



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:53 AM UTC

PDB ID : 4HGJ / pdb_00004hgj
Title : Crystal structure of P450 BM3 5F5 heme domain variant
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Deposited on : 2012-10-08
Resolution : 1.90 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

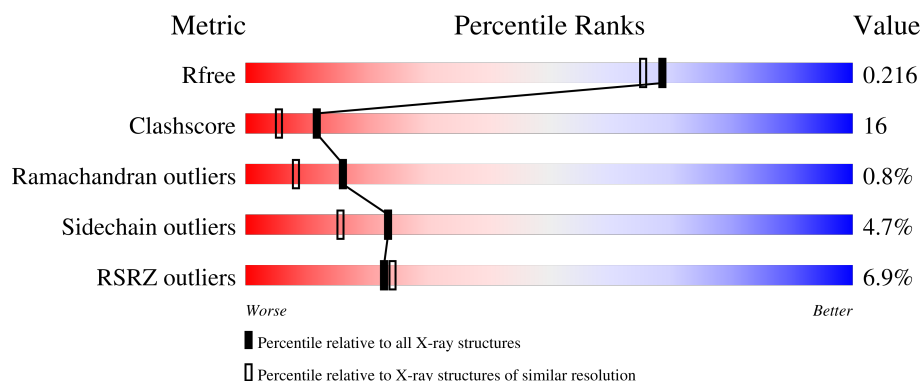
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	7% 75% 21% ..
1	B	455	7% 77% 18% ..

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.

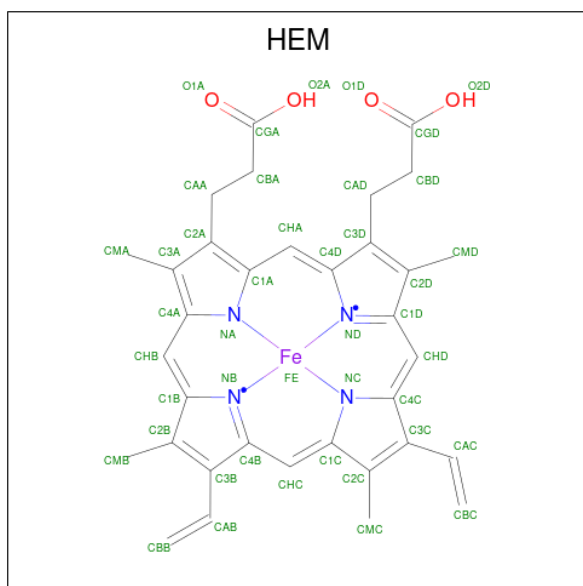
- Molecule 1 is a protein called Bifunctional P-450/NADPH-P450 reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	448	Total 3708	C 2375	N 627	O 688	S 18	0	15	0
1	B	451	Total 3697	C 2366	N 628	O 686	S 17	0	11	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	87	ALA	PHE	engineered mutation	UNP P14779
A	235	ALA	THR	engineered mutation	UNP P14779
B	87	ALA	PHE	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 GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O		
			6	3	3	0	0
3	A	1	Total	C	O		
			6	3	3	0	0
3	A	1	Total	C	O		
			6	3	3	0	0
3	B	1	Total	C	O		
			6	3	3	0	0
3	B	1	Total	C	O		
			6	3	3	0	0

- Molecule 4 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

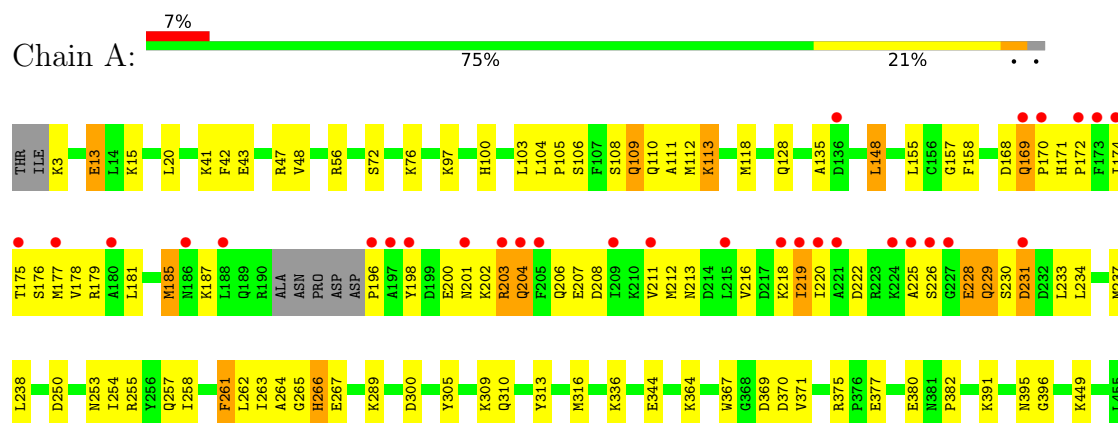
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	391	Total	O	0	0
			391	391		
5	B	412	Total	O	0	0
			412	412		

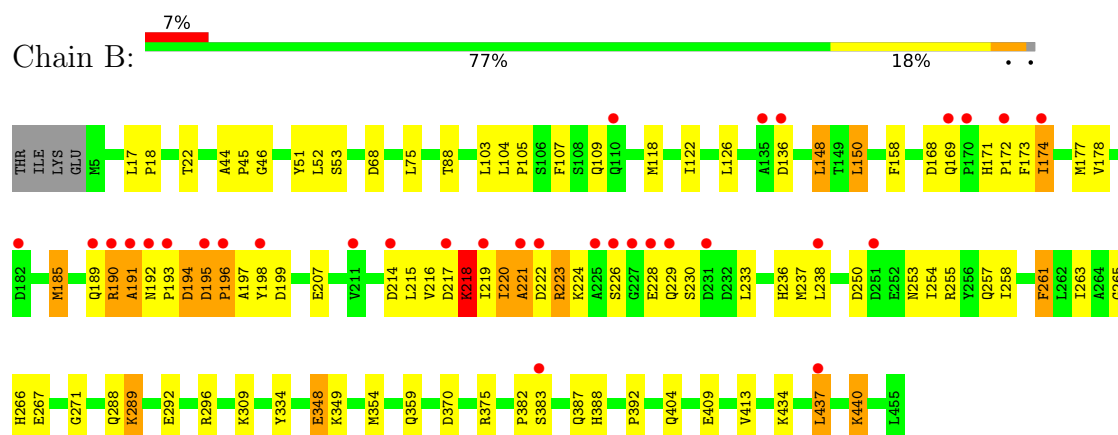
3 Residue-property plots [i](#)

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: Bifunctional P-450/NADPH-P450 reductase



- Molecule 1: Bifunctional P-450/NADPH-P450 reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.01Å 148.37Å 64.15Å 90.00° 98.20° 90.00°	Depositor
Resolution (Å)	19.93 – 1.90 19.93 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.4 (19.93-1.90) 97.4 (19.93-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.69 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.182 , 0.217 0.182 , 0.216	Depositor DCC
R_{free} test set	4174 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	21.8	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 56.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8336	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.97% 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: GOL, MES, 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	0.54	4/3820 (0.1%)	0.81	5/5155 (0.1%)
1	B	0.51	2/3815 (0.1%)	0.81	3/5156 (0.1%)
All	All	0.53	6/7635 (0.1%)	0.81	8/10311 (0.1%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	56	ARG	C-O	-6.57	1.16	1.24
1	B	261[A]	PHE	C-O	-5.52	1.17	1.24
1	B	261[B]	PHE	C-O	-5.52	1.17	1.24
1	A	266	HIS	ND1-CE1	5.46	1.38	1.32
1	A	261[A]	PHE	C-O	-5.05	1.18	1.24
1	A	261[B]	PHE	C-O	-5.05	1.18	1.24

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	195	ASP	N-CA-C	-7.23	101.29	109.60
1	A	266	HIS	CB-CG-CD2	-6.59	122.64	131.20
1	B	194	ASP	N-CA-C	6.04	122.73	113.19
1	A	266	HIS	CB-CG-ND1	5.80	131.41	122.70
1	A	204	GLN	N-CA-C	-5.62	105.06	111.07
1	A	15	LYS	CB-CA-C	-5.45	110.31	116.63
1	B	226	SER	N-CA-C	-5.27	106.69	113.23
1	A	135	ALA	N-CA-C	5.16	119.84	111.37

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	3708	0	3700	110	0
1	B	3697	0	3697	119	0
2	A	43	0	30	4	0
2	B	43	0	30	10	0
3	A	18	0	24	5	0
3	B	12	0	16	0	0
4	A	12	0	12	0	0
5	A	391	0	0	12	0
5	B	412	0	0	22	0
All	All	8336	0	7509	239	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 16.

All (239) 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:177:MET:HG3	1:A:212:MET:CE	1.82	1.09
1:B:437:LEU:HD12	1:B:437:LEU:H	1.17	1.08
1:B:171:HIS:HB3	1:B:174:ILE:CG1	1.90	1.02
1:B:440[A]:LYS:HD2	5:B:999:HOH:O	1.63	0.97
1:B:171:HIS:CG	1:B:172:PRO:HD2	2.00	0.96
1:A:179:ARG:NE	1:A:204:GLN:OE1	1.99	0.96
2:B:502:HEM:HBC2	2:B:502:HEM:HHD	1.48	0.95
1:B:171:HIS:HB3	1:B:174:ILE:HG12	1.49	0.95
1:B:214:ASP:O	1:B:218:LYS:HD3	1.68	0.93
1:A:370:ASP:OD2	1:A:375:ARG:NH1	2.03	0.92
1:B:437:LEU:H	1:B:437:LEU:CD1	1.82	0.91
1:A:234:LEU:HA	1:A:237:MET:HE3	1.54	0.89
1:A:234:LEU:HA	1:A:237:MET:CE	2.02	0.89
1:A:200:GLU:O	1:A:203:ARG:HG3	1.72	0.89
1:B:237:MET:HE1	1:B:258:ILE:CG1	2.02	0.88
1:A:181:LEU:O	1:A:185:MET:HG3	1.72	0.88
1:B:292:GLU:OE2	1:B:296[A]:ARG:NH1	2.07	0.88
1:B:189:GLN:O	1:B:191:ALA:N	2.08	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:237:MET:HE1	1:B:258:ILE:HG13	1.55	0.87
1:A:202:LYS:O	1:A:206:GLN:HG2	1.73	0.86
1:B:177:MET:SD	1:B:263:ILE:HG12	2.16	0.86
1:B:185:MET:CE	1:B:437:LEU:HB3	2.07	0.85
1:A:41:LYS:NZ	1:A:43:GLU:OE2	2.11	0.83
1:B:237:MET:HE3	1:B:254:ILE:HG23	1.60	0.83
1:A:177:MET:HG3	1:A:212:MET:HE2	1.63	0.81
1:A:395:ASN:HA	3:A:502:GOL:H11	1.61	0.81
1:B:158:PHE:HZ	1:B:261[B]:PHE:HD2	1.28	0.81
1:A:128[B]:GLN:NE2	5:A:855:HOH:O	2.05	0.80
1:A:174:ILE:O	1:A:178:VAL:HG23	1.82	0.78
1:B:217:ASP:OD1	5:B:896:HOH:O	2.02	0.78
1:B:216:VAL:HG11	1:B:255:ARG:HB2	1.66	0.77
1:B:440[A]:LYS:CD	5:B:999:HOH:O	2.23	0.77
1:B:189:GLN:O	1:B:190:ARG:C	2.24	0.77
1:A:177:MET:SD	1:A:263:ILE:HG12	2.26	0.76
1:B:267:GLU:OE1	1:B:437:LEU:HD21	1.84	0.76
1:A:177:MET:CG	1:A:212:MET:CE	2.64	0.76
1:A:177:MET:HE1	1:A:266:HIS:HE1	1.52	0.75
1:B:185:MET:HE1	1:B:437:LEU:HB3	1.68	0.75
1:A:177:MET:CB	1:A:212:MET:HE3	2.17	0.74
1:A:226:SER:OG	1:A:228:GLU:HG2	1.87	0.74
1:B:288:GLN:OE1	5:B:909:HOH:O	2.07	0.72
2:B:502:HEM:HBB2	2:B:502:HEM:HMB2	1.71	0.71
2:B:502:HEM:HBC1	5:B:994:HOH:O	1.90	0.71
1:B:309:LYS:NZ	5:B:667:HOH:O	2.22	0.71
2:B:502:HEM:HBB2	2:B:502:HEM:CMB	2.21	0.70
1:A:103:LEU:HD13	1:A:261[B]:PHE:HZ	1.55	0.70
1:B:437:LEU:HD12	1:B:437:LEU:N	2.00	0.69
1:A:109[B]:GLN:O	1:A:112:MET:HB2	1.93	0.69
1:A:174:ILE:HG22	1:A:178:VAL:CG2	2.23	0.68
1:A:177:MET:CA	1:A:212:MET:HE3	2.23	0.68
1:A:267[B]:GLU:OE1	1:A:267[B]:GLU:HA	1.93	0.68
1:A:177:MET:HG3	1:A:212:MET:HE1	1.71	0.68
1:B:185:MET:HE3	1:B:437:LEU:HB3	1.76	0.68
1:A:13:GLU:H	1:A:13:GLU:CD	2.02	0.68
1:B:158:PHE:CZ	1:B:261[B]:PHE:HD2	2.12	0.67
1:B:118:MET:HE2	5:B:735:HOH:O	1.95	0.67
2:B:502:HEM:CBC	5:B:994:HOH:O	2.43	0.67
1:A:229:GLN:HA	1:A:229:GLN:OE1	1.93	0.67
1:A:177:MET:HA	1:A:212:MET:HE3	1.75	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:193:PRO:O	1:B:194:ASP:OD1	2.14	0.65
1:A:168:ASP:N	1:B:168:ASP:OD1	2.24	0.65
1:B:107:PHE:CZ	1:B:261[A]:PHE:CE1	2.84	0.65
1:A:106[B]:SER:HB3	1:A:233:LEU:HD23	1.78	0.64
1:B:171:HIS:ND1	1:B:172:PRO:HD2	2.11	0.64
1:B:292:GLU:CD	1:B:296[A]:ARG:NH1	2.56	0.64
1:B:177:MET:HE1	1:B:266:HIS:HE1	1.63	0.64
1:B:222:ASP:OD2	5:B:809:HOH:O	2.15	0.64
1:A:220:ILE:HD11	1:A:258:ILE:CD1	2.27	0.64
1:B:171:HIS:HB3	1:B:174:ILE:CD1	2.27	0.64
1:A:207:GLU:O	1:A:211:VAL:HG23	1.97	0.63
1:A:336:LYS:NZ	3:A:503:GOL:O2	2.31	0.63
1:B:194:ASP:O	1:B:195:ASP:C	2.42	0.63
1:B:171:HIS:CD2	1:B:172:PRO:HD2	2.33	0.62
1:B:292:GLU:CD	1:B:296[A]:ARG:HH12	2.07	0.62
1:B:109:GLN:CD	1:B:309:LYS:HZ3	2.06	0.62
1:A:203:ARG:O	1:A:207:GLU:HG2	2.01	0.61
1:A:204:GLN:O	1:A:207:GLU:HG3	1.99	0.61
1:A:108[B]:SER:O	1:A:112:MET:HG2	2.00	0.60
1:A:106[A]:SER:HB3	1:A:233:LEU:HD23	1.82	0.60
1:B:289:LYS:HD2	5:B:745:HOH:O	2.00	0.60
1:A:218:LYS:O	1:A:222:ASP:HB2	2.00	0.60
1:A:174:ILE:HG22	1:A:178:VAL:HG23	1.83	0.59
1:B:440[A]:LYS:HE3	5:B:999:HOH:O	2.03	0.59
1:B:237:MET:CE	1:B:258:ILE:HG13	2.31	0.58
1:A:47:ARG:NH2	5:A:940:HOH:O	2.34	0.58
1:A:169:GLN:HB2	1:A:170:PRO:HD2	1.85	0.57
1:A:310:GLN:HG3	5:A:977:HOH:O	2.03	0.57
1:B:171:HIS:CB	1:B:174:ILE:HG12	2.29	0.57
1:A:395:ASN:HA	3:A:502:GOL:C1	2.31	0.57
1:B:193:PRO:C	1:B:194:ASP:OD1	2.48	0.57
1:B:288:GLN:CD	5:B:909:HOH:O	2.47	0.57
1:A:204:GLN:O	1:A:207:GLU:CG	2.53	0.56
1:A:187:LYS:NZ	5:A:820:HOH:O	2.32	0.56
1:A:367:TRP:HB2	1:A:371:VAL:HG12	1.88	0.56
1:B:440[B]:LYS:CE	5:B:1012:HOH:O	2.53	0.56
1:B:109:GLN:CD	1:B:309:LYS:NZ	2.63	0.55
1:B:122:ILE:HG22	1:B:148:LEU:HD22	1.88	0.55
1:B:387:GLN:HG2	1:B:388:HIS:CD2	2.41	0.55
1:A:41:LYS:NZ	1:A:43:GLU:CD	2.64	0.55
1:B:370:ASP:OD2	1:B:375[B]:ARG:NH1	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:MET:HA	1:A:212:MET:CE	2.36	0.55
1:B:267:GLU:OE1	1:B:437:LEU:CD2	2.54	0.55
1:A:316:MET:CE	1:A:377:GLU:HA	2.37	0.55
1:A:174:ILE:CG2	1:A:178:VAL:CG2	2.85	0.54
1:B:192:ASN:O	1:B:198:TYR:CE2	2.60	0.54
1:B:437:LEU:CD1	1:B:437:LEU:N	2.54	0.54
1:A:382:PRO:HG3	5:A:809:HOH:O	2.07	0.54
1:B:215:LEU:O	1:B:219:ILE:HG13	2.08	0.54
1:B:196:PRO:O	1:B:199:ASP:HB2	2.08	0.54
1:A:264:ALA:HA	5:A:770:HOH:O	2.07	0.54
1:A:158[B]:PHE:HZ	1:A:262:LEU:HD21	1.73	0.54
1:B:158:PHE:CE1	1:B:258:ILE:HG12	2.43	0.54
1:B:171:HIS:HB3	1:B:174:ILE:HD11	1.90	0.54
1:B:107:PHE:CZ	1:B:261[A]:PHE:HE1	2.25	0.53
2:B:502:HEM:HMB2	2:B:502:HEM:CBB	2.37	0.53
1:A:108[A]:SER:O	1:A:111[A]:ALA:HB3	2.09	0.53
1:A:266:HIS:C	1:A:266:HIS:CD2	2.87	0.53
1:B:265:GLY:HA2	2:B:502:HEM:HBB2	1.88	0.53
1:A:198:TYR:HD1	1:A:201:ASN:OD1	1.92	0.53
1:B:292:GLU:OE1	1:B:296[A]:ARG:NH1	2.42	0.53
1:A:148:LEU:HD13	1:A:148:LEU:C	2.33	0.53
1:A:198:TYR:CD1	1:A:201:ASN:OD1	2.62	0.53
1:B:150:LEU:HD11	1:B:174:ILE:HD12	1.91	0.53
1:A:169:GLN:CB	1:A:170:PRO:HD2	2.39	0.53
1:A:106[A]:SER:HB3	1:A:233:LEU:CD2	2.38	0.53
1:A:250:ASP:O	1:A:254:ILE:HG13	2.09	0.53
1:B:150:LEU:HD11	1:B:174:ILE:CD1	2.37	0.53
1:B:228:GLU:OE2	1:B:228:GLU:HA	2.08	0.53
1:A:267[A]:GLU:HB3	5:A:761:HOH:O	2.09	0.52
1:A:264:ALA:O	5:A:770:HOH:O	2.18	0.52
1:A:289:LYS:HD3	1:A:313:TYR:CZ	2.44	0.52
1:B:238:LEU:HD23	1:B:254:ILE:HD13	1.92	0.52
1:B:265:GLY:HA2	2:B:502:HEM:CBB	2.40	0.52
2:A:501:HEM:HMC2	2:A:501:HEM:HBC2	1.92	0.52
1:B:223:ARG:HD2	5:B:1007:HOH:O	2.10	0.51
1:B:195:ASP:O	1:B:197:ALA:N	2.43	0.51
1:B:185:MET:HE3	1:B:437:LEU:CB	2.41	0.51
2:A:501:HEM:HBC2	2:A:501:HEM:CMC	2.41	0.51
1:A:106[B]:SER:HB3	1:A:233:LEU:CD2	2.40	0.50
1:A:253:ASN:O	1:A:257:GLN:HG2	2.12	0.50
1:A:174:ILE:HG22	1:A:178:VAL:HG21	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:219:ILE:O	1:B:222:ASP:HB2	2.11	0.50
1:B:174:ILE:O	1:B:178:VAL:HG23	2.11	0.50
1:B:440[A]:LYS:CE	5:B:999:HOH:O	2.53	0.50
1:B:171:HIS:CE1	1:B:172:PRO:HD2	2.47	0.49
1:B:253:ASN:O	1:B:257:GLN:HG2	2.12	0.49
1:B:51:TYR:CE2	1:B:354:MET:HG2	2.46	0.49
1:B:392:PRO:O	2:B:502:HEM:CBC	2.60	0.49
1:B:440[B]:LYS:HE2	5:B:1012:HOH:O	2.10	0.49
1:A:257:GLN:O	1:A:261[A]:PHE:HD1	1.96	0.49
1:B:440[B]:LYS:NZ	5:B:851:HOH:O	2.45	0.49
1:A:174:ILE:CG2	1:A:178:VAL:HG23	2.42	0.49
1:B:104:LEU:N	1:B:105:PRO:HD2	2.28	0.49
1:A:204:GLN:HA	1:A:207:GLU:HG2	1.96	0.48
1:A:157:GLY:C	1:A:158[A]:PHE:CG	2.91	0.48
1:B:434:LYS:HB3	1:B:440[A]:LYS:HD3	1.95	0.48
1:A:158[B]:PHE:CD1	1:A:219:ILE:HD13	2.49	0.48
1:A:177:MET:HB2	1:A:212:MET:HE3	1.95	0.48
1:B:237:MET:O	1:B:254:ILE:HD11	2.14	0.48
1:B:216:VAL:CG1	1:B:255:ARG:HB2	2.42	0.48
1:A:364:LYS:HD3	1:A:369:ASP:HA	1.95	0.47
1:B:404[B]:GLN:OE1	5:B:657:HOH:O	2.20	0.47
1:A:118[A]:MET:SD	1:A:155:LEU:HD21	2.54	0.47
1:B:44:ALA:HB1	1:B:45:PRO:HD2	1.95	0.47
1:A:158[B]:PHE:HB3	1:A:234:LEU:HB2	1.97	0.47
1:A:158[B]:PHE:CE1	1:A:219:ILE:HD13	2.50	0.47
1:B:17:LEU:HB3	1:B:18:PRO:HD3	1.95	0.47
1:B:68:ASP:HB3	1:B:334:TYR:CE1	2.50	0.47
1:B:220:ILE:O	1:B:221:ALA:C	2.58	0.47
1:B:103:LEU:HD21	1:B:237:MET:HG3	1.97	0.47
1:B:148:LEU:HD11	1:B:413:VAL:HG21	1.97	0.47
2:A:501:HEM:CMB	2:A:501:HEM:HBB2	2.45	0.46
1:B:107:PHE:HZ	1:B:261[A]:PHE:CE1	2.33	0.46
1:B:257:GLN:O	1:B:261[A]:PHE:HD2	1.98	0.46
1:B:237:MET:HE1	1:B:258:ILE:CD1	2.46	0.46
1:B:173:PHE:HD2	1:B:174:ILE:HD13	1.80	0.45
1:B:195:ASP:C	1:B:197:ALA:H	2.24	0.45
1:B:173:PHE:CD2	1:B:173:PHE:C	2.95	0.45
1:A:113:LYS:HD2	1:A:305:TYR:CE2	2.51	0.45
1:A:177:MET:CB	1:A:212:MET:CE	2.88	0.45
1:A:179:ARG:CZ	1:A:204:GLN:OE1	2.64	0.45
1:A:204:GLN:CA	1:A:207:GLU:HG2	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:220:ILE:HG23	1:B:238:LEU:HD21	1.99	0.45
1:A:300:ASP:OD2	5:A:926:HOH:O	2.21	0.45
1:A:396:GLY:H	3:A:502:GOL:C1	2.30	0.45
1:A:230[B]:SER:OG	1:A:231:ASP:N	2.48	0.45
1:A:41:LYS:HZ2	1:A:43:GLU:CD	2.24	0.44
2:A:501:HEM:HBB2	2:A:501:HEM:HMB2	1.99	0.44
1:A:213:ASN:OD1	1:A:255:ARG:HD3	2.17	0.44
1:A:196:PRO:C	1:A:198:TYR:N	2.72	0.44
1:A:109[A]:GLN:O	1:A:110[A]:GLN:C	2.60	0.44
1:A:175:THR:CG2	1:A:179:ARG:HH12	2.31	0.44
1:B:171:HIS:C	1:B:174:ILE:HG12	2.43	0.44
1:B:236:HIS:CE1	5:B:1010:HOH:O	2.70	0.44
1:A:171:HIS:CG	1:A:172:PRO:HD2	2.52	0.44
1:A:100:HIS:HE1	3:A:505:GOL:O2	2.01	0.43
1:B:288:GLN:NE2	5:B:909:HOH:O	2.51	0.43
1:B:382:PRO:HD2	5:B:953:HOH:O	2.18	0.43
1:A:43:GLU:HG2	1:A:48:VAL:HG22	2.00	0.43
1:B:171:HIS:CG	1:B:172:PRO:CD	2.87	0.43
1:B:75:LEU:HD22	1:B:88:THR:HG22	2.01	0.43
1:B:107:PHE:CE2	1:B:261[A]:PHE:CE1	3.07	0.43
1:B:171:HIS:HB3	1:B:174:ILE:HG13	1.92	0.43
1:A:200:GLU:O	1:A:203:ARG:CG	2.55	0.43
1:B:46:GLY:N	5:B:913:HOH:O	2.27	0.43
1:B:75:LEU:HD23	1:B:75:LEU:HA	1.79	0.42
1:B:250:ASP:C	1:B:250:ASP:OD1	2.62	0.42
1:A:225:ALA:O	1:A:226:SER:C	2.62	0.42
1:B:237:MET:CE	1:B:254:ILE:HG23	2.39	0.42
1:B:348:GLU:O	1:B:349:LYS:C	2.62	0.42
1:A:104:LEU:N	1:A:105:PRO:CD	2.83	0.42
1:A:104:LEU:HB3	1:A:105:PRO:HD3	2.02	0.42
1:A:266:HIS:N	5:A:759:HOH:O	2.52	0.42
1:B:271:GLY:HA2	1:B:440[B]:LYS:HG3	2.00	0.42
2:B:502:HEM:HHD	2:B:502:HEM:CBC	2.35	0.42
1:A:234:LEU:HA	1:A:237:MET:HE2	1.94	0.42
1:A:237:MET:HE3	1:A:237:MET:HB2	1.83	0.41
1:A:169:GLN:H	1:A:169:GLN:HG2	1.38	0.41
1:B:53:SER:HB3	1:B:359:GLN:HB3	2.01	0.41
1:A:47:ARG:HD2	5:A:942:HOH:O	2.19	0.41
1:A:72:SER:O	1:A:76:LYS:HG3	2.21	0.41
1:B:107:PHE:CE2	1:B:233:LEU:HD21	2.56	0.41
1:A:216:VAL:HG13	1:A:258:ILE:HD13	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:218:LYS:O	1:B:219:ILE:C	2.63	0.41
1:A:255:ARG:HA	1:A:258:ILE:HD12	2.03	0.41
1:A:220:ILE:HG23	1:A:238:LEU:HD21	2.03	0.41
1:B:177:MET:HE1	1:B:266:HIS:CE1	2.50	0.41
1:A:169:GLN:CB	1:A:170:PRO:CD	2.99	0.41
1:B:171:HIS:O	1:B:174:ILE:HG12	2.21	0.41
1:B:216:VAL:O	1:B:220:ILE:HG13	2.21	0.41
1:A:97:LYS:HB2	5:A:740:HOH:O	2.21	0.40
1:A:176:SER:O	1:A:208:ASP:HB3	2.21	0.40
1:B:185:MET:HE3	1:B:437:LEU:CA	2.52	0.40
1:B:158:PHE:HZ	1:B:261[B]:PHE:CD2	2.20	0.40
1:B:197:ALA:C	1:B:199:ASP:H	2.30	0.40
1:A:20:LEU:HG	1:A:42:PHE:CZ	2.56	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	459/455 (101%)	438 (95%)	20 (4%)	1 (0%)	43	36
1	B	460/455 (101%)	436 (95%)	18 (4%)	6 (1%)	9	3
All	All	919/910 (101%)	874 (95%)	38 (4%)	7 (1%)	16	8

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	190	ARG
1	B	191	ALA
1	B	220	ILE
1	B	221	ALA
1	B	218	LYS

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Mol	Chain	Res	Type
1	A	228	GLU
1	B	196	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	405/397 (102%)	388 (96%)	17 (4%)	26	19
1	B	404/397 (102%)	381 (94%)	23 (6%)	18	11
All	All	809/794 (102%)	769 (95%)	40 (5%)	23	14

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	13	GLU
1	A	109[A]	GLN
1	A	109[B]	GLN
1	A	113	LYS
1	A	148	LEU
1	A	169	GLN
1	A	185	MET
1	A	203	ARG
1	A	219	ILE
1	A	229	GLN
1	A	231	ASP
1	A	309	LYS
1	A	344	GLU
1	A	380	GLU
1	A	391	LYS
1	A	449	LYS
1	B	22	THR
1	B	52	LEU
1	B	126	LEU
1	B	136	ASP

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Mol	Chain	Res	Type
1	B	148	LEU
1	B	150	LEU
1	B	169	GLN
1	B	174	ILE
1	B	185	MET
1	B	207	GLU
1	B	218	LYS
1	B	223	ARG
1	B	224	LYS
1	B	229	GLN
1	B	230	SER
1	B	289	LYS
1	B	348	GLU
1	B	383	SER
1	B	409[A]	GLU
1	B	409[B]	GLU
1	B	437	LEU
1	B	440[A]	LYS
1	B	440[B]	LYS

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	359	GLN
1	A	381	ASN
1	B	70	ASN
1	B	204	GLN
1	B	213	ASN
1	B	239	ASN
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 ⓘ

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](#)

8 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	GOL	A	505	-	5,5,5	0.36	0	5,5,5	0.46	0
2	HEM	B	502	5,1	50,50,50	1.86	11 (22%)	67,82,82	1.46	12 (17%)
3	GOL	B	501	-	5,5,5	0.35	0	5,5,5	0.51	0
4	MES	A	504	-	12,12,12	1.92	1 (8%)	15,16,16	2.61	6 (40%)
3	GOL	A	503	-	5,5,5	0.37	0	5,5,5	0.65	0
3	GOL	B	503	-	5,5,5	0.36	0	5,5,5	0.37	0
3	GOL	A	502	-	5,5,5	0.32	0	5,5,5	0.31	0
2	HEM	A	501	5,1	50,50,50	1.98	11 (22%)	67,82,82	1.41	10 (14%)

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	GOL	A	505	-	-	2/4/4/4	-
2	HEM	B	502	5,1	-	3/14/54/54	-
3	GOL	B	501	-	-	3/4/4/4	-
4	MES	A	504	-	-	2/6/14/14	0/1/1/1
3	GOL	A	503	-	-	2/4/4/4	-
3	GOL	B	503	-	-	0/4/4/4	-
3	GOL	A	502	-	-	0/4/4/4	-
2	HEM	A	501	5,1	-	2/14/54/54	-

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	C3D-C2D	7.81	1.53	1.36
2	B	502	HEM	C3D-C2D	7.34	1.52	1.36
2	B	502	HEM	FE-ND	6.03	2.13	1.94
4	A	504	MES	C8-S	-5.92	1.69	1.77
2	A	501	HEM	FE-ND	5.75	2.12	1.94
2	A	501	HEM	FE-NA	4.71	2.10	1.95
2	A	501	HEM	FE-NB	3.39	2.05	1.94
2	B	502	HEM	FE-NA	3.31	2.06	1.95
2	A	501	HEM	CAC-C3C	3.10	1.55	1.47
2	B	502	HEM	CAC-C3C	2.79	1.54	1.47
2	A	501	HEM	CAB-C3B	2.74	1.54	1.47
2	B	502	HEM	FE-NB	2.65	2.03	1.94
2	A	501	HEM	CMD-C2D	2.56	1.56	1.50
2	B	502	HEM	CAB-C3B	2.54	1.54	1.47
2	B	502	HEM	CMB-C2B	2.39	1.55	1.50
2	A	501	HEM	FE-NC	2.38	2.03	1.95
2	B	502	HEM	CMD-C2D	2.33	1.55	1.50
2	B	502	HEM	FE-NC	2.31	2.02	1.95
2	A	501	HEM	CMB-C2B	2.30	1.55	1.50
2	A	501	HEM	CMA-C3A	2.08	1.55	1.50
2	B	502	HEM	CMA-C3A	2.07	1.55	1.50
2	A	501	HEM	CMC-C2C	2.05	1.55	1.50
2	B	502	HEM	CMC-C2C	2.04	1.55	1.50

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	504	MES	C5-N4-C3	5.24	120.13	108.84
2	A	501	HEM	C4D-ND-C1D	5.02	111.15	105.21
2	B	502	HEM	C4D-ND-C1D	4.82	110.92	105.21
4	A	504	MES	C7-N4-C3	4.82	124.09	111.24
4	A	504	MES	O3S-S-O2S	-3.78	101.95	111.40
4	A	504	MES	O2S-S-C8	3.62	112.19	106.73
2	A	501	HEM	CHD-C4C-NC	3.46	128.22	124.45
2	A	501	HEM	CBD-CAD-C3D	-2.75	104.93	112.53
4	A	504	MES	O3S-S-O1S	-2.57	104.97	111.40
2	B	502	HEM	CBA-CAA-C2A	-2.54	105.51	112.53
2	A	501	HEM	C1B-NB-C4B	2.53	108.20	105.21
2	B	502	HEM	CMD-C2D-C1D	2.47	128.89	125.03
2	A	501	HEM	CHC-C4B-NB	2.43	127.04	124.42
2	A	501	HEM	C2A-C1A-NA	-2.37	107.52	110.15
2	B	502	HEM	C1B-NB-C4B	2.35	107.99	105.21
2	B	502	HEM	CHD-C4C-NC	2.31	126.97	124.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	C1D-C2D-C3D	-2.23	104.63	106.98
2	A	501	HEM	CBA-CAA-C2A	-2.22	106.39	112.53
2	B	502	HEM	C3B-C4B-NB	-2.20	107.89	109.47
2	B	502	HEM	CHC-C1C-NC	2.16	126.81	124.45
2	B	502	HEM	O2D-CGD-CBD	2.14	120.77	114.00
2	B	502	HEM	C4C-C3C-C2C	2.14	108.67	106.81
2	B	502	HEM	CBD-CAD-C3D	-2.13	106.63	112.53
2	A	501	HEM	CHA-C4D-ND	2.08	126.94	124.37
2	B	502	HEM	C1D-C2D-C3D	-2.07	104.80	106.98
4	A	504	MES	O1-C2-C3	2.05	116.19	111.77
2	B	502	HEM	C2A-C1A-NA	-2.04	107.89	110.15
2	A	501	HEM	C3B-C4B-NB	-2.02	108.02	109.47

There are no chirality outliers.

All (14) torsion outliers are listed below:

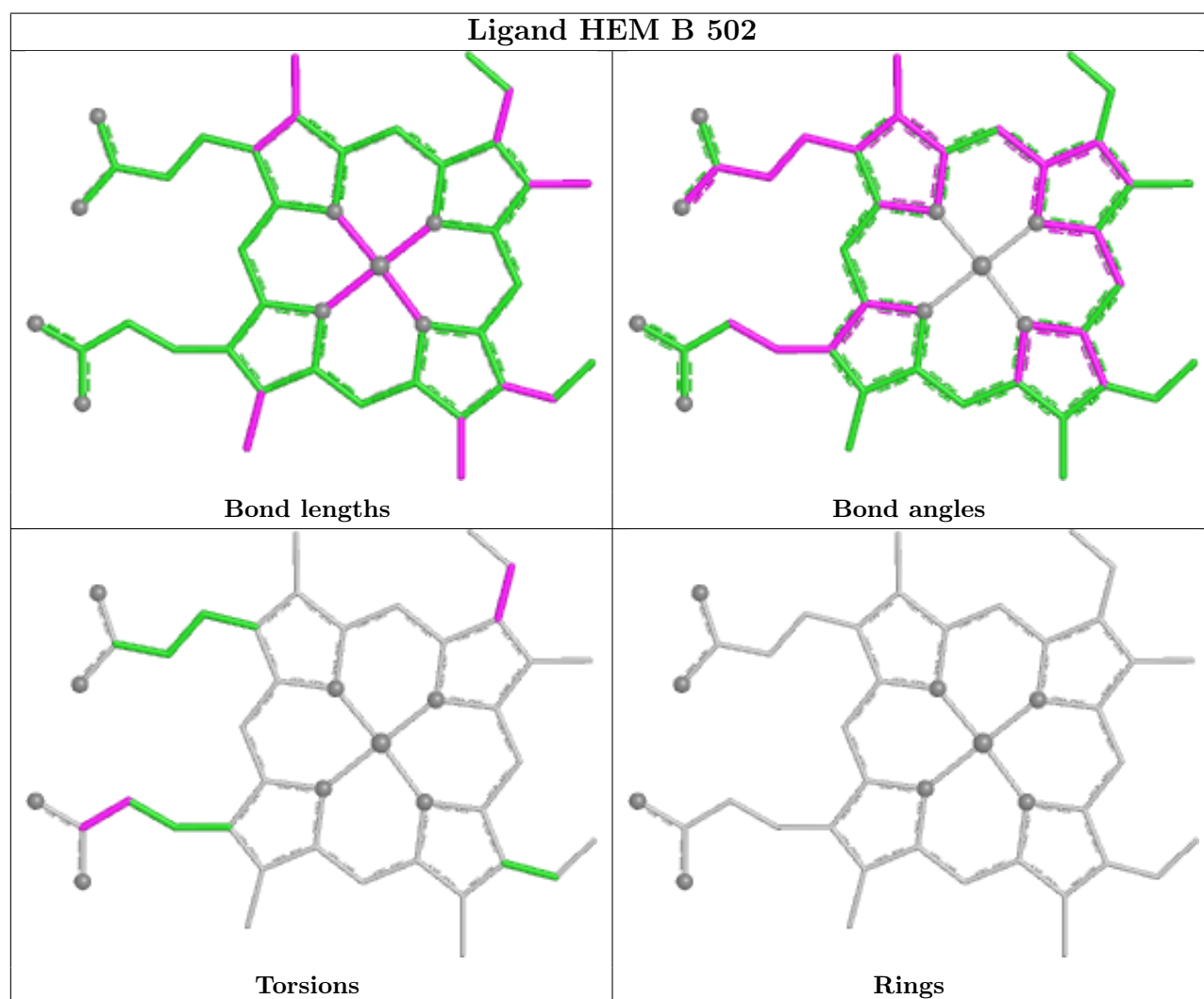
Mol	Chain	Res	Type	Atoms
3	B	501	GOL	O1-C1-C2-C3
4	A	504	MES	C8-C7-N4-C3
3	A	503	GOL	O1-C1-C2-C3
3	A	505	GOL	O1-C1-C2-C3
3	B	501	GOL	O1-C1-C2-O2
4	A	504	MES	N4-C7-C8-S
3	A	505	GOL	O1-C1-C2-O2
3	A	503	GOL	O1-C1-C2-O2
3	B	501	GOL	O2-C2-C3-O3
2	B	502	HEM	C4C-C3C-CAC-CBC
2	B	502	HEM	CAA-CBA-CGA-O2A
2	B	502	HEM	CAA-CBA-CGA-O1A
2	A	501	HEM	CAD-CBD-CGD-O1D
2	A	501	HEM	CAD-CBD-CGD-O2D

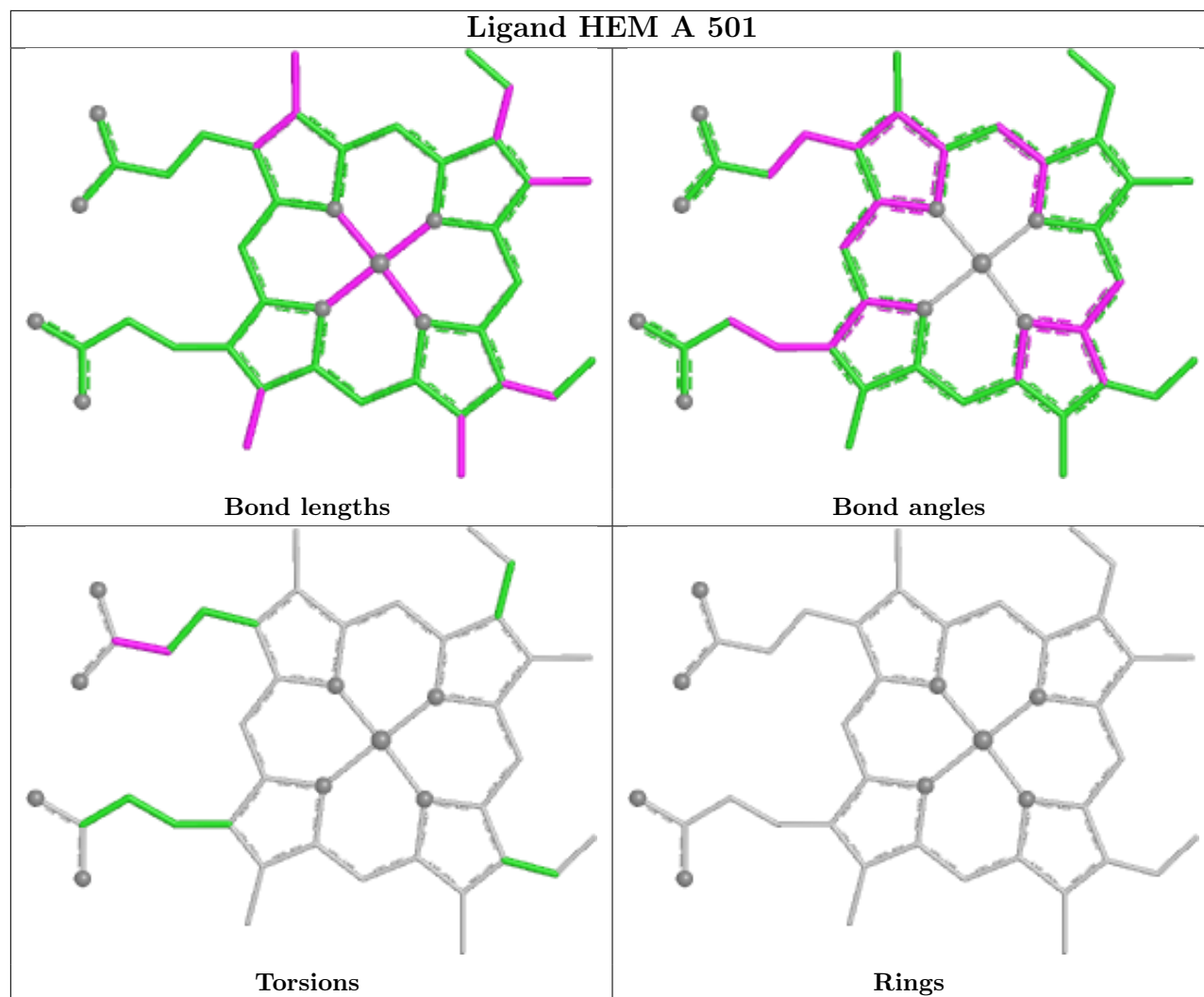
There are no ring outliers.

5 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	505	GOL	1	0
2	B	502	HEM	10	0
3	A	503	GOL	1	0
3	A	502	GOL	3	0
2	A	501	HEM	4	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	448/455 (98%)	0.00	30 (6%) 24 25	10, 22, 59, 82	15 (3%)
1	B	451/455 (99%)	0.03	32 (7%) 22 23	10, 23, 57, 95	11 (2%)
All	All	899/910 (98%)	0.02	62 (6%) 23 24	10, 23, 58, 95	26 (2%)

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	221	ALA	6.8
1	B	191	ALA	5.6
1	A	173	PHE	4.5
1	B	219	ILE	4.4
1	B	225	ALA	4.2
1	A	197	ALA	4.1
1	B	193	PRO	4.0
1	B	198	TYR	4.0
1	A	221	ALA	3.9
1	A	215	LEU	3.9
1	A	211	VAL	3.8
1	B	196	PRO	3.7
1	B	195	ASP	3.6
1	A	198	TYR	3.6
1	A	225	ALA	3.5
1	A	219	ILE	3.4
1	A	180	ALA	3.4
1	A	196	PRO	3.4
1	A	174	ILE	3.3
1	B	192	ASN	3.2
1	B	383	SER	3.1
1	A	204	GLN	3.0
1	A	209	ILE	3.0
1	A	186	ASN	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	170	PRO	2.8
1	B	170	PRO	2.8
1	B	214	ASP	2.6
1	A	227	GLY	2.6
1	B	135	ALA	2.6
1	B	189	GLN	2.5
1	B	226	SER	2.5
1	B	228	GLU	2.5
1	B	136	ASP	2.5
1	B	437	LEU	2.4
1	A	175	THR	2.4
1	A	177	MET	2.4
1	B	238	LEU	2.4
1	A	220	ILE	2.4
1	B	110	GLN	2.3
1	A	172	PRO	2.3
1	B	222	ASP	2.3
1	A	169	GLN	2.3
1	B	229	GLN	2.2
1	A	201	ASN	2.2
1	A	224	LYS	2.2
1	A	226	SER	2.2
1	B	227	GLY	2.2
1	A	205	PHE	2.2
1	A	136	ASP	2.2
1	B	211	VAL	2.2
1	A	231	ASP	2.2
1	A	218	LYS	2.2
1	A	203	ARG	2.1
1	B	169	GLN	2.1
1	B	182	ASP	2.1
1	B	231	ASP	2.1
1	B	172	PRO	2.1
1	A	188	LEU	2.1
1	B	251	ASP	2.1
1	B	190	ARG	2.1
1	B	217	ASP	2.0
1	B	174	ILE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

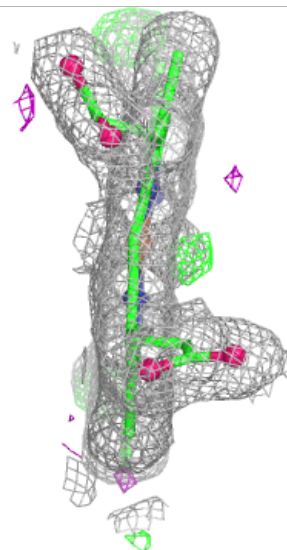
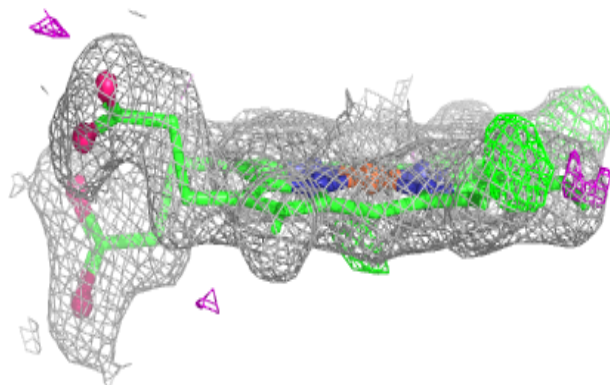
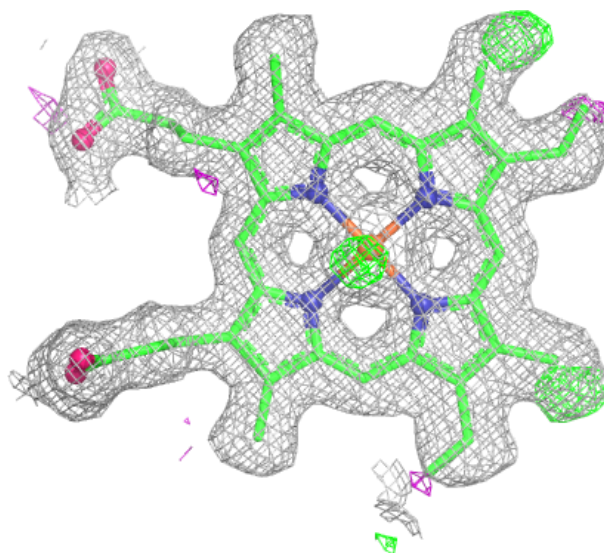
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

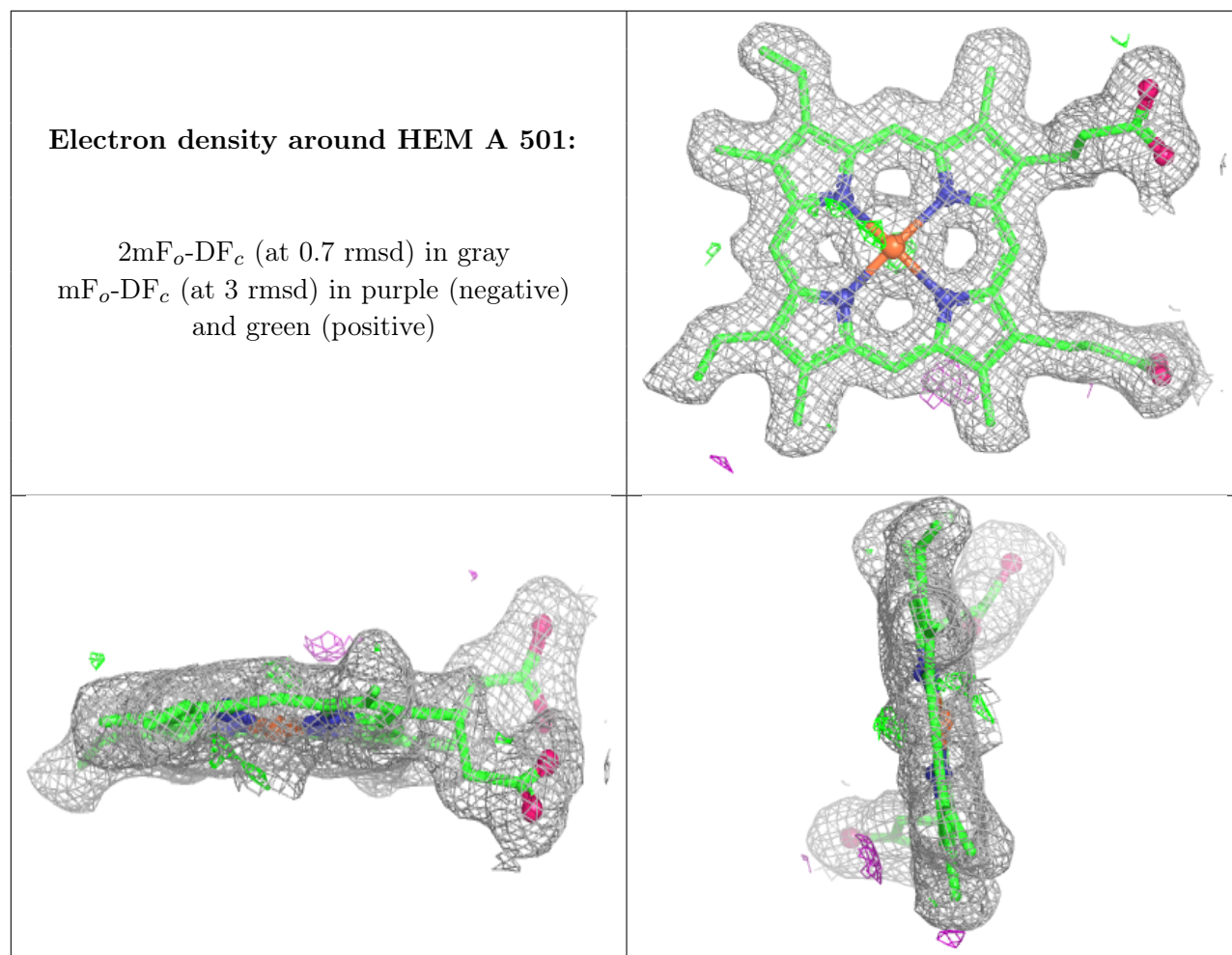
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MES	A	504	12/12	0.73	0.17	34,52,89,93	0
3	GOL	A	502	6/6	0.82	0.15	30,40,41,51	0
3	GOL	A	505	6/6	0.85	0.16	27,31,39,43	0
3	GOL	B	503	6/6	0.86	0.12	39,47,51,51	0
3	GOL	A	503	6/6	0.88	0.13	24,40,43,50	0
3	GOL	B	501	6/6	0.89	0.11	24,37,45,54	0
2	HEM	B	502	43/43	0.98	0.05	9,15,21,31	0
2	HEM	A	501	43/43	0.99	0.05	10,16,22,26	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 502:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.