



# Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 10:22 am BST

PDB ID : 6M7L  
Title : Complex of OxyA with the X-domain from GPA biosynthesis  
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Deposited on : 2018-08-20  
Resolution : 2.65 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](mailto: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.

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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

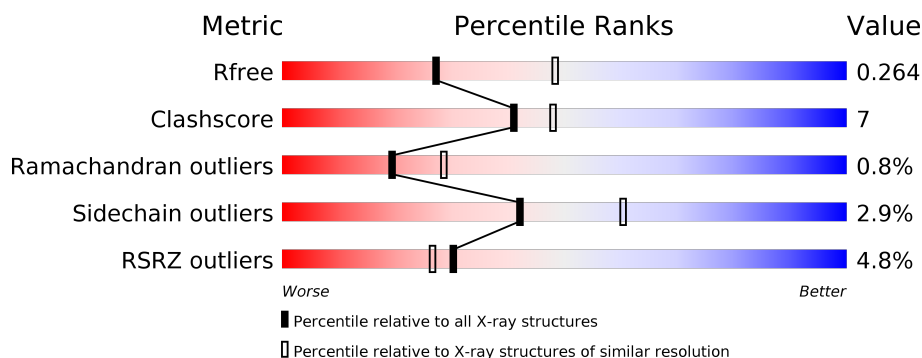
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.65 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	1426 (2.66-2.62)
Clashscore	141614	1472 (2.66-2.62)
Ramachandran outliers	138981	1446 (2.66-2.62)
Sidechain outliers	138945	1446 (2.66-2.62)
RSRZ outliers	127900	1408 (2.66-2.62)

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  $\geq 3$ , 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	B	486	<div> <div>2%</div> <div> <div></div> <div>70%</div> <div>18%</div> <div>•</div> <div>11%</div> </div> </div>
2	A	404	<div> <div>7%</div> <div> <div></div> <div>71%</div> <div>17%</div> <div>11%</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6233 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 Putative non-ribosomal peptide synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	433	Total	C	N	O	S	0	0	0
			3382	2120	627	628	7			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1020	GLY	-	expression tag	UNP A0A2P9IBG7
B	1021	ALA	-	expression tag	UNP A0A2P9IBG7
B	1022	MET	-	expression tag	UNP A0A2P9IBG7
B	1023	ALA	-	expression tag	UNP A0A2P9IBG7
B	1494	LEU	-	expression tag	UNP A0A2P9IBG7
B	1495	GLU	-	expression tag	UNP A0A2P9IBG7
B	1496	SER	-	expression tag	UNP A0A2P9IBG7
B	1497	ALA	-	expression tag	UNP A0A2P9IBG7
B	1498	TRP	-	expression tag	UNP A0A2P9IBG7
B	1499	SER	-	expression tag	UNP A0A2P9IBG7
B	1500	HIS	-	expression tag	UNP A0A2P9IBG7
B	1501	PRO	-	expression tag	UNP A0A2P9IBG7
B	1502	GLN	-	expression tag	UNP A0A2P9IBG7
B	1503	PHE	-	expression tag	UNP A0A2P9IBG7
B	1504	GLU	-	expression tag	UNP A0A2P9IBG7
B	1505	LYS	-	expression tag	UNP A0A2P9IBG7

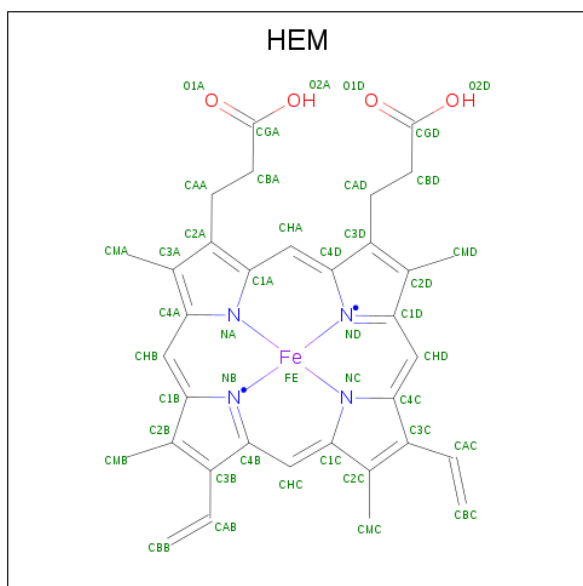
- Molecule 2 is a protein called Putative cytochrome P450 hydroxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	359	Total	C	N	O	S	0	0	0
			2808	1766	520	507	15			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP A0A2P9IBF7
A	-18	GLY	-	expression tag	UNP A0A2P9IBF7
A	-17	SER	-	expression tag	UNP A0A2P9IBF7
A	-16	SER	-	expression tag	UNP A0A2P9IBF7
A	-15	HIS	-	expression tag	UNP A0A2P9IBF7
A	-14	HIS	-	expression tag	UNP A0A2P9IBF7
A	-13	HIS	-	expression tag	UNP A0A2P9IBF7
A	-12	HIS	-	expression tag	UNP A0A2P9IBF7
A	-11	HIS	-	expression tag	UNP A0A2P9IBF7
A	-10	HIS	-	expression tag	UNP A0A2P9IBF7
A	-9	SER	-	expression tag	UNP A0A2P9IBF7
A	-8	SER	-	expression tag	UNP A0A2P9IBF7
A	-7	GLY	-	expression tag	UNP A0A2P9IBF7
A	-6	LEU	-	expression tag	UNP A0A2P9IBF7
A	-5	VAL	-	expression tag	UNP A0A2P9IBF7
A	-4	PRO	-	expression tag	UNP A0A2P9IBF7
A	-3	ARG	-	expression tag	UNP A0A2P9IBF7
A	-2	GLY	-	expression tag	UNP A0A2P9IBF7
A	-1	SER	-	expression tag	UNP A0A2P9IBF7
A	0	HIS	-	expression tag	UNP A0A2P9IBF7

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



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



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	118.83Å 87.34Å 95.95Å 90.00° 95.77° 90.00°	Depositor
Resolution (Å)	47.73 – 2.65 47.73 – 2.65	Depositor EDS
% Data completeness (in resolution range)	99.6 (47.73-2.65) 99.6 (47.73-2.65)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.10 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, $R_{free}$	0.214 , 0.264 0.214 , 0.264	Depositor DCC
$R_{free}$ test set	1415 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	65.0	Xtriage
Anisotropy	0.248	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 48.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6233	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.17% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEM

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	B	0.24	0/3452	0.43	0/4699
2	A	0.24	0/2860	0.41	0/3867
All	All	0.24	0/6312	0.42	0/8566

There are no bond length outliers.

There are no bond angle outliers.

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	B	3382	0	3369	49	0
2	A	2808	0	2835	39	0
3	A	43	0	30	4	0
All	All	6233	0	6234	89	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 7.

All (89) 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)
2:A:190:ARG:NH2	2:A:211:ASP:OD2	2.10	0.85
1:B:1070:ARG:NH2	1:B:1126:GLU:OE2	2.13	0.82
2:A:120:ASP:OD1	2:A:352:ARG:NH2	2.15	0.77
2:A:291:LYS:N	2:A:294:GLU:OE2	2.15	0.76
2:A:90:ARG:NH1	2:A:208:ASP:O	2.20	0.74
3:A:401:HEM:HBC2	3:A:401:HEM:HHD	1.69	0.74
1:B:1178:ASP:OD2	1:B:1182:ARG:NH1	2.23	0.71
1:B:1343:ASP:O	1:B:1356:ARG:NH1	2.27	0.66
1:B:1042:PRO:O	1:B:1107:GLN:NE2	2.30	0.64
1:B:1259:VAL:HG13	1:B:1260:LEU:HD12	1.82	0.61
1:B:1046:ARG:NH2	1:B:1373:GLU:OE2	2.34	0.60
2:A:12:ARG:HD2	2:A:272:GLU:OE2	2.02	0.60
2:A:325:PHE:HB3	2:A:332:CYS:HB3	1.83	0.60
2:A:284:GLU:HG2	2:A:289:VAL:HG12	1.84	0.59
1:B:1041:LEU:HD12	1:B:1042:PRO:HD2	1.84	0.59
1:B:1287:ALA:HB3	1:B:1354:LEU:HD22	1.85	0.58
2:A:334:GLY:HA3	3:A:401:HEM:C2B	2.38	0.58
2:A:48:TYR:CZ	2:A:52:ARG:HD2	2.39	0.57
2:A:190:ARG:NH1	2:A:209:ILE:O	2.32	0.56
1:B:1192:ARG:NH2	2:A:117:ASP:OD1	2.39	0.55
3:A:401:HEM:HMB2	3:A:401:HEM:HBB2	1.89	0.55
1:B:1217:LEU:HD11	1:B:1371:PRO:HD3	1.89	0.55
2:A:30:LEU:HB2	2:A:50:VAL:HG11	1.90	0.53
2:A:114:ILE:HD13	2:A:142:ARG:HG3	1.91	0.53
2:A:96:ALA:HA	2:A:333:LEU:HD11	1.90	0.53
1:B:1139:ARG:HB2	1:B:1150:ARG:NH2	2.24	0.52
1:B:1271:ARG:HG2	1:B:1447:VAL:HG22	1.92	0.52
2:A:31:VAL:HG22	2:A:44:PHE:HB2	1.91	0.52
1:B:1064:HIS:O	1:B:1169:ARG:NH2	2.44	0.51
1:B:1309:ARG:HG2	1:B:1311:SER:H	1.76	0.50
2:A:98:ALA:HA	2:A:200:MET:HE1	1.92	0.50
1:B:1347:ASP:OD1	1:B:1483:ARG:HB3	2.12	0.50
1:B:1347:ASP:OD2	1:B:1485:SER:HB3	2.11	0.50
2:A:80:LEU:HD21	2:A:91:LEU:HB3	1.94	0.49
2:A:108:GLU:HG2	2:A:340:MET:SD	2.52	0.49
1:B:1086:ASP:OD2	1:B:1191:ARG:NE	2.43	0.49
2:A:92:ARG:HD2	2:A:333:LEU:HD13	1.94	0.48
1:B:1261:PRO:HG3	1:B:1385:ILE:O	2.13	0.48
1:B:1459:ARG:NH2	1:B:1463:GLU:OE2	2.40	0.48
1:B:1217:LEU:HG	1:B:1227:ILE:HD12	1.96	0.48
2:A:17:LEU:HG	2:A:21:LEU:HD23	1.96	0.48
1:B:1111:ASP:OD1	1:B:1114:THR:OG1	2.28	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:78:GLY:HA2	2:A:91:LEU:HD21	1.96	0.47
1:B:1405:GLU:HB3	1:B:1417:PRO:HG3	1.95	0.47
1:B:1096:THR:O	1:B:1148:PRO:HD3	2.15	0.47
1:B:1348:PRO:HB2	1:B:1353:ILE:HG13	1.97	0.47
2:A:10:HIS:CE1	2:A:372:SER:HA	2.50	0.46
1:B:1096:THR:HG21	1:B:1107:GLN:HG2	1.95	0.46
1:B:1215:GLU:O	1:B:1217:LEU:N	2.48	0.46
2:A:53:GLU:OE2	2:A:285:ILE:HG23	2.15	0.46
2:A:274:LEU:HD21	2:A:300:LEU:HD11	1.98	0.46
1:B:1389:PRO:O	1:B:1391:PHE:N	2.44	0.46
1:B:1483:ARG:HB2	1:B:1486:GLN:HG2	1.97	0.46
2:A:218:CYS:O	2:A:222:MET:HG2	2.15	0.46
1:B:1101:SER:OG	1:B:1102:ARG:N	2.49	0.46
2:A:262:GLU:OE2	2:A:339:ARG:NE	2.39	0.45
1:B:1136:ARG:O	1:B:1150:ARG:NH2	2.35	0.45
1:B:1072:ARG:NE	1:B:1160:GLU:OE1	2.42	0.45
1:B:1270:LEU:HD22	1:B:1433:LEU:HD21	1.99	0.45
2:A:206:GLY:C	2:A:208:ASP:H	2.20	0.45
2:A:276:PRO:HB2	2:A:295:ALA:HB1	1.99	0.44
1:B:1056:ILE:HD12	1:B:1263:ARG:NH2	2.32	0.44
1:B:1176:SER:HB3	1:B:1331:VAL:HG13	1.99	0.44
1:B:1433:LEU:HD23	1:B:1448:GLY:HA3	1.99	0.44
1:B:1292:PHE:HA	1:B:1397:LEU:HD13	1.98	0.44
2:A:307:ASP:OD2	2:A:321:ARG:NH2	2.50	0.43
1:B:1117:PRO:HG3	1:B:1149:TRP:CZ2	2.53	0.43
1:B:1430:TRP:CD2	1:B:1432:LYS:HE3	2.54	0.43
1:B:1277:HIS:CG	1:B:1443:ALA:HB1	2.54	0.43
2:A:219:ALA:O	2:A:223:ASN:HB2	2.19	0.43
2:A:278:THR:HG22	2:A:295:ALA:HB2	2.01	0.42
1:B:1062:GLY:HA3	1:B:1064:HIS:CE1	2.54	0.42
2:A:277:ARG:HH22	3:A:401:HEM:CGD	2.32	0.42
1:B:1048:ARG:NH2	1:B:1102:ARG:O	2.21	0.42
1:B:1237:ARG:HD3	1:B:1237:ARG:HA	1.87	0.42
1:B:1093:ILE:HD12	1:B:1206:PHE:CD2	2.54	0.42
1:B:1073:GLY:N	1:B:1159:ARG:O	2.32	0.42
2:A:48:TYR:CE1	2:A:52:ARG:HD2	2.55	0.42
1:B:1223:ARG:HB2	1:B:1223:ARG:HE	1.62	0.42
2:A:332:CYS:SG	2:A:335:ALA:N	2.92	0.41
2:A:92:ARG:NH1	2:A:331:HIS:O	2.54	0.41
2:A:22:LYS:O	2:A:26:GLU:HG3	2.20	0.41
1:B:1129:LEU:HD23	1:B:1129:LEU:HA	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1297:ALA:O	1:B:1301:MET:HG3	2.21	0.41
2:A:43:TRP:HB2	2:A:296:VAL:HA	2.03	0.41
1:B:1384:SER:OG	1:B:1386:THR:HG22	2.21	0.40
2:A:368:ARG:HG2	2:A:369:PRO:O	2.21	0.40
2:A:19:ALA:O	2:A:23:LEU:HG	2.22	0.40
1:B:1115:GLY:O	1:B:1147:VAL:HG11	2.22	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	B	429/486 (88%)	407 (95%)	18 (4%)	4 (1%)	17	26
2	A	351/404 (87%)	333 (95%)	16 (5%)	2 (1%)	25	37
All	All	780/890 (88%)	740 (95%)	34 (4%)	6 (1%)	19	28

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1217	LEU
2	A	374	VAL
1	B	1324	GLN
1	B	1127	ASP
1	B	1381	PRO
2	A	370	PRO

### 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	B	350/388 (90%)	337 (96%)	13 (4%)	34	51
2	A	296/333 (89%)	290 (98%)	6 (2%)	55	72
All	All	646/721 (90%)	627 (97%)	19 (3%)	42	60

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	1066	ARG
1	B	1070	ARG
1	B	1107	GLN
1	B	1116	ARG
1	B	1180	LEU
1	B	1192	ARG
1	B	1217	LEU
1	B	1223	ARG
1	B	1227	ILE
1	B	1296	HIS
1	B	1401	ASP
1	B	1465	LEU
1	B	1487	VAL
2	A	92	ARG
2	A	182	MET
2	A	225	SER
2	A	228	GLN
2	A	271	PHE
2	A	374	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

1 ligand is 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	A	401	-	27,50,50	1.88	4 (14%)	17,82,82	1.47	1 (5%)

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	A	401	-	-	0/6/54/54	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	HEM	C3C-C2C	-4.69	1.33	1.40
3	A	401	HEM	C3B-CAB	3.74	1.55	1.47
3	A	401	HEM	C3B-C2B	-3.68	1.35	1.40
3	A	401	HEM	C3C-CAC	3.57	1.55	1.47

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	HEM	CMB-C2B-C3B	2.20	128.79	124.68

There are no chirality outliers.

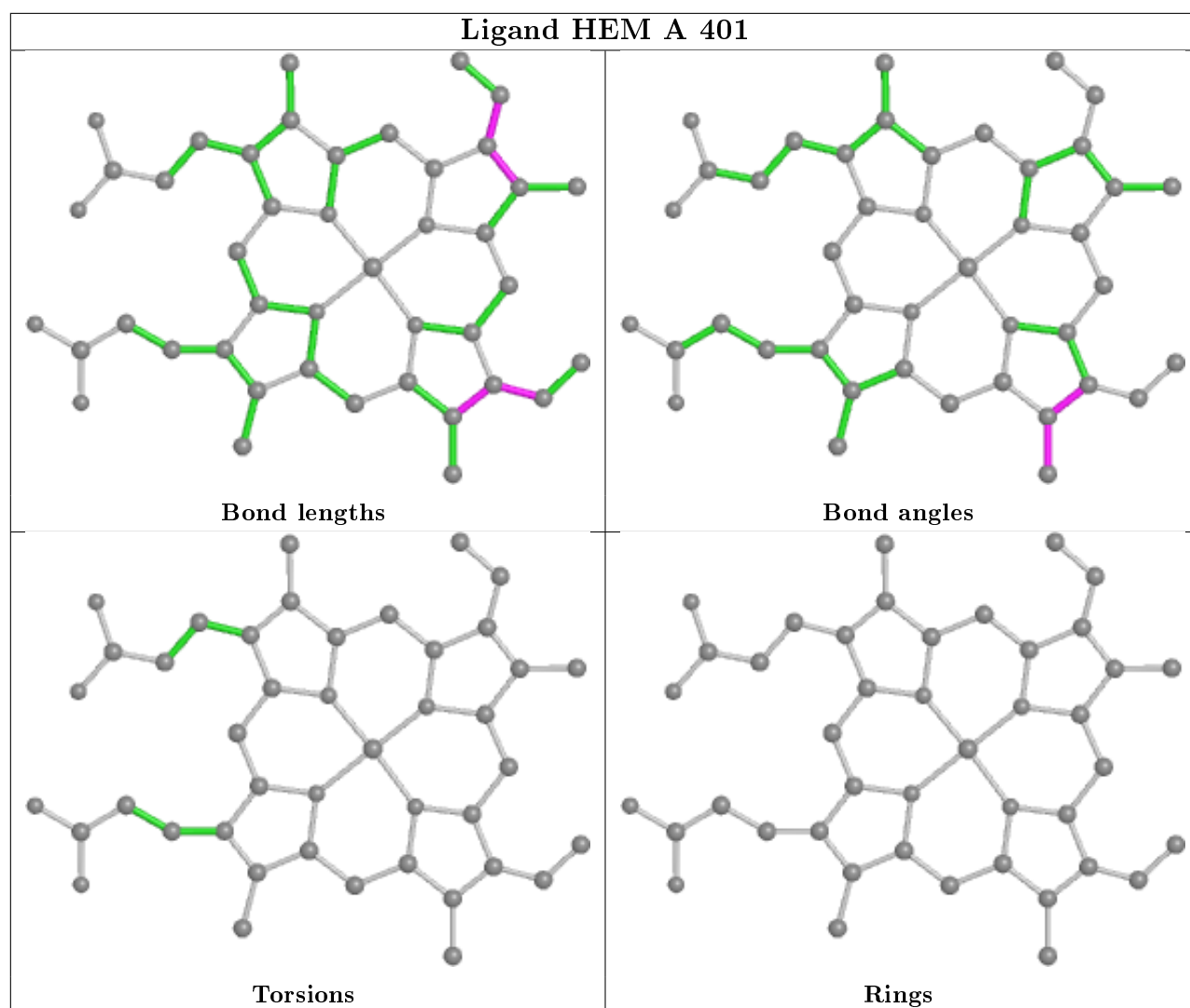
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	HEM	4	0

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



## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	B	433/486 (89%)	0.20	10 (2%) 60 57	47, 67, 96, 118	0
2	A	359/404 (88%)	0.44	28 (7%) 13 10	46, 72, 119, 144	0
All	All	792/890 (88%)	0.31	38 (4%) 30 26	46, 69, 109, 144	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	289	VAL	8.7
2	A	285	ILE	6.2
2	A	280	THR	5.2
2	A	23	LEU	4.3
1	B	1059	GLU	4.3
2	A	284	GLU	4.2
1	B	1489	VAL	4.1
2	A	287	GLY	4.1
2	A	191	ARG	3.7
2	A	290	ILE	3.7
2	A	295	ALA	3.7
2	A	33	ILE	3.6
2	A	24	LEU	3.5
2	A	40	ALA	3.5
2	A	34	PRO	3.5
2	A	43	TRP	3.3
1	B	1220	ALA	3.2
1	B	1204	LEU	2.9
2	A	206	GLY	2.8
2	A	27	ARG	2.7
2	A	42	HIS	2.7
2	A	207	ALA	2.5
2	A	31	VAL	2.5
2	A	283	VAL	2.5

*Continued on next page...*

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Mol	Chain	Res	Type	RSRZ
2	A	7	ARG	2.5
1	B	1487	VAL	2.4
2	A	288	GLN	2.4
2	A	30	LEU	2.3
2	A	297	PHE	2.3
1	B	1478	ALA	2.3
1	B	1488	ASP	2.2
2	A	41	VAL	2.2
2	A	39	SER	2.2
2	A	35	LEU	2.2
1	B	1346	GLY	2.1
1	B	1483	ARG	2.1
1	B	1486	GLN	2.0
2	A	32	ARG	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, 95<sup>th</sup> 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.

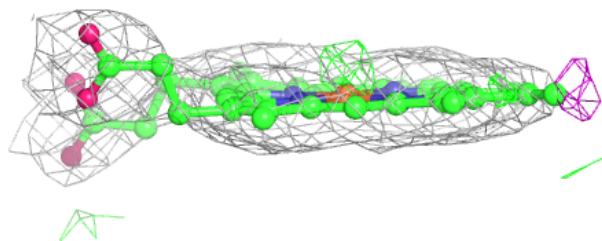
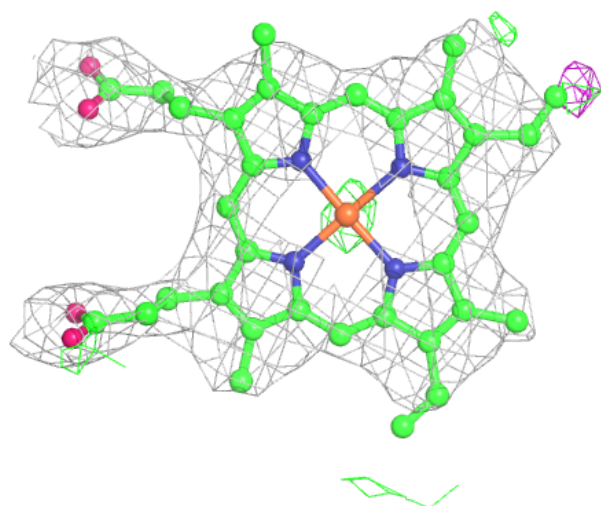
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	HEM	A	401	43/43	0.98	0.21	49,59,74,76	0

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 A 401:**

$2mF_o-DF_c$  (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.