



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 03:58 AM UTC

PDB ID : 7XZN / pdb_00007xzn
Title : Formate-tetrahydrofolate ligase from *Peptostreptococcus anaerobius*
Authors : Fang, C.L.; Zhang, Y.
Deposited on : 2022-06-03
Resolution : 2.04 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 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

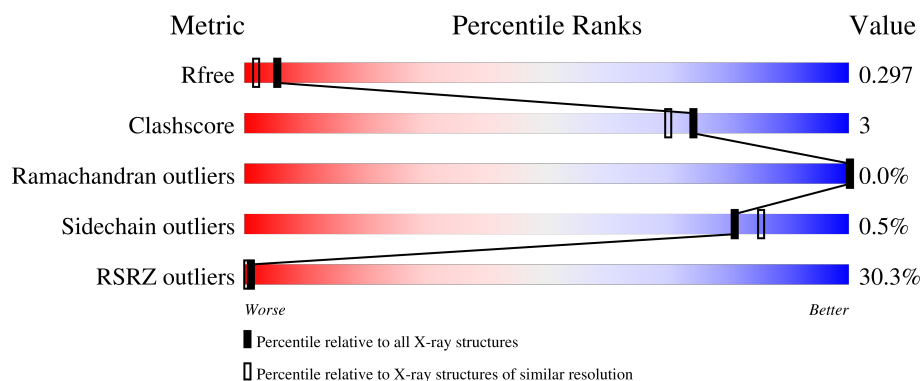
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.04 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	562	30% 91% 7%
1	B	562	34% 92% 7%
1	C	562	34% 90% 9%
1	D	562	23% 93% 6%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	TAM	A	605	-	X	-	-
3	TAM	B	602	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 16831 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 Formate--tetrahydrofolate ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	555	Total	C	N	O	S	0	0	0
			4159	2617	717	804	21			
1	B	556	Total	C	N	O	S	0	0	0
			4162	2619	715	807	21			
1	C	556	Total	C	N	O	S	0	0	0
			4172	2625	719	807	21			
1	D	555	Total	C	N	O	S	0	0	0
			4137	2600	714	802	21			

There are 16 discrepancies between the modelled and reference sequences:

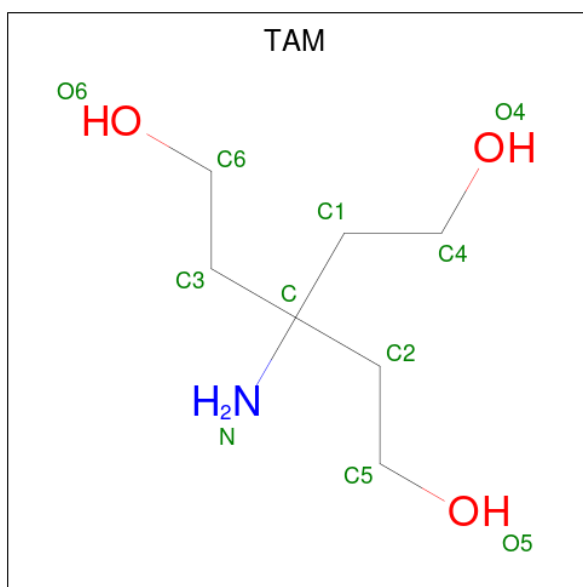
Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP A0A379CIH2
A	-2	SER	-	expression tag	UNP A0A379CIH2
A	-1	HIS	-	expression tag	UNP A0A379CIH2
A	0	MET	-	expression tag	UNP A0A379CIH2
B	-3	GLY	-	expression tag	UNP A0A379CIH2
B	-2	SER	-	expression tag	UNP A0A379CIH2
B	-1	HIS	-	expression tag	UNP A0A379CIH2
B	0	MET	-	expression tag	UNP A0A379CIH2
C	-3	GLY	-	expression tag	UNP A0A379CIH2
C	-2	SER	-	expression tag	UNP A0A379CIH2
C	-1	HIS	-	expression tag	UNP A0A379CIH2
C	0	MET	-	expression tag	UNP A0A379CIH2
D	-3	GLY	-	expression tag	UNP A0A379CIH2
D	-2	SER	-	expression tag	UNP A0A379CIH2
D	-1	HIS	-	expression tag	UNP A0A379CIH2
D	0	MET	-	expression tag	UNP A0A379CIH2

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



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

- Molecule 3 is TRIS(HYDROXYETHYL)AMINOMETHANE (CCD ID: TAM) (formula: $C_7H_{17}NO_3$).



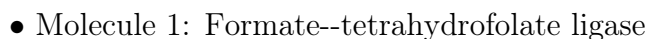
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			11	7	1	3		
3	B	1	Total	C	N	O	0	0
			11	7	1	3		
3	C	1	Total	C	N	O	0	0
			11	7	1	3		
3	D	1	Total	C	N	O	0	0
			11	7	1	3		

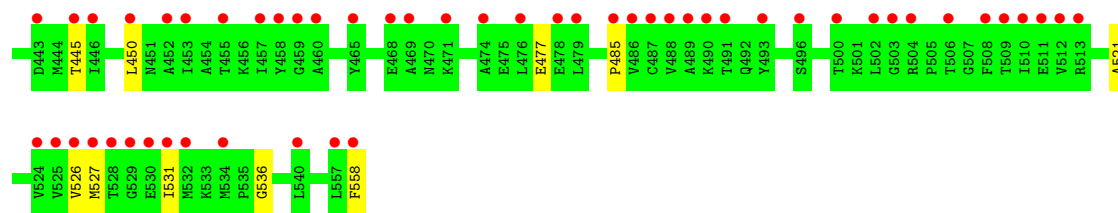
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	28	Total	O	0	0
			28	28		
4	B	36	Total	O	0	0
			36	36		
4	C	28	Total	O	0	0
			28	28		
4	D	25	Total	O	0	0
			25	25		

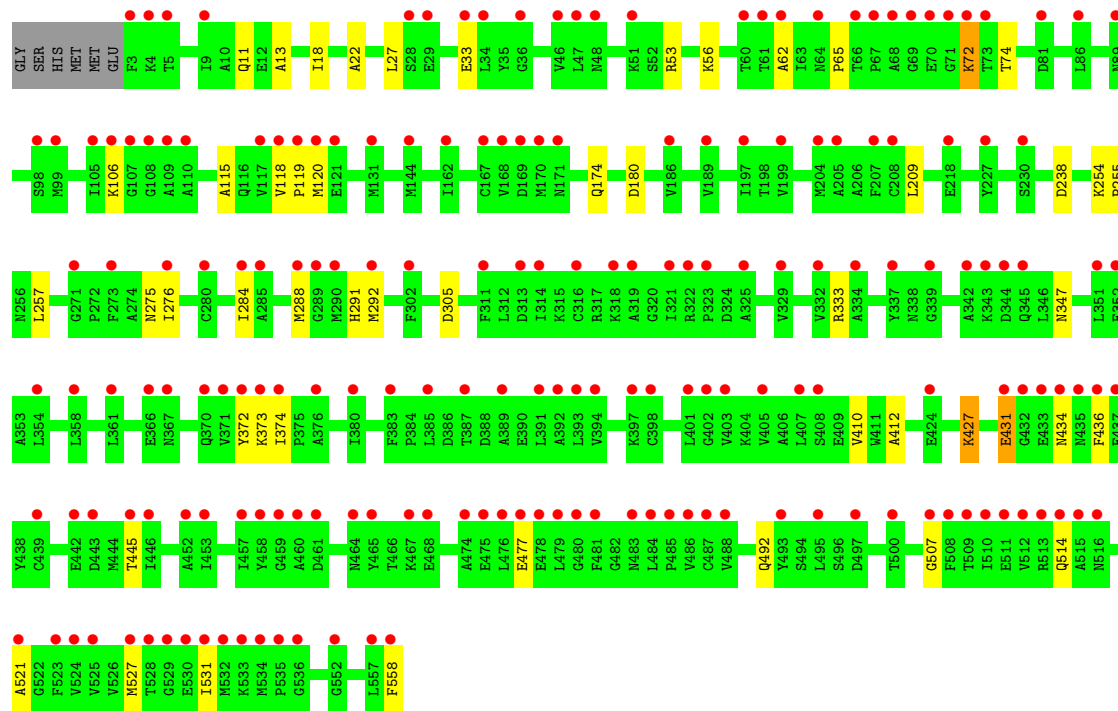
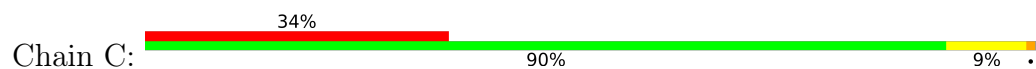
i

- Molecule 1: Formate--tetrahydrofolate ligase

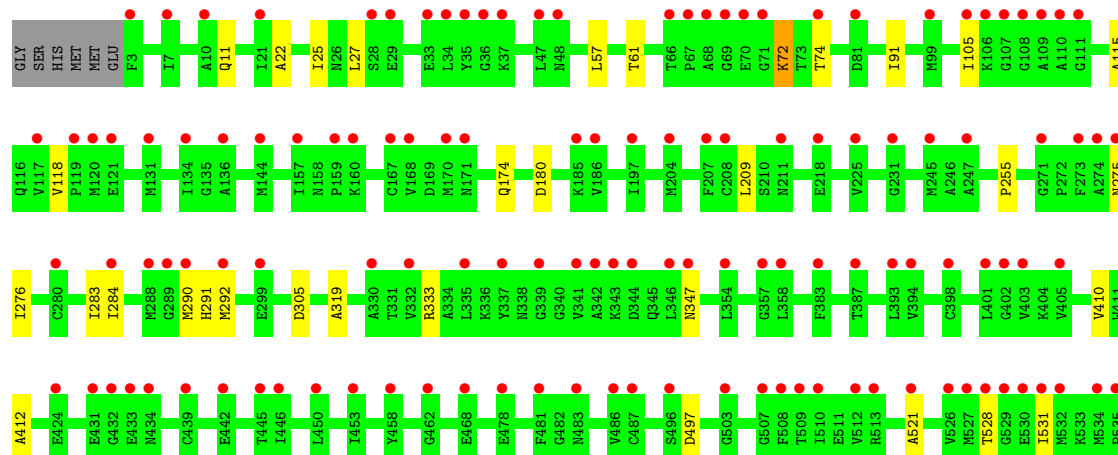




• Molecule 1: Formate--tetrahydrofolate ligase



• Molecule 1: Formate--tetrahydrofolate ligase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	215.87Å 116.30Å 107.38Å 90.00° 92.58° 90.00°	Depositor
Resolution (Å)	30.50 – 2.04 30.50 – 2.04	Depositor EDS
% Data completeness (in resolution range)	99.8 (30.50-2.04) 99.9 (30.50-2.04)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 2.05Å)	Xtriage
Refinement program	PHENIX 1.14 _3260	Depositor
R, R_{free}	0.272 , 0.283 0.286 , 0.297	Depositor DCC
R_{free} test set	2012 reflections (1.20%)	wwPDB-VP
Wilson B-factor (Å ²)	36.0	Xtriage
Anisotropy	0.408	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.55$, $\langle L^2 \rangle = 0.40$	Xtriage
Estimated twinning fraction	0.016 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	16831	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 54.13 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.8167e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: TAM, 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	0.11	0/4218	0.30	0/5711
1	B	0.10	0/4221	0.28	0/5716
1	C	0.10	0/4232	0.29	0/5731
1	D	0.11	0/4195	0.29	0/5685
All	All	0.11	0/16866	0.29	0/22843

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	4159	0	4195	27	0
1	B	4162	0	4190	23	0
1	C	4172	0	4201	32	0
1	D	4137	0	4155	23	0
2	A	16	0	24	2	0
2	B	4	0	6	2	0
2	C	12	0	18	4	0
2	D	8	0	12	1	0
3	A	11	0	17	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	11	0	17	3	0
3	C	11	0	17	3	0
3	D	11	0	17	1	0
4	A	28	0	0	0	0
4	B	36	0	0	1	0
4	C	28	0	0	0	0
4	D	25	0	0	0	0
All	All	16831	0	16869	103	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 3.

The worst 5 of 103 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:105:ILE:HG22	1:A:106:LYS:H	1.46	0.79
1:D:291:HIS:HD2	1:D:292:MET:HE3	1.49	0.76
1:D:11:GLN:O	3:D:603:TAM:N	2.25	0.69
1:D:57:LEU:HB2	1:D:290:MET:HE2	1.76	0.68
1:A:173:ARG:HH21	2:A:603:EDO:H21	1.59	0.68

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	553/562 (98%)	532 (96%)	21 (4%)	0	100	100
1	B	554/562 (99%)	535 (97%)	19 (3%)	0	100	100
1	C	554/562 (99%)	532 (96%)	21 (4%)	1 (0%)	43	37
1	D	553/562 (98%)	531 (96%)	22 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	2214/2248 (98%)	2130 (96%)	83 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	507	GLY

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	434/447 (97%)	432 (100%)	2 (0%)	81	85
1	B	434/447 (97%)	434 (100%)	0	100	100
1	C	436/447 (98%)	432 (99%)	4 (1%)	70	75
1	D	430/447 (96%)	427 (99%)	3 (1%)	76	80
All	All	1734/1788 (97%)	1725 (100%)	9 (0%)	81	85

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	105	ILE
1	D	118	VAL
1	C	106	LYS
1	C	427	LYS
1	C	431	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 15 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	514	GLN
1	D	16	GLN
1	B	516	ASN
1	D	516	ASN

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Mol	Chain	Res	Type
1	C	514	GLN

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 ⓘ

14 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	TAM	D	603	-	10,10,10	0.50	0	12,12,12	1.31	1 (8%)
3	TAM	B	602	-	10,10,10	0.47	0	12,12,12	1.60	3 (25%)
2	EDO	A	603	-	3,3,3	0.43	0	2,2,2	0.35	0
2	EDO	C	603	-	3,3,3	0.39	0	2,2,2	0.48	0
2	EDO	C	602	-	3,3,3	0.42	0	2,2,2	0.34	0
2	EDO	D	602	-	3,3,3	0.43	0	2,2,2	0.35	0
3	TAM	A	605	-	10,10,10	0.52	0	12,12,12	5.39	8 (66%)
2	EDO	A	601	-	3,3,3	0.42	0	2,2,2	0.31	0
2	EDO	B	601	-	3,3,3	0.42	0	2,2,2	0.32	0
2	EDO	C	601	-	3,3,3	0.43	0	2,2,2	0.38	0
3	TAM	C	604	-	10,10,10	0.51	0	12,12,12	1.53	3 (25%)
2	EDO	A	602	-	3,3,3	0.43	0	2,2,2	0.37	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	EDO	D	601	-	3,3,3	0.42	0	2,2,2	0.39	0
2	EDO	A	604	-	3,3,3	0.42	0	2,2,2	0.43	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	TAM	D	603	-	-	7/12/12/12	-
3	TAM	B	602	-	-	11/12/12/12	-
2	EDO	A	603	-	-	0/1/1/1	-
2	EDO	C	603	-	-	0/1/1/1	-
2	EDO	C	602	-	-	1/1/1/1	-
2	EDO	D	602	-	-	1/1/1/1	-
3	TAM	A	605	-	-	7/12/12/12	-
2	EDO	A	601	-	-	1/1/1/1	-
2	EDO	B	601	-	-	0/1/1/1	-
2	EDO	C	601	-	-	0/1/1/1	-
3	TAM	C	604	-	-	6/12/12/12	-
2	EDO	A	602	-	-	0/1/1/1	-
2	EDO	D	601	-	-	0/1/1/1	-
2	EDO	A	604	-	-	0/1/1/1	-

There are no bond length outliers.

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	605	TAM	C1-C-N	-10.21	82.58	108.22
3	A	605	TAM	C2-C-N	-9.19	85.16	108.22
3	A	605	TAM	C3-C-N	-7.30	89.89	108.22
3	A	605	TAM	C2-C-C1	5.73	120.62	110.50
3	A	605	TAM	C3-C-C1	4.95	119.23	110.50

There are no chirality outliers.

5 of 34 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	605	TAM	C3-C-C1-C4
3	A	605	TAM	N-C-C1-C4

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Mol	Chain	Res	Type	Atoms
3	A	605	TAM	C1-C-C2-C5
3	A	605	TAM	N-C-C2-C5
3	A	605	TAM	C1-C-C3-C6

There are no ring outliers.

10 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	603	TAM	1	0
3	B	602	TAM	3	0
2	A	603	EDO	1	0
2	C	603	EDO	2	0
2	C	602	EDO	1	0
2	A	601	EDO	1	0
2	B	601	EDO	2	0
2	C	601	EDO	1	0
3	C	604	TAM	3	0
2	D	601	EDO	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(Å²)	Q<0.9
1	A	555/562 (98%)	1.63	166 (29%)	1 0	31, 54, 88, 102	0
1	B	556/562 (98%)	1.75	189 (33%)	1 0	32, 52, 81, 96	0
1	C	556/562 (98%)	1.68	189 (33%)	1 0	31, 50, 76, 104	0
1	D	555/562 (98%)	1.52	130 (23%)	2 1	31, 50, 78, 106	0
All	All	2222/2248 (98%)	1.65	674 (30%)	1 0	31, 52, 82, 106	0

The worst 5 of 674 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	68	ALA	8.0
1	D	110	ALA	7.1
1	B	105	ILE	5.8
1	C	529	GLY	5.8
1	D	105	ILE	5.7

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($\text{\AA}^2$)	Q<0.9
2	EDO	A	604	4/4	0.40	0.25	56,60,61,63	0
3	TAM	B	602	11/11	0.52	0.28	61,70,76,77	0
3	TAM	C	604	11/11	0.59	0.25	57,71,75,81	0
2	EDO	C	603	4/4	0.61	0.21	55,56,59,60	0
2	EDO	A	602	4/4	0.62	0.24	47,51,51,52	0
3	TAM	D	603	11/11	0.62	0.23	69,74,81,81	0
3	TAM	A	605	11/11	0.64	0.19	75,78,82,83	0
2	EDO	D	602	4/4	0.65	0.24	43,57,60,61	0
2	EDO	A	601	4/4	0.67	0.18	50,52,55,57	0
2	EDO	D	601	4/4	0.74	0.22	39,40,50,55	0
2	EDO	C	601	4/4	0.75	0.21	33,38,45,54	0
2	EDO	C	602	4/4	0.77	0.23	33,42,51,53	0
2	EDO	B	601	4/4	0.78	0.21	35,48,52,58	0
2	EDO	A	603	4/4	0.85	0.16	35,47,51,56	0

6.5 Other polymers [i](#)

There are no such residues in this entry.