



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:35 PM UTC

PDB ID : 7JJ1 / pdb_00007jj1
Title : Crystal structure of the sterol 14alpha-demethylase-ferredoxin (CYP51-fx)
heme domain and architectural comparison to the whole fusion protein
Authors : Zhao, B.; Lamb, D.C.
Deposited on : 2020-07-24
Resolution : 3.00 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	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

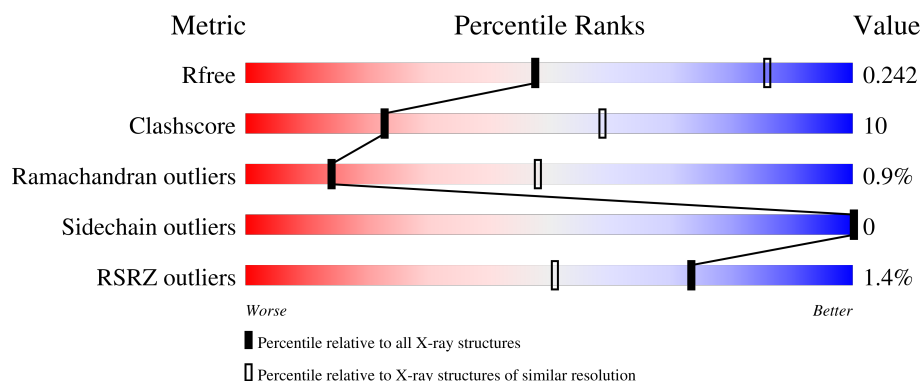
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

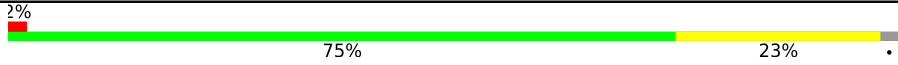
The reported resolution of this entry is 3.00 Å.

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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	450	
1	B	450	
1	C	450	

2 Entry composition [i](#)

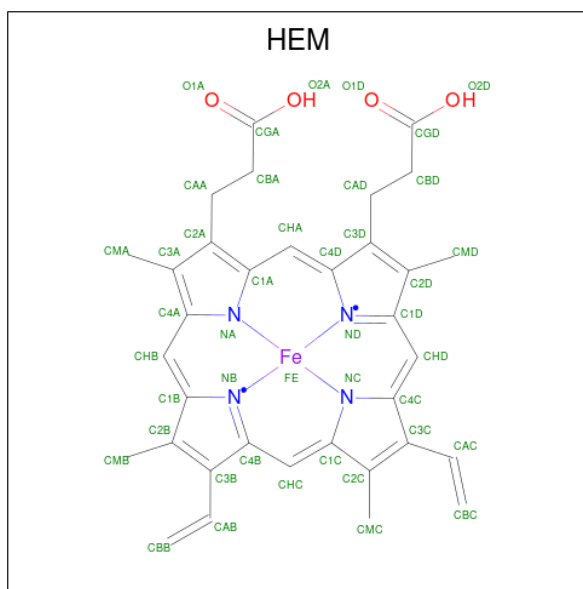
There are 5 unique types of molecules in this entry. The entry contains 10661 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 P450 51.

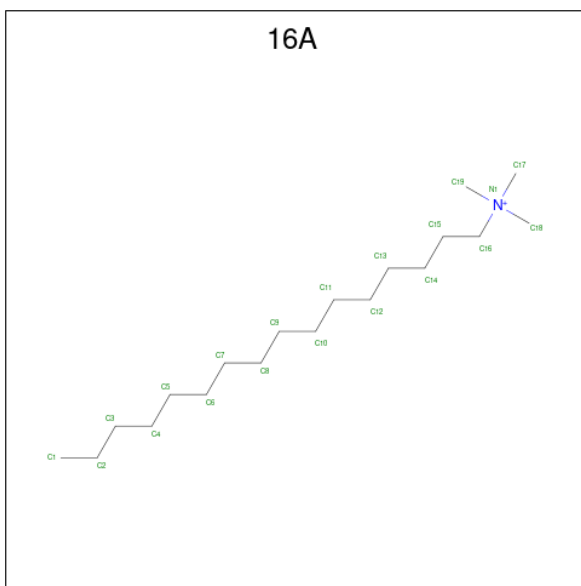
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	0	1	0
			3465	2212	612	618	23			
1	B	434	Total	C	N	O	S	0	1	0
			3400	2182	595	602	21			
1	C	442	Total	C	N	O	S	0	1	0
			3535	2262	628	622	23			

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$) (labeled as "Ligand of Interest" by depositor).



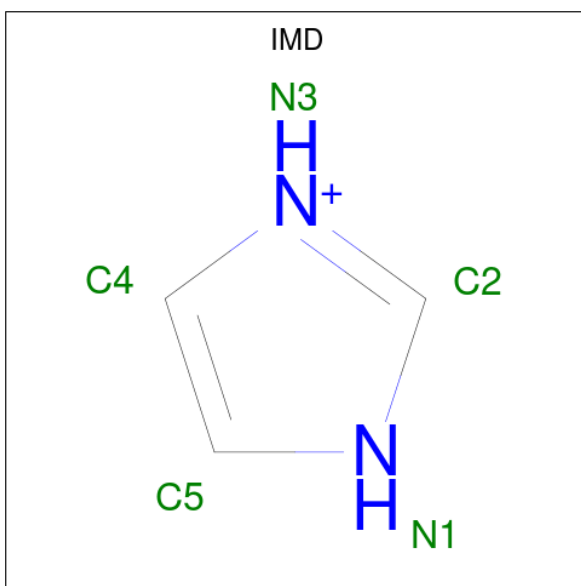
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is CETYL-TRIMETHYL-AMMONIUM (CCD ID: 16A) (formula: $C_{19}H_{42}N$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			20	19	1		
3	B	1	Total	C	N	0	0
			20	19	1		
3	C	1	Total	C	N	0	0
			20	19	1		

- Molecule 4 is IMIDAZOLE (CCD ID: IMD) (formula: $C_3H_5N_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N 5 3 2	0	0
4	A	1	Total C N 5 3 2	0	0
4	A	1	Total C N 5 3 2	0	0
4	B	1	Total C N 5 3 2	0	0
4	C	1	Total C N 5 3 2	0	0
4	C	1	Total C N 5 3 2	0	0

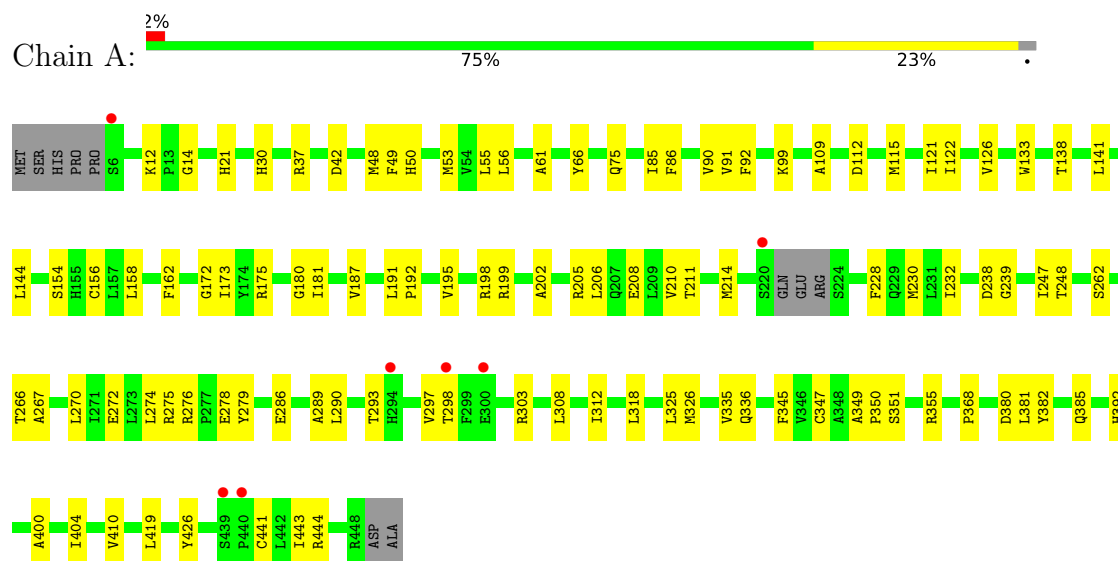
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	19	Total O 19 19	0	0
5	B	14	Total O 14 14	0	0
5	C	9	Total O 9 9	0	0

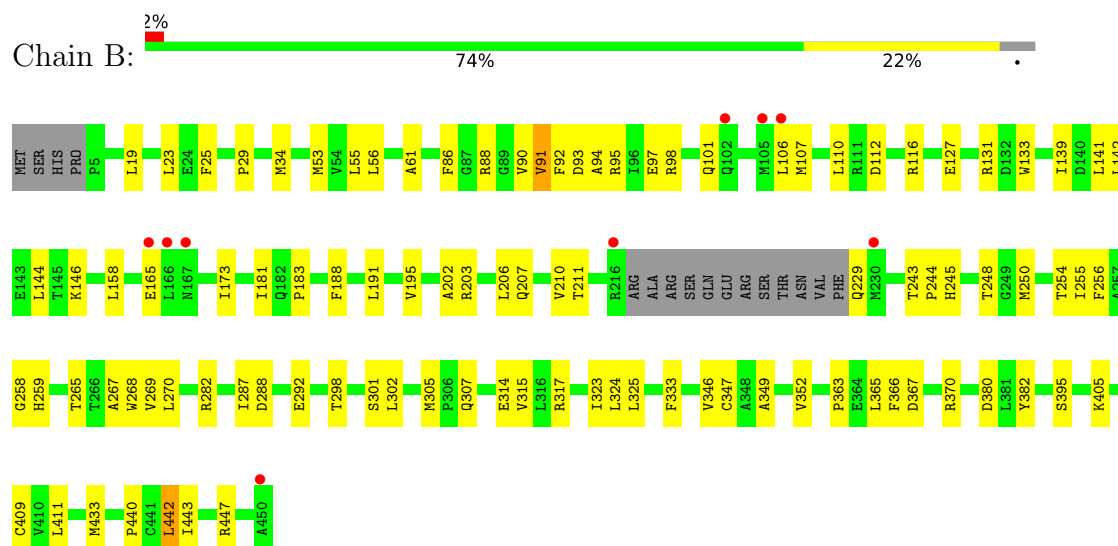
3 Residue-property plots [i](#)

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.

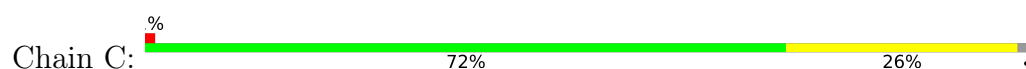
• Molecule 1: Cytochrome P450 51

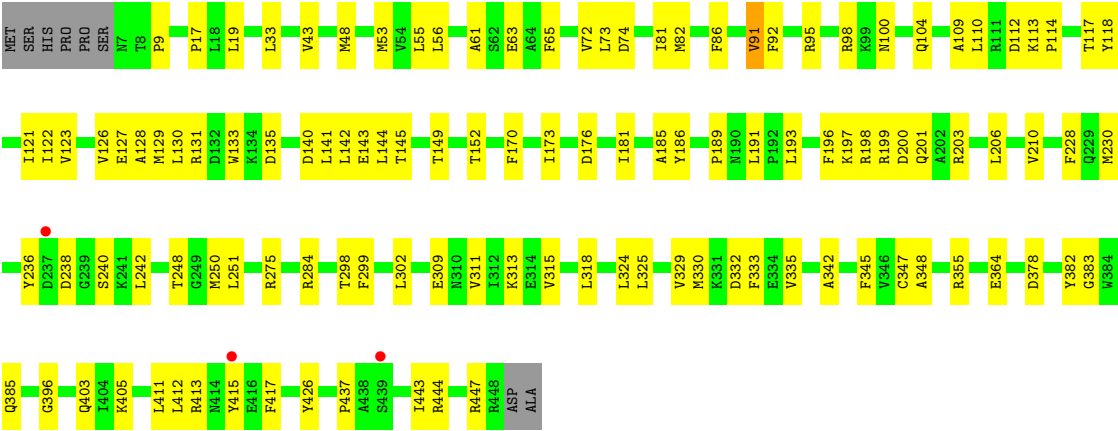


• Molecule 1: Cytochrome P450 51



• Molecule 1: Cytochrome P450 51





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	157.68Å 164.75Å 187.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.52 – 3.00 29.52 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.52-3.00) 99.7 (29.52-3.00)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.81 (at 2.96Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.196 , 0.237 0.198 , 0.242	Depositor DCC
R_{free} test set	5354 reflections (10.81%)	wwPDB-VP
Wilson B-factor (Å ²)	45.6	Xtriage
Anisotropy	0.792	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 39.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.037 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	10661	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.91% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, IMD, 16A

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	0.21	0/3550	0.53	0/4810
1	B	0.19	0/3485	0.52	0/4724
1	C	0.21	0/3622	0.54	0/4904
All	All	0.21	0/10657	0.53	0/14438

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3465	0	3384	62	0
1	B	3400	0	3300	63	0
1	C	3535	0	3515	86	0
2	A	43	0	30	3	0
2	B	43	0	30	5	0
2	C	43	0	30	2	0
3	A	20	0	42	3	0
3	B	20	0	42	2	0
3	C	20	0	42	4	0
4	A	15	0	15	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	5	0	5	0	0
4	C	10	0	10	0	0
5	A	19	0	0	0	0
5	B	14	0	0	0	0
5	C	9	0	0	0	0
All	All	10661	0	10445	218	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 10.

All (218) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:91:VAL:HG13	1:C:92:PHE:H	1.37	0.89
1:C:73:LEU:HD21	1:C:333:PHE:HE2	1.38	0.87
1:C:206:LEU:HD22	1:C:248:THR:HG23	1.61	0.83
1:B:207:GLN:HA	1:B:248:THR:HG21	1.63	0.80
1:A:138:THR:HG22	1:A:444:ARG:HG2	1.63	0.79
1:B:173:ILE:HG21	1:B:206:LEU:HG	1.65	0.78
1:B:302:LEU:HB2	1:B:305:MET:HE2	1.66	0.75
1:B:250:MET:HB3	3:B:502:16A:H13	1.69	0.73
1:B:298:THR:H	1:B:301:SER:HB3	1.52	0.73
1:C:135:ASP:HA	1:C:447:ARG:HE	1.52	0.72
1:C:95:ARG:HH11	1:C:98:ARG:HD3	1.55	0.72
1:C:142:LEU:HD21	1:C:437:PRO:HD2	1.72	0.72
1:C:242:LEU:HD22	1:C:250:MET:HE1	1.72	0.70
1:C:199:ARG:NH1	1:C:200:ASP:OD1	2.24	0.69
1:B:173:ILE:HG23	1:B:202:ALA:HB1	1.76	0.68
1:A:92:PHE:H	1:A:99:LYS:HG3	1.59	0.67
1:C:142:LEU:HA	1:C:145:THR:HG22	1.76	0.67
1:C:298:THR:HG22	1:C:299:PHE:H	1.58	0.66
1:C:81:ILE:HG13	1:C:185:ALA:HB3	1.78	0.65
1:C:100:ASN:O	1:C:104:GLN:HG2	1.99	0.63
1:C:189:PRO:O	1:C:199:ARG:NH2	2.30	0.63
1:C:228:PHE:CD1	1:C:251:LEU:HD11	2.33	0.63
1:A:326:MET:H	4:A:504:IMD:H2	1.64	0.62
1:B:141:LEU:HD12	1:B:443:ILE:HD13	1.80	0.62
1:A:48:MET:HB2	1:A:53:MET:CE	2.29	0.62
1:B:302:LEU:HD12	1:B:405:LYS:HB3	1.82	0.62
1:C:170:PHE:HA	1:C:173:ILE:HG12	1.80	0.62
1:C:141:LEU:HD12	1:C:443:ILE:HG21	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50[B]:HIS:CD2	1:B:195:VAL:HG21	2.34	0.62
2:A:501:HEM:HMD3	3:A:502:16A:H13	1.81	0.62
1:C:82:MET:HE1	1:C:181:ILE:HD13	1.82	0.61
1:B:142:LEU:HG	1:B:146:LYS:HE3	1.82	0.61
1:C:109:ALA:HA	1:C:230:MET:SD	2.41	0.61
1:A:56:LEU:HB3	1:A:61:ALA:HB1	1.83	0.60
1:B:90:VAL:HG12	1:B:91:VAL:HG23	1.84	0.60
1:A:272:GLU:OE1	1:A:276:ARG:NH1	2.35	0.60
1:B:181:ILE:HD11	1:B:256[B]:PHE:HE2	1.68	0.59
1:A:266:THR:HG22	1:A:404:ILE:HG23	1.84	0.59
1:C:82:MET:HE3	3:C:502:16A:H82	1.84	0.58
1:B:255:ILE:O	1:B:259:HIS:HB3	2.03	0.58
1:A:350:PRO:HA	1:A:385:GLN:HE21	1.69	0.57
1:C:284:ARG:NH1	1:C:412:LEU:O	2.37	0.57
1:B:250:MET:O	1:B:254:THR:HG23	2.04	0.57
1:A:228:PHE:O	1:A:232:ILE:HG12	2.04	0.57
1:A:419:LEU:HA	1:A:443:ILE:HD12	1.87	0.57
1:B:34:MET:HG2	1:B:352:VAL:HG22	1.86	0.57
2:C:501:HEM:HMD2	3:C:502:16A:H13	1.87	0.57
1:B:56:LEU:HB3	1:B:61:ALA:HB1	1.87	0.56
2:B:501:HEM:HMD3	3:B:502:16A:H61	1.87	0.56
1:C:48:MET:HE3	1:C:53:MET:HG3	1.86	0.56
1:C:236:TYR:HB2	1:C:240:SER:HB2	1.87	0.56
1:C:332:ASP:OD1	1:C:342:ALA:N	2.38	0.55
1:C:133:TRP:HZ2	1:C:144:LEU:HD13	1.71	0.55
1:B:244:PRO:O	1:B:248:THR:HG23	2.06	0.55
1:C:81:ILE:HB	1:C:186:TYR:CE1	2.42	0.55
1:C:86:PHE:O	1:C:203:ARG:NH2	2.33	0.55
1:A:158:LEU:HB3	1:A:162:PHE:CD2	2.42	0.55
1:C:129:MET:HE2	1:C:133:TRP:HE1	1.73	0.54
1:C:140:ASP:HB3	1:C:143:GLU:HB2	1.88	0.54
1:B:325:LEU:HD22	2:B:501:HEM:HBA2	1.89	0.54
1:C:112:ASP:O	1:C:114:PRO:HD2	2.08	0.54
1:A:121:ILE:HD11	1:A:156:CYS:SG	2.48	0.54
1:C:196:PHE:HA	1:C:199:ARG:HG2	1.88	0.54
1:A:279:TYR:CZ	1:A:368:PRO:HD2	2.43	0.54
1:A:173:ILE:HG23	1:A:202:ALA:HB1	1.89	0.53
1:B:365:LEU:HD22	1:B:367:ASP:HB2	1.89	0.53
1:C:236:TYR:CE2	1:C:242:LEU:HD23	2.44	0.53
1:B:94:ALA:HB3	1:B:98:ARG:HB2	1.90	0.53
1:B:367:ASP:O	1:B:370:ARG:HG3	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:110:LEU:HD13	1:C:396:GLY:HA2	1.91	0.52
1:A:133:TRP:HZ2	1:A:144:LEU:HD13	1.74	0.52
1:B:317:ARG:NH2	1:B:363:PRO:O	2.38	0.52
1:C:48:MET:HE3	1:C:53:MET:HE2	1.91	0.52
1:C:197:LYS:O	1:C:201:GLN:HG3	2.08	0.52
1:C:206:LEU:CD2	1:C:248:THR:HG23	2.36	0.52
1:B:265:THR:O	1:B:269:VAL:HG23	2.09	0.52
1:B:447:ARG:HG3	1:B:447:ARG:HH11	1.74	0.52
1:B:19:LEU:HD13	1:B:23:LEU:HD11	1.91	0.52
1:C:91:VAL:HG13	1:C:92:PHE:N	2.16	0.52
1:C:56:LEU:HD12	1:C:65:PHE:HB2	1.91	0.51
1:A:55:LEU:HD11	1:A:349:ALA:HB2	1.92	0.51
1:A:48:MET:HB2	1:A:53:MET:HE1	1.91	0.51
1:A:335:VAL:HG12	1:A:336:GLN:HG3	1.91	0.51
1:C:112:ASP:C	1:C:114:PRO:HD2	2.36	0.51
1:A:205:ARG:NH1	1:A:208:GLU:OE1	2.44	0.51
1:A:37:ARG:NH1	1:A:42:ASP:OD1	2.43	0.51
1:C:383:GLY:O	1:C:385:GLN:HG3	2.11	0.51
1:B:25:PHE:CZ	1:B:324:LEU:HD11	2.45	0.51
1:B:127:GLU:O	1:B:131:ARG:HG3	2.11	0.51
1:C:56:LEU:HB3	1:C:61:ALA:HB1	1.92	0.51
1:C:43:VAL:HG22	1:C:56:LEU:HD22	1.93	0.51
1:A:180:GLY:HA3	1:A:199:ARG:HD3	1.93	0.50
1:C:309:GLU:HG2	1:C:405:LYS:HE2	1.93	0.50
1:A:49:PHE:CZ	1:A:187:VAL:HG22	2.45	0.50
1:C:173:ILE:HD12	1:C:206:LEU:HB2	1.93	0.50
1:B:183:PRO:HG3	1:B:433:MET:HE3	1.94	0.50
1:B:107:MET:HE1	1:B:395:SER:HB2	1.94	0.50
1:A:262:SER:O	1:A:266:THR:HG23	2.11	0.50
1:A:286:GLU:O	1:A:290:LEU:HD23	2.11	0.50
1:A:266:THR:HG22	1:A:404:ILE:HA	1.94	0.50
1:B:258:GLY:HA2	2:B:501:HEM:HMC2	1.95	0.49
1:C:126:VAL:O	1:C:130:LEU:HD13	2.13	0.49
1:B:207:GLN:O	1:B:211:THR:HG23	2.12	0.48
1:C:122:ILE:HD13	1:C:403:GLN:HA	1.95	0.48
2:C:501:HEM:CMD	3:C:502:16A:H13	2.43	0.48
1:A:232:ILE:HD13	1:A:247:ILE:HD13	1.95	0.48
1:B:243:THR:HG22	1:B:245:HIS:H	1.79	0.48
1:A:109:ALA:HB2	1:A:230:MET:HG3	1.96	0.48
1:C:122:ILE:HG12	1:C:152:THR:OG1	2.14	0.48
1:A:289:ALA:O	1:A:293:THR:HG22	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:112:ASP:O	1:B:116:ARG:HG3	2.14	0.48
1:B:282:ARG:NE	1:B:307:GLN:OE1	2.42	0.48
1:C:236:TYR:O	1:C:238:ASP:N	2.43	0.48
1:A:210:VAL:HG13	1:A:228:PHE:CZ	2.49	0.48
1:A:274:LEU:HD13	1:A:443:ILE:HD11	1.95	0.47
1:B:86:PHE:O	1:B:203:ARG:NE	2.44	0.47
1:B:380:ASP:O	1:B:382:TYR:N	2.42	0.47
1:C:118:TYR:O	1:C:122:ILE:HG13	2.15	0.47
1:C:117:THR:O	1:C:121:ILE:HG13	2.15	0.47
1:A:12:LYS:HD3	1:A:21:HIS:CD2	2.49	0.47
1:A:195:VAL:HG12	1:A:198:ARG:HH21	1.79	0.47
1:A:75:GLN:NE2	1:A:392:HIS:CG	2.82	0.47
1:B:288:ASP:O	1:B:292:GLU:HG2	2.14	0.47
1:C:149:THR:HA	1:C:152:THR:HG22	1.97	0.47
1:C:275:ARG:NH2	1:C:426:TYR:O	2.47	0.47
1:A:211:THR:HA	1:A:214:MET:HE2	1.97	0.47
1:A:297:VAL:HG21	1:A:410:VAL:HG22	1.97	0.47
2:A:501:HEM:HBC2	2:A:501:HEM:CMC	2.45	0.46
1:A:318:LEU:O	1:A:355:ARG:NH2	2.44	0.46
1:B:141:LEU:CD2	1:B:267:ALA:HA	2.45	0.46
1:C:413:ARG:HD2	1:C:413:ARG:O	2.16	0.46
1:B:380:ASP:C	1:B:382:TYR:H	2.23	0.46
1:A:115:MET:HE3	1:A:115:MET:HB3	1.74	0.46
1:C:191:LEU:C	1:C:193:LEU:H	2.24	0.46
1:B:106:LEU:O	1:B:110:LEU:HG	2.16	0.46
1:B:333:PHE:HZ	1:B:346:VAL:HG11	1.81	0.45
1:A:419:LEU:HD22	1:A:426:TYR:CE2	2.51	0.45
1:B:287:ILE:HD13	1:B:409:CYS:HA	1.97	0.45
1:C:63:GLU:HG3	1:C:382:TYR:CE2	2.51	0.45
1:C:95:ARG:HH11	1:C:98:ARG:CD	2.27	0.45
1:C:298:THR:HG22	1:C:299:PHE:N	2.29	0.45
1:C:329:VAL:HG12	1:C:342:ALA:HA	1.98	0.45
1:A:275:ARG:NH2	1:A:426:TYR:O	2.50	0.45
1:B:97:GLU:O	1:B:101:GLN:HG3	2.17	0.45
2:A:501:HEM:HBC2	2:A:501:HEM:HMC2	1.99	0.45
1:C:33:LEU:HD22	1:C:55:LEU:HD22	1.99	0.44
1:C:318:LEU:O	1:C:355:ARG:NH2	2.48	0.44
1:A:380:ASP:C	1:A:382:TYR:H	2.25	0.44
1:A:238:ASP:OD1	1:A:239:GLY:N	2.51	0.44
1:B:88:ARG:O	1:B:93:ASP:HB3	2.17	0.44
1:A:141:LEU:HD22	1:A:267:ALA:HA	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158:LEU:HA	1:B:229:GLN:OE1	2.16	0.44
1:B:55:LEU:HA	1:B:347:CYS:HB2	1.99	0.44
1:C:206:LEU:O	1:C:210:VAL:HG23	2.17	0.44
1:C:313:LYS:NZ	1:C:378:ASP:OD2	2.42	0.44
1:A:53:MET:SD	1:A:345:PHE:HB2	2.58	0.44
1:B:91:VAL:HG12	1:B:92:PHE:N	2.32	0.44
1:B:188:PHE:HB3	1:B:191:LEU:HG	1.99	0.44
1:C:228:PHE:HD1	1:C:251:LEU:HD11	1.81	0.44
1:C:325:LEU:O	1:C:347:CYS:HA	2.18	0.44
1:C:123:VAL:O	1:C:127:GLU:HG2	2.17	0.44
1:A:278:GLU:CD	1:A:278:GLU:H	2.26	0.44
1:A:400:ALA:O	1:A:404:ILE:HG13	2.17	0.44
1:C:9:PRO:HB3	1:C:43:VAL:HG12	2.00	0.44
1:C:309:GLU:CG	1:C:405:LYS:HE2	2.48	0.43
1:A:141:LEU:HD21	1:A:270:LEU:HD12	2.00	0.43
1:B:110:LEU:HD21	2:B:501:HEM:HBC1	1.99	0.43
3:A:502:16A:H192	3:A:502:16A:H151	1.73	0.43
1:C:355:ARG:HD3	1:C:364:GLU:OE2	2.18	0.43
1:A:86:PHE:C	1:A:90:VAL:HG21	2.44	0.43
1:A:308:LEU:O	1:A:312:ILE:HG13	2.18	0.43
1:B:258:GLY:HA2	2:B:501:HEM:CMC	2.48	0.43
1:B:365:LEU:HD23	1:B:366:PHE:N	2.34	0.43
1:B:133:TRP:HZ2	1:B:144:LEU:HD13	1.83	0.43
1:A:66:TYR:HE2	1:A:385:GLN:OE1	2.01	0.43
1:B:53:MET:HE3	1:B:347:CYS:SG	2.59	0.43
1:C:72:VAL:CG1	1:C:333:PHE:HD2	2.32	0.43
1:C:126:VAL:HG12	1:C:130:LEU:HD13	2.01	0.43
1:B:270:LEU:HD11	1:B:411:LEU:HD23	1.99	0.43
1:C:53:MET:SD	1:C:345:PHE:HB2	2.59	0.43
1:C:210:VAL:HG13	1:C:228:PHE:CE1	2.54	0.43
1:A:325:LEU:O	1:A:347:CYS:HA	2.18	0.43
1:B:25:PHE:HZ	1:B:324:LEU:HD11	1.84	0.43
1:A:50[B]:HIS:HD2	1:B:195:VAL:HG21	1.81	0.42
1:B:323:ILE:O	1:B:349:ALA:HA	2.19	0.42
1:C:130:LEU:HG	1:C:415:TYR:CE2	2.54	0.42
1:C:128:ALA:O	1:C:131:ARG:HG2	2.19	0.42
3:C:502:16A:H22	3:C:502:16A:H51	1.83	0.42
1:C:122:ILE:O	1:C:126:VAL:HG23	2.20	0.42
1:A:14:GLY:O	1:A:21:HIS:NE2	2.53	0.42
1:C:228:PHE:HA	1:C:251:LEU:HD21	2.01	0.42
1:C:333:PHE:CE1	1:C:335:VAL:HG23	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:VAL:HB	1:B:248:THR:HG22	2.02	0.42
1:C:74:ASP:HB2	1:C:330:MET:SD	2.60	0.42
1:C:311:VAL:O	1:C:315:VAL:HG23	2.20	0.42
1:A:191:LEU:HD12	1:A:192:PRO:HD2	2.02	0.42
1:A:154:SER:O	1:A:158:LEU:HB2	2.20	0.41
1:B:142:LEU:O	1:B:146:LYS:HG3	2.20	0.41
1:A:122:ILE:O	1:A:126:VAL:HG23	2.20	0.41
1:A:172:GLY:HA2	1:A:175:ARG:NH1	2.35	0.41
1:C:81:ILE:HB	1:C:186:TYR:CD1	2.56	0.41
1:C:324:LEU:HA	1:C:348:ALA:O	2.20	0.41
1:A:30:HIS:HD1	1:A:351:SER:HB2	1.85	0.41
1:B:139:ILE:O	1:B:442:LEU:HA	2.20	0.41
1:A:206:LEU:HD22	1:A:248:THR:OG1	2.21	0.41
1:C:17:PRO:C	1:C:19:LEU:H	2.28	0.41
1:C:130:LEU:HD11	1:C:411:LEU:HD23	2.03	0.41
1:C:176:ASP:OD2	1:C:198:ARG:HG2	2.21	0.41
3:A:502:16A:H81	4:A:505:IMD:H5	2.03	0.40
1:B:29:PRO:HG3	1:B:433:MET:HE1	2.03	0.40
1:B:268:TRP:HB3	1:B:315:VAL:HG13	2.04	0.40
1:C:95:ARG:N	1:C:95:ARG:HD3	2.36	0.40
1:C:298:THR:O	1:C:302:LEU:HG	2.21	0.40
1:C:417:PHE:HA	1:C:444:ARG:O	2.21	0.40
1:A:85:ILE:CD1	1:A:181:ILE:HG12	2.51	0.40
1:A:380:ASP:O	1:A:381:LEU:HB2	2.20	0.40
1:B:314:GLU:HA	1:B:314:GLU:OE2	2.20	0.40
1:A:326:MET:HG2	1:A:347:CYS:SG	2.62	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	437/450 (97%)	415 (95%)	17 (4%)	5 (1%)	11	43
1	B	431/450 (96%)	404 (94%)	22 (5%)	5 (1%)	10	40
1	C	441/450 (98%)	425 (96%)	14 (3%)	2 (0%)	24	60
All	All	1309/1350 (97%)	1244 (95%)	53 (4%)	12 (1%)	14	48

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	91	VAL
1	A	303	ARG
1	B	95	ARG
1	B	165	GLU
1	C	113	LYS
1	A	441	CYS
1	B	91	VAL
1	B	440	PRO
1	C	91	VAL
1	B	442	LEU
1	A	112	ASP
1	A	298	THR

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	357/389 (92%)	357 (100%)	0	100	100
1	B	341/389 (88%)	341 (100%)	0	100	100
1	C	369/389 (95%)	369 (100%)	0	100	100
All	All	1067/1167 (91%)	1067 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	385	GLN
1	B	102	GLN
1	C	28	ASN
1	C	101	GLN
1	C	167	ASN
1	C	260	HIS
1	C	294	HIS
1	C	385	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

12 ligands are modelled in this entry.

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	$\# Z > 2$
4	IMD	C	504	-	5,5,5	0.47	0	5,5,5	0.53	0
2	HEM	C	501	-	50,50,50	1.72	9 (18%)	67,82,82	0.99	6 (8%)
3	16A	C	502	-	19,19,19	0.55	0	21,21,21	0.53	0
4	IMD	A	504	-	5,5,5	0.46	0	5,5,5	0.53	0
4	IMD	A	505	-	5,5,5	0.50	0	5,5,5	0.50	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HEM	A	501	-	50,50,50	1.73	10 (20%)	67,82,82	1.05	7 (10%)
2	HEM	B	501	4	50,50,50	1.73	10 (20%)	67,82,82	1.00	5 (7%)
4	IMD	B	503	2	5,5,5	0.52	0	5,5,5	0.61	0
3	16A	A	502	-	19,19,19	0.54	0	21,21,21	0.66	0
4	IMD	C	503	-	5,5,5	0.48	0	5,5,5	0.46	0
4	IMD	A	503	-	5,5,5	0.49	0	5,5,5	0.57	0
3	16A	B	502	-	19,19,19	0.55	0	21,21,21	0.68	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
4	IMD	C	504	-	-	-	0/1/1/1
2	HEM	C	501	-	-	3/14/54/54	-
3	16A	C	502	-	-	4/17/17/17	-
4	IMD	A	504	-	-	-	0/1/1/1
4	IMD	A	505	-	-	-	0/1/1/1
2	HEM	A	501	-	-	3/14/54/54	-
2	HEM	B	501	4	-	2/14/54/54	-
4	IMD	B	503	2	-	-	0/1/1/1
3	16A	A	502	-	-	3/17/17/17	-
4	IMD	C	503	-	-	-	0/1/1/1
4	IMD	A	503	-	-	-	0/1/1/1
3	16A	B	502	-	-	2/17/17/17	-

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	HEM	CBB-CAB	4.95	1.54	1.30
2	A	501	HEM	CBB-CAB	4.95	1.54	1.30
2	C	501	HEM	CBB-CAB	4.91	1.54	1.30
2	C	501	HEM	CBC-CAC	4.77	1.53	1.30
2	B	501	HEM	CBC-CAC	4.76	1.53	1.30
2	A	501	HEM	CBC-CAC	4.74	1.53	1.30
2	B	501	HEM	FE-NA	3.04	2.05	1.95
2	C	501	HEM	FE-NA	3.04	2.05	1.95
2	A	501	HEM	FE-NA	3.02	2.05	1.95
2	B	501	HEM	CAB-C3B	2.97	1.55	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	HEM	CAB-C3B	2.96	1.55	1.47
2	A	501	HEM	CAB-C3B	2.94	1.55	1.47
2	C	501	HEM	FE-NC	2.91	2.04	1.95
2	B	501	HEM	FE-NC	2.90	2.04	1.95
2	A	501	HEM	FE-NC	2.86	2.04	1.95
2	C	501	HEM	CAC-C3C	2.72	1.54	1.47
2	A	501	HEM	CAC-C3C	2.65	1.54	1.47
2	B	501	HEM	CAC-C3C	2.62	1.54	1.47
2	A	501	HEM	FE-ND	2.46	2.02	1.94
2	C	501	HEM	FE-ND	2.45	2.02	1.94
2	A	501	HEM	FE-NB	2.45	2.02	1.94
2	B	501	HEM	FE-NB	2.41	2.02	1.94
2	C	501	HEM	FE-NB	2.39	2.02	1.94
2	B	501	HEM	FE-ND	2.38	2.02	1.94
2	C	501	HEM	CMC-C2C	2.28	1.55	1.50
2	A	501	HEM	CMC-C2C	2.25	1.55	1.50
2	B	501	HEM	CMC-C2C	2.24	1.55	1.50
2	A	501	HEM	O2A-CGA	-2.07	1.24	1.30
2	B	501	HEM	O2A-CGA	-2.04	1.24	1.30

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	O1A-CGA-CBA	-2.70	114.52	123.09
2	C	501	HEM	O1A-CGA-CBA	-2.59	114.86	123.09
2	B	501	HEM	CBC-CAC-C3C	-2.55	114.80	127.53
2	B	501	HEM	O1A-CGA-CBA	-2.45	115.31	123.09
2	A	501	HEM	CBB-CAB-C3B	-2.36	115.75	127.53
2	C	501	HEM	C2A-C1A-NA	-2.31	107.59	110.15
2	C	501	HEM	C4A-NA-C1A	2.29	109.56	105.82
2	A	501	HEM	CBC-CAC-C3C	-2.28	116.13	127.53
2	A	501	HEM	C4A-NA-C1A	2.26	109.51	105.82
2	A	501	HEM	C2A-C1A-NA	-2.25	107.66	110.15
2	A	501	HEM	CAD-C3D-C4D	2.23	128.59	124.70
2	C	501	HEM	CBB-CAB-C3B	-2.23	116.39	127.53
2	B	501	HEM	CBB-CAB-C3B	-2.21	116.48	127.53
2	C	501	HEM	CBC-CAC-C3C	-2.20	116.53	127.53
2	B	501	HEM	C1B-NB-C4B	2.18	107.79	105.21
2	B	501	HEM	C4A-NA-C1A	2.18	109.38	105.82
2	C	501	HEM	C1B-NB-C4B	2.06	107.65	105.21
2	A	501	HEM	C1B-NB-C4B	2.03	107.61	105.21

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	16A	C14-C15-C16-N1
3	A	502	16A	C12-C13-C14-C15
2	B	501	HEM	C2C-C3C-CAC-CBC
3	C	502	16A	C10-C11-C12-C13
3	C	502	16A	C15-C16-N1-C18
2	C	501	HEM	CAD-CBD-CGD-O1D
2	A	501	HEM	CAD-CBD-CGD-O1D
2	A	501	HEM	CAD-CBD-CGD-O2D
2	B	501	HEM	C4C-C3C-CAC-CBC
3	B	502	16A	C4-C5-C6-C7
2	C	501	HEM	CAD-CBD-CGD-O2D
2	A	501	HEM	CAA-CBA-CGA-O2A
3	B	502	16A	C2-C3-C4-C5
3	C	502	16A	C15-C16-N1-C19
3	C	502	16A	C15-C16-N1-C17
3	A	502	16A	C9-C10-C11-C12
2	C	501	HEM	C2B-C3B-CAB-CBB

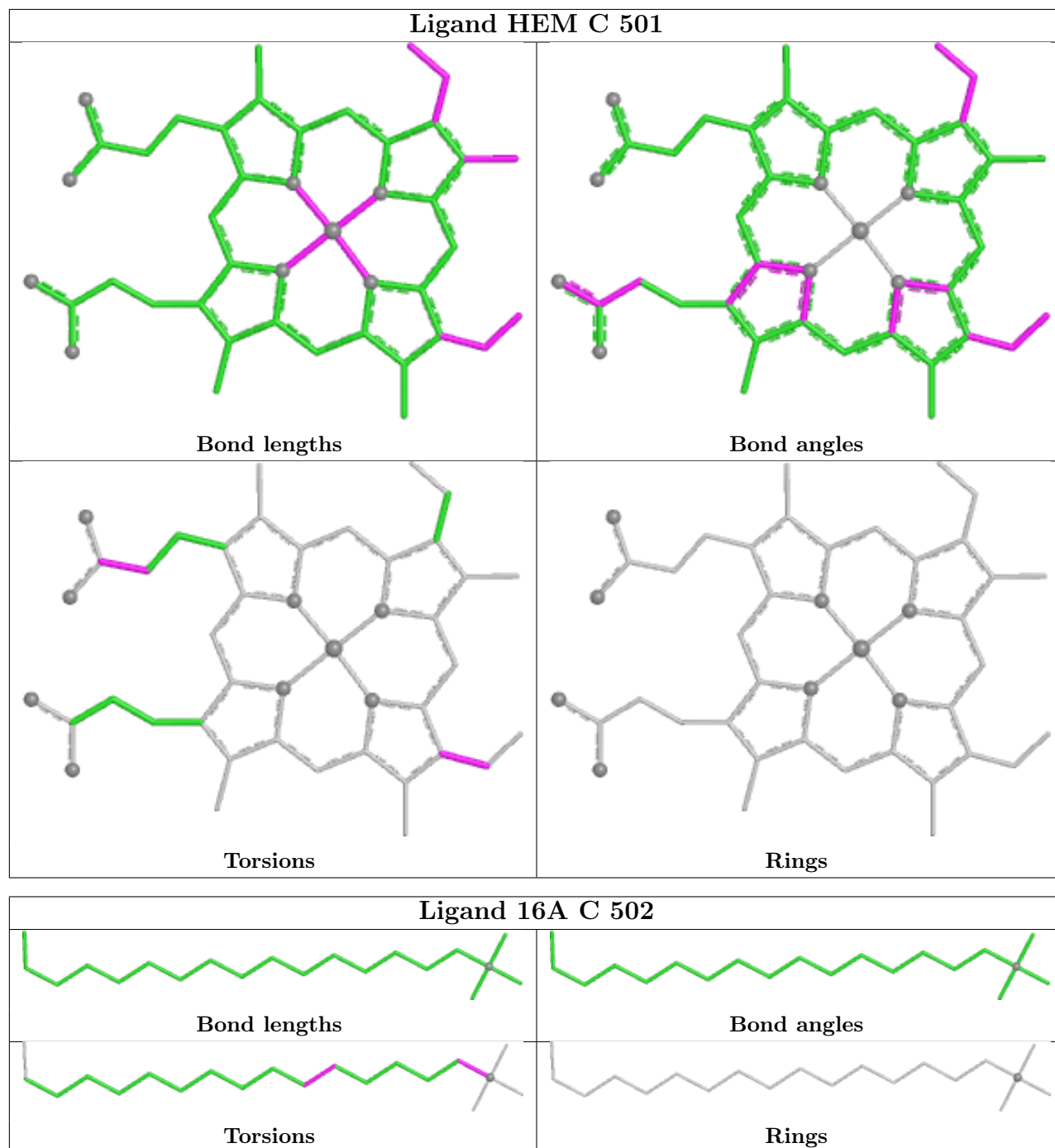
There are no ring outliers.

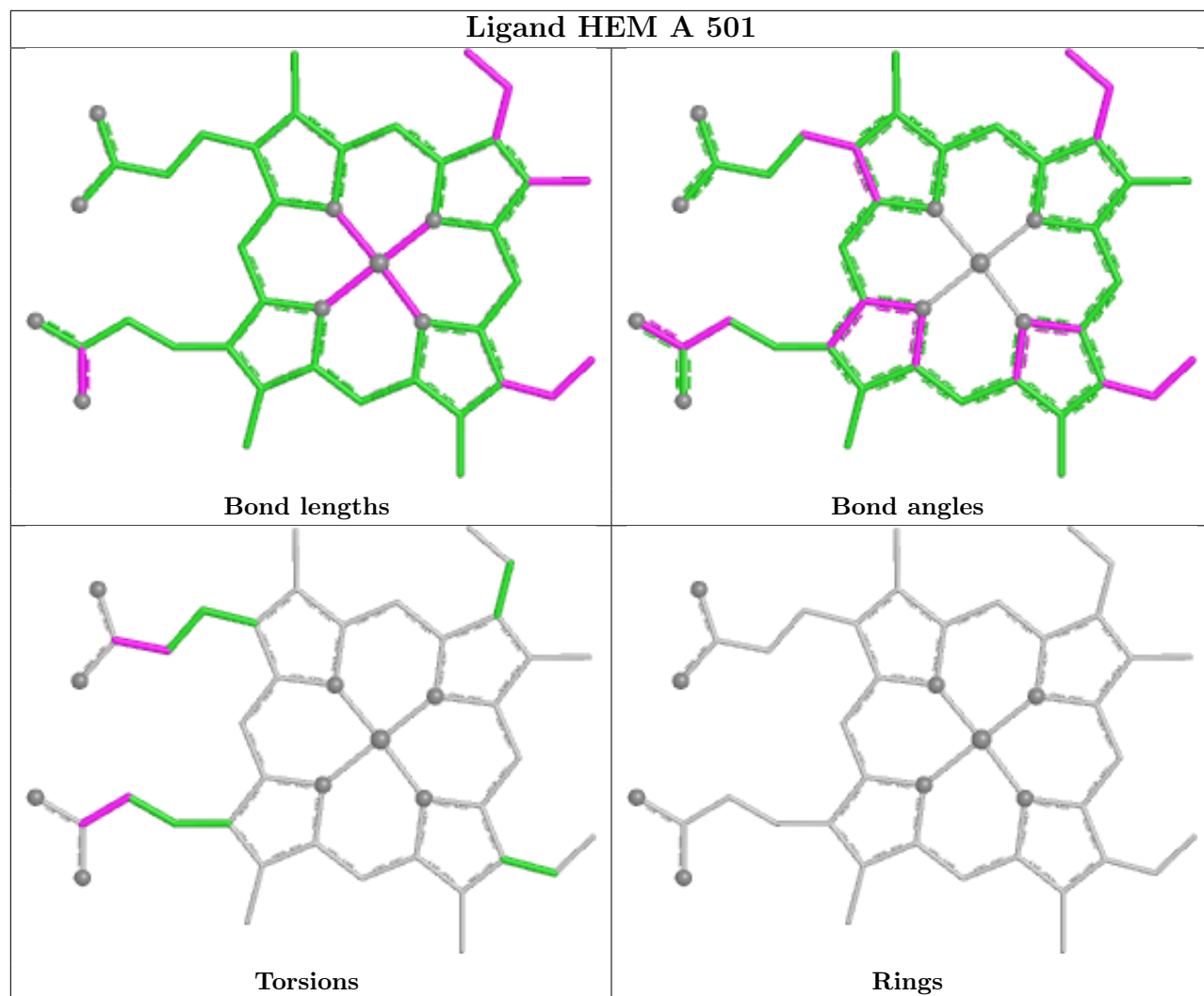
8 monomers are involved in 16 short contacts:

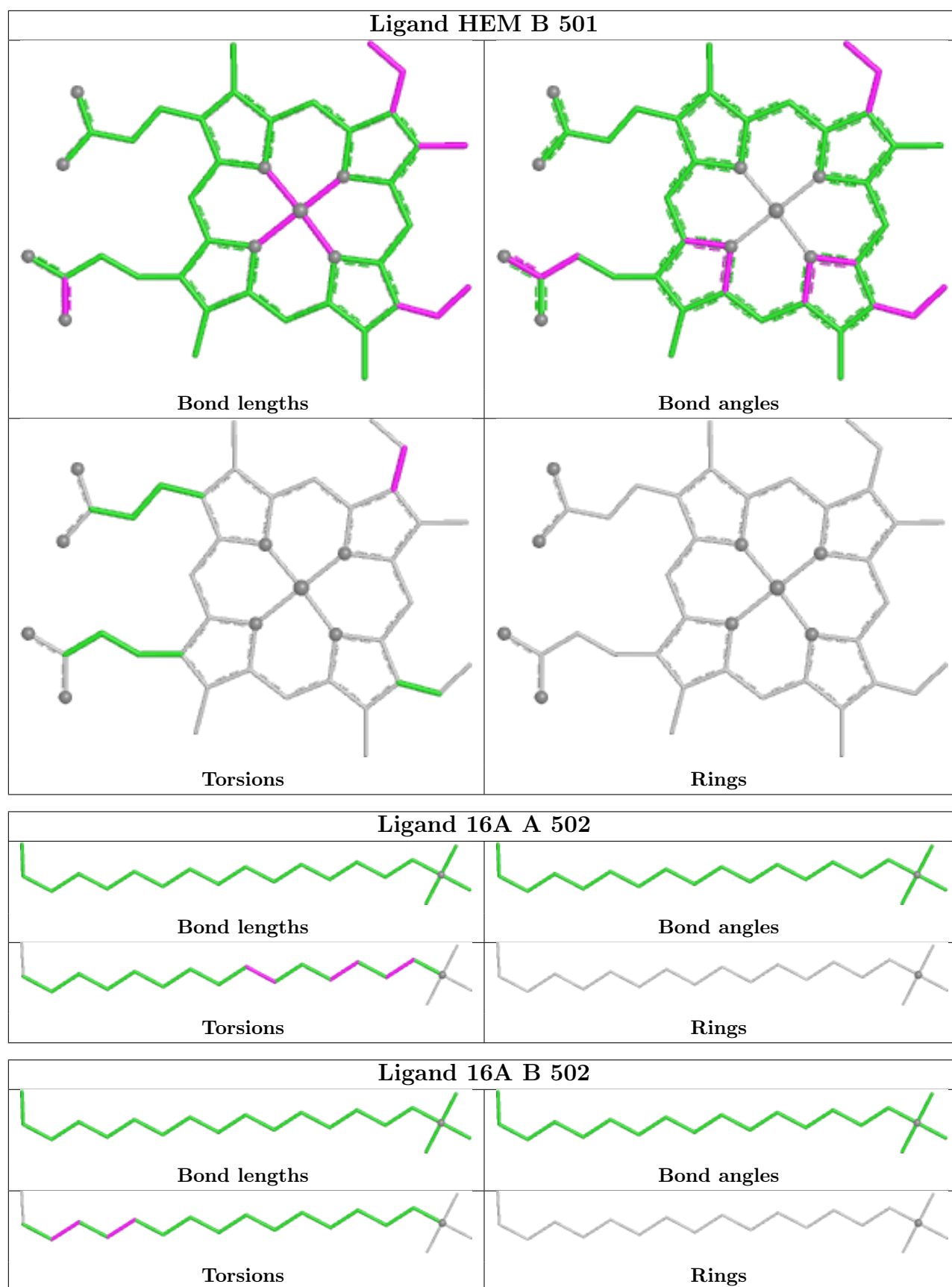
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	501	HEM	2	0
3	C	502	16A	4	0
4	A	504	IMD	1	0
4	A	505	IMD	1	0
2	A	501	HEM	3	0
2	B	501	HEM	5	0
3	A	502	16A	3	0
3	B	502	16A	2	0

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.







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	440/450 (97%)	-0.25	7 (1%) 70 47	24, 47, 79, 102	1 (0%)
1	B	434/450 (96%)	-0.22	9 (2%) 63 40	22, 44, 79, 119	1 (0%)
1	C	442/450 (98%)	-0.37	3 (0%) 84 66	21, 39, 71, 91	1 (0%)
All	All	1316/1350 (97%)	-0.28	19 (1%) 73 51	21, 43, 77, 119	3 (0%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	450	ALA	5.3
1	B	106	LEU	4.4
1	C	415	TYR	3.3
1	A	439	SER	3.0
1	B	165	GLU	2.9
1	B	102	GLN	2.7
1	A	298	THR	2.7
1	B	166	LEU	2.6
1	C	439	SER	2.5
1	B	167	ASN	2.5
1	B	105	MET	2.5
1	B	230	MET	2.4
1	A	300	GLU	2.3
1	A	220	SER	2.3
1	A	6	SER	2.2
1	A	440	PRO	2.2
1	B	216	ARG	2.2
1	A	294	HIS	2.1
1	C	237	ASP	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](#)

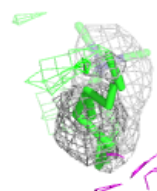
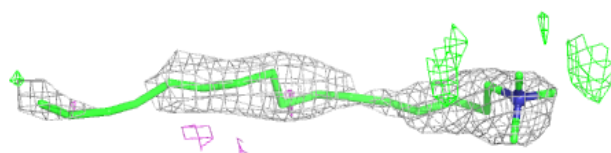
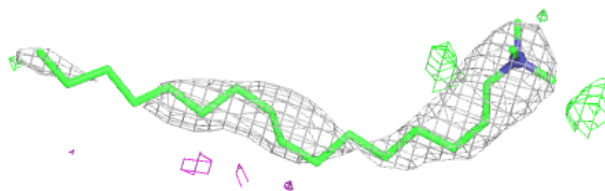
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	IMD	C	504	5/5	0.84	0.23	64,67,78,83	0
4	IMD	A	504	5/5	0.87	0.18	54,60,65,78	0
4	IMD	A	505	5/5	0.88	0.21	71,73,75,82	0
3	16A	B	502	20/20	0.88	0.25	36,56,75,77	0
4	IMD	C	503	5/5	0.91	0.25	24,41,54,59	0
3	16A	C	502	20/20	0.93	0.15	15,35,45,46	0
3	16A	A	502	20/20	0.94	0.12	29,35,49,50	0
4	IMD	A	503	5/5	0.96	0.14	40,41,42,59	0
2	HEM	C	501	43/43	0.98	0.07	18,27,37,48	0
4	IMD	B	503	5/5	0.98	0.09	29,38,47,53	0
2	HEM	A	501	43/43	0.98	0.08	26,41,52,57	0
2	HEM	B	501	43/43	0.98	0.08	19,31,43,60	0

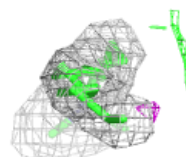
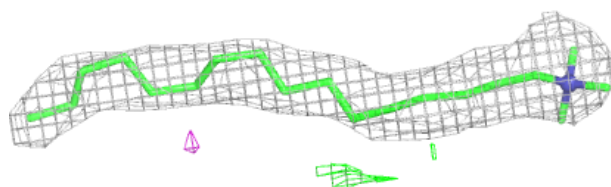
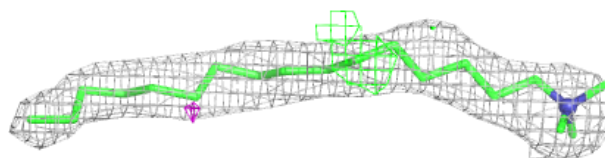
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 16A B 502:

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

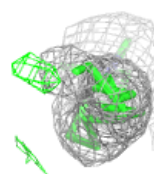
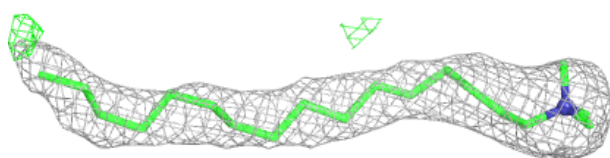
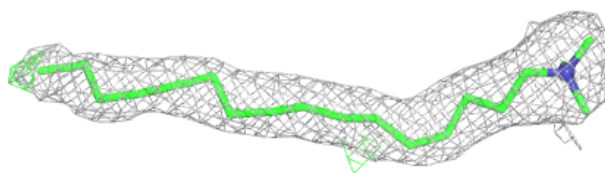
**Electron density around 16A C 502:**

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

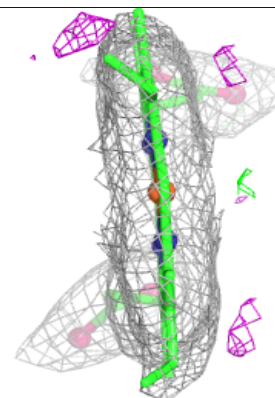
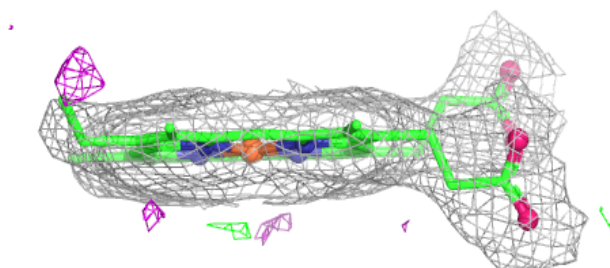
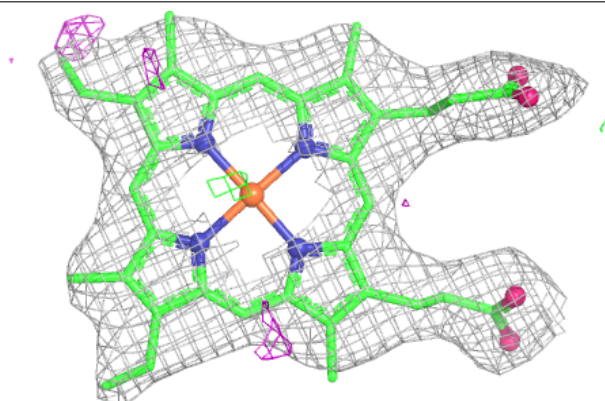


Electron density around 16A A 502:

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

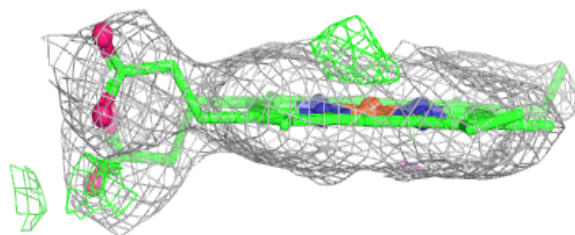
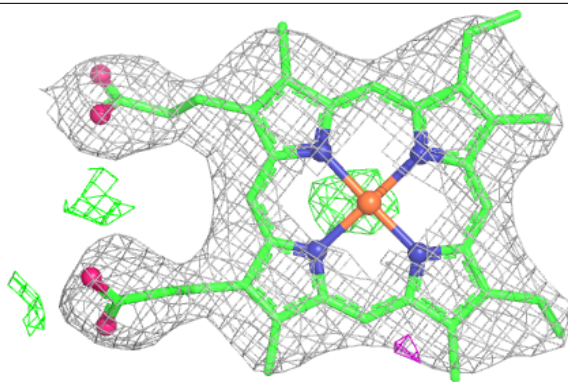
**Electron density around HEM C 501:**

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



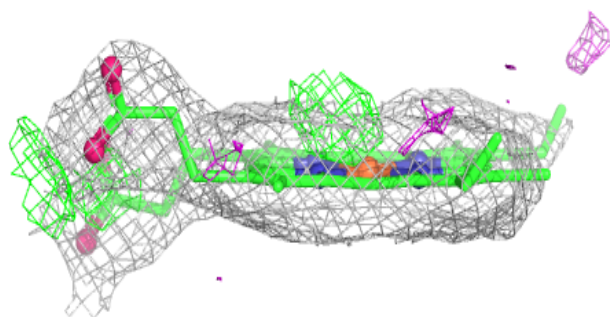
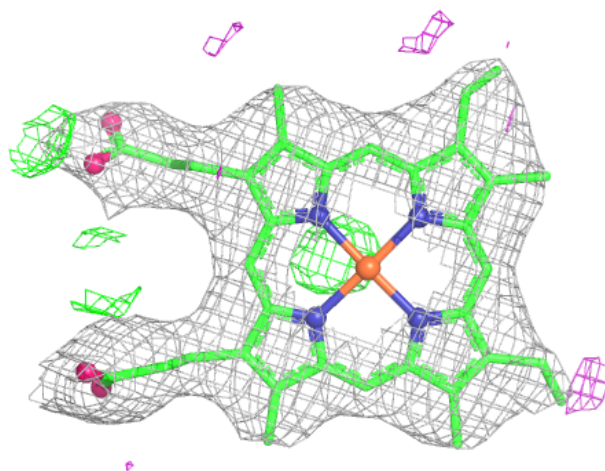
Electron density around HEM A 501:

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



Electron density around HEM B 501:

$2mF_o-DF_c$ (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.