SUPPLEMENTARY MATERIALS

Materials and Methods

Sample Preparation. H-*BEA was prepared from commercial CP814E NH_4-*BEA zeolite by calcination in static air at 550°C, as in refs. (6) and (7). H-CHA zeolite used for the preparation of the Fe-CHA samples was prepared following the CBV720 recipe with N,N,N-trimethyl-1-admantylammonium cations ($TMAda^+$) template from ref. (35). A molar batch composition of 1 Si : 0.067 Al : 0.22 $TMAda^+$: 0.13 Na^+ : 0.35 OH^- : 24.5 H_2O was targeted using zeolite Y (Zeolyst international CBV720) as the Si and Al source. 28.69 g of aqueous N,N,N-trimethyladamantylammonium hydroxide (TMAdaOH) solution (25 wt %, Sachem), 5.29 g of NaOH solution (15 wt%, from >98 wt% NaOH pellets, Sigma Aldrich) and 40.58 g of deionized water (18.2 MΩ cm) was mixed in a 125 ml Teflon lined stainless steel autoclave (Parr Instruments) and homogenized. 11.25 g of the zeolite Y precursor was then added and the mixture was stirred for 2 hours at ambient conditions. The autoclave was then sealed off and oven-heated at 160 °C for 4 days under static conditions.

Fe-CHA and Fe-*BEA materials were prepared by a strategy analogous to the samples in refs (6), (7), and (36): Fe was introduced into dried H-CHA and H-*BEA by diffusion impregnation in a solution of $^{57}Fe(acac)_3$ in toluene (25 mL/g zeolite). The concentration of $^{57}Fe(acac)_3$ in toluene is approximately 0.01 M. All samples were calcined in air with a heating ramp of 2 °C/min to 550 °C for 30 hours to remove organic

material. Each sample is then introduced into a quartz U-tube/flow cell and heated at 10°C/min in a dried 30 ml/min flow of helium (99.999% purity, O_2 filtered with an Agilent 17970 system, H_2O filtered with an Agilent 17971 system) to 900°C where it is kept for one hour. The Fe-*BEA and Fe-CHA materials studied here were compositionally similar with Si/Al ~ 10 and ~0.30 wt.% Fe. Sample specific Si/Al and iron loadings are given in the main manuscript.

N_2O Activation and CH_4/H_2 Reactions. ~0.5g of Fe zeolite is introduced into a quartz U-tube/flow cell. After heating to 900°C in helium (*vide supra*), a 35% N_2O (99%, IJsfabriek Strombeek) in He flow is passed over the sample (20 min at 160°C for Fe-*BEA, 25 min at 180°C for Fe-CHA to achieve maximal α-Fe(IV)=O and minimal spectator). The reactor is flushed in helium at room temperature and a 20 mL/min flow of CH_4 (99.995%, Air Liquide) or H_2 (99.99%, Air Liquide) is passed through the flow cell at room temperature for 5 minutes, followed by another flush in helium. For the isotope labelled experiments CH_4 and H_2 were replaced by 99% labeled $^{13}CH_4$ (99%, 99 atom% ^{13}C, Sigma Aldrich) and D_2 (99%, 99.9 atom% D, Sigma Aldrich) respectively.

Catalyst Cycling Tests. A known amount of Fe-*BEA or Fe-CHA was put into a quartz reactor. The sample was pretreated in a dried helium flow at 900°C, then a 35% N_2O/He atmosphere (large molar excess of N_2O over iron) was inserted into the reactor at room temperature. The reactor was heated in the oven for 30 minutes at 180 °C for Fe-CHA and 160 °C for Fe-*BEA to ensure full oxidation. Then the N_2O/He was removed at room

temperature by a vacuum, and replaced by a 1.5 bar $^{13}CH_4$ atmosphere (99% ^{13}C labeled). After 15 minutes all $^{13}CH_4$ was removed by a vacuum. For the single cycle experiments, methanol was extracted at this point in a 35 ml/min He stream that passed through a bubbler with water at room temperature. The outflow was monitored by in line mass spectrometry using an Omnistar Pfeiffer Vacuum GSD 30102 quadrupole mass spectrometer. For a second cycle experiment, the $^{13}CH_4$ atmosphere was removed by vacuum and replaced by 35%N_2O/He. The reactor was heated in the oven for 15 minutes at 180 °C for Fe-CHA and 160 °C for Fe-*BEA. Then the N_2O/He was removed at room temperature by a vacuum, and replaced by a $^{12}CH_4$ atmosphere. After 15 minutes methanol extraction was started. Different N_2O activation temperatures and durations were tested to optimize for α-Fe(II) conversion and minimize overoxidation of first cycle oxidation products. Quantification of ^{13}C and ^{12}C methanol products was calibrated by a known amount of methanol adsorbed onto Fe-CHA and Fe-*BEA and then extracted using the same procedure as for the reactive cycles. Exemplary mass spectrometry data are shown in Fig. S6. Mass spectrometry during CH_4 reaction did not detect the gas phase release of methanol or other oxidation products. Products other than methanol were also not detected in quantifiable amounts during the steaming procedure. Nevertheless, the methanol yields (both ^{13}C and ^{12}C) are ~30% below what is expected based on Mössbauer spectroscopy. We explain this by the sampling for Mössbauer from the part of the zeolite bed that is least exposed to contaminants. The Mössbauer data therefore does not quantitatively but only relatively represent the whole zeolite bed used for the reactivity tests.

Steam assisted extractions were similar for H and Na exchanged zeolites and for the CHA and BEA topologies (Fig. S7). Although the time required for methanol recovery differs markedly, the final outcome is similar for the purpose of this work.

DR-UV-vis Spectroscopy. Diffuse reflectance spectroscopy in the UV–vis–NIR energy range was performed with a Varian Cary 5000 UV-vis-NIR spectrophotometer at room temperature against a Halon white reflectance standard in the 4000-40000 cm^{-1} energy range. All treatments before UV-vis-NIR spectroscopic measurements were performed in a quartz U-tube/flow cell, equipped with a window for *in situ* measurements.

Mössbauer Spectroscopy. ^{57}Fe Mössbauer spectra were recorded with a See Co. W302 resonant gamma ray spectrometer in horizontal geometry at room temperature with zero external field using a 1.85 GBq source (Be window, Rh matrix). Data were collected from samples enriched with 100% ^{57}Fe. Isomer shifts are given relative to α-iron foil at room temperature. Spectra were collected with 1024 points and summed up to 512 points before analyzing. Spectra were fit to Lorentzian doublets and hyperfine-split multiplets using the Vinda software package for Microsoft Excel. Hyperfine features were fit using a spin Hamiltonian model (SPINHAM), using a vanishingly small external field. A slowly relaxing S=5/2 model was required to fit the data. First, the α-Fe(III)-OH component of H_2-reacted BEA was considered. Its isomer shift, quadrupole splitting, and hyperfine coupling tensor (constrained to be isotropic) were determined from the position of peripheral spectral lines, where there are no contributions from overlapping Lorentzian

components (i.e. above 4 mm/s and below -2 mm/s). The *g* tensor was constrained to be isotropic, with *g=2.*0023. The axial and rhombic zero-field splitting parameters *D* and *E* (respectively) were then floated to properly model the intensity distribution among these peripheral spectral lines, accounting for the 6K sample temperature. For the CH_4 reacted sample, the features of α-Fe(III)-OH were subtracted from the experimental data to facilitate fitting of the remaining α-Fe(III)-OCH_3 component.

Resonance Raman Spectroscopy. Resonance Raman (rR) spectra were recorded using a Spex 1877 CP triple monochromator with 1200, 1800, and 2400 grooves/mm holographic spectrograph gratings and an Andor Newton CCD cooled to –80 °C. Excitation was provided by either a Coherent I90C–K Kr+ ion laser (λ_{ex} = 407) or an Innova Sabre 25/7 Ar+ CW ion laser (λex = 458, 502 nm). The spectral resolution was ~2 cm^{-1}. Spectra were recorded at room temperature, using 20 mW of laser power at the sample. Baseline spectra were collected from samples of α-Fe(IV)=O in BEA and CHA lattice.

Spectroscopic Studies of Fe-CHA. Mössbauer data from the H_2 and CH_4 reactions in CHA are presented in Fig. S1. For the H_2 reaction, α-Fe(IV)=O (initially 84%) is converted into a new majority Fe(III) component (51% of Fe) with hyperfine structure. Based on its resemblance to α-Fe(III)-OH in BEA, as well as rR data (see Fig. S5), we assign this as the CHA α-Fe(III)-OH site (solid black trace, 61% of α-Fe(IV)=O). Some α-Fe(II) is also regenerated (21% of α-Fe(IV)=O). The reaction of α-Fe(IV)=O with H_2 therefore gives a

similar distribution of Fe species in CHA and BEA. For the CH_4 reaction, however, only ~60% of the α-Fe(IV)=O is converted into α-Fe(III)-OH/α-Fe(III)-OCH_3. Unlike BEA, a significant portion (~40% of α-Fe(II)) is regenerated.

DR-UV-vis data from H_2- and CH_4-reacted CHA are presented in Fig. S5. As with BEA, reacting CHA with H_2 at 300 K (Fig. S5d) results in decay of the α-Fe(IV)=O band (17,500 cm^{-1}, gray trace), and leaves a rising slope with somewhat diminished Abs intensity from 20,000-30,000 cm^{-1}. For the CH_4 reaction (Fig. S5e), there is additional Abs intensity at ~22,000 cm^{-1}. Finally, a low energy band appears at 15,000 cm^{-1} that is not present in the BEA sample. However, heating the sample to 450 K results in loss of this absorption feature, but with minimal change to its Mössbauer spectrum, suggesting this is due to a small quantity (<5%) of a strongly absorbing species. Tuning a laser (ν_{ex}=21,800 cm^{-1}) into the Abs shoulder of the H_2-reacted sample resonance enhances a single vibration at 730 cm^{-1} (Fig. S5f, blue trace). We assign this to the α-Fe(III)-OH species. No α-Fe(III)-OH/α-Fe(III)-OCH_3 features are resolved in the CH_4 reacted CHA sample (Fig. S5f, red trace). This is a key difference between CHA and BEA – consistent with Mössbauer data, which show a greatly diminished contribution from Fe(III) species in CH_4-reacted CHA.

Computational Details. Spin-unrestricted DFT calculations were performed using Gaussian 09.(37) Cluster models were generated from crystallographic coordinates of BEA polymorph A (28) and CHA,(38) and dangling Si-O groups were replaced with capping hydrides at 1.42 Å (see supplementary tables 1-17 for coordinates). The B3LYP

functional was used for all calculations. For geometry optimizations and frequency calculations, the 6-311G* basis set was used for Fe, all atoms directly coordinated to iron, and all atoms of the CH_4 substrate. The 6-31G* basis set was used on all other atoms. For geometry optimizations, the six T-sites of the 6-membered ring hosting the iron center, and all atoms of the substrate were allowed to relax. All other atoms were constrained to their crystallographic positions. The stability of all optimized geometries were confirmed by frequency calculations. Guess transition state structures were located through linear transits of the Fe-$OH\cdots CH_3$ distance, and then geometry optimized. Transition state structures were validated through frequency calculations, which showed a single imaginary frequency corresponding to motion along the reaction coordinate. Single point calculations were performed on optimized structures using the 6-311G+* basis set on all atoms, and including the D3 dispersion correction of Grimme in the DFT functional.(39) All single point energies were corrected for thermal contributions to enthalpy and Gibbs free energy at 300K. Mössbauer isomer shifts and quadrupole splittings were calculated according to previously reported methods.(6)

Supplementary Figures

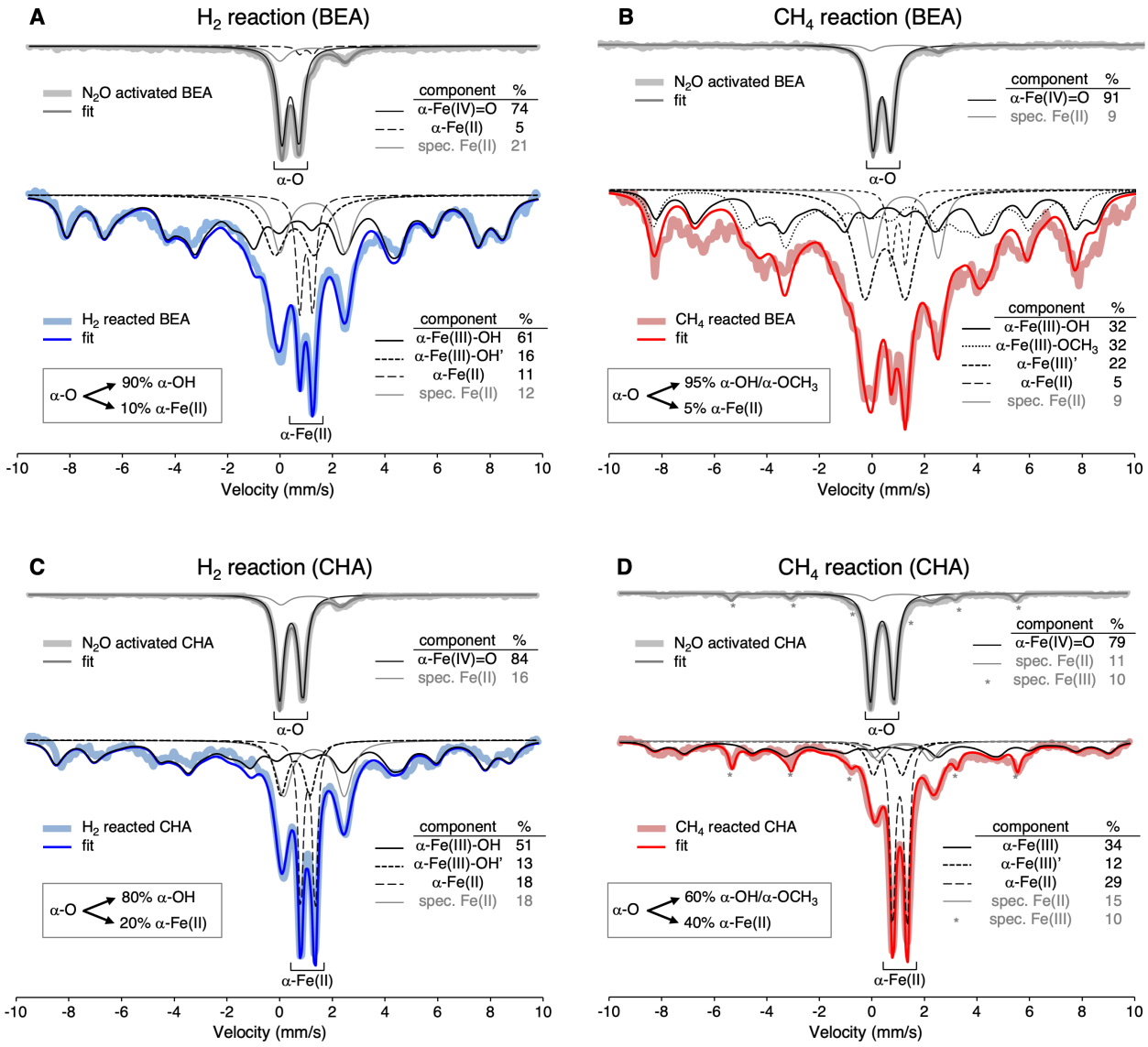
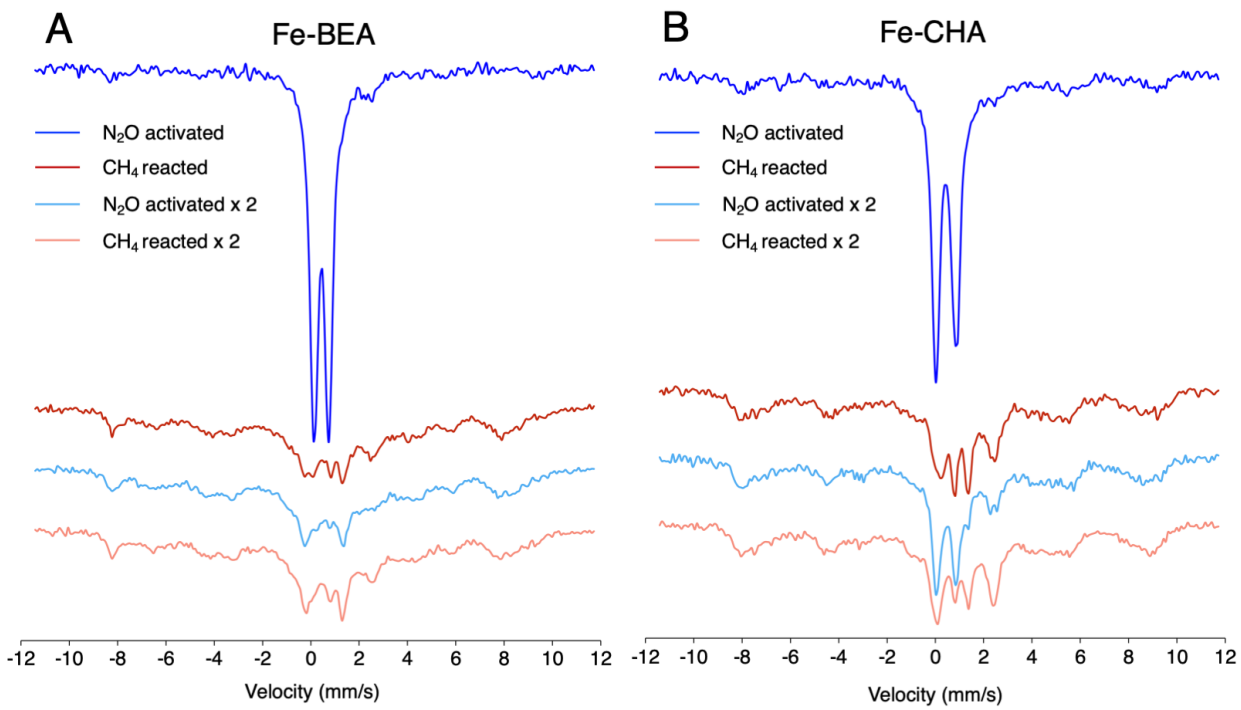


Figure S1. ^{57}Fe Mössbauer spectra of H_2- and CH_4-reacted Fe-CHA and Fe-BEA. Data were collected at 6K. Components of each fit are indicated in the figure legends. Data from the corresponding α-Fe(IV)=O samples are presented above the spectrum of each reacted material (**A** – H_2-reacted BEA, **B** – CH_4-reacted BEA, **C** – H_2-reacted CHA, **D** – CH_4-reacted CHA).



	component	BEA (%)	CHA (%)
N_2O activated	α-Fe(IV)=O	90	74
	α-Fe(III)	0	0
	α-Fe(II)	0	0
CH_4 reacted	α-Fe(IV)=O	0	0
	α-Fe(III)	85	46
	α-Fe(II)	5	28
N_2O activated x 2	α-Fe(IV)=O	0	22
	α-Fe(III)	90	52
	α-Fe(II)	0	0
CH_4 reacted x 2	α-Fe(IV)=O	0	0
	α-Fe(III)	90	54
	α-Fe(II)	0	20

Figure S2. Normalized 6K Mössbauer spectra from **(A)** Fe-*BEA and **(B)** Fe-CHA collected over the course of catalyst cycling. Note regeneration of α-Fe(IV)=O upon reoxidation in Fe-CHA, but not Fe-*BEA (see pale blue traces). **(C)** Active site speciation over the course of catalyst cycling, as quantified from Mössbauer data in panels A and B. These values are associated with errors of ±5%.

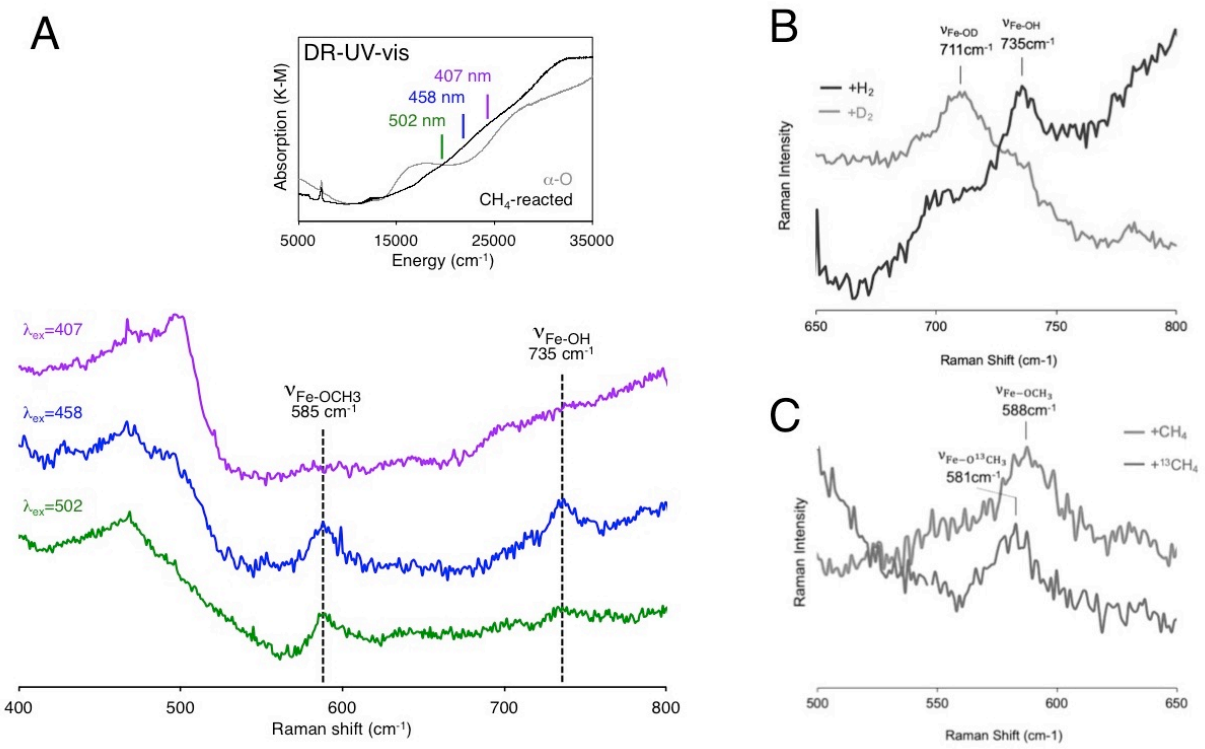
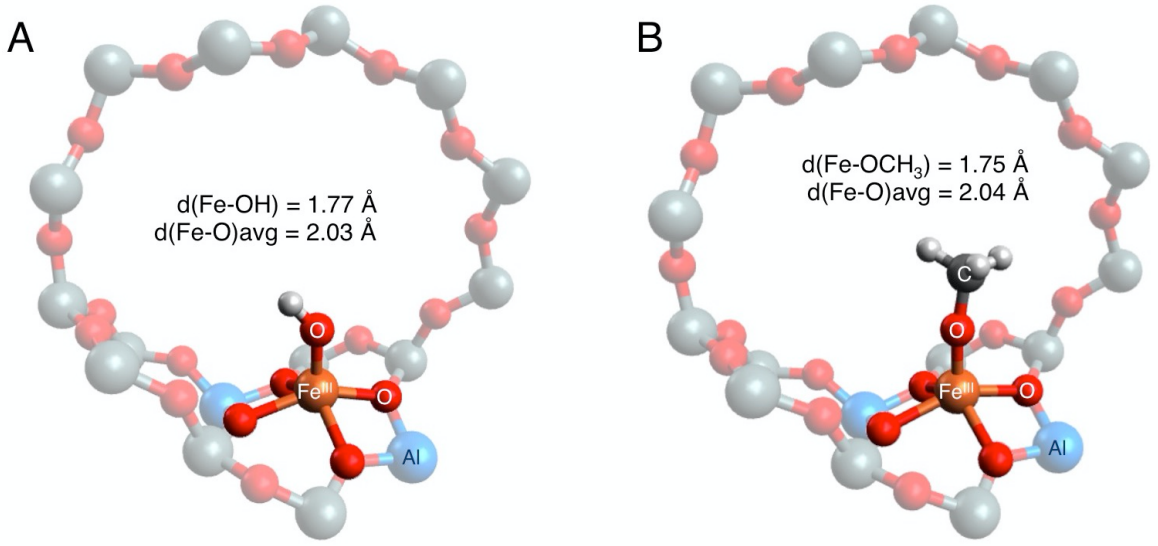


Figure S3. (A) Resonance Raman profile of the characteristic vibrations of α-OH and α-OCH_3 in CH_4-reacted Fe-BEA. The intensities of both features are maximized using an excitation wavelength of 458 nm, indicating these Raman features are in resonance with electronic absorption bands within the 407-502 nm interval. To the right, isotopically perturbed resonance Raman data (λ_{ex}=458 nm) are presented, focusing on the H/D sensitivity of the Fe-OH stretch of α-OH in H_2/D_2-reacted Fe-BEA (**A**), and the $^{12}C/^{13}C$ sensitivity of the Fe-OCH_3 stretch of α-OCH_3 in $^{12}CH_4/^{13}CH_4$-reacted Fe-BEA (**B**). The hydrogen and methane reactions were all performed at 300K.



	Mössbauer parameters		$\nu_{Fe\text{-}OH}$ (ΔH/D)
α-OH	**IS (mm/s)**	**QS (mm/s)**	**(cm^{-1})**
EXP	0.5±0.1	-1.6±0.1	735 (24)
DFT	0.41	-1.41	764 (32)

	Mössbauer parameters		$\nu_{Fe\text{-}OH}$ ($\Delta^{12}C/^{13}C$)
α-OCH$_3$	**IS (mm/s)**	**QS (mm/s)**	**(cm^{-1})**
EXP	0.5±0.1	-1.6±0.1	585 (7)
DFT	0.47	-1.34	577 (6)

Figure S4. S=5/2 DFT models of α-OH **(A)** and α-OCH$_3$ sites **(B)** in Fe-BEA, including a comparison of their experimental and calculated spectroscopic features in Mössbauer (IS = isomer shift, QS = quadrupole splitting) and resonance Raman spectroscopy. Atoms have been omitted for clarity.

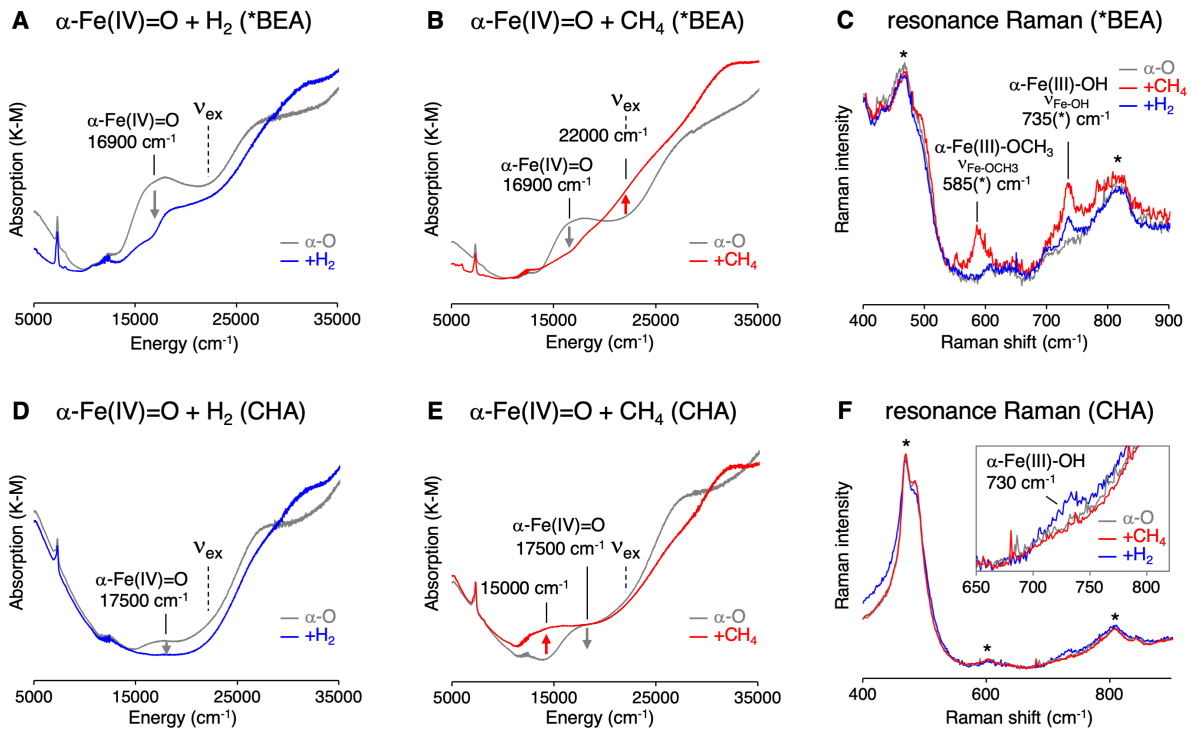


Figure S5. Comparison of DR-UV-vis and rR data (λ_{ex}=458 nm) from H_2/CH_4 reacted BEA (top) and CHA (bottom). All reactions were performed at 300 K. While rR features of the cage escape products α-Fe(III)-OH and α-Fe(III)-OCH_3 are observed in CH_4-reacted BEA (panel C), these species are not detected in CH_4-reacted CHA. Cage escape products are only observed for the smaller substrate H_2 (panel F). Starred features are non-resonant Raman vibrations of the zeolite lattice.

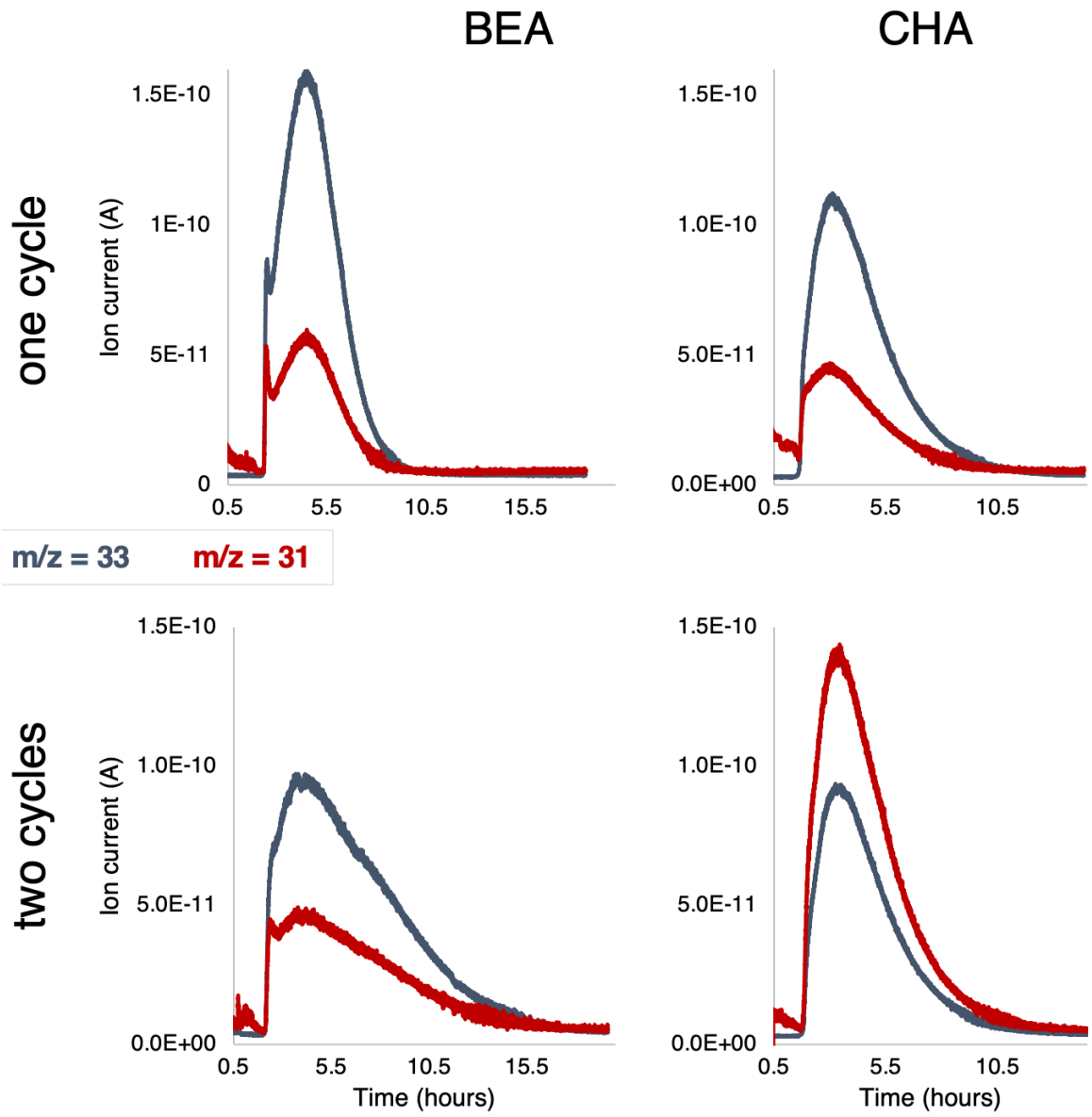


Figure S6. Selected mass spectrometry data used to quantify the steam desorbed $^{13}CH_3OH$ and $^{12}CH_3OH$. m/z=33 was used to quantify $^{13}CH_3OH$. The portion of the m/z=31 signal attributed to $^{13}CH_3OH$ was subtracted from the total signal to quantify $^{12}CH_3OH$.

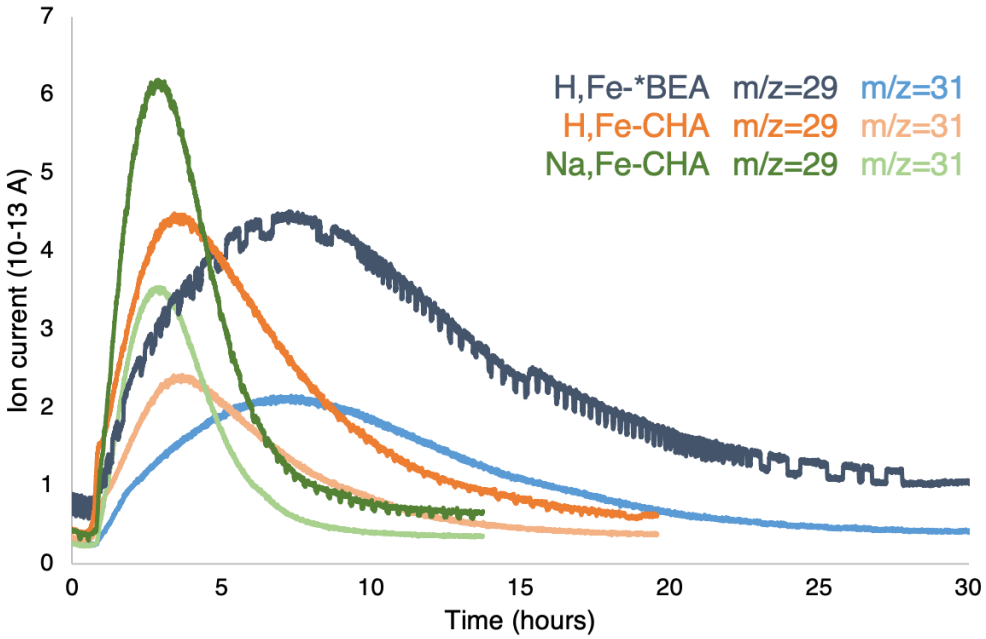


Figure S7. Methanol desorption by a flow of 30 mL/min He saturated with H_2O at room temperature compared for the *BEA vs. CHA topologies and for Na^+ vs H^+ exchanged CHA. Methanol was produced by a reaction of α-O with methane at room temperature.

Supplementary Tables

Table S1 – Coordinates for DFT Model of α-OH···CH_3 Complex (BEA)

element	constrained during optimization?	X	Y	Z
Fe	N	-0.88317713	0.2078925	-0.30780668
O	N	-1.0400598	-0.44184015	1.60450329
O	N	-2.37687441	0.25268539	-1.68318867
O	N	-1.93334343	1.88022371	0.21946534
O	N	-0.84106095	-1.84277563	-0.51339426
O	N	0.75101568	0.71266097	-0.72332064
H	N	1.65230655	0.48644183	-0.43617148
C	N	3.21305879	-0.24578148	0.88023069
H	N	3.13304826	0.65959271	1.46811711
H	N	2.65000536	-1.12165503	1.17712814
H	N	4.03337029	-0.34787691	0.18140014
O	Y	4.87661135	-4.44257331	0.30163545
O	Y	1.27825899	5.77135874	0.64899151
O	Y	6.93006414	-2.79880345	0.15694788
O	Y	6.04196136	4.42530407	0.03027365
O	Y	3.71060695	5.56193854	-0.34005062
O	Y	7.15463405	2.24695049	-0.92669782
O	Y	7.53770541	-0.32180801	-0.50751246
Si	Y	3.32319985	-4.86647248	0.25003353
Si	Y	-0.31975761	5.56885478	0.68192352
Si	Y	5.98373678	-3.79995768	-0.67419749
Si	Y	7.30423865	3.43501077	0.1455109
Si	Y	2.38113727	6.46630161	-0.29602195
Si	Y	7.58505566	0.85172085	-1.60825877
Si	Y	8.11782004	-1.71583324	0.05034822
Si	Y	5.203328	5.41191462	-0.93008396
O	Y	2.57809006	-4.33650714	1.57477511
Si	Y	1.20514996	-4.24318502	2.41140011
H	Y	0.72498441	0.22216818	6.48083059
O	Y	2.63218595	-4.22169495	-1.05123203
H	Y	-0.86362965	5.85729448	-0.59768194
H	Y	7.37260057	2.89143817	1.455571
H	Y	2.99572467	-3.2083064	-3.28320658
H	Y	8.68835307	-1.51596069	1.3352399
H	Y	8.48622223	4.17916665	-0.11053247
H	Y	5.3278827	-3.0748261	-1.70396832
H	Y	1.8632252	6.57139177	-1.61401213

H	Y	6.76597082	-4.83091757	-1.25868583
H	Y	2.7000874	7.75475836	0.20850288
H	Y	9.11683636	-2.20172141	-0.83411566
H	Y	8.90445043	0.97342957	-2.11893871
H	Y	5.15098332	4.8672679	-2.24043
H	Y	6.68520139	0.54664373	-2.66351801
H	Y	5.82687181	6.68694693	-0.97356146
Si	Y	2.02476692	-3.89163911	-2.50423579
O	N	-0.0069791	0.02802504	4.05761858
O	N	-2.84524028	-0.01262259	-4.33628098
H	Y	0.81254472	-1.82960421	5.40400638
H	Y	3.22908059	-6.28209649	0.18842809
H	Y	-4.21283594	4.21609455	-4.11866462
H	Y	1.50081345	-4.13868655	3.79622425
H	Y	-2.61179643	5.31420437	-2.85074504
H	Y	-1.23152514	-0.95737013	6.07445221
H	Y	-5.25708837	-0.3330108	-5.11446855
H	Y	-3.94173059	-1.20644052	5.29758307
H	Y	-6.61689752	-1.55802481	-2.95555335
H	Y	1.66300639	-5.09938037	-3.15758428
H	Y	-0.88434447	6.44110335	1.64980538
O	N	0.78236699	-2.83971473	-2.37288099
O	N	-0.550889	3.99749498	1.05214991
O	N	-0.93329768	1.99026324	2.65089395
O	N	-1.50007281	-1.82494797	-3.06699486
O	N	-2.63925551	0.20452577	3.67742707
O	N	-4.12166885	-1.59107063	-2.61180931
H	Y	-5.1244251	0.2113667	3.88579629
H	Y	-5.73872247	-3.08613362	-1.44034967
O	Y	-4.15970964	-1.85095029	2.91533171
H	Y	-4.80957296	5.15973295	2.809539
O	Y	-5.82528084	-0.70576163	-0.76809888
H	Y	-2.08634563	-6.59763456	-1.71124899
O	N	-4.70897598	1.68675949	-0.4926341
H	Y	-1.75220556	-5.4596028	0.27397778
O	N	-2.24472889	-3.16715043	1.60942187
H	Y	-5.02787416	4.0636545	0.78497986
O	N	-3.00128128	3.00435962	-2.32174871
H	Y	0.43034821	-5.41048192	2.17983227
O	N	0.53591137	-2.83783967	1.94158606
H	Y	-4.71832796	4.91717127	-1.96710834
H	Y	-6.15080497	0.70660986	1.23869362
H	Y	-3.6842613	-5.0945297	-0.96606054

H	Y	-4.62910758	-3.55124479	1.17848673
H	Y	-5.30802453	2.89734291	2.77622182
H	Y	-7.16041854	1.3818953	-0.73616746
O	N	-1.45110768	-4.16661595	-1.81069443
H	Y	-3.77103589	-4.28139872	3.20479645
O	N	-2.96527547	3.60161245	2.07384681
Si	Y	0.05712837	-0.63874217	5.57023429
Si	Y	-4.01817743	0.26379672	-5.46850182
Si	N	-1.17249889	0.42884422	3.02084121
Si	N	-2.74459067	-0.78988245	-2.93008988
Si	Y	-4.01005516	-0.6628577	3.98750595
Si	Y	-5.62852639	-1.76838418	-1.95787707
Al	N	-3.13641855	1.81341683	-1.13164839
Al	N	-0.88302309	-2.21720719	1.25138142
Si	Y	-6.00443507	0.78041402	-0.17182709
Si	Y	-2.29210997	-5.36549847	-1.03598896
Si	Y	-3.72923507	-3.24613987	2.23378664
Si	Y	-4.58582977	3.93410279	2.12818444
Si	N	-1.63279803	2.89898729	1.49627949
Si	N	-0.75964507	-2.71296991	-1.92531086
Si	Y	-3.69769648	4.40016792	-2.80825149
H	Y	-4.16301828	1.67580886	-5.50964355
H	Y	-3.61030004	-0.20749562	-6.74441532

Table S2 – Coordinates for DFT Model of α-Fe(II) (BEA)

element	constrained during optimization?	X	Y	Z
Fe	N	-1.12883888	0.10320115	-0.21656273
O	N	-0.95669034	-0.46232606	1.67806171
O	N	-2.38006755	0.28004486	-1.74770877
O	N	-1.87230785	1.92004893	0.18314947
O	N	-0.72559543	-1.84267712	-0.49485712
O	Y	4.94558774	-4.43295001	0.3203729
O	Y	1.34801784	5.78217523	0.63992601
O	Y	6.99958707	-2.78975642	0.17683658
O	Y	6.11283216	4.43421721	0.03404538
O	Y	3.78244321	5.57045776	-0.34349751
O	Y	7.22722346	2.25382251	-0.91625629
O	Y	7.60901638	-0.31416147	-0.49119815
Si	Y	3.39223206	-4.85672332	0.26628911

Si	Y	-0.25008882	5.57997464	0.66984097
Si	Y	6.05488926	-3.79240787	-0.65435803
Si	Y	7.37471357	3.44397058	0.15392925
Si	Y	2.45301683	6.47509881	-0.30409032
Si	Y	7.65889313	0.85719167	-1.59414893
Si	Y	8.18772954	-1.70716996	0.07064424
Si	Y	5.27640016	5.41905698	-0.9300427
O	Y	2.64438202	-4.32405323	1.58840047
Si	Y	1.26966375	-4.22887428	2.42188553
H	Y	0.78145736	0.24455456	6.48148439
O	Y	2.704092	-4.21439524	-1.03772303
H	Y	-0.7911904	5.8659687	-0.61148552
H	Y	7.44019418	2.90296724	1.46520016
H	Y	3.07254949	-3.20546414	-3.27090739
H	Y	8.75554567	-1.50485016	1.35635652
H	Y	8.55735233	4.18745002	-0.10105195
H	Y	5.40134515	-3.06921069	-1.68695405
H	Y	1.93793779	6.57766828	-1.62338951
H	Y	6.83821786	-4.82462959	-1.23514268
H	Y	2.77107989	7.76449924	0.19857752
H	Y	9.18856242	-2.19494208	-0.81072449
H	Y	8.9793949	0.97770386	-2.10224611
H	Y	5.22677505	4.87183918	-2.23942399
H	Y	6.76125093	0.55016805	-2.6507268
H	Y	5.90022515	6.69391106	-0.97469686
Si	Y	2.09982881	-3.88711335	-2.49267638
O	N	0.06284683	0.10200961	4.0902585
O	N	-2.74467337	-0.04704576	-4.38211772
H	Y	0.87101099	-1.80934599	5.40888958
H	Y	3.29801954	-6.27244536	0.20725802
H	Y	-4.13311217	4.21832532	-4.13638188
H	Y	1.56236278	-4.12169756	3.80713269
H	Y	-2.53461622	5.31869375	-2.86721211
H	Y	-1.1743574	-0.93549513	6.07324611
H	Y	-5.17590393	-0.33258811	-5.12546182
H	Y	-3.8829308	-1.18570719	5.29107787
H	Y	-6.54051057	-1.55314484	-2.96704774
H	Y	1.73928256	-5.09609267	-3.14441136
H	Y	-0.81661821	6.45420081	1.63480057
O	N	0.8586752	-2.82983618	-2.37915712
O	N	-0.4764637	4.00451433	1.03000645
O	N	-0.88270978	2.00231555	2.62543751
O	N	-1.42427486	-1.81339575	-3.04023414

O	N	-2.58215821	0.23701016	3.69301448
O	N	-4.05908255	-1.6092377	-2.650114
H	Y	-5.062398	0.22949071	3.87398002
H	Y	-5.66580375	-3.07839904	-1.44696866
O	Y	-4.09591657	-1.83487079	2.90963848
H	Y	-4.74451271	5.17568389	2.7886595
O	Y	-5.75344016	-0.69669442	-0.77959309
H	Y	-2.01337515	-6.59095677	-1.70314456
O	N	-4.64751236	1.71438462	-0.50138873
H	Y	-1.68330782	-5.44906365	0.2805449
O	N	-2.18808524	-3.15412089	1.58526527
H	Y	-4.95864676	4.07565573	0.76579898
O	N	-2.91396278	3.01176417	-2.36281586
H	Y	0.49517078	-5.39650471	2.1909498
O	N	0.60414085	-2.82216998	1.98361439
H	Y	-4.64309021	4.92370948	-1.98729486
H	Y	-6.0830435	0.71966884	1.22371881
H	Y	-3.61265717	-5.08615517	-0.96433631
H	Y	-4.56185308	-3.5385155	1.17514492
H	Y	-5.24322414	2.91330441	2.75872925
H	Y	-7.08833766	1.39120941	-0.75461774
O	N	-1.37448548	-4.16001761	-1.80276527
H	Y	-3.70822355	-4.26480121	3.20471866
O	N	-2.89643547	3.63232163	2.05714146
Si	Y	0.1154217	-0.61804675	5.57115903
Si	Y	-3.936152	0.26334971	-5.47803034
Si	N	-1.10792523	0.45214418	3.04892912
Si	N	-2.6884559	-0.79859526	-2.962902
Si	Y	-3.9483813	-0.64468503	3.97979586
Si	Y	-5.55430415	-1.76170588	-1.96685747
Al	N	-3.08729187	1.84254904	-1.16560541
Al	N	-0.80153797	-2.21992559	1.28072638
Si	Y	-5.93366132	0.79069511	-0.18661926
Si	Y	-2.22040236	-5.35747053	-1.03075807
Si	Y	-3.66418475	-3.23146486	2.23176469
Si	Y	-4.51950263	3.94869207	2.1102012
Si	N	-1.57262015	2.91939965	1.47686939
Si	N	-0.67700176	-2.71078598	-1.9112984
Si	Y	-3.62074899	4.40491193	-2.82525113
H	Y	-4.08069853	1.67529606	-5.52224516
H	Y	-3.52562135	-0.21050944	-6.75213146

Table S3 – Coordinates for DFT Model of α-O (BEA)

element	constrained during optimization?	X	Y	Z
Fe	N	-0.91671491	0.17603945	-0.28283458
O	N	-1.00631261	-0.44267911	1.6278311
O	N	-2.3666186	0.23921784	-1.67295926
O	N	-1.86718256	1.85971359	0.21941703
O	N	-0.77442756	-1.80711442	-0.50276371
O	N	0.49253807	0.69473342	-0.80528232
O	Y	4.92431485	-4.45223802	0.32936005
O	Y	1.3408724	5.76730625	0.66532759
O	Y	6.98076239	-2.81167126	0.19090354
O	Y	6.10433749	4.41381069	0.06125501
O	Y	3.77590579	5.55407073	-0.3162491
O	Y	7.21653202	2.23368335	-0.89222701
O	Y	7.59430578	-0.33565401	-0.47178009
Si	Y	3.37040808	-4.87372562	0.27302532
Si	Y	-0.2575663	5.56732953	0.69330328
Si	Y	6.03542263	-3.81137918	-0.64310176
Si	Y	7.36470794	3.42155221	0.18039299
Si	Y	2.44772037	6.46052151	-0.27631782
Si	Y	7.64685949	0.83775567	-1.57241807
Si	Y	8.17053022	-1.73056113	0.08790428
Si	Y	5.27019146	5.40169192	-0.90170181
O	Y	2.62208921	-4.34251704	1.59545854
Si	Y	1.24664945	-4.24693411	2.42782899
H	Y	0.76108173	0.21923792	6.4956374
O	Y	2.6843789	-4.22788486	-1.03037751
H	Y	-0.79706307	5.8565604	-0.58789692
H	Y	7.42820736	2.8779253	1.49067615
H	Y	3.05634753	-3.21516319	-3.26126442
H	Y	8.7374367	-1.53152817	1.37453077
H	Y	8.54863091	4.16385109	-0.07205527
H	Y	5.38385997	-3.08526615	-1.67490273
H	Y	1.9340129	6.56635738	-1.59589864
H	Y	6.81783284	-4.84358316	-1.22515437
H	Y	2.7671369	7.74848819	0.22912946
H	Y	9.17149143	-2.21804387	-0.79347852
H	Y	8.96800176	0.95738286	-2.07905875
H	Y	5.2210107	4.85707511	-2.21218437
H	Y	6.7497663	0.53404203	-2.63041758
H	Y	5.89585615	6.67574742	-0.94331467

Si	Y	2.08196929	-3.89694836	-2.4852466
O	N	0.03480766	0.03465076	4.07609949
O	N	-2.77407588	-0.00927613	-4.33499126
H	Y	0.84872652	-1.83270399	5.41915837
H	Y	3.2742533	-6.28917773	0.21116755
H	Y	-4.13800483	4.22046345	-4.11907612
H	Y	1.53831031	-4.142948	3.81358004
H	Y	-2.5391601	5.31611279	-2.8463067
H	Y	-1.19602386	-0.95726345	6.08332217
H	Y	-5.18632007	-0.32705993	-5.11790854
H	Y	-3.90421286	-1.20213952	5.2981505
H	Y	-6.5546532	-1.54984045	-2.96311442
H	Y	1.72029122	-5.10415226	-3.13966181
H	Y	-0.82373802	6.44047945	1.65950386
O	N	0.83979004	-2.84147363	-2.36105765
O	N	-0.49486008	3.99077958	1.04467572
O	N	-0.88378877	1.99159792	2.65998817
O	N	-1.43836757	-1.81712505	-3.05548652
O	N	-2.59942071	0.21608372	3.68679659
O	N	-4.06796105	-1.60401688	-2.63107386
H	Y	-5.08038665	0.21746078	3.88270789
H	Y	-5.68352521	-3.07927502	-1.44517909
O	Y	-4.1159038	-1.84641691	2.91532497
H	Y	-4.75451019	5.16529369	2.80724728
O	Y	-5.76844832	-0.69873186	-0.77333166
H	Y	-2.03579429	-6.59649371	-1.70475958
O	N	-4.65243556	1.70205473	-0.48579717
H	Y	-1.70596774	-5.45890122	0.28143107
O	N	-2.2109951	-3.15414251	1.58595664
H	Y	-4.96833368	4.06947714	0.78206813
O	N	-2.92765629	3.00833732	-2.31948764
H	Y	0.47082236	-5.41312167	2.19395151
O	N	0.56720348	-2.84587645	1.97128005
H	Y	-4.64900616	4.9224383	-1.9691203
H	Y	-6.09790428	0.71419138	1.23244788
H	Y	-3.63365116	-5.09088331	-0.96451501
H	Y	-4.58266724	-3.54601413	1.17707446
H	Y	-5.25639211	2.9036907	2.77246794
H	Y	-7.10035951	1.39097968	-0.74555779
O	N	-1.40187944	-4.14287071	-1.77417077
H	Y	-3.73190616	-4.27741161	3.20603238
O	N	-2.91309445	3.59164476	2.05443655
Si	Y	0.09461771	-0.6406435	5.58302212

Si	Y	-3.94539248	0.2678036	-5.4681443
Si	N	-1.12707209	0.43391823	3.04182609
Si	N	-2.69429523	-0.79728156	-2.93632018
Si	Y	-3.9677119	-0.65841852	3.98784093
Si	Y	-5.56958208	-1.76176385	-1.96242874
Al	N	-3.08557794	1.8198421	-1.13503489
Al	N	-0.83106996	-2.214622	1.27119934
Si	Y	-5.94707744	0.78775679	-0.17744032
Si	Y	-2.24173864	-5.36394586	-1.03022906
Si	Y	-3.68553141	-3.24229701	2.23497095
Si	Y	-4.53048906	3.93929759	2.12659366
Si	N	-1.57039102	2.89516745	1.49936172
Si	N	-0.69530589	-2.70118635	-1.91430514
Si	Y	-3.6266018	4.40364775	-2.80698067
H	Y	-4.08791191	1.68003488	-5.50979579
H	Y	-3.53435056	-0.20417456	-6.74280757

Table S4 – Coordinates for DFT Model of α-OH (BEA)

element	constrained during optimization?	X	Y	Z
Fe	N	-0.84673097	0.20228632	-0.30335058
O	N	-0.99908053	-0.4371057	1.61115894
O	N	-2.34219503	0.24565126	-1.66945215
O	N	-1.87509261	1.87832175	0.22198741
O	N	-0.78982757	-1.83840538	-0.50020977
O	N	0.75391944	0.69714192	-0.86389625
H	N	1.62063145	0.30266454	-0.98581128
O	Y	4.91208494	-4.45498761	0.33673611
O	Y	1.33281809	5.76581926	0.66872748
O	Y	6.96819533	-2.81525728	0.19921073
O	Y	6.09614477	4.41008206	0.06841789
O	Y	3.7685036	5.55132229	-0.31099179
O	Y	7.20809402	2.22933761	-0.88394042
O	Y	7.58445419	-0.34017503	-0.46307391
Si	Y	3.35784438	-4.87629668	0.27931788
Si	Y	-0.26565355	5.56645528	0.69562488
Si	Y	6.02420609	-3.81524059	-0.63503391
Si	Y	7.35598594	3.41728092	0.18863644
Si	Y	2.44069235	6.45834559	-0.27219654
Si	Y	7.63831714	0.83313276	-1.56362275

Si	Y	8.15972029	-1.73529494	0.09782321
Si	Y	5.2631625	5.39820459	-0.89529839
O	Y	2.6087764	-4.34495309	1.60110359
Si	Y	1.23313356	-4.24823247	2.43266567
H	Y	0.74617268	0.21872493	6.4993101
O	Y	2.67322782	-4.23025274	-1.02468351
H	Y	-0.80411637	5.85578241	-0.58613775
H	Y	7.41825645	2.87379856	1.49903674
H	Y	3.04727877	-3.21797295	-3.25541199
H	Y	8.72551232	-1.53612378	1.38448658
H	Y	8.54042549	4.15902337	-0.06301558
H	Y	5.37373104	-3.08894162	-1.66746883
H	Y	1.92802673	6.56425348	-1.59216619
H	Y	6.80652136	-4.84787423	-1.21609666
H	Y	2.76029263	7.74625617	0.23333591
H	Y	9.16106854	-2.2232972	-0.78303637
H	Y	8.95989318	0.95210804	-2.06928191
H	Y	5.21472924	4.85343799	-2.20574425
H	Y	6.74188902	0.52967562	-2.62226042
H	Y	5.88942157	6.6719782	-0.93660641
Si	Y	2.07203909	-3.89932126	-2.48000431
O	N	0.01969752	0.02586564	4.07447106
O	N	-2.78308135	-0.01081375	-4.33014602
H	Y	0.83372217	-1.83341178	5.42317048
H	Y	3.26129072	-6.29164115	0.21765393
H	Y	-4.14314917	4.22069113	-4.11960376
H	Y	1.52351255	-4.14407387	3.81840425
H	Y	-2.54477643	5.31584783	-2.84577908
H	Y	-1.21114092	-0.95696235	6.08566415
H	Y	-5.19269866	-0.32646414	-5.1186344
H	Y	-3.91886512	-1.20073953	5.29849918
H	Y	-6.56318761	-1.54838179	-2.96473637
H	Y	1.71030646	-5.10635063	-3.13457987
H	Y	-0.83226423	6.44003203	1.66114582
O	N	0.83608434	-2.8348601	-2.3521504
O	N	-0.50217742	3.99515225	1.06128894
O	N	-0.89486463	1.99065543	2.66259589
O	N	-1.4452312	-1.82318199	-3.05391205
O	N	-2.608863	0.2071799	3.67674359
O	N	-4.07134424	-1.59459267	-2.61830977
H	Y	-5.09329881	0.21919578	3.88195363
H	Y	-5.69388047	-3.07799643	-1.44595201
O	Y	-4.12891878	-1.84513323	2.91542377

H	Y	-4.7644209	5.16673254	2.80610548
O	Y	-5.77842685	-0.69735751	-0.77442547
H	Y	-2.04754521	-6.59685054	-1.7023138
O	N	-4.6582699	1.69559204	-0.49115668
H	Y	-1.71869341	-5.45914924	0.28399011
O	N	-2.21766884	-3.15906153	1.60415505
H	Y	-4.97719853	4.07073885	0.78091327
O	N	-2.94114345	3.00552284	-2.31876304
H	Y	0.45661454	-5.41402926	2.1981137
O	N	0.55852339	-2.8461634	1.96834617
H	Y	-4.65544797	4.92313546	-1.97011659
H	Y	-6.10859636	0.71604906	1.23088735
H	Y	-3.64526306	-5.09040451	-0.96347727
H	Y	-4.59520382	-3.54485288	1.17719573
H	Y	-5.26728714	2.90534315	2.77126219
H	Y	-7.10926885	1.39302477	-0.74793448
O	N	-1.40221806	-4.1661173	-1.79712555
H	Y	-3.7463547	-4.27638762	3.20689152
O	N	-2.92171567	3.60099292	2.07049065
Si	Y	0.08007292	-0.64105714	5.58641198
Si	Y	-3.9511515	0.26778246	-5.46799418
Si	N	-1.13649622	0.42960376	3.03134226
Si	N	-2.69403362	-0.79132786	-2.92524023
Si	Y	-3.98112716	-0.65728318	3.98814575
Si	Y	-5.57887893	-1.76051173	-1.9632497
Al	N	-3.0847148	1.81762632	-1.12715349
Al	N	-0.84579878	-2.21498176	1.26809327
Si	Y	-5.95635911	0.78927257	-0.1790747
Si	Y	-2.25341202	-5.36422199	-1.02804471
Si	Y	-3.69876662	-3.24150891	2.23594663
Si	Y	-4.54045264	3.94049848	2.12579388
Si	N	-1.58506172	2.89898877	1.50433197
Si	N	-0.7102836	-2.71373844	-1.9107719
Si	Y	-3.63247027	4.40391989	-2.80719698
H	Y	-4.09310483	1.68008184	-5.50997434
H	Y	-3.5394727	-0.20454351	-6.74231845

Table S5 – Coordinates for DFT Model of α-OCH_3 (BEA)

element	constrained during optimization?	X	Y	Z
Fe	N	-0.8592125	0.19339361	-0.29282275
O	N	-1.02665169	-0.45587964	1.62262601
O	N	-2.35571336	0.2527359	-1.67079623
O	N	-1.91023793	1.87569731	0.24556535
O	N	-0.81393149	-1.85066386	-0.50270537
O	N	0.71212542	0.80126276	-0.75806112
C	N	1.89271544	1.55277248	-0.81872581
H	N	1.74691806	2.52455531	-0.33619774
H	N	2.6979614	1.01736257	-0.30582234
H	N	2.17471855	1.7069376	-1.86479575
O	Y	4.85187037	-4.51956881	0.34093809
O	Y	1.3514797	5.72804685	0.69735929
O	Y	6.92206372	-2.8955436	0.21267284
O	Y	6.10526905	4.3372355	0.10183303
O	Y	3.78727529	5.49765916	-0.27844718
O	Y	7.20206881	2.15082576	-0.85509165
O	Y	7.55764422	-0.42273467	-0.44116205
Si	Y	3.29471279	-4.92811915	0.27929173
Si	Y	-0.24859466	5.54113351	0.72057867
Si	Y	5.97076934	-3.88503545	-0.6267278
Si	Y	7.35714039	3.3343232	0.22137885
Si	Y	2.4664686	6.4148542	-0.23935959
Si	Y	7.62273429	0.75339088	-1.53822606
Si	Y	8.12092499	-1.8239426	0.11595495
Si	Y	5.28180717	5.33471813	-0.86041981
O	Y	2.54728138	-4.39463255	1.60131111
Si	Y	1.17048091	-4.2901793	2.43036983
H	Y	0.71077156	0.16801925	6.50989423
O	Y	2.61745815	-4.27289233	-1.02401209
H	Y	-0.78232624	5.83845894	-0.56121217
H	Y	7.41270855	2.78639604	1.53023099
H	Y	3.00362914	-3.25679959	-3.25095003
H	Y	8.68603628	-1.63330044	1.40464048
H	Y	8.5477755	4.06760739	-0.02579177
H	Y	5.32791694	-3.15060941	-1.65810941
H	Y	1.95714154	6.52873717	-1.55996463
H	Y	6.74622627	-4.92193406	-1.20973864
H	Y	2.7950958	7.69869287	0.27068905
H	Y	9.1201861	-2.3170676	-0.76421824

H	Y	8.94614981	0.86363792	-2.04104525
H	Y	5.2316311	4.7943255	-2.17261107
H	Y	6.72598603	0.46012097	-2.59945992
H	Y	5.91800672	6.60371031	-0.89666762
Si	Y	2.02164933	-3.93282448	-2.47951977
O	N	-0.01244916	-0.00124912	4.08818702
O	N	-2.79913337	0.00034769	-4.32823746
H	Y	0.7844363	-1.881438	5.42765304
H	Y	3.18711889	-6.34255139	0.21304979
H	Y	-4.12721258	4.24009962	-4.10591085
H	Y	1.45930153	-4.19260402	3.81714853
H	Y	-2.52281019	5.31888833	-2.8257601
H	Y	-1.2548174	-0.99117248	6.08900077
H	Y	-5.21013176	-0.29575632	-5.12077167
H	Y	-3.96283659	-1.21153437	5.296014
H	Y	-6.59415231	-1.51350784	-2.97315548
H	Y	1.65181063	-5.13510043	-3.1383686
H	Y	-0.81015792	6.41606766	1.68780397
O	N	0.78674115	-2.87036429	-2.36727561
O	N	-0.48934716	3.96395776	1.07105754
O	N	-0.91171541	1.9704297	2.67848771
O	N	-1.47970367	-1.82545486	-3.05355578
O	N	-2.64000709	0.19495069	3.68379611
O	N	-4.10425229	-1.57395828	-2.62142031
H	Y	-5.12357331	0.22177439	3.88164412
H	Y	-5.73961837	-3.05443606	-1.45741569
O	Y	-4.17348831	-1.84713566	2.91076536
H	Y	-4.75429692	5.16986076	2.82150042
O	Y	-5.80678218	-0.67523329	-0.77884292
H	Y	-2.1201833	-6.60071787	-1.71768811
O	N	-4.67586401	1.71339136	-0.49699566
H	Y	-1.78628746	-5.47162894	0.27266537
O	N	-2.27264688	-3.16338923	1.59324518
H	Y	-4.97171774	4.08172158	0.79257373
O	N	-2.92960176	3.0157564	-2.3025749
H	Y	0.38572539	-5.44926628	2.19102076
O	N	0.50160657	-2.88324973	1.97645762
H	Y	-4.63812438	4.93999698	-1.95525839
H	Y	-6.12995195	0.73452714	1.23018536
H	Y	-3.70758706	-5.08418654	-0.97725187
H	Y	-4.64956709	-3.53779944	1.1663401
H	Y	-5.27459957	2.9125618	2.7787948
H	Y	-7.12159339	1.42524684	-0.74845621

O	N	-1.46330123	-4.16926521	-1.80046721
H	Y	-3.81022177	-4.28203142	3.1953827
O	N	-2.9152999	3.61843152	2.09414905
Si	Y	0.03970907	-0.68370607	5.59304082
Si	Y	-3.96345094	0.28990607	-5.46604215
Si	N	-1.16428047	0.40638095	3.04051947
Si	N	-2.71804872	-0.78319476	-2.92303524
Si	Y	-4.01839862	-0.66351522	3.98715087
Si	Y	-5.61351908	-1.73641399	-1.97052749
Al	N	-3.09519279	1.81956129	-1.11913249
Al	N	-0.8878703	-2.23253576	1.2626404
Si	Y	-5.97478897	0.81095351	-0.17910129
Si	Y	-2.31779361	-5.3685133	-1.04011574
Si	Y	-3.75277689	-3.24453857	2.22748192
Si	Y	-4.53855345	3.94405763	2.13788652
Si	N	-1.59719311	2.88557752	1.52237701
Si	N	-0.75100454	-2.72558608	-1.91204167
Si	Y	-3.61780121	4.41528547	-2.79198805
H	Y	-4.09426421	1.70338198	-5.50390997
H	Y	-3.55291993	-0.18172244	-6.74096699

Table S6 – Coordinates for DFT Model of C-O Recombination TS (BEA)

element	constrained during optimization?	X	Y	Z
Fe	N	-0.88618006	0.20844731	-0.29585694
O	N	-1.05297281	-0.43907767	1.62043272
O	N	-2.37610343	0.24640971	-1.67837952
O	N	-1.94760098	1.8832949	0.2292295
O	N	-0.8105164	-1.83401901	-0.50705038
O	N	0.7422407	0.80886032	-0.67061895
H	N	1.28635252	1.5784887	-0.48836409
C	N	3.10023815	-0.46753597	-0.06997848
H	N	3.70511097	0.33058103	0.3393772
H	N	2.59619764	-1.15529723	0.59363855
H	N	3.08430711	-0.63820456	-1.13550146
O	Y	4.88581737	-4.4356887	0.33753518
O	Y	1.2693725	5.77257892	0.66306481
O	Y	6.93713558	-2.78864727	0.19988652
O	Y	6.03771593	4.43390052	0.06442378
O	Y	3.70600621	5.56644199	-0.31606467

O	Y	7.15778324	2.25668951	-0.88650291
O	Y	7.54337856	-0.31113549	-0.46392635
Si	Y	3.33332698	-4.86217118	0.28001801
Si	Y	-0.32842242	5.56748379	0.68973809
Si	Y	5.99578969	-3.79194864	-0.63432077
Si	Y	7.30114196	3.44576333	0.18543635
Si	Y	2.37489106	6.46865433	-0.27801949
Si	Y	7.5932189	0.86167972	-1.5653236
Si	Y	8.12353394	-1.70380646	0.0972643
Si	Y	5.20132513	5.41844472	-0.90000147
O	Y	2.58205508	-4.33248896	1.60138929
Si	Y	1.20561363	-4.24080822	2.43241844
H	Y	0.70181274	0.22667225	6.49665828
O	Y	2.64647612	-4.2194486	-1.02447904
H	Y	-0.86763969	5.85410793	-0.59224156
H	Y	7.36514055	2.90324274	1.49615322
H	Y	3.0172991	-3.20707544	-3.25571285
H	Y	8.68858311	-1.50207563	1.38428722
H	Y	8.48292186	4.19167374	-0.0664065
H	Y	5.3428826	-3.06863386	-1.66723708
H	Y	1.86209564	6.5719482	-1.59815126
H	Y	6.78204908	-4.82204252	-1.21492242
H	Y	2.68970495	7.75799316	0.22684622
H	Y	9.12688374	-2.18868822	-0.78283635
H	Y	8.91444982	0.98518674	-2.07079884
H	Y	5.15512692	4.87277292	-2.21015288
H	Y	6.6981041	0.55436973	-2.62396009
H	Y	5.82294876	6.69446675	-0.94190231
Si	Y	2.044348	-3.89143313	-2.4801519
O	N	-0.02343086	0.03920489	4.07595948
O	N	-2.82483916	-0.01941877	-4.33373095
H	Y	0.79704803	-1.82572498	5.42167873
H	Y	3.24175998	-6.27798447	0.21904953
H	Y	-4.20001242	4.20488974	-4.12543099
H	Y	1.49553102	-4.13483636	3.81834287
H	Y	-2.60586408	5.30653061	-2.8519049
H	Y	-1.25111937	-0.95636258	6.0832937
H	Y	-5.23280366	-0.34664309	-5.12212104
H	Y	-3.95777358	-1.21044176	5.29574857
H	Y	-6.59925364	-1.57232937	-2.96778438
H	Y	1.68718566	-5.10024112	-3.13406378
H	Y	-0.89831844	6.439491	1.65472275
O	N	0.79761179	-2.8436859	-2.36976179

O	N	-0.54858291	3.98626749	1.04266313
O	N	-0.94753953	1.99513975	2.66018707
O	N	-1.48247513	-1.82758495	-3.05635085
O	N	-2.65538772	0.21028054	3.68437438
O	N	-4.10786966	-1.60221467	-2.61837524
H	Y	-5.13712237	0.20441116	3.87820835
H	Y	-5.72465954	-3.09791174	-1.44797033
O	Y	-4.16514592	-1.85701392	2.91310945
H	Y	-4.82606626	5.15251413	2.79963895
O	Y	-5.81780733	-0.71720236	-0.77779748
H	Y	-2.06547571	-6.60360982	-1.70168176
O	N	-4.71262248	1.68421001	-0.50541892
H	Y	-1.74116117	-5.46360386	0.28404055
O	N	-2.24916821	-3.16045109	1.60090859
H	Y	-5.03444968	4.0546292	0.77501468
O	N	-2.98562377	3.00069945	-2.32022377
H	Y	0.43364559	-5.40953258	2.19858163
O	N	0.52667047	-2.84067493	1.96560567
H	Y	-4.71526782	4.90667936	-1.97642634
H	Y	-6.1536893	0.69606895	1.22665443
H	Y	-3.66882717	-5.10259342	-0.96399565
H	Y	-4.62478903	-3.55932385	1.1756293
H	Y	-5.32066917	2.88928592	2.76596111
H	Y	-7.15648527	1.36827798	-0.75272227
O	N	-1.43651059	-4.16531504	-1.79389069
H	Y	-3.77365485	-4.28661301	3.20589118
O	N	-2.97014659	3.62079364	2.07528018
Si	Y	0.0390251	-0.63598232	5.58401608
Si	Y	-3.99346425	0.25194994	-5.47162767
Si	N	-1.18490748	0.43048969	3.0323473
Si	N	-2.73109713	-0.7980431	-2.92708556
Si	Y	-4.02174209	-0.66790228	3.98502069
Si	Y	-5.61454638	-1.78037401	-1.96601106
Al	N	-3.13578021	1.80662946	-1.13347924
Al	N	-0.87928837	-2.21131964	1.26031156
Si	Y	-6.00179919	0.76912145	-0.18331201
Si	Y	-2.27596732	-5.37133561	-1.02814987
Si	Y	-3.72964818	-3.25198684	2.23430777
Si	Y	-4.59759213	3.92677409	2.12007698
Si	N	-1.64554584	2.90063395	1.50404411
Si	N	-0.74062483	-2.71436646	-1.91298029
Si	Y	-3.69043532	4.39075655	-2.81311301
H	Y	-4.14045576	1.6636894	-5.51435624

H	Y	-3.57970558	-0.21958455	-6.7455468

Table S7 – Coordinates for DFT Model of α-MeOH (BEA)

element	constrained during optimization?	X	Y	Z
Fe	N	-0.87493352	0.20179363	-0.30248982
O	N	-1.03191452	-0.4772009	1.65429081
O	N	-2.40134985	0.23750702	-1.69300219
O	N	-1.94130703	1.95823977	0.18720259
O	N	-0.74770353	-1.87526633	-0.50682239
O	N	0.88647998	1.29258603	-0.2116378
H	N	0.7508182	2.23251501	-0.01363312
C	N	2.11106502	0.81270246	0.40400945
H	N	2.01810547	0.83468201	1.4896733
H	N	2.24989768	-0.21056128	0.06425877
H	N	2.94548428	1.42954035	0.07201882
O	Y	4.86222375	-4.50175858	0.32392216
O	Y	1.32971847	5.73640801	0.63172018
O	Y	6.92661469	-2.8718311	0.17822878
O	Y	6.08579872	4.35746043	0.02695722
O	Y	3.76264124	5.50804845	-0.35169763
O	Y	7.18619915	2.16889832	-0.92086971
O	Y	7.55170103	-0.40095558	-0.492804
Si	Y	3.306199	-4.91570932	0.27049828
Si	Y	-0.26963826	5.54440954	0.66203665
Si	Y	5.97546997	-3.86943969	-0.65168013
Si	Y	7.34137065	3.35935241	0.14788813
Si	Y	2.43899864	6.4211711	-0.31322922
Si	Y	7.60890536	0.76874804	-1.59714859
Si	Y	8.12160551	-1.7969487	0.07063202
Si	Y	5.25554136	5.34645529	-0.93821387
O	Y	2.56189504	-4.37672721	1.59205232
Si	Y	1.18790051	-4.27182016	2.42556286
H	Y	0.72859132	0.20943266	6.47990083
O	Y	2.62201567	-4.27056499	-1.03420526
H	Y	-0.80904953	5.83231922	-0.61957338
H	Y	7.40355231	2.81949879	1.45979321
H	Y	2.99663863	-3.26664594	-3.26862207
H	Y	8.6908364	-1.59671944	1.35604583
H	Y	8.52868543	4.09499349	-0.10809451

H	Y	5.32642561	-3.14332613	-1.68506719
H	Y	1.92443889	6.52544955	-1.63259697
H	Y	6.75215575	-4.90730955	-1.23131897
H	Y	2.76530904	7.70911782	0.18787637
H	Y	9.11922062	-2.29211991	-0.81025859
H	Y	8.93009145	0.88025818	-2.10552191
H	Y	5.20229524	4.79801204	-2.24693992
H	Y	6.70921423	0.46618623	-2.6532706
H	Y	5.88745552	6.6172627	-0.98444351
Si	Y	2.01968775	-3.94117256	-2.48948461
O	N	-0.00089567	0.06494848	4.08122606
O	N	-2.81945973	-0.05036195	-4.35043728
H	Y	0.80496602	-1.84626599	5.40973451
H	Y	3.20297956	-6.3308725	0.21315631
H	Y	-4.16176378	4.20178235	-4.14217424
H	Y	1.48142568	-4.1648645	3.81065218
H	Y	-2.55616521	5.29346766	-2.87447259
H	Y	-1.23473192	-0.95863924	6.07326066
H	Y	-5.23358062	-0.34357713	-5.12574895
H	Y	-3.94492642	-1.19254959	5.29166388
H	Y	-6.60568631	-1.55287148	-2.96575053
H	Y	1.65139022	-5.14860652	-3.13974833
H	Y	-0.83049217	6.42336406	1.62601567
O	N	0.7879022	-2.87591182	-2.39889864
O	N	-0.46387488	3.9495379	1.01747896
O	N	-0.92837877	1.9827154	2.62847584
O	N	-1.49288076	-1.8472495	-3.046787
O	N	-2.64027773	0.22864479	3.69694634
O	N	-4.12051997	-1.61724204	-2.63423801
H	Y	-5.11552461	0.2284386	3.87300752
H	Y	-5.74053094	-3.08185358	-1.44395163
O	Y	-4.16229468	-1.84316883	2.91101804
H	Y	-4.76631146	5.17121964	2.78178777
O	Y	-5.81294805	-0.6988504	-0.77939325
H	Y	-2.11054033	-6.61786718	-1.69632964
O	N	-4.69905515	1.70915586	-0.51457936
H	Y	-1.77300292	-5.47574493	0.28597019
O	N	-2.2635486	-3.16825164	1.58748661
H	Y	-4.98765599	4.07017779	0.76025552
O	N	-2.95892751	2.98532139	-2.35889162
H	Y	0.40597332	-5.43477486	2.19609087
O	N	0.51797619	-2.86705513	1.97282015
H	Y	-4.66701242	4.9129428	-1.99387416

H	Y	-6.13332053	0.72195437	1.2222701
H	Y	-3.70014072	-5.10205129	-0.95914543
H	Y	-4.63924355	-3.54587142	1.17859409
H	Y	-5.27940224	2.9120235	2.75459196
H	Y	-7.13453902	1.39752784	-0.75675988
O	N	-1.45064962	-4.19282593	-1.79828309
H	Y	-3.79002911	-4.27516369	3.20894152
O	N	-2.90919754	3.68290215	2.08169025
Si	Y	0.05698504	-0.64999457	5.57066695
Si	Y	-3.99010269	0.24404654	-5.4791496
Si	N	-1.17313151	0.41859019	3.02599014
Si	N	-2.73957246	-0.80994169	-2.92774361
Si	Y	-4.00707782	-0.65267736	3.97974768
Si	Y	-5.62071743	-1.76651353	-1.96541337
Al	N	-3.12104493	1.80439445	-1.15446412
Al	N	-0.88205328	-2.22483776	1.25468125
Si	Y	-5.98364296	0.79035701	-0.18816623
Si	Y	-2.30964666	-5.38229307	-1.02538609
Si	Y	-3.73952608	-3.2432827	2.23475796
Si	Y	-4.54918199	3.9420183	2.10476289
Si	N	-1.62855547	2.90557653	1.48493314
Si	N	-0.74804548	-2.74124449	-1.91014412
Si	Y	-3.64808201	4.38666151	-2.83131772
H	Y	-4.12567267	1.65682999	-5.52502483
H	Y	-3.58273207	-0.2339239	-6.75272911

Table S8 – Coordinates for DFT Model of MeOH-Lattice Complex (BEA)

element	constrained during optimization?	X	Y	Z
O	N	-1.06942049	-0.66917138	1.77002981
O	N	-2.32789161	0.30915243	-1.75601242
O	N	-1.9133366	2.05713545	0.21464748
O	N	-0.86053165	-2.16769033	-0.36864897
O	N	0.07851261	0.68211435	-1.07646296
H	N	-0.2223849	1.16237376	-0.28533347
C	N	0.82140306	1.56988289	-1.9275306
H	N	1.70434428	1.93669941	-1.40094447
H	N	1.13441994	0.98656233	-2.79144607
H	N	0.20725948	2.413284	-2.25485835
O	Y	4.84905553	-4.49380269	0.33423381

O	Y	1.33777062	5.74865296	0.72827339
O	Y	6.91775098	-2.86717533	0.21521131
O	Y	6.09371241	4.36516371	0.13301002
O	Y	3.77501161	5.52478289	-0.24540725
O	Y	7.19393641	2.18381215	-0.83149995
O	Y	7.55162155	-0.39106187	-0.42768271
Si	Y	3.29239766	-4.90366671	0.26894236
Si	Y	-0.26228637	5.55982543	0.74871516
Si	Y	5.96851447	-3.85418166	-0.62942921
Si	Y	7.34645673	3.36304141	0.25000871
Si	Y	2.45321697	6.44049882	-0.20419216
Si	Y	7.61687805	0.78961982	-1.51983261
Si	Y	8.11563814	-1.79396995	0.12438773
Si	Y	5.27043227	5.36574856	-0.82617283
O	Y	2.5427636	-4.3764491	1.59221663
Si	Y	1.16490037	-4.27674899	2.41995591
H	Y	0.69546102	0.164236	6.51717682
O	Y	2.61613724	-4.24386776	-1.03252268
H	Y	-0.79456742	5.8621718	-0.53247421
H	Y	7.40095383	2.80980624	1.55666988
H	Y	3.00401137	-3.21823708	-3.25478717
H	Y	8.678947	-1.60804238	1.4145508
H	Y	8.53664997	4.09854164	0.00733556
H	Y	5.32619944	-3.11618472	-1.6585937
H	Y	1.94542464	6.55925016	-1.52497393
H	Y	6.74575588	-4.88789335	-1.21571587
H	Y	2.77990171	7.72253622	0.3115098
H	Y	9.11649869	-2.28246547	-0.75654943
H	Y	8.94080698	0.90327194	-2.02054028
H	Y	5.22244444	4.83069219	-2.14063165
H	Y	6.72175293	0.49979472	-2.58338012
H	Y	5.90538182	6.63552366	-0.85642199
Si	Y	2.02153652	-3.89819838	-2.48715379
O	N	-0.01725854	0.02791329	4.11523276
O	N	-2.77133179	-0.05197932	-4.38182566
H	Y	0.77259692	-1.88071258	5.42667972
H	Y	3.18633682	-6.31794915	0.1967795
H	Y	-4.13342236	4.27492205	-4.08789035
H	Y	1.45183864	-4.18453976	3.80750018
H	Y	-2.5316979	5.35004184	-2.80125968
H	Y	-1.26837986	-0.99521543	6.08910221
H	Y	-5.2104524	-0.25785138	-5.12274607
H	Y	-3.97520068	-1.21502377	5.2918271

H	Y	-6.5958668	-1.48587541	-2.98185917
H	Y	1.65401278	-5.09831255	-3.151615
H	Y	-0.82574024	6.43047721	1.71885457
O	N	0.78540851	-2.84617173	-2.34855394
O	N	-0.48467331	4.01314419	1.20067597
O	N	-0.92347128	1.89886918	2.62247829
O	N	-1.45639825	-1.80531318	-2.96803884
O	N	-2.65618388	0.20116204	3.71963311
O	N	-4.10397122	-1.54290768	-2.62050849
H	Y	-5.13560581	0.22286032	3.88190977
H	Y	-5.7416357	-3.03209968	-1.47139736
O	Y	-4.18223947	-1.841084	2.90365093
H	Y	-4.77003602	5.17559813	2.84249622
O	Y	-5.81204316	-0.65577754	-0.78304167
H	Y	-2.11829404	-6.57356055	-1.74167605
O	N	-4.67210696	1.69592304	-0.51714954
H	Y	-1.78804991	-5.45237509	0.25375146
O	N	-2.22009021	-3.11570472	1.66059436
H	Y	-4.98389019	4.09561301	0.80885788
O	N	-2.97106471	3.01232318	-2.33461117
H	Y	0.38154096	-5.43561705	2.1749009
O	N	0.37668121	-2.9111599	1.92353512
H	Y	-4.64771362	4.96539136	-1.93498546
H	Y	-6.1391818	0.74544393	1.23128303
H	Y	-3.70817917	-5.06174698	-0.99704588
H	Y	-4.65439432	-3.52507389	1.15175566
H	Y	-5.28803663	2.91800685	2.78993176
H	Y	-7.12911807	1.44325147	-0.74572426
O	N	-1.46711115	-4.23114509	-1.93635162
H	Y	-3.81684499	-4.27679627	3.17877762
O	N	-2.92806895	3.62990408	2.19240348
Si	Y	0.02644828	-0.68442846	5.59575011
Si	Y	-3.96380939	0.33051562	-5.46378732
Si	N	-1.18130503	0.36045416	3.044126
Si	N	-2.71420393	-0.79830674	-2.95971316
Si	Y	-4.02960777	-0.6618972	3.98499767
Si	Y	-5.61631346	-1.71170838	-1.9787642
Si	N	-0.92610165	-2.20475243	1.25617743
Si	Y	-5.98246284	0.82779082	-0.17764699
Si	Y	-2.31794849	-5.34426264	-1.05911882
Si	Y	-3.75927739	-3.23533471	2.21537142
Si	Y	-4.55227438	3.95270909	2.1539684
Si	N	-1.61959664	2.91149279	1.54826463

Si	N	-0.75132602	-2.77806397	-1.86848614
Si	Y	-3.6257276	4.44527522	-2.77274006
H	Y	-4.09617182	1.7439631	-5.49627357
H	Y	-3.55134862	-0.13552178	-6.74033631
H	N	-1.31549439	0.3412137	-1.52681502
Al	N	-3.06823854	1.87640144	-1.06777356

Table S9 – Coordinates for DFT Model of α-OH···CH_3 Complex (CHA)

element	constrained during optimization?	X	Y	Z
Fe	N	0.43956566	0.0002841	0.76787327
O	N	-0.07750667	1.5082483	-0.56037765
O	N	1.65230782	1.69061476	1.20644832
O	N	1.65103493	-1.69017097	1.20759459
O	N	-0.07860585	-1.50806352	-0.55923202
O	N	-0.80588079	0.0013323	1.99839168
H	N	-1.78985265	0.00333617	2.02433995
C	N	-3.63709433	0.00033431	1.51924325
H	N	-3.62152262	0.9302752	0.96785195
H	N	-3.6190251	-0.93042656	0.96930704
H	N	-4.01060369	0.00046853	2.53496541
O	N	-1.73231354	0.00064424	-1.9567264
O	N	-2.33307251	2.43309357	-1.80977368
O	N	-2.3336837	-2.4309627	-1.81074441
O	Y	-4.18054375	3.3137707	-0.17803865
O	Y	-6.0538145	2.43860121	1.44080375
O	Y	-4.17956345	-3.31324777	-0.17943522
O	Y	-7.01118632	0.0001211	1.26158047
O	Y	-6.05294263	-2.43941297	1.44006774
Si	N	-1.06841561	1.46863281	-1.89096819
Si	N	-1.06840732	-1.46732644	-1.89049771
Si	Y	-5.23680809	3.73949866	0.95855558
Si	Y	-3.1226263	-3.73898055	-1.31677176
Si	Y	-7.39841183	1.55121008	1.43477609
Si	Y	-5.23582919	-3.73952417	0.95682368
Si	Y	-7.39771442	-1.55279845	1.43468822
Si	Y	-3.12434191	3.74004679	-1.31571904
O	N	-0.22681486	1.81614298	-3.22097441
O	N	-0.22662535	-1.81620306	-3.21996429
O	N	3.19485237	0.00014075	2.3221407

O	Y	4.14166443	2.00923425	-1.64948624
H	Y	3.83698812	4.24410954	-2.67062225
H	Y	6.26699869	2.33299077	-0.42169885
O	Y	2.33246753	2.46505917	-3.50165081
Si	Y	3.11647452	3.09742839	-2.2458615
Si	Y	5.08140669	1.55199829	-0.42600713
O	N	0.09971126	4.26530763	0.67898206
H	Y	-3.81518242	4.29692816	-2.42410969
H	Y	-0.81761429	6.65542972	0.35126662
O	N	2.06370024	3.49435729	-1.09318275
H	Y	-1.84910778	5.1012945	1.73051037
Al	N	0.94474648	2.9201839	0.06385808
Si	Y	-1.15939967	5.28562628	0.50349914
O	Y	5.46796142	-0.00008637	-0.59846492
O	N	3.24273918	-2.26555687	3.42919496
O	N	4.33162293	-1.95115203	0.98857863
Si	N	3.1102897	-1.56558141	1.97736212
O	Y	4.03602589	0.00020144	4.96026169
O	N	3.24318938	2.26596966	3.42885979
H	Y	4.64269909	-2.35812501	5.40889016
H	Y	4.64354027	2.35889088	5.40680012
O	N	4.33286587	1.95111559	0.98870666
H	Y	2.42491625	-1.73561078	5.6900006
H	Y	2.42538171	1.73740142	5.68917386
Si	N	3.11111482	1.56585638	1.97710364
Si	Y	3.59270897	-1.54593665	4.90484147
Si	Y	3.59345449	1.54707435	4.90487576
O	Y	1.58614901	-0.00103234	-4.014346
H	Y	0.86992122	-1.91231743	-5.41578259
H	Y	0.8702113	1.9106962	-5.41604299
Si	Y	1.1427779	-1.54618426	-4.07120145
Si	Y	1.14281952	1.54484465	-4.07184476
O	Y	4.14107955	-2.01078895	-1.64981393
H	Y	6.26670976	-2.33402304	-0.42035806
H	Y	3.83564198	-4.24592113	-2.66968921
O	Y	2.33287438	-2.46795453	-3.50021219
Si	Y	5.08069599	-1.55301136	-0.42604104
Si	Y	3.11507257	-3.09859301	-2.24452489
O	N	0.0984241	-4.26536378	0.68228785
H	Y	-0.81885808	-6.6546178	0.35358918
O	N	2.06265166	-3.49417041	-1.09093424
H	Y	-1.8510646	-5.09974246	1.73290616
Al	N	0.94310497	-2.9205198	0.0657613

Si	Y	-1.16114676	-5.28541994	0.50534586
O	Y	-2.09901942	4.82783747	-0.72027317
H	Y	-4.54664011	4.29661639	2.06719894
H	Y	-6.14287562	4.69904133	0.43413274
O	Y	-2.09610832	-4.82619466	-0.72184799
H	Y	-8.22785097	1.9557549	0.35575404
H	Y	-3.8134613	-4.29609679	-2.42468826
H	Y	-8.0883219	1.73589162	2.66234083
H	Y	-4.54431648	-4.29740709	2.06449057
H	Y	-8.22726273	-1.9572567	0.35531552
H	Y	-6.14135111	-4.69998669	0.43180572
H	Y	-8.08742606	-1.73811782	2.6616992

Table S10 – Coordinates for DFT Model of α-Fe(II) (CHA)

element	constrained during optimization?	X	Y	Z
Fe	N	0.64406224	0.00033164	0.57190717
O	N	-0.1741793	1.54244769	-0.49332666
O	N	1.5833205	1.7365958	1.25685225
O	N	1.58211543	-1.73628717	1.25798601
O	N	-0.17558635	-1.54210711	-0.491994
O	N	-1.82055954	0.00076919	-1.86089377
O	N	-2.44211283	2.43854763	-1.73589861
O	N	-2.44294202	-2.43622434	-1.73739881
O	Y	-4.25891576	3.31370039	-0.05283248
O	Y	-6.10765882	2.43867248	1.5939389
O	Y	-4.25788521	-3.31308683	-0.05428291
O	Y	-7.06761418	0.00020727	1.4289581
O	Y	-6.10686145	-2.43934992	1.59303196
Si	N	-1.17906326	1.47741305	-1.81405041
Si	N	-1.17921507	-1.47590045	-1.81353535
Si	Y	-5.29777071	3.73960074	1.09972273
Si	Y	-3.2179759	-3.73888301	-1.20717828
Si	Y	-7.4521883	1.55129565	1.60796747
Si	Y	-5.29703992	-3.73945335	1.09763089
Si	Y	-7.4515391	-1.552713	1.60776216
Si	Y	-3.21956249	3.74020812	-1.20592878
O	N	-0.34595561	1.81704696	-3.15185385
O	N	-0.34559619	-1.81708101	-3.15057835
O	N	3.06160646	0.00000761	2.33298365

O	Y	4.04057187	2.00922092	-1.64816452
H	Y	3.72074716	4.24416724	-2.6645497
H	Y	6.18399063	2.33296942	-0.45224735
O	Y	2.20393909	2.4651931	-3.47308869
Si	Y	3.00661261	3.09749241	-2.22910073
Si	Y	4.99853158	1.55201814	-0.43879751
O	N	0.04917566	4.29158107	0.75605052
H	Y	-3.92693426	4.29713982	-2.30382574
H	Y	-0.88820594	6.65545681	0.42656573
O	N	1.9768986	3.50044893	-1.0598991
H	Y	-1.89903472	5.1012751	1.82097761
Al	N	0.86144738	2.95032763	0.10852293
Si	Y	-1.22766356	5.28561286	0.58388588
O	Y	5.38242801	-0.00011681	-0.61714707
O	N	3.22388779	-2.26723277	3.44535878
O	N	4.26944492	-1.92135877	0.99410986
Si	N	3.05124339	-1.57434908	1.9975059
O	Y	4.03364769	-0.0000048	4.96230672
O	N	3.22424626	2.26734017	3.44509671
H	Y	4.64679844	-2.35838428	5.40182903
H	Y	4.64768074	2.35869081	5.39989145
O	N	4.27079749	1.92105162	0.99441378
H	Y	2.43349366	-1.73577045	5.71601527
H	Y	2.43399753	1.73715338	5.7152985
Si	N	3.05202538	1.57433414	1.99725408
Si	Y	3.58939779	-1.54621417	4.9135192
Si	Y	3.59017997	1.54701222	4.91359842
O	Y	1.45002969	-0.0008941	-3.97465119
H	Y	0.71296019	-1.91210126	-5.36534501
H	Y	0.71329482	1.91089582	-5.3654736
Si	Y	1.00576646	-1.5460708	-4.02501555
Si	Y	1.00594547	1.54506331	-4.02557401
O	Y	4.03990549	-2.01072906	-1.64860179
H	Y	6.18364375	-2.33411062	-0.45105658
H	Y	3.71927552	-4.24586527	-2.66386872
O	Y	2.2043244	-2.46781921	-3.47183123
Si	Y	4.99773177	-1.55311287	-0.4389192
Si	Y	3.00513203	-3.09849458	-2.22791028
O	N	0.04772027	-4.2916266	0.7590511
H	Y	-0.8896193	-6.65458944	0.42843848
O	N	1.97559335	-3.49999323	-1.05792076
H	Y	-1.90112295	-5.09975803	1.8230454
Al	N	0.85964675	-2.95058105	0.11034779

Si	Y	-1.22959475	-5.28536285	0.58533141
O	Y	-2.18545062	4.82786775	-0.62580867
H	Y	-4.591252	4.29660756	2.19787181
H	Y	-6.2116639	4.69915972	0.58881468
O	Y	-2.18273224	-4.82611979	-0.62781563
H	Y	-8.29763156	1.95589007	0.54145779
H	Y	-3.92533381	-4.29593019	-2.30475856
H	Y	-8.12370042	1.73594482	2.84569772
H	Y	-4.58910102	-4.29737421	2.19482383
H	Y	-8.297109	-1.95712064	0.54087571
H	Y	-6.21031968	-4.69989681	0.58613888
H	Y	-8.1228668	-1.73806457	2.84492296

Table S11 – Coordinates for DFT Model of α-O (CHA)

element	constrained during optimization?	X	Y	Z
Fe	N	0.43069028	0.00031886	0.72569816
O	N	-0.13936506	1.49426365	-0.53051914
O	N	1.60373886	1.6709757	1.20775213
O	N	1.60240342	-1.67046946	1.20888027
O	N	-0.14055575	-1.49399071	-0.52934538
O	N	-0.6598037	0.00118572	1.87604154
O	N	-1.82178329	0.00074995	-1.91635687
O	N	-2.40530669	2.44332665	-1.76025299
O	N	-2.40654389	-2.44049535	-1.7606462
O	Y	-4.24212824	3.31364969	-0.09477907
O	Y	-6.09492253	2.43865857	1.54744319
O	Y	-4.24105294	-3.31313233	-0.09603779
O	Y	-7.05417896	-0.00015129	1.38019583
O	Y	-6.09410797	-2.4390012	1.54692632
Si	N	-1.16003798	1.46593573	-1.85044767
Si	N	-1.1601781	-1.46449959	-1.84995349
Si	Y	-5.28380669	3.73957633	1.05521532
Si	Y	-3.19829467	-3.73894927	-1.24638447
Si	Y	-7.43943724	1.55127913	1.55823456
Si	Y	-5.2830515	-3.73947306	1.05329095
Si	Y	-7.43893225	-1.55274738	1.55794652
Si	Y	-3.19989665	3.74015868	-1.24536022
O	N	-0.31767842	1.81735801	-3.17732608
O	N	-0.31745724	-1.81802411	-3.17602851

O	N	3.17537792	0.00011283	2.31436653
O	Y	4.06131253	2.00917673	-1.66948803
H	Y	3.74399434	4.24408046	-2.68672864
H	Y	6.20176151	2.33294504	-0.46828316
O	Y	2.22921788	2.46505909	-3.49897936
Si	Y	3.02874439	3.09740264	-2.2529818
Si	Y	5.0162315	1.55201948	-0.4577567
O	N	0.0500761	4.26278177	0.71374154
H	Y	-3.90455485	4.29704353	-2.34494818
H	Y	-0.87261314	6.65543199	0.39288027
O	N	1.99169668	3.48626302	-1.08306239
H	Y	-1.8868873	5.10129862	1.78483981
Al	N	0.88080233	2.92167817	0.08283123
Si	Y	-1.21255552	5.2856853	0.5493707
O	Y	5.40061351	-0.00014354	-0.63510668
O	N	3.21470574	-2.26463463	3.41918341
O	N	4.27659146	-1.95552058	0.96279915
Si	N	3.07530149	-1.56316318	1.97067217
O	Y	4.03804343	0.00011604	4.94097824
O	N	3.21512233	2.26498882	3.41891261
H	Y	4.65009486	-2.35824801	5.38209029
H	Y	4.65096755	2.35881196	5.38002761
O	N	4.27801746	1.95531642	0.96303523
H	Y	2.43601659	-1.73564611	5.69078182
H	Y	2.43651637	1.7372912	5.68998594
Si	N	3.07618905	1.56339895	1.97045485
Si	Y	3.59389473	-1.54609549	4.89115305
Si	Y	3.59467998	1.54707782	4.8912114
O	Y	1.47654726	-0.00102546	-4.00230897
H	Y	0.74293639	-1.91227049	-5.39480732
H	Y	0.74325444	1.91071775	-5.39503773
Si	Y	1.03240811	-1.54622184	-4.05380105
Si	Y	1.03259938	1.54490917	-4.05445237
O	Y	4.06066809	-2.01079272	-1.66983524
H	Y	6.20143199	-2.33412153	-0.46697489
H	Y	3.74256234	-4.24594378	-2.68583889
O	Y	2.22960569	-2.46793012	-3.49758602
Si	Y	5.01547763	-1.55314182	-0.45781499
Si	Y	3.0273083	-3.09860802	-2.25168097
O	N	0.04889539	-4.26297057	0.71672264
H	Y	-0.87398071	-6.65461462	0.39511822
O	N	1.9905508	-3.48596408	-1.08088935
H	Y	-1.88893248	-5.09974758	1.78717579

Al	N	0.87908462	-2.92188439	0.08470401
Si	Y	-1.21434559	-5.2854378	0.55111969
O	Y	-2.16725299	4.82782146	-0.66262729
H	Y	-4.58002337	4.29661521	2.1550853
H	Y	-6.19644624	4.6991181	0.54201001
O	Y	-2.16449843	-4.8261883	-0.66436555
H	Y	-8.28224859	1.95584834	0.48953983
H	Y	-3.90292949	-4.29604772	-2.34564158
H	Y	-8.11408236	1.73595032	2.79421543
H	Y	-4.57783815	-4.29737352	2.15227522
H	Y	-8.28172147	-1.95716435	0.4890667
H	Y	-6.19505371	-4.69992237	0.53955113
H	Y	-8.11326769	-1.73810005	2.79350853

Table S12 – Coordinates for DFT Model of α-OH (CHA)

element	constrained during optimization?	X	Y	Z
Fe	N	0.37108239	-0.00136969	0.78271673
O	N	-0.1451366	1.49590057	-0.5267685
O	N	1.60550984	1.68319114	1.21096759
O	N	1.59896882	-1.67209074	1.21752678
O	N	-0.13953104	-1.50258828	-0.5339549
O	N	-0.85245754	0.00673157	2.05031402
H	N	-1.68753407	0.4239811	2.27662958
O	N	-1.80965153	-0.00407613	-1.92283604
O	N	-2.39679947	2.43482158	-1.76966602
O	N	-2.39973166	-2.44142114	-1.77416873
O	Y	-4.23316298	3.31256021	-0.11719167
O	Y	-6.08900955	2.43985486	1.52261509
O	Y	-4.2318172	-3.3140232	-0.10901817
O	Y	-7.04835097	0.00117932	1.35702122
O	Y	-6.08791643	-2.43798803	1.52896034
Si	N	-1.14446634	1.4606543	-1.85924787
Si	N	-1.14470447	-1.47125831	-1.85739416
Si	Y	-5.2772334	3.74023119	1.02994616
Si	Y	-3.18678772	-3.7415049	-1.25665747
Si	Y	-7.43364455	1.5525947	1.53199289
Si	Y	-5.27581895	-3.73889067	1.03921367
Si	Y	-7.43271452	-1.55159569	1.53650669
Si	Y	-3.18878097	3.7373718	-1.26645921

O	N	-0.30227872	1.8113666	-3.18649742
O	N	-0.30687104	-1.81924996	-3.18897912
O	N	3.17564456	0.00462808	2.31289952
O	Y	4.07327558	2.00638631	-1.67422905
H	Y	3.75778388	4.23979264	-2.69538526
H	Y	6.21145339	2.33204404	-0.46939892
O	Y	2.24468268	2.45947339	-3.50792981
Si	Y	3.04191143	3.09371058	-2.26142128
Si	Y	5.02597735	1.55101325	-0.460054
O	N	0.05016614	4.2556444	0.69944347
H	Y	-3.89131494	4.29277715	-2.36834982
H	Y	-0.86484284	6.6553587	0.37183142
O	N	1.99802742	3.48562136	-1.09766966
H	Y	-1.88167093	5.10320931	1.76413874
Al	N	0.8926704	2.91509517	0.07169203
Si	Y	-1.20487568	5.28570626	0.52965024
O	Y	5.41068153	-0.00132255	-0.6343889
O	N	3.22163368	-2.25868	3.41880993
O	N	4.27980869	-1.95044263	0.96274697
Si	N	3.07397369	-1.55943076	1.96792236
O	Y	4.03739963	0.0070397	4.9392173
O	N	3.22157568	2.27074245	3.41298265
H	Y	4.64889818	-2.35061252	5.38482572
H	Y	4.6494808	2.36640271	5.37583701
O	N	4.28470296	1.95532537	0.95834876
H	Y	2.4341667	-1.7277631	5.68836523
H	Y	2.43445089	1.74524416	5.68246379
Si	N	3.07780423	1.56739243	1.96343859
Si	Y	3.59364454	-1.53919442	4.89065437
Si	Y	3.59423579	1.55380967	4.88616878
O	Y	1.49313847	-0.007389	-4.00917242
H	Y	0.76224085	-1.92074291	-5.40014106
H	Y	0.76236906	1.90226747	-5.40599085
Si	Y	1.04928967	-1.55259056	-4.05897762
Si	Y	1.04916619	1.53834997	-4.06413262
O	Y	4.07290993	-2.01358756	-1.66869602
H	Y	6.2114137	-2.33497251	-0.46123849
H	Y	3.75688059	-4.25023576	-2.6820467
O	Y	2.2453266	-2.47358962	-3.49929891
Si	Y	5.02542138	-1.55401462	-0.45554881
Si	Y	3.04082462	-3.10230439	-2.25101869
O	N	0.05744847	-4.26454685	0.70875652
H	Y	-0.86538835	-6.65467246	0.39360064

O	N	2.0017943	-3.48684483	-1.08141634
H	Y	-1.88308587	-5.09781865	1.78145987
Al	N	0.88800745	-2.91720668	0.08201174
Si	Y	-1.20608343	-5.285223	0.54695913
O	Y	-2.15736912	4.82615035	-0.68344626
H	Y	-4.57546293	4.29888513	2.13042127
H	Y	-6.18881667	4.69891497	0.5136357
O	Y	-2.15400467	-4.82787244	-0.67094649
H	Y	-8.27439576	1.95542022	0.46125454
H	Y	-3.88920432	-4.3004099	-2.35636383
H	Y	-8.11050655	1.73893235	2.76658767
H	Y	-4.57273461	-4.29501994	2.14025755
H	Y	-8.27362877	-1.95755013	0.46653259
H	Y	-6.18683347	-4.70013763	0.52502926
H	Y	-8.10946788	-1.73504266	2.7709989

Table S13 – Coordinates for DFT Model of α-OCH_3 (CHA)

element	constrained during optimization?	X	Y	Z
Fe	N	-0.39157967	-0.0003875	0.75576297
O	N	0.09947851	-1.5110778	-0.5677878
O	N	-1.61634215	-1.69106569	1.21642237
O	N	-1.61517574	1.6903601	1.21769784
O	N	0.10075862	1.51116116	-0.56634329
O	N	0.94520129	-0.0014753	1.89644267
C	N	2.33594573	-0.00419693	2.09969198
H	N	2.78384044	-0.89340365	1.64419161
H	N	2.78602544	0.88769297	1.65164778
H	N	2.53954111	-0.00893839	3.17376884
O	N	1.74963488	-0.00044584	-1.96505294
O	N	2.35123354	-2.4359787	-1.82069557
O	N	2.35199852	2.43429448	-1.82172653
O	Y	4.19614853	-3.31345979	-0.17773241
O	Y	6.06746241	-2.43854465	1.44340497
O	Y	4.19489531	3.31332604	-0.17847148
O	Y	7.02497059	-0.00003192	1.26541803
O	Y	6.06650571	2.43947779	1.44302098
Si	N	1.08998464	-1.46970346	-1.89974815
Si	N	1.09010923	1.46883043	-1.89900081
Si	Y	5.25085583	-3.73944682	0.96030104

Si	Y	3.13914275	3.73919333	-1.31682263
Si	Y	7.41203057	-1.55112686	1.43892822
Si	Y	5.24986663	3.73960742	0.95900827
Si	Y	7.41128512	1.55288178	1.4390576
Si	Y	3.14094508	-3.73990437	-1.31641232
O	N	0.24503359	-1.81806302	-3.22674805
O	N	0.24473046	1.81886671	-3.22525606
O	N	-3.16277284	-0.00032706	2.32506694
O	Y	-4.12468104	-2.0090964	-1.65798
H	Y	-3.8188483	-4.24390658	-2.67892157
H	Y	-6.25132907	-2.33299658	-0.4325227
O	Y	-2.31343359	-2.46478932	-3.50818121
Si	Y	-3.09878881	-3.09726109	-2.25325323
Si	Y	-5.06573322	-1.55202208	-0.43545763
O	N	-0.08456613	-4.26432298	0.6768569
H	Y	3.83308265	-4.2966777	-2.42402107
H	Y	0.83249554	-6.65537204	0.3478833
O	N	-2.04947209	-3.48934168	-1.09536325
H	Y	1.86246355	-5.10131653	1.72835561
Al	N	-0.92635005	-2.91868596	0.05880871
Si	Y	1.17413042	-5.28562889	0.50059915
O	Y	-5.45216056	0.00011334	-0.60820649
O	N	-3.22956227	2.26580714	3.42227758
O	N	-4.3030539	1.94177438	0.9761366
Si	N	-3.08409964	1.56390408	1.97139573
O	Y	-4.02633868	-0.00054451	4.95202876
O	N	-3.22991189	-2.26676285	3.4216664
H	Y	-4.63344743	2.35777079	5.40017921
H	Y	-4.63422117	-2.35929041	5.3977542
O	N	-4.30425748	-1.94193527	0.97602388
H	Y	-2.41597842	1.73522149	5.6836502
H	Y	-2.41639605	-1.73774848	5.68258889
Si	N	-3.08487338	-1.564499	1.97095617
Si	Y	-3.58293668	1.54560596	4.89719398
Si	Y	-3.58363651	-1.54740451	4.89703341
O	Y	-1.56660499	0.0013647	-4.01985207
H	Y	-0.84889654	1.9127397	-5.42041707
H	Y	-0.84911244	-1.91025088	-5.42094206
Si	Y	-1.1231067	1.54657736	-4.07620059
Si	Y	-1.12322979	-1.54454474	-4.0770828
O	Y	-4.1241428	2.01088727	-1.6580169
H	Y	-6.25110919	2.33404061	-0.43085277
H	Y	-3.81763185	4.24611893	-2.67738473

O	Y	-2.31393687	2.46821225	-3.50639308
Si	Y	-5.06506726	1.55308285	-0.43528666
Si	Y	-3.09750582	3.09879499	-2.25149456
O	N	-0.08344757	4.26437308	0.68081576
H	Y	0.83354104	6.65467781	0.35114515
O	N	-2.04841058	3.48927507	-1.09272723
H	Y	1.86426802	5.09972466	1.73146489
Al	N	-0.92472537	2.91904748	0.06108231
Si	Y	1.17565545	5.28549073	0.50316538
O	Y	2.11501817	-4.82764453	-0.72211862
H	Y	4.55961002	-4.29659069	2.06805959
H	Y	6.15762305	-4.69892356	0.43670116
O	Y	2.11197689	4.8263639	-0.72308104
H	Y	8.24265354	-1.95558314	0.36078422
H	Y	3.83122436	4.29639473	-2.4240245
H	Y	8.10060389	-1.73588418	2.66723255
H	Y	4.5571583	4.29739128	2.0659434
H	Y	8.24200519	1.9574276	0.36062
H	Y	6.15595894	4.70013315	0.43503042
H	Y	8.09965503	1.73812525	2.66683391

Table S14 – Coordinates for DFT Model of C-O Recombination TS (CHA)

element	constrained during optimization?	X	Y	Z
Fe	N	0.43565724	-0.00320197	0.76391618
O	N	-0.08792594	1.51364167	-0.55221599
O	N	1.64433117	1.71136807	1.20969844
O	N	1.64308036	-1.69431003	1.20980373
O	N	-0.0904309	-1.52586141	-0.55274544
O	N	-0.79950884	0.00616413	2.03935795
H	N	-1.43131338	0.62815474	2.41497695
C	N	-3.4693545	0.01551726	1.40781789
H	N	-3.53364874	0.93770336	0.84802883
H	N	-3.31298448	-0.91419027	0.88328304
H	N	-3.81846821	-0.01375737	2.43116477
O	N	-1.72983373	-0.00588261	-1.95690732
O	N	-2.33759234	2.42797791	-1.81176791
O	N	-2.34287733	-2.4368044	-1.81439792
O	Y	-4.18356632	3.31146662	-0.18045778
O	Y	-6.05477374	2.43949421	1.44239032

O	Y	-4.18120077	-3.31529555	-0.16787733
O	Y	-7.01187123	0.00046932	1.26928913
O	Y	-6.05300179	-2.43851824	1.45181352
Si	N	-1.07230425	1.46293424	-1.88892608
Si	N	-1.07376498	-1.47777861	-1.88518612
Si	Y	-5.23838667	3.73955945	0.95667608
Si	Y	-3.12535555	-3.74326559	-1.30535084
Si	Y	-7.39919337	1.55184475	1.43967318
Si	Y	-5.2361435	-3.73948167	0.97042397
Si	Y	-7.39792927	-1.55215673	1.44604275
Si	Y	-3.12839453	3.73577736	-1.31996863
O	N	-0.22696936	1.81258038	-3.21645174
O	N	-0.23010841	-1.82519059	-3.21493894
O	N	3.16951652	0.00714733	2.32349412
O	Y	4.13748644	2.00552964	-1.65792968
H	Y	3.83130168	4.2382231	-2.68337618
H	Y	6.26407467	2.3322238	-0.4330929
O	Y	2.32621054	2.45718591	-3.50909056
Si	Y	3.11142242	3.09227573	-2.2554194
Si	Y	5.07860144	1.55101419	-0.43448551
O	N	0.09429443	4.26354975	0.67875462
H	Y	-3.82062721	4.29021418	-2.42873172
H	Y	-0.82049055	6.65498515	0.33848544
O	N	2.05962301	3.49701188	-1.10432796
H	Y	-1.85022245	5.10353882	1.72206329
Al	N	0.94309471	2.92466139	0.05369736
Si	Y	-1.16191407	5.28548529	0.49393581
O	Y	5.465266	-0.00135669	-0.60417041
O	N	3.24492842	-2.25683625	3.43018871
O	N	4.32722841	-1.9372267	0.98952852
Si	N	3.10206192	-1.56011566	1.97722832
O	Y	4.03930905	0.01023947	4.9560954
O	N	3.24405841	2.27493435	3.42260381
H	Y	4.64688885	-2.34704051	5.40898292
H	Y	4.64686793	2.36996089	5.39707331
O	N	4.3285036	1.94515702	0.98117492
H	Y	2.42929617	-1.72434506	5.69117669
H	Y	2.42912836	1.74866241	5.68312637
Si	N	3.10439592	1.5724478	1.97177581
Si	Y	3.5962138	-1.53608839	4.90437996
Si	Y	3.5963863	1.5569062	4.8979766
O	Y	1.57979829	-0.01011692	-4.01582209
H	Y	0.86243068	-1.92443007	-5.4125436

H	Y	0.86200647	1.89856125	-5.42075436
Si	Y	1.13655215	-1.555497	-4.0690478
Si	Y	1.13617005	1.5355727	-4.07614446
O	Y	4.13762943	-2.0144826	-1.64988627
H	Y	6.26463719	-2.33477614	-0.42204633
H	Y	3.83150735	-4.25178771	-2.66478136
O	Y	2.327539	-2.47579904	-3.49738556
Si	Y	5.07848332	-1.55399367	-0.42806672
Si	Y	3.11119942	-3.10371881	-2.24124171
O	N	0.10612091	-4.2768663	0.69189583
H	Y	-0.81931388	-6.65502987	0.36850593
O	N	2.06286453	-3.49823504	-1.08471023
H	Y	-1.85031582	-5.09747253	1.7456843
Al	N	0.93819312	-2.92830206	0.06884865
Si	Y	-1.16165942	-5.28558749	0.51774885
O	Y	-2.10268066	4.82488789	-0.72786734
H	Y	-4.54724376	4.29904436	2.06333462
H	Y	-6.14529428	4.69782971	0.43114176
O	Y	-2.09804686	-4.82909907	-0.709408
H	Y	-8.22986433	1.9539942	0.36070349
H	Y	-3.81734455	-4.30282688	-2.41145138
H	Y	-8.08782051	1.73895637	2.66759105
H	Y	-4.54335543	-4.29492361	2.07849601
H	Y	-8.22856234	-1.95900878	0.36840591
H	Y	-6.14205947	-4.70121001	0.44836827
H	Y	-8.08629113	-1.73504699	2.67417724

Table S15 – Coordinates for DFT Model of α-MeOH (CHA)

element	constrained during optimization?	X	Y	Z
Fe	N	-0.49667952	0.03250597	0.65651902
O	N	0.09353453	-1.51370229	-0.61056201
O	N	-1.59339729	-1.88715759	1.24640659
O	N	-1.58890051	1.75946508	1.24777801
O	N	0.14723964	1.62012919	-0.54861354
O	N	0.51954987	-0.56940842	2.3508367
H	N	0.18999296	-1.42172011	2.66300371
C	N	1.90854677	-0.36253709	2.69093041
H	N	2.54408865	-1.06400162	2.14800576
H	N	2.15119483	0.65611169	2.39816213

H	N	2.04473782	-0.47316586	3.76686606
O	N	1.73889873	0.05651086	-1.94569605
O	N	2.3608952	-2.38152177	-1.87714608
O	N	2.39615066	2.48524307	-1.8256888
O	Y	4.21150022	-3.29468697	-0.26377078
O	Y	6.08218179	-2.45409008	1.37614615
O	Y	4.20978814	3.33063902	-0.12454975
O	Y	7.03958245	-0.01230287	1.2500415
O	Y	6.08088414	2.42285197	1.47878947
Si	N	1.089499	-1.41936196	-1.93157416
Si	N	1.11158347	1.54249125	-1.88378305
Si	Y	5.26583819	-3.74455009	0.86538278
Si	Y	3.15440369	3.780386	-1.25402093
Si	Y	7.42668942	-1.56669192	1.39088947
Si	Y	5.26432647	3.73286327	1.0220537
Si	Y	7.42572695	1.53662131	1.45657721
Si	Y	3.15672677	-3.69705271	-1.41157623
O	N	0.26265208	-1.7316995	-3.28199234
O	N	0.26642805	1.89545612	-3.21253624
O	N	-2.99392807	-0.05806929	2.29400018
O	Y	-4.10886846	-1.9598649	-1.7190703
H	Y	-3.80254474	-4.17259989	-2.78689748
H	Y	-6.23594276	-2.30972929	-0.5015389
O	Y	-2.29696508	-2.37626687	-3.57785955
Si	Y	-3.08271657	-3.0351385	-2.33684521
Si	Y	-5.05046643	-1.52880157	-0.487492
O	N	-0.07163567	-4.28435304	0.58401164
H	Y	3.84929744	-4.23026428	-2.53045498
H	Y	0.84789712	-6.64716603	0.18995122
O	N	-2.03716239	-3.48438253	-1.20034107
H	Y	1.87726824	-5.1225487	1.60330227
Al	N	-0.92118596	-2.94309064	-0.02875452
Si	Y	1.18937801	-5.28092006	0.371682
O	Y	-5.43689464	0.02663846	-0.62767843
O	N	-3.23159663	2.1887346	3.44720544
O	N	-4.28772097	1.8719938	1.00957354
Si	N	-3.04192384	1.52335497	1.9836168
O	Y	-4.01304341	-0.09136673	4.93187033
O	N	-3.25741854	-2.33793306	3.36573874
H	Y	-4.62045494	2.25691578	5.42955744
H	Y	-4.62089623	-2.45904042	5.32750767
O	N	-4.2979403	-1.90748509	0.92758178
H	Y	-2.4030515	1.62863556	5.70060086

H	Y	-2.40322473	-1.8434943	5.62618695
Si	N	-3.0551116	-1.62881502	1.9247309
Si	Y	-3.56967515	1.45571249	4.90991735
Si	Y	-3.57016841	-1.63679214	4.84441691
O	Y	-1.55012598	0.1001865	-4.03707121
H	Y	-0.83205514	2.04076112	-5.39669405
H	Y	-0.83200527	-1.78136627	-5.47796324
Si	Y	-1.10672211	1.64626795	-4.06060586
Si	Y	-1.10662731	-1.44414405	-4.12677709
O	Y	-4.10861244	2.05920938	-1.63420261
H	Y	-6.23604844	2.3562664	-0.40129725
H	Y	-3.80192249	4.3154997	-2.60604426
O	Y	-2.29781245	2.55559056	-3.47188606
Si	Y	-5.05001638	1.57565187	-0.42173455
Si	Y	-3.08186234	3.15946265	-2.20421844
O	N	-0.09204641	4.30718976	0.76981544
H	Y	0.84800948	6.65984586	0.47432851
O	N	-2.04080813	3.55301797	-1.04343385
H	Y	1.87835899	5.07615283	1.82186557
Al	N	-0.90819473	2.97623614	0.0965137
Si	Y	1.19017294	5.28776905	0.59752054
O	Y	2.13066974	-4.79716569	-0.84075523
H	Y	4.57423927	-4.32500913	1.96088313
H	Y	6.17285647	-4.69269713	0.32195719
O	Y	2.12694716	4.85470321	-0.63782208
H	Y	8.25772138	-1.94823531	0.30473809
H	Y	3.84683623	4.3608912	-2.34896789
H	Y	8.11484252	-1.77730809	2.61526147
H	Y	4.57118703	4.26710063	2.14027751
H	Y	8.25679977	1.96390627	0.38721906
H	Y	6.17053582	4.70429787	0.51880209
H	Y	8.11365097	1.69593537	2.68823671

Table S16 – Coordinates for DFT Model of CH_3 Radical Escape TS (CHA)

element	constrained during optimization?	X	Y	Z
Fe	N	-0.42325988	-0.00001559	0.78301668
O	N	0.07037008	-1.49955078	-0.53626773
O	N	-1.66192762	-1.67371466	1.22678969
O	N	-1.66221319	1.67350103	1.22673924

O	N	0.07005806	1.49909166	-0.53660963
O	N	0.8872166	0.00070105	1.96682918
H	N	1.84940885	0.00087394	1.9371646
C	N	4.38574851	-0.04093813	-0.46803989
H	N	4.61381279	-0.75214552	0.3066797
H	N	4.16423679	-0.38627711	-1.46488546
H	N	4.43038397	1.01579209	-0.26634365
O	N	1.69793385	-0.00068867	-1.98289992
O	N	2.34357013	-2.43048714	-1.74548294
O	N	2.343914	2.42826172	-1.74114246
O	Y	4.16960175	-3.31115635	-0.08555197
O	Y	6.01957309	-2.4344541	1.55976545
O	Y	4.16541055	3.31611505	-0.0880205
O	Y	6.9780231	0.00432698	1.39347318
O	Y	6.01617026	2.44343298	1.55785962
Si	N	1.07457348	-1.47817856	-1.86101801
Si	N	1.07463659	1.47662031	-1.86149577
Si	Y	5.20955258	-3.73594301	1.06672216
Si	Y	3.12459539	3.74092385	-1.2404936
Si	Y	7.36350878	-1.54654415	1.57241944
Si	Y	5.20529745	3.74314059	1.06282743
Si	Y	7.36133708	1.55746949	1.5716297
Si	Y	3.1297947	-3.7381368	-1.23752062
O	N	0.2489546	-1.86281098	-3.18906316
O	N	0.24785918	1.86295843	-3.18862078
O	N	-3.24748519	-0.00032465	2.3161802
O	Y	-4.13172738	-2.01081735	-1.6741517
H	Y	-3.81158911	-4.24576407	-2.69033982
H	Y	-6.27398918	-2.3352755	-0.47639148
O	Y	-2.29627935	-2.46627087	-3.50052881
Si	Y	-3.09769949	-3.09856576	-2.25576495
Si	Y	-5.08883481	-1.55370755	-0.4641359
O	N	-0.11919901	-4.2585798	0.71100524
H	Y	3.83636008	-4.29476267	-2.33597219
H	Y	0.80107032	-6.65415456	0.39747564
O	N	-2.06472869	-3.4864235	-1.08115543
H	Y	1.81226062	-5.09922015	1.79075532
Al	N	-0.95670559	-2.91408118	0.08596859
Si	Y	1.14001836	-5.28410806	0.55418852
O	Y	-5.47363452	-0.00185857	-0.64244171
O	N	-3.30005928	2.26499538	3.41487771
O	N	-4.34412082	1.95012297	0.95315026
Si	N	-3.14350899	1.5620349	1.96622435

O	Y	-4.12026126	-0.00017032	4.9359731
O	N	-3.2985142	-2.26539289	3.4156819
H	Y	-4.73431837	2.35796849	5.37541546
H	Y	-4.73289237	-2.35904104	5.37444448
O	N	-4.34370828	-1.95202246	0.95433226
H	Y	-2.52042522	1.73658368	5.68792352
H	Y	-2.51923399	-1.73642779	5.68792479
Si	N	-3.14299342	-1.56278966	1.96682164
Si	Y	-3.6769556	1.54617279	4.88640902
Si	Y	-3.67623731	-1.54684666	4.88717972
O	Y	-1.54405785	0.00013917	-4.00343507
H	Y	-0.80896327	1.91147916	-5.39484378
H	Y	-0.80746682	-1.91160745	-5.39419586
Si	Y	-1.10076688	1.54527937	-4.05401131
Si	Y	-1.09933316	-1.54533413	-4.05384995
O	Y	-4.13305028	2.00926031	-1.67544292
H	Y	-6.2759356	2.33173517	-0.47616462
H	Y	-3.81428863	4.24427116	-2.69143234
O	Y	-2.29901189	2.46697693	-3.50027951
Si	Y	-5.08960942	1.5512809	-0.46491275
Si	Y	-3.09927452	3.09730103	-2.25590695
O	N	-0.12105639	4.25934637	0.71117871
H	Y	0.79596258	6.6558964	0.39662469
O	N	-2.06664039	3.48524595	-1.08085844
H	Y	1.80933575	5.10184518	1.79074072
Al	N	-0.95733947	2.913915	0.0855779
Si	Y	1.13671906	5.2868504	0.55351477
O	Y	2.0965486	-4.82617665	-0.65637376
H	Y	4.5043855	-4.29320744	2.16522974
H	Y	6.12377153	-4.69509793	0.55490139
O	Y	2.08912112	4.82782558	-0.66025741
H	Y	8.20827114	-1.95093689	0.50535064
H	Y	3.83061606	4.29807312	-2.33859526
H	Y	8.03611645	-1.73060238	2.80968944
H	Y	4.4980411	4.30090263	2.16044671
H	Y	8.20583571	1.96207292	0.5039673
H	Y	6.11782802	4.70373314	0.55033667
H	Y	8.03358444	1.74340341	2.80820261

Table S17 – Coordinates for DFT Model of MeOH-lattice complex (BEA)

element	constrained during optimization?	X	Y	Z
O	N	0.23875467	-1.90385893	-0.51721361
O	N	-1.5198645	-2.21068051	1.45656972
O	N	-1.73065459	2.40695958	1.35224349
O	N	0.07406308	1.99968837	-0.5115548
O	N	0.67012203	-1.12027788	2.05485961
H	N	0.90450173	-1.07159549	1.11086445
C	N	0.95515199	0.10677302	2.7376365
H	N	0.40101404	0.93995993	2.30247832
H	N	0.65668452	-0.03159447	3.77577931
H	N	2.02747897	0.31018569	2.7022753
O	N	1.51961619	0.05592869	-1.62331749
O	N	2.45987781	-2.35198205	-1.93358611
O	N	2.36411832	2.46481477	-1.82512686
O	Y	4.22915551	-3.27591937	-0.21362643
O	Y	6.0964446	-2.4054383	1.41452962
O	Y	4.17450178	3.35017582	-0.13469453
O	Y	7.03393322	0.04269235	1.26415211
O	Y	6.0561786	2.47206994	1.47279119
Si	N	1.0892969	-1.50158883	-1.82660282
Si	N	1.02391303	1.57578114	-1.78341679
Si	Y	5.28943811	-3.70700174	0.91728411
Si	Y	3.11319502	3.78110576	-1.2659976
Si	Y	7.43380664	-1.50719474	1.41830679
Si	Y	5.22818577	3.77127705	1.00590243
Si	Y	7.4080491	1.59647973	1.45575486
Si	Y	3.17530681	-3.69719017	-1.35551653
O	N	0.26091687	-1.70512518	-3.21352424
O	N	0.21916682	1.83384097	-3.16743072
O	N	-2.56433286	0.02387973	2.02492867
O	Y	-4.10469942	-2.02117826	-1.66317382
H	Y	-3.78275988	-4.2410113	-2.71144282
H	Y	-6.22629729	-2.37696199	-0.43789189
O	Y	-2.29330312	-2.43997288	-3.52194237
Si	Y	-3.0710556	-3.0937501	-2.27317183
Si	Y	-5.0469567	-1.5864338	-0.43355791
O	N	-0.17691352	-4.5231512	0.66988404
H	Y	3.86981072	-4.23499799	-2.47092361
H	Y	0.89345936	-6.65094065	0.27775559
O	N	-2.09802868	-3.56244945	-1.09604557

H	Y	1.91360767	-5.10541414	1.67499915
Si	Y	1.22426799	-5.28012351	0.44635398
O	Y	-5.44646508	-0.03550315	-0.58683867
O	N	-3.29714147	2.09980697	3.4469828
O	N	-4.31584489	1.76826999	1.03490861
Si	N	-2.99008699	1.59534208	1.94924238
O	Y	-4.01014734	-0.09138108	4.96999909
O	N	-3.28935471	-2.16909323	3.37432725
H	Y	-4.63524649	2.25641369	5.44789401
H	Y	-4.59800537	-2.46017036	5.38876518
O	N	-4.19880959	-1.84135853	0.95378636
H	Y	-2.41232223	1.64844772	5.71989027
H	Y	-2.38475639	-1.82419115	5.67707486
Si	N	-2.96605004	-1.54127409	1.93314276
Si	Y	-3.57907835	1.4588349	4.93350774
Si	Y	-3.55485396	-1.63386663	4.89609485
O	Y	-1.56736635	0.03813303	-4.0054817
H	Y	-0.86774075	1.97191918	-5.38405116
H	Y	-0.83716278	-1.85066583	-5.43053885
Si	Y	-1.13644996	1.58742508	-4.04372889
Si	Y	-1.11161208	-1.50339116	-4.08181594
O	Y	-4.13655611	1.99843278	-1.6148696
H	Y	-6.26370404	2.28962734	-0.38013403
H	Y	-3.84997813	4.24812845	-2.60784358
O	Y	-2.33356268	2.49254027	-3.46086232
Si	Y	-5.07142104	1.51840698	-0.39608516
Si	Y	-3.11979169	3.10157753	-2.19702625
O	N	-0.24274896	4.41509503	0.64135128
H	Y	0.78736313	6.65765559	0.44106939
O	N	-2.05663759	3.55213682	-1.03672211
H	Y	1.83319845	5.09462406	1.80074402
Si	N	-0.98050381	3.09324406	0.08905518
Si	Y	1.14073013	5.28950495	0.57598897
O	Y	2.15922116	-4.80022698	-0.77250045
H	Y	4.60479882	-4.28297412	2.01950834
H	Y	6.20291051	-4.65276467	0.38056317
O	Y	2.07850708	4.85279502	-0.6573162
H	Y	8.26561914	-1.89195121	0.33388733
H	Y	3.79868734	4.35716721	-2.36762779
H	Y	8.12617099	-1.70111838	2.64306214
H	Y	4.53309132	4.31011366	2.12071506
H	Y	8.23343945	2.02064593	0.38076733
H	Y	6.12550395	4.74531503	0.49189284

H	Y	8.09722727	1.77252501	2.68443015
H	N	-0.69475622	-1.73610267	1.86640164
Al	N	-0.88050228	-3.1437323	0.0021482

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