



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2026 – 07:46 AM EDT

PDB ID : 3FYI / pdb_00003fyi
Title : Catalytic core subunits (I and II) of cytochrome C oxidase from Rhodobacter sphaeroides in the reduced state bound with cyanide
Authors : Qin, L.; Mills, D.A.; Proshlyakov, D.A.; Hiser, C.; Ferguson-Miller, S.
Deposited on : 2009-01-22
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

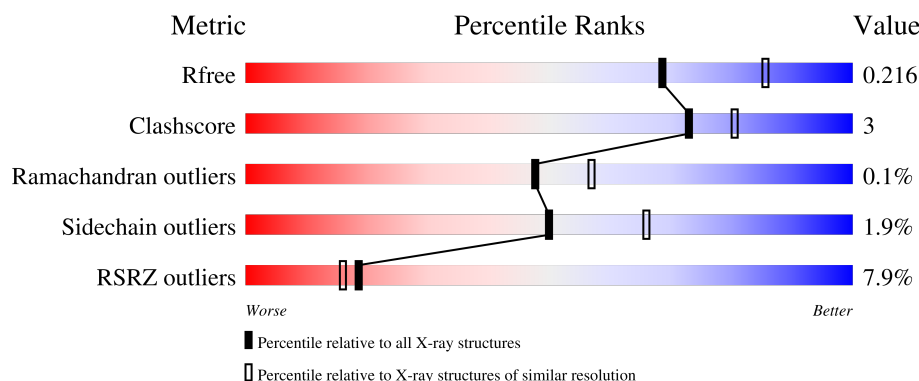
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	566	4% 87% 7% 5%
1	C	566	14% 84% 9% 6%
2	B	262	3% 93% ...
2	D	262	6% 91% 6%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
10	HTO	B	1	X	-	-	-
8	DMU	D	1011	-	-	-	X

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 13601 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 Cytochrome c oxidase subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	535	Total	C	N	O	S	22	0	0
			4212	2822	663	696	31			
1	C	531	Total	C	N	O	S	26	0	0
			4172	2794	656	691	31			

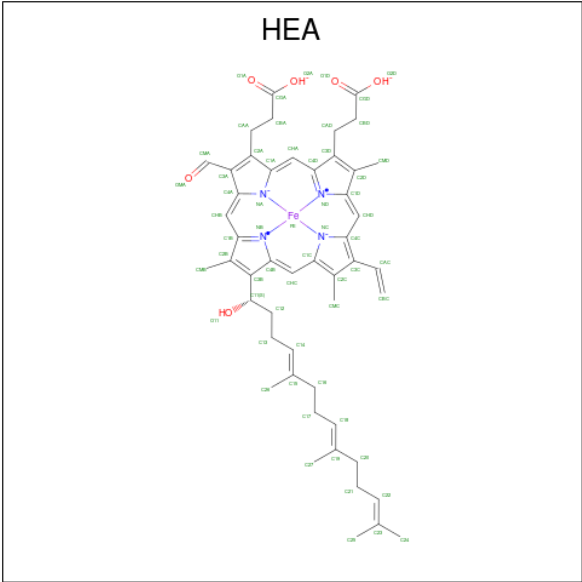
- Molecule 2 is a protein called Cytochrome c oxidase subunit 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	256	Total	C	N	O	S	4	1	0
			2031	1325	333	367	6			
2	D	256	Total	C	N	O	S	9	1	0
			2031	1325	333	367	6			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	282	HIS	-	expression tag	UNP Q03736
B	283	HIS	-	expression tag	UNP Q03736
B	284	HIS	-	expression tag	UNP Q03736
B	285	HIS	-	expression tag	UNP Q03736
B	286	HIS	-	expression tag	UNP Q03736
B	287	HIS	-	expression tag	UNP Q03736
D	282	HIS	-	expression tag	UNP Q03736
D	283	HIS	-	expression tag	UNP Q03736
D	284	HIS	-	expression tag	UNP Q03736
D	285	HIS	-	expression tag	UNP Q03736
D	286	HIS	-	expression tag	UNP Q03736
D	287	HIS	-	expression tag	UNP Q03736

- Molecule 3 is HEME-A (CCD ID: HEA) (formula: $C_{49}H_{56}FeN_4O_6$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	0	0
			60	49	1	4	6		
3	A	1	Total	C	Fe	N	O	0	0
			60	49	1	4	6		
3	C	1	Total	C	Fe	N	O	0	0
			60	49	1	4	6		
3	C	1	Total	C	Fe	N	O	0	0
			60	49	1	4	6		

- Molecule 4 is COPPER (I) ION (CCD ID: CU1) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cu	0	0
			1	1		
4	B	2	Total	Cu	0	0
			2	2		
4	C	1	Total	Cu	0	0
			1	1		
4	D	2	Total	Cu	0	0
			2	2		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		

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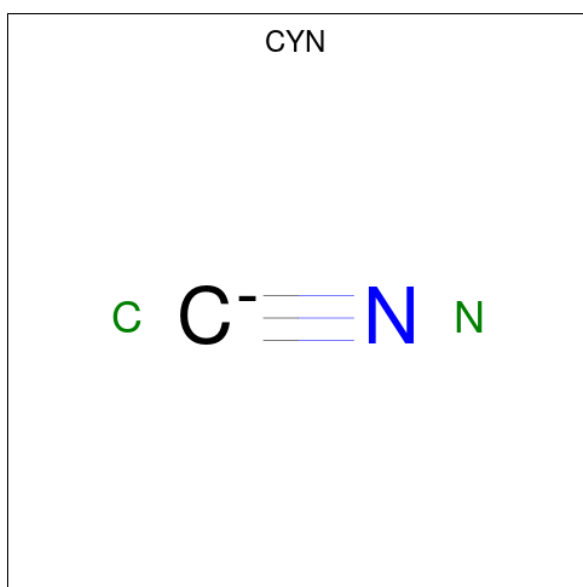
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	1	Total	Mg	0	0
			1	1		

- Molecule 6 is CALCIUM ION (CCD ID: CA) (formula: Ca).

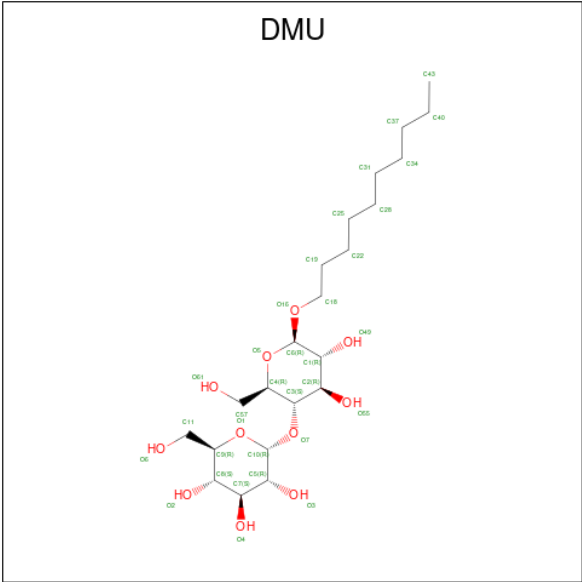
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Ca	0	0
			1	1		
6	C	1	Total	Ca	0	0
			1	1		

- Molecule 7 is CYANIDE ION (CCD ID: CYN) (formula: CN).



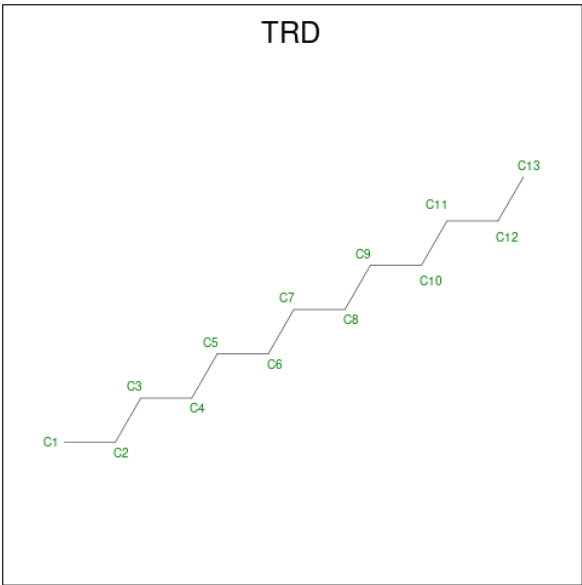
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	N	0	0
			2	1	1		
7	C	1	Total	C	N	0	0
			2	1	1		

- Molecule 8 is DECYL-BETA-D-MALTOPYRANOSIDE (CCD ID: DMU) (formula: C₂₂H₄₂O₁₁).



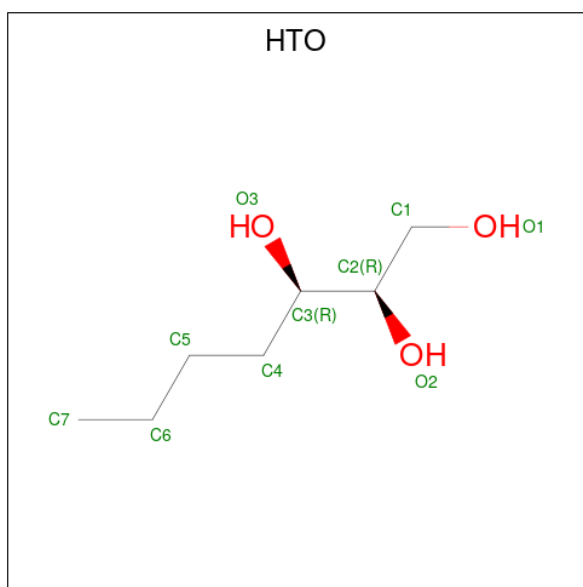
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			33	22	11		
8	A	1	Total	C	O	0	0
			33	22	11		
8	A	1	Total	C	O	0	0
			22	16	6		
8	A	1	Total	C	O	0	0
			33	22	11		
8	B	1	Total	C	O	0	0
			33	22	11		
8	B	1	Total	C	O	0	0
			23	12	11		
8	C	1	Total	C	O	0	0
			23	12	11		
8	C	1	Total	C	O	0	0
			33	22	11		
8	D	1	Total	C	O	0	0
			33	22	11		
8	D	1	Total	C	O	0	0
			23	12	11		

- Molecule 9 is TRIDECANE (CCD ID: TRD) (formula: C₁₃H₂₈).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	1	Total C 13 13	0	0
9	A	1	Total C 13 13	0	0
9	A	1	Total C 7 7	0	0
9	A	1	Total C 7 7	0	0
9	A	1	Total C 13 13	0	0
9	A	1	Total C 13 13	0	0
9	C	1	Total C 13 13	0	0
9	C	1	Total C 13 13	0	0
9	C	1	Total C 7 7	0	0
9	C	1	Total C 9 9	0	0
9	D	1	Total C 13 13	0	0
9	D	1	Total C 7 7	0	0

- Molecule 10 is HEPTANE-1,2,3-TRIOL (CCD ID: HTO) (formula: C₇H₁₆O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	B	1	Total	C	O	0	0
			10	7	3		

- Molecule 11 is CADMIUM ION (CCD ID: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	B	2	Total	Cd	0	0
			2	2		
11	D	2	Total	Cd	0	0
			2	2		

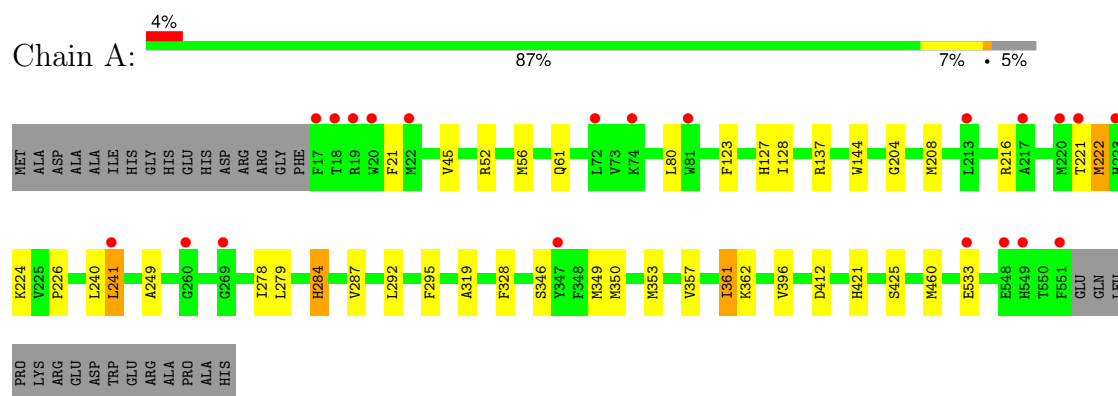
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	126	Total	O	0	0
			126	126		
12	B	133	Total	O	0	0
			133	133		
12	C	102	Total	O	0	0
			102	102		
12	D	109	Total	O	0	0
			109	109		

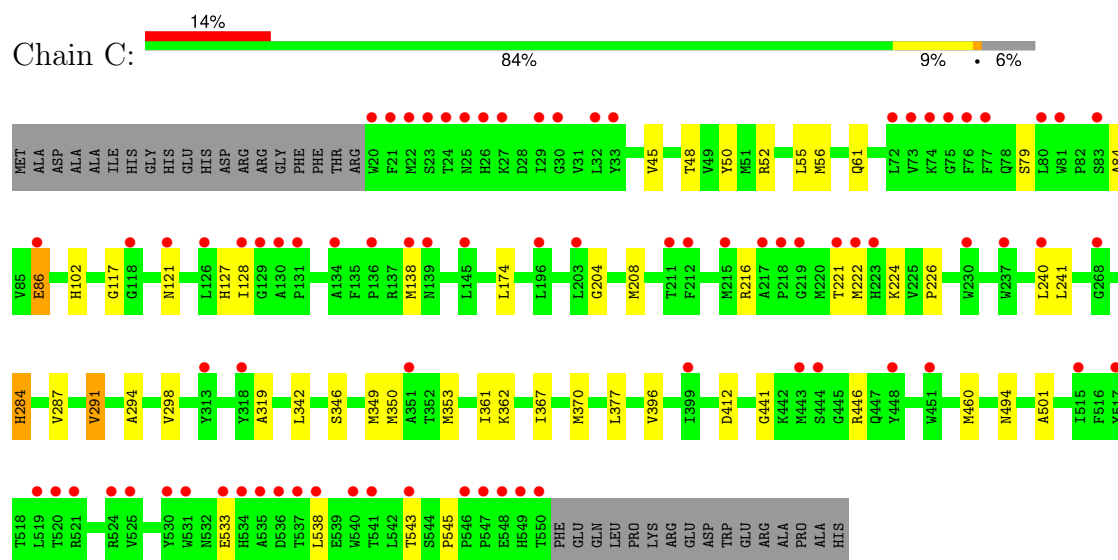
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

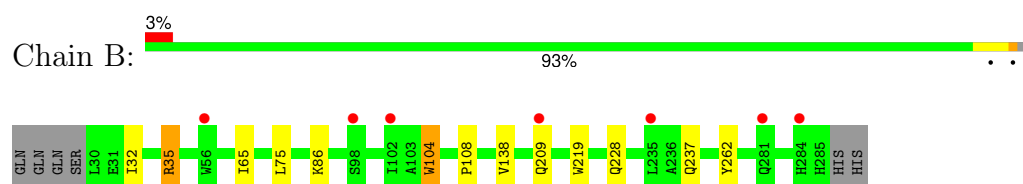
• Molecule 1: Cytochrome c oxidase subunit 1



• Molecule 1: Cytochrome c oxidase subunit 1

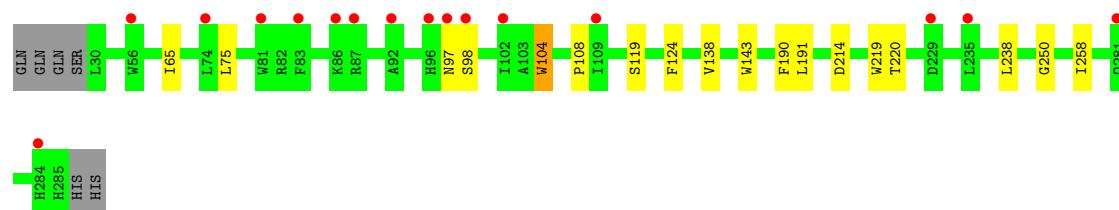


• Molecule 2: Cytochrome c oxidase subunit 2



- Molecule 2: Cytochrome c oxidase subunit 2

Chain D: 6% 91% 6%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	124.34Å 131.88Å 176.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20 50.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.2 (50.00-2.20) 96.2 (50.00-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.194 , 0.219 0.193 , 0.216	Depositor DCC
R_{free} test set	3933 reflections (2.78%)	wwPDB-VP
Wilson B-factor (Å ²)	40.1	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13601	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, CYN, DMU, TRD, CD, CU1, MG, HTO, HEA

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	0.66	0/4368	0.88	2/5961 (0.0%)
1	C	0.56	1/4326 (0.0%)	0.88	2/5905 (0.0%)
2	B	0.62	0/2096	0.78	1/2869 (0.0%)
2	D	0.57	0/2096	0.79	0/2869
All	All	0.60	1/12886 (0.0%)	0.85	5/17604 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	284	HIS	CE1-NE2	5.04	1.37	1.32

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	284	HIS	N-CA-CB	7.13	117.17	110.39
1	A	284	HIS	N-CA-CB	5.64	118.23	110.11
1	C	538	LEU	N-CA-C	5.17	117.66	111.71
2	B	262	TYR	N-CA-C	5.09	117.40	110.88
1	A	123	PHE	N-CA-C	5.00	119.51	113.41

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	4212	0	4134	29	0
1	C	4172	0	4097	32	0
2	B	2031	0	1988	9	0
2	D	2031	0	1988	10	0
3	A	120	0	108	5	0
3	C	120	0	108	8	0
4	A	1	0	0	0	0
4	B	2	0	0	0	0
4	C	1	0	0	0	0
4	D	2	0	0	0	0
5	A	1	0	0	0	0
5	C	1	0	0	0	0
6	A	1	0	0	0	0
6	C	1	0	0	0	0
7	A	2	0	0	0	0
7	C	2	0	0	0	0
8	A	121	0	157	1	0
8	B	56	0	63	1	0
8	C	56	0	63	2	0
8	D	56	0	63	1	0
9	A	66	0	138	0	0
9	C	42	0	86	2	0
9	D	20	0	41	0	0
10	B	10	0	16	0	0
11	B	2	0	0	0	0
11	D	2	0	0	0	0
12	A	126	0	0	0	0
12	B	133	0	0	2	0
12	C	102	0	0	0	0
12	D	109	0	0	0	0
All	All	13601	0	13050	82	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 3.

All (82) 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)
2:B:228:GLN:NE2	12:B:389:HOH:O	1.96	0.95
2:B:32:ILE:HG22	2:B:35:ARG:HD3	1.58	0.86
1:A:21:PHE:HB3	1:A:144:TRP:HZ2	1.45	0.81
1:C:349:MET:HG2	1:C:353:MET:HE2	1.73	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:MET:HE1	8:A:576:DMU:H7	1.76	0.67
1:A:21:PHE:HB3	1:A:144:TRP:CZ2	2.29	0.66
1:C:56:MET:HE1	8:C:1005:DMU:H7	1.76	0.66
1:A:216:ARG:CZ	1:A:222:MET:HE3	2.26	0.65
1:C:84:ALA:HB1	1:C:86:GLU:OE1	1.95	0.65
1:A:349:MET:HG2	1:A:353:MET:HE2	1.79	0.64
1:C:284:HIS:CD2	1:C:284:HIS:C	2.74	0.64
3:C:568:HEA:HMD1	3:C:568:HEA:HBD2	1.79	0.64
1:A:396:VAL:HB	2:B:65:ILE:HB	1.80	0.63
1:C:50:TYR:OH	1:C:79:SER:HB2	2.00	0.62
1:C:204:GLY:O	1:C:208:MET:HG2	1.99	0.62
1:C:350:MET:HA	1:C:353:MET:HE3	1.81	0.61
3:C:567:HEA:HBC1	3:C:567:HEA:HMC3	1.83	0.60
1:A:357:VAL:O	1:A:361:ILE:HG12	2.02	0.59
1:C:396:VAL:HB	2:D:65:ILE:HB	1.84	0.59
1:A:350:MET:HA	1:A:353:MET:HE3	1.86	0.57
1:A:221:THR:CG2	1:A:224:LYS:HG2	2.34	0.57
1:A:284:HIS:C	1:A:284:HIS:CD2	2.83	0.57
1:A:204:GLY:O	1:A:208:MET:HG2	2.07	0.55
1:C:86:GLU:H	1:C:86:GLU:CD	2.16	0.54
1:C:361:ILE:HD11	2:D:108:PRO:HG2	1.90	0.54
1:A:287:VAL:HB	3:A:568:HEA:HAC	1.90	0.53
1:C:287:VAL:HB	3:C:568:HEA:HAC	1.91	0.52
1:A:533:GLU:H	1:A:533:GLU:CD	2.19	0.51
1:A:361:ILE:HD11	2:B:108:PRO:HG2	1.93	0.50
3:A:568:HEA:HBD2	3:A:568:HEA:HMD1	1.93	0.50
1:A:127:HIS:HB3	1:A:226:PRO:HG2	1.93	0.49
1:C:284:HIS:O	1:C:287:VAL:HG22	2.12	0.49
1:A:279:LEU:C	1:A:279:LEU:HD13	2.38	0.49
1:C:221:THR:HG22	1:C:224:LYS:HG3	1.94	0.49
2:B:35:ARG:HD2	12:B:343:HOH:O	2.12	0.49
1:C:128:ILE:HB	1:C:216:ARG:HG2	1.95	0.48
1:A:221:THR:HG22	1:A:224:LYS:CG	2.43	0.48
1:A:221:THR:HG23	1:A:224:LYS:HG2	1.95	0.48
1:A:221:THR:HG22	1:A:224:LYS:HG3	1.96	0.47
1:C:48:THR:CG2	1:C:102:HIS:CE1	2.97	0.46
3:C:568:HEA:H122	3:C:568:HEA:HHC	1.97	0.46
1:A:292:LEU:O	1:A:295:PHE:HB2	2.16	0.46
1:A:460:MET:HE1	3:A:567:HEA:C14	2.46	0.46
1:C:342:LEU:HD21	2:D:124:PHE:CD2	2.50	0.45
1:C:319:ALA:HB3	1:C:362:LYS:HE2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:55:LEU:O	1:C:494:ASN:HB3	2.16	0.45
1:C:52:ARG:HG3	1:C:501:ALA:CB	2.47	0.45
1:C:441:GLY:HA2	1:C:446:ARG:O	2.16	0.45
1:A:249:ALA:HB2	1:A:278:ILE:HG22	1.99	0.45
1:C:367:ILE:HA	1:C:370:MET:HE2	1.99	0.45
1:A:346:SER:O	1:A:350:MET:HG2	2.17	0.44
2:D:104:TRP:CD1	2:D:104:TRP:C	2.96	0.44
1:A:128:ILE:HB	1:A:216:ARG:HG2	2.00	0.44
1:C:543:THR:HG23	1:C:545:PRO:O	2.18	0.44
2:D:75:LEU:HD13	8:D:1003:DMU:H18	2.00	0.43
1:A:319:ALA:HB3	1:A:362:LYS:HE2	1.99	0.43
3:C:567:HEA:HBC1	3:C:567:HEA:CMC	2.48	0.43
1:A:421:HIS:HA	1:A:425:SER:HB2	2.00	0.43
1:C:127:HIS:HB3	1:C:226:PRO:HG2	2.01	0.43
1:C:291:VAL:HB	3:C:568:HEA:HBC2	2.00	0.43
1:C:377:LEU:HD13	9:C:1010:TRD:H31	2.00	0.43
1:C:367:ILE:HA	1:C:370:MET:CE	2.49	0.43
2:B:75:LEU:HD13	8:B:288:DMU:H18	2.01	0.42
2:D:143:TRP:CH2	2:D:258:ILE:HG22	2.53	0.42
2:D:220:THR:O	2:D:250:GLY:HA3	2.20	0.42
1:A:241:LEU:HB3	1:A:328:PHE:CZ	2.55	0.42
1:A:45:VAL:HG21	3:A:567:HEA:H171	2.01	0.42
1:C:460:MET:HE1	3:C:567:HEA:C14	2.49	0.42
1:A:221:THR:CG2	1:A:224:LYS:CG	2.97	0.42
1:C:346:SER:O	1:C:350:MET:HG2	2.20	0.42
2:D:190:PHE:O	2:D:191:LEU:HB2	2.20	0.42
2:B:138:VAL:HG11	2:B:219:TRP:CD1	2.55	0.42
1:C:294:ALA:O	1:C:298:VAL:HG23	2.20	0.41
2:B:104:TRP:CD1	2:B:104:TRP:C	2.98	0.41
2:B:209:GLN:HG2	2:B:237:GLN:HG2	2.02	0.41
2:D:138:VAL:HG11	2:D:219:TRP:CD1	2.56	0.41
1:C:117:GLY:O	1:C:121:ASN:HB2	2.20	0.41
2:D:238:LEU:C	2:D:238:LEU:HD12	2.46	0.41
1:C:45:VAL:HG21	3:C:567:HEA:H171	2.02	0.41
1:C:533:GLU:H	1:C:533:GLU:CD	2.29	0.40
8:C:1005:DMU:H9	9:C:1006:TRD:H81	2.04	0.40
3:A:568:HEA:H122	3:A:568:HEA:HHC	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	533/566 (94%)	523 (98%)	10 (2%)	0	100	100
1	C	529/566 (94%)	524 (99%)	5 (1%)	0	100	100
2	B	255/262 (97%)	249 (98%)	6 (2%)	0	100	100
2	D	255/262 (97%)	249 (98%)	5 (2%)	1 (0%)	30	34
All	All	1572/1656 (95%)	1545 (98%)	26 (2%)	1 (0%)	48	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	97	ASN

5.3.2 Protein sidechains

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	435/459 (95%)	426 (98%)	9 (2%)	47	63
1	C	431/459 (94%)	422 (98%)	9 (2%)	47	63
2	B	216/221 (98%)	213 (99%)	3 (1%)	59	75
2	D	216/221 (98%)	212 (98%)	4 (2%)	50	66
All	All	1298/1360 (95%)	1273 (98%)	25 (2%)	50	66

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	ARG
1	A	61	GLN
1	A	80	LEU
1	A	137	ARG
1	A	222	MET
1	A	240	LEU
1	A	241	LEU
1	A	361	ILE
1	A	412	ASP
2	B	35	ARG
2	B	86	LYS
2	B	104	TRP
1	C	61	GLN
1	C	86	GLU
1	C	138	MET
1	C	174	LEU
1	C	222	MET
1	C	240	LEU
1	C	241	LEU
1	C	291	VAL
1	C	412	ASP
2	D	98	SER
2	D	104	TRP
2	D	119	SER
2	D	214	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	ASN
1	A	139	ASN
1	A	163	ASN
1	A	165	GLN
1	A	276	GLN
1	A	465	ASN
2	B	96	HIS
2	B	170	ASN
2	B	178	GLN
1	C	163	ASN
1	C	344	GLN
1	C	465	ASN
2	D	170	ASN
2	D	251	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 43 ligands modelled in this entry, 14 are monoatomic - leaving 29 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	HEA	C	567	1	67,67,67	1.09	7 (10%)	81,103,103	0.99	3 (3%)
7	CYN	A	572	4	1,1,1	0.28	0	-		
8	DMU	C	1002	-	24,24,34	0.52	0	35,35,45	0.66	1 (2%)
9	TRD	D	1007	-	12,12,12	0.23	0	11,11,11	0.59	0
9	TRD	A	581	-	12,12,12	0.27	0	11,11,11	0.54	0
9	TRD	D	1008	-	6,6,12	0.26	0	5,5,11	0.42	0
8	DMU	A	574	-	34,34,34	0.58	1 (2%)	45,45,45	0.85	2 (4%)
9	TRD	A	580	-	6,6,12	0.25	0	5,5,11	0.40	0
8	DMU	B	288	-	34,34,34	0.54	0	45,45,45	0.72	0
10	HTO	B	1	-	9,9,9	0.37	0	10,10,10	0.94	1 (10%)
9	TRD	C	1009	-	6,6,12	0.27	0	5,5,11	0.35	0
3	HEA	A	567	1	67,67,67	1.08	5 (7%)	81,103,103	1.05	6 (7%)
8	DMU	B	11	-	24,24,34	0.50	0	35,35,45	0.71	0
8	DMU	A	575	-	22,22,34	0.60	1 (4%)	27,27,45	0.86	1 (3%)
9	TRD	A	579	-	6,6,12	0.26	0	5,5,11	0.37	0
8	DMU	C	1005	-	34,34,34	0.60	1 (2%)	45,45,45	0.85	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	TRD	C	1006	-	12,12,12	0.27	0	11,11,11	0.49	0
9	TRD	A	582	-	12,12,12	0.29	0	11,11,11	0.51	0
3	HEA	A	568	1	67,67,67	1.74	8 (11%)	81,103,103	1.89	25 (30%)
8	DMU	A	576	-	34,34,34	0.59	0	45,45,45	0.74	0
8	DMU	D	1011	-	24,24,34	0.51	0	35,35,45	0.85	0
9	TRD	C	1010	-	8,8,12	0.27	0	7,7,11	0.40	0
7	CYN	C	572	4	1,1,1	0.22	0	-		
9	TRD	A	578	-	12,12,12	0.28	0	11,11,11	0.52	0
3	HEA	C	568	1	67,67,67	1.86	8 (11%)	81,103,103	1.91	24 (29%)
8	DMU	D	1003	-	34,34,34	0.53	0	45,45,45	0.62	0
8	DMU	A	573	-	34,34,34	0.50	0	45,45,45	0.96	3 (6%)
9	TRD	A	577	-	12,12,12	0.25	0	11,11,11	0.48	0
9	TRD	C	1001	-	12,12,12	0.25	0	11,11,11	0.52	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	HEA	C	567	1	-	2/36/76/76	-
8	DMU	C	1002	-	-	3/8/48/59	0/2/2/2
9	TRD	D	1007	-	-	6/10/10/10	-
9	TRD	A	581	-	-	6/10/10/10	-
9	TRD	D	1008	-	-	1/4/4/10	-
8	DMU	A	574	-	-	9/19/59/59	0/2/2/2
9	TRD	A	580	-	-	2/4/4/10	-
8	DMU	B	288	-	-	5/19/59/59	0/2/2/2
10	HTO	B	1	-	1/1/2/2	6/10/10/10	-
9	TRD	C	1009	-	-	3/4/4/10	-
3	HEA	A	567	1	-	3/36/76/76	-
8	DMU	B	11	-	-	1/8/48/59	0/2/2/2
8	DMU	A	575	-	-	4/13/33/59	0/1/1/2
9	TRD	A	579	-	-	2/4/4/10	-
8	DMU	C	1005	-	-	11/19/59/59	0/2/2/2
9	TRD	C	1006	-	-	0/10/10/10	-
9	TRD	A	582	-	-	4/10/10/10	-
3	HEA	A	568	1	-	10/36/76/76	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	DMU	A	576	-	-	11/19/59/59	0/2/2/2
8	DMU	D	1011	-	-	4/8/48/59	0/2/2/2
9	TRD	C	1010	-	-	3/6/6/10	-
9	TRD	A	578	-	-	5/10/10/10	-
3	HEA	C	568	1	-	10/36/76/76	-
8	DMU	D	1003	-	-	6/19/59/59	0/2/2/2
8	DMU	A	573	-	-	2/19/59/59	0/2/2/2
9	TRD	A	577	-	-	5/10/10/10	-
9	TRD	C	1001	-	-	2/10/10/10	-

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	568	HEA	FE-NC	8.04	2.21	1.95
3	C	568	HEA	FE-NB	7.94	2.19	1.94
3	A	568	HEA	FE-NA	6.34	2.16	1.95
3	A	568	HEA	FE-NC	6.01	2.15	1.95
3	A	568	HEA	FE-NB	5.77	2.12	1.94
3	C	568	HEA	C11-C3B	4.92	1.57	1.51
3	C	568	HEA	FE-NA	4.57	2.10	1.95
3	A	568	HEA	C11-C3B	4.36	1.56	1.51
3	A	568	HEA	FE-ND	4.30	2.08	1.94
3	C	567	HEA	C11-C3B	3.24	1.55	1.51
3	A	567	HEA	C11-C3B	3.03	1.55	1.51
3	A	567	HEA	CAC-C3C	2.99	1.55	1.47
3	C	568	HEA	FE-ND	2.99	2.04	1.94
3	C	567	HEA	CAC-C3C	2.77	1.54	1.47
3	C	567	HEA	CMA-C3A	2.69	1.51	1.45
3	A	567	HEA	CMA-C3A	2.62	1.51	1.45
3	C	568	HEA	CMD-C2D	2.34	1.55	1.50
3	A	568	HEA	CMA-C3A	2.34	1.50	1.45
3	A	567	HEA	CMD-C2D	2.31	1.55	1.50
3	A	568	HEA	CMD-C2D	2.27	1.55	1.50
3	C	567	HEA	CMD-C2D	2.27	1.55	1.50
3	C	568	HEA	CMA-C3A	2.24	1.50	1.45
8	A	575	DMU	O16-C6	2.21	1.43	1.40
3	C	567	HEA	CMB-C2B	2.20	1.55	1.50
8	A	574	DMU	O16-C6	2.16	1.43	1.40
8	C	1005	DMU	O16-C6	2.13	1.43	1.40
3	C	567	HEA	CMC-C2C	2.13	1.55	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	567	HEA	CMB-C2B	2.10	1.55	1.50
3	A	568	HEA	CMC-C2C	2.08	1.55	1.50
3	C	567	HEA	C3C-C2C	-2.06	1.34	1.41
3	C	568	HEA	CAC-C3C	2.05	1.52	1.47

All (66) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	568	HEA	C3C-C4C-NC	-6.23	104.56	109.80
3	A	568	HEA	C3C-C4C-NC	-6.22	104.56	109.80
3	C	568	HEA	C2A-C1A-NA	-4.60	105.88	110.32
3	A	568	HEA	C2A-C1A-NA	-4.56	105.93	110.32
3	A	568	HEA	C4C-NC-C1C	3.95	112.26	105.82
3	C	568	HEA	C4C-NC-C1C	3.85	112.09	105.82
3	C	568	HEA	C1D-ND-C4D	3.76	109.66	105.21
3	C	568	HEA	C4B-NB-C1B	3.56	109.42	105.21
3	A	568	HEA	C4B-NB-C1B	3.49	109.33	105.21
3	A	568	HEA	C1D-ND-C4D	3.35	109.17	105.21
3	A	568	HEA	C4A-NA-C1A	3.28	111.17	105.82
3	C	568	HEA	C4A-NA-C1A	3.27	111.15	105.82
3	C	568	HEA	C2B-C1B-NB	-3.25	106.15	109.90
3	C	568	HEA	C17-C18-C19	-3.19	120.32	127.62
3	C	568	HEA	C2D-C1D-ND	-3.18	106.19	109.84
8	A	574	DMU	O16-C6-C1	3.09	112.97	108.27
3	C	568	HEA	C3D-C4D-ND	-3.00	107.45	110.35
3	A	568	HEA	C26-C15-C16	2.99	120.43	115.23
8	A	575	DMU	O16-C6-C1	2.89	112.66	108.27
3	A	568	HEA	C3D-C4D-ND	-2.86	107.59	110.35
3	A	568	HEA	C2B-C1B-NB	-2.83	106.63	109.90
3	A	567	HEA	C26-C15-C16	2.83	120.14	115.23
3	A	567	HEA	C17-C18-C19	-2.79	121.23	127.62
3	A	568	HEA	C2D-C1D-ND	-2.77	106.65	109.84
3	C	568	HEA	C2C-C1C-NC	-2.77	105.71	110.14
3	A	568	HEA	CMB-C2B-C3B	-2.75	124.96	130.28
3	C	568	HEA	C26-C15-C16	2.67	119.85	115.23
3	A	568	HEA	C2C-C1C-NC	-2.65	105.90	110.14
3	A	568	HEA	CBA-CAA-C2A	-2.64	105.24	112.53
3	A	568	HEA	OMA-CMA-C3A	-2.62	119.71	125.62
8	A	573	DMU	C10-O7-C3	-2.62	111.77	117.98
3	A	568	HEA	C4A-C3A-C2A	2.57	109.04	106.81
3	A	568	HEA	C17-C18-C19	-2.54	121.82	127.62
3	C	568	HEA	CMB-C2B-C3B	-2.53	125.39	130.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	568	HEA	OMA-CMA-C3A	-2.48	120.02	125.62
3	A	567	HEA	C13-C12-C11	-2.46	110.47	114.39
3	C	568	HEA	C4A-C3A-C2A	2.46	108.95	106.81
3	C	568	HEA	C1B-C2B-C3B	2.44	109.62	106.80
3	C	568	HEA	CBA-CAA-C2A	-2.40	105.91	112.53
3	A	568	HEA	C3B-C4B-NB	-2.32	107.17	109.84
3	C	568	HEA	C25-C23-C24	2.32	119.92	114.59
3	C	568	HEA	CHA-C1A-C2A	2.31	128.51	124.86
3	A	567	HEA	O2D-CGD-CBD	2.31	121.29	114.00
10	B	1	HTO	C5-C4-C3	-2.29	110.35	114.11
3	C	568	HEA	CAD-C3D-C4D	-2.29	120.71	124.70
8	A	573	DMU	O2-C8-C9	2.28	114.95	109.32
3	C	568	HEA	C27-C19-C20	2.28	119.19	115.23
3	A	568	HEA	CAD-C3D-C4D	-2.26	120.75	124.70
8	A	574	DMU	C10-O7-C3	-2.25	112.64	117.98
3	A	567	HEA	C27-C19-C20	2.24	119.12	115.23
3	A	568	HEA	O11-C11-C3B	-2.17	107.29	111.26
3	A	568	HEA	CHA-C1A-C2A	2.16	128.28	124.86
3	A	567	HEA	O2A-CGA-CBA	2.16	120.82	114.00
3	C	567	HEA	C25-C23-C24	2.11	119.45	114.59
3	A	568	HEA	C27-C19-C20	2.10	118.87	115.23
3	C	568	HEA	C3B-C4B-NB	-2.10	107.43	109.84
3	C	568	HEA	CHB-C1B-C2B	2.08	128.32	125.03
3	C	567	HEA	C13-C14-C15	-2.08	122.87	127.62
3	A	568	HEA	CBD-CAD-C3D	2.07	118.27	112.53
3	A	568	HEA	CHA-C4D-ND	2.07	126.65	124.42
3	C	567	HEA	C27-C19-C20	2.06	118.81	115.23
3	A	568	HEA	CBC-CAC-C3C	-2.06	117.25	127.53
3	C	568	HEA	CHA-C4D-ND	2.05	126.63	124.42
8	C	1002	DMU	C10-O7-C3	-2.01	113.20	117.98
3	A	568	HEA	CHB-C4A-C3A	2.01	128.61	125.21
8	A	573	DMU	C7-C8-C9	-2.00	106.60	110.23

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
10	B	1	HTO	C2

All (126) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	568	HEA	C2A-C3A-CMA-OMA

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Mol	Chain	Res	Type	Atoms
3	A	568	HEA	C2D-C3D-CAD-CBD
3	C	568	HEA	C2D-C3D-CAD-CBD
8	A	574	DMU	C19-C18-O16-C6
10	B	1	HTO	O1-C1-C2-O2
10	B	1	HTO	O1-C1-C2-C3
10	B	1	HTO	C2-C3-C4-C5
10	B	1	HTO	O3-C3-C4-C5
8	D	1003	DMU	O6-C11-C9-O1
3	A	568	HEA	C4D-C3D-CAD-CBD
3	C	568	HEA	C4D-C3D-CAD-CBD
8	C	1002	DMU	O6-C11-C9-O1
8	D	1003	DMU	O6-C11-C9-C8
8	A	574	DMU	O5-C6-O16-C18
8	A	575	DMU	O5-C6-O16-C18
8	A	574	DMU	C1-C6-O16-C18
8	A	575	DMU	C1-C6-O16-C18
8	C	1005	DMU	O1-C10-O7-C3
8	D	1011	DMU	O5-C4-C57-O61
8	D	1011	DMU	O6-C11-C9-O1
8	A	576	DMU	O1-C10-O7-C3
9	A	578	TRD	C11-C10-C9-C8
8	A	576	DMU	C19-C18-O16-C6
8	C	1005	DMU	C19-C18-O16-C6
8	A	575	DMU	C19-C22-C25-C28
9	C	1009	TRD	C2-C3-C4-C5
8	B	288	DMU	C22-C25-C28-C31
3	C	568	HEA	C19-C20-C21-C22
8	A	574	DMU	C25-C28-C31-C34
8	A	573	DMU	C31-C34-C37-C40
9	D	1007	TRD	C11-C10-C9-C8
8	D	1003	DMU	C22-C25-C28-C31
8	B	11	DMU	O6-C11-C9-O1
9	A	577	TRD	C2-C3-C4-C5
9	C	1009	TRD	C3-C4-C5-C6
8	A	575	DMU	C28-C31-C34-C37
9	A	578	TRD	C4-C5-C6-C7
9	C	1010	TRD	C3-C4-C5-C6
8	C	1005	DMU	C18-C19-C22-C25
9	D	1007	TRD	C6-C7-C8-C9
8	A	576	DMU	C3-C4-C57-O61
9	C	1001	TRD	C5-C6-C7-C8
8	C	1002	DMU	O6-C11-C9-C8

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Mol	Chain	Res	Type	Atoms
9	D	1008	TRD	C2-C3-C4-C5
9	A	579	TRD	C3-C4-C5-C6
9	D	1007	TRD	C4-C5-C6-C7
8	A	576	DMU	C18-C19-C22-C25
8	B	288	DMU	C25-C28-C31-C34
9	A	581	TRD	C2-C3-C4-C5
8	D	1003	DMU	C25-C28-C31-C34
9	A	579	TRD	C2-C3-C4-C5
9	A	582	TRD	C2-C3-C4-C5
9	D	1007	TRD	C5-C6-C7-C8
3	C	568	HEA	C2A-C3A-CMA-OMA
8	C	1005	DMU	C31-C34-C37-C40
8	A	576	DMU	C2-C3-O7-C10
8	D	1003	DMU	C19-C18-O16-C6
8	A	574	DMU	C34-C37-C40-C43
9	A	577	TRD	C11-C10-C9-C8
8	A	576	DMU	C4-C3-O7-C10
9	D	1007	TRD	C10-C11-C12-C13
9	C	1010	TRD	C4-C5-C6-C7
8	B	288	DMU	C31-C34-C37-C40
9	A	582	TRD	C10-C11-C12-C13
9	C	1009	TRD	C4-C5-C6-C7
9	A	577	TRD	C1-C2-C3-C4
8	C	1005	DMU	C28-C31-C34-C37
8	A	574	DMU	C18-C19-C22-C25
3	A	568	HEA	C19-C20-C21-C22
9	A	581	TRD	C1-C2-C3-C4
9	A	581	TRD	C5-C6-C7-C8
8	D	1003	DMU	C28-C31-C34-C37
8	C	1005	DMU	C19-C22-C25-C28
9	C	1001	TRD	C4-C5-C6-C7
9	A	581	TRD	C4-C5-C6-C7
3	A	568	HEA	C4A-C3A-CMA-OMA
3	C	568	HEA	C4A-C3A-CMA-OMA
8	D	1011	DMU	C3-C4-C57-O61
8	C	1005	DMU	C3-C4-C57-O61
9	A	578	TRD	C5-C6-C7-C8
8	A	574	DMU	C28-C31-C34-C37
8	A	574	DMU	C19-C22-C25-C28
8	A	574	DMU	O6-C11-C9-C8
8	A	576	DMU	C22-C25-C28-C31
8	A	576	DMU	C28-C31-C34-C37

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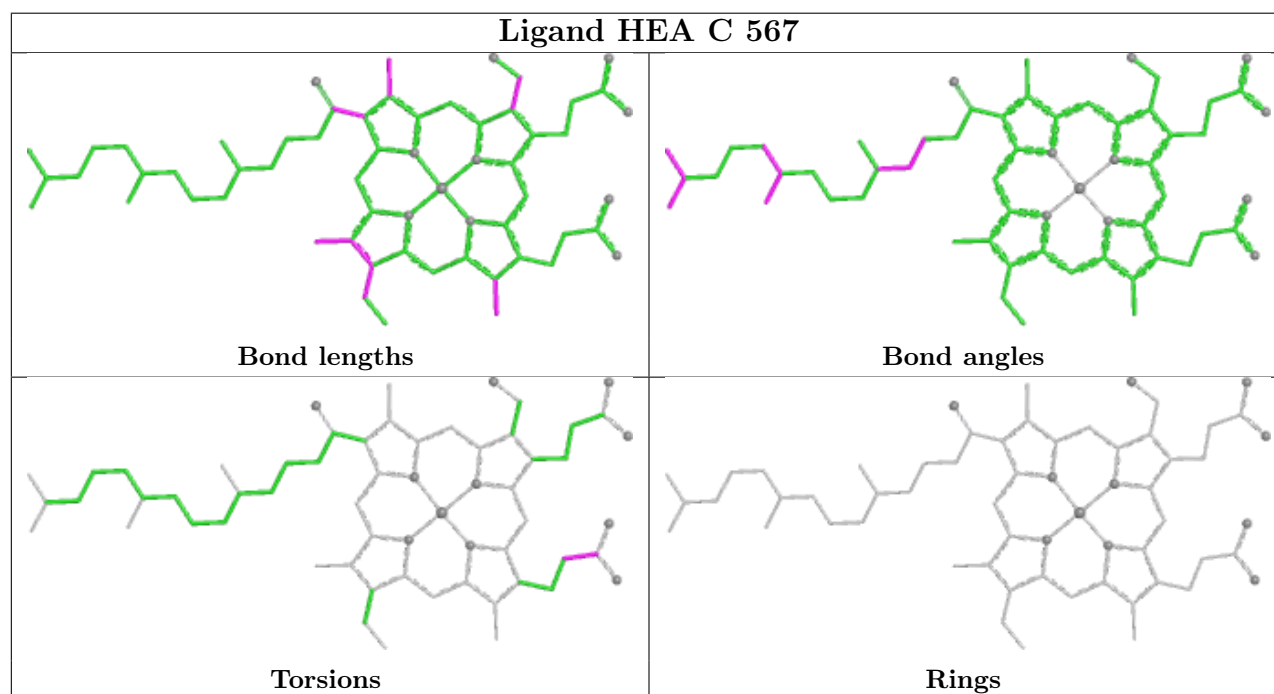
Mol	Chain	Res	Type	Atoms
9	A	577	TRD	C4-C5-C6-C7
8	C	1002	DMU	C3-C4-C57-O61
9	A	582	TRD	C1-C2-C3-C4
9	A	577	TRD	C9-C10-C11-C12
9	A	578	TRD	C6-C7-C8-C9
8	C	1005	DMU	C2-C3-O7-C10
8	B	288	DMU	O6-C11-C9-C8
3	C	567	HEA	CAD-CBD-CGD-O1D
8	C	1005	DMU	C4-C3-O7-C10
8	A	576	DMU	C19-C22-C25-C28
3	A	568	HEA	CAA-CBA-CGA-O1A
3	C	567	HEA	CAD-CBD-CGD-O2D
8	C	1005	DMU	C25-C28-C31-C34
3	C	568	HEA	C2C-C3C-CAC-CBC
9	A	580	TRD	C4-C5-C6-C7
3	C	568	HEA	CAA-CBA-CGA-O2A
8	A	576	DMU	O5-C4-C57-O61
3	A	567	HEA	CAD-CBD-CGD-O2D
3	C	568	HEA	CAA-CBA-CGA-O1A
8	A	573	DMU	C19-C22-C25-C28
3	A	568	HEA	C4C-C3C-CAC-CBC
3	C	568	HEA	C4C-C3C-CAC-CBC
3	A	567	HEA	CAD-CBD-CGD-O1D
10	B	1	HTO	O2-C2-C3-O3
8	B	288	DMU	C3-C4-C57-O61
3	A	568	HEA	CAA-CBA-CGA-O2A
9	A	581	TRD	C6-C7-C8-C9
9	A	578	TRD	C2-C3-C4-C5
9	C	1010	TRD	C5-C6-C7-C8
9	D	1007	TRD	C2-C3-C4-C5
10	B	1	HTO	C1-C2-C3-O3
8	C	1005	DMU	C5-C10-O7-C3
8	A	576	DMU	C5-C10-O7-C3
3	A	568	HEA	CAD-CBD-CGD-O2D
9	A	581	TRD	C9-C10-C11-C12
9	A	580	TRD	C1-C2-C3-C4
3	C	568	HEA	C16-C17-C18-C19
8	D	1011	DMU	O6-C11-C9-C8
3	A	568	HEA	CAD-CBD-CGD-O1D
9	A	582	TRD	C7-C8-C9-C10
3	A	567	HEA	CAA-CBA-CGA-O1A

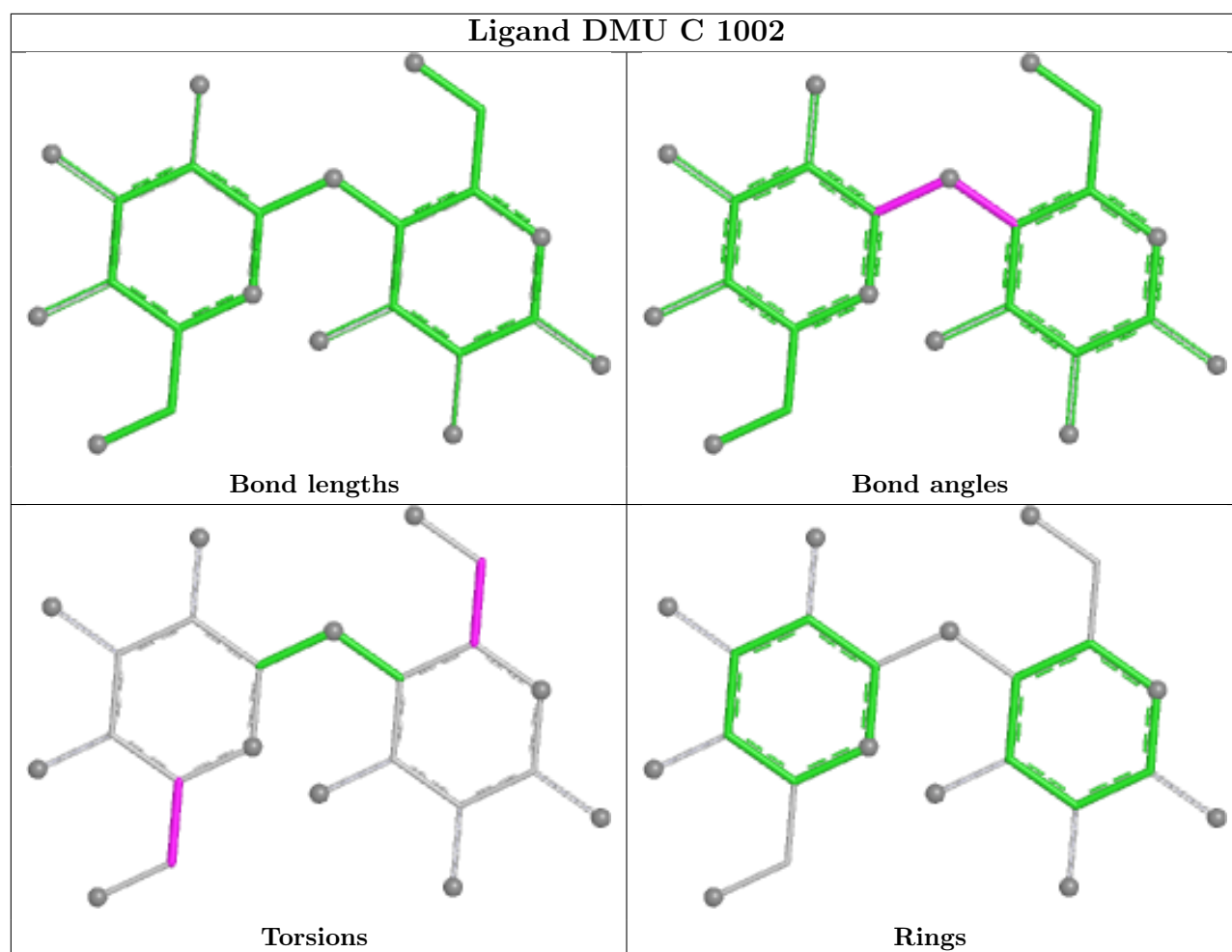
There are no ring outliers.

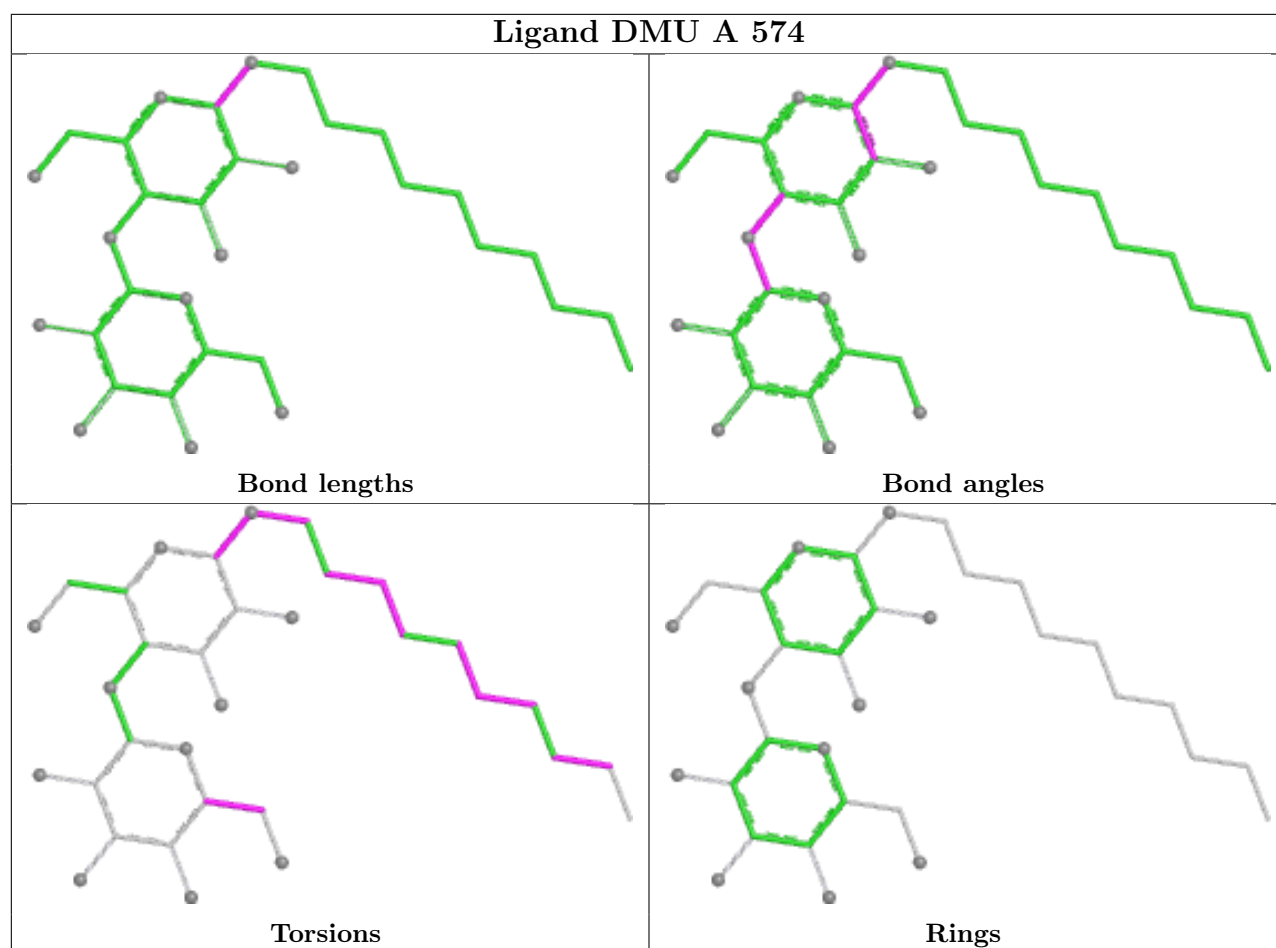
10 monomers are involved in 19 short contacts:

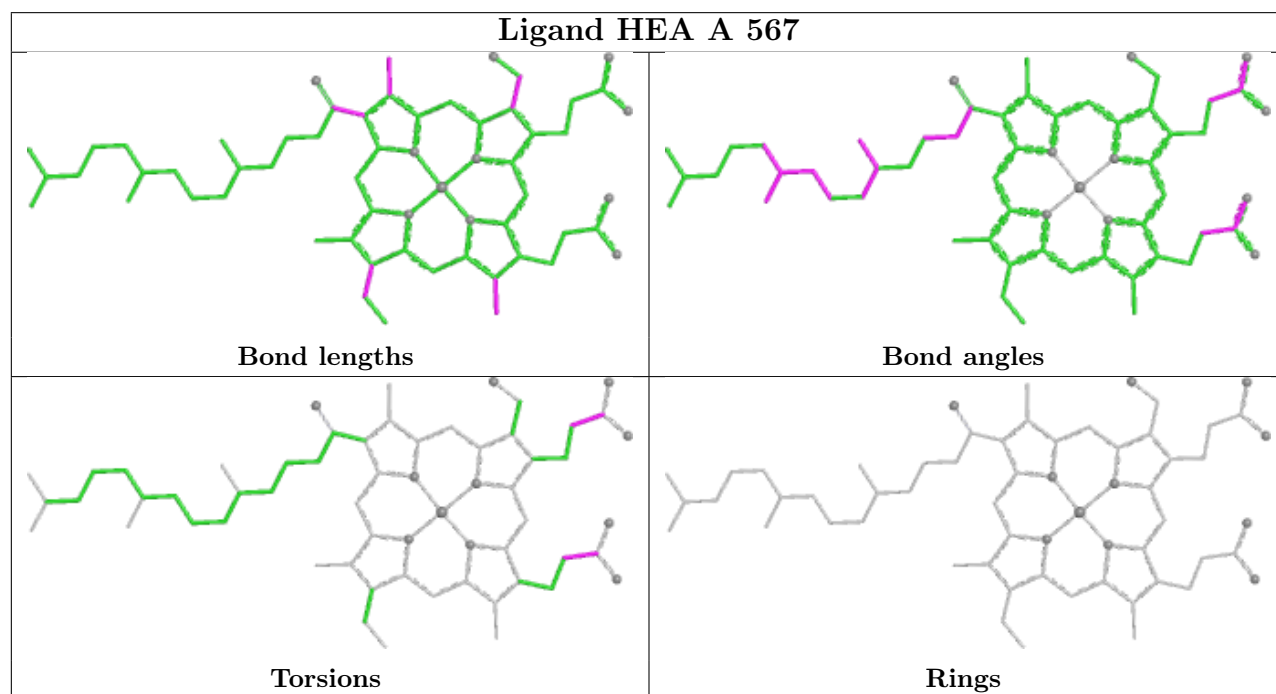
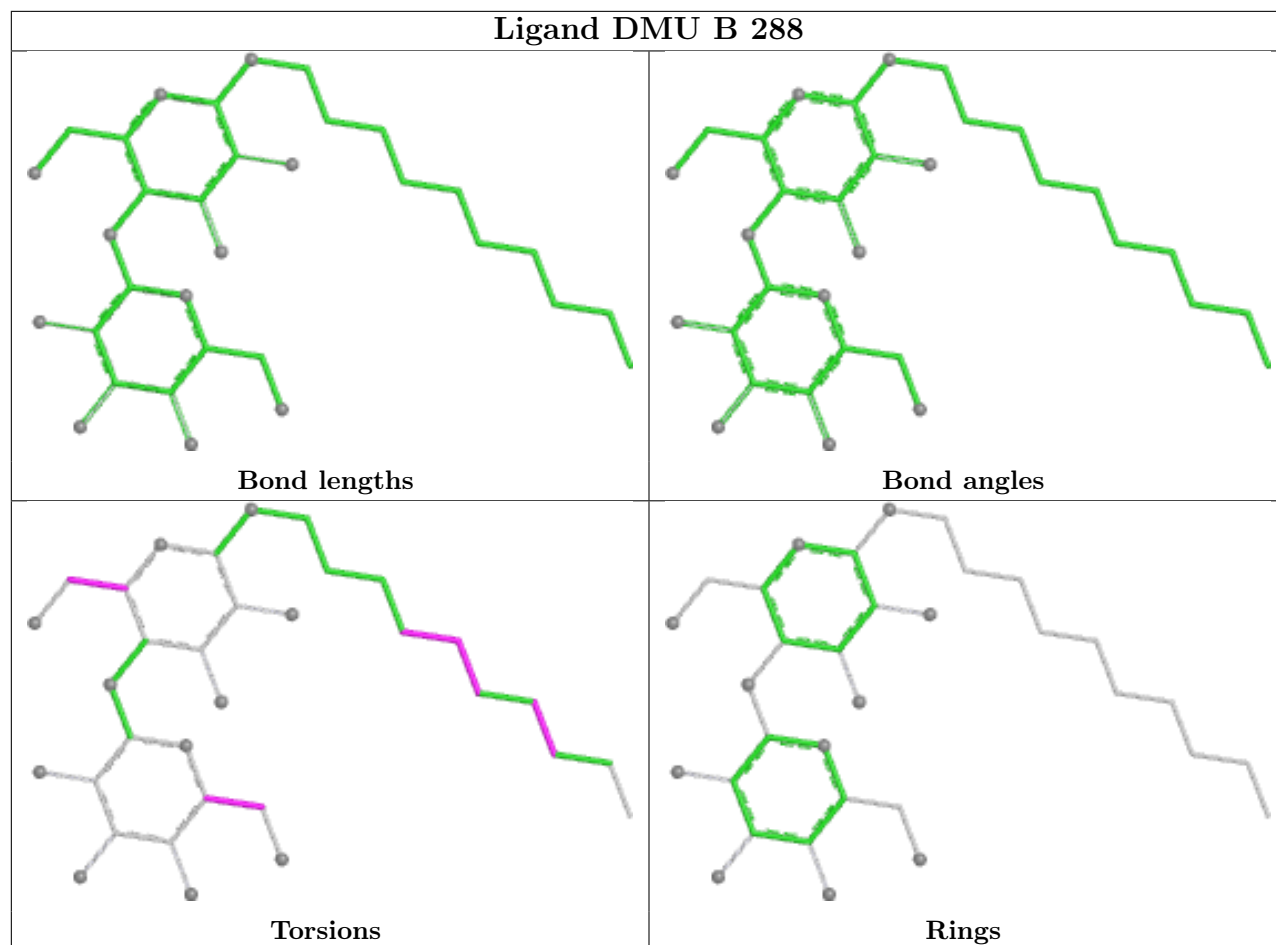
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	567	HEA	4	0
8	B	288	DMU	1	0
3	A	567	HEA	2	0
8	C	1005	DMU	2	0
9	C	1006	TRD	1	0
3	A	568	HEA	3	0
8	A	576	DMU	1	0
9	C	1010	TRD	1	0
3	C	568	HEA	4	0
8	D	1003	DMU	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

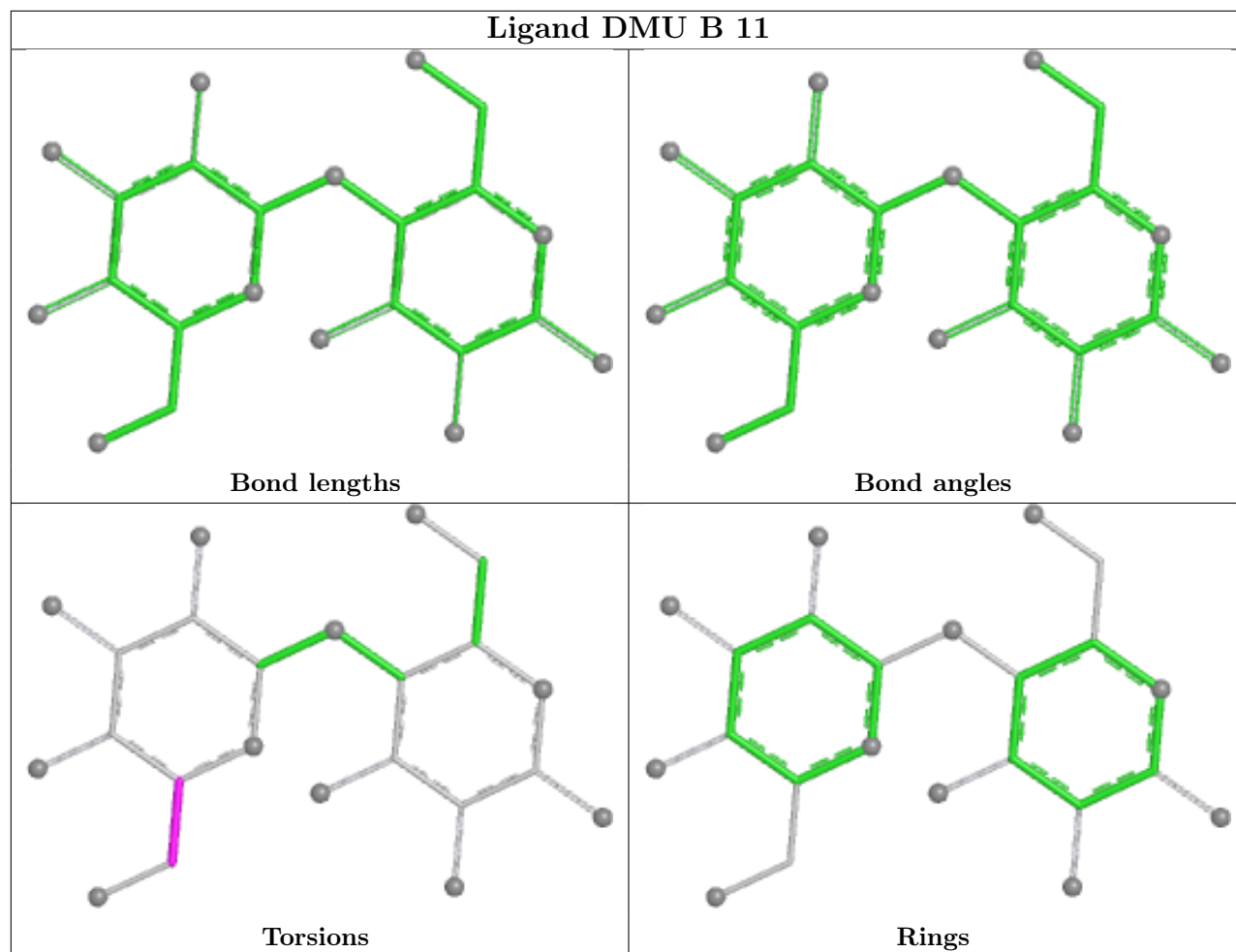




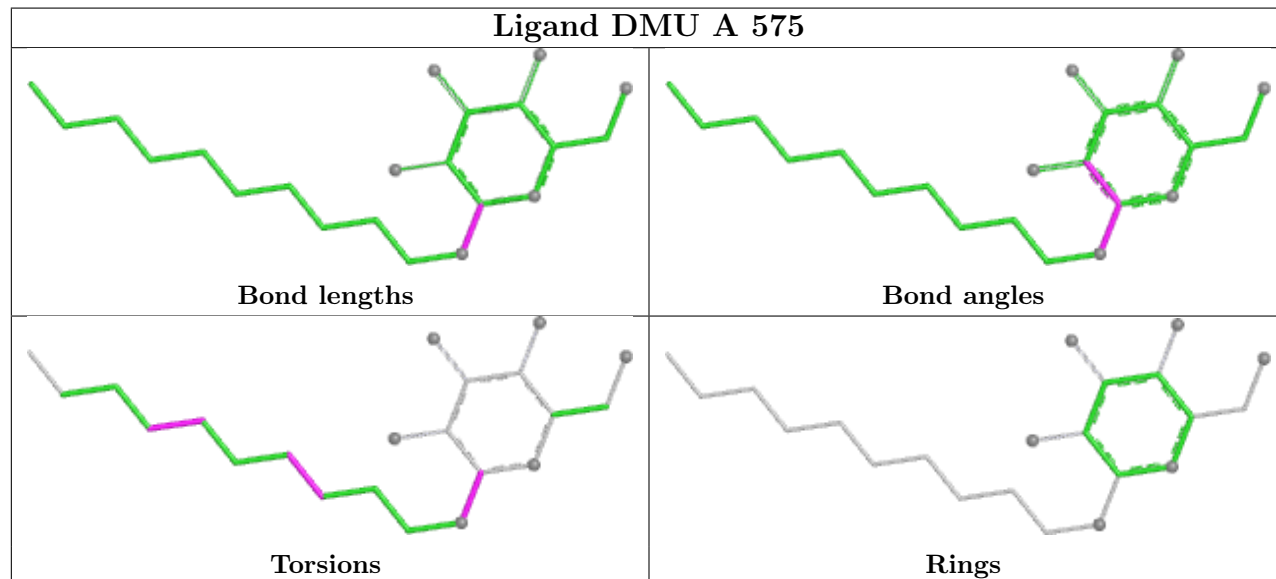




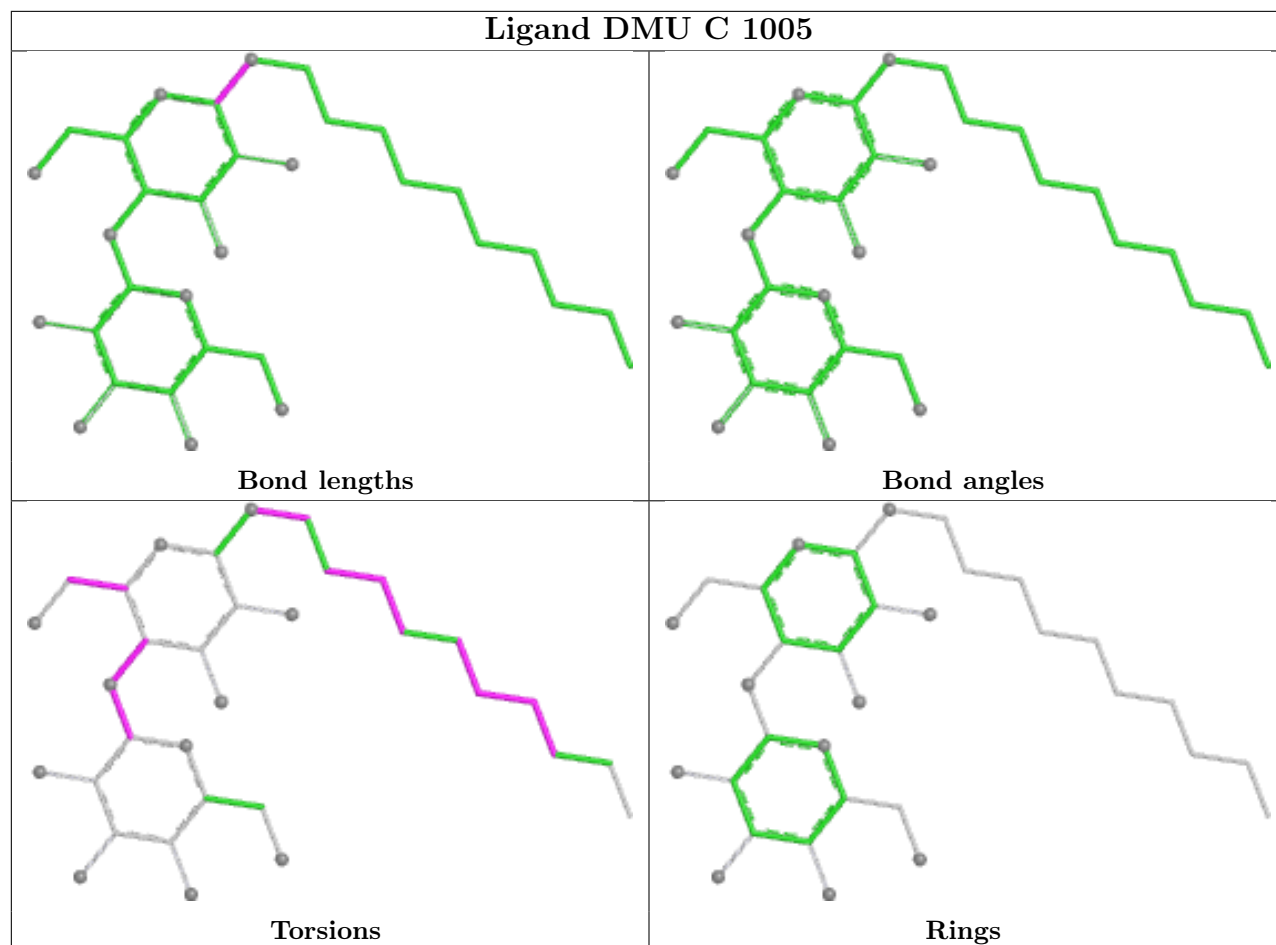
Ligand DMU B 11



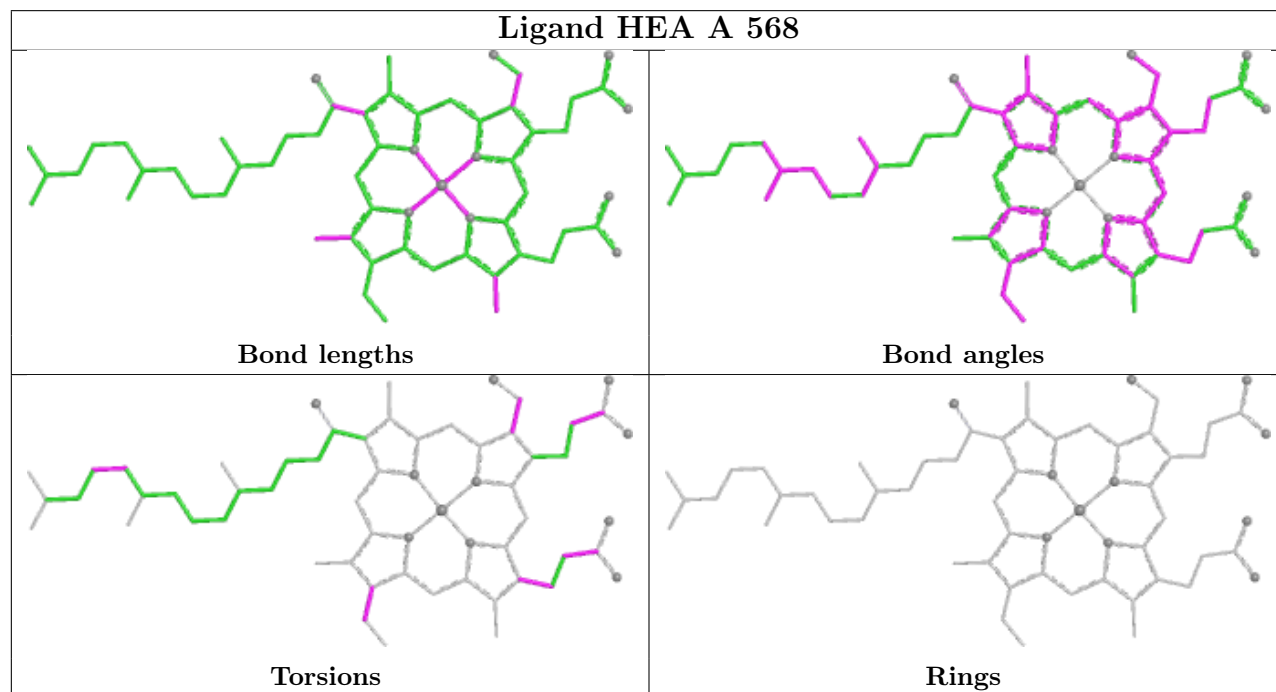
Ligand DMU A 575

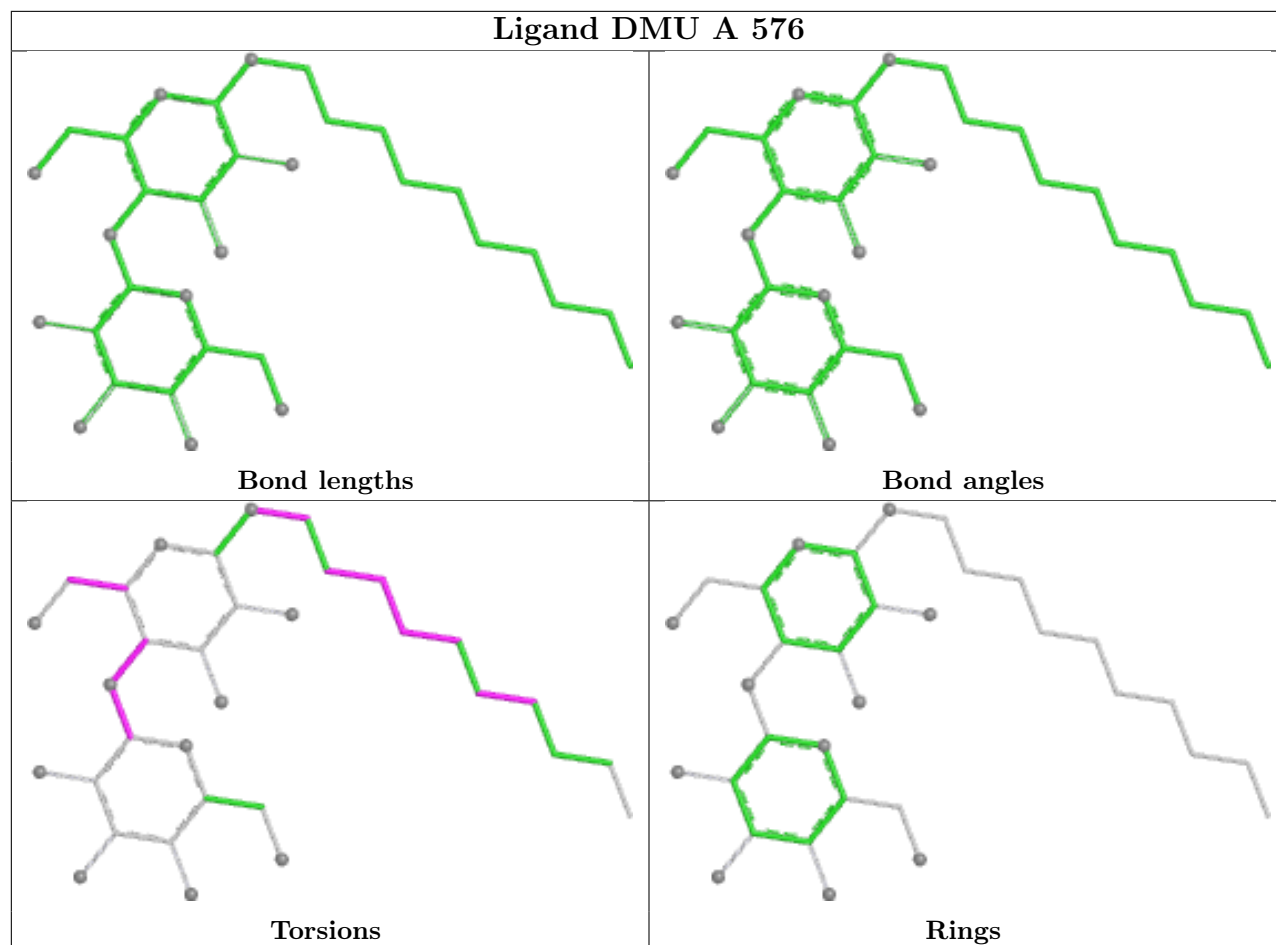


Ligand DMU C 1005

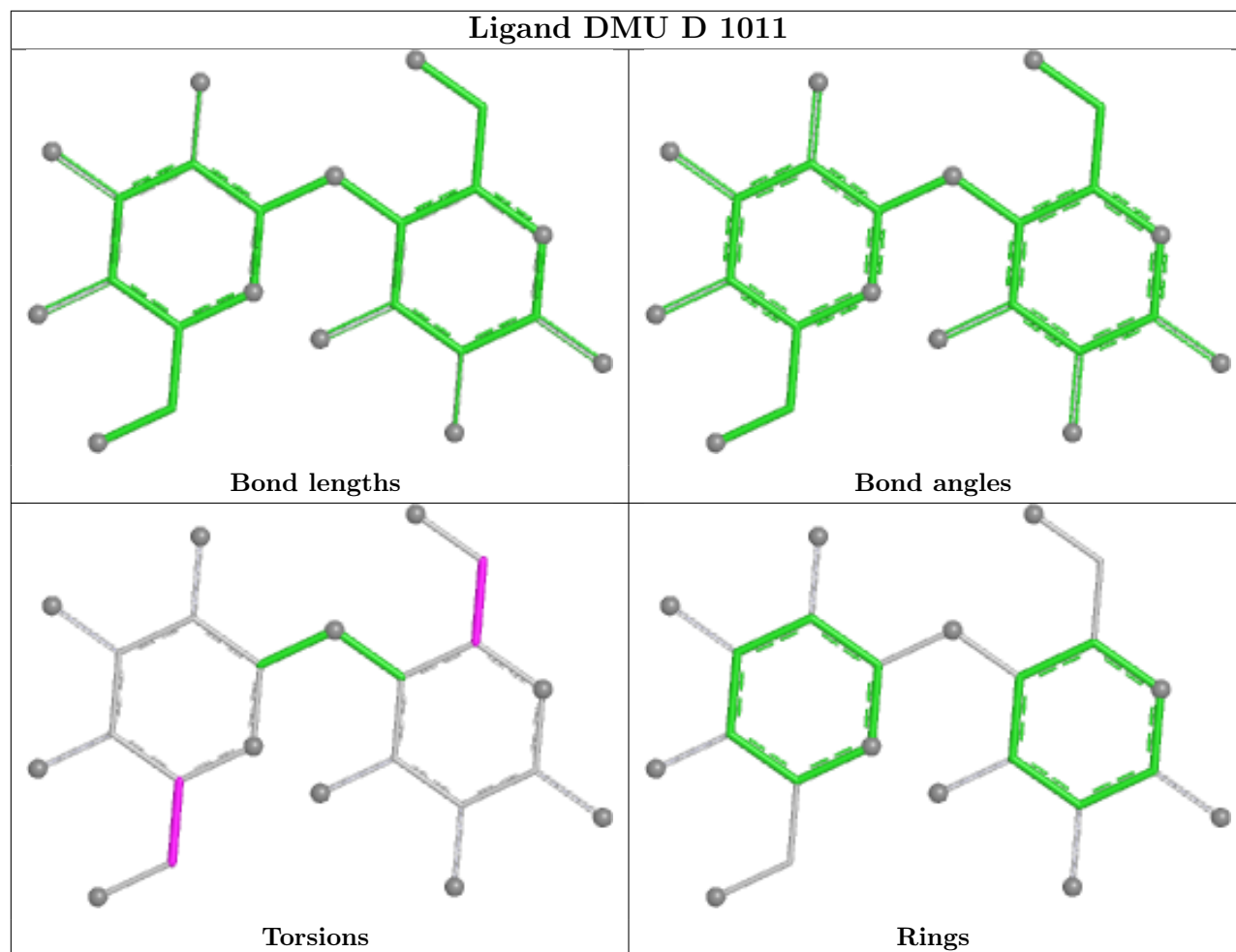


Ligand HEA A 568

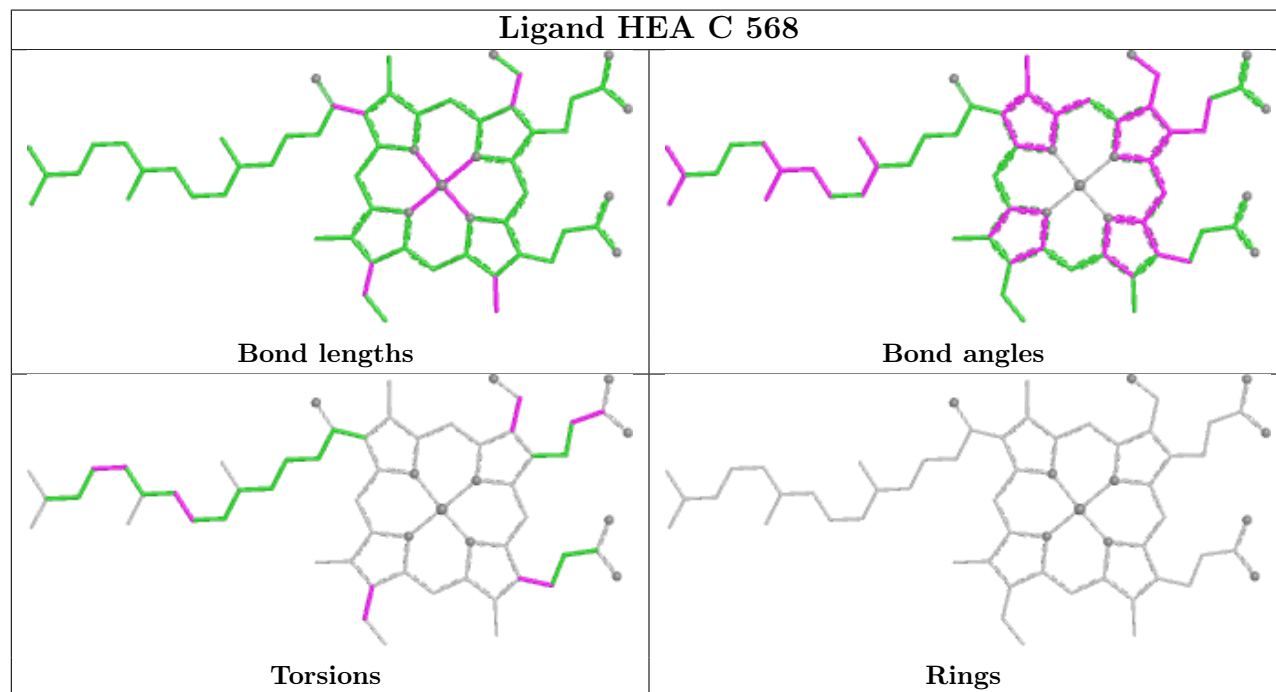


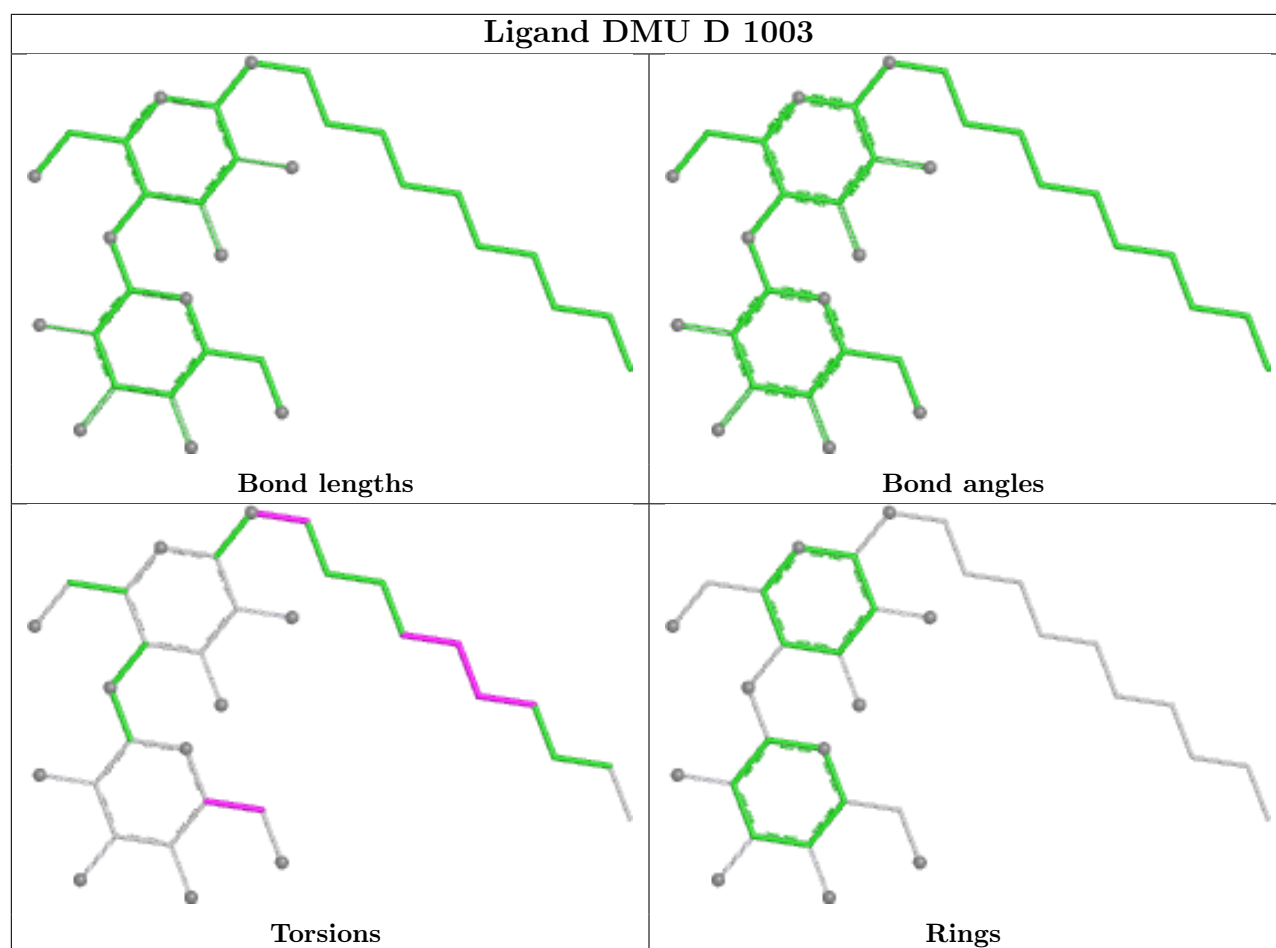


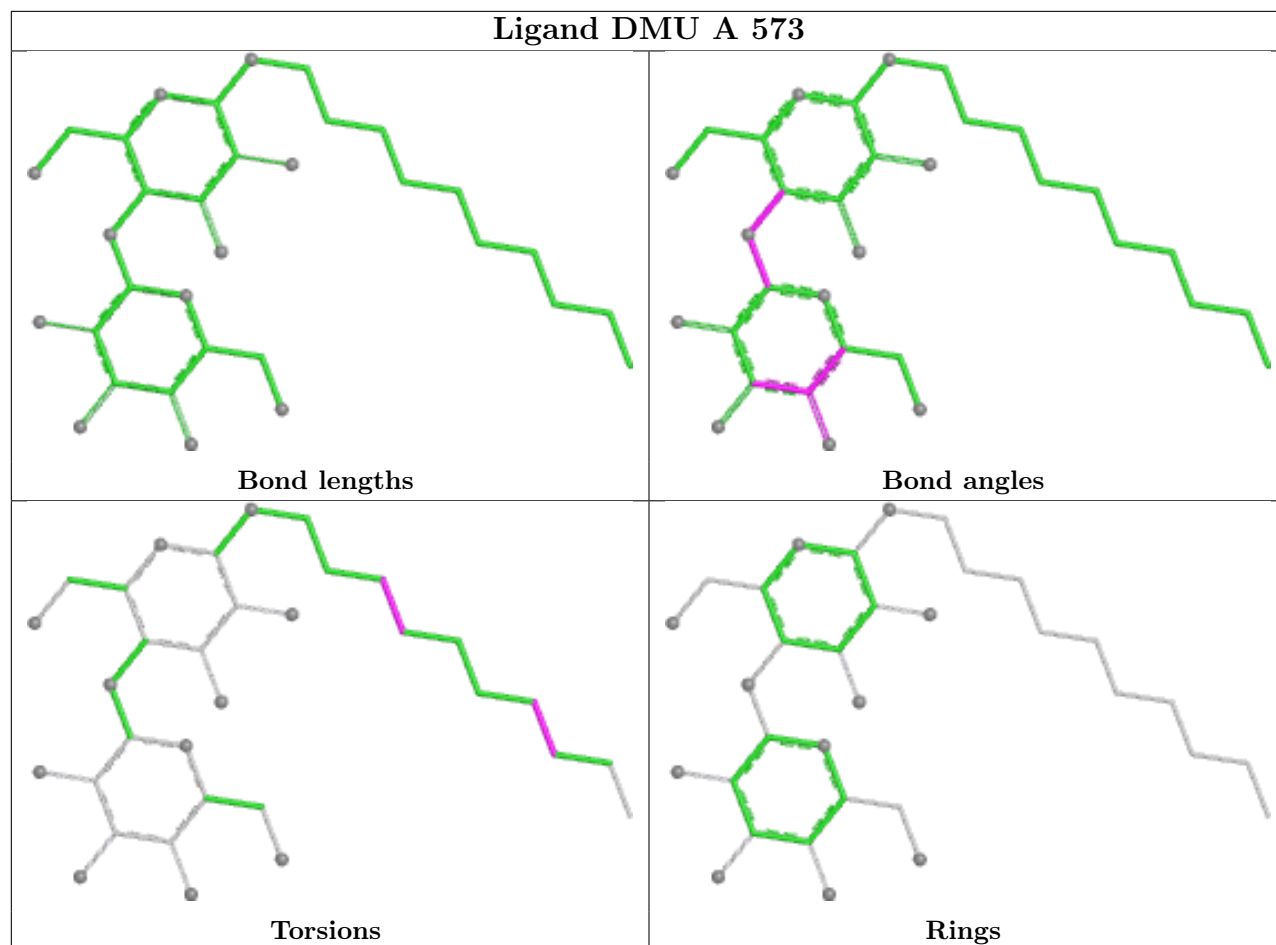
Ligand DMU D 1011



Ligand HEA C 568







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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	535/566 (94%)	0.07	21 (3%) 43 40	23, 34, 53, 72	5 (0%)
1	C	531/566 (93%)	0.96	80 (15%) 5 4	29, 51, 74, 95	5 (0%)
2	B	256/262 (97%)	0.10	7 (2%) 56 53	21, 37, 55, 60	2 (0%)
2	D	256/262 (97%)	0.37	16 (6%) 26 23	22, 42, 62, 71	3 (1%)
All	All	1578/1656 (95%)	0.42	124 (7%) 18 16	21, 41, 67, 95	15 (0%)

All (124) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	56	TRP	6.5
1	A	20	TRP	5.1
1	C	550	THR	4.9
1	C	72	LEU	4.9
1	C	20	TRP	4.8
1	C	222	MET	4.7
1	C	134	ALA	4.5
1	A	22	MET	3.9
1	C	76	PHE	3.9
1	A	81	TRP	3.9
2	D	96	HIS	3.9
1	A	551	PHE	3.8
1	C	533	GLU	3.7
1	C	548	GLU	3.7
1	C	138	MET	3.5
1	C	23	SER	3.5
1	C	77	PHE	3.5
1	C	217	ALA	3.3
1	C	530	TYR	3.1
1	C	24	THR	3.1
1	C	221	THR	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	525	VAL	3.1
2	D	109	ILE	3.1
1	A	18	THR	3.1
1	A	17	PHE	3.1
1	C	22	MET	3.1
1	C	534	HIS	3.0
1	C	543	THR	3.0
1	C	139	ASN	3.0
1	C	32	LEU	3.0
1	C	318	TYR	2.9
1	A	19	ARG	2.9
1	C	531	TRP	2.9
1	A	220	MET	2.9
1	C	541	THR	2.9
1	C	535	ALA	2.8
1	C	126	LEU	2.8
1	C	540	TRP	2.8
1	C	21	PHE	2.8
1	C	547	PRO	2.7
2	B	281	GLN	2.7
1	C	118	GLY	2.7
2	D	284	HIS	2.7
1	C	73	VAL	2.7
1	C	524	ARG	2.6
1	C	81	TRP	2.6
1	C	536	ASP	2.6
1	C	129	GLY	2.6
1	C	223	HIS	2.6
2	D	87	ARG	2.5
1	A	217	ALA	2.5
1	C	215	MET	2.5
1	C	136	PRO	2.5
1	C	80	LEU	2.5
1	C	268	GLY	2.5
1	A	347	TYR	2.5
2	D	235	LEU	2.5
1	A	221	THR	2.5
1	C	83	SER	2.5
2	D	86	LYS	2.5
1	C	351	ALA	2.5
1	C	451	TRP	2.4
1	A	72	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	203	LEU	2.4
1	C	537	THR	2.4
1	C	128	ILE	2.4
2	D	74	LEU	2.4
1	A	241	LEU	2.4
1	C	448	TYR	2.3
1	C	218	PRO	2.3
2	D	92	ALA	2.3
2	D	102	ILE	2.3
1	C	86	GLU	2.3
1	C	145	LEU	2.3
2	D	97	ASN	2.3
1	C	30	GLY	2.3
1	C	75	GLY	2.3
1	C	517	TYR	2.3
1	C	26	HIS	2.3
2	B	284	HIS	2.3
1	A	533	GLU	2.3
1	C	546	PRO	2.3
1	A	260	GLY	2.3
1	C	130	ALA	2.3
1	C	549	HIS	2.3
1	C	25	ASN	2.2
1	C	237	TRP	2.2
2	D	98	SER	2.2
1	C	399	ILE	2.2
2	B	235	LEU	2.2
1	A	548	GLU	2.2
1	C	74	LYS	2.2
1	C	444	SER	2.2
2	B	102	ILE	2.2
1	C	196	LEU	2.2
1	C	538	LEU	2.2
1	C	443	MET	2.2
1	C	313	TYR	2.2
1	C	519	LEU	2.2
1	C	521	ARG	2.1
1	C	240	LEU	2.1
2	D	56	TRP	2.1
2	D	229	ASP	2.1
1	C	212	PHE	2.1
1	A	223	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	549	HIS	2.1
1	C	131	PRO	2.1
1	C	230	TRP	2.1
1	A	269	GLY	2.1
1	C	219	GLY	2.1
1	C	520	THR	2.1
1	A	74	LYS	2.1
1	C	29	ILE	2.1
1	C	121	ASN	2.1
2	B	209	GLN	2.1
2	D	281	GLN	2.1
2	B	98	SER	2.1
1	C	27	LYS	2.1
1	C	33	TYR	2.1
2	D	83	PHE	2.1
1	C	515	ILE	2.0
2	D	81	TRP	2.0
1	C	211	THR	2.0
1	A	213	LEU	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	DMU	C	1002	23/33	0.66	0.22	97,98,99,99	23
9	TRD	C	1001	13/13	0.67	0.39	78,79,79,79	0
9	TRD	A	579	7/13	0.72	0.28	59,59,60,60	0
9	TRD	C	1006	13/13	0.74	0.32	72,74,76,76	0

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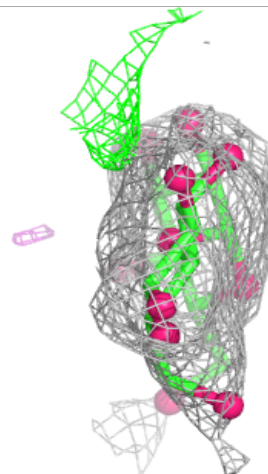
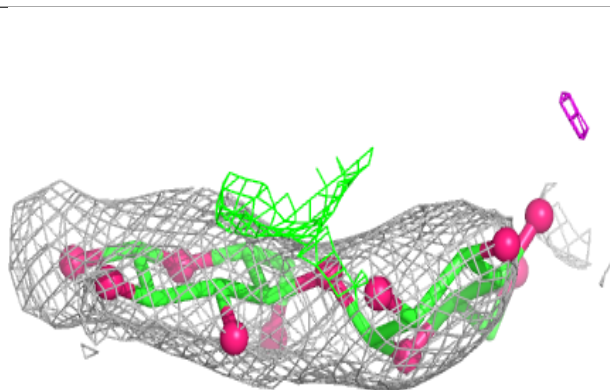
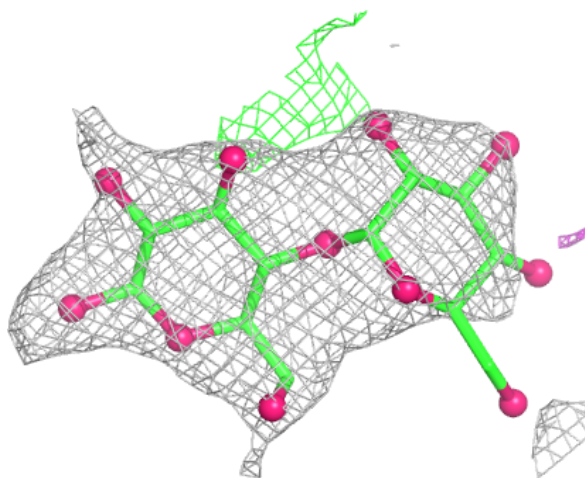
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
9	TRD	A	582	13/13	0.76	0.27	63,66,67,68	0
9	TRD	A	578	13/13	0.77	0.22	43,45,48,50	0
9	TRD	A	581	13/13	0.78	0.27	53,60,67,68	0
8	DMU	B	11	23/33	0.79	0.22	72,73,75,75	23
8	DMU	D	1011	23/33	0.79	0.48	68,69,70,70	23
9	TRD	C	1010	9/13	0.80	0.29	68,69,70,70	0
9	TRD	C	1009	7/13	0.81	0.25	64,65,65,65	0
9	TRD	A	577	13/13	0.81	0.26	66,67,71,71	0
8	DMU	A	574	33/33	0.83	0.18	66,72,74,75	0
8	DMU	A	576	33/33	0.83	0.19	41,50,51,52	33
8	DMU	D	1003	33/33	0.84	0.19	93,99,103,103	0
9	TRD	D	1007	13/13	0.84	0.23	51,53,56,57	0
8	DMU	A	575	22/33	0.85	0.15	41,61,68,70	0
9	TRD	D	1008	7/13	0.86	0.22	61,62,62,63	0
8	DMU	C	1005	33/33	0.87	0.20	59,61,62,63	33
8	DMU	B	288	33/33	0.87	0.17	77,82,86,86	0
9	TRD	A	580	7/13	0.88	0.21	59,59,60,60	0
10	HTO	B	1	10/10	0.90	0.17	54,57,58,59	0
11	CD	D	9	1/1	0.92	0.19	45,45,45,45	1
8	DMU	A	573	33/33	0.94	0.09	24,35,51,53	0
7	CYN	C	572	2/2	0.95	0.09	36,36,36,37	0
3	HEA	C	568	60/60	0.97	0.09	30,37,49,50	0
3	HEA	A	568	60/60	0.98	0.08	26,31,46,47	0
7	CYN	A	572	2/2	0.98	0.07	31,31,31,32	0
11	CD	B	9	1/1	0.98	0.10	47,47,47,47	1
3	HEA	C	567	60/60	0.98	0.08	30,33,45,46	0
6	CA	C	571	1/1	0.99	0.04	40,40,40,40	0
3	HEA	A	567	60/60	0.99	0.05	20,23,29,30	0
11	CD	B	8	1/1	0.99	0.02	40,40,40,40	0
4	CU1	C	569	1/1	0.99	0.03	42,42,42,42	0
11	CD	D	8	1/1	0.99	0.02	40,40,40,40	0
5	MG	C	570	1/1	0.99	0.06	20,20,20,20	0
4	CU1	D	3	1/1	1.00	0.02	33,33,33,33	0
4	CU1	D	4	1/1	1.00	0.02	33,33,33,33	0
5	MG	A	570	1/1	1.00	0.09	17,17,17,17	0
4	CU1	B	3	1/1	1.00	0.01	25,25,25,25	0
6	CA	A	571	1/1	1.00	0.03	27,27,27,27	0
4	CU1	B	4	1/1	1.00	0.03	25,25,25,25	0
4	CU1	A	569	1/1	1.00	0.03	35,35,35,35	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different

orientation to approximate a three-dimensional view.

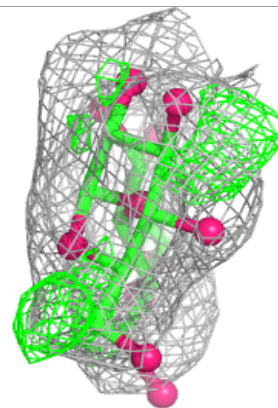
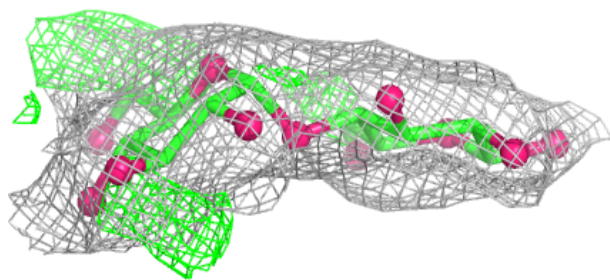
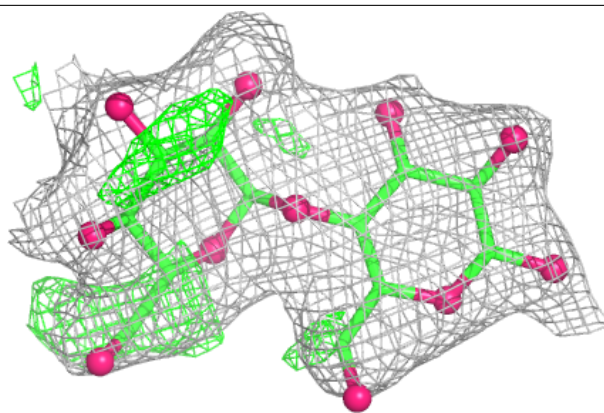
Electron density around DMU C 1002:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



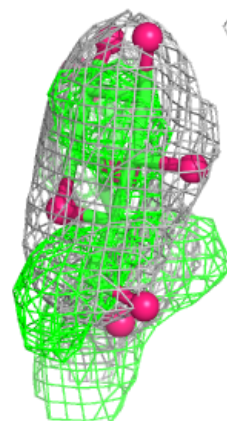
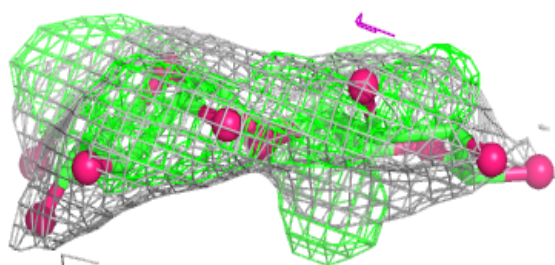
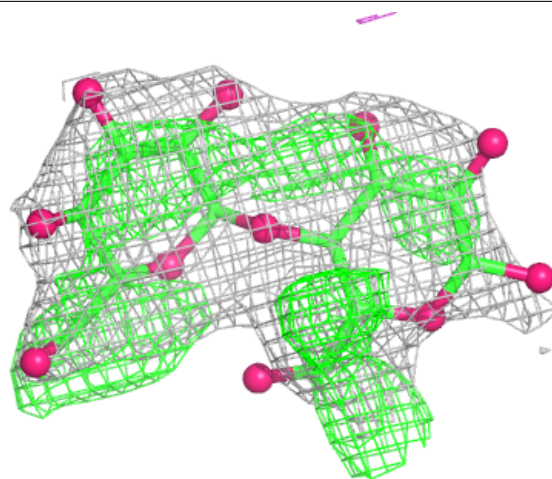
Electron density around DMU B 11:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



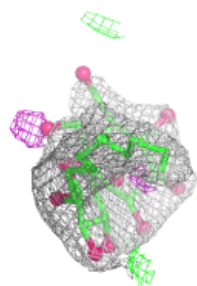
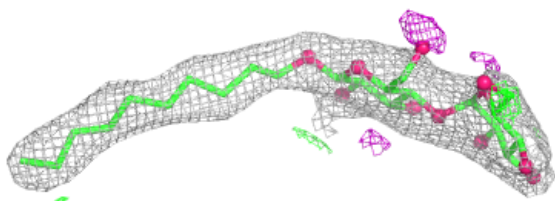
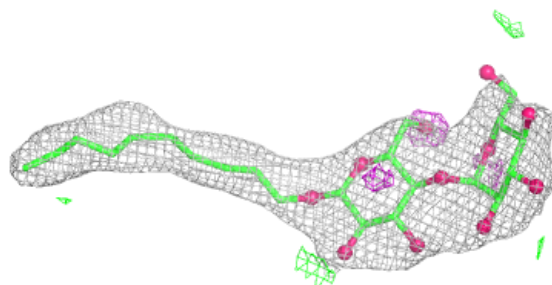
Electron density around DMU D 1011:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

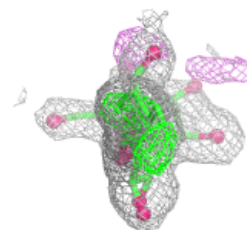
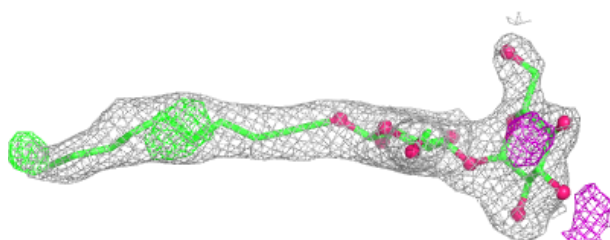
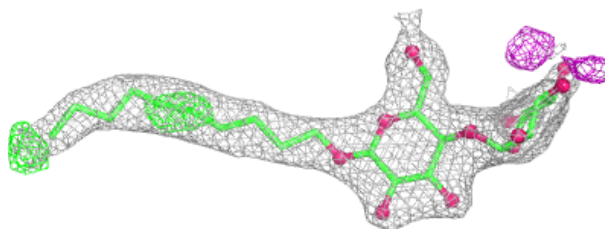


Electron density around DMU A 574:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

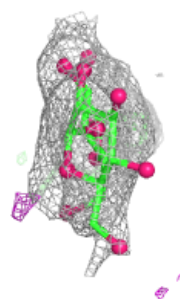
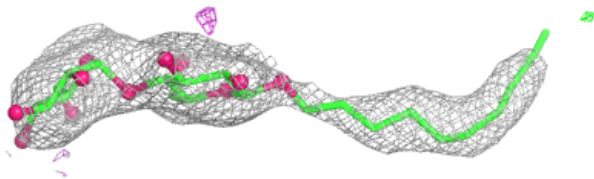
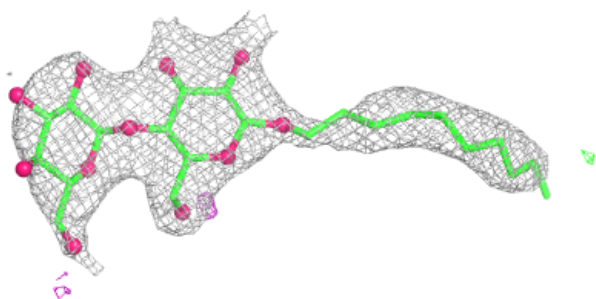
**Electron density around DMU A 576:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

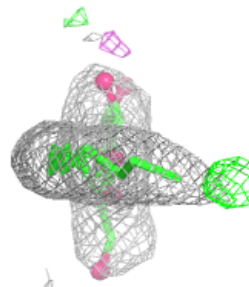
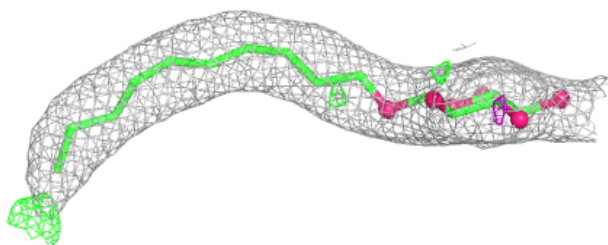
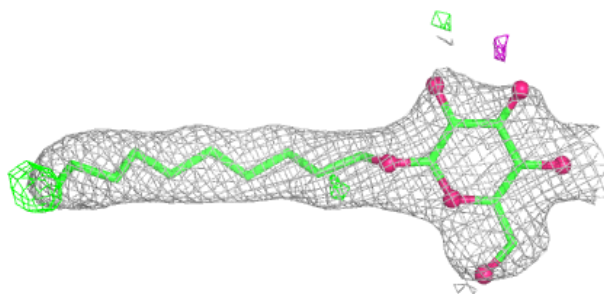


Electron density around DMU D 1003:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

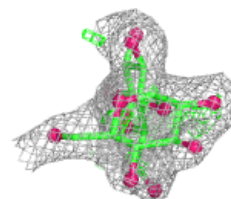
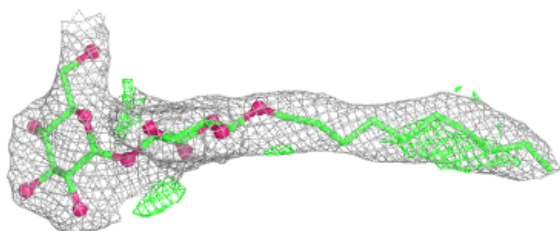
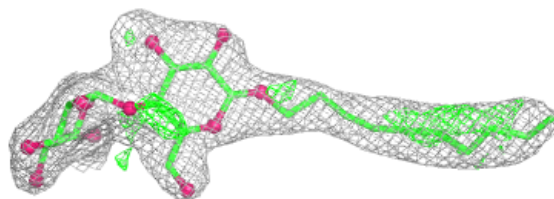
**Electron density around DMU A 575:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

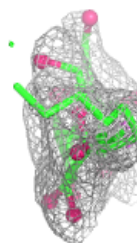
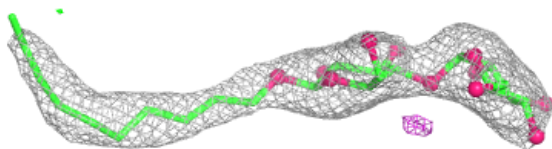
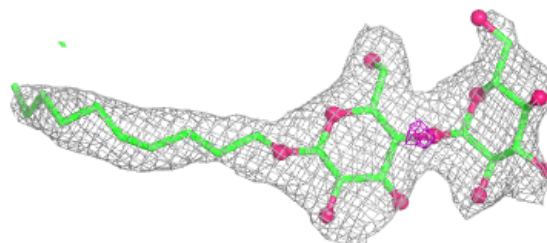


Electron density around DMU C 1005:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

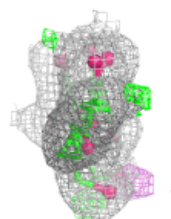
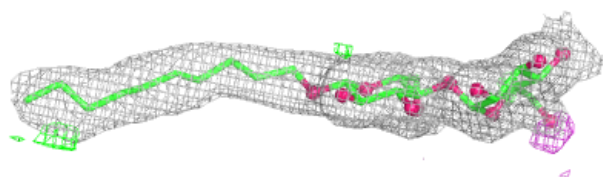
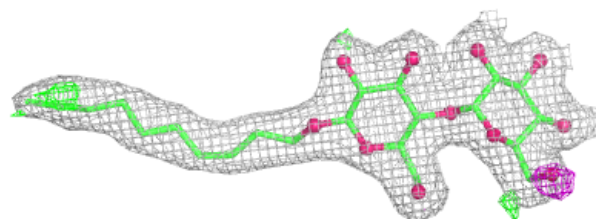
**Electron density around DMU B 288:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

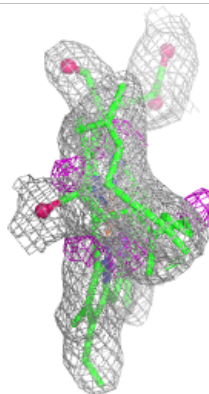
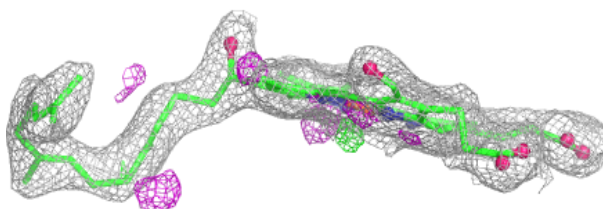
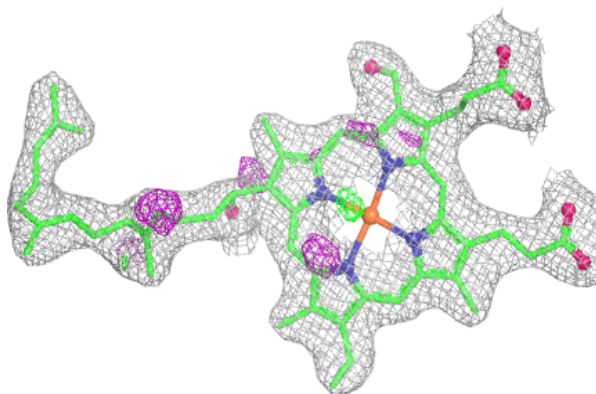


Electron density around DMU A 573:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

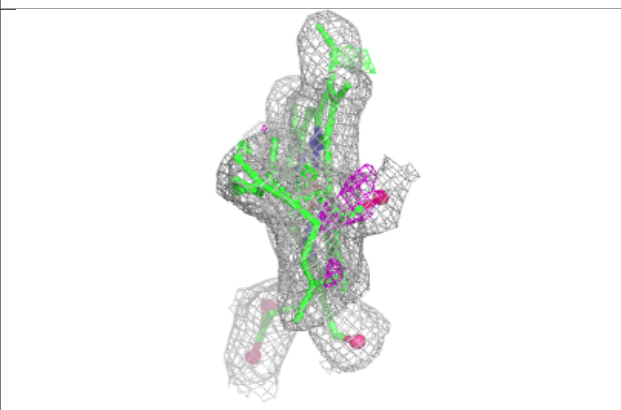
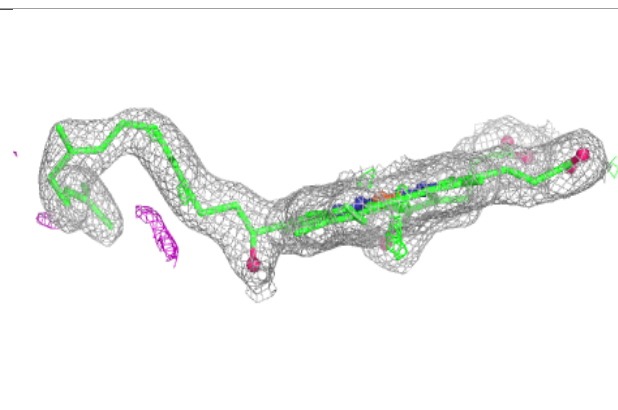
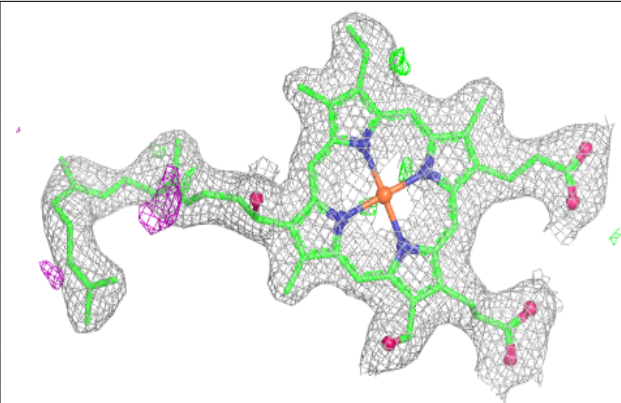
**Electron density around HEA C 568:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

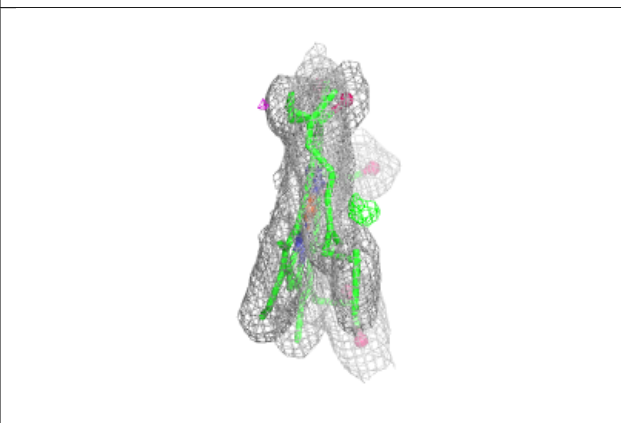
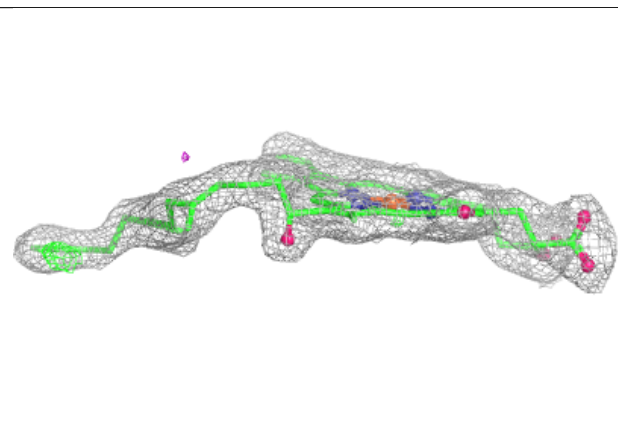
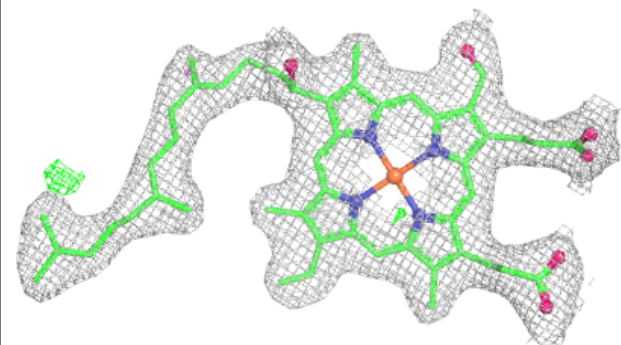


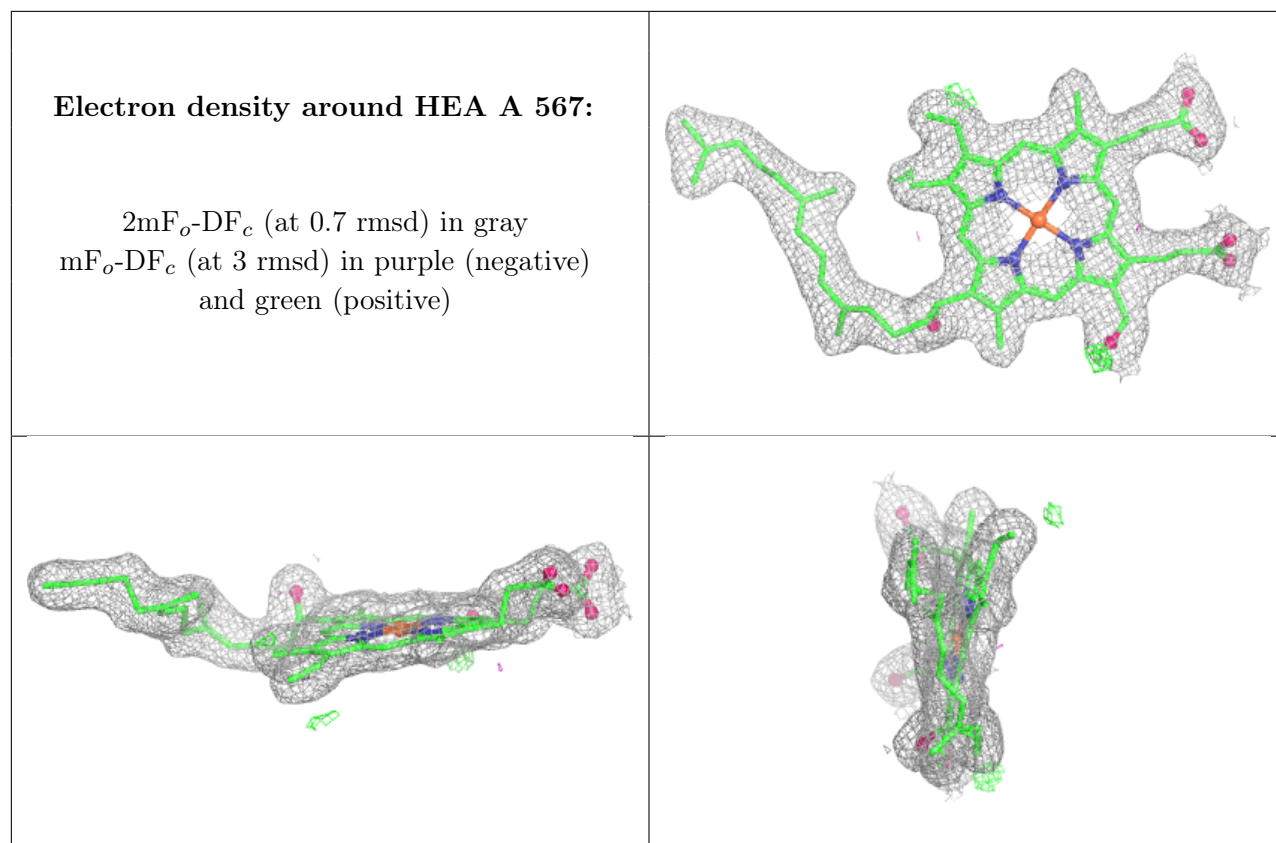
Electron density around HEA A 568:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around HEA C 567:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.