.08	1.73	0.037
169. LP* (6)Fe	67	/559. RY* (11)Cl	66	0.23	1.75	0.065

within unit 5

132. CR (2)Fe	67	/169. LP* (6)Fe	67	1.15	29.67	0.253
133. CR (3)Fe	67	/169. LP* (6)Fe	67	4.34	5.28	0.207
168. LP (5)Fe	67	/169. LP* (6)Fe	67	0.25	0.92	0.021
168. LP (5)Fe	67	/563. RY* (3)Fe	67	0.31	4.55	0.048
168. LP (5)Fe	67	/580. RY* (20)Fe	67	0.34	48.41	0.162
169. LP* (6)Fe	67	/172. LP* (9)Fe	67	0.37	1.16	0.066
169. LP* (6)Fe	67	/561. RY* (1)Fe	67	15.40	0.11	0.133
169. LP* (6)Fe	67	/562. RY* (2)Fe	67	8.07	1.08	0.301
169. LP* (6)Fe	67	/563. RY* (3)Fe	67	32.10	3.63	1.101
169. LP* (6)Fe	67	/564. RY* (4)Fe	67	0.26	0.86	0.049
169. LP* (6)Fe	67	/565. RY* (5)Fe	67	1.41	0.94	0.118
169. LP* (6)Fe	67	/567. RY* (7)Fe	67	10.96	1.54	0.420
169. LP* (6)Fe	67	/569. RY* (9)Fe	67	0.33	8.65	0.174
169. LP* (6)Fe	67	/570. RY* (10)Fe	67	1.58	5.34	0.297
169. LP* (6)Fe	67	/572. RY* (12)Fe	67	0.47	6.19	0.174
169. LP* (6)Fe	67	/573. RY* (13)Fe	67	2.11	6.16	0.368
169. LP* (6)Fe	67	/574. RY* (14)Fe	67	1.05	5.99	0.257
169. LP* (6)Fe	67	/575. RY* (15)Fe	67	10.89	164.85	4.335
169. LP* (6)Fe	67	/579. RY* (19)Fe	67	5.07	729.40	6.221
169. LP* (6)Fe	67	/580. RY* (20)Fe	67	22.26	47.49	3.327