



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 02:16 PM GMT

PDB ID : 3WI3  
Title : Crystal Structure of the Sld3/Treslin domain from yeast Sld3  
Authors : Itou, H.; Araki, H.; Shirakihara, Y.  
Deposited on : 2013-09-05  
Resolution : 2.40 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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  
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

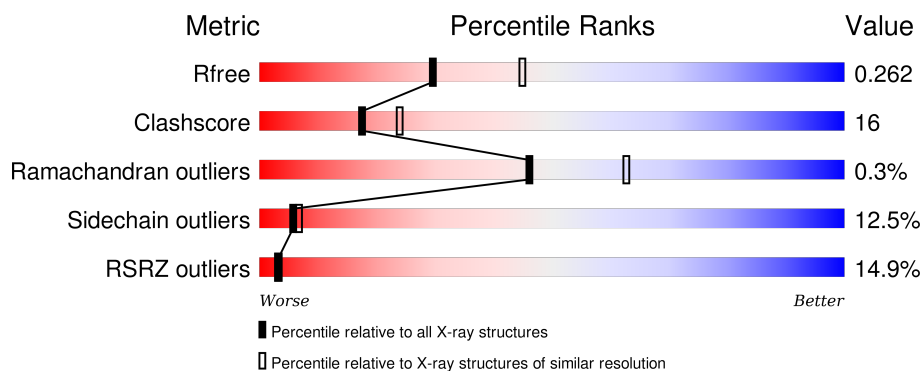
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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))
$R_{free}$	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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. 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	292	<div> <div>3%</div> <div>52%</div> <div>18%</div> <div>•</div> <div>26%</div> </div>
1	B	292	<div> <div>3%</div> <div>49%</div> <div>20%</div> <div>•</div> <div>27%</div> </div>
1	C	292	<div> <div>26%</div> <div>53%</div> <div>14%</div> <div>•</div> <div>30%</div> </div>

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	502	-	-	X	-
3	EDO	A	503	-	-	X	X
3	EDO	A	504	-	-	X	-
3	EDO	A	505	-	-	X	X
3	EDO	A	506	-	-	X	X
3	EDO	A	507	-	-	X	-
3	EDO	A	508	-	-	X	X
3	EDO	A	509	-	-	-	X
3	EDO	A	514	-	-	X	X
3	EDO	A	515	-	X	-	-
3	EDO	A	516	-	-	X	-
3	EDO	A	517	-	-	X	-
3	EDO	B	502	-	-	X	-
3	EDO	B	503	-	-	X	X
3	EDO	B	504	-	-	X	-
3	EDO	B	505	-	-	X	X
3	EDO	B	506	-	-	X	-
3	EDO	B	508	-	-	-	X
3	EDO	B	509	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5473 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 DNA replication regulator SLD3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	215	Total	C	N	O	S	0	0	0
			1766	1145	289	326	6			
1	B	212	Total	C	N	O	S	0	0	0
			1741	1127	287	321	6			
1	C	205	Total	C	N	O	S	0	0	0
			1686	1095	276	309	6			

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	147	MET	-	EXPRESSION TAG	UNP P53135
A	431	LEU	-	EXPRESSION TAG	UNP P53135
A	432	GLU	-	EXPRESSION TAG	UNP P53135
A	433	HIS	-	EXPRESSION TAG	UNP P53135
A	434	HIS	-	EXPRESSION TAG	UNP P53135
A	435	HIS	-	EXPRESSION TAG	UNP P53135
A	436	HIS	-	EXPRESSION TAG	UNP P53135
A	437	HIS	-	EXPRESSION TAG	UNP P53135
A	438	HIS	-	EXPRESSION TAG	UNP P53135
B	147	MET	-	EXPRESSION TAG	UNP P53135
B	431	LEU	-	EXPRESSION TAG	UNP P53135
B	432	GLU	-	EXPRESSION TAG	UNP P53135
B	433	HIS	-	EXPRESSION TAG	UNP P53135
B	434	HIS	-	EXPRESSION TAG	UNP P53135
B	435	HIS	-	EXPRESSION TAG	UNP P53135
B	436	HIS	-	EXPRESSION TAG	UNP P53135
B	437	HIS	-	EXPRESSION TAG	UNP P53135
B	438	HIS	-	EXPRESSION TAG	UNP P53135
C	147	MET	-	EXPRESSION TAG	UNP P53135
C	431	LEU	-	EXPRESSION TAG	UNP P53135
C	432	GLU	-	EXPRESSION TAG	UNP P53135
C	433	HIS	-	EXPRESSION TAG	UNP P53135
C	434	HIS	-	EXPRESSION TAG	UNP P53135

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Chain	Residue	Modelled	Actual	Comment	Reference
C	435	HIS	-	EXPRESSION TAG	UNP P53135
C	436	HIS	-	EXPRESSION TAG	UNP P53135
C	437	HIS	-	EXPRESSION TAG	UNP P53135
C	438	HIS	-	EXPRESSION TAG	UNP P53135

- Molecule 2 is SULFATE ION (three-letter code: SO<sub>4</sub>) (formula: O<sub>4</sub>S).



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

- Molecule 3 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
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		

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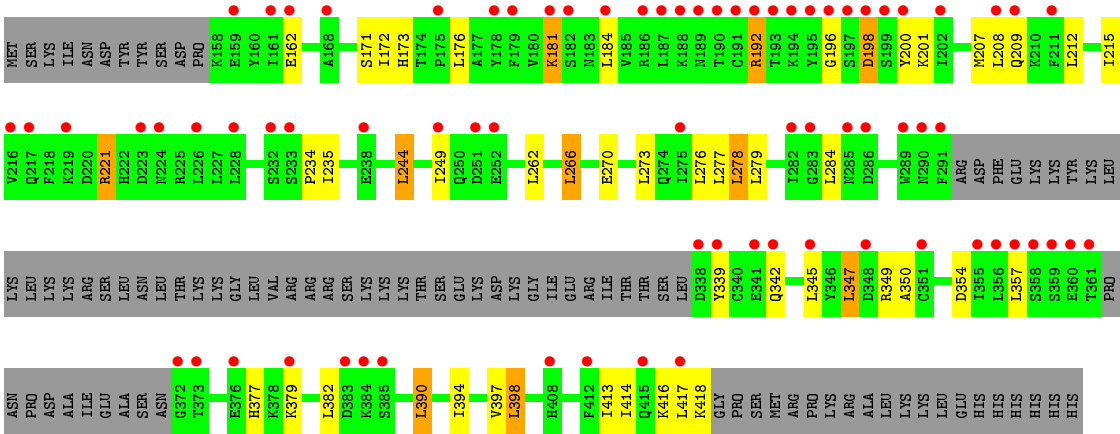
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	68	Total O 68 68	0	0
4	B	82	Total O 82 82	0	0
4	C	19	Total O 19 19	0	0







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	65.26 Å 92.82 Å 160.87 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.64 – 2.40 19.64 – 2.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (19.64-2.40) 88.1 (19.64-2.40)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	8.06 (at 2.41 Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.217 , 0.262 0.218 , 0.262	Depositor DCC
$R_{free}$ test set	1714 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	34.4	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 58.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	1 of 34298 reflections (0.003%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	5473	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 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	A	0.98	0/1799	1.08	6/2424 (0.2%)
1	B	0.94	1/1772 (0.1%)	1.08	11/2386 (0.5%)
1	C	0.62	0/1715	0.88	3/2309 (0.1%)
All	All	0.87	1/5286 (0.0%)	1.02	20/7119 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	172	ILE	N-CA	7.57	1.61	1.46

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	221	ARG	NE-CZ-NH2	-14.82	112.89	120.30
1	A	221	ARG	NE-CZ-NH1	12.14	126.37	120.30
1	B	221	ARG	NE-CZ-NH2	-11.35	114.62	120.30
1	B	221	ARG	NE-CZ-NH1	10.31	125.46	120.30
1	C	221	ARG	NE-CZ-NH1	9.58	125.09	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	171	SER	Peptide

## 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	1766	0	1804	72	0
1	B	1741	0	1783	71	0
1	C	1686	0	1738	23	0
2	A	5	0	0	1	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
3	A	64	0	96	57	0
3	B	32	0	48	45	0
4	A	68	0	0	4	0
4	B	82	0	0	5	0
4	C	19	0	0	3	0
All	All	5473	0	5469	172	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 16.

The worst 5 of 172 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:256:SER:O	3:A:502:EDO:H21	1.04	1.22
1:B:222:HIS:CD2	3:B:506:EDO:H22	1.78	1.19
1:A:256:SER:O	3:A:502:EDO:C2	1.93	1.17
1:B:222:HIS:HD2	3:B:506:EDO:C2	1.74	1.00
1:B:222:HIS:CD2	3:B:506:EDO:C2	2.45	0.99

There are no symmetry-related clashes.

## 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	209/292 (72%)	205 (98%)	4 (2%)	0	100	100
1	B	206/292 (70%)	201 (98%)	4 (2%)	1 (0%)	34	48
1	C	199/292 (68%)	192 (96%)	6 (3%)	1 (0%)	34	48
All	All	614/876 (70%)	598 (97%)	14 (2%)	2 (0%)	46	63

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	198	ASP
1	B	198	ASP

### 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	201/273 (74%)	181 (90%)	20 (10%)	9	14
1	B	198/273 (72%)	169 (85%)	29 (15%)	4	4
1	C	192/273 (70%)	167 (87%)	25 (13%)	5	6
All	All	591/819 (72%)	517 (88%)	74 (12%)	6	7

5 of 74 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	276	LEU

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Mol	Chain	Res	Type
1	B	370	SER
1	C	357	LEU
1	B	277	LEU
1	B	289	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 25 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	285	ASN
1	B	375	GLN
1	C	395	ASN
1	B	371	ASN
1	B	380	ASN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

27 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
2	SO4	A	501	-	4,4,4	2.18	2 (50%)	6,6,6	0.46	0
3	EDO	A	502	-	3,3,3	1.08	0	2,2,2	0.70	0
3	EDO	A	503	-	3,3,3	0.95	0	2,2,2	0.77	0
3	EDO	A	504	-	3,3,3	0.83	0	2,2,2	0.63	0
3	EDO	A	505	-	3,3,3	0.98	0	2,2,2	1.21	0
3	EDO	A	506	-	3,3,3	0.74	0	2,2,2	0.71	0
3	EDO	A	507	-	3,3,3	0.52	0	2,2,2	0.63	0
3	EDO	A	508	-	3,3,3	1.49	0	2,2,2	0.28	0
3	EDO	A	509	-	3,3,3	0.83	0	2,2,2	0.62	0
3	EDO	A	510	-	3,3,3	0.81	0	2,2,2	0.10	0
3	EDO	A	511	-	3,3,3	0.43	0	2,2,2	0.70	0
3	EDO	A	512	-	3,3,3	0.78	0	2,2,2	0.86	0
3	EDO	A	513	-	3,3,3	0.37	0	2,2,2	1.30	0
3	EDO	A	514	-	3,3,3	0.61	0	2,2,2	0.51	0
3	EDO	A	515	-	3,3,3	2.25	2 (66%)	2,2,2	1.49	1 (50%)
3	EDO	A	516	-	3,3,3	0.37	0	2,2,2	0.34	0
3	EDO	A	517	-	3,3,3	0.72	0	2,2,2	0.51	0
2	SO4	B	501	-	4,4,4	1.44	0	6,6,6	0.85	0
3	EDO	B	502	-	3,3,3	0.69	0	2,2,2	0.04	0
3	EDO	B	503	-	3,3,3	1.25	0	2,2,2	1.46	1 (50%)
3	EDO	B	504	-	3,3,3	0.53	0	2,2,2	0.30	0
3	EDO	B	505	-	3,3,3	1.28	0	2,2,2	1.15	0
3	EDO	B	506	-	3,3,3	0.67	0	2,2,2	0.37	0
3	EDO	B	507	-	3,3,3	0.58	0	2,2,2	0.30	0
3	EDO	B	508	-	3,3,3	0.93	0	2,2,2	0.33	0
3	EDO	B	509	-	3,3,3	0.72	0	2,2,2	0.18	0
2	SO4	C	501	-	4,4,4	1.91	2 (50%)	6,6,6	0.31	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
2	SO4	A	501	-	-	0/0/0/0	0/0/0/0
3	EDO	A	502	-	-	0/1/1/1	0/0/0/0
3	EDO	A	503	-	-	0/1/1/1	0/0/0/0
3	EDO	A	504	-	-	0/1/1/1	0/0/0/0
3	EDO	A	505	-	-	0/1/1/1	0/0/0/0
3	EDO	A	506	-	-	0/1/1/1	0/0/0/0
3	EDO	A	507	-	-	0/1/1/1	0/0/0/0
3	EDO	A	508	-	-	0/1/1/1	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	A	509	-	-	0/1/1/1	0/0/0/0
3	EDO	A	510	-	-	0/1/1/1	0/0/0/0
3	EDO	A	511	-	-	0/1/1/1	0/0/0/0
3	EDO	A	512	-	-	0/1/1/1	0/0/0/0
3	EDO	A	513	-	-	0/1/1/1	0/0/0/0
3	EDO	A	514	-	-	0/1/1/1	0/0/0/0
3	EDO	A	515	-	-	0/1/1/1	0/0/0/0
3	EDO	A	516	-	-	0/1/1/1	0/0/0/0
3	EDO	A	517	-	-	0/1/1/1	0/0/0/0
2	SO4	B	501	-	-	0/0/0/0	0/0/0/0
3	EDO	B	502	-	-	0/1/1/1	0/0/0/0
3	EDO	B	503	-	-	0/1/1/1	0/0/0/0
3	EDO	B	504	-	-	0/1/1/1	0/0/0/0
3	EDO	B	505	-	-	0/1/1/1	0/0/0/0
3	EDO	B	506	-	-	0/1/1/1	0/0/0/0
3	EDO	B	507	-	-	0/1/1/1	0/0/0/0
3	EDO	B	508	-	-	0/1/1/1	0/0/0/0
3	EDO	B	509	-	-	0/1/1/1	0/0/0/0
2	SO4	C	501	-	-	0/0/0/0	0/0/0/0

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	SO4	O3-S	2.02	1.54	1.47
2	C	501	SO4	O1-S	2.05	1.54	1.47
3	A	515	EDO	C2-C1	2.19	1.63	1.47
2	A	501	SO4	O3-S	2.45	1.56	1.47
2	A	501	SO4	O2-S	2.64	1.56	1.47

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	503	EDO	O2-C2-C1	2.05	127.22	112.54
3	A	515	EDO	O2-C2-C1	2.09	127.56	112.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

23 monomers are involved in 103 short contacts:



Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	SO4	1	0
3	A	502	EDO	11	0
3	A	503	EDO	5	0
3	A	504	EDO	4	0
3	A	505	EDO	7	0
3	A	506	EDO	4	0
3	A	507	EDO	4	0
3	A	508	EDO	4	0
3	A	509	EDO	1	0
3	A	510	EDO	2	0
3	A	513	EDO	2	0
3	A	514	EDO	5	0
3	A	515	EDO	3	0
3	A	516	EDO	8	0
3	A	517	EDO	4	0
3	B	502	EDO	8	0
3	B	503	EDO	16	0
3	B	504	EDO	7	0
3	B	505	EDO	5	0
3	B	506	EDO	8	0
3	B	507	EDO	1	0
3	B	508	EDO	1	0
3	B	509	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	215/292 (73%)	0.06	10 (4%) 35 36	14, 31, 59, 86	0
1	B	212/292 (72%)	-0.02	9 (4%) 40 41	13, 33, 68, 84	0
1	C	205/292 (70%)	1.86	75 (36%) 0 0	44, 75, 128, 170	0
All	All	632/876 (72%)	0.62	94 (14%) 3 3	13, 42, 104, 170	0

The worst 5 of 94 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	199	SER	9.2
1	C	178	TYR	8.5
1	C	193	THR	8.1
1	C	195	TYR	7.9
1	C	289	TRP	7.7

### 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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	EDO	B	503	4/4	0.90	0.33	9.63	21,28,28,28	0
3	EDO	A	514	4/4	0.94	0.36	8.09	30,32,35,36	0
3	EDO	B	508	4/4	0.75	0.30	5.85	62,62,70,77	0
3	EDO	A	506	4/4	0.90	0.31	5.56	28,31,32,34	0
3	EDO	B	509	4/4	0.88	0.26	5.16	39,44,44,54	0
3	EDO	A	505	4/4	0.93	0.20	2.70	29,31,32,34	0
3	EDO	A	508	4/4	0.90	0.23	2.34	34,43,45,54	0
3	EDO	A	509	4/4	0.79	0.21	2.22	57,60,64,64	0
3	EDO	A	503	4/4	0.90	0.21	2.10	28,35,40,47	0
3	EDO	B	505	4/4	0.87	0.24	2.02	26,30,32,44	0
3	EDO	B	502	4/4	0.96	0.16	1.70	21,22,23,24	0
3	EDO	B	504	4/4	0.92	0.22	1.67	32,33,42,45	0
3	EDO	A	515	4/4	0.92	0.18	1.56	25,28,29,32	0
3	EDO	B	506	4/4	0.91	0.25	1.22	28,33,35,36	0
3	EDO	A	502	4/4	0.96	0.18	1.06	31,35,37,37	0
3	EDO	A	511	4/4	0.93	0.23	0.91	39,46,49,54	0
2	SO4	C	501	5/5	0.62	0.39	0.39	118,121,128,130	0
3	EDO	A	513	4/4	0.79	0.22	0.34	52,54,57,62	0
3	EDO	A	507	4/4	0.91	0.16	0.23	35,39,44,48	0
3	EDO	A	504	4/4	0.96	0.20	0.22	20,21,24,25	0
2	SO4	A	501	5/5	0.83	0.21	-0.21	66,72,92,105	0
3	EDO	B	507	4/4	0.97	0.14	-0.28	29,34,36,37	0
2	SO4	B	501	5/5	1.00	0.12	-1.09	19,19,22,23	0
3	EDO	A	516	4/4	0.95	0.12	-1.69	35,36,36,38	0
3	EDO	A	517	4/4	0.86	0.32	-	44,49,51,52	0
3	EDO	A	510	4/4	0.67	0.24	-	51,51,54,58	0
3	EDO	A	512	4/4	0.83	0.29	-	38,40,45,47	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.