



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:11 AM UTC

PDB ID : 8RP4 / pdb_00008rp4
Title : Alpha-Methylacyl-CoA racemase from Mycobacterium tuberculosis (D156A mutant)
Authors : Mojanaga, O.O.; Acharya, K.R.; Lloyd, M.D.
Deposited on : 2024-01-12
Resolution : 1.64 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

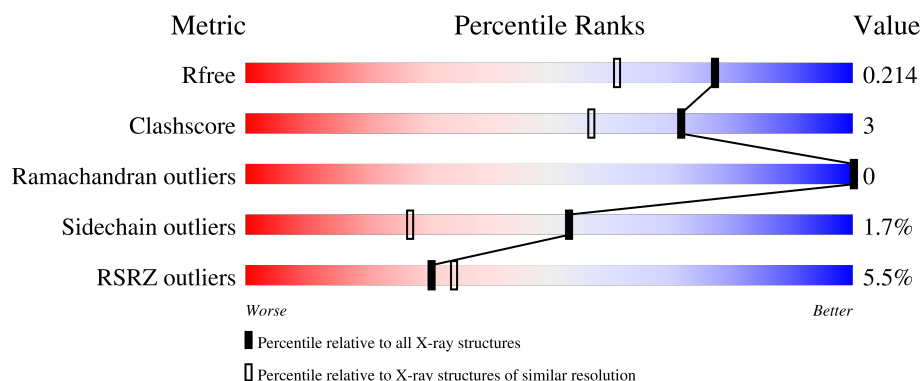
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.64 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}	180053	1141 (1.64-1.64)
Clashscore	190562	1171 (1.64-1.64)
Ramachandran outliers	187476	1151 (1.64-1.64)
Sidechain outliers	187428	1150 (1.64-1.64)
RSRZ outliers	180081	1141 (1.64-1.64)

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 $\leq 5\%$. 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	365	5% 88% 8% ..
1	B	365	5% 88% 8% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11186 atoms, of which 5340 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 Alpha-methylacyl-CoA racemase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	354	Total	C	H	N	O	S	56	4	0
			5337	1690	2645	482	504	16			
1	B	354	Total	C	H	N	O	S	56	3	0
			5329	1688	2641	481	503	16			

There are 12 discrepancies between the modelled and reference sequences:

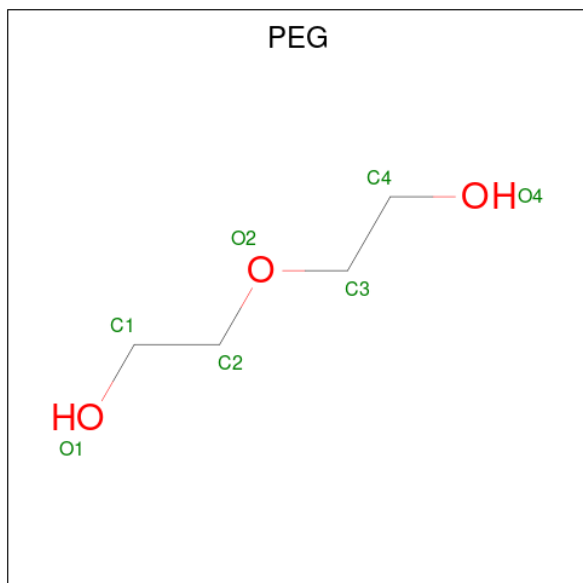
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP O06543
A	156	ALA	ASP	engineered mutation	UNP O06543
A	361	GLY	-	expression tag	UNP O06543
A	362	SER	-	expression tag	UNP O06543
A	363	GLY	-	expression tag	UNP O06543
A	364	CYS	-	expression tag	UNP O06543
B	0	MET	-	initiating methionine	UNP O06543
B	156	ALA	ASP	engineered mutation	UNP O06543
B	361	GLY	-	expression tag	UNP O06543
B	362	SER	-	expression tag	UNP O06543
B	363	GLY	-	expression tag	UNP O06543
B	364	CYS	-	expression tag	UNP O06543

- Molecule 2 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	1	0
			10	2	6	2		
2	A	1	Total	C	H	O	1	0
			10	2	6	2		
2	B	1	Total	C	H	O	1	0
			10	2	6	2		
2	B	1	Total	C	H	O	1	0
			10	2	6	2		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C H O 17 4 10 3	1	0
3	A	1	Total C H O 17 4 10 3	1	0
3	B	1	Total C H O 17 4 10 3	1	0

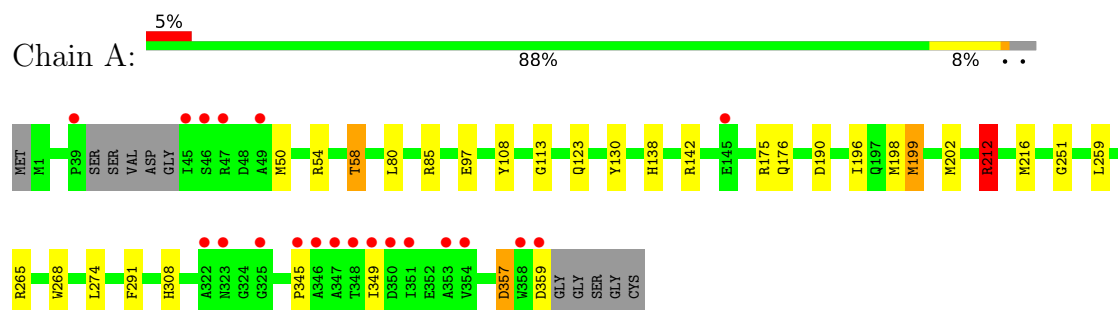
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	207	Total O 207 207	0	0
4	B	222	Total O 222 222	0	0

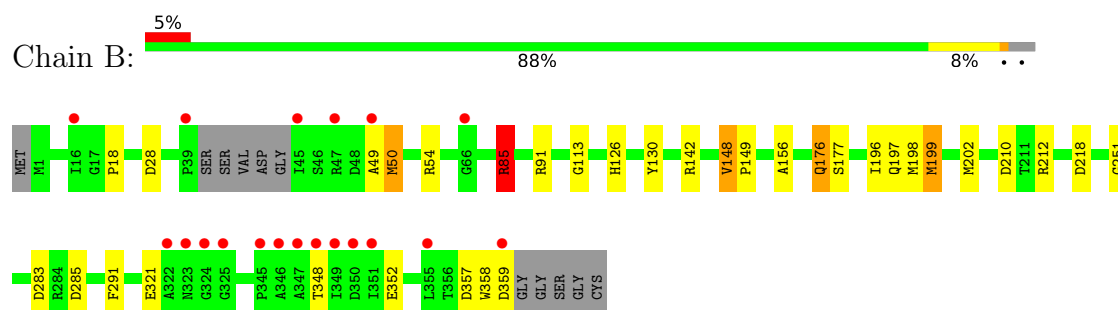
3 Residue-property plots [i](#)

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: Alpha-methylacyl-CoA racemase



- Molecule 1: Alpha-methylacyl-CoA racemase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	180.47Å 78.75Å 58.94Å 90.00° 92.15° 90.00°	Depositor
Resolution (Å)	90.17 – 1.64 90.17 – 1.64	Depositor EDS
% Data completeness (in resolution range)	98.8 (90.17-1.64) 98.7 (90.17-1.64)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.16 (at 1.64Å)	Xtriage
Refinement program	REFMAC 5.8.0419, REFMAC 5.8.0419	Depositor
R, R_{free}	0.179 , 0.211 0.180 , 0.214	Depositor DCC
R_{free} test set	5055 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	23.3	Xtriage
Anisotropy	0.151	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 33.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.116 for -h,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	11186	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.30% 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

5.1 Standard geometry

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	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.99	0/2772	1.34	8/3770 (0.2%)
1	B	1.01	2/2764 (0.1%)	1.38	14/3759 (0.4%)
All	All	1.00	2/5536 (0.0%)	1.36	22/7529 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4
1	B	0	1
All	All	0	5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	212	ARG	NE-CZ	-6.18	1.26	1.33
1	B	126	HIS	ND1-CE1	5.22	1.37	1.32

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	54	ARG	CB-CA-C	-8.54	91.29	110.07
1	B	28	ASP	CA-CB-CG	7.58	120.18	112.60
1	B	148	VAL	CA-C-O	7.11	124.21	119.38
1	B	54	ARG	CD-NE-CZ	7.04	134.25	124.40
1	A	216	MET	CG-SD-CE	6.49	115.18	100.90
1	B	142	ARG	NE-CZ-NH1	-6.42	115.08	121.50
1	A	212	ARG	CD-NE-CZ	-6.26	115.63	124.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	50	MET	N-CA-C	-6.16	104.67	111.82
1	B	283[A]	ASP	CA-CB-CG	5.99	118.59	112.60
1	B	283[B]	ASP	CA-CB-CG	5.99	118.59	112.60
1	A	58	THR	CA-CB-OG1	-5.75	100.98	109.60
1	B	50	MET	CA-C-N	-5.59	113.55	122.66
1	B	50	MET	C-N-CA	-5.59	113.55	122.66
1	B	85	ARG	NE-CZ-NH2	5.53	124.18	119.20
1	A	190	ASP	CA-CB-CG	5.53	118.13	112.60
1	A	97	GLU	CB-CG-CD	5.46	121.88	112.60
1	B	357	ASP	CA-CB-CG	5.34	117.94	112.60
1	A	175	ARG	NE-CZ-NH1	-5.33	116.17	121.50
1	A	357	ASP	CA-CB-CG	5.28	117.88	112.60
1	B	285	ASP	CB-CA-C	5.26	120.36	110.63
1	B	85	ARG	CD-NE-CZ	5.25	131.76	124.40
1	A	308	HIS	CA-CB-CG	-5.20	108.60	113.80

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	142	ARG	Sidechain
1	A	212	ARG	Sidechain
1	A	54	ARG	Sidechain
1	A	85	ARG	Sidechain
1	B	85	ARG	Sidechain

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	2692	2645	2634	21	0
1	B	2688	2641	2632	17	1
2	A	8	12	11	2	0
2	B	8	12	12	0	0
3	A	14	20	19	1	1
3	B	7	10	10	0	0
4	A	207	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	222	0	0	5	0
All	All	5846	5340	5318	35	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 3.

All (35) 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:A:199:MET:HE3	1:A:202:MET:CE	2.22	0.69
1:A:176:GLN:CD	1:B:176:GLN:HG2	2.24	0.62
1:A:123:GLN:NE2	4:A:505:HOH:O	2.37	0.57
1:A:198:MET:HG2	1:A:202:MET:HE2	1.87	0.57
1:A:199:MET:HE3	1:A:202:MET:HE1	1.87	0.56
1:A:259:LEU:HD22	1:A:274:LEU:HD13	1.91	0.53
1:B:91:ARG:NH2	4:B:505:HOH:O	2.43	0.50
1:A:357:ASP:O	1:A:359:ASP:N	2.45	0.49
1:B:113:GLY:HA3	1:B:130:TYR:CZ	2.48	0.49
1:B:352:GLU:HG2	4:B:674:HOH:O	2.13	0.48
1:B:358:TRP:O	1:B:359:ASP:HB2	2.13	0.48
1:A:138:HIS:O	1:A:212:ARG:HD3	2.14	0.48
1:A:212:ARG:NH2	1:A:212:ARG:CB	2.76	0.48
1:B:49:ALA:HB2	4:B:633:HOH:O	2.13	0.47
1:B:198:MET:HG2	1:B:202:MET:HE2	1.96	0.47
1:B:148:VAL:CG2	1:B:149:PRO:HD2	2.44	0.47
1:B:251:GLY:HA3	1:B:291:PHE:CE1	2.49	0.47
1:A:176:GLN:HG3	1:B:176:GLN:NE2	2.31	0.46
1:A:251:GLY:HA3	1:A:291:PHE:CE1	2.51	0.45
1:A:113:GLY:HA3	1:A:130:TYR:CZ	2.52	0.45
1:A:198:MET:HE3	1:B:50:MET:HE1	1.99	0.45
3:A:403:PEG:O1	3:A:403:PEG:H32	2.18	0.43
1:A:80:LEU:HD23	1:A:108:TYR:CE2	2.53	0.43
1:A:138:HIS:HD2	4:A:553:HOH:O	2.00	0.43
1:A:345:PRO:HB3	4:A:630:HOH:O	2.19	0.43
1:B:218:ASP:HB2	4:B:517:HOH:O	2.19	0.43
1:A:196:ILE:HG12	1:A:199:MET:HB2	2.01	0.42
1:A:50:MET:HA	1:B:197:GLN:HE22	1.85	0.41
1:A:265:ARG:HB3	2:A:401:EDO:H22	2.02	0.41
1:B:148:VAL:HA	1:B:149:PRO:HD3	1.95	0.41
1:A:212:ARG:NH2	1:A:212:ARG:HB2	2.36	0.41
1:B:196:ILE:HG12	1:B:199:MET:HB2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:352:GLU:CG	4:B:674:HOH:O	2.68	0.41
1:B:18:PRO:HB3	1:B:156:ALA:O	2.21	0.40
1:A:268:TRP:HB2	2:A:401:EDO:H11	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	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:ASP:H	3:A:404:PEG:H41[4_545]	1.25	0.35

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	354/365 (97%)	342 (97%)	12 (3%)	0	100	100
1	B	353/365 (97%)	343 (97%)	10 (3%)	0	100	100
All	All	707/730 (97%)	685 (97%)	22 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	274/277 (99%)	271 (99%)	3 (1%)	65	44

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	B	273/277 (99%)	267 (98%)	6 (2%)	45 18
All	All	547/554 (99%)	538 (98%)	9 (2%)	53 30

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	THR
1	A	199	MET
1	A	349	ILE
1	B	85	ARG
1	B	176	GLN
1	B	177	SER
1	B	199	MET
1	B	321	GLU
1	B	348	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	123	GLN
1	A	138	HIS
1	B	176	GLN
1	B	282	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides 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	404	-	6,6,6	0.29	0	5,5,5	0.60	0
2	EDO	A	402	-	3,3,3	0.35	0	2,2,2	0.53	0
2	EDO	B	401	-	3,3,3	0.47	0	2,2,2	0.56	0
3	PEG	B	403	-	6,6,6	0.27	0	5,5,5	0.43	0
2	EDO	A	401	-	3,3,3	1.02	0	2,2,2	0.22	0
3	PEG	A	403	-	6,6,6	0.70	0	5,5,5	0.58	0
2	EDO	B	402	-	3,3,3	0.29	0	2,2,2	0.62	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	404	-	-	3/4/4/4	-
2	EDO	A	402	-	-	1/1/1/1	-
2	EDO	B	401	-	-	0/1/1/1	-
3	PEG	B	403	-	-	4/4/4/4	-
2	EDO	A	401	-	-	1/1/1/1	-
3	PEG	A	403	-	-	3/4/4/4	-
2	EDO	B	402	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	403	PEG	C4-C3-O2-C2

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Mol	Chain	Res	Type	Atoms
3	A	404	PEG	O2-C3-C4-O4
3	B	403	PEG	O1-C1-C2-O2
3	B	403	PEG	O2-C3-C4-O4
2	A	401	EDO	O1-C1-C2-O2
3	A	404	PEG	C1-C2-O2-C3
3	A	403	PEG	O1-C1-C2-O2
3	B	403	PEG	C4-C3-O2-C2
3	A	404	PEG	C4-C3-O2-C2
3	B	403	PEG	C1-C2-O2-C3
2	A	402	EDO	O1-C1-C2-O2
3	A	403	PEG	O2-C3-C4-O4

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	404	PEG	0	1
2	A	401	EDO	2	0
3	A	403	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

6.1 Protein, DNA and RNA chains

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	354/365 (96%)	0.07	20 (5%) 30 33	15, 26, 53, 94	4 (1%)
1	B	354/365 (96%)	0.02	19 (5%) 31 35	15, 26, 54, 86	3 (0%)
All	All	708/730 (96%)	0.04	39 (5%) 30 34	15, 26, 53, 94	7 (0%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	324	GLY	5.9
1	A	349	ILE	4.6
1	A	45	ILE	4.3
1	A	359	ASP	4.2
1	B	325	GLY	4.0
1	A	47	ARG	3.9
1	B	39	PRO	3.9
1	A	347	ALA	3.8
1	B	349	ILE	3.7
1	B	45	ILE	3.7
1	A	346	ALA	3.6
1	B	346	ALA	3.4
1	A	39	PRO	3.4
1	B	347	ALA	3.0
1	B	323	ASN	3.0
1	A	350	ASP	2.9
1	A	323	ASN	2.9
1	A	345	PRO	2.8
1	B	359	ASP	2.8
1	B	49	ALA	2.7
1	B	345	PRO	2.7
1	A	351	ILE	2.6
1	B	322	ALA	2.6
1	B	16	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	49	ALA	2.4
1	B	350	ASP	2.3
1	A	358	TRP	2.3
1	B	47	ARG	2.3
1	A	354	VAL	2.2
1	A	322	ALA	2.2
1	B	355	LEU	2.2
1	A	353	ALA	2.2
1	A	325	GLY	2.2
1	A	145	GLU	2.1
1	B	348	THR	2.1
1	B	66	GLY	2.1
1	A	46	SER	2.1
1	A	348	THR	2.0
1	B	351	ILE	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 oligosaccharides 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
3	PEG	A	403	7/7	0.84	0.19	29,52,61,63	1
3	PEG	B	403	7/7	0.86	0.20	29,54,61,62	1
2	EDO	B	402	4/4	0.90	0.11	39,40,43,47	1
2	EDO	A	401	4/4	0.91	0.09	35,36,39,41	1
3	PEG	A	404	7/7	0.92	0.11	26,35,46,47	1
2	EDO	A	402	4/4	0.94	0.10	30,37,46,51	1
2	EDO	B	401	4/4	0.97	0.12	27,33,35,35	1

6.5 Other polymers [i](#)

There are no such residues in this entry.