

Full wwPDB X-ray Structure Validation Report (i)

Jan 24, 2023 – 05:11 PM EST

PDB ID : 3KTD

Title : CRYSTAL STRUCTURE OF A PUTATIVE PREPHENATE DEHYDROGE-

NASE (CGL0226) FROM CORYNEBACTERIUM GLUTAMICUM ATCC

13032 AT 2.60 A RESOLUTION

Authors : Joint Center for Structural Genomics (JCSG)

Deposited on : 2009-11-24

Resolution : 2.60 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 (1)) were used in the production of this report:

 $Mol Probity \quad : \quad 4.02b\text{--}467$

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)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

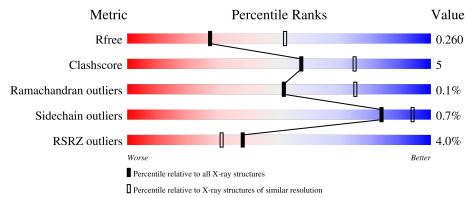
Validation Pipeline (wwPDB-VP) : 2.31.2

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 2.60 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries,\ resolution\ range(\mathring{A})}) \end{array}$
R_{free}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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% 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	341	86%	9%	5%
1	В	341	81%	12%	6%
1	С	341	8%	12%	7%
1	D	341	85%	9%	6%



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 9588 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 Prephenate dehydrogenase.

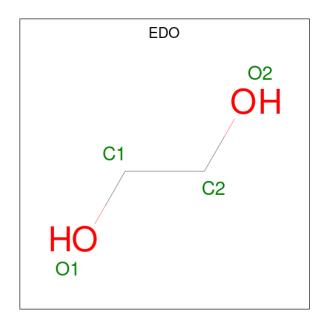
Mol	Chain	Residues		A	Atoms	5			ZeroOcc	AltConf	Trace
1	Λ	323	Total	С	N	О	S	Se	0	3	0
1	A	323	2409	1506	432	463	2	6	U	J	
1	В	319	Total	С	N	О	S	Se	0	3	0
1	Ъ	319	2350	1474	418	450	2	6	0	3	0
1	С	318	Total	С	N	О	S	Se	0	1	0
1		310	2354	1477	424	445	2	6	0	1	0
1	D	321	Total	С	N	О	S	Se	0	0	0
	ע	321	2372	1485	425	454	2	6	U	U	U

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q8NTS6
В	0	GLY	-	expression tag	UNP Q8NTS6
С	0	GLY	-	expression tag	UNP Q8NTS6
D	0	GLY	-	expression tag	UNP Q8NTS6

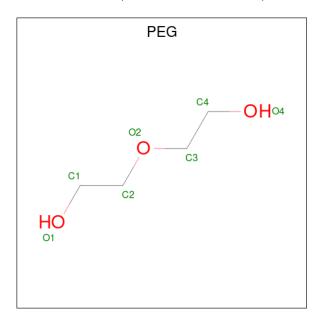
• Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).





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

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



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

• Molecule 4 is water.



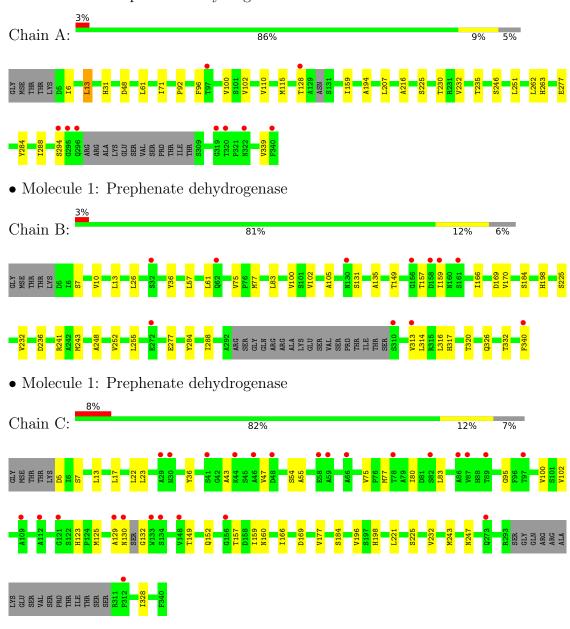
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	30	Total O 30 30	0	0
4	В	19	Total O 19 19	0	0
4	С	22	Total O 22 22	0	0
4	D	21	Total O 21 21	0	0



3 Residue-property plots (i)

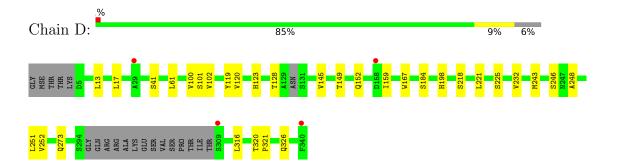
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



• Molecule 1: Prephenate dehydrogenase







4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	99.32Å 107.23Å 146.91Å	Donositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.53 - 2.60	Depositor
Resolution (A)	29.53 - 2.60	EDS
% Data completeness	99.9 (29.53-2.60)	Depositor
(in resolution range)	100.0 (29.53-2.60)	EDS
R_{merge}	0.11	Depositor
R_{sym}	0.11	Depositor
$< I/\sigma(I) > 1$	2.16 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.5.0053, PHENIX	Depositor
D D.	0.213 , 0.254	Depositor
R, R_{free}	0.222 , 0.260	DCC
R_{free} test set	2475 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	48.5	Xtriage
Anisotropy	0.502	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.32, 50.5	EDS
L-test for twinning ²	$ < L > = 0.49, < L^2> = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9588	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.54% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities estimated from amplitudes.

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: EDO, 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		
IVIOI	Chain	$\mid \text{RMSZ} \mid \# Z > 5$		RMSZ	# Z > 5	
1	A	0.42	0/2457	0.61	0/3334	
1	В	0.38	0/2399	0.56	0/3266	
1	С	0.40	0/2392	0.56	0/3251	
1	D	0.40	0/2410	0.58	0/3272	
All	All	0.40	0/9658	0.58	0/13123	

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 (i)

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	2409	0	2354	21	0
1	В	2350	0	2263	29	0
1	С	2354	0	2292	33	0
1	D	2372	0	2317	22	0
2	A	4	0	6	0	0
3	A	7	0	10	0	0
4	A	30	0	0	1	0
4	В	19	0	0	0	0
4	С	22	0	0	0	0



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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	21	0	0	1	0
All	All	9588	0	9242	91	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 (91) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${ m distance} \; ({ m \AA})$	overlap (Å)
1:B:75:VAL:HG21	1:B:83:LEU:HD12	1.70	0.73
1:D:13:LEU:HD21	1:D:61:LEU:HD21	1.76	0.68
1:C:77:MSE:HG2	1:C:100:VAL:HG13	1.77	0.67
1:B:317[B]:HIS:CD2	1:B:340:PHE:HB2	2.30	0.67
1:C:13:LEU:HD23	1:C:83:LEU:HD12	1.77	0.66
1:C:100:VAL:HG12	1:C:102:VAL:H	1.61	0.64
1:C:17:LEU:HD22	1:C:129:ALA:HB2	1.86	0.58
1:C:75:VAL:CG2	1:C:80:ILE:HD13	2.36	0.55
1:B:100:VAL:HG12	1:B:102:VAL:H	1.71	0.55
1:B:149:THR:HA	1:B:184:SER:O	2.06	0.55
1:A:277:GLU:HG3	1:B:241:ARG:CZ	2.37	0.55
1:C:129:ALA:O	1:C:130:ASN:CB	2.55	0.54
1:B:157:THR:HG22	1:B:159:ILE:HG23	1.91	0.53
1:C:43:ALA:O	1:C:47:VAL:HG23	2.08	0.53
1:C:123:HIS:CE1	1:C:125:MSE:SE	3.11	0.53
1:C:225:SER:CB	1:D:232:VAL:HG11	2.39	0.52
1:A:102:VAL:HG21	1:A:246:SER:HB2	1.91	0.52
1:C:130:ASN:O	1:C:132:GLY:N	2.42	0.52
1:C:77:MSE:HG2	1:C:100:VAL:CG1	2.40	0.52
1:C:232:VAL:HG11	1:D:225:SER:HB2	1.91	0.52
1:A:262:LEU:HD22	1:B:255:LEU:HD11	1.93	0.51
1:C:149:THR:HA	1:C:184:SER:O	2.10	0.51
1:C:225:SER:HB3	1:D:232:VAL:HG11	1.92	0.51
1:B:36:TYR:CE2	1:B:57:LEU:HD13	2.45	0.51
1:C:159:ILE:HD12	1:C:159:ILE:O	2.11	0.51
1:C:125:MSE:HE2	1:D:221:LEU:O	2.11	0.50
1:C:129:ALA:O	1:C:130:ASN:HB2	2.11	0.50
1:A:159:ILE:C	1:A:159:ILE:HD12	2.31	0.50
1:B:77:MSE:HE2	1:B:105:ALA:HB1	1.94	0.50
1:B:316:LEU:HD21	1:B:326:GLN:HB3	1.94	0.50
1:D:17:LEU:HD23	1:D:128:THR:O	2.10	0.50
1:C:23:LEU:HD23	1:C:23:LEU:O	2.12	0.50
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Continuea from previ		Interatomic	Clash
Atom-1	Atom-2	${\rm distance}\ (\mathring{\rm A})$	overlap (Å)
1:C:22:LEU:HD12	1:C:177:VAL:HG11	1.93	0.49
1:A:232:VAL:HG11	1:B:225:SER:CB	2.42	0.49
1:A:216:ALA:HB3	1:B:332:THR:O	2.13	0.49
1:A:110:VAL:HG13	1:A:115:MSE:HB2	1.93	0.49
1:C:198:HIS:CD2	1:C:243:MSE:HG2	2.48	0.49
1:B:131:SER:HB3	1:B:135:ALA:HB2	1.96	0.48
1:C:198:HIS:CE1	1:C:247:ASN:HD22	2.31	0.48
1:C:221:LEU:CD2	1:D:145:VAL:HG11	2.44	0.48
1:C:157:THR:HG22	1:C:159:ILE:HG23	1.96	0.47
1:B:316:LEU:HD21	1:B:326:GLN:CB	2.43	0.47
1:D:248:ALA:O	1:D:252:VAL:HG23	2.14	0.47
1:A:100:VAL:HG12	1:A:102:VAL:H	1.79	0.47
1:B:10:VAL:HG21	1:B:26:LEU:HD13	1.97	0.46
1:B:248:ALA:O	1:B:252:VAL:HG23	2.15	0.46
1:D:198:HIS:HB2	1:D:251:LEU:HD13	1.97	0.46
1:B:159:ILE:HD12	1:B:159:ILE:C	2.36	0.46
1:C:7:SER:HB3	1:C:169:ASP:OD1	2.16	0.46
1:C:196:VAL:HG11	1:D:218:SER:O	2.16	0.45
1:D:101:SER:HA	1:D:123:HIS:CD2	2.51	0.45
1:A:207:LEU:HD23	1:A:230:THR:HG23	1.97	0.45
1:D:100:VAL:HG12	1:D:102:VAL:H	1.82	0.45
1:C:43:ALA:CB	1:C:55:ALA:HB2	2.47	0.45
1:A:6:ILE:HB	1:A:31[B]:HIS:CE1	2.53	0.44
1:C:232:VAL:HG11	1:D:225:SER:CB	2.47	0.44
1:A:96:PHE:CZ	1:A:115:MSE:HE2	2.53	0.44
1:B:317[B]:HIS:HD2	1:B:340:PHE:HB2	1.79	0.44
1:D:119:TYR:O	1:D:152:GLN:NE2	2.50	0.44
1:A:194:ALA:HB1	1:A:251:LEU:HA	2.00	0.44
1:A:128:THR:HG21	1:A:235:THR:HG22	1.98	0.44
1:D:159:ILE:C	1:D:159:ILE:HD12	2.38	0.44
1:D:120:VAL:HG21	1:D:167:TRP:HA	2.00	0.44
1:A:339:VAL:HG23	1:A:339:VAL:O	2.18	0.43
1:B:13:LEU:HD21	1:B:61:LEU:HD21	1.99	0.43
1:B:313:VAL:C	1:B:314:LEU:HD12	2.38	0.43
1:A:284:TYR:O	1:A:288:ILE:HG12	2.18	0.43
1:A:263:HIS:HB2	4:A:367:HOH:O	2.17	0.43
1:B:284:TYR:CE1	1:B:288:ILE:HD11	2.54	0.43
1:A:13:LEU:HD21	1:A:61:LEU:HD21	1.99	0.43
1:B:159:ILE:HD12	1:B:159:ILE:O	2.19	0.43
1:D:102:VAL:HG21	1:D:246:SER:HB2	2.01	0.43
1:D:273:GLN:N	4:D:361:HOH:O	2.38	0.42



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Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${ m distance}({ m \AA})$	overlap (Å)
1:D:149:THR:HA	1:D:184:SER:O	2.20	0.42
1:A:225:SER:HB2	1:B:232:VAL:HG11	2.01	0.42
1:C:152:GLN:HB3	1:C:159:ILE:HG21	2.02	0.42
1:D:316:LEU:HD22	1:D:326:GLN:HB2	2.00	0.42
1:A:13:LEU:HD22	1:A:71:ILE:CG2	2.50	0.42
1:B:166:ILE:O	1:B:170:VAL:HG23	2.19	0.42
1:C:95:GLY:HA3	1:C:166:ILE:HD13	2.02	0.42
1:D:320:THR:HG22	1:D:321:PRO:O	2.20	0.41
1:D:198:HIS:CD2	1:D:243:MSE:HB3	2.55	0.41
1:C:75:VAL:CG2	1:C:80:ILE:CD1	2.99	0.41
1:A:232:VAL:HG11	1:B:225:SER:HB2	2.01	0.41
1:B:198:HIS:CD2	1:B:243:MSE:HG2	2.56	0.41
1:C:36:TYR:HA	1:C:54:SER:O	2.21	0.41
1:B:7:SER:OG	1:B:169:ASP:OD2	2.31	0.40
1:C:95:GLY:CA	1:C:166:ILE:HD13	2.51	0.40
1:A:48:ASP:HB3	1:C:328:ILE:HD13	2.03	0.40
1:B:317[A]:HIS:HD2	1:B:320:THR:HG23	1.86	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	Perce	ntiles
1	A	319/341~(94%)	300 (94%)	18 (6%)	1 (0%)	41	64
1	В	$318/341 \ (93\%)$	302 (95%)	16 (5%)	0	100	100
1	С	313/341~(92%)	292 (93%)	21 (7%)	0	100	100
1	D	315/341~(92%)	298 (95%)	17 (5%)	0	100	100
All	All	$1265/1364\ (93\%)$	1192 (94%)	72 (6%)	1 (0%)	51	75

All (1) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	A	294	SER

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	P	erce	ntiles
1	A	247/260~(95%)	245 (99%)	2 (1%)		81	92
1	В	$236/260 \ (91\%)$	234 (99%)	2 (1%)		81	92
1	С	235/260~(90%)	233 (99%)	2 (1%)		78	91
1	D	241/260~(93%)	240 (100%)	1 (0%)		91	97
All	All	959/1040 (92%)	952 (99%)	7 (1%)		84	94

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	92	PRO
1	В	236	ASP
1	В	277	GLU
1	С	5	ASP
1	С	160	ASN
1	D	41	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

\mathbf{Mol}	Chain	Res	Type
1	В	273	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

2 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	Tune	Chain	Res	Link	В	ond leng	${ m gths}$	В	ond ang	gles
MIOI	Mol Type Chain	Chain	nes	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PEG	A	342	-	6,6,6	0.48	0	5,5,5	0.32	0
2	EDO	A	341	-	3,3,3	0.54	0	2,2,2	0.19	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	A	342	-	-	2/4/4/4	-
2	EDO	A	341	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	341	EDO	O1-C1-C2-O2
3	A	342	PEG	C1-C2-O2-C3



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Mol	Chain	Res	Type	Atoms
3	A	342	PEG	O2-C3-C4-O4

There are no ring outliers.

No monomer is involved in short contacts.

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^{th} 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>	$\#\mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q < 0.9
1	A	317/341 (92%)	-0.06	9 (2%) 53 46	28, 37, 49, 68	0
1	В	313/341 (91%)	0.03	11 (3%) 44 36	29, 38, 47, 58	0
1	С	312/341 (91%)	0.31	26 (8%) 11 8	27, 40, 48, 67	0
1	D	315/341 (92%)	-0.13	4 (1%) 77 73	28, 37, 47, 65	0
All	All	1257/1364 (92%)	0.04	50 (3%) 38 31	27, 38, 48, 68	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	С	156	GLY	5.3
1	С	66	ALA	4.8
1	A	296	GLN	4.4
1	A	128	THR	4.0
1	В	130	ASN	3.9
1	С	48	ASP	3.9
1	С	312	PRO	3.8
1	С	29	ALA	3.6
1	С	129	ALA	3.5
1	D	309	SER	3.5
1	D	29	ALA	3.5
1	A	340	PHE	3.4
1	В	158	ASP	3.4
1	A	295	GLY	3.2
1	С	87	VAL	3.2
1	В	310	SER	3.2
1	С	89	THR	3.1
1	С	134	SER	3.1
1	В	340	PHE	3.0
1	С	30	ASN	2.9
1	В	272	GLU	2.9



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Mol	Chain	Res	Type	RSRZ	
1	D	340	PHE	2.9	
1	С	133	TRP	2.8	
1	С	46	ALA	2.8	
1	A	322	ASN	2.7	
1	С	97	THR	2.7	
1	С	86	ALA	2.6	
1	С	130	ASN	2.6	
1	С	41	SER	2.4	
1	C C	148	VAL	2.4	
1	С	44	LYS	2.4	
1	С	121	GLY	2.4	
1	В	62	GLN	2.3	
1	С	112	ALA	2.3	
1	С	273	GLN	2.3	
1	C	58	GLU	2.3	
1	В	313	VAL	2.3	
1	A	320	THR	2.3	
1	В	159	ILE	2.3	
1	С	82	SER	2.2	
1	A	319	GLY	2.2	
1	A	294	SER	2.2	
1	D	158	ASP	2.2	
1	С	59	ALA	2.2	
1	С	109	ALA	2.2	
1	С	78	THR	2.2	
1	В	156	GLY	2.1	
1	A	97	THR	2.1	
1	В	32	SER	2.0	
1	В	161	SER	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 monosaccharides 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, 95^{th} 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
3	PEG	A	342	7/7	0.77	0.20	80,80,83,84	0
2	EDO	A	341	4/4	0.84	0.22	76,76,76,76	0

6.5 Other polymers (i)

There are no such residues in this entry.

