Full wwPDB X-ray Structure Validation Report (i)

Oct 31, 2023 – 06:55 PM JST

PDB ID	:	5GUX
Title	:	Cytochrome c-dependent nitric oxide reductase (cNOR) from Pseudomonas aeruginosa in complex with xenon
Authors	:	Ishii, S.; Terasaka, E.; Sugimoto, H.; Shiro, Y.; Tosha, T.
Deposited on	:	2016-08-31
Resolution	:	3.30 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 (i) 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 (i)) were used in the production of this report:

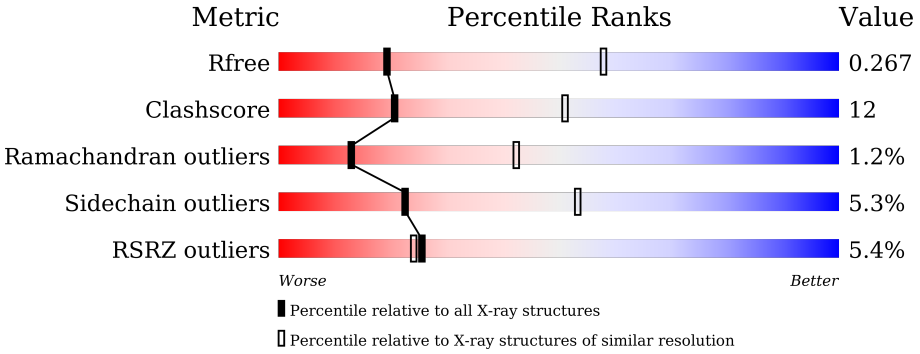
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.30 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}	130704	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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 <=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	L	213	4% 71% 28% •
2	H	225	5% 70% 28% •
3	B	465	6% 70% 23% • •
4	C	146	5% 78% 18% • •

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
8	XE	B	808	-	-	X	-
8	XE	B	809	-	-	X	-
9	10M	B	812	-	-	-	X

2 Entry composition (i)

There are 11 unique types of molecules in this entry. The entry contains 8265 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 Antibody fab fragment light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	L	213	1669	1047	277	338	7	0	0	0

- Molecule 2 is a protein called Antibody fab fragment heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	H	225	1692	1065	280	338	9	0	0	0

- Molecule 3 is a protein called Nitric oxide reductase subunit B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	B	449	3576	2416	563	572	25	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	ARG	deletion	UNP Q59647

- Molecule 4 is a protein called Nitric oxide reductase subunit C.

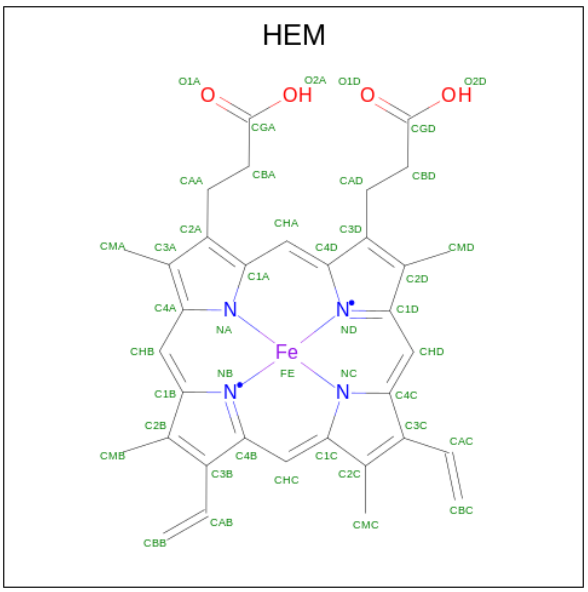
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	C	142	1123	720	195	202	6	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	100	LYS	ASN	conflict	UNP Q59646

- Molecule 5 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (for-

mula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
5	B	1	43	34	1	4	4	0	0
			Total	C	Fe	N	O		
5	B	1	43	34	1	4	4	0	0

- Molecule 6 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Fe		
6	B	1	1	1	0	0

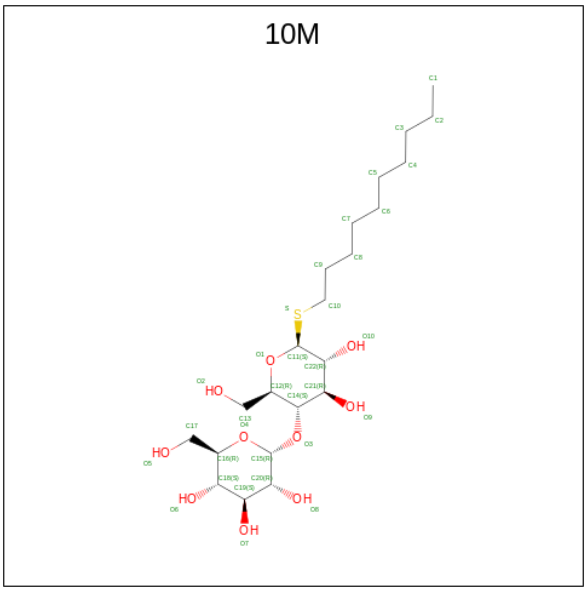
- Molecule 7 is OXYGEN ATOM (three-letter code: O) (formula: O).

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

- Molecule 8 is XENON (three-letter code: XE) (formula: Xe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Xe		
8	B	7	7	7	0	0

- Molecule 9 is decyl 4-O-alpha-D-glucopyranosyl-1-thio-beta-D-glucopyranoside (three-letter code: 10M) (formula: $C_{22}H_{42}O_{10}S$).

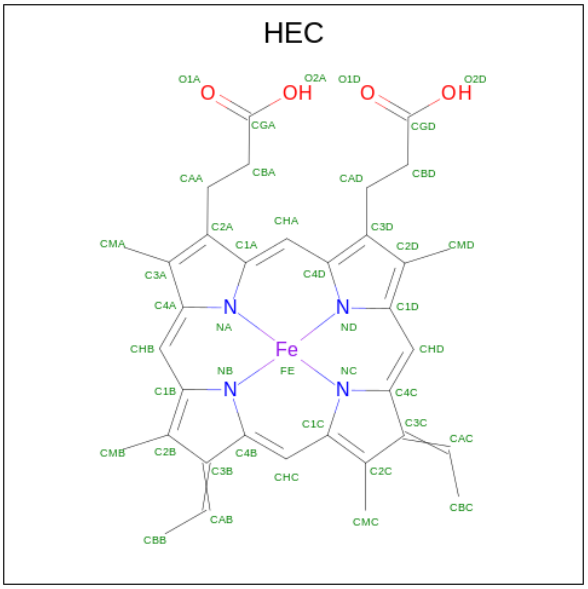


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	S		
9	B	1	33	22	10	1	0	0
			Total	C	O	S		
9	B	1	33	22	10	1	0	0

- Molecule 10 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
10	B	1	1	1	0	0

- Molecule 11 is HEME C (three-letter code: HEC) (formula: $C_{34}H_{34}FeN_4O_4$).



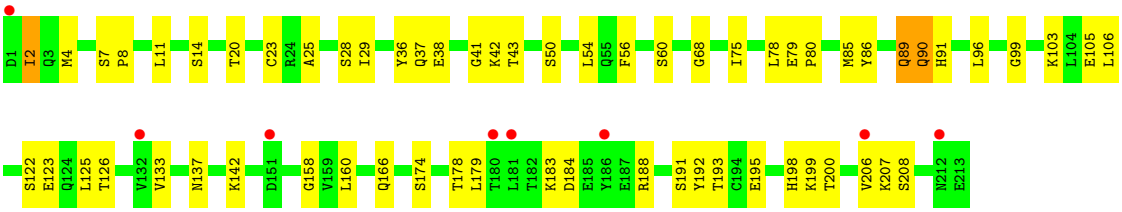
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
11	C	1	43	34	1	4	4	0	0

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.

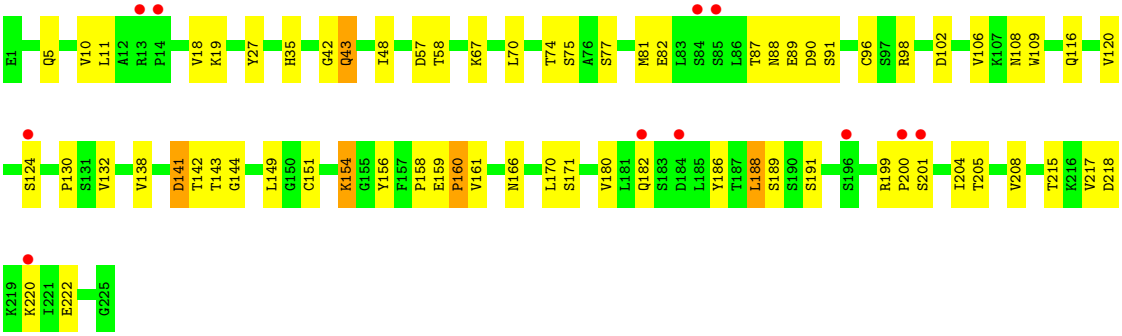
• Molecule 1: Antibody fab fragment light chain

Chain L:



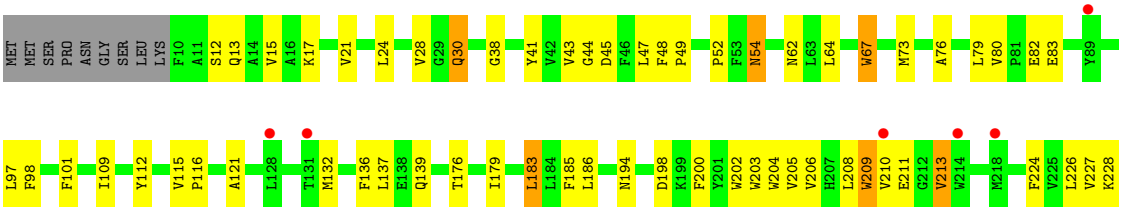
• Molecule 2: Antibody fab fragment heavy chain

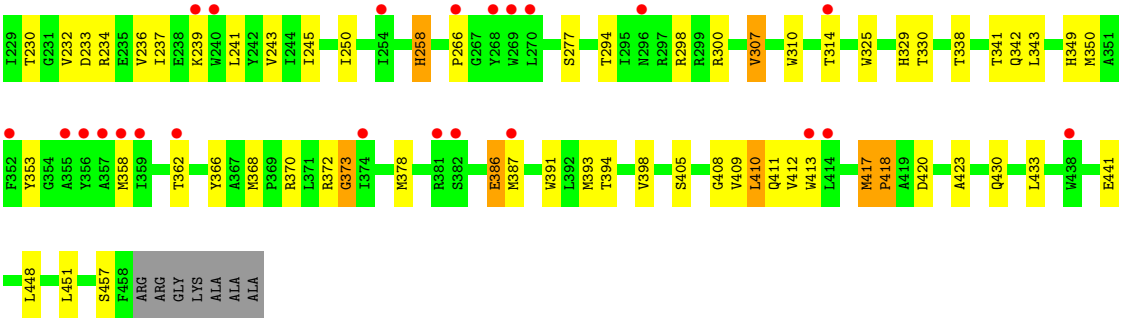
Chain H:



• Molecule 3: Nitric oxide reductase subunit B

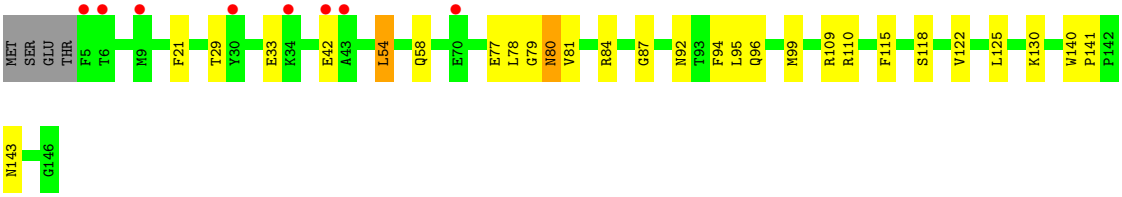
Chain B:





- Molecule 4: Nitric oxide reductase subunit C

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	89.84Å 105.38Å 192.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.39 – 3.30 39.36 – 3.30	Depositor EDS
% Data completeness (in resolution range)	96.9 (39.39-3.30) 97.2 (39.36-3.30)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.37 (at 3.32Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.209 , 0.268 0.208 , 0.267	Depositor DCC
R_{free} test set	1369 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	84.8	Xtriage
Anisotropy	0.083	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 37.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	8265	wwPDB-VP
Average B, all atoms (Å^2)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.00% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: FE, CA, O, 10M, HEM, HEC, XE

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	L	0.57	0/1709	0.77	0/2317
2	H	0.57	0/1735	0.82	0/2367
3	B	0.55	0/3693	0.72	0/5039
4	C	0.54	0/1153	0.74	0/1559
All	All	0.56	0/8290	0.75	0/11282

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	L	1669	0	1606	44	0
2	H	1692	0	1647	56	0
3	B	3576	0	3619	88	0
4	C	1123	0	1092	25	0
5	B	86	0	60	4	0
6	B	1	0	0	0	0
7	B	1	0	0	1	0
8	B	7	0	0	7	0
9	B	66	0	84	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	B	1	0	0	0	0
11	C	43	0	30	4	0
All	All	8265	0	8138	201	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 12.

All (201) 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:H:199:ARG:CG	2:H:200:PRO:HA	1.80	1.10
2:H:199:ARG:HG3	2:H:200:PRO:HA	1.33	1.09
1:L:195:GLU:HG2	1:L:206:VAL:HG22	1.36	1.08
2:H:159:GLU:HG3	2:H:160:PRO:HA	1.46	0.97
3:B:73:MET:CE	8:B:808:XE:XE	2.91	0.95
3:B:230:THR:HG22	3:B:232:VAL:HG23	1.45	0.95
1:L:96:LEU:HD11	2:H:108:ASN:HB3	1.50	0.92
2:H:159:GLU:HG3	2:H:160:PRO:CA	2.06	0.85
2:H:199:ARG:HG2	2:H:200:PRO:HA	1.59	0.84
3:B:83:GLU:OE1	3:B:234:ARG:NH2	2.12	0.81
3:B:73:MET:HE2	8:B:808:XE:XE	2.58	0.80
1:L:37:GLN:HG3	1:L:86:TYR:CE2	2.18	0.79
2:H:149:LEU:HD11	2:H:199:ARG:HD3	1.65	0.79
3:B:73:MET:HE3	8:B:808:XE:XE	2.59	0.79
2:H:199:ARG:CG	2:H:200:PRO:CA	2.62	0.78
3:B:230:THR:CG2	3:B:232:VAL:HG23	2.12	0.78
2:H:199:ARG:HG2	2:H:200:PRO:CA	2.15	0.77
3:B:341:THR:HG22	3:B:343:LEU:H	1.49	0.76
3:B:226:LEU:O	3:B:230:THR:HB	1.86	0.76
1:L:8:PRO:HG3	1:L:11:LEU:HD13	1.70	0.74
2:H:154:LYS:HE2	2:H:182:GLN:HE22	1.51	0.74
1:L:75:ILE:HG21	1:L:78:LEU:HD23	1.69	0.72
3:B:393:MET:HE1	3:B:451:LEU:HD13	1.74	0.70
3:B:405:SER:O	3:B:409:VAL:HG23	1.92	0.69
2:H:130:PRO:HB3	2:H:156:TYR:HB3	1.74	0.69
3:B:298:ARG:HG3	3:B:300:ARG:O	1.93	0.69
3:B:338:THR:O	3:B:341:THR:HB	1.93	0.68
2:H:102:ASP:O	4:C:109:ARG:NH2	2.27	0.67
3:B:121:ALA:CB	4:C:58:GLN:HG2	2.24	0.67
3:B:408:GLY:O	3:B:412:VAL:HG23	1.95	0.67
1:L:125:LEU:O	1:L:183:LYS:HD2	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:109:ILE:O	3:B:112:TYR:HB2	1.96	0.66
3:B:342:GLN:H	3:B:411:GLN:NE2	1.94	0.65
1:L:195:GLU:CG	1:L:206:VAL:HG22	2.19	0.65
3:B:387:MET:O	3:B:391:TRP:HD1	1.79	0.65
3:B:194:ASN:HD22	4:C:42:GLU:HG3	1.61	0.65
1:L:96:LEU:CD1	2:H:108:ASN:HB3	2.26	0.64
3:B:136:PHE:CE1	8:B:809:XE:XE	3.29	0.64
3:B:17:LYS:O	3:B:21:VAL:HG23	1.99	0.63
1:L:198:HIS:HD2	1:L:200:THR:OG1	1.81	0.62
3:B:121:ALA:HA	3:B:132:MET:HE1	1.80	0.62
2:H:48:ILE:HD13	2:H:81:MET:HE1	1.80	0.62
3:B:387:MET:O	3:B:391:TRP:CD1	2.53	0.61
2:H:222:GLU:HA	2:H:222:GLU:OE1	2.00	0.61
2:H:27:TYR:CE2	2:H:98:ARG:HD2	2.36	0.60
3:B:413:TRP:HD1	3:B:417:MET:CE	2.14	0.60
2:H:70:LEU:CD2	2:H:81:MET:HG3	2.32	0.59
2:H:11:LEU:HD21	2:H:158:PRO:HG3	1.84	0.59
4:C:140:TRP:CG	4:C:141:PRO:HA	2.38	0.59
3:B:310:TRP:HB3	3:B:362:THR:OG1	2.02	0.58
4:C:92:ASN:OD1	4:C:130:LYS:HE2	2.04	0.58
3:B:121:ALA:HB1	4:C:58:GLN:HG2	1.86	0.58
3:B:30:GLN:NE2	3:B:67:TRP:HE1	2.02	0.57
3:B:137:LEU:HA	3:B:139:GLN:HE22	1.70	0.57
1:L:29:ILE:HG21	1:L:90:GLN:HG3	1.85	0.56
1:L:184:ASP:HB3	1:L:188:ARG:NH2	2.20	0.56
3:B:372:ARG:HH21	3:B:378:MET:CE	2.18	0.56
3:B:349:HIS:CD2	5:B:801:HEM:NC	2.73	0.56
1:L:193:THR:HA	1:L:208:SER:HB3	1.88	0.56
3:B:239:LYS:O	3:B:243:VAL:HG23	2.06	0.55
2:H:199:ARG:HG2	2:H:200:PRO:N	2.20	0.55
3:B:413:TRP:HD1	3:B:417:MET:HE2	1.71	0.55
1:L:79:GLU:HG3	1:L:80:PRO:HD2	1.89	0.54
1:L:4:MET:HB3	1:L:99:GLY:HA2	1.88	0.54
2:H:199:ARG:HA	2:H:201:SER:N	2.22	0.54
2:H:199:ARG:HA	2:H:201:SER:H	1.73	0.54
4:C:96:GLN:HA	4:C:122:VAL:HG11	1.89	0.54
3:B:13:GLN:HE22	3:B:82:GLU:HG3	1.74	0.53
1:L:96:LEU:HD11	2:H:108:ASN:CB	2.31	0.53
2:H:87:THR:HG22	2:H:89:GLU:H	1.73	0.53
3:B:372:ARG:HH21	3:B:378:MET:HE3	1.73	0.53
3:B:230:THR:HG21	3:B:294:THR:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:417:MET:HB2	3:B:418:PRO:HD3	1.89	0.53
2:H:132:VAL:HG21	2:H:217:VAL:HG21	1.91	0.52
1:L:54:LEU:HD11	1:L:60:SER:HA	1.92	0.52
3:B:49:PRO:O	3:B:52:PRO:HD3	2.10	0.52
3:B:233:ASP:O	3:B:236:VAL:HG22	2.10	0.51
1:L:43:THR:HG23	2:H:116:GLN:HG3	1.93	0.51
2:H:19:LYS:HB2	2:H:82:GLU:HG3	1.92	0.51
3:B:258:HIS:CE1	5:B:802:HEM:CHD	2.93	0.51
3:B:241:LEU:O	3:B:245:ILE:HG12	2.10	0.51
3:B:314:THR:OG1	3:B:358:MET:HB3	2.10	0.51
1:L:193:THR:OG1	1:L:208:SER:HB3	2.11	0.51
3:B:204:TRP:O	3:B:208:LEU:HB2	2.10	0.51
2:H:48:ILE:CD1	2:H:81:MET:CE	2.89	0.51
2:H:48:ILE:HG21	2:H:81:MET:CE	2.41	0.51
3:B:52:PRO:HB2	3:B:54:ASN:ND2	2.26	0.51
2:H:67:LYS:HE3	2:H:90:ASP:OD1	2.11	0.50
1:L:96:LEU:HD13	2:H:108:ASN:HD22	1.76	0.50
3:B:230:THR:CG2	3:B:294:THR:O	2.60	0.50
2:H:161:VAL:CG1	2:H:208:VAL:HG13	2.42	0.50
1:L:85:MET:HE2	1:L:103:LYS:HD2	1.92	0.50
1:L:192:TYR:O	1:L:208:SER:HB2	2.11	0.50
2:H:70:LEU:HD23	2:H:81:MET:HG3	1.94	0.50
2:H:149:LEU:HD11	2:H:199:ARG:CD	2.40	0.50
3:B:307:VAL:HG11	3:B:366:TYR:HB2	1.94	0.50
3:B:62:ASN:ND2	3:B:112:TYR:OH	2.44	0.49
3:B:115:VAL:HG12	3:B:116:PRO:O	2.12	0.49
3:B:202:TRP:HZ3	3:B:203:TRP:CZ2	2.30	0.49
1:L:192:TYR:O	1:L:208:SER:CB	2.60	0.49
3:B:43:VAL:O	3:B:45:ASP:N	2.45	0.49
3:B:205:VAL:O	3:B:209:TRP:HB2	2.12	0.49
2:H:48:ILE:HD13	2:H:81:MET:CE	2.43	0.49
1:L:2:ILE:HD11	1:L:4:MET:SD	2.53	0.49
3:B:410:LEU:HB3	3:B:433:LEU:HD21	1.95	0.49
4:C:140:TRP:CD2	4:C:141:PRO:HA	2.47	0.49
1:L:38:GLU:HB3	1:L:85:MET:HB3	1.95	0.48
3:B:418:PRO:HG2	3:B:423:ALA:HA	1.96	0.48
3:B:82:GLU:OE1	3:B:228:LYS:HE3	2.14	0.48
3:B:353:TYR:O	3:B:358:MET:HG3	2.14	0.48
1:L:50:SER:H	1:L:91:HIS:CE1	2.31	0.48
2:H:87:THR:HG22	2:H:88:ASN:N	2.29	0.48
3:B:411:GLN:HB2	3:B:433:LEU:HD11	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:220:LYS:HE3	2:H:222:GLU:OE2	2.13	0.48
4:C:95:LEU:O	4:C:99:MET:HG2	2.13	0.48
4:C:80:ASN:HD21	4:C:84:ARG:HH12	1.62	0.47
1:L:158:GLY:O	1:L:179:LEU:HA	2.14	0.47
2:H:57:ASP:OD1	2:H:58:THR:N	2.46	0.47
3:B:136:PHE:CD1	8:B:809:XE:XE	3.45	0.47
3:B:211:GLU:OE2	7:B:804:O:O	2.32	0.47
2:H:156:TYR:CE1	2:H:186:TYR:HB2	2.50	0.47
2:H:106:VAL:HG22	2:H:109:TRP:HB2	1.96	0.47
4:C:122:VAL:O	4:C:125:LEU:HB2	2.15	0.47
3:B:64:LEU:HD23	8:B:809:XE:XE	2.92	0.47
3:B:179:ILE:HG22	3:B:183:LEU:HD22	1.97	0.47
4:C:77:GLU:OE1	4:C:79:GLY:HA3	2.15	0.47
3:B:230:THR:CG2	3:B:232:VAL:CG2	2.90	0.46
1:L:106:LEU:O	1:L:166:GLN:NE2	2.38	0.46
3:B:342:GLN:H	3:B:411:GLN:HE22	1.59	0.46
3:B:198:ASP:C	3:B:198:ASP:OD1	2.54	0.46
2:H:42:GLY:O	2:H:43:GLN:HB2	2.14	0.46
3:B:386:GLU:HG3	3:B:457:SER:OG	2.15	0.46
2:H:48:ILE:HD12	2:H:81:MET:HE3	1.97	0.46
3:B:48:PHE:CG	3:B:49:PRO:HA	2.51	0.46
8:B:807:XE:XE	8:B:808:XE:XE	3.90	0.46
3:B:372:ARG:O	3:B:373:GLY:C	2.54	0.45
3:B:393:MET:HE3	3:B:451:LEU:HB2	1.99	0.45
2:H:138:VAL:O	2:H:141:ASP:HB2	2.17	0.45
3:B:97:LEU:O	3:B:98:PHE:C	2.55	0.45
1:L:123:GLU:OE2	1:L:123:GLU:N	2.45	0.45
9:B:812:10M:H10A	9:B:812:10M:H22	1.89	0.45
4:C:115:PHE:CE2	11:C:201:HEC:HBB2	2.52	0.45
1:L:56:PHE:CE1	4:C:94:PHE:CE1	3.05	0.45
1:L:137:ASN:ND2	1:L:174:SER:HB3	2.32	0.44
3:B:200:PHE:HE2	4:C:29:THR:HG23	1.82	0.44
1:L:85:MET:HE2	1:L:85:MET:HB2	1.85	0.44
1:L:4:MET:HG3	1:L:23:CYS:SG	2.57	0.44
1:L:28:SER:HA	1:L:68:GLY:O	2.18	0.44
1:L:160:LEU:HD23	2:H:180:VAL:CG2	2.48	0.44
1:L:133:VAL:HG22	1:L:178:THR:OG1	2.17	0.44
2:H:10:VAL:HG21	2:H:18:VAL:HG21	1.99	0.44
3:B:12:SER:O	3:B:15:VAL:HG22	2.18	0.44
2:H:161:VAL:HG13	2:H:208:VAL:HG13	2.00	0.43
1:L:78:LEU:HD23	1:L:78:LEU:HA	1.76	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:30:GLN:HE21	3:B:67:TRP:HE1	1.66	0.43
3:B:137:LEU:HA	3:B:139:GLN:NE2	2.33	0.43
3:B:227:VAL:HG23	3:B:237:ILE:HG21	2.00	0.43
2:H:48:ILE:CD1	2:H:81:MET:HE3	2.48	0.43
2:H:188:LEU:C	2:H:188:LEU:HD12	2.39	0.43
9:B:813:10M:H13	9:B:813:10M:H16	2.00	0.43
2:H:159:GLU:CG	2:H:160:PRO:HA	2.34	0.43
2:H:222:GLU:OE1	2:H:222:GLU:CA	2.66	0.43
3:B:430:GLN:NE2	4:C:110:ARG:HH22	2.17	0.43
4:C:78:LEU:O	4:C:81:VAL:HG12	2.19	0.43
2:H:205:THR:HG21	2:H:218:ASP:HB3	2.00	0.43
3:B:394:THR:O	3:B:398:VAL:HG23	2.18	0.43
3:B:413:TRP:HA	3:B:417:MET:HG3	2.00	0.43
4:C:115:PHE:CD2	11:C:201:HEC:HBB2	2.53	0.42
3:B:13:GLN:NE2	3:B:82:GLU:HG3	2.33	0.42
1:L:206:VAL:O	1:L:207:LYS:HG2	2.20	0.42
2:H:91:SER:HA	2:H:120:VAL:O	2.18	0.42
2:H:35:HIS:O	2:H:96:CYS:HA	2.19	0.42
3:B:420:ASP:OD1	4:C:143:ASN:HB2	2.19	0.42
3:B:183:LEU:O	3:B:186:LEU:HB2	2.19	0.42
3:B:202:TRP:CZ2	3:B:206:VAL:HG21	2.55	0.42
1:L:41:GLY:O	1:L:42:LYS:HG3	2.19	0.42
3:B:76:ALA:O	3:B:80:VAL:HB	2.20	0.41
3:B:38:GLY:HA3	3:B:441:GLU:OE1	2.20	0.41
2:H:48:ILE:HD12	2:H:81:MET:CE	2.50	0.41
2:H:130:PRO:CB	2:H:156:TYR:HB3	2.48	0.41
1:L:36:TYR:HE2	1:L:89:GLN:HE21	1.69	0.41
3:B:210:VAL:HG11	5:B:802:HEM:C2B	2.55	0.41
3:B:250:ILE:HG23	4:C:21:PHE:CE2	2.55	0.41
3:B:277:SER:HB3	3:B:330:THR:OG1	2.20	0.41
3:B:24:LEU:HD23	3:B:24:LEU:HA	1.85	0.41
3:B:28:VAL:HG22	3:B:451:LEU:HD21	2.03	0.41
1:L:105:GLU:OE1	1:L:142:LYS:HE3	2.21	0.41
3:B:417:MET:HB2	3:B:418:PRO:CD	2.50	0.41
1:L:122:SER:O	1:L:126:THR:HG23	2.21	0.41
2:H:170:LEU:HD12	2:H:171:SER:N	2.35	0.41
1:L:4:MET:HB3	1:L:99:GLY:CA	2.50	0.41
3:B:210:VAL:CG1	5:B:802:HEM:C2B	3.04	0.41
4:C:140:TRP:CD1	4:C:141:PRO:HA	2.56	0.41
2:H:166:ASN:HD21	2:H:204:ILE:HA	1.86	0.40
3:B:79:LEU:HB2	3:B:224:PHE:CE2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:C:80:ASN:HD22	4:C:80:ASN:H	1.70	0.40
1:L:79:GLU:O	1:L:80:PRO:C	2.58	0.40
4:C:54:LEU:HD22	4:C:58:GLN:NE2	2.37	0.40
1:L:2:ILE:HD13	1:L:25:ALA:HB1	2.03	0.40
4:C:109:ARG:HH11	11:C:201:HEC:CGD	2.33	0.40
4:C:109:ARG:NH1	11:C:201:HEC:O2D	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	L	211/213 (99%)	200 (95%)	10 (5%)	1 (0%)	29	61
2	H	223/225 (99%)	202 (91%)	18 (8%)	3 (1%)	12	40
3	B	447/465 (96%)	414 (93%)	26 (6%)	7 (2%)	9	36
4	C	140/146 (96%)	130 (93%)	9 (6%)	1 (1%)	22	54
All	All	1021/1049 (97%)	946 (93%)	63 (6%)	12 (1%)	13	42

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	B	44	GLY
2	H	43	GLN
2	H	143	THR
2	H	144	GLY
3	B	209	TRP
4	C	87	GLY
1	L	199	LYS
3	B	185	PHE
3	B	266	PRO

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Mol	Chain	Res	Type
3	B	47	LEU
3	B	213	VAL
3	B	373	GLY

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	189/189 (100%)	182 (96%)	7 (4%)	34	63
2	H	192/192 (100%)	178 (93%)	14 (7%)	14	41
3	B	360/371 (97%)	340 (94%)	20 (6%)	21	52
4	C	116/120 (97%)	112 (97%)	4 (3%)	37	65
All	All	857/872 (98%)	812 (95%)	45 (5%)	22	53

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	2	ILE
1	L	7	SER
1	L	14	SER
1	L	20	THR
1	L	89	GLN
1	L	90	GLN
1	L	191	SER
2	H	5	GLN
2	H	74	THR
2	H	75	SER
2	H	77	SER
2	H	124	SER
2	H	141	ASP
2	H	142	THR
2	H	151	CYS
2	H	154	LYS
2	H	160	PRO
2	H	188	LEU

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Mol	Chain	Res	Type
2	H	189	SER
2	H	191	SER
2	H	215	THR
3	B	30	GLN
3	B	41	TYR
3	B	54	ASN
3	B	67	TRP
3	B	101	PHE
3	B	176	THR
3	B	183	LEU
3	B	213	VAL
3	B	258	HIS
3	B	307	VAL
3	B	325	TRP
3	B	329	HIS
3	B	350	MET
3	B	368	MET
3	B	370	ARG
3	B	386	GLU
3	B	410	LEU
3	B	417	MET
3	B	418	PRO
3	B	448	LEU
4	C	33	GLU
4	C	54	LEU
4	C	80	ASN
4	C	118	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	89	GLN
1	L	91	HIS
1	L	137	ASN
1	L	161	ASN
1	L	198	HIS
2	H	5	GLN
2	H	108	ASN
2	H	182	GLN
3	B	30	GLN
3	B	54	ASN
3	B	62	ASN

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Mol	Chain	Res	Type
3	B	126	ASN
3	B	139	GLN
3	B	383	GLN
3	B	411	GLN
3	B	430	GLN
4	C	31	HIS
4	C	58	GLN
4	C	60	ASN
4	C	80	ASN
4	C	96	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 15 ligands modelled in this entry, 10 are monoatomic - leaving 5 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	$\#\|Z\| > 2$
9	10M	B	812	-	34,34,34	0.98	1 (2%)	44,45,45	1.13	5 (11%)
5	HEM	B	801	3,10	41,50,50	1.94	8 (19%)	45,82,82	2.95	23 (51%)
11	HEC	C	201	4	32,50,50	2.06	6 (18%)	24,82,82	2.16	6 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
5	HEM	B	802	7,3,10	41,50,50	1.93	5 (12%)	45,82,82	2.26	17 (37%)
9	10M	B	813	-	34,34,34	0.50	0	44,45,45	0.50	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
9	10M	B	812	-	-	11/19/59/59	0/2/2/2
5	HEM	B	801	3,10	-	5/12/54/54	-
11	HEC	C	201	4	-	2/10/54/54	-
5	HEM	B	802	7,3,10	-	4/12/54/54	-
9	10M	B	813	-	-	9/19/59/59	0/2/2/2

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	801	HEM	C3D-C2D	7.89	1.53	1.36
5	B	802	HEM	C3D-C2D	7.78	1.53	1.36
11	C	201	HEC	C3C-C2C	-6.68	1.33	1.40
5	B	802	HEM	C3C-C2C	-4.82	1.33	1.40
11	C	201	HEC	C3D-C2D	4.80	1.51	1.37
5	B	801	HEM	C3C-C2C	-4.02	1.34	1.40
11	C	201	HEC	C2B-C3B	-3.84	1.36	1.40
5	B	802	HEM	C3C-CAC	3.63	1.55	1.47
5	B	801	HEM	C3C-CAC	3.33	1.54	1.47
5	B	802	HEM	CAB-C3B	3.10	1.55	1.47
5	B	801	HEM	CAB-C3B	2.85	1.55	1.47
5	B	801	HEM	CMB-C2B	2.75	1.56	1.50
5	B	801	HEM	CMD-C2D	2.48	1.56	1.50
5	B	801	HEM	CMC-C2C	2.44	1.57	1.51
11	C	201	HEC	CAA-C2A	2.40	1.56	1.52
11	C	201	HEC	C3C-C4C	2.17	1.47	1.43
11	C	201	HEC	CMB-C2B	2.07	1.56	1.51
5	B	801	HEM	C2C-C1C	2.07	1.47	1.42
9	B	812	10M	C10-S	2.01	1.84	1.81
5	B	802	HEM	CAA-C2A	2.00	1.55	1.52

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	B	801	HEM	C2C-C3C-C4C	7.86	112.38	106.90
5	B	801	HEM	C4B-C3B-C2B	5.60	111.56	107.11
11	C	201	HEC	CBD-CAD-C3D	-5.54	103.17	112.62
5	B	802	HEM	C4D-ND-C1D	5.33	110.58	105.07
5	B	801	HEM	CMA-C3A-C4A	-5.26	120.38	128.46
5	B	802	HEM	CBA-CAA-C2A	-4.99	104.11	112.62
5	B	802	HEM	C4C-CHD-C1D	4.93	129.06	122.56
5	B	801	HEM	C4D-ND-C1D	4.90	110.14	105.07
5	B	801	HEM	CHD-C1D-ND	4.71	129.55	124.43
5	B	801	HEM	O1D-CGD-CBD	-4.67	108.09	123.08
5	B	801	HEM	CHC-C4B-NB	4.65	129.48	124.43
5	B	801	HEM	O2D-CGD-CBD	4.57	128.73	114.03
5	B	802	HEM	CHA-C4D-ND	4.52	129.96	124.38
11	C	201	HEC	C1D-C2D-C3D	-4.39	103.94	107.00
5	B	802	HEM	C4B-CHC-C1C	4.29	128.22	122.56
5	B	801	HEM	C4C-CHD-C1D	4.22	128.13	122.56
11	C	201	HEC	CMB-C2B-C1B	-3.97	122.36	128.46
5	B	801	HEM	CAD-C3D-C4D	3.97	131.59	124.66
5	B	801	HEM	C3B-C2B-C1B	-3.91	103.58	106.49
5	B	802	HEM	O2D-CGD-CBD	3.87	126.47	114.03
5	B	802	HEM	CMD-C2D-C1D	3.75	130.75	125.04
5	B	801	HEM	CMB-C2B-C1B	3.70	130.68	125.04
5	B	801	HEM	CBD-CAD-C3D	-3.62	102.57	112.63
5	B	801	HEM	CMA-C3A-C2A	3.58	131.69	124.94
11	C	201	HEC	CMB-C2B-C3B	3.50	129.94	125.82
5	B	801	HEM	C4B-CHC-C1C	3.23	126.82	122.56
5	B	802	HEM	O1D-CGD-CBD	-3.13	113.04	123.08
5	B	802	HEM	CAD-CBD-CGD	-3.00	107.15	113.60
9	B	812	10M	C11-O1-C12	2.94	118.00	112.58
11	C	201	HEC	CMD-C2D-C3D	2.87	130.35	124.94
5	B	801	HEM	C4D-C3D-C2D	-2.86	102.73	106.90
5	B	802	HEM	C1D-C2D-C3D	-2.72	104.09	106.96
11	C	201	HEC	CBA-CAA-C2A	-2.54	108.31	112.60
5	B	802	HEM	CAD-C3D-C4D	2.54	129.10	124.66
9	B	812	10M	C10-S-C11	2.50	104.76	100.09
5	B	801	HEM	CAA-CBA-CGA	-2.43	106.95	113.76
5	B	801	HEM	CHB-C1B-NB	-2.40	121.41	124.38
5	B	801	HEM	O2A-CGA-CBA	2.40	121.73	114.03
5	B	802	HEM	C3B-C2B-C1B	2.36	108.24	106.49
5	B	802	HEM	CHA-C4D-C3D	-2.35	120.91	125.33
5	B	802	HEM	CMC-C2C-C3C	2.30	128.97	124.68
9	B	812	10M	C19-C18-C16	2.24	114.23	110.24
9	B	812	10M	O1-C11-C22	2.24	113.13	110.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	B	801	HEM	CHD-C1D-C2D	-2.23	121.49	124.98
9	B	812	10M	C15-O3-C14	2.22	123.45	117.96
5	B	802	HEM	CHC-C4B-NB	-2.21	122.03	124.43
5	B	801	HEM	CBB-CAB-C3B	-2.17	116.83	127.62
5	B	801	HEM	C3C-C4C-NC	-2.15	106.89	110.94
5	B	802	HEM	C4B-C3B-C2B	-2.11	105.44	107.11
5	B	801	HEM	C1B-NB-C4B	2.10	107.24	105.07
5	B	802	HEM	O2A-CGA-O1A	-2.03	118.23	123.30

There are no chirality outliers.

All (31) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	B	812	10M	C22-C11-S-C10
9	B	812	10M	C9-C10-S-C11
9	B	812	10M	C18-C16-C17-O5
9	B	812	10M	O4-C16-C17-O5
9	B	812	10M	C7-C8-C9-C10
9	B	813	10M	C3-C4-C5-C6
9	B	813	10M	C6-C7-C8-C9
9	B	813	10M	C2-C3-C4-C5
9	B	812	10M	C4-C5-C6-C7
9	B	812	10M	C1-C2-C3-C4
9	B	813	10M	C1-C2-C3-C4
9	B	812	10M	O1-C12-C13-O2
5	B	801	HEM	C2B-C3B-CAB-CBB
9	B	812	10M	C6-C7-C8-C9
9	B	813	10M	C7-C8-C9-C10
9	B	813	10M	O4-C16-C17-O5
11	C	201	HEC	CAD-CBD-CGD-O1D
5	B	802	HEM	CAA-CBA-CGA-O1A
5	B	802	HEM	CAD-CBD-CGD-O1D
11	C	201	HEC	CAD-CBD-CGD-O2D
5	B	802	HEM	CAD-CBD-CGD-O2D
9	B	812	10M	C3-C4-C5-C6
9	B	813	10M	C5-C6-C7-C8
5	B	802	HEM	CAA-CBA-CGA-O2A
9	B	812	10M	C2-C3-C4-C5
9	B	813	10M	O1-C12-C13-O2
9	B	813	10M	C14-C12-C13-O2
5	B	801	HEM	C4B-C3B-CAB-CBB
5	B	801	HEM	CAD-CBD-CGD-O2D

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Mol	Chain	Res	Type	Atoms
5	B	801	HEM	CAA-CBA-CGA-O2A
5	B	801	HEM	CAA-CBA-CGA-O1A

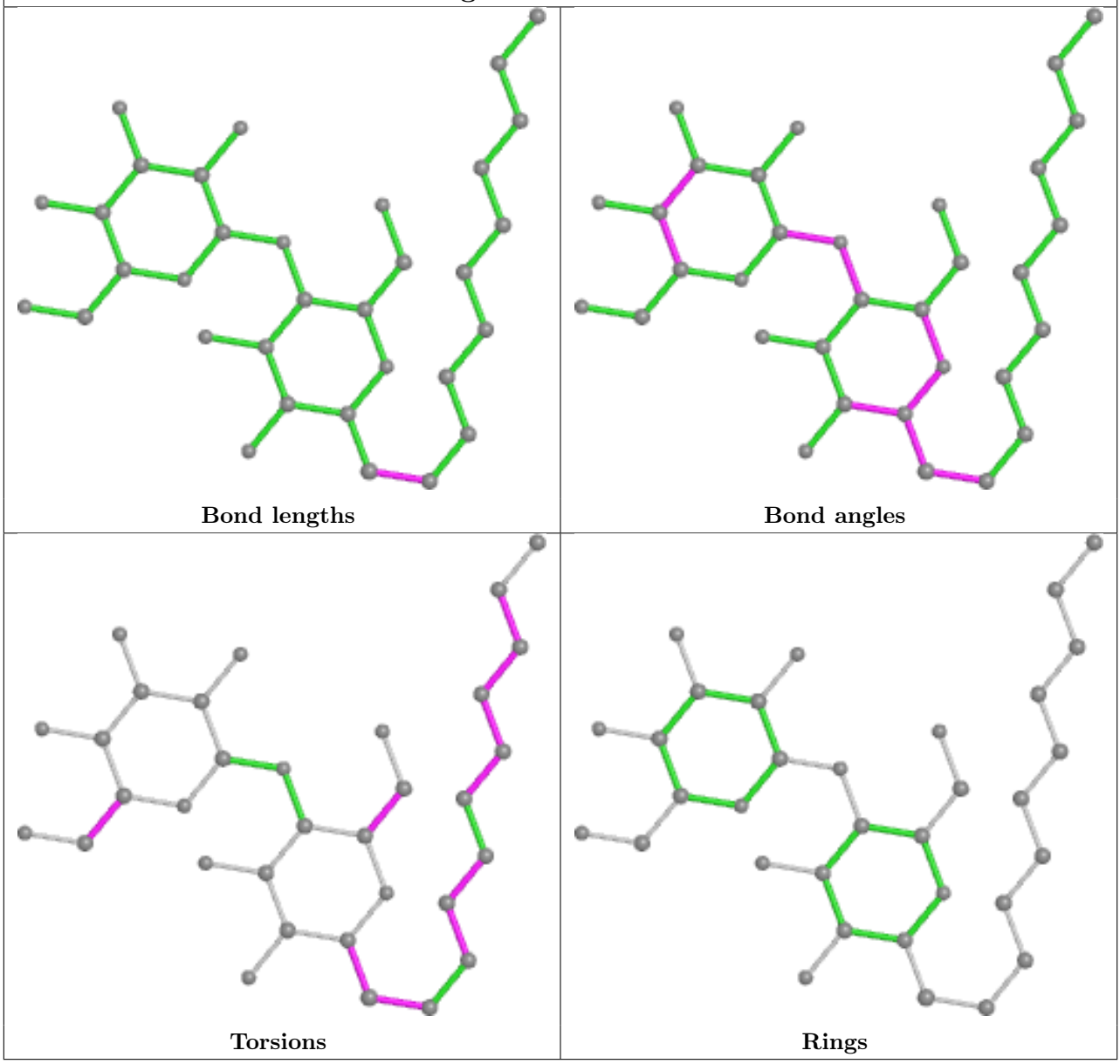
There are no ring outliers.

5 monomers are involved in 10 short contacts:

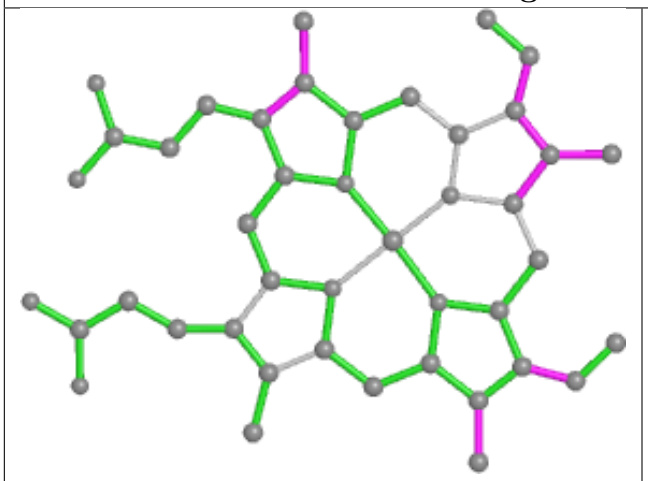
Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	B	812	10M	1	0
5	B	801	HEM	1	0
11	C	201	HEC	4	0
5	B	802	HEM	3	0
9	B	813	10M	1	0

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 then 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.

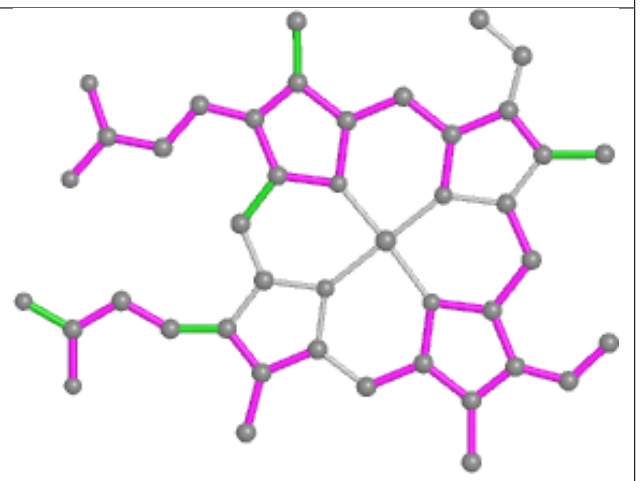
Ligand 10M B 812



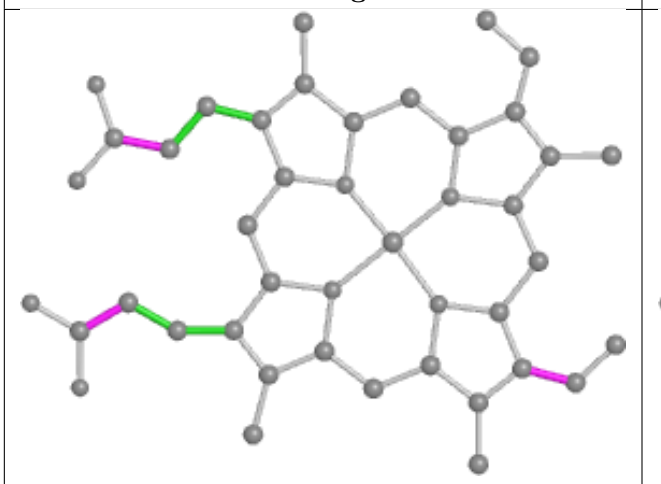
Ligand HEM B 801



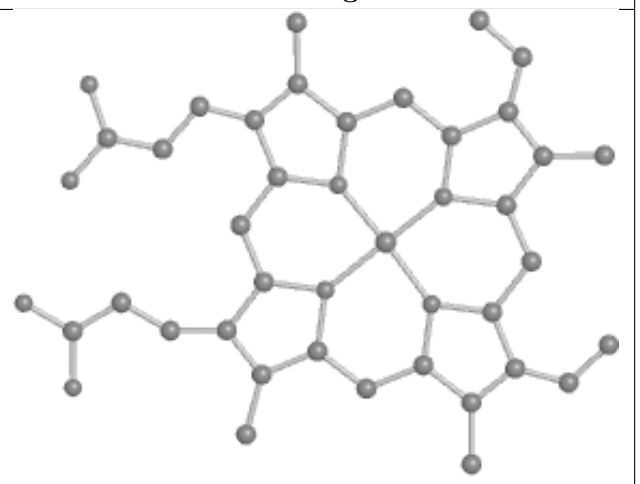
Bond lengths



Bond angles

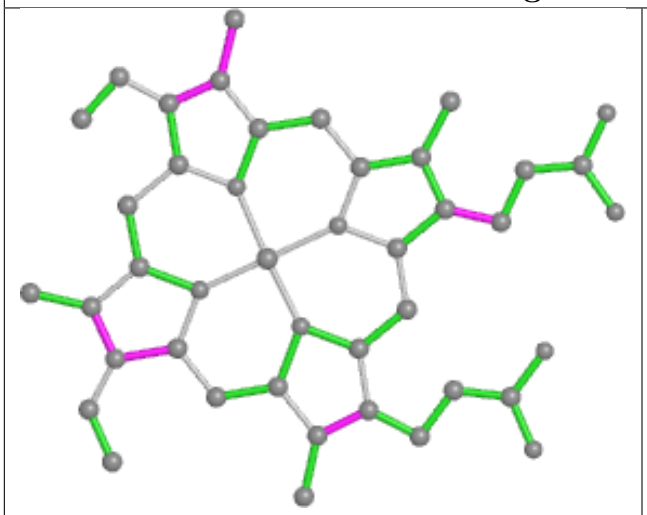


Torsions

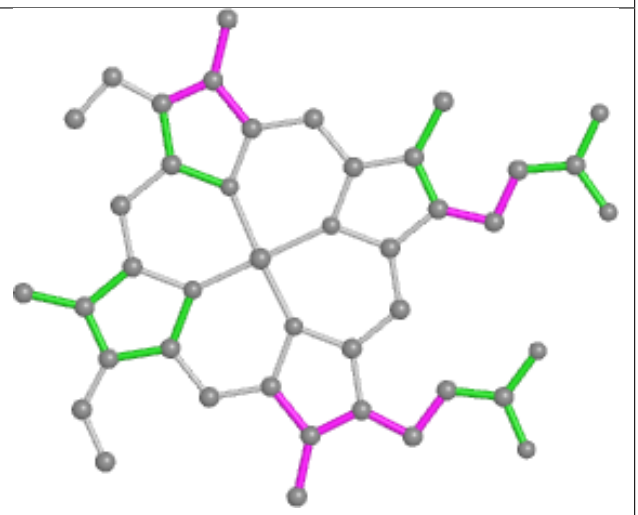


Rings

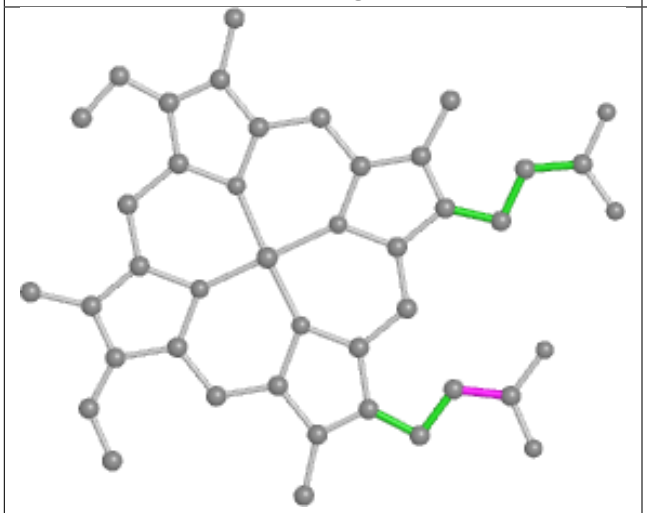
Ligand HEC C 201



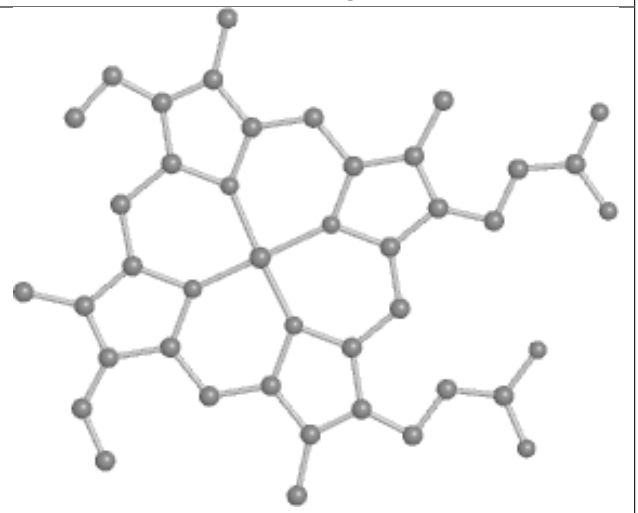
Bond lengths



Bond angles

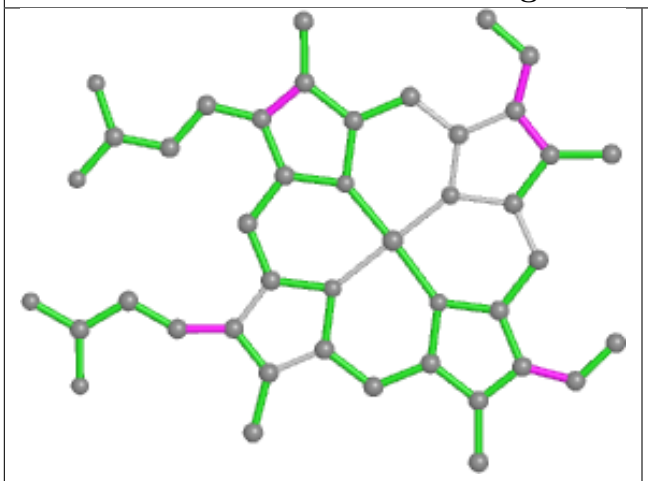


Torsions

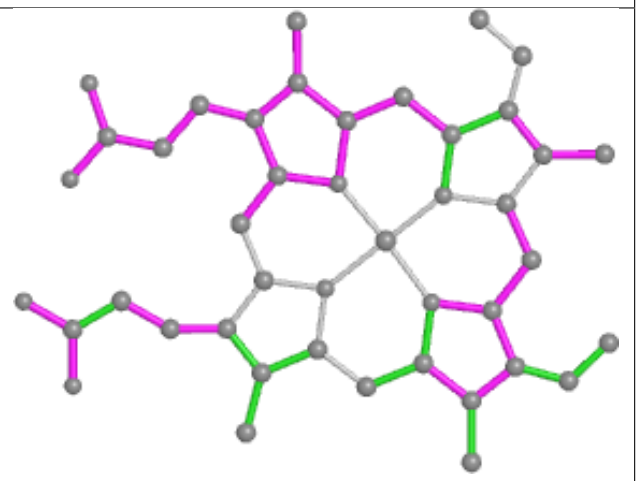


Rings

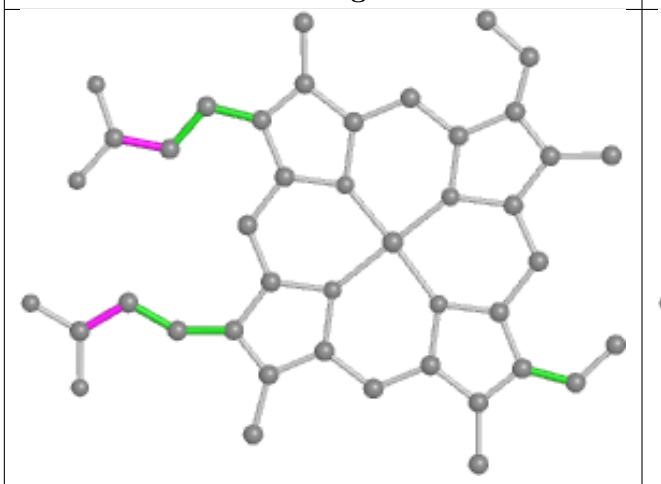
Ligand HEM B 802



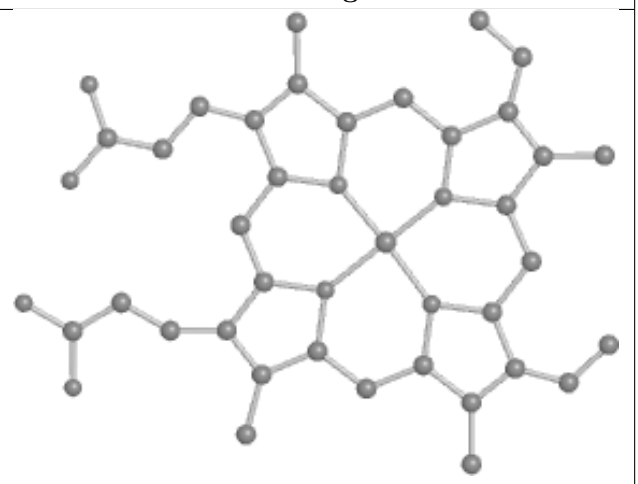
Bond lengths



Bond angles

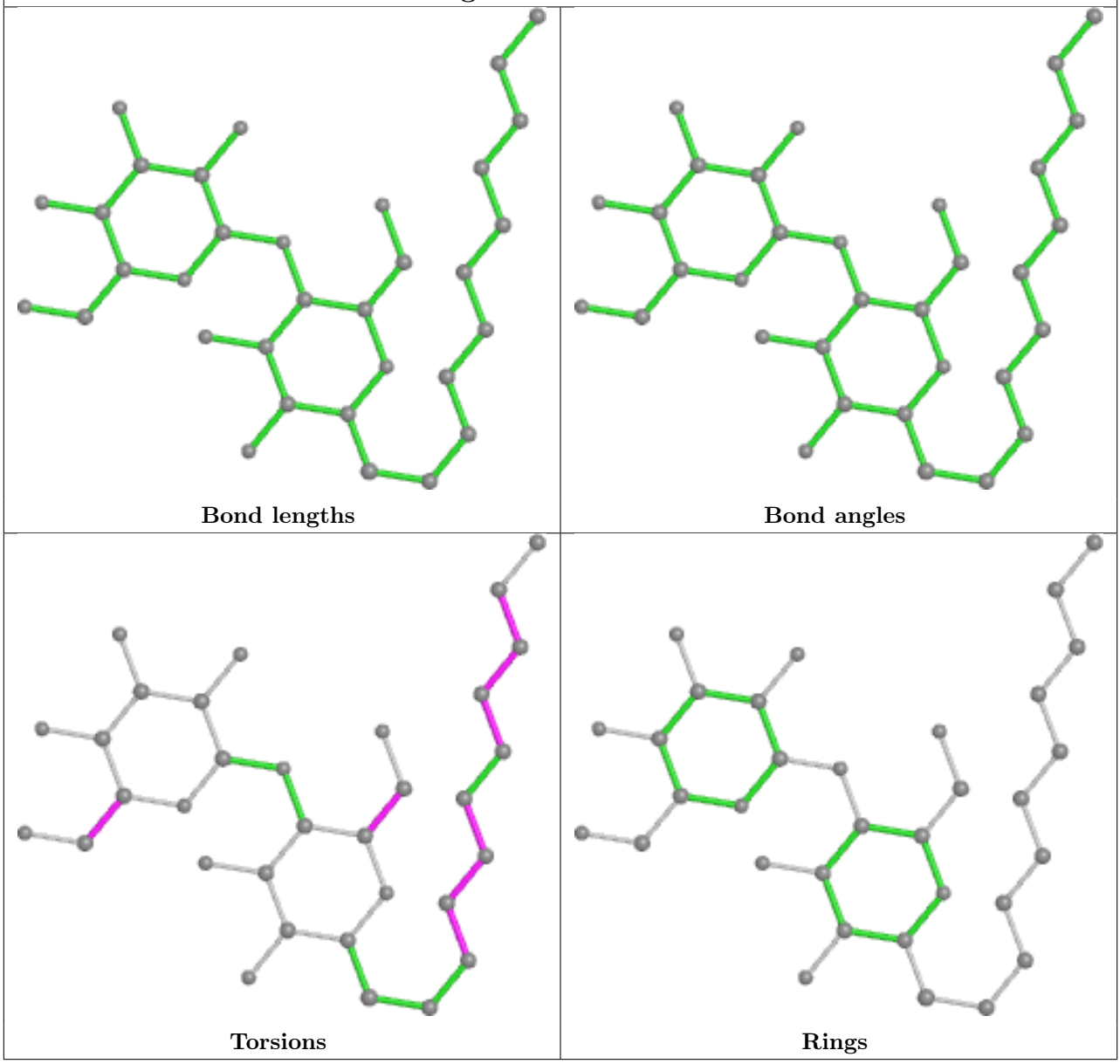


Torsions



Rings

Ligand 10M B 813



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	L	213/213 (100%)	0.20	8 (3%)	40	37	50, 77, 104, 156	0
2	H	225/225 (100%)	0.06	11 (4%)	29	27	55, 76, 124, 187	0
3	B	449/465 (96%)	0.23	29 (6%)	18	18	63, 91, 124, 161	0
4	C	142/146 (97%)	0.07	8 (5%)	24	23	61, 89, 120, 137	0
All	All	1029/1049 (98%)	0.16	56 (5%)	25	24	50, 85, 121, 187	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	B	356	TYR	4.4
4	C	43	ALA	3.9
3	B	355	ALA	3.7
3	B	269	TRP	3.6
2	H	14	PRO	3.5
3	B	254	ILE	3.5
3	B	359	ILE	3.3
4	C	42	GLU	3.3
3	B	210	VAL	3.2
4	C	5	PHE	3.1
2	H	124	SER	3.1
2	H	200	PRO	3.1
1	L	212	ASN	3.0
3	B	382	SER	3.0
3	B	357	ALA	2.9
3	B	374	ILE	2.9
1	L	1	ASP	2.9
2	H	13	ARG	2.8
3	B	270	LEU	2.8
4	C	6	THR	2.8
3	B	266	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
2	H	184	ASP	2.7
3	B	381	ARG	2.7
3	B	128	LEU	2.7
1	L	186	TYR	2.7
1	L	181	LEU	2.7
4	C	70	GLU	2.7
3	B	314	THR	2.7
3	B	239	LYS	2.6
4	C	34	LYS	2.6
1	L	151	ASP	2.5
3	B	268	TYR	2.5
3	B	362	THR	2.5
3	B	438	TRP	2.5
1	L	206	VAL	2.4
3	B	214	TRP	2.4
3	B	413	TRP	2.3
4	C	30	TYR	2.3
4	C	9	MET	2.3
2	H	196	SER	2.3
2	H	84	SER	2.3
2	H	85	SER	2.3
3	B	240	TRP	2.2
2	H	201	SER	2.2
2	H	182	GLN	2.2
1	L	132	VAL	2.2
1	L	180	THR	2.2
2	H	220	LYS	2.1
3	B	414	LEU	2.1
3	B	358	MET	2.1
3	B	218	MET	2.1
3	B	387	MET	2.1
3	B	89	TYR	2.1
3	B	131	THR	2.1
3	B	352	PHE	2.0
3	B	296	ASN	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 monosaccharides in this entry.

6.4 Ligands (i)

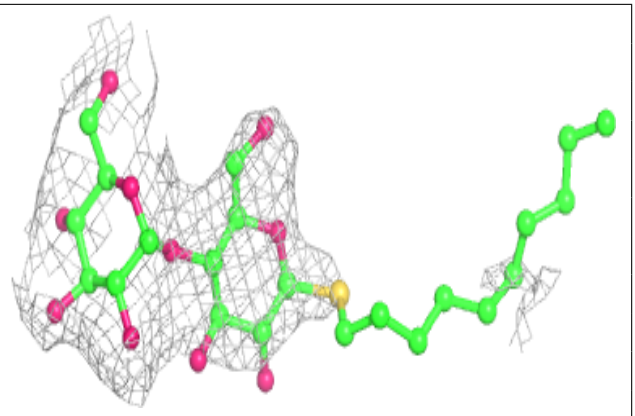
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

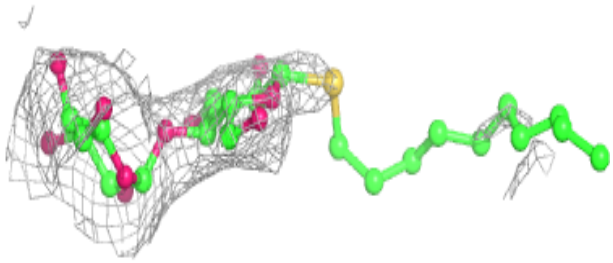
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
9	10M	B	812	33/33	0.74	0.56	112,142,153,157	0
9	10M	B	813	33/33	0.82	0.37	85,156,182,196	0
5	HEM	B	802	43/43	0.96	0.37	71,85,93,102	0
8	XE	B	805	1/1	0.96	0.15	98,98,98,98	1
5	HEM	B	801	43/43	0.97	0.32	62,71,83,96	0
10	CA	B	814	1/1	0.97	0.28	97,97,97,97	0
11	HEC	C	201	43/43	0.97	0.23	56,63,82,87	0
8	XE	B	810	1/1	0.98	0.12	87,87,87,87	1
8	XE	B	811	1/1	0.98	0.31	89,89,89,89	1
7	O	B	804	1/1	0.98	0.32	80,80,80,80	0
6	FE	B	803	1/1	0.98	0.25	80,80,80,80	0
8	XE	B	807	1/1	0.98	0.10	94,94,94,94	1
8	XE	B	808	1/1	0.98	0.10	93,93,93,93	1
8	XE	B	806	1/1	0.99	0.32	91,91,91,91	1
8	XE	B	809	1/1	1.00	0.32	85,85,85,85	1

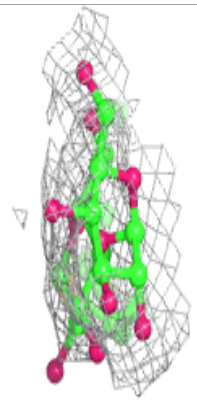
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 10M B 812:

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

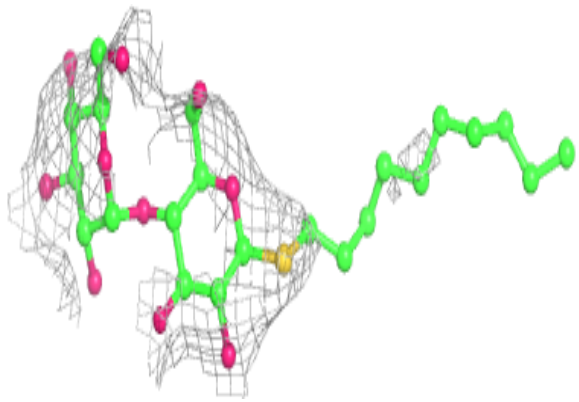


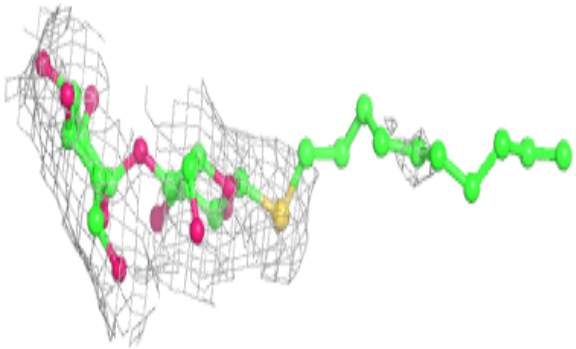


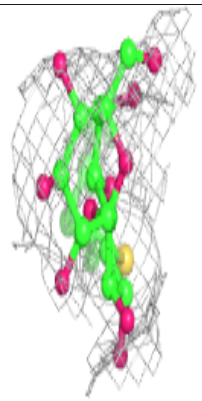


Electron density around 10M B 813:

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

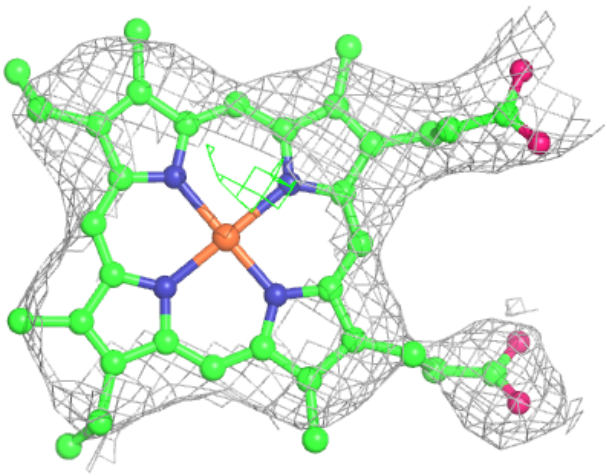


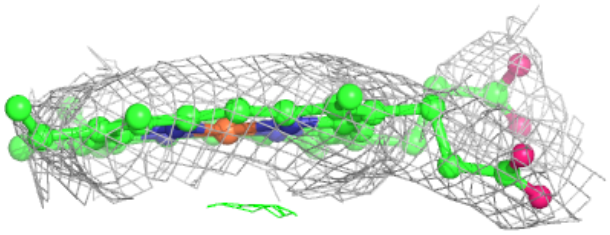


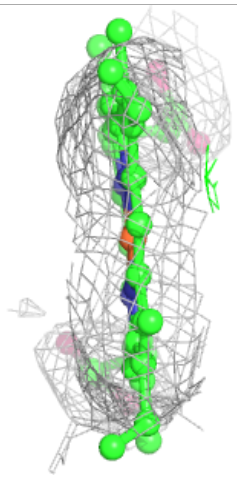


Electron density around HEM B 802:

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

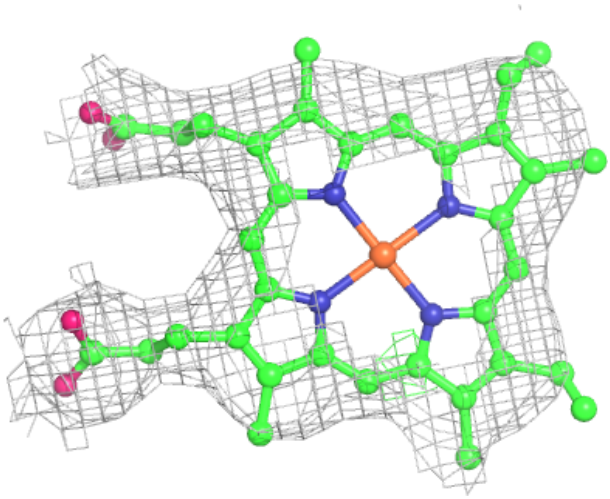


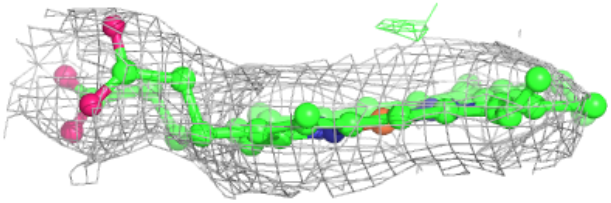


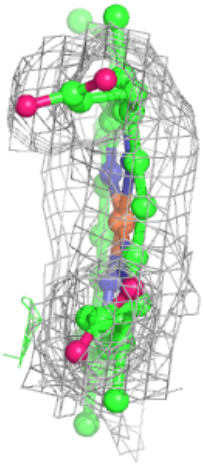


Electron density around HEM B 801:

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

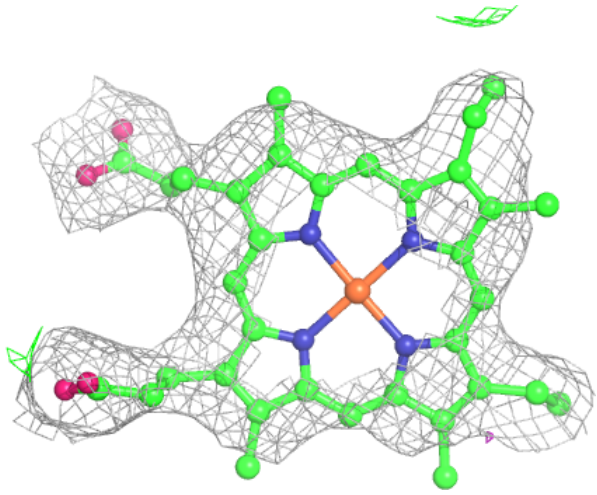


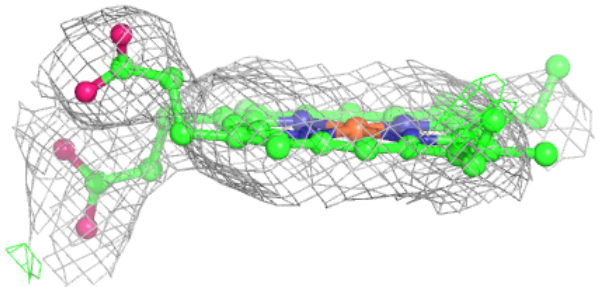


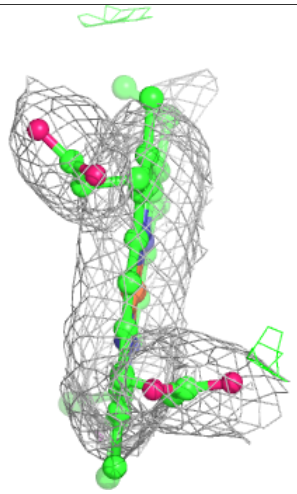


Electron density around HEC C 201:

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.