



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 14, 2024 – 04:41 PM EST

PDB ID : 2GLU
Title : The crystal structure of YcgJ protein from *Bacillus subtilis*
Authors : Burke, T.; Gorman, J.; Shapiro, L.; Burley, S.K.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2006-04-05
Resolution : 2.91 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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.40

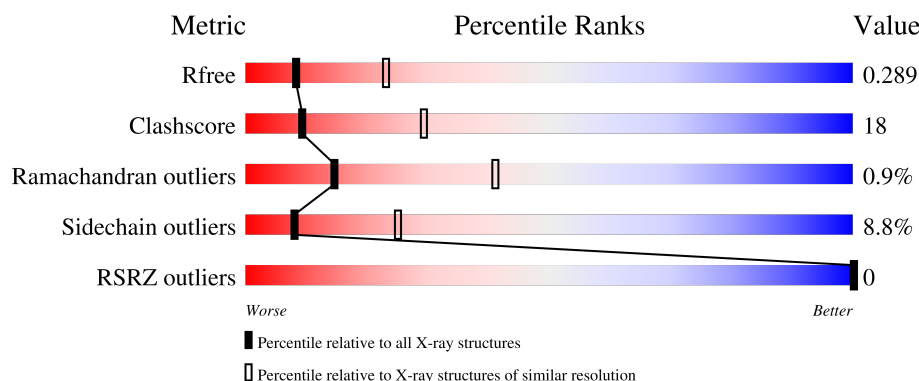
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.91 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}	164625	2797 (2.94-2.90)
Clashscore	180529	3049 (2.94-2.90)
Ramachandran outliers	177936	2981 (2.94-2.90)
Sidechain outliers	177891	2983 (2.94-2.90)
RSRZ outliers	164620	2799 (2.94-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	239	 67% 27% . .
1	B	239	 61% 33% . .

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	A	406	-	-	X	-
2	SO4	B	409	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3923 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 ycgJ.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	234	Total	C	N	O	S	Se	0	0	0
			1861	1166	338	349	5	3			
1	B	234	Total	C	N	O	S	Se	0	0	0
			1861	1166	338	349	5	3			

There are 30 discrepancies between the modelled and reference sequences:

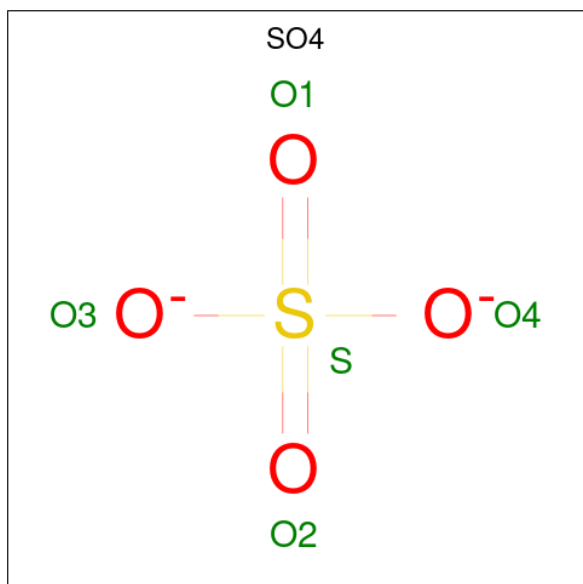
Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	MSE	-	cloning artifact	GB 2632602
A	-7	ALA	-	cloning artifact	GB 2632602
A	-6	HIS	-	expression tag	GB 2632602
A	-5	HIS	-	expression tag	GB 2632602
A	-4	HIS	-	expression tag	GB 2632602
A	-3	HIS	-	expression tag	GB 2632602
A	-2	HIS	-	expression tag	GB 2632602
A	-1	HIS	-	expression tag	GB 2632602
A	0	SER	-	cloning artifact	GB 2632602
A	1	LEU	-	cloning artifact	GB 2632602
A	4	MSE	MET	modified residue	GB 2632602
A	46	MSE	MET	modified residue	GB 2632602
A	151	MSE	MET	modified residue	GB 2632602
A	229	GLU	-	cloning artifact	GB 2632602
A	230	GLY	-	cloning artifact	GB 2632602
B	-8	MSE	-	cloning artifact	GB 2632602
B	-7	ALA	-	cloning artifact	GB 2632602
B	-6	HIS	-	expression tag	GB 2632602
B	-5	HIS	-	expression tag	GB 2632602
B	-4	HIS	-	expression tag	GB 2632602
B	-3	HIS	-	expression tag	GB 2632602
B	-2	HIS	-	expression tag	GB 2632602
B	-1	HIS	-	expression tag	GB 2632602
B	0	SER	-	cloning artifact	GB 2632602
B	1	LEU	-	cloning artifact	GB 2632602

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Chain	Residue	Modelled	Actual	Comment	Reference
B	4	MSE	MET	modified residue	GB 2632602
B	46	MSE	MET	modified residue	GB 2632602
B	151	MSE	MET	modified residue	GB 2632602
B	229	GLU	-	cloning artifact	GB 2632602
B	230	GLY	-	cloning artifact	GB 2632602

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



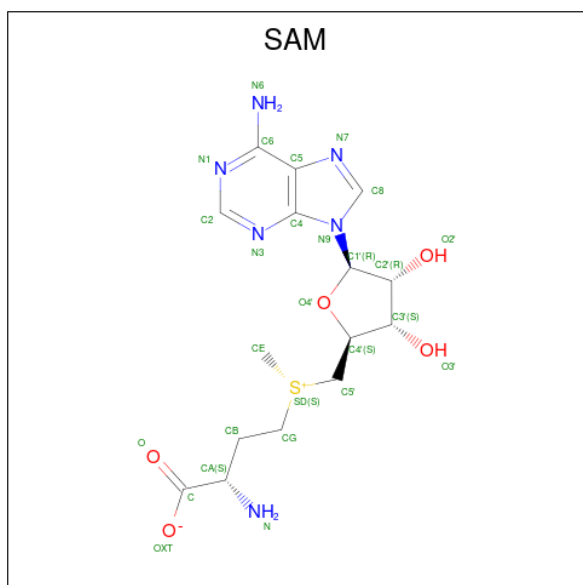
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is S-ADENOSYLMETHIONINE (three-letter code: SAM) (formula: C₁₅H₂₂N₆O₅S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
3	B	1	Total	C	N	O	S	0	0
			27	15	6	5	1		

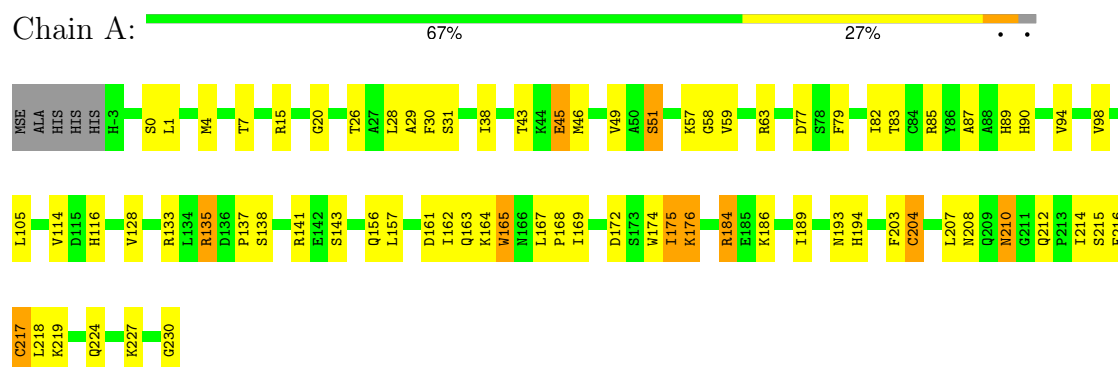
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	47	Total	O	0	0
			47	47		
4	B	45	Total	O	0	0
			45	45		

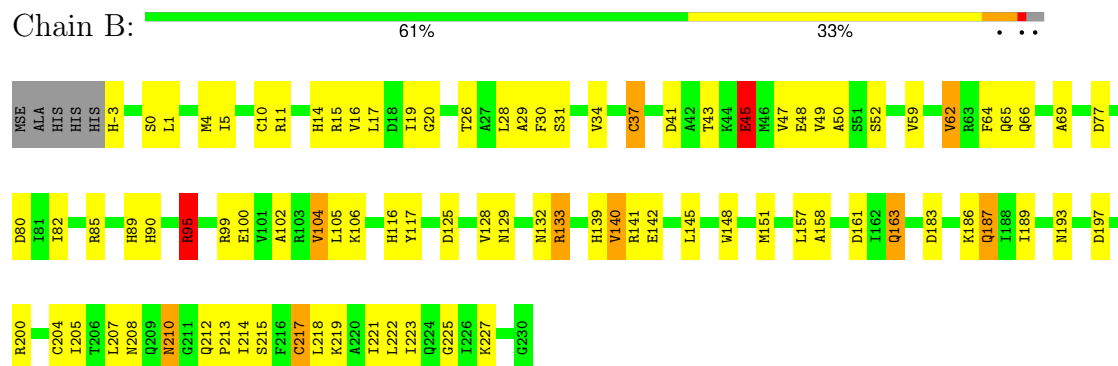
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: ycgJ



• Molecule 1: ycgJ



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	69.90Å 73.64Å 245.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.91 20.00 – 2.91	Depositor EDS
% Data completeness (in resolution range)	94.2 (20.00-2.91) 93.9 (20.00-2.91)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.75 (at 2.93Å)	Xtriage
Refinement program	REFMAC 5.2.0016	Depositor
R, R_{free}	0.192 , 0.283 0.199 , 0.289	Depositor DCC
R_{free} test set	660 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	40.2	Xtriage
Anisotropy	0.802	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 48.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.107 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3923	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.06% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, SAM

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	0.77	2/1900 (0.1%)	0.80	1/2567 (0.0%)
1	B	0.79	2/1900 (0.1%)	0.88	3/2567 (0.1%)
All	All	0.78	4/3800 (0.1%)	0.84	4/5134 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	217	CYS	CB-SG	-6.61	1.71	1.82
1	B	217	CYS	CB-SG	-5.70	1.72	1.81
1	A	204	CYS	CB-SG	-5.46	1.73	1.81
1	B	37	CYS	CB-SG	-5.42	1.73	1.81

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	133	ARG	NE-CZ-NH1	8.06	124.33	120.30
1	B	15	ARG	NE-CZ-NH1	7.69	124.14	120.30
1	B	95	ARG	NE-CZ-NH1	7.66	124.13	120.30
1	A	133	ARG	NE-CZ-NH1	6.00	123.30	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1861	0	1805	63	0
1	B	1861	0	1805	70	0
2	A	20	0	0	6	0
2	B	35	0	0	7	0
3	A	27	0	22	1	0
3	B	27	0	22	1	0
4	A	47	0	0	2	0
4	B	45	0	0	8	0
All	All	3923	0	3654	132	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 18.

All (132) 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:46:MSE:SE	1:A:46:MSE:CE	2.23	1.37
1:B:208:ASN:HD22	1:B:210:ASN:HD21	1.10	0.96
1:B:200:ARG:HA	1:B:205:ILE:HD12	1.47	0.94
1:B:90:HIS:NE2	2:B:409:SO4:O4	2.02	0.93
1:B:89:HIS:NE2	2:B:409:SO4:O2	2.08	0.86
1:B:90:HIS:NE2	2:B:409:SO4:S	2.49	0.85
1:B:89:HIS:NE2	2:B:409:SO4:S	2.50	0.85
1:A:89:HIS:NE2	2:A:406:SO4:S	2.52	0.81
1:B:208:ASN:HD22	1:B:210:ASN:ND2	1.79	0.81
1:A:89:HIS:NE2	2:A:406:SO4:O1	2.16	0.77
1:B:10:CYS:HA	4:B:451:HOH:O	1.86	0.74
1:B:28:LEU:HA	4:B:453:HOH:O	1.89	0.73
1:B:5:ILE:HD11	1:B:29:ALA:HB1	1.71	0.73
1:A:230:GLY:C	1:B:200:ARG:HD3	2.10	0.72
1:B:148:TRP:HA	1:B:151:MSE:HE3	1.70	0.72
1:A:1:LEU:HD13	1:A:85:ARG:NH1	2.08	0.68
1:A:4:MSE:HE3	1:A:114:VAL:HG23	1.76	0.67
1:A:45:GLU:OE2	1:A:45:GLU:N	2.28	0.66
1:B:208:ASN:ND2	1:B:210:ASN:HD21	1.88	0.65
1:B:204:CYS:HG	1:B:217:CYS:HG	1.44	0.63
1:A:45:GLU:H	1:A:45:GLU:CD	2.02	0.62
1:B:99:ARG:HD3	4:B:454:HOH:O	1.99	0.62
1:B:210:ASN:HD21	1:B:212:GLN:HB3	1.66	0.60
1:A:164:LYS:O	1:A:165:TRP:HB3	2.01	0.59
1:B:95:ARG:HH11	1:B:95:ARG:HG3	1.66	0.59
1:B:140:VAL:HG22	1:B:141:ARG:N	2.16	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:ILE:HG13	1:A:105:LEU:HD13	1.84	0.58
1:B:31:SER:HB2	4:B:453:HOH:O	2.03	0.58
1:A:51:SER:HB3	4:A:424:HOH:O	2.04	0.58
1:A:193:ASN:ND2	1:A:207:LEU:HD13	2.18	0.58
1:A:20:GLY:O	3:A:301:SAM:N	2.35	0.58
1:B:132:ASN:ND2	1:B:142:GLU:OE2	2.38	0.56
1:A:43:THR:OG1	1:A:46:MSE:HB2	2.05	0.56
1:A:1:LEU:HD13	1:A:85:ARG:CZ	2.36	0.56
1:B:82:ILE:HD12	1:B:105:LEU:HD13	1.86	0.55
1:A:89:HIS:NE2	2:A:406:SO4:O2	2.38	0.54
1:B:20:GLY:O	3:B:302:SAM:HG2	2.07	0.54
1:B:50:ALA:HB1	1:B:64:PHE:CE1	2.42	0.54
1:A:89:HIS:CD2	2:A:406:SO4:O1	2.60	0.54
1:A:210:ASN:H	1:A:210:ASN:ND2	2.06	0.54
1:B:43:THR:CB	1:B:45:GLU:HG2	2.38	0.54
1:A:163:GLN:HG2	1:A:164:LYS:O	2.08	0.54
1:A:210:ASN:HD22	1:A:210:ASN:N	2.06	0.54
1:A:169:ILE:HG21	1:A:216:PHE:CZ	2.43	0.54
1:A:168:PRO:O	1:A:169:ILE:HD12	2.09	0.53
1:B:208:ASN:HB2	1:B:210:ASN:ND2	2.23	0.53
1:A:1:LEU:HD11	1:A:26:THR:HA	1.91	0.53
1:A:175:ILE:HG22	1:A:176:LYS:N	2.22	0.53
1:B:43:THR:HB	1:B:45:GLU:HG2	1.89	0.52
1:B:82:ILE:CD1	1:B:105:LEU:HD13	2.39	0.52
1:A:30:PHE:CE1	1:A:83:THR:HG21	2.45	0.52
1:B:208:ASN:ND2	1:B:212:GLN:HB3	2.25	0.52
1:B:158:ALA:O	1:B:225:GLY:HA2	2.10	0.52
1:B:30:PHE:O	1:B:34:VAL:HG22	2.10	0.51
1:B:183:ASP:O	1:B:187:GLN:HG2	2.11	0.51
1:B:210:ASN:ND2	1:B:212:GLN:H	2.08	0.51
1:B:210:ASN:HD22	1:B:210:ASN:C	2.13	0.50
1:B:89:HIS:CD2	2:B:409:SO4:O2	2.64	0.50
1:A:45:GLU:O	1:A:49:VAL:HG23	2.11	0.50
1:A:214:ILE:HG22	1:A:215:SER:HB3	1.93	0.50
1:B:129:ASN:OD1	1:B:142:GLU:HG2	2.11	0.50
1:B:19:ILE:N	1:B:19:ILE:HD13	2.25	0.50
1:B:125:ASP:C	1:B:125:ASP:OD2	2.51	0.49
1:B:210:ASN:HD22	1:B:212:GLN:H	1.61	0.49
1:B:117:TYR:CG	1:B:145:LEU:HD21	2.47	0.49
1:B:0:SER:O	1:B:4:MSE:HG3	2.13	0.49
1:A:193:ASN:HD22	1:A:207:LEU:HD13	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:10:CYS:C	4:B:451:HOH:O	2.50	0.49
1:B:189:ILE:HD13	1:B:213:PRO:CD	2.42	0.49
1:A:204:CYS:HG	1:A:217:CYS:HG	1.59	0.48
1:A:208:ASN:HB2	1:A:212:GLN:HB3	1.96	0.48
1:A:1:LEU:HD21	1:A:29:ALA:HB2	1.95	0.48
1:B:11:ARG:HD3	1:B:14:HIS:CE1	2.48	0.48
1:A:164:LYS:NZ	4:A:414:HOH:O	2.32	0.48
1:B:99:ARG:HD2	4:B:421:HOH:O	2.14	0.47
1:B:62:VAL:HG11	4:B:453:HOH:O	2.13	0.47
1:B:161:ASP:OD1	1:B:161:ASP:C	2.52	0.47
1:A:156:GLN:N	1:A:156:GLN:CD	2.67	0.47
1:A:210:ASN:HD21	1:A:212:GLN:HB3	1.79	0.47
1:A:208:ASN:N	1:A:214:ILE:HD13	2.30	0.47
1:A:116:HIS:HA	1:A:219:LYS:O	2.15	0.47
1:A:89:HIS:HA	1:A:141:ARG:O	2.14	0.47
1:B:116:HIS:HA	1:B:219:LYS:O	2.15	0.47
1:B:17:LEU:HG	1:B:19:ILE:HD11	1.97	0.47
1:B:41:ASP:O	1:B:66:GLN:HA	2.14	0.47
1:B:80:ASP:OD2	1:B:106:LYS:HD2	2.16	0.46
1:A:210:ASN:ND2	1:A:210:ASN:N	2.62	0.46
1:B:95:ARG:HH11	1:B:95:ARG:CG	2.29	0.46
1:A:135:ARG:HG3	1:A:174:TRP:CZ2	2.50	0.46
1:A:208:ASN:CG	1:A:214:ILE:HD11	2.36	0.46
1:B:222:LEU:HD12	1:B:223:ILE:N	2.31	0.46
1:B:100:GLU:O	1:B:104:VAL:HG13	2.15	0.45
1:B:105:LEU:HD23	1:B:227:LYS:HB2	1.98	0.45
1:A:214:ILE:HD12	1:A:214:ILE:N	2.30	0.45
1:B:90:HIS:NE2	2:B:409:SO4:O2	2.48	0.45
1:B:16:VAL:CG1	1:B:34:VAL:HG21	2.47	0.45
1:B:221:ILE:O	1:B:221:ILE:HG23	2.17	0.45
1:B:19:ILE:HG21	1:B:69:ALA:HB2	1.98	0.45
1:A:0:SER:O	1:A:4:MSE:HG3	2.18	0.44
1:A:90:HIS:HD2	1:A:138:SER:O	2.01	0.44
1:A:167:LEU:O	1:A:217:CYS:HA	2.18	0.44
1:A:87:ALA:O	1:A:90:HIS:HB2	2.16	0.44
1:A:207:LEU:C	1:A:214:ILE:HD13	2.38	0.44
1:B:90:HIS:CD2	2:B:409:SO4:O4	2.68	0.44
1:A:157:LEU:HD23	1:A:227:LYS:HA	1.99	0.44
1:B:128:VAL:HG13	1:B:129:ASN:N	2.32	0.44
1:A:208:ASN:HB3	1:A:210:ASN:ND2	2.33	0.43
1:B:37:CYS:HB2	1:B:62:VAL:HB	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:VAL:HB	1:A:203:PHE:CZ	2.53	0.43
1:A:90:HIS:CE1	2:A:406:SO4:O2	2.72	0.43
1:A:137:PRO:HB3	1:A:184:ARG:NH2	2.32	0.43
1:B:157:LEU:HD23	1:B:227:LYS:HA	2.01	0.43
1:B:48:GLU:O	1:B:49:VAL:C	2.58	0.42
1:A:210:ASN:H	1:A:210:ASN:HD22	1.65	0.42
1:B:-3:HIS:HB3	1:B:85:ARG:HH12	1.85	0.42
1:A:208:ASN:N	1:A:214:ILE:CD1	2.82	0.42
1:A:208:ASN:OD1	1:A:214:ILE:HD11	2.19	0.42
1:B:102:ALA:O	1:B:227:LYS:NZ	2.41	0.42
1:B:214:ILE:HG22	1:B:215:SER:HB3	2.00	0.42
1:A:90:HIS:CE1	2:A:406:SO4:S	3.12	0.42
1:A:214:ILE:N	1:A:214:ILE:CD1	2.82	0.42
1:A:204:CYS:CB	1:A:217:CYS:HG	2.32	0.42
1:B:14:HIS:CE1	4:B:451:HOH:O	2.74	0.41
1:A:38:ILE:HD12	1:A:79:PHE:CZ	2.54	0.41
1:B:133:ARG:HG3	1:B:139:HIS:O	2.20	0.41
1:A:28:LEU:HD23	1:A:59:VAL:HG21	2.03	0.41
1:B:19:ILE:CG2	1:B:69:ALA:HB2	2.51	0.41
1:A:7:THR:O	1:A:7:THR:HG22	2.19	0.41
1:A:162:ILE:HD11	1:A:224:GLN:HB3	2.02	0.41
1:A:186:LYS:HA	1:A:189:ILE:HD12	2.03	0.41
1:A:94:VAL:O	1:A:98:VAL:HG23	2.21	0.40
1:B:47:VAL:HG11	1:B:65:GLN:HA	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	232/239 (97%)	213 (92%)	17 (7%)	2 (1%)	14 41

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	232/239 (97%)	215 (93%)	15 (6%)	2 (1%)	14	41
All	All	464/478 (97%)	428 (92%)	32 (7%)	4 (1%)	14	41

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	58	GLY
1	A	165	TRP
1	B	45	GLU
1	B	163	GLN

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	199/199 (100%)	182 (92%)	17 (8%)	8	26
1	B	199/199 (100%)	181 (91%)	18 (9%)	8	24
All	All	398/398 (100%)	363 (91%)	35 (9%)	8	25

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	ARG
1	A	31	SER
1	A	45	GLU
1	A	51	SER
1	A	57	LYS
1	A	63	ARG
1	A	77	ASP
1	A	135	ARG
1	A	143	SER
1	A	161	ASP
1	A	172	ASP
1	A	175	ILE

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Mol	Chain	Res	Type
1	A	176	LYS
1	A	184	ARG
1	A	194	HIS
1	A	210	ASN
1	A	218	LEU
1	B	1	LEU
1	B	26	THR
1	B	45	GLU
1	B	52	SER
1	B	59	VAL
1	B	62	VAL
1	B	77	ASP
1	B	95	ARG
1	B	104	VAL
1	B	140	VAL
1	B	163	GLN
1	B	186	LYS
1	B	187	GLN
1	B	193	ASN
1	B	197	ASP
1	B	207	LEU
1	B	210	ASN
1	B	218	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	HIS
1	A	107	GLN
1	A	193	ASN
1	A	210	ASN
1	B	107	GLN
1	B	187	GLN
1	B	210	ASN

5.3.3 RNA ⓘ

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](#)

13 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	SO4	B	408	-	4,4,4	0.20	0	6,6,6	0.26	0
2	SO4	A	406	-	4,4,4	0.27	0	6,6,6	0.88	0
2	SO4	A	410	-	4,4,4	0.25	0	6,6,6	0.29	0
2	SO4	A	404	-	4,4,4	0.28	0	6,6,6	0.42	0
2	SO4	B	407	-	4,4,4	0.29	0	6,6,6	0.53	0
2	SO4	B	401	-	4,4,4	0.16	0	6,6,6	0.27	0
2	SO4	B	405	-	4,4,4	0.22	0	6,6,6	0.19	0
3	SAM	A	301	-	23,29,29	1.25	2 (8%)	20,42,42	2.22	6 (30%)
2	SO4	B	411	-	4,4,4	0.25	0	6,6,6	0.51	0
2	SO4	A	402	-	4,4,4	0.26	0	6,6,6	0.25	0
2	SO4	B	409	-	4,4,4	0.27	0	6,6,6	0.59	0
3	SAM	B	302	-	23,29,29	1.26	2 (8%)	20,42,42	2.31	6 (30%)
2	SO4	B	403	-	4,4,4	0.25	0	6,6,6	0.12	0

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	SAM	A	301	-	-	4/13/33/33	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SAM	B	302	-	-	6/13/33/33	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	302	SAM	C2-N3	4.09	1.38	1.32
3	A	301	SAM	C2-N3	3.59	1.37	1.32
3	B	302	SAM	C2-N1	2.44	1.38	1.33
3	A	301	SAM	O4'-C4'	-2.08	1.40	1.45

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	301	SAM	N3-C2-N1	-6.83	119.40	128.67
3	B	302	SAM	N3-C2-N1	-6.11	120.38	128.67
3	B	302	SAM	O4'-C1'-N9	-5.53	101.41	108.75
3	A	301	SAM	O4'-C1'-N9	-4.43	102.87	108.75
3	B	302	SAM	OXT-C-O	-2.74	117.86	124.08
3	B	302	SAM	C4-C5-N7	-2.69	106.50	109.34
3	A	301	SAM	C4-C5-N7	-2.59	106.60	109.34
3	A	301	SAM	OXT-C-O	-2.53	118.33	124.08
3	A	301	SAM	C1'-N9-C4	-2.50	122.25	126.64
3	B	302	SAM	O4'-C4'-C3'	2.31	109.75	105.15
3	B	302	SAM	O2'-C2'-C3'	-2.22	104.70	111.82
3	A	301	SAM	C4'-O4'-C1'	-2.22	107.89	109.92

There are no chirality outliers.

All (10) torsion outliers are listed below:

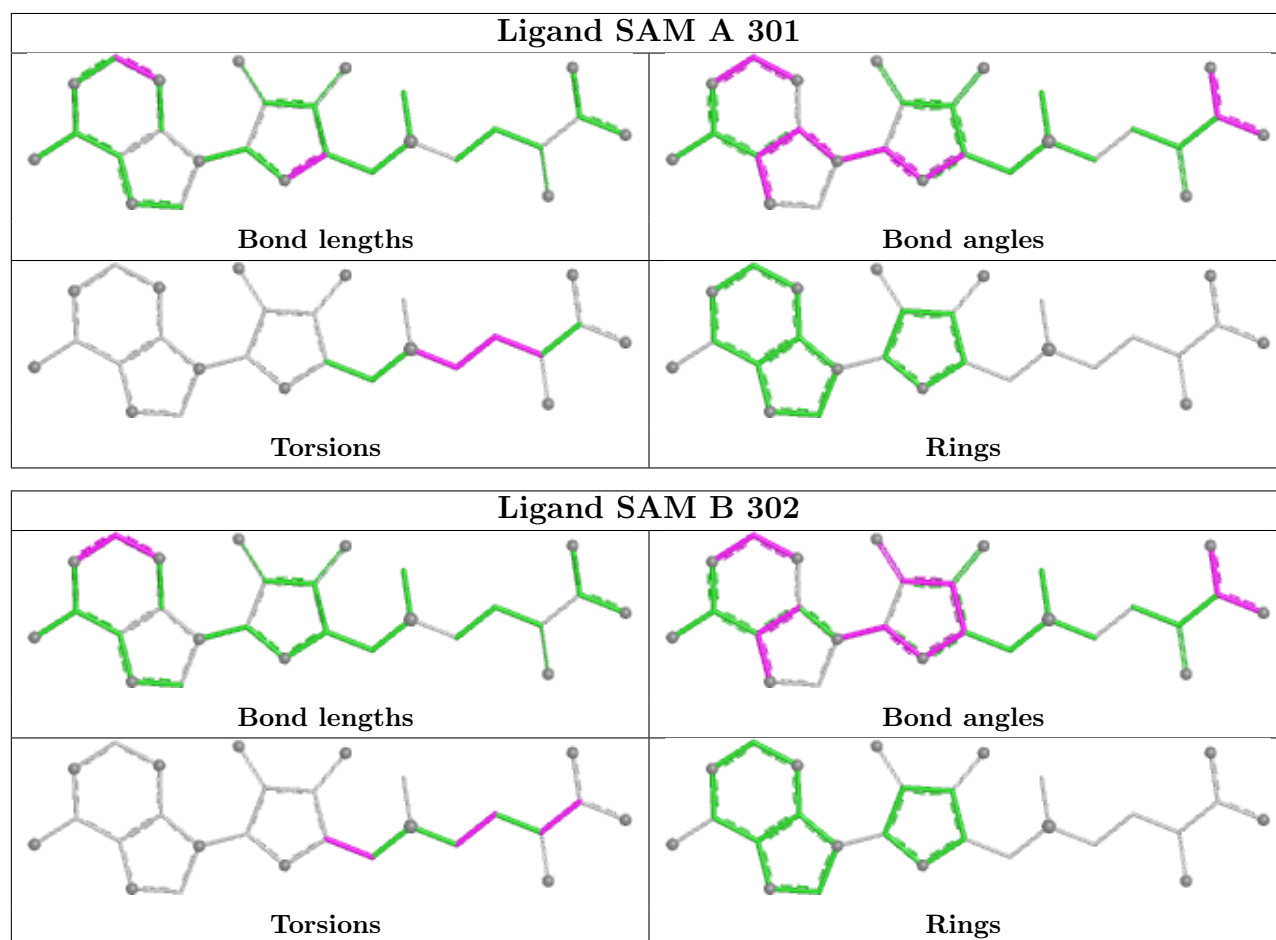
Mol	Chain	Res	Type	Atoms
3	A	301	SAM	N-CA-CB-CG
3	A	301	SAM	CA-CB-CG-SD
3	A	301	SAM	CB-CG-SD-CE
3	B	302	SAM	CA-CB-CG-SD
3	B	302	SAM	OXT-C-CA-CB
3	B	302	SAM	OXT-C-CA-N
3	B	302	SAM	O-C-CA-CB
3	B	302	SAM	O-C-CA-N
3	B	302	SAM	O4'-C4'-C5'-SD
3	A	301	SAM	C-CA-CB-CG

There are no ring outliers.

4 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	406	SO4	6	0
3	A	301	SAM	1	0
2	B	409	SO4	7	0
3	B	302	SAM	1	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	231/239 (96%)	-0.51	0 100 100	20, 32, 47, 51	0
1	B	231/239 (96%)	-0.57	0 100 100	20, 29, 42, 50	0
All	All	462/478 (96%)	-0.54	0 100 100	20, 31, 44, 51	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	B	403	5/5	0.85	0.15	118,118,119,119	0
2	SO4	B	411	5/5	0.85	0.22	83,85,86,86	0
2	SO4	A	410	5/5	0.86	0.14	67,69,70,71	0
2	SO4	B	408	5/5	0.88	0.09	75,75,76,76	0
2	SO4	B	405	5/5	0.88	0.13	73,73,74,75	0
2	SO4	B	409	5/5	0.91	0.10	66,68,69,69	0
2	SO4	A	406	5/5	0.93	0.08	65,65,66,66	0

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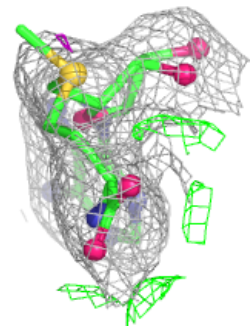
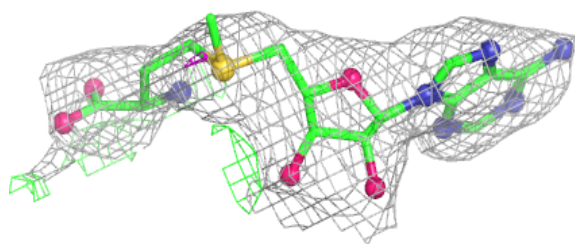
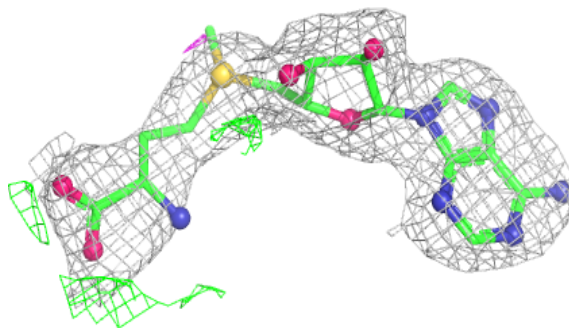
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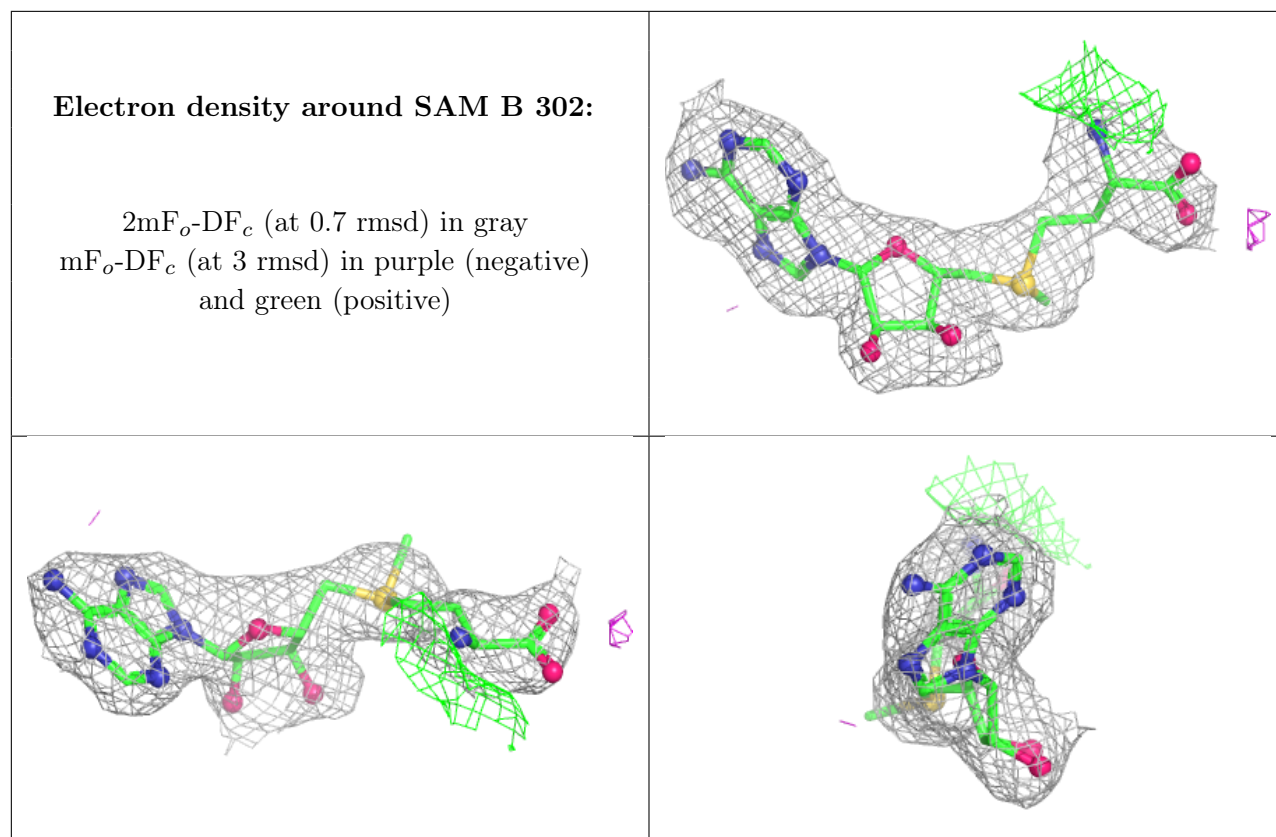
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	SO4	A	402	5/5	0.94	0.13	84,84,84,85	0
3	SAM	A	301	27/27	0.94	0.09	30,33,47,48	0
3	SAM	B	302	27/27	0.95	0.08	23,26,39,40	0
2	SO4	B	407	5/5	0.98	0.07	55,56,57,57	0
2	SO4	A	404	5/5	0.98	0.06	33,34,35,36	0
2	SO4	B	401	5/5	0.99	0.09	35,36,37,38	0

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 SAM A 301:

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.