



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:52 PM UTC

PDB ID : 9GWO / pdb_00009gwo
Title : SARS-CoV-2 methyltransferase nsp10-16 in complex with SAM and theophylline derivative LAS 54571126
Authors : Kremling, V.; Sprenger, J.; Oberthuer, D.; Kiene, A.
Deposited on : 2024-09-27
Resolution : 2.00 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

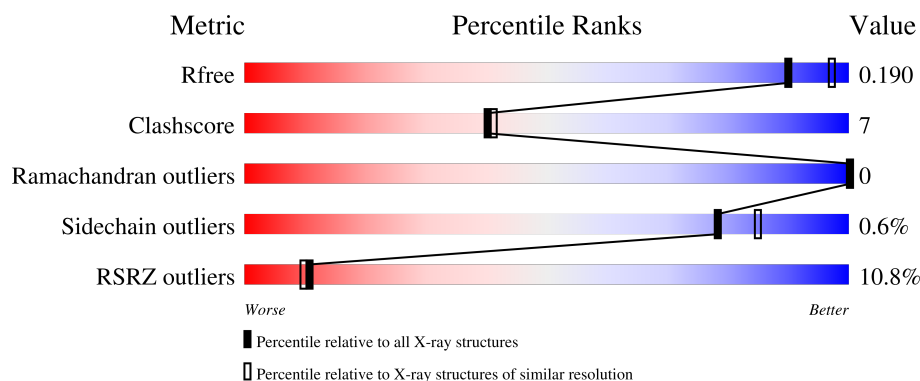
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	304	8% 86% 12%
2	B	140	14% 77% 5% 18%

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
6	A1IQS	A	7209	-	-	-	X

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 3494 atoms, of which 23 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 2'-O-methyltransferase nsp16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	300	Total	C	N	O	S	0	1	0
			2368	1511	397	443	17			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	7097	GLU	-	expression tag	UNP P0DTD1
A	7098	ASN	-	expression tag	UNP P0DTD1
A	7099	LEU	-	expression tag	UNP P0DTD1
A	7100	TYR	-	expression tag	UNP P0DTD1
A	7101	PHE	-	expression tag	UNP P0DTD1
A	7102	GLN	-	expression tag	UNP P0DTD1

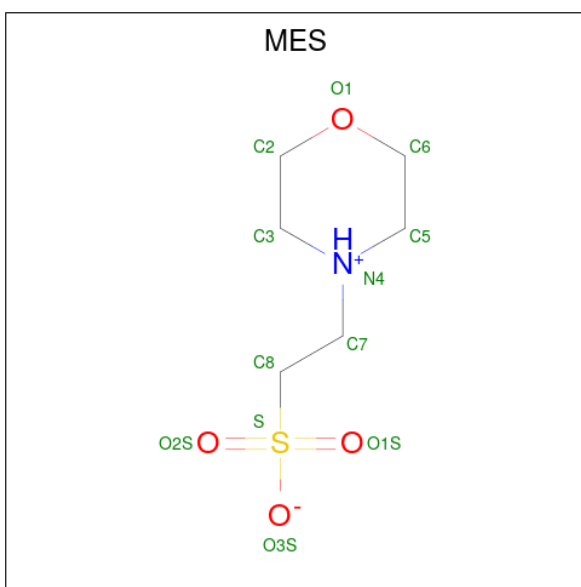
- Molecule 2 is a protein called Non-structural protein 10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	115	Total	C	N	O	S	0	0	0
			848	527	143	163	15			

There is a discrepancy between the modelled and reference sequences:

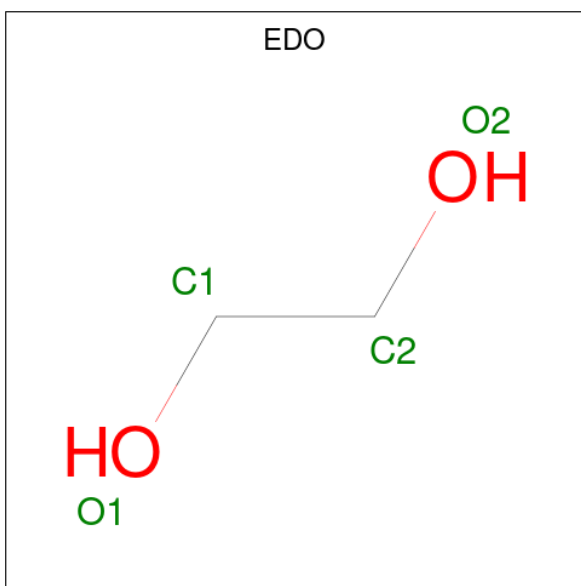
Chain	Residue	Modelled	Actual	Comment	Reference
B	4252	GLY	-	expression tag	UNP P0DTD1

- Molecule 3 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

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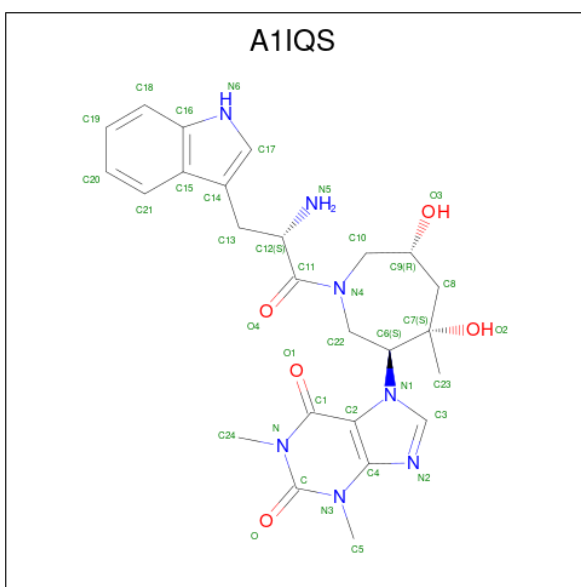
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0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

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

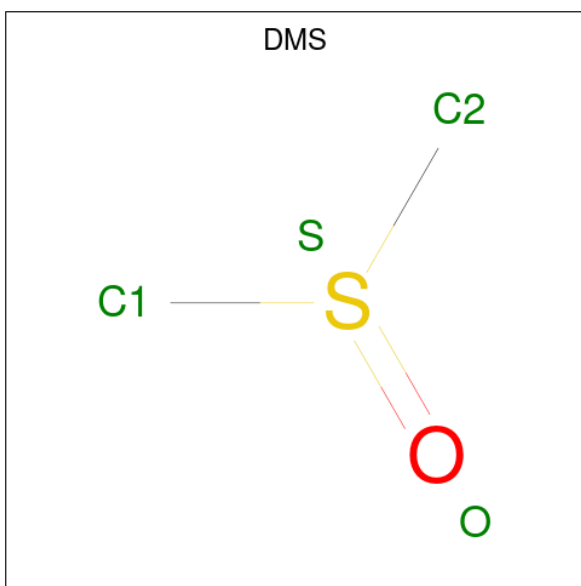
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	2	Total Na 2 2	0	0

- Molecule 6 is 7-[(3 {S},4 {S},6 {R})-1-[(2 {S})-2-azanyl-3-(1 {H}-indol-3-yl)propanoyl]-4-methyl-4,6-bis(oxidanyl)azepan-3-yl]-1,3-dimethyl-purine-2,6-dione (CCD ID: A1IQS) (formula: C₂₅H₃₁N₇O₅) (labeled as "Ligand of Interest" by depositor).



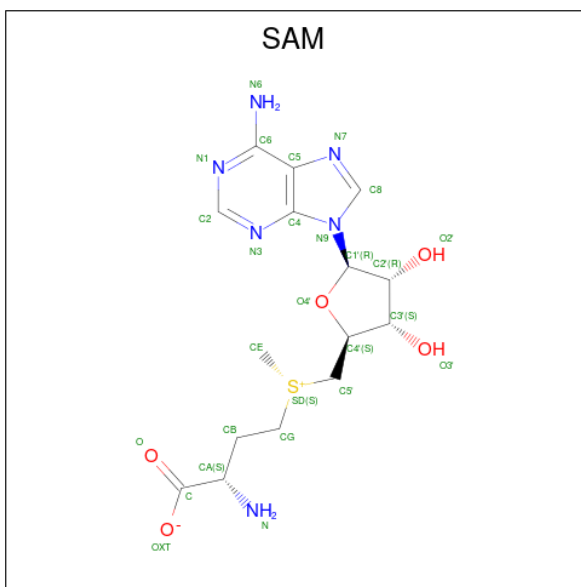
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			37	25	7	5		

- Molecule 7 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C_2H_6OS).



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

- Molecule 8 is S-ADENOSYLMETHIONINE (CCD ID: SAM) (formula: $C_{15}H_{22}N_6O_5S$).

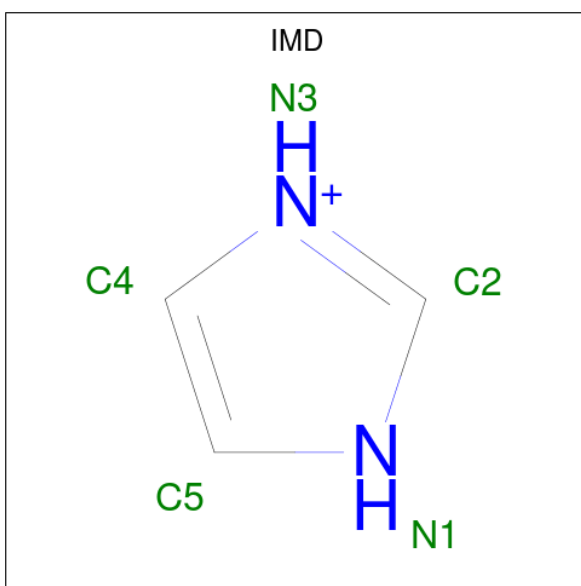


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	A	1	Total	C	H	N	O	S	
			50	15	23	6	5	1	

- Molecule 9 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	B	2	Total	Zn		
			2	2	0	0

- Molecule 10 is IMIDAZOLE (CCD ID: IMD) (formula: C₃H₅N₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	B	1	Total	C	N	0	0
			5	3	2		

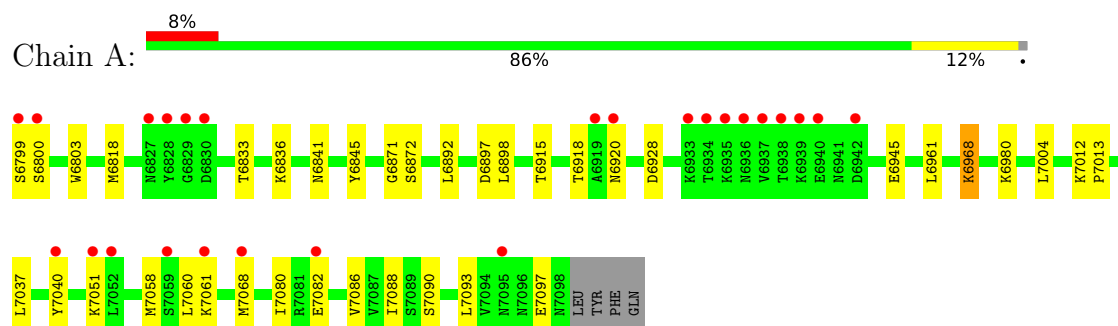
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	89	Total	O	0	0
			89	89		
11	B	21	Total	O	0	0
			21	21		

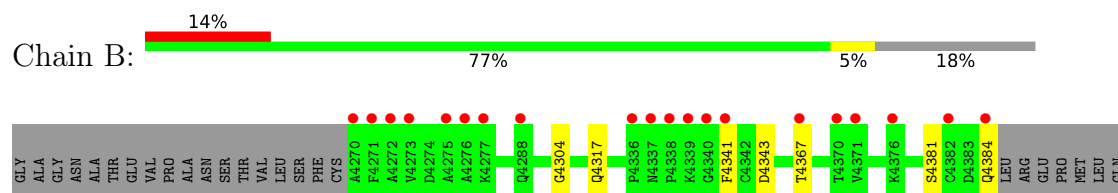
3 Residue-property plots [i](#)

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: 2'-O-methyltransferase nsp16



• Molecule 2: Non-structural protein 10



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	168.08Å 168.08Å 51.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.74 – 2.00 48.74 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.74-2.00) 99.8 (48.74-2.00)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.10 (at 1.52Å)	Xtriage
Refinement program	PHENIX 1.21.1-5286	Depositor
R, R_{free}	0.189 , 0.203 0.186 , 0.190	Depositor DCC
R_{free} test set	1028 reflections (1.05%)	wwPDB-VP
Wilson B-factor (Å ²)	37.8	Xtriage
Anisotropy	0.018	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 28.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3494	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.65% 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: IMD, DMS, EDO, SAM, MES, ZN, A1IQS, NA

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	0.34	0/2420	0.50	0/3283
2	B	0.28	0/867	0.45	0/1176
All	All	0.33	0/3287	0.49	0/4459

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 [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	2368	0	2349	41	0
2	B	848	0	805	3	0
3	A	12	0	13	0	0
4	A	44	0	66	12	0
4	B	12	0	18	0	0
5	A	2	0	0	0	0
6	A	37	0	0	0	0
7	A	4	0	6	0	0
8	A	27	23	22	5	0
9	B	2	0	0	0	0
10	B	5	0	5	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
11	A	89	0	0	2	0
11	B	21	0	0	0	0
All	All	3471	23	3284	44	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 7.

All (44) 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:7082:GLU:HB2	4:A:7208:EDO:H12	1.45	0.98
1:A:6980:LYS:HD3	4:A:7206:EDO:H22	1.46	0.97
1:A:7051:LYS:HD2	1:A:7051:LYS:H	1.40	0.86
1:A:7097:GLU:OE1	1:A:7097:GLU:N	2.08	0.86
1:A:6915:THR:HG22	4:A:7215:EDO:H12	1.62	0.79
1:A:6833:THR:H	4:A:7212:EDO:H11	1.51	0.75
1:A:7082:GLU:HB2	4:A:7208:EDO:C1	2.17	0.74
1:A:6928:ASP:OD2	1:A:6968:LYS:HE2	1.88	0.72
1:A:7060:LEU:C	1:A:7061:LYS:HD3	2.17	0.69
1:A:6799:SER:OG	1:A:6800:SER:N	2.27	0.68
1:A:7058:MET:CE	1:A:7060:LEU:HD11	2.32	0.59
1:A:7061:LYS:HD3	1:A:7061:LYS:N	2.18	0.58
1:A:6818:MET:HE3	11:A:7375:HOH:O	2.06	0.54
1:A:7082:GLU:HB2	4:A:7208:EDO:C2	2.37	0.54
1:A:6961:LEU:HB2	1:A:7080:ILE:HB	1.89	0.54
1:A:6803:TRP:HD1	1:A:7040[B]:TYR:CZ	2.26	0.54
1:A:6833:THR:H	4:A:7212:EDO:C1	2.19	0.53
1:A:7058:MET:HE1	1:A:7060:LEU:HD11	1.91	0.53
1:A:7093:LEU:HB2	4:A:7215:EDO:H21	1.92	0.51
2:B:4341:PHE:HE2	2:B:4343:ASP:OD1	1.95	0.50
2:B:4381:SER:O	2:B:4384:GLN:HG2	2.12	0.49
1:A:6803:TRP:HA	1:A:7040[B]:TYR:CD2	2.50	0.47
1:A:6872:SER:HB3	8:A:7216:SAM:N	2.29	0.47
1:A:6918:THR:HG22	1:A:7088:ILE:HG22	1.96	0.47
1:A:6892:LEU:HD12	1:A:7086:VAL:HG21	1.97	0.46
1:A:6897:ASP:OD1	8:A:7216:SAM:O2'	2.32	0.46
1:A:6897:ASP:OD1	1:A:6898:LEU:N	2.46	0.46
1:A:6803:TRP:HD1	1:A:7040[B]:TYR:CE1	2.34	0.45
1:A:7012:LYS:HB2	1:A:7013:PRO:HD2	1.98	0.45
1:A:6836:LYS:HZ3	4:A:7211:EDO:H21	1.82	0.45
1:A:6845:TYR:OH	1:A:6928:ASP:HB2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6915:THR:O	1:A:7090:SER:HA	2.17	0.45
1:A:6803:TRP:HD1	1:A:7040[B]:TYR:CE2	2.36	0.44
1:A:6945:GLU:O	4:A:7213:EDO:H22	2.18	0.44
1:A:6928:ASP:OD1	8:A:7216:SAM:HG1	2.19	0.42
1:A:7068:MET:HE3	1:A:7068:MET:HB3	1.95	0.42
1:A:6836:LYS:NZ	4:A:7211:EDO:H21	2.35	0.42
1:A:6871:GLY:O	8:A:7216:SAM:HB1	2.20	0.42
1:A:7004:LEU:HD23	1:A:7004:LEU:HA	1.95	0.41
1:A:6920:ASN:ND2	1:A:6920:ASN:H	2.19	0.41
1:A:7037:LEU:HD23	11:A:7368:HOH:O	2.21	0.41
1:A:6833:THR:N	4:A:7212:EDO:H11	2.29	0.40
1:A:6841:ASN:ND2	8:A:7216:SAM:N	2.70	0.40
2:B:4304:GLY:HA2	2:B:4317:GLN:OE1	2.21	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	299/304 (98%)	291 (97%)	8 (3%)	0	100	100
2	B	113/140 (81%)	110 (97%)	3 (3%)	0	100	100
All	All	412/444 (93%)	401 (97%)	11 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	261/264 (99%)	260 (100%)	1 (0%)	84	89
2	B	93/113 (82%)	92 (99%)	1 (1%)	65	73
All	All	354/377 (94%)	352 (99%)	2 (1%)	78	85

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6968	LYS
2	B	4367	THR

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	6804	GLN
1	A	6841	ASN
1	A	6917	HIS
1	A	6920	ASN
1	A	7008	ASN
2	B	4305	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](#)

Of 23 ligands modelled in this entry, 4 are monoatomic - leaving 19 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	EDO	A	7204	-	3,3,3	0.31	0	2,2,2	0.13	0
10	IMD	B	8205	-	5,5,5	0.31	0	5,5,5	0.82	0
4	EDO	A	7211	-	3,3,3	0.26	0	2,2,2	0.35	0
4	EDO	A	7212	-	3,3,3	0.23	0	2,2,2	0.42	0
4	EDO	A	7213	-	3,3,3	0.20	0	2,2,2	0.48	0
4	EDO	A	7215	-	3,3,3	0.25	0	2,2,2	0.16	0
3	MES	A	7201	-	12,12,12	1.41	1 (8%)	15,16,16	0.94	0
4	EDO	A	7203	-	3,3,3	0.29	0	2,2,2	0.11	0
4	EDO	B	8201	-	3,3,3	0.32	0	2,2,2	0.27	0
4	EDO	A	7206	-	3,3,3	0.29	0	2,2,2	0.26	0
4	EDO	A	7210	-	3,3,3	0.33	0	2,2,2	0.28	0
7	DMS	A	7214	-	3,3,3	0.61	0	3,3,3	0.16	0
8	SAM	A	7216	-	27,29,29	0.46	0	34,42,42	0.95	2 (5%)
4	EDO	A	7208	-	3,3,3	0.27	0	2,2,2	0.39	0
4	EDO	B	8206	-	3,3,3	0.28	0	2,2,2	0.56	0
4	EDO	B	8204	-	3,3,3	0.24	0	2,2,2	0.42	0
4	EDO	A	7202	-	3,3,3	0.27	0	2,2,2	0.51	0
6	A1IQS	A	7209	-	37,41,41	0.21	0	50,63,63	0.58	1 (2%)
4	EDO	A	7207	-	3,3,3	0.28	0	2,2,2	0.15	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	MES	A	7201	-	-	3/6/14/14	0/1/1/1
4	EDO	A	7204	-	-	1/1/1/1	-
4	EDO	A	7203	-	-	1/1/1/1	-
10	IMD	B	8205	-	-	-	0/1/1/1
4	EDO	A	7210	-	-	0/1/1/1	-
4	EDO	A	7211	-	-	1/1/1/1	-
4	EDO	A	7212	-	-	1/1/1/1	-
4	EDO	B	8201	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	8206	-	-	0/1/1/1	-
8	SAM	A	7216	-	-	5/17/33/33	0/3/3/3
4	EDO	B	8204	-	-	0/1/1/1	-
4	EDO	A	7202	-	-	1/1/1/1	-
4	EDO	A	7208	-	-	1/1/1/1	-
6	A1IQS	A	7209	-	-	5/12/35/35	0/5/5/5
4	EDO	A	7206	-	-	1/1/1/1	-
4	EDO	A	7213	-	-	0/1/1/1	-
4	EDO	A	7215	-	-	0/1/1/1	-
4	EDO	A	7207	-	-	0/1/1/1	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	7201	MES	C8-S	3.91	1.83	1.77

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	7216	SAM	CG-SD-C5'	3.81	112.74	103.43
6	A	7209	A1IQS	C22-C6-N1	2.31	112.63	109.72
8	A	7216	SAM	C4'-O4'-C1'	-2.07	104.90	109.47

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	7201	MES	C7-C8-S-O1S
6	A	7209	A1IQS	C22-C6-N1-C3
6	A	7209	A1IQS	C11-C12-C13-C14
6	A	7209	A1IQS	N5-C12-C13-C14
6	A	7209	A1IQS	O4-C11-N4-C10
6	A	7209	A1IQS	O4-C11-N4-C22
8	A	7216	SAM	CB-CG-SD-CE
3	A	7201	MES	C7-C8-S-O3S
4	A	7202	EDO	O1-C1-C2-O2
4	A	7206	EDO	O1-C1-C2-O2
4	A	7203	EDO	O1-C1-C2-O2
4	A	7204	EDO	O1-C1-C2-O2
4	A	7211	EDO	O1-C1-C2-O2
3	A	7201	MES	C7-C8-S-O2S

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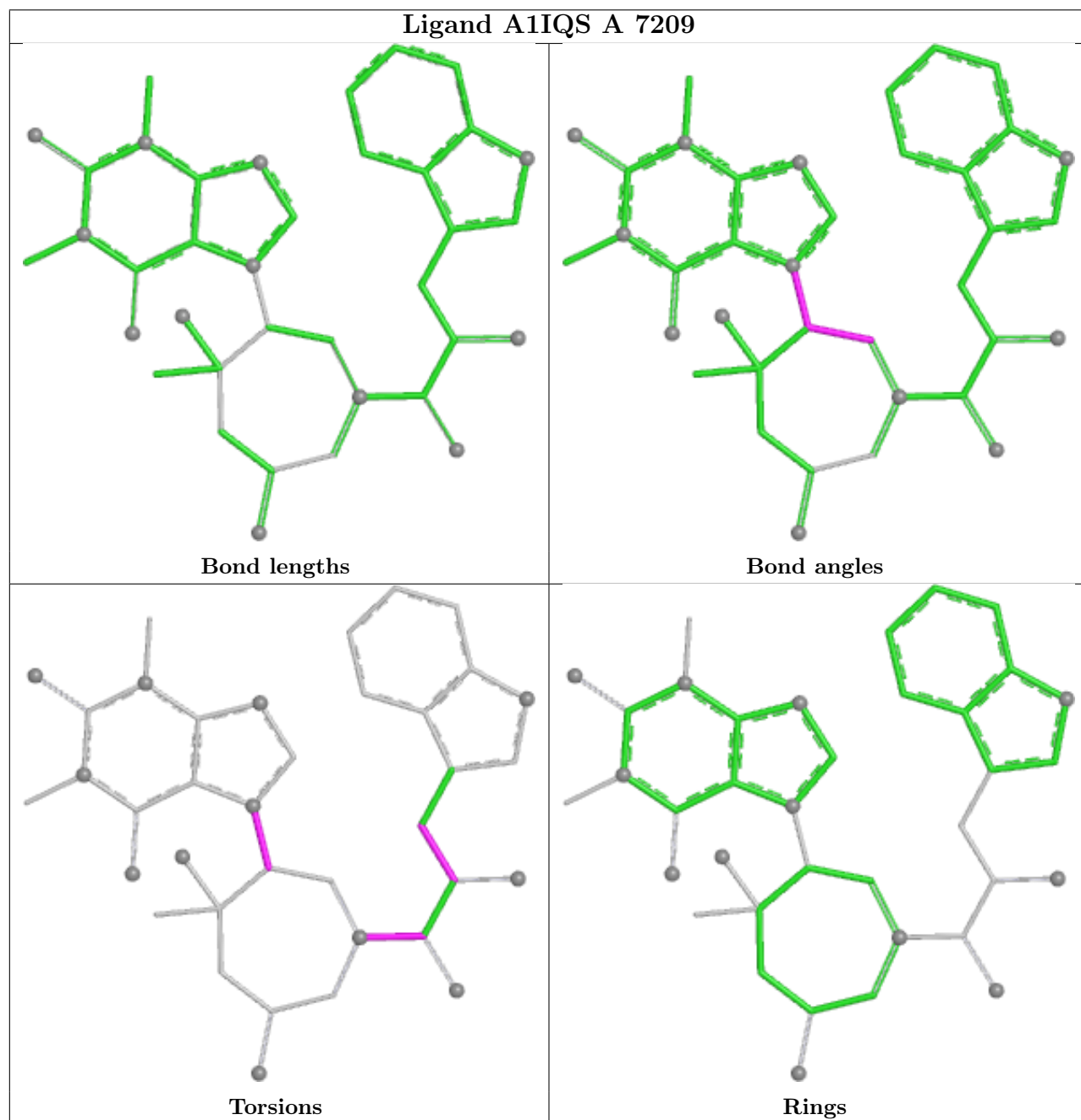
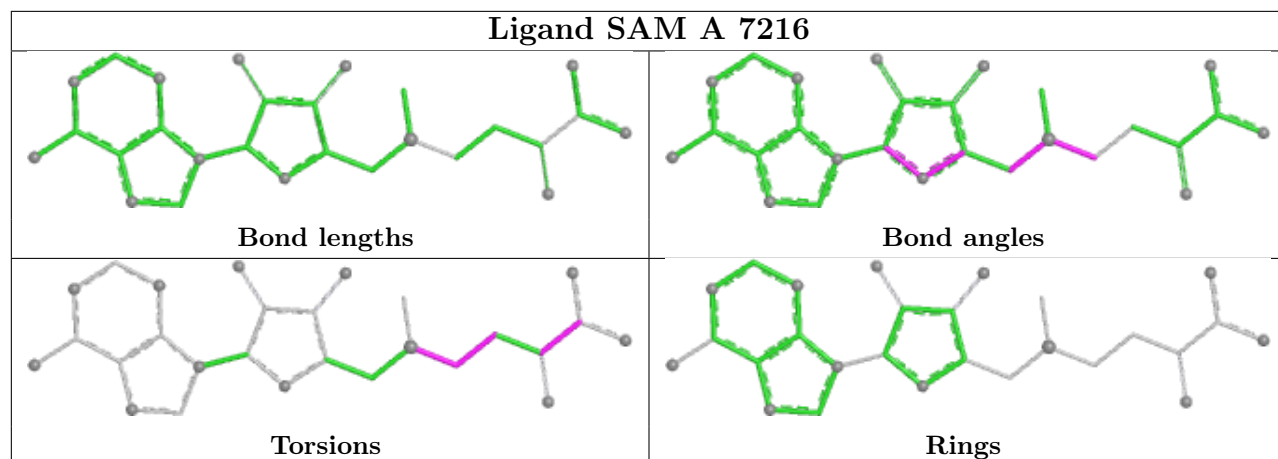
Mol	Chain	Res	Type	Atoms
8	A	7216	SAM	CA-CB-CG-SD
8	A	7216	SAM	O-C-CA-CB
8	A	7216	SAM	OXT-C-CA-CB
4	B	8201	EDO	O1-C1-C2-O2
4	A	7212	EDO	O1-C1-C2-O2
4	A	7208	EDO	O1-C1-C2-O2
8	A	7216	SAM	CB-CG-SD-C5'

There are no ring outliers.

7 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	7211	EDO	2	0
4	A	7212	EDO	3	0
4	A	7213	EDO	1	0
4	A	7215	EDO	2	0
4	A	7206	EDO	1	0
8	A	7216	SAM	5	0
4	A	7208	EDO	3	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

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	300/304 (98%)	0.20	25 (8%) 17 16	16, 35, 66, 91	1 (0%)
2	B	115/140 (82%)	0.89	20 (17%) 4 3	28, 45, 78, 89	0
All	All	415/444 (93%)	0.39	45 (10%) 11 10	16, 38, 70, 91	1 (0%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	4270	ALA	5.4
2	B	4338	PRO	5.3
2	B	4341	PHE	5.2
1	A	6828	TYR	4.6
1	A	6937	VAL	4.5
2	B	4340	GLY	4.5
1	A	6938	THR	4.3
1	A	6829	GLY	4.1
1	A	6800	SER	4.1
1	A	6799	SER	3.9
1	A	6939	LYS	3.6
2	B	4339	LYS	3.5
1	A	7052	LEU	3.5
1	A	6936	ASN	3.4
2	B	4336	PRO	3.4
2	B	4337	ASN	3.0
2	B	4370	THR	3.0
2	B	4384	GLN	2.9
1	A	6830	ASP	2.8
2	B	4271	PHE	2.8
1	A	6942	ASP	2.7
1	A	7095	ASN	2.6
2	B	4371	VAL	2.6
1	A	6935	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	4382	CYS	2.5
1	A	6934	THR	2.5
1	A	6933	LYS	2.5
1	A	7059	SER	2.4
1	A	7051	LYS	2.4
1	A	7061	LYS	2.4
2	B	4273	VAL	2.4
1	A	7040[A]	TYR	2.3
2	B	4272	ALA	2.3
2	B	4277	LYS	2.3
1	A	7082	GLU	2.2
2	B	4288	GLN	2.2
1	A	6919	ALA	2.2
1	A	7068	MET	2.1
2	B	4367	THR	2.1
1	A	6920	ASN	2.1
2	B	4276	ALA	2.1
1	A	6940	GLU	2.0
2	B	4275	ALA	2.1
1	A	6827	ASN	2.0
2	B	4376	LYS	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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
6	A1IQS	A	7209	37/37	0.70	0.42	63,103,132,134	37
4	EDO	A	7204	4/4	0.74	0.26	50,61,64,68	0

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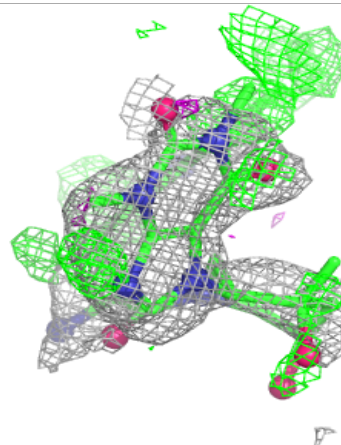
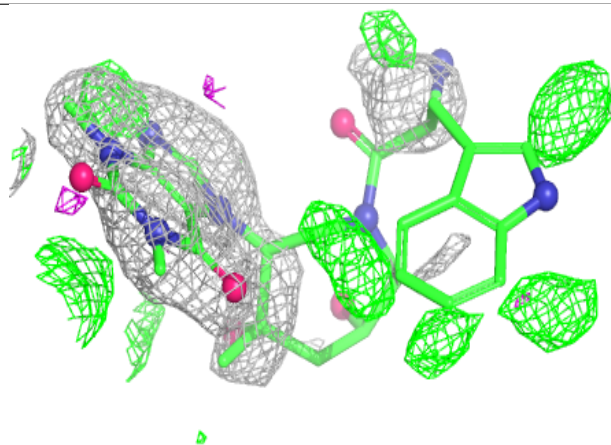
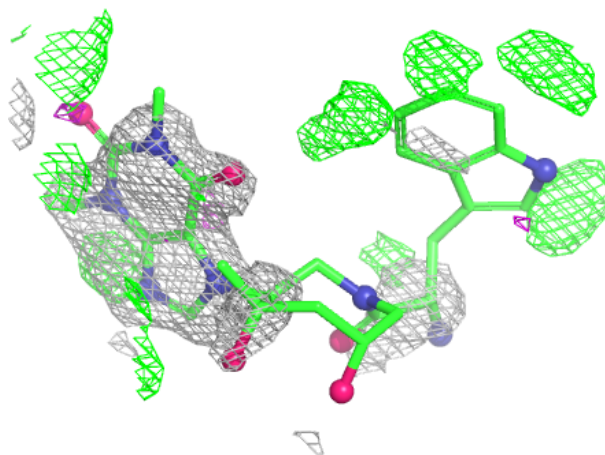
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	EDO	B	8204	4/4	0.76	0.26	64,67,71,82	0
7	DMS	A	7214	4/4	0.78	0.30	76,80,91,103	0
4	EDO	A	7212	4/4	0.82	0.19	45,47,54,66	0
4	EDO	A	7211	4/4	0.83	0.22	54,59,59,61	0
4	EDO	A	7210	4/4	0.85	0.22	47,48,52,60	0
4	EDO	A	7208	4/4	0.85	0.20	45,56,56,57	0
4	EDO	A	7206	4/4	0.86	0.21	55,58,62,66	0
8	SAM	A	7216	27/27	0.87	0.16	32,48,65,68	50
4	EDO	B	8206	4/4	0.89	0.26	44,48,48,59	0
4	EDO	A	7207	4/4	0.90	0.25	47,49,53,54	0
3	MES	A	7201	12/12	0.90	0.13	40,44,67,67	0
4	EDO	A	7215	4/4	0.90	0.20	50,55,61,65	0
4	EDO	A	7203	4/4	0.90	0.20	53,59,60,62	0
10	IMD	B	8205	5/5	0.90	0.19	59,63,68,70	0
4	EDO	A	7213	4/4	0.92	0.15	39,45,48,55	0
4	EDO	A	7202	4/4	0.92	0.15	51,52,58,63	0
4	EDO	B	8201	4/4	0.94	0.31	38,41,43,43	0
5	NA	A	7217	1/1	0.96	0.13	56,56,56,56	0
9	ZN	B	8203	1/1	0.97	0.06	69,69,69,69	0
5	NA	A	7205	1/1	0.98	0.11	41,41,41,41	0
9	ZN	B	8202	1/1	0.99	0.03	37,37,37,37	1

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.

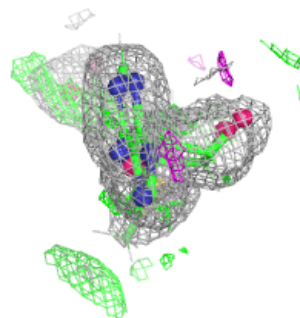
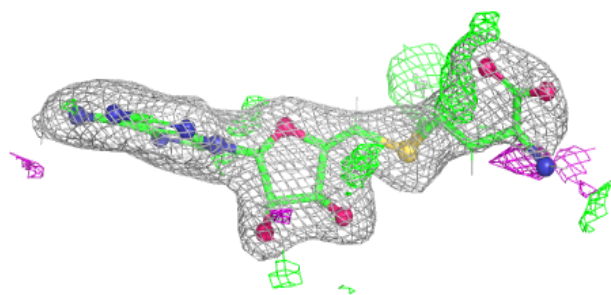
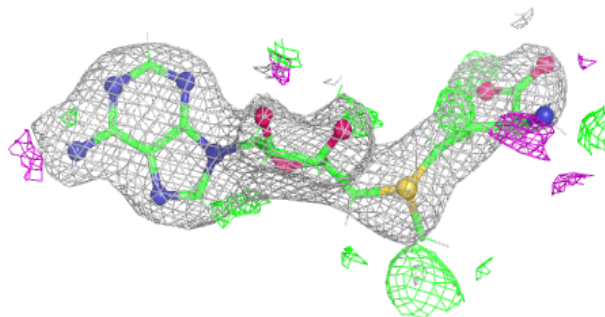
Electron density around A1IQS A 7209:

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



Electron density around SAM A 7216:

$2mF_o-DF_c$ (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.