



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:42 PM UTC

PDB ID : 3EKD / pdb_00003ekd
Title : Crystal structure of the A264M heme domain of cytochrome P450 BM3
Authors : Toogood, H.S.; Leys, D.
Deposited on : 2008-09-19
Resolution : 2.50 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

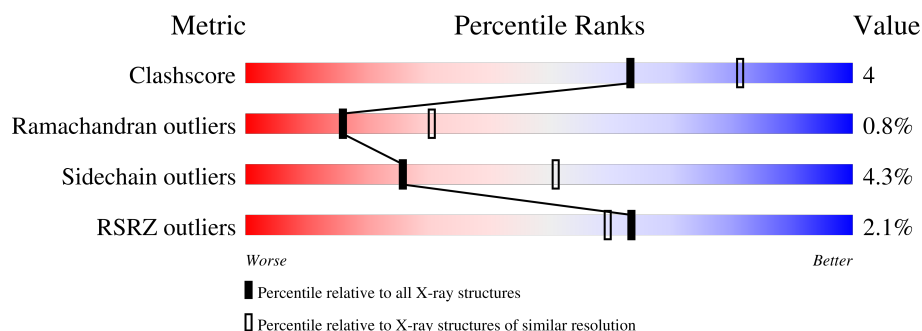
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	470	% 84% 12% . .
1	B	470	3% 80% 15% . .

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7523 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

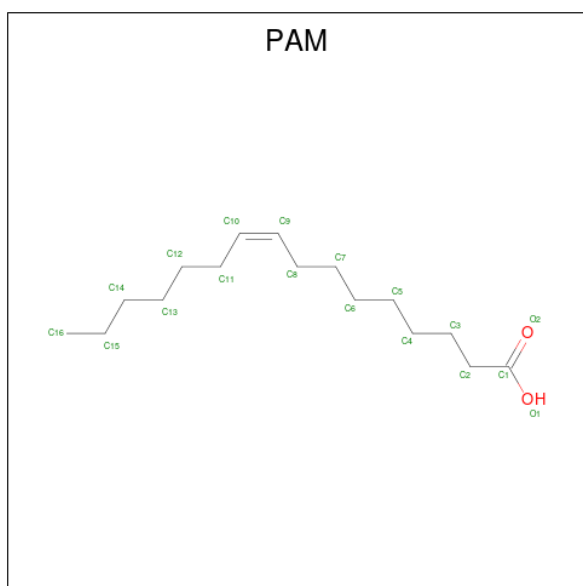
- Molecule 1 is a protein called Cytochrome P450(BM-3).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	454	Total	C	N	O	S	0	0	0
			3623	2320	610	675	18			
1	B	455	Total	C	N	O	S	0	0	0
			3624	2321	612	673	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	264	MET	ALA	engineered mutation	UNP P14779
B	264	MET	ALA	engineered mutation	UNP P14779

- Molecule 2 is PALMITOLEIC ACID (CCD ID: PAM) (formula: C₁₆H₃₀O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			18	16	2		

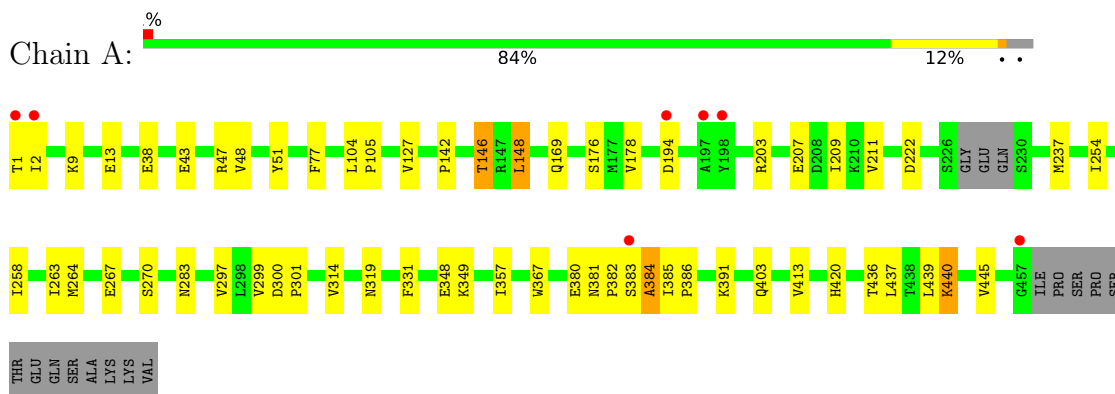
- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



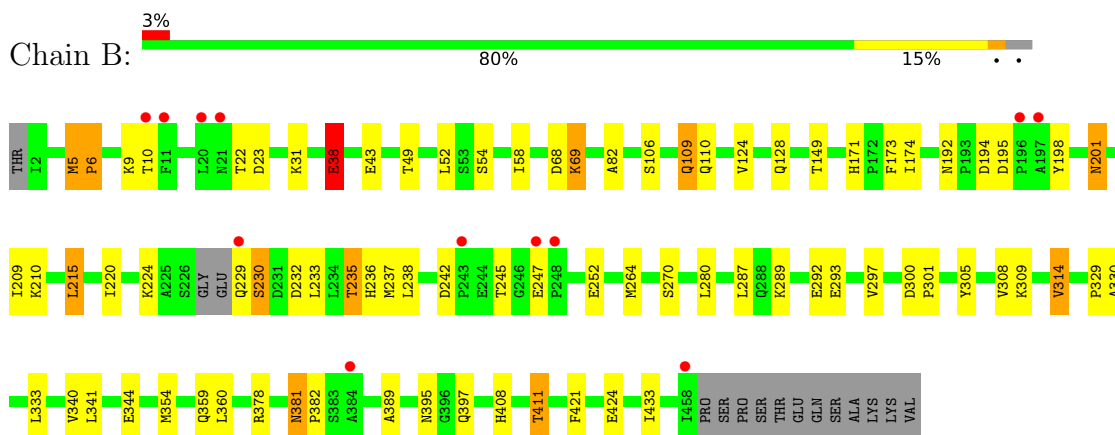
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Cytochrome P450(BM-3)



• Molecule 1: Cytochrome P450(BM-3)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.77Å 125.17Å 153.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 20.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.4 (20.00-2.50) 99.1 (20.00-2.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.217 , 0.252 0.221 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	34.5	Xtriage
Anisotropy	0.252	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 51.5	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	0.93	EDS
Total number of atoms	7523	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% 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: PAM, 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	1.64	16/3707 (0.4%)	1.25	11/5017 (0.2%)
1	B	1.52	20/3708 (0.5%)	1.26	17/5020 (0.3%)
All	All	1.58	36/7415 (0.5%)	1.26	28/10037 (0.3%)

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	357	ILE	CA-CB	11.73	1.61	1.53
1	B	149	THR	CA-CB	8.15	1.66	1.53
1	A	391	LYS	CA-C	8.02	1.62	1.53
1	B	314	VAL	CA-CB	7.21	1.63	1.54
1	B	424	GLU	CA-C	6.51	1.60	1.52
1	B	52	LEU	CA-C	5.98	1.60	1.52
1	B	270	SER	CA-C	5.94	1.60	1.52
1	B	411	THR	CA-CB	5.90	1.62	1.53
1	B	421	PHE	N-CA	5.70	1.53	1.45
1	A	314	VAL	CA-CB	5.70	1.61	1.54
1	B	174	ILE	CA-CB	5.61	1.61	1.54
1	A	209	ILE	CA-CB	5.52	1.61	1.54
1	B	397	GLN	CA-C	5.50	1.60	1.52
1	B	69	LYS	N-CA	5.47	1.52	1.46
1	A	77	PHE	CA-C	5.46	1.60	1.52
1	B	381	ASN	N-CA	5.42	1.51	1.46
1	A	283	ASN	CA-C	5.39	1.58	1.52
1	A	267	GLU	CA-C	5.37	1.59	1.52
1	A	439	LEU	N-CA	5.35	1.52	1.46
1	B	408	HIS	CA-C	5.35	1.59	1.52
1	B	245	THR	CA-CB	5.33	1.61	1.53
1	B	308	VAL	CA-CB	5.32	1.61	1.54
1	B	174	ILE	CA-C	5.28	1.59	1.52
1	A	2	ILE	CA-C	5.28	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	51	TYR	CA-C	5.26	1.59	1.52
1	A	331	PHE	CA-C	5.17	1.59	1.52
1	A	299	VAL	CA-C	5.14	1.58	1.52
1	A	367	TRP	CA-C	5.14	1.59	1.52
1	B	389	ALA	CA-C	5.12	1.59	1.52
1	A	127	VAL	CA-CB	5.08	1.60	1.54
1	B	247	GLU	N-CA	5.08	1.51	1.45
1	B	242	ASP	CA-C	5.06	1.58	1.52
1	A	445	VAL	N-CA	5.04	1.51	1.46
1	B	82	ALA	N-CA	5.02	1.52	1.46
1	A	297	VAL	CA-CB	5.01	1.61	1.54
1	B	340	VAL	N-CA	5.01	1.51	1.46

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	381	ASN	CA-C-N	8.48	130.44	119.84
1	B	381	ASN	C-N-CA	8.48	130.44	119.84
1	B	329	PRO	N-CA-C	8.23	123.24	114.68
1	B	344	GLU	N-CA-C	8.07	120.16	111.36
1	A	420	HIS	N-CA-C	7.23	122.11	113.28
1	B	5	MET	CA-C-N	6.71	126.89	120.31
1	B	5	MET	C-N-CA	6.71	126.89	120.31
1	A	319	ASN	N-CA-C	6.45	118.31	111.28
1	A	47	ARG	N-CA-C	6.42	118.83	109.07
1	B	10	THR	N-CA-C	5.81	118.37	110.35
1	A	380	GLU	N-CA-C	5.75	118.01	111.11
1	A	386	PRO	N-CA-C	5.70	120.18	111.34
1	A	403	GLN	N-CA-C	5.69	117.94	111.11
1	A	38	GLU	N-CA-C	5.68	119.31	112.38
1	B	38	GLU	N-CA-C	5.68	119.47	112.54
1	A	391	LYS	N-CA-C	5.45	119.18	110.73
1	B	174	ILE	N-CA-C	5.39	116.77	110.62
1	B	6	PRO	N-CA-C	5.39	119.76	111.57
1	B	43	GLU	N-CA-C	5.36	118.14	109.40
1	B	238	LEU	N-CA-C	5.27	117.03	111.28
1	B	421	PHE	N-CA-C	5.20	117.09	109.24
1	A	385	ILE	CB-CA-C	-5.17	101.94	111.36
1	B	395	ASN	N-CA-C	5.08	117.43	108.75
1	A	178	VAL	N-CA-C	5.07	117.39	111.05
1	A	43	GLU	N-CA-C	5.06	117.65	109.40
1	B	341	LEU	N-CA-C	5.06	117.15	108.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	171	HIS	CA-C-N	5.01	126.10	119.84
1	B	171	HIS	C-N-CA	5.01	126.10	119.84

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	3623	0	3556	19	0
1	B	3624	0	3550	34	0
2	A	18	0	29	0	0
3	A	43	0	30	1	0
3	B	43	0	30	4	0
4	A	111	0	0	2	0
4	B	61	0	0	3	0
All	All	7523	0	7195	58	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (58) 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:173:PHE:HD1	1:B:215:LEU:HD21	1.21	1.04
1:B:330:ALA:HB1	1:B:354:MET:CE	1.94	0.97
1:A:237:MET:HE1	1:A:258:ILE:CD1	1.96	0.96
1:B:209:ILE:HG22	4:B:506:HOH:O	1.67	0.94
1:B:173:PHE:CD1	1:B:215:LEU:HD21	2.11	0.85
1:B:58:ILE:HD12	1:B:360:LEU:HD13	1.63	0.81
1:B:330:ALA:HB1	1:B:354:MET:HE2	1.63	0.79
3:B:471:HEM:HMC1	3:B:471:HEM:HBC2	1.65	0.78
1:A:237:MET:HE1	1:A:258:ILE:HD13	1.66	0.77
3:B:471:HEM:HMB2	3:B:471:HEM:HBB2	1.70	0.74
1:B:220:ILE:HG22	1:B:224:LYS:HE3	1.71	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:MET:HE1	1:A:258:ILE:HD11	1.68	0.71
1:A:237:MET:CE	1:A:254:ILE:HG23	2.22	0.70
1:B:201:ASN:HD22	1:B:201:ASN:H	1.39	0.70
1:A:237:MET:HE3	1:A:254:ILE:HG23	1.76	0.68
1:A:146:THR:HG22	1:A:270:SER:OG	1.98	0.63
1:B:68:ASP:CG	4:B:505:HOH:O	2.42	0.62
1:B:330:ALA:HB1	1:B:354:MET:HE1	1.82	0.62
1:B:201:ASN:H	1:B:201:ASN:ND2	1.98	0.61
3:B:471:HEM:HBC2	3:B:471:HEM:CMC	2.33	0.59
1:B:230:SER:O	1:B:235:THR:HG21	2.02	0.59
1:B:220:ILE:HG22	1:B:224:LYS:CE	2.34	0.58
1:B:232:ASP:OD1	1:B:235:THR:HG22	2.03	0.57
1:A:207:GLU:O	1:A:211:VAL:HG23	2.05	0.57
1:A:381:ASN:O	1:A:384:ALA:HB3	2.05	0.56
1:B:124:VAL:O	1:B:128:GLN:HG3	2.05	0.56
1:B:229:GLN:O	1:B:230:SER:O	2.23	0.55
3:B:471:HEM:HBB2	3:B:471:HEM:CMB	2.39	0.52
1:B:232:ASP:H	1:B:235:THR:CG2	2.24	0.51
1:A:258:ILE:HA	4:A:573:HOH:O	2.11	0.51
1:B:280:LEU:HB3	1:B:287:LEU:HD13	1.92	0.50
1:A:237:MET:CE	1:A:258:ILE:HD11	2.38	0.50
1:A:148:LEU:HD21	1:A:413:VAL:HG21	1.93	0.50
1:B:305:TYR:C	1:B:305:TYR:CD1	2.92	0.48
1:B:293:GLU:O	1:B:297:VAL:HG23	2.14	0.48
1:B:38:GLU:HB3	1:B:54:SER:HB3	1.95	0.47
1:B:300:ASP:HB3	1:B:301:PRO:CD	2.45	0.47
1:A:436:THR:O	1:A:437:LEU:HB2	2.15	0.47
1:A:142:PRO:O	1:A:146:THR:HG23	2.14	0.46
1:B:109:GLN:NE2	1:B:109:GLN:HA	2.30	0.46
1:A:382:PRO:O	1:A:384:ALA:N	2.49	0.45
1:A:440:LYS:HD3	4:A:575:HOH:O	2.16	0.45
1:B:5:MET:HA	1:B:6:PRO:HD3	1.78	0.45
1:B:381:ASN:HA	1:B:382:PRO:HD2	1.73	0.45
3:A:472:HEM:HBC2	3:A:472:HEM:HMC2	1.99	0.45
1:B:233:LEU:HG	1:B:237:MET:HE3	2.00	0.44
1:B:235:THR:HB	4:B:495:HOH:O	2.18	0.44
1:A:348:GLU:O	1:A:349:LYS:C	2.61	0.44
1:B:106:SER:HG	1:B:236:HIS:HD1	1.64	0.44
1:A:104:LEU:N	1:A:105:PRO:CD	2.82	0.43
1:B:314:VAL:HG12	1:B:411:THR:HG23	2.01	0.43
1:B:359:GLN:O	1:B:360:LEU:C	2.62	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:232:ASP:O	1:B:235:THR:HG23	2.20	0.41
1:B:192:ASN:O	1:B:195:ASP:HB2	2.19	0.41
1:A:300:ASP:HB3	1:A:301:PRO:CD	2.50	0.41
1:A:237:MET:CE	1:A:258:ILE:CD1	2.83	0.41
1:B:192:ASN:O	1:B:195:ASP:CB	2.68	0.41
1:B:58:ILE:CD1	1:B:360:LEU:HD13	2.43	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	450/470 (96%)	425 (94%)	22 (5%)	3 (1%)	18	34
1	B	451/470 (96%)	404 (90%)	43 (10%)	4 (1%)	14	27
All	All	901/940 (96%)	829 (92%)	65 (7%)	7 (1%)	16	31

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	384	ALA
1	B	230	SER
1	A	383	SER
1	B	378	ARG
1	B	9	LYS
1	B	198	TYR
1	A	263	ILE

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	386/412 (94%)	373 (97%)	13 (3%)	32	60
1	B	385/412 (93%)	365 (95%)	20 (5%)	21	42
All	All	771/824 (94%)	738 (96%)	33 (4%)	26	51

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	THR
1	A	9	LYS
1	A	13	GLU
1	A	48	VAL
1	A	146	THR
1	A	148	LEU
1	A	169	GLN
1	A	176	SER
1	A	194	ASP
1	A	203	ARG
1	A	222	ASP
1	A	264	MET
1	A	440	LYS
1	B	22	THR
1	B	23	ASP
1	B	31	LYS
1	B	38	GLU
1	B	49	THR
1	B	69	LYS
1	B	109	GLN
1	B	110	GLN
1	B	194	ASP
1	B	201	ASN
1	B	210	LYS
1	B	215	LEU
1	B	235	THR
1	B	252	GLU

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Mol	Chain	Res	Type
1	B	264	MET
1	B	289	LYS
1	B	292	GLU
1	B	309	LYS
1	B	333	LEU
1	B	433	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	70	ASN
1	A	73	GLN
1	A	116	HIS
1	A	163	ASN
1	A	189	GLN
1	A	201	ASN
1	A	239	ASN
1	A	359	GLN
1	A	388	HIS
1	A	426	HIS
1	B	100	HIS
1	B	109	GLN
1	B	116	HIS
1	B	201	ASN
1	B	253	ASN
1	B	257	GLN
1	B	266	HIS
1	B	283	ASN
1	B	359	GLN
1	B	388	HIS
1	B	397	GLN
1	B	403	GLN
1	B	420	HIS
1	B	426	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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	HEM	B	471	1	50,50,50	2.05	11 (22%)	67,82,82	1.49	13 (19%)
2	PAM	A	471	-	17,17,17	1.32	2 (11%)	17,17,17	0.86	0
3	HEM	A	472	1	50,50,50	2.67	11 (22%)	67,82,82	1.97	17 (25%)

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	HEM	B	471	1	-	0/14/54/54	-
2	PAM	A	471	-	-	5/15/15/15	-
3	HEM	A	472	1	-	2/14/54/54	-

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	472	HEM	FE-ND	10.37	2.26	1.94
3	A	472	HEM	C3D-C2D	7.91	1.53	1.36
3	B	471	HEM	C3D-C2D	7.56	1.53	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	472	HEM	FE-NC	5.90	2.14	1.95
3	B	471	HEM	FE-NC	5.48	2.13	1.95
3	B	471	HEM	FE-NA	5.26	2.12	1.95
3	A	472	HEM	FE-NB	4.97	2.10	1.94
3	A	472	HEM	FE-NA	4.45	2.09	1.95
2	A	471	PAM	C10-C9	3.91	1.53	1.31
3	A	472	HEM	CAC-C3C	3.86	1.57	1.47
3	B	471	HEM	CAB-C3B	3.23	1.56	1.47
3	A	472	HEM	CAB-C3B	3.01	1.55	1.47
3	B	471	HEM	CAC-C3C	2.84	1.55	1.47
3	A	472	HEM	CMC-C2C	2.75	1.56	1.50
3	B	471	HEM	CMD-C2D	2.66	1.56	1.50
3	A	472	HEM	CAA-C2A	2.43	1.57	1.51
3	B	471	HEM	CMC-C2C	2.36	1.55	1.50
3	B	471	HEM	C1C-C2C	2.36	1.50	1.45
3	A	472	HEM	CMB-C2B	2.35	1.55	1.50
3	B	471	HEM	CAA-C2A	2.33	1.57	1.51
3	A	472	HEM	CMD-C2D	2.16	1.55	1.50
3	B	471	HEM	CMA-C3A	2.16	1.55	1.50
2	A	471	PAM	O2-C1	2.14	1.29	1.22
3	B	471	HEM	CMB-C2B	2.08	1.55	1.50

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	472	HEM	C4D-ND-C1D	7.18	113.71	105.21
3	A	472	HEM	C3B-C4B-NB	-5.22	105.72	109.47
3	A	472	HEM	C1B-NB-C4B	4.60	110.66	105.21
3	B	471	HEM	C4D-ND-C1D	4.02	109.97	105.21
3	B	471	HEM	CHD-C1D-ND	3.80	128.51	124.42
3	A	472	HEM	C2B-C1B-NB	-3.74	105.54	109.84
3	A	472	HEM	C3B-C2B-C1B	3.51	109.05	106.41
3	A	472	HEM	C4C-C3C-C2C	3.43	109.79	106.81
3	A	472	HEM	C4A-C3A-C2A	3.17	110.44	106.82
3	B	471	HEM	CHC-C4B-NB	2.99	127.64	124.42
3	B	471	HEM	O2A-CGA-CBA	2.97	123.38	114.00
3	A	472	HEM	C2D-C1D-ND	-2.78	106.69	109.90
3	B	471	HEM	CHB-C1B-NB	2.55	127.53	124.37
3	A	472	HEM	CMD-C2D-C1D	2.52	128.97	125.03
3	B	471	HEM	CHD-C1D-C2D	-2.52	121.06	125.03
3	B	471	HEM	O1A-CGA-CBA	-2.51	115.12	123.09
3	B	471	HEM	CHA-C4D-ND	2.51	127.47	124.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	472	HEM	C3A-C4A-NA	-2.48	106.16	110.14
3	B	471	HEM	C3B-C2B-C1B	2.47	108.26	106.41
3	A	472	HEM	CHD-C4C-NC	2.40	127.07	124.45
3	B	471	HEM	CAD-C3D-C4D	2.35	128.80	124.70
3	B	471	HEM	CAA-C2A-C1A	2.33	129.50	124.94
3	A	472	HEM	CHD-C1D-ND	2.28	126.87	124.42
3	A	472	HEM	C4A-CHB-C1B	2.24	131.52	126.25
3	A	472	HEM	CAA-C2A-C1A	2.20	129.23	124.94
3	A	472	HEM	CBA-CAA-C2A	2.14	118.44	112.53
3	B	471	HEM	C1D-C2D-C3D	-2.10	104.78	106.98
3	B	471	HEM	CAA-CBA-CGA	2.04	119.07	113.67
3	A	472	HEM	C3D-C4D-ND	-2.02	107.95	110.17
3	A	472	HEM	C4A-NA-C1A	2.02	109.11	105.82

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	471	PAM	C13-C14-C15-C16
2	A	471	PAM	C11-C10-C9-C8
2	A	471	PAM	C12-C13-C14-C15
2	A	471	PAM	C1-C2-C3-C4
3	A	472	HEM	CAD-CBD-CGD-O1D
3	A	472	HEM	CAD-CBD-CGD-O2D
2	A	471	PAM	C7-C8-C9-C10

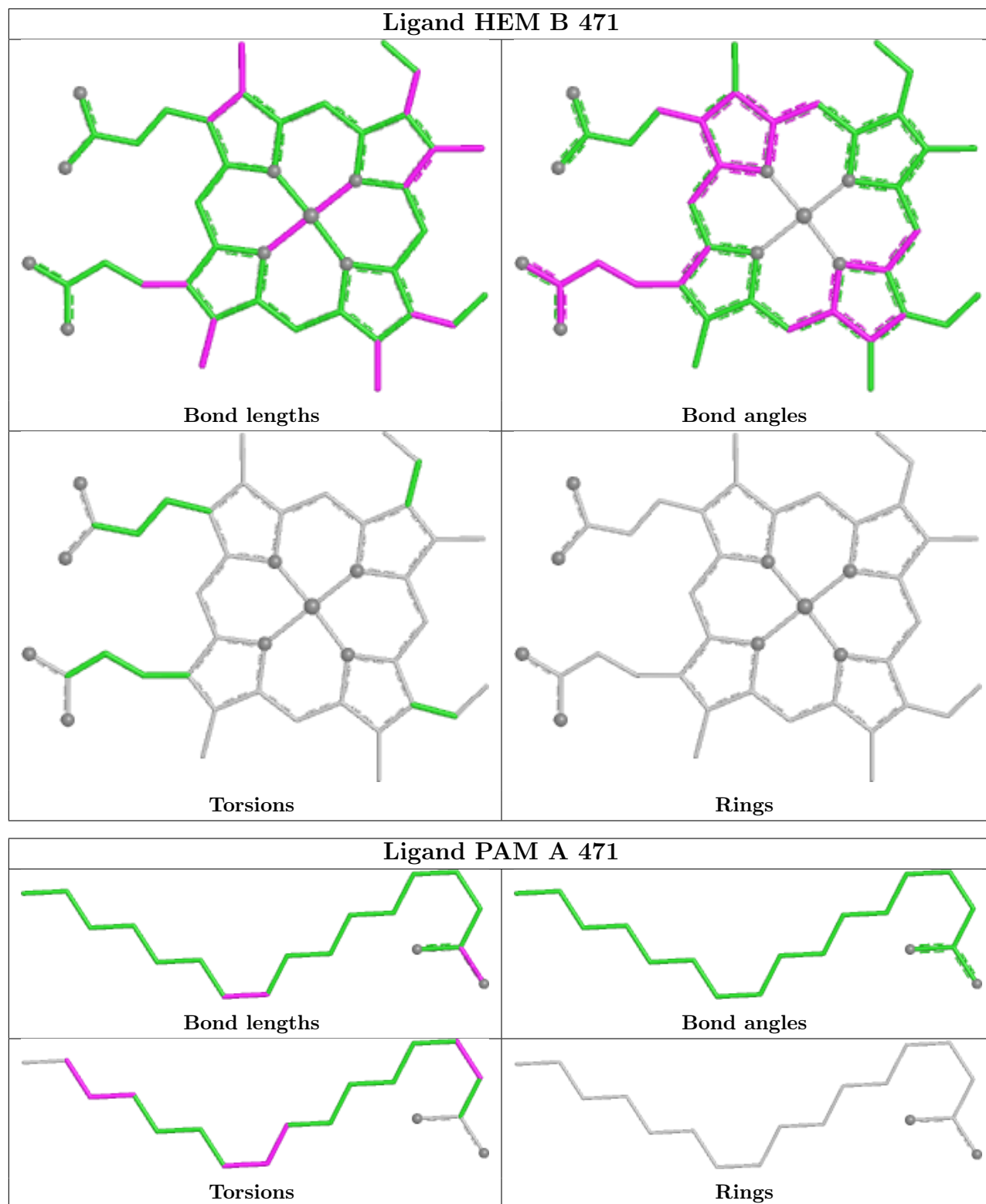
There are no ring outliers.

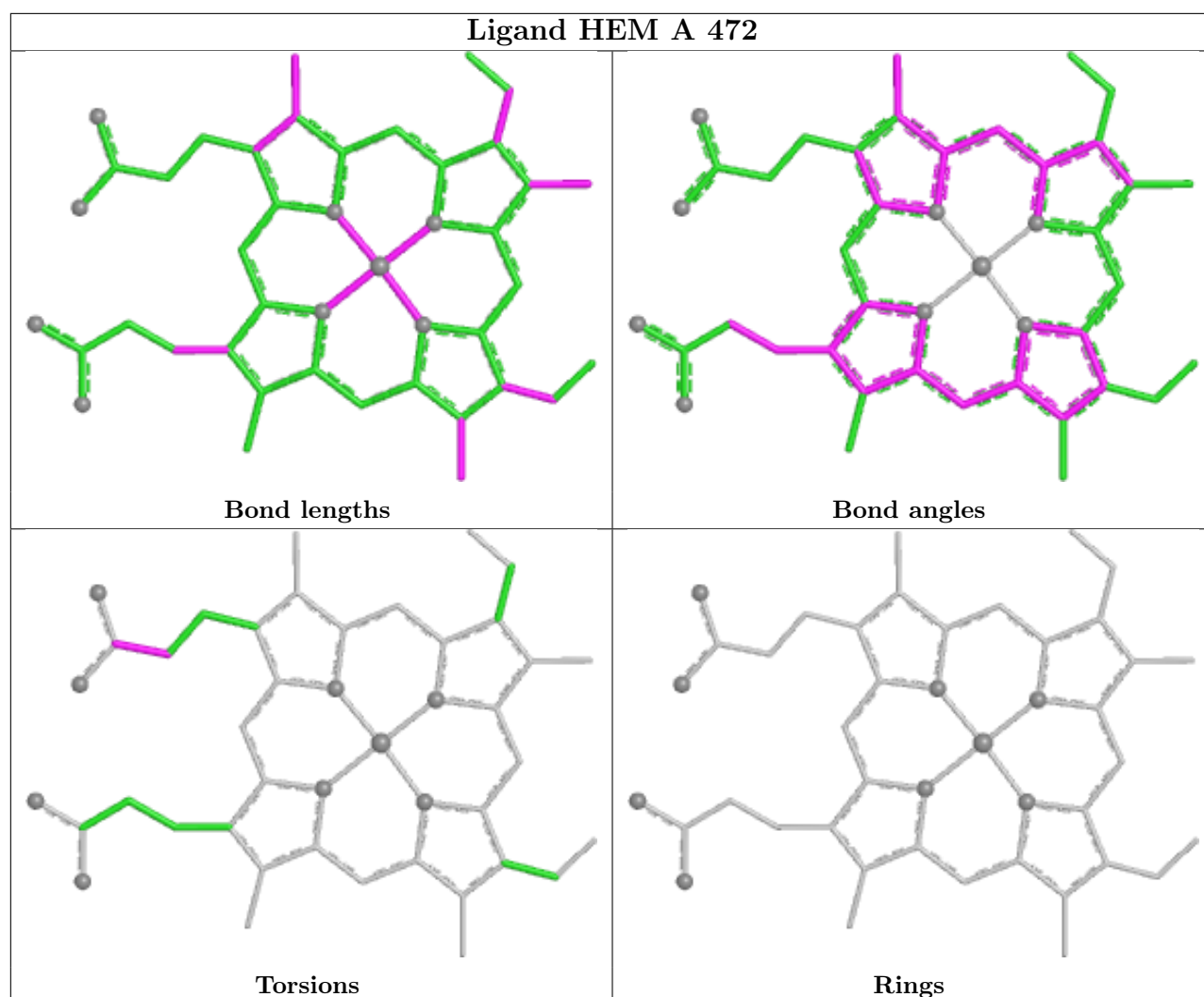
2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	471	HEM	4	0
3	A	472	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	454/470 (96%)	-0.34	7 (1%) 72 68	13, 27, 52, 73	0
1	B	455/470 (96%)	0.21	12 (2%) 57 52	21, 39, 66, 72	0
All	All	909/940 (96%)	-0.07	19 (2%) 63 59	13, 33, 61, 73	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	458	ILE	3.8
1	B	243	PRO	3.0
1	B	196	PRO	2.8
1	A	457	GLY	2.7
1	A	383	SER	2.6
1	B	248	PRO	2.6
1	A	194	ASP	2.4
1	B	197	ALA	2.3
1	B	229	GLN	2.2
1	B	247	GLU	2.2
1	A	197	ALA	2.2
1	B	10	THR	2.2
1	B	20	LEU	2.2
1	A	1	THR	2.1
1	A	198	TYR	2.1
1	A	2	ILE	2.1
1	B	384	ALA	2.0
1	B	21	ASN	2.0
1	B	11	PHE	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](#)

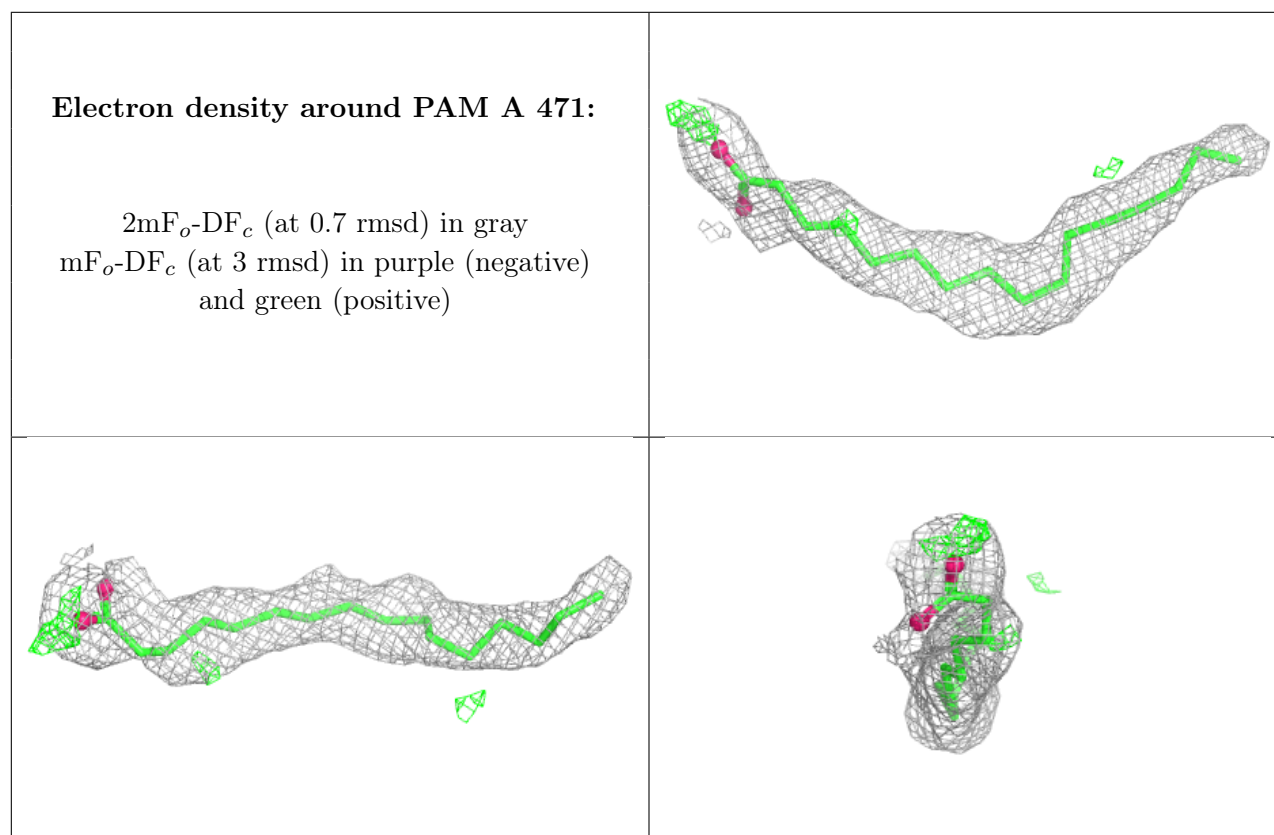
There are no oligosaccharides in this entry.

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.

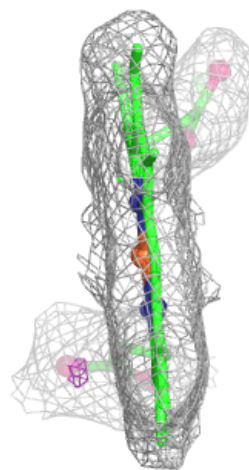
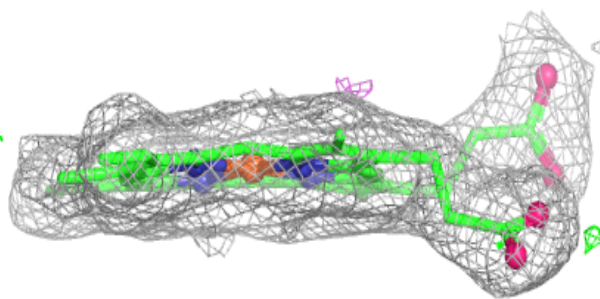
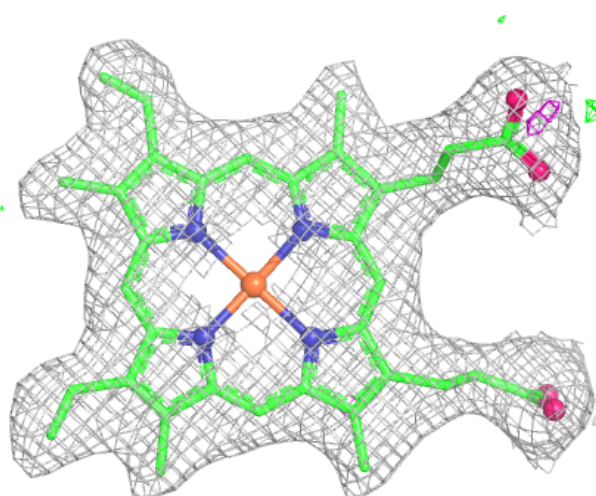
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PAM	A	471	18/18	0.79	0.18	47,50,60,61	0
3	HEM	B	471	43/43	0.97	0.06	18,23,28,30	0
3	HEM	A	472	43/43	0.98	0.05	6,14,17,20	0

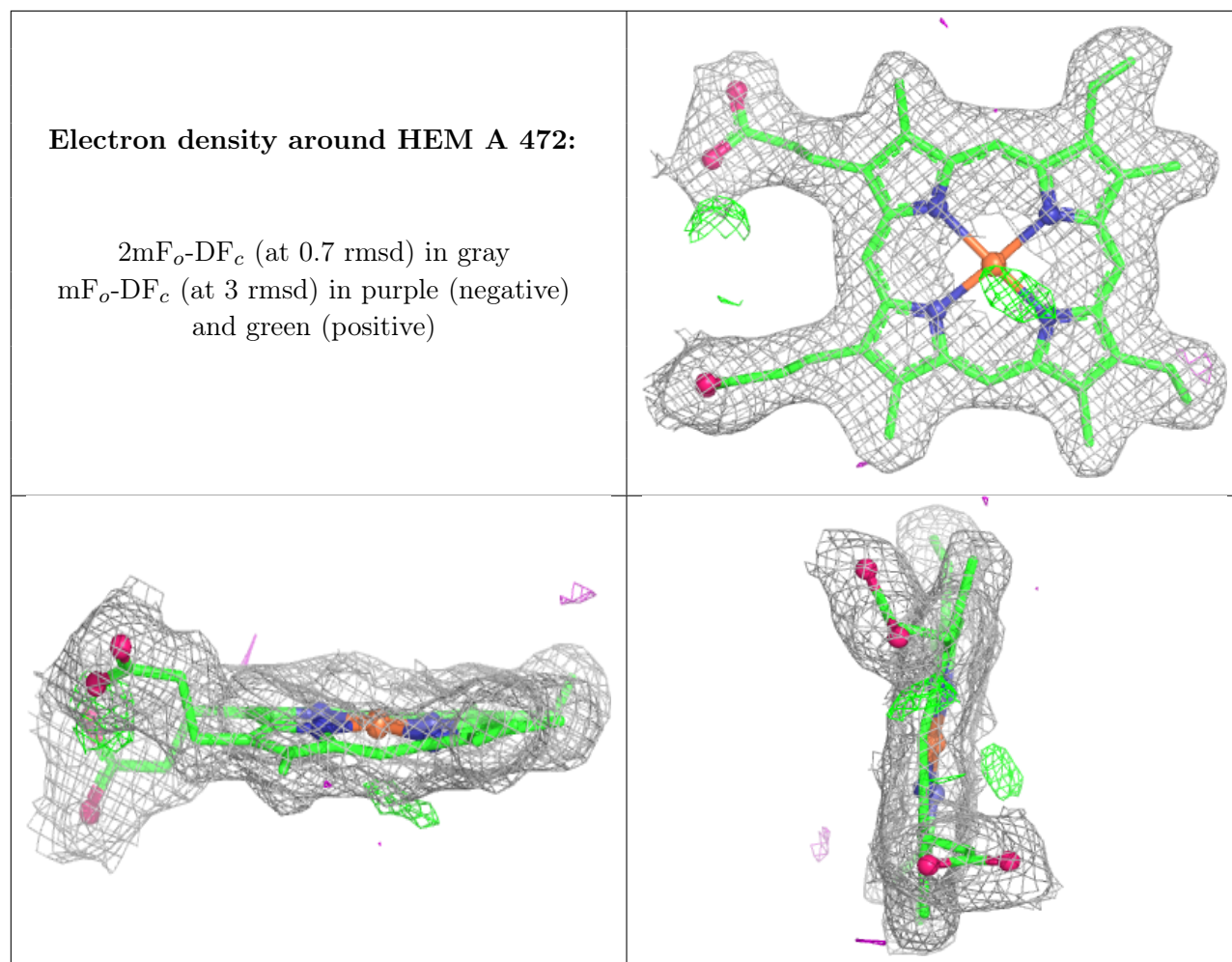
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around HEM B 471:

$2mF_o-DF_c$ (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.