



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:01 PM UTC

PDB ID : 6GRY / pdb_00006gry
Title : Glucuronoyl Esterase from Solibacter usitatus.
Authors : Lo Leggio, L.; Larsbrink, J.; Meland Knudsen, R.; Mazurkewich, S.; Navarro Poulsen, J.C.
Deposited on : 2018-06-12
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
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

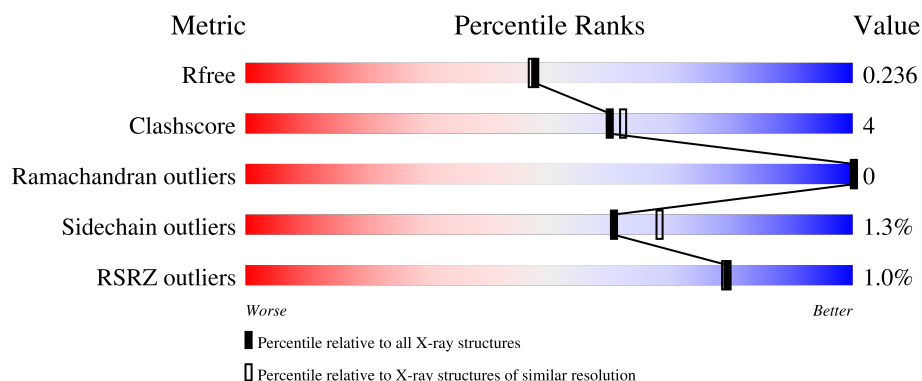
1 Overall quality at a glance

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	397	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6651 atoms, of which 3202 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.

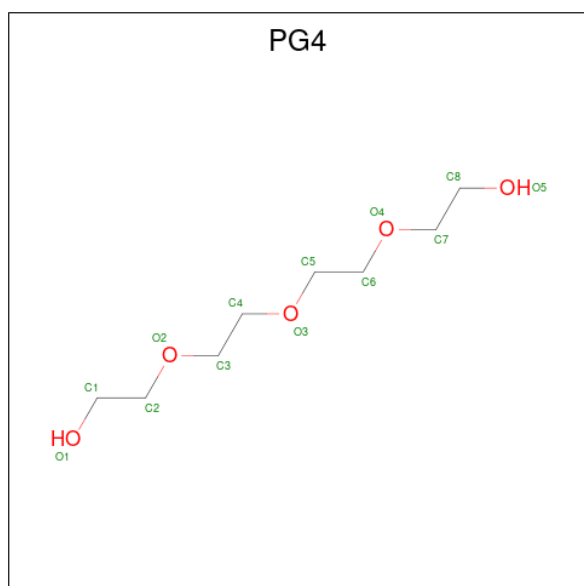
- Molecule 1 is a protein called Putative acetyl xylan esterase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	392	6111	1977	3014	541	570	9	0	6	0

There are 2 discrepancies between the modelled and reference sequences:

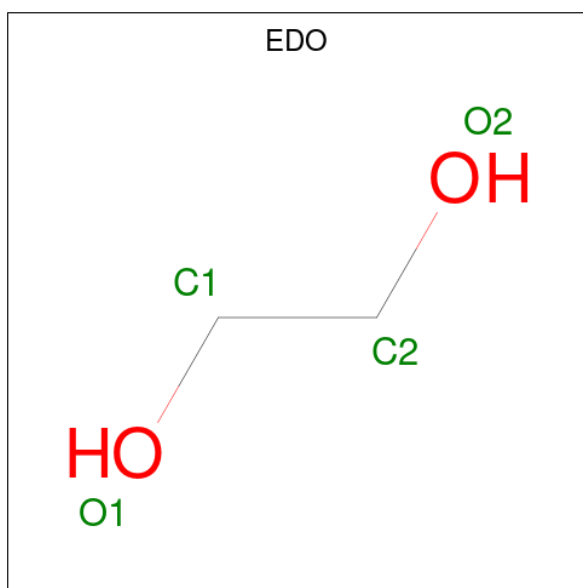
Chain	Residue	Modelled	Actual	Comment	Reference
A	21	GLY	-	expression tag	UNP Q01YM8
A	22	HIS	-	expression tag	UNP Q01YM8

- Molecule 2 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



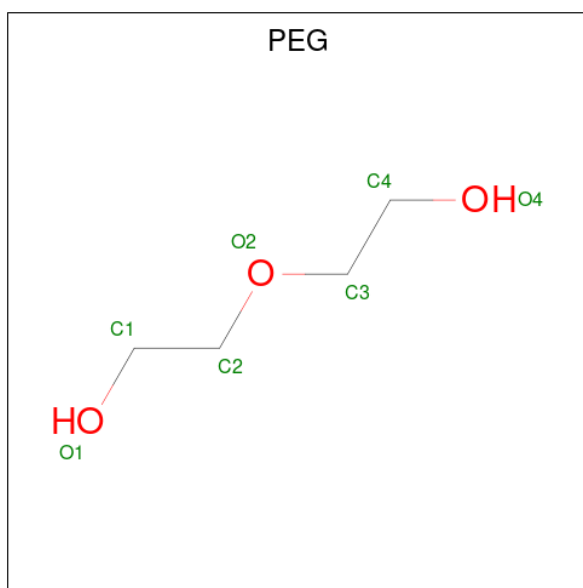
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
2	A	1	31	8	18	5	0	0
2	A	1	31	8	18	5	0	0

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



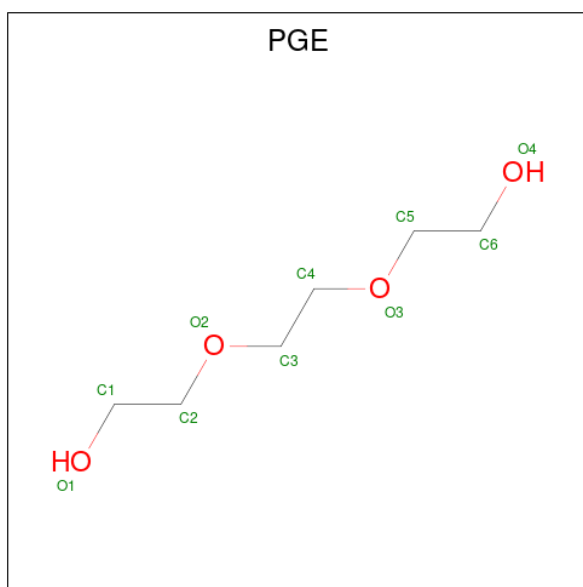
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			10	2	6	2		
3	A	1	Total	C	H	O	0	0
			10	2	6	2		
3	A	1	Total	C	H	O	0	0
			10	2	6	2		
3	A	1	Total	C	H	O	0	0
			10	2	6	2		
3	A	1	Total	C	H	O	0	0
			10	2	6	2		
3	A	1	Total	C	H	O	0	0
			10	2	6	2		
3	A	1	Total	C	H	O	0	0
			10	2	6	2		
3	A	1	Total	C	H	O	0	0
			10	2	6	2		
3	A	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		
4	A	1	Total	C	H	O	0	0
			17	4	10	3		
4	A	1	Total	C	H	O	0	0
			17	4	10	3		
4	A	1	Total	C	H	O	0	0
			17	4	10	3		
4	A	1	Total	C	H	O	0	0
			17	4	10	3		

- Molecule 5 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O	0	0
			24	6	14	4		

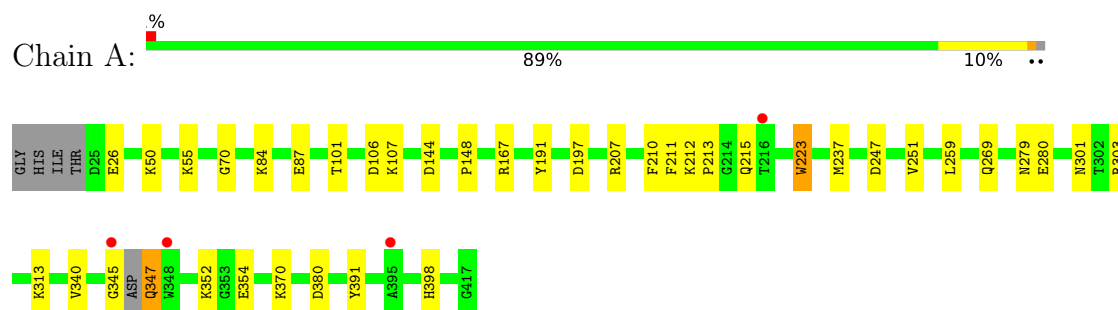
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	222	Total	O	0	0
			222	222		

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: Putative acetyl xylan esterase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	54.58Å 54.58Å 284.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.29 – 2.00 47.29 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.4 (47.29-2.00) 98.4 (47.29-2.00)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.188 , 0.239 0.189 , 0.236	Depositor DCC
R_{free} test set	967 reflections (3.18%)	wwPDB-VP
Wilson B-factor (Å ²)	32.6	Xtriage
Anisotropy	0.084	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 47.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6651	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.39% 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, PGE, PG4, 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	0.56	0/3196	0.66	0/4329

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	3097	3014	3021	28	0
2	A	26	36	36	1	0
3	A	52	78	78	0	2
4	A	42	60	60	3	0
5	A	10	14	14	1	0
6	A	222	0	0	5	0
All	All	3449	3202	3209	28	2

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 4.

All (28) 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:345:GLY:O	1:A:347:GLN:HG2	1.85	0.77
1:A:354:GLU:OE2	6:A:601:HOH:O	2.07	0.71
1:A:269:GLN:OE1	6:A:602:HOH:O	2.15	0.65
1:A:345:GLY:C	1:A:347:GLN:HG2	2.23	0.63
1:A:87:GLU:OE2	6:A:603:HOH:O	2.17	0.60
1:A:280:GLU:OE1	1:A:398:HIS:ND1	2.37	0.57
1:A:279:ASN:OD1	1:A:280:GLU:HG3	2.04	0.56
1:A:26:GLU:OE1	1:A:352:LYS:HE2	2.07	0.55
1:A:380[A]:ASP:OD2	6:A:604:HOH:O	2.18	0.54
1:A:280:GLU:OE1	1:A:398:HIS:HA	2.11	0.51
1:A:106:ASP:OD1	1:A:107:LYS:N	2.43	0.51
1:A:191:TYR:CD2	1:A:259:LEU:HD21	2.45	0.50
1:A:144:ASP:O	1:A:167:ARG:NH1	2.45	0.49
1:A:340:VAL:O	1:A:391:TYR:HA	2.12	0.49
1:A:55:LYS:NZ	6:A:617:HOH:O	2.48	0.46
1:A:84:LYS:HE2	1:A:101:THR:OG1	2.16	0.46
1:A:211:PHE:HB3	1:A:215:GLN:HB3	1.97	0.45
1:A:370:LYS:HZ2	4:A:519:PEG:H42	1.81	0.45
1:A:237[A]:MET:HE1	1:A:251:VAL:HB	1.98	0.44
1:A:210:PHE:HB3	2:A:502:PG4:H52	2.00	0.44
1:A:148:PRO:HG2	4:A:520:PEG:C3	2.49	0.43
1:A:197:ASP:O	5:A:522:PGE:H1	2.20	0.42
1:A:148:PRO:HG2	4:A:520:PEG:H32	2.02	0.42
1:A:301:ASN:HB3	1:A:313:LYS:HD2	2.00	0.42
1:A:70:GLY:HA3	1:A:223:TRP:HA	2.02	0.42
1:A:207:ARG:HG2	1:A:223:TRP:CE2	2.55	0.41
1:A:247:ASP:OD1	1:A:247:ASP:C	2.63	0.41
1:A:212:LYS:CG	1:A:213:PRO:HD2	2.50	0.41

All (2) 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)
3:A:504:EDO:O1	3:A:504:EDO:O1[7_556]	1.76	0.44
3:A:504:EDO:O1	3:A:504:EDO:HO1[7_556]	1.51	0.09

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	394/397 (99%)	376 (95%)	18 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	309/308 (100%)	305 (99%)	4 (1%)	61	68

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	LYS
1	A	223	TRP
1	A	303	ARG
1	A	347	GLN

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	80	ASN
1	A	143	ASN
1	A	407	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

22 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	$\# Z > 2$
3	EDO	A	507	-	3,3,3	0.45	0	2,2,2	0.38	0
3	EDO	A	508	-	3,3,3	0.63	0	2,2,2	0.18	0
4	PEG	A	518	-	6,6,6	0.50	0	5,5,5	0.42	0
4	PEG	A	519	-	6,6,6	0.59	0	5,5,5	0.83	0
2	PG4	A	501	-	12,12,12	0.66	0	11,11,11	0.89	0
3	EDO	A	506	-	3,3,3	0.50	0	2,2,2	0.24	0
3	EDO	A	513	-	3,3,3	0.48	0	2,2,2	0.29	0
3	EDO	A	505	-	3,3,3	0.42	0	2,2,2	0.21	0
4	PEG	A	521	-	6,6,6	0.55	0	5,5,5	0.46	0
3	EDO	A	510	-	3,3,3	0.60	0	2,2,2	0.19	0
4	PEG	A	517	-	6,6,6	0.65	0	5,5,5	1.27	0
3	EDO	A	512	-	3,3,3	0.64	0	2,2,2	0.19	0
2	PG4	A	502	-	12,12,12	0.56	0	11,11,11	0.51	0
3	EDO	A	511	-	3,3,3	0.63	0	2,2,2	0.12	0
4	PEG	A	520	-	6,6,6	0.44	0	5,5,5	0.42	0
3	EDO	A	515	-	3,3,3	0.44	0	2,2,2	0.29	0
5	PGE	A	522	-	9,9,9	0.39	0	8,8,8	0.32	0
3	EDO	A	509	-	3,3,3	0.45	0	2,2,2	0.34	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	PEG	A	516	-	6,6,6	0.48	0	5,5,5	0.32	0
3	EDO	A	504	-	3,3,3	0.38	0	2,2,2	0.45	0
3	EDO	A	514	-	3,3,3	0.36	0	2,2,2	0.28	0
3	EDO	A	503	-	3,3,3	0.34	0	2,2,2	0.45	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	507	-	-	0/1/1/1	-
3	EDO	A	508	-	-	1/1/1/1	-
4	PEG	A	518	-	-	2/4/4/4	-
4	PEG	A	519	-	-	2/4/4/4	-
2	PG4	A	501	-	-	7/10/10/10	-
3	EDO	A	506	-	-	0/1/1/1	-
3	EDO	A	513	-	-	0/1/1/1	-
3	EDO	A	505	-	-	1/1/1/1	-
4	PEG	A	521	-	-	1/4/4/4	-
3	EDO	A	510	-	-	0/1/1/1	-
4	PEG	A	517	-	-	4/4/4/4	-
3	EDO	A	512	-	-	1/1/1/1	-
2	PG4	A	502	-	-	5/10/10/10	-
3	EDO	A	511	-	-	0/1/1/1	-
4	PEG	A	520	-	-	3/4/4/4	-
3	EDO	A	515	-	-	1/1/1/1	-
5	PGE	A	522	-	-	6/7/7/7	-
3	EDO	A	509	-	-	1/1/1/1	-
4	PEG	A	516	-	-	3/4/4/4	-
3	EDO	A	504	-	-	1/1/1/1	-
3	EDO	A	514	-	-	0/1/1/1	-
3	EDO	A	503	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (40) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	502	PG4	O3-C5-C6-O4
4	A	516	PEG	O2-C3-C4-O4
5	A	522	PGE	O3-C5-C6-O4
5	A	522	PGE	O2-C3-C4-O3
2	A	501	PG4	O4-C7-C8-O5
4	A	517	PEG	O2-C3-C4-O4
4	A	518	PEG	O1-C1-C2-O2
4	A	520	PEG	O2-C3-C4-O4
3	A	503	EDO	O1-C1-C2-O2
3	A	508	EDO	O1-C1-C2-O2
3	A	512	EDO	O1-C1-C2-O2
2	A	502	PG4	C8-C7-O4-C6
2	A	501	PG4	O2-C3-C4-O3
4	A	516	PEG	O1-C1-C2-O2
3	A	509	EDO	O1-C1-C2-O2
5	A	522	PGE	C1-C2-O2-C3
4	A	517	PEG	C1-C2-O2-C3
4	A	520	PEG	C1-C2-O2-C3
4	A	519	PEG	O2-C3-C4-O4
5	A	522	PGE	O1-C1-C2-O2
2	A	502	PG4	C4-C3-O2-C2
2	A	501	PG4	C6-C5-O3-C4
4	A	518	PEG	C4-C3-O2-C2
2	A	501	PG4	C1-C2-O2-C3
4	A	517	PEG	O1-C1-C2-O2
2	A	501	PG4	C4-C3-O2-C2
4	A	516	PEG	C1-C2-O2-C3
2	A	501	PG4	O3-C5-C6-O4
3	A	505	EDO	O1-C1-C2-O2
4	A	517	PEG	C4-C3-O2-C2
5	A	522	PGE	C3-C4-O3-C5
3	A	504	EDO	O1-C1-C2-O2
4	A	520	PEG	O1-C1-C2-O2
2	A	502	PG4	O2-C3-C4-O3
5	A	522	PGE	C4-C3-O2-C2
2	A	501	PG4	C3-C4-O3-C5
3	A	515	EDO	O1-C1-C2-O2
4	A	521	PEG	C4-C3-O2-C2
4	A	519	PEG	O1-C1-C2-O2
2	A	502	PG4	C6-C5-O3-C4

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	519	PEG	1	0
2	A	502	PG4	1	0
4	A	520	PEG	2	0
5	A	522	PGE	1	0
3	A	504	EDO	0	2

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	392/397 (98%)	-0.19	4 (1%) 79 79	17, 38, 72, 108	4 (1%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	345	GLY	4.4
1	A	348	TRP	3.2
1	A	395	ALA	3.1
1	A	216	THR	2.3

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(Å ²)	Q<0.9
3	EDO	A	508	4/4	0.68	0.16	54,68,81,96	0
2	PG4	A	502	13/13	0.74	0.15	55,84,110,113	0
4	PEG	A	518	7/7	0.74	0.16	56,81,97,100	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EDO	A	515	4/4	0.75	0.13	67,80,89,89	0
4	PEG	A	521	7/7	0.78	0.16	60,81,101,101	0
4	PEG	A	519	7/7	0.79	0.14	37,67,80,80	0
3	EDO	A	506	4/4	0.79	0.12	53,64,71,72	0
5	PGE	A	522	10/10	0.80	0.14	39,67,97,98	0
2	PG4	A	501	13/13	0.81	0.14	39,61,72,75	0
3	EDO	A	507	4/4	0.82	0.14	51,62,67,76	0
3	EDO	A	513	4/4	0.83	0.12	77,92,99,110	0
3	EDO	A	510	4/4	0.85	0.20	32,64,77,78	0
4	PEG	A	520	7/7	0.85	0.13	58,70,97,97	0
3	EDO	A	511	4/4	0.86	0.12	57,72,90,90	0
4	PEG	A	516	7/7	0.88	0.12	45,58,74,74	0
3	EDO	A	509	4/4	0.88	0.12	43,51,57,63	0
4	PEG	A	517	7/7	0.89	0.13	29,53,73,73	0
3	EDO	A	512	4/4	0.89	0.14	47,60,76,76	0
3	EDO	A	505	4/4	0.90	0.18	49,63,79,79	0
3	EDO	A	514	4/4	0.90	0.20	40,59,71,71	0
3	EDO	A	503	4/4	0.91	0.09	46,55,60,70	0
3	EDO	A	504	4/4	0.94	0.09	39,52,64,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.