



Full wwPDB X-ray Structure Validation Report ⓘ

May 27, 2020 – 11:21 pm BST

PDB ID : 3LI2
Title : Closed Conformation of HtsA Complexed with Staphyloferrin A
Authors : Grigg, J.C.; Murphy, M.E.P.
Deposited on : 2010-01-23
Resolution : 2.20 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

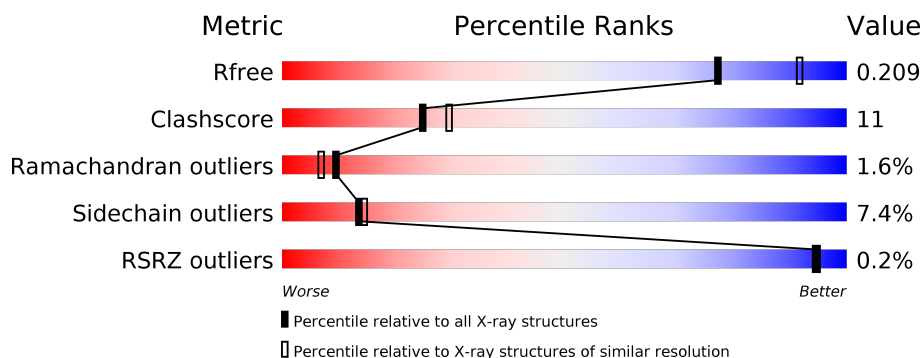
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	296	 71% 24% . .
1	B	296	 71% 24% . .

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4945 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 Ferrichrome ABC transporter lipoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	0	4	0
			2316	1466	408	439	3			
1	B	289	Total	C	N	O	S	0	5	0
			2333	1478	410	442	3			

There are 14 discrepancies between the modelled and reference sequences:

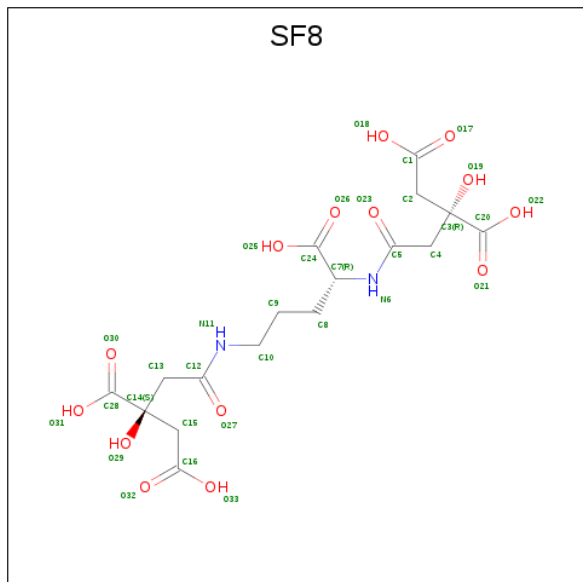
Chain	Residue	Modelled	Actual	Comment	Reference
A	32	GLY	-	EXPRESSION TAG	UNP A6QJ18
A	33	SER	-	EXPRESSION TAG	UNP A6QJ18
A	34	HIS	-	EXPRESSION TAG	UNP A6QJ18
A	35	MET	-	EXPRESSION TAG	UNP A6QJ18
A	36	ALA	-	EXPRESSION TAG	UNP A6QJ18
A	37	SER	-	EXPRESSION TAG	UNP A6QJ18
A	38	THR	-	EXPRESSION TAG	UNP A6QJ18
B	32	GLY	-	EXPRESSION TAG	UNP A6QJ18
B	33	SER	-	EXPRESSION TAG	UNP A6QJ18
B	34	HIS	-	EXPRESSION TAG	UNP A6QJ18
B	35	MET	-	EXPRESSION TAG	UNP A6QJ18
B	36	ALA	-	EXPRESSION TAG	UNP A6QJ18
B	37	SER	-	EXPRESSION TAG	UNP A6QJ18
B	38	THR	-	EXPRESSION TAG	UNP A6QJ18

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Fe	0	0
			1	1		
2	A	1	Total	Fe	0	0
			1	1		

- Molecule 3 is (2R)-2-(2-{[(1R)-1-carboxy-4-{[(3S)-3,4-dicarboxy-3-hydroxybutanoyl]amin

o}butyl]amino}-2-oxoethyl)-2-hydroxybutanedioic acid (three-letter code: SF8) (formula: $C_{17}H_{24}N_2O_{14}$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			33	17	2	14		
3	B	1	Total	C	N	O	0	0
			33	17	2	14		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		

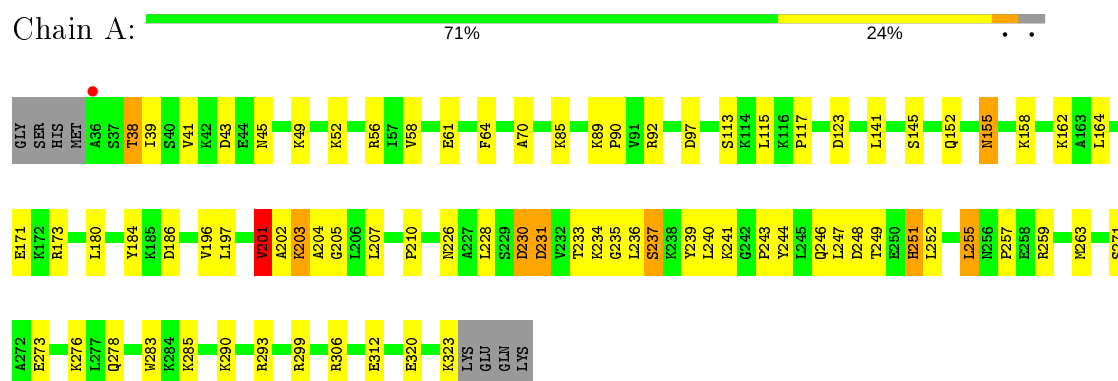
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	102	Total	O	0	0
			102	102		
6	B	118	Total	O	0	0
			118	118		

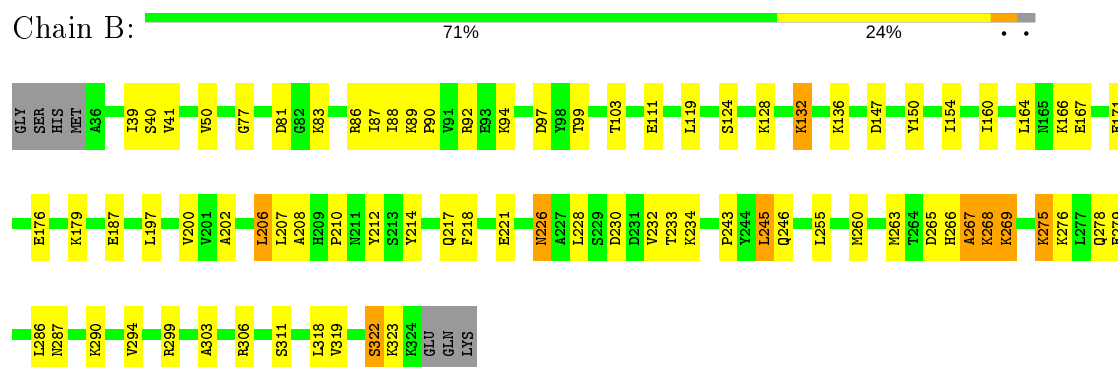
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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: Ferrichrome ABC transporter lipoprotein



- Molecule 1: Ferrichrome ABC transporter lipoprotein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.28Å 148.60Å 52.27Å 90.00° 117.14° 90.00°	Depositor
Resolution (Å)	33.15 – 2.20 33.15 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.0 (33.15-2.20) 99.1 (33.15-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.75 (at 2.20Å)	Xtriage
Refinement program	REFMAC refmac _5.5.0102	Depositor
R, R_{free}	0.165 , 0.216 0.164 , 0.209	Depositor DCC
R_{free} test set	1785 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	28.4	Xtriage
Anisotropy	0.590	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 23.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.017 for l,k,-h-l 0.017 for -h-l,k,h 0.023 for -h-l,-k,l 0.022 for h,-k,-h-l 0.477 for l,-k,h	Xtriage
Reported twinning fraction	0.509 for H, K, L 0.491 for L, -K, H	Depositor
Outliers	0 of 35627 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4945	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.19% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: ZN, FE, SF8, ACT

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	0.61	0/2349	0.69	0/3154
1	B	0.59	0/2366	0.70	0/3173
All	All	0.60	0/4715	0.70	0/6327

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
1	B	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	201	VAL	Peptide
1	B	269	LYS	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	2316	0	2408	50	0
1	B	2333	0	2433	54	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	33	0	17	4	0
3	B	33	0	17	4	0
4	A	4	0	0	0	0
5	A	4	0	3	0	0
6	A	102	0	0	3	0
6	B	118	0	0	1	0
All	All	4945	0	4878	104	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 11.

All (104) 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:202:ALA:HB2	1:A:207:LEU:HD23	1.32	1.10
1:B:263:MET:CE	1:B:303:ALA:HB2	1.87	1.04
1:B:263:MET:HE1	1:B:303:ALA:HB2	1.46	0.98
1:B:268:LYS:HA	1:B:269:LYS:HB3	1.60	0.83
1:A:248:ASP:OD2	1:A:251:HIS:N	2.13	0.79
1:B:200:VAL:O	1:B:206:LEU:O	2.02	0.78
1:B:263:MET:HE3	1:B:303:ALA:HB2	1.67	0.75
1:A:123:ASP:OD1	1:A:145:SER:HB3	1.88	0.73
1:B:132[A]:LYS:CB	1:B:132[A]:LYS:NZ	2.50	0.73
1:B:132[A]:LYS:HB3	1:B:132[A]:LYS:NZ	2.04	0.72
1:A:202:ALA:HB2	1:A:207:LEU:CD2	2.16	0.71
1:B:202:ALA:HB1	3:B:501:SF8:O25	1.94	0.68
1:B:226:ASN:HD22	1:B:228[A]:LEU:H	1.43	0.67
1:B:132[A]:LYS:CB	1:B:132[A]:LYS:HZ2	2.08	0.67
1:B:132[A]:LYS:HB2	1:B:132[A]:LYS:HZ2	1.61	0.66
1:B:81:ASP:HA	1:B:103:THR:HA	1.78	0.65
1:B:150:TYR:CZ	1:B:154:ILE:HD11	2.32	0.64
1:A:201:VAL:HG23	1:A:201:VAL:O	1.97	0.64
1:A:226:ASN:ND2	1:A:228:LEU:HB2	2.12	0.64
1:B:275:LYS:HA	1:B:278:GLN:HE21	1.62	0.64
1:B:132[A]:LYS:HB3	1:B:132[A]:LYS:HZ3	1.62	0.64
1:B:88:ILE:HG12	1:B:90:PRO:HD2	1.79	0.64
1:A:197:LEU:HD13	1:A:228:LEU:HD13	1.81	0.63
1:B:318:LEU:O	1:B:322:SER:HB2	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:VAL:CG2	1:A:201:VAL:O	2.48	0.61
1:A:58:VAL:HG23	1:A:117:PRO:HB3	1.82	0.61
1:A:278:GLN:HA	1:A:283:TRP:CD1	2.36	0.61
1:B:260:MET:HB2	1:B:294:VAL:HG22	1.84	0.59
1:A:255:LEU:HD23	1:A:257:PRO:HD3	1.84	0.59
1:A:202:ALA:HB1	3:A:501:SF8:O25	2.01	0.59
1:A:39:ILE:HG13	1:A:52:LYS:HG2	1.84	0.59
1:B:50:VAL:HG13	1:B:119:LEU:HD13	1.83	0.58
1:A:180:LEU:HD22	1:A:184:TYR:CE2	2.38	0.58
1:A:226:ASN:HD22	1:A:228:LEU:HB2	1.69	0.58
1:A:38:THR:HG21	1:A:49:LYS:HD3	1.84	0.58
1:B:226:ASN:HD22	1:B:228[B]:LEU:H	1.50	0.58
1:A:239:TYR:CZ	1:A:240:LEU:HD23	2.39	0.57
1:B:299:ARG:NH2	3:B:501:SF8:H8A	2.20	0.57
1:A:162:LYS:HE2	6:A:369:HOH:O	2.04	0.57
1:B:207:LEU:HD13	3:B:501:SF8:H10	1.87	0.57
1:B:228[B]:LEU:HD13	1:B:255:LEU:HD11	1.87	0.57
1:A:89:LYS:HB3	1:A:90:PRO:HD3	1.86	0.56
1:B:228[A]:LEU:CD2	1:B:255:LEU:HD11	2.36	0.56
1:A:259:ARG:NH2	1:A:293:ARG:HD2	2.21	0.56
1:B:83:LYS:HB2	1:B:86:ARG:HD2	1.88	0.55
1:B:228[A]:LEU:HD21	1:B:255:LEU:HD11	1.87	0.55
1:B:319:VAL:O	1:B:323:LYS:HG2	2.08	0.54
1:B:111:GLU:HA	1:B:111:GLU:OE1	2.08	0.52
1:B:212:TYR:OH	1:B:233:THR:HG23	2.10	0.52
1:B:160:ILE:O	1:B:164:LEU:HG	2.10	0.51
1:A:152:GLN:HG2	6:A:359:HOH:O	2.12	0.50
1:B:187[A]:GLU:HB3	1:B:319:VAL:HG13	1.93	0.50
1:A:202:ALA:HB1	3:A:501:SF8:H8A	1.94	0.50
1:B:210:PRO:HB3	1:B:228[B]:LEU:HD23	1.93	0.50
1:A:202:ALA:CB	1:A:207:LEU:HD23	2.23	0.50
1:A:233:THR:O	1:A:235:GLY:O	2.30	0.50
1:A:210:PRO:HD2	1:A:243:PRO:O	2.11	0.50
1:B:77:GLY:HA2	1:B:99:THR:O	2.12	0.50
1:B:268:LYS:HG2	1:B:269:LYS:HD2	1.94	0.49
1:A:61:GLU:OE2	1:A:306:ARG:NH2	2.46	0.49
1:A:41:VAL:CG1	1:A:141:LEU:HD21	2.43	0.48
1:A:202:ALA:O	1:A:203:LYS:CB	2.61	0.48
1:A:196:VAL:HA	1:A:259:ARG:O	2.14	0.48
1:B:166:LYS:HE3	6:B:359:HOH:O	2.13	0.48
1:A:70:ALA:HB1	1:A:173[B]:ARG:HG2	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:GLU:HB2	1:A:64:PHE:CE2	2.49	0.47
1:A:164:LEU:HD23	1:A:164:LEU:HA	1.72	0.47
1:B:275:LYS:O	1:B:279:GLU:HB2	2.15	0.47
1:B:214:TYR:CZ	1:B:306:ARG:HA	2.50	0.47
1:B:124:SER:O	1:B:128:LYS:HG3	2.15	0.46
1:B:87:ILE:HB	1:B:92:ARG:HB2	1.98	0.46
1:B:230:ASP:HA	1:B:233:THR:HB	1.97	0.46
1:B:81:ASP:OD2	1:B:86:ARG:NH1	2.48	0.46
1:A:202:ALA:H	1:A:205:GLY:CA	2.28	0.46
1:A:239:TYR:CE1	3:A:501:SF8:H8	2.51	0.46
1:A:237:SER:HB3	1:A:246[A]:GLN:NE2	2.31	0.45
1:B:263:MET:HE1	1:B:303:ALA:CB	2.32	0.45
1:B:217:GLN:O	1:B:221:GLU:HG2	2.17	0.44
1:A:56:ARG:NH2	1:A:115:LEU:O	2.46	0.44
1:B:286:LEU:O	1:B:287:ASN:C	2.56	0.44
1:A:306:ARG:NH1	3:A:501:SF8:O33	2.40	0.44
1:B:207:LEU:HD13	3:B:501:SF8:C10	2.48	0.44
1:A:320:GLU:O	1:A:323:LYS:HD2	2.18	0.43
1:B:197:LEU:HD11	1:B:208:ALA:HB1	1.99	0.43
1:A:285:LYS:HA	1:A:290:LYS:HE3	2.01	0.43
1:B:207:LEU:HA	1:B:245:LEU:O	2.18	0.43
1:A:173[A]:ARG:NH2	1:A:312:GLU:OE2	2.44	0.42
1:B:176:GLU:HA	1:B:179:LYS:HE2	2.01	0.42
1:A:155:ASN:HD22	1:A:158:LYS:HD3	1.85	0.42
1:A:43:ASP:OD1	1:A:45:ASN:N	2.53	0.42
1:A:155:ASN:HA	1:A:158:LYS:HD3	2.02	0.42
1:B:167:GLU:O	1:B:171:GLU:HG3	2.20	0.42
1:B:267:ALA:O	1:B:268:LYS:HG3	2.20	0.42
1:B:210:PRO:HD2	1:B:243:PRO:O	2.19	0.42
1:B:266:HIS:O	1:B:268:LYS:N	2.53	0.41
1:A:155:ASN:HA	1:A:155:ASN:HD22	1.60	0.41
1:A:263:MET:HB3	1:A:299:ARG:HG3	2.02	0.41
1:A:52:LYS:HA	6:A:348:HOH:O	2.21	0.41
1:A:252:LEU:O	1:A:255:LEU:HD22	2.21	0.40
1:B:218:PHE:CZ	1:B:311:SER:HB2	2.56	0.40
1:A:247:LEU:HG	1:A:251:HIS:HB3	2.03	0.40
1:B:164:LEU:O	1:B:166:LYS:HG2	2.21	0.40
1:A:239:TYR:HB3	1:A:244:TYR:HB2	2.03	0.40
1:A:203:LYS:HA	1:A:204:ALA:HA	1.92	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	290/296 (98%)	271 (93%)	13 (4%)	6 (2%)	7	4
1	B	292/296 (99%)	275 (94%)	14 (5%)	3 (1%)	15	14
All	All	582/592 (98%)	546 (94%)	27 (5%)	9 (2%)	9	8

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	203	LYS
1	A	236	LEU
1	A	271	SER
1	B	267	ALA
1	A	230	ASP
1	A	231	ASP
1	B	268	LYS
1	A	234	LYS
1	B	290	LYS

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	254/257 (99%)	236 (93%)	18 (7%)	14	16
1	B	256/257 (100%)	236 (92%)	20 (8%)	12	13
All	All	510/514 (99%)	472 (92%)	38 (8%)	13	14

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	THR
1	A	85	LYS
1	A	92	ARG
1	A	97	ASP
1	A	113	SER
1	A	155	ASN
1	A	171	GLU
1	A	186	ASP
1	A	201	VAL
1	A	230	ASP
1	A	231	ASP
1	A	237	SER
1	A	241	LYS
1	A	249	THR
1	A	251	HIS
1	A	255	LEU
1	A	273	GLU
1	A	276	LYS
1	B	39	ILE
1	B	40	SER
1	B	41	VAL
1	B	89	LYS
1	B	94	LYS
1	B	97	ASP
1	B	132[A]	LYS
1	B	132[B]	LYS
1	B	136	LYS
1	B	147	ASP
1	B	206	LEU
1	B	226	ASN
1	B	232	VAL
1	B	234	LYS
1	B	245	LEU
1	B	246	GLN
1	B	265	ASP
1	B	275	LYS
1	B	276	LYS
1	B	322	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	135	ASN
1	A	155	ASN
1	A	226	ASN
1	B	131	ASN
1	B	135	ASN
1	B	182	ASN
1	B	226	ASN
1	B	256	ASN
1	B	278	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](#)

Of 9 ligands modelled in this entry, 6 are monoatomic - leaving 3 for Mogul analysis.

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	SF8	B	501	2	17,32,32	0.97	2 (11%)	20,45,45	1.97	7 (35%)
3	SF8	A	501	2	17,32,32	0.47	0	20,45,45	1.18	1 (5%)
5	ACT	A	14	-	1,3,3	1.62	0	0,3,3	0.00	-

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	SF8	B	501	2	-	5/27/47/47	-
3	SF8	A	501	2	-	6/27/47/47	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	501	SF8	C13-C12	2.53	1.54	1.51
3	B	501	SF8	C8-C7	2.17	1.56	1.53

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	501	SF8	C8-C7-N6	4.83	117.22	110.19
3	B	501	SF8	C13-C12-N11	3.37	119.04	115.57
3	A	501	SF8	O23-C5-C4	-2.92	118.30	122.14
3	B	501	SF8	O27-C12-C13	-2.52	118.83	122.14
3	B	501	SF8	C7-N6-C5	-2.41	119.48	123.33
3	B	501	SF8	O23-C5-C4	-2.21	119.24	122.14
3	B	501	SF8	C3-C4-C5	-2.18	108.39	113.76
3	B	501	SF8	C10-N11-C12	2.12	126.77	122.84

There are no chirality outliers.

All (11) torsion outliers are listed below:

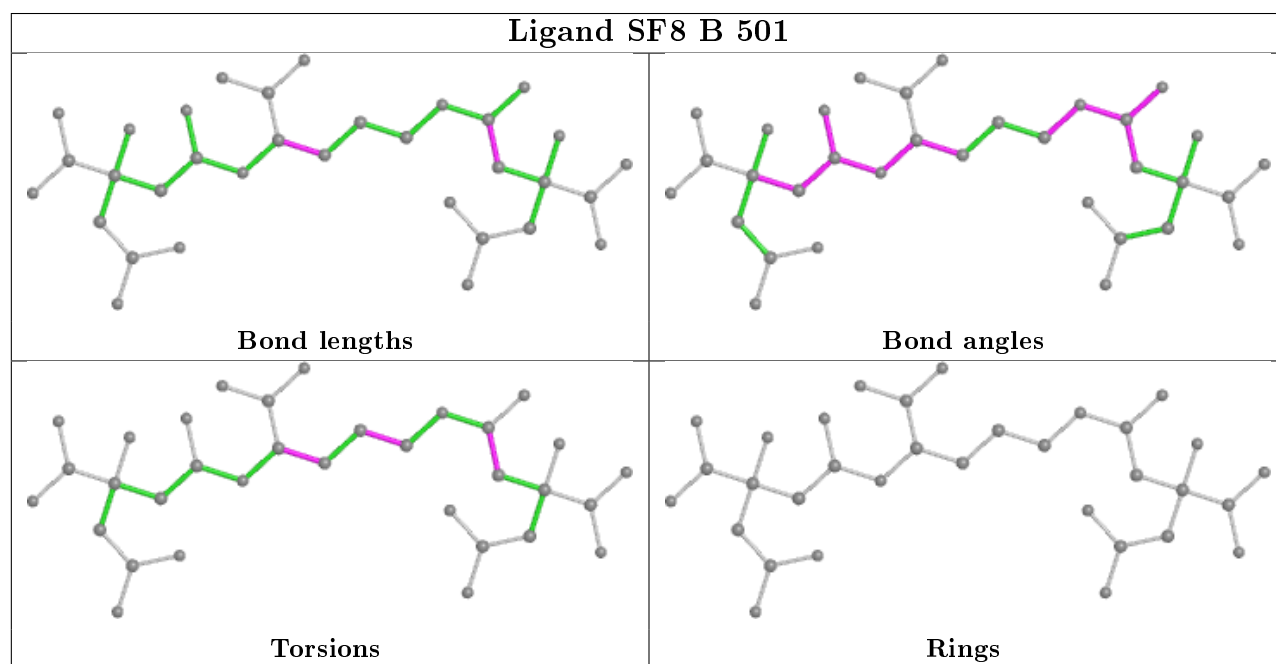
Mol	Chain	Res	Type	Atoms
3	B	501	SF8	N6-C7-C8-C9
3	B	501	SF8	C24-C7-C8-C9
3	B	501	SF8	N11-C10-C9-C8
3	A	501	SF8	N11-C10-C9-C8
3	B	501	SF8	N11-C12-C13-C14
3	B	501	SF8	O27-C12-C13-C14
3	A	501	SF8	N11-C12-C13-C14
3	A	501	SF8	O19-C3-C4-C5
3	A	501	SF8	C2-C3-C4-C5
3	A	501	SF8	O27-C12-C13-C14
3	A	501	SF8	C20-C3-C4-C5

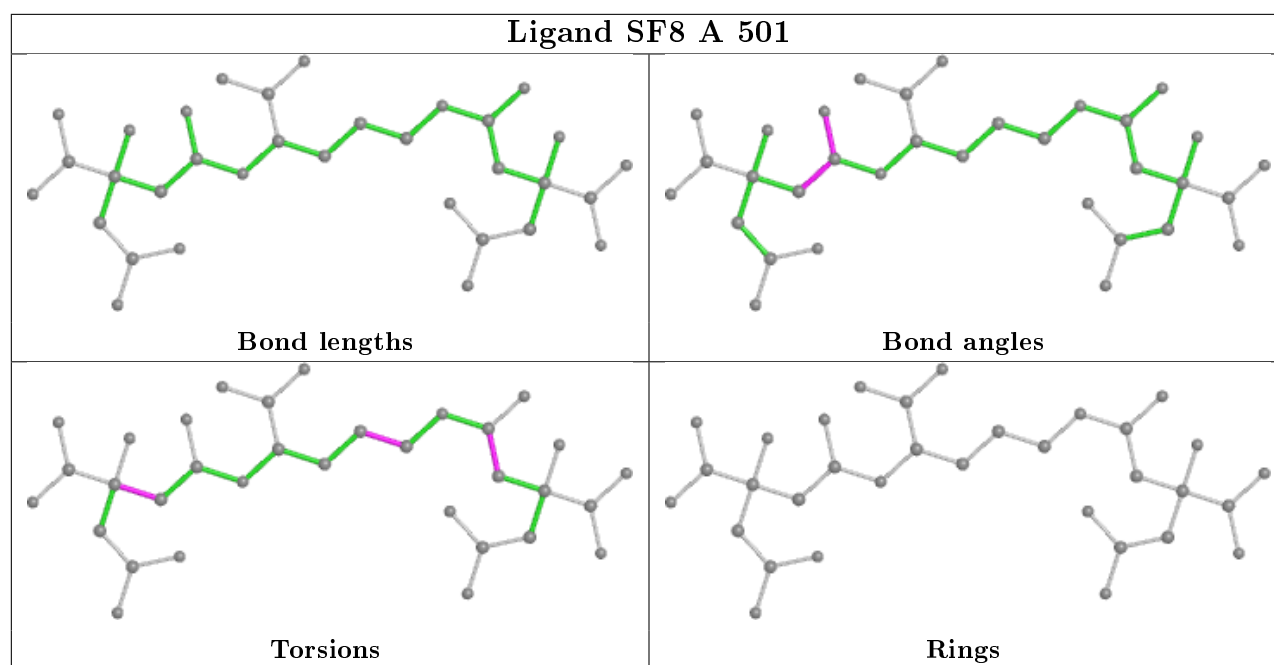
There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	501	SF8	4	0
3	A	501	SF8	4	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	288/296 (97%)	-0.64	1 (0%) 94 93	17, 35, 57, 70	0
1	B	289/296 (97%)	-0.61	0 100 100	17, 36, 54, 61	0
All	All	577/592 (97%)	-0.62	1 (0%) 95 94	17, 36, 56, 70	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	36	ALA	3.9

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(Å ²)	Q<0.9
3	SF8	A	501	33/33	0.96	0.10	15,24,36,37	0
3	SF8	B	501	33/33	0.96	0.10	14,25,34,37	0
5	ACT	A	14	4/4	0.97	0.06	30,30,31,31	0
4	ZN	A	4	1/1	0.99	0.08	34,34,34,34	0

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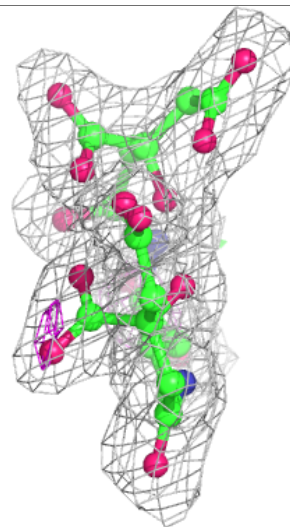
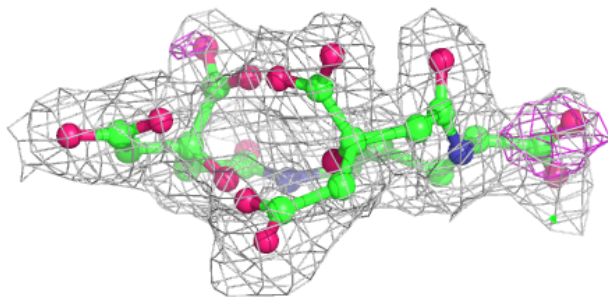
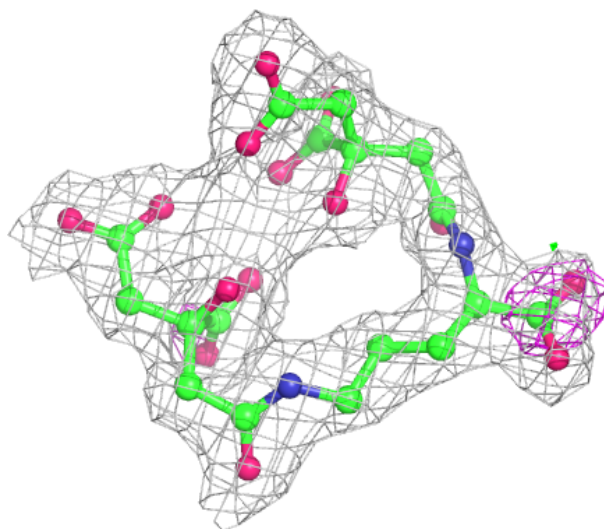
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	ZN	A	2	1/1	1.00	0.10	26,26,26,26	0
2	FE	A	500	1/1	1.00	0.09	21,21,21,21	0
4	ZN	A	1	1/1	1.00	0.06	33,33,33,33	0
4	ZN	A	3	1/1	1.00	0.09	23,23,23,23	0
2	FE	B	500	1/1	1.00	0.10	20,20,20,20	0

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.

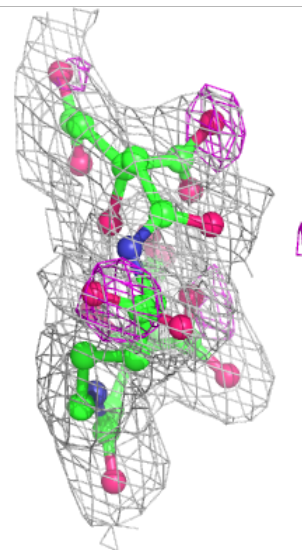
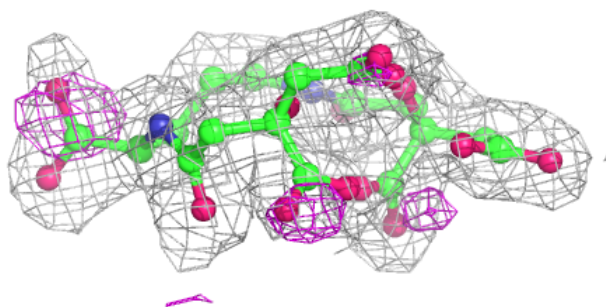
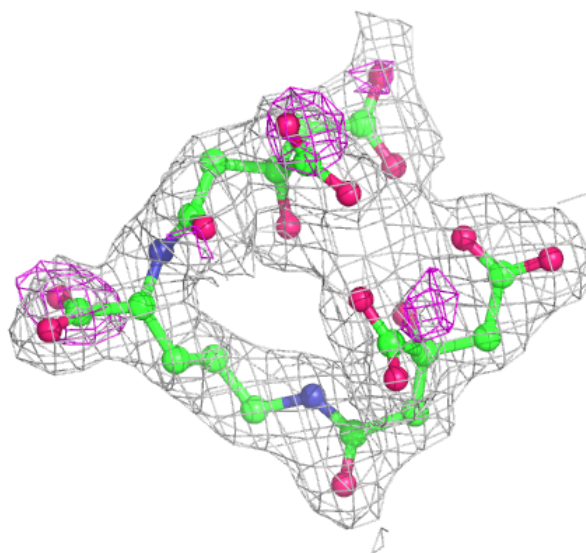
Electron density around SF8 A 501:

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 SF8 B 501:

$2mF_o-DF_c$ (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.