



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:15 PM UTC

PDB ID : 1JME / pdb_00001jme
Title : Crystal Structure of Phe393His Cytochrome P450 BM3
Authors : Ost, T.W.B.; Munro, A.W.; Mowat, C.G.; Pesseguiro, A.; Fulco, A.J.; Cho, A.K.; Cheesman, M.A.; Walkinshaw, M.D.; Chapman, S.K.
Deposited on : 2001-07-18
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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

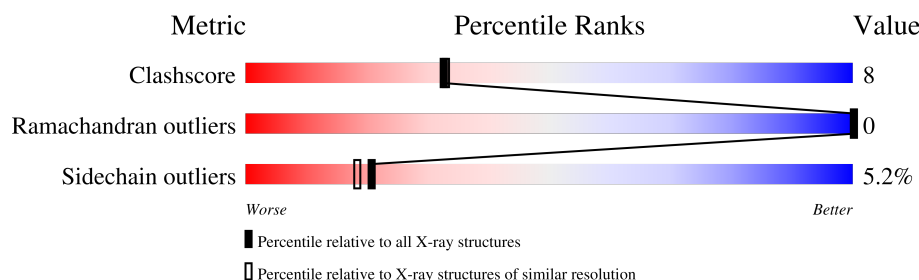
1 Overall quality at a glance

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))
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	455	 80%14% . .
1	B	455	 81%14% . .

2 Entry composition ⓘ

There are 3 unique types of molecules in this entry. The entry contains 8278 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.

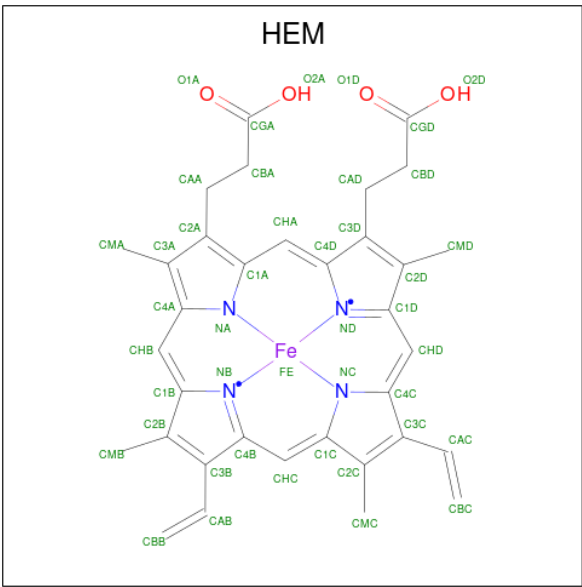
- Molecule 1 is a protein called BIFUNCTIONAL P-450:NADPH-P450 REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	441	Total	C	N	O	S	0	0	0
			3496	2241	595	643	17			
1	B	441	Total	C	N	O	S	0	0	0
			3514	2252	596	649	17			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	393	HIS	PHE	engineered mutation	UNP P14779
B	393	HIS	PHE	engineered mutation	UNP P14779

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: C₃₄H₃₂FeN₄O₄).



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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	569	Total	O	0	0
			569	569		
3	B	613	Total	O	0	0
			613	613		

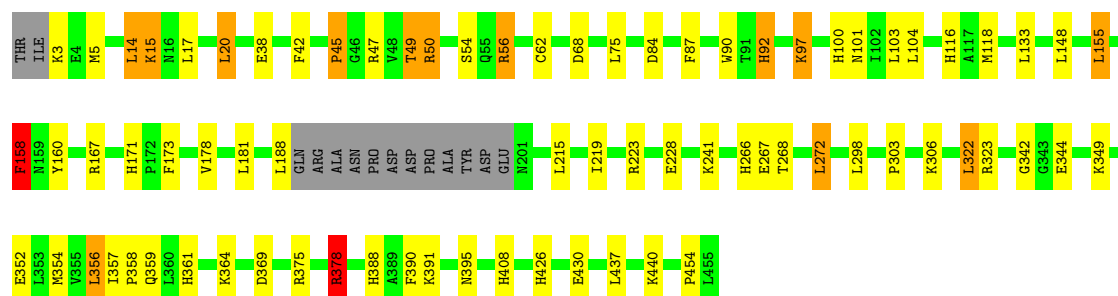
3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

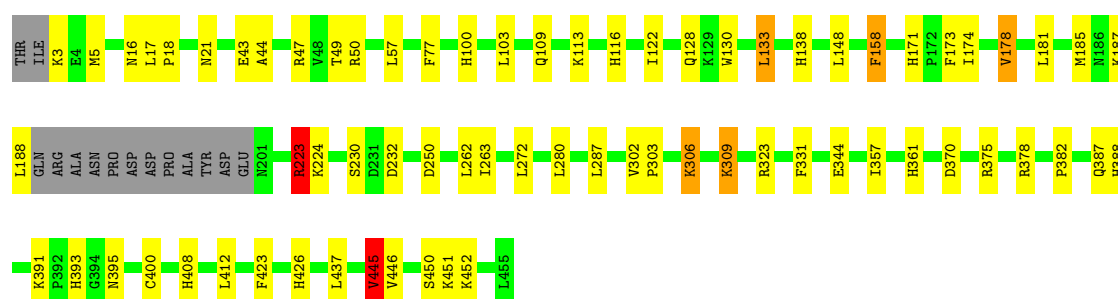
• Molecule 1: BIFUNCTIONAL P-450:NADPH-P450 REDUCTASE

Chain A: 



• Molecule 1: BIFUNCTIONAL P-450:NADPH-P450 REDUCTASE

Chain B: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.84Å 153.10Å 61.48Å 90.00° 94.67° 90.00°	Depositor
Resolution (Å)	40.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (40.00-2.00)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.179 , 0.231	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	8278	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

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	0.43	0/3576	1.25	18/4841 (0.4%)
1	B	0.43	0/3595	1.25	10/4863 (0.2%)
All	All	0.43	0/7171	1.25	28/9704 (0.3%)

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	167	ARG	CG-CD-NE	8.18	130.00	112.00
1	A	87	PHE	CA-CB-CG	7.67	121.47	113.80
1	B	223	ARG	CD-NE-CZ	7.40	134.76	124.40
1	B	47	ARG	N-CA-C	6.93	118.31	108.74
1	B	393	HIS	CA-CB-CG	-6.57	107.23	113.80
1	B	250	ASP	CA-CB-CG	6.54	119.14	112.60
1	A	357	ILE	CB-CA-C	-6.20	107.78	114.35
1	A	84	ASP	CA-CB-CG	6.03	118.63	112.60
1	A	378	ARG	CA-CB-CG	5.97	126.04	114.10
1	B	445	VAL	N-CA-CB	-5.91	101.75	111.45
1	B	331	PHE	CA-CB-CG	5.75	119.55	113.80
1	A	158	PHE	CA-CB-CG	5.73	119.53	113.80
1	B	44	ALA	O-C-N	5.71	126.00	121.55
1	A	56	ARG	CD-NE-CZ	5.69	132.36	124.40
1	A	68	ASP	CA-CB-CG	5.58	118.18	112.60
1	A	388	HIS	CA-CB-CG	-5.54	108.25	113.80
1	A	454	PRO	CA-C-N	5.51	131.62	121.70
1	A	454	PRO	C-N-CA	5.51	131.62	121.70
1	B	357	ILE	CB-CA-C	-5.40	108.63	114.35
1	A	181	LEU	CA-C-N	5.20	127.20	120.44
1	A	181	LEU	C-N-CA	5.20	127.20	120.44
1	A	47	ARG	N-CA-C	5.17	117.04	109.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	50	ARG	NE-CZ-NH2	5.15	123.83	119.20
1	B	445	VAL	CB-CA-C	5.08	119.50	110.71
1	A	15	LYS	CB-CA-C	-5.04	110.78	116.63
1	A	45	PRO	CA-C-N	-5.02	113.79	122.83
1	A	45	PRO	C-N-CA	-5.02	113.79	122.83
1	B	158	PHE	CA-CB-CG	5.01	118.81	113.80

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	3496	0	3423	54	0
1	B	3514	0	3471	57	0
2	A	43	0	30	1	0
2	B	43	0	30	1	0
3	A	569	0	0	12	0
3	B	613	0	0	16	0
All	All	8278	0	6954	111	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 8.

All (111) 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:B:280:LEU:HB3	1:B:287:LEU:HD23	1.57	0.85
1:B:16:ASN:HD22	1:B:43:GLU:H	1.22	0.85
1:B:116:HIS:HD2	1:B:408:HIS:HE2	1.24	0.83
1:A:356:LEU:HD13	1:A:358:PRO:HD2	1.65	0.78
1:A:171:HIS:HD2	1:A:173:PHE:H	1.32	0.78
1:A:97:LYS:HB2	3:A:926:HOH:O	1.84	0.77
1:B:174:ILE:O	1:B:178:VAL:HG13	1.84	0.76
1:B:171:HIS:HD2	1:B:173:PHE:H	1.33	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:361:HIS:HE1	1:B:391:LYS:H	1.34	0.74
1:B:128:GLN:HG2	3:B:1060:HOH:O	1.89	0.73
1:A:272:LEU:HG	1:A:322:LEU:HD13	1.72	0.70
1:A:361:HIS:HE1	1:A:391:LYS:H	1.39	0.70
1:B:57:LEU:HD21	3:B:984:HOH:O	1.91	0.69
1:A:375:ARG:O	1:A:378:ARG:HG2	1.94	0.67
1:A:158:PHE:HE2	1:A:219:ILE:HD13	1.61	0.66
1:B:382:PRO:HD2	3:B:790:HOH:O	1.97	0.65
1:B:223:ARG:HG2	1:B:223:ARG:HH11	1.63	0.63
1:A:356:LEU:HD12	1:A:359:GLN:HG2	1.81	0.63
1:B:116:HIS:CD2	1:B:408:HIS:HE2	2.14	0.62
1:A:158:PHE:CE2	1:A:219:ILE:HD13	2.35	0.61
1:B:130:TRP:HA	1:B:133:LEU:HD22	1.83	0.61
1:A:241:LYS:HD2	3:A:802:HOH:O	2.01	0.59
1:B:450:SER:O	1:B:452:LYS:HE2	2.03	0.59
1:A:97:LYS:HE2	1:A:101:ASN:OD1	2.03	0.58
1:B:100:HIS:HD2	3:B:710:HOH:O	1.86	0.58
1:B:21:ASN:HB2	3:B:741:HOH:O	2.03	0.58
1:A:352:GLU:OE2	1:A:354:MET:HE1	2.04	0.58
1:B:223:ARG:HD2	3:B:1032:HOH:O	2.03	0.58
1:B:100:HIS:HE1	3:B:1014:HOH:O	1.87	0.58
1:A:352:GLU:HB2	3:A:806:HOH:O	2.04	0.57
1:B:451:LYS:C	1:B:452:LYS:HE3	2.30	0.57
1:B:5:MET:HE1	1:B:50:ARG:HD3	1.88	0.56
1:B:391:LYS:HZ3	1:B:395:ASN:HD22	1.51	0.56
1:A:100:HIS:HD2	3:A:686:HOH:O	1.90	0.55
1:A:223:ARG:HD3	1:A:228:GLU:HG2	1.87	0.55
1:B:109:GLN:HE22	1:B:309:LYS:NZ	2.04	0.55
1:B:223:ARG:HH22	1:B:230:SER:HB2	1.72	0.55
1:B:17:LEU:HB3	1:B:18:PRO:HD3	1.90	0.53
1:B:185:MET:HE1	1:B:437:LEU:HD13	1.90	0.53
1:B:185:MET:CE	1:B:437:LEU:HD13	2.39	0.53
1:B:423:PHE:HD1	3:B:1055:HOH:O	1.91	0.53
1:A:267:GLU:HG2	3:A:970:HOH:O	2.09	0.52
1:B:280:LEU:CB	1:B:287:LEU:HD23	2.34	0.52
1:A:171:HIS:CD2	1:A:173:PHE:H	2.19	0.52
1:B:116:HIS:HE1	1:B:303:PRO:O	1.93	0.52
1:B:171:HIS:CD2	1:B:173:PHE:H	2.21	0.51
1:B:122:ILE:HG22	1:B:148:LEU:HD22	1.92	0.51
1:B:16:ASN:ND2	1:B:43:GLU:H	1.99	0.50
1:A:116:HIS:HD2	1:A:408:HIS:NE2	2.08	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:364:LYS:HD3	1:A:369:ASP:HA	1.93	0.50
1:A:342:GLY:O	1:A:344:GLU:HG3	2.12	0.49
1:A:178:VAL:HG22	3:A:792:HOH:O	2.11	0.49
1:A:266:HIS:CE1	1:A:267:GLU:HG3	2.48	0.49
1:B:323:ARG:HA	1:B:361:HIS:CD2	2.48	0.48
1:A:5:MET:SD	1:A:50:ARG:HG2	2.53	0.48
1:A:20:LEU:HD22	1:A:42:PHE:HZ	1.77	0.48
1:A:118:MET:SD	1:A:155:LEU:HD13	2.53	0.48
1:B:109:GLN:HE22	1:B:309:LYS:HZ2	1.61	0.48
1:A:391:LYS:NZ	1:A:395:ASN:HD22	2.12	0.48
1:B:287:LEU:HD11	3:B:1055:HOH:O	2.13	0.48
1:A:356:LEU:HD12	1:A:359:GLN:CG	2.45	0.47
1:A:306:LYS:HG3	3:A:568:HOH:O	2.15	0.46
1:B:138:HIS:CD2	1:B:445:VAL:HG22	2.50	0.46
1:A:49:THR:HG23	1:A:352:GLU:HB3	1.97	0.46
1:A:361:HIS:CE1	1:A:390:PHE:HA	2.50	0.46
1:A:364:LYS:NZ	3:A:1004:HOH:O	2.48	0.46
1:B:375:ARG:O	1:B:378:ARG:HG3	2.15	0.46
1:B:391:LYS:NZ	1:B:395:ASN:HD22	2.12	0.46
1:B:426:HIS:HE1	3:B:676:HOH:O	1.98	0.46
1:A:437:LEU:HD23	3:A:1027:HOH:O	2.16	0.45
1:B:223:ARG:NH2	1:B:232:ASP:OD2	2.49	0.45
1:A:3:LYS:HA	1:A:3:LYS:HD3	1.73	0.45
1:A:298:LEU:HD22	1:A:303:PRO:HB3	1.99	0.45
1:A:426:HIS:HE1	3:A:690:HOH:O	2.00	0.45
1:B:50:ARG:NH1	3:B:643:HOH:O	2.50	0.45
1:B:50:ARG:NE	3:B:727:HOH:O	2.48	0.45
1:B:77:PHE:CE2	1:B:187:LYS:HB3	2.51	0.45
1:B:113:LYS:NZ	3:B:974:HOH:O	2.50	0.44
1:A:272:LEU:HG	1:A:322:LEU:CD1	2.43	0.44
1:B:370:ASP:OD2	1:B:375:ARG:NH2	2.49	0.44
1:A:215:LEU:O	1:A:219:ILE:HD12	2.18	0.44
1:A:20:LEU:HD22	1:A:42:PHE:CZ	2.52	0.44
1:A:116:HIS:HE1	1:A:303:PRO:O	2.00	0.44
1:B:303:PRO:HG2	1:B:412:LEU:CD1	2.48	0.44
1:A:323:ARG:HA	1:A:361:HIS:CD2	2.53	0.43
1:A:62:CYS:HB3	1:A:395:ASN:ND2	2.34	0.43
1:A:241:LYS:HE3	3:A:542:HOH:O	2.18	0.43
1:B:387:GLN:HG2	1:B:388:HIS:ND1	2.34	0.43
1:B:370:ASP:OD2	1:B:375:ARG:NH1	2.50	0.43
1:B:391:LYS:HE2	1:B:395:ASN:HB2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:426:HIS:H	1:B:426:HIS:CD2	2.36	0.43
1:A:361:HIS:CE1	1:A:391:LYS:H	2.28	0.43
1:B:173:PHE:HE1	1:B:262:LEU:HD13	1.83	0.42
1:B:306:LYS:NZ	3:B:921:HOH:O	2.51	0.42
1:A:90:TRP:HB3	1:A:92:HIS:CE1	2.55	0.42
1:A:378:ARG:HH11	1:A:378:ARG:HD3	1.69	0.42
1:A:14:LEU:HD12	1:A:14:LEU:HA	1.86	0.42
1:A:268:THR:HB	2:A:460:HEM:C3B	2.55	0.42
1:A:17:LEU:HD22	1:A:45:PRO:CD	2.50	0.42
1:A:160:TYR:OH	1:A:171:HIS:HE1	2.03	0.42
1:B:50:ARG:NH2	3:B:817:HOH:O	2.52	0.42
1:B:138:HIS:HD2	1:B:445:VAL:HG22	1.83	0.42
1:A:118:MET:HB3	1:A:155:LEU:HD13	2.01	0.41
1:B:128:GLN:NE2	3:B:491:HOH:O	2.52	0.41
1:B:223:ARG:HH22	1:B:230:SER:CB	2.33	0.41
1:A:38:GLU:HB2	1:A:54:SER:HB3	2.03	0.41
1:A:356:LEU:HD12	1:A:359:GLN:HB2	2.02	0.41
1:B:388:HIS:HD2	1:B:391:LYS:NZ	2.19	0.41
1:A:3:LYS:HE2	1:A:344:GLU:HA	2.02	0.40
1:B:400:CYS:HA	2:B:460:HEM:CHA	2.52	0.40
1:A:440:LYS:NZ	3:A:1017:HOH:O	2.53	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	437/455 (96%)	423 (97%)	14 (3%)	0	100	100
1	B	437/455 (96%)	427 (98%)	10 (2%)	0	100	100
All	All	874/910 (96%)	850 (97%)	24 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	369/399 (92%)	348 (94%)	21 (6%)	18	15
1	B	377/399 (94%)	359 (95%)	18 (5%)	23	21
All	All	746/798 (94%)	707 (95%)	39 (5%)	21	18

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	LEU
1	A	15	LYS
1	A	20	LEU
1	A	49	THR
1	A	56	ARG
1	A	75	LEU
1	A	92	HIS
1	A	97	LYS
1	A	103	LEU
1	A	104	LEU
1	A	133	LEU
1	A	148	LEU
1	A	155	LEU
1	A	158	PHE
1	A	188	LEU
1	A	272	LEU
1	A	322	LEU
1	A	349	LYS
1	A	356	LEU
1	A	378	ARG
1	A	430	GLU
1	B	3	LYS
1	B	49	THR
1	B	103	LEU
1	B	133	LEU

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Mol	Chain	Res	Type
1	B	158	PHE
1	B	178	VAL
1	B	181	LEU
1	B	188	LEU
1	B	223	ARG
1	B	224	LYS
1	B	263	ILE
1	B	272	LEU
1	B	302	VAL
1	B	306	LYS
1	B	309	LYS
1	B	344	GLU
1	B	445	VAL
1	B	446	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (41) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	GLN
1	A	21	ASN
1	A	27	GLN
1	A	95	ASN
1	A	100	HIS
1	A	110	GLN
1	A	116	HIS
1	A	159	ASN
1	A	169	GLN
1	A	171	HIS
1	A	213	ASN
1	A	229	GLN
1	A	239	ASN
1	A	266	HIS
1	A	319	ASN
1	A	361	HIS
1	A	381	ASN
1	A	395	ASN
1	A	426	HIS
1	B	7	GLN
1	B	16	ASN
1	B	27	GLN
1	B	70	ASN
1	B	95	ASN

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Mol	Chain	Res	Type
1	B	100	HIS
1	B	109	GLN
1	B	110	GLN
1	B	116	HIS
1	B	128	GLN
1	B	138	HIS
1	B	171	HIS
1	B	201	ASN
1	B	236	HIS
1	B	266	HIS
1	B	288	GLN
1	B	319	ASN
1	B	361	HIS
1	B	387	GLN
1	B	388	HIS
1	B	395	ASN
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	A	460	3,1	50,50,50	1.31	5 (10%)	67,82,82	0.90	2 (2%)
2	HEM	B	460	3,1	50,50,50	1.35	7 (14%)	67,82,82	0.83	2 (2%)

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	A	460	3,1	-	2/14/54/54	-
2	HEM	B	460	3,1	-	2/14/54/54	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	460	HEM	FE-NA	3.32	2.06	1.95
2	B	460	HEM	FE-ND	3.27	2.05	1.94
2	A	460	HEM	FE-ND	3.18	2.04	1.94
2	A	460	HEM	CAC-C3C	-3.14	1.39	1.47
2	A	460	HEM	CAB-C3B	-3.09	1.39	1.47
2	B	460	HEM	CBB-CAB	3.09	1.45	1.30
2	B	460	HEM	CAC-C3C	-3.07	1.39	1.47
2	B	460	HEM	CBC-CAC	3.04	1.45	1.30
2	A	460	HEM	CBB-CAB	2.92	1.44	1.30
2	B	460	HEM	CAB-C3B	-2.84	1.39	1.47
2	A	460	HEM	CBC-CAC	2.81	1.43	1.30
2	B	460	HEM	FE-NB	2.06	2.01	1.94

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	460	HEM	CHC-C1C-NC	-2.08	122.18	124.45
2	A	460	HEM	C3B-C4B-NB	2.06	110.95	109.47
2	B	460	HEM	C3B-C4B-NB	2.03	110.92	109.47
2	B	460	HEM	CHA-C4D-ND	-2.02	121.87	124.37

There are no chirality outliers.

All (4) torsion outliers are listed below:

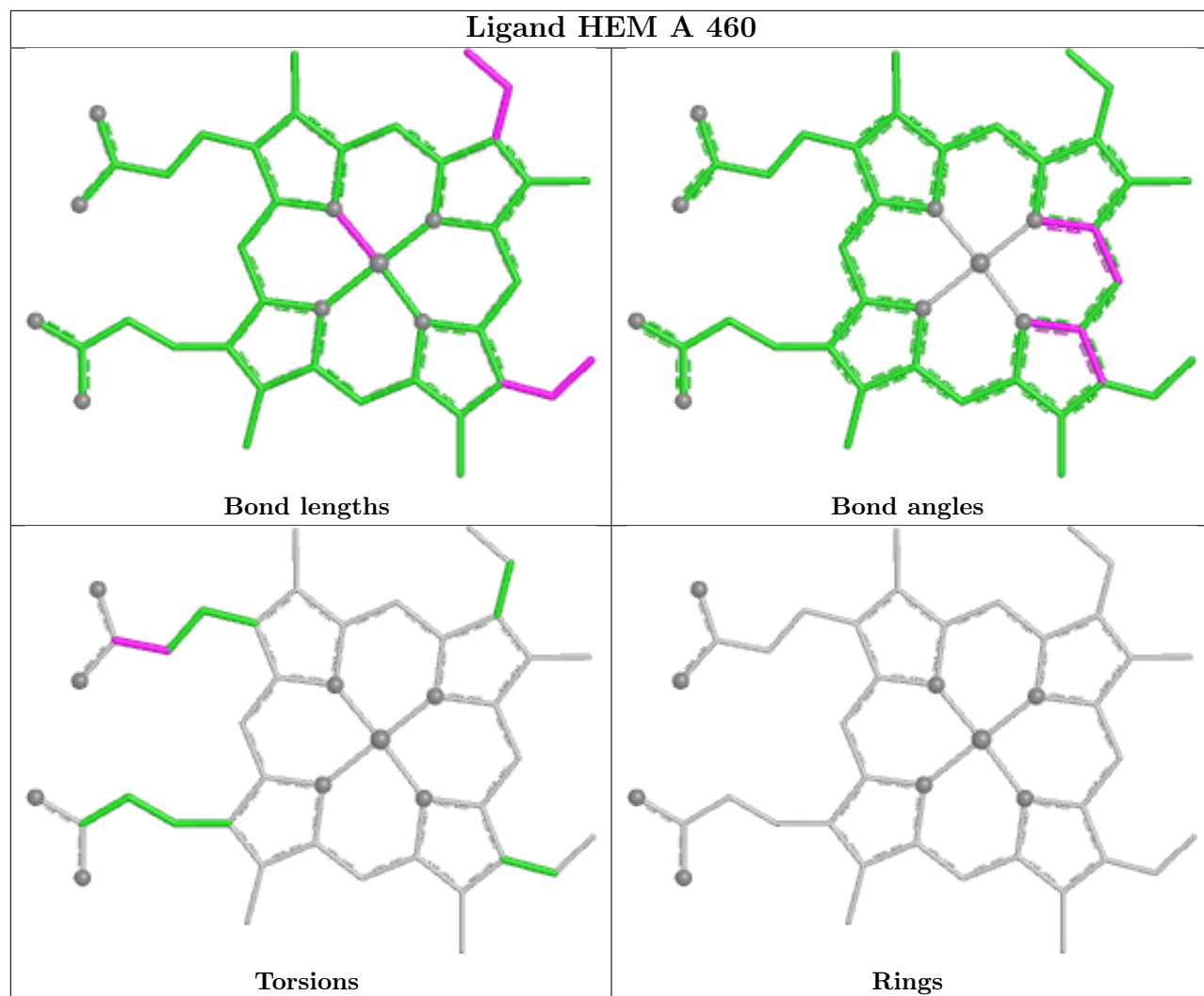
Mol	Chain	Res	Type	Atoms
2	B	460	HEM	CAD-CBD-CGD-O1D
2	B	460	HEM	CAD-CBD-CGD-O2D
2	A	460	HEM	CAD-CBD-CGD-O1D
2	A	460	HEM	CAD-CBD-CGD-O2D

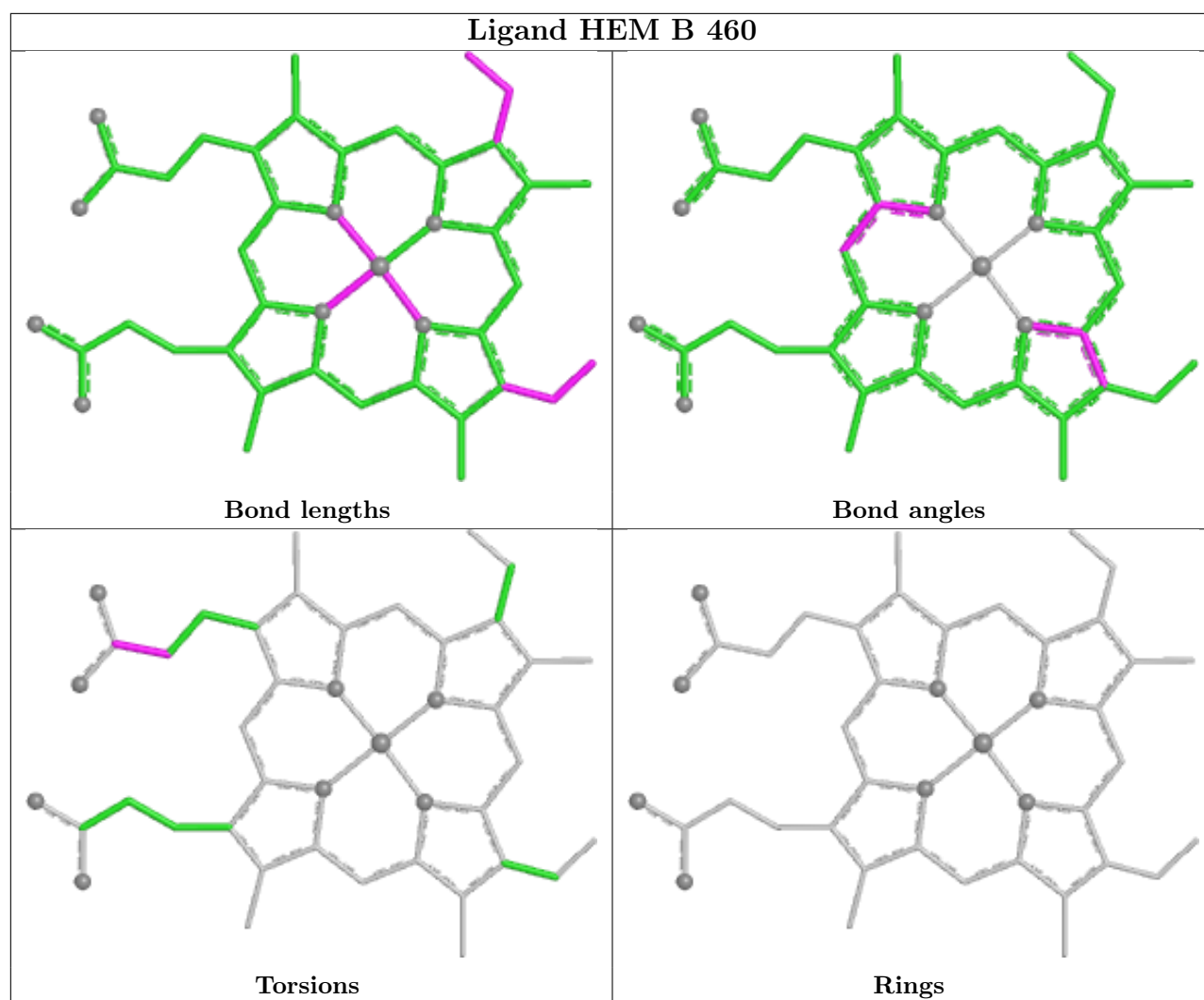
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	460	HEM	1	0
2	B	460	HEM	1	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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.