



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 11:38 PM UTC

PDB ID : 7Q9E / pdb_00007q9e
Title : CYP106A1
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Deposited on : 2021-11-12
Resolution : 1.70 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

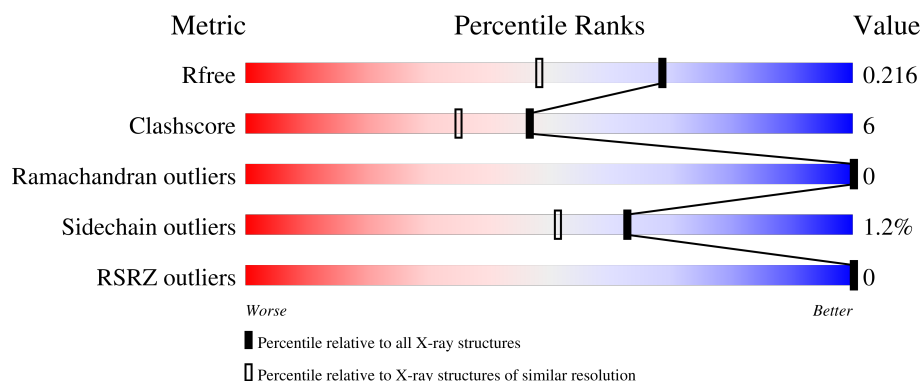
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	416	 81% 11% 8%
1	B	416	 82% 10% 8%
1	C	416	 81% 13% 5%
1	D	416	 83% 12% 5%

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

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	A	502	-	-	X	-
4	DMS	B	504	-	-	X	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 14013 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	381	Total	C	N	O	S	0	8	0
			3086	1981	513	581	11			
1	B	383	Total	C	N	O	S	0	9	0
			3111	1998	517	584	12			
1	C	397	Total	C	N	O	S	0	4	0
			3246	2087	541	606	12			
1	D	397	Total	C	N	O	S	0	4	0
			3263	2096	547	609	11			

There are 24 discrepancies between the modelled and reference sequences:

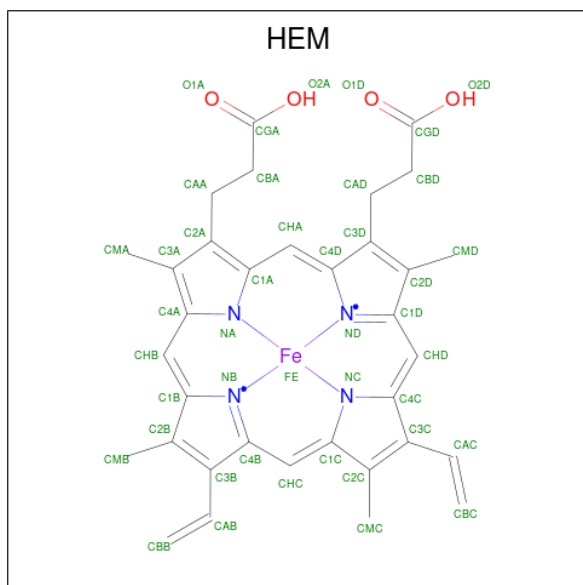
Chain	Residue	Modelled	Actual	Comment	Reference
A	411	HIS	-	expression tag	UNP D5DF35
A	412	HIS	-	expression tag	UNP D5DF35
A	413	HIS	-	expression tag	UNP D5DF35
A	414	HIS	-	expression tag	UNP D5DF35
A	415	HIS	-	expression tag	UNP D5DF35
A	416	HIS	-	expression tag	UNP D5DF35
B	411	HIS	-	expression tag	UNP D5DF35
B	412	HIS	-	expression tag	UNP D5DF35
B	413	HIS	-	expression tag	UNP D5DF35
B	414	HIS	-	expression tag	UNP D5DF35
B	415	HIS	-	expression tag	UNP D5DF35
B	416	HIS	-	expression tag	UNP D5DF35
C	411	HIS	-	expression tag	UNP D5DF35
C	412	HIS	-	expression tag	UNP D5DF35
C	413	HIS	-	expression tag	UNP D5DF35
C	414	HIS	-	expression tag	UNP D5DF35
C	415	HIS	-	expression tag	UNP D5DF35
C	416	HIS	-	expression tag	UNP D5DF35
D	411	HIS	-	expression tag	UNP D5DF35
D	412	HIS	-	expression tag	UNP D5DF35
D	413	HIS	-	expression tag	UNP D5DF35

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Chain	Residue	Modelled	Actual	Comment	Reference
D	414	HIS	-	expression tag	UNP D5DF35
D	415	HIS	-	expression tag	UNP D5DF35
D	416	HIS	-	expression tag	UNP D5DF35

- 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 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



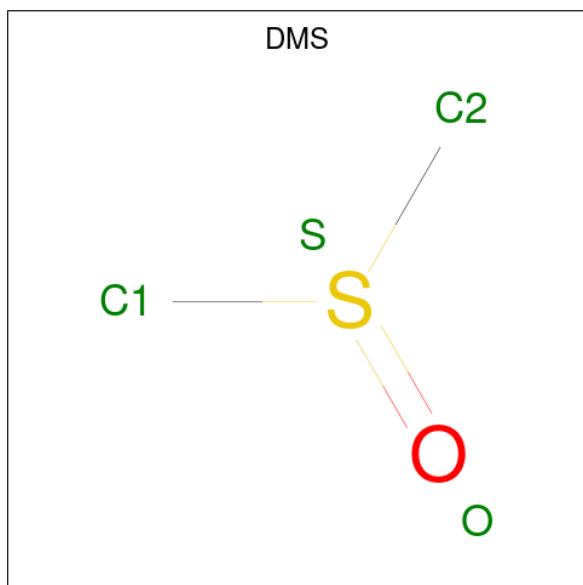
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		

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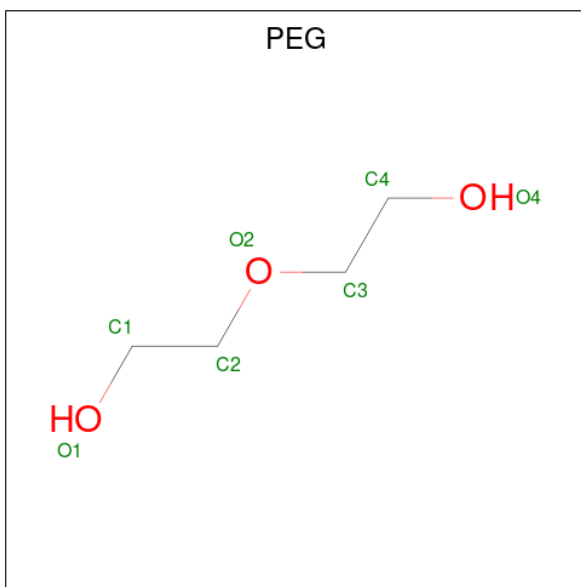
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C_2H_6OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	O	S	0	0
			4	2	1	1		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).

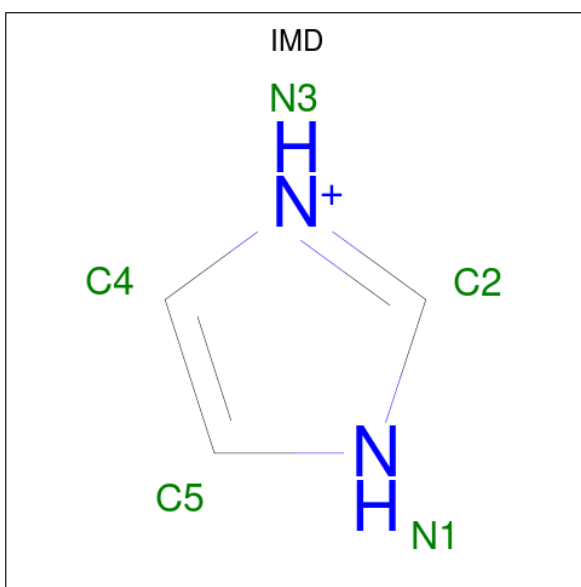


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			7	4	3		
5	C	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is COBALT (II) ION (CCD ID: CO) (formula: Co).

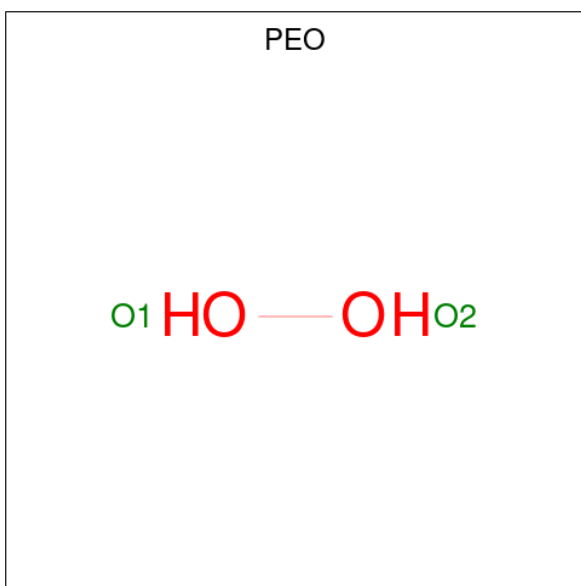
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	2	Total	Co	0	0
			2	2		
6	D	1	Total	Co	0	0
			1	1		

- Molecule 7 is IMIDAZOLE (CCD ID: IMD) (formula: C₃H₅N₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	C	1	Total	C	N	0	0
			5	3	2		

- Molecule 8 is HYDROGEN PEROXIDE (CCD ID: PEO) (formula: H_2O_2).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	C	1	Total	O	0	0
			2	2		

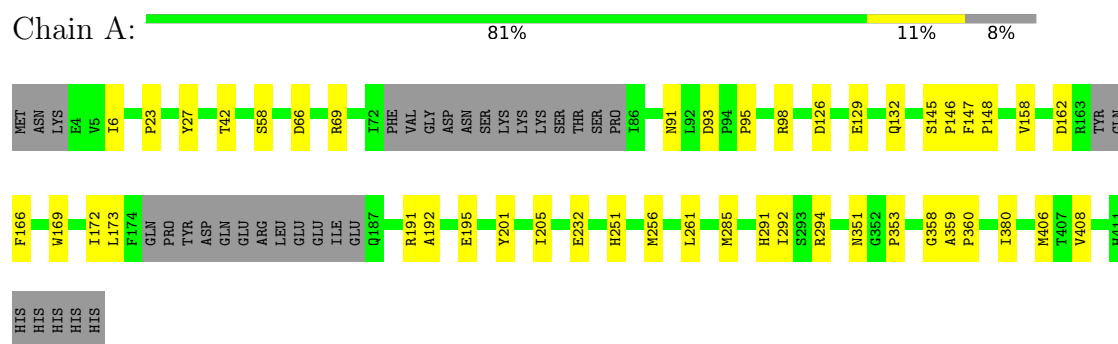
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	274	Total 274	O 274	0	0
9	B	275	Total 275	O 275	0	0
9	C	266	Total 266	O 266	0	0
9	D	207	Total 207	O 207	0	0

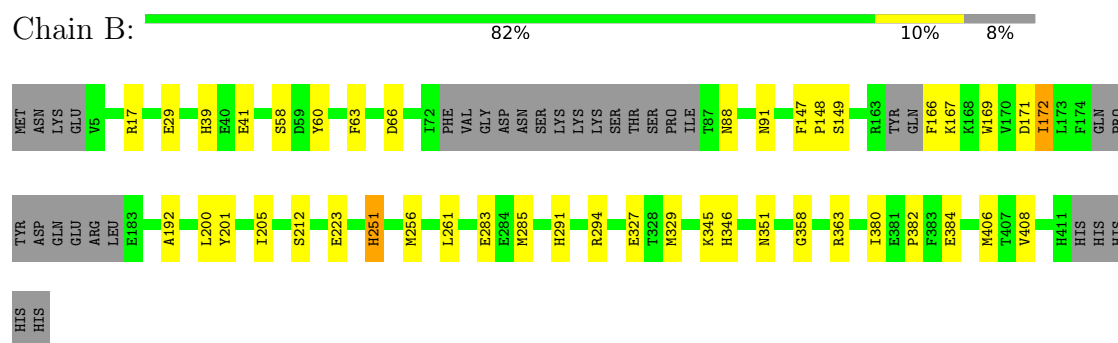
3 Residue-property plots [i](#)

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.

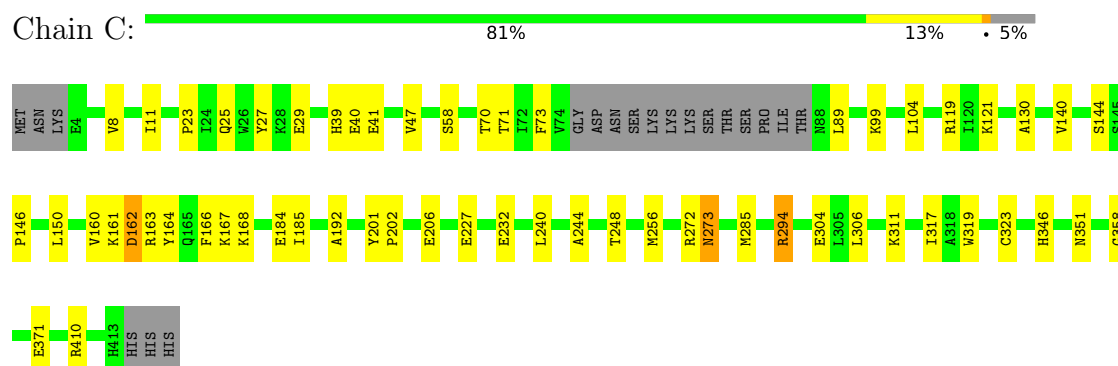
• Molecule 1: Cytochrome P450



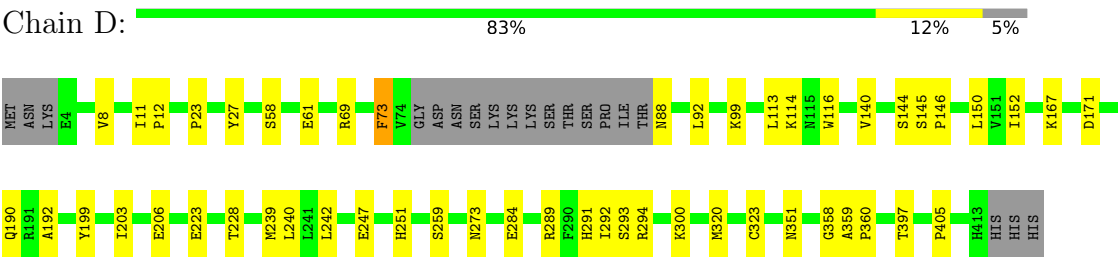
• Molecule 1: Cytochrome P450



• Molecule 1: Cytochrome P450



● Molecule 1: Cytochrome P450



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	63.69Å 82.87Å 84.86Å 95.36° 90.00° 90.00°	Depositor
Resolution (Å)	25.82 – 1.70 25.82 – 1.70	Depositor EDS
% Data completeness (in resolution range)	88.9 (25.82-1.70) 88.9 (25.82-1.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.169 , 0.206 0.180 , 0.216	Depositor DCC
R_{free} test set	8441 reflections (4.44%)	wwPDB-VP
Wilson B-factor (Å ²)	18.3	Xtriage
Anisotropy	0.023	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 27.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.389 for h,-k,-l 0.013 for -h,l,k 0.013 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	14013	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.15% 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: DMS, SO4, HEM, PEG, CO, IMD, PEO

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	1.08	2/3186 (0.1%)	1.21	1/4324 (0.0%)
1	B	1.09	2/3214 (0.1%)	1.23	2/4359 (0.0%)
1	C	1.09	2/3343 (0.1%)	1.30	7/4537 (0.2%)
1	D	1.09	2/3360 (0.1%)	1.27	5/4559 (0.1%)
All	All	1.09	8/13103 (0.1%)	1.25	15/17779 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	251	HIS	CE1-NE2	6.39	1.39	1.32
1	B	251	HIS	CE1-NE2	6.04	1.38	1.32
1	C	346	HIS	CE1-NE2	5.83	1.38	1.32
1	B	346	HIS	CE1-NE2	5.79	1.38	1.32
1	A	251	HIS	CE1-NE2	5.55	1.38	1.32
1	C	346	HIS	C-O	5.52	1.31	1.23
1	D	405	PRO	C-O	-5.39	1.17	1.23
1	A	6	ILE	N-CA	5.10	1.50	1.46

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	273	ASN	CA-CB-CG	-5.93	106.67	112.60
1	C	40	GLU	CB-CG-CD	5.83	122.51	112.60
1	D	190	GLN	CA-C-N	5.78	128.50	120.29
1	D	190	GLN	C-N-CA	5.78	128.50	120.29
1	C	40	GLU	CB-CA-C	5.76	119.93	110.88
1	C	192	ALA	CA-C-N	5.76	126.37	119.98
1	C	192	ALA	C-N-CA	5.76	126.37	119.98
1	C	206	GLU	CB-CG-CD	5.52	121.98	112.60
1	A	93	ASP	CA-CB-CG	5.47	118.07	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	273	ASN	CA-CB-CG	-5.37	107.23	112.60
1	B	171	ASP	CA-C-N	5.31	127.35	120.56
1	B	171	ASP	C-N-CA	5.31	127.35	120.56
1	C	70	THR	CB-CA-C	5.25	122.37	110.67
1	D	192	ALA	CA-C-N	5.10	125.64	119.98
1	D	192	ALA	C-N-CA	5.10	125.64	119.98

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3086	0	2970	24	1
1	B	3111	0	2998	31	0
1	C	3246	0	3129	50	1
1	D	3263	0	3160	35	0
2	A	43	0	30	1	0
2	B	43	0	30	1	0
2	C	43	0	30	3	0
2	D	43	0	30	2	0
3	A	30	0	0	3	0
3	B	25	0	0	1	0
3	C	20	0	0	1	0
3	D	10	0	0	0	0
4	B	4	0	6	6	0
5	B	7	0	10	0	0
5	C	7	0	10	1	0
6	B	2	0	0	0	0
6	D	1	0	0	0	0
7	C	5	0	5	0	0
8	C	2	0	0	0	0
9	A	274	0	0	2	0
9	B	275	0	0	4	0
9	C	266	0	0	9	0
9	D	207	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	14013	0	12408	144	1

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 6.

All (144) 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:163:ARG:O	1:C:167:LYS:N	2.08	0.85
1:B:39:HIS:CE1	1:B:41:GLU:HG2	2.18	0.79
1:C:163:ARG:O	1:C:167:LYS:CB	2.30	0.79
1:A:66:ASP:HB3	1:A:91:ASN:O	1.86	0.76
1:D:58:SER:HB3	1:D:351:ASN:HD22	1.50	0.74
1:C:371:GLU:OE1	9:C:601:HOH:O	2.06	0.73
1:D:199:TYR:CZ	1:D:203:ILE:HD11	2.23	0.73
1:C:163:ARG:O	1:C:167:LYS:HB3	1.89	0.72
1:B:29:GLU:HG3	9:B:618:HOH:O	1.90	0.70
1:C:121:LYS:HE2	9:C:601:HOH:O	1.91	0.70
1:C:304:GLU:OE1	9:C:602:HOH:O	2.09	0.70
1:D:293[A]:SER:OG	1:D:320:MET:HE2	1.92	0.70
1:C:162:ASP:O	1:C:166:PHE:N	2.28	0.65
1:D:199:TYR:CZ	1:D:203:ILE:CD1	2.79	0.65
1:C:73:PHE:HZ	1:C:319:TRP:CZ3	2.16	0.64
1:A:98:ARG:HD3	3:A:502:SO4:O2	1.98	0.63
1:D:167:LYS:NZ	1:D:171:ASP:OD2	2.30	0.63
4:B:504:DMS:C1	9:B:764:HOH:O	2.46	0.63
1:C:58[B]:SER:HB3	1:C:351:ASN:HD22	1.63	0.63
1:A:95:PRO:HB3	1:B:327[A]:GLU:HG3	1.81	0.63
1:C:244:ALA:HB2	5:C:505:PEG:H32	1.80	0.62
1:C:25:GLN:NE2	9:C:605:HOH:O	2.31	0.62
1:B:329[A]:MET:HE2	1:B:345:LYS:HD3	1.81	0.62
1:A:172:ILE:HG22	1:A:192:ALA:HB2	1.81	0.61
1:B:251:HIS:HE1	1:B:291:HIS:O	1.82	0.61
4:B:504:DMS:H13	9:B:764:HOH:O	1.99	0.61
1:B:149:SER:CB	1:B:167:LYS:HG2	2.31	0.60
1:B:172:ILE:HG22	1:B:192:ALA:HB1	1.84	0.60
1:D:73:PHE:CE2	1:D:397:THR:C	2.80	0.60
1:C:273:ASN:ND2	3:C:507:SO4:O1	2.34	0.59
1:B:149:SER:HB2	1:B:167:LYS:HG2	1.85	0.58
1:B:58:SER:HB3	1:B:351:ASN:HD22	1.68	0.58
1:C:73:PHE:CZ	1:C:319:TRP:CZ3	2.91	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:160:VAL:HG12	1:C:161:LYS:HG2	1.86	0.57
1:B:384:GLU:H	4:B:504:DMS:H23	1.69	0.57
1:D:150:LEU:HD21	1:D:167:LYS:HD3	1.85	0.57
1:C:73:PHE:CZ	1:C:294:ARG:HG2	2.40	0.56
1:D:69:ARG:NH1	9:D:604:HOH:O	2.37	0.56
1:A:58:SER:HB3	1:A:351:ASN:HD22	1.70	0.56
1:A:380:ILE:HG22	1:A:408:VAL:HG12	1.88	0.55
1:C:8:VAL:HG23	1:C:71:THR:HB	1.88	0.55
1:D:199:TYR:CE2	1:D:203:ILE:CD1	2.90	0.55
1:C:8:VAL:CG2	1:C:71:THR:HB	2.37	0.55
1:C:150:LEU:O	1:C:163:ARG:HD3	2.08	0.54
1:B:256:MET:HG3	1:B:285:MET:HG3	1.90	0.53
1:C:358:GLY:HA3	2:C:501:HEM:C3C	2.43	0.53
1:A:358:GLY:HA3	2:A:501:HEM:C3C	2.44	0.52
2:C:501:HEM:HMC1	2:C:501:HEM:HBC2	1.90	0.52
1:B:17:ARG:NH2	3:B:502:SO4:O4	2.41	0.52
1:C:119:ARG:NH2	9:C:613:HOH:O	2.43	0.52
1:A:126:ASP:OD2	9:A:601:HOH:O	2.18	0.52
1:A:145[A]:SER:OG	1:A:146:PRO:HD3	2.09	0.52
1:C:89:LEU:HD22	1:C:104:LEU:HD12	1.92	0.51
1:D:11:ILE:HB	1:D:12:PRO:HD3	1.92	0.51
1:C:163:ARG:O	1:C:167:LYS:CA	2.58	0.51
1:D:247:GLU:HG2	9:D:750:HOH:O	2.08	0.51
1:C:162:ASP:N	1:C:162:ASP:OD1	2.44	0.51
1:C:184:GLU:HG3	1:C:185:ILE:HG23	1.93	0.51
1:C:23:PRO:HG2	1:C:27:TYR:CZ	2.46	0.50
1:C:99:LYS:HD2	1:C:227:GLU:OE1	2.11	0.50
1:C:8:VAL:HG13	1:C:11:ILE:HD12	1.93	0.50
1:A:98:ARG:CD	3:A:502:SO4:O2	2.60	0.50
1:A:166:PHE:HA	1:A:169:TRP:CD1	2.46	0.50
1:B:384:GLU:H	4:B:504:DMS:C2	2.25	0.50
1:B:201:TYR:CE2	1:B:205:ILE:HD11	2.47	0.49
1:D:23:PRO:HG2	1:D:27:TYR:CZ	2.48	0.49
1:A:158:VAL:CG1	1:A:162:ASP:HB2	2.43	0.49
1:D:358:GLY:HA3	2:D:501:HEM:C3C	2.48	0.49
1:A:261:LEU:HD11	1:A:406:MET:HE2	1.95	0.48
1:B:172:ILE:HG22	1:B:192:ALA:CB	2.43	0.48
1:A:256:MET:HG3	1:A:285:MET:HG3	1.94	0.48
1:B:358:GLY:HA3	2:B:501:HEM:C3C	2.48	0.48
1:D:150:LEU:CD2	1:D:167:LYS:HD3	2.42	0.48
1:C:164:TYR:CD1	1:C:167:LYS:HE2	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:382:PRO:O	4:B:504:DMS:H11	2.13	0.48
1:B:251:HIS:CE1	1:B:291:HIS:O	2.66	0.48
1:B:261:LEU:HD11	1:B:406:MET:HE2	1.96	0.47
1:D:8:VAL:CG2	1:D:73:PHE:HB3	2.43	0.47
1:D:359:ALA:HB3	1:D:360:PRO:HD3	1.96	0.47
1:D:259:SER:HB2	1:D:289:ARG:HG3	1.97	0.47
1:C:47:VAL:HG11	1:C:306:LEU:HD13	1.97	0.47
1:D:223:GLU:HG3	1:D:228:THR:HG22	1.96	0.47
1:D:293[A]:SER:OG	1:D:320:MET:CE	2.63	0.47
1:C:164:TYR:O	1:C:168:LYS:HB2	2.14	0.47
1:C:227:GLU:HA	9:C:756:HOH:O	2.15	0.47
1:C:311:LYS:NZ	9:C:619:HOH:O	2.47	0.47
1:B:384:GLU:HB2	4:B:504:DMS:H12	1.97	0.46
1:C:240:LEU:HD23	1:C:240:LEU:C	2.41	0.46
1:D:88:ASN:O	1:D:92:LEU:HG	2.15	0.46
1:B:380:ILE:HG22	1:B:408:VAL:HG12	1.98	0.46
1:B:283:GLU:OE2	1:B:363[B]:ARG:CD	2.64	0.46
1:A:359:ALA:HB3	1:A:360:PRO:HD3	1.98	0.45
1:C:25:GLN:O	1:C:29:GLU:HG3	2.16	0.45
1:D:73:PHE:HE2	1:D:397:THR:O	2.00	0.45
1:C:8:VAL:O	1:C:8:VAL:CG1	2.64	0.45
1:C:184:GLU:HG3	1:C:185:ILE:N	2.31	0.45
1:D:113:LEU:HD22	1:D:116:TRP:CZ3	2.52	0.45
1:B:223:GLU:HG2	9:B:613:HOH:O	2.16	0.45
1:D:73:PHE:HE2	1:D:397:THR:C	2.25	0.45
1:C:39:HIS:CE1	1:C:41:GLU:CG	3.00	0.45
1:B:166:PHE:HA	1:B:169:TRP:CD1	2.53	0.44
1:C:130:ALA:HB3	9:C:682:HOH:O	2.16	0.44
1:A:23:PRO:HG2	1:A:27:TYR:CZ	2.52	0.44
1:A:129:GLU:O	1:A:132:GLN:HG2	2.18	0.44
1:D:140:VAL:HA	1:D:144:SER:OG	2.18	0.44
1:D:223:GLU:HB3	9:D:657:HOH:O	2.18	0.44
1:D:284:GLU:HA	1:D:284:GLU:OE1	2.17	0.44
1:A:42:THR:O	1:A:69:ARG:NH2	2.50	0.43
2:D:501:HEM:HMC1	2:D:501:HEM:HBC2	2.00	0.43
1:D:61:GLU:O	1:D:300:LYS:HD3	2.18	0.43
1:C:248:THR:HG23	9:C:797:HOH:O	2.18	0.43
1:C:71:THR:HG21	1:C:317:ILE:HD11	2.00	0.43
1:B:283:GLU:OE2	1:B:363[B]:ARG:HD3	2.19	0.43
1:D:152:ILE:HG21	1:D:242:LEU:CD2	2.49	0.43
1:C:73:PHE:CE1	1:C:294:ARG:HD3	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:ARG:O	1:A:195:GLU:HG3	2.19	0.43
1:A:129:GLU:HG3	9:A:809:HOH:O	2.18	0.43
1:D:240:LEU:C	1:D:240:LEU:HD23	2.44	0.43
1:A:201:TYR:CE2	1:A:205:ILE:HD11	2.53	0.42
1:B:39:HIS:CE1	1:B:41:GLU:CG	2.96	0.42
1:C:164:TYR:O	1:C:168:LYS:N	2.49	0.42
1:C:272:ARG:HH11	1:C:272:ARG:HG3	1.83	0.42
1:B:201:TYR:CZ	1:B:205:ILE:HD11	2.54	0.42
1:C:58[A]:SER:HB2	1:C:351:ASN:HD22	1.84	0.42
1:C:164:TYR:CE1	1:C:167:LYS:HE2	2.54	0.42
1:D:73:PHE:CE2	1:D:292:ILE:HB	2.54	0.42
1:D:114:LYS:HA	1:D:114:LYS:HD2	1.88	0.42
1:A:147:PHE:HB3	1:A:148:PRO:HD3	2.01	0.42
1:B:66:ASP:O	1:B:66:ASP:CG	2.62	0.42
1:A:291:HIS:HA	1:A:292:ILE:HA	1.86	0.42
1:C:73:PHE:HE1	1:C:319:TRP:CH2	2.38	0.42
1:B:88:ASN:OD1	1:B:91:ASN:HB2	2.20	0.41
1:C:201:TYR:N	1:C:202:PRO:HD2	2.35	0.41
1:D:291:HIS:HA	1:D:292:ILE:HA	1.88	0.41
1:A:353:PRO:HB3	3:A:502:SO4:O1	2.20	0.41
1:C:146:PRO:HB3	1:C:167:LYS:NZ	2.35	0.41
1:D:145[A]:SER:OG	1:D:146:PRO:HD3	2.19	0.41
1:C:140:VAL:HA	1:C:144:SER:OG	2.21	0.41
1:D:73:PHE:CZ	1:D:397:THR:HA	2.56	0.41
1:C:256:MET:HG3	1:C:285:MET:HG3	2.02	0.40
2:C:501:HEM:HBC2	2:C:501:HEM:CMC	2.51	0.40
1:B:60:TYR:HA	1:B:63:PHE:O	2.21	0.40
1:B:147:PHE:HB3	1:B:148:PRO:HD3	2.02	0.40
1:D:8:VAL:HG22	1:D:73:PHE:HB3	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:GLU:OE2	1:C:232:GLU:OE2[1_545]	2.12	0.08

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	381/416 (92%)	369 (97%)	12 (3%)	0	100	100
1	B	384/416 (92%)	371 (97%)	13 (3%)	0	100	100
1	C	397/416 (95%)	382 (96%)	15 (4%)	0	100	100
1	D	397/416 (95%)	383 (96%)	14 (4%)	0	100	100
All	All	1559/1664 (94%)	1505 (96%)	54 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	334/377 (89%)	332 (99%)	2 (1%)	78	72
1	B	336/377 (89%)	332 (99%)	4 (1%)	63	51
1	C	352/377 (93%)	348 (99%)	4 (1%)	65	54
1	D	356/377 (94%)	350 (98%)	6 (2%)	53	38
All	All	1378/1508 (91%)	1362 (99%)	16 (1%)	63	51

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	173	LEU
1	A	294	ARG

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Mol	Chain	Res	Type
1	B	172	ILE
1	B	200	LEU
1	B	212	SER
1	B	294	ARG
1	C	162	ASP
1	C	294	ARG
1	C	323	CYS
1	C	410	ARG
1	D	73	PHE
1	D	99	LYS
1	D	206	GLU
1	D	239	MET
1	D	294	ARG
1	D	323	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	ASN
1	A	198	GLN
1	A	221	GLN
1	A	236	HIS
1	A	332	ASN
1	A	351	ASN
1	B	198	GLN
1	B	221	GLN
1	B	251	HIS
1	B	332	ASN
1	B	351	ASN
1	C	15	GLN
1	C	25	GLN
1	C	33	ASN
1	C	55	GLN
1	C	175	GLN
1	C	187	GLN
1	C	198	GLN
1	C	221	GLN
1	C	301	GLN
1	C	351	ASN
1	D	33	ASN
1	D	175	GLN
1	D	179	GLN

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Mol	Chain	Res	Type
1	D	187	GLN
1	D	198	GLN
1	D	255	ASN
1	D	301	GLN
1	D	351	ASN

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](#)

Of 29 ligands modelled in this entry, 3 are monoatomic - leaving 26 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	SO4	C	507	-	4,4,4	0.29	0	6,6,6	0.26	0
8	PEO	C	508	-	1,1,1	1.16	0	-		
3	SO4	B	505	-	4,4,4	0.33	0	6,6,6	0.09	0
3	SO4	A	504	-	4,4,4	0.31	0	6,6,6	0.07	0
3	SO4	A	505	-	4,4,4	0.26	0	6,6,6	0.13	0
3	SO4	C	503	-	4,4,4	0.33	0	6,6,6	0.09	0
4	DMS	B	504	-	3,3,3	0.19	0	3,3,3	0.33	0
3	SO4	B	510	-	4,4,4	0.29	0	6,6,6	0.14	0
2	HEM	C	501	1,9	50,50,50	1.58	8 (16%)	67,82,82	1.57	13 (19%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	D	502	-	4,4,4	0.27	0	6,6,6	0.11	0
3	SO4	C	506	-	4,4,4	0.32	0	6,6,6	0.16	0
3	SO4	A	503	-	4,4,4	0.29	0	6,6,6	0.11	0
3	SO4	B	502	-	4,4,4	0.26	0	6,6,6	0.15	0
3	SO4	D	503	-	4,4,4	0.35	0	6,6,6	0.12	0
5	PEG	C	505	-	6,6,6	0.27	0	5,5,5	0.16	0
2	HEM	A	501	1,9	50,50,50	1.49	10 (20%)	67,82,82	1.68	12 (17%)
5	PEG	B	506	-	6,6,6	0.18	0	5,5,5	0.13	0
2	HEM	B	501	1,9	50,50,50	1.48	8 (16%)	67,82,82	1.51	12 (17%)
3	SO4	B	503	-	4,4,4	0.31	0	6,6,6	0.13	0
2	HEM	D	501	1,9	50,50,50	1.65	10 (20%)	67,82,82	1.77	14 (20%)
3	SO4	A	506	-	4,4,4	0.31	0	6,6,6	0.13	0
7	IMD	C	504	-	5,5,5	0.15	0	5,5,5	0.62	0
3	SO4	B	509	-	4,4,4	0.32	0	6,6,6	0.12	0
3	SO4	A	502	-	4,4,4	0.34	0	6,6,6	0.22	0
3	SO4	A	507	-	4,4,4	0.32	0	6,6,6	0.11	0
3	SO4	C	502	-	4,4,4	0.29	0	6,6,6	0.21	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
2	HEM	D	501	1,9	-	0/14/54/54	-
7	IMD	C	504	-	-	-	0/1/1/1
2	HEM	C	501	1,9	-	1/14/54/54	-
5	PEG	C	505	-	-	3/4/4/4	-
2	HEM	A	501	1,9	-	2/14/54/54	-
5	PEG	B	506	-	-	3/4/4/4	-
2	HEM	B	501	1,9	-	1/14/54/54	-

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	HEM	FE-NC	4.26	2.09	1.95
2	C	501	HEM	FE-NC	4.20	2.09	1.95
2	C	501	HEM	C4D-ND	-4.18	1.33	1.40
2	B	501	HEM	C1B-NB	-4.05	1.33	1.40
2	A	501	HEM	FE-NC	3.96	2.08	1.95
2	A	501	HEM	FE-NB	3.85	2.06	1.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	HEM	FE-NB	3.73	2.06	1.94
2	C	501	HEM	FE-NB	3.72	2.06	1.94
2	B	501	HEM	C4D-ND	-3.70	1.33	1.40
2	A	501	HEM	C4D-ND	-3.48	1.34	1.40
2	D	501	HEM	C4D-ND	-3.40	1.34	1.40
2	B	501	HEM	FE-NB	3.26	2.04	1.94
2	C	501	HEM	C4D-C3D	3.24	1.50	1.45
2	B	501	HEM	C4A-NA	-3.20	1.33	1.39
2	D	501	HEM	C1A-NA	-3.05	1.33	1.39
2	B	501	HEM	FE-NC	2.92	2.04	1.95
2	A	501	HEM	C1B-NB	-2.78	1.35	1.40
2	B	501	HEM	FE-NA	2.74	2.04	1.95
2	B	501	HEM	C4B-NB	-2.72	1.33	1.38
2	D	501	HEM	CHB-C1B	2.69	1.43	1.38
2	D	501	HEM	C1B-NB	-2.65	1.35	1.40
2	D	501	HEM	C1C-NC	-2.47	1.35	1.39
2	A	501	HEM	CAA-C2A	2.42	1.57	1.51
2	D	501	HEM	C3B-C4B	2.34	1.49	1.44
2	C	501	HEM	CMD-C2D	2.31	1.55	1.50
2	A	501	HEM	C4D-C3D	2.29	1.48	1.45
2	C	501	HEM	C1D-C2D	2.28	1.49	1.44
2	A	501	HEM	C1C-C2C	-2.23	1.40	1.45
2	D	501	HEM	C4B-NB	-2.19	1.34	1.38
2	D	501	HEM	C4C-NC	-2.14	1.35	1.39
2	C	501	HEM	CMC-C2C	2.14	1.55	1.50
2	B	501	HEM	C4C-NC	-2.09	1.35	1.39
2	C	501	HEM	CBA-CGA	2.06	1.55	1.50
2	A	501	HEM	C3C-C4C	-2.04	1.42	1.46
2	A	501	HEM	CAC-C3C	2.02	1.52	1.47
2	A	501	HEM	FE-NA	2.01	2.01	1.95

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	CHC-C4B-NB	5.46	130.29	124.42
2	D	501	HEM	CHD-C4C-NC	5.10	130.00	124.45
2	A	501	HEM	C1B-NB-C4B	4.72	110.80	105.21
2	A	501	HEM	O2A-CGA-CBA	4.00	126.65	114.00
2	C	501	HEM	C1B-NB-C4B	3.99	109.93	105.21
2	C	501	HEM	O2A-CGA-CBA	3.93	126.43	114.00
2	B	501	HEM	C1B-NB-C4B	3.88	109.80	105.21
2	B	501	HEM	CHB-C1B-NB	3.82	129.09	124.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	HEM	CHD-C1D-ND	3.76	128.47	124.42
2	D	501	HEM	C1B-NB-C4B	3.64	109.51	105.21
2	D	501	HEM	CHA-C4D-ND	3.63	128.85	124.37
2	A	501	HEM	C3B-C4B-NB	-3.61	106.87	109.47
2	D	501	HEM	O2A-CGA-CBA	3.52	125.11	114.00
2	D	501	HEM	CHC-C4B-NB	3.49	128.18	124.42
2	D	501	HEM	CHA-C4D-C3D	-3.44	118.88	125.23
2	C	501	HEM	CHB-C1B-NB	3.43	128.61	124.37
2	B	501	HEM	CHC-C4B-NB	3.41	128.09	124.42
2	B	501	HEM	CHD-C1D-C2D	-3.35	119.74	125.03
2	A	501	HEM	CHB-C1B-NB	3.32	128.47	124.37
2	D	501	HEM	CMD-C2D-C1D	3.26	130.13	125.03
2	B	501	HEM	C4B-C3B-C2B	-3.24	104.30	107.28
2	A	501	HEM	CHD-C1D-ND	3.24	127.91	124.42
2	D	501	HEM	C4B-C3B-C2B	-3.16	104.38	107.28
2	C	501	HEM	O1A-CGA-CBA	-3.14	113.12	123.09
2	D	501	HEM	C4C-NC-C1C	2.97	110.67	105.82
2	A	501	HEM	O1D-CGD-CBD	-2.93	113.80	123.09
2	C	501	HEM	CHC-C4B-NB	2.93	127.58	124.42
2	C	501	HEM	CHD-C1D-ND	2.80	127.44	124.42
2	A	501	HEM	C3B-C2B-C1B	2.77	108.49	106.41
2	A	501	HEM	O1A-CGA-CBA	-2.74	114.41	123.09
2	D	501	HEM	O1A-CGA-CBA	-2.65	114.70	123.09
2	C	501	HEM	C3B-C2B-C1B	2.58	108.35	106.41
2	B	501	HEM	C3B-C2B-C1B	2.55	108.33	106.41
2	B	501	HEM	O2A-CGA-O1A	-2.47	116.99	123.33
2	C	501	HEM	C3B-C4B-NB	-2.46	107.70	109.47
2	A	501	HEM	CAA-C2A-C1A	2.46	129.74	124.94
2	D	501	HEM	C3C-C2C-C1C	-2.41	104.77	107.05
2	D	501	HEM	CAD-C3D-C4D	2.38	128.85	124.70
2	C	501	HEM	CHC-C1C-NC	2.31	126.96	124.45
2	D	501	HEM	CHC-C1C-NC	2.30	126.96	124.45
2	C	501	HEM	CHD-C1D-C2D	-2.24	121.50	125.03
2	C	501	HEM	C4C-C3C-C2C	2.22	108.74	106.81
2	B	501	HEM	CAD-C3D-C4D	2.16	128.47	124.70
2	A	501	HEM	CHD-C1D-C2D	-2.15	121.64	125.03
2	C	501	HEM	CMA-C3A-C4A	2.14	128.68	125.42
2	B	501	HEM	O1D-CGD-CBD	-2.10	116.44	123.09
2	D	501	HEM	C1A-CHA-C4D	-2.08	121.35	126.25
2	A	501	HEM	CAB-C3B-C2B	2.05	135.11	128.43
2	B	501	HEM	O2D-CGD-CBD	2.05	120.47	114.00
2	C	501	HEM	CHA-C4D-ND	2.02	126.86	124.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	HEM	C3D-C4D-ND	2.01	112.37	110.17

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	505	PEG	O1-C1-C2-O2
5	B	506	PEG	O2-C3-C4-O4
5	C	505	PEG	C4-C3-O2-C2
5	B	506	PEG	C4-C3-O2-C2
5	C	505	PEG	O2-C3-C4-O4
2	A	501	HEM	CAA-CBA-CGA-O1A
5	B	506	PEG	O1-C1-C2-O2
2	A	501	HEM	CAA-CBA-CGA-O2A
2	B	501	HEM	CAD-CBD-CGD-O2D
2	C	501	HEM	CAA-CBA-CGA-O2A

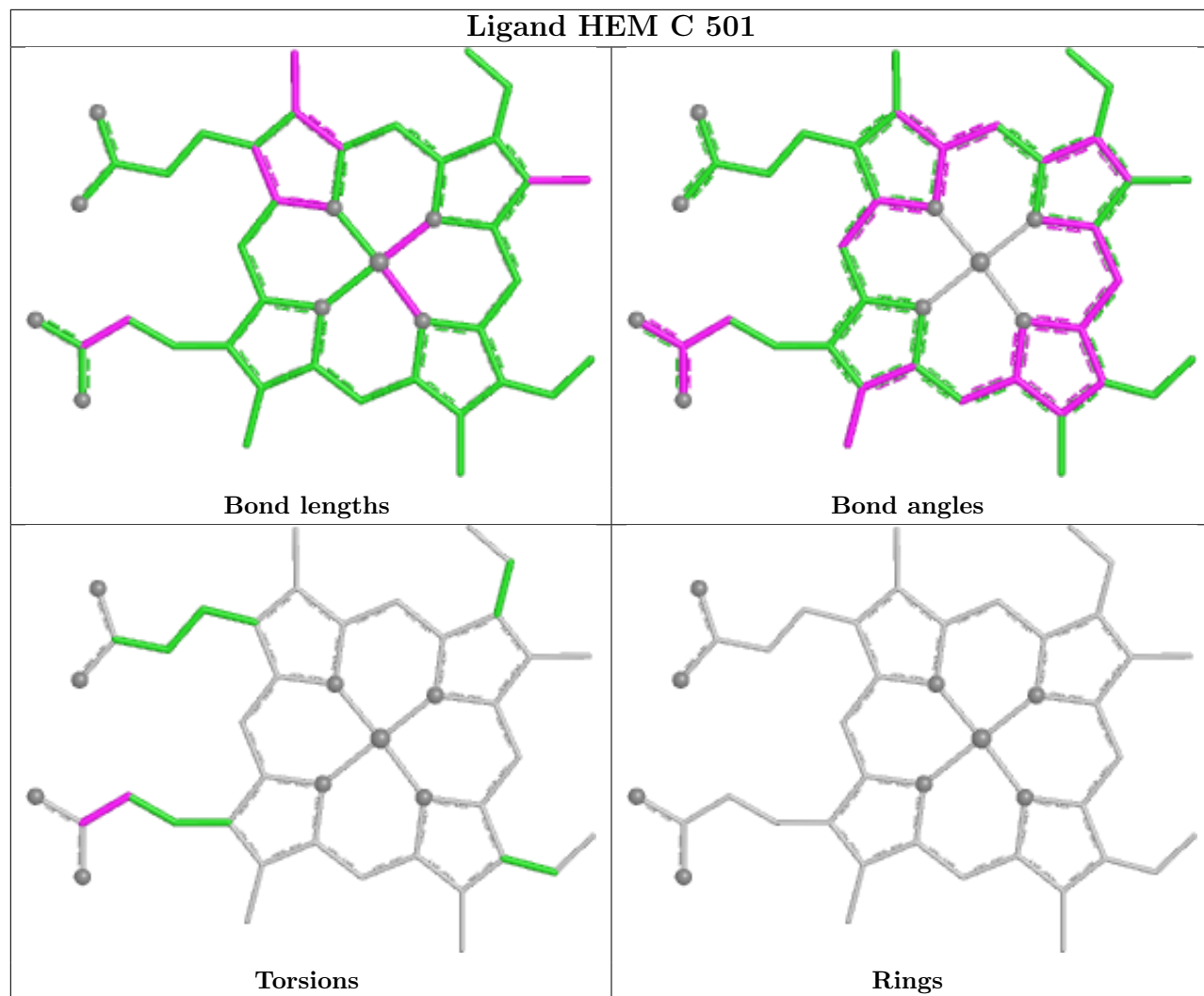
There are no ring outliers.

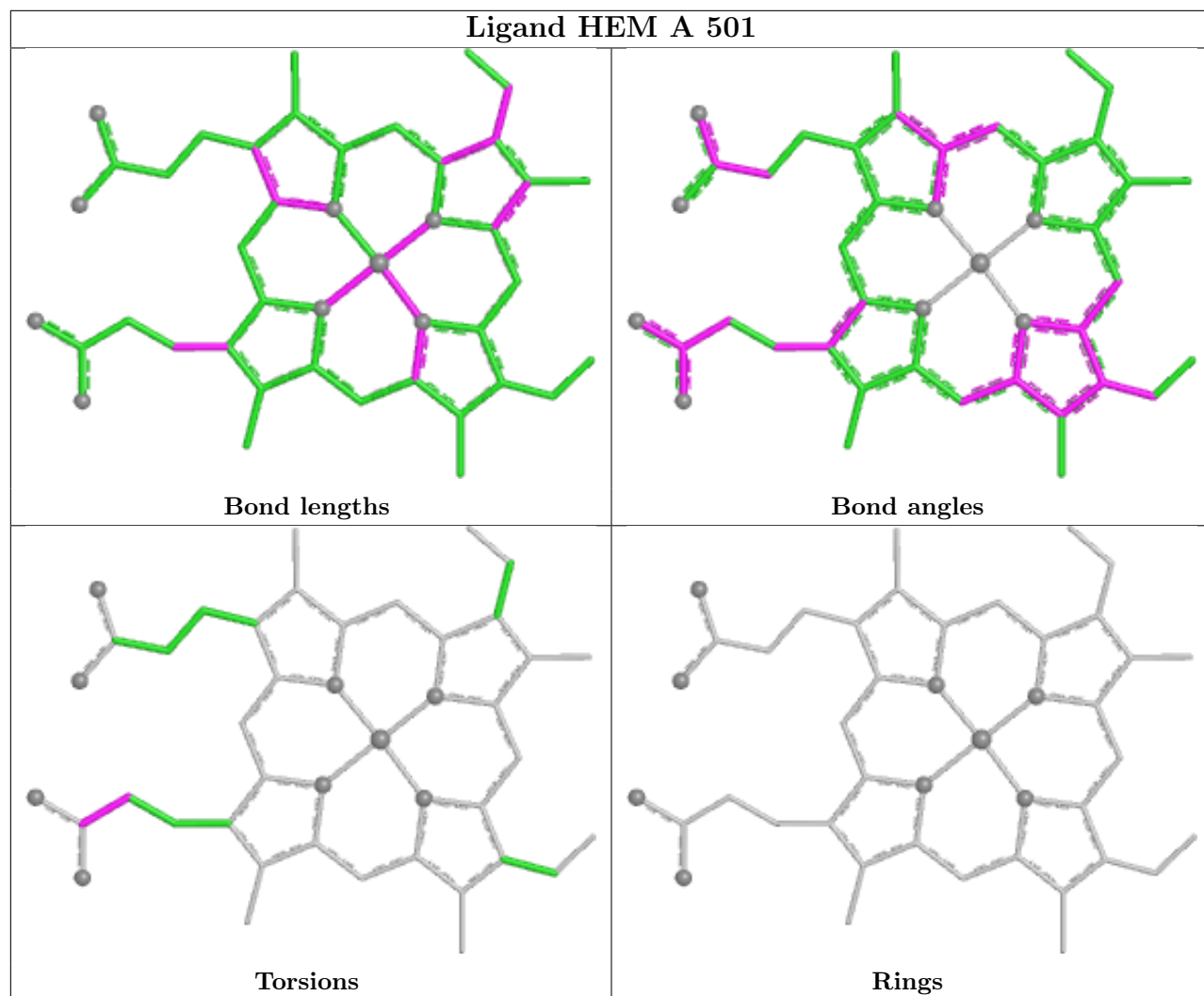
9 monomers are involved in 19 short contacts:

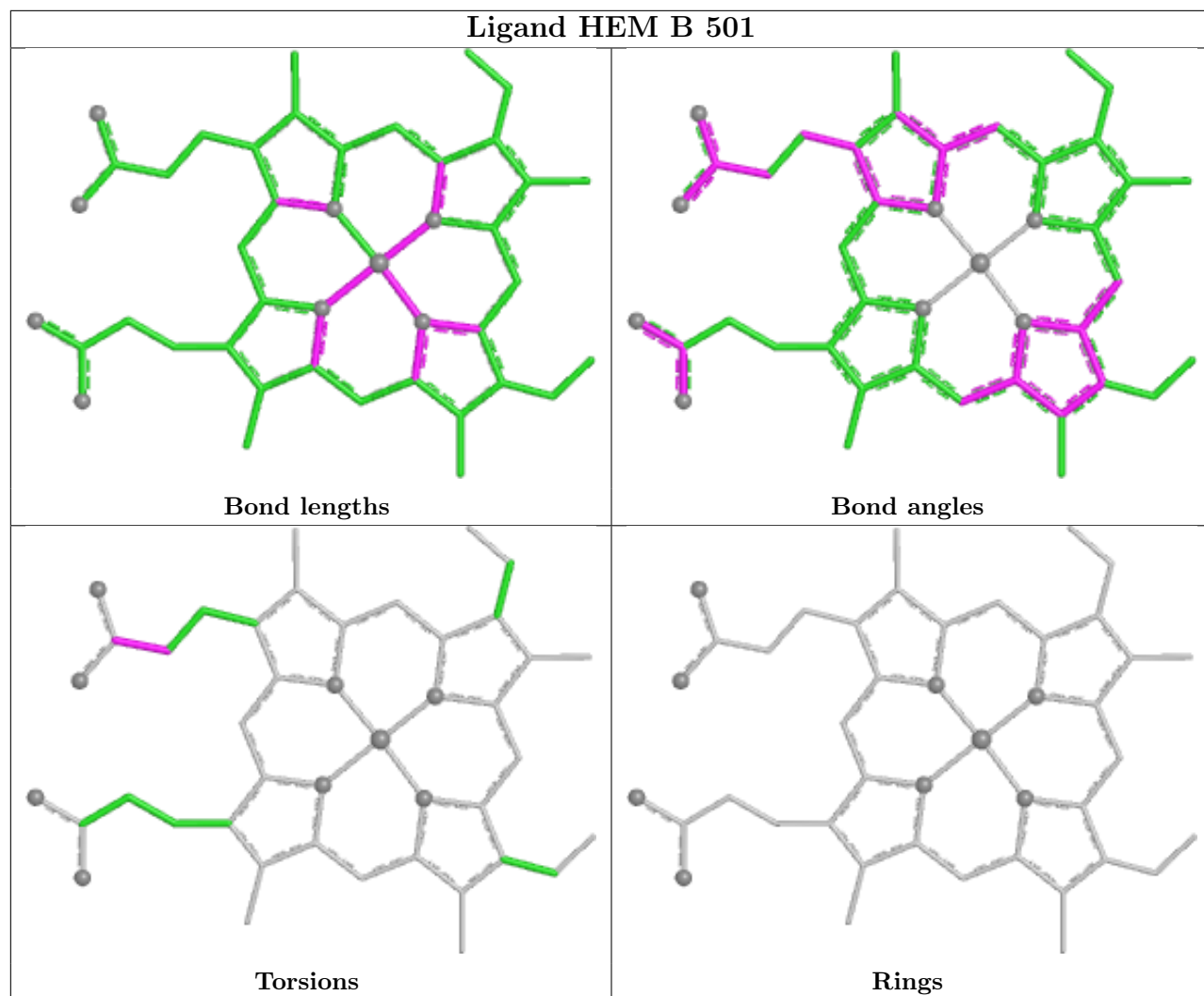
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	507	SO4	1	0
4	B	504	DMS	6	0
2	C	501	HEM	3	0
3	B	502	SO4	1	0
5	C	505	PEG	1	0
2	A	501	HEM	1	0
2	B	501	HEM	1	0
2	D	501	HEM	2	0
3	A	502	SO4	3	0

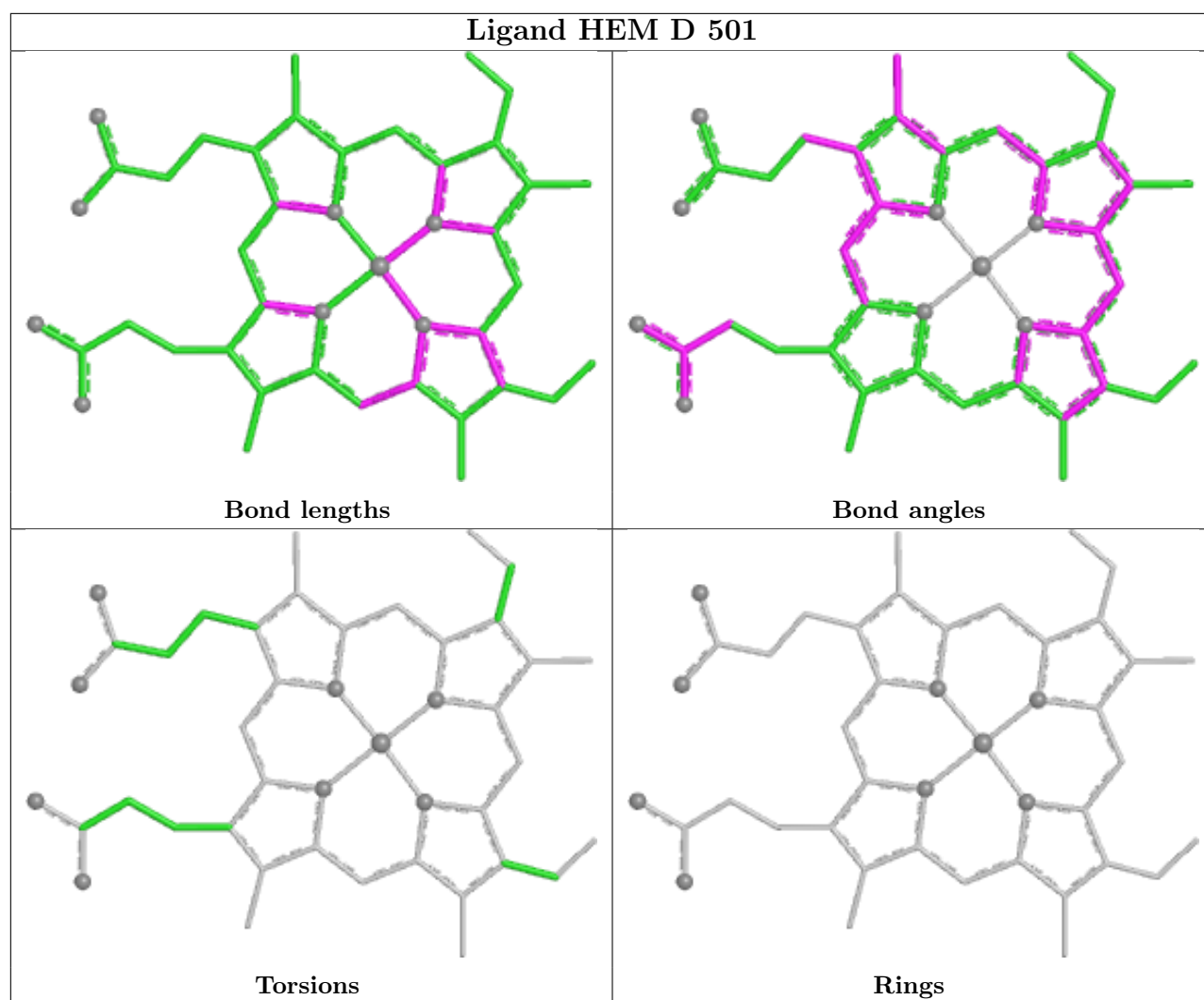
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 [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	A	381/416 (91%)	-1.15	0 100 100	8, 19, 52, 72	8 (2%)
1	B	383/416 (92%)	-1.17	0 100 100	8, 19, 50, 77	9 (2%)
1	C	397/416 (95%)	-1.09	0 100 100	10, 20, 52, 86	4 (1%)
1	D	397/416 (95%)	-1.08	0 100 100	10, 21, 52, 64	4 (1%)
All	All	1558/1664 (93%)	-1.12	0 100 100	8, 20, 52, 86	25 (1%)

There are no RSRZ outliers to report.

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(Å ²)	Q<0.9
3	SO4	C	506	5/5	0.96	0.06	57,63,64,67	0
5	PEG	C	505	7/7	0.96	0.07	49,51,52,53	0
3	SO4	A	507	5/5	0.97	0.05	65,67,71,72	0
3	SO4	B	509	5/5	0.97	0.06	56,61,63,64	0

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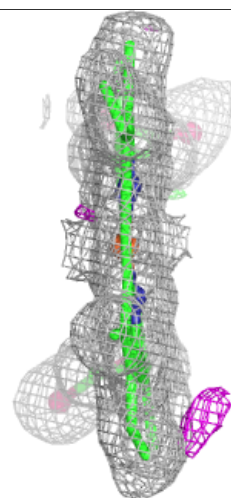
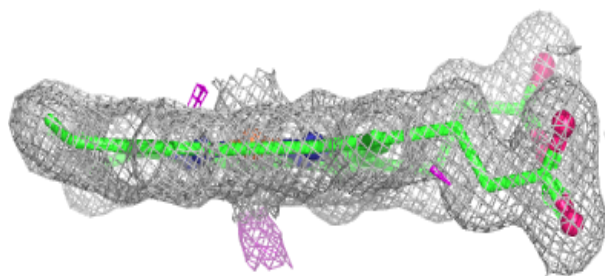
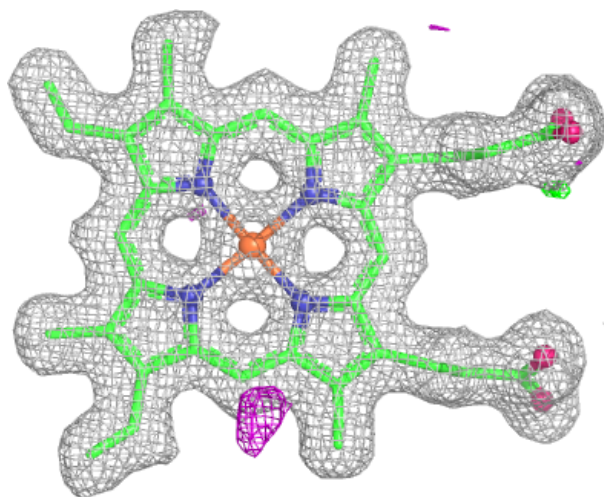
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	IMD	C	504	5/5	0.97	0.05	42,43,45,47	0
3	SO4	B	505	5/5	0.98	0.05	49,50,56,61	0
3	SO4	A	505	5/5	0.98	0.05	50,56,58,60	0
3	SO4	B	510	5/5	0.98	0.05	55,60,67,69	0
3	SO4	C	503	5/5	0.98	0.04	53,57,61,66	0
3	SO4	A	506	5/5	0.98	0.05	55,56,62,66	0
3	SO4	C	507	5/5	0.98	0.04	30,41,44,45	0
4	DMS	B	504	4/4	0.98	0.06	28,38,39,41	0
5	PEG	B	506	7/7	0.98	0.06	50,52,56,58	0
3	SO4	A	502	5/5	0.98	0.06	34,44,47,53	0
6	CO	B	507	1/1	0.98	0.07	58,58,58,58	0
6	CO	B	508	1/1	0.98	0.03	69,69,69,69	0
6	CO	D	504	1/1	0.98	0.04	71,71,71,71	0
3	SO4	B	502	5/5	0.98	0.04	43,45,57,58	0
8	PEO	C	508	2/2	0.98	0.04	36,36,36,36	0
3	SO4	A	504	5/5	0.99	0.04	31,50,56,58	0
3	SO4	B	503	5/5	0.99	0.05	32,55,58,60	0
3	SO4	D	502	5/5	0.99	0.04	37,38,39,41	0
3	SO4	D	503	5/5	0.99	0.03	43,44,47,52	0
3	SO4	C	502	5/5	0.99	0.03	34,39,40,40	0
3	SO4	A	503	5/5	0.99	0.03	49,54,57,58	0
2	HEM	B	501	43/43	1.00	0.02	9,10,14,19	0
2	HEM	C	501	43/43	1.00	0.02	12,14,17,19	0
2	HEM	D	501	43/43	1.00	0.02	12,13,16,22	0
2	HEM	A	501	43/43	1.00	0.02	10,12,15,21	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.

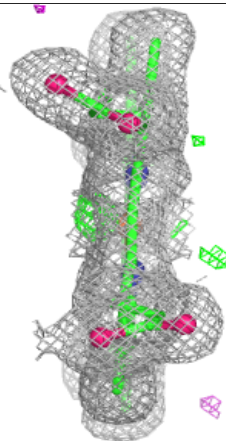
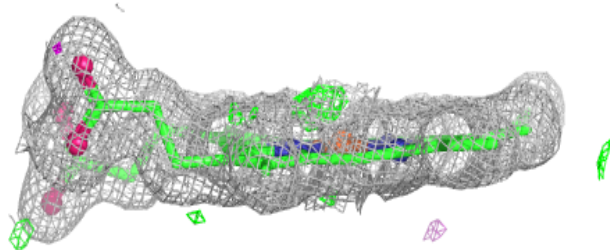
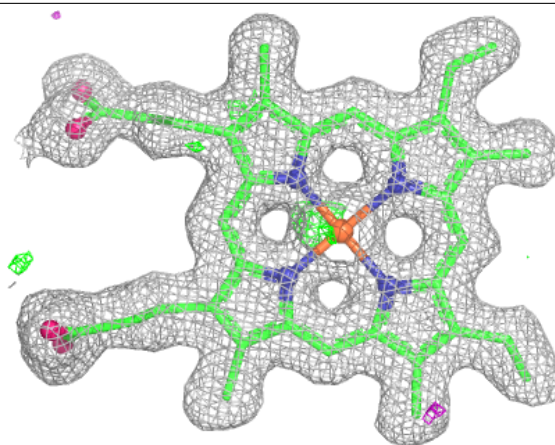
Electron density around HEM B 501:

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



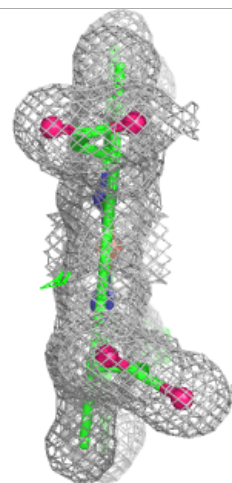
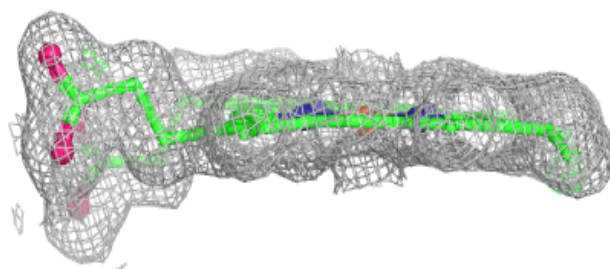
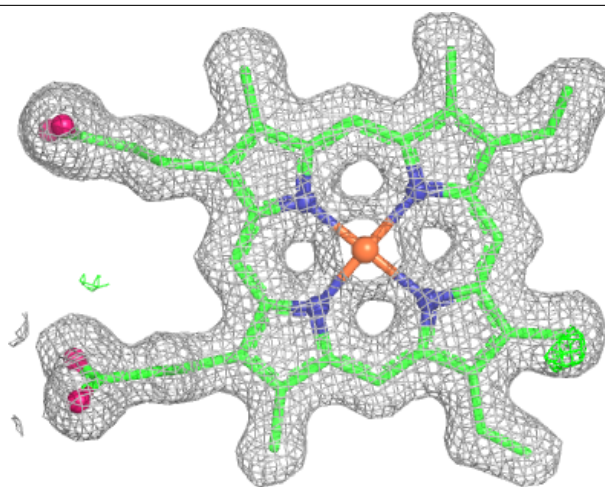
Electron density around HEM C 501:

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



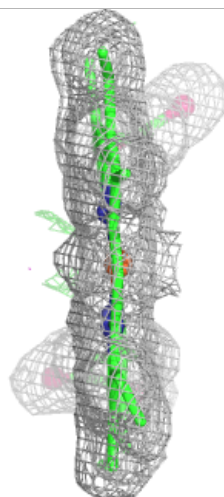
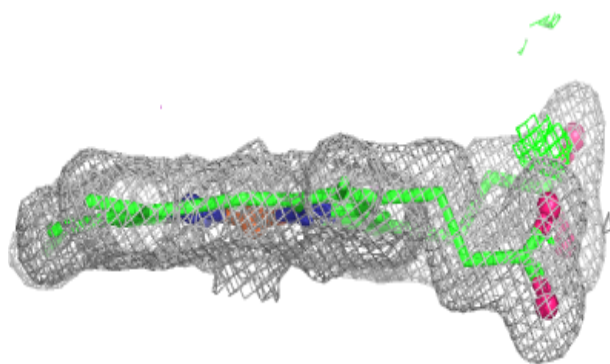
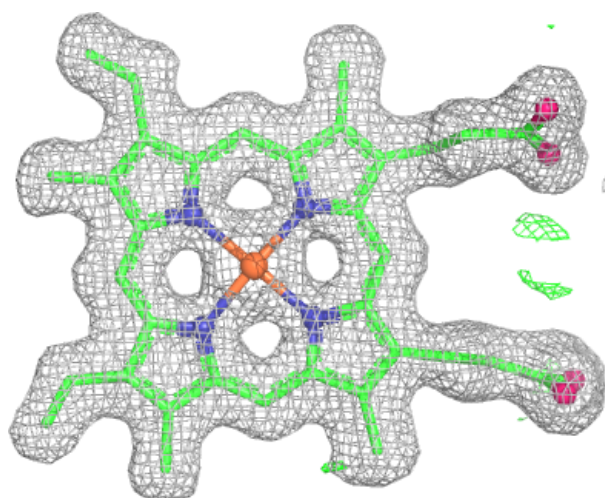
Electron density around HEM D 501:

$2mF_o-DF_c$ (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)



6.5 Other polymers [i](#)

There are no such residues in this entry.