



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 11:46 PM UTC

PDB ID : 3ZG2 / pdb_00003zg2
Title : Sterol 14 alpha-demethylase (CYP51) from Trypanosoma cruzi in complex with the pyridine inhibitor (S)-2-(4-chlorophenyl)-2-(pyridin-3-yl)-1-(4-(4-(trifluoromethyl)phenyl)piperazin-1-yl)ethanone (EPL-BS1246,UDO)
Authors : Hargrove, T.Y.; Wawrzak, Z.; Keenan, M.; Chatelain, E.; Lepesheva, G.I.
Deposited on : 2012-12-14
Resolution : 2.80 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

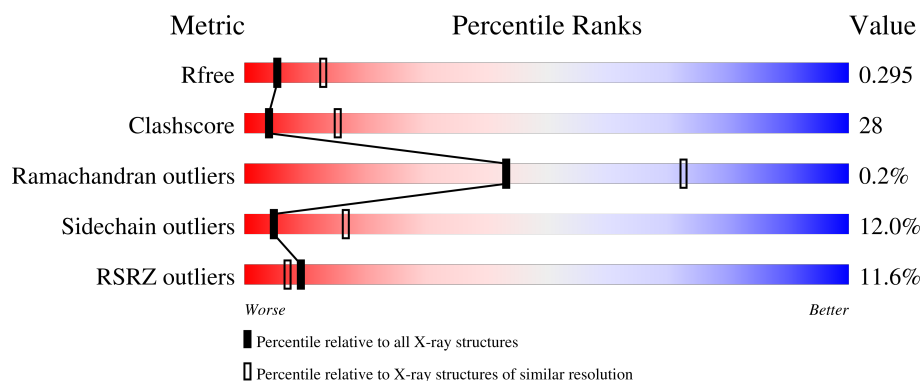
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	460	11% 50% 40% 7%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3677 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
			Total	C	N	O	S			
1	A	449	3589	2291	630	640	28	0	0	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	28	ALA	ARG	engineered mutation	UNP Q7Z1V1
A	29	LYS	PRO	engineered mutation	UNP Q7Z1V1
A	30	LYS	THR	engineered mutation	UNP Q7Z1V1
A	31	THR	ASP	engineered mutation	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 (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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- The chemical structure is a complex molecule with the following features:
- Top Group:** A central carbon atom is bonded to two fluorine atoms (FAB, FAD) and two other fluorine atoms (FAC, FAF) via green bonds. This central carbon is also bonded to a green oxygen atom (OFB).
 - Phenyl Ring 1:** A benzene ring with a green oxygen atom (OEB) at the top position. Other positions are labeled with green atoms: CAO, CAP, CAN, CAM, and CBA.
 - Piperazine Ring:** A six-membered ring with two nitrogen atoms (NBE, NED) in blue. The ring is substituted with green atoms: CAT, CAU, CAS, and CAV.
 - Pyridine Ring:** A six-membered ring with a nitrogen atom (NRV) in blue. It is substituted with green atoms: CAG, CAZ, CAN, and CAX.
 - Carboxylic Acid Group:** A carboxylic acid group (COA) is attached to the piperazine ring via a green carbon atom (CAW). The oxygen atom is red (ORA).
 - Phenyl Ring 2:** A benzene ring with a green oxygen atom (OEB) at the top position. Other positions are labeled with green atoms: CAY, CAL, CLJ, CLK, and CLA.
 - Chlorine Atom:** A green chlorine atom (CLA) is attached to the bottom of the second phenyl ring.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	Cl	F	N	O	0	0
			32	24	1	3	3	1		

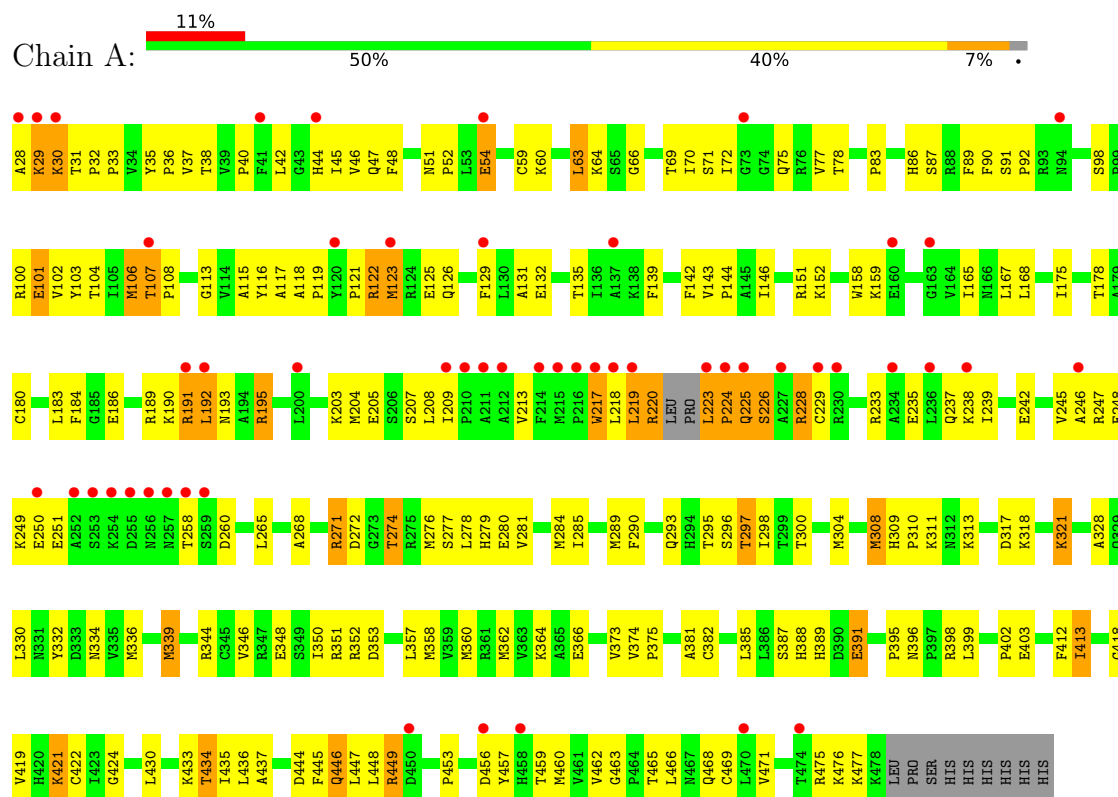
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	13	Total 13	O 13	0	0

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: 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	63.70Å 63.70Å 223.26Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	27.66 – 2.80 27.66 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.2 (27.66-2.80) 94.2 (27.66-2.80)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.92 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.268 , 0.282 0.265 , 0.295	Depositor DCC
R_{free} test set	672 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	40.6	Xtriage
Anisotropy	0.202	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 50.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.054 for -h,-k,l	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	3677	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.79% 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: UDO, HEM

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.33	0/3672	0.91	7/4960 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	223	LEU	CA-C-N	-7.63	111.65	120.12
1	A	223	LEU	C-N-CA	-7.63	111.65	120.12
1	A	224	PRO	N-CA-C	-6.96	105.85	114.20
1	A	29	LYS	N-CA-C	-6.79	104.05	112.47
1	A	226	SER	N-CA-C	-6.11	104.62	111.28
1	A	117	ALA	N-CA-C	-5.21	105.61	111.28
1	A	228	ARG	N-CA-C	-5.00	108.21	114.56

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3589	0	3618	202	0
2	A	43	0	30	4	0
3	A	32	0	21	5	0
4	A	13	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3677	0	3669	203	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 28.

All (203) 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:33:PRO:HB2	1:A:63:LEU:CD1	1.69	1.20
1:A:293:GLN:O	1:A:297:THR:HG22	1.48	1.11
1:A:223:LEU:HG	1:A:224:PRO:CD	1.85	1.05
1:A:223:LEU:HG	1:A:224:PRO:HD2	1.03	1.03
1:A:191:ARG:HH22	1:A:242:GLU:HB3	1.22	1.01
1:A:237:GLN:OE1	1:A:279:HIS:HA	1.62	1.00
1:A:271:ARG:HG3	1:A:271:ARG:HH11	1.23	1.00
1:A:33:PRO:HB2	1:A:63:LEU:HD11	1.40	1.00
1:A:30:LYS:HE2	1:A:31:THR:HG23	1.43	0.97
1:A:308:MET:HE2	1:A:445:PHE:HB3	1.48	0.95
1:A:60:LYS:HG3	1:A:66:GLY:HA2	1.48	0.95
1:A:33:PRO:HB2	1:A:63:LEU:HD12	1.47	0.94
1:A:33:PRO:CB	1:A:63:LEU:CD1	2.50	0.90
1:A:33:PRO:CB	1:A:63:LEU:HD12	2.02	0.89
1:A:100:ARG:HH22	1:A:119:PRO:HA	1.35	0.89
1:A:271:ARG:HG3	1:A:271:ARG:NH1	1.87	0.87
1:A:277:SER:O	1:A:281:VAL:HG23	1.75	0.87
1:A:308:MET:HE2	1:A:445:PHE:CB	2.05	0.86
1:A:203:LYS:HG2	1:A:228:ARG:HD3	1.58	0.85
1:A:223:LEU:CG	1:A:224:PRO:HD2	1.99	0.84
1:A:309:HIS:HD2	1:A:311:LYS:H	1.27	0.83
1:A:193:ASN:HD21	1:A:195:ARG:HD2	1.44	0.82
1:A:102:VAL:O	1:A:213:VAL:CG2	2.29	0.80
1:A:51:ASN:HD22	1:A:54:GLU:H	1.30	0.79
1:A:33:PRO:CG	1:A:63:LEU:HD12	2.11	0.79
1:A:235:GLU:O	1:A:239:ILE:HG13	1.84	0.78
1:A:293:GLN:O	1:A:297:THR:CG2	2.29	0.78
1:A:225:GLN:OE1	1:A:228:ARG:HB3	1.84	0.78
1:A:102:VAL:O	1:A:213:VAL:HG23	1.85	0.76
1:A:203:LYS:CG	1:A:228:ARG:HD3	2.15	0.75
1:A:246:ALA:C	1:A:250:GLU:OE2	2.29	0.75
1:A:116:TYR:HD1	1:A:123:MET:HE1	1.51	0.75
1:A:300:THR:O	1:A:304:MET:HG3	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:ARG:NH1	1:A:328:ALA:O	2.22	0.72
1:A:237:GLN:OE1	1:A:279:HIS:CA	2.37	0.72
1:A:360:MET:HE1	3:A:1490:UDO:FAC	1.80	0.71
1:A:126:GLN:HA	1:A:129:PHE:CD2	2.25	0.71
1:A:191:ARG:HH22	1:A:242:GLU:CB	2.01	0.71
1:A:352:ARG:NH1	1:A:353:ASP:OD2	2.23	0.70
1:A:205:GLU:O	1:A:208:LEU:HG	1.91	0.70
1:A:388:HIS:HE1	1:A:413:ILE:H	1.40	0.70
1:A:271:ARG:HH11	1:A:271:ARG:CG	2.04	0.69
1:A:90:PHE:O	1:A:419:VAL:HG23	1.93	0.69
1:A:122:ARG:HD3	1:A:126:GLN:NE2	2.08	0.69
1:A:281:VAL:O	1:A:285:ILE:HD12	1.93	0.69
1:A:360:MET:CE	3:A:1490:UDO:FAC	2.30	0.69
1:A:389:HIS:HE1	1:A:398:ARG:HH11	1.39	0.69
1:A:121:PRO:O	1:A:125:GLU:HG3	1.91	0.69
1:A:446:GLN:NE2	1:A:448:LEU:HD23	2.07	0.69
1:A:223:LEU:C	1:A:225:GLN:H	1.99	0.67
1:A:219:LEU:HB2	4:A:2003:HOH:O	1.93	0.67
1:A:100:ARG:HH22	1:A:119:PRO:CA	2.08	0.67
1:A:228:ARG:HG2	1:A:228:ARG:O	1.95	0.66
1:A:106:MET:HE2	1:A:290:PHE:CE2	2.31	0.66
1:A:102:VAL:O	1:A:213:VAL:HG21	1.96	0.65
1:A:217:TRP:CE3	1:A:217:TRP:HA	2.31	0.65
1:A:385:LEU:O	1:A:389:HIS:HD2	1.80	0.64
1:A:309:HIS:CD2	1:A:310:PRO:HD2	2.32	0.63
1:A:102:VAL:HG11	1:A:360:MET:HB2	1.80	0.63
1:A:203:LYS:HG2	1:A:228:ARG:CD	2.27	0.63
1:A:143:VAL:HG21	1:A:332:TYR:HA	1.80	0.62
1:A:63:LEU:O	1:A:64:LYS:HB2	2.00	0.62
1:A:389:HIS:CE1	1:A:398:ARG:HH11	2.18	0.61
1:A:193:ASN:ND2	1:A:195:ARG:HD2	2.13	0.61
1:A:204:MET:SD	1:A:233:ARG:HG2	2.40	0.61
1:A:446:GLN:HE22	1:A:448:LEU:CD2	2.13	0.61
1:A:35:TYR:HD2	1:A:44:HIS:CE1	2.17	0.61
1:A:444:ASP:OD2	1:A:476:LYS:HE3	2.00	0.60
1:A:72:ILE:O	1:A:75:GLN:HB3	2.01	0.60
1:A:465:THR:HG21	1:A:468:GLN:HE21	1.67	0.60
1:A:446:GLN:NE2	1:A:448:LEU:CD2	2.65	0.59
1:A:35:TYR:CD2	1:A:44:HIS:CE1	2.91	0.59
1:A:100:ARG:NH2	1:A:119:PRO:HA	2.13	0.59
1:A:118:ALA:HB1	1:A:122:ARG:CD	2.33	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:309:HIS:NE2	1:A:311:LYS:HE3	2.17	0.59
1:A:385:LEU:O	1:A:389:HIS:CD2	2.56	0.58
1:A:32:PRO:HD3	1:A:373:VAL:O	2.04	0.58
1:A:192:LEU:HD12	1:A:193:ASN:N	2.18	0.58
1:A:298:ILE:HG12	1:A:463:GLY:HA3	1.85	0.58
1:A:51:ASN:HB3	1:A:54:GLU:HB2	1.86	0.58
1:A:339:MET:HE1	1:A:433:LYS:O	2.03	0.58
1:A:465:THR:CB	1:A:468:GLN:HE21	2.15	0.58
1:A:70:ILE:HG13	1:A:77:VAL:HB	1.86	0.57
1:A:391:GLU:O	1:A:395:PRO:HD3	2.04	0.57
1:A:223:LEU:C	1:A:225:GLN:N	2.55	0.57
1:A:205:GLU:O	1:A:207:SER:N	2.38	0.57
1:A:245:VAL:HG12	1:A:249:LYS:HE2	1.86	0.57
1:A:422:CYS:HA	2:A:1480:HEM:CHA	2.34	0.57
1:A:346:VAL:O	1:A:350:ILE:HG13	2.05	0.56
1:A:101:GLU:HG3	1:A:362:MET:CB	2.35	0.56
1:A:104:THR:O	1:A:107:THR:HG23	2.05	0.56
1:A:456:ASP:O	1:A:462:VAL:HG13	2.05	0.56
1:A:246:ALA:HB1	1:A:250:GLU:OE2	2.06	0.56
1:A:277:SER:OG	1:A:280:GLU:HG3	2.06	0.56
1:A:387:SER:C	1:A:389:HIS:H	2.14	0.55
1:A:113:GLY:O	1:A:118:ALA:HB2	2.07	0.55
1:A:31:THR:HG22	1:A:375:PRO:HD3	1.90	0.54
1:A:396:ASN:HB3	1:A:399:LEU:HB3	1.90	0.54
1:A:276:MET:HE1	1:A:284:MET:HG3	1.90	0.54
1:A:183:LEU:HD12	1:A:289:MET:HE1	1.89	0.54
1:A:217:TRP:HZ3	1:A:220:ARG:HH11	1.54	0.54
1:A:430:LEU:O	1:A:434:THR:OG1	2.25	0.54
1:A:189:ARG:O	1:A:192:LEU:O	2.27	0.53
1:A:357:LEU:HD11	1:A:462:VAL:HG21	1.91	0.53
1:A:89:PHE:CD2	1:A:382:CYS:HB2	2.44	0.53
1:A:118:ALA:HB1	1:A:122:ARG:CG	2.38	0.52
1:A:186:GLU:O	1:A:190:LYS:HG2	2.09	0.52
1:A:45:ILE:HG23	1:A:46:VAL:HG23	1.91	0.52
1:A:98:SER:CB	1:A:364:LYS:HE3	2.38	0.52
1:A:246:ALA:O	1:A:250:GLU:OE2	2.26	0.52
1:A:330:LEU:HA	1:A:334:ASN:HD22	1.75	0.52
1:A:104:THR:C	1:A:106:MET:H	2.17	0.52
1:A:113:GLY:C	1:A:118:ALA:HB2	2.34	0.52
1:A:344:ARG:HD2	4:A:2007:HOH:O	2.10	0.51
1:A:465:THR:CB	1:A:468:GLN:NE2	2.73	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:LEU:CB	4:A:2003:HOH:O	2.55	0.51
1:A:118:ALA:HB1	1:A:122:ARG:HG3	1.92	0.51
1:A:233:ARG:HH12	1:A:279:HIS:CD2	2.29	0.50
1:A:28:ALA:C	1:A:30:LYS:N	2.66	0.50
1:A:42:LEU:HB3	1:A:45:ILE:CG2	2.41	0.50
1:A:118:ALA:HB1	1:A:122:ARG:HD2	1.94	0.50
1:A:465:THR:CG2	1:A:468:GLN:HE21	2.23	0.50
1:A:223:LEU:O	1:A:225:GLN:N	2.46	0.49
1:A:69:THR:HA	1:A:78:THR:HA	1.94	0.49
1:A:59:CYS:O	1:A:63:LEU:HB2	2.12	0.49
1:A:107:THR:N	1:A:108:PRO:CD	2.75	0.49
1:A:139:PHE:CD1	1:A:142:PHE:CD2	3.01	0.49
1:A:348:GLU:HA	1:A:348:GLU:OE1	2.12	0.49
1:A:213:VAL:HG22	1:A:213:VAL:O	2.13	0.48
1:A:465:THR:HB	1:A:468:GLN:NE2	2.27	0.48
1:A:33:PRO:HG2	1:A:63:LEU:HD12	1.95	0.48
1:A:203:LYS:HG3	1:A:228:ARG:HD3	1.94	0.47
1:A:51:ASN:HD22	1:A:54:GLU:HB2	1.78	0.47
1:A:272:ASP:OD1	1:A:274:THR:OG1	2.25	0.47
1:A:126:GLN:HA	1:A:129:PHE:HD2	1.77	0.47
1:A:106:MET:C	1:A:108:PRO:HD2	2.40	0.47
1:A:100:ARG:HE	1:A:100:ARG:HB2	1.60	0.47
1:A:295:THR:HG21	3:A:1490:UDO:HAG	1.97	0.46
1:A:152:LYS:HB3	1:A:152:LYS:HE2	1.69	0.46
1:A:217:TRP:HA	1:A:217:TRP:HE3	1.79	0.46
1:A:146:ILE:HG12	1:A:178:THR:HB	1.95	0.46
1:A:339:MET:CE	1:A:437:ALA:HB2	2.45	0.46
1:A:395:PRO:O	1:A:396:ASN:HB2	2.14	0.46
1:A:98:SER:HB2	1:A:364:LYS:HE3	1.96	0.46
1:A:205:GLU:C	1:A:207:SER:N	2.74	0.46
1:A:366:GLU:HA	1:A:374:VAL:O	2.16	0.46
2:A:1480:HEM:HHA	2:A:1480:HEM:HBD2	1.96	0.46
1:A:203:LYS:HG2	1:A:228:ARG:CG	2.46	0.45
1:A:387:SER:C	1:A:389:HIS:N	2.74	0.45
1:A:418:GLY:O	1:A:421:LYS:HG2	2.16	0.45
1:A:360:MET:HE3	3:A:1490:UDO:FAC	2.04	0.45
1:A:180:CYS:O	1:A:184:PHE:HB2	2.16	0.45
1:A:100:ARG:NH2	1:A:118:ALA:O	2.50	0.45
1:A:131:ALA:O	1:A:135:THR:HG23	2.16	0.45
1:A:158:TRP:O	1:A:475:ARG:NH2	2.50	0.45
1:A:42:LEU:HB3	1:A:45:ILE:HG22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:LEU:HD23	1:A:466:LEU:HD13	1.99	0.45
1:A:218:LEU:C	1:A:220:ARG:H	2.25	0.44
1:A:449:ARG:NH2	1:A:468:GLN:OE1	2.51	0.44
1:A:309:HIS:CD2	1:A:311:LYS:HB2	2.52	0.44
1:A:30:LYS:HE2	1:A:31:THR:H	1.82	0.44
1:A:44:HIS:O	1:A:48:PHE:HB2	2.17	0.44
1:A:107:THR:N	1:A:108:PRO:HD2	2.33	0.44
1:A:310:PRO:O	1:A:313:LYS:HG3	2.18	0.44
1:A:446:GLN:HE22	1:A:448:LEU:HD23	1.76	0.44
1:A:293:GLN:HA	1:A:293:GLN:OE1	2.18	0.44
1:A:469:CYS:O	1:A:471:VAL:HG23	2.18	0.44
1:A:344:ARG:HB3	1:A:402:PRO:O	2.17	0.43
1:A:351:ARG:O	1:A:388:HIS:CD2	2.71	0.43
1:A:248:GLU:C	1:A:250:GLU:H	2.27	0.43
1:A:321:LYS:HA	1:A:321:LYS:HD3	1.87	0.43
1:A:332:TYR:CZ	1:A:336:MET:HG3	2.52	0.43
1:A:358:MET:HE3	1:A:381:ALA:HB1	2.00	0.43
1:A:412:PHE:CD1	1:A:412:PHE:C	2.95	0.43
1:A:422:CYS:HA	2:A:1480:HEM:C4D	2.53	0.43
1:A:91:SER:N	1:A:92:PRO:CD	2.82	0.43
1:A:115:ALA:O	1:A:118:ALA:HB3	2.17	0.43
1:A:218:LEU:C	1:A:220:ARG:N	2.77	0.43
1:A:265:LEU:O	1:A:268:ALA:HB3	2.18	0.43
1:A:122:ARG:HD3	1:A:126:GLN:HE21	1.80	0.43
1:A:226:SER:O	1:A:229:CYS:N	2.49	0.43
1:A:35:TYR:CD1	1:A:36:PRO:HD2	2.53	0.43
1:A:103:TYR:CE1	3:A:1490:UDO:HAM	2.54	0.43
1:A:318:LYS:NZ	1:A:403:GLU:OE2	2.52	0.43
1:A:143:VAL:HB	1:A:144:PRO:HD3	2.01	0.42
1:A:52:PRO:HD2	1:A:457:TYR:CG	2.53	0.42
1:A:424:GLY:HA3	2:A:1480:HEM:C2C	2.54	0.42
1:A:48:PHE:CE1	1:A:52:PRO:HB3	2.54	0.42
1:A:476:LYS:O	1:A:477:LYS:HG3	2.20	0.42
1:A:435:ILE:O	1:A:436:LEU:C	2.62	0.42
1:A:116:TYR:HD1	1:A:123:MET:CE	2.27	0.41
1:A:83:PRO:HA	1:A:86:HIS:CE1	2.55	0.41
1:A:260:ASP:C	1:A:260:ASP:OD1	2.64	0.41
1:A:167:LEU:HD12	1:A:167:LEU:HA	1.80	0.41
1:A:203:LYS:HG3	1:A:228:ARG:HH11	1.85	0.41
1:A:205:GLU:C	1:A:207:SER:H	2.28	0.41
1:A:91:SER:N	1:A:92:PRO:HD3	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:HIS:HD2	1:A:71:SER:O	2.04	0.41
1:A:116:TYR:CD1	1:A:123:MET:SD	3.13	0.41
1:A:86:HIS:CD2	1:A:382:CYS:SG	3.14	0.41
1:A:225:GLN:NE2	1:A:225:GLN:O	2.54	0.41
1:A:48:PHE:CD1	1:A:52:PRO:HA	2.55	0.41
1:A:40:PRO:C	1:A:42:LEU:H	2.29	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	445/460 (97%)	412 (93%)	32 (7%)	1 (0%)	43 72

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	453	PRO

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	391/402 (97%)	344 (88%)	47 (12%)	5 17

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	LYS
1	A	30	LYS
1	A	37	VAL
1	A	38	THR
1	A	47	GLN
1	A	54	GLU
1	A	63	LEU
1	A	87	SER
1	A	101	GLU
1	A	106	MET
1	A	107	THR
1	A	122	ARG
1	A	123	MET
1	A	132	GLU
1	A	159	LYS
1	A	165	ILE
1	A	175	ILE
1	A	191	ARG
1	A	192	LEU
1	A	195	ARG
1	A	209	ILE
1	A	217	TRP
1	A	219	LEU
1	A	220	ARG
1	A	225	GLN
1	A	238	LYS
1	A	247	ARG
1	A	251	GLU
1	A	258	THR
1	A	271	ARG
1	A	274	THR
1	A	278	LEU
1	A	296	SER
1	A	297	THR
1	A	308	MET
1	A	317	ASP
1	A	321	LYS
1	A	339	MET
1	A	391	GLU
1	A	413	ILE
1	A	421	LYS
1	A	434	THR

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Mol	Chain	Res	Type
1	A	446	GLN
1	A	447	LEU
1	A	449	ARG
1	A	459	THR
1	A	460	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	HIS
1	A	47	GLN
1	A	51	ASN
1	A	140	GLN
1	A	193	ASN
1	A	306	HIS
1	A	309	HIS
1	A	329	GLN
1	A	334	ASN
1	A	388	HIS
1	A	389	HIS
1	A	446	GLN
1	A	458	HIS
1	A	468	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 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$
2	HEM	A	1480	3,1	50,50,50	1.47	7 (14%)	67,82,82	2.04	25 (37%)
3	UDO	A	1490	2	35,35,35	2.66	5 (14%)	48,50,50	1.37	6 (12%)

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	3,1	-	4/14/54/54	-
3	UDO	A	1490	2	-	2/26/36/36	0/4/4/4

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1490	UDO	CAZ-CBC	-10.19	1.39	1.53
3	A	1490	UDO	CAY-CBC	-10.05	1.39	1.53
3	A	1490	UDO	CBF-CBB	-5.00	1.39	1.49
2	A	1480	HEM	FE-NB	4.68	2.09	1.94
2	A	1480	HEM	FE-NC	3.76	2.07	1.95
2	A	1480	HEM	C1B-NB	-3.43	1.34	1.40
2	A	1480	HEM	C4D-ND	-3.38	1.34	1.40
2	A	1480	HEM	C4B-NB	-2.52	1.33	1.38
3	A	1490	UDO	CAQ-NAV	2.45	1.39	1.34
2	A	1480	HEM	C1D-ND	-2.41	1.34	1.38
2	A	1480	HEM	C3C-C4C	-2.08	1.42	1.46
3	A	1490	UDO	CAG-NAV	2.07	1.39	1.33

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1480	HEM	CHD-C1D-C2D	-4.44	118.02	125.03
2	A	1480	HEM	CAD-CBD-CGD	-4.32	102.20	113.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1480	HEM	C2D-C1D-ND	4.27	114.84	109.90
2	A	1480	HEM	CAD-C3D-C4D	4.05	131.76	124.70
3	A	1490	UDO	CAT-NBE-CAU	4.03	120.62	111.57
3	A	1490	UDO	CAR-NBD-CAS	3.83	120.50	112.68
2	A	1480	HEM	C3D-C4D-ND	3.47	113.97	110.17
2	A	1480	HEM	C2B-C1B-NB	3.44	113.80	109.84
2	A	1480	HEM	CHC-C4B-NB	3.25	127.92	124.42
2	A	1480	HEM	CHA-C1A-NA	3.12	129.52	123.86
2	A	1480	HEM	C1A-CHA-C4D	-3.11	118.94	126.25
2	A	1480	HEM	C2A-C1A-NA	-2.88	106.96	110.15
2	A	1480	HEM	C4C-NC-C1C	2.82	110.41	105.82
2	A	1480	HEM	C1D-C2D-C3D	-2.81	104.02	106.98
3	A	1490	UDO	CAZ-CAQ-NAV	-2.76	119.75	124.05
2	A	1480	HEM	CHA-C4D-C3D	-2.70	120.25	125.23
2	A	1480	HEM	CAA-CBA-CGA	-2.57	106.84	113.67
2	A	1480	HEM	CHD-C1D-ND	2.51	127.12	124.42
2	A	1480	HEM	C4B-C3B-C2B	-2.48	105.00	107.28
2	A	1480	HEM	C4D-ND-C1D	-2.45	102.31	105.21
2	A	1480	HEM	C4C-CHD-C1D	-2.43	120.86	126.02
3	A	1490	UDO	CAH-CAZ-CAQ	2.43	119.39	116.89
2	A	1480	HEM	CAD-C3D-C2D	-2.39	123.39	127.87
3	A	1490	UDO	CAM-CBA-NBE	-2.38	118.09	121.39
2	A	1480	HEM	O2D-CGD-CBD	2.34	121.41	114.00
2	A	1480	HEM	C3B-C4B-NB	2.34	111.15	109.47
2	A	1480	HEM	CHD-C4C-NC	2.17	126.82	124.45
3	A	1490	UDO	CAG-NAV-CAQ	2.17	120.65	116.85
2	A	1480	HEM	C3B-C2B-C1B	-2.09	104.84	106.41
2	A	1480	HEM	CHC-C4B-C3B	-2.08	120.91	125.07
2	A	1480	HEM	C3A-C4A-NA	-2.01	106.93	110.14

There are no chirality outliers.

All (6) torsion outliers are listed below:

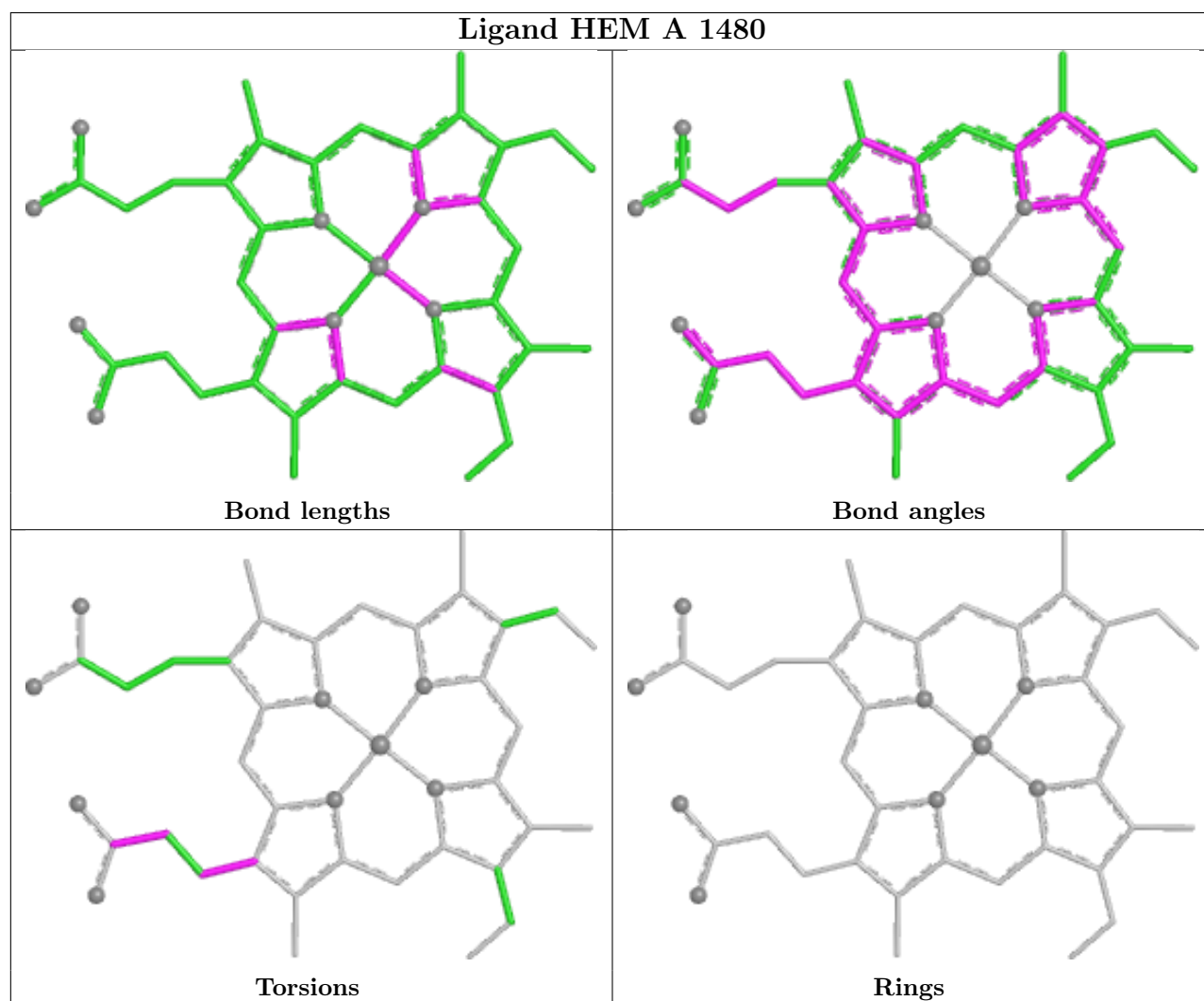
Mol	Chain	Res	Type	Atoms
2	A	1480	HEM	C4D-C3D-CAD-CBD
2	A	1480	HEM	C2D-C3D-CAD-CBD
2	A	1480	HEM	CAD-CBD-CGD-O1D
2	A	1480	HEM	CAD-CBD-CGD-O2D
3	A	1490	UDO	CAH-CAZ-CBC-CAW
3	A	1490	UDO	CAQ-CAZ-CBC-CAW

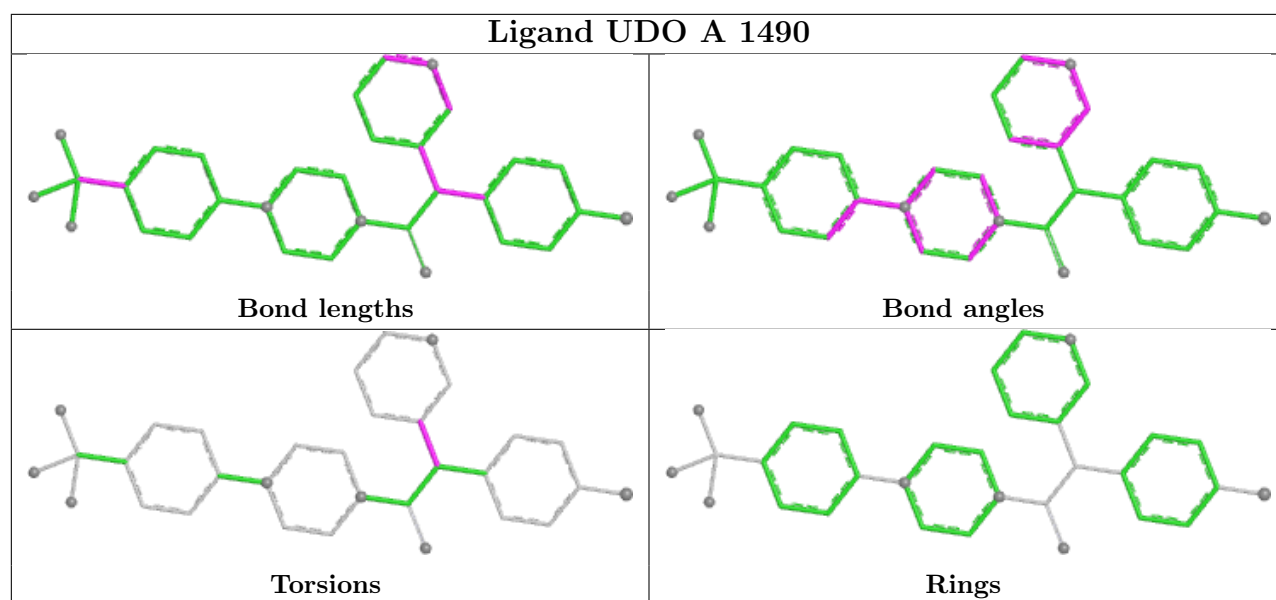
There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1480	HEM	4	0
3	A	1490	UDO	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	449/460 (97%)	0.86	52 (11%) 9 7	20, 56, 101, 122	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	253	SER	9.4
1	A	28	ALA	5.2
1	A	219	LEU	5.1
1	A	223	LEU	5.1
1	A	252	ALA	4.3
1	A	210	PRO	4.2
1	A	470	LEU	4.2
1	A	211	ALA	4.2
1	A	200	LEU	4.0
1	A	217	TRP	3.9
1	A	218	LEU	3.9
1	A	230	ARG	3.8
1	A	258	THR	3.7
1	A	256	ASN	3.7
1	A	214	PHE	3.6
1	A	224	PRO	3.5
1	A	255	ASP	3.4
1	A	225	GLN	3.3
1	A	229	CYS	3.0
1	A	216	PRO	3.0
1	A	246	ALA	3.0
1	A	456	ASP	3.0
1	A	215	MET	2.9
1	A	227	ALA	2.9
1	A	458	HIS	2.7
1	A	73	GLY	2.7
1	A	254	LYS	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	209	ILE	2.6
1	A	212	ALA	2.6
1	A	107	THR	2.5
1	A	192	LEU	2.5
1	A	30	LYS	2.5
1	A	257	ASN	2.5
1	A	450	ASP	2.4
1	A	163	GLY	2.3
1	A	29	LYS	2.3
1	A	120	TYR	2.3
1	A	137	ALA	2.2
1	A	236	LEU	2.2
1	A	94	ASN	2.2
1	A	191	ARG	2.1
1	A	250	GLU	2.1
1	A	44	HIS	2.1
1	A	123	MET	2.1
1	A	238	LYS	2.1
1	A	41	PHE	2.1
1	A	54	GLU	2.1
1	A	259	SER	2.1
1	A	474	THR	2.1
1	A	129	PHE	2.0
1	A	234	ALA	2.0
1	A	160	GLU	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 ⓘ

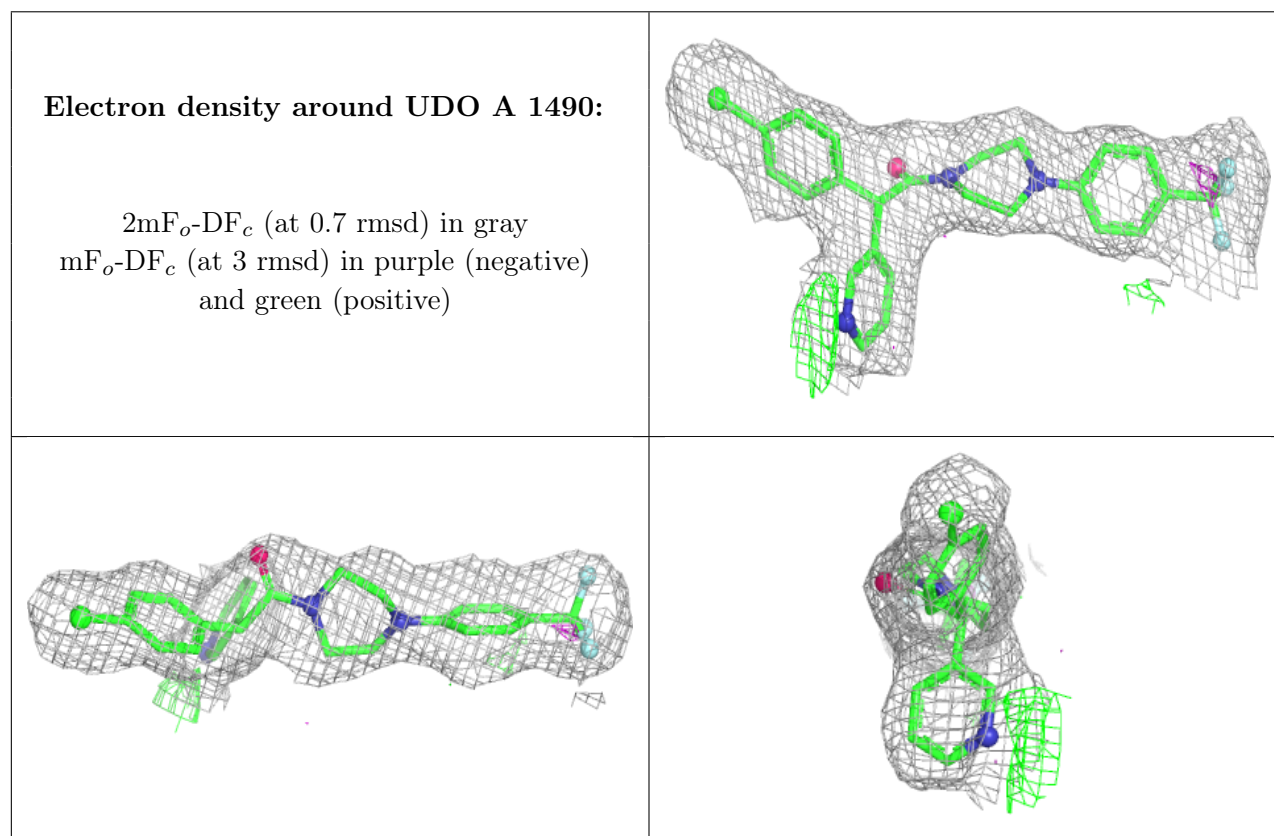
There are no oligosaccharides in this entry.

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.

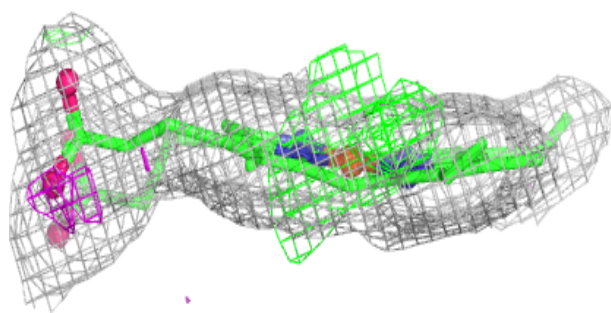
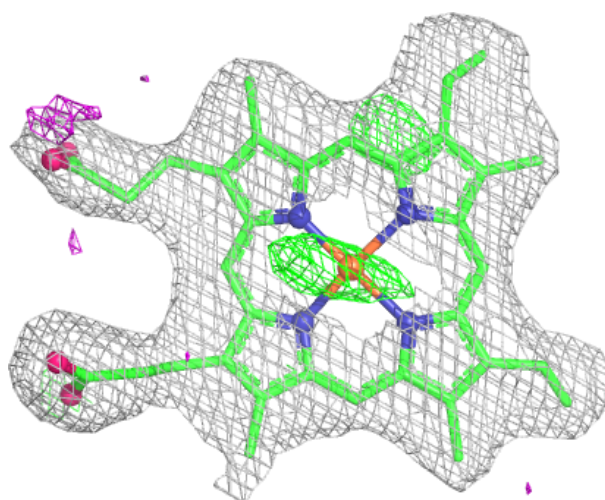
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	UDO	A	1490	32/32	0.91	0.11	26,36,51,54	0
2	HEM	A	1480	43/43	0.96	0.11	22,26,30,34	0

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 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.