



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 09:26 AM UTC

PDB ID : 2CJ9 / pdb_00002cj9
Title : Crystal structure of Methanosarcina barkeri seryl-tRNA synthetase complexed with an analog of seryladenylate
Authors : Bilokapic, S.; Maier, T.; Ahel, D.; Gruic-Sovulj, I.; Soll, D.; Weygand-Durasevic, I.; Ban, N.
Deposited on : 2006-03-29
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

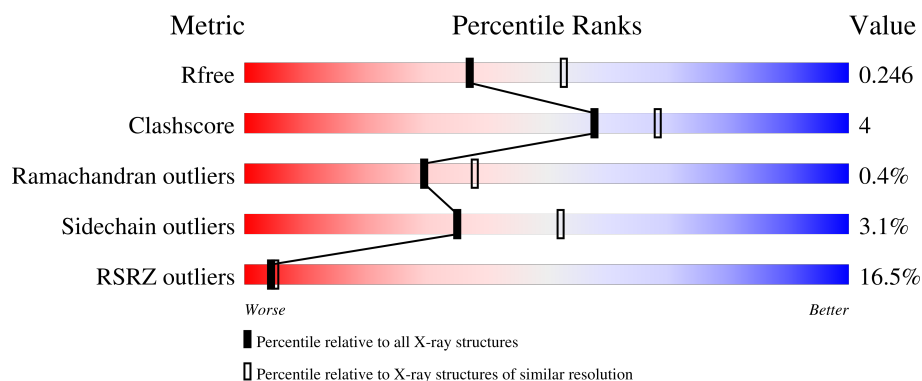
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	522	
1	B	522	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8307 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	477	Total	C	N	O	S	0	1	0
			3900	2505	663	711	21			
1	B	493	Total	C	N	O	S	0	1	0
			4015	2577	681	736	21			

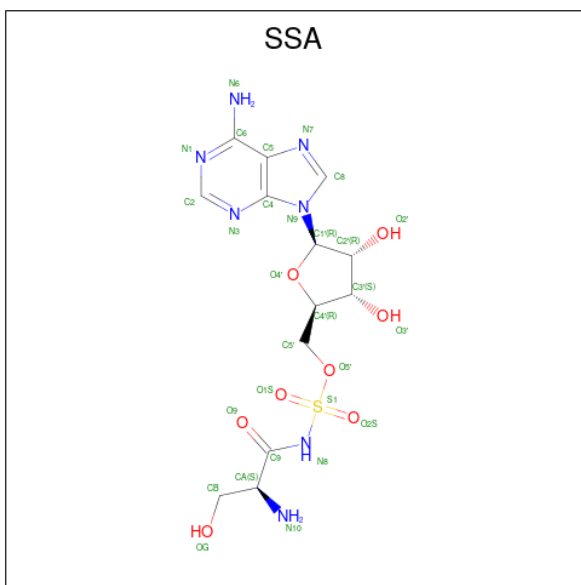
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	B	1	Total	Cl	0	0
			1	1		

- Molecule 4 is 5'-O-(N-(L-SERYL)-SULFAMOYL)ADENOSINE (CCD ID: SSA) (formula: C₁₃H₁₉N₇O₈S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 29	C 13	N 7	O 8	S 1	0	0
4	B	1	Total 29	C 13	N 7	O 8	S 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	149	Total O 149 149	0	0
5	B	181	Total O 181 181	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	96.71Å 96.71Å 270.30Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	14.93 – 2.30 14.93 – 2.30	Depositor EDS
% Data completeness (in resolution range)	78.8 (14.93-2.30) 78.4 (14.93-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.72 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.206 , 0.248 0.208 , 0.246	Depositor DCC
R_{free} test set	2567 reflections (3.90%)	wwPDB-VP
Wilson B-factor (Å ²)	49.0	Xtriage
Anisotropy	0.090	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 59.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.037 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	8307	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% 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: ZN, SSA, CL

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	0.48	0/4003	0.76	1/5411 (0.0%)
1	B	0.47	0/4123	0.79	4/5579 (0.1%)
All	All	0.47	0/8126	0.78	5/10990 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	332	GLY	N-CA-C	5.85	118.76	110.74
1	B	143	ARG	CA-C-N	5.81	123.90	120.24
1	B	143	ARG	C-N-CA	5.81	123.90	120.24
1	B	307	PRO	N-CA-C	5.77	117.74	110.70
1	A	307	PRO	N-CA-C	5.76	117.73	110.70

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	3900	0	3816	41	0
1	B	4015	0	3929	40	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	29	0	19	0	0
4	B	29	0	19	0	0
5	A	149	0	0	4	0
5	B	181	0	0	1	0
All	All	8307	0	7783	71	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 4.

All (71) 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:301[B]:MET:HE2	1:B:301[B]:MET:HE2	1.34	1.07
1:B:16:ASP:HB3	1:B:56:ASP:HB2	1.65	0.78
1:A:197:ARG:HH12	1:A:203:GLN:H	1.39	0.71
1:A:301[B]:MET:CE	1:B:301[B]:MET:HE2	2.18	0.70
1:A:232:ARG:HD3	5:A:2049:HOH:O	1.90	0.70
1:A:450:LYS:NZ	5:A:2126:HOH:O	2.29	0.65
1:B:131:LEU:HD11	1:B:152:LEU:HD22	1.79	0.63
1:B:259:TYR:HB3	1:B:301[A]:MET:HE2	1.83	0.61
1:B:10:TYR:HB3	1:B:59:GLU:HB3	1.82	0.61
1:A:233:GLU:OE2	1:A:330:ARG:HD2	2.02	0.59
1:B:18:THR:HB	1:B:19:PRO:CD	2.33	0.59
1:B:18:THR:HB	1:B:19:PRO:HD2	1.86	0.58
1:B:306:CYS:SG	1:B:307:PRO:HD3	2.44	0.58
1:A:397:PHE:O	1:A:401:GLU:HG2	2.02	0.57
1:A:40:ALA:HB3	1:A:41:PRO:HD3	1.86	0.57
1:A:101:PHE:HE2	1:A:103:ILE:HD11	1.69	0.57
1:A:301[B]:MET:HE2	1:B:301[B]:MET:CE	2.22	0.56
1:A:8:LYS:HB3	1:A:99:GLU:HB3	1.88	0.55
1:A:3:LEU:HD11	1:A:149:LEU:HD11	1.88	0.54
1:A:110:GLU:HB2	1:A:124:ASN:HD22	1.73	0.53
1:B:325:VAL:O	1:B:358:TRP:HA	2.09	0.53
1:B:1:MET:HE3	5:B:2005:HOH:O	2.08	0.52
1:A:243:GLU:HA	1:A:246:MET:HE3	1.92	0.52
1:A:116:VAL:HB	1:A:119:ILE:HD12	1.93	0.50
1:B:225:LEU:HD21	1:B:372:LEU:CD2	2.41	0.50
1:A:369:ALA:HB2	1:A:438:ILE:HD11	1.93	0.50
1:B:121:SER:HB3	1:B:132:GLU:HB3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:VAL:O	1:A:49:THR:HG23	2.13	0.49
1:A:187:GLN:NE2	5:A:2032:HOH:O	2.46	0.49
1:A:258:ILE:O	1:B:264:PRO:HD3	2.13	0.49
1:B:320:ASN:HD21	1:B:453:SER:H	1.60	0.49
1:A:8:LYS:HD2	1:A:61:THR:HG23	1.94	0.48
1:A:186:THR:HA	1:A:189:MET:HE3	1.96	0.48
1:A:118:TYR:HB2	1:A:148:ILE:HD11	1.96	0.47
1:A:306:CYS:SG	1:A:307:PRO:HD3	2.54	0.47
1:B:173:GLN:HG3	1:B:389:ARG:HB3	1.97	0.47
1:A:395:PRO:HG3	1:A:413:GLY:HA2	1.96	0.46
1:A:103:ILE:HG22	1:A:105:VAL:HG23	1.98	0.46
1:B:243:GLU:HG2	1:B:311:MET:HE3	1.97	0.45
1:B:380:PHE:CZ	1:B:433:PHE:HE1	2.34	0.45
1:A:237:PRO:HB3	1:B:203:GLN:CD	2.41	0.45
1:B:366:LEU:HD21	1:B:411:THR:OG1	2.16	0.45
1:A:285:VAL:HG11	1:B:242:TRP:HB3	1.99	0.45
1:A:125:ILE:HG13	1:A:128:GLY:H	1.82	0.44
1:B:19:PRO:HA	1:B:53:LEU:HD22	1.99	0.44
1:A:173:GLN:HG3	1:A:389:ARG:HB3	2.00	0.44
1:B:420:CYS:HB3	1:B:430:TRP:CE2	2.53	0.44
1:A:225:LEU:HD21	1:A:372:LEU:CD2	2.48	0.44
1:A:49:THR:HG21	1:A:61:THR:O	2.18	0.43
1:A:49:THR:OG1	1:A:62:LEU:HA	2.18	0.43
1:B:5:PHE:HB3	1:B:70:VAL:HG22	2.00	0.43
1:B:259:TYR:CB	1:B:301[A]:MET:HE2	2.47	0.43
1:A:339:SER:HB3	1:B:267:ARG:NH1	2.34	0.43
1:B:335:HIS:HA	1:B:349:ASP:O	2.18	0.43
1:A:206:HIS:HB2	1:B:235:ILE:HB	2.01	0.42
1:A:74:ILE:HD12	1:A:96:ILE:HG21	2.01	0.42
1:B:12:LYS:HG2	1:B:95:GLY:O	2.19	0.42
1:B:395:PRO:HG3	1:B:413:GLY:HA2	2.01	0.42
1:B:4:GLN:HG2	1:B:104:LYS:HB2	2.02	0.42
1:B:135:VAL:HG13	1:B:139:GLU:HB2	2.02	0.42
1:B:243:GLU:HG2	1:B:311:MET:CE	2.50	0.42
1:A:147:ARG:NE	1:B:281:VAL:O	2.53	0.41
1:A:330:ARG:NH1	5:A:2087:HOH:O	2.53	0.41
1:A:306:CYS:N	1:A:307:PRO:CD	2.82	0.41
1:A:369:ALA:CB	1:A:438:ILE:HD11	2.50	0.41
1:B:451:LEU:HB2	1:B:457:LEU:HG	2.02	0.41
1:A:36:LEU:HD23	1:A:76:ARG:HD2	2.02	0.41
1:A:279:TYR:CE1	1:B:246:MET:HG2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:353:ARG:HA	1:B:464:VAL:O	2.21	0.41
1:B:433:PHE:HA	1:B:468:ARG:HB3	2.03	0.41
1:B:254:VAL:HG13	1:B:258:ILE:HD13	2.02	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	468/522 (90%)	445 (95%)	21 (4%)	2 (0%)	30	38
1	B	488/522 (94%)	472 (97%)	14 (3%)	2 (0%)	30	38
All	All	956/1044 (92%)	917 (96%)	35 (4%)	4 (0%)	30	38

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	253	GLY
1	B	253	GLY
1	A	127	GLY
1	B	18	THR

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	410/444 (92%)	396 (97%)	14 (3%)	32	49
1	B	422/444 (95%)	410 (97%)	12 (3%)	38	56
All	All	832/888 (94%)	806 (97%)	26 (3%)	35	52

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	24	ILE
1	A	49	THR
1	A	74	ILE
1	A	81	LEU
1	A	110	GLU
1	A	197	ARG
1	A	215	ARG
1	A	239	LEU
1	A	285	VAL
1	A	327	VAL
1	A	330	ARG
1	A	410	ASN
1	A	489	ARG
1	B	7	LEU
1	B	16	ASP
1	B	48	VAL
1	B	60	LEU
1	B	99	GLU
1	B	112	ARG
1	B	135	VAL
1	B	170	LEU
1	B	363	GLU
1	B	367	LYS
1	B	393	VAL
1	B	494	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	GLN
1	A	71	HIS
1	A	80	GLN
1	A	167	HIS

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Mol	Chain	Res	Type
1	A	283	HIS
1	B	320	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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
4	SSA	B	1505	2	30,31,31	1.46	5 (16%)	41,46,46	2.09	9 (21%)
4	SSA	A	1505	2	30,31,31	1.48	4 (13%)	41,46,46	2.27	11 (26%)

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
4	SSA	B	1505	2	-	4/20/37/37	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SSA	A	1505	2	-	2/20/37/37	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1505	SSA	C5-C4	4.87	1.47	1.39
4	B	1505	SSA	C5-C4	4.64	1.47	1.39
4	A	1505	SSA	C5-C6	2.78	1.48	1.41
4	B	1505	SSA	C5-C6	2.56	1.48	1.41
4	A	1505	SSA	C8-N7	2.46	1.36	1.31
4	B	1505	SSA	C4-N9	-2.34	1.32	1.37
4	A	1505	SSA	C9-N8	-2.29	1.33	1.37
4	B	1505	SSA	C5-N7	-2.10	1.35	1.39
4	B	1505	SSA	C8-N7	2.07	1.35	1.31

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1505	SSA	O2S-S1-O1S	-8.41	108.26	120.85
4	B	1505	SSA	O2S-S1-O1S	-7.21	110.05	120.85
4	A	1505	SSA	C5-C4-N3	-5.60	119.01	126.72
4	B	1505	SSA	C5-C4-N3	-5.04	119.77	126.72
4	A	1505	SSA	N3-C4-N9	4.58	134.95	127.17
4	B	1505	SSA	N3-C4-N9	4.37	134.60	127.17
4	B	1505	SSA	N3-C2-N1	-3.77	122.88	128.58
4	A	1505	SSA	C2-N3-C4	3.66	120.77	111.83
4	B	1505	SSA	C2-N3-C4	3.56	120.51	111.83
4	A	1505	SSA	N3-C2-N1	-3.49	123.30	128.58
4	B	1505	SSA	C4-N9-C8	3.26	109.16	105.74
4	A	1505	SSA	C4-C5-N7	-3.22	106.90	110.58
4	B	1505	SSA	C4-C5-N7	-2.81	107.37	110.58
4	A	1505	SSA	C4-N9-C8	2.77	108.64	105.74
4	A	1505	SSA	C5-N7-C8	2.52	107.41	103.45
4	B	1505	SSA	C2-N1-C6	2.38	122.64	118.73
4	A	1505	SSA	C6-C5-N7	2.16	136.26	132.09
4	B	1505	SSA	C5-N7-C8	2.09	106.73	103.45
4	A	1505	SSA	N9-C8-N7	-2.06	111.01	113.94
4	A	1505	SSA	C2'-C1'-N9	-2.04	108.23	113.30

There are no chirality outliers.

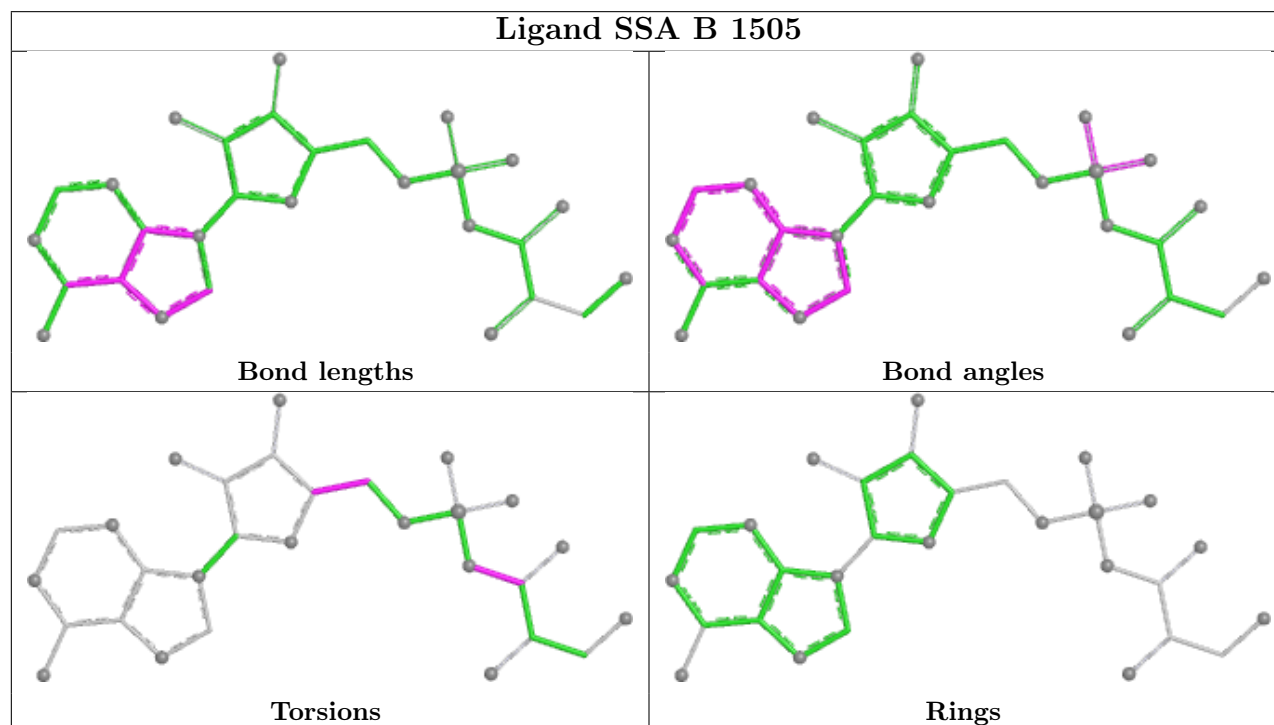
All (6) torsion outliers are listed below:

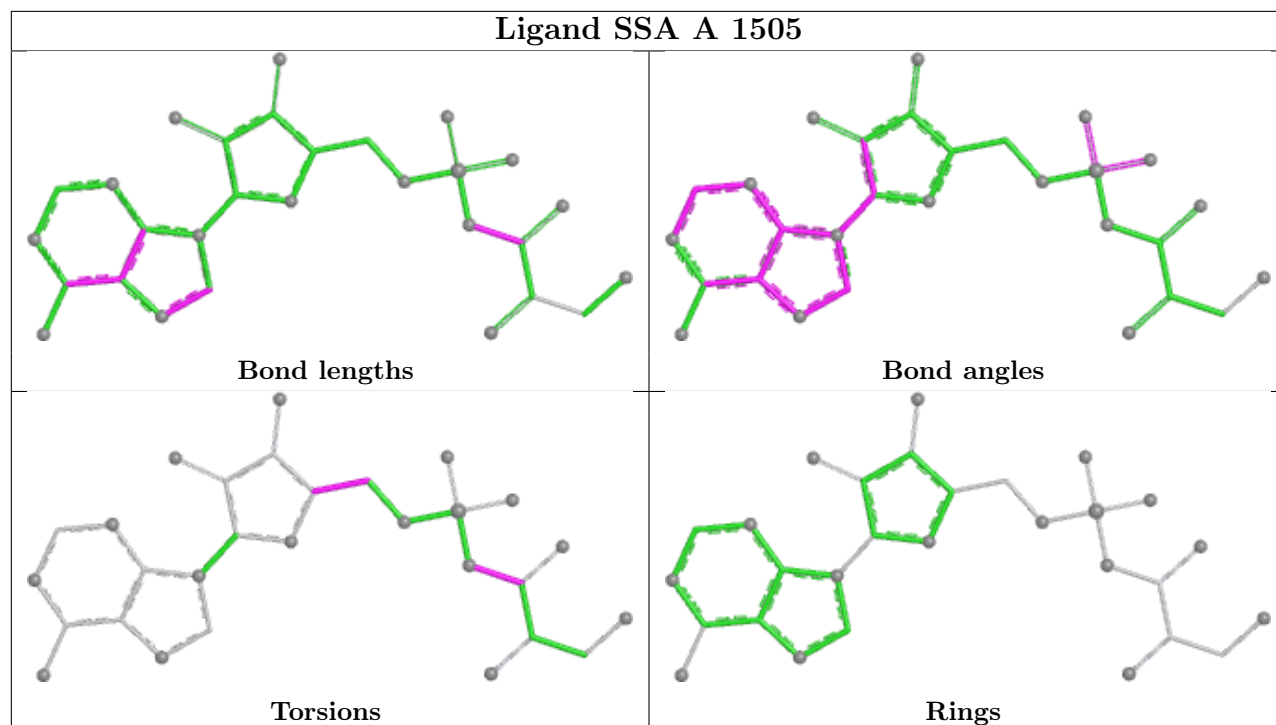
Mol	Chain	Res	Type	Atoms
4	A	1505	SSA	O9-C9-N8-S1
4	B	1505	SSA	CA-C9-N8-S1
4	B	1505	SSA	O9-C9-N8-S1
4	B	1505	SSA	O4'-C4'-C5'-O5'
4	B	1505	SSA	C3'-C4'-C5'-O5'
4	A	1505	SSA	C3'-C4'-C5'-O5'

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	477/522 (91%)	1.41	128 (26%) 1 1	30, 66, 137, 145	1 (0%)
1	B	493/522 (94%)	0.78	32 (6%) 25 27	31, 64, 95, 104	1 (0%)
All	All	970/1044 (92%)	1.09	160 (16%) 4 5	30, 65, 118, 145	2 (0%)

All (160) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	107	ALA	9.0
1	A	68	VAL	7.3
1	A	159	ALA	6.4
1	A	32	ASN	6.2
1	A	73	ALA	5.9
1	A	158	ALA	5.9
1	A	23	ALA	5.7
1	A	20	ALA	5.7
1	A	126	GLU	5.4
1	A	43	GLY	5.2
1	B	161	TYR	5.0
1	A	65	GLY	4.8
1	A	95	GLY	4.6
1	A	128	GLY	4.6
1	A	67	TYR	4.6
1	A	26	ALA	4.5
1	A	31	ALA	4.5
1	A	46	ALA	4.5
1	A	109	HIS	4.5
1	A	165	ALA	4.5
1	A	51	TRP	4.4
1	B	165	ALA	4.4
1	B	482	ALA	4.3
1	A	3	LEU	4.3

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Mol	Chain	Res	Type	RSRZ
1	A	22	ASP	4.3
1	A	35	LEU	4.3
1	A	82	ALA	4.2
1	A	24	ILE	4.1
1	A	83	GLU	4.1
1	A	150	THR	4.0
1	A	72	ASP	4.0
1	A	-1	HIS	4.0
1	A	149	LEU	3.9
1	A	103	ILE	3.9
1	A	61	THR	3.9
1	A	71	HIS	3.8
1	A	-2	SER	3.8
1	A	70	VAL	3.8
1	A	144	VAL	3.8
1	A	106	PRO	3.8
1	A	131	LEU	3.8
1	B	166	GLU	3.7
1	B	123	GLU	3.6
1	A	111	LEU	3.6
1	A	96	ILE	3.6
1	A	76	ARG	3.6
1	A	138	ALA	3.6
1	A	129	ILE	3.5
1	B	404	LEU	3.5
1	A	405	GLY	3.4
1	B	167	HIS	3.4
1	A	44	GLN	3.4
1	A	275	VAL	3.3
1	A	151	LEU	3.3
1	B	27	LEU	3.3
1	A	58	ILE	3.3
1	A	102	ILE	3.3
1	A	9	ALA	3.3
1	B	406	LEU	3.3
1	A	119	ILE	3.2
1	B	14	SER	3.2
1	A	21	LYS	3.1
1	A	28	PHE	3.1
1	A	1	MET	3.1
1	A	48	VAL	3.1
1	A	135	VAL	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	42	GLU	3.0
1	A	167	HIS	2.9
1	A	2	LYS	2.9
1	B	410	ASN	2.9
1	A	10	TYR	2.9
1	A	80	GLN	2.9
1	B	53	LEU	2.8
1	A	166	GLU	2.8
1	A	34	THR	2.8
1	A	157	GLU	2.8
1	A	301[A]	MET	2.8
1	A	60	LEU	2.7
1	A	114	LEU	2.7
1	A	75	PHE	2.7
1	A	394	THR	2.7
1	A	272	TRP	2.7
1	A	40	ALA	2.7
1	A	156	ILE	2.7
1	A	33	SER	2.7
1	A	121	SER	2.7
1	A	395	PRO	2.7
1	A	39	GLY	2.6
1	A	404	LEU	2.6
1	A	396	TRP	2.6
1	A	101	PHE	2.6
1	A	104	LYS	2.6
1	A	276	ALA	2.6
1	A	489	ARG	2.6
1	B	289	LEU	2.6
1	B	138	ALA	2.6
1	A	41	PRO	2.6
1	A	143	ARG	2.5
1	B	20	ALA	2.5
1	B	82	ALA	2.5
1	A	49	THR	2.5
1	A	130	GLN	2.5
1	A	281	VAL	2.5
1	A	27	LEU	2.5
1	A	133	LEU	2.5
1	B	405	GLY	2.5
1	A	64	SER	2.5
1	B	403	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	56	ASP	2.5
1	B	75	PHE	2.4
1	A	385	ASP	2.4
1	A	407	ALA	2.4
1	B	19	PRO	2.4
1	A	125	ILE	2.4
1	A	77	LEU	2.4
1	B	109	HIS	2.4
1	A	178	GLU	2.4
1	A	100	SER	2.4
1	A	105	VAL	2.4
1	B	160	GLN	2.4
1	A	45	GLY	2.3
1	B	253	GLY	2.3
1	A	57	ARG	2.3
1	A	406	LEU	2.3
1	A	262	CYS	2.3
1	A	5	PHE	2.3
1	B	-3	GLY	2.3
1	B	408	GLU	2.3
1	A	7	LEU	2.2
1	B	396	TRP	2.2
1	A	141	LYS	2.2
1	A	11	PHE	2.2
1	A	108	ASP	2.2
1	A	277	ASP	2.2
1	A	63	GLN	2.2
1	B	40	ALA	2.2
1	A	116	VAL	2.2
1	A	4	GLN	2.2
1	A	482	ALA	2.2
1	B	425	GLY	2.1
1	B	394	THR	2.1
1	A	118	TYR	2.1
1	A	81	LEU	2.1
1	A	403	LEU	2.1
1	A	50	GLU	2.1
1	A	74	ILE	2.1
1	A	124	ASN	2.1
1	A	320	ASN	2.1
1	A	299	GLY	2.1
1	A	321	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	6	ASN	2.1
1	A	146	ASP	2.1
1	B	383	ILE	2.1
1	A	297	PRO	2.1
1	B	70	VAL	2.1
1	A	148	ILE	2.0
1	A	294	ILE	2.0
1	B	275	VAL	2.0
1	A	78	ARG	2.0
1	A	500	ARG	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](#)

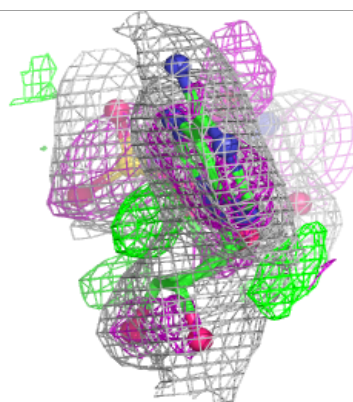
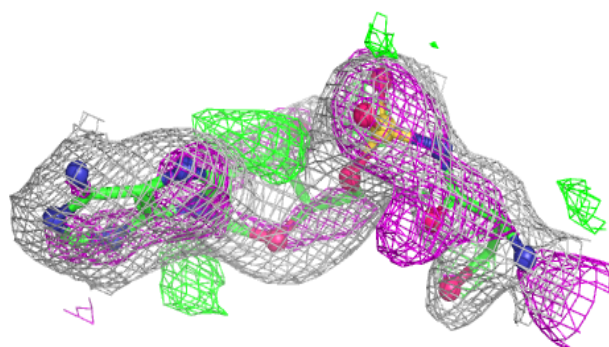
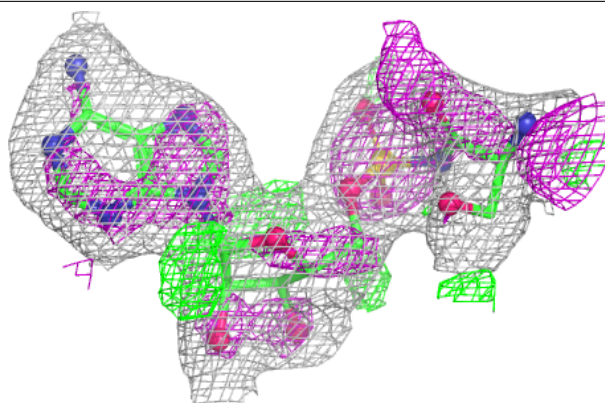
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	SSA	A	1505	29/29	0.93	0.10	37,39,42,42	0
4	SSA	B	1505	29/29	0.94	0.10	34,37,41,41	0
3	CL	B	1504	1/1	0.96	0.19	37,37,37,37	0
2	ZN	A	1503	1/1	0.97	0.18	51,51,51,51	0
3	CL	A	1504	1/1	0.98	0.15	32,32,32,32	0
2	ZN	B	1503	1/1	0.99	0.17	51,51,51,51	0

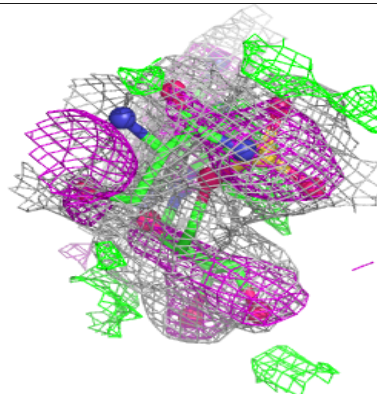
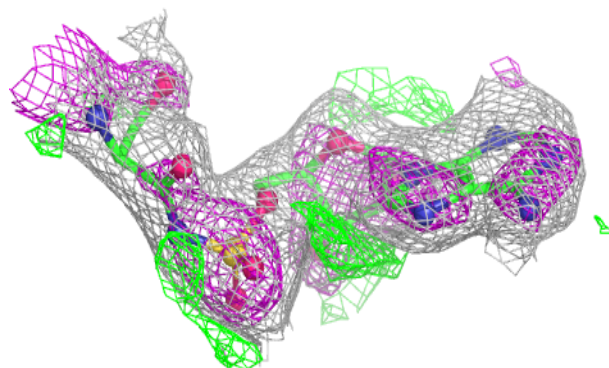
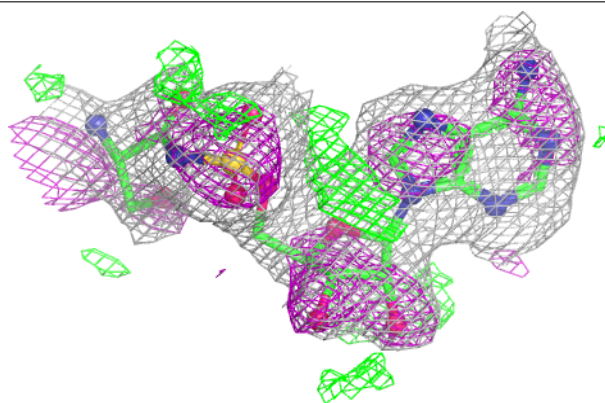
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around SSA A 1505:

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

**Electron density around SSA B 1505:**

$2mF_o-DF_c$ (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.