



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 1, 2025 – 10:05 PM EDT

PDB ID : 9Q7W / pdb_00009q7w
Title : One of a series of engineered variants of I-OnuI meganuclease targeting altered DNA target sequence
Authors : Werther, R.; Stoddard, B.L.
Deposited on : 2025-08-25
Resolution : 2.54 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 Phenix2.0rc1
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.45.1

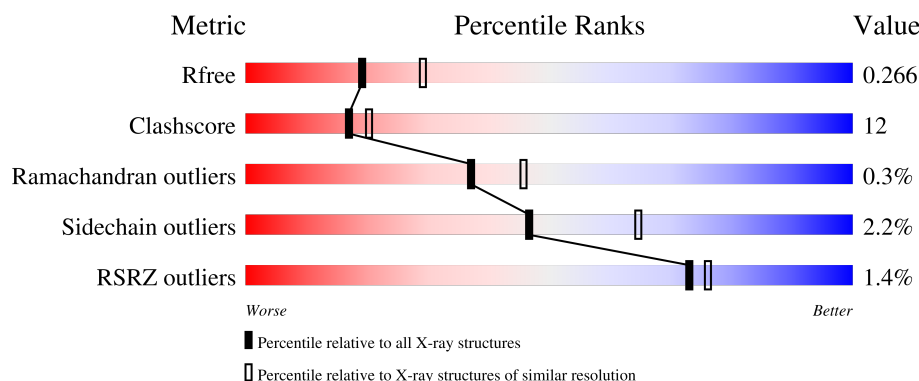
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.54 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	1004 (2.54-2.54)
Clashscore	180529	1055 (2.54-2.54)
Ramachandran outliers	177936	1048 (2.54-2.54)
Sidechain outliers	177891	1048 (2.54-2.54)
RSRZ outliers	164620	1004 (2.54-2.54)

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	300	 <div>2%</div> <div>71%</div> <div>25%</div> <div>..</div>
2	B	26	 <div>46%</div> <div>54%</div>
3	C	26	 <div>50%</div> <div>50%</div>

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 3364 atoms, of which 28 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 Ribosomal protein 3/homing endonuclease-like protein fusion.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	293	Total	C	N	O	S	0	0	0
			2239	1442	375	417	5			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	MET	-	initiating methionine	UNP Q4VWW5
A	3	ALA	-	expression tag	UNP Q4VWW5
A	4	SER	-	expression tag	UNP Q4VWW5
A	108	LYS	MET	conflict	UNP Q4VWW5
A	115	SER	CYS	conflict	UNP Q4VWW5
A	125	GLU	ILE	conflict	UNP Q4VWW5
A	153	ASN	ILE	conflict	UNP Q4VWW5
A	180	SER	CYS	conflict	UNP Q4VWW5
A	182	ILE	PHE	conflict	UNP Q4VWW5
A	199	ARG	VAL	conflict	UNP Q4VWW5
A	201	THR	SER	conflict	UNP Q4VWW5
A	225	VAL	LYS	conflict	UNP Q4VWW5
A	227	SER	LYS	conflict	UNP Q4VWW5
A	236	THR	ASP	conflict	UNP Q4VWW5

- Molecule 2 is a DNA chain called DNA (26-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	26	Total	C	N	O	P	0	0	0
			528	256	86	161	25			

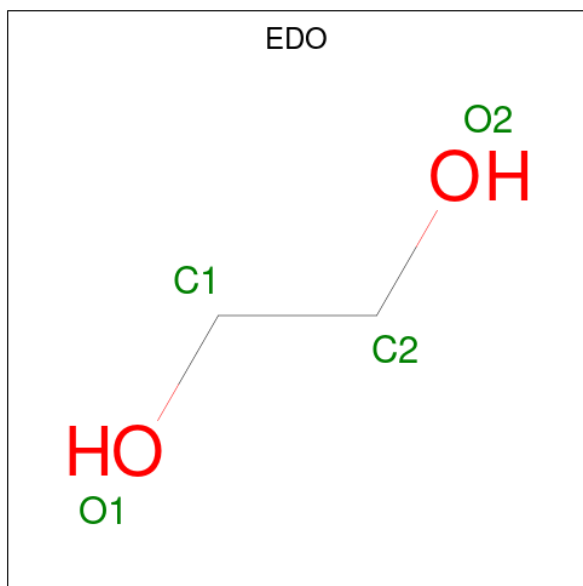
- Molecule 3 is a DNA chain called DNA (26-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	26	Total	C	N	O	P	0	0	0
			532	255	105	147	25			

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Ca	0	0
			2	2		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

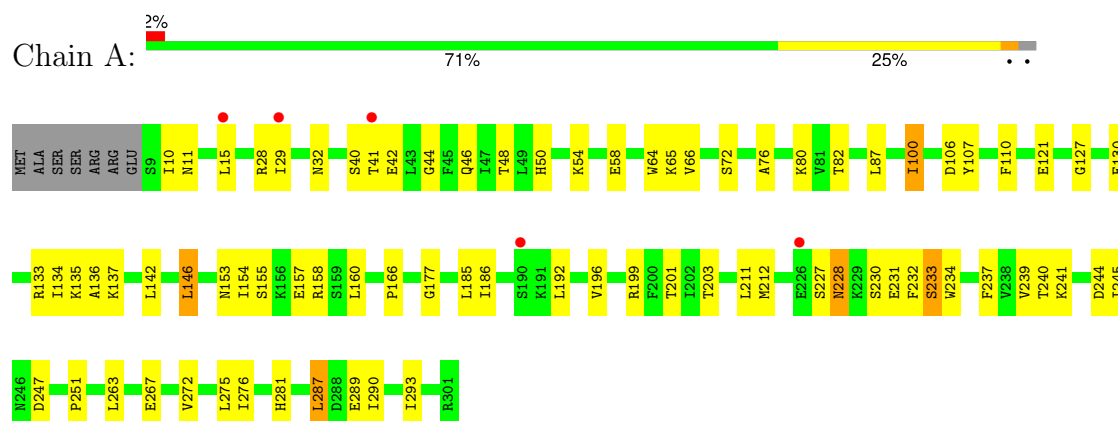
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	8	Total	O	0	0
			8	8		
7	B	4	Total	O	0	0
			4	4		
7	C	3	Total	O	0	0
			3	3		

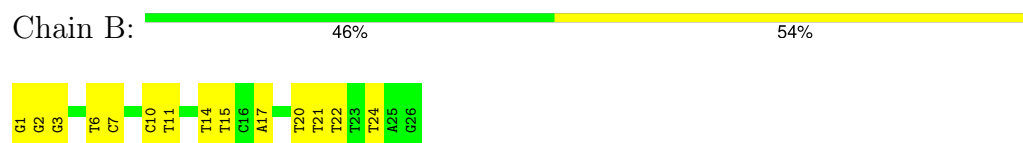
3 Residue-property plots [i](#)

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: Ribosomal protein 3/homing endonuclease-like protein fusion



- Molecule 2: DNA (26-MER)



- Molecule 3: DNA (26-MER)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	39.22Å 70.49Å 166.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.66 – 2.54 43.66 – 2.54	Depositor EDS
% Data completeness (in resolution range)	98.0 (43.66-2.54) 98.0 (43.66-2.54)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.17 (at 2.54Å)	Xtriage
Refinement program	PHENIX 1.11.1 _2575	Depositor
R, R_{free}	0.213 , 0.266 0.214 , 0.266	Depositor DCC
R_{free} test set	1579 reflections (9.83%)	wwPDB-VP
Wilson B-factor (Å ²)	55.6	Xtriage
Anisotropy	0.716	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 35.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3364	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.56% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, CA, GOL

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	0.06	0/2282	0.22	0/3097
2	B	0.20	0/589	0.40	0/908
3	C	0.19	0/599	0.37	0/922
All	All	0.12	0/3470	0.29	0/4927

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

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	2239	0	2150	62	0
2	B	528	0	300	11	0
3	C	532	0	293	11	0
4	A	2	0	0	0	0
5	A	4	6	6	0	0
5	B	4	6	6	1	0
6	B	12	16	16	0	0
7	A	8	0	0	3	0
7	B	4	0	0	0	0
7	C	3	0	0	0	0
All	All	3336	28	2771	76	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 12.

All (76) 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:10:ILE:HD11	1:A:15:LEU:HD12	1.51	0.92
1:A:137:LYS:HG2	1:A:157:GLU:HA	1.51	0.91
1:A:212:MET:HA	1:A:212:MET:HE2	1.52	0.91
1:A:276:ILE:HD13	1:A:281:HIS:HB3	1.59	0.84
1:A:106:ASP:OD1	1:A:158:ARG:NH2	2.20	0.74
1:A:275:LEU:HD12	1:A:290:ILE:HG22	1.70	0.72
1:A:10:ILE:CD1	1:A:15:LEU:HB2	2.23	0.69
1:A:48:THR:O	3:C:15:DT:H5''	1.96	0.66
1:A:72:SER:O	1:A:76:ALA:HB3	1.96	0.66
1:A:272:VAL:O	1:A:276:ILE:HG12	1.97	0.65
1:A:121:GLU:O	1:A:127:GLY:HA3	1.99	0.62
1:A:201:THR:HA	1:A:237:PHE:O	1.98	0.62
1:A:287:LEU:O	1:A:290:ILE:HG13	1.99	0.62
1:A:230:SER:OG	1:A:231:GLU:N	2.28	0.59
3:C:22:DA:H2''	3:C:23:DA:C8	2.39	0.58
2:B:21:DT:H2'	2:B:22:DT:H71	1.85	0.58
1:A:29:ILE:HB	1:A:146:LEU:HD21	1.86	0.56
1:A:10:ILE:HD11	1:A:15:LEU:HB2	1.87	0.56
1:A:137:LYS:HG2	1:A:157:GLU:CA	2.31	0.56
3:C:11:DT:H2''	3:C:12:DG:C8	2.40	0.56
1:A:240:THR:OG1	3:C:6:DA:OP2	2.21	0.56
1:A:199:ARG:NH2	2:B:20:DT:O4	2.38	0.55
1:A:203:THR:HG21	2:B:17:DA:C8	2.43	0.54
1:A:241:LYS:HE2	1:A:244:ASP:OD2	2.07	0.54
1:A:185:LEU:HD21	1:A:290:ILE:CD1	2.38	0.54
1:A:137:LYS:CG	1:A:157:GLU:HA	2.33	0.53
2:B:24:DT:H5''	5:B:101:EDO:O1	2.07	0.53
1:A:212:MET:HA	1:A:212:MET:CE	2.34	0.53
3:C:24:DA:H1'	3:C:25:DC:C6	2.44	0.53
1:A:133:ARG:HG2	1:A:153:ASN:OD1	2.10	0.52
7:A:508:HOH:O	3:C:16:DA:H5'	2.10	0.52
1:A:154:ILE:O	1:A:155:SER:HB2	2.10	0.51
3:C:17:DA:H2''	3:C:18:DG:O5'	2.09	0.51
2:B:2:DG:H1'	2:B:3:DG:C8	2.45	0.51
1:A:32:ASN:ND2	1:A:40:SER:OG	2.43	0.51
1:A:177:GLY:HA2	7:A:504:HOH:O	2.11	0.51
1:A:10:ILE:HD12	1:A:11:ASN:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:TRP:HB3	1:A:66:VAL:HG22	1.92	0.51
1:A:107:TYR:O	1:A:110:PHE:HB3	2.11	0.50
3:C:22:DA:H2''	3:C:23:DA:N7	2.27	0.50
1:A:186:ILE:HG23	2:B:22:DT:H73	1.94	0.49
1:A:281:HIS:NE2	3:C:4:DA:OP1	2.40	0.49
1:A:135:LYS:HG2	1:A:142:LEU:HD23	1.94	0.49
1:A:196:VAL:HG11	1:A:276:ILE:HD11	1.95	0.49
3:C:23:DA:H2''	3:C:24:DA:H5'	1.95	0.48
1:A:46:GLN:OE1	1:A:80:LYS:NZ	2.43	0.47
1:A:227:SER:HB2	1:A:234:TRP:CE2	2.49	0.47
1:A:192:LEU:HD13	1:A:287:LEU:HD13	1.97	0.47
1:A:136:ALA:HA	1:A:142:LEU:HG	1.97	0.46
1:A:42:GLU:OE2	1:A:82:THR:HG21	2.14	0.46
1:A:64:TRP:O	1:A:65:LYS:HB2	2.16	0.46
2:B:1:DG:H2'	2:B:2:DG:C8	2.50	0.46
1:A:28:ARG:O	1:A:41:THR:HA	2.16	0.46
1:A:228:ASN:HA	1:A:232:PHE:O	2.16	0.46
1:A:100:ILE:HD11	1:A:211:LEU:CD2	2.46	0.45
2:B:10:DC:C6	2:B:11:DT:H72	2.52	0.45
1:A:44:GLY:HA2	1:A:87:LEU:HD21	1.99	0.45
1:A:247:ASP:O	1:A:251:PRO:HG2	2.17	0.45
1:A:228:ASN:HB3	1:A:233:SER:HB3	1.99	0.44
1:A:290:ILE:C	1:A:290:ILE:HD12	2.43	0.44
1:A:137:LYS:HD2	1:A:155:SER:HB3	1.98	0.44
1:A:130:GLU:O	1:A:134:ILE:HG13	2.18	0.44
1:A:136:ALA:HB3	7:A:506:HOH:O	2.17	0.44
1:A:10:ILE:HD11	1:A:15:LEU:CD1	2.34	0.44
1:A:54:LYS:HE2	1:A:58:GLU:OE2	2.18	0.44
1:A:199:ARG:NH2	2:B:20:DT:H71	2.34	0.43
1:A:10:ILE:HD12	1:A:10:ILE:C	2.42	0.43
1:A:263:LEU:O	1:A:267:GLU:HG3	2.19	0.43
1:A:160:LEU:C	1:A:160:LEU:HD23	2.45	0.42
1:A:239:VAL:HG12	1:A:245:ILE:HG13	2.01	0.42
1:A:289:GLU:O	1:A:293:ILE:HG13	2.20	0.42
2:B:6:DT:H2''	2:B:7:DC:O5'	2.20	0.41
1:A:50:HIS:HB2	3:C:14:DA:H3'	2.03	0.41
1:A:130:GLU:HG3	1:A:133:ARG:NH1	2.36	0.41
1:A:146:LEU:HD12	1:A:146:LEU:HA	1.79	0.40
2:B:14:DT:H4'	2:B:15:DT:OP1	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	291/300 (97%)	276 (95%)	14 (5%)	1 (0%)	37	46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	166	PRO

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	231/272 (85%)	226 (98%)	5 (2%)	47	65

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	100	ILE
1	A	146	LEU
1	A	228	ASN
1	A	233	SER
1	A	287	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	254	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 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	$\# Z > 2$
5	EDO	A	403	-	3,3,3	0.43	0	2,2,2	0.37	0
5	EDO	B	101	-	3,3,3	0.43	0	2,2,2	0.36	0
6	GOL	B	102	-	5,5,5	0.37	0	5,5,5	0.29	0
6	GOL	B	103	-	5,5,5	0.39	0	5,5,5	0.20	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
5	EDO	A	403	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	B	101	-	-	0/1/1/1	-
6	GOL	B	102	-	-	2/4/4/4	-
6	GOL	B	103	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	103	GOL	O1-C1-C2-C3
6	B	102	GOL	C1-C2-C3-O3
6	B	103	GOL	O1-C1-C2-O2
6	B	102	GOL	O2-C2-C3-O3
5	A	403	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	101	EDO	1	0

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	293/300 (97%)	0.45	5 (1%) 69 72	41, 55, 75, 80	0
2	B	26/26 (100%)	-0.26	0 100 100	46, 64, 73, 79	0
3	C	26/26 (100%)	-0.21	0 100 100	52, 63, 78, 81	0
All	All	345/352 (98%)	0.34	5 (1%) 73 76	41, 56, 75, 81	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	41	THR	2.8
1	A	29	ILE	2.8
1	A	190	SER	2.3
1	A	226	GLU	2.2
1	A	15	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
6	GOL	B	103	6/6	0.65	0.13	71,85,94,94	0
5	EDO	B	101	4/4	0.68	0.17	63,75,86,91	0
6	GOL	B	102	6/6	0.75	0.15	64,77,85,93	0
5	EDO	A	403	4/4	0.82	0.17	56,68,82,82	0
4	CA	A	402	1/1	0.96	0.08	54,54,54,54	0
4	CA	A	401	1/1	0.98	0.05	46,46,46,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.