Full wwPDB X-ray Structure Validation Report (i)

Jan 24, 2023 – 12:50 PM EST

PDB ID	:	3KTC
Title	:	Crystal structure of Putative sugar isomerase (YP_050048.1) from ERWINIA CAROTOVORA ATROSEPTICA SCRI1043 at 1.54 A resolution
Authors	:	Joint Center for Structural Genomics (JCSG)
Deposited on	:	2009-11-24
Resolution	:	1.54 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 (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:

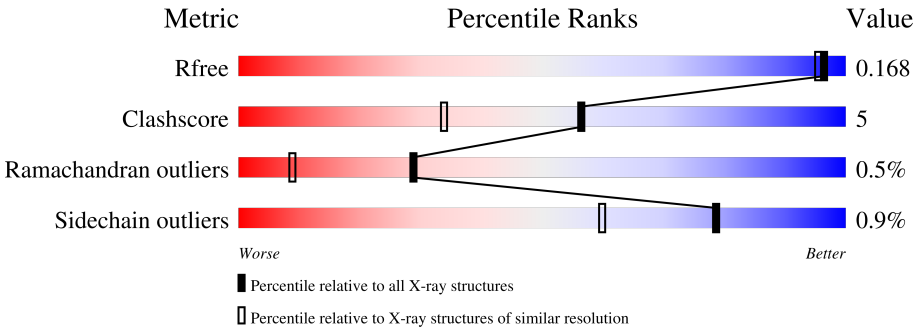
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.31.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.2

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 1.54 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}	130704	2556 (1.56-1.52)
Clashscore	141614	2634 (1.56-1.52)
Ramachandran outliers	138981	2580 (1.56-1.52)
Sidechain outliers	138945	2577 (1.56-1.52)

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%

Mol	Chain	Length	Quality of chain
1	A	333	92% 7% .
1	B	333	92% 6% ..

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
5	UNL	A	337	-	-	X	-
5	UNL	A	338	-	-	X	-
5	UNL	B	337	-	-	X	-

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 6418 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 Xylose isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total 2813	C 1797	N 481	O 528	Se 7	0	25	0
1	B	330	Total 2796	C 1789	N 477	O 524	Se 6	0	24	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q6D5T7
B	0	GLY	-	expression tag	UNP Q6D5T7

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total 3	Fe 3	0	1
2	B	2	Total 3	Fe 3	0	1

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Mg 1	0	0
3	B	1	Total 1	Mg 1	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Cl 1	0	0

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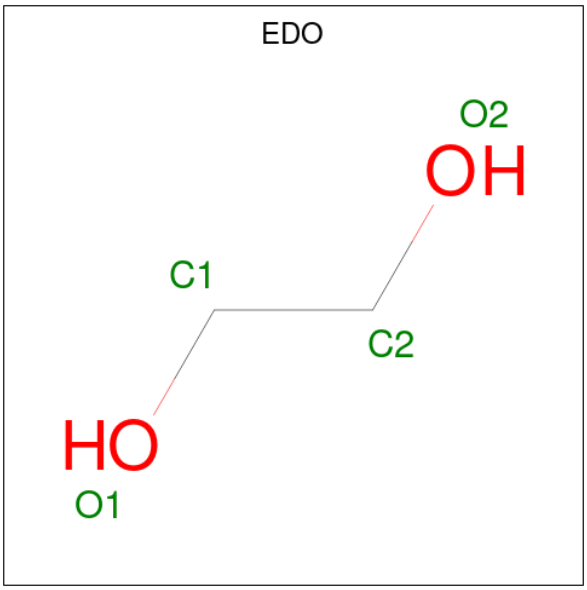
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Cl 1 1	0	0

- Molecule 5 is UNKNOWN LIGAND (three-letter code: UNL) (formula:).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	2	Total O 18 18	0	0
5	B	1	Total O 9 9	0	0

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

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

- Molecule 7 is water.

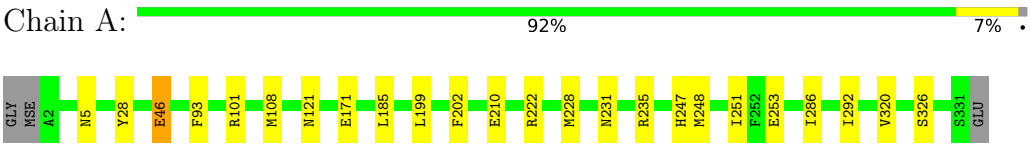
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	342	Total O 344 344	0	2
7	B	367	Total O 368 368	0	1

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.

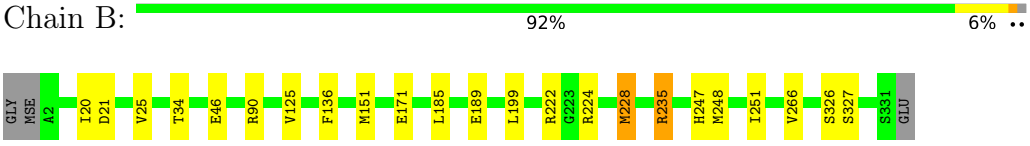
- Molecule 1: Xylose isomerase

Chain A:



- Molecule 1: Xylose isomerase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	79.65Å 79.65Å 194.41Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.32 – 1.54 29.33 – 1.54	Depositor EDS
% Data completeness (in resolution range)	99.3 (29.32-1.54) 99.3 (29.33-1.54)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.72 (at 1.54Å)	Xtriage
Refinement program	REFMAC 5.2.0019, PHENIX	Depositor
R, R_{free}	0.136 , 0.157 0.151 , 0.168	Depositor DCC
R_{free} test set	5292 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	15.2	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 42.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.025 for -h,-k,l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	6418	wwPDB-VP
Average B, all atoms (Å^2)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.73% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CL, UNL, MG, EDO, FE

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	0.64	1/2886 (0.0%)	0.76	1/3916 (0.0%)
1	B	0.73	2/2869 (0.1%)	0.78	3/3895 (0.1%)
All	All	0.69	3/5755 (0.1%)	0.77	4/7811 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	228	MSE	CG-SE	-18.37	1.32	1.95
1	A	228	MSE	CG-SE	-10.78	1.58	1.95
1	B	228	MSE	CB-CG	7.26	1.74	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	101	ARG	NE-CZ-NH1	5.91	123.25	120.30
1	B	235[A]	ARG	NE-CZ-NH1	5.39	123.00	120.30
1	B	235[B]	ARG	NE-CZ-NH1	5.39	123.00	120.30
1	B	90	ARG	NE-CZ-NH2	-5.26	117.67	120.30

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	2813	0	2697	18	0
1	B	2796	0	2682	23	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	18	0	0	8	0
5	B	9	0	0	7	0
6	A	24	0	36	3	0
6	B	36	0	54	4	0
7	A	344	0	0	5	0
7	B	368	0	0	3	0
All	All	6418	0	5469	56	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 5.

All (56) 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:B:228:MSE:SE	1:B:228:MSE:CG	1.33	1.52
5:A:338:UNL:O1	5:A:338:UNL:O2	1.53	1.26
5:A:338:UNL:O6	5:A:338:UNL:O7	1.60	1.19
5:A:338:UNL:O6	5:A:338:UNL:O5	1.59	1.16
1:B:228:MSE:SE	1:B:228:MSE:HG3	1.87	1.11
5:A:338:UNL:O2	5:A:338:UNL:O4	1.68	1.09
1:B:228:MSE:SE	1:B:228:MSE:HG2	1.87	1.08
5:A:337:UNL:O8	5:A:337:UNL:O9	1.70	1.07
5:A:337:UNL:O1	5:A:337:UNL:O3	1.75	1.04
5:B:337:UNL:O4	5:B:337:UNL:O5	1.78	0.99
1:A:251[B]:ILE:HG21	7:A:379:HOH:O	1.63	0.97
5:B:337:UNL:O2	5:B:337:UNL:O3	1.81	0.97
1:B:228:MSE:SE	1:B:228:MSE:CB	2.63	0.95
5:B:337:UNL:O3	5:B:337:UNL:O1	1.88	0.91
5:A:337:UNL:O8	5:A:337:UNL:O7	1.89	0.89
1:A:202:PHE:CZ	1:A:251[B]:ILE:HG22	2.17	0.79
5:A:337:UNL:O1	5:A:337:UNL:O2	1.99	0.79
1:B:228:MSE:CG	1:B:228:MSE:CE	2.62	0.77
5:B:337:UNL:O5	5:B:337:UNL:O6	2.03	0.77
1:A:210:GLU:HB3	6:A:342:EDO:H11	1.71	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:337:UNL:O5	5:B:337:UNL:O3	2.09	0.70
1:A:235[A]:ARG:HD2	7:A:496:HOH:O	1.96	0.64
1:B:34:THR:OG1	6:B:346:EDO:H21	1.98	0.64
1:A:251[B]:ILE:HD11	1:A:292:ILE:HG13	1.81	0.61
1:A:185:LEU:HD13	1:A:222[A]:ARG:HD2	1.82	0.59
1:A:202:PHE:CE2	1:A:251[B]:ILE:HG22	2.37	0.58
1:A:5[A]:ASN:ND2	7:A:566:HOH:O	2.36	0.58
1:B:248[B]:MSE:HA	1:B:251[B]:ILE:HG22	1.88	0.55
1:A:46:GLU:HB3	1:A:286:ILE:HD13	1.88	0.54
1:B:20[A]:ILE:HD11	1:B:25:VAL:O	2.08	0.53
1:B:185:LEU:HD13	1:B:222[A]:ARG:HD2	1.91	0.53
1:B:248[B]:MSE:HA	1:B:251[B]:ILE:CG2	2.38	0.53
1:B:326[B]:SER:O	1:B:327[B]:SER:HB2	2.09	0.52
1:A:247:HIS:O	1:A:251[B]:ILE:HG23	2.10	0.51
5:B:337:UNL:O7	5:B:337:UNL:O8	2.29	0.49
1:B:189[B]:GLU:OE2	1:B:222[B]:ARG:NH2	2.45	0.49
1:B:125:VAL:HG11	1:B:151:MSE:HE1	1.94	0.48
1:A:231:ASN:ND2	7:B:733[B]:HOH:O	2.46	0.48
1:A:251[B]:ILE:HD12	1:A:251[B]:ILE:C	2.34	0.48
5:B:337:UNL:O8	5:B:337:UNL:O9	2.33	0.46
6:A:342:EDO:H21	6:B:341:EDO:H12	1.98	0.45
1:B:247:HIS:O	1:B:251[B]:ILE:HG22	2.16	0.45
1:A:235[B]:ARG:HD3	7:A:484:HOH:O	2.15	0.45
1:B:20[B]:ILE:HG22	1:B:21:ASP:N	2.32	0.45
1:B:235[A]:ARG:HD3	7:B:422:HOH:O	2.18	0.43
1:B:248[B]:MSE:CA	1:B:251[B]:ILE:HG22	2.49	0.43
1:A:253:GLU:HA	1:B:136:PHE:CZ	2.55	0.42
1:A:28[B]:TYR:O	6:A:343:EDO:O1	2.25	0.42
1:A:93:PHE:CE1	1:A:108:MSE:HE1	2.54	0.42
1:A:320:VAL:HG21	1:B:136:PHE:CG	2.56	0.41
1:B:224:ARG:HE	6:B:338:EDO:H21	1.85	0.41
1:A:248[A]:MSE:SE	7:A:485:HOH:O	2.88	0.41
1:B:248[B]:MSE:O	1:B:251[B]:ILE:HG22	2.21	0.41
1:B:228:MSE:HE3	1:B:266:VAL:O	2.20	0.41
1:B:248[B]:MSE:SE	7:B:526:HOH:O	2.88	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	353/333 (106%)	341 (97%)	9 (2%)	3 (1%)	19 4
1	B	352/333 (106%)	333 (95%)	18 (5%)	1 (0%)	41 19
All	All	705/666 (106%)	674 (96%)	27 (4%)	4 (1%)	29 7

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	171	GLU
1	A	326[A]	SER
1	A	326[B]	SER
1	B	171	GLU

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	291/264 (110%)	288 (99%)	3 (1%)	76 55
1	B	288/264 (109%)	286 (99%)	2 (1%)	84 68
All	All	579/528 (110%)	574 (99%)	5 (1%)	78 60

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

Mol	Chain	Res	Type
1	A	46	GLU
1	A	121	ASN

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Mol	Chain	Res	Type
1	A	199	LEU
1	B	46	GLU
1	B	199	LEU

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 28 ligands modelled in this entry, 10 are monoatomic and 3 are unknown - leaving 15 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	$\#\|Z\| > 2$
6	EDO	A	341	-	3,3,3	0.41	0	2,2,2	0.42	0
6	EDO	B	343	-	3,3,3	0.49	0	2,2,2	0.22	0
6	EDO	B	342	-	3,3,3	0.46	0	2,2,2	0.23	0
6	EDO	A	342	-	3,3,3	0.61	0	2,2,2	0.88	0
6	EDO	A	340	-	3,3,3	0.47	0	2,2,2	0.34	0
6	EDO	B	344	-	3,3,3	0.40	0	2,2,2	0.41	0
6	EDO	A	343	-	3,3,3	0.51	0	2,2,2	0.18	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
6	EDO	A	339	-	3,3,3	0.55	0	2,2,2	0.21	0
6	EDO	B	340	-	3,3,3	0.50	0	2,2,2	0.34	0
6	EDO	B	345	-	3,3,3	0.45	0	2,2,2	0.34	0
6	EDO	A	344	-	3,3,3	0.60	0	2,2,2	0.12	0
6	EDO	B	338	-	3,3,3	0.48	0	2,2,2	0.10	0
6	EDO	B	341	-	3,3,3	0.60	0	2,2,2	0.22	0
6	EDO	B	339	-	3,3,3	0.51	0	2,2,2	0.40	0
6	EDO	B	346	-	3,3,3	0.62	0	2,2,2	0.30	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
6	EDO	A	341	-	-	1/1/1/1	-
6	EDO	B	343	-	-	0/1/1/1	-
6	EDO	B	342	-	-	0/1/1/1	-
6	EDO	A	342	-	-	1/1/1/1	-
6	EDO	A	340	-	-	0/1/1/1	-
6	EDO	B	344	-	-	0/1/1/1	-
6	EDO	A	343	-	-	0/1/1/1	-
6	EDO	A	339	-	-	1/1/1/1	-
6	EDO	B	340	-	-	1/1/1/1	-
6	EDO	B	345	-	-	0/1/1/1	-
6	EDO	A	344	-	-	0/1/1/1	-
6	EDO	B	338	-	-	1/1/1/1	-
6	EDO	B	341	-	-	0/1/1/1	-
6	EDO	B	339	-	-	0/1/1/1	-
6	EDO	B	346	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	338	EDO	O1-C1-C2-O2
6	B	346	EDO	O1-C1-C2-O2
6	A	341	EDO	O1-C1-C2-O2
6	A	342	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
6	B	340	EDO	O1-C1-C2-O2
6	A	339	EDO	O1-C1-C2-O2

There are no ring outliers.

6 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	343	EDO	1	0
6	A	342	EDO	2	0
6	A	343	EDO	1	0
6	B	338	EDO	1	0
6	B	341	EDO	1	0
6	B	346	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

6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.