



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 11:54 pm BST

PDB ID : 4CKA
Title : STEROL 14-ALPHA DEMETHYLASE (CYP51)FROM TRYPANOSOMA CRUZI IN COMPLEX WITH (S)-1-(4-fluorophenyl)-2-(1H-imidazol-1-yl)ethyl 4- isopropylphenylcarbamate (LFS)
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Deposited on : 2013-12-31
Resolution : 2.70 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

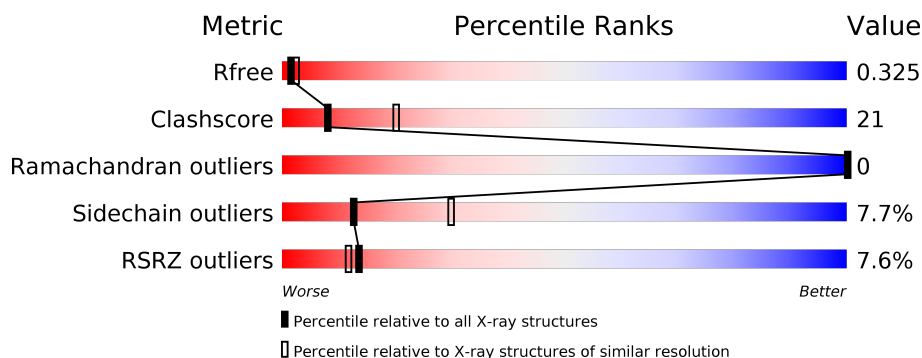
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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

Mol	Chain	Length	Quality of chain
1	A	460	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	LFS	A	1490	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3541 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.

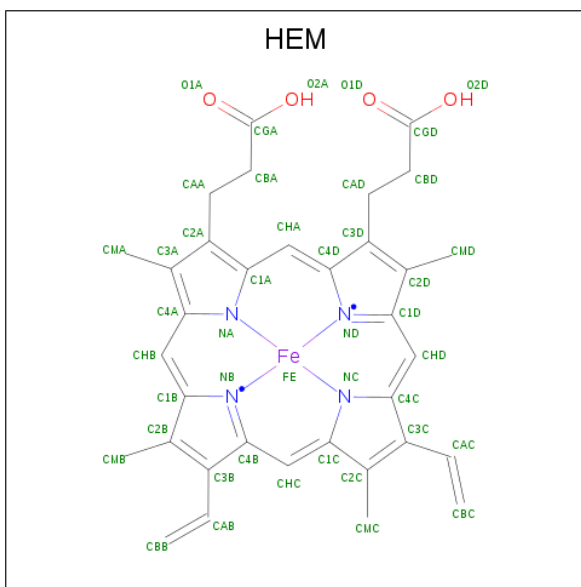
- Molecule 1 is a protein called STEROL 14-ALPHA DEMETHYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	434	Total	C	N	O	S	0	0	0
			3467	2215	607	617	28			

There are 10 discrepancies between the modelled and reference sequences:

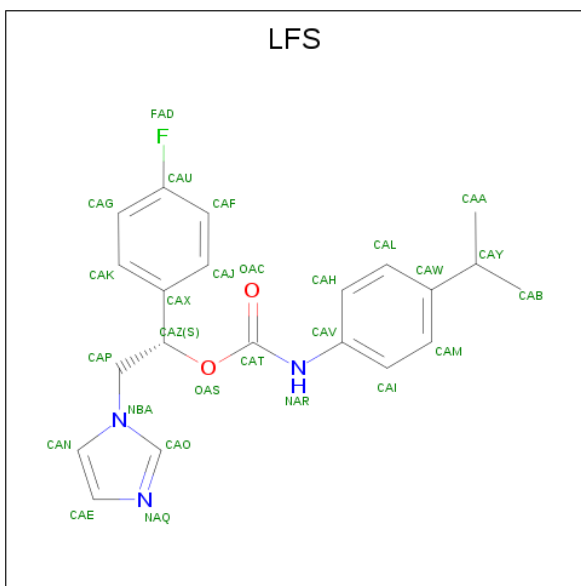
Chain	Residue	Modelled	Actual	Comment	Reference
A	28	ALA	-	expression tag	UNP Q7Z1V1
A	29	LYS	-	expression tag	UNP Q7Z1V1
A	30	LYS	-	expression tag	UNP Q7Z1V1
A	31	THR	-	expression tag	UNP Q7Z1V1
A	482	HIS	-	expression tag	UNP Q7Z1V1
A	483	HIS	-	expression tag	UNP Q7Z1V1
A	484	HIS	-	expression tag	UNP Q7Z1V1
A	485	HIS	-	expression tag	UNP Q7Z1V1
A	486	HIS	-	expression tag	UNP Q7Z1V1
A	487	HIS	-	expression tag	UNP Q7Z1V1

- 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	0	0
			43	34	1	4	4		

- Molecule 3 is (1S)-1-(4-fluorophenyl)-2-(1H-imidazol-1-yl)ethyl 4-isopropylphenylcarbamate (three-letter code: LFS) (formula: $C_{21}H_{22}FN_3O_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			27	21	1	3	2		

- Molecule 4 is water.

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



- Molecule 1: STEROL 14-ALPHA DEMETHYLASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	62.58Å 62.58Å 221.07Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.80 – 2.70 28.80 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.4 (28.80-2.70) 99.5 (28.80-2.70)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.20 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.263 , 0.284 0.259 , 0.325	Depositor DCC
R_{free} test set	732 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	79.3	Xtriage
Anisotropy	0.075	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 57.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.049 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3541	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.31% 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, LFS

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	0.58	0/3546	0.75	2/4787 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	143	VAL	C-N-CD	6.09	141.20	128.40
1	A	359	VAL	N-CA-C	-5.14	97.11	111.00

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	3467	0	3495	145	0
2	A	43	0	30	10	0
3	A	27	0	22	10	0
4	A	4	0	0	1	0
All	All	3541	0	3547	151	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 21.

All (151) 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:103:TYR:CD2	3:A:1490:LFS:HAF	1.92	1.03
1:A:136:ILE:H	1:A:136:ILE:HD13	1.19	1.02
1:A:183:LEU:HD12	1:A:289:MET:CE	2.04	0.86
1:A:352:ARG:NH1	1:A:353:ASP:OD2	2.12	0.82
1:A:281:VAL:O	1:A:285:ILE:HD12	1.80	0.82
1:A:33:PRO:CG	1:A:63:LEU:HD21	2.13	0.78
1:A:361:ARG:HH21	1:A:420:HIS:CD2	2.07	0.73
1:A:146:ILE:CG2	1:A:182:CYS:SG	2.76	0.72
1:A:424:GLY:HA3	2:A:1480:HEM:C2C	2.24	0.72
1:A:303:SER:O	1:A:307:LEU:HD12	1.90	0.72
1:A:103:TYR:CE2	3:A:1490:LFS:HAF	2.25	0.71
1:A:29:LYS:HE3	4:A:2001:HOH:O	1.90	0.71
1:A:295:THR:HB	2:A:1480:HEM:HAB	1.73	0.70
1:A:145:ALA:HB1	1:A:181:GLN:OE1	1.92	0.70
2:A:1480:HEM:HBC2	2:A:1480:HEM:CMC	2.21	0.69
1:A:225:GLN:HA	1:A:228:ARG:HB3	1.75	0.68
1:A:449:ARG:NH1	1:A:468:GLN:OE1	2.26	0.67
1:A:124:ARG:NH1	1:A:124:ARG:HG2	2.08	0.67
1:A:309:HIS:CE1	1:A:451:GLU:HA	2.29	0.67
1:A:418:GLY:O	1:A:421:LYS:HG3	1.96	0.66
1:A:424:GLY:HA3	2:A:1480:HEM:C3C	2.31	0.65
1:A:33:PRO:HG2	1:A:63:LEU:HD21	1.79	0.65
1:A:158:TRP:O	1:A:475:ARG:NH2	2.29	0.65
1:A:109:VAL:O	1:A:233:ARG:HD2	1.97	0.64
1:A:103:TYR:CD2	3:A:1490:LFS:CAF	2.75	0.64
1:A:146:ILE:HD11	1:A:434:THR:HB	1.80	0.64
1:A:130:LEU:HD23	3:A:1490:LFS:HAB	1.79	0.63
1:A:33:PRO:CB	1:A:63:LEU:HD21	2.28	0.63
1:A:102:VAL:HG13	1:A:213:VAL:HG13	1.80	0.63
1:A:183:LEU:HD12	1:A:289:MET:HE2	1.78	0.63
1:A:102:VAL:HG11	1:A:360:MET:HB2	1.80	0.63
1:A:358:MET:HE3	1:A:381:ALA:HB1	1.80	0.63
1:A:339:MET:HE2	1:A:433:LYS:HB3	1.80	0.63
1:A:107:THR:OG1	1:A:108:PRO:HD3	1.99	0.62
1:A:126:GLN:HB3	3:A:1490:LFS:HAAB	1.80	0.62
1:A:105:ILE:H	1:A:105:ILE:HD12	1.64	0.62
1:A:124:ARG:HH11	1:A:124:ARG:HG2	1.62	0.62
1:A:136:ILE:H	1:A:136:ILE:CD1	1.93	0.61
1:A:33:PRO:HB2	1:A:63:LEU:HD21	1.81	0.61
1:A:43:GLY:O	1:A:71:SER:O	2.18	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:SER:O	1:A:230:ARG:HB2	2.00	0.61
2:A:1480:HEM:C4D	3:A:1490:LFS:HAO	2.36	0.60
1:A:43:GLY:O	1:A:44:HIS:HB2	2.01	0.60
1:A:98:SER:HG	1:A:120:TYR:HH	1.44	0.59
1:A:63:LEU:HD22	1:A:68:PHE:HB3	1.84	0.59
2:A:1480:HEM:HBC2	2:A:1480:HEM:HMC2	1.84	0.59
1:A:144:PRO:HG2	1:A:145:ALA:H	1.67	0.59
1:A:127:LEU:HD23	3:A:1490:LFS:HABB	1.85	0.58
1:A:124:ARG:CG	1:A:124:ARG:HH11	2.15	0.58
1:A:105:ILE:HG22	1:A:208:LEU:HD22	1.85	0.58
1:A:107:THR:HG22	1:A:117:ALA:HB2	1.84	0.58
1:A:144:PRO:HG2	1:A:145:ALA:N	2.19	0.58
1:A:76:ARG:NH2	1:A:378:ASP:OD2	2.37	0.57
1:A:43:GLY:HA3	1:A:73:GLY:H	1.68	0.57
1:A:358:MET:CE	1:A:381:ALA:HB1	2.35	0.57
1:A:72:ILE:O	1:A:75:GLN:HB3	2.05	0.56
1:A:132:GLU:HA	1:A:135:THR:CG2	2.36	0.55
1:A:358:MET:SD	1:A:383:SER:HB2	2.46	0.55
2:A:1480:HEM:CBC	2:A:1480:HEM:HMC2	2.36	0.55
1:A:357:LEU:HD21	1:A:462:VAL:HG21	1.88	0.55
1:A:158:TRP:HE3	1:A:443:TYR:CE2	2.24	0.54
1:A:147:GLN:OE1	1:A:330:LEU:HB2	2.07	0.54
1:A:165:ILE:HG23	1:A:170:ASP:OD2	2.07	0.54
1:A:146:ILE:HG23	1:A:182:CYS:SG	2.48	0.54
1:A:309:HIS:HE1	1:A:451:GLU:HA	1.70	0.54
1:A:272:ASP:OD1	1:A:274:THR:OG1	2.20	0.54
1:A:146:ILE:HG22	1:A:182:CYS:SG	2.47	0.53
1:A:442:GLU:OE2	1:A:443:TYR:CE2	2.61	0.53
1:A:339:MET:N	1:A:340:PRO:CD	2.72	0.53
1:A:281:VAL:O	1:A:285:ILE:CD1	2.55	0.53
1:A:76:ARG:HH22	1:A:375:PRO:HG2	1.74	0.52
1:A:98:SER:OG	1:A:120:TYR:OH	2.21	0.52
1:A:351:ARG:HD2	1:A:388:HIS:HB3	1.91	0.52
2:A:1480:HEM:C1A	3:A:1490:LFS:HAO	2.44	0.51
1:A:389:HIS:CD2	1:A:397:PRO:HB2	2.46	0.51
1:A:58:ARG:HA	1:A:61:ARG:HH12	1.76	0.51
1:A:122:ARG:NH2	1:A:280:GLU:OE2	2.44	0.51
1:A:31:THR:HG23	1:A:32:PRO:HD2	1.91	0.51
1:A:158:TRP:CE3	1:A:443:TYR:CE2	3.00	0.50
1:A:295:THR:HB	2:A:1480:HEM:CAB	2.41	0.49
1:A:189:ARG:O	1:A:192:LEU:N	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:PRO:O	1:A:228:ARG:N	2.39	0.49
1:A:245:VAL:O	1:A:249:LYS:HG2	2.12	0.49
1:A:160:GLU:O	1:A:475:ARG:NH1	2.46	0.49
1:A:100:ARG:HH22	1:A:119:PRO:HA	1.78	0.49
1:A:361:ARG:NH2	1:A:420:HIS:CD2	2.77	0.49
1:A:357:LEU:CD2	1:A:462:VAL:HG21	2.43	0.48
1:A:63:LEU:HD23	1:A:65:SER:HB3	1.95	0.48
1:A:90:PHE:O	1:A:419:VAL:HG22	2.12	0.48
1:A:300:THR:O	1:A:304:MET:HG3	2.14	0.48
1:A:43:GLY:O	1:A:72:ILE:HA	2.14	0.47
1:A:281:VAL:HG12	1:A:285:ILE:HD11	1.96	0.47
1:A:309:HIS:HD2	1:A:311:LYS:CB	2.27	0.47
1:A:395:PRO:O	1:A:404:ARG:NH2	2.47	0.47
1:A:452:VAL:HG13	1:A:453:PRO:HD2	1.96	0.47
1:A:91:SER:N	1:A:92:PRO:HD2	2.30	0.47
1:A:147:GLN:NE2	1:A:330:LEU:HD12	2.30	0.46
1:A:104:THR:O	1:A:107:THR:HG23	2.15	0.46
1:A:130:LEU:CD2	3:A:1490:LFS:HAB	2.45	0.46
1:A:105:ILE:HD11	1:A:213:VAL:HA	1.96	0.46
1:A:82:ASP:OD2	1:A:85:GLU:HG2	2.16	0.45
1:A:58:ARG:HA	1:A:61:ARG:NH1	2.30	0.45
1:A:161:ASP:OD1	1:A:161:ASP:N	2.49	0.45
1:A:438:THR:O	1:A:441:ARG:O	2.34	0.45
1:A:51:ASN:HB3	1:A:54:GLU:HB2	1.99	0.45
1:A:236:LEU:HA	1:A:239:ILE:HD12	1.98	0.45
1:A:366:GLU:HG3	1:A:374:VAL:O	2.15	0.45
1:A:448:LEU:CD2	1:A:470:LEU:HB3	2.47	0.45
1:A:143:VAL:HG21	1:A:332:TYR:HA	1.98	0.45
1:A:32:PRO:CD	1:A:373:VAL:O	2.65	0.44
1:A:260:ASP:OD1	1:A:263:GLY:N	2.36	0.44
1:A:332:TYR:CE2	1:A:336:MET:HG3	2.52	0.44
1:A:69:THR:HA	1:A:78:THR:HA	1.99	0.44
1:A:292:GLY:O	1:A:296:SER:OG	2.20	0.44
1:A:173:ALA:HA	1:A:176:ILE:HG22	2.00	0.44
1:A:168:LEU:HD23	1:A:466:LEU:HA	1.99	0.44
1:A:98:SER:O	1:A:362:MET:N	2.48	0.44
1:A:144:PRO:CG	1:A:145:ALA:N	2.81	0.44
1:A:442:GLU:OE2	1:A:443:TYR:CZ	2.71	0.44
1:A:412:PHE:C	1:A:412:PHE:CD1	2.91	0.43
1:A:235:GLU:O	1:A:239:ILE:HG13	2.19	0.43
2:A:1480:HEM:CHA	3:A:1490:LFS:HAO	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:LYS:HG3	1:A:66:GLY:HA2	2.01	0.43
1:A:249:LYS:O	1:A:249:LYS:HG3	2.19	0.43
1:A:163:GLY:HA3	1:A:473:TYR:CD2	2.53	0.43
1:A:180:CYS:O	1:A:184:PHE:HB2	2.19	0.43
1:A:260:ASP:C	1:A:260:ASP:OD1	2.58	0.43
1:A:175:ILE:HG23	1:A:431:GLN:HE22	1.84	0.43
1:A:395:PRO:O	1:A:396:ASN:HB2	2.18	0.43
1:A:42:LEU:HA	1:A:42:LEU:HD23	1.86	0.42
1:A:98:SER:OG	1:A:101:GLU:OE1	2.36	0.42
1:A:405:ASP:O	1:A:406:GLU:C	2.56	0.42
1:A:432:VAL:O	1:A:436:LEU:HG	2.20	0.42
1:A:32:PRO:HD3	1:A:373:VAL:O	2.19	0.41
1:A:400:TRP:CE3	1:A:402:PRO:HD3	2.54	0.41
1:A:150:VAL:CG1	1:A:438:THR:HG21	2.50	0.41
1:A:146:ILE:HD11	1:A:434:THR:CB	2.47	0.41
1:A:146:ILE:HD11	1:A:434:THR:CG2	2.50	0.41
1:A:100:ARG:HE	1:A:100:ARG:HB2	1.72	0.41
1:A:39:VAL:HG13	1:A:42:LEU:HB2	2.02	0.41
1:A:89:PHE:CD2	1:A:382:CYS:HB2	2.55	0.41
1:A:163:GLY:O	1:A:472:LYS:HG2	2.21	0.41
1:A:105:ILE:H	1:A:105:ILE:CD1	2.33	0.41
1:A:339:MET:N	1:A:340:PRO:HD3	2.35	0.41
1:A:76:ARG:NH2	1:A:375:PRO:HG2	2.35	0.41
1:A:441:ARG:O	1:A:442:GLU:HB3	2.21	0.41
1:A:101:GLU:OE1	1:A:101:GLU:N	2.49	0.41
1:A:348:GLU:HG3	1:A:400:TRP:CD1	2.56	0.40
1:A:107:THR:N	1:A:108:PRO:CD	2.83	0.40
1:A:144:PRO:CG	1:A:145:ALA:H	2.34	0.40
1:A:150:VAL:HG11	1:A:438:THR:HG21	2.03	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	428/460 (93%)	390 (91%)	38 (9%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	378/402 (94%)	349 (92%)	29 (8%)	13	30

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	LYS
1	A	47	GLN
1	A	63	LEU
1	A	87	SER
1	A	124	ARG
1	A	136	ILE
1	A	146	ILE
1	A	161	ASP
1	A	206	SER
1	A	209	ILE
1	A	214	PHE
1	A	223	LEU
1	A	229	CYS
1	A	233	ARG
1	A	269	VAL
1	A	274	THR
1	A	278	LEU
1	A	293	GLN
1	A	307	LEU
1	A	339	MET
1	A	371	SER
1	A	376	LYS
1	A	441	ARG

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Mol	Chain	Res	Type
1	A	442	GLU
1	A	446	GLN
1	A	460	MET
1	A	474	THR
1	A	476	LYS
1	A	477	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	237	GLN
1	A	309	HIS
1	A	420	HIS
1	A	446	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 carbohydrates 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
2	HEM	A	1480	1,3	27,50,50	1.89	6 (22%)	17,82,82	1.48	3 (17%)
3	LFS	A	1490	2	27,29,29	3.67	8 (29%)	35,39,39	1.85	9 (25%)

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
2	HEM	A	1480	1,3	-	0/6/54/54	-
3	LFS	A	1490	2	-	3/20/20/20	0/3/3/3

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1490	LFS	FAD-CAU	-13.18	1.04	1.36
3	A	1490	LFS	CAX-CAZ	-7.67	1.38	1.51
2	A	1480	HEM	C3B-C2B	-5.43	1.32	1.40
3	A	1490	LFS	CAW-CAY	-5.35	1.39	1.53
3	A	1490	LFS	OAS-CAZ	-5.02	1.39	1.46
3	A	1490	LFS	CAV-NAR	-4.92	1.31	1.41
3	A	1490	LFS	CAN-NBA	-4.50	1.29	1.37
2	A	1480	HEM	C3C-C2C	-3.94	1.34	1.40
3	A	1490	LFS	CAP-NBA	-3.43	1.44	1.48
2	A	1480	HEM	C4A-CHB	-2.44	1.34	1.41
2	A	1480	HEM	CAA-C2A	-2.31	1.48	1.52
2	A	1480	HEM	C3D-C2D	-2.29	1.30	1.37
2	A	1480	HEM	CAD-C3D	-2.23	1.48	1.52
3	A	1490	LFS	OAS-CAT	-2.02	1.31	1.35

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1490	LFS	OAS-CAT-NAR	5.02	115.44	109.07
3	A	1490	LFS	CAP-NBA-CAO	4.18	133.99	125.76
3	A	1490	LFS	CAP-NBA-CAN	-3.78	118.13	125.92
2	A	1480	HEM	CAD-CBD-CGD	-3.33	107.08	112.67
2	A	1480	HEM	CBA-CAA-C2A	-3.30	106.39	112.49
3	A	1490	LFS	CAK-CAG-CAU	3.06	121.53	118.36
3	A	1490	LFS	CAP-CAZ-CAX	-3.00	106.56	111.46
3	A	1490	LFS	OAS-CAT-OAC	-2.63	120.57	124.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1480	HEM	C1D-C2D-C3D	-2.53	105.24	107.00
3	A	1490	LFS	CAF-CAJ-CAX	2.46	123.68	121.20
3	A	1490	LFS	CAG-CAK-CAX	-2.24	118.94	121.20
3	A	1490	LFS	CAV-NAR-CAT	-2.22	122.70	126.36

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1490	LFS	CAL-CAW-CAY-CAA
3	A	1490	LFS	CAP-CAZ-OAS-CAT
3	A	1490	LFS	CAM-CAW-CAY-CAA

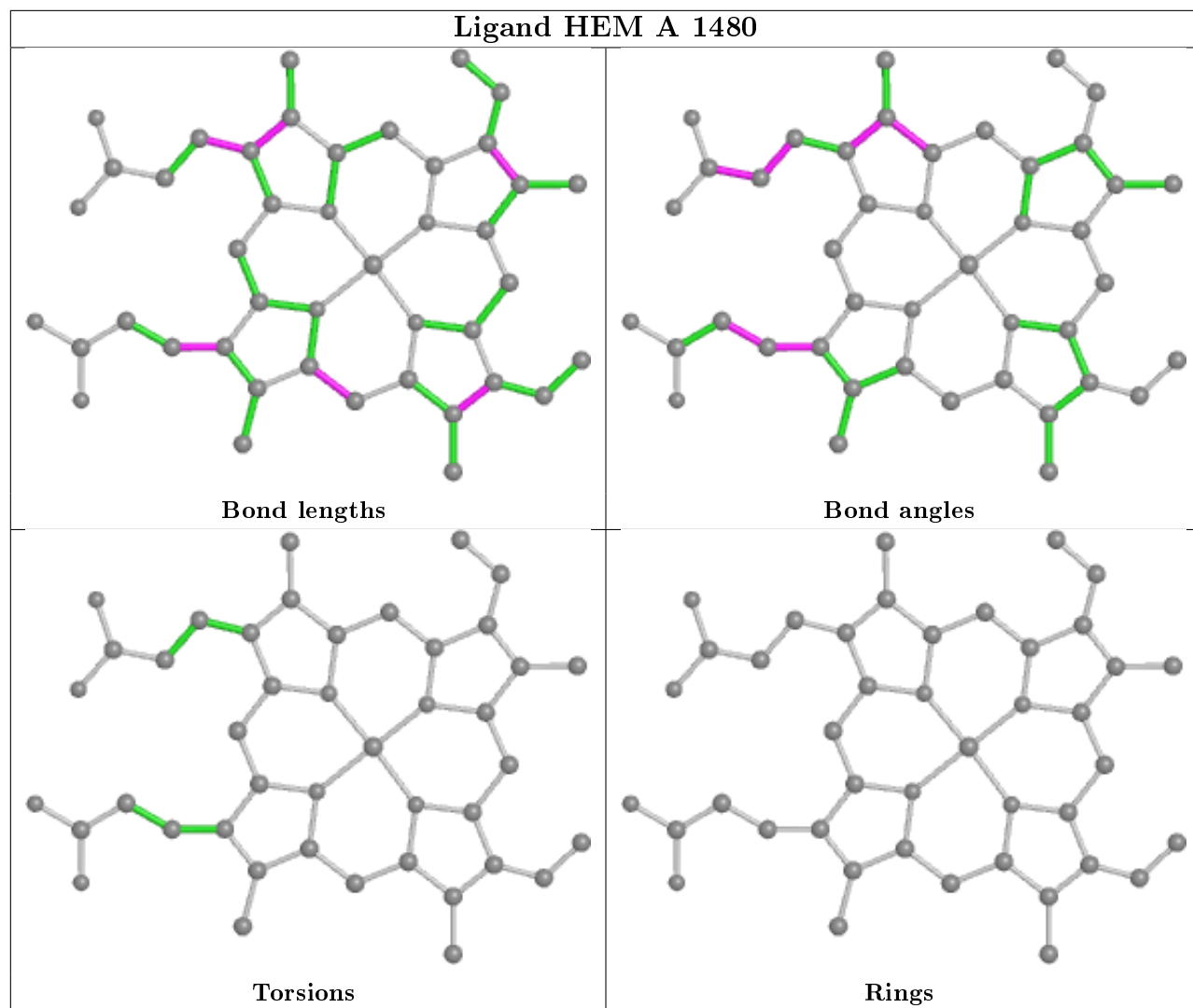
There are no ring outliers.

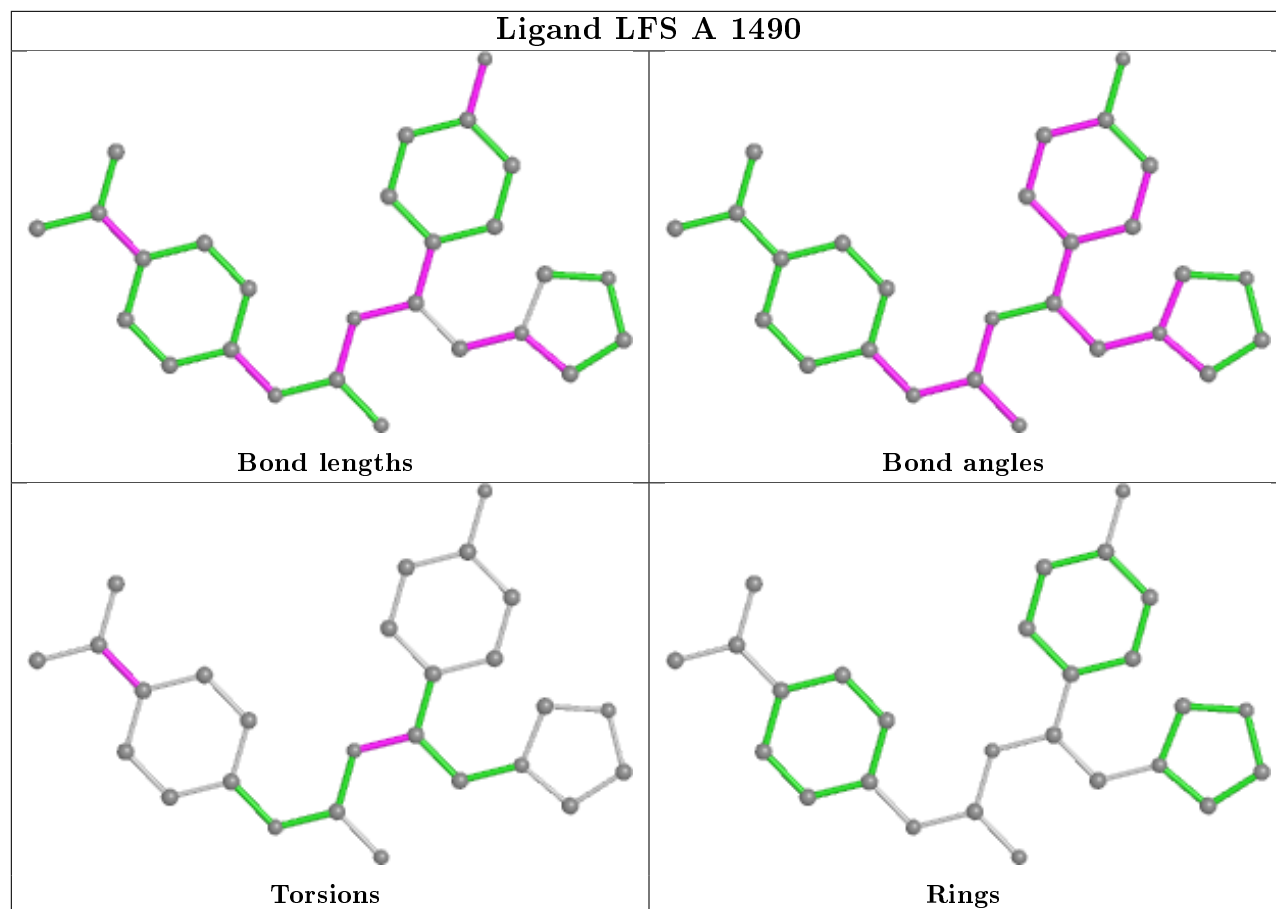
2 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1480	HEM	10	0
3	A	1490	LFS	10	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.

Ligand HEM A 1480





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	434/460 (94%)	0.46	33 (7%) 13 12	47, 89, 134, 292	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	223	LEU	24.2
1	A	224	PRO	10.9
1	A	225	GLN	7.0
1	A	248	GLU	4.0
1	A	214	PHE	3.8
1	A	160	GLU	3.8
1	A	38	THR	3.8
1	A	159	LYS	3.7
1	A	275	ARG	3.7
1	A	271	ARG	3.5
1	A	365	ALA	3.4
1	A	73	GLY	3.2
1	A	41	PHE	3.1
1	A	285	ILE	3.0
1	A	247	ARG	3.0
1	A	112	GLU	2.9
1	A	44	HIS	2.9
1	A	270	TYR	2.8
1	A	288	ALA	2.8
1	A	362	MET	2.7
1	A	475	ARG	2.6
1	A	195	ARG	2.5
1	A	274	THR	2.5
1	A	207	SER	2.5
1	A	230	ARG	2.4
1	A	435	ILE	2.3
1	A	209	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	286	VAL	2.2
1	A	269	VAL	2.2
1	A	281	VAL	2.1
1	A	215	MET	2.1
1	A	291	ALA	2.1
1	A	289	MET	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

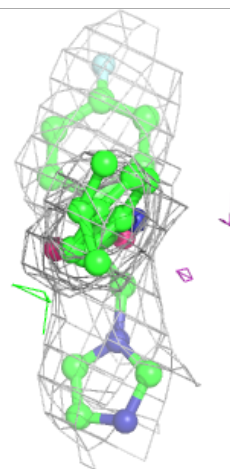
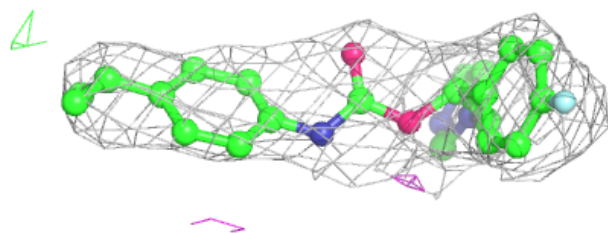
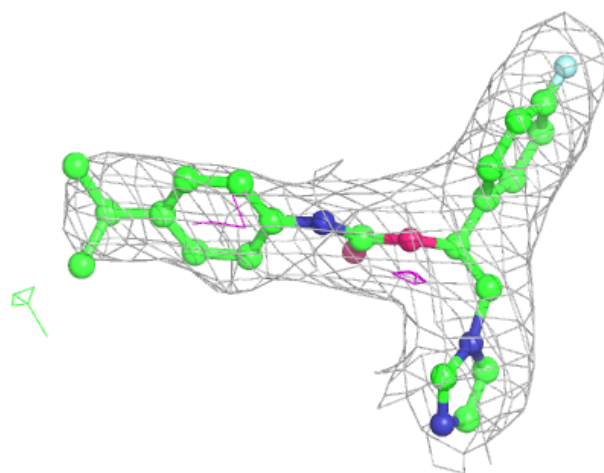
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	LFS	A	1490	27/27	0.95	0.22	50,74,84,89	0
2	HEM	A	1480	43/43	0.97	0.18	46,54,64,75	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.

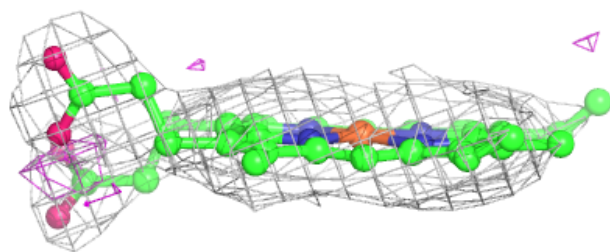
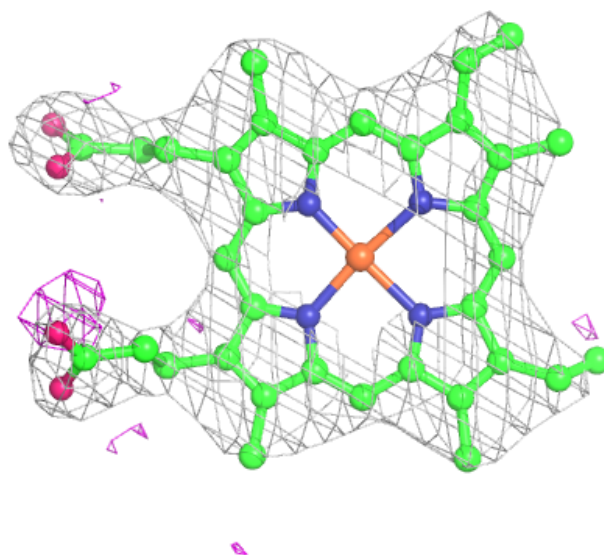
Electron density around LFS A 1490:

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



Electron density around HEM A 1480:

$2mF_o-DF_c$ (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.