



Full wwPDB X-ray Structure Validation Report i

Feb 28, 2014 – 09:39 AM GMT

PDB ID : 3KTD
Title : CRYSTAL STRUCTURE OF A PUTATIVE PREPHENATE DEHYDROGENASE (CGL0226) FROM CORYNEBACTERIUM GLUTAMICUM ATCC 13032 AT 2.60 Å RESOLUTION
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2009-11-24
Resolution : 2.60 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

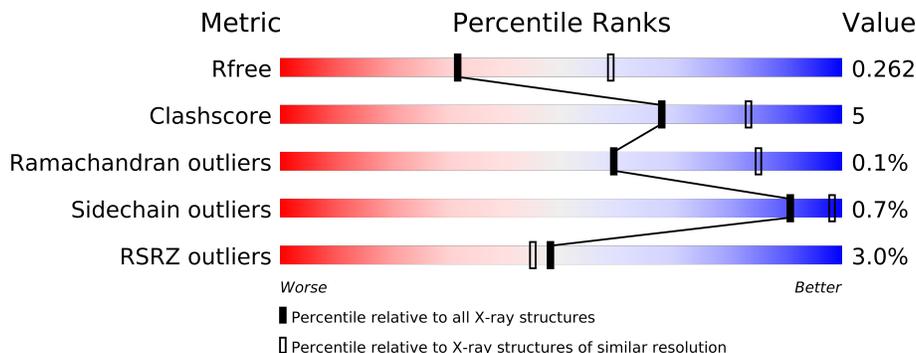
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	1718 (2.60-2.60)
Clashscore	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)
RSRZ outliers	66119	1718 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	341	
1	B	341	
1	C	341	
1	D	341	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	PEG	A	342	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9588 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

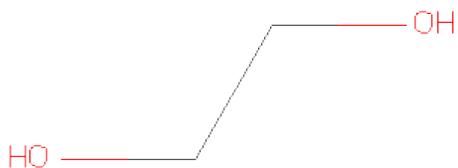
- Molecule 1 is a protein called Prephenate dehydrogenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	323	Total 2409	C 1506	N 432	O 463	S 2	Se 6	0	3	0
1	B	319	Total 2350	C 1474	N 418	O 450	S 2	Se 6	0	3	0
1	C	318	Total 2354	C 1477	N 424	O 445	S 2	Se 6	0	1	0
1	D	321	Total 2372	C 1485	N 425	O 454	S 2	Se 6	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

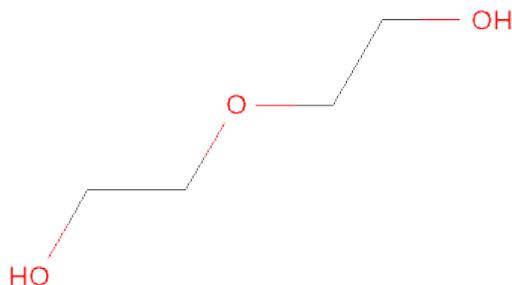
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	leader sequence	UNP Q8NTS6
B	0	GLY	-	leader sequence	UNP Q8NTS6
C	0	GLY	-	leader sequence	UNP Q8NTS6
D	0	GLY	-	leader sequence	UNP Q8NTS6

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

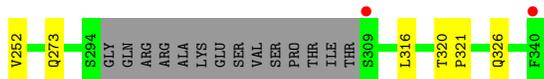
- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



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

- Molecule 4 is water.

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



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	99.32Å 107.23Å 146.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.53 – 2.60 29.53 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.53-2.60) 100.0 (29.53-2.60)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.16 (at 2.61Å)	Xtrriage
Refinement program	REFMAC 5.5.0053	Depositor
R, R_{free}	0.213 , 0.254 0.222 , 0.262	Depositor DCC
R_{free} test set	2475 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	48.5	Xtrriage
Anisotropy	0.502	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 27.5	EDS
Estimated twinning fraction	No twinning to report.	Xtrriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Outliers	0 of 48915 reflections	Xtrriage
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

Xtrriage'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.*

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: PEG, 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	0.42	0/2457	0.61	0/3334
1	B	0.38	0/2399	0.56	0/3266
1	C	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 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2409	0	2354	21	1
1	B	2350	0	2263	29	0
1	C	2354	0	2292	33	1
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	B	19	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	22	0	0	0	0
4	D	21	0	0	1	0
All	All	9588	0	9242	91	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 5.

All (91) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(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:225:SER:HB3	1:D:232:VAL:HG11	1.92	0.51
1:C:149:THR:HA	1:C:184:SER:O	2.10	0.51
1:B:36:TYR:CE2	1:B:57:LEU:HD13	2.45	0.51
1:C:159:ILE:O	1:C:159:ILE:HD12	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:316:LEU:HD21	1:B:326:GLN:HB3	1.94	0.50
1:B:77:MSE:HE2	1:B:105:ALA:HB1	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
1:C:22:LEU:HD12	1:C:177:VAL:HG11	1.93	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:248:ALA:O	1:B:252:VAL:HG23	2.15	0.46
1:B:10:VAL:HG21	1:B:26:LEU:HD13	1.97	0.46
1:D:198:HIS:HB2	1:D:251:LEU:HD13	1.97	0.46
1:B:159:ILE:C	1:B:159:ILE:HD12	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:C:232:VAL:HG11	1:D:225:SER:CB	2.47	0.44
1:A:6:ILE:HB	1:A:31[B]:HIS:CE1	2.53	0.44
1:B:317[B]:HIS:HD2	1:B:340:PHE:HB2	1.79	0.44
1:A:96:PHE:CZ	1:A:115:MSE:HE2	2.53	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:B:313:VAL:C	1:B:314:LEU:HD12	2.38	0.43
1:B:13:LEU:HD21	1:B:61:LEU:HD21	1.99	0.43
1:A:339:VAL:HG23	1:A:339:VAL:O	2.18	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
1:D:149:THR:HA	1:D:184:SER:O	2.20	0.42
1:C:152:GLN:HB3	1:C:159:ILE:HG21	2.02	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:225:SER:HB2	1:B:232:VAL:HG11	2.01	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:C:95:GLY:CA	1:C:166:ILE:HD13	2.51	0.40
1:B:7:SER:OG	1:B:169:ASP:OD2	2.31	0.40
1:B:317[A]:HIS:HD2	1:B:320:THR:HG23	1.86	0.40
1:A:48:ASP:HB3	1:C:328:ILE:HD13	2.03	0.40

All (1) 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	Distance(Å)	Clash(Å)
1:A:281[B]:ASP:OD2	1:C:236[B]:ASP:OD1[2_574]	1.82	0.38

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	319/341 (94%)	300 (94%)	18 (6%)	1 (0%)	50 77
1	B	318/341 (93%)	302 (95%)	16 (5%)	0	100 100
1	C	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%)	59 85

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	Percentiles	
1	A	247/260 (95%)	245 (99%)	2 (1%)	89	97
1	B	236/260 (91%)	234 (99%)	2 (1%)	89	97
1	C	235/260 (90%)	233 (99%)	2 (1%)	87	97
1	D	241/260 (93%)	240 (100%)	1 (0%)	95	99
All	All	959/1040 (92%)	952 (99%)	7 (1%)	91	98

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

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

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

Mol	Chain	Res	Type
1	B	273	GLN

5.3.3 RNA [i](#)

There are no RNA chains 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

There are no carbohydrates in this entry.

5.6 Ligand geometry

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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	EDO	A	341	-	3,3,3	0.62	0	2,2,2	0.23	0
3	PEG	A	342	-	6,6,6	0.48	0	5,5,5	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
2	EDO	A	341	-	-	0/1/1/1	0/0/0/0
3	PEG	A	342	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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

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	323/341 (94%)	-0.16	6 (1%) 64 61	28, 37, 50, 68	0
1	B	319/341 (93%)	-0.09	8 (2%) 54 52	29, 38, 47, 58	0
1	C	318/341 (93%)	0.19	21 (6%) 18 15	27, 40, 49, 67	0
1	D	321/341 (94%)	-0.18	3 (0%) 81 82	28, 38, 48, 65	0
All	All	1281/1364 (93%)	-0.06	38 (2%) 48 45	27, 38, 49, 68	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	156	GLY	5.1
1	A	296	GLN	4.0
1	C	66	ALA	4.0
1	B	130	ASN	3.9
1	C	312	PRO	3.9
1	C	29	ALA	3.9
1	C	48	ASP	3.8
1	D	309	SER	3.8
1	A	128	THR	3.7
1	B	340	PHE	3.5
1	C	129	ALA	3.4
1	A	340	PHE	3.3
1	B	310	SER	3.2
1	D	29	ALA	3.0
1	C	134	SER	2.9
1	C	30	ASN	2.9
1	A	295	GLY	2.9
1	C	89	THR	2.9
1	B	158	ASP	2.8
1	D	340	PHE	2.7
1	B	272	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	35	GLY	2.6
1	C	41	SER	2.6
1	C	133	TRP	2.6
1	A	322	ASN	2.5
1	B	32	SER	2.5
1	C	87	VAL	2.4
1	C	58	GLU	2.4
1	C	90	HIS	2.2
1	B	62	GLN	2.2
1	C	130	ASN	2.1
1	C	59	ALA	2.1
1	C	148	VAL	2.1
1	A	294	SER	2.1
1	C	273	GLN	2.1
1	B	156	GLY	2.1
1	C	86	ALA	2.0
1	C	46	ALA	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 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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	PEG	A	342	7/7	0.22	3.88	80,80,83,84	0
2	EDO	A	341	4/4	0.21	1.15	76,76,76,76	0

6.5 Other polymers

There are no such residues in this entry.