



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:25 PM UTC

PDB ID : 4DUE / pdb_00004due
Title : cytochrome P450 BM3h-2G9C6 MRI sensor bound to serotonin
Authors : Brustad, E.M.; Lelyveld, V.S.; Snow, C.D.; Crook, N.; Martinez, F.M.; Scholl, T.J.; Jasanoff, A.; Arnold, F.H.
Deposited on : 2012-02-21
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

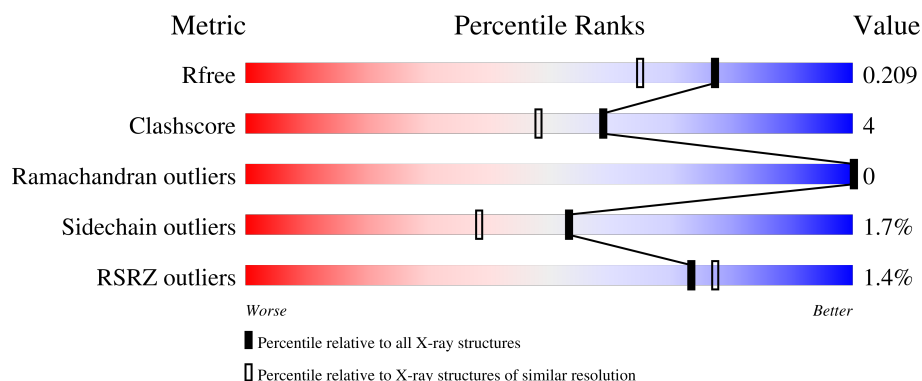
1 Overall quality at a glance

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	471	% 82% 13% ••
1	B	471	% 83% 12% •

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	SRO	A	501	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8229 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 BM3 variant 2G9C6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	453	Total	C	N	O	S	0	5	0
			3650	2333	623	674	20			
1	B	450	Total	C	N	O	S	0	5	0
			3637	2327	616	674	20			

There are 26 discrepancies between the modelled and reference sequences:

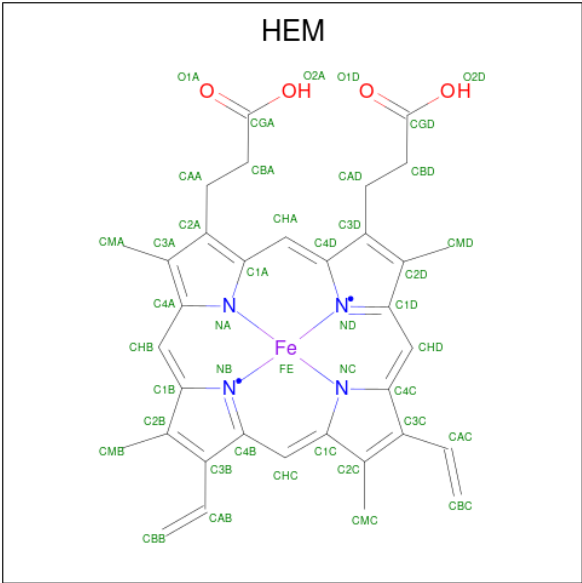
Chain	Residue	Modelled	Actual	Comment	Reference
A	50	CYS	ARG	engineered mutation	UNP P14779
A	87	LEU	PHE	engineered mutation	UNP P14779
A	268	SER	THR	engineered mutation	UNP P14779
A	437	GLN	LEU	engineered mutation	UNP P14779
A	438	LEU	THR	engineered mutation	UNP P14779
A	464	LEU	-	expression tag	UNP P14779
A	465	GLU	-	expression tag	UNP P14779
A	466	HIS	-	expression tag	UNP P14779
A	467	HIS	-	expression tag	UNP P14779
A	468	HIS	-	expression tag	UNP P14779
A	469	HIS	-	expression tag	UNP P14779
A	470	HIS	-	expression tag	UNP P14779
A	471	HIS	-	expression tag	UNP P14779
B	50	CYS	ARG	engineered mutation	UNP P14779
B	87	LEU	PHE	engineered mutation	UNP P14779
B	268	SER	THR	engineered mutation	UNP P14779
B	437	GLN	LEU	engineered mutation	UNP P14779
B	438	LEU	THR	engineered mutation	UNP P14779
B	464	LEU	-	expression tag	UNP P14779
B	465	GLU	-	expression tag	UNP P14779
B	466	HIS	-	expression tag	UNP P14779
B	467	HIS	-	expression tag	UNP P14779
B	468	HIS	-	expression tag	UNP P14779
B	469	HIS	-	expression tag	UNP P14779
B	470	HIS	-	expression tag	UNP P14779

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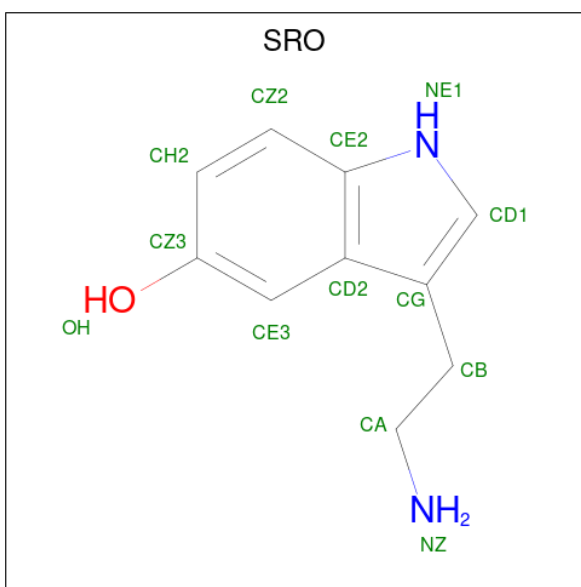
Chain	Residue	Modelled	Actual	Comment	Reference
B	471	HIS	-	expression tag	UNP P14779

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



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

- Molecule 3 is SEROTONIN (CCD ID: SRO) (formula: $C_{10}H_{12}N_2O$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			13	10	2	1		
3	B	1	Total	C	N	O	0	0
			13	10	2	1		

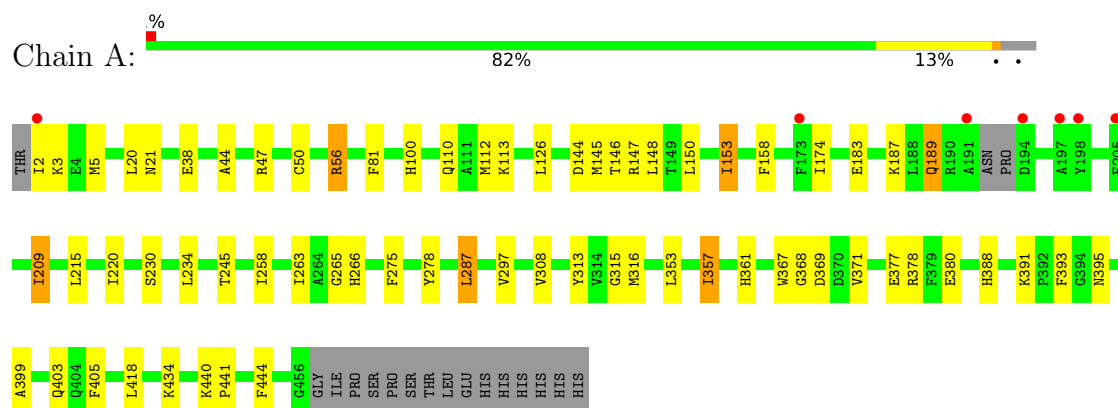
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	431	Total	O	0	0
			431	431		
4	B	399	Total	O	0	0
			399	399		

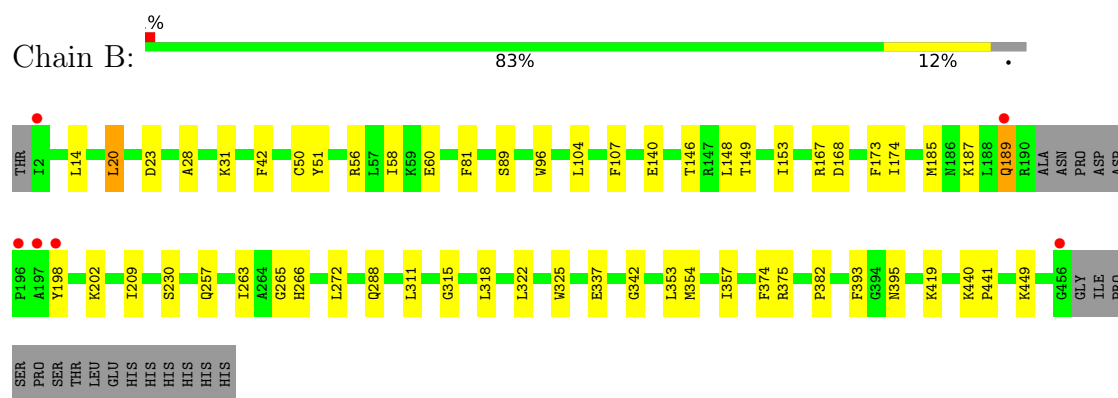
3 Residue-property plots

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 BM3 variant 2G9C6



- Molecule 1: cytochrome P450 BM3 variant 2G9C6



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.61Å 148.42Å 64.03Å 90.00° 98.88° 90.00°	Depositor
Resolution (Å)	38.41 – 1.70 38.41 – 1.70	Depositor EDS
% Data completeness (in resolution range)	96.4 (38.41-1.70) 96.4 (38.41-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.168 , 0.211 0.167 , 0.209	Depositor DCC
R_{free} test set	5899 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	24.3	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 41.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8229	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.34% 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: SRO, HEM

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	1.54	19/3744 (0.5%)	1.28	9/5058 (0.2%)
1	B	1.53	16/3733 (0.4%)	1.26	7/5041 (0.1%)
All	All	1.54	35/7477 (0.5%)	1.27	16/10099 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	357	ILE	CA-C	7.77	1.59	1.52
1	A	393	PHE	N-CA	7.49	1.55	1.46
1	B	441	PRO	C-O	7.29	1.31	1.23
1	A	357	ILE	CA-CB	-6.80	1.50	1.54
1	B	140	GLU	N-CA	-6.69	1.38	1.46
1	B	374	PHE	CA-C	6.67	1.61	1.52
1	A	357	ILE	CA-C	6.56	1.58	1.52
1	B	257	GLN	C-O	-6.48	1.16	1.24
1	B	318	LEU	CA-C	6.16	1.60	1.52
1	B	58	ILE	CA-CB	6.05	1.62	1.54
1	B	393	PHE	CA-CB	6.03	1.61	1.53
1	A	220	ILE	C-O	6.02	1.30	1.24
1	B	325	TRP	N-CA	6.00	1.53	1.46
1	A	113	LYS	N-CA	5.78	1.53	1.46
1	A	308	VAL	N-CA	5.78	1.53	1.46
1	A	215	LEU	CA-C	5.77	1.60	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	287	LEU	N-CA	5.70	1.53	1.46
1	A	174	ILE	CA-CB	-5.68	1.47	1.54
1	A	275	PHE	N-CA	5.58	1.53	1.46
1	A	245	THR	N-CA	5.54	1.53	1.46
1	B	28	ALA	CA-CB	5.39	1.61	1.53
1	A	378	ARG	CA-C	-5.28	1.45	1.52
1	B	104	LEU	CA-C	5.28	1.59	1.52
1	A	399	ALA	CA-CB	5.20	1.61	1.53
1	A	153	ILE	N-CA	5.19	1.52	1.46
1	B	311	LEU	CA-CB	5.19	1.59	1.52
1	A	297	VAL	CA-CB	5.18	1.62	1.54
1	A	315	GLY	N-CA	5.14	1.52	1.45
1	A	3	LYS	N-CA	5.13	1.51	1.45
1	B	107	PHE	CA-C	-5.10	1.46	1.52
1	A	395	ASN	CA-C	5.08	1.58	1.52
1	B	395	ASN	C-O	-5.07	1.17	1.23
1	B	315	GLY	N-CA	5.06	1.52	1.45
1	A	403	GLN	CA-C	5.01	1.59	1.52
1	B	96	TRP	CA-CB	5.00	1.61	1.53

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	189	GLN	N-CA-C	6.83	118.46	110.41
1	B	263	ILE	CB-CG1-CD1	-6.07	101.06	113.80
1	B	89	SER	CB-CA-C	6.02	119.67	109.50
1	A	380	GLU	N-CA-C	6.01	117.83	111.28
1	B	357	ILE	CB-CA-C	-5.63	108.39	114.35
1	A	209	ILE	N-CA-C	-5.62	105.04	110.72
1	B	419	LYS	N-CA-C	-5.45	104.98	111.69
1	A	357	ILE	CA-C-N	-5.43	113.31	119.28
1	A	357	ILE	C-N-CA	-5.43	113.31	119.28
1	A	377	GLU	N-CA-C	-5.42	106.51	113.01
1	A	263	ILE	CB-CG1-CD1	-5.21	102.85	113.80
1	A	378	ARG	CB-CA-C	-5.11	100.54	110.46
1	A	100	HIS	N-CA-C	-5.05	105.67	111.07
1	A	189	GLN	N-CA-C	5.03	119.50	112.90
1	B	375	ARG	N-CA-CB	-5.03	104.81	111.40
1	B	14	LEU	N-CA-C	-5.03	107.14	113.28

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	441	PRO	Mainchain

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	3650	0	3623	33	0
1	B	3637	0	3620	28	0
2	A	43	0	30	2	0
2	B	43	0	30	3	0
3	A	13	0	11	0	0
3	B	13	0	10	0	0
4	A	431	0	0	5	0
4	B	399	0	0	7	0
All	All	8229	0	7324	64	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 4.

All (64) 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:313:TYR:HA	1:A:316:MET:HE3	1.55	0.88
1:B:23:ASP:OD2	4:B:976:HOH:O	1.96	0.82
1:B:173:PHE:HD1	1:B:174:ILE:HD12	1.52	0.74
1:A:21:ASN:HA	1:A:189:GLN:HG3	1.77	0.67
1:A:434:LYS:HE3	4:A:674:HOH:O	1.95	0.67
1:A:50[A]:CYS:SG	1:A:353:LEU:HD23	2.38	0.64
1:A:110:GLN:NE2	4:A:908:HOH:O	2.30	0.64
1:B:187:LYS:C	1:B:189:GLN:H	2.06	0.63
2:B:500:HEM:HMC2	2:B:500:HEM:HBC2	1.82	0.62
1:A:146:THR:HG23	1:A:266:HIS:CE1	2.36	0.60
1:A:367:TRP:HB2	1:A:371:VAL:HG12	1.84	0.59
1:A:150:LEU:HB2	1:A:266:HIS:CE1	2.38	0.59
1:B:51:TYR:CE1	1:B:354[B]:MET:HG2	2.38	0.58
1:A:81:PHE:HB3	1:A:209:ILE:HG12	1.87	0.56
1:B:56[A]:ARG:NH2	4:B:979:HOH:O	2.31	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:81:PHE:HB3	1:B:209:ILE:HG12	1.88	0.54
1:B:266:HIS:HE1	4:B:833:HOH:O	1.90	0.54
1:A:50[A]:CYS:SG	1:A:353:LEU:CD2	2.96	0.54
1:A:368:GLY:O	1:A:371:VAL:HG13	2.09	0.53
1:A:146:THR:HG23	1:A:266:HIS:HE1	1.74	0.51
1:B:337[B]:GLU:CD	4:B:916:HOH:O	2.52	0.51
1:B:20:LEU:HD22	1:B:42:PHE:CZ	2.46	0.51
1:A:126:LEU:HD21	1:A:145:MET:HE1	1.92	0.51
1:B:146:THR:HG23	1:B:266:HIS:HE2	1.77	0.50
1:A:316:MET:HE1	4:A:1014:HOH:O	2.12	0.49
1:B:174:ILE:HD12	1:B:174:ILE:N	2.27	0.49
1:B:265:GLY:HA2	2:B:500:HEM:C2C	2.48	0.48
2:A:500:HEM:HMC2	2:A:500:HEM:HBC2	1.94	0.48
1:B:440:LYS:HG2	4:B:654:HOH:O	2.14	0.47
1:A:440:LYS:HD3	4:A:976:HOH:O	2.14	0.47
1:B:288:GLN:HG3	4:B:899:HOH:O	2.14	0.47
1:A:313:TYR:HA	1:A:316:MET:CE	2.36	0.47
1:A:388:HIS:HA	1:A:391:LYS:HE3	1.98	0.46
1:B:174:ILE:N	1:B:174:ILE:CD1	2.79	0.46
2:B:500:HEM:HBC2	2:B:500:HEM:CMC	2.45	0.46
1:B:187:LYS:O	1:B:189:GLN:N	2.49	0.45
1:A:287:LEU:CD1	1:A:418:LEU:CD2	2.94	0.45
1:B:50[A]:CYS:SG	1:B:353:LEU:HD23	2.57	0.45
1:B:173:PHE:CD1	1:B:174:ILE:HD12	2.41	0.45
1:B:50[A]:CYS:SG	1:B:353:LEU:CD2	3.05	0.45
1:B:153:ILE:HD12	1:B:153:ILE:HA	1.91	0.44
1:A:265:GLY:HA2	2:A:500:HEM:C2C	2.53	0.44
1:A:5:MET:HE3	1:A:5:MET:HB2	1.56	0.44
1:B:272:LEU:HD13	1:B:322:LEU:HG	1.98	0.44
1:B:187:LYS:C	1:B:189:GLN:N	2.70	0.44
1:A:153:ILE:HD12	1:A:153:ILE:HA	1.71	0.43
1:B:60:GLU:OE2	1:B:342:GLY:HA2	2.18	0.43
1:A:278:TYR:HA	1:A:444:PHE:CZ	2.54	0.43
1:A:234:LEU:HD13	1:A:258:ILE:HD11	2.00	0.43
1:B:198:TYR:O	1:B:202:LYS:HG3	2.19	0.43
1:B:354[A]:MET:HA	1:B:354[A]:MET:HE3	2.01	0.43
1:A:144:ASP:OD1	1:A:147:ARG:NH1	2.52	0.42
1:A:20:LEU:HB2	1:A:189:GLN:NE2	2.34	0.42
1:A:110:GLN:NE2	1:A:110:GLN:HA	2.35	0.42
1:A:112:MET:SD	1:A:405:PHE:HA	2.60	0.42
1:A:44:ALA:HB3	1:A:47:ARG:HG2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:434:LYS:CE	4:A:674:HOH:O	2.62	0.41
1:B:382:PRO:HD2	4:B:638:HOH:O	2.21	0.41
1:A:183:GLU:O	1:A:187:LYS:HG3	2.21	0.41
1:B:149:THR:OG1	1:B:266:HIS:HD2	2.04	0.41
1:A:20:LEU:HB2	1:A:189:GLN:HE21	1.87	0.40
1:B:148:LEU:HD23	1:B:148:LEU:C	2.46	0.40
1:A:357:ILE:HG22	1:A:361:HIS:CE1	2.56	0.40
1:A:38:GLU:OE2	1:A:56[B]:ARG:NH1	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	454/471 (96%)	442 (97%)	12 (3%)	0	100	100
1	B	451/471 (96%)	437 (97%)	14 (3%)	0	100	100
All	All	905/942 (96%)	879 (97%)	26 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	394/413 (95%)	387 (98%)	7 (2%)	51	36

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	395/413 (96%)	388 (98%)	7 (2%)	51	36
All	All	789/826 (96%)	775 (98%)	14 (2%)	53	36

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ILE
1	A	56[A]	ARG
1	A	56[B]	ARG
1	A	148	LEU
1	A	158	PHE
1	A	230	SER
1	A	369	ASP
1	B	20	LEU
1	B	31	LYS
1	B	167	ARG
1	B	168	ASP
1	B	185	MET
1	B	230	SER
1	B	449	LYS

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	70	ASN
1	A	229	GLN
1	A	239	ASN
1	A	266	HIS
1	A	404	GLN
1	B	70	ASN
1	B	109	GLN
1	B	310	GLN
1	B	359	GLN

5.3.3 RNA ⓘ

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](#)

4 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	SRO	A	501	2	14,14,14	2.03	6 (42%)	19,19,19	2.34	10 (52%)
3	SRO	B	501	2	14,14,14	1.63	4 (28%)	19,19,19	2.30	8 (42%)
2	HEM	B	500	3,1	50,50,50	1.84	12 (24%)	67,82,82	1.88	19 (28%)
2	HEM	A	500	3,1	50,50,50	2.05	16 (32%)	67,82,82	2.42	21 (31%)

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	SRO	A	501	2	-	1/3/3/3	0/2/2/2
3	SRO	B	501	2	-	1/3/3/3	0/2/2/2
2	HEM	B	500	3,1	-	2/14/54/54	-
2	HEM	A	500	3,1	-	2/14/54/54	-

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	HEM	C3D-C2D	5.78	1.49	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	HEM	FE-NA	5.42	2.13	1.95
2	A	500	HEM	FE-NC	4.89	2.11	1.95
3	A	501	SRO	CD2-CE2	4.17	1.46	1.41
2	A	500	HEM	FE-NA	4.06	2.08	1.95
2	B	500	HEM	C3D-C2D	3.72	1.44	1.36
2	A	500	HEM	CMD-C2D	3.64	1.58	1.50
2	A	500	HEM	FE-NB	3.55	2.05	1.94
2	A	500	HEM	CAB-C3B	3.45	1.56	1.47
2	B	500	HEM	FE-NB	3.39	2.05	1.94
2	A	500	HEM	CMA-C3A	3.24	1.57	1.50
2	B	500	HEM	FE-ND	3.15	2.04	1.94
3	A	501	SRO	CD2-CG	3.07	1.49	1.44
2	B	500	HEM	CAB-C3B	3.07	1.55	1.47
3	B	501	SRO	CD2-CE2	2.97	1.45	1.41
3	B	501	SRO	CE3-CD2	2.88	1.44	1.39
2	A	500	HEM	CMB-C2B	2.87	1.56	1.50
2	B	500	HEM	O2D-CGD	-2.80	1.21	1.30
2	A	500	HEM	O1D-CGD	2.76	1.31	1.22
2	B	500	HEM	CAC-C3C	2.75	1.54	1.47
2	B	500	HEM	C4C-NC	-2.74	1.34	1.39
2	B	500	HEM	FE-NC	2.73	2.04	1.95
3	A	501	SRO	CZ2-CH2	2.67	1.43	1.38
2	A	500	HEM	FE-ND	2.67	2.03	1.94
3	B	501	SRO	CD1-CG	2.62	1.40	1.36
2	A	500	HEM	C3B-C2B	-2.59	1.31	1.37
2	A	500	HEM	C2A-C3A	-2.58	1.32	1.38
2	B	500	HEM	CMC-C2C	2.49	1.55	1.50
2	A	500	HEM	CAD-C3D	2.46	1.57	1.51
3	B	501	SRO	CD2-CG	2.45	1.48	1.44
2	B	500	HEM	CMA-C3A	2.43	1.55	1.50
2	B	500	HEM	O1A-CGA	2.37	1.29	1.22
2	A	500	HEM	CAC-C3C	2.28	1.53	1.47
3	A	501	SRO	CE3-CZ3	2.19	1.42	1.39
3	A	501	SRO	CB-CG	2.11	1.54	1.50
2	A	500	HEM	O2D-CGD	-2.09	1.23	1.30
2	A	500	HEM	CBD-CGD	2.02	1.55	1.50
3	A	501	SRO	CD1-NE1	-2.00	1.34	1.37

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	HEM	C3B-C4B-NB	-6.48	104.81	109.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	HEM	CHC-C4B-NB	5.96	130.84	124.42
2	A	500	HEM	CHD-C4C-NC	5.75	130.72	124.45
2	A	500	HEM	C1B-NB-C4B	5.70	111.96	105.21
2	A	500	HEM	CHA-C4D-ND	5.65	131.36	124.37
2	A	500	HEM	C4D-ND-C1D	5.50	111.72	105.21
2	B	500	HEM	C4D-ND-C1D	4.58	110.63	105.21
2	B	500	HEM	CHD-C1D-ND	4.43	129.19	124.42
2	A	500	HEM	CBD-CAD-C3D	-4.32	100.58	112.53
2	B	500	HEM	CBD-CAD-C3D	-4.31	100.63	112.53
2	A	500	HEM	CHD-C1D-ND	4.26	129.00	124.42
3	B	501	SRO	CZ2-CE2-CD2	-4.18	118.14	122.19
3	A	501	SRO	CH2-CZ3-CE3	4.03	124.61	120.19
2	A	500	HEM	C2B-C1B-NB	-3.90	105.35	109.84
3	B	501	SRO	CB-CG-CD1	3.72	134.66	127.18
3	B	501	SRO	CD2-CE3-CZ3	-3.67	115.14	120.11
2	A	500	HEM	C3B-C2B-C1B	3.56	109.09	106.41
3	A	501	SRO	CB-CG-CD1	3.56	134.34	127.18
3	B	501	SRO	CB-CG-CD2	-3.53	119.39	126.18
2	B	500	HEM	C1B-NB-C4B	3.49	109.34	105.21
2	B	500	HEM	C3B-C2B-C1B	3.43	108.99	106.41
3	A	501	SRO	CD2-CG-CD1	-3.40	102.68	106.16
2	B	500	HEM	CHB-C1B-NB	3.39	128.56	124.37
2	B	500	HEM	CMA-C3A-C4A	-3.36	120.30	125.42
2	A	500	HEM	C2A-C1A-NA	-3.24	106.56	110.15
3	A	501	SRO	CE2-NE1-CD1	3.14	111.87	109.08
2	B	500	HEM	C2B-C1B-NB	-3.12	106.25	109.84
3	B	501	SRO	CE2-NE1-CD1	3.12	111.85	109.08
3	A	501	SRO	CZ2-CE2-CD2	-3.09	119.19	122.19
3	A	501	SRO	CD2-CE3-CZ3	-3.06	115.96	120.11
2	A	500	HEM	CMB-C2B-C1B	-3.05	120.27	125.03
2	B	500	HEM	CBA-CAA-C2A	-3.03	104.15	112.53
2	B	500	HEM	CMA-C3A-C2A	2.99	131.96	125.62
2	A	500	HEM	C2D-C1D-ND	-2.94	106.50	109.90
3	B	501	SRO	CH2-CZ3-CE3	2.93	123.41	120.19
2	B	500	HEM	C3B-C4B-NB	-2.88	107.40	109.47
3	B	501	SRO	CE3-CD2-CE2	2.87	122.19	119.39
2	A	500	HEM	C1A-C2A-C3A	2.87	111.31	106.87
2	B	500	HEM	C2D-C1D-ND	-2.86	106.59	109.90
3	A	501	SRO	CZ2-CE2-NE1	2.80	135.23	130.74
2	B	500	HEM	CHA-C4D-ND	2.72	127.74	124.37
2	B	500	HEM	O1A-CGA-CBA	-2.55	114.99	123.09
2	B	500	HEM	CHC-C4B-NB	2.49	127.10	124.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	HEM	C4C-NC-C1C	2.47	109.85	105.82
2	B	500	HEM	CHC-C1C-NC	2.44	127.11	124.45
2	A	500	HEM	C3D-C4D-ND	-2.44	107.50	110.17
2	A	500	HEM	CMA-C3A-C2A	2.37	130.65	125.62
3	A	501	SRO	CD2-CE2-NE1	-2.36	105.55	107.52
3	A	501	SRO	CE3-CD2-CE2	2.26	121.59	119.39
2	A	500	HEM	CHA-C4D-C3D	-2.26	121.06	125.23
2	B	500	HEM	C2A-C1A-NA	-2.21	107.70	110.15
3	B	501	SRO	CZ2-CE2-NE1	2.19	134.25	130.74
2	A	500	HEM	CHB-C1B-NB	2.17	127.05	124.37
2	B	500	HEM	CAD-C3D-C4D	2.10	128.36	124.70
2	A	500	HEM	C4C-CHD-C1D	-2.09	121.58	126.02
2	B	500	HEM	CHA-C1A-NA	2.07	127.62	123.86
3	A	501	SRO	CG-CD1-NE1	2.06	112.56	110.31
2	A	500	HEM	CAD-CBD-CGD	-2.03	108.27	113.67

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	501	SRO	NZ-CA-CB-CG
3	B	501	SRO	NZ-CA-CB-CG
2	B	500	HEM	CAD-CBD-CGD-O2D
2	B	500	HEM	CAD-CBD-CGD-O1D
2	A	500	HEM	CAD-CBD-CGD-O1D
2	A	500	HEM	CAD-CBD-CGD-O2D

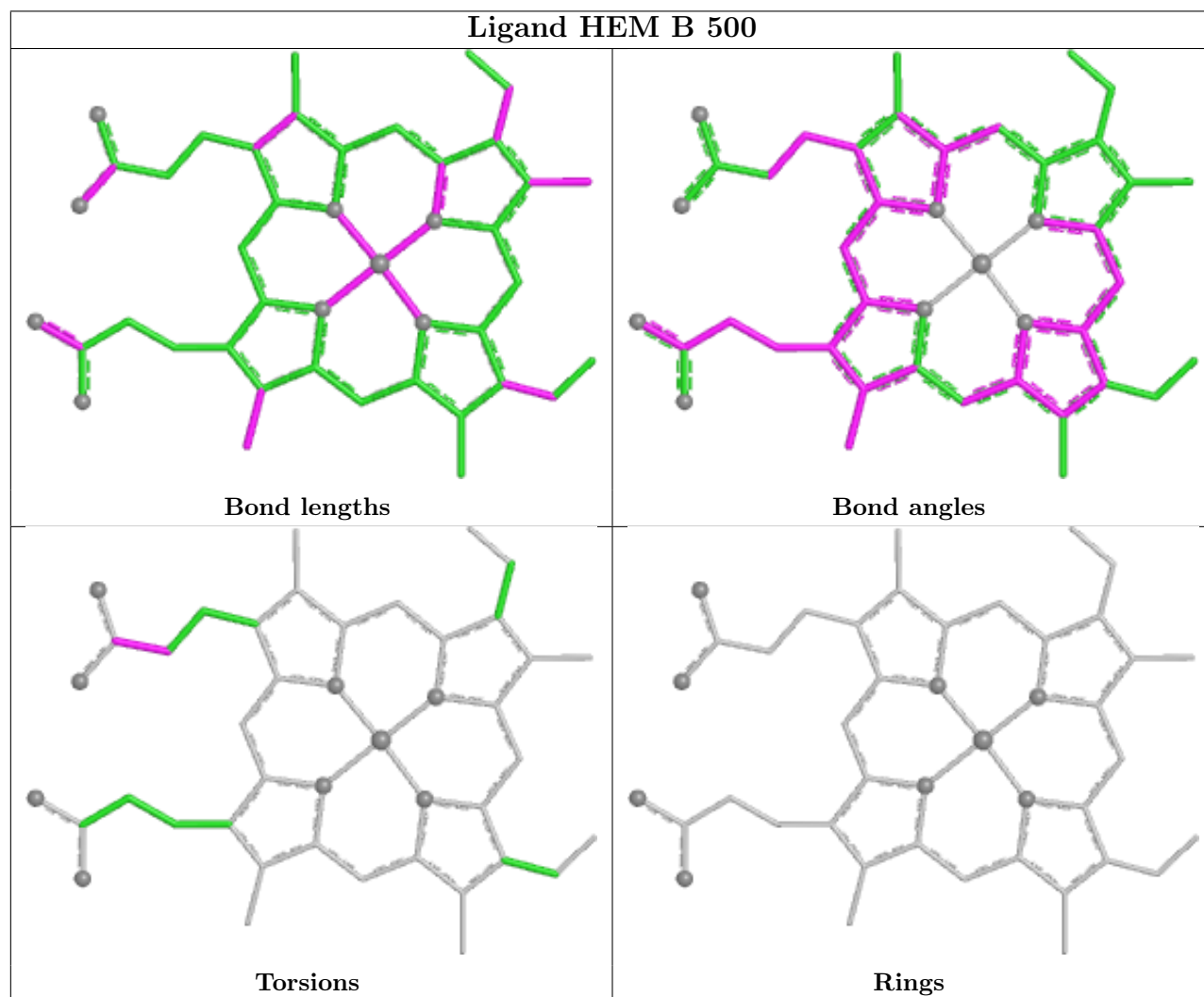
There are no ring outliers.

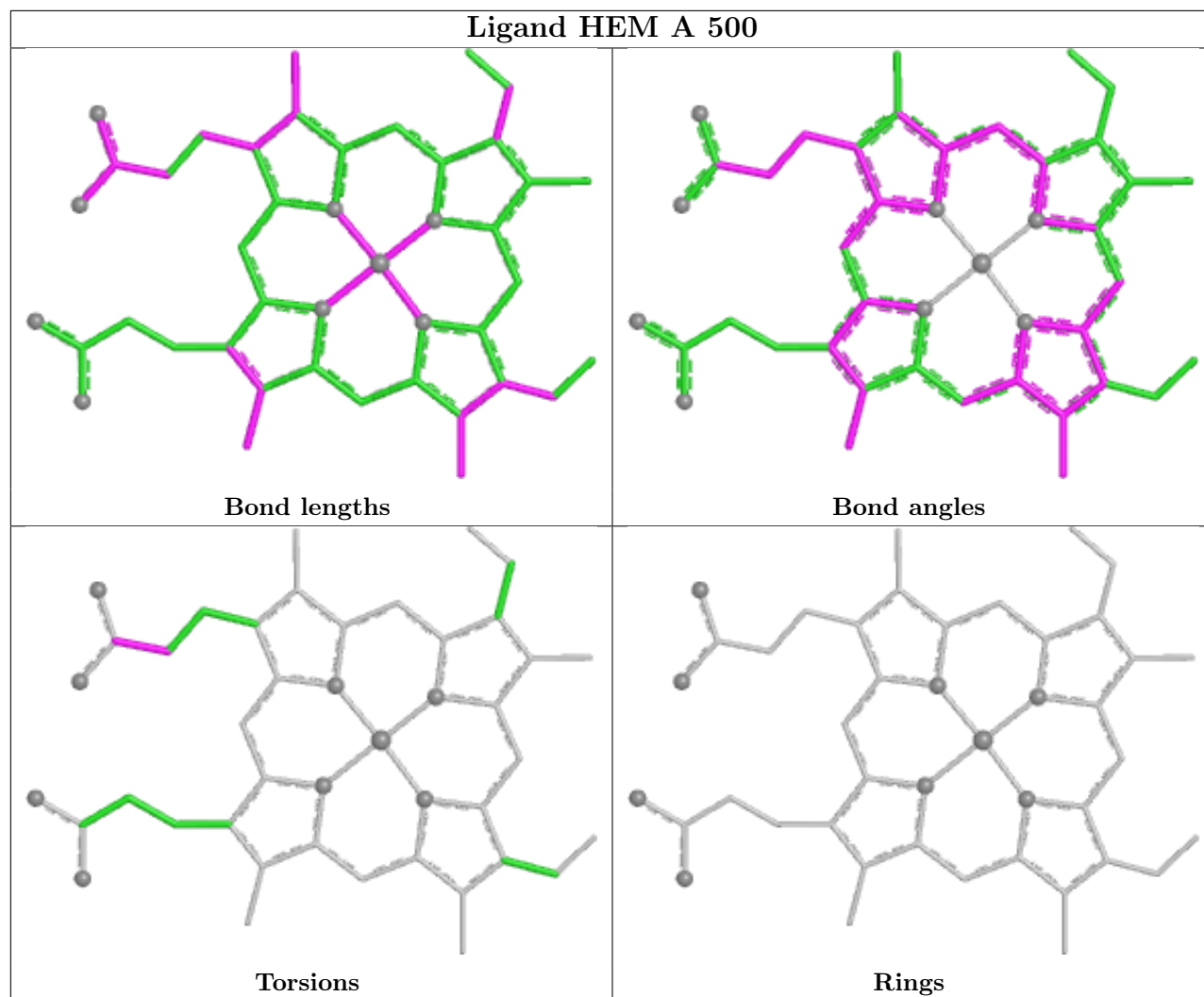
2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	500	HEM	3	0
2	A	500	HEM	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 [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	453/471 (96%)	-0.22	7 (1%) 72 76	9, 23, 46, 64	5 (1%)
1	B	450/471 (95%)	-0.22	6 (1%) 75 79	10, 24, 44, 59	5 (1%)
All	All	903/942 (95%)	-0.22	13 (1%) 73 77	9, 23, 45, 64	10 (1%)

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	2	ILE	5.5
1	B	196	PRO	4.3
1	A	194	ASP	4.1
1	A	2	ILE	3.9
1	A	191	ALA	3.7
1	A	197	ALA	3.3
1	A	198	TYR	3.0
1	B	198	TYR	2.8
1	B	197	ALA	2.7
1	A	173	PHE	2.3
1	B	456	GLY	2.2
1	B	189	GLN	2.2
1	A	205	PHE	2.1

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](#)

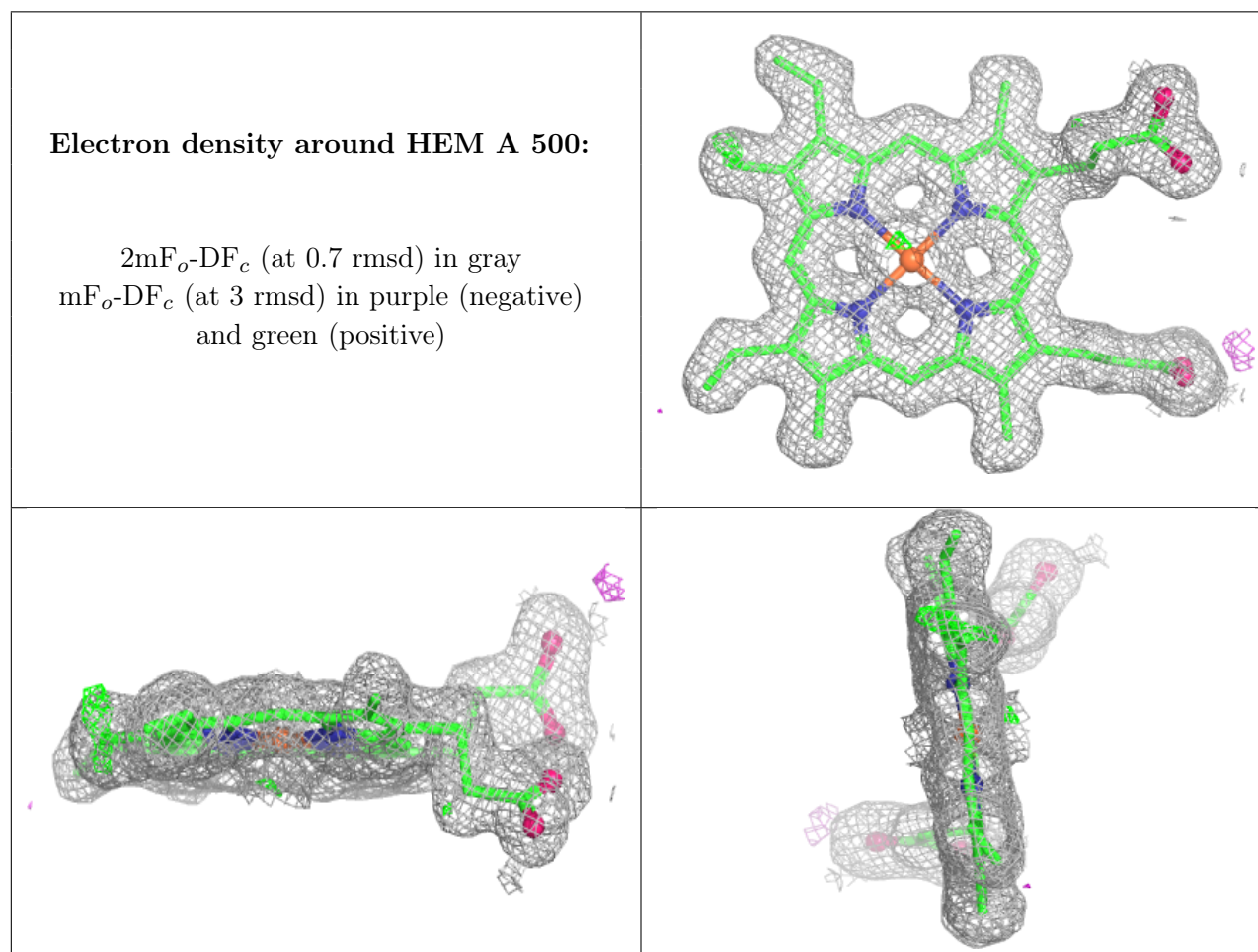
There are no oligosaccharides in this entry.

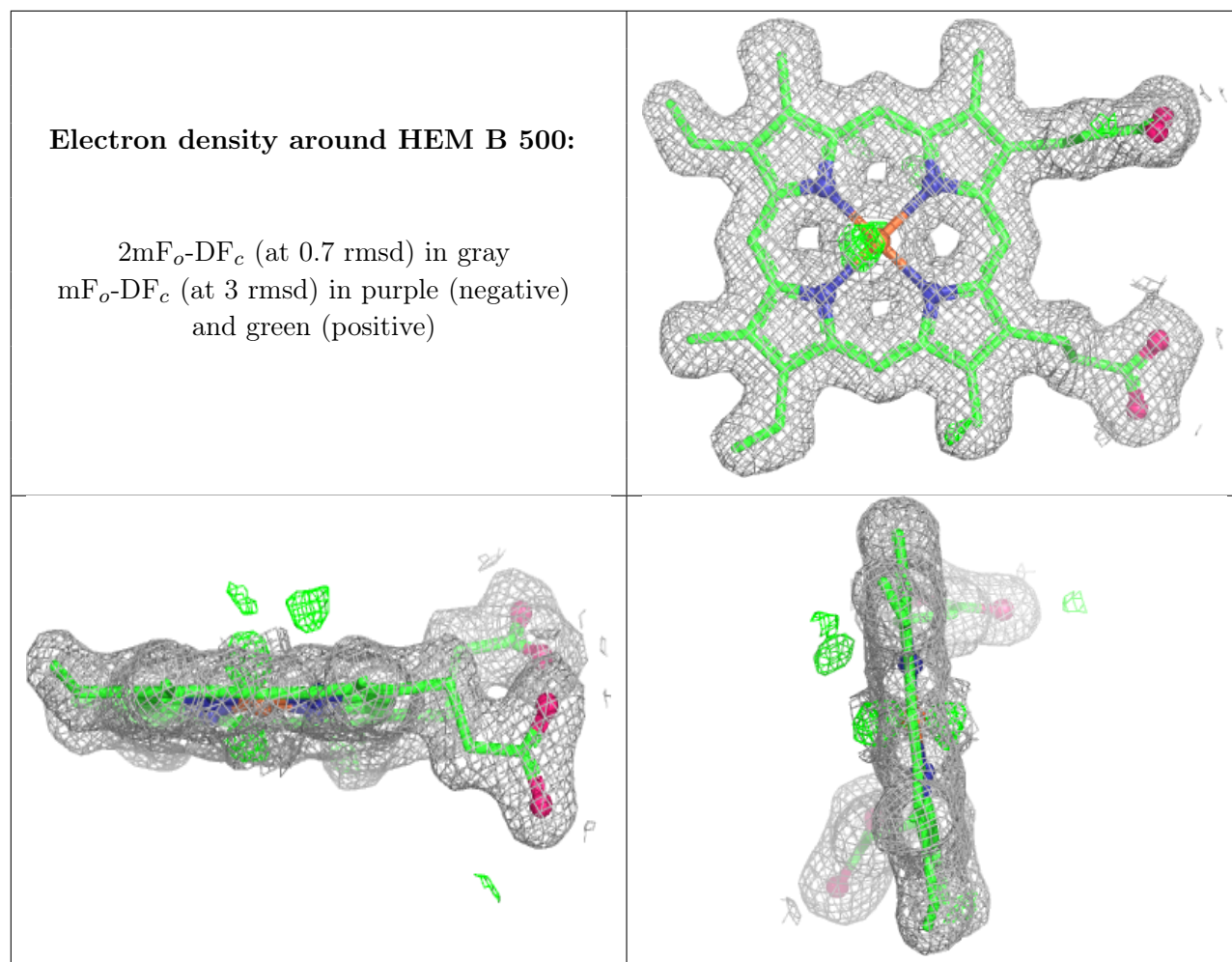
6.4 Ligands [i](#)

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
3	SRO	A	501	13/13	0.92	0.09	17,26,30,31	0
3	SRO	B	501	13/13	0.95	0.07	20,25,29,30	0
2	HEM	A	500	43/43	0.99	0.04	10,15,18,27	0
2	HEM	B	500	43/43	0.99	0.04	10,14,19,27	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.