



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 6, 2025 – 02:07 PM JST

PDB ID : 9IQL / pdb_00009iql
Title : SacM in complex with L6S
Authors : Zhang, B.; Ge, H.M.
Deposited on : 2024-07-12
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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.46

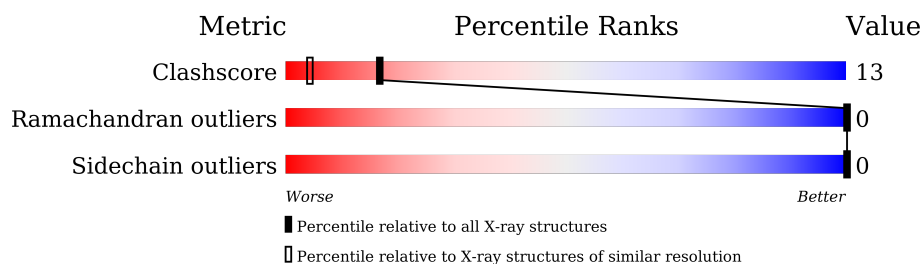
1 Overall quality at a glance

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))
Clashscore	180529	2515 (1.66-1.66)
Ramachandran outliers	177936	2475 (1.66-1.66)
Sidechain outliers	177891	2475 (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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	133	
1	B	133	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 2250 atoms, of which 88 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 Nuclear transport factor 2 family protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	121	Total	C	N	O	S	0	2	0
			960	608	164	185	3			
1	B	125	Total	C	N	O	S	0	3	0
			991	630	169	189	3			

There are 26 discrepancies between the modelled and reference sequences:

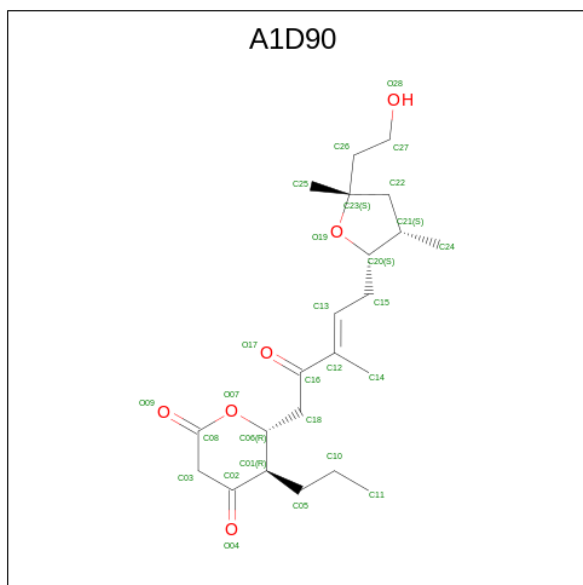
Chain	Residue	Modelled	Actual	Comment	Reference
A	121	LYS	-	expression tag	UNP A0A5Q0H2K6
A	122	LEU	-	expression tag	UNP A0A5Q0H2K6
A	123	ALA	-	expression tag	UNP A0A5Q0H2K6
A	124	ALA	-	expression tag	UNP A0A5Q0H2K6
A	125	ALA	-	expression tag	UNP A0A5Q0H2K6
A	126	LEU	-	expression tag	UNP A0A5Q0H2K6
A	127	GLU	-	expression tag	UNP A0A5Q0H2K6
A	128	HIS	-	expression tag	UNP A0A5Q0H2K6
A	129	HIS	-	expression tag	UNP A0A5Q0H2K6
A	130	HIS	-	expression tag	UNP A0A5Q0H2K6
A	131	HIS	-	expression tag	UNP A0A5Q0H2K6
A	132	HIS	-	expression tag	UNP A0A5Q0H2K6
A	133	HIS	-	expression tag	UNP A0A5Q0H2K6
B	121	LYS	-	expression tag	UNP A0A5Q0H2K6
B	122	LEU	-	expression tag	UNP A0A5Q0H2K6
B	123	ALA	-	expression tag	UNP A0A5Q0H2K6
B	124	ALA	-	expression tag	UNP A0A5Q0H2K6
B	125	ALA	-	expression tag	UNP A0A5Q0H2K6
B	126	LEU	-	expression tag	UNP A0A5Q0H2K6
B	127	GLU	-	expression tag	UNP A0A5Q0H2K6
B	128	HIS	-	expression tag	UNP A0A5Q0H2K6
B	129	HIS	-	expression tag	UNP A0A5Q0H2K6
B	130	HIS	-	expression tag	UNP A0A5Q0H2K6
B	131	HIS	-	expression tag	UNP A0A5Q0H2K6
B	132	HIS	-	expression tag	UNP A0A5Q0H2K6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	133	HIS	-	expression tag	UNP A0A5Q0H2K6

- Molecule 2 is (5 {R},6 {R})-6-[({E})-5-[(2 {S},3 {S},5 {S})-5-(2-hydroxyethyl)-3,5-dimethyl-oxolan-2-yl]-3-methyl-2-oxidanylidene-pent-3-enyl]-5-propyl-oxane-2,4-dione (CCD ID: A1D90) (formula: C₂₂H₃₄O₆) (labeled as "Ligand of Interest" by depositor).



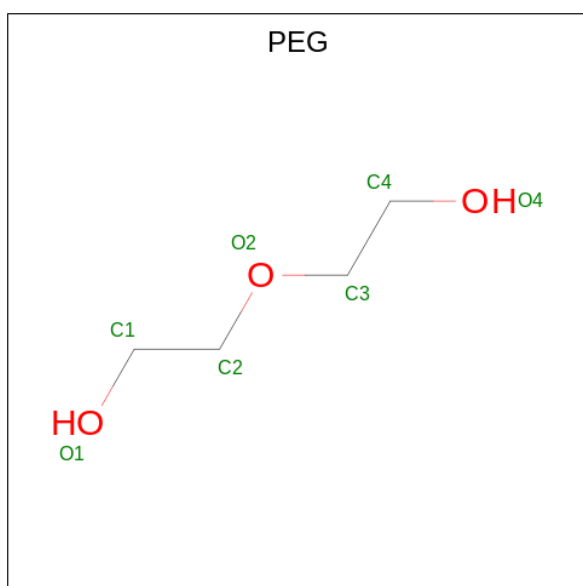
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			61	22	33	6		
2	B	1	Total	C	H	O	0	0
			61	22	33	6		

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			10	2	6	2		
3	B	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			17	4	10	3		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



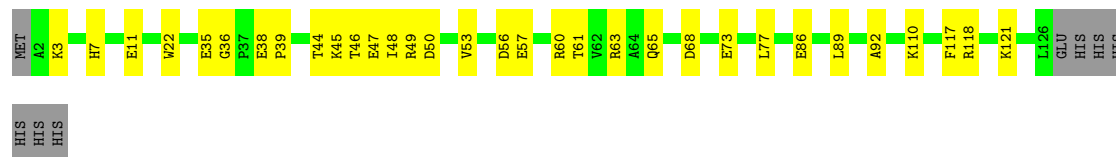
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	70	Total	O	0	0
			70	70		
6	B	65	Total	O	0	0
			65	65		

Note EDS was not executed.

- Chain A: 77% 14% 9%



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	52.24Å 52.24Å 203.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.34 – 1.65	Depositor
% Data completeness (in resolution range)	92.7 (36.34-1.65)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.209 , 0.225	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2250	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

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, A1D90, EDO, PEG

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	0.35	0/978	0.56	0/1322
1	B	0.29	0/1009	0.52	0/1365
All	All	0.32	0/1987	0.54	0/2687

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	960	0	940	17	2
1	B	991	0	978	38	2
2	A	28	33	0	1	0
2	B	28	33	0	3	0
3	A	4	6	6	0	1
3	B	4	6	6	0	0
4	A	7	10	10	0	0
5	B	5	0	0	1	0
6	A	70	0	0	2	1
6	B	65	0	0	8	1
All	All	2162	88	1940	50	4

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 13.

All (50) 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:39:PRO:HB3	1:B:121:LYS:HG2	1.29	1.12
1:A:97:VAL:O	1:A:121:LYS:NZ	2.10	0.83
1:B:56:ASP:O	6:B:301:HOH:O	2.00	0.79
5:B:203:SO4:O3	6:B:302:HOH:O	2.02	0.78
1:B:73:GLU:CD	6:B:303:HOH:O	2.26	0.78
1:B:73:GLU:OE2	6:B:303:HOH:O	2.03	0.75
1:A:91:LYS:O	6:A:301:HOH:O	2.05	0.74
1:B:45:LYS:O	1:B:48[B]:ILE:HG12	1.88	0.74
1:B:57:GLU:O	1:B:60:ARG:HG3	1.90	0.70
1:B:63:ARG:HH11	1:B:63:ARG:HG3	1.57	0.69
1:A:122:LEU:O	6:A:302:HOH:O	2.12	0.67
1:A:41:ARG:HD2	1:A:43:TYR:CZ	2.31	0.66
1:B:118:ARG:HH22	2:B:201:A1D90:C16	2.08	0.66
1:B:50:ASP:O	1:B:53:VAL:HG22	1.97	0.65
1:B:86:GLU:OE2	6:B:304:HOH:O	2.14	0.64
1:B:46[B]:THR:HG23	1:B:47:GLU:HG3	1.79	0.64
1:A:57:GLU:O	1:A:61[B]:THR:HG23	1.99	0.62
1:B:73:GLU:CG	6:B:303:HOH:O	2.47	0.62
1:B:63:ARG:HG3	1:B:63:ARG:NH1	2.15	0.62
1:B:45:LYS:HG3	1:B:48[B]:ILE:HD11	1.81	0.61
1:A:77:LEU:CD2	1:B:77:LEU:HD22	2.31	0.61
1:B:22:TRP:CD1	1:B:48[B]:ILE:HG13	2.37	0.59
1:A:121:LYS:HA	1:B:36:GLY:O	2.03	0.59
1:B:49:ARG:O	1:B:53:VAL:HG13	2.03	0.59
1:B:118:ARG:NH1	2:B:201:A1D90:O07	2.35	0.59
1:A:121:LYS:HB2	1:B:39:PRO:HB3	1.85	0.59
1:B:44:THR:O	1:B:48[B]:ILE:HG23	2.02	0.58
1:B:63:ARG:NH2	6:B:309:HOH:O	2.37	0.56
1:B:110:LYS:HE3	6:B:311:HOH:O	2.07	0.53
1:B:45:LYS:HA	1:B:48[B]:ILE:CD1	2.40	0.52
1:A:61[B]:THR:O	1:A:65:GLN:HG3	2.10	0.51
1:B:45:LYS:HA	1:B:48[B]:ILE:HD13	1.93	0.51
1:B:61:THR:O	1:B:65:GLN:HG3	2.11	0.50
1:B:60:ARG:HD3	1:B:61:THR:N	2.26	0.50
1:A:61[A]:THR:O	1:A:65:GLN:HG3	2.11	0.49
1:A:56:ASP:O	1:A:60:ARG:HG3	2.14	0.48
1:B:22:TRP:HD1	1:B:48[B]:ILE:HG13	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:VAL:HB	1:A:121:LYS:HZ2	1.81	0.45
1:B:7:HIS:NE2	1:B:73:GLU:CD	2.75	0.44
1:A:33:ALA:HA	2:A:201:A1D90:O09	2.17	0.44
1:B:68:ASP:OD1	1:B:68:ASP:N	2.51	0.44
1:A:117:PHE:HB3	1:B:117:PHE:HB3	2.00	0.44
1:B:46[B]:THR:HG23	1:B:47:GLU:N	2.33	0.44
1:B:89:LEU:HD13	2:B:201:A1D90:C24	2.48	0.43
1:A:90:ARG:HA	1:A:95:GLU:O	2.19	0.42
1:B:63:ARG:HA	1:B:63:ARG:HD2	1.84	0.42
1:A:120:ALA:O	1:B:35:GLU:HA	2.20	0.41
1:B:3:LYS:O	1:B:3:LYS:HG2	2.20	0.41
1:B:60:ARG:HD3	1:B:60:ARG:C	2.47	0.40
1:B:38:GLU:N	1:B:39:PRO:HA	2.37	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:319:HOH:O	6:B:311:HOH:O[1_455]	1.57	0.63
3:A:202:EDO:O2	3:A:202:EDO:O2[7_555]	1.77	0.43
1:A:41:ARG:NH1	1:B:11:GLU:OE2[5_544]	2.15	0.05
1:A:121:LYS:NZ	1:B:92:ALA:O[5_544]	2.18	0.02

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	121/133 (91%)	119 (98%)	2 (2%)	0	100	100
1	B	126/133 (95%)	123 (98%)	3 (2%)	0	100	100
All	All	247/266 (93%)	242 (98%)	5 (2%)	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	100/108 (93%)	100 (100%)	0	100	100
1	B	102/108 (94%)	102 (100%)	0	100	100
All	All	202/216 (94%)	202 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	B	8	GLN
1	B	54	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](#)

6 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
4	PEG	A	203	-	6,6,6	0.21	0	5,5,5	0.13	0
3	EDO	A	202	-	3,3,3	0.39	0	2,2,2	0.39	0
2	A1D90	B	201	-	29,29,29	2.79	8 (27%)	24,41,41	2.83	9 (37%)
3	EDO	B	202	-	3,3,3	0.50	0	2,2,2	0.32	0
2	A1D90	A	201	-	29,29,29	2.82	8 (27%)	24,41,41	2.90	9 (37%)
5	SO4	B	203	-	4,4,4	0.21	0	6,6,6	0.17	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
4	PEG	A	203	-	-	2/4/4/4	-
3	EDO	A	202	-	-	0/1/1/1	-
2	A1D90	B	201	-	-	5/20/50/50	0/2/2/2
3	EDO	B	202	-	-	0/1/1/1	-
2	A1D90	A	201	-	-	2/20/50/50	0/2/2/2

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	201	A1D90	O07-C08	9.75	1.49	1.34
2	A	201	A1D90	O07-C06	-9.58	1.33	1.46
2	A	201	A1D90	O19-C23	-6.32	1.39	1.46
2	A	201	A1D90	C22-C23	5.02	1.61	1.53
2	B	201	A1D90	C26-C27	4.48	1.60	1.52
2	B	201	A1D90	O19-C23	-4.25	1.41	1.46
2	B	201	A1D90	C05-C01	-4.17	1.46	1.54
2	B	201	A1D90	C14-C12	-4.16	1.40	1.50
2	B	201	A1D90	C01-C06	3.92	1.58	1.53
2	B	201	A1D90	C22-C23	3.78	1.59	1.53
2	A	201	A1D90	C26-C27	3.68	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	201	A1D90	O07-C08	3.61	1.40	1.34
2	A	201	A1D90	C14-C12	-3.33	1.42	1.50
2	B	201	A1D90	C01-C02	3.06	1.56	1.51
2	A	201	A1D90	C03-C02	2.87	1.56	1.51
2	A	201	A1D90	C01-C02	2.40	1.55	1.51

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	201	A1D90	C15-C13-C12	-10.38	114.12	126.59
2	B	201	A1D90	C15-C13-C12	-8.24	116.69	126.59
2	B	201	A1D90	C20-C15-C13	5.57	122.63	113.02
2	B	201	A1D90	C14-C12-C16	-4.55	108.39	116.02
2	A	201	A1D90	O19-C23-C25	-4.38	102.72	108.44
2	A	201	A1D90	C20-C15-C13	3.87	119.70	113.02
2	B	201	A1D90	O07-C06-C01	3.68	115.80	110.23
2	B	201	A1D90	C24-C21-C22	-3.52	108.60	113.76
2	B	201	A1D90	C14-C12-C13	3.24	130.83	123.64
2	B	201	A1D90	C10-C05-C01	-3.15	107.37	114.02
2	A	201	A1D90	C05-C01-C02	-2.94	109.07	112.82
2	A	201	A1D90	O09-C08-C03	2.60	124.30	118.37
2	A	201	A1D90	O07-C06-C01	2.56	114.10	110.23
2	B	201	A1D90	O09-C08-C03	2.36	123.76	118.37
2	B	201	A1D90	O19-C23-C25	-2.24	105.52	108.44
2	A	201	A1D90	O19-C20-C21	2.08	106.77	103.45
2	A	201	A1D90	C14-C12-C13	2.07	128.24	123.64
2	A	201	A1D90	C14-C12-C16	-2.07	112.56	116.02

There are no chirality outliers.

All (9) torsion outliers are listed below:

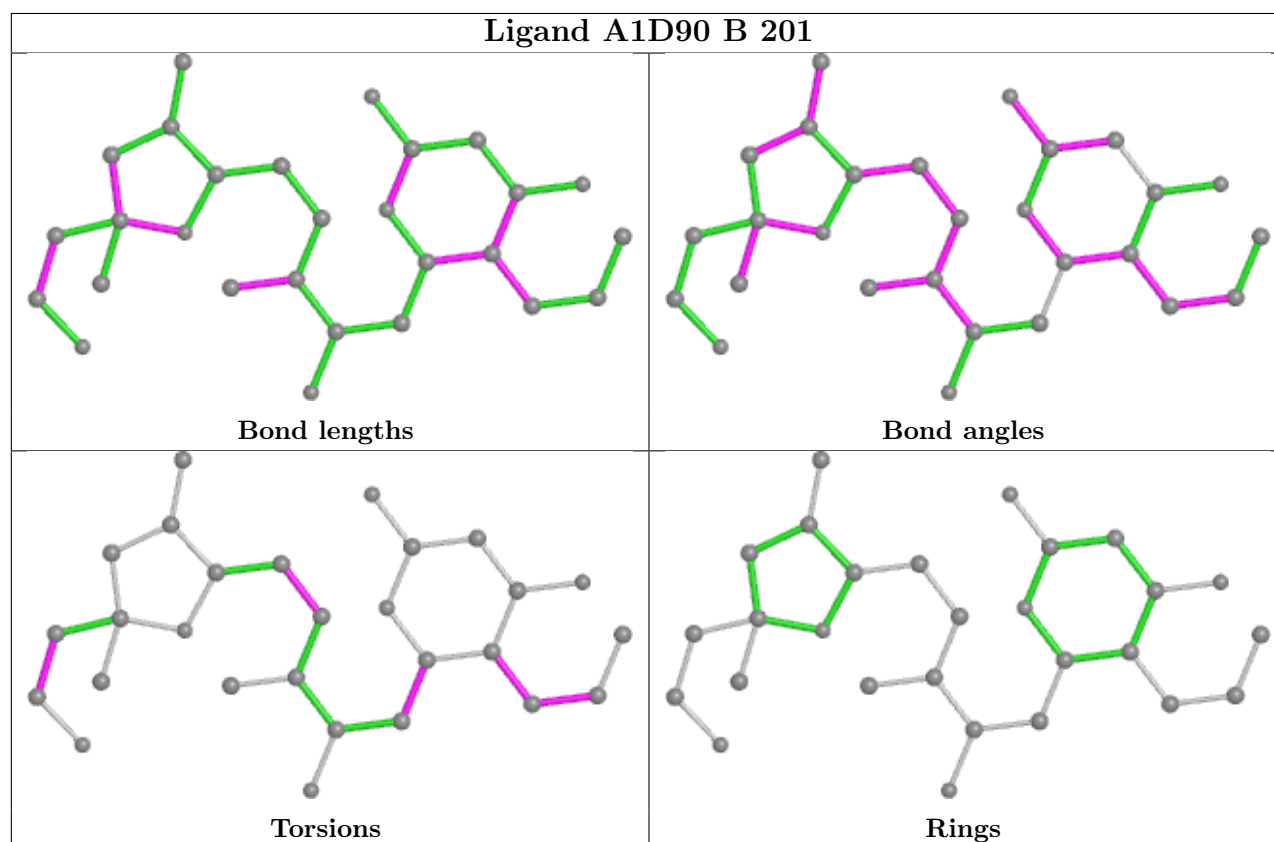
Mol	Chain	Res	Type	Atoms
2	A	201	A1D90	C23-C26-C27-O28
2	B	201	A1D90	C12-C13-C15-C20
2	B	201	A1D90	C23-C26-C27-O28
4	A	203	PEG	O2-C3-C4-O4
2	A	201	A1D90	C01-C05-C10-C11
4	A	203	PEG	O1-C1-C2-O2
2	B	201	A1D90	C01-C05-C10-C11
2	B	201	A1D90	C06-C01-C05-C10
2	B	201	A1D90	C01-C06-C18-C16

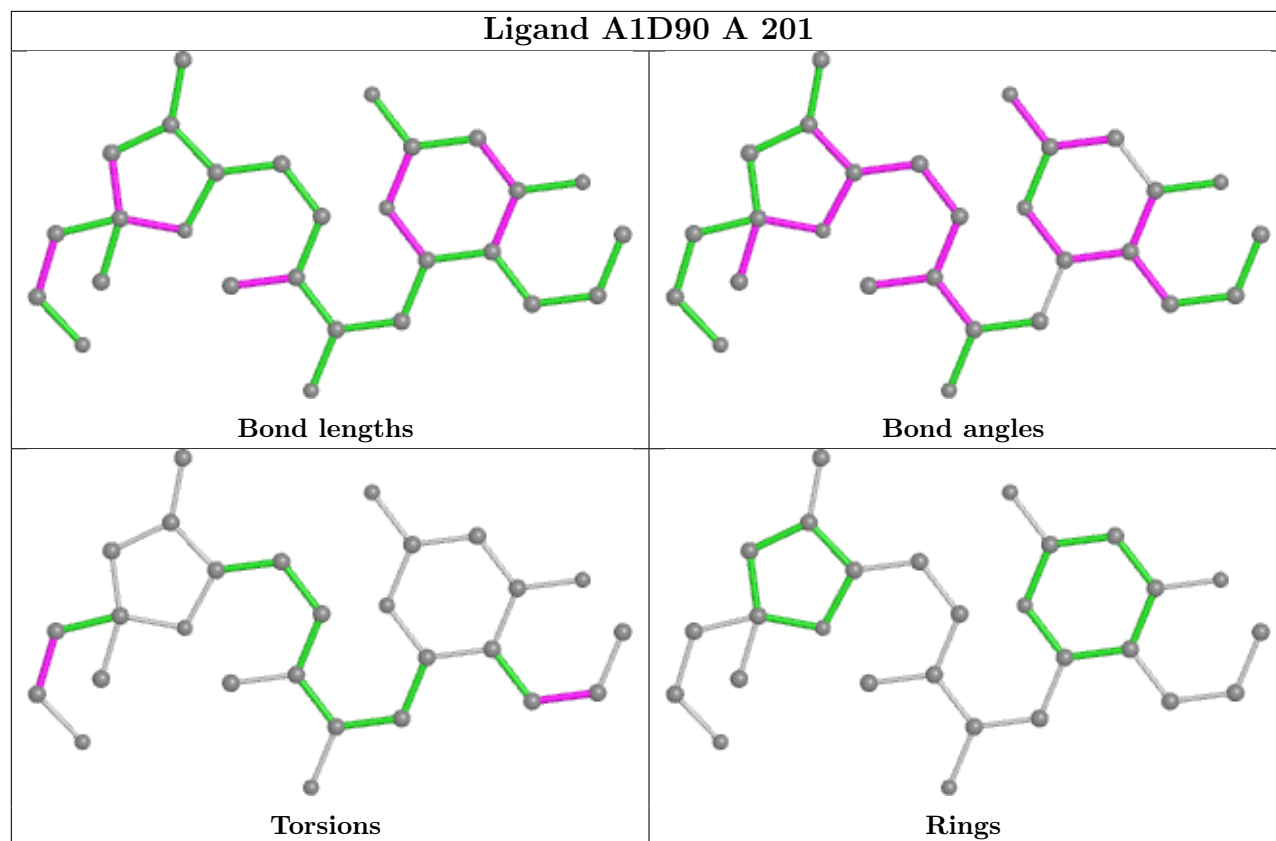
There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	202	EDO	0	1
2	B	201	A1D90	3	0
2	A	201	A1D90	1	0
5	B	203	SO4	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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.