



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:27 AM UTC

PDB ID : 1SMI / pdb_00001smi
Title : A single mutation of P450 BM3 induces the conformational rearrangement seen upon substrate-binding in wild-type enzyme
Authors : Joyce, M.G.; Girvan, H.M.; Munro, A.W.; Leys, D.
Deposited on : 2004-03-09
Resolution : 2.00 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

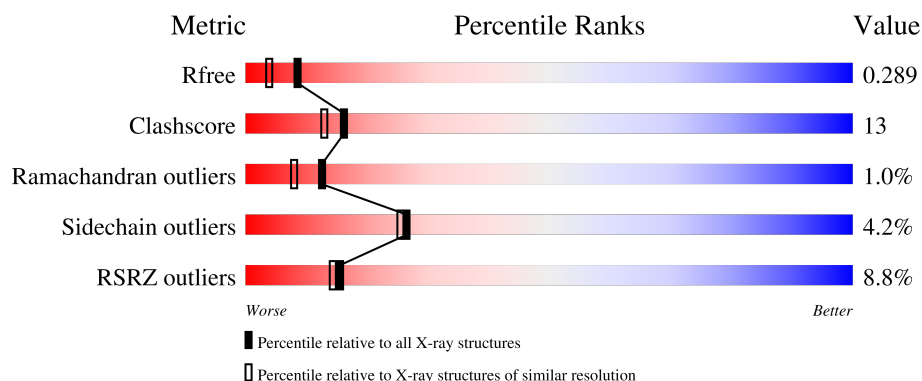
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	 6% 72% 19% • 7%
1	B	471	 11% 72% 21% • •

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7702 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.

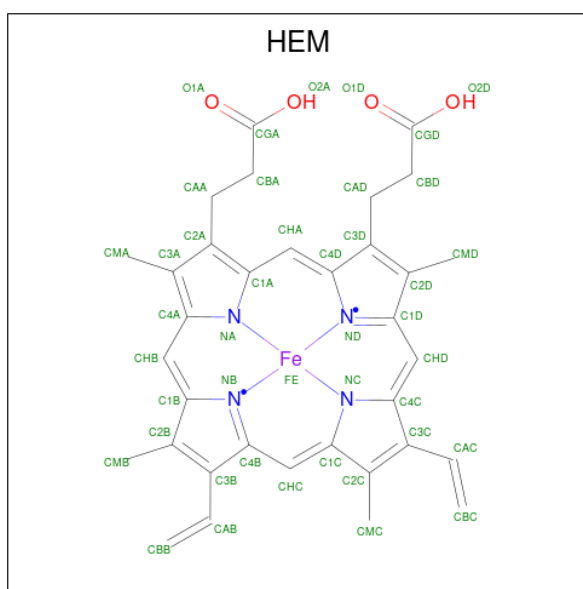
- Molecule 1 is a protein called Bifunctional P-450:NADPH-P450 reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	439	Total	C	N	O	S	0	0	0
			3500	2244	593	646	17			
1	B	455	Total	C	N	O	S	0	0	0
			3624	2310	616	681	17			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	264	GLU	ALA	engineered mutation	UNP P14779
B	264	GLU	ALA	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	0	0
			43	34	1	4	4		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

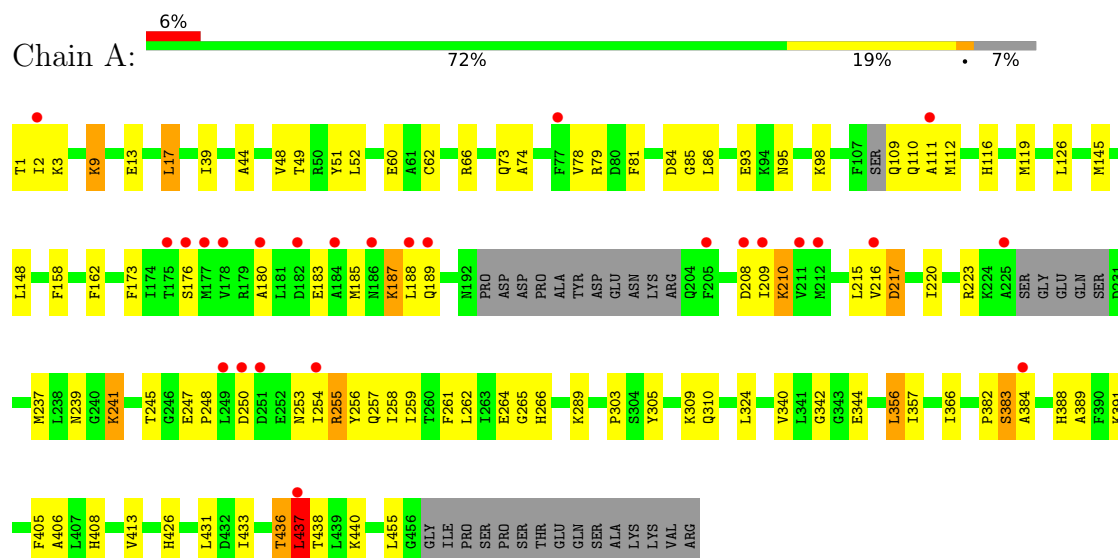
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	292	Total	O	0	0
			292	292		
3	B	200	Total	O	0	0
			200	200		

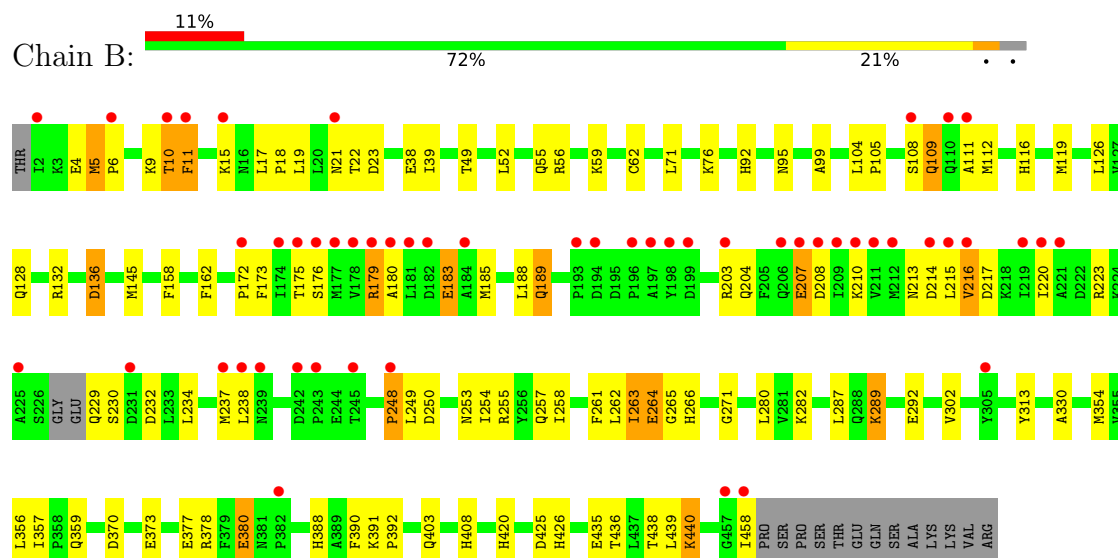
3 Residue-property plots

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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.21Å 118.93Å 146.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.92 – 2.00 14.92 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (14.92-2.00) 93.8 (14.92-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.55 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.1.9999	Depositor
R, R_{free}	0.224 , 0.288 0.224 , 0.289	Depositor DCC
R_{free} test set	3440 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	35.9	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 55.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7702	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.67% 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: 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	1.08	1/3579 (0.0%)	1.15	10/4839 (0.2%)
1	B	0.95	1/3708 (0.0%)	1.11	7/5018 (0.1%)
All	All	1.02	2/7287 (0.0%)	1.13	17/9857 (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 (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	357	ILE	CA-CB	7.29	1.58	1.54
1	A	406	ALA	N-CA	5.11	1.52	1.46

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	10	THR	N-CA-C	7.61	119.77	110.41
1	A	437	LEU	CA-C-N	7.45	134.54	121.80
1	A	437	LEU	C-N-CA	7.45	134.54	121.80
1	B	22	THR	N-CA-C	-7.13	98.24	109.07
1	A	111	ALA	N-CA-C	-7.03	103.67	111.82
1	A	436	THR	N-CA-C	-6.80	102.28	110.44
1	A	357	ILE	CB-CA-C	-6.74	107.20	114.35
1	A	255	ARG	N-CA-C	-6.53	104.16	111.28
1	A	438	THR	N-CA-C	-5.90	99.98	108.60

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	216	VAL	N-CA-C	-5.88	103.56	111.44
1	B	207	GLU	N-CA-C	-5.67	104.73	111.03
1	B	11	PHE	N-CA-C	-5.39	103.60	111.30
1	B	5	MET	N-CA-C	5.24	116.02	109.57
1	B	38	GLU	N-CA-C	5.24	117.89	111.82
1	A	324	LEU	N-CA-C	5.13	116.68	111.14
1	A	405	PHE	N-CA-C	-5.09	105.64	111.14
1	A	86	LEU	N-CA-C	5.07	117.46	111.33

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	437	LEU	Peptide

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	3500	0	3448	71	0
1	B	3624	0	3530	110	0
2	A	43	0	30	3	0
2	B	43	0	30	2	0
3	A	292	0	0	10	0
3	B	200	0	0	19	0
All	All	7702	0	7038	186	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 13.

All (186) 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:437:LEU:HA	3:A:576:HOH:O	1.33	1.23
1:B:179:ARG:NH1	1:B:180:ALA:HB2	1.59	1.16

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:173:PHE:HA	3:B:663:HOH:O	1.49	1.07
1:B:263:ILE:HG22	3:B:530:HOH:O	1.55	1.06
1:B:108:SER:O	1:B:111:ALA:HB3	1.58	1.04
1:B:172:PRO:HA	3:B:548:HOH:O	1.62	0.99
1:B:179:ARG:HD2	1:B:180:ALA:N	1.82	0.93
1:A:237:MET:HE3	1:A:257:GLN:HB2	1.52	0.91
1:B:179:ARG:HD2	1:B:180:ALA:CA	2.00	0.90
1:B:271:GLY:CA	1:B:440:LYS:HE3	2.03	0.87
1:B:21:ASN:HB3	3:B:608:HOH:O	1.74	0.86
1:B:179:ARG:CZ	1:B:208:ASP:HB3	2.08	0.82
1:B:248:PRO:HD3	3:B:586:HOH:O	1.78	0.81
1:B:179:ARG:HD2	1:B:180:ALA:HA	1.61	0.80
1:B:179:ARG:CZ	1:B:180:ALA:HB2	2.12	0.79
1:A:180:ALA:HB2	3:A:759:HOH:O	1.85	0.77
1:B:176:SER:C	1:B:179:ARG:HH21	1.93	0.77
1:A:162:PHE:HE1	1:A:215:LEU:HD11	1.51	0.75
1:B:179:ARG:CD	1:B:180:ALA:N	2.50	0.75
1:B:17:LEU:HD11	1:B:189:GLN:HA	1.69	0.74
1:B:158:PHE:CE1	1:B:258:ILE:HG12	2.21	0.74
1:B:56:ARG:HA	1:B:59:LYS:HE2	1.68	0.74
1:B:216:VAL:O	1:B:220:ILE:HD12	1.89	0.73
1:B:271:GLY:N	1:B:440:LYS:HE3	2.05	0.72
1:B:55:GLN:O	1:B:59:LYS:HG2	1.88	0.72
1:B:108:SER:O	1:B:111:ALA:CB	2.37	0.72
1:B:330:ALA:HB1	1:B:354:MET:HE1	1.72	0.71
1:B:116:HIS:HD2	1:B:408:HIS:NE2	1.89	0.69
1:A:436:THR:O	1:A:437:LEU:C	2.37	0.68
1:A:210:LYS:HE2	1:A:210:LYS:HA	1.76	0.67
1:B:112:MET:HE1	1:B:119:MET:HE1	1.76	0.67
1:B:420:HIS:NE2	3:B:474:HOH:O	2.27	0.67
1:B:11:PHE:CD1	1:B:18:PRO:HD2	2.31	0.66
1:B:11:PHE:HE1	1:B:19:LEU:HG	1.60	0.66
1:B:185:MET:HA	1:B:188:LEU:HD12	1.77	0.66
1:A:237:MET:HE1	1:A:258:ILE:CD1	2.25	0.66
1:B:207:GLU:O	1:B:210:LYS:HB2	1.94	0.66
2:B:472:HEM:HMC2	2:B:472:HEM:HBC2	1.78	0.66
1:A:210:LYS:HA	1:A:210:LYS:CE	2.26	0.66
1:B:179:ARG:NH2	1:B:208:ASP:HB3	2.09	0.66
2:A:472:HEM:HBC2	2:A:472:HEM:HMC2	1.78	0.65
1:B:435:GLU:HG2	1:B:439:LEU:CD2	2.27	0.65
1:B:207:GLU:HG3	3:B:615:HOH:O	1.97	0.64

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:253:ASN:O	1:A:256:TYR:HB2	1.97	0.64
1:B:214:ASP:O	1:B:217:ASP:HB2	1.97	0.64
1:A:85:GLY:HA2	1:A:257:GLN:NE2	2.12	0.64
1:B:92:HIS:HB3	3:B:634:HOH:O	1.98	0.62
1:B:11:PHE:CE1	1:B:19:LEU:HG	2.34	0.62
1:B:173:PHE:CA	3:B:663:HOH:O	2.22	0.61
1:A:241:LYS:HG3	1:A:248:PRO:HB3	1.82	0.61
1:B:229:GLN:HA	3:B:662:HOH:O	2.00	0.60
1:B:173:PHE:CD1	1:B:215:LEU:HD11	2.37	0.60
1:A:289:LYS:NZ	3:A:684:HOH:O	2.33	0.60
1:A:162:PHE:CE1	1:A:215:LEU:HD11	2.35	0.59
1:B:330:ALA:HB1	1:B:354:MET:CE	2.31	0.59
1:B:330:ALA:CB	1:B:354:MET:HE1	2.32	0.59
1:B:261:PHE:O	1:B:265:GLY:N	2.34	0.59
1:A:255:ARG:O	1:A:259:ILE:HD12	2.04	0.57
1:B:263:ILE:CG2	3:B:530:HOH:O	2.32	0.57
1:A:382:PRO:C	1:A:384:ALA:H	2.11	0.57
1:A:66:ARG:NH2	1:A:340:VAL:O	2.37	0.57
1:B:179:ARG:HH11	1:B:180:ALA:HB2	1.60	0.57
1:A:223:ARG:HB3	3:A:722:HOH:O	2.05	0.57
1:B:234:LEU:HD13	1:B:258:ILE:HD11	1.88	0.56
1:A:216:VAL:HG21	1:A:259:ILE:CG1	2.35	0.56
1:B:204:GLN:NE2	1:B:207:GLU:OE2	2.40	0.55
1:A:261:PHE:O	1:A:265:GLY:HA3	2.05	0.55
1:B:179:ARG:CZ	1:B:208:ASP:CB	2.84	0.55
1:B:426:HIS:HE1	3:B:577:HOH:O	1.89	0.55
1:B:213:ASN:HB3	1:B:255:ARG:HH21	1.72	0.55
1:B:377:GLU:O	1:B:380:GLU:HB2	2.07	0.55
1:B:238:LEU:HD23	1:B:254:ILE:HD13	1.87	0.54
1:A:51:TYR:HD2	1:A:356:LEU:HD13	1.71	0.54
1:B:176:SER:C	1:B:179:ARG:NH2	2.64	0.54
1:B:176:SER:O	1:B:179:ARG:NE	2.41	0.53
1:A:74:ALA:O	1:A:78:VAL:HG23	2.08	0.53
1:A:183:GLU:HA	3:A:694:HOH:O	2.08	0.53
1:B:204:GLN:NE2	1:B:208:ASP:OD1	2.42	0.53
1:B:179:ARG:NH1	1:B:180:ALA:CB	2.53	0.52
1:B:249:LEU:HD22	1:B:253:ASN:OD1	2.09	0.52
1:A:112:MET:HE1	1:A:119:MET:HE1	1.91	0.52
1:B:261:PHE:O	1:B:265:GLY:HA3	2.10	0.52
1:A:237:MET:HE2	1:A:254:ILE:HG23	1.91	0.51
1:B:282:LYS:NZ	1:B:425:ASP:OD2	2.40	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158:PHE:HE1	1:B:258:ILE:HG12	1.73	0.51
1:B:390:PHE:CZ	1:B:392:PRO:HG3	2.45	0.51
1:A:116:HIS:HE1	1:A:303:PRO:O	1.93	0.51
2:A:472:HEM:HBC2	2:A:472:HEM:CMC	2.40	0.51
1:A:62:CYS:SG	1:A:391:LYS:HE2	2.51	0.51
1:B:280:LEU:HB3	1:B:287:LEU:HD13	1.93	0.51
1:B:179:ARG:HD2	1:B:179:ARG:C	2.37	0.50
1:B:262:LEU:O	1:B:266:HIS:HD2	1.95	0.50
1:A:73:GLN:HB2	1:A:188:LEU:HD21	1.94	0.50
1:A:245:THR:OG1	1:A:247:GLU:HG2	2.11	0.50
1:B:173:PHE:HD1	1:B:215:LEU:HD21	1.77	0.49
1:B:264:GLU:HB3	3:B:496:HOH:O	2.12	0.49
1:B:289:LYS:HD2	1:B:313:TYR:CZ	2.47	0.49
1:A:3:LYS:HD2	1:A:344:GLU:OE1	2.13	0.49
1:B:62:CYS:SG	1:B:391:LYS:HE2	2.52	0.49
1:A:216:VAL:HG21	1:A:259:ILE:HG13	1.93	0.49
1:A:426:HIS:CD2	1:A:426:HIS:H	2.31	0.49
1:B:39:ILE:HD12	1:B:52:LEU:CD2	2.43	0.49
1:A:310:GLN:HG2	3:A:560:HOH:O	2.12	0.48
1:B:271:GLY:HA2	1:B:440:LYS:HE3	1.93	0.48
1:B:216:VAL:HG12	1:B:220:ILE:HD11	1.96	0.48
1:A:158:PHE:CE1	1:A:258:ILE:HD12	2.49	0.48
1:B:223:ARG:NH2	1:B:232:ASP:OD2	2.30	0.48
1:B:391:LYS:HA	3:B:524:HOH:O	2.12	0.48
1:A:17:LEU:O	1:A:17:LEU:HD12	2.13	0.48
1:A:81:PHE:HB3	1:A:209:ILE:HG12	1.96	0.48
1:A:431:LEU:HD21	1:A:433:ILE:HD11	1.96	0.47
1:A:208:ASP:HB2	3:A:759:HOH:O	2.13	0.47
1:B:220:ILE:HG23	3:B:570:HOH:O	2.14	0.47
1:A:237:MET:HE1	1:A:258:ILE:HD13	1.93	0.47
1:B:216:VAL:HG12	1:B:220:ILE:CD1	2.45	0.47
1:A:17:LEU:HD22	1:A:44:ALA:HB1	1.96	0.47
1:A:98:LYS:HE3	1:A:247:GLU:HG3	1.97	0.47
2:B:472:HEM:HBC2	2:B:472:HEM:CMC	2.43	0.47
1:B:179:ARG:CD	1:B:179:ARG:C	2.88	0.46
1:A:109:GLN:HA	1:A:112:MET:HB2	1.98	0.46
1:B:179:ARG:O	1:B:183:GLU:N	2.35	0.46
1:A:9:LYS:HE2	1:A:9:LYS:HB2	1.73	0.45
1:A:220:ILE:HA	3:A:722:HOH:O	2.16	0.45
1:B:99:ALA:HB2	1:B:249:LEU:HD21	1.99	0.45
1:B:261:PHE:O	1:B:265:GLY:CA	2.65	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:391:LYS:N	1:B:392:PRO:CD	2.80	0.45
1:A:3:LYS:CD	1:A:344:GLU:OE1	2.64	0.45
1:B:388:HIS:HD2	1:B:391:LYS:NZ	2.14	0.45
1:B:112:MET:HE1	1:B:119:MET:CE	2.43	0.45
1:A:148:LEU:HD21	1:A:413:VAL:HG21	1.98	0.44
1:A:254:ILE:O	1:A:258:ILE:HG12	2.17	0.44
1:B:217:ASP:OD1	1:B:255:ARG:NH1	2.50	0.44
1:A:185:MET:O	1:A:188:LEU:HB2	2.18	0.44
1:B:162:PHE:HE1	1:B:215:LEU:HD13	1.82	0.44
1:A:180:ALA:CB	3:A:759:HOH:O	2.56	0.44
1:B:216:VAL:O	1:B:220:ILE:CD1	2.62	0.44
1:A:366:ILE:HG21	1:A:389:ALA:HB1	2.00	0.44
1:A:183:GLU:O	1:A:187:LYS:HB2	2.18	0.43
1:A:79:ARG:NH2	1:A:93:GLU:OE2	2.40	0.43
1:A:388:HIS:HD2	1:A:391:LYS:NZ	2.15	0.43
1:B:179:ARG:NE	1:B:208:ASP:HB3	2.32	0.43
1:A:1:THR:HB	1:A:2:ILE:H	1.65	0.43
1:A:84:ASP:OD2	1:A:95:ASN:OD1	2.36	0.43
1:B:354:MET:HE3	1:B:354:MET:HB3	1.82	0.43
1:A:60:GLU:OE2	1:A:342:GLY:HA2	2.18	0.43
1:B:370:ASP:HB2	1:B:373:GLU:OE1	2.18	0.43
1:B:403:GLN:NE2	3:B:529:HOH:O	2.47	0.43
1:B:237:MET:HB3	1:B:254:ILE:HG12	2.01	0.43
1:B:257:GLN:O	1:B:261:PHE:CD1	2.72	0.43
1:A:162:PHE:HE1	1:A:215:LEU:CD1	2.27	0.43
1:A:39:ILE:HD12	1:A:52:LEU:HD21	2.00	0.43
1:B:126:LEU:HD21	1:B:145:MET:HE1	1.99	0.43
1:A:262:LEU:O	1:A:266:HIS:ND1	2.50	0.43
2:A:472:HEM:CMB	2:A:472:HEM:HBB2	2.49	0.43
1:A:116:HIS:HD2	1:A:408:HIS:NE2	2.17	0.42
1:A:305:TYR:O	1:A:309:LYS:HG2	2.19	0.42
1:B:71:LEU:O	1:B:76:LYS:HE3	2.20	0.42
1:B:237:MET:SD	1:B:254:ILE:HG23	2.59	0.42
1:B:253:ASN:O	1:B:257:GLN:HG2	2.20	0.42
1:B:257:GLN:O	1:B:261:PHE:HD1	2.02	0.42
1:B:271:GLY:HA2	1:B:440:LYS:HG3	2.02	0.42
1:A:126:LEU:HD21	1:A:145:MET:HE1	2.01	0.42
1:A:237:MET:HE1	1:A:258:ILE:HG12	2.01	0.42
1:A:239:ASN:HB2	3:A:688:HOH:O	2.20	0.42
1:B:175:THR:HB	3:B:548:HOH:O	2.19	0.42
1:B:213:ASN:O	1:B:255:ARG:NH2	2.52	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:436:THR:C	1:B:438:THR:H	2.28	0.42
1:A:73:GLN:CB	1:A:188:LEU:HD21	2.50	0.42
1:A:237:MET:O	1:A:254:ILE:HD11	2.20	0.42
1:B:458:ILE:HG23	3:B:650:HOH:O	2.20	0.42
1:A:257:GLN:O	1:A:261:PHE:HD1	2.03	0.41
1:B:302:VAL:HG23	3:B:508:HOH:O	2.19	0.41
1:B:5:MET:HA	1:B:6:PRO:HD3	1.80	0.41
1:B:95:ASN:HD22	1:B:95:ASN:HA	1.61	0.41
1:B:176:SER:HB3	1:B:179:ARG:HH21	1.86	0.41
1:B:356:LEU:HD23	1:B:359:GLN:HG3	2.01	0.41
1:A:342:GLY:O	1:A:344:GLU:HG3	2.21	0.41
1:B:128:GLN:O	1:B:132:ARG:HG3	2.21	0.41
1:A:382:PRO:C	1:A:384:ALA:N	2.78	0.41
1:B:109:GLN:C	1:B:111:ALA:N	2.76	0.41
1:A:217:ASP:OD1	1:A:255:ARG:NH1	2.46	0.40
1:A:173:PHE:HB2	1:A:215:LEU:HD13	2.03	0.40
1:B:439:LEU:O	1:B:440:LYS:NZ	2.46	0.40
1:B:104:LEU:N	1:B:105:PRO:HD2	2.36	0.40
1:A:183:GLU:O	1:A:187:LYS:NZ	2.55	0.40
1:B:10:THR:HA	1:B:15:LYS:O	2.21	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	431/471 (92%)	412 (96%)	17 (4%)	2 (0%)	24	21
1	B	451/471 (96%)	419 (93%)	25 (6%)	7 (2%)	7	3
All	All	882/942 (94%)	831 (94%)	42 (5%)	9 (1%)	12	8

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	250	ASP
1	A	110	GLN
1	A	383	SER
1	B	4	GLU
1	B	263	ILE
1	B	9	LYS
1	B	378	ARG
1	B	136	ASP
1	B	248	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	374/413 (91%)	356 (95%)	18 (5%)	23	21
1	B	387/413 (94%)	373 (96%)	14 (4%)	31	31
All	All	761/826 (92%)	729 (96%)	32 (4%)	26	25

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	LYS
1	A	13	GLU
1	A	17	LEU
1	A	48	VAL
1	A	49	THR
1	A	176	SER
1	A	187	LYS
1	A	189	GLN
1	A	210	LYS
1	A	217	ASP
1	A	241	LYS
1	A	250	ASP
1	A	264	GLU
1	A	356	LEU
1	A	383	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	437	LEU
1	A	440	LYS
1	A	455	LEU
1	B	23	ASP
1	B	49	THR
1	B	109	GLN
1	B	136	ASP
1	B	179	ARG
1	B	183	GLU
1	B	189	GLN
1	B	203	ARG
1	B	230	SER
1	B	264	GLU
1	B	289	LYS
1	B	292	GLU
1	B	380	GLU
1	B	440	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	92	HIS
1	A	95	ASN
1	A	116	HIS
1	A	186	ASN
1	A	239	ASN
1	A	359	GLN
1	A	381	ASN
1	A	388	HIS
1	A	403	GLN
1	A	426	HIS
1	B	70	ASN
1	B	95	ASN
1	B	101	ASN
1	B	116	HIS
1	B	204	GLN
1	B	388	HIS
1	B	403	GLN
1	B	426	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	$\# Z > 2$
2	HEM	B	472	1,3	50,50,50	1.92	10 (20%)	67,82,82	1.55	14 (20%)
2	HEM	A	472	1	50,50,50	2.03	13 (26%)	67,82,82	1.71	9 (13%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	B	472	1,3	-	2/14/54/54	-
2	HEM	A	472	1	-	2/14/54/54	-

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	472	HEM	C3D-C2D	7.22	1.52	1.36

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	472	HEM	FE-NB	5.95	2.13	1.94
2	A	472	HEM	C3D-C2D	5.91	1.49	1.36
2	A	472	HEM	FE-ND	5.85	2.12	1.94
2	B	472	HEM	FE-NC	4.41	2.09	1.95
2	B	472	HEM	FE-NA	3.83	2.07	1.95
2	A	472	HEM	CMB-C2B	3.47	1.57	1.50
2	A	472	HEM	CAC-C3C	3.32	1.56	1.47
2	B	472	HEM	CAB-C3B	3.02	1.55	1.47
2	B	472	HEM	CMA-C3A	2.80	1.56	1.50
2	B	472	HEM	CMD-C2D	2.69	1.56	1.50
2	A	472	HEM	O2D-CGD	-2.41	1.22	1.30
2	A	472	HEM	CMA-C3A	2.37	1.55	1.50
2	A	472	HEM	CMD-C2D	2.28	1.55	1.50
2	A	472	HEM	C4D-ND	-2.27	1.36	1.40
2	A	472	HEM	C4B-NB	-2.25	1.34	1.38
2	A	472	HEM	FE-NA	2.24	2.02	1.95
2	B	472	HEM	CAC-C3C	2.23	1.53	1.47
2	B	472	HEM	C2A-C3A	-2.23	1.33	1.38
2	B	472	HEM	C1C-NC	-2.16	1.35	1.39
2	A	472	HEM	CMC-C2C	2.13	1.55	1.50
2	A	472	HEM	CAB-C3B	2.08	1.52	1.47
2	B	472	HEM	CMB-C2B	2.05	1.55	1.50

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	472	HEM	C4D-ND-C1D	5.60	111.83	105.21
2	A	472	HEM	C1B-NB-C4B	5.59	111.83	105.21
2	B	472	HEM	CAD-C3D-C4D	3.79	131.31	124.70
2	A	472	HEM	C3B-C4B-NB	-3.73	106.79	109.47
2	A	472	HEM	CHC-C4B-NB	3.54	128.23	124.42
2	A	472	HEM	C2B-C1B-NB	-3.39	105.94	109.84
2	B	472	HEM	CMD-C2D-C1D	3.23	130.09	125.03
2	A	472	HEM	CBD-CAD-C3D	-3.09	103.99	112.53
2	B	472	HEM	CMA-C3A-C4A	-2.97	120.90	125.42
2	A	472	HEM	CHD-C1D-ND	2.78	127.42	124.42
2	B	472	HEM	CHC-C4B-NB	2.77	127.40	124.42
2	B	472	HEM	CAD-C3D-C2D	-2.61	122.98	127.87
2	B	472	HEM	O2A-CGA-CBA	2.60	122.21	114.00
2	B	472	HEM	CMA-C3A-C2A	2.57	131.08	125.62
2	A	472	HEM	CHC-C1C-NC	-2.53	121.70	124.45
2	B	472	HEM	C3B-C2B-C1B	2.45	108.25	106.41

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	472	HEM	C2A-C1A-NA	-2.43	107.46	110.15
2	B	472	HEM	C4A-NA-C1A	2.37	109.68	105.82
2	B	472	HEM	CBD-CAD-C3D	-2.32	106.11	112.53
2	B	472	HEM	CHD-C1D-ND	2.11	126.70	124.42
2	A	472	HEM	C1D-C2D-C3D	-2.10	104.77	106.98
2	B	472	HEM	CHB-C4A-NA	2.08	127.64	123.86
2	B	472	HEM	C4D-ND-C1D	2.08	107.67	105.21

There are no chirality outliers.

All (4) torsion outliers are listed below:

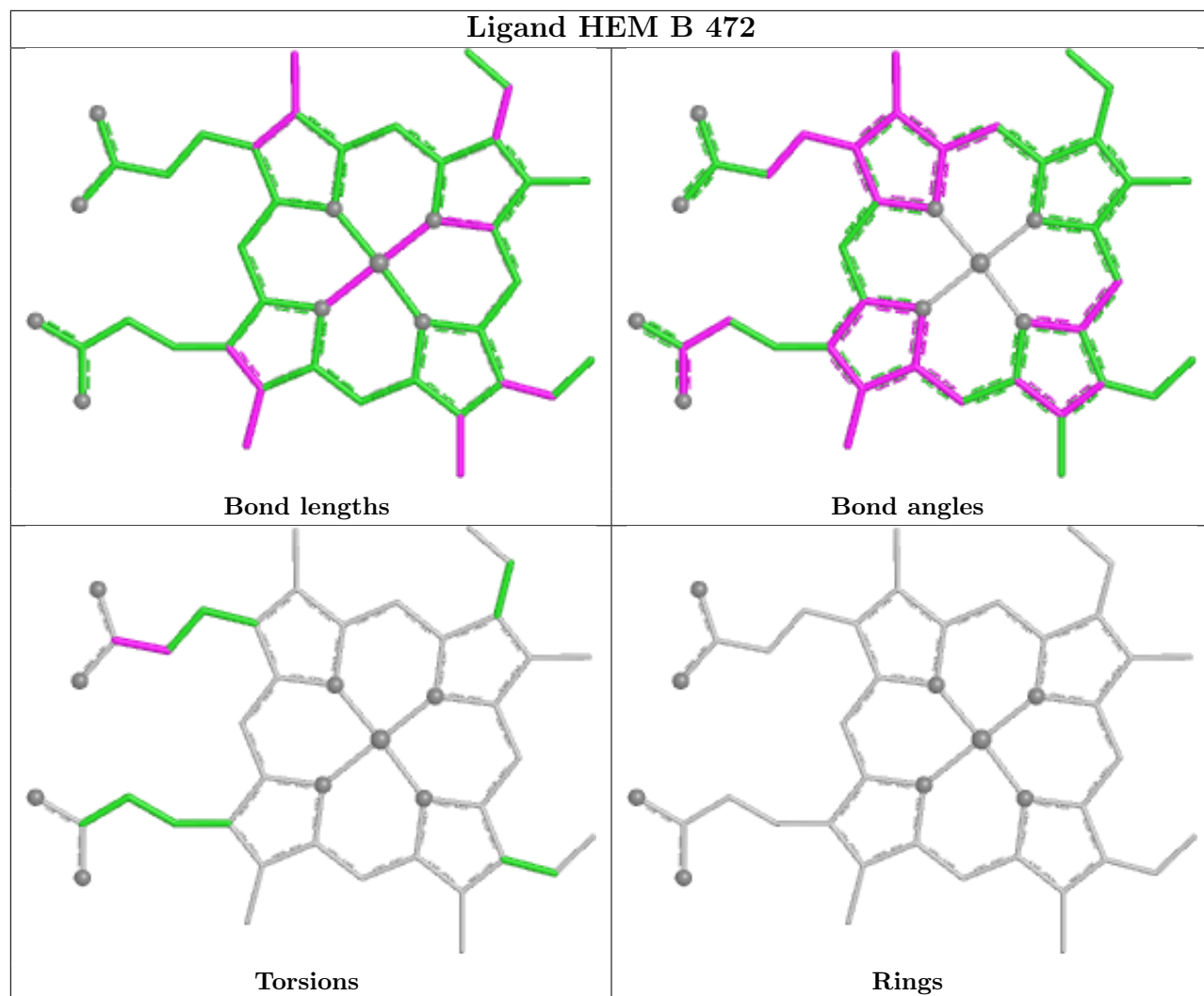
Mol	Chain	Res	Type	Atoms
2	A	472	HEM	CAD-CBD-CGD-O1D
2	A	472	HEM	CAD-CBD-CGD-O2D
2	B	472	HEM	CAD-CBD-CGD-O2D
2	B	472	HEM	CAD-CBD-CGD-O1D

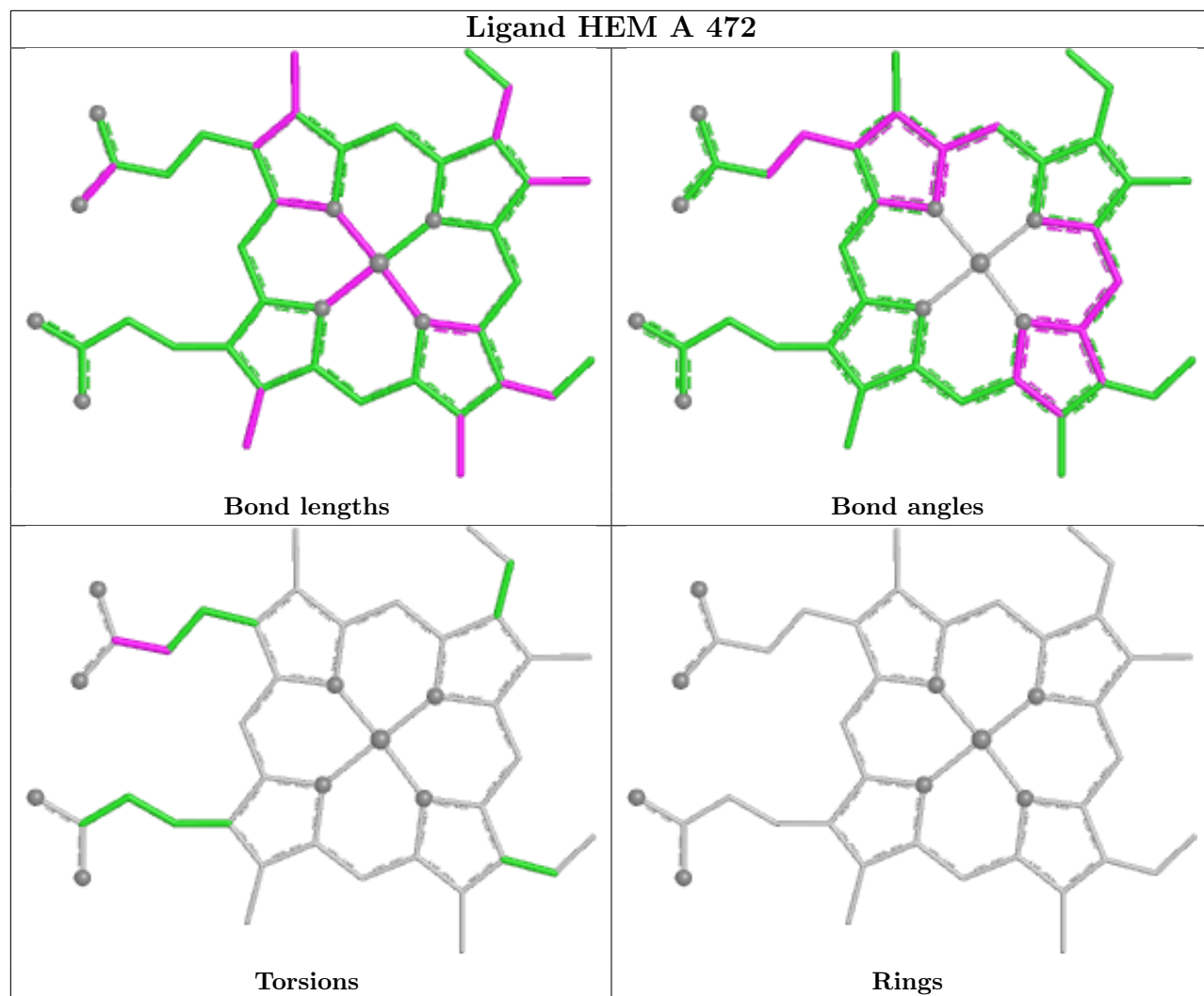
There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	472	HEM	2	0
2	A	472	HEM	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [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	439/471 (93%)	0.22	26 (5%) 28 27	24, 39, 66, 76	0
1	B	455/471 (96%)	0.70	53 (11%) 9 8	32, 49, 75, 82	0
All	All	894/942 (94%)	0.47	79 (8%) 15 14	24, 44, 72, 82	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	211	VAL	5.2
1	A	211	VAL	4.9
1	A	209	ILE	4.6
1	A	225	ALA	4.1
1	A	212	MET	4.1
1	B	209	ILE	4.0
1	B	181	LEU	3.9
1	B	458	ILE	3.7
1	B	242	ASP	3.7
1	A	250	ASP	3.7
1	B	214	ASP	3.5
1	A	184	ALA	3.5
1	B	208	ASP	3.5
1	B	174	ILE	3.5
1	A	205	PHE	3.5
1	B	11	PHE	3.4
1	A	254	ILE	3.3
1	B	176	SER	3.2
1	A	384	ALA	3.1
1	A	180	ALA	3.0
1	B	180	ALA	3.0
1	B	198	TYR	2.9
1	B	221	ALA	2.9
1	B	457	GLY	2.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	110	GLN	2.9
1	B	220	ILE	2.9
1	B	184	ALA	2.9
1	A	2	ILE	2.8
1	B	2	ILE	2.8
1	B	219	ILE	2.8
1	B	245	THR	2.8
1	B	177	MET	2.7
1	B	203	ARG	2.7
1	B	10	THR	2.7
1	B	182	ASP	2.7
1	B	111	ALA	2.7
1	B	212	MET	2.7
1	A	176	SER	2.7
1	B	179	ARG	2.7
1	A	251	ASP	2.6
1	B	15	LYS	2.6
1	B	206	GLN	2.6
1	B	215	LEU	2.6
1	A	208	ASP	2.6
1	A	437	LEU	2.5
1	A	178	VAL	2.5
1	B	194	ASP	2.5
1	B	6	PRO	2.5
1	B	238	LEU	2.5
1	B	199	ASP	2.4
1	B	243	PRO	2.4
1	B	197	ALA	2.4
1	B	108	SER	2.4
1	A	216	VAL	2.4
1	B	305	TYR	2.3
1	B	210	LYS	2.3
1	B	231	ASP	2.3
1	B	216	VAL	2.3
1	A	175	THR	2.3
1	B	21	ASN	2.2
1	B	196	PRO	2.2
1	B	248	PRO	2.2
1	B	239	ASN	2.2
1	B	172	PRO	2.2
1	B	178	VAL	2.2
1	B	237	MET	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	189	GLN	2.1
1	A	177	MET	2.1
1	A	182	ASP	2.1
1	A	188	LEU	2.1
1	A	249	LEU	2.1
1	A	186	ASN	2.1
1	A	77	PHE	2.1
1	B	225	ALA	2.1
1	B	207	GLU	2.0
1	B	175	THR	2.0
1	B	382	PRO	2.0
1	A	111	ALA	2.0
1	B	193	PRO	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](#)

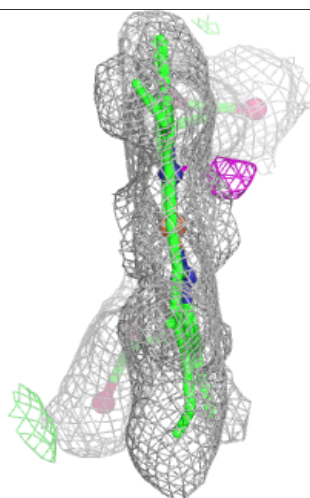
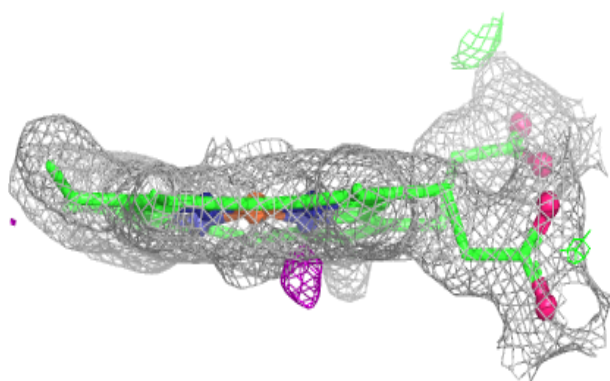
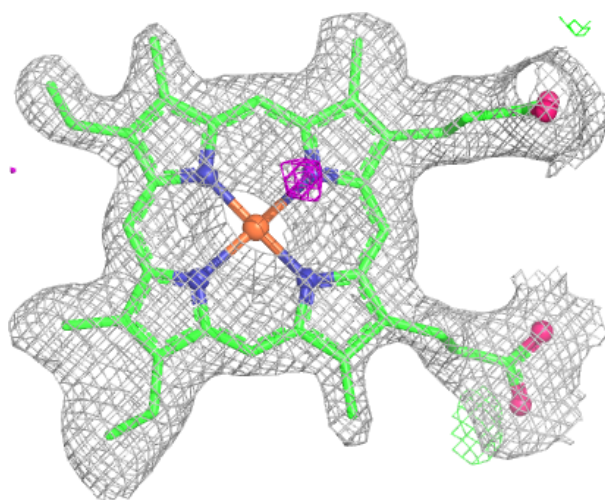
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

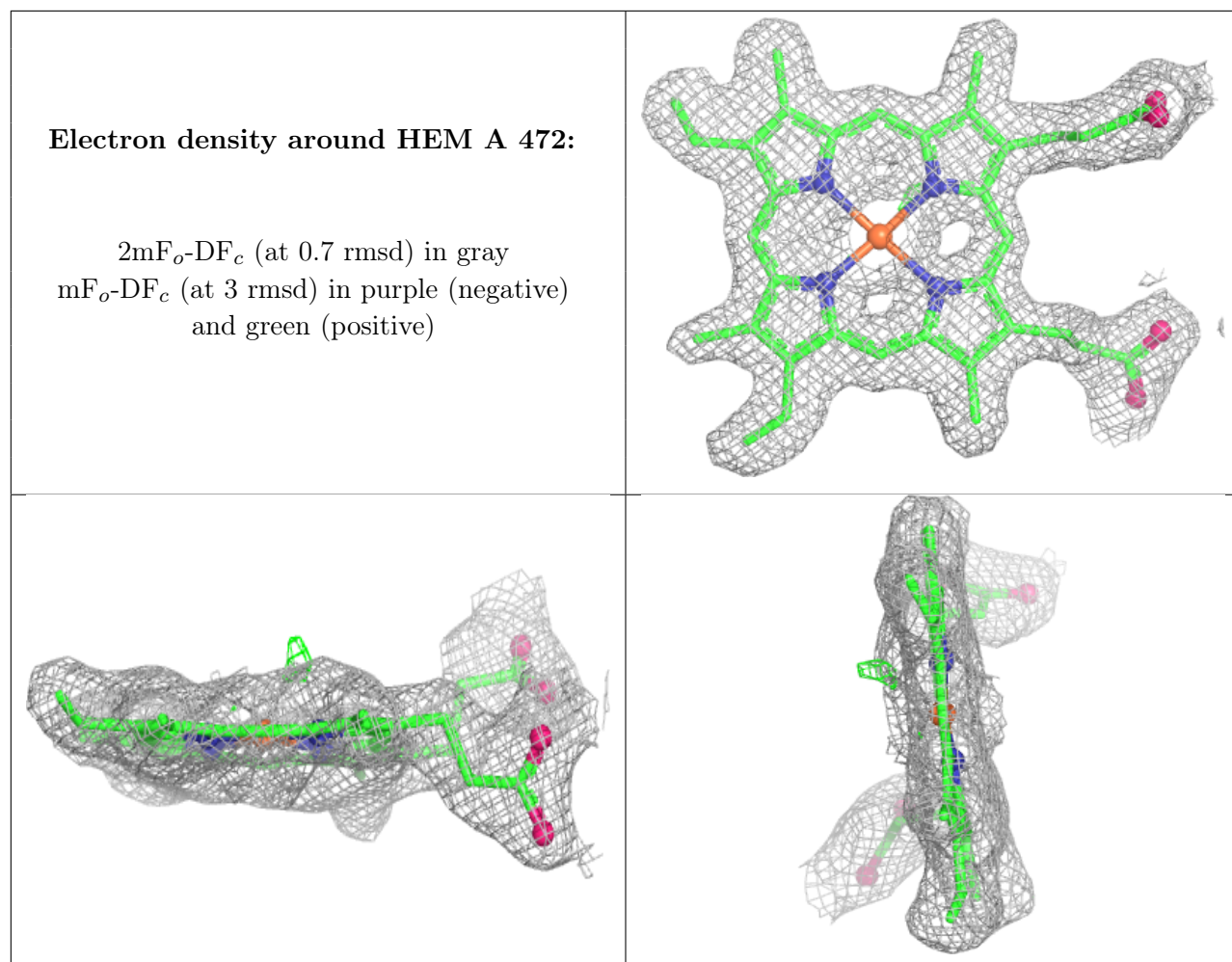
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	HEM	B	472	43/43	0.97	0.07	29,34,37,41	0
2	HEM	A	472	43/43	0.98	0.06	21,27,29,36	0

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 472:

$2mF_o-DF_c$ (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.