



Full wwPDB X-ray Structure Validation Report

Dec 12, 2023 – 12:06 pm GMT

PDB ID : 3ZG3
Title : STEROL 14-ALPHA DEMETHYLASE (CYP51)FROM TRYPANOSOMA CRUZI IN COMPLEX WITH THE PYRIDINE INHIBITOR N-(1-(5-(trifluoromethyl)(pyridin-2-yl)) piperidin-4yl)-N-(4-(trifluoromethyl)phenyl)pyridin-3-amine (EPL- BS967, UDD)
Authors : Hargrove, T.Y.; Wawrzak, Z.; Keenan, M.; Chatelain, E.; Lepesheva, G.I.
Deposited on : 2012-12-14
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](#) ①) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

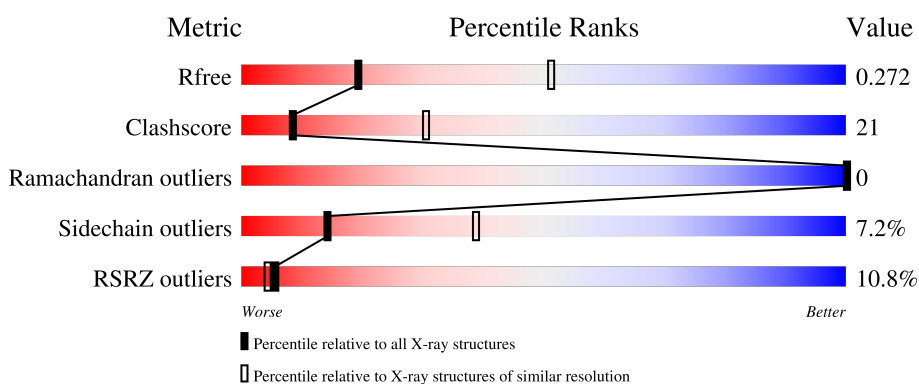
1 Overall quality at a glance

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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (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	460	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3639 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	445	3557	2270	623	636	28	0	0	0

There are 10 discrepancies between the modelled and reference sequences:

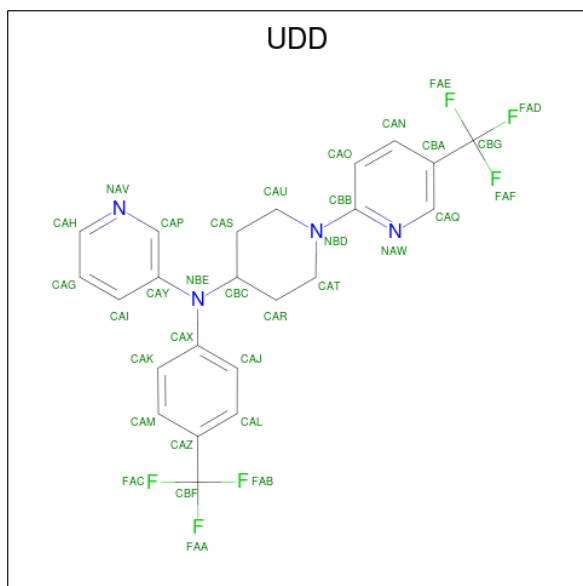
Chain	Residue	Modelled	Actual	Comment	Reference
A	28	ALA	ARG	engineered mutation	UNP Q7Z1V1
A	29	LYS	PRO	engineered mutation	UNP Q7Z1V1
A	30	LYS	THR	engineered mutation	UNP Q7Z1V1
A	31	THR	ASP	engineered mutation	UNP Q7Z1V1
A	482	HIS	-	expression tag	UNP Q7Z1V1
A	483	HIS	-	expression tag	UNP Q7Z1V1
A	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 (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

- Molecule 3 is N-[4-(trifluoromethyl)phenyl]-N-[1-[5-(trifluoromethyl)pyridin-2-yl]piperidin-4-yl]pyridin-3-amine (three-letter code: UDD) (formula: C₂₃H₂₀F₆N₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	F	N		
3	A	1	33	23	6	4	0	0

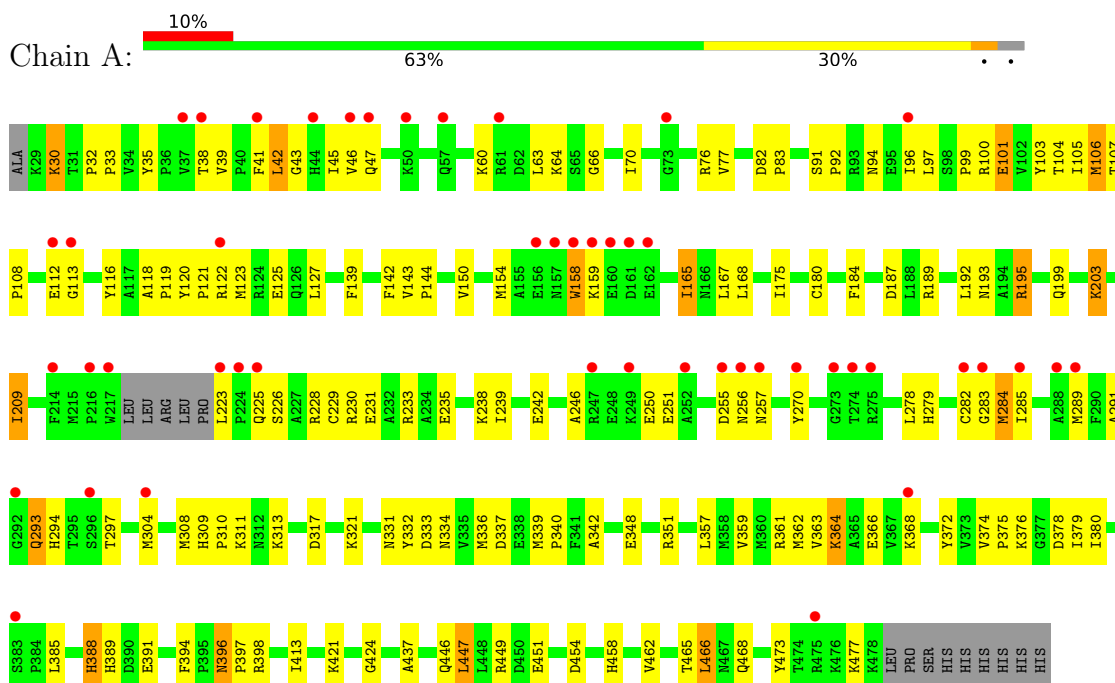
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	6	Total O 6 6	0	0

3 Residue-property plots

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 i

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	62.85Å 62.85Å 222.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.90 27.44 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.5 (30.00-2.90) 99.6 (27.44-2.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 2.89Å)	Xtrriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.248 , 0.289 0.249 , 0.272	Depositor DCC
R_{free} test set	567 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	93.1	Xtrriage
Anisotropy	0.041	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 69.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	0.054 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3639	wwPDB-VP
Average B, all atoms (Å ²)	114.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.69% 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: UDD, 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	0.23	0/3640	0.40	0/4917

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3557	0	3578	139	0
2	A	43	0	30	13	0
3	A	33	0	20	8	0
4	A	6	0	0	2	0
All	All	3639	0	3628	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)
2:A:490:HEM:HBC2	2:A:490:HEM:HMC2	1.35	1.03
1:A:293:GLN:O	1:A:297:THR:HG22	1.58	1.03
2:A:490:HEM:HMB2	2:A:490:HEM:HBB2	1.50	0.93
1:A:458:HIS:CE1	4:A:2006:HOH:O	2.27	0.88
1:A:105:ILE:HG13	1:A:106:MET:SD	2.13	0.87
1:A:366:GLU:HB3	1:A:374:VAL:O	1.76	0.85
1:A:42:LEU:HD22	1:A:45:ILE:HG22	1.57	0.84
2:A:490:HEM:HBB2	2:A:490:HEM:CMB	2.08	0.82
1:A:42:LEU:CD2	1:A:45:ILE:HG22	2.09	0.82
1:A:357:LEU:HD22	1:A:385:LEU:HD22	1.61	0.82
2:A:490:HEM:HBC2	2:A:490:HEM:CMC	2.09	0.80
1:A:103:TYR:HB3	1:A:116:TYR:HE2	1.50	0.76
1:A:42:LEU:HD22	1:A:42:LEU:O	1.87	0.75
1:A:168:LEU:HD23	1:A:466:LEU:HD22	1.66	0.75
1:A:103:TYR:HB3	1:A:116:TYR:CE2	2.22	0.74
1:A:30:LYS:HA	1:A:30:LYS:NZ	2.03	0.73
1:A:293:GLN:O	1:A:297:THR:CG2	2.36	0.73
1:A:103:TYR:CB	1:A:116:TYR:CE2	2.73	0.72
1:A:143:VAL:HB	1:A:144:PRO:HD3	1.72	0.71
1:A:351:ARG:O	1:A:351:ARG:HD2	1.90	0.71
1:A:388:HIS:HE1	1:A:413:ILE:H	1.39	0.71
1:A:189:ARG:HA	1:A:192:LEU:O	1.92	0.70
1:A:359:VAL:HG13	2:A:490:HEM:O2A	1.92	0.70
1:A:38:THR:HG23	1:A:39:VAL:H	1.57	0.69
1:A:105:ILE:CG1	1:A:106:MET:SD	2.81	0.68
1:A:64:LYS:HA	4:A:2001:HOH:O	1.93	0.67
1:A:255:ASP:O	1:A:257:ASN:ND2	2.27	0.66
1:A:30:LYS:HA	1:A:30:LYS:HZ3	1.61	0.66
1:A:203:LYS:HG3	1:A:228:ARG:HD3	1.79	0.65
1:A:103:TYR:CB	1:A:116:TYR:HE2	2.09	0.65
1:A:351:ARG:HD2	1:A:351:ARG:C	2.17	0.64
1:A:42:LEU:HD13	1:A:43:GLY:C	2.18	0.64
1:A:77:VAL:HA	1:A:379:ILE:HG23	1.80	0.64
1:A:154:MET:O	1:A:158:TRP:HB2	1.98	0.63
1:A:209:ILE:N	1:A:209:ILE:HD13	2.14	0.62
1:A:363:VAL:HG22	1:A:378:ASP:O	2.00	0.62
1:A:366:GLU:CB	1:A:374:VAL:O	2.47	0.62
1:A:310:PRO:O	1:A:313:LYS:HG3	2.00	0.62
3:A:491:UDD:HAK	3:A:491:UDD:HAS1	1.83	0.60
1:A:293:GLN:HG3	1:A:294:HIS:N	2.16	0.59
1:A:389:HIS:HE1	1:A:398:ARG:HH11	1.51	0.59
1:A:91:SER:N	1:A:92:PRO:CD	2.66	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:ARG:NH1	1:A:116:TYR:O	2.36	0.58
2:A:490:HEM:HMB2	2:A:490:HEM:CBB	2.29	0.58
1:A:99:PRO:HD2	1:A:120:TYR:OH	2.03	0.58
1:A:447:LEU:HD22	1:A:449:ARG:HB2	1.85	0.58
3:A:491:UDD:HAS1	3:A:491:UDD:CAK	2.34	0.58
1:A:317:ASP:O	1:A:321:LYS:HG2	2.02	0.58
2:A:490:HEM:HBD2	2:A:490:HEM:HHA	1.86	0.57
1:A:116:TYR:HD1	1:A:123:MET:SD	2.27	0.57
1:A:348:GLU:HA	1:A:348:GLU:OE1	2.02	0.57
1:A:42:LEU:HD21	1:A:46:VAL:H	1.69	0.57
1:A:76:ARG:O	1:A:379:ILE:HG22	2.05	0.57
1:A:60:LYS:HG3	1:A:66:GLY:HA2	1.87	0.56
1:A:38:THR:HG23	1:A:39:VAL:N	2.20	0.56
1:A:309:HIS:HD2	1:A:311:LYS:HB2	1.71	0.56
1:A:100:ARG:HH22	1:A:119:PRO:HA	1.68	0.56
1:A:363:VAL:HG21	1:A:375:PRO:O	2.04	0.56
1:A:42:LEU:HD13	1:A:42:LEU:C	2.26	0.55
1:A:424:GLY:HA3	2:A:490:HEM:C2C	2.41	0.55
2:A:490:HEM:HMC2	2:A:490:HEM:CBC	2.25	0.54
2:A:490:HEM:C4D	3:A:491:UDD:HAP	2.42	0.54
1:A:375:PRO:O	1:A:378:ASP:HB2	2.07	0.54
1:A:388:HIS:CE1	1:A:413:ILE:H	2.21	0.53
1:A:107:THR:N	1:A:108:PRO:HD2	2.23	0.53
1:A:235:GLU:O	1:A:239:ILE:HG13	2.07	0.53
1:A:180:CYS:O	1:A:184:PHE:HB2	2.08	0.53
1:A:33:PRO:HD3	1:A:372:TYR:CE2	2.45	0.52
1:A:291:ALA:HB2	3:A:491:UDD:HAJ	1.92	0.52
1:A:389:HIS:HA	1:A:397:PRO:HB3	1.90	0.52
1:A:175:ILE:HG21	1:A:293:GLN:HA	1.92	0.51
1:A:357:LEU:HD21	1:A:462:VAL:HG21	1.92	0.51
1:A:35:TYR:HD1	1:A:63:LEU:HD21	1.75	0.51
1:A:339:MET:CE	1:A:437:ALA:HB2	2.41	0.51
1:A:379:ILE:HG23	1:A:379:ILE:O	2.11	0.51
1:A:230:ARG:HG3	1:A:231:GLU:N	2.27	0.50
1:A:91:SER:H	1:A:92:PRO:HD3	1.77	0.50
1:A:396:ASN:N	1:A:397:PRO:HD3	2.26	0.50
1:A:103:TYR:HB2	1:A:116:TYR:CD2	2.47	0.49
1:A:447:LEU:HD22	1:A:449:ARG:H	1.77	0.49
1:A:105:ILE:O	1:A:108:PRO:HD2	2.12	0.49
1:A:101:GLU:HG3	1:A:362:MET:HG2	1.93	0.48
1:A:113:GLY:O	1:A:118:ALA:HB2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195:ARG:HH11	1:A:195:ARG:HB3	1.78	0.48
1:A:233:ARG:NH2	1:A:282:CYS:SG	2.86	0.48
1:A:125:GLU:OE1	1:A:270:TYR:HD1	1.96	0.48
1:A:119:PRO:HB2	1:A:121:PRO:HD2	1.96	0.48
1:A:339:MET:HG2	1:A:342:ALA:HB3	1.96	0.48
1:A:256:ASN:O	1:A:257:ASN:OD1	2.32	0.47
1:A:91:SER:N	1:A:92:PRO:HD3	2.29	0.47
1:A:96:ILE:N	1:A:96:ILE:HD12	2.30	0.47
2:A:490:HEM:ND	3:A:491:UDD:HAP	2.29	0.47
3:A:491:UDD:CAK	3:A:491:UDD:CAS	2.91	0.47
1:A:42:LEU:HD13	1:A:43:GLY:N	2.30	0.47
1:A:256:ASN:ND2	1:A:256:ASN:H	2.12	0.47
1:A:103:TYR:CB	1:A:116:TYR:CD2	2.98	0.47
1:A:116:TYR:CD1	1:A:116:TYR:N	2.77	0.47
1:A:167:LEU:HD21	1:A:304:MET:HB2	1.95	0.47
1:A:284:MET:CE	1:A:284:MET:HA	2.45	0.46
1:A:332:TYR:CZ	1:A:336:MET:HG3	2.50	0.46
1:A:424:GLY:HA3	2:A:490:HEM:C3C	2.49	0.46
1:A:70:ILE:HG13	1:A:77:VAL:HB	1.97	0.46
1:A:94:ASN:HA	1:A:97:LEU:O	2.16	0.46
1:A:226:SER:HA	1:A:229:CYS:HB3	1.98	0.46
1:A:112:GLU:H	1:A:279:HIS:HE1	1.62	0.46
1:A:285:ILE:O	1:A:289:MET:HG2	2.16	0.46
1:A:282:CYS:SG	1:A:283:GLY:N	2.89	0.46
1:A:465:THR:HB	1:A:468:GLN:NE2	2.31	0.46
1:A:165:ILE:HD12	1:A:473:TYR:HD2	1.80	0.45
1:A:331:ASN:H	1:A:334:ASN:ND2	2.15	0.45
1:A:339:MET:HE1	1:A:437:ALA:HB2	1.99	0.45
1:A:103:TYR:HB2	1:A:116:TYR:CE2	2.50	0.45
1:A:223:LEU:HD12	1:A:223:LEU:N	2.31	0.45
1:A:454:ASP:H	1:A:468:GLN:HE22	1.63	0.45
1:A:127:LEU:HD21	3:A:491:UDD:FAB	2.07	0.45
1:A:150:VAL:O	1:A:154:MET:HG3	2.17	0.45
1:A:333:ASP:O	1:A:337:ASP:HB2	2.16	0.45
1:A:465:THR:HB	1:A:468:GLN:HE21	1.82	0.45
1:A:120:TYR:N	1:A:121:PRO:CD	2.80	0.44
1:A:364:LYS:O	1:A:376:LYS:HG3	2.18	0.43
1:A:105:ILE:HD11	3:A:491:UDD:HAQ	2.00	0.43
1:A:389:HIS:CE1	1:A:398:ARG:HH11	2.34	0.43
1:A:42:LEU:HD23	1:A:46:VAL:HG23	1.99	0.43
1:A:150:VAL:HG12	1:A:154:MET:HE2	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:ALA:O	1:A:250:GLU:HG3	2.18	0.43
1:A:32:PRO:HA	1:A:372:TYR:CD2	2.54	0.43
1:A:309:HIS:CD2	1:A:311:LYS:H	2.36	0.43
1:A:304:MET:O	1:A:308:MET:HB2	2.18	0.43
1:A:139:PHE:HA	1:A:142:PHE:CD2	2.54	0.42
1:A:332:TYR:CE1	1:A:336:MET:HG3	2.54	0.42
1:A:361:ARG:HB2	1:A:380:ILE:CG2	2.48	0.42
1:A:309:HIS:CD2	1:A:311:LYS:HB2	2.52	0.42
1:A:348:GLU:OE1	1:A:348:GLU:CA	2.67	0.42
1:A:143:VAL:HG21	1:A:332:TYR:HA	2.00	0.42
1:A:187:ASP:N	1:A:187:ASP:OD1	2.52	0.42
1:A:238:LYS:HE3	1:A:242:GLU:OE2	2.20	0.41
1:A:107:THR:N	1:A:108:PRO:CD	2.84	0.41
1:A:339:MET:N	1:A:340:PRO:CD	2.84	0.41
1:A:193:ASN:HD21	1:A:195:ARG:HH11	1.68	0.41
1:A:104:THR:O	1:A:105:ILE:C	2.58	0.41
1:A:363:VAL:HG23	1:A:376:LYS:HA	2.01	0.41
1:A:42:LEU:HD22	1:A:42:LEU:C	2.41	0.41
1:A:77:VAL:HA	1:A:379:ILE:CG2	2.48	0.41
1:A:359:VAL:HG13	2:A:490:HEM:CGA	2.50	0.41
1:A:394:PHE:O	1:A:397:PRO:HG3	2.21	0.41
1:A:42:LEU:HD23	1:A:45:ILE:HG22	2.00	0.41
1:A:82:ASP:HA	1:A:83:PRO:HD2	1.94	0.41
1:A:199:GLN:O	1:A:203:LYS:HB2	2.22	0.40
1:A:368:LYS:H	1:A:368:LYS:HG2	1.63	0.40
1:A:351:ARG:C	1:A:351:ARG:CD	2.87	0.40
1:A:396:ASN:HD22	1:A:396:ASN:HA	1.63	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	441/460 (96%)	421 (96%)	20 (4%)	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	388/402 (96%)	360 (93%)	28 (7%)	14	39

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	LYS
1	A	41	PHE
1	A	42	LEU
1	A	47	GLN
1	A	101	GLU
1	A	106	MET
1	A	122	ARG
1	A	158	TRP
1	A	159	LYS
1	A	165	ILE
1	A	195	ARG
1	A	203	LYS
1	A	209	ILE
1	A	225	GLN
1	A	251	GLU
1	A	278	LEU
1	A	284	MET
1	A	293	GLN
1	A	364	LYS
1	A	388	HIS
1	A	391	GLU
1	A	396	ASN
1	A	421	LYS

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Mol	Chain	Res	Type
1	A	446	GLN
1	A	447	LEU
1	A	451	GLU
1	A	466	LEU
1	A	477	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	51	ASN
1	A	193	ASN
1	A	256	ASN
1	A	309	HIS
1	A	329	GLN
1	A	334	ASN
1	A	388	HIS
1	A	389	HIS
1	A	396	ASN
1	A	431	GLN
1	A	468	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 monosaccharides 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	490	1,3	41,50,50	1.91	10 (24%)	45,82,82	2.32	14 (31%)
3	UDD	A	491	2	36,36,36	1.56	8 (22%)	52,53,53	1.50	5 (9%)

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	490	1,3	-	4/12/54/54	-
3	UDD	A	491	2	-	8/28/38/38	0/4/4/4

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	490	HEM	C1B-NB	-5.51	1.30	1.40
3	A	491	UDD	CBG-CBA	-4.77	1.39	1.49
2	A	490	HEM	C4D-ND	-4.31	1.32	1.40
3	A	491	UDD	CBF-CAZ	-4.22	1.40	1.49
2	A	490	HEM	C1D-ND	-3.70	1.31	1.38
2	A	490	HEM	C4B-NB	-3.51	1.31	1.38
2	A	490	HEM	FE-ND	-3.14	1.81	1.96
3	A	491	UDD	CAQ-NAW	2.94	1.40	1.34
2	A	490	HEM	C3C-C2C	-2.93	1.36	1.40
3	A	491	UDD	CAP-NAV	2.91	1.40	1.34
2	A	490	HEM	C1B-C2B	-2.76	1.39	1.44
2	A	490	HEM	C3D-C2D	-2.73	1.30	1.36
3	A	491	UDD	CAY-NBE	-2.55	1.33	1.42
3	A	491	UDD	CAH-NAV	2.40	1.40	1.33
2	A	490	HEM	C4A-CHB	-2.26	1.34	1.41
3	A	491	UDD	CAX-NBE	-2.24	1.34	1.42
2	A	490	HEM	O2A-CGA	-2.09	1.23	1.30
3	A	491	UDD	CAU-NBD	2.05	1.49	1.46

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	490	HEM	CHC-C4B-NB	6.18	131.14	124.43
2	A	490	HEM	CHD-C1D-ND	5.53	130.44	124.43
2	A	490	HEM	CAD-C3D-C4D	5.36	134.02	124.66
3	A	491	UDD	CAT-NBD-CAU	4.57	121.61	111.52
3	A	491	UDD	CAX-NBE-CBC	4.51	124.85	119.37
2	A	490	HEM	CAD-CBD-CGD	-3.99	105.01	113.60
2	A	490	HEM	CHD-C1D-C2D	-3.92	118.85	124.98
2	A	490	HEM	CAD-C3D-C2D	-3.90	120.62	127.88
2	A	490	HEM	CHB-C1B-NB	3.83	129.11	124.38
2	A	490	HEM	C1B-NB-C4B	3.72	108.91	105.07
3	A	491	UDD	CBA-CAQ-NAW	-3.45	120.04	123.34
3	A	491	UDD	CAY-CAP-NAV	-2.72	119.92	122.92
2	A	490	HEM	C4C-CHD-C1D	-2.69	119.01	122.56
2	A	490	HEM	CHA-C4D-ND	2.45	127.40	124.38
2	A	490	HEM	CHB-C1B-C2B	-2.41	120.06	126.72
2	A	490	HEM	CHC-C4B-C3B	-2.32	121.02	124.57
2	A	490	HEM	CMA-C3A-C4A	-2.30	124.93	128.46
3	A	491	UDD	CAN-CAO-CBB	2.15	120.49	117.53
2	A	490	HEM	O2A-CGA-CBA	2.05	120.62	114.03

There are no chirality outliers.

All (12) torsion outliers are listed below:

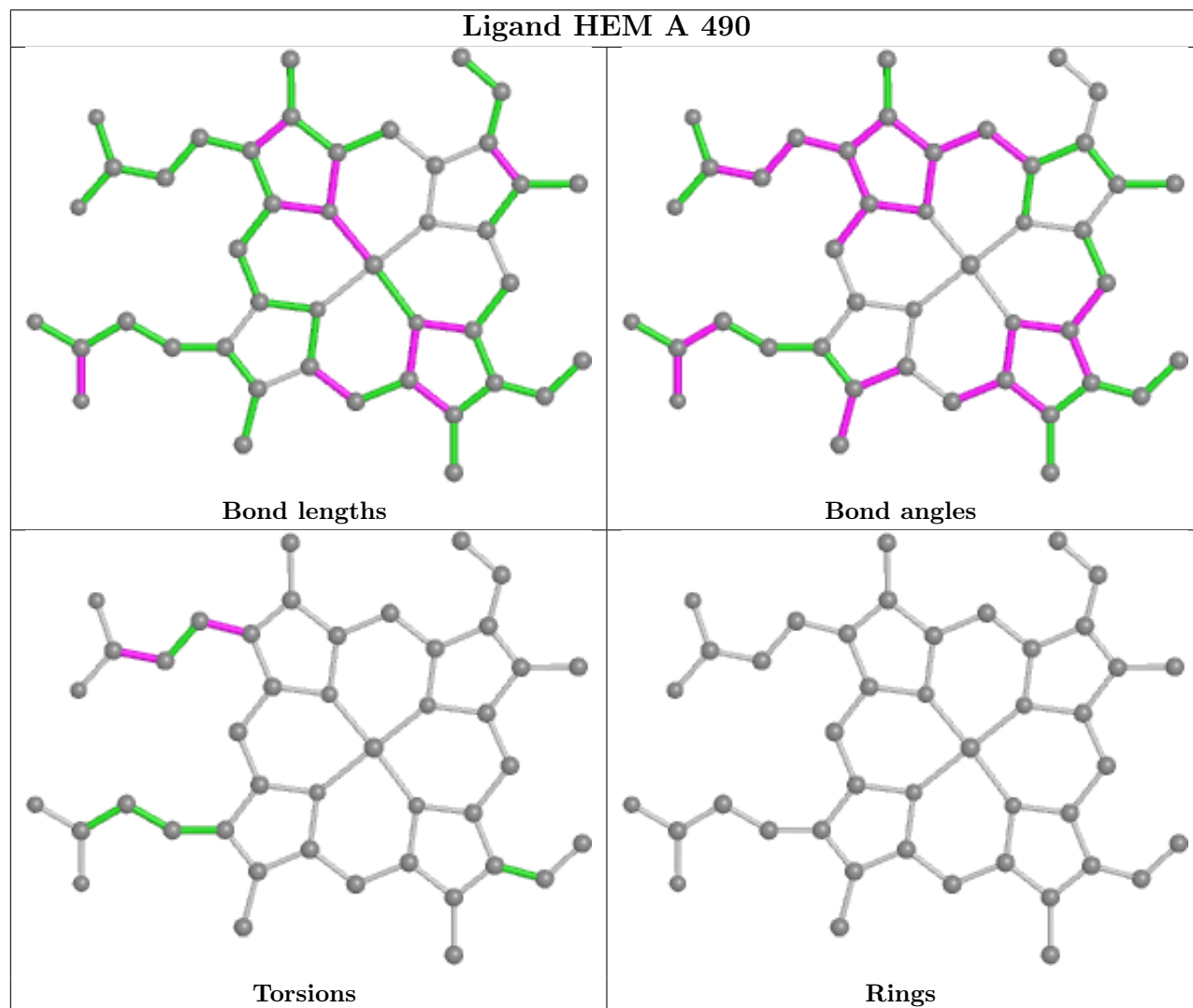
Mol	Chain	Res	Type	Atoms
2	A	490	HEM	C2D-C3D-CAD-CBD
3	A	491	UDD	CAJ-CAX-NBE-CBC
3	A	491	UDD	CAK-CAX-NBE-CBC
2	A	490	HEM	C4D-C3D-CAD-CBD
3	A	491	UDD	CAS-CBC-NBE-CAY
3	A	491	UDD	CAQ-CBA-CBG-FAF
2	A	490	HEM	CAD-CBD-CGD-O1D
2	A	490	HEM	CAD-CBD-CGD-O2D
3	A	491	UDD	CAQ-CBA-CBG-FAE
3	A	491	UDD	CAK-CAX-NBE-CAY
3	A	491	UDD	CAQ-CBA-CBG-FAD
3	A	491	UDD	CAI-CAY-NBE-CBC

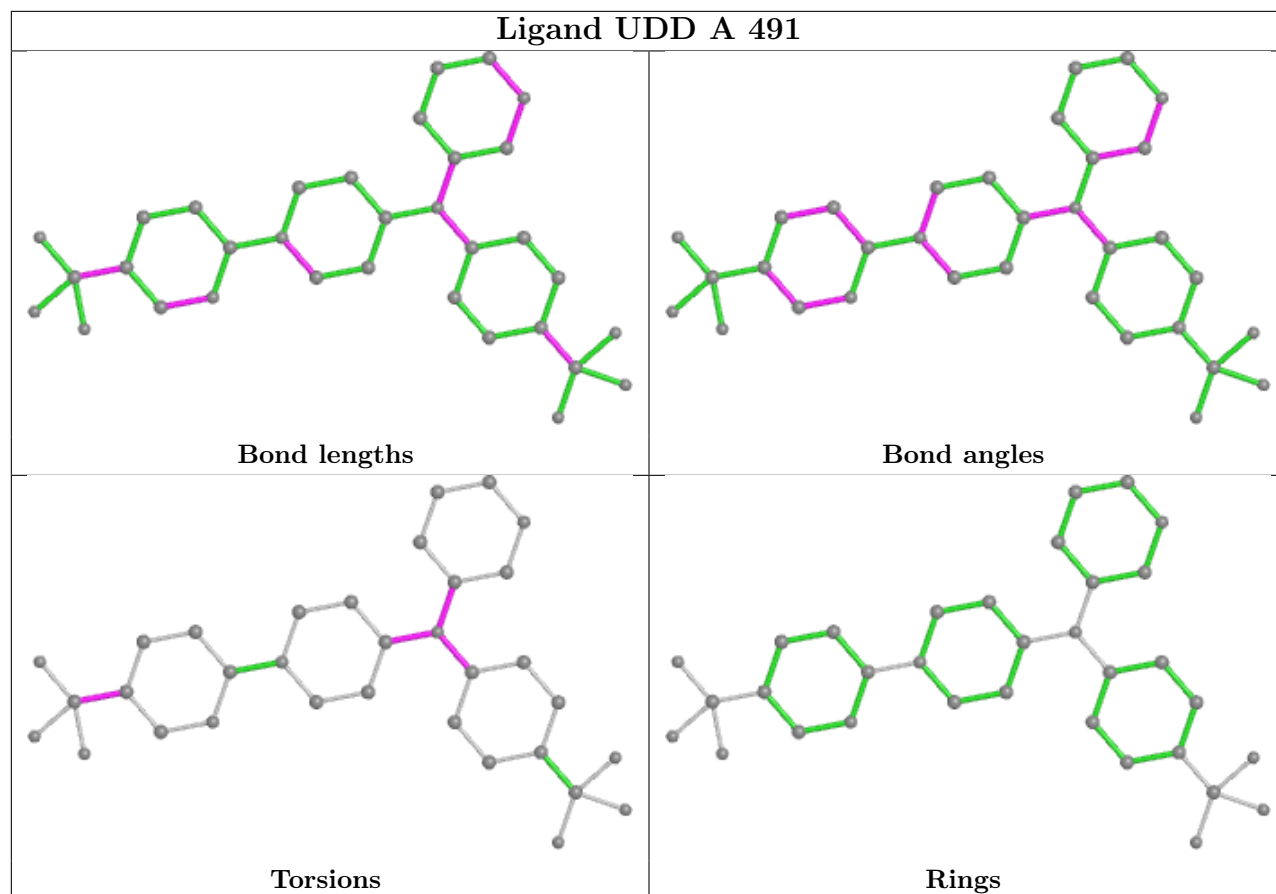
There are no ring outliers.

2 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	490	HEM	13	0
3	A	491	UDD	8	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	445/460 (96%)	0.42	48 (10%) 5 4	57, 108, 182, 252	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	224	PRO	10.5
1	A	217	TRP	8.0
1	A	223	LEU	6.7
1	A	161	ASP	5.4
1	A	274	THR	5.2
1	A	475	ARG	4.9
1	A	159	LYS	4.4
1	A	44	HIS	4.3
1	A	247	ARG	4.2
1	A	275	ARG	3.8
1	A	282	CYS	3.8
1	A	273	GLY	3.6
1	A	255	ASP	3.6
1	A	256	ASN	3.6
1	A	257	ASN	3.5
1	A	156	GLU	3.5
1	A	73	GLY	3.4
1	A	160	GLU	3.3
1	A	225	GLN	3.2
1	A	214	PHE	3.1
1	A	96	ILE	3.0
1	A	112	GLU	3.0
1	A	216	PRO	2.8
1	A	61	ARG	2.8
1	A	37	VAL	2.8
1	A	113	GLY	2.7
1	A	252	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	122	ARG	2.7
1	A	50	LYS	2.6
1	A	57	GLN	2.6
1	A	46	VAL	2.5
1	A	158	TRP	2.5
1	A	283	GLY	2.4
1	A	285	ILE	2.3
1	A	288	ALA	2.3
1	A	383	SER	2.3
1	A	249	LYS	2.3
1	A	368	LYS	2.2
1	A	47	GLN	2.2
1	A	38	THR	2.2
1	A	304	MET	2.2
1	A	41	PHE	2.1
1	A	292	GLY	2.1
1	A	270	TYR	2.1
1	A	296	SER	2.0
1	A	162	GLU	2.0
1	A	289	MET	2.0
1	A	157	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](#)

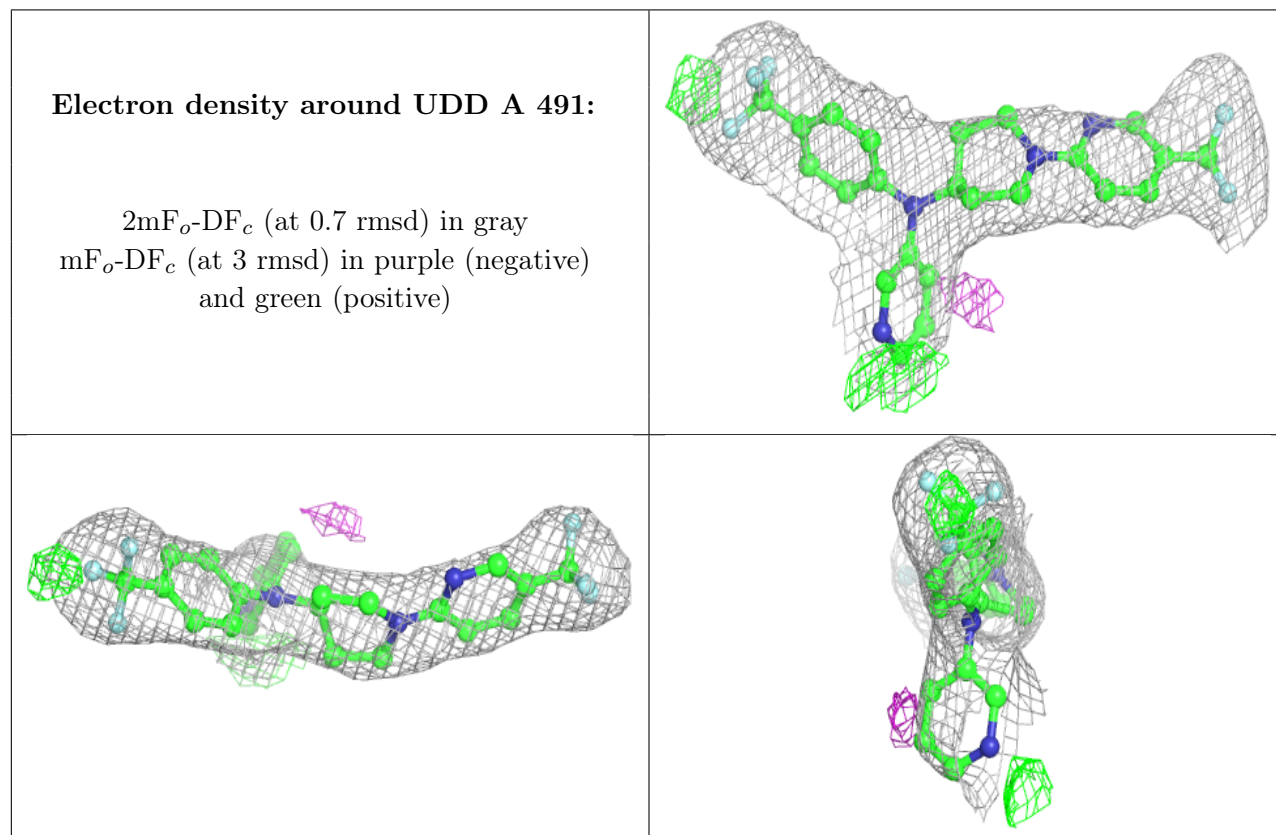
There are no monosaccharides in this entry.

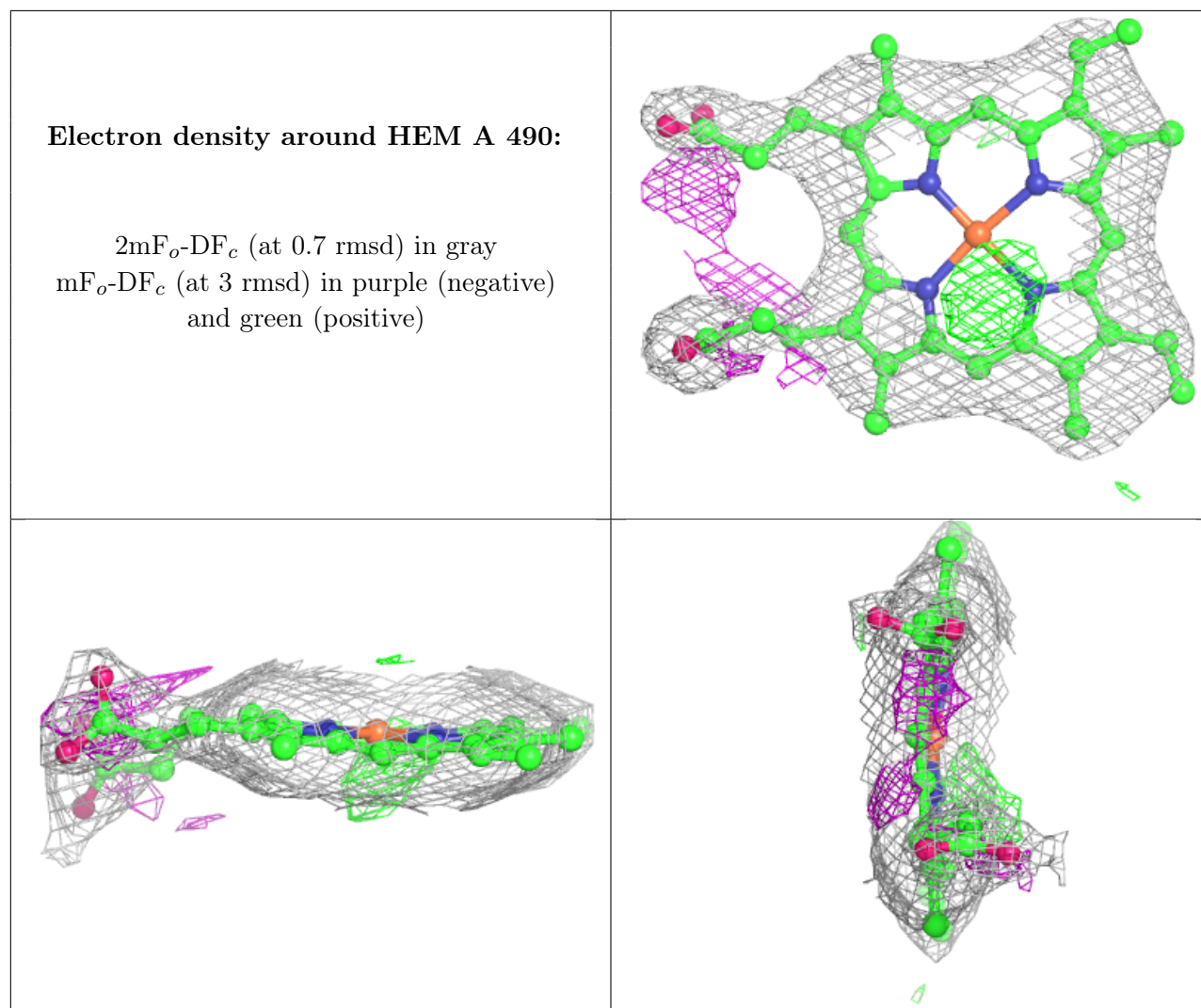
6.4 Ligands [i](#)

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	UDD	A	491	33/33	0.95	0.26	65,85,133,136	0
2	HEM	A	490	43/43	0.96	0.24	48,64,77,85	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.