



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:46 PM UTC

PDB ID : 1WLE / pdb_00001wle
Title : Crystal Structure of mammalian mitochondrial seryl-tRNA synthetase complexed with seryl-adenylate
Authors : Chimnaronk, S.; Jeppesen, M.G.; Suzuki, T.; Nyborg, J.; Watanabe, K.
Deposited on : 2004-06-25
Resolution : 1.65 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

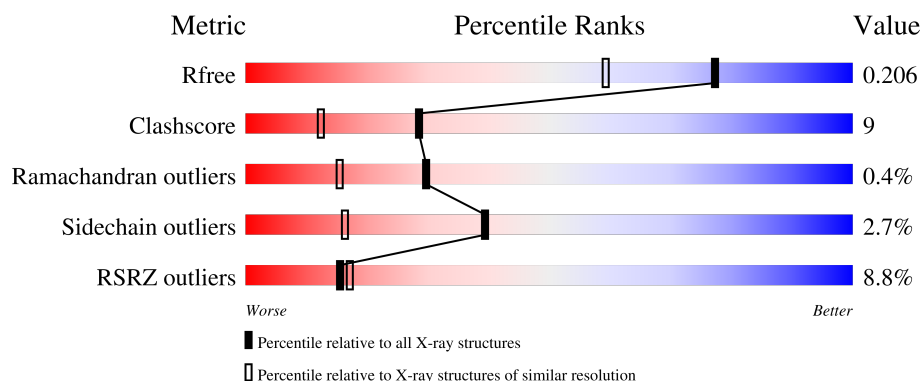
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	501	6% 77% 14% • 7%
1	B	501	10% 73% 18% • 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8106 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 Seryl-tRNA synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	465	Total	C	N	O	S	0	0	0
			3701	2337	661	687	16			
1	B	469	Total	C	N	O	S	0	0	0
			3736	2357	667	696	16			

There are 34 discrepancies between the modelled and reference sequences:

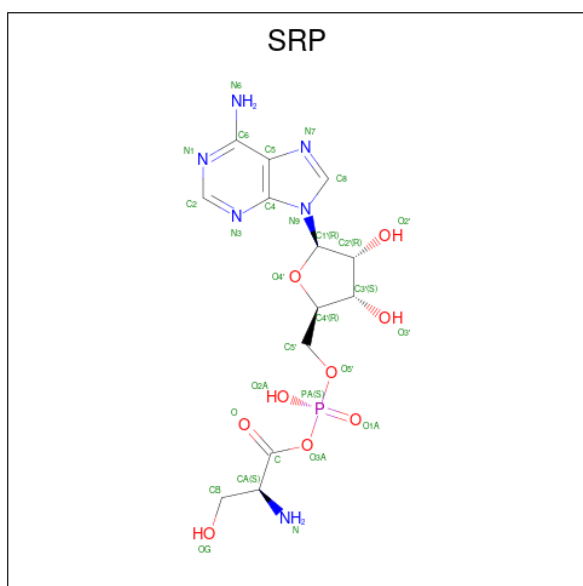
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q9N0F3
A	2	GLY	-	expression tag	UNP Q9N0F3
A	3	HIS	-	expression tag	UNP Q9N0F3
A	4	HIS	-	expression tag	UNP Q9N0F3
A	5	HIS	-	expression tag	UNP Q9N0F3
A	6	HIS	-	expression tag	UNP Q9N0F3
A	7	HIS	-	expression tag	UNP Q9N0F3
A	8	HIS	-	expression tag	UNP Q9N0F3
A	9	SER	-	expression tag	UNP Q9N0F3
A	10	SER	-	expression tag	UNP Q9N0F3
A	11	GLY	-	expression tag	UNP Q9N0F3
A	12	LEU	-	expression tag	UNP Q9N0F3
A	13	VAL	-	expression tag	UNP Q9N0F3
A	14	PRO	-	expression tag	UNP Q9N0F3
A	15	ARG	-	expression tag	UNP Q9N0F3
A	16	GLY	-	expression tag	UNP Q9N0F3
A	17	SER	-	expression tag	UNP Q9N0F3
B	1	MET	-	expression tag	UNP Q9N0F3
B	2	GLY	-	expression tag	UNP Q9N0F3
B	3	HIS	-	expression tag	UNP Q9N0F3
B	4	HIS	-	expression tag	UNP Q9N0F3
B	5	HIS	-	expression tag	UNP Q9N0F3
B	6	HIS	-	expression tag	UNP Q9N0F3
B	7	HIS	-	expression tag	UNP Q9N0F3
B	8	HIS	-	expression tag	UNP Q9N0F3

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	9	SER	-	expression tag	UNP Q9N0F3
B	10	SER	-	expression tag	UNP Q9N0F3
B	11	GLY	-	expression tag	UNP Q9N0F3
B	12	LEU	-	expression tag	UNP Q9N0F3
B	13	VAL	-	expression tag	UNP Q9N0F3
B	14	PRO	-	expression tag	UNP Q9N0F3
B	15	ARG	-	expression tag	UNP Q9N0F3
B	16	GLY	-	expression tag	UNP Q9N0F3
B	17	SER	-	expression tag	UNP Q9N0F3

- Molecule 2 is SERYL ADENYLATE (CCD ID: SRP) (formula: $C_{13}H_{19}N_6O_9P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			29	13	6	9	1		
2	B	1	Total	C	N	O	P	0	0
			29	13	6	9	1		

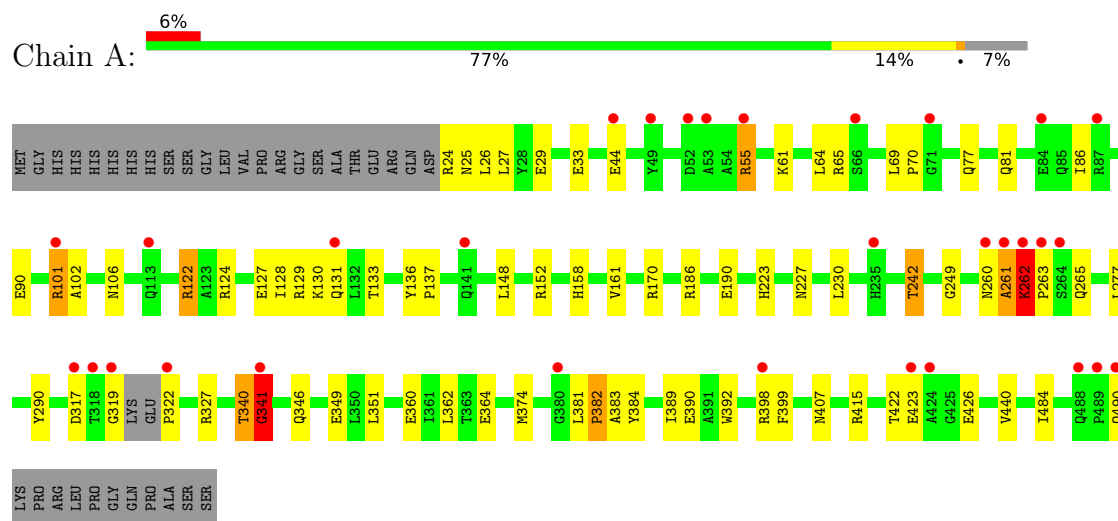
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	260	Total	O	0	0
			260	260		
3	B	351	Total	O	0	0
			351	351		

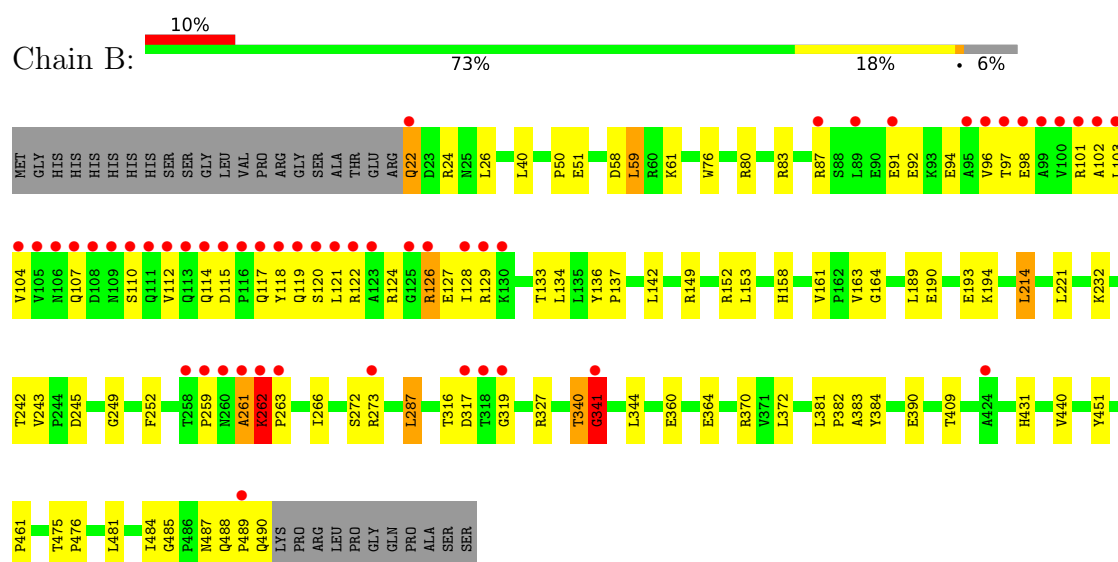
3 Residue-property plots [i](#)

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: Seryl-tRNA synthetase



• Molecule 1: Seryl-tRNA synthetase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	79.86Å 230.35Å 135.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.65 10.00 – 1.65	Depositor EDS
% Data completeness (in resolution range)	96.2 (10.00-1.65) 96.1 (10.00-1.65)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 1.64Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.210 , 0.226 0.212 , 0.206	Depositor DCC
R_{free} test set	14487 reflections (9.96%)	wwPDB-VP
Wilson B-factor (Å ²)	19.9	Xtriage
Anisotropy	0.247	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.47 , 52.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8106	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% 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: SRP

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	0/3783	0.89	11/5127 (0.2%)
1	B	0.36	0/3819	1.04	19/5177 (0.4%)
All	All	0.34	0/7602	0.97	30/10304 (0.3%)

There are no bond length outliers.

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	262	LYS	CA-C-N	-20.81	93.83	119.84
1	B	262	LYS	C-N-CA	-20.81	93.83	119.84
1	B	261	ALA	N-CA-C	-11.80	92.05	109.63
1	A	261	ALA	N-CA-C	-9.29	96.11	109.96
1	B	440	VAL	N-CA-C	8.47	122.52	112.35
1	B	104	VAL	N-CA-C	-8.26	105.86	113.71
1	A	440	VAL	N-CA-C	8.04	121.99	112.35
1	A	340	THR	N-CA-C	7.96	120.08	108.86
1	B	340	THR	N-CA-C	6.92	118.62	108.86
1	A	262	LYS	CA-C-N	-6.76	111.39	119.84
1	A	262	LYS	C-N-CA	-6.76	111.39	119.84
1	B	262	LYS	C-N-CD	6.43	151.38	125.00
1	A	374	MET	CA-C-N	6.17	126.13	120.21
1	A	374	MET	C-N-CA	6.17	126.13	120.21
1	B	341	GLY	N-CA-C	5.84	124.25	112.34
1	B	164	GLY	N-CA-C	5.80	119.20	110.97
1	B	263	PRO	N-CA-C	-5.79	100.55	112.47
1	A	341	GLY	N-CA-C	5.76	124.08	112.34
1	B	485	GLY	CA-C-N	5.74	125.11	118.85
1	B	485	GLY	C-N-CA	5.74	125.11	118.85
1	B	262	LYS	CB-CA-C	5.71	121.42	110.17
1	B	384	TYR	N-CA-C	-5.61	106.56	113.41

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	384	TYR	N-CA-C	-5.43	106.33	113.12
1	B	242	THR	N-CA-C	-5.38	100.25	109.24
1	B	383	ALA	N-CA-C	5.31	118.21	109.76
1	B	488	GLN	CA-C-N	5.28	125.53	119.83
1	B	488	GLN	C-N-CA	5.28	125.53	119.83
1	A	383	ALA	N-CA-C	5.15	118.34	110.20
1	B	344	LEU	N-CA-C	5.14	116.88	111.28
1	A	382	PRO	N-CA-C	5.09	121.25	114.18

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	3701	0	3670	67	0
1	B	3736	0	3701	70	0
2	A	29	0	18	0	0
2	B	29	0	18	0	0
3	A	260	0	0	4	0
3	B	351	0	0	4	0
All	All	8106	0	7407	131	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 9.

All (131) 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:130:LYS:O	1:A:133:THR:HG22	1.67	0.94
1:A:101:ARG:HH21	1:A:101:ARG:HB3	1.40	0.87
1:A:122:ARG:HB3	1:A:122:ARG:HH21	1.46	0.79
1:B:370:ARG:HD2	1:B:390:GLU:OE1	1.83	0.78
1:B:261:ALA:C	1:B:262:LYS:HD3	2.08	0.78
1:B:118:TYR:O	1:B:122:ARG:HB3	1.87	0.74

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:101:ARG:HB3	1:A:101:ARG:NH2	2.02	0.73
1:A:262:LYS:HD2	1:A:262:LYS:C	2.13	0.73
1:A:242:THR:HB	1:B:214:LEU:HB2	1.73	0.69
1:A:422:THR:OG1	1:A:426:GLU:HG2	1.94	0.68
1:B:58:ASP:O	1:B:61:LYS:HD2	1.96	0.66
1:B:94:GLU:O	1:B:98:GLU:HG2	1.95	0.66
1:B:120:SER:C	1:B:121:LEU:HD22	2.20	0.66
1:A:55:ARG:O	1:A:55:ARG:HD3	1.96	0.65
1:A:319:GLY:HA2	1:A:327:ARG:HD2	1.80	0.64
1:A:186:ARG:HD3	1:A:190:GLU:OE2	1.97	0.63
1:A:61:LYS:HB3	1:A:61:LYS:NZ	2.14	0.63
1:B:243:VAL:HG11	1:B:287:LEU:HD13	1.81	0.62
1:A:262:LYS:HB3	1:A:263:PRO:CD	2.30	0.62
1:B:340:THR:OG1	1:B:341:GLY:N	2.32	0.61
1:B:129:ARG:O	1:B:133:THR:HG23	2.00	0.60
1:B:26:LEU:HD13	1:B:249:GLY:HA3	1.84	0.60
1:A:263:PRO:HG2	3:A:1150:HOH:O	2.03	0.59
1:B:136:TYR:HB2	1:B:137:PRO:HD3	1.84	0.59
1:A:262:LYS:HB3	1:A:263:PRO:HD3	1.84	0.59
1:B:262:LYS:HD3	1:B:262:LYS:N	2.14	0.59
1:B:149:ARG:HH11	1:B:152:ARG:CZ	2.17	0.58
1:A:136:TYR:HB2	1:A:137:PRO:HD3	1.85	0.58
1:A:101:ARG:HD2	1:A:101:ARG:C	2.29	0.58
1:B:232:LYS:HE2	3:B:1238:HOH:O	2.03	0.58
1:B:122:ARG:C	1:B:122:ARG:HD3	2.29	0.58
1:A:340:THR:OG1	1:A:341:GLY:N	2.35	0.57
1:A:122:ARG:HH21	1:A:122:ARG:CB	2.16	0.56
1:B:107:GLN:O	1:B:112:VAL:HG23	2.04	0.56
1:B:259:PRO:HA	1:B:266:ILE:HD12	1.87	0.56
1:B:117:GLN:OE1	1:B:121:LEU:HG	2.06	0.56
1:B:252:PHE:HB2	1:B:259:PRO:HG3	1.87	0.56
1:B:272:SER:C	1:B:273:ARG:HD2	2.30	0.56
1:B:489:PRO:O	1:B:490:GLN:HG3	2.06	0.55
1:B:119:GLN:O	1:B:120:SER:HB2	2.07	0.54
1:B:94:GLU:O	1:B:97:THR:HG22	2.09	0.53
1:B:124:ARG:O	1:B:124:ARG:HD3	2.08	0.53
1:A:102:ALA:O	1:A:106:ASN:HB2	2.08	0.53
1:B:190:GLU:HG2	1:B:194:LYS:HE2	1.91	0.53
1:A:170:ARG:HD2	3:A:1076:HOH:O	2.08	0.53
1:A:341:GLY:HA3	1:A:346:GLN:HG2	1.91	0.53
1:B:149:ARG:HE	1:B:152:ARG:NH1	2.07	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:381:LEU:H	1:A:382:PRO:HD2	1.74	0.52
1:B:262:LYS:N	1:B:262:LYS:CD	2.73	0.52
1:A:24:ARG:HD3	1:A:29:GLU:OE1	2.11	0.51
1:B:94:GLU:C	1:B:97:THR:HG22	2.35	0.51
1:A:265:GLN:HB3	3:A:1037:HOH:O	2.11	0.51
1:A:290:TYR:CE2	1:B:481:LEU:HD11	2.46	0.51
1:A:86:ILE:O	1:A:90:GLU:HG3	2.11	0.50
1:B:26:LEU:HD12	1:B:259:PRO:HG2	1.92	0.50
1:B:381:LEU:H	1:B:382:PRO:HD2	1.76	0.50
1:A:55:ARG:HD3	1:A:55:ARG:C	2.37	0.50
1:A:64:LEU:HD22	1:A:64:LEU:N	2.27	0.50
1:B:80:ARG:NH2	1:B:83:ARG:HH21	2.10	0.50
1:B:149:ARG:HE	1:B:152:ARG:HH11	1.59	0.50
1:A:346:GLN:O	1:A:349:GLU:HG2	2.12	0.49
1:A:223:HIS:CE1	1:A:227:ASN:HD21	2.30	0.49
1:A:124:ARG:O	1:A:128:ILE:HG13	2.12	0.49
1:B:360:GLU:O	1:B:364:GLU:HG3	2.11	0.49
1:A:360:GLU:O	1:A:364:GLU:HG3	2.12	0.49
1:B:92:GLU:O	1:B:96:VAL:HG23	2.13	0.49
1:B:101:ARG:HD3	1:B:101:ARG:C	2.38	0.48
1:A:262:LYS:HE2	1:A:263:PRO:HD3	1.94	0.48
1:A:158:HIS:HB3	1:A:161:VAL:HG23	1.95	0.48
1:A:242:THR:CG2	1:B:214:LEU:HB2	2.44	0.48
1:A:381:LEU:N	1:A:382:PRO:HD2	2.29	0.48
1:B:124:ARG:O	1:B:128:ILE:HG13	2.13	0.47
1:A:262:LYS:HD2	1:A:263:PRO:N	2.28	0.47
1:A:262:LYS:CE	1:A:263:PRO:HD3	2.45	0.47
1:B:451:TYR:CE2	1:B:461:PRO:HG3	2.50	0.47
1:B:76:TRP:O	1:B:80:ARG:HG2	2.15	0.47
1:A:129:ARG:HG3	1:A:129:ARG:HH21	1.79	0.47
1:A:261:ALA:O	1:A:262:LYS:C	2.58	0.47
1:A:322:PRO:HG2	3:A:1064:HOH:O	2.13	0.46
1:A:26:LEU:HG	1:A:260:ASN:HD21	1.80	0.46
1:B:262:LYS:HZ3	1:B:262:LYS:H	1.62	0.46
1:A:242:THR:CB	1:B:214:LEU:HB2	2.43	0.45
1:A:260:ASN:O	1:A:261:ALA:C	2.59	0.45
1:A:44:GLU:OE1	1:A:44:GLU:HA	2.16	0.45
1:A:242:THR:HG21	3:B:1002:HOH:O	2.16	0.45
1:A:242:THR:HG22	3:B:1007:HOH:O	2.17	0.45
1:B:87:ARG:O	1:B:91:GLU:HG3	2.17	0.45
1:A:64:LEU:O	1:A:65:ARG:HD3	2.17	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:319:GLY:HA3	1:B:327:ARG:NH2	2.32	0.44
1:A:242:THR:CG2	3:B:1007:HOH:O	2.66	0.44
1:B:484:ILE:HD12	1:B:484:ILE:C	2.43	0.44
1:B:59:LEU:HG	1:B:163:VAL:HG13	2.01	0.43
1:A:77:GLN:O	1:A:81:GLN:HG3	2.18	0.43
1:A:415:ARG:HG3	1:A:415:ARG:HH21	1.82	0.43
1:B:126:ARG:HD2	1:B:127:GLU:N	2.33	0.43
1:A:351:LEU:HD22	1:A:407:ASN:HB2	1.99	0.43
1:A:362:LEU:CD1	1:A:389:ILE:HG21	2.49	0.43
1:B:126:ARG:HG2	1:B:126:ARG:HH21	1.84	0.43
1:A:340:THR:O	1:A:341:GLY:O	2.36	0.43
1:A:484:ILE:C	1:A:484:ILE:HD12	2.44	0.43
1:B:50:PRO:HG2	1:B:51:GLU:OE2	2.19	0.43
1:B:316:THR:O	1:B:317:ASP:CG	2.62	0.43
1:A:127:GLU:HG3	1:A:131:GLN:HE21	1.84	0.43
1:B:124:ARG:NH2	1:B:128:ILE:HG12	2.34	0.43
1:A:490:GLN:O	1:A:490:GLN:HG3	2.19	0.42
1:A:26:LEU:HD13	1:A:249:GLY:HA3	2.01	0.42
1:A:148:LEU:O	1:A:152:ARG:HD2	2.19	0.42
1:B:158:HIS:HB3	1:B:161:VAL:HG23	2.00	0.42
1:A:25:ASN:OD1	1:A:27:LEU:HB2	2.19	0.42
1:B:189:LEU:O	1:B:193:GLU:HG3	2.19	0.42
1:B:102:ALA:C	1:B:103:LEU:HD12	2.44	0.42
1:B:381:LEU:N	1:B:382:PRO:HD2	2.34	0.42
1:A:29:GLU:O	1:A:33:GLU:HG2	2.20	0.42
1:B:22:GLN:OE1	1:B:24:ARG:HG3	2.19	0.42
1:B:370:ARG:HD3	1:B:372:LEU:HD21	2.01	0.42
1:B:272:SER:O	1:B:273:ARG:HD2	2.20	0.41
1:B:59:LEU:HD12	1:B:59:LEU:HA	1.91	0.41
1:A:242:THR:HG21	1:B:214:LEU:HD22	2.01	0.41
1:A:392:TRP:CZ3	1:A:398:ARG:HA	2.55	0.41
1:A:242:THR:HB	1:B:214:LEU:CB	2.48	0.41
1:B:98:GLU:HA	1:B:98:GLU:OE2	2.19	0.41
1:B:134:LEU:C	1:B:137:PRO:HD2	2.45	0.41
1:A:69:LEU:HB2	1:A:70:PRO:HD3	2.03	0.41
1:A:319:GLY:HA2	1:A:327:ARG:CD	2.49	0.41
1:B:340:THR:O	1:B:341:GLY:O	2.39	0.41
1:B:451:TYR:CZ	1:B:461:PRO:HG3	2.55	0.41
1:B:475:THR:HA	1:B:476:PRO:HD3	1.86	0.41
1:B:262:LYS:N	1:B:262:LYS:HZ3	2.19	0.40
1:B:409:THR:O	1:B:431:HIS:HA	2.22	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:487:ASN:N	1:B:487:ASN:HD22	2.18	0.40
1:A:390:GLU:HB3	1:A:399:PHE:HB3	2.03	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	461/501 (92%)	448 (97%)	11 (2%)	2 (0%)	30	15
1	B	467/501 (93%)	442 (95%)	23 (5%)	2 (0%)	30	15
All	All	928/1002 (93%)	890 (96%)	34 (4%)	4 (0%)	30	15

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	341	GLY
1	B	341	GLY
1	B	110	SER
1	A	262	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.

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	400/430 (93%)	391 (98%)	9 (2%)	44	21
1	B	404/430 (94%)	391 (97%)	13 (3%)	34	12
All	All	804/860 (94%)	782 (97%)	22 (3%)	39	16

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	ARG
1	A	101	ARG
1	A	122	ARG
1	A	230	LEU
1	A	242	THR
1	A	262	LYS
1	A	277	LEU
1	A	317	ASP
1	A	423	GLU
1	B	22	GLN
1	B	40	LEU
1	B	59	LEU
1	B	114	GLN
1	B	115	ASP
1	B	126	ARG
1	B	142	LEU
1	B	153	LEU
1	B	214	LEU
1	B	221	LEU
1	B	245	ASP
1	B	262	LYS
1	B	287	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	85	GLN
1	A	111	GLN
1	A	131	GLN
1	A	141	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	145	GLN
1	A	205	HIS
1	A	223	HIS
1	A	235	HIS
1	A	260	ASN
1	A	329	HIS
1	A	434	ASN
1	B	77	GLN
1	B	81	GLN
1	B	85	GLN
1	B	113	GLN
1	B	119	GLN
1	B	141	GLN
1	B	145	GLN
1	B	184	GLN
1	B	222	GLN
1	B	223	HIS
1	B	377	GLN
1	B	482	GLN
1	B	487	ASN
1	B	490	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 oligosaccharides 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	SRP	B	901	-	29,31,31	1.73	5 (17%)	41,46,46	2.96	9 (21%)
2	SRP	A	900	-	29,31,31	1.74	5 (17%)	41,46,46	2.97	10 (24%)

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	SRP	B	901	-	-	4/19/37/37	0/3/3/3
2	SRP	A	900	-	-	4/19/37/37	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	900	SRP	C4-N3	4.77	1.43	1.34
2	B	901	SRP	O5'-C5'	-4.75	1.26	1.44
2	A	900	SRP	O5'-C5'	-4.73	1.26	1.44
2	B	901	SRP	C4-N3	4.48	1.42	1.34
2	B	901	SRP	PA-O5'	-3.49	1.45	1.59
2	A	900	SRP	PA-O5'	-3.34	1.46	1.59
2	A	900	SRP	C8-N9	2.15	1.41	1.37
2	B	901	SRP	PA-O3A	-2.13	1.56	1.60
2	B	901	SRP	O4'-C1'	2.12	1.46	1.42
2	A	900	SRP	PA-O1A	-2.02	1.43	1.50

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	901	SRP	O5'-C5'-C4'	13.23	154.03	108.99
2	A	900	SRP	O5'-C5'-C4'	13.07	153.49	108.99
2	A	900	SRP	O5'-PA-O1A	-6.73	82.24	108.94
2	B	901	SRP	O5'-PA-O1A	-6.57	82.91	108.94
2	A	900	SRP	PA-O5'-C5'	6.18	156.78	121.35
2	B	901	SRP	PA-O5'-C5'	6.00	155.71	121.35
2	A	900	SRP	C5'-C4'-C3'	-5.38	95.85	115.21

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	901	SRP	C5'-C4'-C3'	-4.95	97.40	115.21
2	B	901	SRP	O2A-PA-O3A	3.84	117.35	104.94
2	A	900	SRP	O2A-PA-O3A	3.77	117.12	104.94
2	B	901	SRP	O4'-C4'-C5'	-3.18	99.14	109.33
2	A	900	SRP	O4'-C4'-C5'	-3.13	99.30	109.33
2	B	901	SRP	O3A-PA-O5'	-3.06	93.80	103.00
2	A	900	SRP	O3A-PA-O1A	2.98	119.54	109.95
2	B	901	SRP	O3A-PA-O1A	2.83	119.05	109.95
2	A	900	SRP	O3A-PA-O5'	-2.80	94.60	103.00
2	B	901	SRP	O2A-PA-O1A	2.32	123.23	112.44
2	A	900	SRP	O2A-PA-O1A	2.28	123.04	112.44
2	A	900	SRP	C3'-C2'-C1'	2.15	105.53	101.46

There are no chirality outliers.

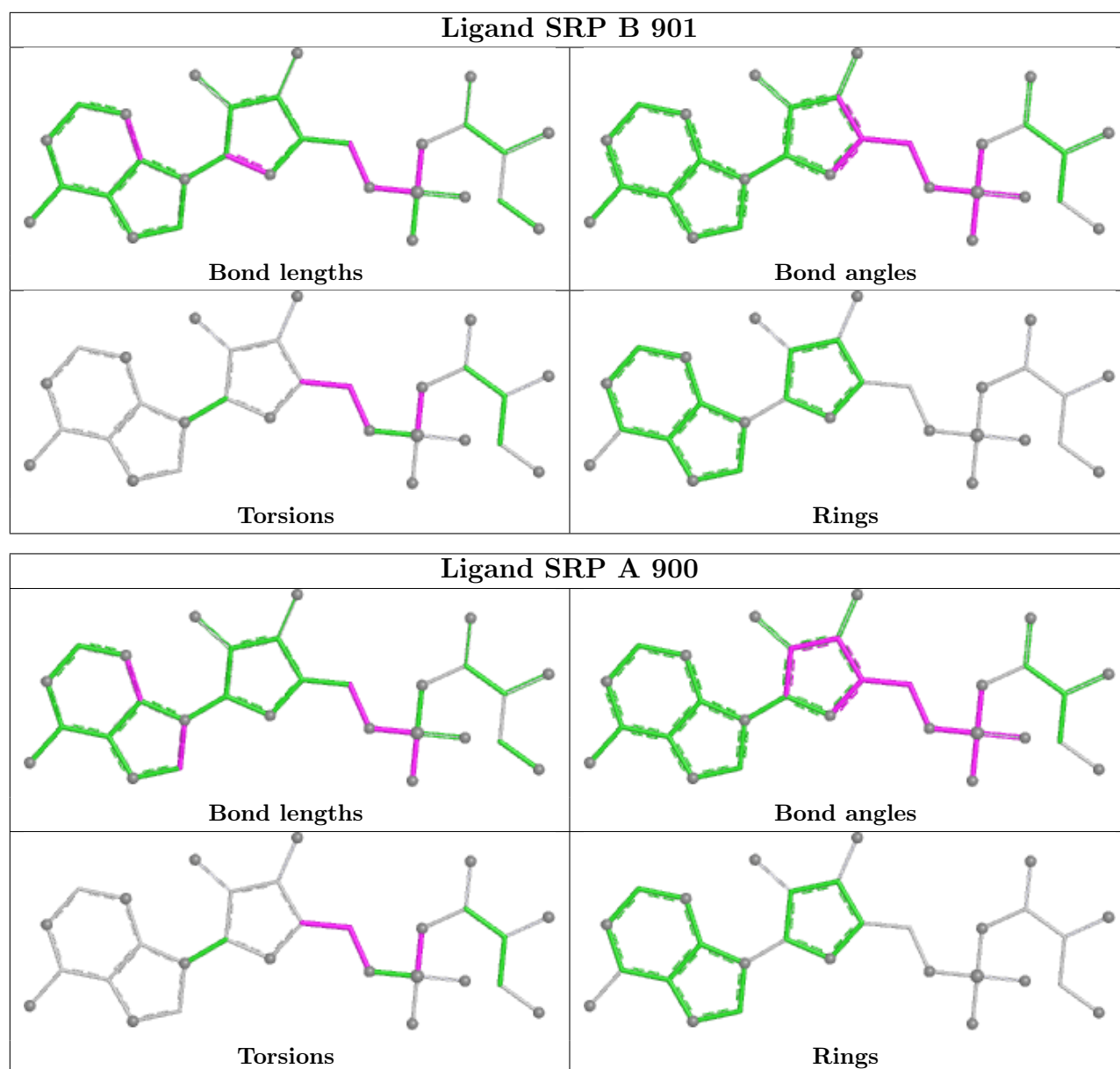
All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	901	SRP	C3'-C4'-C5'-O5'
2	A	900	SRP	O4'-C4'-C5'-O5'
2	A	900	SRP	C3'-C4'-C5'-O5'
2	B	901	SRP	O4'-C4'-C5'-O5'
2	A	900	SRP	C4'-C5'-O5'-PA
2	B	901	SRP	C4'-C5'-O5'-PA
2	B	901	SRP	C-O3A-PA-O1A
2	A	900	SRP	C-O3A-PA-O1A

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, 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	465/501 (92%)	0.35	31 (6%) 24 27	14, 25, 47, 62	0
1	B	469/501 (93%)	0.28	51 (10%) 10 11	12, 20, 86, 100	0
All	All	934/1002 (93%)	0.32	82 (8%) 15 17	12, 23, 55, 100	0

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	105	VAL	11.5
1	B	112	VAL	11.5
1	B	104	VAL	8.1
1	B	341	GLY	7.7
1	B	261	ALA	7.0
1	B	121	LEU	6.7
1	B	99	ALA	6.5
1	B	95	ALA	6.3
1	B	102	ALA	6.2
1	B	118	TYR	5.8
1	B	116	PRO	5.8
1	A	261	ALA	5.7
1	B	103	LEU	5.7
1	A	341	GLY	5.6
1	A	319	GLY	5.4
1	B	100	VAL	5.3
1	A	318	THR	5.3
1	B	489	PRO	5.1
1	B	109	ASN	4.9
1	B	120	SER	4.8
1	A	262	LYS	4.7
1	B	262	LYS	4.5
1	B	110	SER	4.5
1	B	111	GLN	4.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	263	PRO	4.1
1	B	97	THR	4.0
1	A	489	PRO	3.9
1	B	317	ASP	3.9
1	B	96	VAL	3.8
1	A	490	GLN	3.7
1	B	126	ARG	3.7
1	B	319	GLY	3.5
1	B	273	ARG	3.5
1	A	317	ASP	3.4
1	B	117	GLN	3.3
1	B	101	ARG	3.2
1	B	106	ASN	3.2
1	B	107	GLN	3.2
1	A	44	GLU	3.2
1	A	101	ARG	3.2
1	B	318	THR	3.2
1	B	119	GLN	3.1
1	B	130	LYS	3.1
1	A	87	ARG	3.1
1	B	260	ASN	3.1
1	B	123	ALA	3.0
1	B	115	ASP	3.0
1	B	108	ASP	3.0
1	B	259	PRO	3.0
1	A	264	SER	2.9
1	A	260	ASN	2.9
1	A	235	HIS	2.9
1	B	424	ALA	2.9
1	B	263	PRO	2.8
1	B	258	THR	2.8
1	B	113	GLN	2.8
1	B	122	ARG	2.7
1	A	398	ARG	2.7
1	A	49	TYR	2.6
1	A	113	GLN	2.6
1	A	66	SER	2.6
1	B	128	ILE	2.5
1	B	129	ARG	2.5
1	B	125	GLY	2.5
1	A	423	GLU	2.5
1	B	22	GLN	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	322	PRO	2.4
1	A	131	GLN	2.4
1	A	424	ALA	2.4
1	A	52	ASP	2.4
1	A	141	GLN	2.3
1	B	114	GLN	2.3
1	A	84	GLU	2.3
1	A	71	GLY	2.3
1	A	380	GLY	2.3
1	B	91	GLU	2.2
1	B	87	ARG	2.2
1	A	53	ALA	2.2
1	A	55	ARG	2.2
1	B	89	LEU	2.2
1	A	488	GLN	2.1
1	B	98	GLU	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

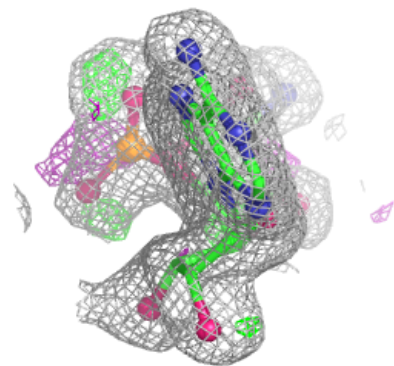
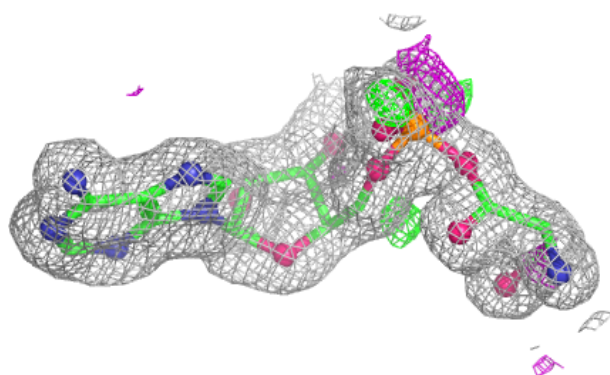
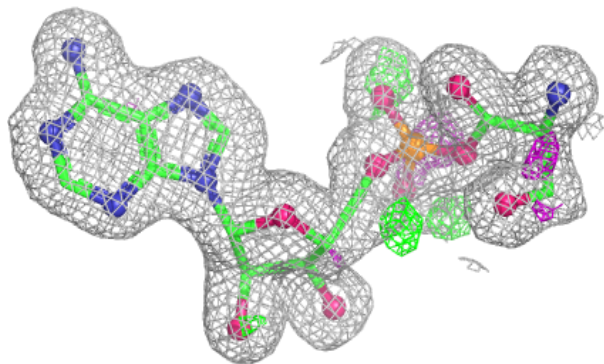
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SRP	A	900	29/29	0.93	0.07	16,21,25,27	0
2	SRP	B	901	29/29	0.94	0.06	14,17,22,22	0

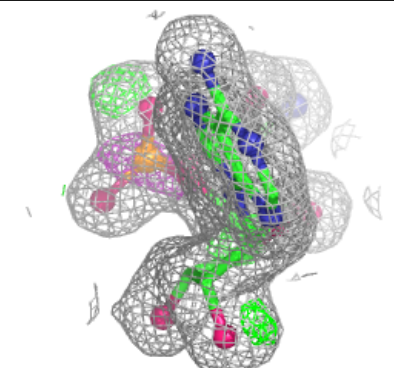
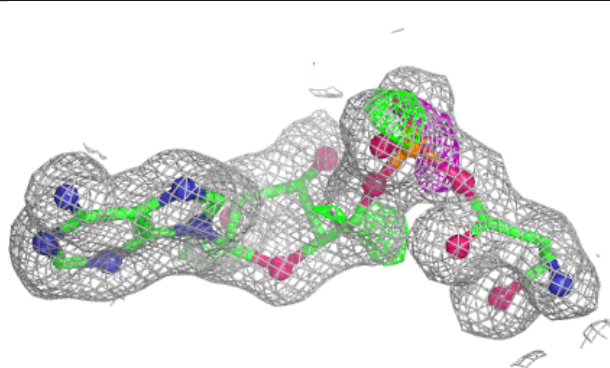
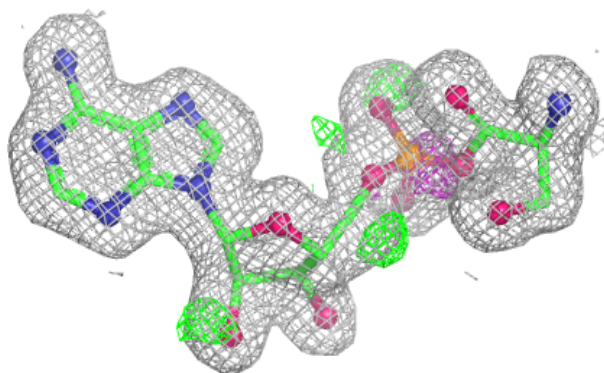
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around SRP A 900:

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

**Electron density around SRP 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.