



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:33 AM UTC

PDB ID : 4HGF / pdb_00004hgf
Title : Crystal structure of P450 BM3 5F5K heme domain variant complexed with styrene
Authors : Shehzad, A.; Panneerselvam, S.; Bocola, M.; Mueller-Dieckmann, J.; Wilmanns, M.; Schwaneberg, U.
Deposited on : 2012-10-08
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

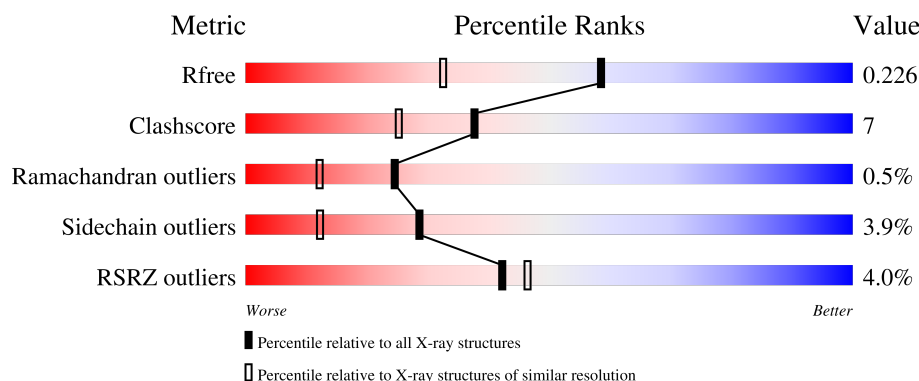
1 Overall quality at a glance

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	455	4% 77% 16% ...
1	B	455	3% 79% 16% ..

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	SYN	B	502	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 8178 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.

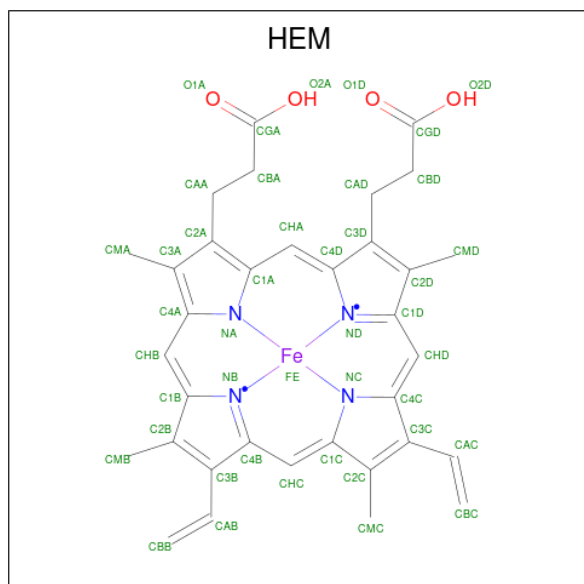
- Molecule 1 is a protein called Bifunctional P-450/NADPH-P450 reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	442	Total	C	N	O	S	0	9	0
			3607	2314	608	668	17			
1	B	441	Total	C	N	O	S	0	8	0
			3599	2303	609	669	18			

There are 6 discrepancies between the modelled and reference sequences:

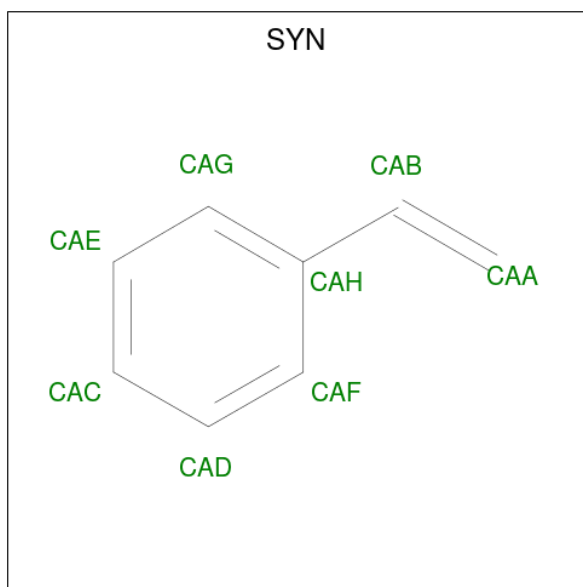
Chain	Residue	Modelled	Actual	Comment	Reference
A	87	ALA	PHE	engineered mutation	UNP P14779
A	184	LYS	ALA	engineered mutation	UNP P14779
A	235	ALA	THR	engineered mutation	UNP P14779
B	87	ALA	PHE	engineered mutation	UNP P14779
B	184	LYS	ALA	engineered mutation	UNP P14779
B	235	ALA	THR	engineered mutation	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	C	Fe	N	O	
			43	34	1	4	4	
2	B	1	Total	C	Fe	N	O	
			43	34	1	4	4	

- Molecule 3 is ethenylbenzene (CCD ID: SYN) (formula: C₈H₈).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	C		
			8	8	0	0
3	B	1	Total	C		
			8	8	0	0

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl		
			1	1	0	0

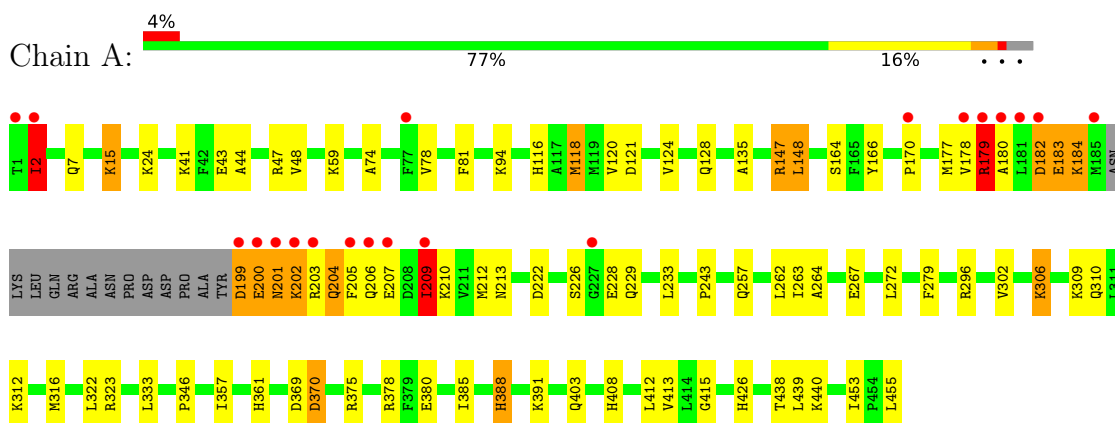
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	423	Total	O		
			423	423	0	0
5	B	446	Total	O		
			446	446	0	0

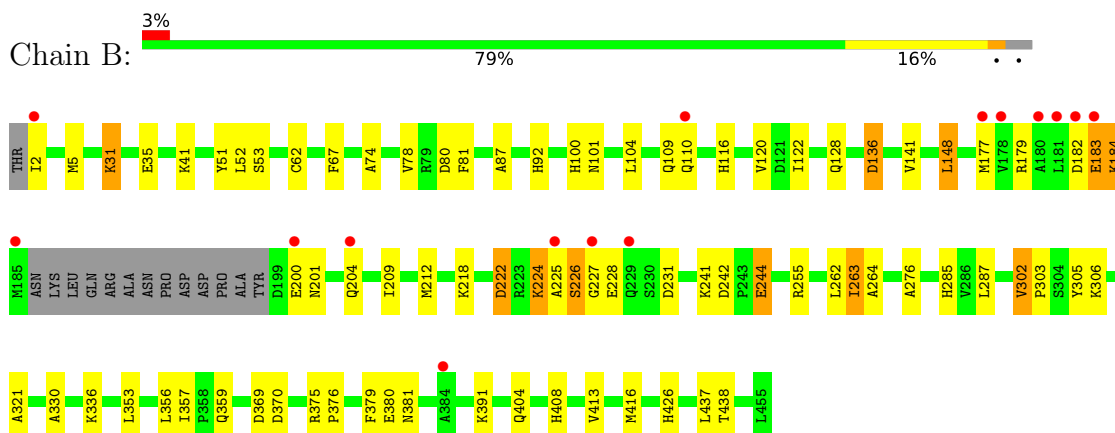
3 Residue-property plots

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: Bifunctional P-450/NADPH-P450 reductase



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.14Å 148.12Å 64.07Å 90.00° 98.16° 90.00°	Depositor
Resolution (Å)	19.51 – 1.70 19.51 – 1.70	Depositor EDS
% Data completeness (in resolution range)	98.3 (19.51-1.70) 98.3 (19.51-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.189 , 0.228 0.189 , 0.226	Depositor DCC
R_{free} test set	824 reflections (0.69%)	wwPDB-VP
Wilson B-factor (Å ²)	17.1	Xtriage
Anisotropy	0.032	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 51.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8178	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.86% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: CL, HEM, SYN

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	1.56	29/3714 (0.8%)	1.29	16/5012 (0.3%)
1	B	1.54	24/3700 (0.6%)	1.27	10/4993 (0.2%)
All	All	1.55	53/7414 (0.7%)	1.28	26/10005 (0.3%)

All (53) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	361	HIS	ND1-CE1	9.32	1.41	1.32
1	A	408	HIS	ND1-CE1	7.48	1.40	1.32
1	A	388	HIS	ND1-CE1	7.06	1.39	1.32
1	B	67	PHE	N-CA	6.87	1.54	1.46
1	B	263	ILE	C-O	6.84	1.30	1.24
1	B	116	HIS	CE1-NE2	6.76	1.39	1.32
1	A	135	ALA	C-O	-6.74	1.16	1.23
1	B	255	ARG	C-O	-6.50	1.16	1.24
1	A	453	ILE	N-CA	6.47	1.51	1.46
1	B	336	LYS	N-CA	6.44	1.54	1.46
1	B	302	VAL	CA-C	6.42	1.57	1.52
1	A	357	ILE	CA-C	6.25	1.57	1.52
1	A	378	ARG	N-CA	6.22	1.54	1.46
1	B	285	HIS	N-CA	6.22	1.53	1.46
1	B	379	PHE	N-CA	6.20	1.53	1.46
1	B	357	ILE	CA-CB	6.20	1.58	1.53
1	A	361	HIS	CE1-NE2	6.17	1.38	1.32
1	B	141	VAL	CA-C	6.13	1.59	1.52
1	A	116	HIS	C-O	-6.09	1.17	1.24
1	A	385	ILE	N-CA	6.07	1.53	1.46
1	A	426	HIS	CE1-NE2	5.89	1.38	1.32
1	B	241	LYS	N-CA	5.86	1.53	1.46
1	A	415	GLY	N-CA	5.79	1.53	1.45
1	B	408	HIS	CG-ND1	-5.77	1.31	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	403	GLN	C-O	5.74	1.30	1.24
1	B	262	LEU	C-O	5.65	1.30	1.24
1	B	438	THR	CA-C	5.60	1.60	1.52
1	A	178	VAL	CA-CB	-5.59	1.47	1.54
1	A	24	LYS	N-CA	5.59	1.52	1.46
1	B	321	ALA	C-O	5.55	1.30	1.24
1	A	408	HIS	N-CA	5.50	1.52	1.46
1	A	44	ALA	N-CA	5.49	1.50	1.45
1	B	416	MET	N-CA	5.43	1.52	1.46
1	A	120	VAL	N-CA	5.42	1.52	1.46
1	B	100	HIS	CE1-NE2	5.36	1.38	1.32
1	B	426	HIS	CE1-NE2	5.33	1.37	1.32
1	A	306	LYS	CA-C	5.32	1.59	1.52
1	A	412	LEU	N-CA	5.32	1.52	1.46
1	A	243	PRO	C-O	-5.31	1.17	1.24
1	B	437	LEU	N-CA	5.28	1.53	1.46
1	B	242	ASP	C-O	5.21	1.30	1.24
1	A	408	HIS	C-O	5.17	1.30	1.24
1	B	330	ALA	C-O	5.14	1.30	1.23
1	B	92	HIS	ND1-CE1	5.14	1.37	1.32
1	A	147	ARG	C-O	-5.13	1.18	1.24
1	A	279	PHE	CA-C	5.13	1.59	1.52
1	A	257	GLN	CD-OE1	5.12	1.33	1.23
1	A	408	HIS	CG-ND1	-5.12	1.32	1.38
1	A	118	MET	C-O	-5.11	1.17	1.24
1	A	357	ILE	CA-CB	5.11	1.57	1.53
1	B	376	PRO	C-O	5.10	1.30	1.24
1	B	104	LEU	CA-C	5.06	1.59	1.52
1	A	333	LEU	C-O	5.06	1.29	1.23

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	180	ALA	CA-C-N	8.53	132.41	120.29
1	A	180	ALA	C-N-CA	8.53	132.41	120.29
1	A	200	GLU	CA-C-N	-8.18	110.10	122.21
1	A	200	GLU	C-N-CA	-8.18	110.10	122.21
1	B	381	ASN	CA-C-N	-7.99	111.49	119.56
1	B	381	ASN	C-N-CA	-7.99	111.49	119.56
1	A	178	VAL	N-CA-C	7.29	117.42	110.42
1	B	222	ASP	N-CA-C	7.16	119.08	111.28
1	A	323	ARG	N-CA-C	-6.62	103.99	111.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	122	ILE	N-CA-C	6.56	117.35	110.72
1	A	202	LYS	N-CA-C	-6.35	107.38	114.62
1	B	244	GLU	CB-CG-CD	6.20	123.13	112.60
1	A	262	LEU	N-CA-C	-6.05	104.80	111.82
1	A	213	ASN	N-CA-C	5.92	117.53	111.14
1	B	302	VAL	N-CA-CB	-5.87	106.14	112.37
1	A	385	ILE	N-CA-C	5.63	113.47	108.63
1	A	209	ILE	N-CA-C	-5.61	105.06	110.72
1	B	136[A]	ASP	N-CA-C	5.61	119.70	112.86
1	B	136[B]	ASP	N-CA-C	5.61	119.70	112.86
1	B	276	ALA	N-CA-C	-5.47	105.22	111.07
1	A	121	ASP	N-CA-C	-5.36	105.34	111.07
1	A	179	ARG	N-CA-C	5.26	119.41	112.89
1	A	15	LYS	CA-CB-CG	-5.12	103.86	114.10
1	A	2	ILE	N-CA-C	5.09	115.61	108.89
1	A	166	TYR	N-CA-C	-5.05	106.70	112.92
1	B	369	ASP	N-CA-C	5.04	117.66	111.82

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	3607	0	3636	51	0
1	B	3599	0	3603	49	0
2	A	43	0	30	0	0
2	B	43	0	30	0	0
3	A	8	0	8	3	0
3	B	8	0	7	4	0
4	A	1	0	0	0	0
5	A	423	0	0	10	0
5	B	446	0	0	13	0
All	All	8178	0	7314	103	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 7.

All (103) 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:370:ASP:OD2	1:A:375:ARG:NH1	1.91	1.03
1:A:128:GLN:HG3	5:B:604:HOH:O	1.59	1.01
1:B:128:GLN:NE2	5:B:1003:HOH:O	1.92	0.98
1:B:183:GLU:OE1	1:B:201:ASN:HB3	1.71	0.90
1:A:267[B]:GLU:HG2	1:A:438:THR:HG21	1.54	0.89
1:A:199:ASP:O	1:A:202:LYS:HD3	1.74	0.87
1:A:118:MET:HG2	5:A:941:HOH:O	1.73	0.87
1:B:128:GLN:HG2	5:B:1004:HOH:O	1.77	0.85
1:A:310:GLN:HG3	5:A:899:HOH:O	1.78	0.82
1:A:43:GLU:HG2	1:A:48[B]:VAL:HG12	1.61	0.82
1:B:177:MET:HE2	1:B:263:ILE:HD12	1.64	0.80
1:A:267[B]:GLU:CG	1:A:438:THR:HG21	2.13	0.78
1:B:204:GLN:NE2	1:B:204:GLN:HA	2.00	0.75
1:A:267[B]:GLU:HG2	1:A:438:THR:CG2	2.16	0.75
1:B:222:ASP:OD2	5:B:885:HOH:O	2.05	0.75
1:B:226:SER:C	1:B:228:GLU:H	1.95	0.75
1:B:183:GLU:O	1:B:183:GLU:HG3	1.86	0.75
1:B:182:ASP:C	1:B:184:LYS:H	1.95	0.75
1:A:179:ARG:CD	1:A:204:GLN:OE1	2.35	0.74
1:A:316:MET:HE1	1:A:380:GLU:HG3	1.71	0.70
1:A:229:GLN:HA	1:A:229:GLN:OE1	1.93	0.67
1:A:309:LYS:O	1:A:312[B]:LYS:HE3	1.95	0.67
1:B:31[A]:LYS:HD2	5:B:811:HOH:O	1.95	0.66
1:A:179:ARG:HD2	1:A:204:GLN:OE1	1.95	0.66
1:A:212:MET:SD	5:A:1005:HOH:O	2.54	0.66
1:B:204:GLN:HA	1:B:204:GLN:HE21	1.59	0.64
3:B:502:SYN:H8	5:B:1046:HOH:O	1.97	0.64
1:A:183:GLU:HG2	1:A:205:PHE:HB2	1.79	0.64
1:A:183:GLU:O	1:A:183:GLU:HG3	1.98	0.64
1:A:147:ARG:HG2	1:A:164[B]:SER:OG	1.98	0.63
1:A:199:ASP:OD2	1:A:199:ASP:N	2.30	0.63
1:A:177:MET:HE2	1:A:263:ILE:HD12	1.81	0.62
1:B:177:MET:HB2	1:B:212:MET:HE3	1.82	0.62
1:B:183:GLU:OE1	1:B:201:ASN:CB	2.47	0.61
3:A:502:SYN:H8	5:A:874:HOH:O	2.01	0.60
1:B:303:PRO:HG3	5:B:883:HOH:O	2.01	0.60
1:B:404:GLN:HG3	5:B:1024:HOH:O	2.02	0.60
1:A:7:GLN:HB2	1:A:41[A]:LYS:HG3	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:ALA:O	3:A:502:SYN:H9	2.02	0.59
1:A:182:ASP:C	1:A:184:LYS:H	2.11	0.58
1:A:388:HIS:HA	1:A:391:LYS:HD3	1.86	0.58
1:B:224:LYS:NZ	1:B:225:ALA:HB2	2.19	0.57
1:B:74:ALA:O	1:B:78:VAL:HG23	2.05	0.57
1:B:80:ASP:HB2	1:B:184:LYS:NZ	2.20	0.55
1:A:272:LEU:HD13	1:A:322:LEU:HG	1.87	0.55
1:A:316:MET:CE	1:A:380:GLU:HG3	2.36	0.55
1:B:306:LYS:HG2	5:B:844:HOH:O	2.06	0.55
1:B:109:GLN:OE1	1:B:305:TYR:OH	2.18	0.55
1:B:101:ASN:HD21	1:B:244:GLU:CD	2.15	0.54
1:B:87:ALA:HB1	3:B:502:SYN:H4	1.90	0.54
1:A:201:ASN:OD1	1:A:201:ASN:N	2.40	0.53
1:B:35:GLU:HG3	5:B:979:HOH:O	2.08	0.53
1:A:296:ARG:NH2	5:A:849:HOH:O	2.38	0.53
1:B:287:LEU:C	1:B:287:LEU:HD23	2.34	0.52
1:B:177:MET:HE2	1:B:263:ILE:CD1	2.37	0.52
1:B:224:LYS:HZ3	1:B:225:ALA:HB2	1.75	0.52
1:A:267[B]:GLU:CG	1:A:438:THR:CG2	2.84	0.51
1:B:148:LEU:HD21	1:B:413:VAL:HG21	1.91	0.51
1:A:148:LEU:HD21	1:A:413:VAL:HG21	1.93	0.51
1:B:136[A]:ASP:O	1:B:136[A]:ASP:OD2	2.29	0.50
1:B:182:ASP:C	1:B:184:LYS:N	2.59	0.50
1:B:182:ASP:O	1:B:184:LYS:N	2.39	0.50
1:B:226:SER:O	1:B:228:GLU:N	2.45	0.50
1:B:53:SER:HB3	1:B:359:GLN:HB3	1.94	0.49
1:B:370:ASP:OD2	1:B:375[B]:ARG:NH1	2.41	0.49
1:A:94[B]:LYS:HE3	5:A:784:HOH:O	2.12	0.49
1:A:7:GLN:HB2	1:A:41[A]:LYS:HE2	1.95	0.49
1:A:204:GLN:HG3	1:A:205:PHE:N	2.27	0.49
1:B:226:SER:C	1:B:228:GLU:N	2.62	0.48
3:B:502:SYN:H8	5:B:821:HOH:O	2.14	0.48
1:A:177:MET:HG2	1:A:212:MET:CE	2.44	0.47
1:A:177:MET:CB	1:A:212:MET:HE3	2.45	0.47
1:A:81:PHE:HE2	1:A:209:ILE:HD11	1.80	0.47
1:B:209:ILE:HD13	1:B:209:ILE:HA	1.70	0.47
1:A:47:ARG:NH2	5:A:767:HOH:O	2.48	0.46
5:A:855:HOH:O	1:B:218:LYS:HE2	2.15	0.46
1:A:264:ALA:O	3:A:502:SYN:CAA	2.63	0.46
1:A:222:ASP:OD2	5:A:824:HOH:O	2.21	0.46
1:A:205:PHE:HD2	1:A:206:GLN:HG2	1.79	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:GLU:O	1:A:202:LYS:N	2.47	0.45
1:A:170:PRO:HA	5:A:873:HOH:O	2.18	0.44
1:A:177:MET:HG2	1:A:212:MET:HE1	2.00	0.44
1:B:31[A]:LYS:HB3	1:B:31[A]:LYS:HE2	1.40	0.44
1:A:41[B]:LYS:NZ	1:A:41[B]:LYS:CB	2.81	0.44
1:B:81:PHE:HE1	1:B:209:ILE:HD11	1.83	0.43
1:A:182:ASP:O	1:A:184:LYS:N	2.52	0.43
1:A:74:ALA:O	1:A:78:VAL:HG23	2.18	0.43
1:A:124:VAL:HG13	1:A:455:LEU:HD13	2.01	0.42
1:B:51:TYR:HD2	1:B:356:LEU:HD13	1.85	0.42
1:B:62:CYS:SG	1:B:391:LYS:HE2	2.59	0.42
1:A:205:PHE:CD2	1:A:205:PHE:C	2.97	0.42
1:A:2:ILE:HG13	1:A:346:PRO:HG3	2.02	0.42
1:B:31[A]:LYS:NZ	5:B:716:HOH:O	2.52	0.42
1:B:5:MET:HE3	1:B:5:MET:HB2	1.94	0.42
1:B:264:ALA:O	3:B:502:SYN:CAA	2.68	0.42
1:B:370:ASP:OD2	1:B:375[B]:ARG:NH2	2.49	0.41
1:A:439:LEU:O	1:A:440[A]:LYS:HG2	2.20	0.41
1:B:52:LEU:HD11	1:B:353:LEU:HD13	2.02	0.41
1:B:380:GLU:HB3	5:B:769:HOH:O	2.20	0.41
1:B:179:ARG:HH11	1:B:179:ARG:HG3	1.86	0.41
1:A:182:ASP:C	1:A:184:LYS:N	2.78	0.41
1:B:80:ASP:HB2	1:B:184:LYS:HZ2	1.85	0.40
1:B:120:VAL:HG11	1:B:302:VAL:CG1	2.52	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	447/455 (98%)	436 (98%)	9 (2%)	2 (0%)	30 16

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	445/455 (98%)	429 (96%)	14 (3%)	2 (0%)	30	16
All	All	892/910 (98%)	865 (97%)	23 (3%)	4 (0%)	24	16

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	183	GLU
1	B	183	GLU
1	B	227	GLY
1	A	370	ASP

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	396/398 (100%)	376 (95%)	20 (5%)	21	7
1	B	394/398 (99%)	383 (97%)	11 (3%)	38	21
All	All	790/796 (99%)	759 (96%)	31 (4%)	28	12

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ILE
1	A	15	LYS
1	A	59	LYS
1	A	148	LEU
1	A	179	ARG
1	A	182	ASP
1	A	184	LYS
1	A	199	ASP
1	A	201	ASN
1	A	203	ARG
1	A	204	GLN
1	A	207	GLU

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Mol	Chain	Res	Type
1	A	209	ILE
1	A	210	LYS
1	A	226	SER
1	A	228	GLU
1	A	233	LEU
1	A	302	VAL
1	A	306	LYS
1	A	369	ASP
1	B	2	ILE
1	B	31[A]	LYS
1	B	31[B]	LYS
1	B	41	LYS
1	B	110	GLN
1	B	148	LEU
1	B	184	LYS
1	B	200	GLU
1	B	224	LYS
1	B	226	SER
1	B	231	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	ASN
1	A	239	ASN
1	A	359	GLN
1	A	403	GLN
1	B	70	ASN
1	B	101	ASN
1	B	110	GLN
1	B	236	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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	SYN	B	502	-	8,8,8	1.88	2 (25%)	9,9,9	1.09	0
2	HEM	B	501	1	50,50,50	2.44	14 (28%)	67,82,82	2.71	21 (31%)
3	SYN	A	502	-	8,8,8	2.19	3 (37%)	9,9,9	0.97	1 (11%)
2	HEM	A	501	1	50,50,50	2.35	16 (32%)	67,82,82	2.58	18 (26%)

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	SYN	B	502	-	-	2/2/2/2	0/1/1/1
2	HEM	B	501	1	-	2/14/54/54	-
3	SYN	A	502	-	-	2/2/2/2	0/1/1/1
2	HEM	A	501	1	-	3/14/54/54	-

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	FE-NA	8.70	2.23	1.95
2	B	501	HEM	FE-NA	6.94	2.18	1.95
2	B	501	HEM	C3D-C2D	6.51	1.50	1.36
2	A	501	HEM	FE-ND	5.99	2.13	1.94
2	B	501	HEM	FE-NB	5.98	2.13	1.94
2	A	501	HEM	FE-NB	5.34	2.11	1.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	C3D-C2D	5.11	1.47	1.36
2	B	501	HEM	C4A-NA	-4.98	1.30	1.39
2	B	501	HEM	FE-ND	4.81	2.09	1.94
2	B	501	HEM	FE-NC	4.75	2.10	1.95
3	A	502	SYN	CAA-CAB	4.43	1.57	1.29
3	B	502	SYN	CAA-CAB	4.19	1.55	1.29
2	B	501	HEM	C1B-NB	-3.89	1.33	1.40
2	A	501	HEM	C4A-NA	-3.52	1.33	1.39
2	B	501	HEM	O2D-CGD	-3.01	1.20	1.30
2	A	501	HEM	C4C-NC	-3.00	1.34	1.39
2	B	501	HEM	O2A-CGA	-2.95	1.21	1.30
2	A	501	HEM	C1B-NB	-2.86	1.35	1.40
2	A	501	HEM	CAC-C3C	2.84	1.55	1.47
2	A	501	HEM	C3B-C2B	-2.76	1.31	1.37
3	A	502	SYN	CAE-CAG	2.73	1.43	1.38
2	B	501	HEM	CAC-C3C	2.71	1.54	1.47
2	B	501	HEM	C1C-NC	-2.57	1.34	1.39
2	A	501	HEM	O2A-CGA	-2.55	1.22	1.30
2	A	501	HEM	C1D-ND	-2.54	1.33	1.38
2	A	501	HEM	C1C-NC	-2.37	1.35	1.39
2	B	501	HEM	C1A-NA	-2.31	1.35	1.39
2	B	501	HEM	C4C-NC	-2.28	1.35	1.39
2	B	501	HEM	C4D-ND	-2.21	1.36	1.40
2	A	501	HEM	O2D-CGD	-2.21	1.23	1.30
2	A	501	HEM	CMC-C2C	2.11	1.55	1.50
2	A	501	HEM	C4B-NB	-2.10	1.34	1.38
3	A	502	SYN	CAH-CAB	-2.06	1.41	1.49
3	B	502	SYN	CAH-CAB	-2.05	1.41	1.49
2	A	501	HEM	C4D-ND	-2.05	1.36	1.40

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	C4D-ND-C1D	8.82	115.65	105.21
2	B	501	HEM	C1B-NB-C4B	8.56	115.34	105.21
2	B	501	HEM	C4D-ND-C1D	6.69	113.14	105.21
2	A	501	HEM	CHD-C4C-NC	6.52	131.55	124.45
2	B	501	HEM	C2B-C1B-NB	-6.39	102.50	109.84
2	B	501	HEM	C3B-C4B-NB	-6.32	104.93	109.47
2	A	501	HEM	C2A-C1A-NA	-6.10	103.38	110.15
2	A	501	HEM	C4A-NA-C1A	5.80	115.28	105.82
2	B	501	HEM	C3B-C2B-C1B	5.76	110.74	106.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	C3D-C4D-ND	-5.39	104.26	110.17
2	B	501	HEM	C4A-NA-C1A	5.18	114.26	105.82
2	B	501	HEM	C4C-NC-C1C	5.14	114.20	105.82
2	A	501	HEM	CHB-C1B-NB	5.06	130.62	124.37
2	B	501	HEM	CHD-C1D-ND	4.82	129.61	124.42
2	B	501	HEM	CHD-C4C-NC	4.73	129.60	124.45
2	A	501	HEM	C3B-C2B-C1B	4.66	109.91	106.41
2	A	501	HEM	C1B-NB-C4B	4.40	110.41	105.21
2	A	501	HEM	CHA-C4D-ND	4.10	129.44	124.37
2	B	501	HEM	C2A-C1A-NA	-3.94	105.78	110.15
2	B	501	HEM	CHA-C4D-ND	3.81	129.08	124.37
2	A	501	HEM	C4C-NC-C1C	3.80	112.02	105.82
2	A	501	HEM	C2B-C1B-NB	-3.59	105.71	109.84
2	B	501	HEM	C3A-C4A-NA	-3.55	104.45	110.14
2	A	501	HEM	C3A-C4A-NA	-3.54	104.46	110.14
2	B	501	HEM	CMB-C2B-C1B	-3.43	119.67	125.03
2	A	501	HEM	CHC-C1C-NC	3.37	128.12	124.45
2	B	501	HEM	CHC-C4B-NB	3.29	127.97	124.42
2	B	501	HEM	CHB-C1B-NB	3.21	128.34	124.37
2	B	501	HEM	CHC-C1C-NC	3.21	127.94	124.45
2	B	501	HEM	C2C-C1C-NC	-2.90	104.28	109.64
2	A	501	HEM	C3B-C4B-NB	-2.75	107.50	109.47
2	A	501	HEM	CBD-CAD-C3D	-2.54	105.51	112.53
2	A	501	HEM	CHA-C1A-NA	2.37	128.17	123.86
2	B	501	HEM	O1D-CGD-CBD	-2.36	115.61	123.09
2	A	501	HEM	CHD-C1D-ND	2.35	126.95	124.42
2	B	501	HEM	C3D-C4D-ND	-2.30	107.65	110.17
2	B	501	HEM	CHD-C1D-C2D	-2.26	121.46	125.03
2	A	501	HEM	C2D-C1D-ND	-2.25	107.30	109.90
3	A	502	SYN	CAH-CAB-CAA	-2.14	114.49	125.87
2	B	501	HEM	C4A-C3A-C2A	2.08	109.20	106.82

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	502	SYN	CAA-CAB-CAH-CAG
3	B	502	SYN	CAA-CAB-CAH-CAF
3	A	502	SYN	CAA-CAB-CAH-CAF
3	A	502	SYN	CAA-CAB-CAH-CAG
2	B	501	HEM	CAD-CBD-CGD-O1D
2	A	501	HEM	CAD-CBD-CGD-O2D

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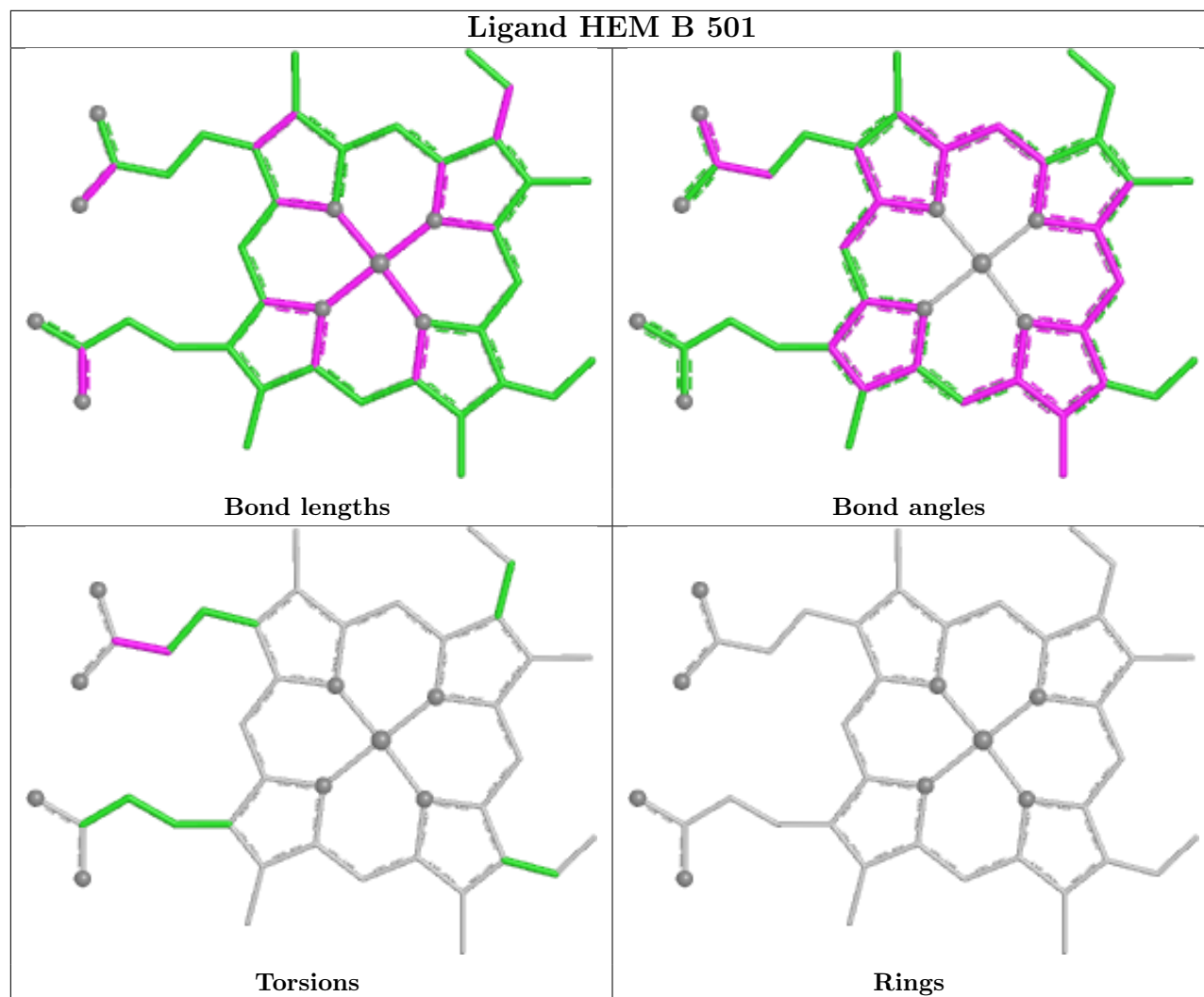
Mol	Chain	Res	Type	Atoms
2	A	501	HEM	CAD-CBD-CGD-O1D
2	B	501	HEM	CAD-CBD-CGD-O2D
2	A	501	HEM	CAA-CBA-CGA-O1A

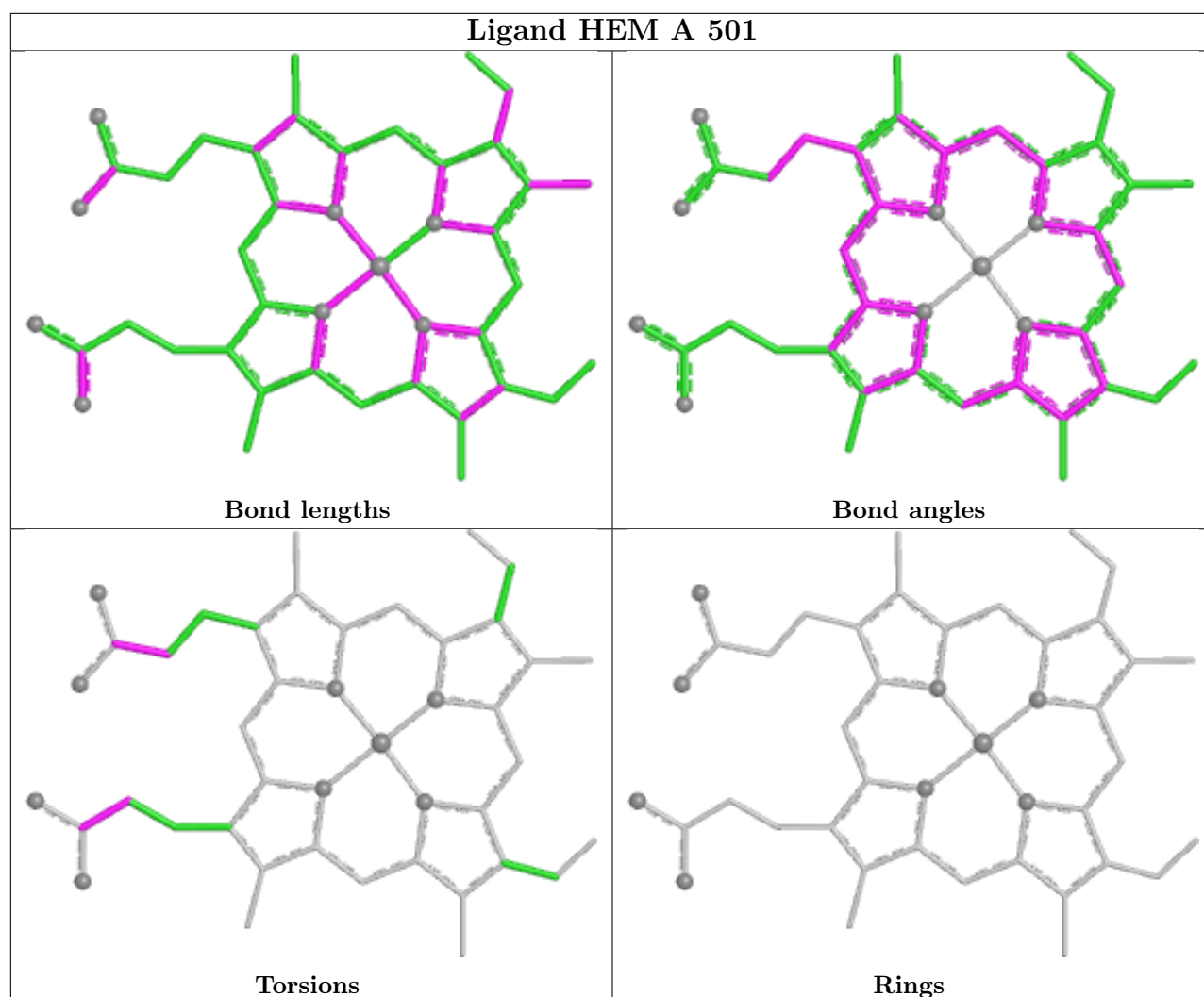
There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	502	SYN	4	0
3	A	502	SYN	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

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	442/455 (97%)	-0.24	20 (4%) 38 42	6, 15, 50, 125	10 (2%)
1	B	441/455 (96%)	-0.24	15 (3%) 48 52	7, 16, 44, 104	8 (1%)
All	All	883/910 (97%)	-0.24	35 (3%) 42 46	6, 16, 49, 125	18 (2%)

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	2	ILE	6.3
1	A	180	ALA	5.5
1	A	199	ASP	4.4
1	A	181	LEU	4.1
1	A	205	PHE	4.1
1	A	77	PHE	4.0
1	B	200	GLU	3.8
1	B	185	MET	3.8
1	A	200	GLU	3.7
1	A	178	VAL	3.4
1	A	2	ILE	3.3
1	A	202	LYS	3.2
1	B	227	GLY	3.2
1	A	182	ASP	3.1
1	A	185	MET	3.0
1	A	170	PRO	3.0
1	B	181	LEU	2.9
1	A	201	ASN	2.9
1	A	209	ILE	2.6
1	B	225	ALA	2.6
1	A	227	GLY	2.6
1	B	229	GLN	2.4
1	A	207	GLU	2.4
1	A	1	THR	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	177	MET	2.3
1	B	384	ALA	2.2
1	A	206	GLN	2.2
1	B	178	VAL	2.2
1	B	183	GLU	2.2
1	B	180	ALA	2.2
1	B	204	GLN	2.1
1	A	203	ARG	2.1
1	B	110	GLN	2.1
1	A	179	ARG	2.1
1	B	182	ASP	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

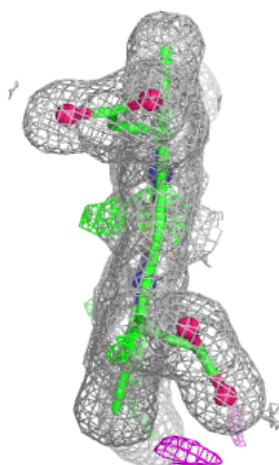
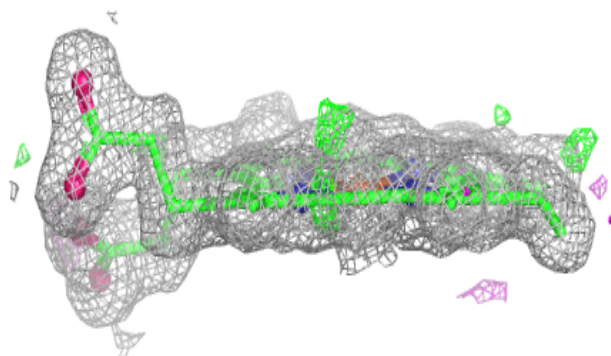
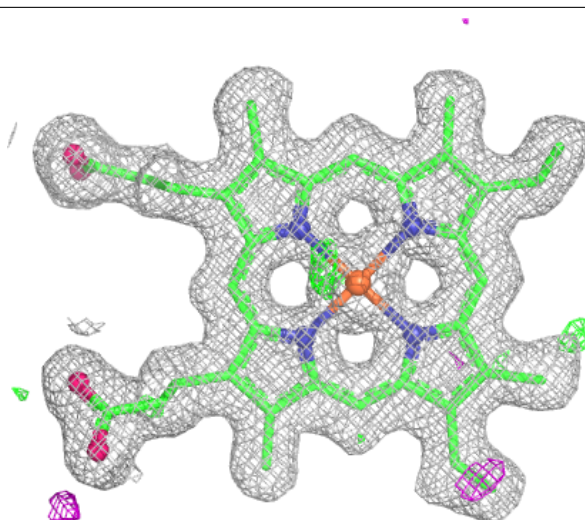
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	SYN	A	502	8/8	0.87	0.13	19,30,38,41	0
3	SYN	B	502	8/8	0.87	0.13	21,32,37,38	0
4	CL	A	503	1/1	0.90	0.11	53,53,53,53	0
2	HEM	B	501	43/43	0.98	0.05	8,9,12,15	0
2	HEM	A	501	43/43	0.98	0.05	8,10,13,15	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.

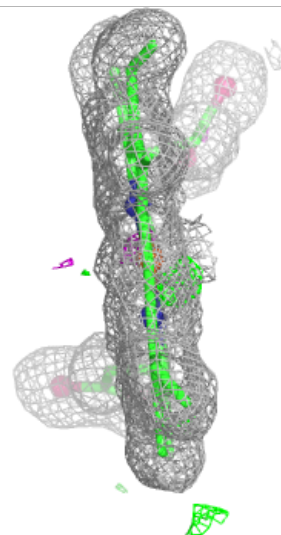
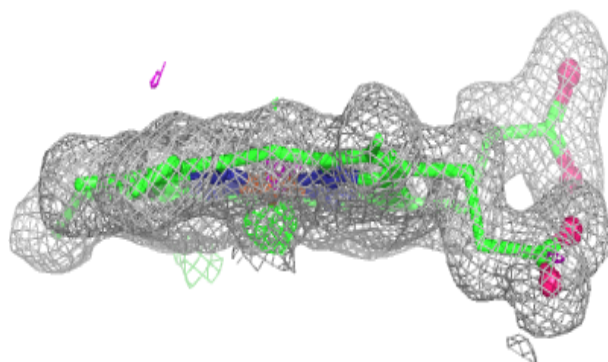
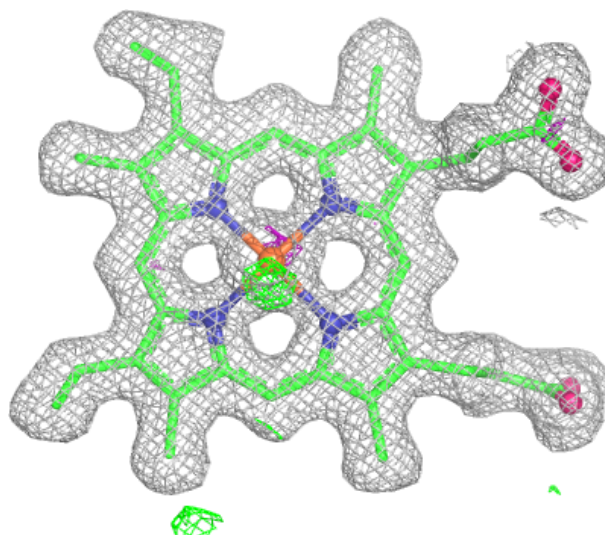
Electron density around HEM B 501:

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



Electron density around HEM 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.