Full wwPDB X-ray Structure Validation Report (i)

Jun 23, 2024 – 05:54 AM EDT

PDB ID	:	5MZN
Title	:	Helicase Sen1
Authors	:	Leonaite, B.; Basquin, J.; Conti, E.
Deposited on	:	2017-02-01
Resolution	:	1.79 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 (i) 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 (i)) were used in the production of this report:

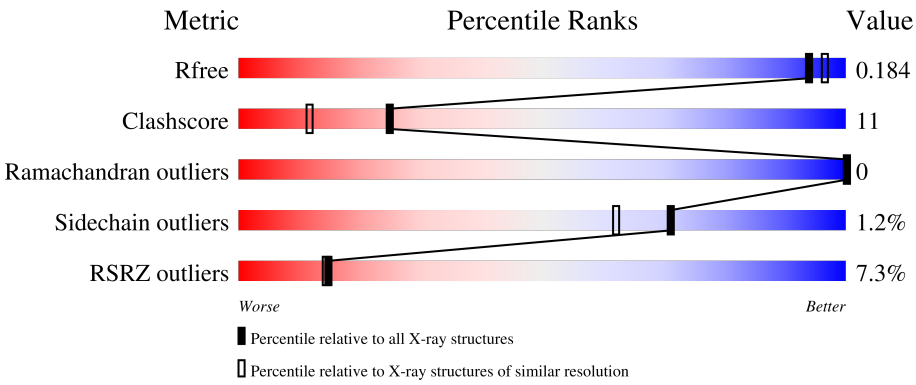
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.79 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}	130704	9185 (1.80-1.76)
Clashscore	141614	10184 (1.80-1.76)
Ramachandran outliers	138981	10051 (1.80-1.76)
Sidechain outliers	138945	10050 (1.80-1.76)
RSRZ outliers	127900	9032 (1.80-1.76)

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 <=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	749	7% 79% 12% 9%

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
3	EDO	A	2022	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	EDO	A	2028	-	-	X	X
3	EDO	A	2032	-	-	-	X
3	EDO	A	2036	-	-	-	X
3	EDO	A	2038	-	-	-	X
3	EDO	A	2039	-	-	-	X
4	GOL	A	2044	-	-	-	X
4	GOL	A	2046	-	-	-	X
4	GOL	A	2047	-	-	X	X
4	GOL	A	2048	-	-	-	X
4	GOL	A	2052	-	-	-	X
5	PEG	A	2054	-	-	X	-
5	PEG	A	2056	-	-	X	X
5	PEG	A	2057	-	-	X	-
5	PEG	A	2059	-	-	X	-
5	PEG	A	2062	-	-	-	X
5	PEG	A	2063	-	-	-	X
5	PEG	A	2064	-	-	-	X

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 12935 atoms, of which 5965 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.

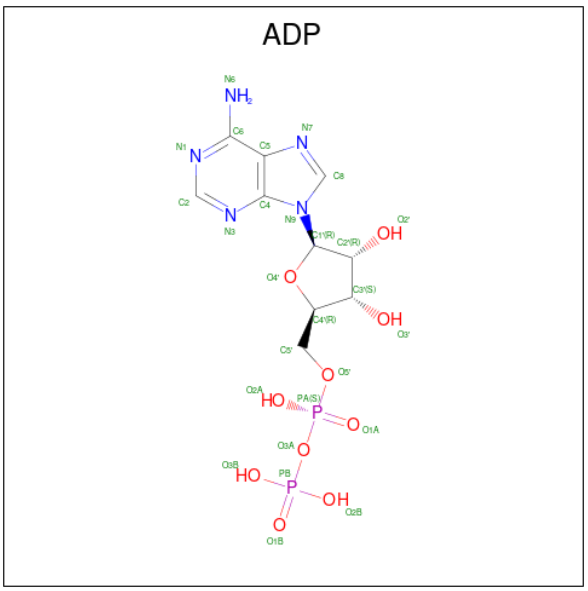
- Molecule 1 is a protein called Helicase SEN1,Helicase SEN1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	682	11057	3513	5514	957	1041	32	0	20	0

There are 7 discrepancies between the modelled and reference sequences:

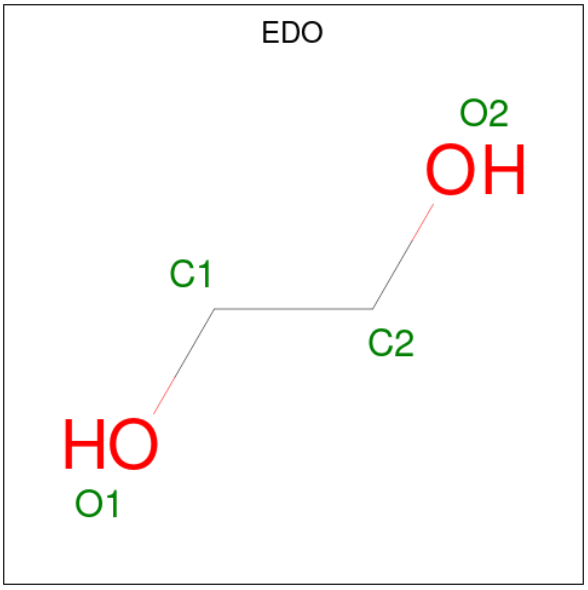
Chain	Residue	Modelled	Actual	Comment	Reference
A	1092	MET	-	initiating methionine	UNP Q00416
A	1093	ASN	-	expression tag	UNP Q00416
A	1094	SER	-	expression tag	UNP Q00416
A	1535	GLY	-	linker	UNP Q00416
A	1536	SER	-	linker	UNP Q00416
A	1537	GLY	-	linker	UNP Q00416
A	1538	SER	-	linker	UNP Q00416

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	H	N	O	P		
2	A	1	39	10	12	5	10	2	0	0

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	8	2	4	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0

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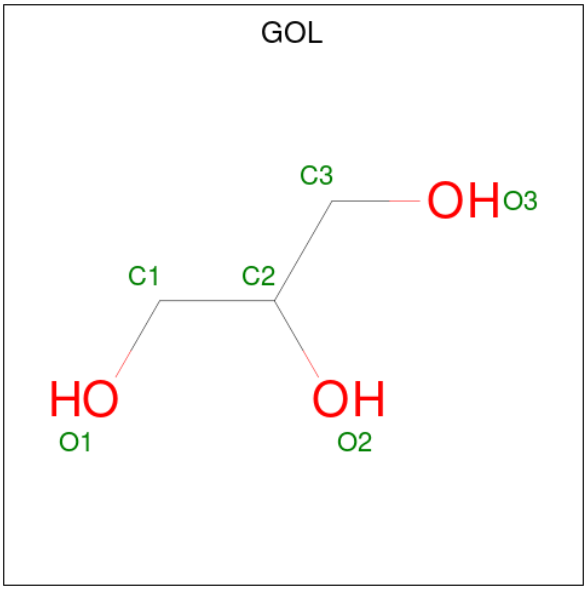
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total	C	H	O		
3	A	1	10	2	6	2	0	0
			Total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

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
4	A	1	14	3	8	3	0	0
			Total	C	H	O		
4	A	1	14	3	8	3	0	0
			Total	C	H	O		
4	A	1	12	3	6	3	0	0

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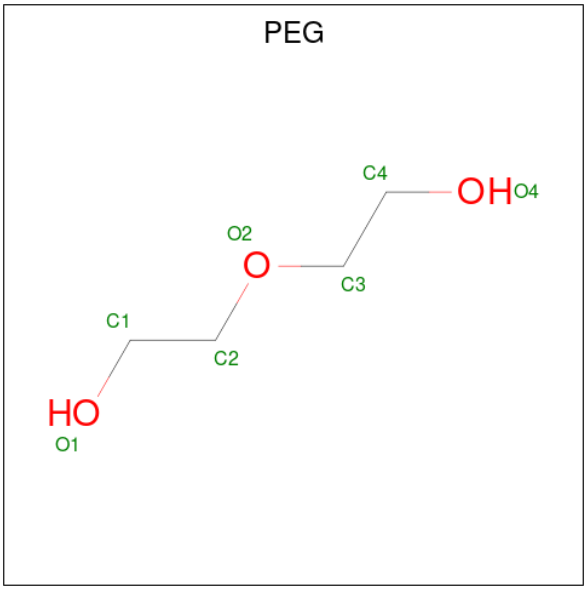
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
4	A	1	14	3	8	3	0	0
			Total	C	H	O		
4	A	1	13	3	7	3	0	0
			Total	C	H	O		
4	A	1	13	3	7	3	0	0
			Total	C	H	O		
4	A	1	14	3	8	3	0	0
			Total	C	H	O		
4	A	1	14	3	8	3	0	0
			Total	C	H	O		
4	A	1	14	3	8	3	0	0
			Total	C	H	O		
4	A	1	14	3	8	3	0	0
			Total	C	H	O		
4	A	1	14	3	8	3	0	0
			Total	C	H	O		
4	A	1	13	3	7	3	0	0

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



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

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

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
6	A	1	1	1	0	0

- Molecule 7 is water.

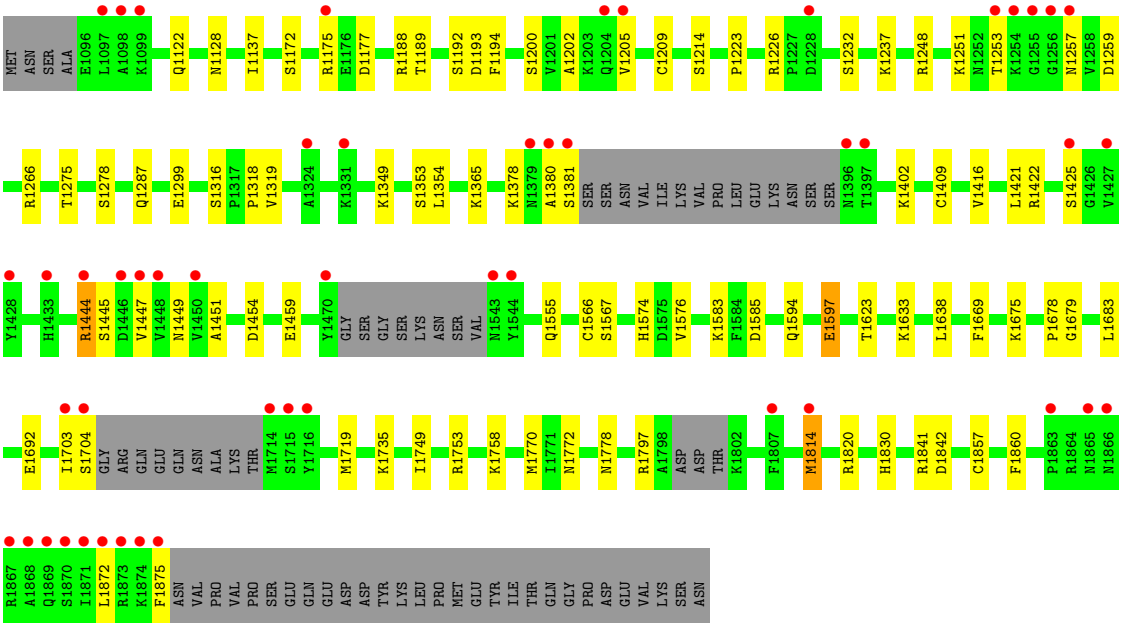
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
7	A	1090	1090	1090	0	0

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: Helicase SEN1,Helicase SEN1

Chain A: 7% 79% 12% 9%



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	90.28Å 171.94Å 69.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.39 – 1.79 48.39 – 1.79	Depositor EDS
% Data completeness (in resolution range)	98.3 (48.39-1.79) 98.3 (48.39-1.79)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.01 (at 1.78Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.153 , 0.184 0.154 , 0.184	Depositor DCC
R_{free} test set	5079 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å^2)	29.8	Xtriage
Anisotropy	0.272	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 62.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	12935	wwPDB-VP
Average B, all atoms (Å^2)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.56% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: ADP, EDO, GOL, PEG, MG

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	0.52	1/5654 (0.0%)	0.66	2/7628 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1597	GLU	CB-CG	-5.12	1.42	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	1814[A]	MET	CG-SD-CE	6.92	111.27	100.20
1	A	1814[B]	MET	CG-SD-CE	6.92	111.27	100.20

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	5543	5514	5501	109	1
2	A	27	12	12	0	0
3	A	160	238	238	22	1
4	A	72	91	91	15	0
5	A	77	110	110	33	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	1	0	0	0	0
7	A	1090	0	0	54	22
All	All	6970	5965	5952	128	23

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 11.

All (128) 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:1683[A]:LEU:HB2	4:A:2047:GOL:O3	1.60	1.01
1:A:1257:ASN:OD1	7:A:2101:HOH:O	1.81	0.98
1:A:1857:CYS:SG	7:A:2735:HOH:O	2.26	0.94
1:A:1209:CYS:O	7:A:2102:HOH:O	1.89	0.91
1:A:1683[A]:LEU:CB	4:A:2047:GOL:O3	2.19	0.89
4:A:2050:GOL:O2	7:A:2104:HOH:O	1.92	0.87
1:A:1692:GLU:OE2	7:A:2103:HOH:O	1.91	0.87
1:A:1192:SER:O	1:A:1266[B]:ARG:NH2	2.08	0.85
1:A:1402:LYS:NZ	7:A:2116:HOH:O	2.12	0.82
1:A:1299:GLU:O	3:A:2038:EDO:O1	1.97	0.80
1:A:1365[A]:LYS:NZ	3:A:2008:EDO:O2	2.15	0.80
1:A:1266[A]:ARG:NH1	7:A:2118:HOH:O	2.14	0.80
4:A:2049:GOL:O3	7:A:2105:HOH:O	1.99	0.79
1:A:1841[B]:ARG:NH2	7:A:2120:HOH:O	2.16	0.78
4:A:2048:GOL:O2	7:A:2106:HOH:O	2.00	0.77
1:A:1193:ASP:OD2	3:A:2039:EDO:O1	2.03	0.77
1:A:1425:SER:O	7:A:2107:HOH:O	2.03	0.76
1:A:1319:VAL:O	7:A:2108:HOH:O	2.04	0.76
4:A:2052:GOL:O2	7:A:2109:HOH:O	2.05	0.74
4:A:2053:GOL:O1	7:A:2110:HOH:O	2.06	0.74
1:A:1175:ARG:NH1	7:A:2124:HOH:O	2.20	0.72
1:A:1253:THR:O	7:A:2111:HOH:O	2.08	0.72
1:A:1259:ASP:OD1	7:A:2113:HOH:O	2.09	0.71
3:A:2035:EDO:O1	7:A:2112:HOH:O	2.08	0.71
1:A:1583:LYS:NZ	7:A:2130:HOH:O	2.24	0.69
3:A:2014:EDO:O2	7:A:2114:HOH:O	2.09	0.69
1:A:1683[A]:LEU:N	4:A:2047:GOL:O3	2.29	0.65
1:A:1266[B]:ARG:NH2	7:A:2135:HOH:O	2.29	0.65
1:A:1623:THR:O	7:A:2119:HOH:O	2.15	0.63
1:A:1555:GLN:OE1	3:A:2023:EDO:O1	2.13	0.62
1:A:1122[A]:GLN:NE2	7:A:2143:HOH:O	2.32	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1194:PHE:O	1:A:1266[B]:ARG:NH1	2.34	0.61
1:A:1830:HIS:HB2	3:A:2029:EDO:C2	2.31	0.60
1:A:1679:GLY:O	4:A:2047:GOL:O3	2.15	0.60
3:A:2002:EDO:O2	5:A:2064:PEG:O4	2.02	0.60
5:A:2057:PEG:H32	7:A:2432:HOH:O	2.03	0.59
1:A:1449:ASN:OD1	1:A:1451:ALA:N	2.35	0.59
3:A:2018:EDO:H11	5:A:2054:PEG:H21	1.84	0.58
1:A:1128:ASN:HB2	5:A:2059:PEG:C4	2.34	0.58
1:A:1772:ASN:ND2	7:A:2140:HOH:O	2.31	0.57
1:A:1226:ARG:NH1	7:A:2146:HOH:O	2.34	0.56
1:A:1449:ASN:HB3	5:A:2054:PEG:O2	2.05	0.56
1:A:1422:ARG:HH21	3:A:2014:EDO:H21	1.71	0.55
5:A:2056:PEG:O1	7:A:2115:HOH:O	2.09	0.55
1:A:1444:ARG:NE	7:A:2160:HOH:O	2.38	0.55
1:A:1758:LYS:NZ	7:A:2163:HOH:O	2.39	0.55
1:A:1841[B]:ARG:NH2	7:A:2164:HOH:O	2.39	0.55
1:A:1128:ASN:HA	5:A:2059:PEG:H31	1.88	0.55
1:A:1449:ASN:HB2	5:A:2054:PEG:H12	1.90	0.54
1:A:1719:MET:HG2	1:A:1758:LYS:HE2	1.89	0.54
5:A:2061:PEG:O1	5:A:2062:PEG:O4	2.09	0.54
1:A:1200:SER:O	7:A:2123:HOH:O	2.18	0.54
1:A:1237:LYS:HE3	5:A:2059:PEG:H32	1.89	0.53
1:A:1820:ARG:HH12	5:A:2057:PEG:H41	1.74	0.53
1:A:1669:PHE:CZ	1:A:1814[A]:MET:HE3	2.44	0.53
1:A:1365[B]:LYS:NZ	3:A:2008:EDO:O1	2.34	0.53
1:A:1669:PHE:CZ	1:A:1814[A]:MET:CE	2.93	0.52
1:A:1683[A]:LEU:CA	4:A:2047:GOL:O3	2.58	0.52
1:A:1237:LYS:CE	5:A:2059:PEG:H41	2.40	0.52
1:A:1422:ARG:HH21	3:A:2014:EDO:C2	2.23	0.52
1:A:1633:LYS:NZ	7:A:2175:HOH:O	2.42	0.51
1:A:1669:PHE:HZ	1:A:1814[A]:MET:CE	2.24	0.51
1:A:1703:ILE:O	7:A:2122:HOH:O	2.17	0.51
1:A:1841[B]:ARG:NE	7:A:2177:HOH:O	2.43	0.51
1:A:1223:PRO:O	3:A:2036:EDO:O2	2.17	0.50
1:A:1353:SER:HA	3:A:2037:EDO:H11	1.94	0.50
1:A:1237:LYS:HE3	5:A:2059:PEG:C3	2.42	0.49
5:A:2063:PEG:C3	7:A:2286:HOH:O	2.60	0.49
1:A:1192:SER:O	1:A:1266[B]:ARG:CZ	2.60	0.49
1:A:1354:LEU:H	3:A:2037:EDO:C1	2.25	0.49
1:A:1841[B]:ARG:NH1	7:A:2184:HOH:O	2.46	0.48
1:A:1830:HIS:HB2	3:A:2029:EDO:C1	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1574:HIS:CE1	1:A:1576:VAL:HG13	2.49	0.47
1:A:1597:GLU:HG3	1:A:1638:LEU:HB2	1.97	0.47
1:A:1237:LYS:HE3	5:A:2059:PEG:H41	1.97	0.47
1:A:1447:VAL:HB	5:A:2054:PEG:H21	1.95	0.47
1:A:1749:ILE:HA	1:A:1778:ASN:O	2.15	0.47
5:A:2056:PEG:O2	7:A:2125:HOH:O	2.20	0.47
1:A:1237:LYS:HE3	5:A:2059:PEG:C4	2.45	0.46
1:A:1416:VAL:HG13	1:A:1566[B]:CYS:SG	2.55	0.46
1:A:1319:VAL:HG22	7:A:2621:HOH:O	2.16	0.46
3:A:2028:EDO:H22	7:A:2181:HOH:O	2.16	0.46
3:A:2006:EDO:C1	7:A:2436:HOH:O	2.62	0.46
1:A:1820:ARG:HH22	5:A:2057:PEG:H31	1.80	0.46
1:A:1449:ASN:CB	5:A:2054:PEG:O2	2.64	0.45
1:A:1128:ASN:HB2	5:A:2059:PEG:H42	1.97	0.45
1:A:1675:LYS:HG2	5:A:2064:PEG:O2	2.17	0.45
1:A:1287:GLN:O	4:A:2053:GOL:H31	2.16	0.45
1:A:1214[B]:SER:OG	7:A:2126:HOH:O	2.21	0.45
1:A:1683[A]:LEU:HB2	4:A:2047:GOL:O2	2.17	0.45
1:A:1421:LEU:HD12	7:A:2765:HOH:O	2.17	0.45
1:A:1449:ASN:HD22	5:A:2054:PEG:H12	1.82	0.45
1:A:1814[B]:MET:HG2	7:A:2793:HOH:O	2.17	0.45
1:A:1447:VAL:O	5:A:2054:PEG:O4	2.13	0.44
1:A:1675:LYS:HG2	5:A:2064:PEG:C3	2.47	0.44
1:A:1841[B]:ARG:CZ	7:A:2184:HOH:O	2.65	0.44
1:A:1275:THR:HG1	1:A:1278:SER:HG	1.65	0.44
1:A:1797:ARG:CD	7:A:2242:HOH:O	2.65	0.44
1:A:1202:ALA:N	5:A:2056:PEG:H32	2.33	0.43
1:A:1454[A]:ASP:HA	1:A:1459:GLU:HG2	1.99	0.43
1:A:1188:ARG:HG3	7:A:2448:HOH:O	2.18	0.43
1:A:1585:ASP:OD2	4:A:2046:GOL:H32	2.18	0.43
1:A:1128:ASN:HB2	5:A:2059:PEG:O4	2.18	0.43
1:A:1860:PHE:HB2	1:A:1872:LEU:HD11	2.01	0.42
1:A:1316[A]:SER:OG	1:A:1349:LYS:HD2	2.19	0.42
1:A:1449:ASN:HD22	5:A:2054:PEG:C1	2.32	0.42
1:A:1594:GLN:NE2	7:A:2221:HOH:O	2.53	0.42
1:A:1248:ARG:HD3	7:A:2872:HOH:O	2.19	0.41
1:A:1380:ALA:O	1:A:1381:SER:C	2.58	0.41
1:A:1444:ARG:HG2	1:A:1444:ARG:HH11	1.85	0.41
1:A:1205:VAL:HG21	5:A:2056:PEG:H12	2.02	0.41
1:A:1137:ILE:HD12	7:A:2436:HOH:O	2.20	0.41
1:A:1251:LYS:HG2	7:A:2351:HOH:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1409:CYS:HA	1:A:1567:SER:O	2.21	0.41
3:A:2032:EDO:H11	5:A:2056:PEG:H11	2.03	0.41
1:A:1214[A]:SER:OG	1:A:1287:GLN:HG3	2.21	0.41
1:A:1128:ASN:OD1	5:A:2059:PEG:H21	2.21	0.41
1:A:1189:THR:HG23	4:A:2052:GOL:H2	2.03	0.41
1:A:1678:PRO:HD3	3:A:2015:EDO:H21	2.03	0.41
1:A:1318:PRO:HB3	7:A:2129:HOH:O	2.21	0.41
3:A:2028:EDO:H21	7:A:2487:HOH:O	2.21	0.41
1:A:1820:ARG:NH1	5:A:2057:PEG:H41	2.35	0.40
1:A:1445[B]:SER:HB2	5:A:2054:PEG:C3	2.52	0.40
1:A:1692:GLU:HG2	4:A:2048:GOL:C3	2.52	0.40
3:A:2028:EDO:C2	7:A:2487:HOH:O	2.69	0.40
1:A:1735:LYS:NZ	7:A:2150:HOH:O	2.35	0.40

All (23) 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)
7:A:2894:HOH:O	7:A:3004:HOH:O[1_556]	1.80	0.40
7:A:3020:HOH:O	7:A:3053:HOH:O[4_454]	1.83	0.37
7:A:2264:HOH:O	7:A:2285:HOH:O[4_455]	1.84	0.36
7:A:2375:HOH:O	7:A:2985:HOH:O[4_455]	1.85	0.35
7:A:2610:HOH:O	7:A:2653:HOH:O[4_555]	1.86	0.34
7:A:2649:HOH:O	7:A:2968:HOH:O[4_555]	1.86	0.34
7:A:2686:HOH:O	7:A:2767:HOH:O[1_554]	1.88	0.32
7:A:2120:HOH:O	7:A:2137:HOH:O[2_455]	1.90	0.30
7:A:2331:HOH:O	7:A:3080:HOH:O[4_454]	1.93	0.27
7:A:2611:HOH:O	7:A:2649:HOH:O[4_455]	1.94	0.26
7:A:2693:HOH:O	7:A:3031:HOH:O[4_555]	1.94	0.26
7:A:2996:HOH:O	7:A:3010:HOH:O[4_455]	1.94	0.26
7:A:2752:HOH:O	7:A:2754:HOH:O[1_556]	1.95	0.25
7:A:2895:HOH:O	7:A:2912:HOH:O[2_455]	1.96	0.24
7:A:2686:HOH:O	7:A:2987:HOH:O[1_554]	1.98	0.22
1:A:1842:ASP:OD1	3:A:2028:EDO:HO1[2_455]	1.41	0.19
7:A:2429:HOH:O	7:A:2507:HOH:O[1_554]	2.03	0.17
7:A:2106:HOH:O	7:A:2752:HOH:O[1_554]	2.04	0.16
7:A:3023:HOH:O	7:A:3091:HOH:O[4_455]	2.12	0.08
7:A:2285:HOH:O	7:A:3023:HOH:O[4_555]	2.17	0.03
7:A:2605:HOH:O	7:A:2757:HOH:O[4_455]	2.18	0.02
7:A:2693:HOH:O	7:A:2759:HOH:O[4_555]	2.18	0.02
7:A:2868:HOH:O	7:A:2870:HOH:O[4_554]	2.19	0.01

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	692/749 (92%)	678 (98%)	14 (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	612/670 (91%)	605 (99%)	7 (1%)	73 65

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1232	SER
1	A	1378	LYS
1	A	1444	ARG
1	A	1704	SER
1	A	1753	ARG
1	A	1770	MET
1	A	1875	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 65 ligands modelled in this entry, 1 is monoatomic - leaving 64 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	$\#\|Z\| > 2$
3	EDO	A	2009	-	3,3,3	0.46	0	2,2,2	0.38	0
5	PEG	A	2056	-	6,6,6	0.50	0	5,5,5	0.83	0
5	PEG	A	2063	-	6,6,6	0.49	0	5,5,5	0.38	0
4	GOL	A	2044	3	5,5,5	0.42	0	5,5,5	0.25	0
3	EDO	A	2011	-	3,3,3	0.51	0	2,2,2	0.06	0
3	EDO	A	2039	-	3,3,3	0.38	0	2,2,2	0.46	0
3	EDO	A	2006	-	3,3,3	0.56	0	2,2,2	0.39	0
4	GOL	A	2042	-	5,5,5	0.42	0	5,5,5	1.06	0
5	PEG	A	2055	-	6,6,6	0.44	0	5,5,5	0.80	0
3	EDO	A	2032	-	3,3,3	0.41	0	2,2,2	0.27	0
4	GOL	A	2043	-	5,5,5	0.45	0	5,5,5	0.54	0
3	EDO	A	2025	-	3,3,3	0.48	0	2,2,2	0.26	0
3	EDO	A	2037	-	3,3,3	0.27	0	2,2,2	0.54	0
5	PEG	A	2054	-	6,6,6	0.59	0	5,5,5	0.42	0
5	PEG	A	2058	-	6,6,6	0.53	0	5,5,5	0.43	0
3	EDO	A	2029	-	3,3,3	0.46	0	2,2,2	0.22	0
3	EDO	A	2016	-	3,3,3	0.46	0	2,2,2	0.29	0
3	EDO	A	2017	-	3,3,3	0.41	0	2,2,2	0.37	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	GOL	A	2045	-	5,5,5	0.46	0	5,5,5	0.35	0
3	EDO	A	2027	-	3,3,3	0.39	0	2,2,2	0.39	0
3	EDO	A	2013	-	3,3,3	0.46	0	2,2,2	0.41	0
3	EDO	A	2036	-	3,3,3	0.45	0	2,2,2	0.26	0
3	EDO	A	2028	-	3,3,3	0.34	0	2,2,2	0.42	0
4	GOL	A	2049	-	5,5,5	0.39	0	5,5,5	0.23	0
4	GOL	A	2048	-	5,5,5	0.35	0	5,5,5	0.30	0
3	EDO	A	2005	-	3,3,3	0.48	0	2,2,2	0.37	0
2	ADP	A	2001	6	24,29,29	0.98	0	29,45,45	1.28	2 (6%)
3	EDO	A	2010	-	3,3,3	0.41	0	2,2,2	0.70	0
3	EDO	A	2030	-	3,3,3	0.51	0	2,2,2	0.21	0
3	EDO	A	2021	4	3,3,3	0.46	0	2,2,2	0.39	0
4	GOL	A	2052	-	5,5,5	0.38	0	5,5,5	0.20	0
3	EDO	A	2035	-	3,3,3	0.46	0	2,2,2	0.30	0
4	GOL	A	2051	-	5,5,5	0.38	0	5,5,5	0.53	0
3	EDO	A	2003	-	3,3,3	0.38	0	2,2,2	0.28	0
3	EDO	A	2015	-	3,3,3	0.52	0	2,2,2	0.31	0
4	GOL	A	2053	1	5,5,5	0.62	0	5,5,5	0.70	0
5	PEG	A	2057	-	6,6,6	0.51	0	5,5,5	0.62	0
5	PEG	A	2062	-	6,6,6	0.49	0	5,5,5	0.35	0
3	EDO	A	2034	-	3,3,3	0.49	0	2,2,2	0.05	0
3	EDO	A	2024	-	3,3,3	0.39	0	2,2,2	0.38	0
5	PEG	A	2060	-	6,6,6	0.49	0	5,5,5	0.31	0
3	EDO	A	2019	-	3,3,3	0.52	0	2,2,2	0.08	0
3	EDO	A	2008	-	3,3,3	0.43	0	2,2,2	0.34	0
3	EDO	A	2020	-	3,3,3	0.45	0	2,2,2	0.25	0
3	EDO	A	2002	-	3,3,3	0.48	0	2,2,2	0.20	0
3	EDO	A	2038	-	3,3,3	0.42	0	2,2,2	0.43	0
3	EDO	A	2040	-	3,3,3	0.40	0	2,2,2	0.45	0
4	GOL	A	2050	-	5,5,5	0.38	0	5,5,5	0.10	0
5	PEG	A	2059	-	6,6,6	0.52	0	5,5,5	0.35	0
5	PEG	A	2061	-	6,6,6	0.54	0	5,5,5	0.46	0
3	EDO	A	2022	-	3,3,3	0.42	0	2,2,2	0.56	0
3	EDO	A	2018	-	3,3,3	0.48	0	2,2,2	0.41	0
3	EDO	A	2031	-	3,3,3	0.48	0	2,2,2	0.21	0
3	EDO	A	2007	-	3,3,3	0.58	0	2,2,2	0.30	0
3	EDO	A	2041	-	3,3,3	0.42	0	2,2,2	0.27	0
3	EDO	A	2004	-	3,3,3	0.46	0	2,2,2	0.35	0
3	EDO	A	2033	-	3,3,3	0.50	0	2,2,2	0.22	0
5	PEG	A	2064	-	6,6,6	0.53	0	5,5,5	0.50	0
3	EDO	A	2026	-	3,3,3	0.57	0	2,2,2	0.24	0
3	EDO	A	2014	-	3,3,3	0.41	0	2,2,2	0.35	0
3	EDO	A	2023	-	3,3,3	0.45	0	2,2,2	0.33	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	GOL	A	2047	1	5,5,5	0.32	0	5,5,5	0.26	0
4	GOL	A	2046	1	5,5,5	0.33	0	5,5,5	0.23	0
3	EDO	A	2012	-	3,3,3	0.36	0	2,2,2	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
3	EDO	A	2009	-	-	1/1/1/1	-
5	PEG	A	2056	-	-	1/4/4/4	-
5	PEG	A	2063	-	-	2/4/4/4	-
4	GOL	A	2044	3	-	0/4/4/4	-
3	EDO	A	2011	-	-	0/1/1/1	-
3	EDO	A	2039	-	-	0/1/1/1	-
3	EDO	A	2006	-	-	0/1/1/1	-
4	GOL	A	2042	-	-	4/4/4/4	-
5	PEG	A	2055	-	-	2/4/4/4	-
3	EDO	A	2032	-	-	1/1/1/1	-
4	GOL	A	2043	-	-	2/4/4/4	-
3	EDO	A	2025	-	-	0/1/1/1	-
3	EDO	A	2037	-	-	1/1/1/1	-
5	PEG	A	2054	-	-	2/4/4/4	-
5	PEG	A	2058	-	-	2/4/4/4	-
3	EDO	A	2029	-	-	1/1/1/1	-
3	EDO	A	2016	-	-	1/1/1/1	-
3	EDO	A	2017	-	-	0/1/1/1	-
4	GOL	A	2045	-	-	2/4/4/4	-
3	EDO	A	2027	-	-	0/1/1/1	-
3	EDO	A	2013	-	-	0/1/1/1	-
3	EDO	A	2036	-	-	0/1/1/1	-
3	EDO	A	2028	-	-	0/1/1/1	-
4	GOL	A	2049	-	-	2/4/4/4	-
4	GOL	A	2048	-	-	2/4/4/4	-
3	EDO	A	2005	-	-	0/1/1/1	-
2	ADP	A	2001	6	-	0/12/32/32	0/3/3/3
3	EDO	A	2010	-	-	1/1/1/1	-
3	EDO	A	2030	-	-	1/1/1/1	-
3	EDO	A	2021	4	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	A	2052	-	-	4/4/4/4	-
3	EDO	A	2035	-	-	0/1/1/1	-
4	GOL	A	2051	-	-	3/4/4/4	-
3	EDO	A	2003	-	-	0/1/1/1	-
3	EDO	A	2015	-	-	0/1/1/1	-
4	GOL	A	2053	1	-	4/4/4/4	-
5	PEG	A	2057	-	-	2/4/4/4	-
5	PEG	A	2062	-	-	1/4/4/4	-
3	EDO	A	2034	-	-	0/1/1/1	-
3	EDO	A	2024	-	-	0/1/1/1	-
5	PEG	A	2060	-	-	1/4/4/4	-
3	EDO	A	2019	-	-	0/1/1/1	-
3	EDO	A	2008	-	-	0/1/1/1	-
3	EDO	A	2020	-	-	1/1/1/1	-
3	EDO	A	2002	-	-	0/1/1/1	-
3	EDO	A	2038	-	-	1/1/1/1	-
3	EDO	A	2040	-	-	0/1/1/1	-
4	GOL	A	2050	-	-	3/4/4/4	-
5	PEG	A	2059	-	-	1/4/4/4	-
5	PEG	A	2061	-	-	0/4/4/4	-
3	EDO	A	2022	-	-	1/1/1/1	-
3	EDO	A	2018	-	-	0/1/1/1	-
3	EDO	A	2031	-	-	0/1/1/1	-
3	EDO	A	2007	-	-	0/1/1/1	-
3	EDO	A	2041	-	-	0/1/1/1	-
3	EDO	A	2004	-	-	1/1/1/1	-
3	EDO	A	2033	-	-	1/1/1/1	-
5	PEG	A	2064	-	-	4/4/4/4	-
3	EDO	A	2026	-	-	1/1/1/1	-
3	EDO	A	2014	-	-	0/1/1/1	-
3	EDO	A	2023	-	-	0/1/1/1	-
4	GOL	A	2047	1	-	2/4/4/4	-
4	GOL	A	2046	1	-	3/4/4/4	-
3	EDO	A	2012	-	-	0/1/1/1	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	2001	ADP	N6-C6-N1	3.13	125.06	118.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	2001	ADP	N3-C2-N1	-2.56	124.67	128.68

There are no chirality outliers.

All (62) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	2042	GOL	O1-C1-C2-C3
4	A	2042	GOL	C1-C2-C3-O3
4	A	2043	GOL	C1-C2-C3-O3
4	A	2046	GOL	O1-C1-C2-C3
4	A	2050	GOL	C1-C2-C3-O3
4	A	2051	GOL	C1-C2-C3-O3
4	A	2052	GOL	C1-C2-C3-O3
4	A	2053	GOL	O1-C1-C2-C3
5	A	2064	PEG	O2-C3-C4-O4
5	A	2059	PEG	O1-C1-C2-O2
5	A	2063	PEG	O2-C3-C4-O4
3	A	2037	EDO	O1-C1-C2-O2
5	A	2054	PEG	O2-C3-C4-O4
4	A	2045	GOL	C1-C2-C3-O3
4	A	2047	GOL	C1-C2-C3-O3
4	A	2048	GOL	C1-C2-C3-O3
4	A	2053	GOL	C1-C2-C3-O3
4	A	2042	GOL	O1-C1-C2-O2
4	A	2051	GOL	O2-C2-C3-O3
4	A	2053	GOL	O1-C1-C2-O2
3	A	2004	EDO	O1-C1-C2-O2
3	A	2010	EDO	O1-C1-C2-O2
3	A	2026	EDO	O1-C1-C2-O2
3	A	2029	EDO	O1-C1-C2-O2
3	A	2033	EDO	O1-C1-C2-O2
3	A	2038	EDO	O1-C1-C2-O2
5	A	2064	PEG	C1-C2-O2-C3
4	A	2042	GOL	O2-C2-C3-O3
4	A	2043	GOL	O2-C2-C3-O3
4	A	2046	GOL	O1-C1-C2-O2
4	A	2048	GOL	O2-C2-C3-O3
4	A	2052	GOL	O1-C1-C2-O2
4	A	2052	GOL	O2-C2-C3-O3
4	A	2045	GOL	O2-C2-C3-O3
4	A	2047	GOL	O2-C2-C3-O3
4	A	2049	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
4	A	2050	GOL	O1-C1-C2-O2
5	A	2062	PEG	O2-C3-C4-O4
5	A	2058	PEG	C1-C2-O2-C3
5	A	2063	PEG	C1-C2-O2-C3
4	A	2050	GOL	O2-C2-C3-O3
4	A	2053	GOL	O2-C2-C3-O3
5	A	2058	PEG	O2-C3-C4-O4
5	A	2064	PEG	O1-C1-C2-O2
5	A	2056	PEG	C1-C2-O2-C3
3	A	2016	EDO	O1-C1-C2-O2
3	A	2022	EDO	O1-C1-C2-O2
3	A	2030	EDO	O1-C1-C2-O2
3	A	2032	EDO	O1-C1-C2-O2
5	A	2055	PEG	O1-C1-C2-O2
5	A	2064	PEG	C4-C3-O2-C2
5	A	2057	PEG	C1-C2-O2-C3
4	A	2051	GOL	O1-C1-C2-C3
5	A	2054	PEG	O1-C1-C2-O2
3	A	2009	EDO	O1-C1-C2-O2
3	A	2020	EDO	O1-C1-C2-O2
4	A	2046	GOL	C1-C2-C3-O3
4	A	2049	GOL	O1-C1-C2-C3
4	A	2052	GOL	O1-C1-C2-C3
5	A	2057	PEG	C4-C3-O2-C2
5	A	2060	PEG	O2-C3-C4-O4
5	A	2055	PEG	O2-C3-C4-O4

There are no ring outliers.

30 monomers are involved in 68 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	2056	PEG	5	0
5	A	2063	PEG	1	0
3	A	2039	EDO	1	0
3	A	2006	EDO	1	0
3	A	2032	EDO	1	0
3	A	2037	EDO	2	0
5	A	2054	PEG	9	0
3	A	2029	EDO	2	0
3	A	2036	EDO	1	0
3	A	2028	EDO	3	1
4	A	2049	GOL	1	0

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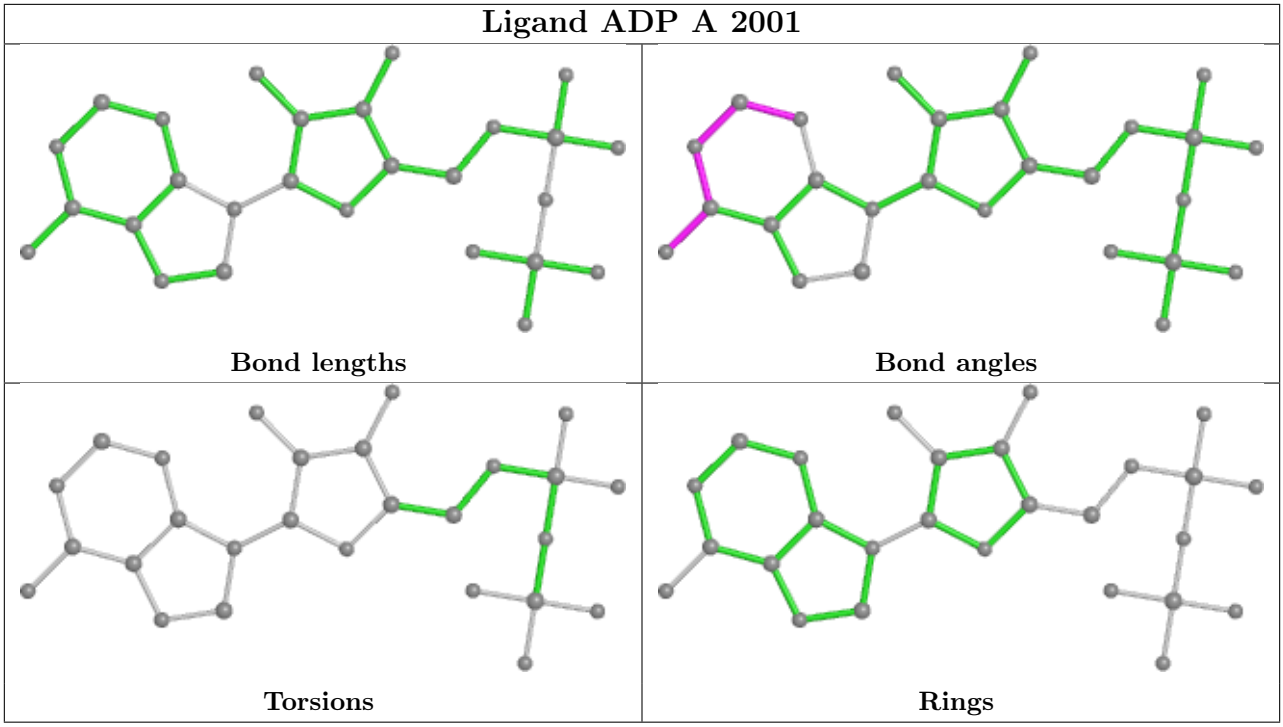
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	2048	GOL	2	0
4	A	2052	GOL	2	0
3	A	2035	EDO	1	0
3	A	2015	EDO	1	0
4	A	2053	GOL	2	0
5	A	2057	PEG	4	0
5	A	2062	PEG	1	0
3	A	2008	EDO	2	0
3	A	2002	EDO	1	0
3	A	2038	EDO	1	0
4	A	2050	GOL	1	0
5	A	2059	PEG	10	0
5	A	2061	PEG	1	0
3	A	2018	EDO	1	0
5	A	2064	PEG	3	0
3	A	2014	EDO	3	0
3	A	2023	EDO	1	0
4	A	2047	GOL	6	0
4	A	2046	GOL	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 then 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	682/749 (91%)	0.20	50 (7%) 15 14	21, 36, 84, 121	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1448	VAL	8.4
1	A	1450	VAL	7.5
1	A	1871	ILE	6.8
1	A	1866	ASN	6.1
1	A	1870	SER	6.0
1	A	1873	ARG	5.5
1	A	1544	TYR	5.4
1	A	1704	SER	5.2
1	A	1872	LEU	5.1
1	A	1396	ASN	5.1
1	A	1875	PHE	5.0
1	A	1381	SER	5.0
1	A	1380	ALA	4.7
1	A	1470	TYR	4.6
1	A	1447	VAL	4.6
1	A	1256	GLY	4.4
1	A	1397	THR	4.1
1	A	1868	ALA	4.0
1	A	1255	GLY	3.9
1	A	1379	ASN	3.6
1	A	1205	VAL	3.6
1	A	1175	ARG	3.2
1	A	1714	MET	3.1
1	A	1253	THR	3.0
1	A	1099	LYS	2.9
1	A	1097	LEU	2.9
1	A	1869	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	1425	SER	2.8
1	A	1863	PRO	2.7
1	A	1228	ASP	2.7
1	A	1254	LYS	2.6
1	A	1444	ARG	2.5
1	A	1716	TYR	2.5
1	A	1807	PHE	2.5
1	A	1543	ASN	2.4
1	A	1428	TYR	2.4
1	A	1257	ASN	2.3
1	A	1703	ILE	2.3
1	A	1867	ARG	2.3
1	A	1427	VAL	2.2
1	A	1446	ASP	2.2
1	A	1433	HIS	2.2
1	A	1098	ALA	2.2
1	A	1331	LYS	2.2
1	A	1865	ASN	2.1
1	A	1715	SER	2.1
1	A	1324	ALA	2.1
1	A	1874	LYS	2.0
1	A	1814[A]	MET	2.0
1	A	1204	GLN	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 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.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	EDO	A	2041	4/4	0.30	0.37	177,213,221,223	0
4	GOL	A	2046	6/6	0.32	0.52	135,154,177,185	0
4	GOL	A	2047	6/6	0.34	0.63	100,124,143,149	0
5	PEG	A	2064	7/7	0.34	0.72	187,232,279,289	0
5	PEG	A	2060	7/7	0.42	0.28	136,163,188,188	0
4	GOL	A	2050	6/6	0.44	0.29	134,160,171,173	0
5	PEG	A	2063	7/7	0.49	0.60	161,193,229,229	0
4	GOL	A	2052	6/6	0.49	0.50	190,228,235,242	0
5	PEG	A	2058	7/7	0.51	0.33	99,139,195,195	0
4	GOL	A	2044	6/6	0.52	0.62	136,153,180,184	0
5	PEG	A	2056	7/7	0.52	0.51	107,142,175,175	0
4	GOL	A	2048	6/6	0.55	0.47	153,184,186,186	0
3	EDO	A	2039	4/4	0.55	0.48	143,172,184,189	0
5	PEG	A	2062	7/7	0.59	0.59	131,158,183,184	0
3	EDO	A	2030	4/4	0.61	0.37	70,98,116,126	0
4	GOL	A	2049	6/6	0.61	0.30	139,167,172,174	0
5	PEG	A	2057	7/7	0.63	0.34	76,101,122,128	0
3	EDO	A	2036	4/4	0.63	0.44	116,139,142,143	0
3	EDO	A	2018	4/4	0.64	0.28	65,79,81,95	0
3	EDO	A	2022	4/4	0.65	0.68	116,139,142,142	0
3	EDO	A	2033	4/4	0.67	0.37	81,103,111,123	0
3	EDO	A	2021	4/4	0.67	0.40	109,127,140,153	0
4	GOL	A	2053	6/6	0.69	0.34	77,109,147,177	0
5	PEG	A	2054	7/7	0.70	0.26	73,93,121,121	0
3	EDO	A	2028	4/4	0.72	0.55	96,121,135,145	0
3	EDO	A	2008	4/4	0.72	0.24	81,97,105,105	0
3	EDO	A	2032	4/4	0.73	0.53	139,167,169,170	0
3	EDO	A	2004	4/4	0.74	0.19	80,96,101,101	0
5	PEG	A	2055	7/7	0.75	0.21	48,84,114,114	0
3	EDO	A	2024	4/4	0.75	0.27	77,100,111,124	0
3	EDO	A	2027	4/4	0.76	0.28	87,104,110,118	0
3	EDO	A	2031	4/4	0.76	0.21	88,106,120,120	0
3	EDO	A	2023	4/4	0.76	0.21	74,97,115,117	0
3	EDO	A	2029	4/4	0.78	0.15	69,83,85,94	0
3	EDO	A	2038	4/4	0.79	0.58	128,153,161,161	0
3	EDO	A	2014	4/4	0.79	0.24	50,65,82,82	0
3	EDO	A	2013	4/4	0.80	0.18	63,76,84,99	0
4	GOL	A	2045	6/6	0.80	0.26	44,74,97,97	0
3	EDO	A	2009	4/4	0.80	0.16	63,76,87,93	0
3	EDO	A	2034	4/4	0.81	0.25	80,96,110,112	0

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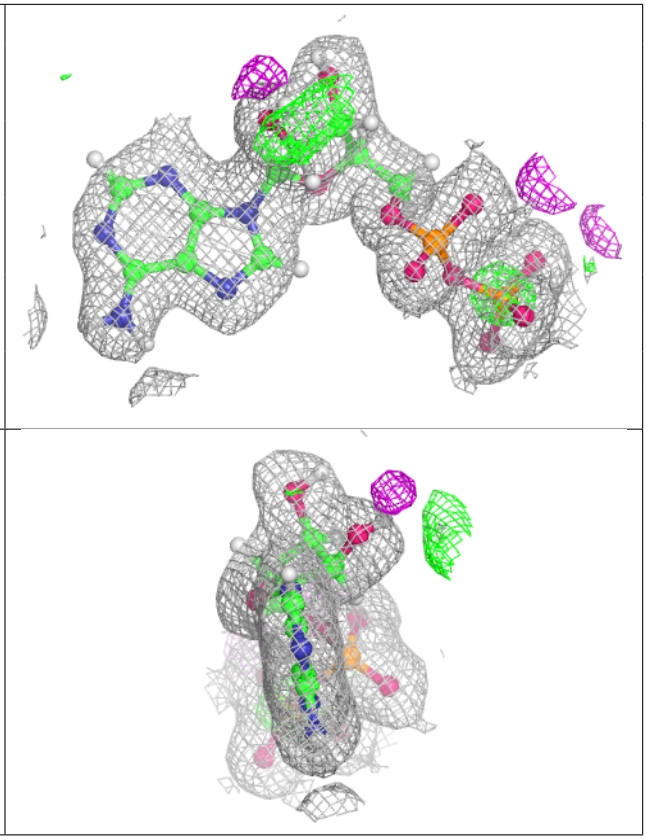
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	EDO	A	2040	4/4	0.83	0.89	100,139,164,180	0
3	EDO	A	2017	4/4	0.83	0.30	73,87,92,92	0
5	PEG	A	2059	7/7	0.84	0.42	161,193,225,228	0
5	PEG	A	2061	7/7	0.85	0.58	99,125,138,144	0
3	EDO	A	2035	4/4	0.85	0.32	88,105,116,122	0
3	EDO	A	2015	4/4	0.86	0.20	60,72,79,79	0
4	GOL	A	2051	6/6	0.86	0.33	93,118,154,173	0
4	GOL	A	2043	6/6	0.86	0.18	53,83,100,104	0
3	EDO	A	2011	4/4	0.88	0.22	44,68,80,86	0
3	EDO	A	2016	4/4	0.88	0.29	48,88,100,108	0
3	EDO	A	2010	4/4	0.89	0.65	74,88,96,96	0
3	EDO	A	2005	4/4	0.89	0.30	62,75,89,89	0
4	GOL	A	2042	6/6	0.89	0.17	56,72,90,90	0
3	EDO	A	2007	4/4	0.89	0.10	48,71,76,91	0
3	EDO	A	2025	4/4	0.89	0.31	81,97,100,103	0
3	EDO	A	2026	4/4	0.91	0.32	50,72,89,101	0
3	EDO	A	2019	4/4	0.92	0.20	51,73,84,93	0
3	EDO	A	2020	4/4	0.92	0.11	61,79,85,97	0
3	EDO	A	2006	4/4	0.92	0.16	46,62,75,75	0
3	EDO	A	2037	4/4	0.93	0.24	32,56,71,86	0
3	EDO	A	2003	4/4	0.93	0.11	45,54,62,62	0
3	EDO	A	2002	4/4	0.94	0.20	36,82,108,113	0
3	EDO	A	2012	4/4	0.97	0.11	36,47,57,59	0
2	ADP	A	2001	27/27	0.99	0.08	22,28,38,42	0
6	MG	A	2065	1/1	0.99	0.09	23,23,23,23	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 ADP A 2001:

$2mF_o$-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.