



Full wwPDB X-ray Structure Validation Report ⓘ

May 7, 2026 – 07:42 AM EDT

PDB ID : 5A5J / pdb_00005a5j
Title : Cytochrome 2C9 P450 inhibitor complex
Authors : Skerratt, S.E.; de Groot, M.J.; Phillips, C.
Deposited on : 2015-06-18
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 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 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

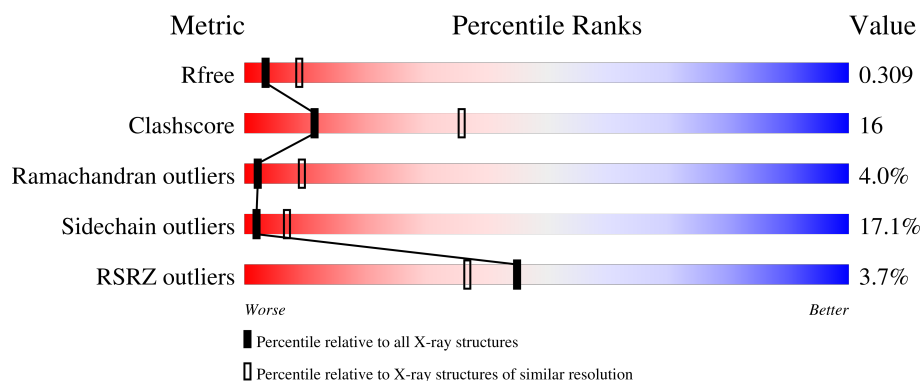
1 Overall quality at a glance

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}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (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 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	477	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3717 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 2C9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	455	3646	2353	614	656	23	0	0	0

There are 10 discrepancies between the modelled and reference sequences:

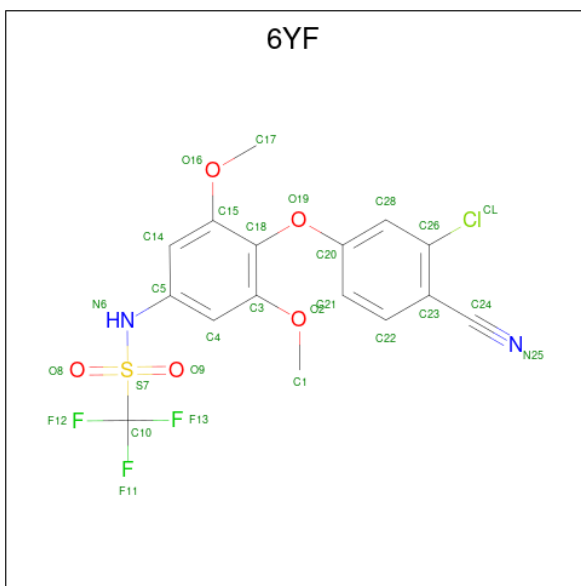
Chain	Residue	Modelled	Actual	Comment	Reference
A	18	MET	-	expression tag	UNP P11712
A	19	ALA	-	expression tag	UNP P11712
A	20	LYS	-	expression tag	UNP P11712
A	21	LYS	-	expression tag	UNP P11712
A	22	THR	-	expression tag	UNP P11712
A	490	ILE	-	expression tag	UNP P11712
A	491	HIS	-	expression tag	UNP P11712
A	492	HIS	-	expression tag	UNP P11712
A	493	HIS	-	expression tag	UNP P11712
A	494	HIS	-	expression tag	UNP P11712

- 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-[4-(3-chloranyl-4-cyano-phenoxy)-3,5-dimethoxy-phenyl]-1,1,1-tris(fluoranyl)methanesulfonamide (CCD ID: 6YF) (formula: $\text{C}_{16}\text{H}_{12}\text{ClF}_3\text{N}_2\text{O}_5\text{S}$).

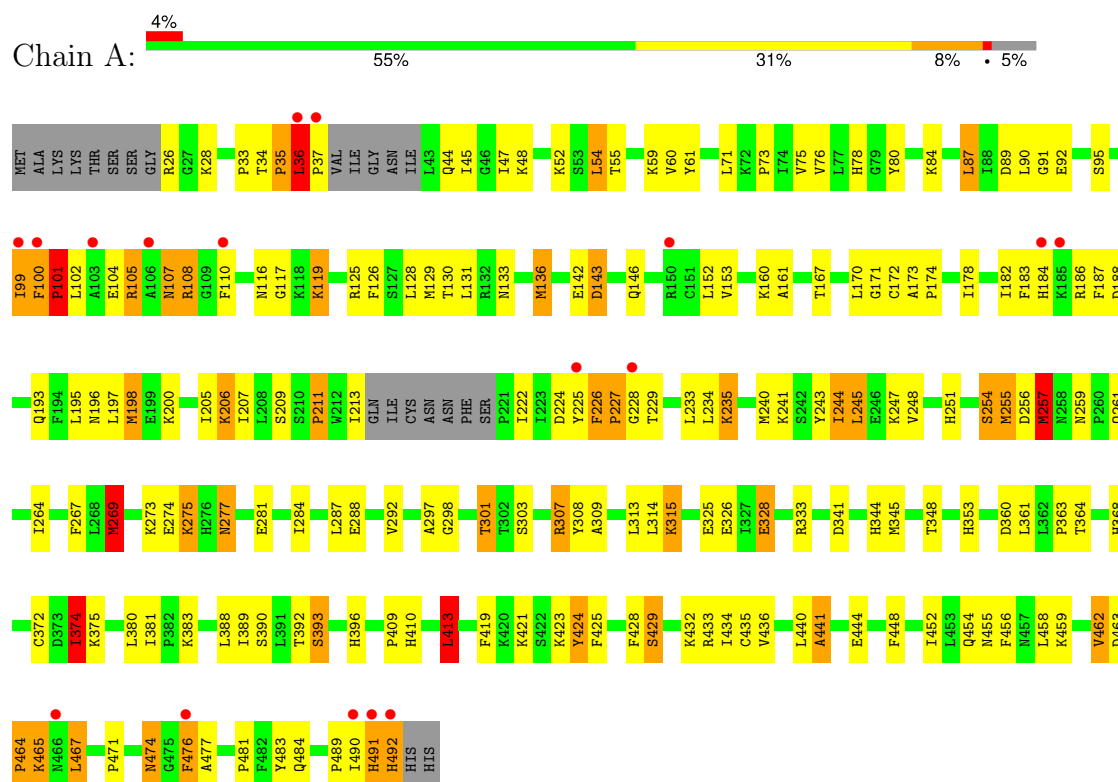


Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
3	A	1	Total	C	Cl	F	N	O	S	0	0
			28	16	1	3	2	5	1		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: CYTOCHROME P450 2C9



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	91.29Å 91.29Å 169.37Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	71.74 – 2.90 71.64 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.7 (71.74-2.90) 99.7 (71.64-2.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.213 , 0.288 0.225 , 0.309	Depositor DCC
R_{free} test set	556 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	61.1	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 33.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.55$, $\langle L^2 \rangle = 0.39$	Xtriage
Estimated twinning fraction	0.009 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3717	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.37% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, 6YF

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.29	8/3735 (0.2%)	1.30	19/5047 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	269	MET	SD-CE	10.69	2.06	1.79
1	A	257	MET	SD-CE	10.55	2.06	1.79
1	A	255	MET	SD-CE	8.70	2.01	1.79
1	A	492	HIS	CE1-NE2	7.39	1.40	1.32
1	A	101	PRO	CA-C	6.56	1.61	1.52
1	A	413	LEU	CA-C	6.03	1.59	1.52
1	A	424	TYR	CA-C	5.13	1.62	1.52
1	A	73	PRO	C-O	5.03	1.29	1.23

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	374	ILE	N-CA-C	7.55	118.31	108.35
1	A	490	ILE	N-CA-C	7.45	118.63	110.21
1	A	206	LYS	N-CA-C	-7.13	102.55	111.11
1	A	235	LYS	N-CA-C	-6.93	103.65	111.07
1	A	75	VAL	CB-CA-C	-6.68	102.39	111.33

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	60	VAL	N-CA-C	6.39	116.43	110.42
1	A	452	ILE	N-CA-C	-6.39	104.10	110.62
1	A	477	ALA	N-CA-C	6.24	118.78	108.99
1	A	60	VAL	CB-CA-C	-6.19	103.88	111.87
1	A	211	PRO	CA-C-O	-5.83	114.66	121.95
1	A	116	ASN	N-CA-C	5.47	115.76	108.38
1	A	229	THR	N-CA-C	-5.36	108.51	114.62
1	A	474	ASN	CA-C-N	5.33	125.75	119.94
1	A	474	ASN	C-N-CA	5.33	125.75	119.94
1	A	424	TYR	N-CA-C	5.26	119.41	113.15
1	A	254	SER	N-CA-C	5.19	119.11	112.26
1	A	441	ALA	N-CA-C	-5.17	105.54	111.07
1	A	372	CYS	CA-CB-SG	-5.17	102.52	114.40
1	A	227	PRO	N-CA-C	5.13	120.51	113.84

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	489	PRO	Peptide

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	3646	0	3678	117	0
2	A	43	0	30	7	0
3	A	28	0	0	1	0
All	All	3717	0	3708	119	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 16.

All (119) 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:257:MET:CE	1:A:257:MET:SD	2.05	1.43
1:A:269:MET:CE	1:A:269:MET:SD	2.06	1.42
1:A:172:CYS:SG	1:A:198:MET:HE1	2.15	0.87
1:A:245:LEU:HD22	1:A:288:GLU:OE2	1.80	0.81
1:A:353:HIS:HD2	1:A:425:PHE:CZ	1.99	0.80
1:A:193:GLN:NE2	1:A:247:LYS:NZ	2.30	0.79
1:A:491:HIS:C	1:A:492:HIS:ND1	2.42	0.78
1:A:47:ILE:HG23	1:A:363:PRO:HG2	1.67	0.77
1:A:209:SER:HB2	1:A:474:ASN:HD21	1.49	0.76
1:A:36:LEU:N	1:A:37:PRO:HD3	2.05	0.72
1:A:193:GLN:NE2	1:A:247:LYS:HZ2	1.88	0.72
1:A:35:PRO:C	1:A:37:PRO:HD3	2.15	0.71
1:A:178:ILE:HG22	1:A:182:ILE:HD12	1.70	0.71
1:A:209:SER:HB2	1:A:474:ASN:ND2	2.06	0.70
1:A:90:LEU:HD22	1:A:374:ILE:HD11	1.77	0.67
1:A:92:GLU:HG2	1:A:432:LYS:NZ	2.10	0.67
1:A:193:GLN:HE21	1:A:247:LYS:HZ1	1.43	0.66
1:A:353:HIS:HD2	1:A:425:PHE:CE2	2.15	0.65
1:A:368:HIS:HD2	1:A:389:ILE:HD13	1.61	0.65
1:A:119:LYS:HG3	1:A:281:GLU:OE1	1.97	0.64
1:A:454:GLN:HB3	1:A:455:ASN:ND2	2.13	0.63
1:A:211:PRO:O	1:A:224:ASP:HB2	1.99	0.63
1:A:309:ALA:O	1:A:313:LEU:HG	1.99	0.62
3:A:1494:6YF:C14	3:A:1494:6YF:O8	2.47	0.62
1:A:267:PHE:CG	1:A:287:LEU:HD13	2.36	0.61
1:A:78:HIS:CD2	1:A:393:SER:HB2	2.36	0.61
1:A:251:HIS:CD2	1:A:264:ILE:HB	2.37	0.60
1:A:275:LYS:C	1:A:277:ASN:H	2.12	0.58
1:A:389:ILE:HD12	1:A:389:ILE:N	2.18	0.58
1:A:100:PHE:CG	1:A:101:PRO:HD2	2.38	0.58
1:A:193:GLN:NE2	1:A:247:LYS:HZ1	1.99	0.58
1:A:463:ASP:O	1:A:464:PRO:C	2.47	0.57
1:A:492:HIS:ND1	1:A:492:HIS:N	2.52	0.57
1:A:90:LEU:HD22	1:A:374:ILE:CD1	2.35	0.56
1:A:183:PHE:O	1:A:184:HIS:HB3	2.06	0.56
1:A:364:THR:HA	1:A:389:ILE:O	2.05	0.56
1:A:125:ARG:O	1:A:129:MET:HG2	2.06	0.56
1:A:183:PHE:O	1:A:184:HIS:CB	2.54	0.56
1:A:80:TYR:HB2	1:A:424:TYR:CE2	2.42	0.55
2:A:1493:HEM:HMB2	2:A:1493:HEM:HBB2	1.89	0.55
1:A:173:ALA:HB3	1:A:174:PRO:HD3	1.88	0.54
1:A:207:ILE:HG21	1:A:233:LEU:HD13	1.89	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:353:HIS:CD2	1:A:425:PHE:CE2	2.96	0.54
1:A:244:ILE:O	1:A:248:VAL:HG23	2.08	0.54
1:A:130:THR:O	1:A:136:MET:HE2	2.08	0.53
1:A:47:ILE:HD13	1:A:476:PHE:HA	1.90	0.53
1:A:78:HIS:NE2	1:A:393:SER:HB2	2.24	0.53
1:A:171:GLY:HA2	1:A:303:SER:HB2	1.89	0.53
1:A:251:HIS:HD2	1:A:264:ILE:HB	1.73	0.53
1:A:187:PHE:O	1:A:188:ASP:C	2.49	0.52
1:A:108:ARG:CZ	1:A:108:ARG:HB3	2.40	0.51
1:A:187:PHE:C	1:A:188:ASP:O	2.53	0.51
1:A:226:PHE:N	1:A:227:PRO:CD	2.74	0.51
1:A:36:LEU:N	1:A:37:PRO:CD	2.73	0.51
1:A:193:GLN:HE21	1:A:247:LYS:NZ	2.02	0.51
1:A:207:ILE:HD13	1:A:233:LEU:HD13	1.93	0.50
1:A:307:ARG:O	1:A:483:TYR:OH	2.27	0.50
1:A:142:GLU:O	1:A:143:ASP:C	2.56	0.49
1:A:462:VAL:CG2	1:A:467:LEU:HB2	2.43	0.49
1:A:100:PHE:CD1	1:A:101:PRO:HD2	2.48	0.48
1:A:454:GLN:HB3	1:A:455:ASN:HD22	1.77	0.48
1:A:78:HIS:CD2	1:A:393:SER:CB	2.97	0.48
1:A:92:GLU:HG2	1:A:432:LYS:HZ1	1.75	0.48
1:A:410:HIS:HA	1:A:413:LEU:HB2	1.94	0.48
1:A:259:ASN:HD21	1:A:261:GLN:HE21	1.60	0.48
1:A:99:ILE:O	1:A:102:LEU:O	2.33	0.47
1:A:245:LEU:HD22	1:A:288:GLU:CD	2.39	0.47
1:A:259:ASN:HD21	1:A:261:GLN:NE2	2.12	0.47
1:A:368:HIS:CD2	1:A:389:ILE:HD13	2.46	0.47
1:A:245:LEU:CD2	1:A:288:GLU:OE2	2.59	0.47
1:A:463:ASP:O	1:A:465:LYS:N	2.48	0.47
1:A:281:GLU:OE1	1:A:281:GLU:HA	2.14	0.47
1:A:196:ASN:O	1:A:200:LYS:HG2	2.15	0.47
1:A:234:LEU:O	1:A:235:LYS:C	2.58	0.46
1:A:178:ILE:HG22	1:A:182:ILE:CD1	2.42	0.46
1:A:428:PHE:HB3	1:A:435:CYS:HB3	1.97	0.46
1:A:315:LYS:O	1:A:315:LYS:HG3	2.17	0.45
1:A:243:TYR:CD1	1:A:243:TYR:C	2.95	0.45
1:A:297:ALA:O	1:A:301:THR:OG1	2.34	0.45
1:A:374:ILE:HG13	1:A:375:LYS:N	2.31	0.45
1:A:275:LYS:C	1:A:277:ASN:N	2.71	0.45
1:A:128:LEU:HD23	1:A:129:MET:SD	2.58	0.44
1:A:54:LEU:N	1:A:54:LEU:HD23	2.33	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:GLY:HA2	2:A:1493:HEM:C2C	2.53	0.44
1:A:462:VAL:HG21	1:A:467:LEU:HB2	2.00	0.44
1:A:160:LYS:O	1:A:161:ALA:HB3	2.17	0.44
1:A:390:SER:HB3	1:A:393:SER:HB3	2.00	0.44
1:A:471:PRO:HA	1:A:481:PRO:HD3	2.00	0.44
1:A:440:LEU:O	1:A:441:ALA:C	2.58	0.43
1:A:170:LEU:O	1:A:174:PRO:HD3	2.19	0.43
1:A:434:ILE:O	1:A:435:CYS:C	2.61	0.43
1:A:267:PHE:CD2	1:A:287:LEU:HD13	2.53	0.43
1:A:172:CYS:HA	1:A:198:MET:SD	2.58	0.43
1:A:325:GLU:O	1:A:328:GLU:N	2.52	0.43
1:A:428:PHE:O	1:A:429:SER:HB3	2.18	0.43
1:A:107:ASN:ND2	1:A:110:PHE:CD2	2.87	0.42
1:A:314:LEU:HD21	1:A:458:LEU:HB3	2.01	0.42
1:A:226:PHE:CD1	1:A:227:PRO:N	2.86	0.42
1:A:436:VAL:HG12	2:A:1493:HEM:C2D	2.54	0.42
1:A:436:VAL:HG12	2:A:1493:HEM:CMD	2.50	0.42
1:A:209:SER:HB2	1:A:474:ASN:CG	2.44	0.42
1:A:301:THR:HG22	1:A:361:LEU:HD21	2.02	0.42
1:A:105:ARG:C	1:A:107:ASN:H	2.27	0.42
1:A:171:GLY:HA2	1:A:303:SER:CB	2.49	0.42
1:A:76:VAL:HG22	1:A:388:LEU:HB3	2.02	0.42
1:A:167:THR:HA	1:A:307:ARG:HD2	2.01	0.42
1:A:126:PHE:CZ	1:A:267:PHE:HA	2.55	0.42
1:A:153:VAL:HG13	1:A:456:PHE:CE2	2.54	0.42
1:A:436:VAL:HG12	2:A:1493:HEM:HMD2	2.02	0.42
1:A:353:HIS:CD2	1:A:425:PHE:CZ	2.92	0.41
1:A:444:GLU:HB3	1:A:448:PHE:CE2	2.56	0.41
1:A:33:PRO:HA	1:A:61:TYR:OH	2.21	0.41
1:A:87:LEU:O	1:A:91:GLY:HA2	2.21	0.41
1:A:360:ASP:O	1:A:361:LEU:C	2.64	0.41
1:A:392:THR:O	1:A:396:HIS:HB2	2.21	0.41
1:A:433:ARG:O	2:A:1493:HEM:HBA2	2.20	0.41
1:A:436:VAL:CG1	2:A:1493:HEM:HMD2	2.51	0.41
1:A:419:PHE:CD2	1:A:419:PHE:C	2.99	0.41
1:A:226:PHE:H	1:A:227:PRO:CD	2.34	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	449/477 (94%)	384 (86%)	47 (10%)	18 (4%)	2 9

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	35	PRO
1	A	36	LEU
1	A	226	PHE
1	A	429	SER
1	A	104	GLU
1	A	107	ASN
1	A	133	ASN
1	A	257	MET
1	A	256	ASP
1	A	326	GLU
1	A	467	LEU
1	A	100	PHE
1	A	101	PRO
1	A	186	ARG
1	A	228	GLY
1	A	464	PRO
1	A	383	LYS
1	A	117	GLY

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	410/430 (95%)	340 (83%)	70 (17%)	2 7

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	ARG
1	A	28	LYS
1	A	34	THR
1	A	36	LEU
1	A	44	GLN
1	A	45	ILE
1	A	48	LYS
1	A	52	LYS
1	A	54	LEU
1	A	55	THR
1	A	59	LYS
1	A	71	LEU
1	A	84	LYS
1	A	87	LEU
1	A	89	ASP
1	A	95	SER
1	A	99	ILE
1	A	105	ARG
1	A	108	ARG
1	A	119	LYS
1	A	131	LEU
1	A	136	MET
1	A	143	ASP
1	A	146	GLN
1	A	152	LEU
1	A	195	LEU
1	A	197	LEU
1	A	198	MET
1	A	205	ILE
1	A	206	LYS
1	A	213	ILE
1	A	222	ILE
1	A	225	TYR
1	A	240	MET
1	A	241	LYS
1	A	244	ILE
1	A	245	LEU
1	A	254	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	255	MET
1	A	269	MET
1	A	273	LYS
1	A	274	GLU
1	A	275	LYS
1	A	277	ASN
1	A	284	ILE
1	A	292	VAL
1	A	301	THR
1	A	307	ARG
1	A	308	TYR
1	A	315	LYS
1	A	328	GLU
1	A	333	ARG
1	A	341	ASP
1	A	344	HIS
1	A	345	MET
1	A	348	THR
1	A	374	ILE
1	A	380	LEU
1	A	381	ILE
1	A	393	SER
1	A	409	PRO
1	A	413	LEU
1	A	421	LYS
1	A	423	LYS
1	A	459	LYS
1	A	462	VAL
1	A	465	LYS
1	A	476	PHE
1	A	484	GLN
1	A	491	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	107	ASN
1	A	192	GLN
1	A	193	GLN
1	A	251	HIS
1	A	259	ASN
1	A	261	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	353	HIS
1	A	457	ASN

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 oligosaccharides 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	6YF	A	1494	-	29,29,29	1.68	3 (10%)	40,43,43	2.39	11 (27%)
2	HEM	A	1493	1	50,50,50	2.02	15 (30%)	67,82,82	1.33	10 (14%)

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	6YF	A	1494	-	-	15/24/24/24	0/2/2/2

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	A	1493	1	-	6/14/54/54	-

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1494	6YF	S7-N6	7.13	1.69	1.59
2	A	1493	HEM	C3D-C2D	6.30	1.50	1.36
2	A	1493	HEM	FE-NA	5.51	2.13	1.95
2	A	1493	HEM	CMA-C3A	3.22	1.57	1.50
2	A	1493	HEM	FE-ND	-3.12	1.85	1.94
2	A	1493	HEM	O1D-CGD	3.08	1.32	1.22
2	A	1493	HEM	CAB-C3B	2.62	1.54	1.47
2	A	1493	HEM	O1A-CGA	2.59	1.30	1.22
2	A	1493	HEM	FE-NB	2.43	2.02	1.94
2	A	1493	HEM	FE-NC	-2.34	1.87	1.95
3	A	1494	6YF	F11-C10	2.29	1.42	1.33
2	A	1493	HEM	CHA-C1A	-2.21	1.34	1.39
2	A	1493	HEM	C3C-C2C	-2.18	1.32	1.37
2	A	1493	HEM	CBA-CGA	2.18	1.55	1.50
2	A	1493	HEM	C3B-C2B	-2.16	1.32	1.37
2	A	1493	HEM	CAC-C3C	2.13	1.53	1.47
3	A	1494	6YF	F13-C10	2.08	1.41	1.33
2	A	1493	HEM	C4D-ND	-2.04	1.36	1.40

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1494	6YF	O9-S7-C10	5.92	114.33	104.08
3	A	1494	6YF	O16-C15-C18	5.55	124.64	115.14
3	A	1494	6YF	O16-C15-C14	-5.01	115.44	124.08
3	A	1494	6YF	O8-S7-C10	4.91	112.58	104.08
3	A	1494	6YF	C20-O19-C18	4.58	125.63	118.40
3	A	1494	6YF	O8-S7-N6	-4.39	101.47	109.42
3	A	1494	6YF	O2-C3-C18	4.00	122.00	115.14
3	A	1494	6YF	O2-C3-C4	-3.63	117.82	124.08
2	A	1493	HEM	C4D-ND-C1D	3.39	109.23	105.21
3	A	1494	6YF	C17-O16-C15	2.85	121.69	117.51
2	A	1493	HEM	CHC-C4B-NB	2.82	127.46	124.42
2	A	1493	HEM	CHD-C1D-ND	2.81	127.45	124.42
2	A	1493	HEM	CHB-C1B-NB	2.75	127.76	124.37
2	A	1493	HEM	CHA-C4D-ND	2.55	127.52	124.37
2	A	1493	HEM	CBD-CAD-C3D	-2.44	105.79	112.53

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1493	HEM	C3B-C4B-NB	2.39	111.19	109.47
2	A	1493	HEM	CBC-CAC-C3C	-2.32	115.92	127.53
3	A	1494	6YF	C28-C26-CL	2.11	121.91	118.45
3	A	1494	6YF	F11-C10-S7	2.05	115.74	110.58
2	A	1493	HEM	CHA-C4D-C3D	-2.03	121.48	125.23
2	A	1493	HEM	CBA-CAA-C2A	2.00	118.08	112.53

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1493	HEM	C2C-C3C-CAC-CBC
3	A	1494	6YF	C5-N6-S7-O9
3	A	1494	6YF	F11-C10-S7-N6
3	A	1494	6YF	F12-C10-S7-N6
3	A	1494	6YF	F13-C10-S7-N6
3	A	1494	6YF	F11-C10-S7-O8
3	A	1494	6YF	F12-C10-S7-O8
3	A	1494	6YF	F13-C10-S7-O8
3	A	1494	6YF	F11-C10-S7-O9
3	A	1494	6YF	F12-C10-S7-O9
3	A	1494	6YF	F13-C10-S7-O9
3	A	1494	6YF	C18-C15-O16-C17
3	A	1494	6YF	C14-C15-O16-C17
3	A	1494	6YF	C4-C3-O2-C1
3	A	1494	6YF	C18-C3-O2-C1
3	A	1494	6YF	C5-N6-S7-C10
2	A	1493	HEM	C4C-C3C-CAC-CBC
2	A	1493	HEM	CAA-CBA-CGA-O1A
2	A	1493	HEM	CAA-CBA-CGA-O2A
2	A	1493	HEM	C2B-C3B-CAB-CBB
2	A	1493	HEM	CAD-CBD-CGD-O2D

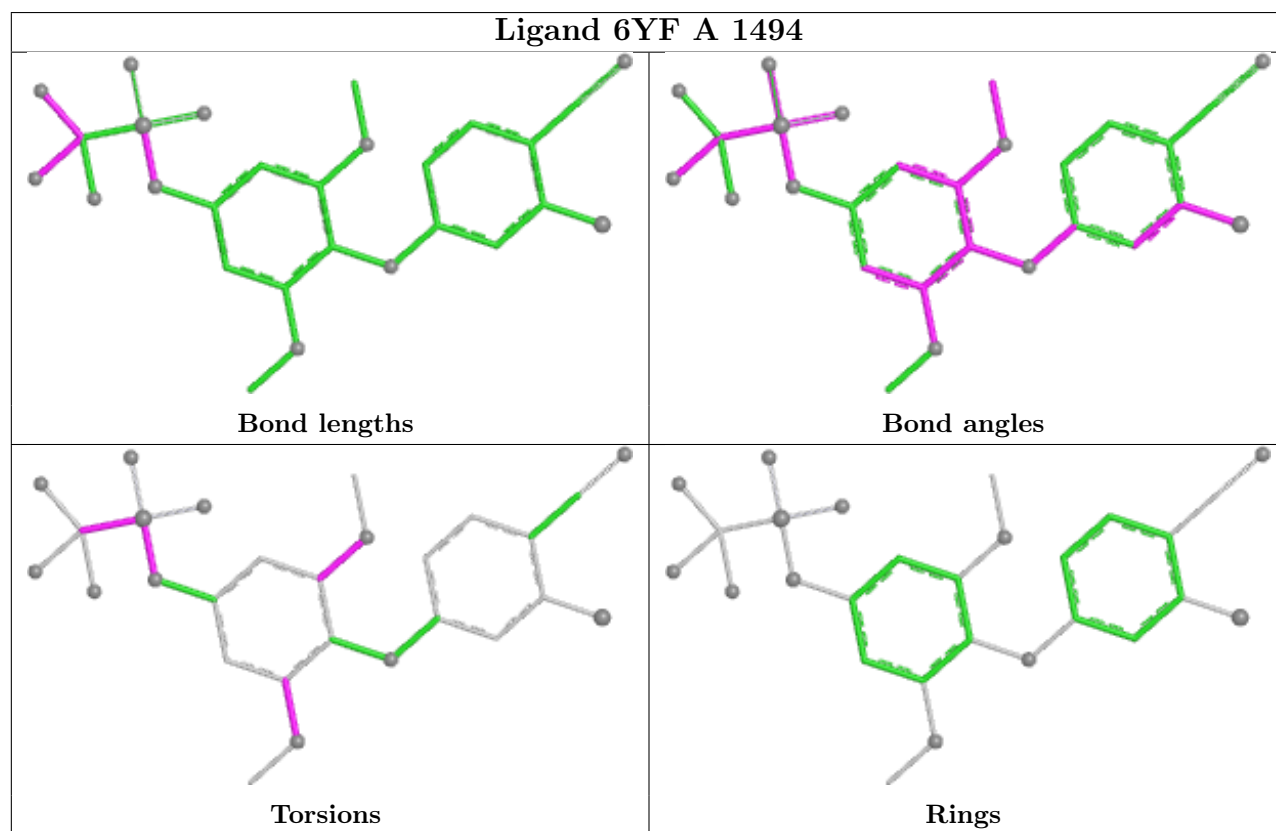
There are no ring outliers.

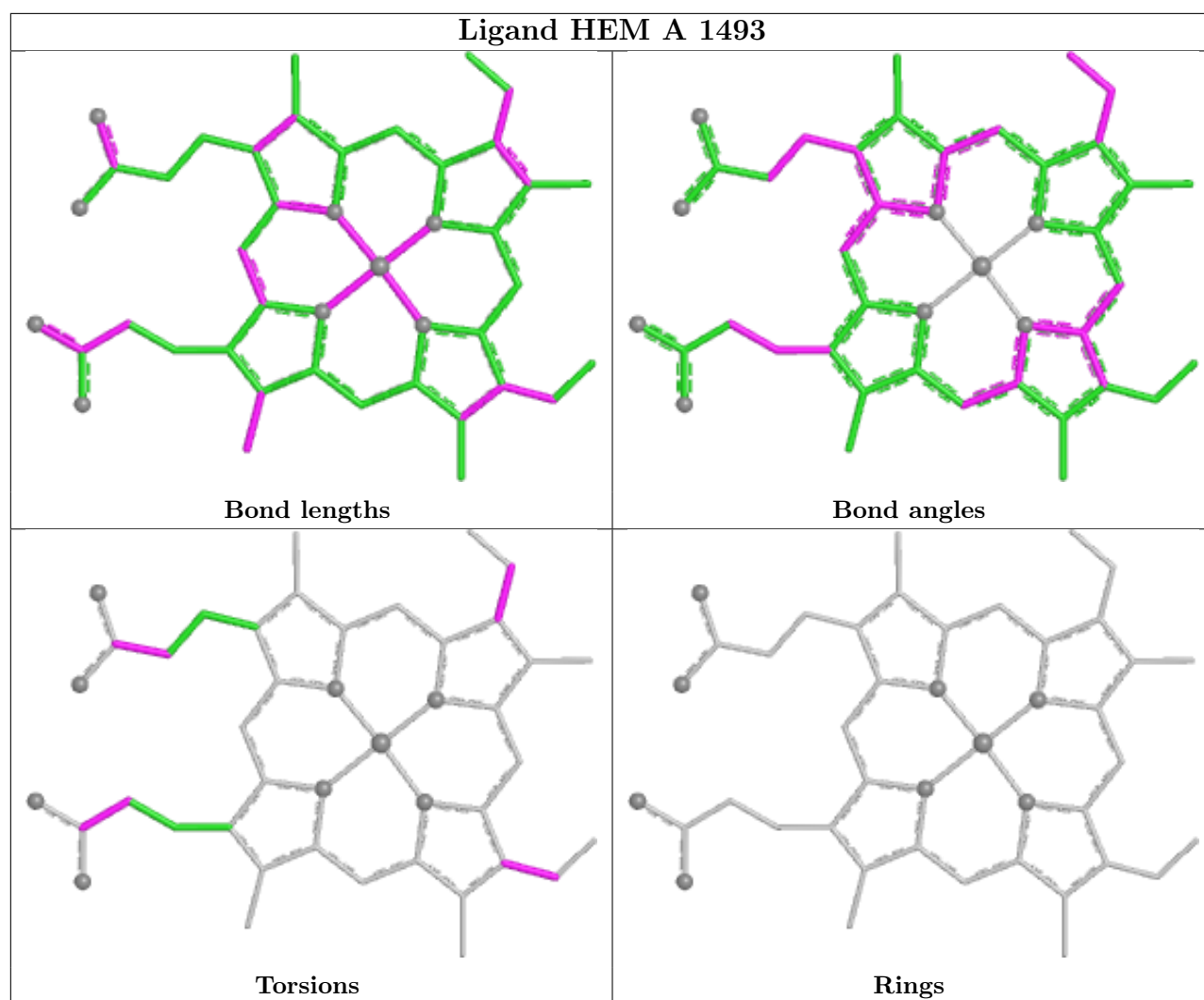
2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1494	6YF	1	0
2	A	1493	HEM	7	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	455/477 (95%)	0.22	17 (3%) 45 37	15, 39, 66, 74	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	492	HIS	4.4
1	A	100	PHE	3.8
1	A	491	HIS	3.5
1	A	490	ILE	3.4
1	A	99	ILE	3.1
1	A	106	ALA	3.0
1	A	225	TYR	2.8
1	A	150	ARG	2.8
1	A	37	PRO	2.7
1	A	228	GLY	2.6
1	A	185	LYS	2.6
1	A	36	LEU	2.6
1	A	110	PHE	2.5
1	A	184	HIS	2.2
1	A	476	PHE	2.1
1	A	103	ALA	2.1
1	A	466	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](#)

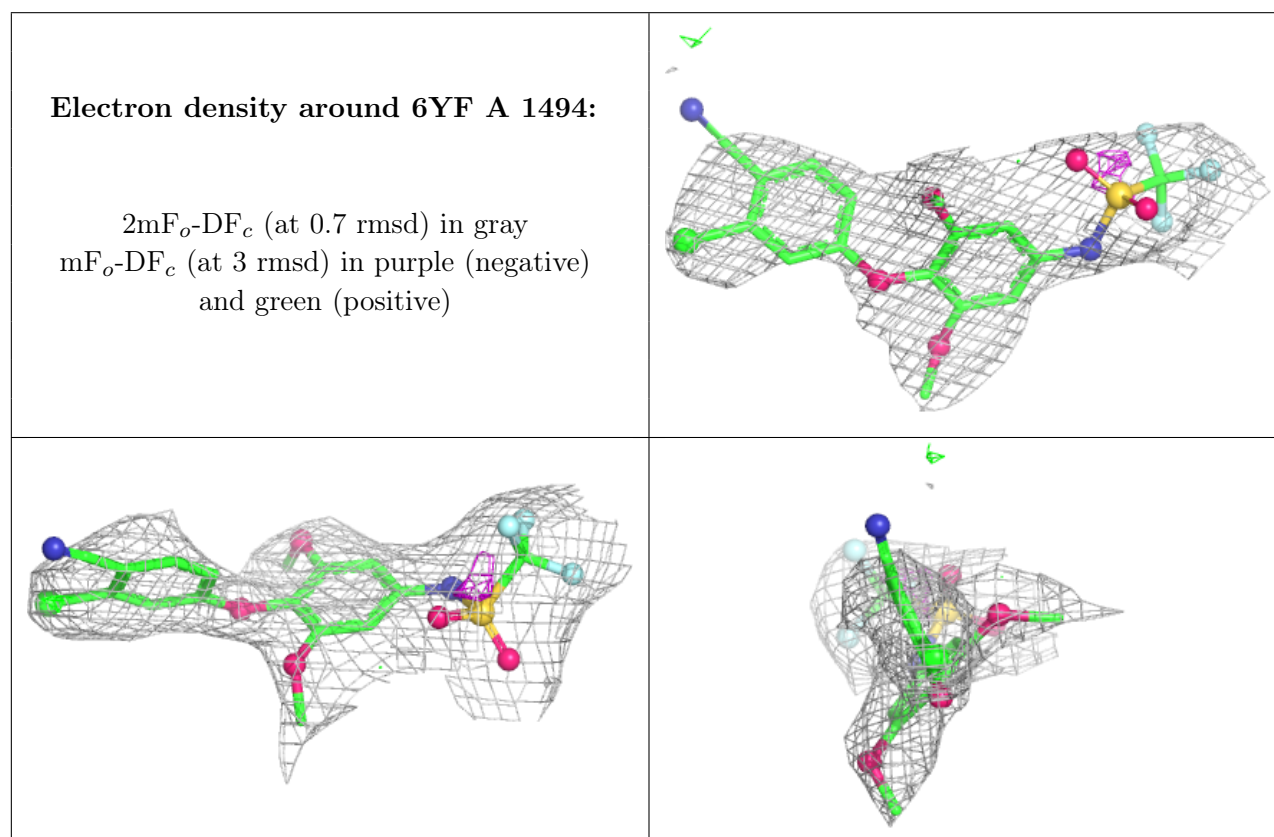
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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.

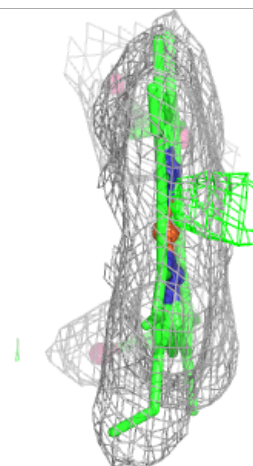
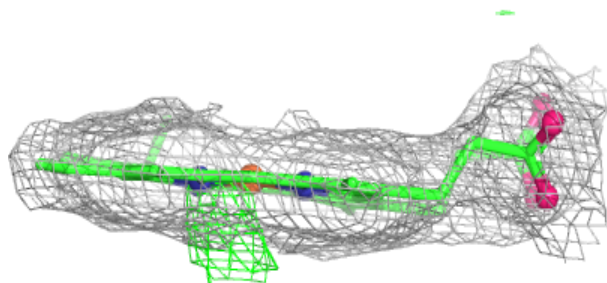
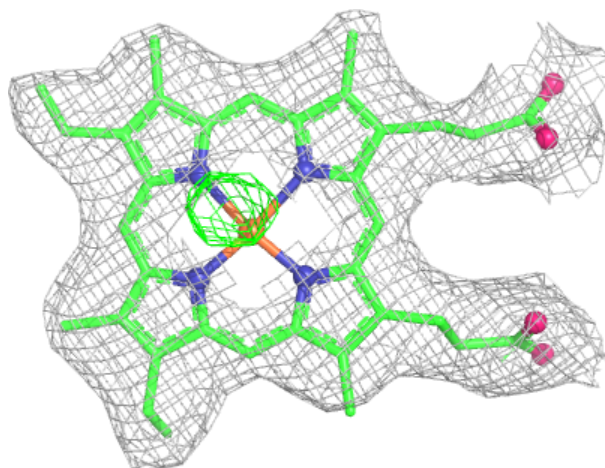
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	6YF	A	1494	28/28	0.78	0.14	52,57,67,68	0
2	HEM	A	1493	43/43	0.97	0.07	8,16,20,23	0

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 1493:

$2mF_o-DF_c$ (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.