



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 02:06 pm BST

PDB ID : 4WYR
Title : Crystal structure of thiolase mutation (V77Q,N153Y,A286K) from *Clostridium acetobutylicum*
Authors : Kim, S.; Ha, S.C.; Ahn, J.W.; Kim, E.J.; Lim, J.H.; Kim, K.J.
Deposited on : 2014-11-18
Resolution : 2.30 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 ⓘ 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

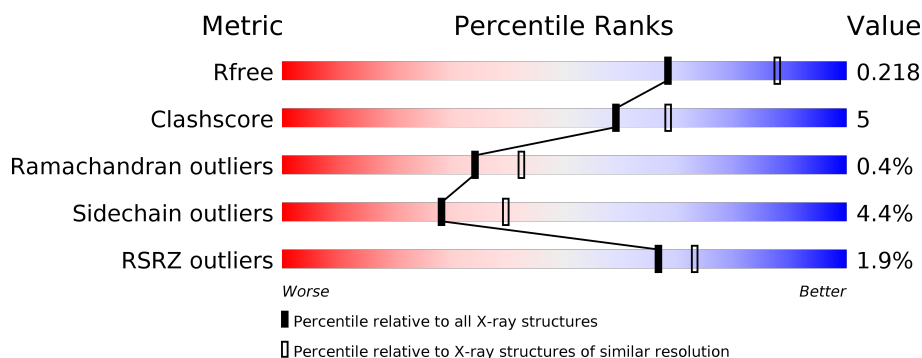
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	400	<div> <div>2%</div> <div> <div></div> <div>85%</div> <div>12%</div> <div>••</div> </div> </div>
1	B	400	<div> <div>2%</div> <div> <div></div> <div>85%</div> <div>13%</div> <div>•</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6112 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 Acetyl-CoA acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	392	Total	C	N	O	S	0	0	0
			2902	1832	502	553	15			
1	B	392	Total	C	N	O	S	0	0	0
			2902	1832	502	553	15			

There are 22 discrepancies between the modelled and reference sequences:

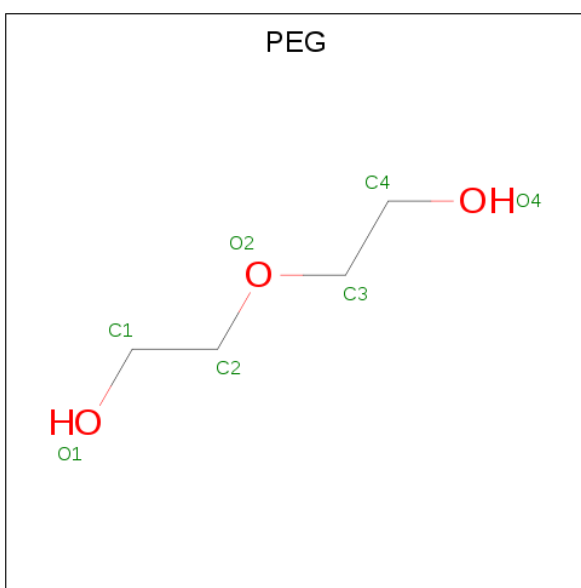
Chain	Residue	Modelled	Actual	Comment	Reference
A	77	GLN	VAL	engineered mutation	UNP P45359
A	153	TYR	ASN	engineered mutation	UNP P45359
A	286	LYS	ALA	engineered mutation	UNP P45359
A	393	LEU	-	expression tag	UNP P45359
A	394	GLU	-	expression tag	UNP P45359
A	395	HIS	-	expression tag	UNP P45359
A	396	HIS	-	expression tag	UNP P45359
A	397	HIS	-	expression tag	UNP P45359
A	398	HIS	-	expression tag	UNP P45359
A	399	HIS	-	expression tag	UNP P45359
A	400	HIS	-	expression tag	UNP P45359
B	77	GLN	VAL	engineered mutation	UNP P45359
B	153	TYR	ASN	engineered mutation	UNP P45359
B	286	LYS	ALA	engineered mutation	UNP P45359
B	393	LEU	-	expression tag	UNP P45359
B	394	GLU	-	expression tag	UNP P45359
B	395	HIS	-	expression tag	UNP P45359
B	396	HIS	-	expression tag	UNP P45359
B	397	HIS	-	expression tag	UNP P45359
B	398	HIS	-	expression tag	UNP P45359
B	399	HIS	-	expression tag	UNP P45359
B	400	HIS	-	expression tag	UNP P45359

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		

- 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	0	0
			7	4	3		
3	A	1	Total	C	O	0	0
			7	4	3		

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

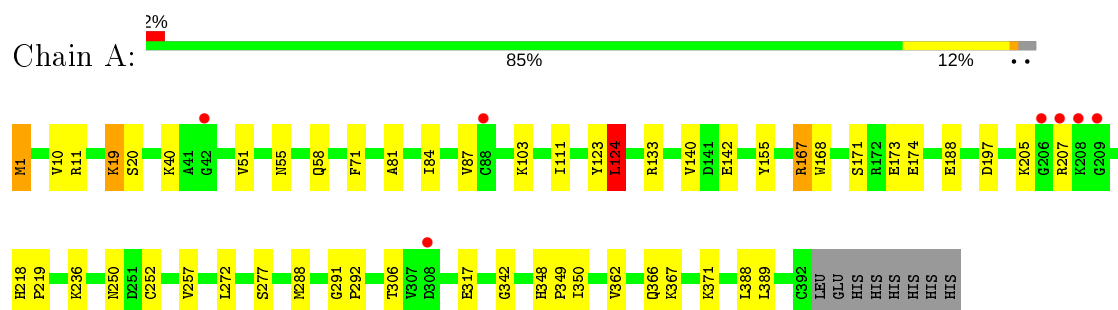
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	142	Total	O	0	0
			142	142		
4	B	119	Total	O	0	0
			119	119		

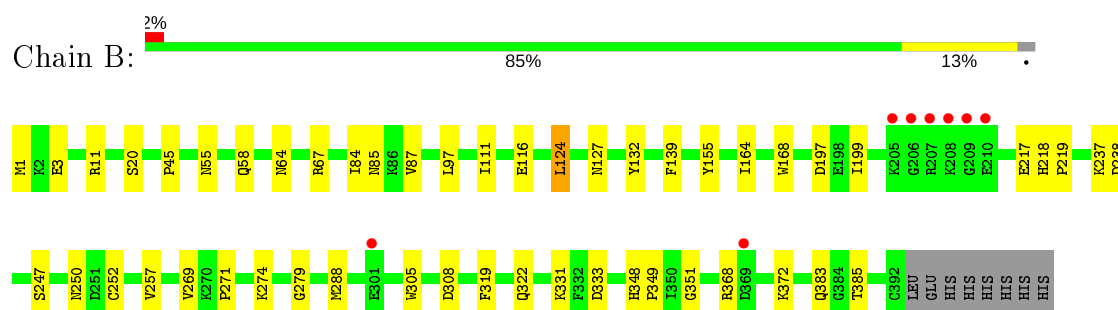
3 Residue-property plots [i](#)

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: Acetyl-CoA acetyltransferase



- Molecule 1: Acetyl-CoA acetyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	116.43 Å 131.20 Å 54.12 Å 90.00° 110.30° 90.00°	Depositor
Resolution (Å)	50.00 – 2.30 35.08 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.7 (50.00-2.30) 98.7 (35.08-2.30)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.01 (at 2.29 Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.151 , 0.213 0.162 , 0.218	Depositor DCC
R_{free} test set	1693 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	28.4	Xtriage
Anisotropy	0.759	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 49.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6112	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.13% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: GOL, 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	0.78	0/2944	0.83	1/3971 (0.0%)
1	B	0.78	0/2944	0.81	3/3971 (0.1%)
All	All	0.78	0/5888	0.82	4/7942 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	124	LEU	CA-CB-CG	6.73	130.78	115.30
1	B	333	ASP	N-CA-C	-5.33	96.59	111.00
1	B	139	PHE	N-CA-C	-5.24	96.86	111.00
1	B	124	LEU	CA-CB-CG	5.12	127.08	115.30

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	2902	0	2973	29	0
1	B	2902	0	2973	31	0
2	A	12	0	16	2	0
3	A	14	0	20	0	0
3	B	21	0	30	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	142	0	0	2	0
4	B	119	0	0	1	0
All	All	6112	0	6012	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:1:MET:CE	1:B:1:MET:HA	1.99	0.93
1:A:250:ASN:HD22	1:A:348:HIS:H	1.36	0.74
1:B:1:MET:HA	1:B:1:MET:HE3	1.71	0.73
1:B:1:MET:HA	1:B:1:MET:HE2	1.75	0.67
1:A:84:ILE:HD12	1:B:84:ILE:HD12	1.77	0.66
1:A:133:ARG:NH2	4:A:601:HOH:O	2.30	0.65
1:B:252:CYS:HB3	1:B:349:PRO:HG3	1.79	0.63
1:B:164:ILE:HG23	1:B:168:TRP:CE3	2.35	0.61
1:A:250:ASN:ND2	1:A:348:HIS:H	1.99	0.61
1:A:252:CYS:HB3	1:A:349:PRO:HG3	1.84	0.59
1:B:111:ILE:CD1	1:B:257:VAL:HG22	2.32	0.59
1:B:250:ASN:HD22	1:B:348:HIS:H	1.52	0.57
1:B:111:ILE:HD13	1:B:257:VAL:HG22	1.86	0.57
1:A:58:GLN:NE2	1:B:85:ASN:HD22	2.04	0.56
1:B:305:TRP:CE2	1:B:372:LYS:HE2	2.42	0.55
1:A:1:MET:HB2	1:A:103:LYS:HG2	1.87	0.55
1:B:11:ARG:O	1:B:199:ILE:HA	2.07	0.55
1:B:319:PHE:H	1:B:322:GLN:HE21	1.55	0.53
3:B:501:PEG:H11	4:B:653:HOH:O	2.08	0.53
1:B:237:LYS:O	1:B:238:ASP:HB2	2.08	0.52
1:A:218:HIS:N	1:A:219:PRO:CD	2.74	0.51
1:B:64:ASN:HB3	1:B:67:ARG:HB3	1.92	0.51
1:A:111:ILE:CD1	1:A:257:VAL:HG22	2.41	0.51
1:B:250:ASN:ND2	1:B:348:HIS:H	2.09	0.50
1:A:317:GLU:CD	1:A:342:GLY:HA3	2.33	0.49
1:A:291:GLY:N	1:A:292:PRO:CD	2.76	0.48
1:A:58:GLN:HE22	1:B:85:ASN:HD22	1.62	0.47
1:A:197:ASP:OD2	1:A:367:LYS:HE2	2.14	0.47
1:A:167:ARG:NH1	4:A:602:HOH:O	2.31	0.47
1:A:123:TYR:HA	1:A:140:VAL:O	2.15	0.46
1:B:20:SER:OG	1:B:217:GLU:OE2	2.23	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:LEU:HD13	1:A:142:GLU:HB2	1.96	0.46
1:B:55:ASN:HD22	1:B:58:GLN:HG2	1.81	0.45
1:A:81:ALA:H	1:B:383:GLN:HE22	1.65	0.45
1:A:10:VAL:HG12	1:A:11:ARG:N	2.32	0.45
1:B:305:TRP:CZ2	1:B:372:LYS:HE2	2.52	0.45
1:B:3:GLU:HG2	1:B:274:LYS:NZ	2.32	0.45
1:B:247:SER:O	3:B:502:PEG:H41	2.17	0.44
1:A:19:LYS:HB3	1:A:20:SER:H	1.52	0.44
1:A:252:CYS:HB3	1:A:349:PRO:CG	2.48	0.43
1:B:218:HIS:N	1:B:219:PRO:CD	2.81	0.43
1:B:1:MET:CA	1:B:1:MET:HE3	2.43	0.42
1:A:84:ILE:HD12	1:B:84:ILE:CD1	2.46	0.42
1:A:171:SER:OG	1:A:173:GLU:HG2	2.19	0.42
1:B:279:GLY:O	1:B:385:THR:HA	2.19	0.42
1:A:277:SER:OG	1:A:388:LEU:HD12	2.19	0.42
1:B:127:ASN:HB2	1:B:132:TYR:CE2	2.54	0.42
1:A:219:PRO:O	2:A:502:GOL:H11	2.19	0.41
1:B:269:VAL:O	1:B:271:PRO:HD3	2.21	0.41
1:A:51:VAL:O	1:A:81:ALA:HA	2.20	0.41
1:B:55:ASN:ND2	1:B:58:GLN:HG2	2.36	0.41
1:A:219:PRO:O	2:A:502:GOL:C1	2.69	0.40
1:A:362:VAL:HG12	1:A:389:LEU:HD13	2.02	0.40
1:A:55:ASN:HD22	1:A:58:GLN:HG2	1.86	0.40
1:A:167:ARG:HB3	1:A:168:TRP:CE3	2.56	0.40
1:B:116:GLU:OE1	1:B:351:GLY:N	2.53	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	390/400 (98%)	381 (98%)	7 (2%)	2 (0%)	29	35

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	390/400 (98%)	380 (97%)	9 (2%)	1 (0%)	41	50
All	All	780/800 (98%)	761 (98%)	16 (2%)	3 (0%)	34	42

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	87	VAL
1	B	87	VAL
1	A	350	ILE

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	295/303 (97%)	278 (94%)	17 (6%)	20	27
1	B	295/303 (97%)	286 (97%)	9 (3%)	40	55
All	All	590/606 (97%)	564 (96%)	26 (4%)	28	39

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	19	LYS
1	A	40	LYS
1	A	71	PHE
1	A	124	LEU
1	A	155	TYR
1	A	167	ARG
1	A	174	GLU
1	A	188	GLU
1	A	205	LYS
1	A	207	ARG
1	A	236	LYS
1	A	272	LEU

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Mol	Chain	Res	Type
1	A	288	MET
1	A	306	THR
1	A	366	GLN
1	A	371	LYS
1	B	45	PRO
1	B	97	LEU
1	B	124	LEU
1	B	155	TYR
1	B	197	ASP
1	B	288	MET
1	B	308	ASP
1	B	331	LYS
1	B	368	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	ASN
1	A	58	GLN
1	A	250	ASN
1	A	322	GLN
1	B	55	ASN
1	B	58	GLN
1	B	250	ASN
1	B	322	GLN
1	B	383	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

7 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	$\# Z > 2$
3	PEG	A	504	-	6,6,6	0.69	0	5,5,5	1.09	1 (20%)
3	PEG	A	503	-	6,6,6	0.32	0	5,5,5	0.28	0
2	GOL	A	502	-	5,5,5	0.64	0	5,5,5	0.81	0
3	PEG	B	501	-	6,6,6	0.30	0	5,5,5	0.29	0
2	GOL	A	501	-	5,5,5	0.18	0	5,5,5	0.33	0
3	PEG	B	502	-	6,6,6	0.31	0	5,5,5	0.40	0
3	PEG	B	503	-	6,6,6	0.44	0	5,5,5	0.64	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	504	-	-	3/4/4/4	-
3	PEG	A	503	-	-	3/4/4/4	-
2	GOL	A	502	-	-	2/4/4/4	-
3	PEG	B	501	-	-	3/4/4/4	-
2	GOL	A	501	-	-	2/4/4/4	-
3	PEG	B	502	-	-	2/4/4/4	-
3	PEG	B	503	-	-	2/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	504	PEG	O2-C2-C1	2.18	119.64	110.07

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	502	GOL	C1-C2-C3-O3
2	A	501	GOL	O1-C1-C2-C3
3	B	501	PEG	O2-C3-C4-O4
3	B	503	PEG	O2-C3-C4-O4
3	A	504	PEG	O2-C3-C4-O4
3	A	503	PEG	O1-C1-C2-O2
2	A	502	GOL	O2-C2-C3-O3
2	A	501	GOL	O1-C1-C2-O2
3	B	501	PEG	O1-C1-C2-O2
3	B	502	PEG	C4-C3-O2-C2
3	B	503	PEG	C4-C3-O2-C2
3	A	504	PEG	C1-C2-O2-C3
3	A	503	PEG	C4-C3-O2-C2
3	A	504	PEG	C4-C3-O2-C2
3	B	502	PEG	C1-C2-O2-C3
3	B	501	PEG	C1-C2-O2-C3
3	A	503	PEG	C1-C2-O2-C3

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	502	GOL	2	0
3	B	501	PEG	1	0
3	B	502	PEG	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, 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	392/400 (98%)	-0.13	7 (1%) 68 74	17, 29, 50, 111	0
1	B	392/400 (98%)	-0.11	8 (2%) 65 71	18, 29, 49, 104	0
All	All	784/800 (98%)	-0.12	15 (1%) 66 73	17, 29, 50, 111	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	207	ARG	6.8
1	A	208	LYS	6.0
1	B	206	GLY	5.9
1	B	207	ARG	5.7
1	B	209	GLY	5.4
1	B	208	LYS	5.3
1	A	209	GLY	3.7
1	B	210	GLU	3.5
1	B	205	LYS	2.7
1	B	369	ASP	2.2
1	B	301	GLU	2.2
1	A	88	CYS	2.1
1	A	308	ASP	2.1
1	A	206	GLY	2.1
1	A	42	GLY	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. 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(Å ²)	Q<0.9
2	GOL	A	501	6/6	0.63	0.24	62,66,69,81	0
2	GOL	A	502	6/6	0.83	0.19	53,55,56,62	0
3	PEG	A	503	7/7	0.84	0.28	52,54,56,57	0
3	PEG	A	504	7/7	0.85	0.24	43,52,58,61	0
3	PEG	B	503	7/7	0.85	0.17	54,55,57,58	0
3	PEG	B	502	7/7	0.88	0.21	46,50,60,65	0
3	PEG	B	501	7/7	0.88	0.17	43,46,49,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.