



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2026 – 03:47 PM EDT

PDB ID : 8EWN / pdb_00008ewn
Title : Crystal structure of CYP3A4 bound to an inhibitor
Authors : Sevrioukova, I.F.
Deposited on : 2022-10-24
Resolution : 2.10 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

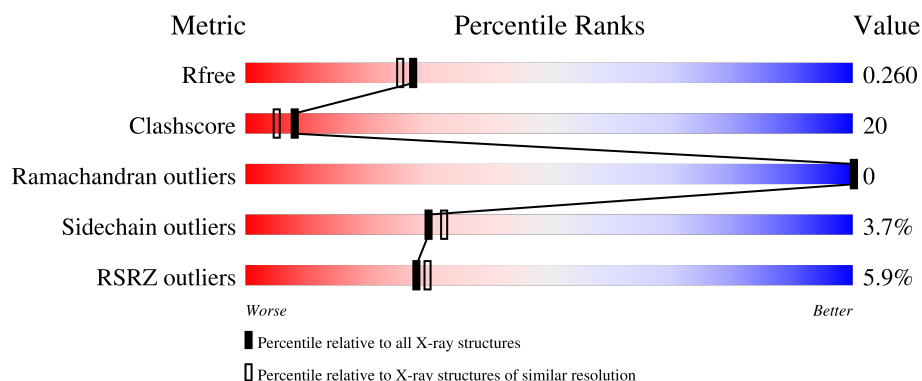
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	487	5% 56% 34% 9%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3663 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.

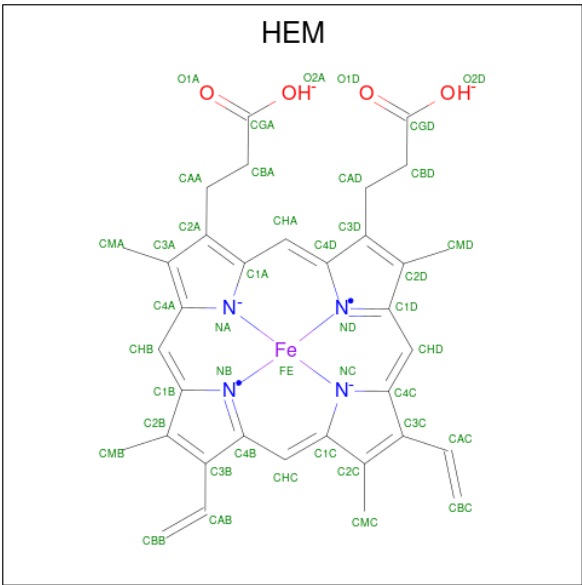
- Molecule 1 is a protein called Cytochrome P450 3A4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	442	Total	C	N	O	S	0	0	0
			3552	2317	582	629	24			

There are 24 discrepancies between the modelled and reference sequences:

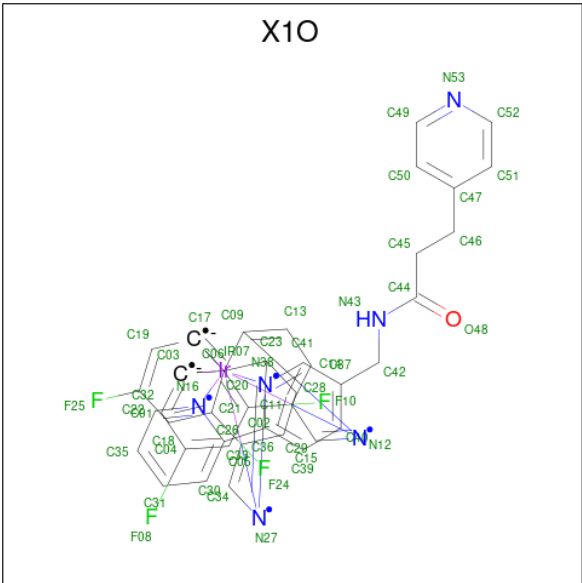
Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	LEU	deletion	UNP P08684
A	?	-	ILE	deletion	UNP P08684
A	?	-	PRO	deletion	UNP P08684
A	?	-	ASP	deletion	UNP P08684
A	?	-	LEU	deletion	UNP P08684
A	?	-	ALA	deletion	UNP P08684
A	?	-	MET	deletion	UNP P08684
A	?	-	GLU	deletion	UNP P08684
A	?	-	THR	deletion	UNP P08684
A	?	-	TRP	deletion	UNP P08684
A	?	-	LEU	deletion	UNP P08684
A	?	-	LEU	deletion	UNP P08684
A	?	-	LEU	deletion	UNP P08684
A	?	-	ALA	deletion	UNP P08684
A	?	-	VAL	deletion	UNP P08684
A	?	-	SER	deletion	UNP P08684
A	?	-	LEU	deletion	UNP P08684
A	?	-	VAL	deletion	UNP P08684
A	?	-	LEU	deletion	UNP P08684
A	?	-	LEU	deletion	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 (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

- Molecule 3 is {N-[(2,2'-bipyridin]-5-yl-kappa 2 N 1 ,N 1')methyl]-3-(pyridin-4-yl)propanamide}bis[3,5-difluoro-2-(pyridin-2-yl-kappaN)phenyl-kappaC 1]iridium(1+) (CCD ID: X1O) (formula: C₄₁H₃₀F₄IrN₆O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	F	Ir	N	O	0	0
			53	41	4	1	6	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	15	Total	O	0	0
			15	15		

4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	75.52Å 99.29Å 125.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.39 – 2.10 39.39 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.1 (39.39-2.10) 99.1 (39.39-2.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.30 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.237 , 0.257 0.242 , 0.260	Depositor DCC
R_{free} test set	1375 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	64.0	Xtriage
Anisotropy	0.214	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 68.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3663	wwPDB-VP
Average B, all atoms (Å ²)	96.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.36% 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: HEM, X1O

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	0.48	0/3639	0.66	1/4920 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	344	PRO	N-CA-C	5.41	117.31	110.70

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	3552	0	3643	136	0
2	A	43	0	30	5	0
3	A	53	0	0	6	0
4	A	15	0	0	0	0
All	All	3663	0	3673	143	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 20.

All (143) 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:188:SER:HA	1:A:271:PHE:HB2	1.62	0.81
1:A:210:LEU:HD11	1:A:245:VAL:HG11	1.62	0.80
1:A:472:GLN:HE21	1:A:476:LYS:H	1.29	0.80
1:A:255:ARG:NH1	1:A:256:MET:HG2	1.97	0.79
1:A:152:TYR:HA	1:A:155:VAL:HG12	1.64	0.78
1:A:255:ARG:HH12	1:A:256:MET:HG2	1.46	0.77
1:A:317:ILE:HD13	1:A:363:THR:HG21	1.67	0.76
1:A:87:PRO:HG3	1:A:431:ILE:HD11	1.69	0.75
1:A:132:LEU:HD13	1:A:290:LEU:HG	1.69	0.74
1:A:140:GLY:HA2	1:A:143:LYS:HD3	1.70	0.73
1:A:59:MET:O	1:A:63:GLU:HG2	1.89	0.71
1:A:247:ASN:O	1:A:251:LYS:HG3	1.91	0.70
1:A:458:ARG:NH1	1:A:461:GLN:OE1	2.25	0.70
1:A:256:MET:SD	1:A:272:LEU:HD13	2.32	0.69
1:A:181:MET:HE1	1:A:207:THR:HB	1.75	0.67
1:A:487:LYS:HD3	1:A:488:PRO:HD2	1.76	0.67
1:A:117:ALA:HB1	1:A:301:ILE:HG13	1.81	0.61
1:A:334:GLU:HG2	1:A:353:MET:HE3	1.82	0.61
1:A:106:ARG:HD2	1:A:108:PHE:CE2	2.36	0.60
1:A:171:THR:HA	1:A:490:VAL:HG12	1.82	0.60
1:A:335:ILE:CG1	1:A:353:MET:HE1	2.32	0.60
1:A:153:GLY:O	1:A:157:VAL:HG13	2.03	0.59
1:A:299:SER:O	1:A:303:ILE:HD12	2.02	0.59
1:A:420:SER:OG	1:A:422:LYS:HD2	2.01	0.59
1:A:172:LEU:HB3	1:A:176:PHE:CD1	2.38	0.59
1:A:229:LEU:HA	1:A:232:ILE:HD12	1.84	0.58
1:A:71:VAL:HG11	1:A:386:MET:HE1	1.86	0.58
1:A:141:LYS:N	1:A:141:LYS:HD2	2.19	0.58
1:A:210:LEU:H	1:A:210:LEU:HD12	1.69	0.58
1:A:335:ILE:HG12	1:A:353:MET:HE1	1.85	0.58
1:A:383:ILE:O	1:A:384:ASN:C	2.46	0.57
1:A:154:ASP:O	1:A:157:VAL:HG22	2.04	0.57
1:A:187:THR:O	1:A:271:PHE:N	2.37	0.57
1:A:232:ILE:O	1:A:236:LEU:HG	2.05	0.57
1:A:169:PRO:HG3	1:A:468:CYS:SG	2.45	0.57
3:A:602:X1O:C09	3:A:602:X1O:C19	2.83	0.57
1:A:183:VAL:O	1:A:187:THR:HG22	2.06	0.56
1:A:148:ILE:H	1:A:148:ILE:HD12	1.70	0.56
1:A:206:ASN:HD21	1:A:245:VAL:HA	1.70	0.56
1:A:296:VAL:O	1:A:300:ILE:HG13	2.06	0.56
1:A:305:ALA:O	1:A:309:THR:OG1	2.21	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:ASP:O	1:A:158:ARG:HG3	2.05	0.55
1:A:442:CYS:HB2	2:A:601:HEM:NA	2.22	0.54
1:A:159:ASN:ND2	1:A:197:ASN:HD21	2.05	0.54
3:A:602:X1O:C19	3:A:602:X1O:N12	2.69	0.54
3:A:602:X1O:C01	3:A:602:X1O:N27	2.68	0.54
1:A:249:LEU:O	1:A:253:VAL:HG23	2.08	0.54
2:A:601:HEM:C1B	3:A:602:X1O:C49	2.91	0.54
1:A:458:ARG:O	1:A:462:ASN:ND2	2.39	0.53
1:A:298:GLN:N	1:A:298:GLN:OE1	2.41	0.53
3:A:602:X1O:C01	3:A:602:X1O:C23	2.87	0.53
3:A:602:X1O:C09	3:A:602:X1O:C32	2.86	0.53
1:A:152:TYR:CD1	1:A:182:ASP:HB3	2.44	0.53
1:A:226:PHE:HA	1:A:228:PHE:CE1	2.44	0.53
1:A:194:ASP:HB3	1:A:198:ASN:OD1	2.09	0.52
1:A:119:SER:HB3	2:A:601:HEM:HAD1	1.91	0.52
1:A:345:PRO:HB2	1:A:454:LEU:HD23	1.92	0.52
1:A:410:GLU:HB3	1:A:413:LYS:HG3	1.91	0.52
1:A:114:MET:HE1	1:A:301:ILE:HG12	1.93	0.51
1:A:159:ASN:HD22	1:A:197:ASN:HD21	1.58	0.51
1:A:442:CYS:HB2	2:A:601:HEM:C4A	2.45	0.51
1:A:99:TYR:CE1	1:A:127:LYS:HE2	2.45	0.51
1:A:162:ARG:HD3	1:A:163:GLU:N	2.26	0.51
1:A:253:VAL:HG21	1:A:296:VAL:HG22	1.93	0.51
1:A:345:PRO:HB2	1:A:454:LEU:CD2	2.41	0.51
1:A:176:PHE:CD2	1:A:314:LEU:HD23	2.47	0.50
1:A:160:LEU:O	1:A:164:ALA:N	2.42	0.50
1:A:155:VAL:HA	1:A:158:ARG:HE	1.76	0.50
1:A:172:LEU:HB3	1:A:176:PHE:CE1	2.48	0.49
1:A:271:PHE:O	1:A:275:MET:HG3	2.12	0.48
1:A:275:MET:HB2	1:A:295:LEU:HD11	1.93	0.48
1:A:438:GLY:O	1:A:441:ASN:ND2	2.47	0.48
2:A:601:HEM:HBB2	2:A:601:HEM:HMB2	1.96	0.47
1:A:209:LYS:HA	1:A:209:LYS:HE2	1.96	0.47
1:A:59:MET:N	1:A:59:MET:HE2	2.30	0.47
1:A:180:SER:OG	1:A:307:TYR:HA	2.15	0.47
1:A:133:LEU:HD13	1:A:133:LEU:HA	1.75	0.46
1:A:226:PHE:N	1:A:227:PRO:HD3	2.30	0.46
1:A:339:LEU:HG	1:A:343:ALA:HB3	1.96	0.46
1:A:253:VAL:HG11	1:A:295:LEU:HD23	1.96	0.46
1:A:349:THR:HA	1:A:352:GLN:OE1	2.16	0.46
1:A:471:THR:OG1	1:A:490:VAL:O	2.29	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:ILE:HG21	1:A:461:GLN:HG3	1.99	0.45
1:A:143:LYS:HG2	1:A:144:GLU:N	2.30	0.45
1:A:236:LEU:HB2	1:A:238:ILE:HG13	1.99	0.45
1:A:251:LYS:HA	1:A:254:LYS:HG2	1.99	0.45
1:A:421:LYS:HA	1:A:424:LYS:HD3	1.98	0.45
1:A:255:ARG:HH11	1:A:255:ARG:HG3	1.82	0.45
1:A:252:SER:O	1:A:256:MET:HG3	2.16	0.45
1:A:339:LEU:HD21	1:A:345:PRO:HA	1.99	0.45
1:A:456:LEU:HD23	1:A:456:LEU:HA	1.77	0.45
1:A:483:LEU:HD12	1:A:483:LEU:HA	1.73	0.45
1:A:91:LYS:HG3	1:A:92:THR:N	2.32	0.45
1:A:159:ASN:ND2	1:A:197:ASN:ND2	2.65	0.45
1:A:379:LYS:O	1:A:381:VAL:HG13	2.17	0.45
1:A:110:PRO:HB3	1:A:233:LEU:HB2	1.99	0.44
1:A:184:ILE:HG12	1:A:303:ILE:HA	1.98	0.44
1:A:181:MET:HG3	1:A:195:SER:OG	2.17	0.44
1:A:193:ILE:HD12	1:A:204:VAL:HG23	1.98	0.44
1:A:88:ASP:HA	1:A:91:LYS:HG2	2.00	0.44
1:A:149:ILE:HG22	1:A:183:VAL:HG13	1.98	0.44
1:A:226:PHE:HB3	1:A:229:LEU:HG	2.00	0.44
1:A:466:LYS:HB3	1:A:466:LYS:HE3	1.60	0.44
1:A:125:GLU:OE2	1:A:129:LEU:HG	2.18	0.44
1:A:338:VAL:CG1	1:A:349:THR:HG23	2.48	0.44
1:A:179:TYR:O	1:A:183:VAL:HG23	2.17	0.44
1:A:357:ASP:O	1:A:360:VAL:HG22	2.18	0.44
1:A:226:PHE:HD1	1:A:228:PHE:CZ	2.36	0.43
1:A:348:ASP:O	1:A:352:GLN:HG3	2.18	0.43
1:A:146:VAL:HG13	1:A:454:LEU:HD11	2.00	0.43
1:A:335:ILE:HG13	1:A:353:MET:HE1	1.99	0.43
1:A:112:GLY:O	1:A:115:LYS:HG2	2.19	0.43
1:A:133:LEU:HD11	1:A:275:MET:SD	2.59	0.43
1:A:206:ASN:OD1	1:A:245:VAL:HG13	2.18	0.43
1:A:189:PHE:O	1:A:256:MET:HE1	2.20	0.42
1:A:194:ASP:CG	1:A:197:ASN:HD22	2.25	0.42
1:A:197:ASN:C	1:A:199:PRO:HD3	2.43	0.42
1:A:124:GLU:H	1:A:124:GLU:CD	2.26	0.42
1:A:126:TRP:CZ2	1:A:130:ARG:HD3	2.54	0.42
1:A:142:LEU:HA	1:A:145:MET:HE3	2.01	0.42
1:A:479:LEU:HD12	1:A:479:LEU:HA	1.79	0.42
1:A:477:LEU:HD23	1:A:477:LEU:HA	1.87	0.42
1:A:172:LEU:HB2	1:A:489:VAL:HG23	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:248:PHE:CD2	1:A:248:PHE:C	2.97	0.42
1:A:159:ASN:HB3	1:A:196:LEU:HD23	2.01	0.42
1:A:321:LEU:HD23	1:A:321:LEU:HA	1.86	0.42
1:A:369:ILE:HD12	1:A:370:ALA:N	2.34	0.42
1:A:136:THR:HG23	1:A:274:LEU:HD13	2.02	0.41
1:A:339:LEU:HD12	1:A:339:LEU:HA	1.76	0.41
1:A:351:LEU:CD2	1:A:450:MET:HE1	2.51	0.41
1:A:487:LYS:HD3	1:A:487:LYS:HA	1.87	0.41
1:A:141:LYS:HD2	1:A:141:LYS:H	1.85	0.41
1:A:172:LEU:HB3	1:A:176:PHE:HD1	1.85	0.41
1:A:224:THR:C	1:A:227:PRO:HD3	2.45	0.41
1:A:101:VAL:HG11	1:A:381:VAL:HG11	2.02	0.41
1:A:181:MET:HE1	1:A:207:THR:CB	2.45	0.41
1:A:290:LEU:HD23	1:A:290:LEU:HA	1.96	0.40
1:A:410:GLU:CB	1:A:413:LYS:HG3	2.50	0.40
1:A:146:VAL:N	1:A:147:PRO:HD2	2.36	0.40
1:A:146:VAL:HA	1:A:149:ILE:HG12	2.03	0.40
1:A:142:LEU:O	1:A:146:VAL:HG23	2.22	0.40
1:A:191:VAL:HG21	1:A:203:PHE:CD2	2.57	0.40
1:A:356:LEU:HD21	1:A:453:LYS:HB3	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	434/487 (89%)	408 (94%)	26 (6%)	0	100	100

There are no Ramachandran outliers to report.

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	403/443 (91%)	388 (96%)	15 (4%)	30	33

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	118	ILE
1	A	139	SER
1	A	165	GLU
1	A	168	LYS
1	A	191	VAL
1	A	198	ASN
1	A	204	VAL
1	A	210	LEU
1	A	276	ILE
1	A	339	LEU
1	A	342	LYS
1	A	384	ASN
1	A	459	VAL
1	A	479	LEU
1	A	482	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	197	ASN
1	A	206	ASN
1	A	472	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	X1O	A	602	2	55,62,62	3.05	14 (25%)	36,100,100	2.36	10 (27%)
2	HEM	A	601	3,1	50,50,50	1.42	7 (14%)	67,82,82	1.41	12 (17%)

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	X1O	A	602	2	-	1/9/160/160	0/10/10/10
2	HEM	A	601	3,1	-	2/14/54/54	-

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	X1O	IR07-N27	-10.02	1.91	2.15
3	A	602	X1O	IR07-N12	-9.97	1.91	2.15
3	A	602	X1O	IR07-N38	-7.37	1.97	2.15
3	A	602	X1O	IR07-N16	-6.57	1.99	2.15
3	A	602	X1O	C36-C33	-6.31	1.32	1.45
3	A	602	X1O	C44-N43	5.84	1.47	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	X1O	C36-N38	-5.46	1.33	1.40
3	A	602	X1O	C18-C21	5.19	1.41	1.35
3	A	602	X1O	C33-N16	-4.16	1.34	1.40
2	A	601	HEM	FE-NC	3.52	2.06	1.95
2	A	601	HEM	FE-ND	3.23	2.04	1.94
3	A	602	X1O	C03-IR07	-3.22	1.92	2.00
3	A	602	X1O	C41-N38	3.03	1.36	1.29
2	A	601	HEM	CAC-C3C	2.93	1.55	1.47
3	A	602	X1O	C17-IR07	-2.91	1.93	2.00
2	A	601	HEM	CMC-C2C	2.80	1.56	1.50
3	A	602	X1O	C09-N12	-2.75	1.36	1.40
2	A	601	HEM	CAB-C3B	2.74	1.54	1.47
2	A	601	HEM	FE-NB	2.71	2.03	1.94
2	A	601	HEM	FE-NA	2.65	2.03	1.95
3	A	602	X1O	C32-N16	2.36	1.35	1.29

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	X1O	O48-C44-N43	-7.49	108.32	123.03
3	A	602	X1O	C45-C44-N43	6.62	128.41	116.34
3	A	602	X1O	C06-C09-N12	3.49	114.91	108.76
3	A	602	X1O	C41-N38-C36	-3.13	114.57	118.40
3	A	602	X1O	C40-C39-C36	-3.03	118.24	122.59
3	A	602	X1O	C39-C40-C37	-2.95	117.34	122.83
2	A	601	HEM	C4D-ND-C1D	2.75	108.46	105.21
2	A	601	HEM	CHD-C1D-ND	2.56	127.18	124.42
2	A	601	HEM	C1B-NB-C4B	2.54	108.21	105.21
2	A	601	HEM	CMA-C3A-C4A	-2.50	121.61	125.42
2	A	601	HEM	CBD-CAD-C3D	-2.45	105.75	112.53
2	A	601	HEM	CBA-CAA-C2A	-2.44	105.77	112.53
2	A	601	HEM	C2A-C1A-NA	-2.43	107.46	110.15
2	A	601	HEM	C3B-C2B-C1B	2.33	108.16	106.41
2	A	601	HEM	O1A-CGA-CBA	-2.31	115.76	123.09
3	A	602	X1O	F24-C21-C18	-2.20	115.43	118.67
2	A	601	HEM	CHC-C4B-NB	2.17	126.76	124.42
3	A	602	X1O	C01-C03-C06	-2.15	119.35	123.08
3	A	602	X1O	C20-C23-N27	2.14	116.45	110.19
2	A	601	HEM	CMA-C3A-C2A	2.13	130.15	125.62
3	A	602	X1O	C52-N53-C49	2.05	121.56	116.86
2	A	601	HEM	CHA-C4D-ND	2.01	126.85	124.37

There are no chirality outliers.

All (3) torsion outliers are listed below:

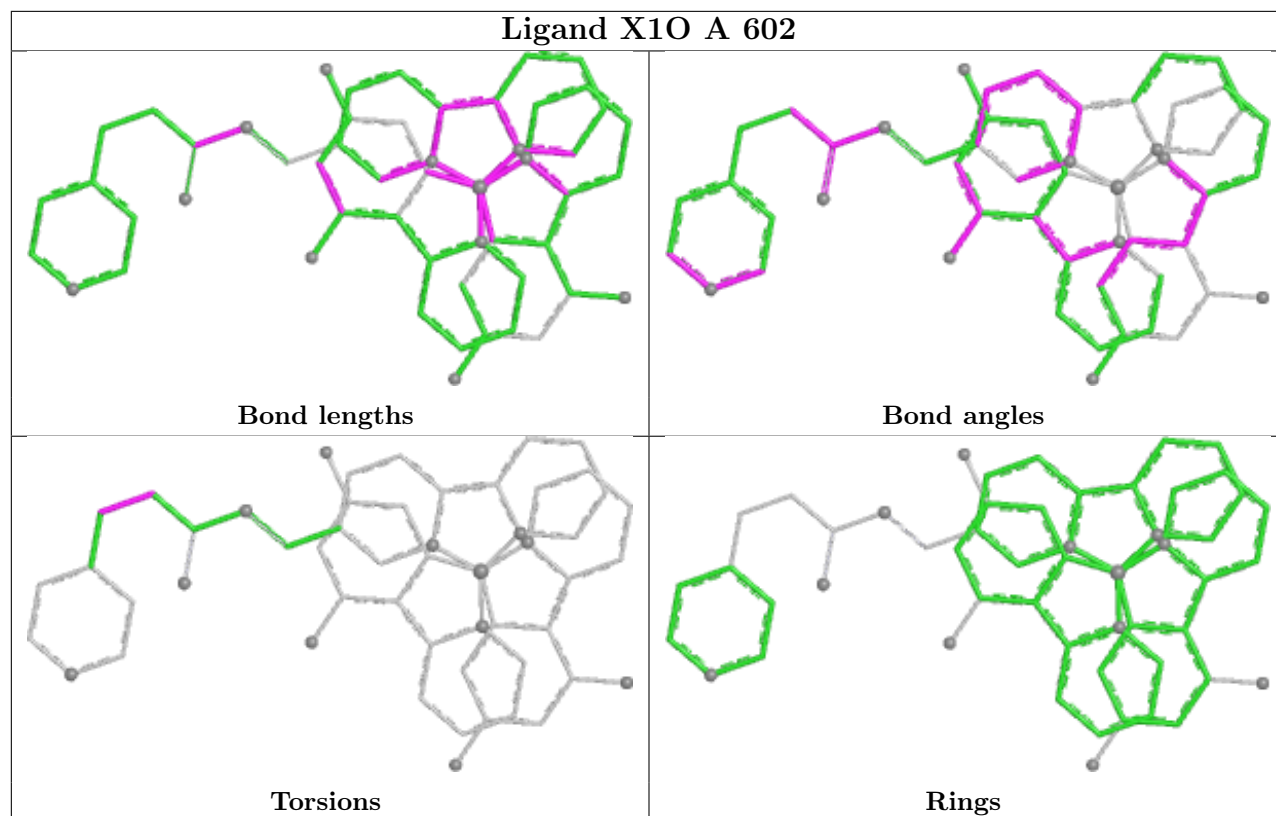
Mol	Chain	Res	Type	Atoms
3	A	602	X1O	C44-C45-C46-C47
2	A	601	HEM	CAD-CBD-CGD-O2D
2	A	601	HEM	CAD-CBD-CGD-O1D

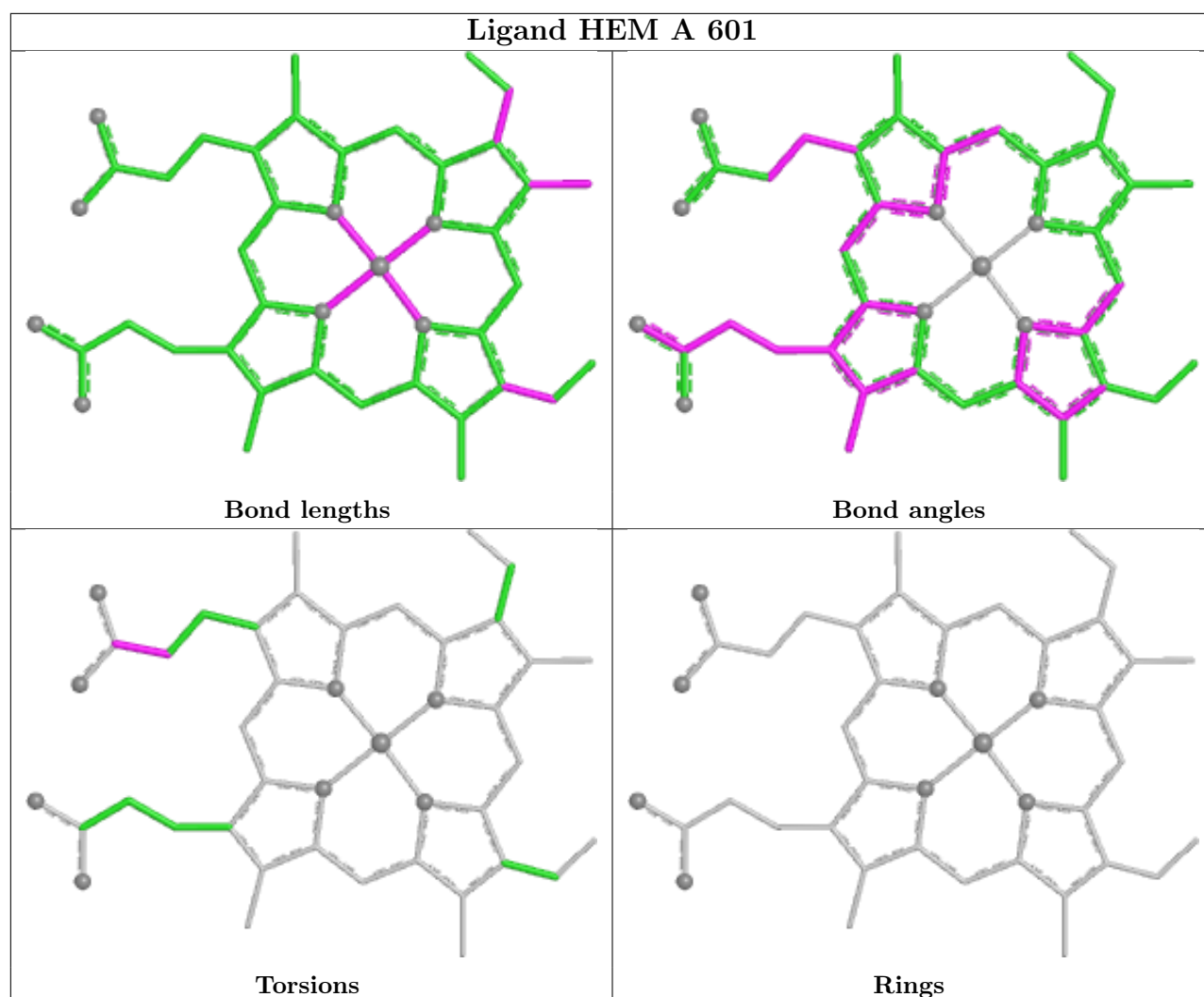
There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	X1O	6	0
2	A	601	HEM	5	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, 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	442/487 (90%)	0.56	26 (5%)	28 30	53, 92, 151, 169	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	241	PHE	3.9
1	A	218	PRO	3.7
1	A	203	PHE	2.7
1	A	322	ALA	2.7
1	A	221	LEU	2.6
1	A	207	THR	2.5
1	A	177	GLY	2.5
1	A	149	ILE	2.5
1	A	46	PHE	2.5
1	A	210	LEU	2.4
1	A	226	PHE	2.3
1	A	271	PHE	2.3
1	A	308	GLU	2.2
1	A	228	PHE	2.1
1	A	183	VAL	2.1
1	A	109	GLY	2.1
1	A	240	VAL	2.1
1	A	297	ALA	2.1
1	A	222	SER	2.1
1	A	156	LEU	2.1
1	A	274	LEU	2.1
1	A	290	LEU	2.1
1	A	242	PRO	2.1
1	A	108	PHE	2.1
1	A	191	VAL	2.1
1	A	384	ASN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

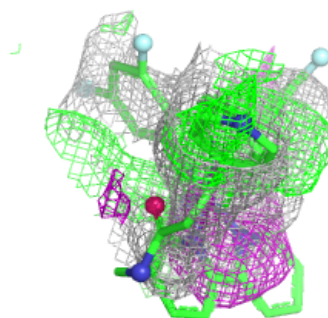
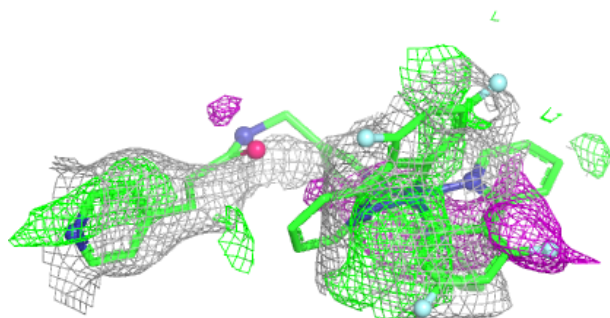
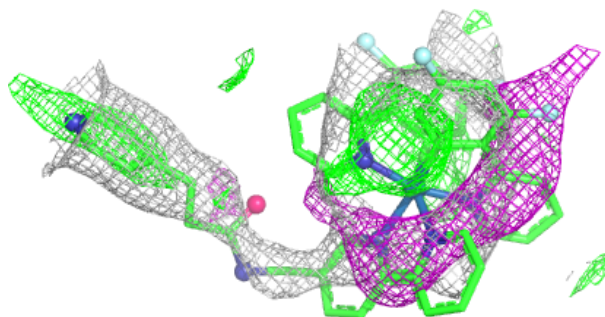
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	X1O	A	602	53/53	0.86	0.19	108,164,181,184	0
2	HEM	A	601	43/43	0.98	0.08	53,67,74,79	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.

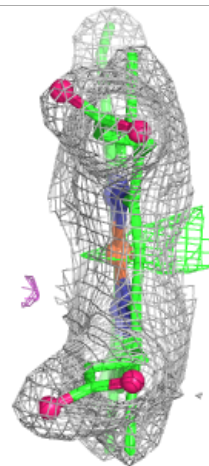
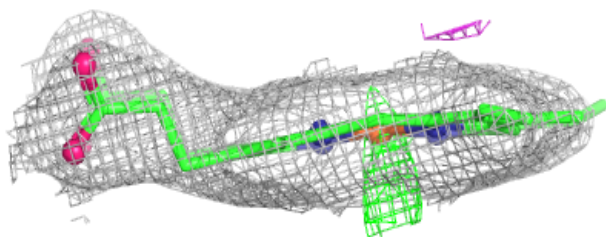
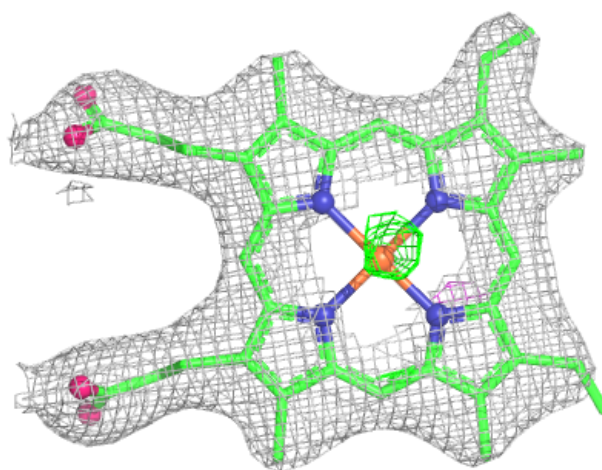
Electron density around X1O A 602:

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



Electron density around HEM A 601:

$2mF_o-DF_c$ (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.