



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 04:13 pm BST

PDB ID : 6UNE  
Title : Human CYP3A4 bound to an inhibitor  
Authors : Sevrioukova, I.F.  
Deposited on : 2019-10-11  
Resolution : 2.55 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

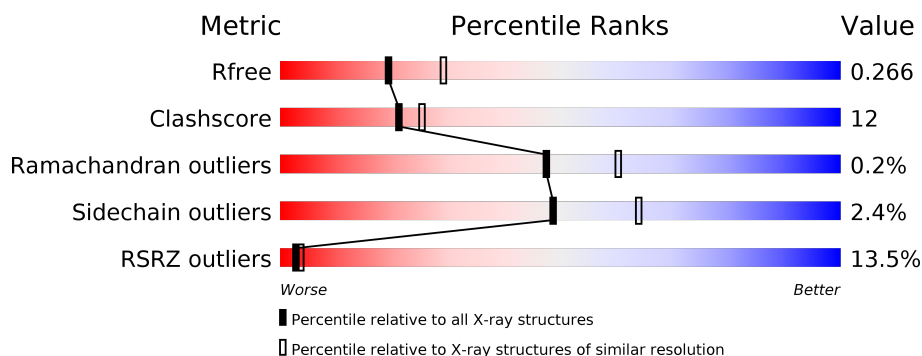
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.55 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	1284 (2.56-2.52)
Clashscore	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)
RSRZ outliers	127900	1272 (2.56-2.52)

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	A	487	

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3750 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 3A4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	452	3667	2392	596	653	26	0	7	0

There are 6 discrepancies between the modelled and reference sequences:

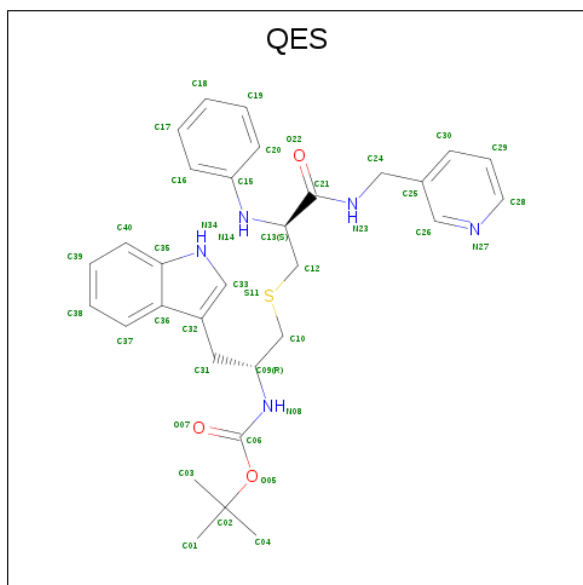
Chain	Residue	Modelled	Actual	Comment	Reference
A	21	MET	-	initiating methionine	UNP P08684
A	22	ALA	-	expression tag	UNP P08684
A	504	HIS	-	expression tag	UNP P08684
A	505	HIS	-	expression tag	UNP P08684
A	506	HIS	-	expression tag	UNP P08684
A	507	HIS	-	expression tag	UNP P08684

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).



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

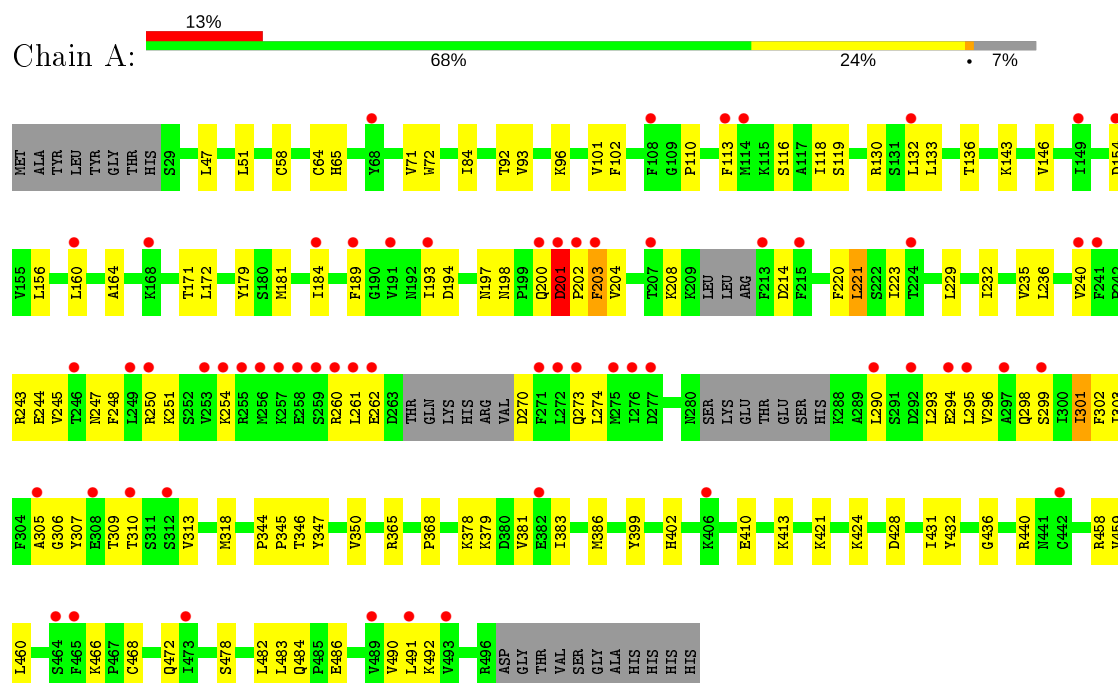
- Molecule 3 is tert-butyl [(2R)-1-(1H-indol-3-yl)-3-{[(2S)-3-oxo-2-(phenylamino)-3-{[(pyridin-3-yl)methyl]amino}propyl]sulfanyl}propan-2-yl]carbamate (three-letter code: QES) (formula: C<sub>31</sub>H<sub>37</sub>N<sub>5</sub>O<sub>3</sub>S) (labeled as "Ligand of Interest" by author).



### 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 P450 3A4



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.09Å 101.51Å 126.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.55 – 2.55 39.55 – 2.55	Depositor EDS
% Data completeness (in resolution range)	96.9 (39.55-2.55) 96.9 (39.55-2.55)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.06 (at 2.54Å)	Xtriage
Refinement program	PHENIX 1.11.1 _2575	Depositor
R, $R_{free}$	0.211 , 0.264 0.216 , 0.266	Depositor DCC
$R_{free}$ test set	771 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	90.2	Xtriage
Anisotropy	0.099	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 87.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3750	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	128.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.91% 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 [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.32	1/3775 (0.0%)	0.48	0/5102

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	201	ASP	C-N	5.59	1.44	1.34

There are no bond angle outliers.

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	3667	0	3771	89	0
2	A	43	0	30	6	0
3	A	40	0	0	0	0
All	All	3750	0	3801	91	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 12.

All (91) 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:198:ASN:HB3	1:A:201:ASP:OD2	1.76	0.86
1:A:202:PRO:HG2	1:A:203:PHE:CD2	2.14	0.83
1:A:201:ASP:HB3	1:A:202:PRO:HD2	1.66	0.77
1:A:260:ARG:HD3	1:A:273:GLN:OE1	1.89	0.73
1:A:201:ASP:HB3	1:A:202:PRO:CD	2.19	0.72
1:A:101:VAL:HG21	1:A:381[A]:VAL:HG11	1.73	0.71
2:A:601:HEM:HBC2	2:A:601:HEM:HHD	1.72	0.69
1:A:156:LEU:HD13	1:A:179:TYR:HB2	1.77	0.67
1:A:202:PRO:HG2	1:A:203:PHE:CE2	2.32	0.65
1:A:119:SER:HB3	2:A:601:HEM:HAD1	1.78	0.64
1:A:421:LYS:HA	1:A:424:LYS:HG2	1.80	0.64
1:A:208:LYS:HG2	1:A:208:LYS:O	1.99	0.62
1:A:200:GLN:HG2	1:A:200:GLN:O	2.00	0.61
1:A:132:LEU:HD13	1:A:290:LEU:HG	1.84	0.60
1:A:172:LEU:HD11	1:A:491:LEU:HD12	1.86	0.58
1:A:171:THR:HG22	1:A:490:VAL:HG12	1.87	0.56
1:A:214:ASP:HA	1:A:482:LEU:HD21	1.89	0.54
1:A:92:THR:HA	1:A:96:LYS:HB2	1.89	0.53
2:A:601:HEM:HMB2	2:A:601:HEM:HBB2	1.88	0.53
1:A:113:PHE:HB2	1:A:293:LEU:HD21	1.91	0.52
1:A:101:VAL:O	1:A:378:LYS:HG2	2.10	0.51
1:A:136:THR:HG23	1:A:274:LEU:HD12	1.92	0.51
1:A:260:ARG:NH2	1:A:270:ASP:OD2	2.43	0.51
1:A:189:PHE:HE2	1:A:303:ILE:HD11	1.76	0.51
1:A:146:VAL:HG21	1:A:347:TYR:HB2	1.92	0.51
1:A:101:VAL:HG11	1:A:381[B]:VAL:HG21	1.93	0.50
1:A:189:PHE:CZ	1:A:299:SER:HB3	2.47	0.50
1:A:482:LEU:HB2	1:A:484:GLN:HG2	1.93	0.50
1:A:189:PHE:HZ	1:A:299:SER:HB3	1.76	0.50
1:A:71[A]:VAL:HG21	1:A:386:MET:HE1	1.94	0.50
1:A:383:ILE:O	1:A:383:ILE:HG13	2.12	0.49
1:A:410:GLU:HB3	1:A:413:LYS:HD3	1.93	0.49
1:A:154:ASP:CG	1:A:458:ARG:HH12	2.17	0.48
1:A:483:LEU:HA	1:A:483:LEU:HD12	1.61	0.48
1:A:466:LYS:HB2	1:A:492:LYS:HB3	1.95	0.48
1:A:184:ILE:HD11	1:A:306:GLY:C	2.34	0.48
1:A:116:SER:O	1:A:298:GLN:NE2	2.40	0.48
1:A:378:LYS:HG3	1:A:379:LYS:HG2	1.93	0.48
1:A:203:PHE:HD1	1:A:248:PHE:CE2	2.32	0.48
1:A:309:THR:OG1	1:A:310:THR:N	2.46	0.48
1:A:478:SER:HB2	1:A:484:GLN:HG3	1.96	0.47
1:A:172:LEU:HD13	1:A:318:MET:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:PHE:O	1:A:260:ARG:NH2	2.48	0.47
1:A:193:ILE:HD13	1:A:201:ASP:OD1	2.15	0.47
1:A:235:VAL:HG13	1:A:236:LEU:HD23	1.96	0.46
1:A:93:VAL:HG13	1:A:102:PHE:CG	2.51	0.46
1:A:428:ASP:HB3	1:A:431:ILE:HG12	1.98	0.46
1:A:270:ASP:O	1:A:274:LEU:HG	2.16	0.46
1:A:160:LEU:O	1:A:164:ALA:N	2.49	0.46
1:A:345:PRO:HD2	1:A:458:ARG:HD3	1.97	0.46
1:A:305:ALA:HB1	2:A:601:HEM:C3C	2.50	0.45
1:A:118:ILE:HD11	1:A:130:ARG:HD2	1.99	0.45
1:A:232:ILE:O	1:A:235:VAL:HG12	2.18	0.44
1:A:101:VAL:HG11	1:A:381[A]:VAL:HG21	2.00	0.44
1:A:143:LYS:HD3	1:A:347:TYR:CD2	2.53	0.44
1:A:201:ASP:CB	1:A:202:PRO:CD	2.93	0.44
1:A:204:VAL:O	1:A:204:VAL:HG12	2.18	0.44
1:A:58:CYS:HB3	1:A:399:TYR:CE2	2.53	0.43
1:A:436:GLY:HA3	2:A:601:HEM:HBA1	2.00	0.43
1:A:133:LEU:HD11	1:A:298:GLN:HG3	2.00	0.43
1:A:251:LYS:HA	1:A:254:LYS:HG2	2.01	0.43
1:A:305:ALA:HB1	2:A:601:HEM:C4C	2.54	0.43
1:A:208:LYS:O	1:A:208:LYS:CG	2.67	0.42
1:A:220:PHE:CD2	1:A:223:ILE:HD12	2.54	0.42
1:A:47:LEU:HD13	1:A:51:LEU:HD21	2.01	0.42
1:A:92:THR:HG21	1:A:383:ILE:HB	2.01	0.42
1:A:250:ARG:HG3	1:A:296:VAL:HG11	2.00	0.42
1:A:346:THR:O	1:A:350:VAL:HG23	2.20	0.42
1:A:110:PRO:HD2	1:A:240:VAL:HG13	2.01	0.42
1:A:184:ILE:HD11	1:A:307:TYR:N	2.34	0.42
1:A:64:CYS:HB3	1:A:72:TRP:CE2	2.54	0.42
1:A:309:THR:O	1:A:313:VAL:HG23	2.20	0.42
1:A:294:GLU:O	1:A:298:GLN:HG2	2.20	0.42
1:A:402:HIS:CD2	1:A:432:TYR:HA	2.55	0.42
1:A:261:LEU:HA	1:A:262:GLU:HA	1.70	0.41
1:A:101:VAL:HG21	1:A:381[B]:VAL:HG21	2.02	0.41
1:A:202:PRO:CG	1:A:203:PHE:CE2	3.03	0.41
1:A:244:GLU:HG3	1:A:245:VAL:H	1.86	0.41
1:A:365:ARG:O	1:A:368:PRO:HD3	2.21	0.41
1:A:243:ARG:O	1:A:247:ASN:ND2	2.54	0.41
1:A:156:LEU:HD12	1:A:156:LEU:HA	1.89	0.41
1:A:194:ASP:OD2	1:A:197:ASN:ND2	2.46	0.41
1:A:221:LEU:HD23	1:A:221:LEU:HA	1.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:460:LEU:HA	1:A:460:LEU:HD23	1.96	0.40
1:A:71[B]:VAL:HG12	1:A:84:ILE:HG22	2.03	0.40
1:A:201:ASP:HB2	1:A:204:VAL:CG2	2.51	0.40
1:A:229:LEU:HA	1:A:232:ILE:HG13	2.03	0.40
1:A:248:PHE:HA	1:A:251:LYS:HE3	2.04	0.40
1:A:344:PRO:HA	1:A:345:PRO:HD2	2.00	0.40
1:A:472:GLN:HE22	1:A:486:GLU:HB3	1.86	0.40
1:A:301:ILE:HG22	1:A:302:PHE:CD1	2.57	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	451/487 (93%)	425 (94%)	25 (6%)	1 (0%)	47 60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	201	ASP

### 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	420/443 (95%)	410 (98%)	10 (2%)	49 64

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	HIS
1	A	181	MET
1	A	201	ASP
1	A	203	PHE
1	A	221	LEU
1	A	295	LEU
1	A	301	ILE
1	A	440	ARG
1	A	459	VAL
1	A	468	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	192	ASN
1	A	461	GLN

### 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 ⓘ

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	QES	A	602	2	41,43,43	1.77	8 (19%)	50,58,58	2.30	19 (38%)
2	HEM	A	601	1,3	27,50,50	1.91	4 (14%)	17,82,82	1.52	5 (29%)

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	QES	A	602	2	-	10/31/32/32	0/4/4/4
2	HEM	A	601	1,3	-	0/6/54/54	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	HEM	C3C-C2C	-4.99	1.33	1.40
3	A	602	QES	C33-N34	3.86	1.44	1.36
2	A	601	HEM	C3B-CAB	3.81	1.55	1.47
3	A	602	QES	C15-N14	3.68	1.46	1.39
2	A	601	HEM	C3B-C2B	-3.60	1.35	1.40
2	A	601	HEM	C3C-CAC	3.57	1.55	1.47
3	A	602	QES	C06-N08	3.38	1.43	1.34
3	A	602	QES	O05-C06	3.31	1.41	1.34
3	A	602	QES	C13-N14	3.25	1.49	1.45
3	A	602	QES	O07-C06	2.85	1.26	1.21
3	A	602	QES	C21-N23	2.39	1.38	1.33
3	A	602	QES	C10-S11	2.02	1.89	1.80

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	QES	O05-C06-N08	7.78	123.09	110.02
3	A	602	QES	C12-S11-C10	5.82	112.02	102.13
3	A	602	QES	O07-C06-N08	-4.70	117.14	124.85
3	A	602	QES	O05-C06-O07	-4.06	118.23	125.62
3	A	602	QES	C25-C24-N23	3.84	121.28	113.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	QES	C09-N08-C06	3.46	127.63	122.30
3	A	602	QES	C31-C09-N08	3.23	116.60	110.39
3	A	602	QES	C13-C21-N23	2.81	122.20	116.54
3	A	602	QES	C24-N23-C21	2.80	126.37	122.34
3	A	602	QES	C31-C32-C33	-2.72	124.61	127.97
2	A	601	HEM	CBD-CAD-C3D	-2.50	107.87	112.48
3	A	602	QES	C17-C16-C15	2.47	122.68	119.72
2	A	601	HEM	CMD-C2D-C1D	-2.45	124.70	128.46
3	A	602	QES	C31-C32-C36	2.41	129.99	126.25
3	A	602	QES	C20-C15-C16	-2.30	115.89	119.03
3	A	602	QES	C28-N27-C26	2.26	120.75	116.85
3	A	602	QES	O05-C02-C01	2.25	116.11	107.20
2	A	601	HEM	CBA-CAA-C2A	-2.23	108.37	112.49
3	A	602	QES	O22-C21-N23	-2.23	118.21	122.99
2	A	601	HEM	C4A-C3A-C2A	2.12	108.47	107.00
3	A	602	QES	C24-C25-C30	2.12	125.31	120.91
3	A	602	QES	C03-C02-C01	-2.12	105.61	111.16
3	A	602	QES	C04-C02-C01	-2.10	105.66	111.16
2	A	601	HEM	CMB-C2B-C3B	2.03	128.47	124.68

There are no chirality outliers.

All (10) torsion outliers are listed below:

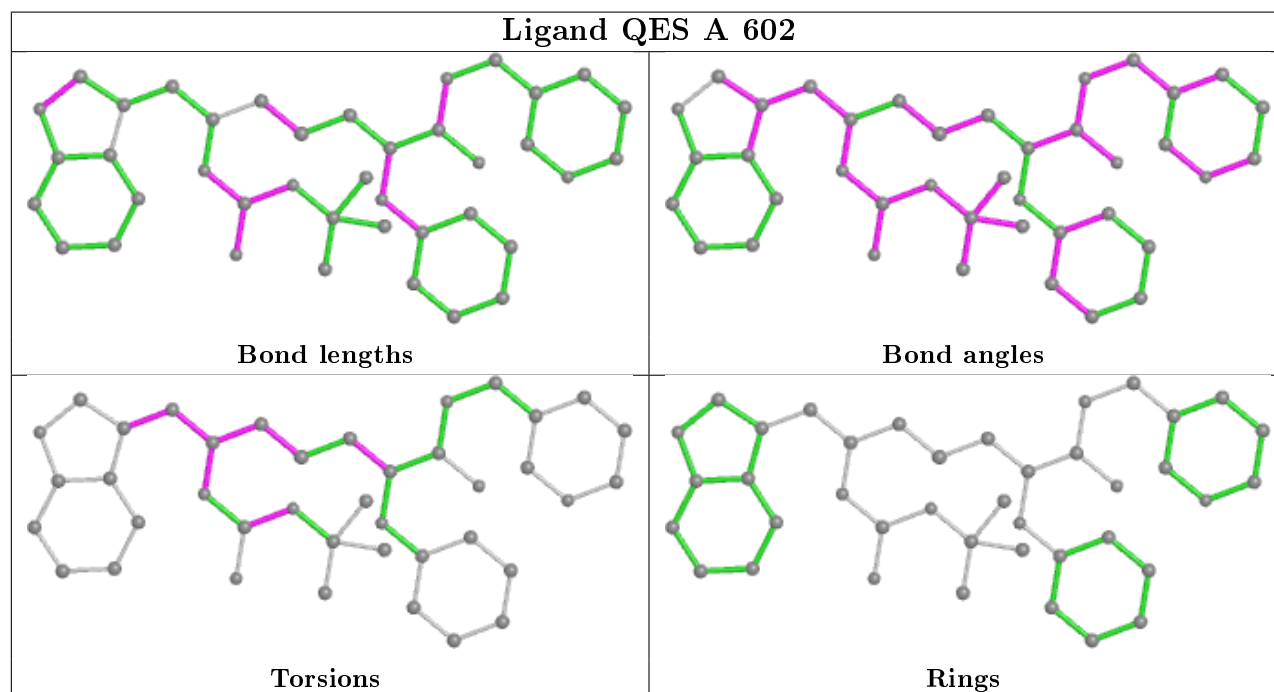
Mol	Chain	Res	Type	Atoms
3	A	602	QES	O07-C06-O05-C02
3	A	602	QES	C31-C09-C10-S11
3	A	602	QES	N08-C09-C10-S11
3	A	602	QES	C10-C09-C31-C32
3	A	602	QES	C31-C09-N08-C06
3	A	602	QES	C09-C10-S11-C12
3	A	602	QES	C09-C31-C32-C33
3	A	602	QES	N08-C06-O05-C02
3	A	602	QES	N08-C09-C31-C32
3	A	602	QES	S11-C12-C13-N14

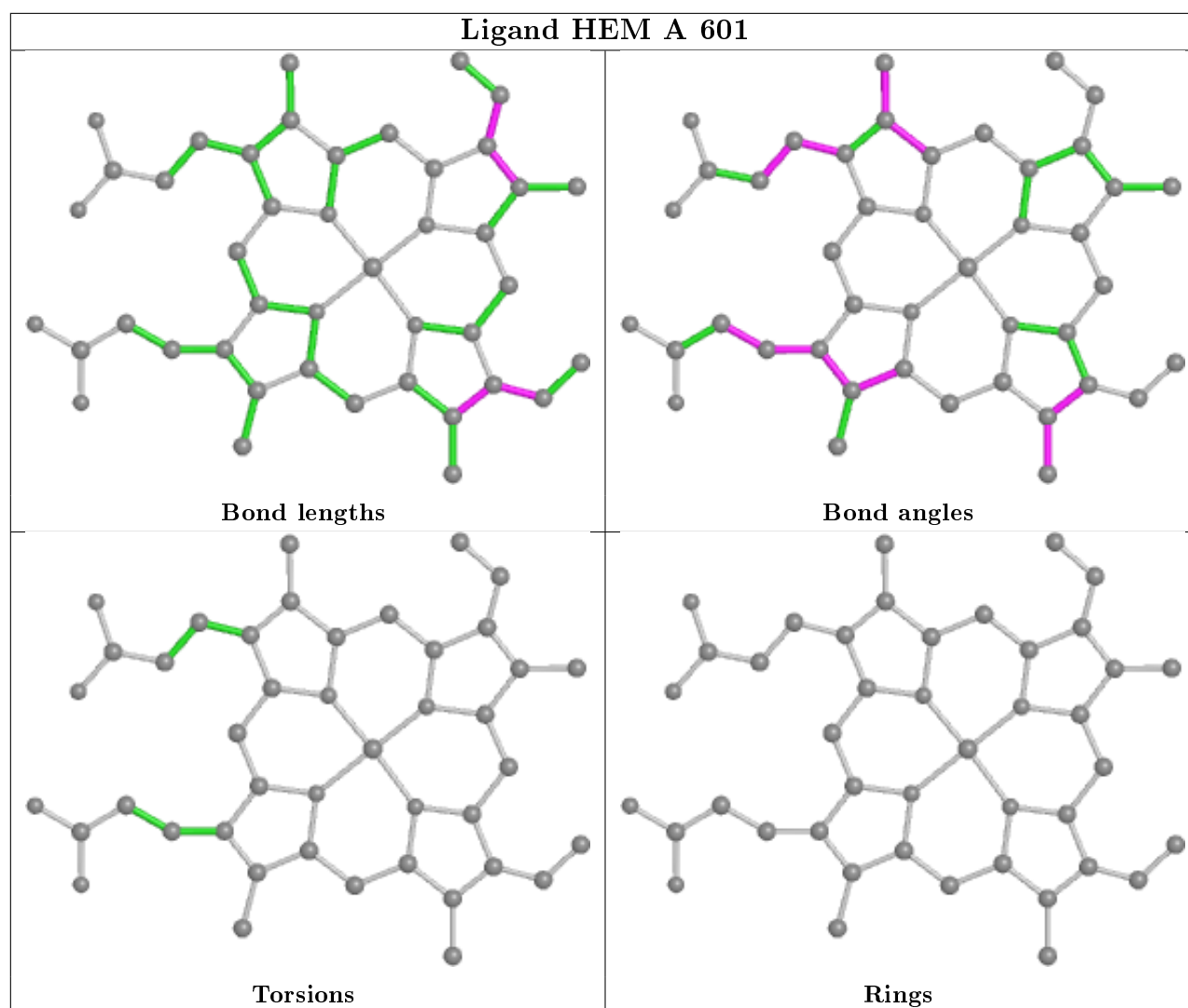
There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	HEM	6	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	A	452/487 (92%)	0.83	61 (13%) 3 4	71, 123, 196, 291	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	189	PHE	9.2
1	A	258	GLU	7.5
1	A	213	PHE	7.1
1	A	241	PHE	6.7
1	A	257	LYS	6.6
1	A	262	GLU	6.5
1	A	290	LEU	6.3
1	A	261	LEU	6.3
1	A	256	MET	5.9
1	A	272	LEU	5.5
1	A	254	LYS	5.4
1	A	203	PHE	5.2
1	A	191	VAL	4.8
1	A	259	SER	4.6
1	A	202	PRO	4.5
1	A	260	ARG	4.3
1	A	295	LEU	4.3
1	A	240	VAL	4.1
1	A	473	ILE	3.9
1	A	271	PHE	3.8
1	A	292	ASP	3.7
1	A	294	GLU	3.6
1	A	246	THR	3.6
1	A	312	SER	3.5
1	A	201	ASP	3.5
1	A	215	PHE	3.3
1	A	255	ARG	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	154	ASP	3.3
1	A	489	VAL	3.2
1	A	114	MET	3.2
1	A	275	MET	3.1
1	A	207	THR	3.1
1	A	491	LEU	3.1
1	A	168	LYS	3.1
1	A	382	GLU	2.9
1	A	305	ALA	2.8
1	A	493	VAL	2.8
1	A	464	SER	2.7
1	A	200	GLN	2.6
1	A	193	ILE	2.6
1	A	160	LEU	2.6
1	A	406	LYS	2.5
1	A	253	VAL	2.5
1	A	149	ILE	2.5
1	A	249	LEU	2.4
1	A	224	THR	2.4
1	A	250	ARG	2.4
1	A	442	CYS	2.4
1	A	68	TYR	2.4
1	A	276	ILE	2.4
1	A	297	ALA	2.3
1	A	299	SER	2.2
1	A	308	GLU	2.2
1	A	108	PHE	2.2
1	A	113	PHE	2.2
1	A	277	ASP	2.2
1	A	132	LEU	2.2
1	A	310	THR	2.1
1	A	465	PHE	2.0
1	A	273	GLN	2.0
1	A	184	ILE	2.0

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

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

### 6.3 Carbohydrates

There are no carbohydrates in this entry.

### 6.4 Ligands

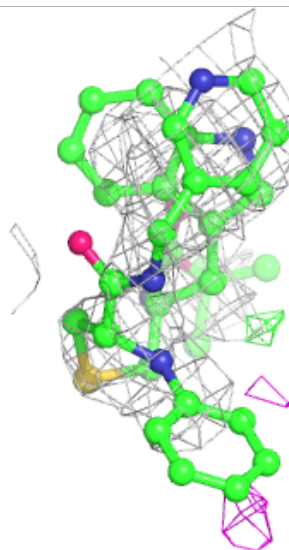
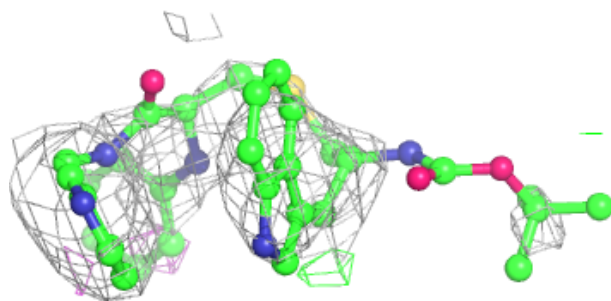
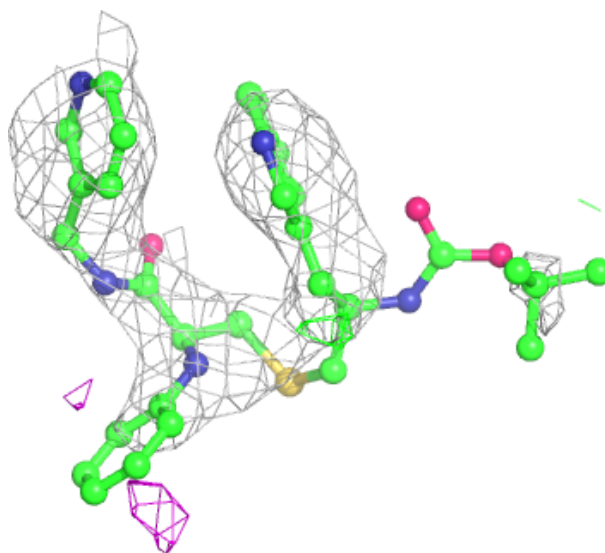
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

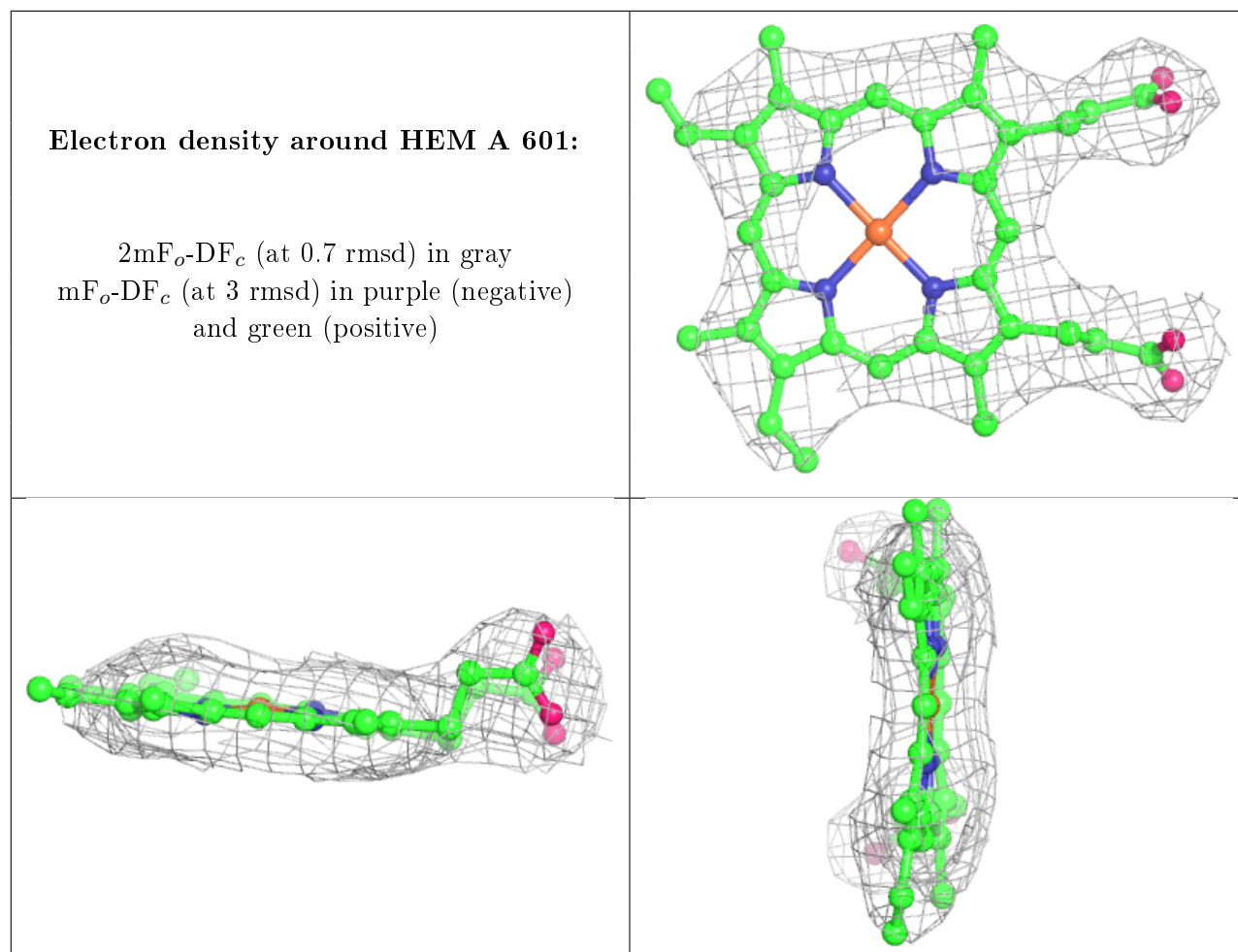
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	QES	A	602	40/40	0.87	0.42	61,167,208,234	0
2	HEM	A	601	43/43	0.98	0.26	63,82,94,104	0

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 QES A 602:**

$2mF_o-DF_c$  (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.