



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 25, 2022 – 10:28 AM EDT

PDB ID : 8D6T  
Title : Rana catesbeiana saxiphilin mutant - Y558I:STX (co-crystal)  
Authors : Chen, Z.; Zakrzewska, S.; Minor, D.L.  
Deposited on : 2022-06-06  
Resolution : 2.15 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](mailto: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>.

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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 : 2.31.2  
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) : 2.31.2

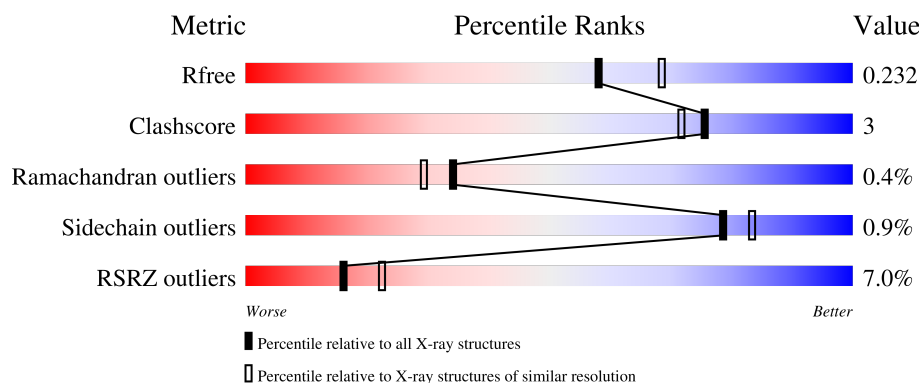
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.15 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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  $\geq 3$ , 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	851	<div> <div>6%</div> <div>89%</div> <div>7%</div> <div>.</div> </div>
1	B	851	<div> <div>8%</div> <div>87%</div> <div>9%</div> <div>.</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13488 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.

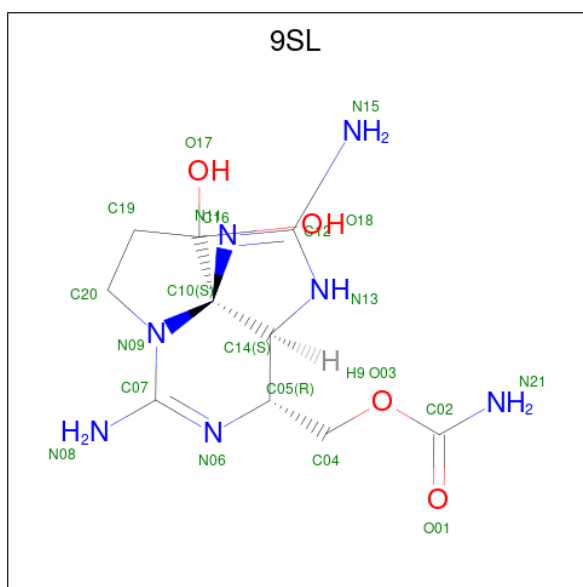
- Molecule 1 is a protein called Saxiphilin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	818	Total	C	N	O	S	0	0	0
			6311	3942	1090	1220	59			
1	A	818	Total	C	N	O	S	0	0	0
			6311	3942	1090	1220	59			

There are 16 discrepancies between the modelled and reference sequences:

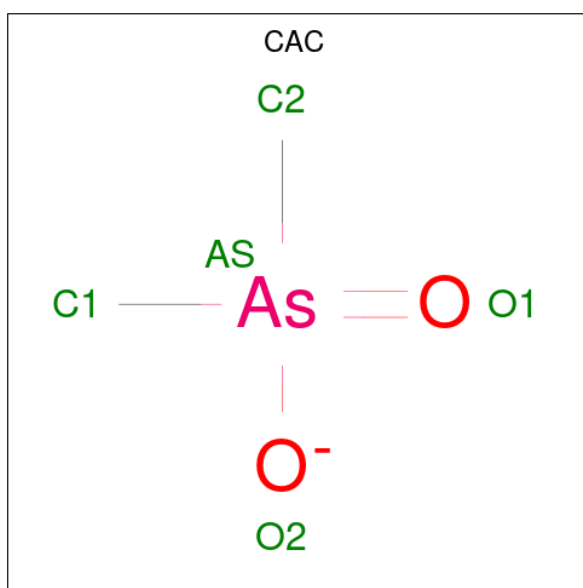
Chain	Residue	Modelled	Actual	Comment	Reference
B	558	ILE	TYR	engineered mutation	UNP P31226
B	826	SER	-	expression tag	UNP P31226
B	827	ASN	-	expression tag	UNP P31226
B	828	SER	-	expression tag	UNP P31226
B	829	LEU	-	expression tag	UNP P31226
B	830	GLU	-	expression tag	UNP P31226
B	831	VAL	-	expression tag	UNP P31226
B	832	LEU	-	expression tag	UNP P31226
A	558	ILE	TYR	engineered mutation	UNP P31226
A	826	SER	-	expression tag	UNP P31226
A	827	ASN	-	expression tag	UNP P31226
A	828	SER	-	expression tag	UNP P31226
A	829	LEU	-	expression tag	UNP P31226
A	830	GLU	-	expression tag	UNP P31226
A	831	VAL	-	expression tag	UNP P31226
A	832	LEU	-	expression tag	UNP P31226

- Molecule 2 is [(3aS,4R,10aS)-2,6-diamino-10,10-dihydroxy-3a,4,9,10-tetrahydro-3H,8H-pyrrolo[1,2-c]purin-4-yl]methyl carbamate (three-letter code: 9SL) (formula: C<sub>10</sub>H<sub>17</sub>N<sub>7</sub>O<sub>4</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total 21	C 10	N 7	O 4	0	0
2	A	1	Total 21	C 10	N 7	O 4	0	0

- Molecule 3 is CACODYLATE ION (three-letter code: CAC) (formula:  $\text{C}_2\text{H}_6\text{AsO}_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total 5	As 1	C 2	O 2	0	0
3	B	1	Total 5	As 1	C 2	O 2	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	As	C	O	0	0
			5	1	2	2		
3	A	1	Total	As	C	O	0	0
			5	1	2	2		
3	A	1	Total	As	C	O	0	0
			5	1	2	2		
3	A	1	Total	As	C	O	0	0
			5	1	2	2		

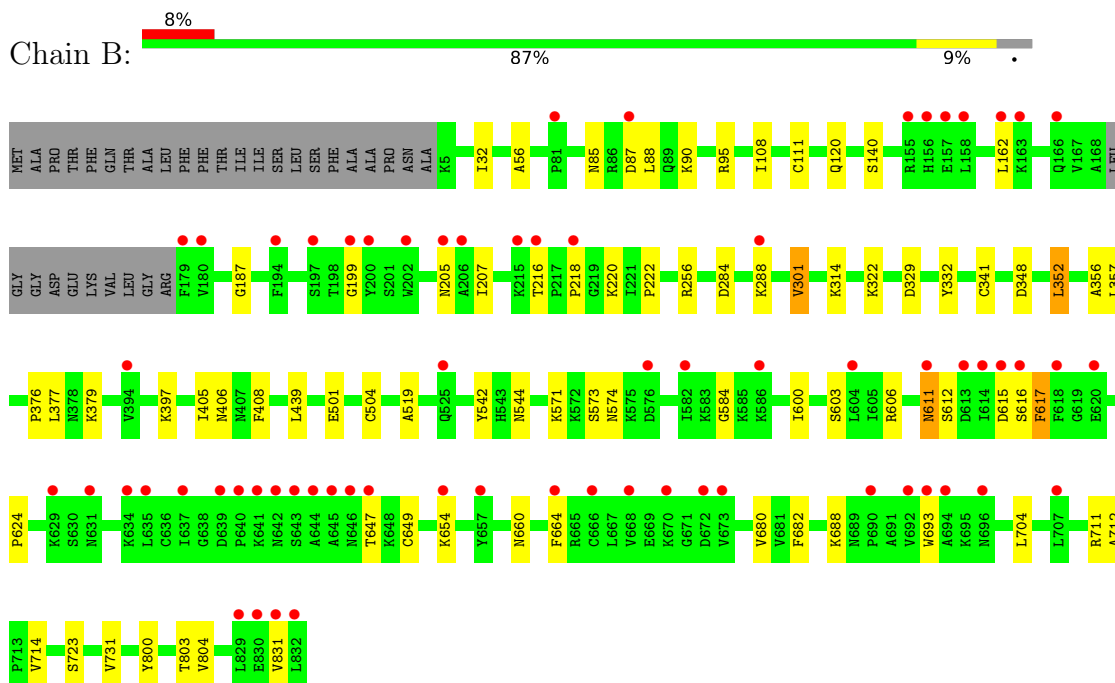
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	379	Total	O	0	0
			379	379		
4	A	415	Total	O	0	0
			415	415		

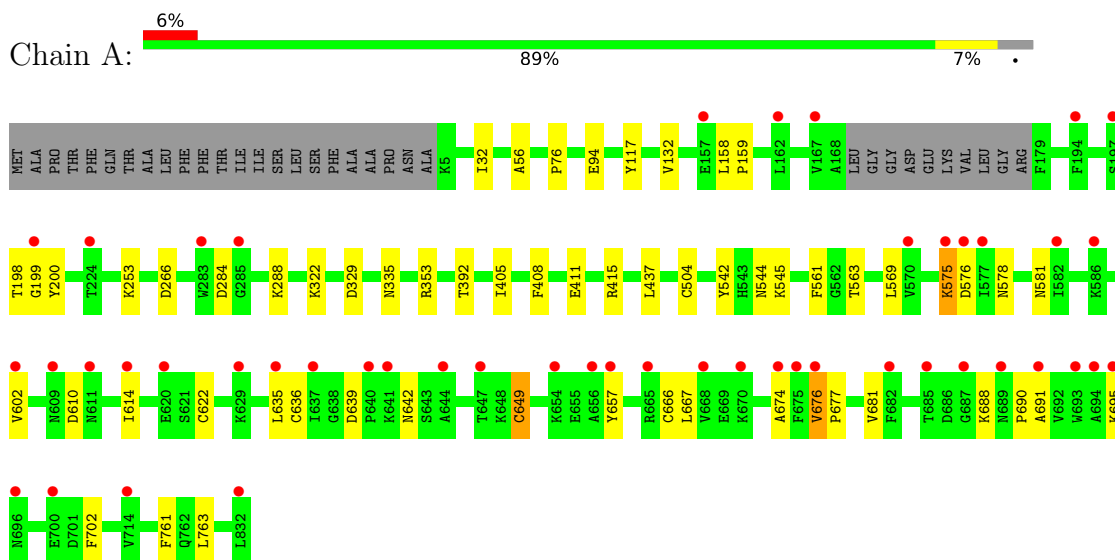
### 3 Residue-property plots [i](#)

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: Saxiphilin



#### • Molecule 1: Saxiphilin



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	96.03Å 107.81Å 253.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.50 – 2.15 47.50 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.4 (47.50-2.15) 99.4 (47.50-2.15)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.59 (at 2.16Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, $R_{free}$	0.207 , 0.234 0.206 , 0.232	Depositor DCC
$R_{free}$ test set	7099 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	44.0	Xtriage
Anisotropy	0.712	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 44.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	13488	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CAC, 9SL

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	0.27	0/6436	0.49	0/8686
1	B	0.27	0/6436	0.48	0/8686
All	All	0.27	0/12872	0.48	0/17372

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

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	6311	0	6183	32	0
1	B	6311	0	6183	46	0
2	A	21	0	0	0	0
2	B	21	0	0	0	0
3	A	15	0	0	0	0
3	B	15	0	0	0	0
4	A	415	0	0	8	0
4	B	379	0	0	7	0
All	All	13488	0	12366	78	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 3.



All (78) 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:636:CYS:SG	4:A:1328:HOH:O	2.36	0.83
1:A:602:VAL:HG23	1:A:614:ILE:HG13	1.68	0.75
1:A:649:CYS:SG	4:A:1328:HOH:O	2.46	0.73
1:B:314:LYS:NZ	4:B:1004:HOH:O	2.27	0.68
1:B:501:GLU:OE1	4:B:1001:HOH:O	2.13	0.67
1:A:253:LYS:NZ	4:A:1002:HOH:O	2.16	0.65
1:B:140:SER:HB3	1:B:831:VAL:HG12	1.78	0.64
1:B:600:ILE:HD13	1:B:723:SER:HB3	1.79	0.64
1:A:411:GLU:OE1	1:A:415:ARG:NH1	2.32	0.63
1:A:94:GLU:OE1	4:A:1001:HOH:O	2.15	0.63
1:B:504:CYS:SG	4:B:1286:HOH:O	2.55	0.60
1:A:32:ILE:HG21	1:A:408:PHE:HB2	1.83	0.60
1:B:205:ASN:O	1:B:207:ILE:N	2.32	0.60
1:A:284:ASP:HB3	1:A:288:LYS:HD2	1.83	0.59
1:B:199:GLY:HA3	1:B:218:PRO:HB3	1.85	0.58
1:A:76:PRO:HB3	1:A:392:THR:HG21	1.86	0.58
1:B:85:ASN:HA	1:B:88:LEU:HD13	1.86	0.58
1:B:647:THR:HB	1:B:654:LYS:HB3	1.86	0.57
1:B:322:LYS:HD3	1:B:329:ASP:HB3	1.86	0.57
1:B:301:VAL:HG12	1:B:314:LYS:HD2	1.86	0.56
1:A:667:LEU:HB2	1:A:674:ALA:HB2	1.88	0.56
1:A:691:ALA:HB1	1:A:695:LYS:HE3	1.88	0.55
1:A:335:ASN:ND2	4:A:1013:HOH:O	2.38	0.54
1:B:120:GLN:NE2	4:B:1015:HOH:O	2.42	0.53
1:B:87:ASP:HA	1:B:90:LYS:HB2	1.91	0.53
1:B:256:ARG:NH1	1:B:348:ASP:OD1	2.41	0.52
1:B:284:ASP:HB3	1:B:288:LYS:HD2	1.92	0.51
1:B:603:SER:OG	1:B:803:THR:O	2.27	0.51
1:A:569:LEU:HD22	1:A:681:VAL:HG22	1.91	0.50
1:B:216:THR:HG21	1:B:222:PRO:HA	1.93	0.50
1:A:504:CYS:SG	4:A:1343:HOH:O	2.59	0.50
1:A:437:LEU:HD13	4:A:1354:HOH:O	2.11	0.49
1:B:612:SER:HB2	1:B:616:SER:HB2	1.93	0.49
1:A:94:GLU:OE1	1:A:117:TYR:OH	2.30	0.49
1:B:571:LYS:O	1:B:573:SER:N	2.41	0.49
1:B:574:ASN:O	1:B:711:ARG:NH2	2.46	0.49
1:B:664:PHE:HD2	1:B:693:TRP:HZ3	1.61	0.49
1:A:691:ALA:HA	1:A:695:LYS:HG3	1.95	0.49
1:B:32:ILE:HG21	1:B:408:PHE:HB2	1.95	0.48
1:A:761:PHE:HE2	1:A:763:LEU:HD23	1.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:712:ALA:HB3	4:B:1278:HOH:O	2.13	0.47
1:A:622:CYS:HB2	1:A:635:LEU:HD12	1.96	0.47
1:A:676:VAL:HG12	1:A:677:PRO:HD2	1.96	0.47
1:B:357:LEU:HG	1:B:377:LEU:HG	1.98	0.46
1:B:624:PRO:HA	1:B:649:CYS:HA	1.98	0.46
1:A:569:LEU:HD11	1:A:702:PHE:HB3	1.98	0.45
1:A:322:LYS:HD2	1:A:329:ASP:HB3	1.99	0.45
1:B:584:GLY:N	1:B:617:PHE:O	2.47	0.44
1:A:159:PRO:HA	4:A:1286:HOH:O	2.17	0.44
1:B:332:TYR:CG	1:B:341:CYS:HB2	2.53	0.44
1:A:575:LYS:H	1:A:575:LYS:HG3	1.62	0.44
1:B:704:LEU:HD13	1:B:714:VAL:HG12	2.00	0.43
1:B:542:TYR:CZ	1:B:544:ASN:HB3	2.53	0.43
1:A:578:ASN:HB2	1:A:581:ASN:ND2	2.33	0.43
1:A:198:THR:O	1:A:200:TYR:N	2.52	0.42
1:B:647:THR:HG21	1:B:654:LYS:HE2	2.01	0.42
1:B:660:ASN:HB3	1:B:680:VAL:HG22	2.01	0.42
1:A:266:ASP:OD2	1:A:353:ARG:HB2	2.20	0.42
1:B:56:ALA:HB2	1:B:405:ILE:HD13	2.02	0.42
1:B:220:LYS:HE3	1:B:220:LYS:HB3	1.85	0.42
1:B:606:ARG:HG2	1:B:611:ASN:HA	2.02	0.42
1:B:95:ARG:HG3	1:B:108:ILE:HG22	2.02	0.41
1:B:406:ASN:OD1	1:B:439:LEU:HB2	2.21	0.41
1:B:501:GLU:HG3	4:B:1162:HOH:O	2.20	0.41
1:A:542:TYR:CZ	1:A:544:ASN:HB3	2.56	0.41
1:B:376:PRO:HD2	1:B:379:LYS:HG3	2.02	0.41
1:B:162:LEU:HD13	1:B:187:GLY:HA2	2.02	0.41
1:B:397:LYS:HE2	4:B:1321:HOH:O	2.19	0.41
1:B:800:TYR:CE1	1:B:804:VAL:HG21	2.56	0.41
1:A:561:PHE:HB3	1:A:563:THR:HG23	2.03	0.41
1:B:352:LEU:HD12	1:B:356:ALA:HB3	2.03	0.41
1:B:603:SER:OG	1:B:804:VAL:HA	2.21	0.41
1:A:545:LYS:HB3	1:A:545:LYS:HE2	1.84	0.41
1:B:682:PHE:O	1:B:688:LYS:HD2	2.21	0.40
1:B:519:ALA:HA	1:B:731:VAL:O	2.22	0.40
1:B:654:LYS:HD2	1:B:654:LYS:HA	1.95	0.40
1:A:657:TYR:CG	1:A:666:CYS:HB2	2.56	0.40
1:A:56:ALA:HB2	1:A:405:ILE:HD13	2.02	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	814/851 (96%)	778 (96%)	33 (4%)	3 (0%)	34	29
1	B	814/851 (96%)	774 (95%)	37 (4%)	3 (0%)	34	29
All	All	1628/1702 (96%)	1552 (95%)	70 (4%)	6 (0%)	34	29

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	639	ASP
1	B	301	VAL
1	A	199	GLY
1	B	611	ASN
1	B	617	PHE
1	A	690	PRO

### 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	704/729 (97%)	695 (99%)	9 (1%)	69	74
1	B	704/729 (97%)	701 (100%)	3 (0%)	91	93
All	All	1408/1458 (97%)	1396 (99%)	12 (1%)	78	83

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	111	CYS
1	B	352	LEU
1	B	615	ASP
1	A	132	VAL
1	A	158	LEU
1	A	575	LYS
1	A	576	ASP
1	A	610	ASP
1	A	642	ASN
1	A	649	CYS
1	A	676	VAL
1	A	688	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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](#)

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	CAC	A	903	-	0,4,4	-	-	0,6,6	-	-
3	CAC	B	904	-	0,4,4	-	-	0,6,6	-	-
3	CAC	A	904	-	0,4,4	-	-	0,6,6	-	-
3	CAC	B	902	-	0,4,4	-	-	0,6,6	-	-
2	9SL	B	901	-	17,23,23	3.41	8 (47%)	13,37,37	3.01	5 (38%)
3	CAC	A	902	-	0,4,4	-	-	0,6,6	-	-
3	CAC	B	903	-	0,4,4	-	-	0,6,6	-	-
2	9SL	A	901	-	17,23,23	3.36	7 (41%)	13,37,37	2.78	4 (30%)

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	9SL	A	901	-	-	2/5/53/53	0/3/3/3
2	9SL	B	901	-	-	4/5/53/53	0/3/3/3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	901	9SL	C12-N13	6.81	1.45	1.35
2	A	901	9SL	C12-N13	6.77	1.45	1.35
2	B	901	9SL	C02-N21	6.62	1.45	1.33
2	A	901	9SL	C02-N21	6.58	1.45	1.33
2	A	901	9SL	C07-N08	6.20	1.45	1.34
2	B	901	9SL	C07-N08	6.20	1.45	1.34
2	B	901	9SL	C12-N15	4.82	1.45	1.34
2	A	901	9SL	C12-N15	4.67	1.45	1.34
2	B	901	9SL	O03-C02	4.04	1.41	1.35
2	A	901	9SL	O03-C02	3.71	1.40	1.35
2	B	901	9SL	C05-C14	2.78	1.59	1.52
2	A	901	9SL	C05-C14	2.76	1.58	1.52
2	B	901	9SL	C07-N09	2.30	1.39	1.35
2	A	901	9SL	C07-N09	2.17	1.39	1.35
2	B	901	9SL	O01-C02	-2.01	1.19	1.21

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	901	9SL	O03-C02-N21	7.81	120.77	111.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	901	9SL	O03-C02-N21	7.15	119.96	111.08
2	B	901	9SL	N09-C07-N06	-4.27	119.48	125.42
2	A	901	9SL	N09-C07-N06	-4.05	119.79	125.42
2	B	901	9SL	O01-C02-N21	-3.77	119.28	125.51
2	A	901	9SL	O01-C02-N21	-3.36	119.97	125.51
2	B	901	9SL	O03-C02-O01	-3.34	119.92	123.07
2	A	901	9SL	O03-C02-O01	-3.17	120.08	123.07
2	B	901	9SL	C04-C05-N06	-2.28	105.09	108.57

There are no chirality outliers.

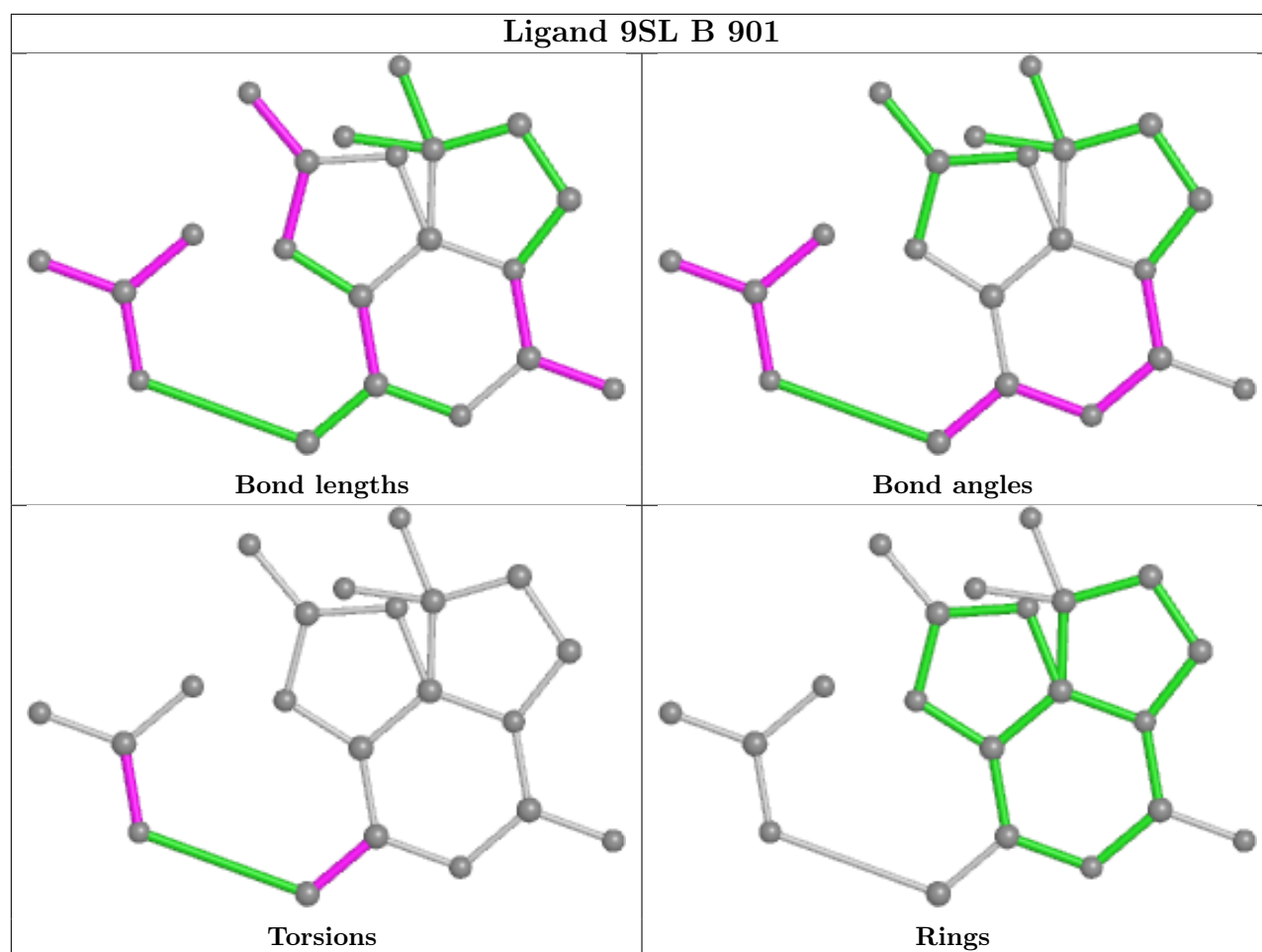
All (6) torsion outliers are listed below:

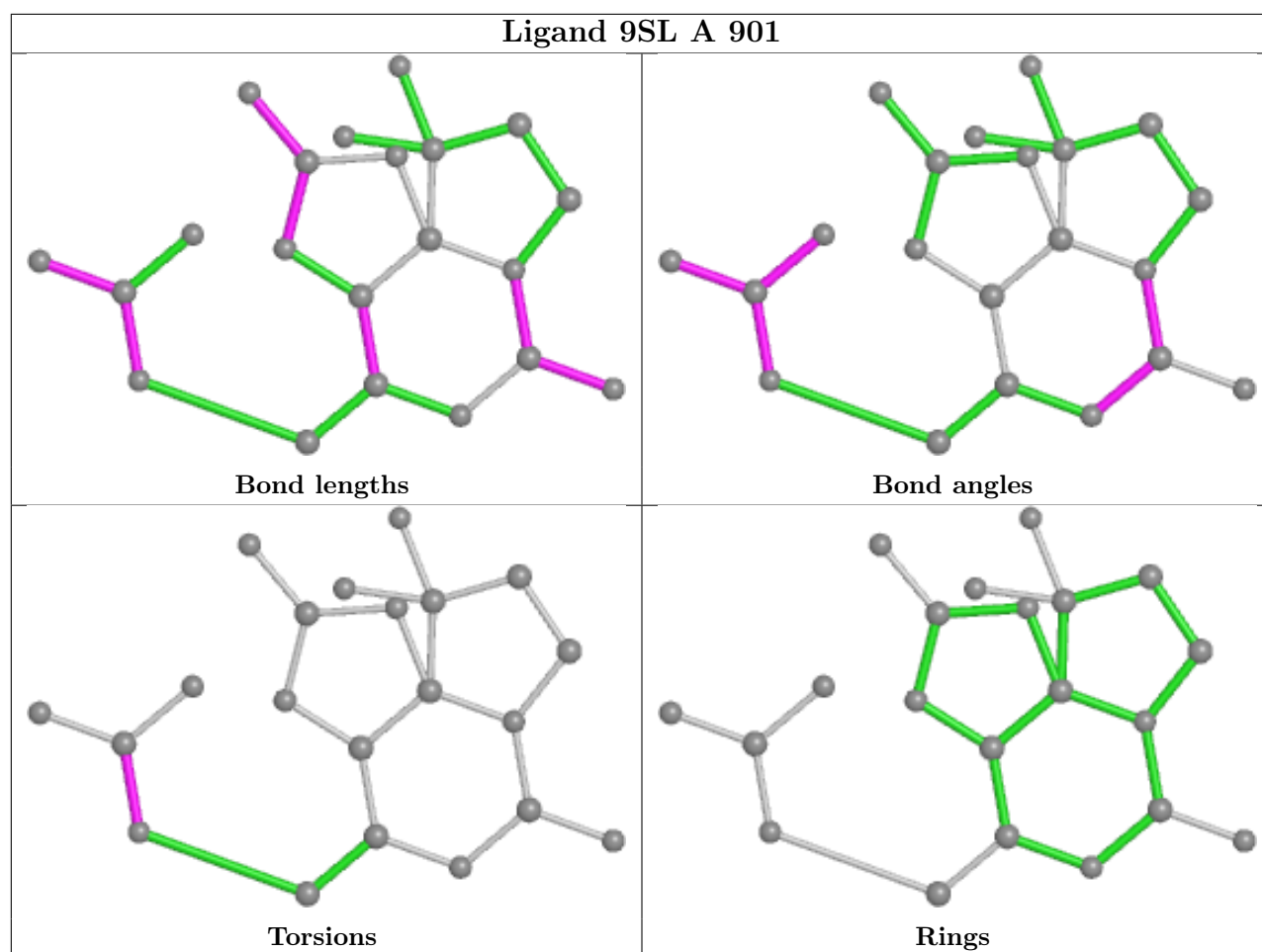
Mol	Chain	Res	Type	Atoms
2	B	901	9SL	O01-C02-O03-C04
2	B	901	9SL	N21-C02-O03-C04
2	B	901	9SL	O03-C04-C05-N06
2	B	901	9SL	O03-C04-C05-C14
2	A	901	9SL	O01-C02-O03-C04
2	A	901	9SL	N21-C02-O03-C04

There are no ring outliers.

No monomer is involved in short contacts.

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	818/851 (96%)	0.14	48 (5%)	22 30	36, 56, 115, 166	0
1	B	818/851 (96%)	0.44	67 (8%)	11 16	39, 64, 133, 191	0
All	All	1636/1702 (96%)	0.29	115 (7%)	16 22	36, 60, 124, 191	0

All (115) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	179	PHE	8.1
1	B	644	ALA	6.2
1	B	657	TYR	5.5
1	A	637	ILE	5.4
1	A	641	LYS	5.2
1	B	635	LEU	5.1
1	B	162	LEU	5.1
1	B	157	GLU	4.9
1	A	693	TRP	4.9
1	B	641	LYS	4.8
1	B	694	ALA	4.8
1	B	668	VAL	4.6
1	B	634	LYS	4.6
1	B	830	GLU	4.6
1	B	156	HIS	4.6
1	B	163	LYS	4.5
1	B	194	PHE	4.5
1	B	642	ASN	4.4
1	B	643	SER	4.4
1	B	646	ASN	4.2
1	A	167	VAL	4.1
1	A	644	ALA	4.1
1	B	199	GLY	4.0
1	B	202	TRP	4.0

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Mol	Chain	Res	Type	RSRZ
1	B	197	SER	4.0
1	B	166	GLN	3.9
1	B	631	ASN	3.8
1	A	582	ILE	3.8
1	A	197	SER	3.8
1	B	616	SER	3.7
1	B	629	LYS	3.7
1	A	691	ALA	3.6
1	B	831	VAL	3.5
1	B	829	LEU	3.5
1	B	664	PHE	3.4
1	B	620	GLU	3.4
1	B	640	PRO	3.4
1	B	613	ASP	3.3
1	B	611	ASN	3.3
1	B	707	LEU	3.3
1	A	685	THR	3.3
1	B	618	PHE	3.2
1	B	693	TRP	3.2
1	A	629	LYS	3.2
1	B	158	LEU	3.2
1	B	582	ILE	3.2
1	A	700	GLU	3.1
1	A	656	ALA	3.1
1	A	577	ILE	3.1
1	B	615	ASP	3.1
1	A	695	LYS	3.0
1	A	694	ALA	3.0
1	A	194	PHE	3.0
1	A	647	THR	3.0
1	B	654	LYS	2.9
1	A	602	VAL	2.9
1	B	205	ASN	2.9
1	B	670	LYS	2.9
1	B	832	LEU	2.9
1	B	200	TYR	2.9
1	A	157	GLU	2.9
1	A	575	LYS	2.9
1	A	635	LEU	2.8
1	B	215	LYS	2.8
1	A	832	LEU	2.8
1	A	586	LYS	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	199	GLY	2.7
1	B	666	CYS	2.7
1	A	611	ASN	2.7
1	A	620	GLU	2.7
1	B	614	ILE	2.7
1	B	218	PRO	2.7
1	A	665	ARG	2.7
1	B	180	VAL	2.7
1	A	689	ASN	2.7
1	A	675	PHE	2.7
1	B	586	LYS	2.7
1	B	639	ASP	2.7
1	B	692	VAL	2.7
1	A	714	VAL	2.7
1	A	576	ASP	2.7
1	A	162	LEU	2.5
1	A	670	LYS	2.5
1	A	676	VAL	2.5
1	A	640	PRO	2.5
1	A	674	ALA	2.4
1	A	657	TYR	2.4
1	A	285	GLY	2.4
1	A	570	VAL	2.3
1	A	654	LYS	2.3
1	A	696	ASN	2.3
1	B	288	LYS	2.3
1	B	637	ILE	2.3
1	B	690	PRO	2.3
1	B	647	THR	2.2
1	A	283	TRP	2.2
1	B	604	LEU	2.2
1	A	687	GLY	2.2
1	B	672	ASP	2.2
1	B	155	ARG	2.2
1	B	576	ASP	2.2
1	A	614	ILE	2.1
1	A	609	ASN	2.1
1	A	682	PHE	2.1
1	B	206	ALA	2.1
1	A	668	VAL	2.1
1	B	394	VAL	2.1
1	A	224	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	87	ASP	2.1
1	B	216	THR	2.1
1	B	525	GLN	2.0
1	B	696	ASN	2.0
1	B	81	PRO	2.0
1	B	645	ALA	2.0
1	B	673	VAL	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](#)

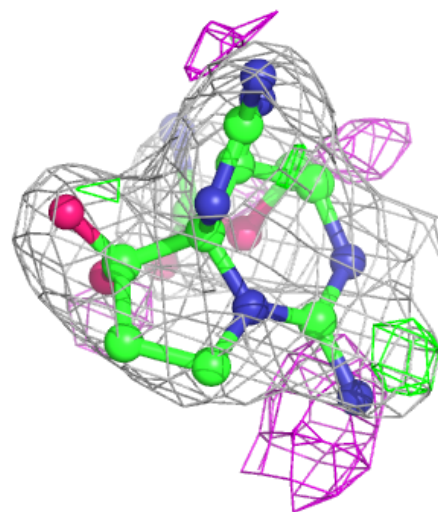
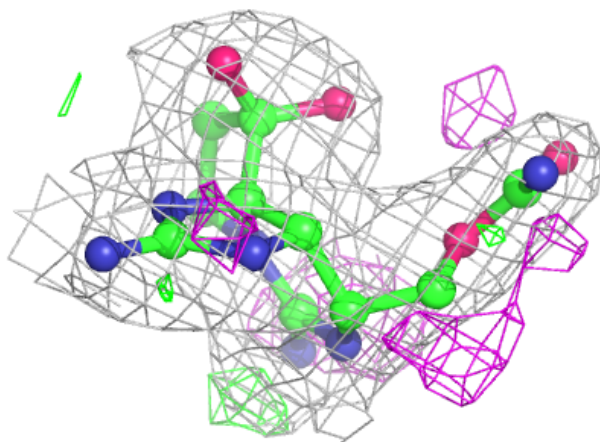
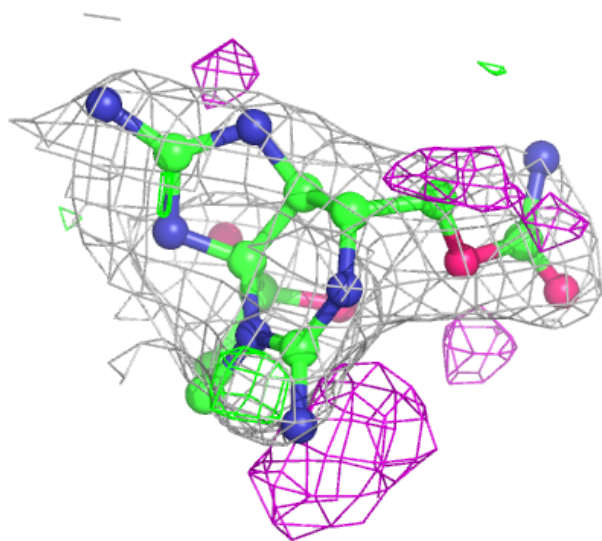
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	CAC	B	903	5/5	0.73	0.20	69,73,93,173	0
3	CAC	A	904	5/5	0.80	0.34	107,109,116,186	0
3	CAC	A	903	5/5	0.84	0.27	66,85,109,162	0
3	CAC	B	902	5/5	0.87	0.18	85,94,128,154	0
3	CAC	A	902	5/5	0.88	0.13	86,89,129,158	0
2	9SL	A	901	21/21	0.91	0.16	60,68,76,79	0
2	9SL	B	901	21/21	0.93	0.14	61,70,75,76	0
3	CAC	B	904	5/5	0.95	0.27	108,114,124,177	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.

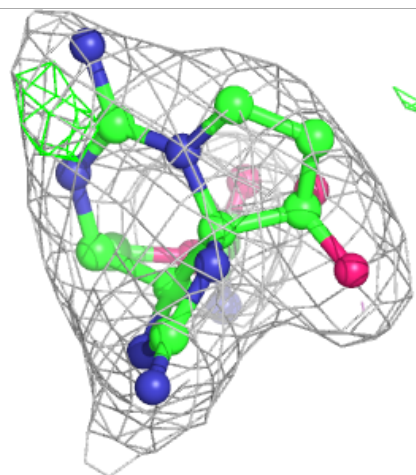
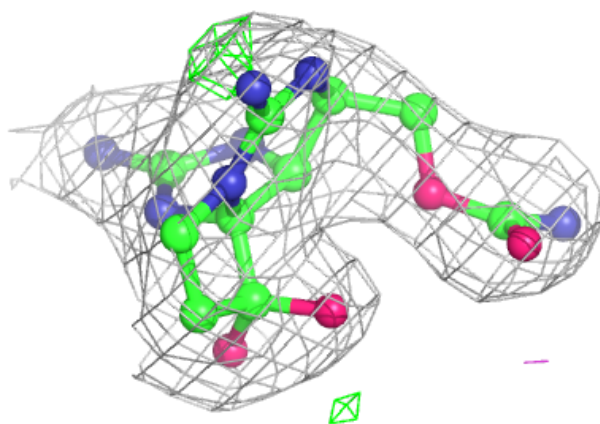
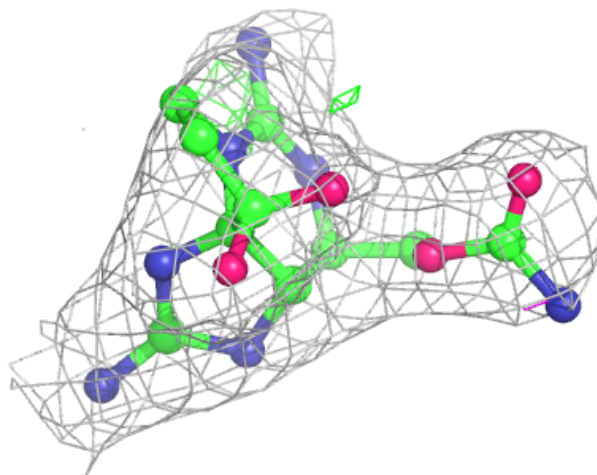
**Electron density around 9SL A 901:**

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



**Electron density around 9SL B 901:**

$2mF_o-DF_c$  (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.