Full wwPDB X-ray Structure Validation Report ⓘ

Sep 3, 2023 – 11:39 AM EDT

PDB ID	:	3S3A
Title	:	Structure of Thermus thermophilus cytochrome ba3 oxidase 120s after Xe de-pressurization
Authors	:	Luna, V.M.; Fee, J.A.; Deniz, A.A.; Stout, C.D.
Deposited on	:	2011-05-18
Resolution	:	4.25 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:

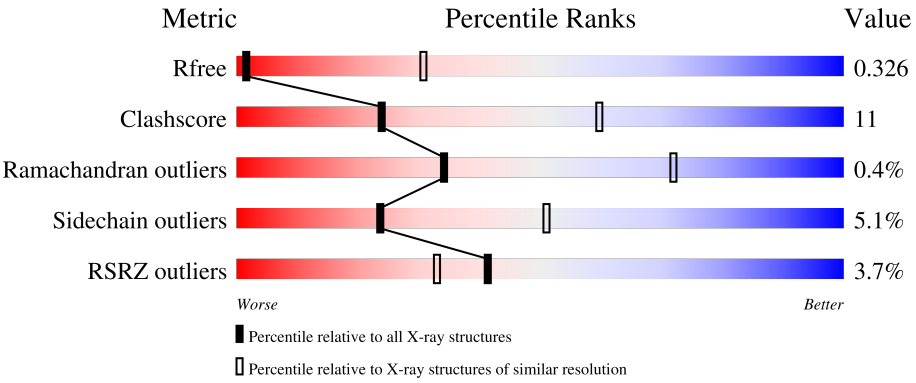
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 4.25 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}	130704	1017 (4.72-3.78)
Clashscore	141614	1059 (4.72-3.80)
Ramachandran outliers	138981	1014 (4.72-3.80)
Sidechain outliers	138945	1018 (4.72-3.78)
RSRZ outliers	127900	1072 (4.80-3.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	568	4% 73% 21%
2	B	166	4% 67% 25% 7%
3	C	33	3% 61% 36%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
7	XE	A	563	-	-	-	X

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 6056 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 c oxidase subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	554	Total	C	N	O	S	0	0	0
			4384	2971	704	693	16			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	MET	-	expression tag	UNP Q5SJ79
A	-4	HIS	-	expression tag	UNP Q5SJ79
A	-3	HIS	-	expression tag	UNP Q5SJ79
A	-2	HIS	-	expression tag	UNP Q5SJ79
A	-1	HIS	-	expression tag	UNP Q5SJ79
A	0	HIS	-	expression tag	UNP Q5SJ79
A	1	HIS	-	expression tag	UNP Q5SJ79

- Molecule 2 is a protein called Cytochrome c oxidase subunit 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	166	Total	C	N	O	S	0	0	0
			1298	844	216	234	4			

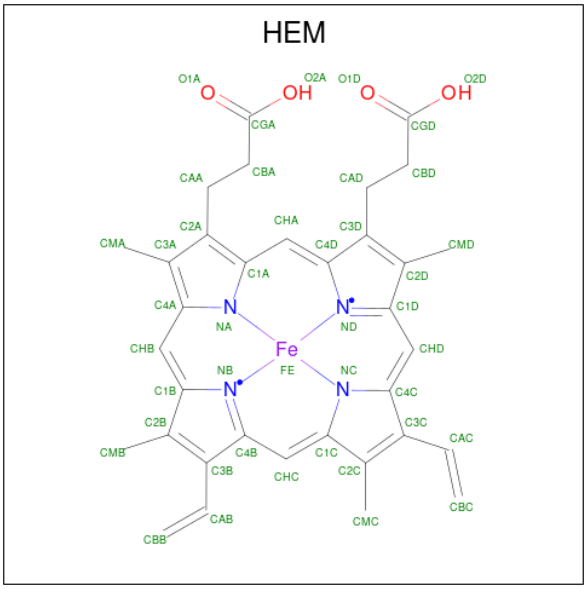
- Molecule 3 is a protein called Cytochrome c oxidase polypeptide 2A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	33	Total	C	N	O	0	0	0
			259	179	39	41			

- Molecule 4 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

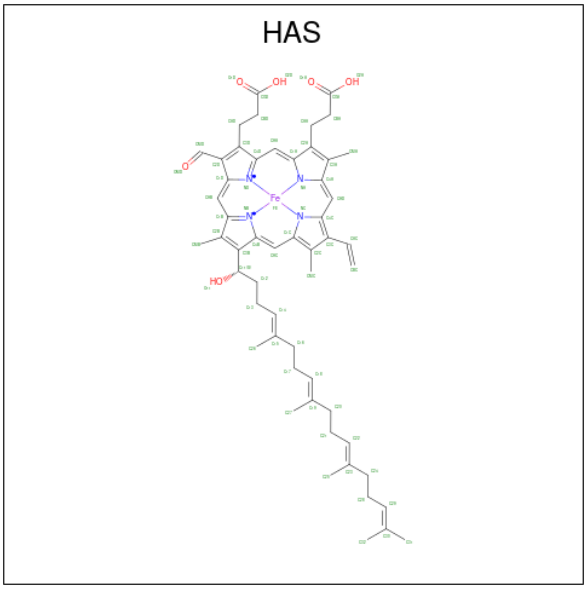
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cu	0	0
			1	1		

- Molecule 5 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

- Molecule 6 is HEME-AS (three-letter code: HAS) (formula: $C_{54}H_{64}FeN_4O_6$).

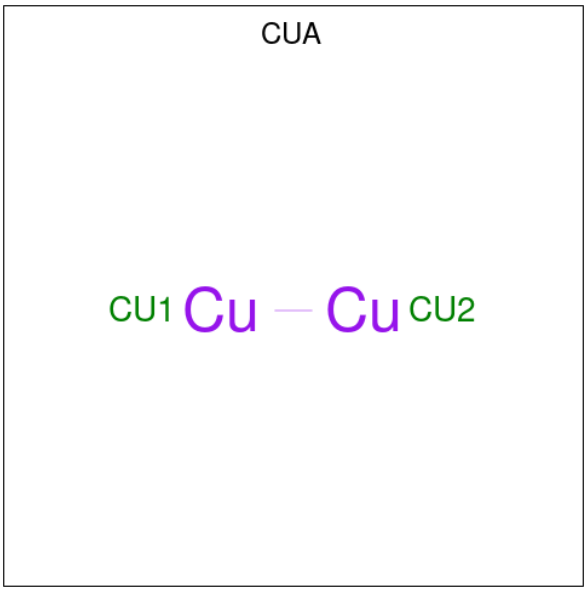


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	Fe	N	O	0	0
			65	54	1	4	6		

- Molecule 7 is XENON (three-letter code: XE) (formula: Xe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	4	Total Xe 4 4	0	0

- Molecule 8 is DINUCLEAR COPPER ION (three-letter code: CUA) (formula: Cu_2).

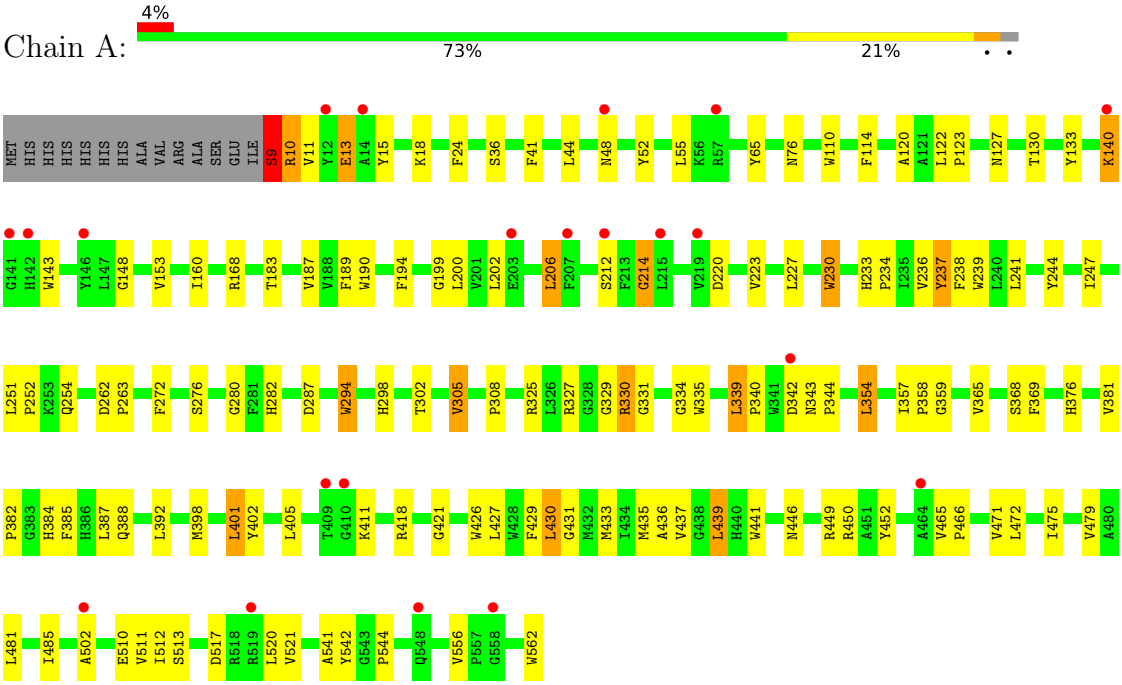


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	B	1	Total Cu 2 2	0	0

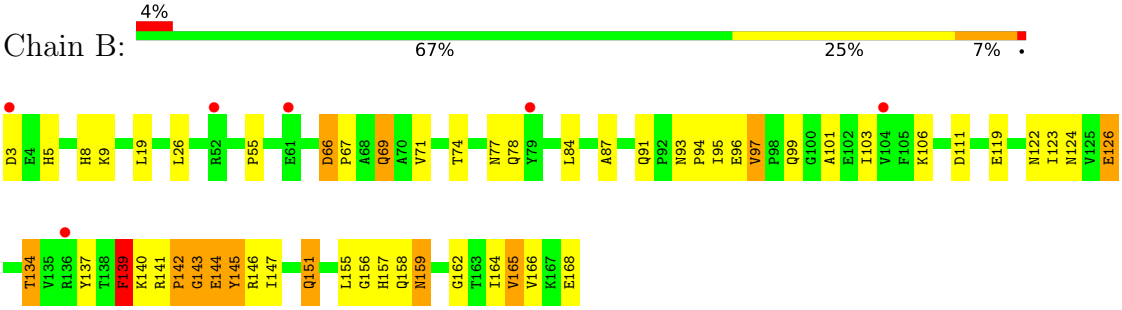
3 Residue-property plots (i)

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: Cytochrome c oxidase subunit 1



• Molecule 2: Cytochrome c oxidase subunit 2



• Molecule 3: Cytochrome c oxidase polypeptide 2A

E2
E3
K4
P5
K6
L9
I12
L13
V14
L15
L20
W23
L24
G25
V26
R33
G34

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α, β, γ	110.29Å 110.29Å 170.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	92.57 – 4.25 55.15 – 4.25	Depositor EDS
% Data completeness (in resolution range)	99.7 (92.57-4.25) 99.7 (55.15-4.25)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.17	Depositor
$< I/\sigma(I) >$ [1]	6.79 (at 4.29Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.279 , 0.330 0.272 , 0.326	Depositor DCC
R_{free} test set	361 reflections (4.60%)	wwPDB-VP
Wilson B-factor (Å^2)	146.8	Xtriage
Anisotropy	0.104	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.27 , 35.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.83	EDS
Total number of atoms	6056	wwPDB-VP
Average B, all atoms (Å^2)	123.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.23% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: HAS, XE, CUA, HEM, CU

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	1.04	11/4541 (0.2%)	0.71	9/6232 (0.1%)
2	B	1.16	14/1335 (1.0%)	0.80	8/1822 (0.4%)
3	C	0.63	0/265	0.55	0/359
All	All	1.06	25/6141 (0.4%)	0.73	17/8413 (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	2

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	13	GLU	CD-OE1	32.90	1.61	1.25
1	A	510	GLU	CD-OE1	28.14	1.56	1.25
2	B	165	VAL	CB-CG1	16.12	1.86	1.52
1	A	140	LYS	CE-NZ	13.79	1.83	1.49
2	B	168	GLU	CD-OE2	10.60	1.37	1.25
1	A	214	GLY	C-O	9.73	1.39	1.23
1	A	9	SER	C-N	9.72	1.56	1.34
2	B	143	GLY	N-CA	9.43	1.60	1.46
2	B	144	GLU	CD-OE2	9.13	1.35	1.25
2	B	144	GLU	C-O	8.77	1.40	1.23
2	B	145	TYR	CZ-OH	7.79	1.51	1.37
1	A	15	TYR	CZ-OH	6.55	1.49	1.37
2	B	146	ARG	CZ-NH2	6.37	1.41	1.33
1	A	510	GLU	CG-CD	6.29	1.61	1.51
2	B	144	GLU	N-CA	6.12	1.58	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	237	TYR	CE2-CZ	6.08	1.46	1.38
2	B	151	GLN	CD-NE2	-5.83	1.18	1.32
1	A	9	SER	CA-C	5.79	1.68	1.52
1	A	9	SER	CB-OG	5.61	1.49	1.42
2	B	144	GLU	CA-C	5.18	1.66	1.52
2	B	144	GLU	CG-CD	5.18	1.59	1.51
2	B	126	GLU	CD-OE1	5.15	1.31	1.25
2	B	140	LYS	C-O	5.10	1.33	1.23
2	B	3	ASP	CG-OD1	5.08	1.37	1.25
1	A	411	LYS	CE-NZ	5.05	1.61	1.49

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	327	ARG	NE-CZ-NH1	-14.15	113.22	120.30
1	A	206	LEU	CB-CG-CD2	8.95	126.22	111.00
1	A	9	SER	O-C-N	-8.74	108.71	122.70
1	A	10	ARG	NE-CZ-NH2	7.85	124.22	120.30
1	A	327	ARG	NH1-CZ-NH2	6.86	126.95	119.40
1	A	9	SER	C-N-CA	6.82	138.76	121.70
1	A	294	TRP	CD1-NE1-CE2	6.57	114.91	109.00
2	B	144	GLU	N-CA-C	6.35	128.14	111.00
2	B	143	GLY	N-CA-C	6.14	128.46	113.10
2	B	3	ASP	CB-CG-OD1	-6.09	112.82	118.30
2	B	165	VAL	CA-CB-CG1	6.00	119.90	110.90
2	B	146	ARG	NE-CZ-NH1	-5.78	117.41	120.30
1	A	13	GLU	OE1-CD-OE2	-5.53	116.67	123.30
2	B	139	PHE	CB-CG-CD2	-5.41	117.01	120.80
1	A	114	PHE	CG-CD2-CE2	-5.40	114.86	120.80
2	B	165	VAL	CA-CB-CG2	-5.31	102.93	110.90
2	B	66	ASP	CB-CG-OD1	5.27	123.05	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	214	GLY	Mainchain
1	A	9	SER	Mainchain

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4384	0	4494	95	24
2	B	1298	0	1280	37	24
3	C	259	0	279	13	0
4	A	1	0	0	0	0
5	A	43	0	30	2	0
6	A	65	0	62	8	0
7	A	4	0	0	0	0
8	B	2	0	0	0	0
All	All	6056	0	6145	135	24

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 11.

All (135) 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:357:ILE:CD1	1:A:357:ILE:CG1	1.75	1.63
6:A:801:HAS:C13	6:A:801:HAS:C12	1.75	1.60
2:B:165:VAL:CG1	2:B:165:VAL:CB	1.86	1.54
1:A:140:LYS:CE	1:A:140:LYS:NZ	1.83	1.41
1:A:262:ASP:HB2	1:A:511:VAL:HG11	1.43	1.00
1:A:562:TRP:HA	2:B:155:LEU:HG	1.55	0.87
6:A:801:HAS:C12	6:A:801:HAS:C14	2.55	0.84
3:C:2:GLU:HG3	3:C:3:GLU:H	1.45	0.81
1:A:233:HIS:NE2	1:A:237:TYR:HE2	1.80	0.80
1:A:233:HIS:NE2	1:A:237:TYR:CE2	2.51	0.77
1:A:357:ILE:HG23	3:C:15:LEU:HD12	1.70	0.73
1:A:233:HIS:HE2	1:A:237:TYR:HE2	1.38	0.72
6:A:801:HAS:C13	6:A:801:HAS:C11	2.67	0.70
1:A:376:HIS:HB3	2:B:124:ASN:ND2	2.06	0.70
1:A:449:ARG:HH12	6:A:801:HAS:CGA	2.04	0.70
1:A:18:LYS:HE2	1:A:502:ALA:O	1.91	0.69
2:B:97:VAL:HG23	2:B:166:VAL:HG12	1.73	0.69
1:A:357:ILE:CD1	1:A:357:ILE:CB	2.71	0.68
1:A:263:PRO:HB3	1:A:520:LEU:HD23	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:365:VAL:HG11	3:C:26:VAL:HG21	1.77	0.66
2:B:165:VAL:CG1	2:B:165:VAL:CG2	2.70	0.66
1:A:52:TYR:CZ	1:A:65:TYR:HD1	2.13	0.66
2:B:69:GLN:HE21	2:B:69:GLN:HA	1.61	0.65
1:A:446:ASN:OD1	2:B:119:GLU:HG3	1.97	0.64
2:B:106:LYS:HG2	2:B:134:THR:HG23	1.80	0.64
1:A:354:LEU:HA	1:A:357:ILE:HD12	1.80	0.63
2:B:69:GLN:HA	2:B:69:GLN:NE2	2.14	0.63
1:A:160:ILE:HG12	1:A:190:TRP:HB3	1.81	0.61
1:A:241:LEU:HA	1:A:244:TYR:HB2	1.83	0.60
1:A:302:THR:O	1:A:305:VAL:HG12	2.02	0.60
1:A:220:ASP:HB3	1:A:223:VAL:HG22	1.84	0.60
1:A:227:LEU:O	1:A:230:TRP:HB3	2.03	0.59
2:B:69:GLN:HE21	2:B:69:GLN:CA	2.15	0.58
1:A:325:ARG:HA	1:A:329:GLY:H	1.67	0.58
1:A:359:GLY:HA3	1:A:388:GLN:NE2	2.19	0.57
2:B:122:ASN:HD21	3:C:33:ARG:HB2	1.69	0.57
1:A:122:LEU:HB2	1:A:123:PRO:HD3	1.87	0.57
2:B:9:LYS:HE2	2:B:9:LYS:HA	1.86	0.57
2:B:147:ILE:HD11	2:B:164:ILE:HG13	1.86	0.57
1:A:233:HIS:O	1:A:236:VAL:HG22	2.06	0.55
1:A:294:TRP:CZ2	1:A:544:PRO:HB2	2.42	0.55
3:C:2:GLU:HG3	3:C:3:GLU:N	2.20	0.55
1:A:357:ILE:HG23	3:C:15:LEU:CD1	2.37	0.54
5:A:800:HEM:HBC2	5:A:800:HEM:HMC2	1.89	0.54
1:A:398:MET:O	1:A:401:LEU:HB2	2.08	0.54
1:A:52:TYR:CZ	1:A:65:TYR:CD1	2.96	0.53
1:A:272:PHE:CZ	1:A:308:PRO:HB2	2.43	0.53
1:A:127:ASN:HD21	1:A:562:TRP:HD1	1.56	0.53
2:B:123:ILE:HD11	2:B:137:TYR:CD1	2.44	0.53
1:A:230:TRP:C	1:A:230:TRP:CD1	2.83	0.53
1:A:272:PHE:O	1:A:276:SER:OG	2.19	0.52
3:C:4:LYS:HE3	3:C:6:LYS:HG2	1.91	0.52
1:A:44:LEU:HD23	1:A:471:VAL:HA	1.92	0.52
1:A:365:VAL:HG22	3:C:23:TRP:HA	1.91	0.52
1:A:450:ARG:NH2	2:B:151:GLN:OE1	2.41	0.52
1:A:140:LYS:HE3	1:A:212:SER:HA	1.92	0.51
1:A:335:TRP:O	1:A:339:LEU:HD22	2.10	0.51
1:A:298:HIS:CE1	1:A:541:ALA:HB1	2.46	0.51
1:A:127:ASN:ND2	1:A:562:TRP:HD1	2.09	0.51
1:A:381:VAL:HB	1:A:382:PRO:HD3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:VAL:HG11	6:A:801:HAS:HHD	1.94	0.50
1:A:189:PHE:CZ	1:A:238:PHE:CD2	3.00	0.50
1:A:282:HIS:CD2	6:A:801:HAS:CMD	2.95	0.50
1:A:153:VAL:HG11	1:A:200:LEU:HD23	1.93	0.50
1:A:450:ARG:O	2:B:157:HIS:CD2	2.64	0.49
3:C:20:LEU:O	3:C:24:LEU:HB2	2.13	0.49
1:A:24:PHE:CE1	1:A:110:TRP:HB2	2.47	0.49
1:A:280:GLY:HA3	1:A:542:TYR:OH	2.12	0.49
5:A:800:HEM:HBC2	5:A:800:HEM:CMC	2.43	0.49
1:A:354:LEU:HA	1:A:357:ILE:CD1	2.43	0.48
1:A:441:TRP:HB3	1:A:466:PRO:HB3	1.95	0.48
1:A:465:VAL:HG23	1:A:466:PRO:HD3	1.95	0.48
1:A:160:ILE:HD13	1:A:194:PHE:HB2	1.95	0.48
1:A:247:ILE:HD11	1:A:392:LEU:HD11	1.96	0.48
1:A:387:LEU:HD21	1:A:436:ALA:HB2	1.96	0.48
2:B:74:THR:HG23	2:B:78:GLN:OE1	2.13	0.47
2:B:84:LEU:HB3	2:B:91:GLN:HB3	1.97	0.47
2:B:157:HIS:C	2:B:159:ASN:H	2.18	0.47
3:C:2:GLU:CG	3:C:3:GLU:H	2.20	0.47
1:A:183:THR:O	1:A:187:VAL:HG23	2.15	0.46
1:A:202:LEU:O	1:A:206:LEU:HB2	2.14	0.46
2:B:139:PHE:CD2	2:B:145:TYR:CE1	3.03	0.46
1:A:120:ALA:HB2	1:A:148:GLY:HA3	1.98	0.46
2:B:122:ASN:ND2	3:C:33:ARG:HB2	2.30	0.46
1:A:330:ARG:HB2	1:A:334:GLY:HA3	1.98	0.46
1:A:384:HIS:O	1:A:388:GLN:HG2	2.16	0.45
2:B:141:ARG:HG3	2:B:142:PRO:HD2	1.98	0.45
1:A:130:THR:HB	1:A:562:TRP:CG	2.51	0.45
1:A:294:TRP:HZ2	1:A:544:PRO:HB2	1.80	0.45
1:A:251:LEU:N	1:A:252:PRO:CD	2.80	0.45
3:C:9:LEU:HD23	3:C:12:ILE:HD11	1.99	0.45
1:A:254:GLN:NE2	1:A:343:ASN:HD21	2.14	0.44
1:A:325:ARG:HA	1:A:329:GLY:N	2.31	0.44
1:A:199:GLY:HA3	1:A:230:TRP:CG	2.53	0.44
1:A:357:ILE:HB	1:A:358:PRO:HD3	1.98	0.44
1:A:429:PHE:O	1:A:433:MET:HG2	2.17	0.44
2:B:66:ASP:HA	2:B:67:PRO:HD3	1.77	0.44
1:A:287:ASP:OD2	2:B:126:GLU:HG3	2.18	0.44
2:B:147:ILE:HB	2:B:162:GLY:HA3	2.00	0.44
1:A:437:VAL:O	1:A:441:TRP:HB2	2.18	0.43
1:A:233:HIS:H	1:A:234:PRO:HD2	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:382:PRO:HA	1:A:385:PHE:CE2	2.53	0.43
2:B:93:ASN:HA	2:B:94:PRO:HA	1.75	0.43
1:A:41:PHE:CZ	1:A:55:LEU:HD13	2.53	0.43
1:A:401:LEU:HG	1:A:405:LEU:HD22	2.00	0.43
1:A:143:TRP:HA	1:A:212:SER:HB2	2.01	0.43
1:A:402:TYR:OH	1:A:421:GLY:HA3	2.18	0.43
2:B:139:PHE:HD2	2:B:145:TYR:CE1	2.37	0.43
1:A:426:TRP:O	1:A:430:LEU:HB2	2.19	0.43
1:A:343:ASN:HA	1:A:344:PRO:HD2	1.94	0.42
1:A:435:MET:HG2	1:A:439:LEU:CD2	2.50	0.42
2:B:96:GLU:HA	2:B:165:VAL:O	2.19	0.42
1:A:262:ASP:HB2	1:A:511:VAL:CG1	2.32	0.42
2:B:69:GLN:NE2	2:B:69:GLN:CA	2.78	0.42
2:B:156:GLY:HA2	2:B:158:GLN:HE22	1.84	0.42
1:A:384:HIS:CE1	6:A:801:HAS:CHC	3.03	0.42
2:B:97:VAL:O	2:B:166:VAL:HA	2.20	0.42
1:A:339:LEU:HB3	1:A:340:PRO:CD	2.49	0.42
1:A:41:PHE:CE1	1:A:55:LEU:HD13	2.55	0.41
1:A:254:GLN:NE2	1:A:343:ASN:ND2	2.68	0.41
1:A:517:ASP:O	1:A:521:VAL:HG23	2.20	0.41
2:B:165:VAL:CG1	2:B:165:VAL:HB	2.26	0.41
1:A:385:PHE:CE2	6:A:801:HAS:HAA1	2.56	0.41
1:A:368:SER:OG	3:C:23:TRP:NE1	2.48	0.41
1:A:427:LEU:O	1:A:431:GLY:N	2.43	0.41
1:A:481:LEU:O	1:A:485:ILE:HG13	2.20	0.41
1:A:512:ILE:CG2	2:B:8:HIS:HB2	2.50	0.41
2:B:95:ILE:O	2:B:165:VAL:N	2.39	0.41
1:A:236:VAL:HA	1:A:239:TRP:CE3	2.57	0.40
1:A:325:ARG:HA	1:A:329:GLY:HA3	2.03	0.40
1:A:342:ASP:O	1:A:418:ARG:NH2	2.52	0.40
1:A:475:ILE:O	1:A:479:VAL:HG23	2.21	0.40
2:B:99:GLN:C	2:B:101:ALA:H	2.23	0.40
2:B:97:VAL:HG11	2:B:103:ILE:HG12	2.03	0.40
1:A:556:VAL:CG1	2:B:55:PRO:HG3	2.52	0.40

All (24) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:SER:C	2:B:143:GLY:N[4_445]	1.17	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:SER:O	2:B:143:GLY:CA[4_445]	1.44	0.76
1:A:9:SER:C	2:B:143:GLY:CA[4_445]	1.45	0.75
1:A:10:ARG:N	2:B:143:GLY:N[4_445]	1.57	0.63
1:A:9:SER:CA	2:B:143:GLY:CA[4_445]	1.59	0.61
1:A:13:GLU:OE2	2:B:165:VAL:CG1[4_445]	1.65	0.55
1:A:9:SER:O	2:B:143:GLY:C[4_445]	1.78	0.42
1:A:9:SER:CB	2:B:143:GLY:CA[4_445]	1.79	0.41
1:A:11:VAL:N	2:B:142:PRO:CB[4_445]	1.83	0.37
1:A:9:SER:CA	2:B:143:GLY:N[4_445]	1.85	0.35
1:A:9:SER:O	2:B:143:GLY:O[4_445]	1.85	0.35
1:A:9:SER:O	2:B:143:GLY:N[4_445]	1.89	0.31
1:A:10:ARG:NH1	2:B:139:PHE:CB[4_445]	2.02	0.18
1:A:13:GLU:CD	2:B:165:VAL:CG1[4_445]	2.02	0.18
1:A:9:SER:C	2:B:143:GLY:C[4_445]	2.03	0.17
1:A:9:SER:OG	2:B:142:PRO:O[4_445]	2.03	0.17
1:A:10:ARG:N	2:B:142:PRO:C[4_445]	2.04	0.16
1:A:13:GLU:OE2	2:B:165:VAL:CB[4_445]	2.08	0.12
1:A:9:SER:N	2:B:144:GLU:N[4_445]	2.11	0.09
1:A:9:SER:N	2:B:143:GLY:C[4_445]	2.13	0.07
1:A:13:GLU:OE2	2:B:165:VAL:CA[4_445]	2.15	0.05
1:A:9:SER:CA	2:B:142:PRO:C[4_445]	2.17	0.03
1:A:9:SER:CA	2:B:142:PRO:O[4_445]	2.18	0.02
1:A:13:GLU:OE2	2:B:165:VAL:C[4_445]	2.19	0.01

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	552/568 (97%)	516 (94%)	34 (6%)	2 (0%)	34 72
2	B	164/166 (99%)	152 (93%)	11 (7%)	1 (1%)	25 65
3	C	31/33 (94%)	30 (97%)	1 (3%)	0	100 100
All	All	747/767 (97%)	698 (93%)	46 (6%)	3 (0%)	34 72

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	331	GLY
1	A	330	ARG
2	B	87	ALA

5.3.2 Protein sidechains (i)

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	450/462 (97%)	433 (96%)	17 (4%)	33	58
2	B	136/136 (100%)	124 (91%)	12 (9%)	10	34
3	C	26/26 (100%)	24 (92%)	2 (8%)	13	39
All	All	612/624 (98%)	581 (95%)	31 (5%)	24	51

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

Mol	Chain	Res	Type
1	A	9	SER
1	A	36	SER
1	A	48	ASN
1	A	76	ASN
1	A	133	TYR
1	A	168	ARG
1	A	230	TRP
1	A	305	VAL
1	A	339	LEU
1	A	354	LEU
1	A	369	PHE
1	A	401	LEU
1	A	430	LEU
1	A	439	LEU
1	A	452	TYR
1	A	472	LEU
1	A	513	SER
2	B	5	HIS

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Mol	Chain	Res	Type
2	B	19	LEU
2	B	26	LEU
2	B	69	GLN
2	B	71	VAL
2	B	77	ASN
2	B	97	VAL
2	B	111	ASP
2	B	134	THR
2	B	139	PHE
2	B	142	PRO
2	B	159	ASN
3	C	3	GLU
3	C	13	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	GLN
1	A	48	ASN
1	A	76	ASN
1	A	254	GLN
1	A	298	HIS
1	A	554	ASN
2	B	69	GLN
2	B	77	ASN
2	B	122	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 8 ligands modelled in this entry, 5 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
8	CUA	B	802	2	0,1,1	-	-	-		
6	HAS	A	801	1	57,72,72	4.62	19 (33%)	48,109,109	3.52	13 (27%)
5	HEM	A	800	1	41,50,50	1.97	7 (17%)	45,82,82	1.56	5 (11%)

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
6	HAS	A	801	1	-	10/40/122/122	-
5	HEM	A	800	1	-	3/12/54/54	-

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	801	HAS	C2D-C3D	18.02	1.53	1.36
6	A	801	HAS	CHA-C4D	14.44	1.62	1.37
6	A	801	HAS	C3C-C2C	10.79	1.55	1.40
6	A	801	HAS	CMC-C2C	9.05	1.73	1.51
6	A	801	HAS	CHB-C1D	8.90	1.56	1.39
6	A	801	HAS	C3B-C2B	8.33	1.53	1.34
5	A	800	HEM	C3D-C2D	7.78	1.53	1.36
6	A	801	HAS	CHB-C1B	7.35	1.53	1.39
6	A	801	HAS	C12-C13	6.73	1.75	1.53
6	A	801	HAS	O1A-CGA	6.01	1.42	1.22
6	A	801	HAS	C2A-C3A	5.97	1.55	1.37
6	A	801	HAS	C3C-C4C	5.17	1.53	1.40
6	A	801	HAS	CHC-C4B	-5.06	1.29	1.37
5	A	800	HEM	C3C-C2C	-4.28	1.34	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	801	HAS	CMD-C2D	4.11	1.51	1.44
6	A	801	HAS	O11-C11	4.05	1.51	1.42
5	A	800	HEM	C3C-CAC	3.79	1.55	1.47
6	A	801	HAS	CHD-C4C	-3.43	1.32	1.39
6	A	801	HAS	C3C-CAC	2.97	1.53	1.47
5	A	800	HEM	CAB-C3B	2.71	1.54	1.47
6	A	801	HAS	CHA-C1A	2.62	1.46	1.40
6	A	801	HAS	O1D-CGD	2.42	1.30	1.22
5	A	800	HEM	CMD-C2D	2.25	1.55	1.50
5	A	800	HEM	FE-ND	2.20	2.07	1.96
6	A	801	HAS	C14-C15	2.18	1.38	1.33
5	A	800	HEM	CMB-C2B	2.13	1.55	1.50

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	A	801	HAS	C4B-C3B-C2B	-18.00	90.30	108.79
6	A	801	HAS	C4B-CHC-C1C	9.90	145.46	118.67
5	A	800	HEM	C4D-ND-C1D	5.51	110.77	105.07
6	A	801	HAS	C12-C13-C14	-4.17	101.22	112.23
6	A	801	HAS	CBA-CAA-C2A	-4.11	105.68	112.60
6	A	801	HAS	CMC-C2C-C3C	-4.03	117.15	124.68
6	A	801	HAS	C13-C12-C11	-3.87	108.53	114.35
6	A	801	HAS	CAD-CBD-CGD	-3.52	106.03	113.60
6	A	801	HAS	C25-C23-C24	3.49	121.14	115.27
6	A	801	HAS	CMA-C3A-C2A	3.38	131.32	124.94
6	A	801	HAS	OMD-CMD-C2D	-3.14	119.99	125.03
6	A	801	HAS	CBD-CAD-C3D	2.73	120.13	112.62
5	A	800	HEM	C1B-NB-C4B	2.62	107.78	105.07
6	A	801	HAS	C27-C19-C20	2.43	119.36	115.27
5	A	800	HEM	C4C-CHD-C1D	2.38	125.69	122.56
5	A	800	HEM	CAD-C3D-C4D	2.08	128.29	124.66
5	A	800	HEM	CBB-CAB-C3B	-2.06	117.38	127.62
6	A	801	HAS	C16-C17-C18	-2.03	105.22	111.88

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	800	HEM	C2B-C3B-CAB-CBB
5	A	800	HEM	C4B-C3B-CAB-CBB
6	A	801	HAS	O11-C11-C3B-C2B

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Mol	Chain	Res	Type	Atoms
6	A	801	HAS	C12-C11-C3B-C2B
6	A	801	HAS	O11-C11-C12-C13
6	A	801	HAS	C25-C23-C24-C28
6	A	801	HAS	CAA-CBA-CGA-O1A
6	A	801	HAS	CAD-CBD-CGD-O1D
6	A	801	HAS	CAD-CBD-CGD-O2D
6	A	801	HAS	CAA-CBA-CGA-O2A
6	A	801	HAS	C22-C23-C24-C28
6	A	801	HAS	C2D-C3D-CAD-CBD
5	A	800	HEM	CAA-CBA-CGA-O2A

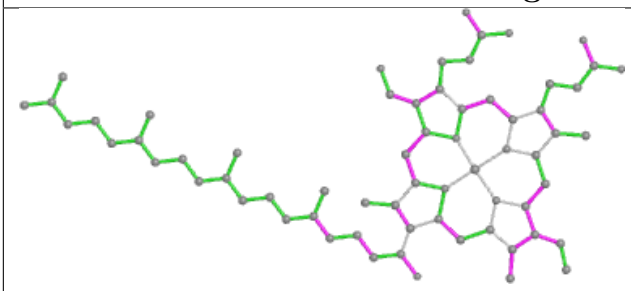
There are no ring outliers.

2 monomers are involved in 10 short contacts:

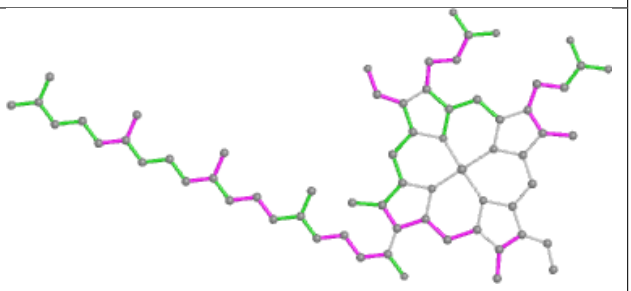
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	801	HAS	8	0
5	A	800	HEM	2	0

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 then 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.

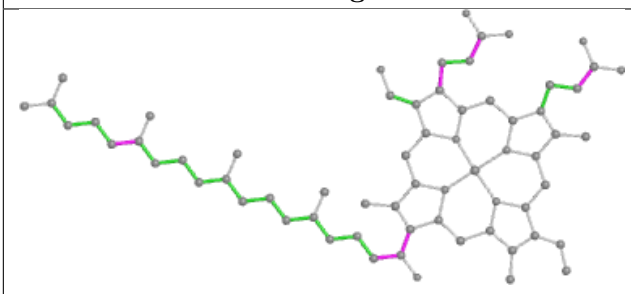
Ligand HAS A 801



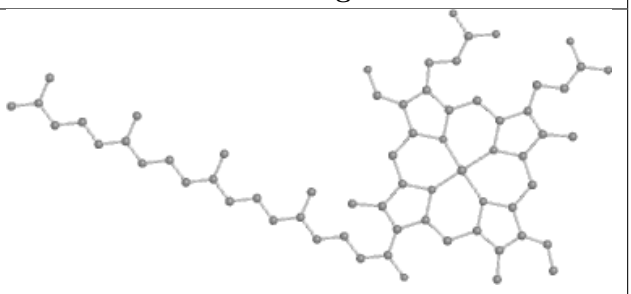
Bond lengths



Bond angles

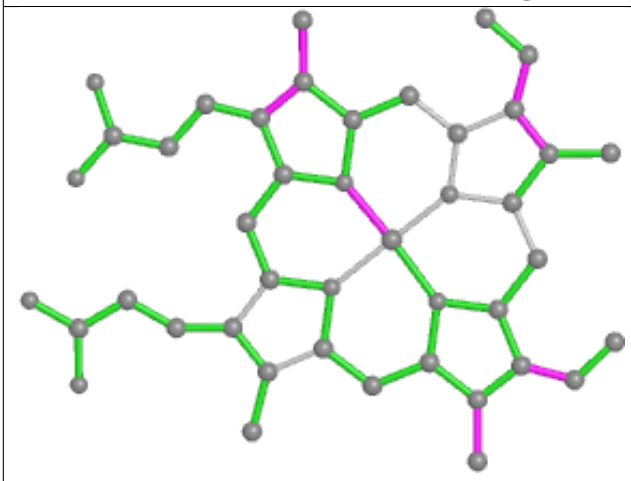


Torsions

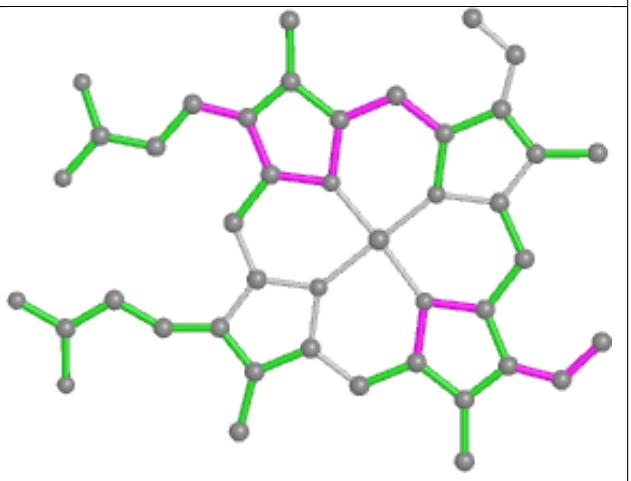


Rings

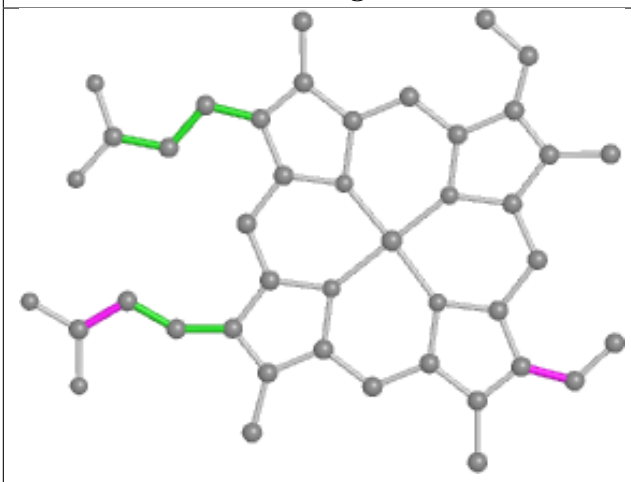
Ligand HEM A 800



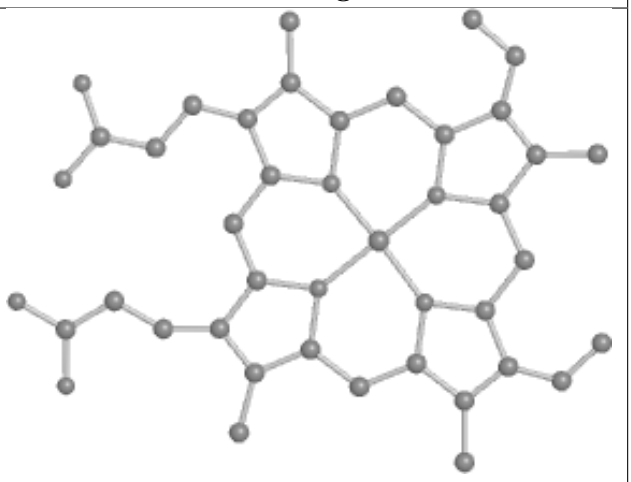
Bond lengths



Bond angles



Torsions



Rings

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, 95^{th} 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	554/568 (97%)	0.10	21 (3%) 40 32	123, 123, 123, 123	0
2	B	166/166 (100%)	0.13	6 (3%) 42 34	123, 123, 123, 123	0
3	C	33/33 (100%)	-0.42	1 (3%) 50 39	123, 123, 123, 123	0
All	All	753/767 (98%)	0.08	28 (3%) 41 33	123, 123, 123, 123	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	141	GLY	4.5
1	A	140	LYS	3.9
2	B	136	ARG	3.6
1	A	142	HIS	3.4
1	A	212	SER	3.1
1	A	342	ASP	3.0
1	A	57	ARG	2.9
1	A	502	ALA	2.9
1	A	409	THR	2.7
2	B	52	ARG	2.7
2	B	61	GLU	2.5
2	B	104	VAL	2.5
1	A	215	LEU	2.4
1	A	203	GLU	2.3
1	A	219	VAL	2.3
1	A	410	GLY	2.3
3	C	2	GLU	2.3
2	B	3	ASP	2.3
1	A	146	TYR	2.2
1	A	558	GLY	2.2
1	A	519	ARG	2.2
1	A	207	PHE	2.1
1	A	464	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	14	ALA	2.1
1	A	48	ASN	2.0
1	A	548	GLN	2.0
2	B	79	TYR	2.0
1	A	12	TYR	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 monosaccharides in this entry.

6.4 Ligands (i)

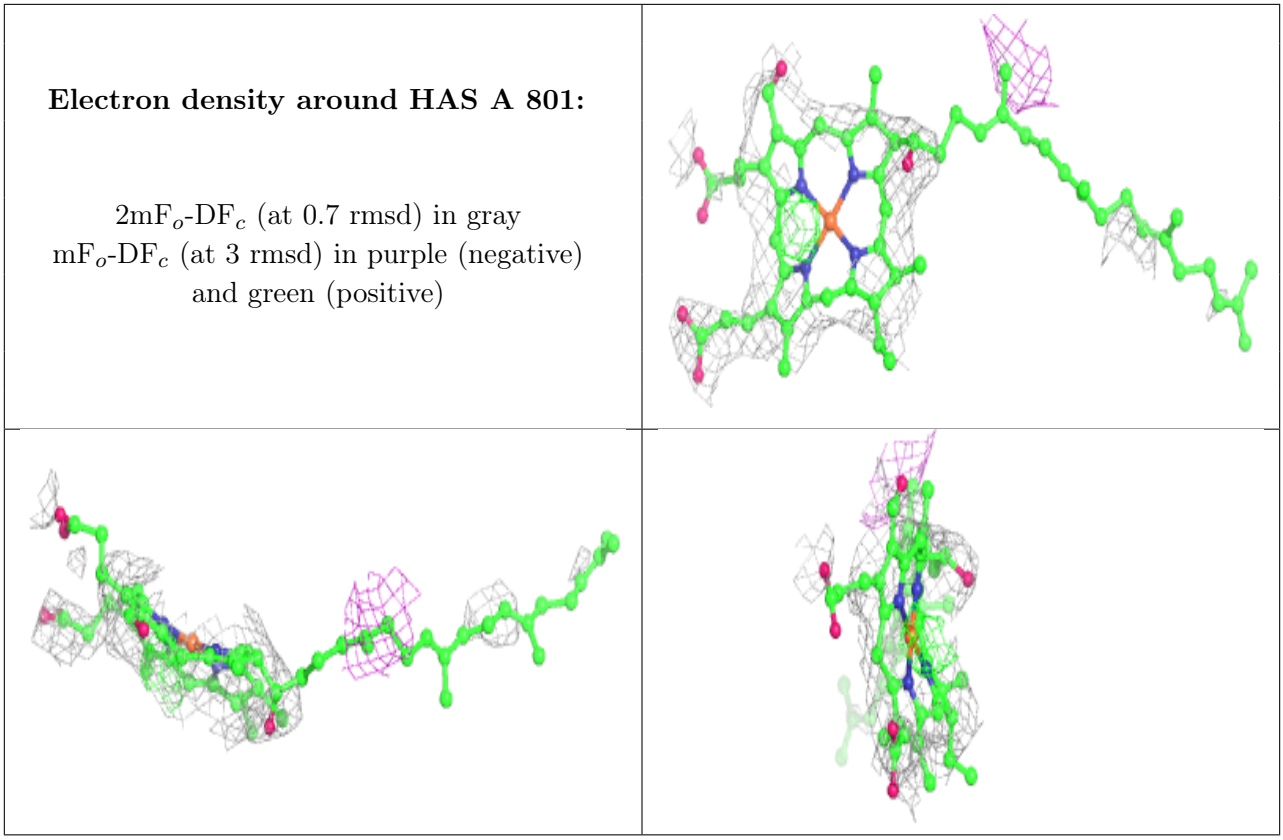
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	CU	A	803	1/1	0.75	0.21	123,123,123,123	0
7	XE	A	563	1/1	0.77	0.75	123,123,123,123	1
6	HAS	A	801	65/65	0.91	0.38	123,123,123,123	0
5	HEM	A	800	43/43	0.91	0.34	123,123,123,123	0
7	XE	A	564	1/1	0.93	0.33	123,123,123,123	1
7	XE	A	565	1/1	0.94	1.30	123,123,123,123	1
8	CUA	B	802	2/2	0.95	0.14	123,123,123,123	0
7	XE	A	566	1/1	0.96	0.94	123,123,123,123	1

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.

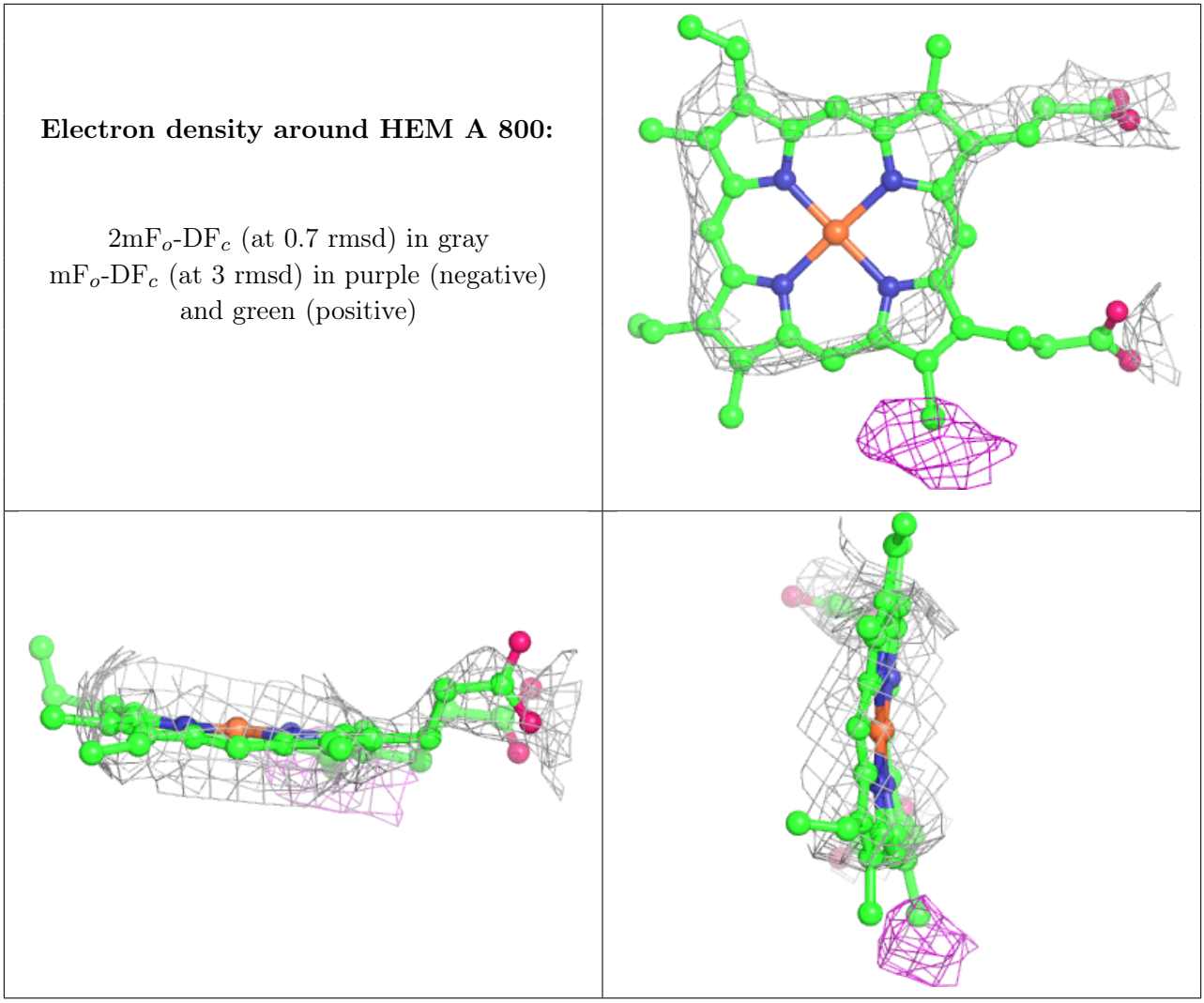
Electron density around HAS A 801:

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



Electron density around HEM A 800:

$2mF_o$-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.