



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 14, 2020 – 03:39 PM BST

PDB ID : 6YQ6  
Title : Promiscuous Reductase LugOII Catalyzes Keto-reduction at C1 during Lugdunomycin Biosynthesis  
Authors : Xiao, X.; Elsayed, S.S.; Wu, C.; van der Heul, H.; Protá, A.; Huang, J.; Guo, R.; Abrahams, J.P.; van Wezel, G.P.  
Deposited on : 2020-04-16  
Resolution : 2.08 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) 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	:	<b>FAILED</b>
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.14.4.dev1

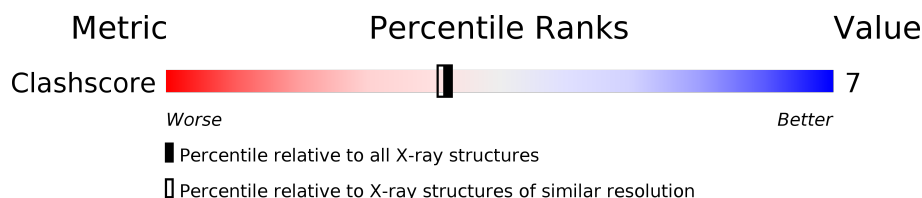
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.08 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	141614	6738 (2.10-2.06)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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 failed to run properly.

Mol	Chain	Length	Quality of chain
1	AAA	255	 82% 9% 9%
1	BBB	255	 86% 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	PEG	AAA	404	-	-	X	-
3	PEG	AAA	405	-	-	X	-
3	PEG	BBB	501	-	-	X	-

## 2 Entry composition [i](#)

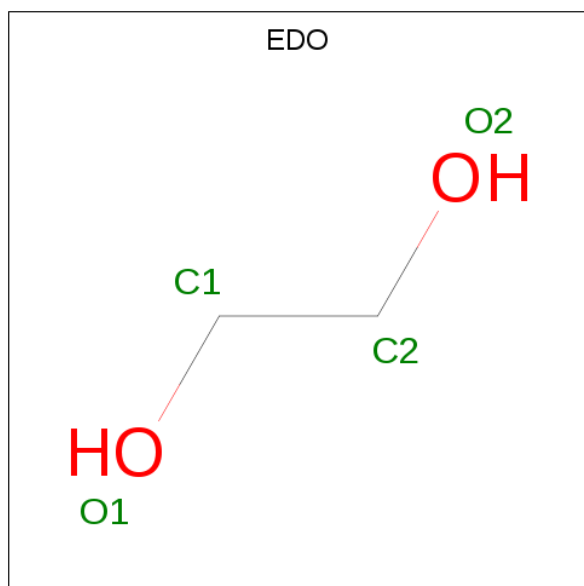
There are 5 unique types of molecules in this entry. The entry contains 4027 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	231	Total	C	N	O	S	0	4	0
			1731	1077	316	332	6			
1	BBB	244	Total	C	N	O	S	0	4	0
			1825	1138	327	353	7			

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



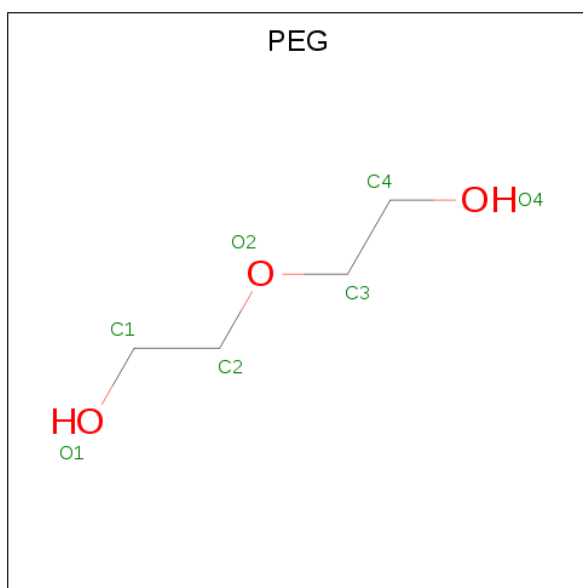
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	AAA	1	Total	C	O	0	0
			4	2	2		
2	AAA	1	Total	C	O	0	0
			4	2	2		
2	AAA	1	Total	C	O	0	0
			4	2	2		
2	AAA	1	Total	C	O	0	0
			4	2	2		

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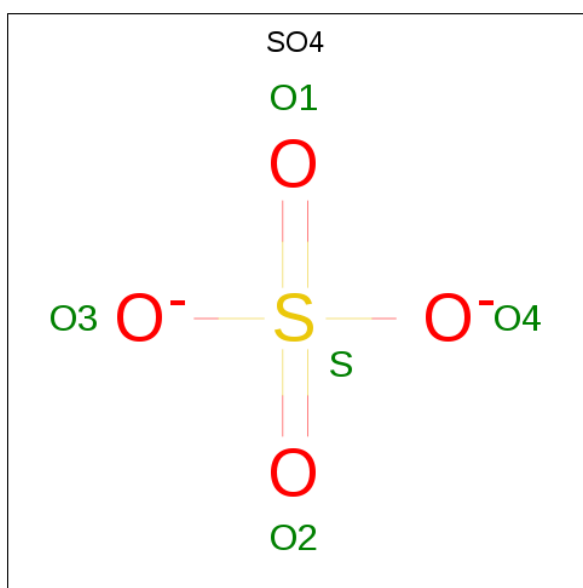
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	AAA	1	Total	C	O	0	0
			4	2	2		
2	AAA	1	Total	C	O	0	0
			4	2	2		
2	AAA	1	Total	C	O	0	0
			4	2	2		
2	AAA	1	Total	C	O	0	0
			4	2	2		
2	AAA	1	Total	C	O	0	0
			4	2	2		
2	BBB	1	Total	C	O	0	0
			4	2	2		
2	BBB	1	Total	C	O	0	0
			4	2	2		
2	BBB	1	Total	C	O	0	0
			4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

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula:  $C_4H_{10}O_3$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	AAA	1	Total C O 7 4 3	0	0
3	AAA	1	Total C O 7 4 3	0	0
3	BBB	1	Total C O 7 4 3	0	0
3	BBB	1	Total C O 7 4 3	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

- Molecule 5 is water.

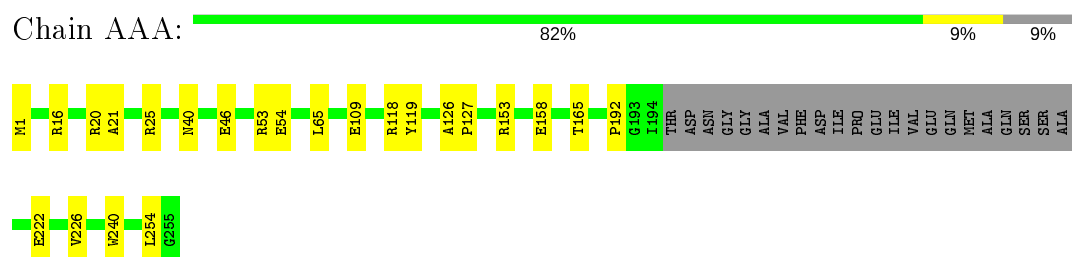
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	AAA	194	Total O 194 194	0	0
5	BBB	170	Total O 170 170	0	0

### 3 Residue-property plots [i](#)

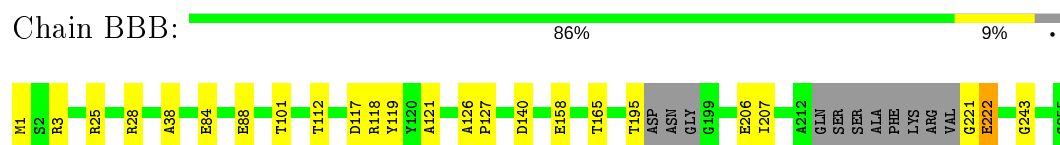
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS failed to run properly.

- Molecule 1: Monooxygenase



- Molecule 1: Monooxygenase



## 4 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	185.99 Å 185.99 Å 75.64 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.47 – 2.08	Depositor
% Data completeness (in resolution range)	99.9 (47.47-2.08)	Depositor
$R_{merge}$	0.40	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.53 (at 2.08 Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, $R_{free}$	0.168 , 0.203	Depositor
Wilson B-factor (Å <sup>2</sup> )	25.3	Xtriage
Anisotropy	0.058	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4027	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.64% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

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

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PEG, SO4, EDO

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	AAA	0.97	6/1764 (0.3%)	0.98	4/2389 (0.2%)
1	BBB	0.91	3/1859 (0.2%)	0.95	3/2518 (0.1%)
All	All	0.94	9/3623 (0.2%)	0.96	7/4907 (0.1%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AAA	158	GLU	CD-OE1	11.09	1.37	1.25
1	BBB	88	GLU	CD-OE2	7.95	1.34	1.25
1	BBB	84	GLU	CD-OE2	6.90	1.33	1.25
1	BBB	88	GLU	CD-OE1	6.48	1.32	1.25
1	AAA	222	GLU	CD-OE1	5.98	1.32	1.25
1	AAA	46	GLU	CD-OE2	-5.80	1.19	1.25
1	AAA	109	GLU	CD-OE1	-5.70	1.19	1.25
1	AAA	54	GLU	CD-OE1	5.35	1.31	1.25
1	AAA	158	GLU	CD-OE2	5.32	1.31	1.25

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BBB	118[A]	ARG	NE-CZ-NH1	-7.50	116.55	120.30
1	BBB	118[B]	ARG	NE-CZ-NH1	-7.50	116.55	120.30
1	AAA	53	ARG	NE-CZ-NH1	5.61	123.10	120.30
1	AAA	118[A]	ARG	NE-CZ-NH2	-5.53	117.53	120.30
1	AAA	118[B]	ARG	NE-CZ-NH2	-5.53	117.53	120.30
1	BBB	222	GLU	CB-CA-C	5.28	120.95	110.40
1	AAA	222	GLU	CB-CA-C	5.20	120.80	110.40

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	AAA	1731	0	1760	32	0
1	BBB	1825	0	1846	23	0
2	AAA	40	0	60	5	0
2	BBB	24	0	36	7	0
3	AAA	14	0	20	16	0
3	BBB	14	0	20	4	0
4	AAA	5	0	0	1	0
4	BBB	10	0	0	0	0
5	AAA	194	0	0	0	0
5	BBB	170	0	0	2	0
All	All	4027	0	3742	51	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (51) 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:BBB:112:THR:HA	2:BBB:507:EDO:H12	1.46	0.96
1:AAA:65:LEU:HB2	3:AAA:405:PEG:H42	1.59	0.84
1:BBB:117:ASP:HB3	2:BBB:505:EDO:H12	1.58	0.82
1:AAA:20:ARG:HG2	3:AAA:404:PEG:H32	1.64	0.77
1:BBB:25:ARG:HH12	3:BBB:501:PEG:C2	2.02	0.72
1:AAA:25:ARG:HH12	3:AAA:404:PEG:H11	1.53	0.71
1:BBB:25:ARG:HH12	3:BBB:501:PEG:H22	1.57	0.70
1:AAA:1:MET:H1	1:BBB:1:MET:HE3	1.61	0.66
1:AAA:65:LEU:H	3:AAA:405:PEG:H42	1.61	0.65
1:AAA:153:ARG:HH22	2:AAA:412:EDO:H22	1.61	0.65
1:BBB:121:ALA:HB2	2:BBB:505:EDO:H22	1.77	0.64
1:AAA:21:ALA:CA	3:AAA:404:PEG:H31	2.29	0.63
1:AAA:126:ALA:HA	2:AAA:410:EDO:H22	1.82	0.60
1:AAA:16:ARG:HH22	1:AAA:40:ASN:HD22	1.50	0.59
1:BBB:195:THR:OG1	1:BBB:207:ILE:HD11	2.05	0.56
1:AAA:1:MET:N	1:BBB:1:MET:CE	2.68	0.56
1:AAA:16:ARG:NH1	4:AAA:413:SO4:O3	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AAA:65:LEU:CB	3:AAA:405:PEG:H42	2.32	0.55
1:BBB:121:ALA:CB	2:BBB:505:EDO:H22	2.36	0.55
1:AAA:21:ALA:N	3:AAA:404:PEG:H31	2.22	0.54
2:AAA:412:EDO:H12	5:BBB:709:HOH:O	2.06	0.54
1:BBB:1:MET:HE2	1:BBB:3:ARG:HD3	1.91	0.53
1:BBB:25:ARG:HH12	3:BBB:501:PEG:H21	1.73	0.52
1:AAA:1:MET:H1	1:BBB:1:MET:CE	2.21	0.52
1:BBB:222:GLU:HG2	2:BBB:503:EDO:O2	2.10	0.52
1:BBB:38:ALA:HB2	2:BBB:502:EDO:H11	1.90	0.51
1:BBB:101:THR:HB	1:BBB:206:GLU:HG3	1.93	0.51
1:BBB:158:GLU:HG3	5:BBB:714:HOH:O	2.11	0.50
1:AAA:65:LEU:N	3:AAA:405:PEG:H42	2.27	0.50
1:BBB:119:TYR:HB3	1:BBB:165:THR:HG21	1.92	0.50
1:AAA:65:LEU:HB2	3:AAA:405:PEG:C4	2.37	0.49
1:AAA:25:ARG:HH12	3:AAA:404:PEG:C1	2.21	0.48
1:AAA:21:ALA:HA	3:AAA:404:PEG:H31	1.94	0.48
1:BBB:126:ALA:HB3	1:BBB:127:PRO:HD3	1.96	0.47
1:AAA:126:ALA:HB2	2:AAA:410:EDO:H11	1.97	0.46
1:AAA:65:LEU:H	3:AAA:405:PEG:C4	2.27	0.46
1:AAA:254:LEU:HD21	1:BBB:243:GLY:HA3	1.97	0.46
1:AAA:21:ALA:HB2	3:AAA:404:PEG:H42	1.97	0.46
1:BBB:222:GLU:HA	2:BBB:503:EDO:H12	1.98	0.45
1:AAA:16:ARG:HH22	1:AAA:40:ASN:ND2	2.14	0.45
1:AAA:126:ALA:HB3	1:AAA:127:PRO:HD3	1.98	0.45
1:AAA:126:ALA:CA	2:AAA:410:EDO:H22	2.46	0.44
1:BBB:28:ARG:HH22	3:BBB:501:PEG:H32	1.84	0.43
1:AAA:119:TYR:HB3	1:AAA:165:THR:HG21	2.00	0.42
1:AAA:25:ARG:NH1	3:AAA:404:PEG:H11	2.26	0.42
1:AAA:20:ARG:HG2	3:AAA:404:PEG:C3	2.41	0.41
1:AAA:240:TRP:HE1	1:BBB:221:GLY:HA2	1.85	0.41
1:AAA:192:PRO:HB3	1:AAA:226:VAL:HG21	2.02	0.41
1:AAA:16:ARG:NH2	1:AAA:40:ASN:HD22	2.16	0.41
1:AAA:25:ARG:HH12	3:AAA:404:PEG:C2	2.34	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

### 5.3.2 Protein sidechains [i](#)

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

### 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](#)

23 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	PEG	BBB	501	-	6,6,6	0.66	0	5,5,5	0.47	0
2	EDO	BBB	505	-	3,3,3	0.30	0	2,2,2	1.01	0
2	EDO	BBB	508	-	3,3,3	0.73	0	2,2,2	0.83	0
2	EDO	AAA	411	-	3,3,3	0.41	0	2,2,2	0.62	0
2	EDO	AAA	410	-	3,3,3	0.26	0	2,2,2	0.14	0
2	EDO	BBB	507	-	3,3,3	0.43	0	2,2,2	0.64	0
4	SO4	BBB	509	-	4,4,4	0.37	0	6,6,6	0.29	0
2	EDO	BBB	504	-	3,3,3	0.30	0	2,2,2	0.11	0
2	EDO	AAA	401	-	3,3,3	0.56	0	2,2,2	0.24	0
3	PEG	BBB	506	-	6,6,6	0.66	0	5,5,5	0.67	0
2	EDO	BBB	502	-	3,3,3	0.38	0	2,2,2	0.23	0
4	SO4	BBB	510	-	4,4,4	0.36	0	6,6,6	0.10	0
2	EDO	AAA	409	-	3,3,3	0.37	0	2,2,2	0.51	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	SO4	AAA	413	-	4,4,4	0.41	0	6,6,6	0.51	0
3	PEG	AAA	405	-	6,6,6	0.43	0	5,5,5	0.30	0
2	EDO	AAA	406	-	3,3,3	0.16	0	2,2,2	0.14	0
2	EDO	AAA	407	-	3,3,3	0.09	0	2,2,2	0.15	0
2	EDO	AAA	402	-	3,3,3	0.36	0	2,2,2	0.37	0
2	EDO	AAA	403	-	3,3,3	0.38	0	2,2,2	0.37	0
3	PEG	AAA	404	-	6,6,6	0.93	0	5,5,5	0.76	0
2	EDO	AAA	412	-	3,3,3	0.63	0	2,2,2	0.93	0
2	EDO	AAA	408	-	3,3,3	0.48	0	2,2,2	0.53	0
2	EDO	BBB	503	-	3,3,3	0.47	0	2,2,2	0.70	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	PEG	AAA	405	-	-	1/4/4/4	-
2	EDO	AAA	401	-	-	1/1/1/1	-
3	PEG	BBB	501	-	-	2/4/4/4	-
2	EDO	AAA	406	-	-	1/1/1/1	-
2	EDO	BBB	505	-	-	1/1/1/1	-
2	EDO	BBB	508	-	-	1/1/1/1	-
2	EDO	AAA	407	-	-	1/1/1/1	-
2	EDO	AAA	411	-	-	0/1/1/1	-
3	PEG	BBB	506	-	-	3/4/4/4	-
2	EDO	AAA	410	-	-	1/1/1/1	-
2	EDO	AAA	402	-	-	0/1/1/1	-
2	EDO	AAA	408	-	-	1/1/1/1	-
2	EDO	BBB	507	-	-	0/1/1/1	-
2	EDO	BBB	502	-	-	1/1/1/1	-
2	EDO	BBB	503	-	-	1/1/1/1	-
2	EDO	AAA	403	-	-	1/1/1/1	-
3	PEG	AAA	404	-	-	1/4/4/4	-
2	EDO	AAA	409	-	-	0/1/1/1	-
2	EDO	BBB	504	-	-	1/1/1/1	-
2	EDO	AAA	412	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	BBB	506	PEG	C1-C2-O2-C3
2	AAA	403	EDO	O1-C1-C2-O2
3	BBB	506	PEG	O1-C1-C2-O2
2	BBB	508	EDO	O1-C1-C2-O2
2	BBB	503	EDO	O1-C1-C2-O2
3	BBB	506	PEG	O2-C3-C4-O4
3	AAA	405	PEG	C4-C3-O2-C2
2	BBB	505	EDO	O1-C1-C2-O2
2	BBB	502	EDO	O1-C1-C2-O2
2	AAA	406	EDO	O1-C1-C2-O2
3	AAA	404	PEG	O2-C3-C4-O4
2	BBB	504	EDO	O1-C1-C2-O2
2	AAA	401	EDO	O1-C1-C2-O2
2	AAA	410	EDO	O1-C1-C2-O2
2	AAA	407	EDO	O1-C1-C2-O2
3	BBB	501	PEG	O2-C3-C4-O4
3	BBB	501	PEG	O1-C1-C2-O2
2	AAA	408	EDO	O1-C1-C2-O2

There are no ring outliers.

10 monomers are involved in 33 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	BBB	501	PEG	4	0
2	BBB	505	EDO	3	0
2	AAA	410	EDO	3	0
2	BBB	507	EDO	1	0
2	BBB	502	EDO	1	0
4	AAA	413	SO4	1	0
3	AAA	405	PEG	6	0
3	AAA	404	PEG	10	0
2	AAA	412	EDO	2	0
2	BBB	503	EDO	2	0

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

EDS failed to run properly - this section is therefore empty.

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

EDS failed to run properly - this section is therefore empty.

### 6.3 Carbohydrates [i](#)

EDS failed to run properly - this section is therefore empty.

### 6.4 Ligands [i](#)

EDS failed to run properly - this section is therefore empty.

### 6.5 Other polymers [i](#)

EDS failed to run properly - this section is therefore empty.