



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 27, 2017 – 12:44 PM JST

PDB ID : 5XUA
Title : The ligand-free dimer of chemoreceptor MCP2201 ligand binding domain
Deposited on : 2017-06-23
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report.

This report is produced by the wwPDB biocuration pipeline after annotation of the structure.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

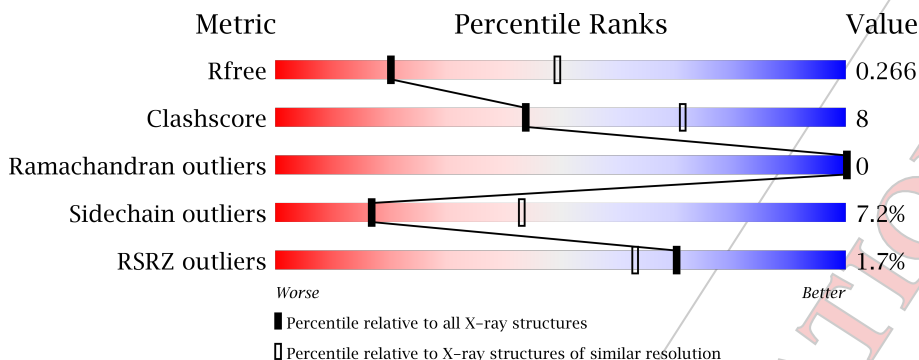
MolProbity : 4.02b-467
Xtrriage (Phenix) : 1.9-1692
EDS : rb-20029077
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20029077

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



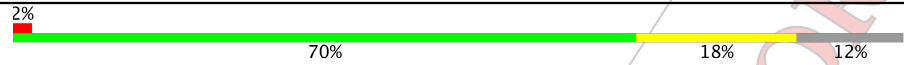


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	2583 (2.80-2.80)
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)
RSRZ outliers	101464	2610 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	154	 .% 68% 20% 5% 8%
1	B	154	 74% 14% • 11%
1	C	154	 68% 19% • 10%
1	D	154	 5% 73% 13% • 11%
1	E	154	 .% 68% 18% • 11%

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Mol	Chain	Length	Quality of chain
1	F	154	
1	G	154	
1	H	154	

CONFIDENTIAL VALIDATION REPORT

2 Entry composition i

There is only 1 type of molecule in this entry. The entry contains 8363 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Methyl-accepting chemotaxis sensory transducer.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	142	Total 1082	C 675	N 185	O 218	S 4	0	0	0
1	B	137	Total 1040	C 649	N 179	O 208	S 4	0	0	0
1	C	138	Total 1047	C 654	N 180	O 209	S 4	0	0	0
1	D	137	Total 1040	C 649	N 179	O 208	S 4	0	0	0
1	E	137	Total 1040	C 649	N 179	O 208	S 4	0	0	0
1	F	136	Total 1032	C 644	N 178	O 207	S 3	0	0	0
1	G	139	Total 1056	C 659	N 181	O 212	S 4	0	0	0
1	H	135	Total 1026	C 641	N 177	O 204	S 4	0	0	0

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	50	MET	-	expression tag	UNP D0IVL9
A	51	HIS	-	expression tag	UNP D0IVL9
A	52	HIS	-	expression tag	UNP D0IVL9
A	53	HIS	-	expression tag	UNP D0IVL9
A	54	HIS	-	expression tag	UNP D0IVL9
A	55	HIS	-	expression tag	UNP D0IVL9
A	56	HIS	-	expression tag	UNP D0IVL9
B	50	MET	-	expression tag	UNP D0IVL9
B	51	HIS	-	expression tag	UNP D0IVL9
B	52	HIS	-	expression tag	UNP D0IVL9
B	53	HIS	-	expression tag	UNP D0IVL9
B	54	HIS	-	expression tag	UNP D0IVL9
B	55	HIS	-	expression tag	UNP D0IVL9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	56	HIS	-	expression tag	UNP D0IVL9
C	50	MET	-	expression tag	UNP D0IVL9
C	51	HIS	-	expression tag	UNP D0IVL9
C	52	HIS	-	expression tag	UNP D0IVL9
C	53	HIS	-	expression tag	UNP D0IVL9
C	54	HIS	-	expression tag	UNP D0IVL9
C	55	HIS	-	expression tag	UNP D0IVL9
C	56	HIS	-	expression tag	UNP D0IVL9
D	50	MET	-	expression tag	UNP D0IVL9
D	51	HIS	-	expression tag	UNP D0IVL9
D	52	HIS	-	expression tag	UNP D0IVL9
D	53	HIS	-	expression tag	UNP D0IVL9
D	54	HIS	-	expression tag	UNP D0IVL9
D	55	HIS	-	expression tag	UNP D0IVL9
D	56	HIS	-	expression tag	UNP D0IVL9
E	50	MET	-	expression tag	UNP D0IVL9
E	51	HIS	-	expression tag	UNP D0IVL9
E	52	HIS	-	expression tag	UNP D0IVL9
E	53	HIS	-	expression tag	UNP D0IVL9
E	54	HIS	-	expression tag	UNP D0IVL9
E	55	HIS	-	expression tag	UNP D0IVL9
E	56	HIS	-	expression tag	UNP D0IVL9
F	50	MET	-	expression tag	UNP D0IVL9
F	51	HIS	-	expression tag	UNP D0IVL9
F	52	HIS	-	expression tag	UNP D0IVL9
F	53	HIS	-	expression tag	UNP D0IVL9
F	54	HIS	-	expression tag	UNP D0IVL9
F	55	HIS	-	expression tag	UNP D0IVL9
F	56	HIS	-	expression tag	UNP D0IVL9
G	50	MET	-	expression tag	UNP D0IVL9
G	51	HIS	-	expression tag	UNP D0IVL9
G	52	HIS	-	expression tag	UNP D0IVL9
G	53	HIS	-	expression tag	UNP D0IVL9
G	54	HIS	-	expression tag	UNP D0IVL9
G	55	HIS	-	expression tag	UNP D0IVL9
G	56	HIS	-	expression tag	UNP D0IVL9
H	50	MET	-	expression tag	UNP D0IVL9
H	51	HIS	-	expression tag	UNP D0IVL9
H	52	HIS	-	expression tag	UNP D0IVL9
H	53	HIS	-	expression tag	UNP D0IVL9
H	54	HIS	-	expression tag	UNP D0IVL9
H	55	HIS	-	expression tag	UNP D0IVL9

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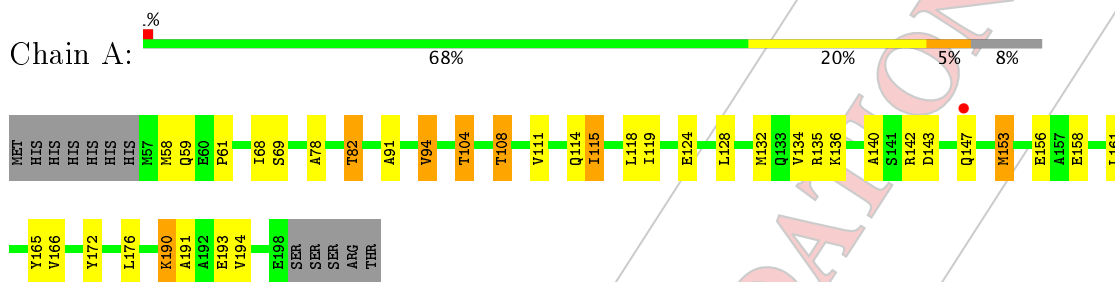
Chain	Residue	Modelled	Actual	Comment	Reference
H	56	HIS	-	expression tag	UNP D0IVL9

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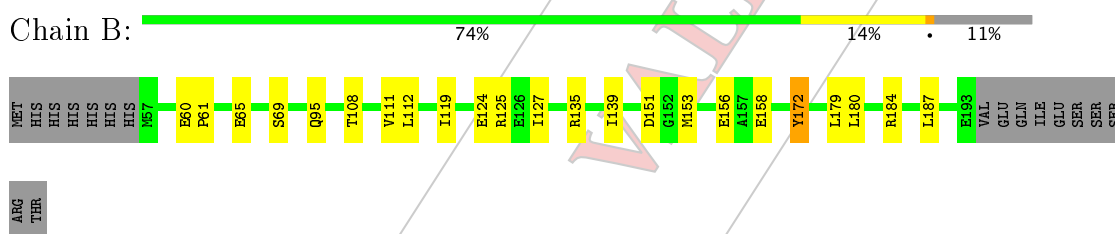
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

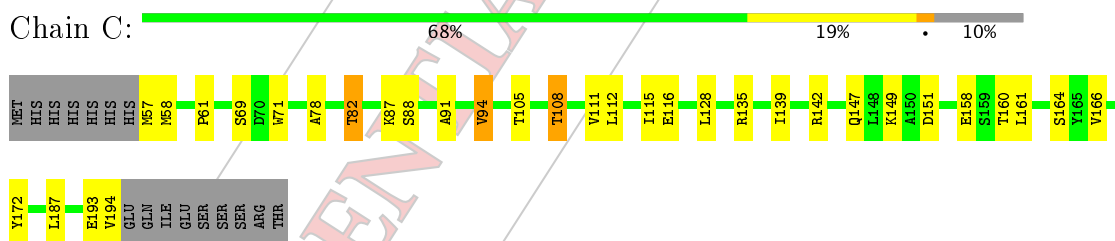
- Molecule 1: Methyl-accepting chemotaxis sensory transducer



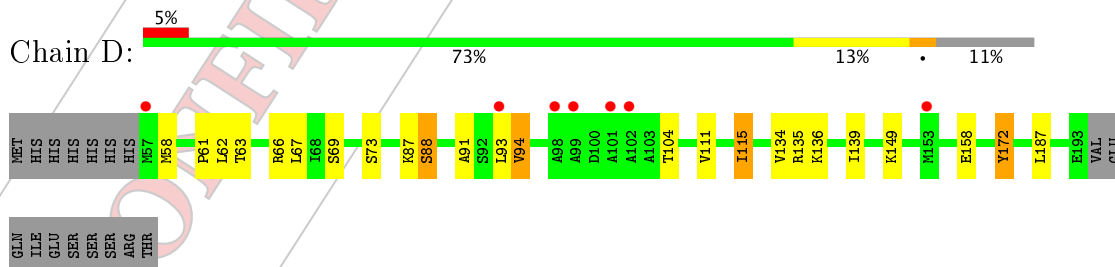
- Molecule 1: Methyl-accepting chemotaxis sensory transducer



- Molecule 1: Methyl-accepting chemotaxis sensory transducer

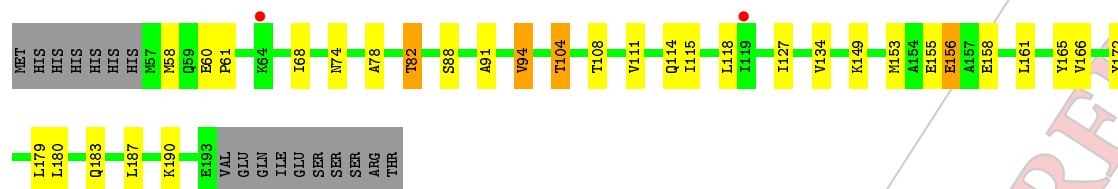


- Molecule 1: Methyl-accepting chemotaxis sensory transducer



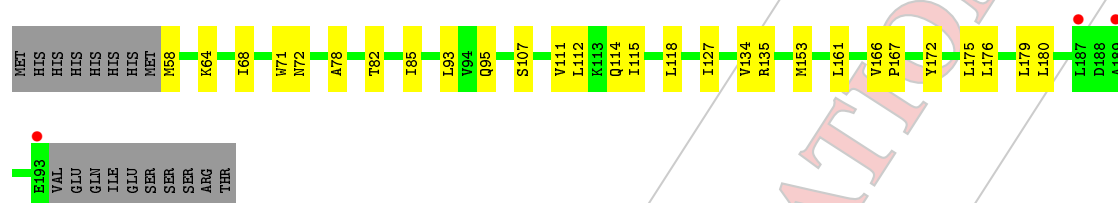
- Molecule 1: Methyl-accepting chemotaxis sensory transducer

Chain E: 



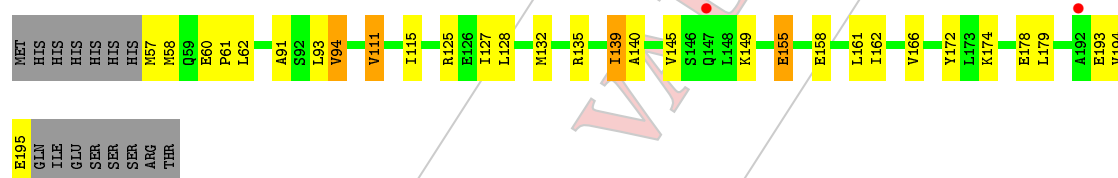
- Molecule 1: Methyl-accepting chemotaxis sensory transducer

Chain F: 



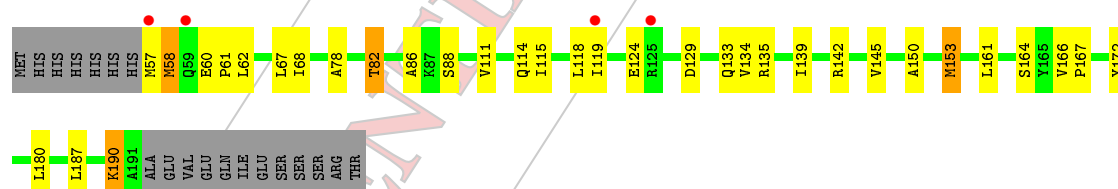
- Molecule 1: Methyl-accepting chemotaxis sensory transducer

Chain G: 



- Molecule 1: Methyl-accepting chemotaxis sensory transducer

Chain H: 



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.78Å 105.03Å 91.07Å 90.00° 90.95° 90.00°	Depositor
Resolution (Å)	53.20 – 2.80 55.13 – 2.80	Depositor EDS
% Data completeness (in resolution range)	95.1 (53.20-2.80) 94.8 (55.13-2.80)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 2.81Å)	Xtrriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.222 , 0.266 0.221 , 0.266	Depositor DCC
R_{free} test set	1448 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	55.6	Xtrriage
Anisotropy	0.320	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.024 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8363	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.40% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.27	0/1091	0.50	0/1476
1	B	0.26	0/1049	0.47	0/1419
1	C	0.26	0/1056	0.46	0/1429
1	D	0.26	0/1049	0.46	0/1419
1	E	0.27	0/1049	0.47	0/1419
1	F	0.26	0/1041	0.48	0/1409
1	G	0.28	0/1065	0.45	0/1441
1	H	0.26	0/1035	0.47	0/1400
All	All	0.26	0/8435	0.47	0/11412

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1082	0	1117	24	0
1	B	1040	0	1077	14	0
1	C	1047	0	1086	19	0
1	D	1040	0	1077	15	0
1	E	1040	0	1077	18	0
1	F	1032	0	1068	13	0
1	G	1056	0	1092	18	0
1	H	1026	0	1066	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	8363	0	8660	129	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (129) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:MET:SD	1:A:135:ARG:NH2	2.49	0.86
1:C:105:THR:HG21	1:H:150:ALA:HB1	1.66	0.78
1:B:61:PRO:HB2	1:B:187:LEU:HD13	1.72	0.72
1:E:78:ALA:O	1:E:82:THR:OG1	2.13	0.67
1:E:161:LEU:O	1:E:166:VAL:HG23	1.96	0.65
1:B:119:ILE:O	1:B:125:ARG:NH2	2.30	0.65
1:D:88:SER:O	1:D:149:LYS:NZ	2.28	0.64
1:E:61:PRO:HB2	1:E:187:LEU:HD23	1.80	0.64
1:D:61:PRO:HB2	1:D:187:LEU:HD13	1.79	0.63
1:F:166:VAL:HG22	1:F:167:PRO:HD3	1.79	0.63
1:C:88:SER:HB2	1:D:93:LEU:HD11	1.79	0.63
1:E:153:MET:HB3	1:E:156:GLU:HG3	1.81	0.62
1:G:161:LEU:O	1:G:166:VAL:HG23	1.99	0.62
1:C:161:LEU:O	1:C:166:VAL:HG23	1.98	0.62
1:C:87:LYS:NZ	1:C:158:GLU:OE1	2.33	0.60
1:B:151:ASP:OD2	1:F:135:ARG:NH2	2.33	0.59
1:F:71:TRP:HB2	1:F:111:VAL:HG21	1.84	0.59
1:A:161:LEU:O	1:A:166:VAL:HG23	2.04	0.58
1:F:161:LEU:O	1:F:166:VAL:HG13	2.04	0.57
1:C:78:ALA:O	1:C:82:THR:OG1	2.18	0.57
1:F:72:ASN:HA	1:F:176:LEU:HD13	1.87	0.57
1:H:166:VAL:HG22	1:H:167:PRO:HD3	1.86	0.57
1:C:116:GLU:HG2	1:C:128:LEU:HD21	1.88	0.56
1:G:194:VAL:HG21	1:H:57:MET:HG3	1.87	0.56
1:G:91:ALA:O	1:G:94:VAL:HG23	2.05	0.56
1:C:151:ASP:OD1	1:H:135:ARG:NH2	2.39	0.55
1:A:78:ALA:O	1:A:82:THR:OG1	2.22	0.55
1:F:127:ILE:HG21	1:F:179:LEU:HB2	1.89	0.55
1:B:184:ARG:HG2	1:B:184:ARG:HH11	1.71	0.55
1:B:60:GLU:HB3	1:B:61:PRO:HD3	1.89	0.54
1:G:93:LEU:HD11	1:H:88:SER:HB2	1.87	0.54
1:D:139:ILE:HG13	1:E:155:GLU:HB3	1.90	0.54
1:B:135:ARG:O	1:B:139:ILE:HG13	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:174:LYS:HE2	1:G:178:GLU:OE2	2.08	0.54
1:G:60:GLU:HB3	1:G:61:PRO:HD3	1.90	0.54
1:D:91:ALA:O	1:D:94:VAL:HG23	2.07	0.54
1:E:104:THR:O	1:E:108:THR:HG23	2.08	0.54
1:E:88:SER:O	1:E:149:LYS:NZ	2.39	0.53
1:D:88:SER:OG	1:D:93:LEU:HD12	2.08	0.53
1:G:125:ARG:HH11	1:G:125:ARG:HG3	1.73	0.53
1:H:114:GLN:O	1:H:118:LEU:HD13	2.08	0.53
1:A:136:LYS:HD2	1:G:158:GLU:HG3	1.90	0.52
1:D:135:ARG:NH1	1:D:172:TYR:OH	2.42	0.52
1:H:82:THR:HG21	1:H:142:ARG:HG3	1.92	0.52
1:B:135:ARG:NH1	1:B:172:TYR:OH	2.41	0.52
1:D:62:LEU:HD21	1:D:66:ARG:CZ	2.39	0.52
1:D:67:LEU:HB3	1:D:115:ILE:CD1	2.40	0.52
1:H:62:LEU:HA	1:H:187:LEU:HD11	1.92	0.52
1:E:61:PRO:HG2	1:E:190:LYS:HD3	1.93	0.51
1:H:68:ILE:HG22	1:H:115:ILE:HD13	1.91	0.51
1:C:61:PRO:HB2	1:C:187:LEU:HD13	1.92	0.51
1:D:136:LYS:HD2	1:E:158:GLU:HG3	1.93	0.51
1:H:58:MET:O	1:H:61:PRO:HD2	2.11	0.50
1:G:128:LEU:O	1:G:132:MET:HG2	2.12	0.50
1:A:104:THR:O	1:A:108:THR:OG1	2.26	0.50
1:E:91:ALA:O	1:E:94:VAL:HG23	2.11	0.50
1:H:111:VAL:O	1:H:115:ILE:HG13	2.12	0.50
1:H:161:LEU:O	1:H:166:VAL:HG13	2.12	0.50
1:A:58:MET:O	1:A:61:PRO:HD2	2.12	0.49
1:F:68:ILE:HD11	1:F:180:LEU:HA	1.93	0.49
1:A:153:MET:N	1:A:153:MET:SD	2.86	0.48
1:C:88:SER:O	1:C:149:LYS:NZ	2.38	0.48
1:E:74:ASN:HB3	1:E:108:THR:HG22	1.95	0.48
1:G:135:ARG:O	1:G:139:ILE:HG23	2.14	0.48
1:G:139:ILE:HG13	1:G:140:ALA:N	2.28	0.47
1:C:139:ILE:HG23	1:C:142:ARG:NH2	2.29	0.47
1:B:180:LEU:O	1:B:184:ARG:HG3	2.14	0.47
1:C:135:ARG:O	1:C:139:ILE:HG13	2.14	0.47
1:C:160:THR:O	1:C:164:SER:OG	2.29	0.46
1:F:85:ILE:HA	1:F:93:LEU:HD12	1.97	0.46
1:E:114:GLN:O	1:E:118:LEU:HD13	2.16	0.46
1:C:147:GLN:HG2	1:H:139:ILE:HD12	1.98	0.46
1:D:63:THR:O	1:D:67:LEU:HD13	2.15	0.46
1:C:91:ALA:O	1:C:94:VAL:HG23	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:68:ILE:HD12	1:E:180:LEU:HA	1.97	0.46
1:A:91:ALA:O	1:A:94:VAL:HG23	2.15	0.45
1:B:153:MET:HB3	1:B:156:GLU:OE1	2.15	0.45
1:B:65:GLU:OE2	1:B:184:ARG:NH1	2.49	0.45
1:A:58:MET:HB2	1:A:190:LYS:HE2	1.97	0.45
1:H:60:GLU:HB2	1:H:61:PRO:HD3	1.99	0.45
1:H:153:MET:H	1:H:153:MET:HG2	1.39	0.45
1:G:125:ARG:NH1	1:G:125:ARG:HG3	2.32	0.45
1:D:67:LEU:HB3	1:D:115:ILE:HD11	1.99	0.45
1:A:69:SER:OG	1:B:69:SER:OG	2.34	0.45
1:C:69:SER:OG	1:D:69:SER:OG	2.35	0.45
1:F:114:GLN:O	1:F:118:LEU:HD13	2.17	0.45
1:G:127:ILE:HG21	1:G:179:LEU:HB2	1.98	0.45
1:A:82:THR:HB	1:A:165:TYR:OH	2.17	0.44
1:G:111:VAL:O	1:G:115:ILE:HG13	2.18	0.44
1:H:129:ASP:O	1:H:133:GLN:HG2	2.17	0.44
1:B:119:ILE:HG23	1:B:124:GLU:HB2	2.00	0.44
1:E:60:GLU:HB2	1:E:61:PRO:HD3	1.99	0.44
1:C:58:MET:HG3	1:C:58:MET:O	2.18	0.43
1:A:128:LEU:O	1:A:132:MET:HG2	2.17	0.43
1:D:135:ARG:O	1:D:139:ILE:HG12	2.19	0.43
1:H:68:ILE:HD11	1:H:180:LEU:HA	2.01	0.43
1:A:68:ILE:HG13	1:A:115:ILE:HG21	2.00	0.43
1:C:108:THR:HA	1:C:111:VAL:HG12	2.01	0.43
1:A:69:SER:HG	1:B:69:SER:HG	1.66	0.43
1:A:143:ASP:O	1:A:147:GLN:HG2	2.18	0.42
1:C:71:TRP:CZ2	1:C:112:LEU:HD13	2.54	0.42
1:H:119:ILE:HG23	1:H:124:GLU:HB3	2.01	0.42
1:G:57:MET:HB2	1:H:190:LYS:NZ	2.34	0.42
1:A:191:ALA:O	1:A:194:VAL:HG12	2.20	0.42
1:H:164:SER:C	1:H:167:PRO:HD2	2.40	0.42
1:E:156:GLU:HG2	1:E:156:GLU:H	1.53	0.42
1:A:82:THR:CG2	1:A:142:ARG:HG3	2.50	0.42
1:D:87:LYS:HE2	1:D:158:GLU:OE2	2.20	0.42
1:A:114:GLN:O	1:A:118:LEU:HD13	2.20	0.41
1:E:82:THR:HB	1:E:165:TYR:OH	2.20	0.41
1:H:166:VAL:CG2	1:H:167:PRO:HD3	2.49	0.41
1:A:190:LYS:HA	1:A:193:GLU:HG2	2.01	0.41
1:A:156:GLU:N	1:A:156:GLU:OE1	2.41	0.41
1:B:127:ILE:HG21	1:B:179:LEU:HB2	2.01	0.41
1:F:134:VAL:HG11	1:F:175:LEU:HD12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:64:LYS:O	1:F:68:ILE:HG23	2.20	0.41
1:A:58:MET:CB	1:A:190:LYS:HE2	2.50	0.41
1:C:111:VAL:O	1:C:115:ILE:HG22	2.21	0.41
1:E:127:ILE:HG21	1:E:179:LEU:HB2	2.02	0.41
1:F:107:SER:O	1:F:111:VAL:HG13	2.21	0.41
1:F:78:ALA:O	1:F:82:THR:HG23	2.20	0.41
1:G:155:GLU:H	1:G:155:GLU:HG3	1.55	0.41
1:A:140:ALA:HB1	1:G:162:ILE:HD12	2.03	0.41
1:H:78:ALA:O	1:H:82:THR:OG1	2.36	0.41
1:H:86:ALA:HB1	1:H:145:VAL:HG11	2.03	0.40
1:A:119:ILE:HG23	1:A:124:GLU:CB	2.51	0.40
1:A:176:LEU:HA	1:A:176:LEU:HD23	1.95	0.40
1:E:68:ILE:HD11	1:E:183:GLN:HB3	2.04	0.40
1:G:145:VAL:O	1:G:149:LYS:HG3	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	140/154 (91%)	137 (98%)	3 (2%)	0	100	100
1	B	135/154 (88%)	129 (96%)	6 (4%)	0	100	100
1	C	136/154 (88%)	135 (99%)	1 (1%)	0	100	100
1	D	135/154 (88%)	129 (96%)	6 (4%)	0	100	100
1	E	135/154 (88%)	131 (97%)	4 (3%)	0	100	100
1	F	134/154 (87%)	130 (97%)	4 (3%)	0	100	100
1	G	137/154 (89%)	133 (97%)	4 (3%)	0	100	100
1	H	133/154 (86%)	128 (96%)	5 (4%)	0	100	100
All	All	1085/1232 (88%)	1052 (97%)	33 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	117/129 (91%)	105 (90%)	12 (10%)	8	24
1	B	112/129 (87%)	106 (95%)	6 (5%)	26	58
1	C	113/129 (88%)	106 (94%)	7 (6%)	21	52
1	D	112/129 (87%)	103 (92%)	9 (8%)	14	38
1	E	112/129 (87%)	103 (92%)	9 (8%)	14	38
1	F	111/129 (86%)	105 (95%)	6 (5%)	26	58
1	G	114/129 (88%)	105 (92%)	9 (8%)	14	38
1	H	111/129 (86%)	104 (94%)	7 (6%)	21	51
All	All	902/1032 (87%)	837 (93%)	65 (7%)	17	43

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	GLN
1	A	82	THR
1	A	94	VAL
1	A	104	THR
1	A	108	THR
1	A	111	VAL
1	A	115	ILE
1	A	134	VAL
1	A	153	MET
1	A	158	GLU
1	A	172	TYR
1	A	190	LYS
1	B	95	GLN
1	B	108	THR
1	B	111	VAL
1	B	112	LEU

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Mol	Chain	Res	Type
1	B	158	GLU
1	B	172	TYR
1	C	57	MET
1	C	82	THR
1	C	94	VAL
1	C	108	THR
1	C	172	TYR
1	C	193	GLU
1	C	194	VAL
1	D	58	MET
1	D	73	SER
1	D	88	SER
1	D	94	VAL
1	D	104	THR
1	D	111	VAL
1	D	115	ILE
1	D	134	VAL
1	D	172	TYR
1	E	58	MET
1	E	82	THR
1	E	94	VAL
1	E	104	THR
1	E	111	VAL
1	E	115	ILE
1	E	134	VAL
1	E	156	GLU
1	E	172	TYR
1	F	58	MET
1	F	95	GLN
1	F	112	LEU
1	F	115	ILE
1	F	153	MET
1	F	172	TYR
1	G	58	MET
1	G	62	LEU
1	G	94	VAL
1	G	111	VAL
1	G	139	ILE
1	G	155	GLU
1	G	172	TYR
1	G	193	GLU
1	G	195	GLU

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Mol	Chain	Res	Type
1	H	58	MET
1	H	67	LEU
1	H	82	THR
1	H	134	VAL
1	H	153	MET
1	H	172	TYR
1	H	190	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	95	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data i

6.1 Protein, DNA and RNA chains i

In the following table, the column labelled '#RSRZ > 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q < 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å ²)	Q < 0.9
1	A	142/154 (92%)	-0.08	1 (0%) 87 83	35, 58, 94, 121	0
1	B	137/154 (88%)	-0.13	0 100 100	36, 54, 97, 116	0
1	C	138/154 (89%)	-0.12	0 100 100	41, 56, 91, 102	0
1	D	137/154 (88%)	0.17	7 (5%) 29 19	41, 63, 108, 129	0
1	E	137/154 (88%)	0.04	2 (1%) 74 67	48, 72, 108, 119	0
1	F	136/154 (88%)	0.17	3 (2%) 62 52	41, 61, 102, 133	0
1	G	139/154 (90%)	0.13	2 (1%) 75 69	44, 63, 95, 124	0
1	H	135/154 (87%)	0.24	4 (2%) 51 39	44, 67, 115, 126	0
All	All	1101/1232 (89%)	0.05	19 (1%) 70 63	35, 62, 103, 133	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	119	ILE	4.3
1	H	57	MET	4.0
1	D	102	ALA	3.8
1	G	192	ALA	3.7
1	D	99	ALA	3.2
1	A	147	GLN	2.8
1	D	98	ALA	2.6
1	H	59	GLN	2.5
1	H	125	ARG	2.4
1	D	57	MET	2.4
1	F	187	LEU	2.3
1	F	193	GLU	2.3
1	E	119	ILE	2.2
1	D	93	LEU	2.1
1	D	153	MET	2.1
1	F	189	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	101	ALA	2.1
1	G	147	GLN	2.0
1	E	64	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.

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