



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 09:31 am BST

PDB ID : 3QJQ
Title : The structure of and photolytic induced changes of carbon monoxide binding to the cytochrome ba3-oxidase from *Thermus thermophilus*
Authors : Liu, B.; Zhang, Y.; Sage, J.T.; Doukov, T.; Chen, Y.; Stout, C.D.; Fee, J.A.
Deposited on : 2011-01-30
Resolution : 2.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The 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 : 2.11
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) : 2.11

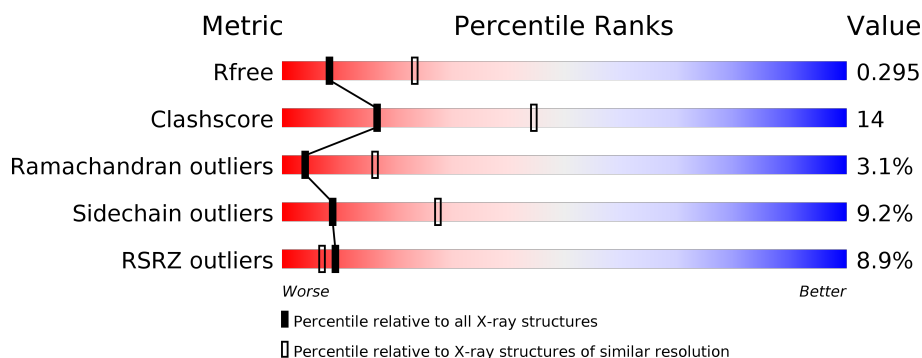
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on 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	568	<div> <div>8%</div> <div> <div></div> <div>65%</div> <div>28%</div> <div>• •</div> </div> </div>
2	B	168	<div> <div>11%</div> <div> <div></div> <div>63%</div> <div>32%</div> <div>• ••</div> </div> </div>
3	C	34	<div> <div>12%</div> <div> <div></div> <div>44%</div> <div>35%</div> <div>18%</div> <div>•</div> </div> </div>

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6037 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	552	Total	C	N	O	S	0	0	0
			4367	2962	699	690	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 (I) ION (three-letter code: CU1) (formula: Cu).

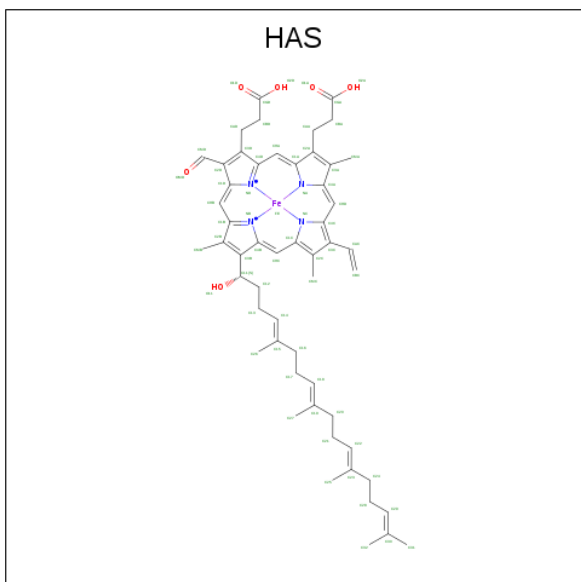
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cu	0	0
			1	1		

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



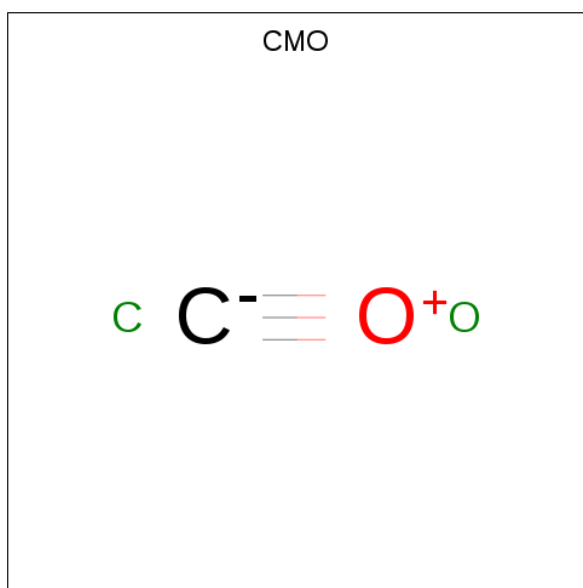
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 6 is HEME-AS (three-letter code: HAS) (formula: $\text{C}_{54}\text{H}_{64}\text{FeN}_4\text{O}_6$).



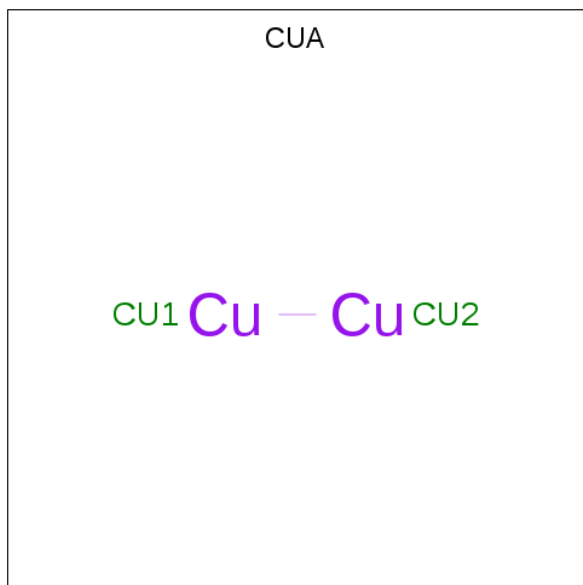
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	Fe	N	O	0	0
			65	54	1	4	6		

- Molecule 7 is CARBON MONOXIDE (three-letter code: CMO) (formula: CO).



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

- Molecule 8 is DINUCLEAR COPPER ION (three-letter code: CUA) (formula: Cu₂).

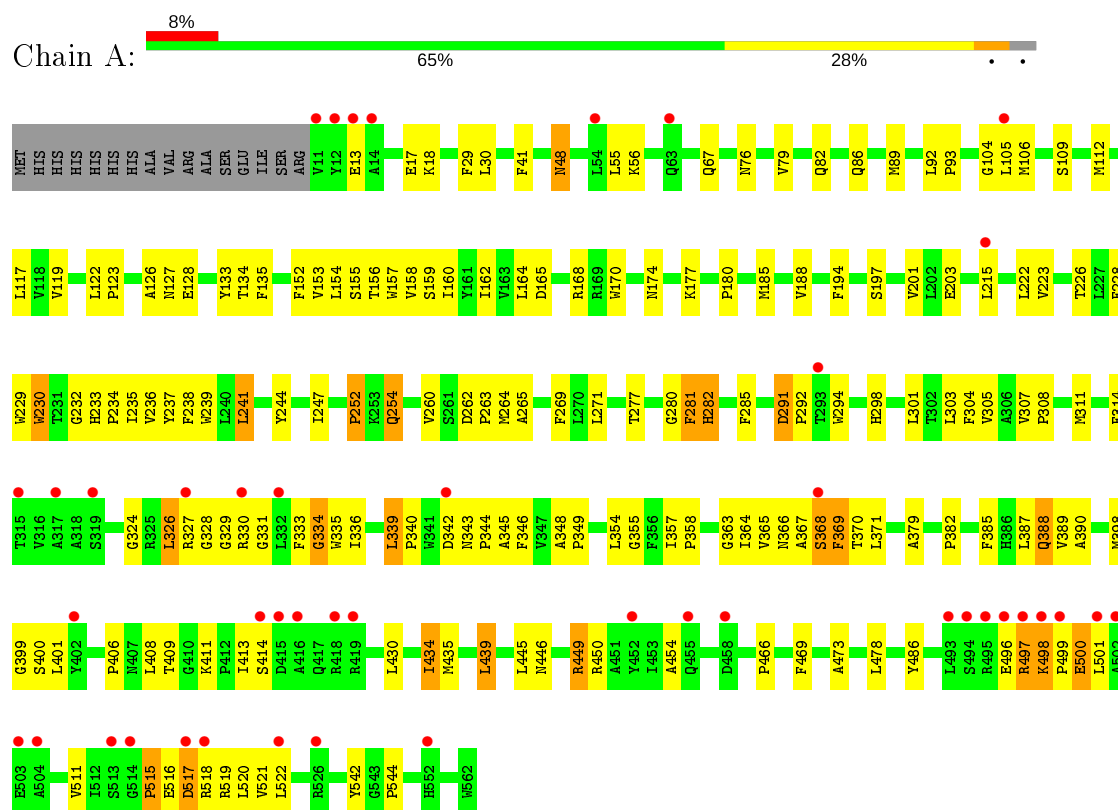


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	1	Total	Cu	0	0
			2	2		

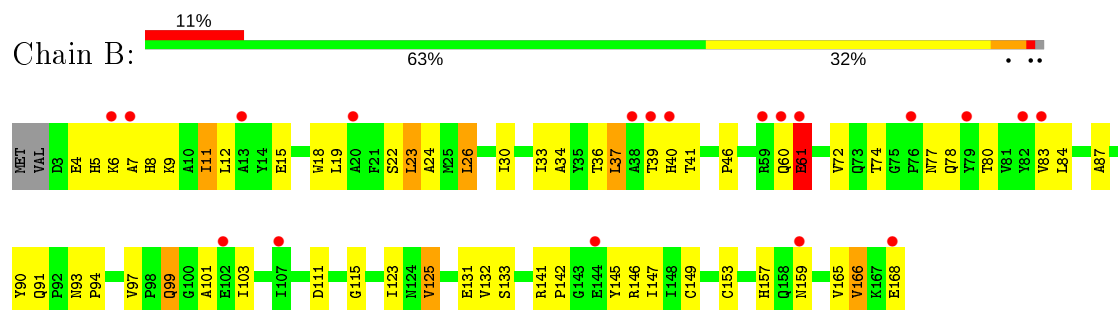
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

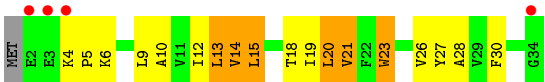
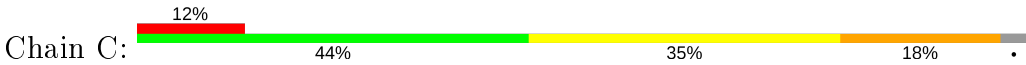
• Molecule 1: Cytochrome c oxidase subunit 1



• Molecule 2: Cytochrome c oxidase subunit 2



• Molecule 3: Cytochrome c oxidase polypeptide 2A



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	108.77Å 108.77Å 164.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.88 – 2.90 19.88 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.5 (19.88-2.90) 99.5 (19.88-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.02	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 2.88Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.229 , 0.284 0.240 , 0.295	Depositor DCC
R_{free} test set	1144 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	90.5	Xtriage
Anisotropy	0.333	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 75.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6037	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CMO, HEM, HAS, CUA, CU1

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	0.64	0/4524	0.76	0/6210
2	B	0.63	1/1335 (0.1%)	0.74	2/1822 (0.1%)
3	C	0.86	0/265	0.89	0/359
All	All	0.65	1/6124 (0.0%)	0.76	2/8391 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	153	CYS	CB-SG	-6.01	1.72	1.82

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	26	LEU	CB-CG-CD2	-9.56	94.75	111.00
2	B	26	LEU	CD1-CG-CD2	-5.30	94.61	110.50

There are no chirality outliers.

There are no planarity outliers.

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	4367	0	4476	132	1
2	B	1298	0	1280	39	1

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	259	0	279	13	0
4	A	1	0	0	0	0
5	A	43	0	30	1	0
6	A	65	0	62	9	0
7	A	2	0	0	1	0
8	B	2	0	0	0	0
All	All	6037	0	6127	171	1

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 14.

All (171) 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:233:HIS:NE2	1:A:237:TYR:HE2	0.93	1.41
1:A:233:HIS:NE2	1:A:237:TYR:CE2	1.79	1.24
1:A:233:HIS:CE1	1:A:237:TYR:HE2	1.81	0.97
1:A:233:HIS:CE1	1:A:237:TYR:CE2	2.54	0.95
2:B:26:LEU:O	2:B:30:ILE:HG13	1.69	0.93
2:B:101:ALA:O	2:B:103:ILE:HD12	1.75	0.87
1:A:449:ARG:HD2	1:A:450:ARG:HG3	1.60	0.81
1:A:368:SER:HB3	2:B:33:ILE:HD11	1.64	0.80
1:A:282:HIS:HA	1:A:285:PHE:CE1	2.16	0.80
1:A:106:MET:O	1:A:109:SER:OG	2.04	0.76
1:A:160:ILE:HD13	1:A:194:PHE:HB2	1.69	0.75
1:A:388:GLN:HE21	1:A:388:GLN:HA	1.50	0.74
1:A:364:ILE:O	1:A:368:SER:OG	2.06	0.74
1:A:233:HIS:NE2	1:A:237:TYR:CD2	2.55	0.74
6:A:801:HAS:ND	7:A:563:CMO:O	2.23	0.71
1:A:388:GLN:HA	1:A:388:GLN:NE2	2.06	0.71
2:B:5:HIS:C	2:B:7:ALA:H	1.95	0.70
1:A:236:VAL:HG12	1:A:239:TRP:CZ3	2.27	0.70
1:A:158:VAL:HG12	1:A:162:ILE:HD12	1.74	0.70
1:A:366:ASN:HB3	6:A:801:HAS:CMD	2.22	0.70
1:A:303:LEU:HB3	2:B:26:LEU:HG	1.73	0.69
1:A:230:TRP:O	1:A:230:TRP:CD1	2.46	0.68
1:A:156:THR:O	1:A:160:ILE:HG13	1.94	0.68
1:A:388:GLN:HE21	1:A:388:GLN:CA	2.06	0.68
1:A:105:LEU:C	1:A:162:ILE:HD11	2.14	0.67
1:A:48:ASN:ND2	1:A:454:ALA:HA	2.10	0.66
1:A:236:VAL:HB	6:A:801:HAS:C3C	2.26	0.66

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:336:ILE:HD13	6:A:801:HAS:H311	1.76	0.66
2:B:142:PRO:HA	2:B:166:VAL:CG2	2.27	0.65
3:C:13:LEU:O	3:C:14:VAL:C	2.35	0.65
2:B:18:TRP:CE3	3:C:12:ILE:HD12	2.32	0.64
2:B:34:ALA:O	2:B:37:LEU:HB2	1.97	0.64
1:A:382:PRO:HA	1:A:385:PHE:CZ	2.34	0.63
1:A:330:ARG:HB2	1:A:334:GLY:HA3	1.81	0.63
1:A:262:ASP:HB3	1:A:263:PRO:HD3	1.80	0.62
1:A:230:TRP:O	1:A:230:TRP:HD1	1.81	0.62
1:A:29:PHE:CE1	1:A:401:LEU:HD11	2.34	0.62
1:A:449:ARG:HH12	6:A:801:HAS:CGA	2.13	0.61
2:B:5:HIS:CE1	2:B:8:HIS:CD2	2.88	0.61
1:A:197:SER:O	1:A:201:VAL:HG23	2.00	0.61
1:A:29:PHE:CZ	1:A:401:LEU:HD11	2.36	0.61
1:A:122:LEU:HB2	1:A:123:PRO:HD3	1.82	0.60
1:A:385:PHE:O	1:A:389:VAL:HG12	2.01	0.60
1:A:367:ALA:HA	6:A:801:HAS:OMD	2.02	0.60
1:A:366:ASN:HB3	6:A:801:HAS:C2D	2.32	0.59
2:B:60:GLN:O	2:B:61:GLU:HB3	2.02	0.59
2:B:99:GLN:OE1	2:B:168:GLU:HG3	2.01	0.59
1:A:357:ILE:HB	1:A:358:PRO:HD3	1.85	0.59
1:A:233:HIS:O	1:A:236:VAL:HG22	2.03	0.58
1:A:339:LEU:HB3	1:A:340:PRO:HD2	1.86	0.58
2:B:5:HIS:CE1	2:B:8:HIS:HD2	2.22	0.58
1:A:230:TRP:C	1:A:230:TRP:CD1	2.77	0.58
1:A:262:ASP:HB2	1:A:511:VAL:HG11	1.86	0.58
1:A:370:THR:HB	3:C:30:PHE:CZ	2.39	0.57
1:A:241:LEU:HA	1:A:244:TYR:HB2	1.84	0.57
1:A:234:PRO:HG3	1:A:277:THR:HA	1.85	0.57
1:A:92:LEU:HB2	1:A:93:PRO:HD3	1.86	0.57
3:C:6:LYS:O	3:C:10:ALA:N	2.24	0.56
1:A:280:GLY:C	1:A:282:HIS:H	2.10	0.55
2:B:94:PRO:HB2	2:B:165:VAL:HG23	1.86	0.55
1:A:260:VAL:HB	2:B:11:ILE:HD13	1.88	0.55
2:B:141:ARG:O	2:B:145:TYR:OH	2.17	0.54
1:A:371:LEU:HG	3:C:26:VAL:HG12	1.90	0.54
1:A:498:LYS:HD2	1:A:500:GLU:HG2	1.90	0.54
1:A:18:LYS:HG2	1:A:408:LEU:HD23	1.89	0.53
1:A:498:LYS:HG3	1:A:501:LEU:H	1.74	0.53
1:A:254:GLN:O	1:A:340:PRO:HG3	2.09	0.53
1:A:112:MET:HB3	1:A:155:SER:HB2	1.91	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:388:GLN:CA	1:A:388:GLN:NE2	2.69	0.52
2:B:5:HIS:C	2:B:7:ALA:N	2.62	0.52
1:A:379:ALA:C	1:A:382:PRO:HD2	2.30	0.52
1:A:365:VAL:HG22	3:C:23:TRP:HA	1.92	0.52
1:A:119:VAL:O	1:A:123:PRO:HD3	2.11	0.51
1:A:326:LEU:C	1:A:328:GLY:H	2.13	0.51
1:A:229:TRP:HZ3	1:A:232:GLY:O	1.94	0.51
1:A:281:PHE:H	1:A:298:HIS:CD2	2.29	0.51
2:B:19:LEU:HA	2:B:22:SER:HB2	1.92	0.50
2:B:9:LYS:HE2	2:B:9:LYS:HA	1.92	0.50
1:A:304:PHE:O	1:A:307:VAL:HG23	2.12	0.50
1:A:86:GLN:HG2	1:A:159:SER:OG	2.11	0.50
1:A:285:PHE:CD2	1:A:369:PHE:CD2	2.99	0.50
1:A:406:PRO:HG3	1:A:413:ILE:HD12	1.94	0.50
1:A:262:ASP:HB2	1:A:511:VAL:CG1	2.42	0.49
1:A:409:THR:O	1:A:497:ARG:NH1	2.45	0.49
1:A:222:LEU:O	1:A:226:THR:OG1	2.26	0.49
2:B:166:VAL:CG2	2:B:166:VAL:O	2.60	0.49
1:A:153:VAL:O	1:A:156:THR:OG1	2.21	0.49
1:A:382:PRO:HA	1:A:385:PHE:CE2	2.48	0.49
1:A:434:ILE:HG22	1:A:473:ALA:HB2	1.93	0.49
1:A:515:PRO:O	1:A:517:ASP:N	2.46	0.49
1:A:466:PRO:HA	1:A:469:PHE:HD2	1.77	0.49
1:A:105:LEU:O	1:A:162:ILE:HD11	2.13	0.48
2:B:125:VAL:HG21	2:B:133:SER:HB3	1.95	0.48
3:C:9:LEU:HA	3:C:12:ILE:HG12	1.95	0.48
2:B:115:GLY:O	2:B:149:CYS:HA	2.13	0.48
1:A:346:PHE:O	1:A:349:PRO:HD2	2.13	0.48
1:A:398:MET:C	1:A:400:SER:H	2.17	0.48
3:C:20:LEU:O	3:C:21:VAL:C	2.52	0.48
1:A:498:LYS:HG2	1:A:501:LEU:HB3	1.95	0.47
2:B:40:HIS:CE1	2:B:41:THR:HG23	2.49	0.47
1:A:154:LEU:O	1:A:157:TRP:HB2	2.15	0.47
1:A:48:ASN:HD22	1:A:454:ALA:HA	1.78	0.47
1:A:294:TRP:CZ2	1:A:544:PRO:HG2	2.50	0.47
1:A:370:THR:HG21	2:B:36:THR:HB	1.97	0.47
2:B:83:VAL:HG12	2:B:84:LEU:N	2.29	0.46
1:A:134:THR:O	1:A:135:PHE:C	2.53	0.46
1:A:281:PHE:H	1:A:298:HIS:HD2	1.63	0.46
1:A:357:ILE:HG23	3:C:15:LEU:HD13	1.97	0.46
1:A:230:TRP:C	1:A:230:TRP:HD1	2.18	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:241:LEU:HD12	1:A:241:LEU:HA	1.78	0.46
1:A:389:VAL:HG13	1:A:390:ALA:N	2.30	0.46
1:A:126:ALA:O	1:A:128:GLU:HG2	2.16	0.46
1:A:280:GLY:C	1:A:282:HIS:N	2.70	0.46
2:B:166:VAL:HG23	2:B:166:VAL:O	2.16	0.45
1:A:76:ASN:HB3	5:A:800:HEM:CAC	2.47	0.45
1:A:247:ILE:O	1:A:252:PRO:HD3	2.16	0.45
1:A:348:ALA:CB	1:A:399:GLY:HA3	2.46	0.45
1:A:435:MET:HG2	1:A:439:LEU:HD22	1.98	0.45
1:A:185:MET:CE	1:A:265:ALA:HB1	2.47	0.45
1:A:188:VAL:HG21	1:A:269:PHE:HB3	1.99	0.45
1:A:235:ILE:O	1:A:238:PHE:HB3	2.16	0.45
2:B:97:VAL:O	2:B:166:VAL:HA	2.16	0.45
1:A:345:ALA:HA	1:A:399:GLY:O	2.16	0.45
3:C:13:LEU:HD22	3:C:13:LEU:HA	1.49	0.45
1:A:41:PHE:CE2	1:A:55:LEU:HB2	2.52	0.44
1:A:517:ASP:N	1:A:517:ASP:OD2	2.49	0.44
1:A:308:PRO:O	1:A:311:MET:HB2	2.18	0.44
2:B:23:LEU:O	2:B:24:ALA:C	2.51	0.44
2:B:34:ALA:HA	2:B:37:LEU:HD22	1.98	0.44
1:A:354:LEU:HA	1:A:357:ILE:HD12	1.99	0.44
2:B:74:THR:HG21	2:B:80:THR:OG1	2.16	0.44
1:A:371:LEU:HD11	3:C:27:TYR:HA	2.00	0.44
1:A:329:GLY:C	1:A:330:ARG:HG3	2.38	0.44
2:B:142:PRO:HA	2:B:166:VAL:HG21	1.99	0.44
1:A:170:TRP:CH2	1:A:180:PRO:HD3	2.53	0.44
1:A:450:ARG:O	2:B:157:HIS:CD2	2.70	0.44
1:A:241:LEU:HG	1:A:269:PHE:CE2	2.52	0.44
1:A:280:GLY:HA3	1:A:542:TYR:OH	2.17	0.44
1:A:41:PHE:CE1	1:A:55:LEU:HD13	2.52	0.44
3:C:4:LYS:HE3	3:C:9:LEU:HD13	2.00	0.44
1:A:174:ASN:HB3	1:A:177:LYS:HB2	2.01	0.43
1:A:67:GLN:HE22	1:A:127:ASN:H	1.66	0.43
2:B:146:ARG:HG2	2:B:147:ILE:N	2.34	0.43
1:A:324:GLY:HA3	1:A:335:TRP:HB2	2.01	0.43
1:A:264:MET:SD	2:B:15:GLU:HG2	2.59	0.42
1:A:388:GLN:HE21	1:A:388:GLN:N	2.17	0.42
1:A:445:LEU:O	1:A:446:ASN:HB2	2.19	0.42
1:A:389:VAL:HB	6:A:801:HAS:HBC2	2.01	0.42
1:A:369:PHE:O	1:A:370:THR:C	2.58	0.42
1:A:79:VAL:HG13	1:A:152:PHE:CZ	2.54	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:5:HIS:O	2:B:7:ALA:N	2.52	0.42
1:A:301:LEU:HA	1:A:304:PHE:HB2	2.01	0.42
1:A:343:ASN:HA	1:A:344:PRO:HD2	1.96	0.41
1:A:370:THR:HG21	2:B:36:THR:CB	2.50	0.41
1:A:203:GLU:OE2	1:A:228:PHE:HB2	2.20	0.41
1:A:339:LEU:HB3	1:A:340:PRO:CD	2.50	0.41
1:A:363:GLY:HA2	1:A:366:ASN:HB2	2.01	0.41
1:A:291:ASP:HA	1:A:292:PRO:HD2	1.83	0.41
1:A:411:LYS:HG2	1:A:496:GLU:O	2.21	0.41
2:B:131:GLU:HG3	2:B:132:VAL:N	2.35	0.41
2:B:74:THR:HG23	2:B:78:GLN:OE1	2.20	0.41
1:A:517:ASP:O	1:A:521:VAL:HB	2.21	0.41
1:A:119:VAL:O	1:A:123:PRO:CD	2.68	0.41
1:A:355:GLY:O	1:A:387:LEU:HB3	2.21	0.41
1:A:271:LEU:HA	1:A:271:LEU:HD23	1.83	0.40
1:A:314:PHE:HB2	3:C:12:ILE:HG21	2.03	0.40
2:B:90:TYR:O	2:B:93:ASN:HB2	2.22	0.40
2:B:94:PRO:HB2	2:B:165:VAL:CG2	2.49	0.40
1:A:335:TRP:O	1:A:339:LEU:HD22	2.21	0.40
1:A:79:VAL:HG21	1:A:117:LEU:HD13	2.04	0.40
1:A:449:ARG:NH1	6:A:801:HAS:CGA	2.82	0.40

All (1) 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:13:GLU:OE1	2:B:166:VAL:N[4_545]	2.00	0.20

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	550/568 (97%)	470 (86%)	70 (13%)	10 (2%)	8	29
2	B	164/168 (98%)	135 (82%)	23 (14%)	6 (4%)	3	13
3	C	31/34 (91%)	12 (39%)	12 (39%)	7 (23%)	0	0
All	All	745/770 (97%)	617 (83%)	105 (14%)	23 (3%)	4	16

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	6	LYS
3	C	21	VAL
1	A	327	ARG
1	A	331	GLY
1	A	515	PRO
2	B	4	GLU
2	B	87	ALA
3	C	14	VAL
1	A	281	PHE
1	A	500	GLU
1	A	516	GLU
2	B	61	GLU
1	A	414	SER
2	B	46	PRO
3	C	23	TRP
3	C	28	ALA
3	C	5	PRO
3	C	20	LEU
1	A	104	GLY
1	A	334	GLY
3	C	19	ILE
1	A	499	PRO
2	B	123	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	448/462 (97%)	409 (91%)	39 (9%)	10	30
2	B	136/138 (99%)	122 (90%)	14 (10%)	7	22
3	C	26/27 (96%)	23 (88%)	3 (12%)	5	17
All	All	610/627 (97%)	554 (91%)	56 (9%)	9	27

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	GLU
1	A	30	LEU
1	A	48	ASN
1	A	56	LYS
1	A	82	GLN
1	A	89	MET
1	A	133	TYR
1	A	164	LEU
1	A	165	ASP
1	A	168	ARG
1	A	215	LEU
1	A	223	VAL
1	A	230	TRP
1	A	241	LEU
1	A	252	PRO
1	A	254	GLN
1	A	282	HIS
1	A	291	ASP
1	A	305	VAL
1	A	326	LEU
1	A	333	PHE
1	A	339	LEU
1	A	342	ASP
1	A	368	SER
1	A	369	PHE
1	A	388	GLN
1	A	430	LEU
1	A	434	ILE
1	A	439	LEU
1	A	449	ARG
1	A	478	LEU
1	A	486	TYR
1	A	497	ARG
1	A	498	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	517	ASP
1	A	518	ARG
1	A	519	ARG
1	A	520	LEU
1	A	522	LEU
2	B	11	ILE
2	B	12	LEU
2	B	23	LEU
2	B	37	LEU
2	B	39	THR
2	B	61	GLU
2	B	72	VAL
2	B	77	ASN
2	B	91	GLN
2	B	99	GLN
2	B	111	ASP
2	B	125	VAL
2	B	159	ASN
2	B	166	VAL
3	C	13	LEU
3	C	15	LEU
3	C	18	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	GLN
1	A	48	ASN
1	A	63	GLN
1	A	76	ASN
1	A	254	GLN
1	A	298	HIS
1	A	388	GLN
1	A	446	ASN
2	B	5	HIS
2	B	8	HIS
2	B	69	GLN
2	B	73	GLN
2	B	117	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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	0.00	-	-		
7	CMO	A	563	4	0,1,1	0.00	-	-		
5	HEM	A	800	1	27,50,50	2.30	6 (22%)	17,82,82	1.31	2 (11%)
6	HAS	A	801	1	56,72,72	4.20	16 (28%)	50,109,109	4.77	23 (46%)

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	-	5/35/122/122	-
5	HEM	A	800	1	-	0/6/54/54	-

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	801	HAS	C1C-C2C	10.52	1.52	1.38
6	A	801	HAS	C1A-C2A	10.50	1.52	1.38
6	A	801	HAS	C4D-ND	-10.35	1.35	1.49
6	A	801	HAS	C4B-NB	-10.19	1.35	1.49
6	A	801	HAS	C3D-C2D	10.12	1.46	1.34
6	A	801	HAS	C1B-NB	-9.77	1.36	1.49
6	A	801	HAS	C1D-ND	-8.46	1.38	1.49
6	A	801	HAS	CHD-C4A	-7.08	1.43	1.51
6	A	801	HAS	C3C-C2C	6.97	1.50	1.40
6	A	801	HAS	C4A-C3A	6.67	1.47	1.38
6	A	801	HAS	CHD-C4C	-6.23	1.44	1.51
6	A	801	HAS	C2B-C3B	6.09	1.39	1.34
5	A	800	HEM	C3B-C2B	-5.57	1.32	1.40
5	A	800	HEM	C3D-C2D	5.49	1.53	1.37
5	A	800	HEM	C3C-C2C	-4.30	1.34	1.40
5	A	800	HEM	C3C-CAC	3.61	1.55	1.47
6	A	801	HAS	C3C-CAC	3.15	1.54	1.47
6	A	801	HAS	CMD-C2D	3.10	1.50	1.44
5	A	800	HEM	C3B-CAB	3.03	1.54	1.47
6	A	801	HAS	C2A-C3A	3.00	1.46	1.37
6	A	801	HAS	CHA-C4D	-2.82	1.48	1.53
5	A	800	HEM	C1C-C2C	2.09	1.47	1.42

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	801	HAS	CHB-C1D-ND	15.65	130.40	110.94
6	A	801	HAS	CHB-C1B-NB	15.19	129.83	110.94
6	A	801	HAS	CHD-C4C-C3C	-10.77	115.49	129.61
6	A	801	HAS	CMC-C2C-C3C	-10.57	104.90	124.68
6	A	801	HAS	CHA-C4D-ND	8.97	127.66	110.75
6	A	801	HAS	CHC-C4B-NB	7.97	125.79	110.75
6	A	801	HAS	CHC-C1C-C2C	-7.76	116.01	129.45
6	A	801	HAS	C4A-C3A-C2A	-5.45	100.01	105.81
6	A	801	HAS	C4C-CHD-C4A	4.96	124.97	112.87
6	A	801	HAS	CAD-CBD-CGD	-4.84	104.55	112.67
6	A	801	HAS	CHD-C4A-C3A	-4.83	121.45	129.53
6	A	801	HAS	C25-C23-C24	4.22	122.37	115.27
6	A	801	HAS	C4C-C3C-C2C	-3.58	98.90	104.41
6	A	801	HAS	CBA-CAA-C2A	-3.39	106.24	112.49
6	A	801	HAS	C21-C22-C23	-3.10	120.20	127.66
6	A	801	HAS	C13-C12-C11	-3.07	109.74	114.35
6	A	801	HAS	CAA-CBA-CGA	-2.79	107.99	112.67

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	801	HAS	CMA-C3A-C2A	2.71	130.05	124.94
5	A	800	HEM	CMD-C2D-C3D	2.67	129.97	124.94
6	A	801	HAS	OMD-CMD-C2D	-2.66	121.02	124.39
5	A	800	HEM	CMD-C2D-C1D	-2.63	124.42	128.46
6	A	801	HAS	C26-C15-C16	2.59	119.63	115.27
6	A	801	HAS	C27-C19-C20	2.50	119.47	115.27
6	A	801	HAS	C13-C14-C15	-2.13	122.54	127.66
6	A	801	HAS	C25-C23-C22	-2.13	118.22	123.68

There are no chirality outliers.

All (5) torsion outliers are listed below:

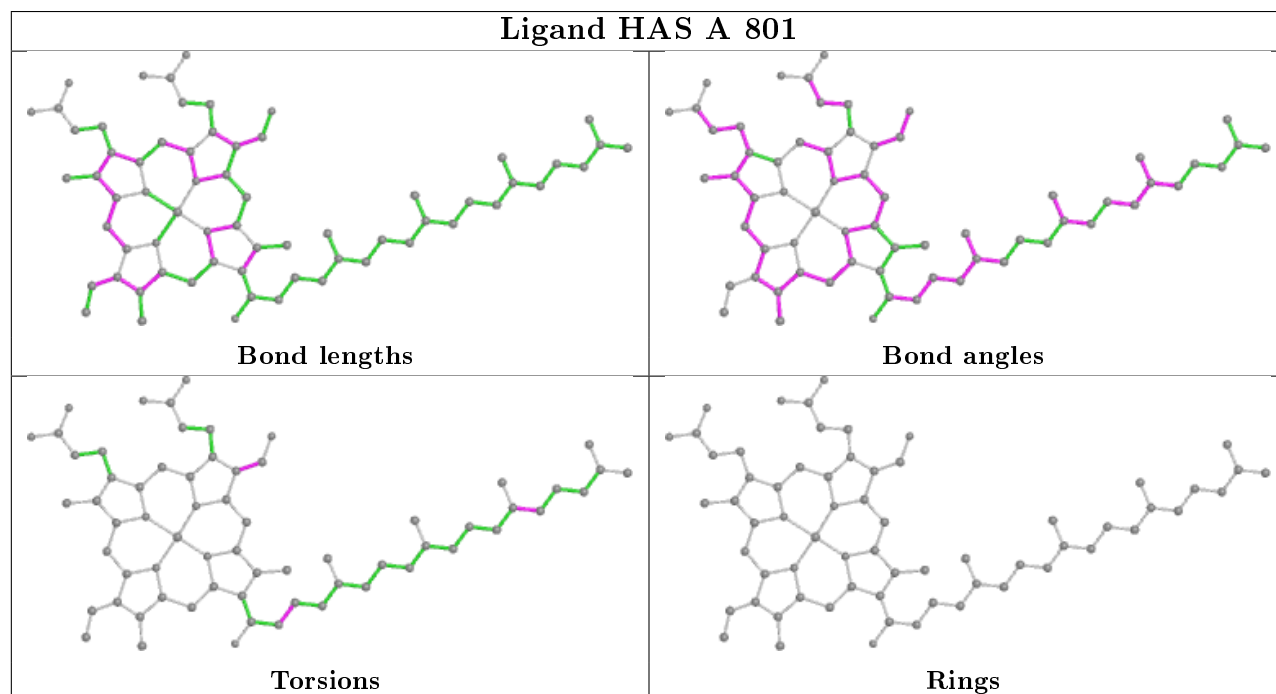
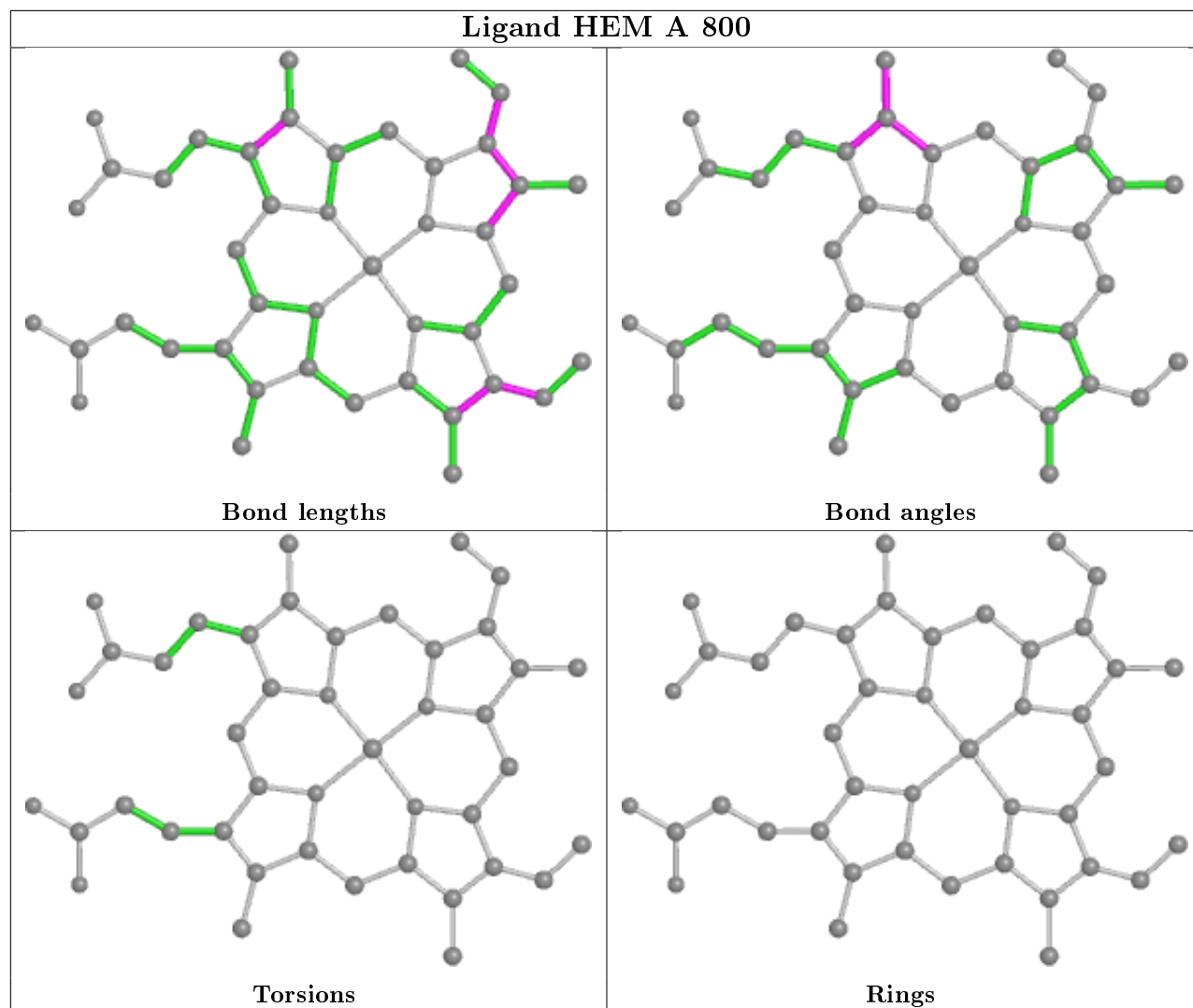
Mol	Chain	Res	Type	Atoms
6	A	801	HAS	C1D-C2D-CMD-OMD
6	A	801	HAS	C22-C23-C24-C28
6	A	801	HAS	C25-C23-C24-C28
6	A	801	HAS	C3D-C2D-CMD-OMD
6	A	801	HAS	C11-C12-C13-C14

There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	563	CMO	1	0
5	A	800	HEM	1	0
6	A	801	HAS	9	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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	552/568 (97%)	0.18	44 (7%) 12 9	33, 43, 49, 58	0
2	B	166/168 (98%)	0.33	19 (11%) 5 3	37, 44, 48, 52	0
3	C	33/34 (97%)	0.19	4 (12%) 4 3	45, 56, 74, 92	0
All	All	751/770 (97%)	0.22	67 (8%) 9 7	33, 43, 51, 92	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	12	TYR	9.4
1	A	14	ALA	9.2
1	A	11	VAL	7.8
1	A	502	ALA	7.0
1	A	13	GLU	7.0
1	A	503	GLU	6.2
1	A	501	LEU	6.0
3	C	2	GLU	5.8
2	B	168	GLU	5.6
1	A	495	ARG	5.4
1	A	416	ALA	4.6
1	A	493	LEU	4.6
1	A	342	ASP	4.6
1	A	504	ALA	4.5
2	B	7	ALA	4.3
1	A	514	GLY	4.1
2	B	60	GLN	4.1
1	A	497	ARG	4.0
1	A	499	PRO	4.0
2	B	39	THR	3.9
1	A	415	ASP	3.9
2	B	6	LYS	3.9
3	C	3	GLU	3.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	498	LYS	3.6
1	A	552	HIS	3.6
1	A	414	SER	3.6
1	A	330	ARG	3.5
1	A	418	ARG	3.5
3	C	34	GLY	3.4
1	A	458	ASP	3.3
2	B	144	GLU	3.3
2	B	76	PRO	3.2
2	B	82	TYR	3.1
1	A	518	ARG	3.1
2	B	38	ALA	3.0
1	A	513	SER	3.0
3	C	4	LYS	3.0
1	A	317	ALA	2.9
1	A	494	SER	2.8
2	B	40	HIS	2.8
1	A	319	SER	2.7
2	B	13	ALA	2.7
2	B	20	ALA	2.7
1	A	327	ARG	2.7
1	A	419	ARG	2.7
1	A	368	SER	2.5
1	A	215	LEU	2.5
2	B	107	ILE	2.5
1	A	455	GLN	2.5
1	A	402	TYR	2.4
2	B	83	VAL	2.4
1	A	63	GLN	2.4
1	A	526	ARG	2.4
2	B	159	ASN	2.4
1	A	105	LEU	2.4
1	A	522	LEU	2.4
1	A	293	THR	2.3
1	A	496	GLU	2.2
1	A	452	TYR	2.2
2	B	102	GLU	2.2
2	B	79	TYR	2.1
2	B	61	GLU	2.1
1	A	54	LEU	2.1
1	A	332	LEU	2.1
1	A	315	THR	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	B	59	ARG	2.1
1	A	517	ASP	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 carbohydrates in this entry.

6.4 Ligands [i](#)

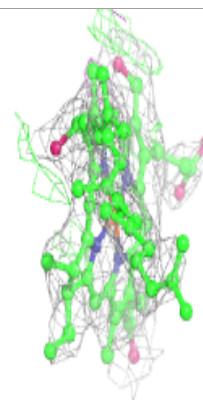
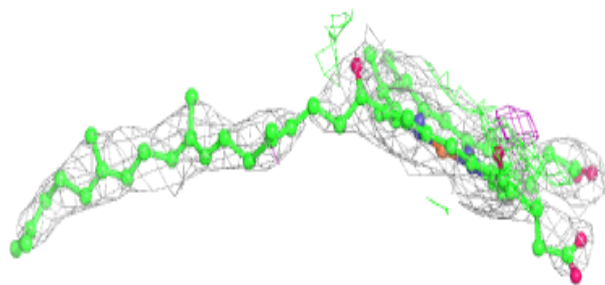
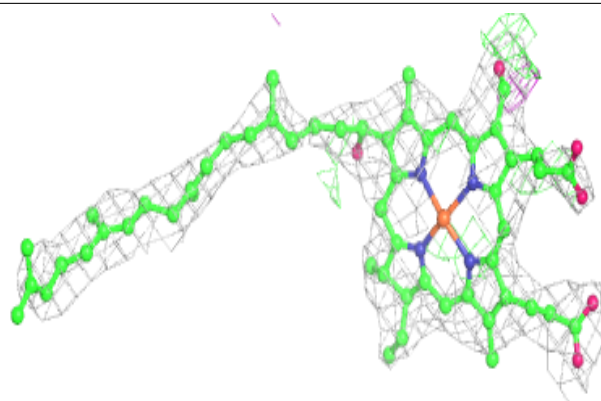
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	HAS	A	801	65/65	0.95	0.16	26,43,46,47	0
5	HEM	A	800	43/43	0.97	0.18	39,44,46,48	0
8	CUA	B	802	2/2	0.99	0.10	44,44,44,47	0
7	CMO	A	563	2/2	0.99	0.23	81,81,81,85	0
4	CU1	A	803	1/1	0.99	0.04	46,46,46,46	0

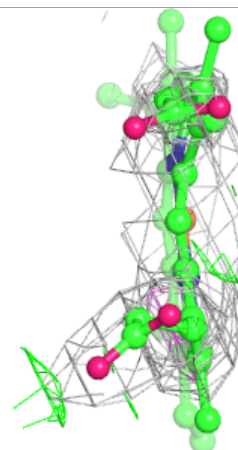
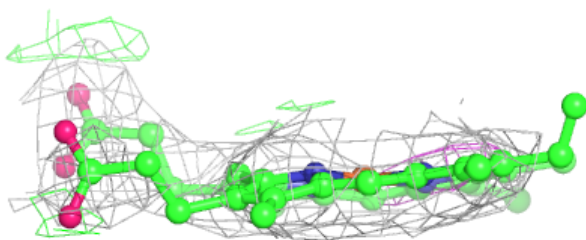
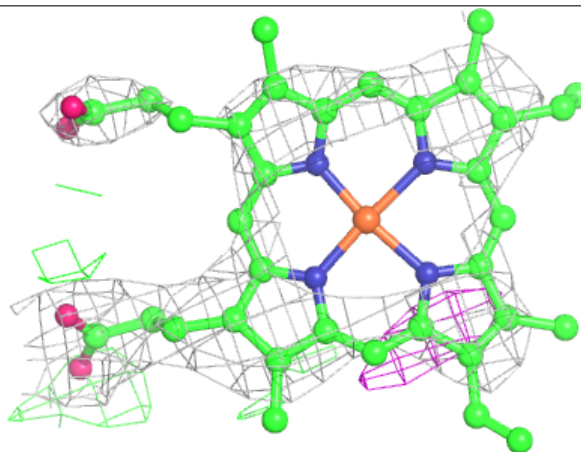
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 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.