



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 02:36 pm BST

PDB ID : 1IZO
Title : Cytochrome P450 BS beta Complexed with Fatty Acid
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Initiative (RSGI)
Deposited on : 2002-10-10
Resolution : 2.10 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.11
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) : 2.11

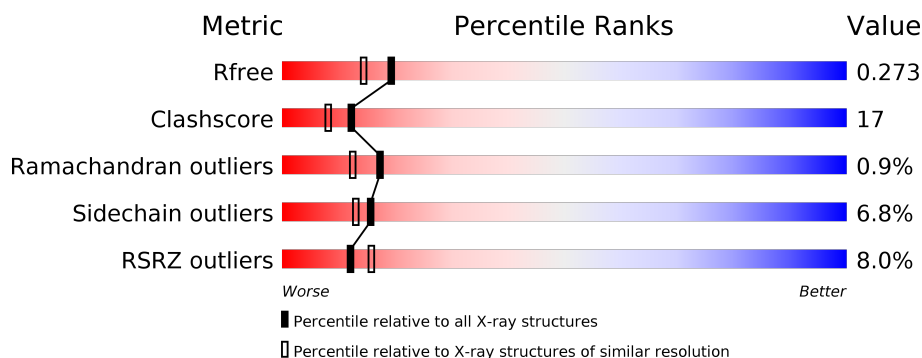
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria 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	417	
1	B	417	
1	C	417	

2 Entry composition [i](#)

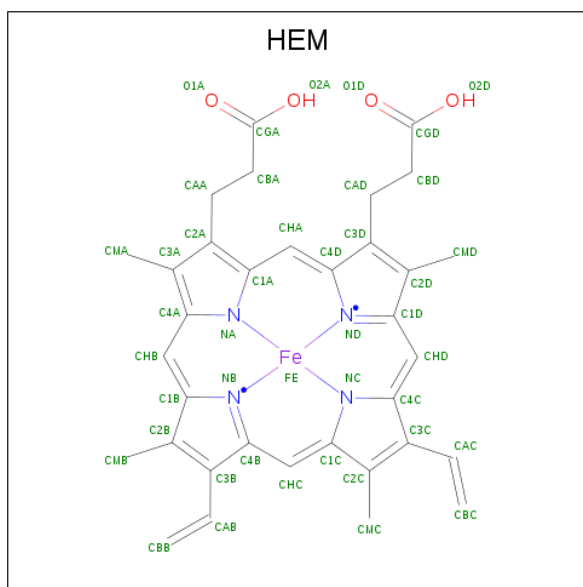
There are 4 unique types of molecules in this entry. The entry contains 10495 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 152A1.

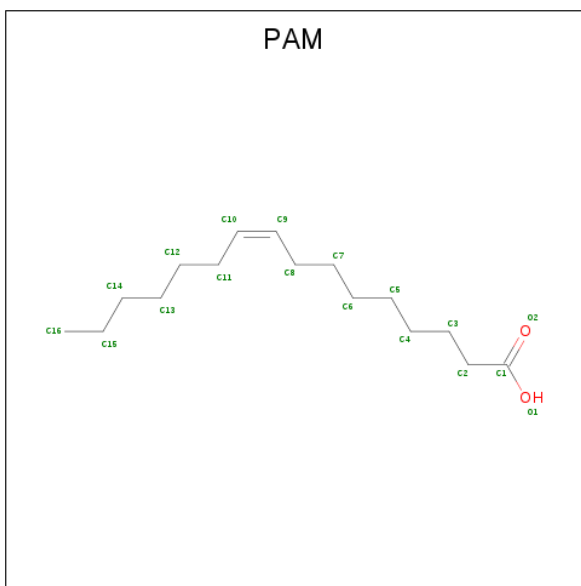
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	411	Total	C	N	O	S	0	0	0
			3339	2125	594	601	19			
1	B	411	Total	C	N	O	S	0	0	0
			3339	2125	594	601	19			
1	C	411	Total	C	N	O	S	0	0	0
			3339	2125	594	601	19			

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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 PALMITOLEIC ACID (three-letter code: PAM) (formula: $C_{16}H_{30}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			18	16	2		
3	B	1	Total	C	O	0	0
			18	16	2		
3	C	1	Total	C	O	0	0
			18	16	2		

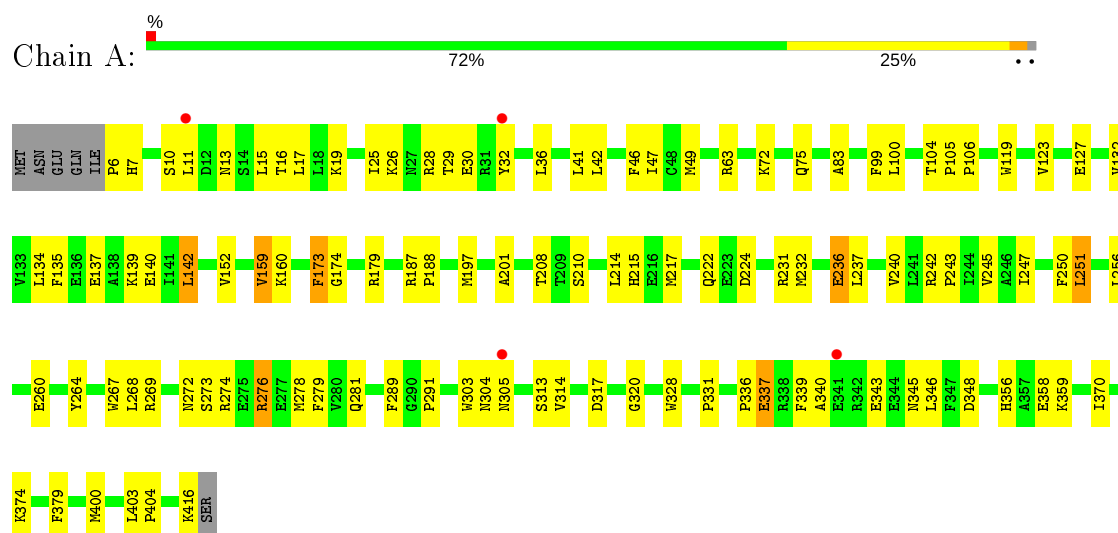
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	153	Total	O	0	0
			153	153		
4	B	84	Total	O	0	0
			84	84		
4	C	58	Total	O	0	0
			58	58		

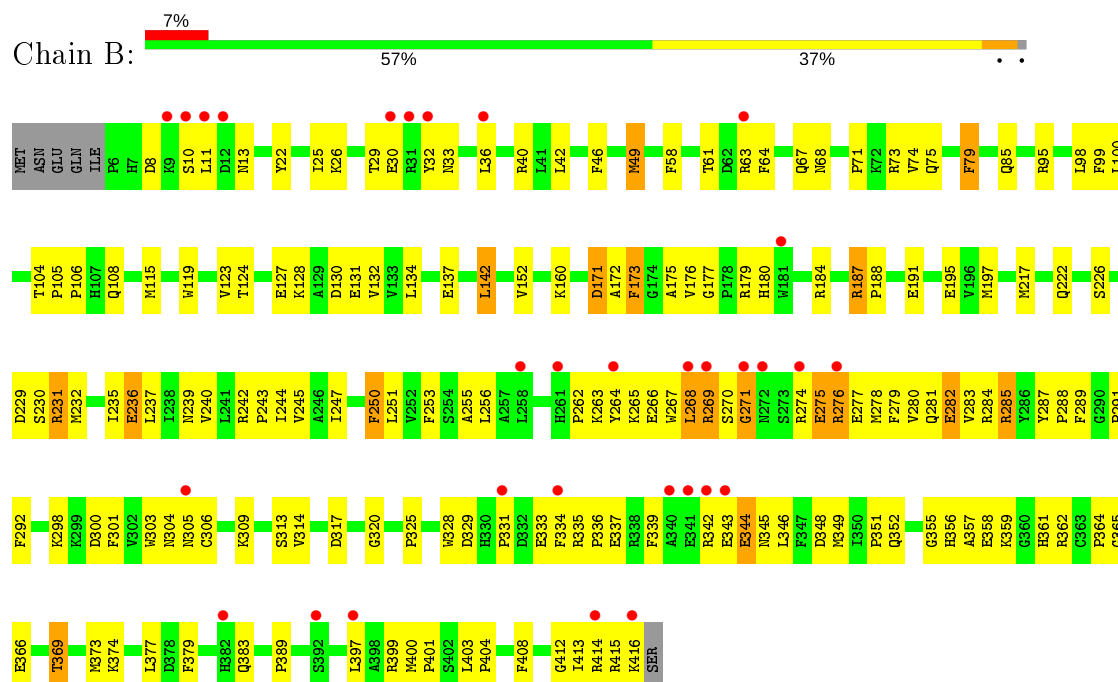
3 Residue-property plots

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

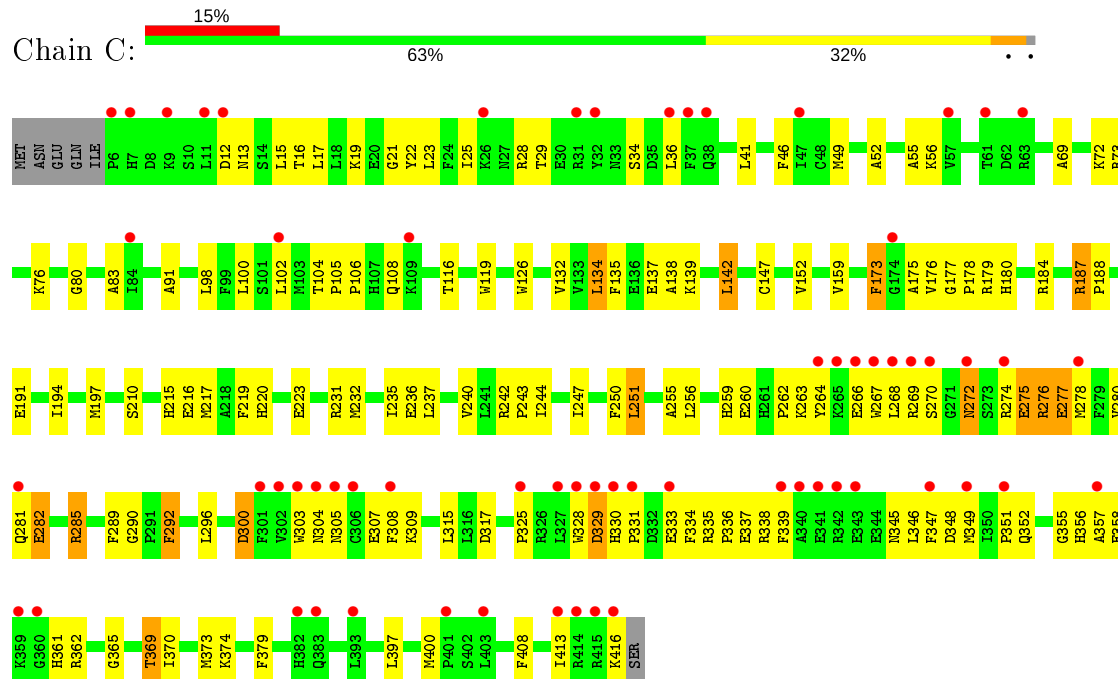
• Molecule 1: Cytochrome P450 152A1



• Molecule 1: Cytochrome P450 152A1



● Molecule 1: Cytochrome P450 152A1



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	155.79Å 155.79Å 111.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.10 46.40 – 2.10	Depositor EDS
% Data completeness (in resolution range)	95.8 (20.00-2.10) 95.7 (46.40-2.10)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.33	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.37 (at 2.10Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.246 , 0.280 0.240 , 0.273	Depositor DCC
R_{free} test set	4432 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	38.7	Xtriage
Anisotropy	0.042	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 55.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.022 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10495	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.63% 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, PAM

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	0.47	0/3422	0.67	0/4614
1	B	0.42	0/3422	0.65	1/4614 (0.0%)
1	C	0.36	0/3422	0.59	2/4614 (0.0%)
All	All	0.42	0/10266	0.64	3/13842 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	271	GLY	N-CA-C	7.78	132.56	113.10
1	C	270	SER	N-CA-C	-6.69	92.95	111.00
1	C	80	GLY	N-CA-C	-5.09	100.36	113.10

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	3339	0	3290	87	0
1	B	3339	0	3290	141	0
1	C	3339	0	3290	117	0
2	A	43	0	30	3	0
2	B	43	0	30	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	43	0	30	5	0
3	A	18	0	29	0	0
3	B	18	0	29	0	0
3	C	18	0	29	2	0
4	A	153	0	0	3	0
4	B	84	0	0	5	0
4	C	58	0	0	4	0
All	All	10495	0	10047	345	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 17.

All (345) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:LEU:HD23	1:A:49:MET:HB3	1.29	1.15
1:A:6:PRO:HB3	1:A:32:TYR:HD2	1.38	0.89
1:C:36:LEU:HD23	1:C:49:MET:HB3	1.54	0.89
1:C:307:GLU:OE2	1:C:309:LYS:HE2	1.78	0.83
1:B:79:PHE:HB3	1:B:85:GLN:NE2	1.93	0.83
1:C:268:LEU:HD12	1:C:276:ARG:HG2	1.57	0.83
1:A:36:LEU:CD2	1:A:49:MET:HB3	2.08	0.83
1:C:328:TRP:O	1:C:331:PRO:HD3	1.84	0.77
1:B:349:MET:HG2	1:B:351:PRO:HD3	1.67	0.77
1:C:83:ALA:HA	1:C:232:MET:HE1	1.66	0.77
1:A:6:PRO:HB3	1:A:32:TYR:CD2	2.23	0.74
1:C:285:ARG:HB2	1:C:349:MET:SD	2.29	0.73
1:C:73:ARG:NH2	1:C:175:ALA:O	2.21	0.73
1:B:265:LYS:HE2	1:B:269:ARG:HH21	1.54	0.73
1:C:187:ARG:HB2	1:C:188:PRO:HD3	1.70	0.73
1:C:36:LEU:CD2	1:C:49:MET:HB3	2.19	0.72
1:C:349:MET:HG2	1:C:351:PRO:HD3	1.70	0.72
1:B:279:PHE:O	1:B:283:VAL:HG23	1.89	0.72
1:B:250:PHE:HB3	1:B:373:MET:CE	2.19	0.72
1:C:282:GLU:HA	1:C:282:GLU:OE1	1.87	0.71
1:B:282:GLU:HG2	1:B:336:PRO:HA	1.74	0.70
1:B:333:GLU:OE1	1:B:335:ARG:NH2	2.22	0.70
1:A:264:TYR:O	1:A:268:LEU:HD23	1.91	0.70
1:A:210:SER:HA	1:A:215:HIS:CD2	2.26	0.70
1:B:236:GLU:OE1	1:B:236:GLU:HA	1.92	0.69
1:A:63:ARG:HH11	1:A:63:ARG:HG3	1.56	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:142:LEU:HG	1:C:244:ILE:O	1.93	0.69
1:B:36:LEU:HD11	1:B:303:TRP:CZ3	2.28	0.69
1:A:41:LEU:HB3	1:A:42:LEU:HD22	1.75	0.68
1:B:231:ARG:O	1:B:235:ILE:HG13	1.93	0.68
1:B:285:ARG:NH2	1:B:331:PRO:O	2.25	0.68
1:B:73:ARG:NH2	1:B:175:ALA:O	2.27	0.66
1:A:140:GLU:OE2	1:A:160:LYS:HG3	1.95	0.66
1:C:231:ARG:O	1:C:235:ILE:HG13	1.96	0.66
1:B:345:ASN:HB3	1:B:348:ASP:OD1	1.95	0.66
1:A:317:ASP:OD2	1:A:320:GLY:HA3	1.96	0.65
1:B:328:TRP:O	1:B:331:PRO:HG3	1.97	0.65
1:C:277:GLU:O	1:C:281:GLN:HG2	1.97	0.65
1:B:280:VAL:HG21	1:B:374:LYS:CG	2.27	0.64
1:C:280:VAL:HG21	1:C:374:LYS:HG2	1.79	0.64
1:B:179:ARG:NH2	1:B:400:MET:O	2.31	0.63
1:A:42:LEU:HD22	1:A:42:LEU:N	2.14	0.63
1:A:268:LEU:HD12	1:A:276:ARG:HG2	1.80	0.63
1:A:179:ARG:NH2	1:A:400:MET:O	2.31	0.63
1:A:273:SER:HA	1:A:276:ARG:NH1	2.12	0.63
1:B:64:PHE:CE1	1:B:314:VAL:HG21	2.34	0.63
1:C:282:GLU:HB3	1:C:334:PHE:CE1	2.33	0.62
1:A:99:PHE:CZ	1:A:236:GLU:HG3	2.34	0.62
1:C:187:ARG:CB	1:C:188:PRO:HD3	2.28	0.62
1:B:255:ALA:HB1	1:B:408:PHE:CE2	2.34	0.62
1:C:36:LEU:HD11	1:C:303:TRP:CZ3	2.35	0.62
1:B:280:VAL:HG21	1:B:374:LYS:HG3	1.79	0.61
1:C:119:TRP:HB3	1:C:379:PHE:CD1	2.35	0.61
1:A:75:GLN:HG3	4:A:707:HOH:O	1.98	0.61
1:B:265:LYS:HE2	1:B:269:ARG:NH2	2.15	0.61
1:C:23:LEU:HD11	1:C:397:LEU:HD13	1.81	0.61
1:B:399:ARG:HG2	1:B:401:PRO:O	2.01	0.61
1:C:17:LEU:C	1:C:17:LEU:HD23	2.21	0.60
1:C:236:GLU:HA	1:C:236:GLU:OE1	2.00	0.60
1:B:369:THR:O	1:B:373:MET:HG3	2.01	0.60
1:C:242:ARG:HB3	1:C:243:PRO:HD3	1.81	0.60
1:B:285:ARG:HB2	1:B:349:MET:SD	2.41	0.60
1:C:91:ALA:HA	1:C:223:GLU:OE2	2.00	0.60
1:C:370:ILE:O	1:C:374:LYS:HG3	2.01	0.60
1:A:6:PRO:CB	1:A:32:TYR:HD2	2.11	0.60
1:B:268:LEU:CD1	1:B:275:GLU:HB2	2.33	0.59
1:C:98:LEU:HD22	1:C:220:HIS:CG	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:187:ARG:CB	1:B:188:PRO:HD3	2.33	0.59
1:B:415:ARG:NH1	4:B:1628:HOH:O	2.35	0.58
1:B:250:PHE:HB3	1:B:373:MET:HE3	1.85	0.58
1:A:6:PRO:CB	1:A:32:TYR:CD2	2.85	0.58
1:C:256:LEU:O	1:C:260:GLU:HG3	2.03	0.58
1:A:152:VAL:HG13	1:A:197:MET:SD	2.43	0.58
1:C:282:GLU:HG2	1:C:336:PRO:HA	1.85	0.58
1:A:100:LEU:HD13	1:A:358:GLU:HG2	1.86	0.58
1:B:352:GLN:HE22	1:B:369:THR:HG21	1.69	0.58
1:C:21:GLY:HA2	1:C:400:MET:CE	2.34	0.57
1:B:61:THR:O	1:B:298:LYS:HE2	2.04	0.57
1:A:370:ILE:HG22	1:A:374:LYS:HE2	1.87	0.57
1:A:274:ARG:HG2	1:A:278:MET:CE	2.34	0.57
1:C:264:TYR:O	1:C:268:LEU:HD23	2.05	0.57
1:C:365:GLY:O	1:C:369:THR:HG22	2.03	0.57
1:A:328:TRP:O	1:A:331:PRO:HD3	2.05	0.57
1:C:345:ASN:HB3	1:C:348:ASP:OD1	2.05	0.57
1:B:100:LEU:HD13	1:B:358:GLU:HG2	1.87	0.57
1:B:282:GLU:HA	1:B:282:GLU:OE1	2.04	0.57
1:B:232:MET:HE2	1:B:236:GLU:OE2	2.05	0.56
1:C:300:ASP:OD2	1:C:309:LYS:HD3	2.05	0.56
1:C:361:HIS:HE1	4:C:2636:HOH:O	1.87	0.56
1:A:123:VAL:O	1:A:127:GLU:HG3	2.05	0.56
1:A:41:LEU:HB3	1:A:42:LEU:CD2	2.36	0.55
1:B:180:HIS:NE2	1:B:184:ARG:NH1	2.54	0.55
1:A:160:LYS:HD3	4:A:634:HOH:O	2.07	0.55
1:A:15:LEU:O	1:A:19:LYS:HG3	2.07	0.55
1:B:240:VAL:HA	2:B:501:HEM:HBC2	1.88	0.55
1:B:282:GLU:CG	1:B:336:PRO:HA	2.36	0.55
1:C:100:LEU:HD13	1:C:358:GLU:HG2	1.88	0.55
1:A:214:LEU:C	1:A:214:LEU:HD13	2.27	0.55
1:B:187:ARG:HB2	1:B:188:PRO:HD3	1.89	0.55
1:B:63:ARG:HG3	1:B:63:ARG:HH11	1.72	0.55
1:C:72:LYS:HD3	4:C:2615:HOH:O	2.06	0.54
1:B:344:GLU:OE2	1:B:346:LEU:HD21	2.08	0.54
1:C:256:LEU:O	1:C:256:LEU:HD13	2.08	0.54
1:B:247:ILE:HA	2:B:501:HEM:HBB1	1.89	0.54
1:B:175:ALA:HB2	1:B:179:ARG:CZ	2.38	0.54
1:B:99:PHE:CZ	1:B:236:GLU:HG3	2.43	0.54
1:A:83:ALA:HA	1:A:232:MET:HE2	1.90	0.53
1:B:242:ARG:HB3	1:B:243:PRO:HD3	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:361:HIS:HD2	2:C:501:HEM:O2D	1.91	0.53
1:A:242:ARG:HB3	1:A:243:PRO:HD3	1.90	0.53
1:B:195:GLU:HG3	1:B:230:SER:HB3	1.90	0.53
1:C:134:LEU:HD12	1:C:408:PHE:CD2	2.43	0.53
1:C:255:ALA:HB1	1:C:408:PHE:CE2	2.43	0.53
1:A:187:ARG:HB3	1:A:188:PRO:HD3	1.91	0.53
1:C:281:GLN:HB3	1:C:339:PHE:CE2	2.44	0.53
1:C:116:THR:HG23	4:C:2641:HOH:O	2.09	0.53
1:B:100:LEU:CD1	1:B:358:GLU:HG2	2.39	0.52
1:C:126:TRP:CE3	1:C:413:ILE:HD13	2.43	0.52
1:A:281:GLN:HA	1:A:281:GLN:OE1	2.08	0.52
1:B:268:LEU:C	1:B:270:SER:H	2.11	0.52
1:B:26:LYS:O	1:B:30:GLU:HG3	2.09	0.52
1:C:104:THR:OG1	1:C:106:PRO:HD2	2.08	0.52
1:B:152:VAL:HG13	1:B:197:MET:SD	2.49	0.52
1:B:264:TYR:CE1	1:B:336:PRO:HG2	2.44	0.52
1:A:26:LYS:O	1:A:30:GLU:HG3	2.10	0.52
1:C:369:THR:O	1:C:373:MET:HG3	2.10	0.52
1:A:247:ILE:HA	2:A:501:HEM:HBB1	1.92	0.52
1:A:345:ASN:HB3	1:A:348:ASP:OD1	2.09	0.52
1:B:104:THR:OG1	1:B:106:PRO:HD2	2.10	0.51
1:A:16:THR:HG21	1:A:28:ARG:NH2	2.24	0.51
1:B:357:ALA:HA	1:B:362:ARG:HB3	1.93	0.51
1:A:17:LEU:C	1:A:17:LEU:HD23	2.31	0.51
1:A:240:VAL:O	1:A:243:PRO:HD2	2.11	0.51
1:A:63:ARG:NH1	1:A:63:ARG:HG3	2.24	0.51
1:B:268:LEU:HD13	1:B:275:GLU:HB2	1.92	0.50
1:C:132:VAL:CG1	1:C:137:GLU:HG3	2.41	0.50
1:C:275:GLU:HA	1:C:278:MET:HE2	1.92	0.50
1:A:269:ARG:CZ	1:A:416:LYS:HZ1	2.24	0.50
1:B:46:PHE:HA	1:B:313:SER:O	2.12	0.50
1:C:346:LEU:HD11	1:C:356:HIS:CE1	2.46	0.50
1:A:135:PHE:CE2	1:A:139:LYS:HD2	2.47	0.50
1:C:280:VAL:HG21	1:C:374:LYS:CG	2.42	0.50
1:B:46:PHE:CE1	1:B:313:SER:HB3	2.46	0.50
1:A:274:ARG:HG2	1:A:278:MET:HE2	1.93	0.49
1:B:266:GLU:HG2	4:B:1671:HOH:O	2.12	0.49
1:C:52:ALA:O	1:C:55:ALA:HB3	2.12	0.49
1:A:46:PHE:CE1	1:A:313:SER:HB3	2.47	0.49
1:C:281:GLN:HA	1:C:281:GLN:OE1	2.12	0.49
1:A:343:GLU:HA	1:A:343:GLU:OE1	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:239:ASN:O	1:B:243:PRO:HG2	2.12	0.49
1:A:119:TRP:HB3	1:A:379:PHE:CD1	2.47	0.49
1:B:71:PRO:HB2	1:B:73:ARG:HG2	1.93	0.49
1:B:63:ARG:HD3	1:B:301:PHE:CE2	2.47	0.49
1:B:300:ASP:OD2	1:B:309:LYS:HD3	2.12	0.49
1:B:373:MET:O	4:B:1613:HOH:O	2.19	0.49
1:B:280:VAL:HG21	1:B:374:LYS:HG2	1.94	0.49
1:C:300:ASP:HA	1:C:308:PHE:O	2.13	0.49
1:A:28:ARG:HG3	1:A:28:ARG:HH11	1.77	0.48
1:B:415:ARG:HG2	1:B:416:LYS:H	1.78	0.48
1:A:269:ARG:NH2	1:A:416:LYS:NZ	2.61	0.48
1:B:240:VAL:HG13	2:B:501:HEM:HBC1	1.95	0.48
1:B:115:MET:O	1:B:119:TRP:CD1	2.66	0.48
1:B:355:GLY:O	1:B:362:ARG:HD2	2.14	0.48
1:C:132:VAL:HG13	1:C:137:GLU:HG3	1.96	0.48
1:C:361:HIS:O	2:C:501:HEM:HBA2	2.13	0.48
1:A:28:ARG:HG3	1:A:28:ARG:NH1	2.29	0.48
1:C:180:HIS:NE2	1:C:184:ARG:NH1	2.60	0.48
1:B:123:VAL:O	1:B:127:GLU:HG3	2.14	0.48
1:C:25:ILE:HG21	1:C:317:ASP:HB2	1.95	0.48
1:B:264:TYR:O	1:B:267:TRP:HB3	2.14	0.48
1:C:138:ALA:O	1:C:142:LEU:HB2	2.14	0.48
1:C:255:ALA:HB1	1:C:408:PHE:CZ	2.49	0.48
1:B:124:THR:HG22	1:B:128:LYS:NZ	2.29	0.47
1:A:173:PHE:C	1:A:173:PHE:CD1	2.88	0.47
1:A:264:TYR:OH	1:A:337:GLU:OE2	2.29	0.47
1:C:275:GLU:HA	1:C:278:MET:CE	2.45	0.47
1:A:72:LYS:CE	1:A:75:GLN:HE22	2.26	0.47
1:B:276:ARG:HH11	1:B:276:ARG:HB2	1.79	0.47
1:B:412:GLY:O	1:B:414:ARG:HG3	2.14	0.47
1:A:370:ILE:O	1:A:374:LYS:HG3	2.14	0.47
1:A:236:GLU:HA	1:A:236:GLU:OE1	2.13	0.47
1:B:187:ARG:O	1:B:191:GLU:HG3	2.14	0.47
1:B:173:PHE:CD1	1:B:173:PHE:C	2.88	0.47
1:B:58:PHE:HE1	1:B:314:VAL:HG23	1.79	0.47
1:C:105:PRO:N	1:C:106:PRO:CD	2.78	0.47
1:B:334:PHE:CE2	1:B:336:PRO:HD3	2.50	0.47
1:B:356:HIS:HB2	1:B:359:LYS:HB2	1.96	0.47
1:C:325:PRO:HA	1:C:328:TRP:O	2.15	0.47
1:B:282:GLU:OE1	1:B:339:PHE:HE1	1.98	0.47
1:A:140:GLU:HG2	1:A:159:VAL:HG12	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:361:HIS:HA	2:B:501:HEM:O1D	2.16	0.46
1:C:173:PHE:C	1:C:173:PHE:CD1	2.88	0.46
1:A:356:HIS:CD2	1:A:359:LYS:HD2	2.50	0.46
1:C:267:TRP:HE3	1:C:268:LEU:HD22	1.81	0.46
1:B:365:GLY:O	1:B:369:THR:CG2	2.63	0.46
1:A:346:LEU:HD11	1:A:356:HIS:CE1	2.50	0.46
1:B:342:ARG:HG3	1:B:343:GLU:H	1.81	0.46
1:C:266:GLU:HA	1:C:269:ARG:HB3	1.97	0.46
1:A:276:ARG:HB2	1:A:276:ARG:HH11	1.80	0.46
1:B:250:PHE:HB3	1:B:373:MET:HE1	1.93	0.46
1:C:104:THR:C	1:C:108:GLN:HE21	2.18	0.46
1:C:330:HIS:N	1:C:331:PRO:CD	2.79	0.46
1:C:357:ALA:HA	1:C:362:ARG:HB3	1.98	0.46
1:C:76:LYS:O	1:C:187:ARG:HG3	2.16	0.46
1:C:356:HIS:HA	4:C:2647:HOH:O	2.15	0.46
1:A:132:VAL:HG13	1:A:137:GLU:HG3	1.98	0.46
1:B:171:ASP:OD2	1:B:399:ARG:NH2	2.49	0.46
1:B:317:ASP:OD2	1:B:320:GLY:HA3	2.15	0.46
1:B:130:ASP:OD1	1:B:131:GLU:HG3	2.17	0.45
1:C:194:ILE:O	1:C:197:MET:HB2	2.15	0.45
1:B:267:TRP:CZ2	1:B:275:GLU:HA	2.51	0.45
1:C:176:VAL:HG12	1:C:177:GLY:N	2.31	0.45
1:C:338:ARG:HG3	1:C:339:PHE:CD1	2.51	0.45
1:B:75:GLN:HG3	4:B:1650:HOH:O	2.16	0.45
1:C:259:HIS:O	1:C:262:PRO:HD3	2.17	0.45
1:C:272:ASN:HA	1:C:272:ASN:HD22	1.64	0.45
1:B:253:PHE:CE1	1:B:404:PRO:HD3	2.51	0.45
1:B:379:PHE:HA	1:B:383:GLN:HG2	1.99	0.45
1:C:352:GLN:HE22	1:C:369:THR:HG21	1.82	0.45
1:A:28:ARG:O	1:A:32:TYR:HD1	2.00	0.45
1:A:243:PRO:HB2	2:A:501:HEM:C3C	2.51	0.45
1:C:285:ARG:NH2	1:C:331:PRO:O	2.50	0.45
1:B:142:LEU:HB3	1:B:245:VAL:HA	1.99	0.45
1:B:63:ARG:HD3	1:B:301:PHE:CD2	2.52	0.45
1:B:46:PHE:CZ	1:B:313:SER:HB3	2.51	0.45
1:C:16:THR:HG21	1:C:28:ARG:NH2	2.32	0.45
1:C:292:PHE:HE1	3:C:2601:PAM:H132	1.82	0.45
1:C:264:TYR:CE2	1:C:336:PRO:HD2	2.52	0.44
1:B:342:ARG:HG3	1:B:343:GLU:N	2.32	0.44
1:C:247:ILE:HG22	1:C:251:LEU:HD13	2.00	0.44
1:A:264:TYR:O	1:A:268:LEU:CD2	2.64	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:PHE:HA	1:A:336:PRO:HB3	1.98	0.44
1:B:32:TYR:O	1:B:33:ASN:C	2.55	0.44
1:C:274:ARG:HG2	1:C:278:MET:CE	2.47	0.44
1:A:201:ALA:HB1	1:A:208:THR:HG21	1.99	0.44
1:B:276:ARG:CD	1:B:377:LEU:HD12	2.48	0.44
1:B:267:TRP:CE2	1:B:275:GLU:HB3	2.53	0.44
1:B:274:ARG:HB3	1:B:278:MET:HE2	2.00	0.44
1:B:279:PHE:HA	1:B:336:PRO:HB3	1.98	0.44
1:B:36:LEU:CD2	1:B:49:MET:HB3	2.48	0.44
1:C:335:ARG:O	1:C:338:ARG:HG2	2.17	0.44
1:A:256:LEU:HD13	1:A:260:GLU:HG3	2.00	0.44
1:C:152:VAL:HG13	1:C:197:MET:SD	2.58	0.44
1:C:269:ARG:CZ	1:C:416:LYS:HZ1	2.31	0.44
1:A:173:PHE:CD1	1:A:174:GLY:N	2.86	0.43
1:A:36:LEU:HD11	1:A:303:TRP:CZ3	2.53	0.43
1:B:36:LEU:HD23	1:B:49:MET:HB3	1.98	0.43
1:C:187:ARG:CB	1:C:188:PRO:CD	2.95	0.43
1:C:355:GLY:O	1:C:362:ARG:HD2	2.19	0.43
1:B:160:LYS:HB3	4:B:1630:HOH:O	2.18	0.43
1:B:364:PRO:HD2	2:B:501:HEM:C2D	2.53	0.43
1:C:256:LEU:C	1:C:256:LEU:HD13	2.37	0.43
1:C:46:PHE:CE1	1:C:315:LEU:HG	2.53	0.43
1:A:47:ILE:O	1:A:314:VAL:HA	2.18	0.43
1:A:281:GLN:HB3	1:A:339:PHE:CZ	2.53	0.43
1:B:172:ALA:HB1	1:B:180:HIS:HA	2.01	0.43
1:B:13:ASN:HB2	1:B:40:ARG:O	2.19	0.43
1:C:22:TYR:CE1	1:C:400:MET:HG2	2.52	0.43
1:B:22:TYR:CE1	1:B:400:MET:HG2	2.52	0.43
1:C:328:TRP:HZ2	1:C:349:MET:HB2	1.84	0.43
1:A:240:VAL:C	1:A:243:PRO:HD2	2.39	0.43
1:A:251:LEU:HD12	1:A:251:LEU:HA	1.83	0.43
1:B:105:PRO:HB2	1:B:106:PRO:HD3	1.99	0.43
1:C:333:GLU:OE1	1:C:335:ARG:NH2	2.52	0.43
1:B:328:TRP:HZ2	1:B:349:MET:HB2	1.83	0.43
1:C:135:PHE:CE2	1:C:139:LYS:HD2	2.54	0.43
1:C:240:VAL:O	1:C:243:PRO:HD2	2.18	0.43
1:A:269:ARG:NH2	1:A:416:LYS:HZ2	2.17	0.43
1:C:104:THR:HB	1:C:105:PRO:CD	2.49	0.43
1:C:361:HIS:HA	2:C:501:HEM:O1D	2.19	0.43
1:B:268:LEU:O	1:B:270:SER:N	2.48	0.42
1:C:102:LEU:HD21	1:C:216:GLU:HG3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:210:SER:HA	1:C:215:HIS:CD2	2.54	0.42
1:B:264:TYR:HD1	1:B:267:TRP:CE3	2.38	0.42
1:B:365:GLY:O	1:B:369:THR:HG22	2.19	0.42
1:B:95:ARG:O	1:B:98:LEU:HB3	2.19	0.42
1:C:329:ASP:O	1:C:330:HIS:C	2.58	0.42
1:A:222:GLN:HB2	1:A:224:ASP:OD1	2.19	0.42
1:A:243:PRO:HB2	2:A:501:HEM:C2C	2.55	0.42
1:B:276:ARG:HH11	1:B:276:ARG:CB	2.32	0.42
1:B:67:GLN:HG2	1:B:68:ASN:ND2	2.34	0.42
1:C:21:GLY:HA2	1:C:400:MET:HE1	2.01	0.42
1:B:176:VAL:HG12	1:B:177:GLY:N	2.34	0.42
1:B:240:VAL:HA	2:B:501:HEM:CBC	2.49	0.42
1:B:277:GLU:O	1:B:281:GLN:HG2	2.20	0.42
1:B:325:PRO:HA	1:B:331:PRO:HG3	2.01	0.42
1:B:25:ILE:HG21	1:B:317:ASP:HB2	2.02	0.42
1:B:403:LEU:HD12	1:B:404:PRO:HD2	2.01	0.42
1:C:15:LEU:O	1:C:19:LYS:HG3	2.20	0.42
1:B:282:GLU:HB3	1:B:334:PHE:CE1	2.54	0.42
1:C:264:TYR:HA	1:C:267:TRP:CB	2.49	0.42
1:C:243:PRO:HB2	2:C:501:HEM:C3C	2.54	0.42
1:C:69:ALA:HA	1:C:296:LEU:HG	2.02	0.42
1:C:175:ALA:HB2	1:C:179:ARG:CZ	2.50	0.42
1:A:105:PRO:N	1:A:106:PRO:CD	2.83	0.41
1:A:142:LEU:HB3	1:A:245:VAL:HA	2.02	0.41
1:B:284:ARG:HH12	1:B:366:GLU:HG3	1.85	0.41
1:B:276:ARG:HD3	1:B:377:LEU:HD12	2.01	0.41
1:B:413:ILE:HG13	1:B:413:ILE:H	1.68	0.41
1:C:13:ASN:HB3	1:C:41:LEU:CD2	2.50	0.41
1:C:29:THR:HB	1:C:34:SER:O	2.21	0.41
1:A:36:LEU:HD21	1:A:303:TRP:HZ3	1.85	0.41
1:A:42:LEU:CD2	1:A:42:LEU:N	2.83	0.41
1:B:240:VAL:O	1:B:244:ILE:HG13	2.21	0.41
1:C:191:GLU:OE1	1:C:231:ARG:NH1	2.52	0.41
1:A:132:VAL:CG1	1:A:137:GLU:HG3	2.49	0.41
1:A:99:PHE:CE2	1:A:236:GLU:HG3	2.55	0.41
1:A:264:TYR:HA	1:A:267:TRP:HB3	2.02	0.41
1:B:132:VAL:CG1	1:B:137:GLU:HG3	2.50	0.41
1:C:147:CYS:HB3	1:C:152:VAL:HB	2.02	0.41
1:A:25:ILE:O	1:A:29:THR:HG23	2.20	0.41
1:A:46:PHE:HA	1:A:313:SER:O	2.19	0.41
1:B:255:ALA:HB1	1:B:408:PHE:CZ	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:280:VAL:O	1:B:283:VAL:N	2.54	0.41
1:A:374:LYS:NZ	4:A:729:HOH:O	2.46	0.41
1:B:119:TRP:HB3	1:B:379:PHE:CD1	2.56	0.41
1:B:222:GLN:HG3	1:B:226:SER:O	2.20	0.41
1:B:25:ILE:O	1:B:29:THR:HG23	2.20	0.41
1:B:79:PHE:CB	1:B:85:GLN:NE2	2.75	0.41
1:B:287:TYR:HA	1:B:288:PRO:HD3	1.80	0.41
1:A:403:LEU:O	1:A:404:PRO:C	2.57	0.41
1:B:267:TRP:CH2	1:B:275:GLU:HA	2.55	0.41
1:B:284:ARG:NH1	1:B:366:GLU:HG3	2.36	0.41
1:B:8:ASP:C	1:B:10:SER:H	2.24	0.41
1:C:247:ILE:HA	2:C:501:HEM:HBB1	2.03	0.41
1:A:256:LEU:HD11	1:A:260:GLU:OE1	2.21	0.41
1:B:74:VAL:HG12	1:B:79:PHE:CE1	2.56	0.41
1:C:36:LEU:HD11	1:C:303:TRP:CE3	2.55	0.41
1:B:276:ARG:O	1:B:280:VAL:HG23	2.21	0.40
1:B:397:LEU:N	1:B:397:LEU:HD22	2.36	0.40
1:C:215:HIS:O	1:C:219:PHE:HD2	2.04	0.40
1:B:187:ARG:CB	1:B:188:PRO:CD	2.99	0.40
1:B:303:TRP:O	1:B:306:CYS:HB3	2.21	0.40
1:C:290:GLY:HA2	3:C:2601:PAM:H62	2.03	0.40
1:C:56:LYS:HG2	1:C:347:PHE:CD2	2.55	0.40
1:B:104:THR:C	1:B:108:GLN:HE21	2.24	0.40
1:B:364:PRO:HD2	2:B:501:HEM:C1D	2.56	0.40
1:C:134:LEU:HB2	1:C:408:PHE:HB3	2.03	0.40
1:C:177:GLY:HA3	1:C:178:PRO:HD2	1.97	0.40
1:A:104:THR:CB	1:A:106:PRO:HD2	2.52	0.40
1:A:267:TRP:HE3	1:A:268:LEU:HD22	1.86	0.40
1:B:240:VAL:C	1:B:243:PRO:HD2	2.42	0.40
1:B:282:GLU:OE1	1:B:285:ARG:HD3	2.21	0.40
1:C:268:LEU:CD1	1:C:276:ARG:HG2	2.41	0.40
1:C:264:TYR:HA	1:C:267:TRP:HB2	2.01	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	409/417 (98%)	393 (96%)	13 (3%)	3 (1%)	22	18
1	B	409/417 (98%)	376 (92%)	26 (6%)	7 (2%)	9	4
1	C	409/417 (98%)	377 (92%)	31 (8%)	1 (0%)	47	49
All	All	1227/1251 (98%)	1146 (93%)	70 (6%)	11 (1%)	17	12

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	263	LYS
1	B	269	ARG
1	A	340	ALA
1	B	291	PRO
1	B	344	GLU
1	A	291	PRO
1	B	262	PRO
1	B	271	GLY
1	A	159	VAL
1	C	159	VAL
1	B	389	PRO

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	352/358 (98%)	333 (95%)	19 (5%)	22	20
1	B	352/358 (98%)	323 (92%)	29 (8%)	11	8
1	C	352/358 (98%)	328 (93%)	24 (7%)	16	13
All	All	1056/1074 (98%)	984 (93%)	72 (7%)	16	13

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	HIS
1	A	10	SER
1	A	11	LEU
1	A	13	ASN
1	A	134	LEU
1	A	142	LEU
1	A	173	PHE
1	A	217	MET
1	A	231	ARG
1	A	236	GLU
1	A	237	LEU
1	A	250	PHE
1	A	251	LEU
1	A	272	ASN
1	A	276	ARG
1	A	289	PHE
1	A	304	ASN
1	A	305	ASN
1	A	337	GLU
1	B	11	LEU
1	B	42	LEU
1	B	49	MET
1	B	79	PHE
1	B	134	LEU
1	B	142	LEU
1	B	171	ASP
1	B	173	PHE
1	B	187	ARG
1	B	217	MET
1	B	229	ASP
1	B	231	ARG
1	B	236	GLU
1	B	237	LEU
1	B	250	PHE
1	B	251	LEU
1	B	256	LEU
1	B	268	LEU
1	B	275	GLU
1	B	276	ARG
1	B	282	GLU
1	B	285	ARG
1	B	289	PHE
1	B	292	PHE

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Mol	Chain	Res	Type
1	B	304	ASN
1	B	305	ASN
1	B	329	ASP
1	B	337	GLU
1	B	369	THR
1	C	12	ASP
1	C	134	LEU
1	C	142	LEU
1	C	173	PHE
1	C	187	ARG
1	C	217	MET
1	C	237	LEU
1	C	250	PHE
1	C	251	LEU
1	C	263	LYS
1	C	272	ASN
1	C	275	GLU
1	C	276	ARG
1	C	277	GLU
1	C	282	GLU
1	C	285	ARG
1	C	289	PHE
1	C	292	PHE
1	C	300	ASP
1	C	304	ASN
1	C	305	ASN
1	C	329	ASP
1	C	337	GLU
1	C	369	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	ASN
1	A	227	GLN
1	A	272	ASN
1	A	304	ASN
1	A	305	ASN
1	A	356	HIS
1	A	383	GLN
1	B	13	ASN
1	B	68	ASN

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Mol	Chain	Res	Type
1	B	92	HIS
1	B	108	GLN
1	B	304	ASN
1	B	305	ASN
1	B	356	HIS
1	B	361	HIS
1	C	13	ASN
1	C	68	ASN
1	C	82	ASN
1	C	108	GLN
1	C	272	ASN
1	C	304	ASN
1	C	305	ASN
1	C	356	HIS
1	C	361	HIS
1	C	383	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 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
3	PAM	C	2601	-	14,17,17	1.79	4 (28%)	13,17,17	0.98	1 (7%)
2	HEM	B	501	1,4	27,50,50	1.95	7 (25%)	17,82,82	1.37	2 (11%)
2	HEM	C	501	1	27,50,50	1.91	7 (25%)	17,82,82	1.25	3 (17%)
3	PAM	B	1601	-	14,17,17	1.82	4 (28%)	13,17,17	1.04	1 (7%)
3	PAM	A	601	-	14,17,17	1.76	4 (28%)	13,17,17	0.96	1 (7%)
2	HEM	A	501	1,4	27,50,50	1.91	7 (25%)	17,82,82	1.75	4 (23%)

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	PAM	C	2601	-	-	5/13/15/15	-
2	HEM	B	501	1,4	-	0/6/54/54	-
2	HEM	C	501	1	-	0/6/54/54	-
3	PAM	B	1601	-	-	4/13/15/15	-
3	PAM	A	601	-	-	6/13/15/15	-
2	HEM	A	501	1,4	-	0/6/54/54	-

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	HEM	C1D-CHD	4.83	1.54	1.41
2	C	501	HEM	C1D-CHD	4.80	1.54	1.41
2	A	501	HEM	C1D-CHD	4.72	1.54	1.41
2	B	501	HEM	C4A-CHB	4.28	1.52	1.41
2	C	501	HEM	C4A-CHB	4.11	1.52	1.41
2	A	501	HEM	C4B-NB	3.90	1.44	1.36
3	A	601	PAM	C10-C9	3.73	1.53	1.31
2	A	501	HEM	C4A-CHB	3.59	1.51	1.41
3	B	1601	PAM	C10-C9	3.50	1.52	1.31
3	C	2601	PAM	C7-C8	-3.48	1.38	1.52
2	B	501	HEM	C4B-NB	3.39	1.43	1.36
3	B	1601	PAM	C7-C8	-3.37	1.38	1.52
3	C	2601	PAM	C10-C9	3.35	1.51	1.31
2	A	501	HEM	C3C-CAC	-3.28	1.41	1.47
3	A	601	PAM	C12-C11	-3.23	1.39	1.52
3	B	1601	PAM	C12-C11	-3.21	1.39	1.52
2	C	501	HEM	C4B-NB	3.16	1.42	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	2601	PAM	C12-C11	-3.06	1.40	1.52
2	B	501	HEM	C3C-CAC	-2.91	1.41	1.47
2	B	501	HEM	C3C-C2C	-2.85	1.36	1.40
3	A	601	PAM	C7-C8	-2.82	1.41	1.52
2	B	501	HEM	C1D-ND	2.55	1.41	1.36
2	C	501	HEM	C3C-CAC	-2.49	1.42	1.47
2	C	501	HEM	C1D-ND	2.42	1.41	1.36
2	A	501	HEM	C1D-ND	2.28	1.40	1.36
2	C	501	HEM	C3B-C2B	-2.19	1.37	1.40
3	C	2601	PAM	C5-C4	-2.14	1.39	1.51
3	B	1601	PAM	C5-C4	-2.11	1.39	1.51
2	A	501	HEM	CAD-C3D	2.11	1.55	1.52
2	A	501	HEM	CBB-CAB	2.11	1.43	1.29
2	B	501	HEM	CMB-C2B	2.10	1.56	1.51
2	C	501	HEM	C3B-CAB	-2.06	1.43	1.47
3	A	601	PAM	C5-C4	-2.04	1.40	1.51

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	C4C-C3C-C2C	-3.83	104.23	106.90
2	A	501	HEM	C1D-C2D-C3D	3.15	109.19	107.00
2	B	501	HEM	C4C-C3C-C2C	-2.93	104.85	106.90
2	B	501	HEM	C1D-C2D-C3D	2.38	108.65	107.00
3	B	1601	PAM	C7-C8-C9	2.37	126.00	112.43
2	C	501	HEM	C4C-C3C-C2C	-2.31	105.28	106.90
2	A	501	HEM	CMC-C2C-C3C	2.12	128.65	124.68
2	C	501	HEM	C1D-C2D-C3D	2.09	108.45	107.00
2	A	501	HEM	CMD-C2D-C1D	-2.07	125.28	128.46
3	A	601	PAM	C7-C8-C9	2.04	124.12	112.43
2	C	501	HEM	CBD-CAD-C3D	2.03	116.22	112.48
3	C	2601	PAM	C7-C8-C9	2.01	123.97	112.43

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	2601	PAM	C1-C2-C3-C4
3	C	2601	PAM	C11-C10-C9-C8
3	B	1601	PAM	C1-C2-C3-C4
3	A	601	PAM	C1-C2-C3-C4
3	B	1601	PAM	C11-C10-C9-C8

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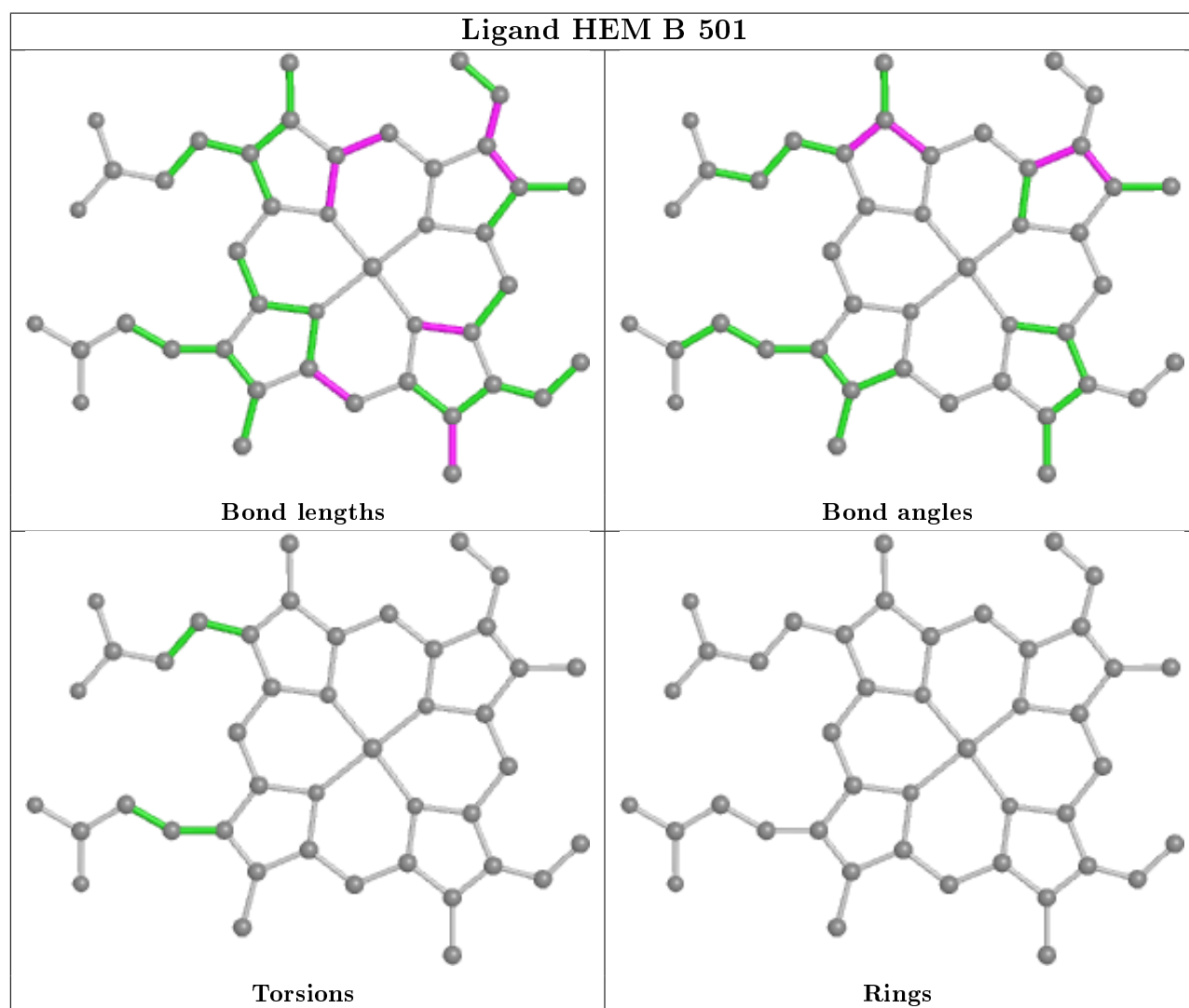
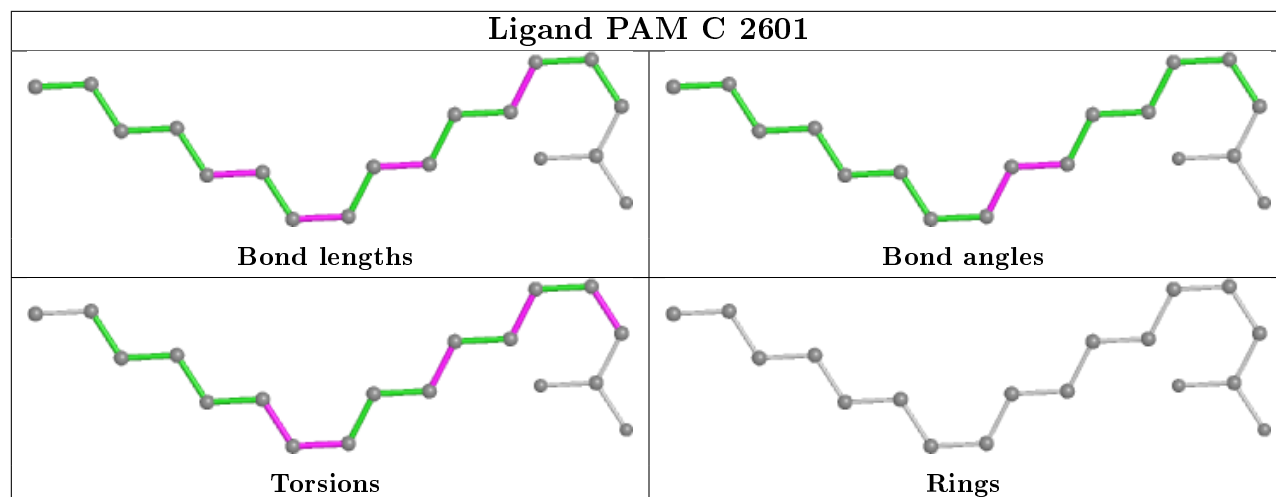
Mol	Chain	Res	Type	Atoms
3	A	601	PAM	C11-C10-C9-C8
3	B	1601	PAM	C3-C4-C5-C6
3	A	601	PAM	C3-C4-C5-C6
3	C	2601	PAM	C3-C4-C5-C6
3	A	601	PAM	C10-C11-C12-C13
3	A	601	PAM	C5-C6-C7-C8
3	C	2601	PAM	C5-C6-C7-C8
3	A	601	PAM	C9-C10-C11-C12
3	B	1601	PAM	C9-C10-C11-C12
3	C	2601	PAM	C9-C10-C11-C12

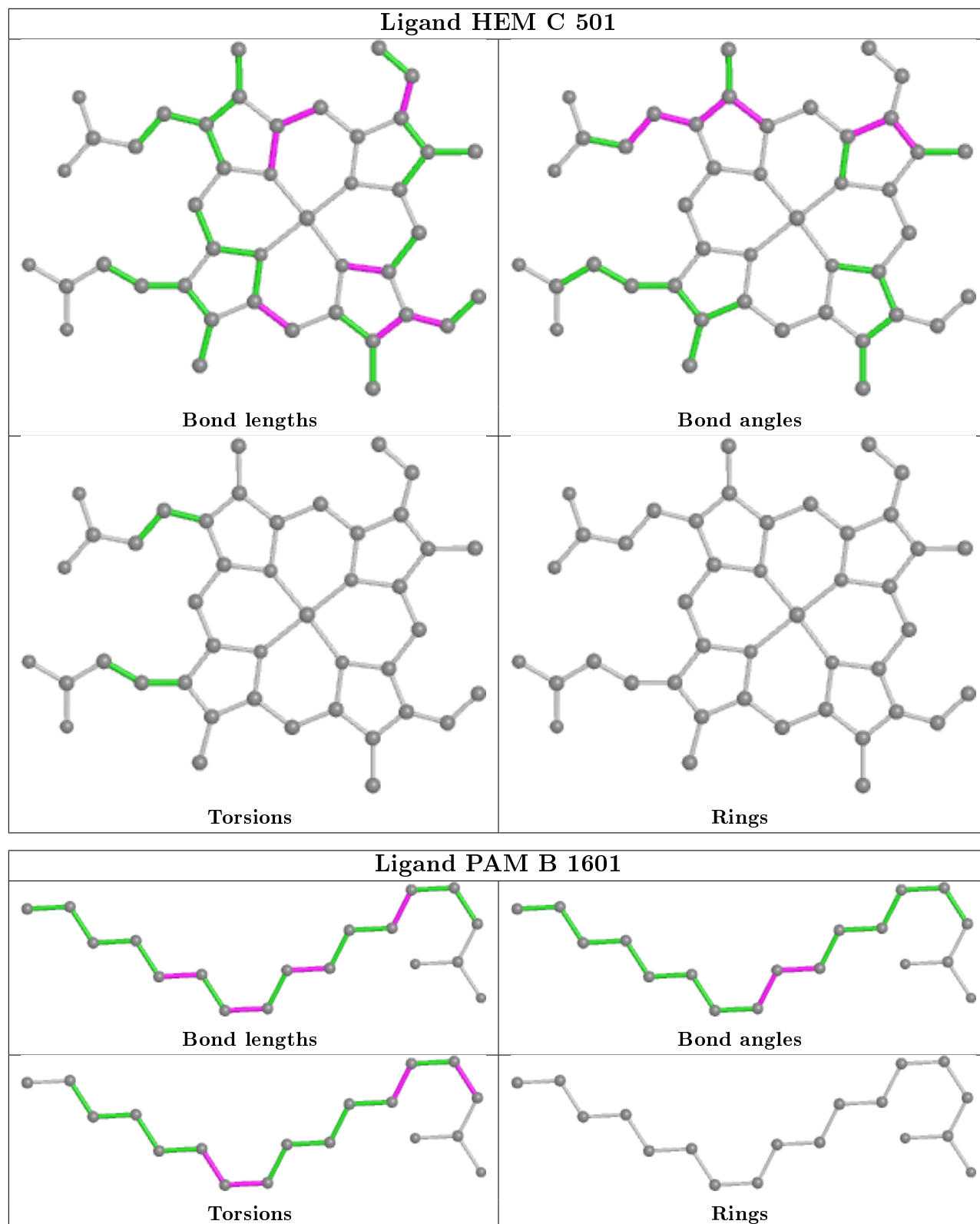
There are no ring outliers.

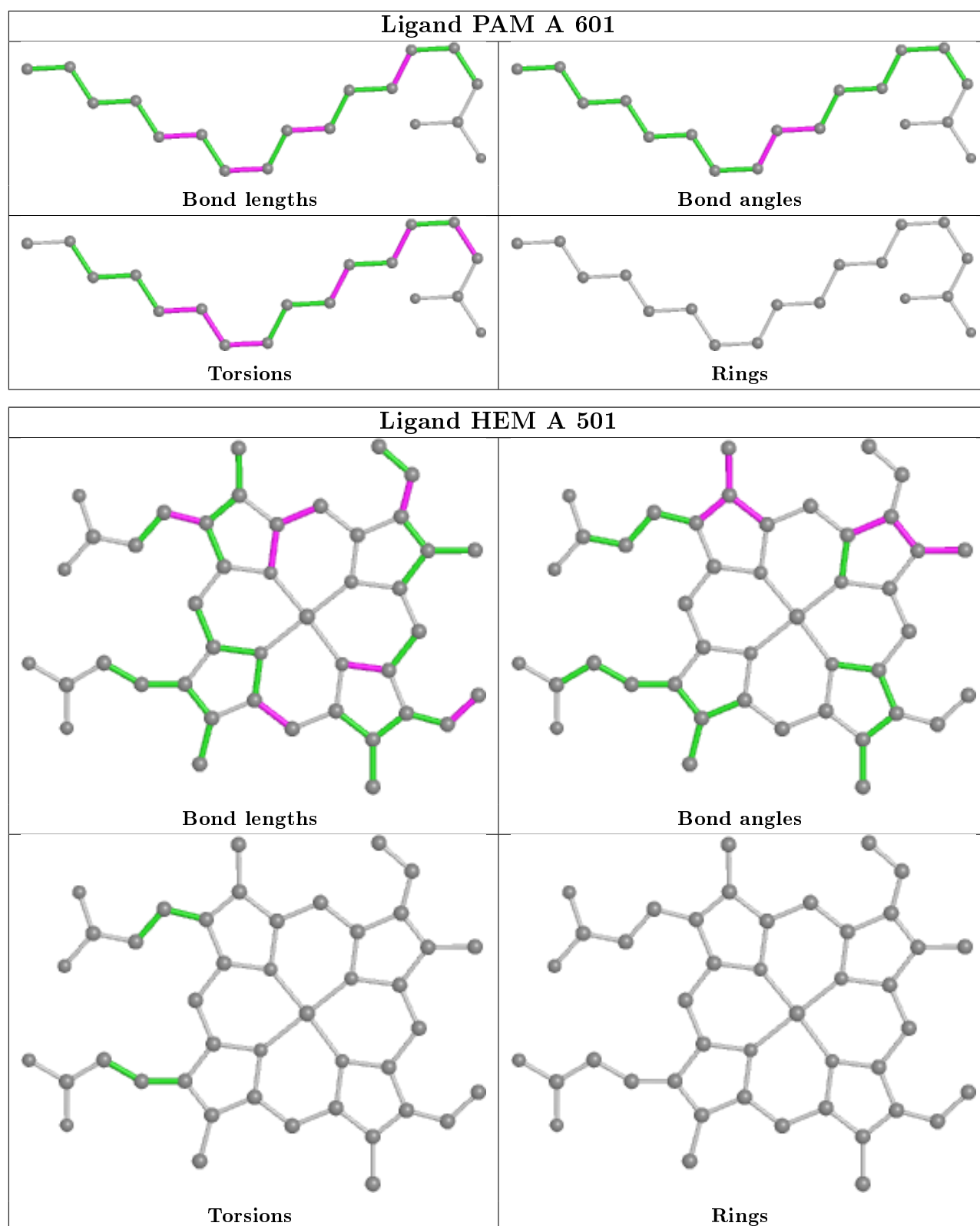
4 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	2601	PAM	2	0
2	B	501	HEM	7	0
2	C	501	HEM	5	0
2	A	501	HEM	3	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	411/417 (98%)	0.07	4 (0%) 82 85	20, 36, 57, 85	0
1	B	411/417 (98%)	0.49	31 (7%) 14 18	27, 50, 80, 107	0
1	C	411/417 (98%)	0.95	64 (15%) 2 2	36, 62, 91, 107	0
All	All	1233/1251 (98%)	0.51	99 (8%) 12 16	20, 49, 85, 107	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	267	TRP	11.2
1	C	306	CYS	7.0
1	C	341	GLU	6.8
1	C	9	LYS	6.4
1	C	303	TRP	6.0
1	C	268	LEU	5.3
1	C	7	HIS	5.1
1	C	413	ILE	5.1
1	C	266	GLU	4.7
1	C	269	ARG	4.5
1	C	11	LEU	4.4
1	B	341	GLU	4.4
1	C	36	LEU	4.3
1	B	276	ARG	4.3
1	C	305	ASN	4.3
1	C	31	ARG	4.2
1	C	330	HIS	4.1
1	C	359	LYS	4.0
1	C	84	ILE	3.9
1	B	340	ALA	3.9
1	B	305	ASN	3.9
1	B	269	ARG	3.8
1	C	272	ASN	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	11	LEU	3.7
1	B	274	ARG	3.7
1	B	334	PHE	3.6
1	C	343	GLU	3.5
1	B	30	GLU	3.5
1	C	416	LYS	3.5
1	B	268	LEU	3.4
1	C	301	PHE	3.4
1	B	331	PRO	3.4
1	C	6	PRO	3.2
1	C	347	PHE	3.2
1	C	357	ALA	3.2
1	C	414	ARG	3.2
1	B	32	TYR	3.1
1	B	342	ARG	3.1
1	C	38	GLN	3.1
1	C	57	VAL	3.1
1	C	32	TYR	3.0
1	B	261	HIS	2.9
1	B	272	ASN	2.9
1	C	383	GLN	2.9
1	C	393	LEU	2.9
1	C	278	MET	2.9
1	A	305	ASN	2.9
1	C	342	ARG	2.8
1	C	327	LEU	2.8
1	B	31	ARG	2.8
1	B	10	SER	2.8
1	C	102	LEU	2.8
1	B	36	LEU	2.7
1	C	37	PHE	2.7
1	C	333	GLU	2.6
1	C	339	PHE	2.6
1	B	271	GLY	2.6
1	C	264	TYR	2.6
1	C	360	GLY	2.6
1	C	308	PHE	2.5
1	C	302	VAL	2.5
1	C	12	ASP	2.5
1	C	47	ILE	2.5
1	C	415	ARG	2.5
1	C	274	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	12	ASP	2.4
1	C	281	GLN	2.4
1	B	382	HIS	2.4
1	C	328	TRP	2.4
1	C	340	ALA	2.4
1	C	61	THR	2.4
1	C	403	LEU	2.4
1	C	331	PRO	2.4
1	C	109	LYS	2.4
1	C	270	SER	2.3
1	C	351	PRO	2.3
1	B	258	LEU	2.3
1	B	397	LEU	2.3
1	C	63	ARG	2.2
1	A	32	TYR	2.2
1	C	329	ASP	2.2
1	B	9	LYS	2.2
1	B	392	SER	2.2
1	C	26	LYS	2.2
1	B	181	TRP	2.1
1	B	264	TYR	2.1
1	A	11	LEU	2.1
1	B	416	LYS	2.1
1	C	304	ASN	2.1
1	A	341	GLU	2.1
1	C	382	HIS	2.1
1	C	349	MET	2.1
1	C	401	PRO	2.1
1	C	265	LYS	2.0
1	B	343	GLU	2.0
1	B	414	ARG	2.0
1	C	325	PRO	2.0
1	C	174	GLY	2.0
1	B	63	ARG	2.0

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

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

6.3 Carbohydrates [i](#)

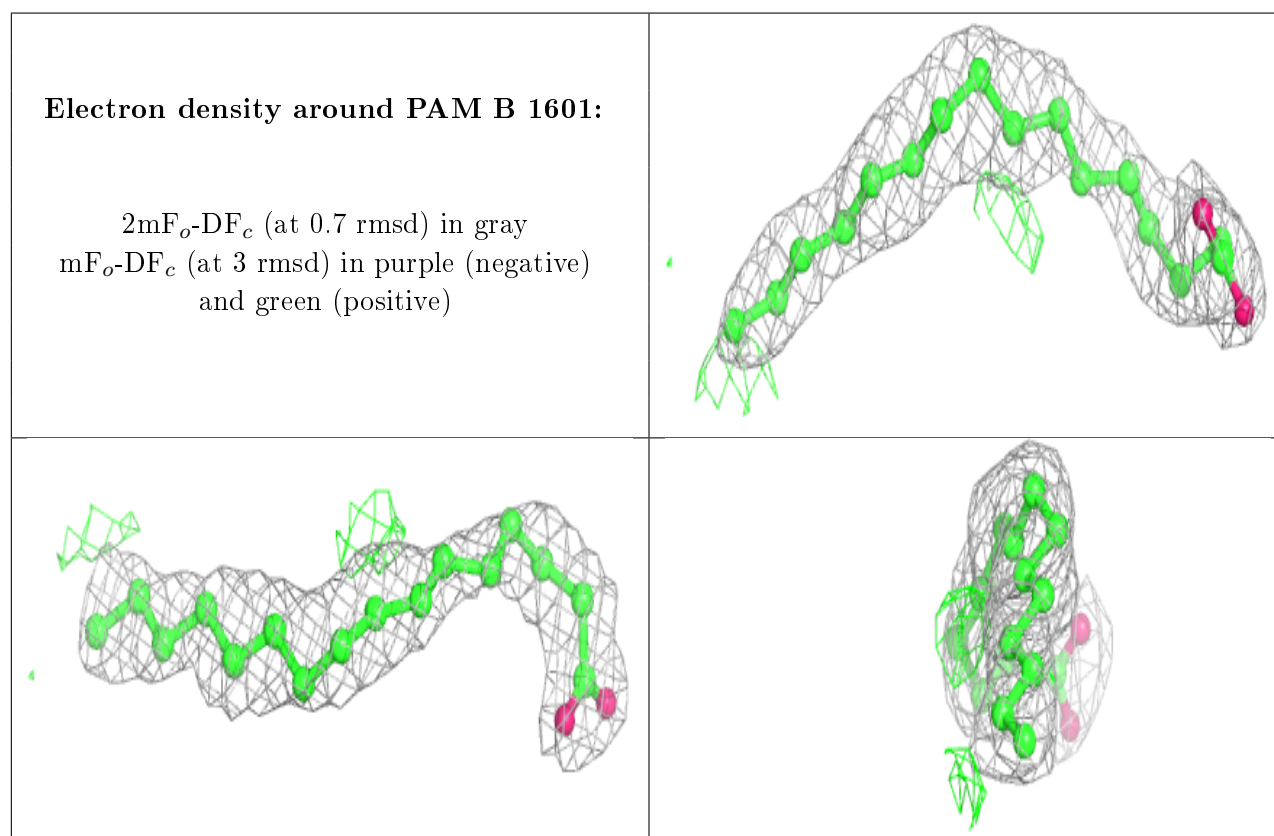
There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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.

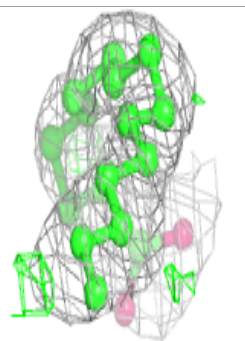
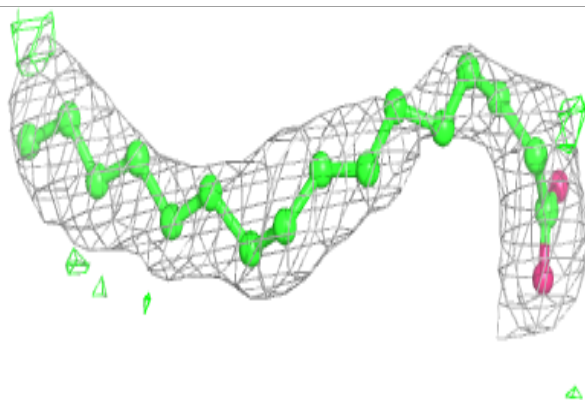
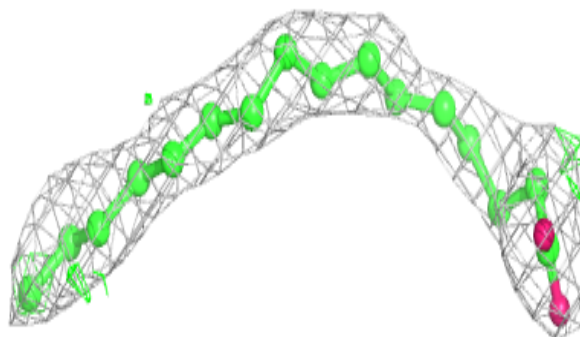
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	PAM	B	1601	18/18	0.76	0.18	56,64,70,71	0
3	PAM	C	2601	18/18	0.84	0.27	68,72,74,74	0
3	PAM	A	601	18/18	0.88	0.19	30,52,61,66	0
2	HEM	C	501	43/43	0.94	0.16	28,41,51,53	0
2	HEM	B	501	43/43	0.96	0.16	26,39,49,51	0
2	HEM	A	501	43/43	0.97	0.15	15,25,28,34	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.

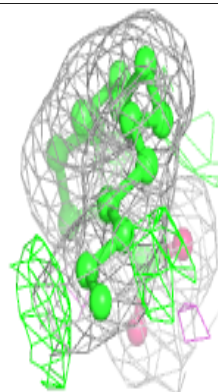
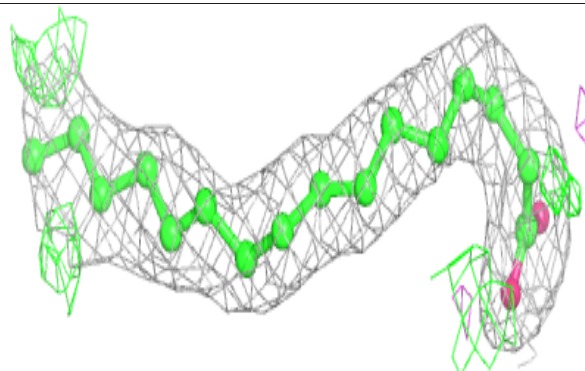
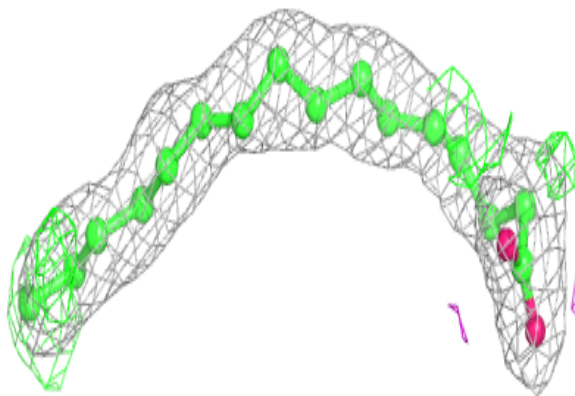


Electron density around PAM C 2601:

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

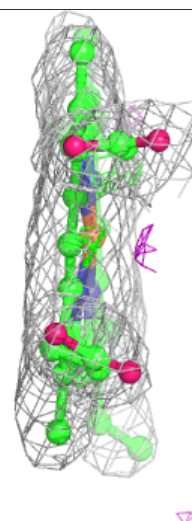
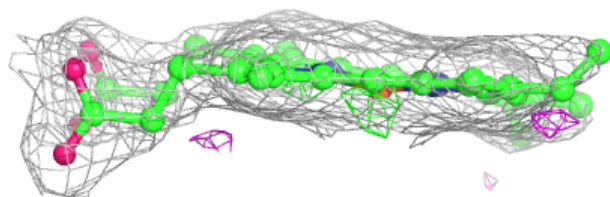
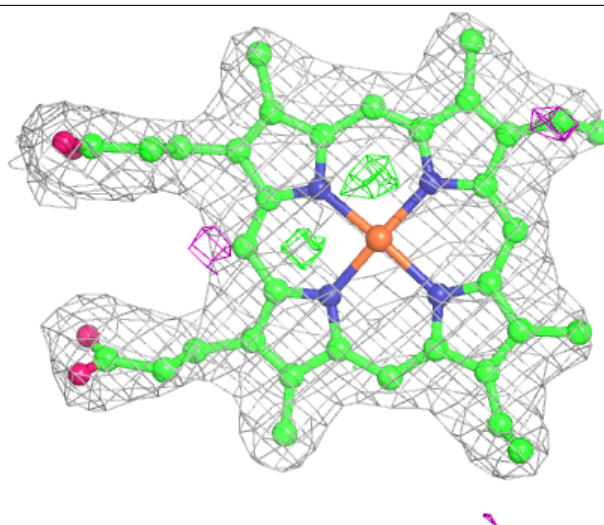
**Electron density around PAM A 601:**

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



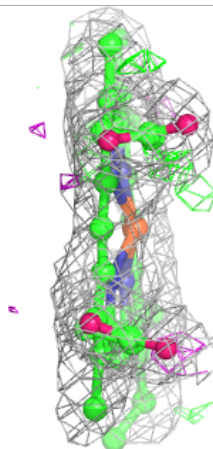
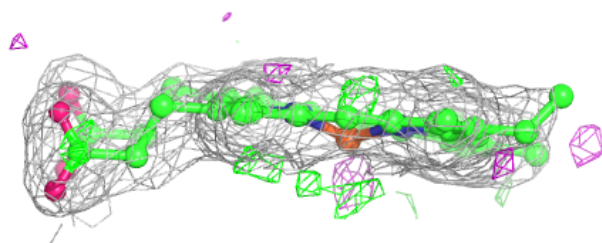
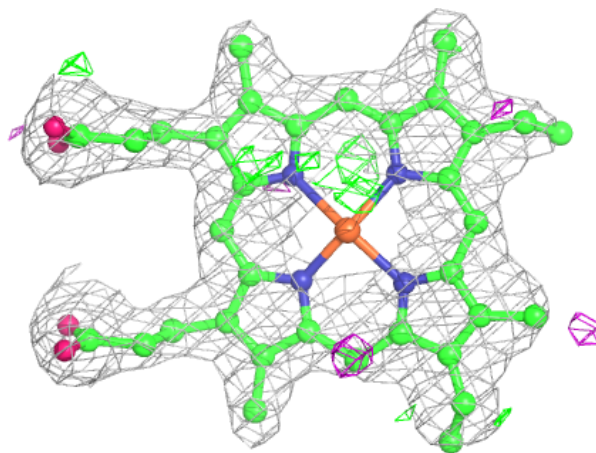
Electron density around HEM C 501:

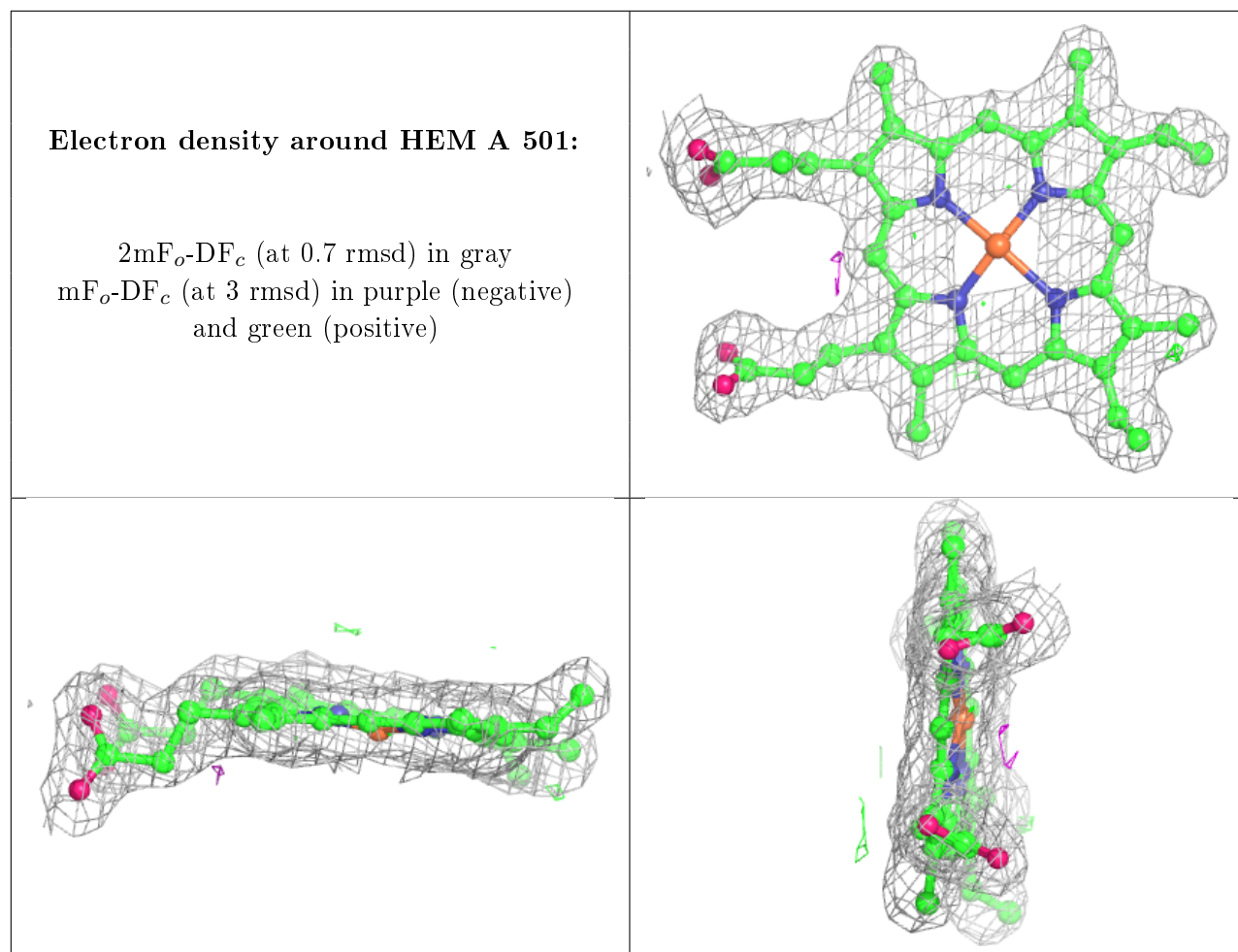
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM 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.