



wwPDB X-ray Structure Validation Summary Report ⓘ

May 25, 2020 – 11:24 pm BST

PDB ID : 3NRE
Title : Crystal structure of a Putative aldose 1-epimerase (b2544) from ES-CHERICHIA COLI K12 at 1.59 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2010-06-30
Resolution : 1.59 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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	:	2.11
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)	:	2.11

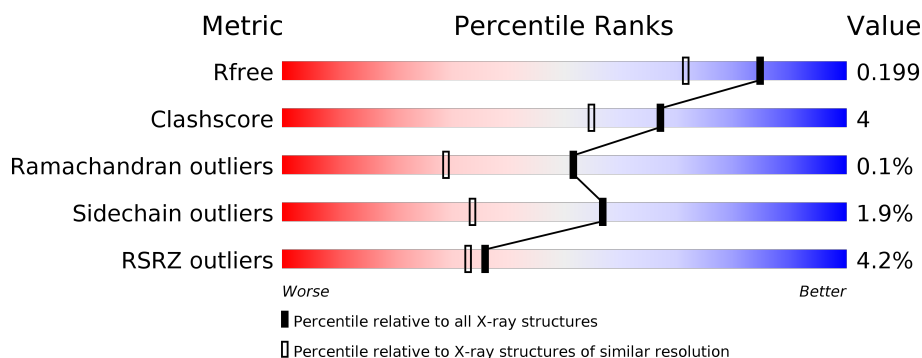
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.59 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	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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 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 $\leq 5\%$. 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	291	<div> <div>4%</div> <div> <div></div> <div>92%</div> <div>8%</div> </div> </div>
1	B	291	<div> <div>4%</div> <div> <div></div> <div>94%</div> <div>6%</div> </div> </div>
1	C	291	<div> <div>4%</div> <div> <div></div> <div>90%</div> <div>10%</div> </div> </div>
1	D	291	<div> <div>4%</div> <div> <div></div> <div>92%</div> <div>7%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10985 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.

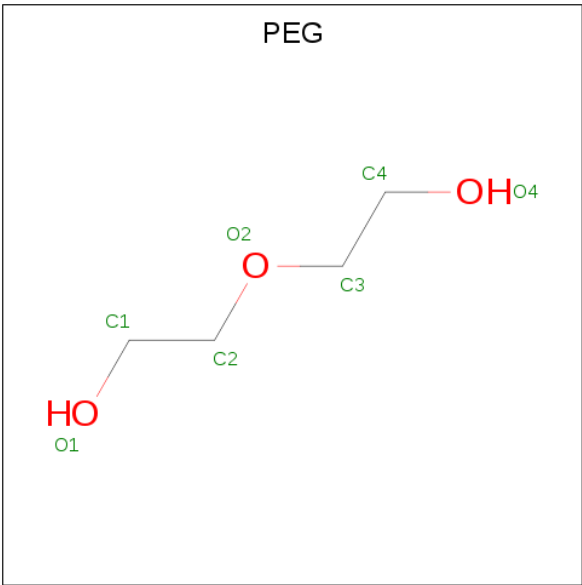
- Molecule 1 is a protein called Aldose 1-epimerase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	291	Total	C	N	O	S	Se	0	22	0
			2409	1549	404	443	7	6			
1	B	291	Total	C	N	O	S	Se	0	23	0
			2415	1558	402	444	7	4			
1	C	291	Total	C	N	O	S	Se	0	31	0
			2454	1600	403	439	7	5			
1	D	291	Total	C	N	O	S	Se	0	26	0
			2446	1575	406	454	7	4			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	leader sequence	UNP C9QPM9
B	0	GLY	-	leader sequence	UNP C9QPM9
C	0	GLY	-	leader sequence	UNP C9QPM9
D	0	GLY	-	leader sequence	UNP C9QPM9

- Molecule 2 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



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

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

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total	Mg	0	0
			1	1		

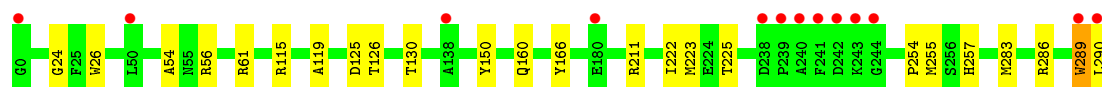
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	237	Total	O	0	13
			248	248		
4	B	259	Total	O	0	11
			268	268		
4	C	319	Total	O	0	12
			330	330		
4	D	279	Total	O	0	9
			288	288		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

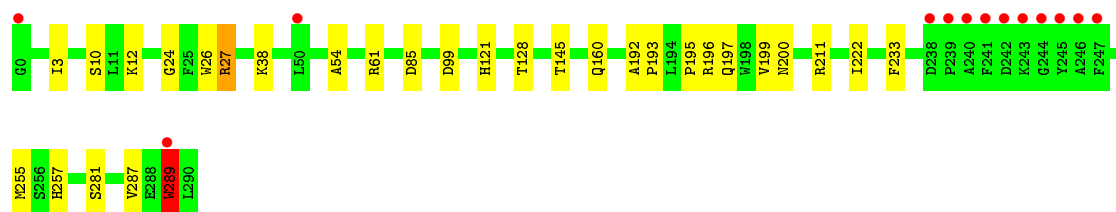
- Molecule 1: Aldose 1-epimerase



- Molecule 1: Aldose 1-epimerase



- Molecule 1: Aldose 1-epimerase



- Molecule 1: Aldose 1-epimerase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	73.73 Å 72.00 Å 272.79 Å 90.00° 90.70° 90.00°	Depositor
Resolution (Å)	28.54 – 1.59 28.54 – 1.59	Depositor EDS
% Data completeness (in resolution range)	98.7 (28.54-1.59) 98.7 (28.54-1.59)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 1.59 Å)	Xtriage
Refinement program	REFMAC 5.5.0110, PHENIX	Depositor
R, R_{free}	0.161 , 0.190 0.172 , 0.199	Depositor DCC
R_{free} test set	9534 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	18.2	Xtriage
Anisotropy	0.559	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.033 for k,h,-l 0.030 for -k,-h,-l 0.034 for -h,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10985	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 4.34% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² 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: MG, 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	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.81	0/2549	0.80	0/3472
1	B	0.81	0/2559	0.81	1/3488 (0.0%)
1	C	0.96	1/2623 (0.0%)	0.91	6/3571 (0.2%)
1	D	0.84	0/2595	0.82	4/3535 (0.1%)
All	All	0.86	1/10326 (0.0%)	0.84	11/14066 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	289	TRP	CB-CG	5.62	1.60	1.50

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	99	ASP	CB-CG-OD1	8.98	126.38	118.30
1	D	17	ASP	CB-CG-OD1	7.28	124.85	118.30
1	B	27	ARG	NE-CZ-NH1	6.63	123.61	120.30
1	C	196[A]	ARG	NE-CZ-NH2	-6.04	117.28	120.30
1	C	196[B]	ARG	NE-CZ-NH2	-6.04	117.28	120.30

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	2285	23	1
1	B	2415	0	2307	17	0
1	C	2454	0	2376	25	0
1	D	2446	0	2321	10	1
2	A	42	0	60	4	0
2	B	28	0	40	4	0
2	C	21	0	30	1	0
2	D	35	0	50	2	0
3	D	1	0	0	0	0
4	A	248	0	0	1	0
4	B	268	0	0	3	0
4	C	330	0	0	4	0
4	D	288	0	0	3	0
All	All	10985	0	9469	75	1

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 4.

The worst 5 of 75 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:C:255[B]:MSE:HE1	4:C:410:HOH:O	1.61	1.00
1:B:287[B]:VAL:HG22	2:B:304:PEG:H22	1.52	0.90
1:C:211[B]:ARG:NH1	1:C:222[B]:ILE:HD11	1.91	0.85
1:A:223[B]:MSE:HE1	1:A:254:PRO:HG3	1.65	0.79
1:D:106[B]:GLU:OE1	1:D:115[B]:ARG:NH1	2.18	0.75

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	Interatomic distance (Å)	Clash overlap (Å)
1:A:115[B]:ARG:NH2	1:D:215:PRO:O[3_455]	2.15	0.05

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	310/291 (106%)	299 (96%)	11 (4%)	0	100	100
1	B	311/291 (107%)	298 (96%)	12 (4%)	1 (0%)	41	21
1	C	319/291 (110%)	311 (98%)	8 (2%)	0	100	100
1	D	314/291 (108%)	306 (98%)	8 (2%)	0	100	100
All	All	1254/1164 (108%)	1214 (97%)	39 (3%)	1 (0%)	51	29

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	239	PRO

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	258/241 (107%)	254 (98%)	4 (2%)	62	41
1	B	262/241 (109%)	257 (98%)	5 (2%)	57	34
1	C	265/241 (110%)	260 (98%)	5 (2%)	57	34
1	D	265/241 (110%)	259 (98%)	6 (2%)	50	25
All	All	1050/964 (109%)	1030 (98%)	20 (2%)	57	34

5 of 20 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	61	ARG
1	C	85	ASP
1	D	217	GLU
1	B	196[B]	ARG
1	B	241	PHE

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

sidechains are listed below:

Mol	Chain	Res	Type
1	A	136	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 1 is monoatomic - leaving 18 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	$\# Z > 2$
2	PEG	D	314	-	6,6,6	0.66	0	5,5,5	0.44	0
2	PEG	D	306	-	6,6,6	0.90	0	5,5,5	0.99	0
2	PEG	B	303	-	6,6,6	0.65	0	5,5,5	0.96	0
2	PEG	A	309	-	6,6,6	0.52	0	5,5,5	0.62	0
2	PEG	A	316	-	6,6,6	0.97	0	5,5,5	1.06	0
2	PEG	C	301	-	6,6,6	0.48	0	5,5,5	0.71	0
2	PEG	A	308	-	6,6,6	0.61	0	5,5,5	0.50	0
2	PEG	B	304	-	6,6,6	0.66	0	5,5,5	0.62	0
2	PEG	D	313	-	6,6,6	0.98	0	5,5,5	0.92	0
2	PEG	B	312	-	6,6,6	0.58	0	5,5,5	0.58	0
2	PEG	A	318	-	6,6,6	0.70	0	5,5,5	0.49	0
2	PEG	C	311	-	6,6,6	0.44	0	5,5,5	0.67	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PEG	D	317	-	6,6,6	0.70	0	5,5,5	0.57	0
2	PEG	A	307	-	6,6,6	0.70	0	5,5,5	0.55	0
2	PEG	A	310	-	6,6,6	0.95	0	5,5,5	1.07	1 (20%)
2	PEG	B	305	-	6,6,6	0.39	0	5,5,5	0.39	0
2	PEG	D	302	-	6,6,6	0.52	0	5,5,5	0.44	0
2	PEG	C	315	-	6,6,6	0.75	0	5,5,5	0.12	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	PEG	D	314	-	-	2/4/4/4	-
2	PEG	D	306	-	-	2/4/4/4	-
2	PEG	B	303	-	-	1/4/4/4	-
2	PEG	A	309	-	-	1/4/4/4	-
2	PEG	A	316	-	-	2/4/4/4	-
2	PEG	C	301	-	-	3/4/4/4	-
2	PEG	A	308	-	-	3/4/4/4	-
2	PEG	B	304	-	-	3/4/4/4	-
2	PEG	D	313	-	-	4/4/4/4	-
2	PEG	B	312	-	-	0/4/4/4	-
2	PEG	A	318	-	-	3/4/4/4	-
2	PEG	C	311	-	-	3/4/4/4	-
2	PEG	D	317	-	-	0/4/4/4	-
2	PEG	A	307	-	-	1/4/4/4	-
2	PEG	A	310	-	-	4/4/4/4	-
2	PEG	B	305	-	-	1/4/4/4	-
2	PEG	D	302	-	-	1/4/4/4	-
2	PEG	C	315	-	-	0/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	310	PEG	O2-C3-C4	2.23	119.89	110.07

There are no chirality outliers.

5 of 34 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	308	PEG	O2-C3-C4-O4
2	D	313	PEG	O1-C1-C2-O2
2	A	307	PEG	O2-C3-C4-O4
2	A	310	PEG	O2-C3-C4-O4
2	D	306	PEG	O2-C3-C4-O4

There are no ring outliers.

6 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	303	PEG	1	0
2	A	316	PEG	1	0
2	B	304	PEG	3	0
2	D	313	PEG	2	0
2	C	311	PEG	1	0
2	A	310	PEG	3	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, 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	287/291 (98%)	0.28	13 (4%) 33 30	17, 29, 47, 78	0
1	B	287/291 (98%)	0.08	11 (3%) 40 37	15, 26, 47, 67	0
1	C	287/291 (98%)	0.05	13 (4%) 33 30	12, 19, 33, 57	0
1	D	287/291 (98%)	0.06	11 (3%) 40 37	16, 25, 43, 58	0
All	All	1148/1164 (98%)	0.12	48 (4%) 36 33	12, 24, 44, 78	0

The worst 5 of 48 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	241	PHE	11.1
1	B	241	PHE	9.3
1	C	240	ALA	8.8
1	D	241	PHE	8.7
1	A	290	LEU	8.6

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. 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	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	PEG	D	313	7/7	0.73	0.16	31,34,46,48	0
2	PEG	A	318	7/7	0.79	0.28	34,43,47,48	0
2	PEG	A	307	7/7	0.84	0.19	36,40,46,47	0
2	PEG	C	301	7/7	0.85	0.25	36,41,45,51	0
2	PEG	A	310	7/7	0.85	0.13	28,34,44,47	0
2	PEG	D	314	7/7	0.86	0.20	37,40,42,44	0
2	PEG	C	311	7/7	0.86	0.16	34,38,48,48	0
2	PEG	D	302	7/7	0.86	0.20	35,38,46,49	0
2	PEG	D	306	7/7	0.87	0.13	36,41,52,53	0
2	PEG	B	304	7/7	0.87	0.25	34,40,46,51	0
2	PEG	D	317	7/7	0.88	0.24	30,38,47,49	0
2	PEG	A	316	7/7	0.88	0.12	30,37,42,43	0
2	PEG	A	308	7/7	0.89	0.27	39,44,45,46	0
2	PEG	B	305	7/7	0.89	0.27	37,39,45,48	0
2	PEG	B	312	7/7	0.89	0.12	33,34,41,42	0
2	PEG	C	315	7/7	0.89	0.23	29,33,45,45	0
2	PEG	B	303	7/7	0.90	0.14	34,37,45,47	0
2	PEG	A	309	7/7	0.91	0.17	39,41,43,45	0
3	MG	D	300	1/1	0.94	0.25	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.