



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 27, 2026 – 11:08 AM UTC

PDB ID : 4PZU / pdb_00004pzu
Title : Crystal structure of a putative uncharacterize protein Rv3404c and likely sugar N-formyltransferase from Mycobacterium tuberculosis
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2014-03-31
Resolution : 2.10 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

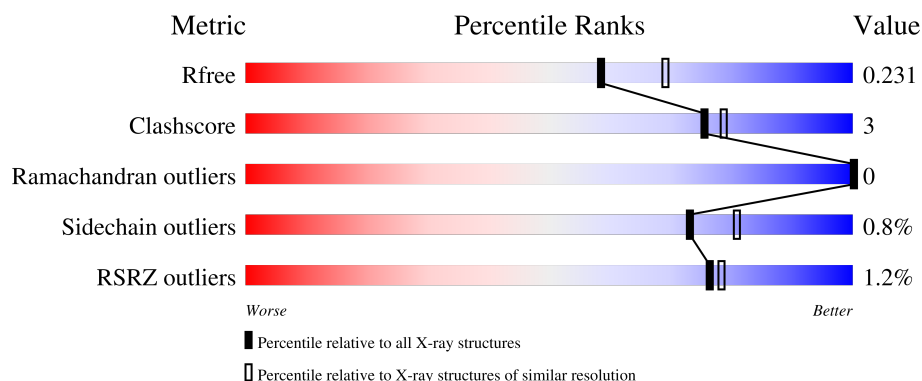
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

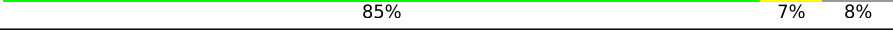
The reported resolution of this entry is 2.10 Å.

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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	256	 84% 8% 8%
1	B	256	 87% 9% 4%
1	C	256	 86% 6% 8% 1%
1	D	256	 85% 7% 8%
1	E	256	 86% 5% 9%

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Mol	Chain	Length	Quality of chain
1	F	256	86%6%8%
1	G	256	4% 86%5%9%
1	H	256	2% 84%6%10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 15706 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 Uncharacterized protein Rv3404c/MT3512.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	235	Total	C	N	O	S	0	0	0
			1836	1160	329	341	6			
1	B	232	Total	C	N	O	S	0	1	0
			1837	1162	333	336	6			
1	C	235	Total	C	N	O	S	0	1	0
			1853	1171	337	339	6			
1	D	235	Total	C	N	O	S	0	1	0
			1853	1171	336	340	6			
1	E	232	Total	C	N	O	S	0	1	0
			1825	1154	323	341	7			
1	F	235	Total	C	N	O	S	0	0	0
			1840	1164	332	338	6			
1	G	232	Total	C	N	O	S	0	0	0
			1791	1135	317	333	6			
1	H	231	Total	C	N	O	S	0	0	0
			1810	1144	323	337	6			

There are 176 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	expression tag	UNP P65073
A	-19	ALA	-	expression tag	UNP P65073
A	-18	HIS	-	expression tag	UNP P65073
A	-17	HIS	-	expression tag	UNP P65073
A	-16	HIS	-	expression tag	UNP P65073
A	-15	HIS	-	expression tag	UNP P65073
A	-14	HIS	-	expression tag	UNP P65073
A	-13	HIS	-	expression tag	UNP P65073
A	-12	MET	-	expression tag	UNP P65073
A	-11	GLY	-	expression tag	UNP P65073
A	-10	THR	-	expression tag	UNP P65073
A	-9	LEU	-	expression tag	UNP P65073
A	-8	GLU	-	expression tag	UNP P65073

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	ALA	-	expression tag	UNP P65073
A	-6	GLN	-	expression tag	UNP P65073
A	-5	THR	-	expression tag	UNP P65073
A	-4	GLN	-	expression tag	UNP P65073
A	-3	GLY	-	expression tag	UNP P65073
A	-2	PRO	-	expression tag	UNP P65073
A	-1	GLY	-	expression tag	UNP P65073
A	0	SER	-	expression tag	UNP P65073
A	2	VAL	-	expression tag	UNP P65073
B	-20	MET	-	expression tag	UNP P65073
B	-19	ALA	-	expression tag	UNP P65073
B	-18	HIS	-	expression tag	UNP P65073
B	-17	HIS	-	expression tag	UNP P65073
B	-16	HIS	-	expression tag	UNP P65073
B	-15	HIS	-	expression tag	UNP P65073
B	-14	HIS	-	expression tag	UNP P65073
B	-13	HIS	-	expression tag	UNP P65073
B	-12	MET	-	expression tag	UNP P65073
B	-11	GLY	-	expression tag	UNP P65073
B	-10	THR	-	expression tag	UNP P65073
B	-9	LEU	-	expression tag	UNP P65073
B	-8	GLU	-	expression tag	UNP P65073
B	-7	ALA	-	expression tag	UNP P65073
B	-6	GLN	-	expression tag	UNP P65073
B	-5	THR	-	expression tag	UNP P65073
B	-4	GLN	-	expression tag	UNP P65073
B	-3	GLY	-	expression tag	UNP P65073
B	-2	PRO	-	expression tag	UNP P65073
B	-1	GLY	-	expression tag	UNP P65073
B	0	SER	-	expression tag	UNP P65073
B	2	VAL	-	expression tag	UNP P65073
C	-20	MET	-	expression tag	UNP P65073
C	-19	ALA	-	expression tag	UNP P65073
C	-18	HIS	-	expression tag	UNP P65073
C	-17	HIS	-	expression tag	UNP P65073
C	-16	HIS	-	expression tag	UNP P65073
C	-15	HIS	-	expression tag	UNP P65073
C	-14	HIS	-	expression tag	UNP P65073
C	-13	HIS	-	expression tag	UNP P65073
C	-12	MET	-	expression tag	UNP P65073
C	-11	GLY	-	expression tag	UNP P65073
C	-10	THR	-	expression tag	UNP P65073

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-9	LEU	-	expression tag	UNP P65073
C	-8	GLU	-	expression tag	UNP P65073
C	-7	ALA	-	expression tag	UNP P65073
C	-6	GLN	-	expression tag	UNP P65073
C	-5	THR	-	expression tag	UNP P65073
C	-4	GLN	-	expression tag	UNP P65073
C	-3	GLY	-	expression tag	UNP P65073
C	-2	PRO	-	expression tag	UNP P65073
C	-1	GLY	-	expression tag	UNP P65073
C	0	SER	-	expression tag	UNP P65073
C	2	VAL	-	expression tag	UNP P65073
D	-20	MET	-	expression tag	UNP P65073
D	-19	ALA	-	expression tag	UNP P65073
D	-18	HIS	-	expression tag	UNP P65073
D	-17	HIS	-	expression tag	UNP P65073
D	-16	HIS	-	expression tag	UNP P65073
D	-15	HIS	-	expression tag	UNP P65073
D	-14	HIS	-	expression tag	UNP P65073
D	-13	HIS	-	expression tag	UNP P65073
D	-12	MET	-	expression tag	UNP P65073
D	-11	GLY	-	expression tag	UNP P65073
D	-10	THR	-	expression tag	UNP P65073
D	-9	LEU	-	expression tag	UNP P65073
D	-8	GLU	-	expression tag	UNP P65073
D	-7	ALA	-	expression tag	UNP P65073
D	-6	GLN	-	expression tag	UNP P65073
D	-5	THR	-	expression tag	UNP P65073
D	-4	GLN	-	expression tag	UNP P65073
D	-3	GLY	-	expression tag	UNP P65073
D	-2	PRO	-	expression tag	UNP P65073
D	-1	GLY	-	expression tag	UNP P65073
D	0	SER	-	expression tag	UNP P65073
D	2	VAL	-	expression tag	UNP P65073
E	-20	MET	-	expression tag	UNP P65073
E	-19	ALA	-	expression tag	UNP P65073
E	-18	HIS	-	expression tag	UNP P65073
E	-17	HIS	-	expression tag	UNP P65073
E	-16	HIS	-	expression tag	UNP P65073
E	-15	HIS	-	expression tag	UNP P65073
E	-14	HIS	-	expression tag	UNP P65073
E	-13	HIS	-	expression tag	UNP P65073
E	-12	MET	-	expression tag	UNP P65073

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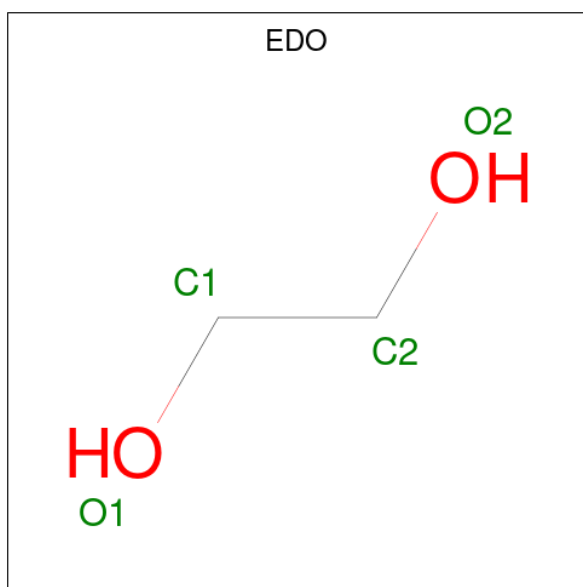
Chain	Residue	Modelled	Actual	Comment	Reference
E	-11	GLY	-	expression tag	UNP P65073
E	-10	THR	-	expression tag	UNP P65073
E	-9	LEU	-	expression tag	UNP P65073
E	-8	GLU	-	expression tag	UNP P65073
E	-7	ALA	-	expression tag	UNP P65073
E	-6	GLN	-	expression tag	UNP P65073
E	-5	THR	-	expression tag	UNP P65073
E	-4	GLN	-	expression tag	UNP P65073
E	-3	GLY	-	expression tag	UNP P65073
E	-2	PRO	-	expression tag	UNP P65073
E	-1	GLY	-	expression tag	UNP P65073
E	0	SER	-	expression tag	UNP P65073
E	2	VAL	-	expression tag	UNP P65073
F	-20	MET	-	expression tag	UNP P65073
F	-19	ALA	-	expression tag	UNP P65073
F	-18	HIS	-	expression tag	UNP P65073
F	-17	HIS	-	expression tag	UNP P65073
F	-16	HIS	-	expression tag	UNP P65073
F	-15	HIS	-	expression tag	UNP P65073
F	-14	HIS	-	expression tag	UNP P65073
F	-13	HIS	-	expression tag	UNP P65073
F	-12	MET	-	expression tag	UNP P65073
F	-11	GLY	-	expression tag	UNP P65073
F	-10	THR	-	expression tag	UNP P65073
F	-9	LEU	-	expression tag	UNP P65073
F	-8	GLU	-	expression tag	UNP P65073
F	-7	ALA	-	expression tag	UNP P65073
F	-6	GLN	-	expression tag	UNP P65073
F	-5	THR	-	expression tag	UNP P65073
F	-4	GLN	-	expression tag	UNP P65073
F	-3	GLY	-	expression tag	UNP P65073
F	-2	PRO	-	expression tag	UNP P65073
F	-1	GLY	-	expression tag	UNP P65073
F	0	SER	-	expression tag	UNP P65073
F	2	VAL	-	expression tag	UNP P65073
G	-20	MET	-	expression tag	UNP P65073
G	-19	ALA	-	expression tag	UNP P65073
G	-18	HIS	-	expression tag	UNP P65073
G	-17	HIS	-	expression tag	UNP P65073
G	-16	HIS	-	expression tag	UNP P65073
G	-15	HIS	-	expression tag	UNP P65073
G	-14	HIS	-	expression tag	UNP P65073

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-13	HIS	-	expression tag	UNP P65073
G	-12	MET	-	expression tag	UNP P65073
G	-11	GLY	-	expression tag	UNP P65073
G	-10	THR	-	expression tag	UNP P65073
G	-9	LEU	-	expression tag	UNP P65073
G	-8	GLU	-	expression tag	UNP P65073
G	-7	ALA	-	expression tag	UNP P65073
G	-6	GLN	-	expression tag	UNP P65073
G	-5	THR	-	expression tag	UNP P65073
G	-4	GLN	-	expression tag	UNP P65073
G	-3	GLY	-	expression tag	UNP P65073
G	-2	PRO	-	expression tag	UNP P65073
G	-1	GLY	-	expression tag	UNP P65073
G	0	SER	-	expression tag	UNP P65073
G	2	VAL	-	expression tag	UNP P65073
H	-20	MET	-	expression tag	UNP P65073
H	-19	ALA	-	expression tag	UNP P65073
H	-18	HIS	-	expression tag	UNP P65073
H	-17	HIS	-	expression tag	UNP P65073
H	-16	HIS	-	expression tag	UNP P65073
H	-15	HIS	-	expression tag	UNP P65073
H	-14	HIS	-	expression tag	UNP P65073
H	-13	HIS	-	expression tag	UNP P65073
H	-12	MET	-	expression tag	UNP P65073
H	-11	GLY	-	expression tag	UNP P65073
H	-10	THR	-	expression tag	UNP P65073
H	-9	LEU	-	expression tag	UNP P65073
H	-8	GLU	-	expression tag	UNP P65073
H	-7	ALA	-	expression tag	UNP P65073
H	-6	GLN	-	expression tag	UNP P65073
H	-5	THR	-	expression tag	UNP P65073
H	-4	GLN	-	expression tag	UNP P65073
H	-3	GLY	-	expression tag	UNP P65073
H	-2	PRO	-	expression tag	UNP P65073
H	-1	GLY	-	expression tag	UNP P65073
H	0	SER	-	expression tag	UNP P65073
H	2	VAL	-	expression tag	UNP P65073

- 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	B	1	Total	C	O	0	0
			4	2	2		
2	B	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		
2	D	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	E	1	Total C O 4 2 2	0	0
2	E	1	Total C O 4 2 2	0	0
2	F	1	Total C O 4 2 2	0	0
2	F	1	Total C O 4 2 2	0	0
2	F	1	Total C O 4 2 2	0	0
2	G	1	Total C O 4 2 2	0	0
2	G	1	Total C O 4 2 2	0	0
2	H	1	Total C O 4 2 2	0	0

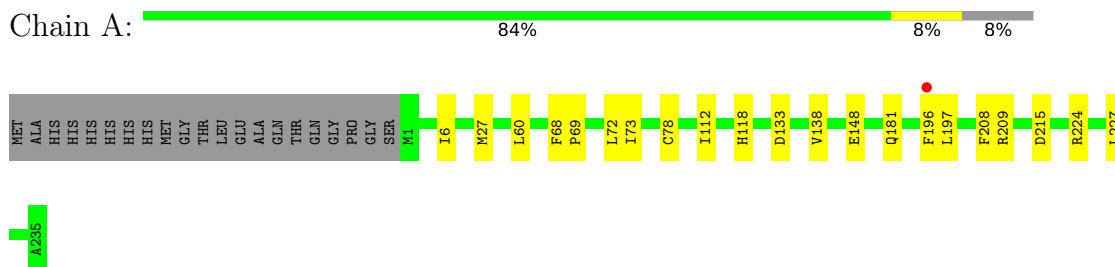
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	153	Total O 153 153	0	0
3	B	136	Total O 136 136	0	0
3	C	158	Total O 158 158	0	0
3	D	139	Total O 139 139	0	0
3	E	156	Total O 156 156	0	0
3	F	92	Total O 92 92	0	0
3	G	90	Total O 90 90	0	0
3	H	49	Total O 49 49	0	0

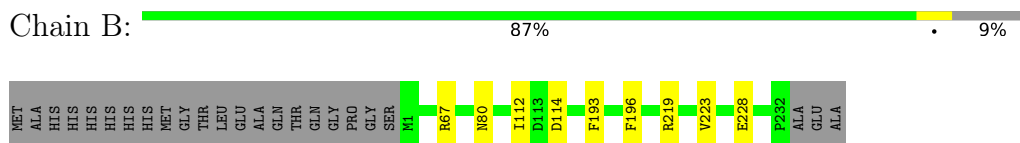
3 Residue-property plots [i](#)

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.

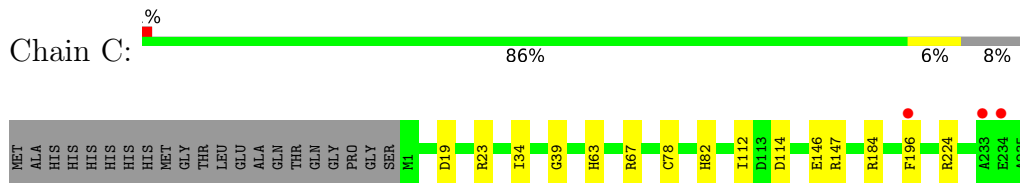
- Molecule 1: Uncharacterized protein Rv3404c/MT3512



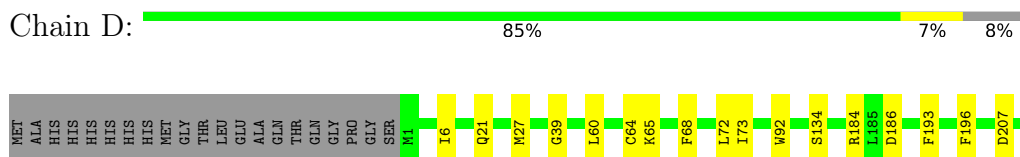
- Molecule 1: Uncharacterized protein Rv3404c/MT3512



- Molecule 1: Uncharacterized protein Rv3404c/MT3512



- Molecule 1: Uncharacterized protein Rv3404c/MT3512

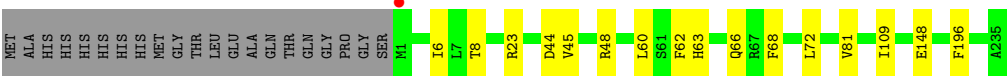
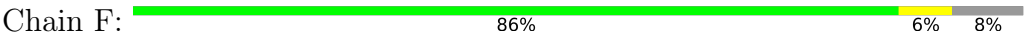


- Molecule 1: Uncharacterized protein Rv3404c/MT3512

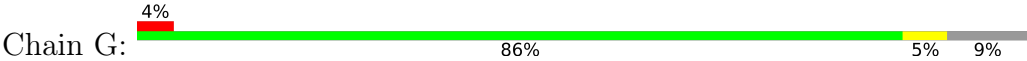




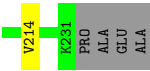
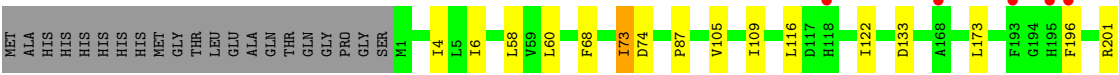
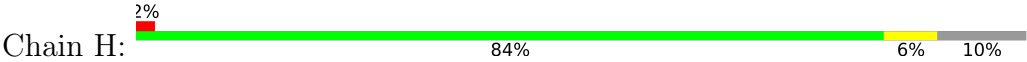
• Molecule 1: Uncharacterized protein Rv3404c/MT3512



• Molecule 1: Uncharacterized protein Rv3404c/MT3512



• Molecule 1: Uncharacterized protein Rv3404c/MT3512



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	89.08Å 136.45Å 99.97Å 90.00° 103.74° 90.00°	Depositor
Resolution (Å)	20.00 – 2.10 20.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.8 (20.00-2.10) 99.8 (20.00-2.10)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 2.09Å)	Xtriage
Refinement program	REFMAC 5.8.0069	Depositor
R, R_{free}	0.187 , 0.231 0.187 , 0.231	Depositor DCC
R_{free} test set	6535 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	33.6	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 40.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15706	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% 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

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	0/1879	0.94	1/2558 (0.0%)
1	B	0.80	0/1883	0.91	0/2560
1	C	0.83	0/1899	0.93	2/2581 (0.1%)
1	D	0.81	0/1899	0.89	0/2581
1	E	0.85	0/1871	0.92	1/2548 (0.0%)
1	F	0.74	0/1883	0.87	0/2561
1	G	0.74	0/1834	0.92	0/2502
1	H	0.72	0/1852	0.91	1/2522 (0.0%)
All	All	0.79	0/15000	0.91	5/20413 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	74	ASP	N-CA-C	6.28	118.12	111.28
1	A	118	HIS	N-CA-C	-5.52	100.33	108.99
1	E	118	HIS	N-CA-C	-5.28	100.71	108.99
1	C	82	HIS	CA-C-N	-5.19	116.41	121.65
1	C	82	HIS	C-N-CA	-5.19	116.41	121.65

There are no chirality outliers.

There are no planarity outliers.

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	1836	0	1750	19	0
1	B	1837	0	1776	10	0
1	C	1853	0	1790	19	0
1	D	1853	0	1788	18	0
1	E	1825	0	1740	12	0
1	F	1840	0	1770	14	0
1	G	1791	0	1684	14	0
1	H	1810	0	1724	13	0
2	A	12	0	18	1	0
2	B	12	0	18	0	0
2	C	12	0	18	0	0
2	D	16	0	24	3	0
2	E	12	0	18	0	0
2	F	12	0	18	0	0
2	G	8	0	12	0	0
2	H	4	0	6	0	0
3	A	153	0	0	3	0
3	B	136	0	0	0	0
3	C	158	0	0	5	0
3	D	139	0	0	1	0
3	E	156	0	0	2	0
3	F	92	0	0	0	0
3	G	90	0	0	1	0
3	H	49	0	0	2	0
All	All	15706	0	14154	86	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.

All (86) 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:196:PHE:HD2	1:B:196:PHE:CD2	1.44	1.32
1:E:196:PHE:HD2	1:F:196:PHE:CD2	1.47	1.31
1:E:196:PHE:CD2	1:F:196:PHE:HD2	1.56	1.23
1:A:196:PHE:CD2	1:B:196:PHE:CD2	2.31	1.17
1:A:196:PHE:CD2	1:B:196:PHE:HD2	1.61	1.17
1:C:196:PHE:CD2	1:D:196:PHE:CD2	2.34	1.14
1:C:196:PHE:CD2	1:D:196:PHE:HD2	1.65	1.14
1:E:196:PHE:CD2	1:F:196:PHE:CD2	2.33	1.10
1:C:196:PHE:HD2	1:D:196:PHE:CD2	1.71	1.03
1:E:146:GLU:OE1	3:E:492:HOH:O	2.05	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:196:PHE:HD2	1:D:196:PHE:CE2	2.05	0.74
1:C:196:PHE:CE2	1:D:196:PHE:HD2	2.05	0.71
1:D:134:SER:OG	2:D:303:EDO:H22	1.92	0.68
1:C:19:ASP:OD2	1:C:147[B]:ARG:NH2	2.27	0.66
1:C:23:ARG:NH1	1:E:25:GLY:O	2.27	0.66
1:A:196:PHE:CE2	1:B:196:PHE:HD2	2.12	0.65
1:G:193:PHE:HA	1:H:196:PHE:CE2	2.32	0.65
1:H:133:ASP:O	1:H:201:ARG:HD2	1.97	0.65
1:A:69:PRO:HD3	3:A:460:HOH:O	1.97	0.63
1:G:193:PHE:HA	1:H:196:PHE:HE2	1.63	0.63
1:H:87:PRO:HD3	3:H:409:HOH:O	2.00	0.62
1:E:196:PHE:HD2	1:F:196:PHE:CE2	2.14	0.61
1:A:73:ILE:HG23	2:A:302:EDO:H22	1.83	0.60
1:G:193:PHE:CG	1:H:196:PHE:HE2	2.20	0.60
1:G:32:SER:HB2	1:G:33:PRO:HD2	1.85	0.59
1:A:209:ARG:NH2	1:B:228:GLU:OE2	2.27	0.57
1:E:196:PHE:CE2	1:F:196:PHE:HD2	2.18	0.55
1:G:193:PHE:CD1	1:H:196:PHE:HE2	2.24	0.55
1:B:80:ASN:HB2	1:B:112:ILE:HD11	1.89	0.55
1:C:224:ARG:NE	3:C:523:HOH:O	2.29	0.54
1:C:184:ARG:NH1	3:C:541:HOH:O	2.41	0.54
1:D:184[A]:ARG:HH12	1:D:186:ASP:HB2	1.74	0.52
1:G:193:PHE:CG	1:H:196:PHE:CE2	2.98	0.51
1:A:227:LEU:HD23	1:B:223:VAL:HG22	1.91	0.50
1:D:73:ILE:HG23	2:D:302:EDO:H21	1.93	0.50
1:A:208:PHE:N	1:A:208:PHE:CD1	2.79	0.49
1:H:109:ILE:HG22	1:H:122:ILE:HD12	1.95	0.49
1:C:67:ARG:HD2	1:C:114:ASP:O	2.13	0.49
1:D:68:PHE:CD2	1:D:72:LEU:HD23	2.47	0.49
1:G:193:PHE:CD1	1:H:196:PHE:CE2	3.01	0.48
1:C:147[B]:ARG:HD3	1:E:39:GLY:O	2.13	0.48
1:A:68:PHE:CD2	1:A:72:LEU:HD23	2.48	0.48
1:A:196:PHE:CE1	1:B:193:PHE:CD2	3.01	0.48
1:A:215:ASP:HB2	3:A:531:HOH:O	2.13	0.47
1:A:6:ILE:HD12	1:A:27:MET:SD	2.55	0.46
1:H:116:LEU:HD23	1:H:173:LEU:HD11	1.98	0.46
1:E:148:GLU:HG3	1:G:39:GLY:HA2	1.96	0.46
1:E:195:HIS:HD2	3:E:429:HOH:O	1.97	0.46
1:F:44:ASP:O	1:F:48:ARG:HG2	2.16	0.46
1:H:68:PHE:HB2	1:H:73:ILE:HD11	1.98	0.46
1:D:21:GLN:HE21	1:F:23:ARG:HH21	1.63	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:116:LEU:HD23	1:G:173:LEU:HD11	1.97	0.45
1:C:146:GLU:OE1	3:C:471:HOH:O	2.21	0.45
1:A:78:CYS:HB3	1:A:112:ILE:HD12	1.98	0.45
1:G:193:PHE:HE2	3:H:429:HOH:O	1.99	0.44
1:D:64:CYS:SG	2:D:304:EDO:H11	2.58	0.44
1:H:6:ILE:HA	1:H:60:LEU:O	2.18	0.44
1:A:197:LEU:HG	1:B:196:PHE:CZ	2.52	0.44
1:F:8:THR:HB	1:F:62:PHE:HB2	2.00	0.43
1:A:224:ARG:NH2	3:A:532:HOH:O	2.40	0.43
1:G:195:HIS:CD2	3:G:408:HOH:O	2.71	0.43
1:C:224:ARG:NH2	3:C:523:HOH:O	2.43	0.43
1:F:81:VAL:HG22	1:F:109:ILE:HG12	2.01	0.43
1:C:196:PHE:HE1	1:D:193:PHE:CD2	2.36	0.43
1:D:65:LYS:HD3	1:D:92:TRP:CZ2	2.54	0.43
1:D:215:ASP:HB2	3:D:454:HOH:O	2.18	0.43
1:C:63:HIS:CD2	1:C:63:HIS:N	2.87	0.42
1:D:6:ILE:HA	1:D:60:LEU:O	2.19	0.42
1:D:39:GLY:HA2	1:F:148:GLU:HG3	2.02	0.42
1:C:196:PHE:CE1	1:D:193:PHE:CD2	3.07	0.42
1:E:143:MET:HE2	1:E:143:MET:HA	2.01	0.42
1:E:231:LYS:HA	1:E:232:PRO:HD3	1.80	0.42
1:F:6:ILE:HA	1:F:60:LEU:O	2.20	0.42
1:G:45:VAL:CG1	1:G:72:LEU:HD22	2.49	0.42
1:A:148:GLU:HG3	1:C:39:GLY:HA2	2.00	0.42
1:B:67:ARG:HD2	1:B:114:ASP:O	2.19	0.42
1:C:147[B]:ARG:NH1	3:C:498:HOH:O	2.53	0.42
1:H:4:ILE:HG12	1:H:58:LEU:HD23	2.01	0.41
1:A:133:ASP:HB3	1:A:138:VAL:HG23	2.01	0.41
1:D:27:MET:O	1:F:23:ARG:NH2	2.53	0.41
1:F:63:HIS:N	1:F:63:HIS:CD2	2.88	0.41
1:A:6:ILE:HA	1:A:60:LEU:O	2.21	0.41
1:F:68:PHE:CD2	1:F:72:LEU:HD23	2.56	0.40
1:G:208:PHE:N	1:G:208:PHE:CD1	2.88	0.40
1:C:78:CYS:HB3	1:C:112:ILE:HD12	2.03	0.40
1:G:63:HIS:CD2	1:G:63:HIS:N	2.87	0.40

There are no symmetry-related clashes.

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	233/256 (91%)	230 (99%)	3 (1%)	0	100	100
1	B	231/256 (90%)	224 (97%)	7 (3%)	0	100	100
1	C	234/256 (91%)	229 (98%)	5 (2%)	0	100	100
1	D	234/256 (91%)	231 (99%)	3 (1%)	0	100	100
1	E	231/256 (90%)	226 (98%)	5 (2%)	0	100	100
1	F	233/256 (91%)	229 (98%)	4 (2%)	0	100	100
1	G	230/256 (90%)	226 (98%)	4 (2%)	0	100	100
1	H	229/256 (90%)	227 (99%)	2 (1%)	0	100	100
All	All	1855/2048 (91%)	1822 (98%)	33 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	191/217 (88%)	190 (100%)	1 (0%)	81	88
1	B	194/217 (89%)	193 (100%)	1 (0%)	81	88
1	C	194/217 (89%)	193 (100%)	1 (0%)	81	88
1	D	194/217 (89%)	193 (100%)	1 (0%)	81	88
1	E	193/217 (89%)	191 (99%)	2 (1%)	68	76
1	F	192/217 (88%)	190 (99%)	2 (1%)	68	76

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	184/217 (85%)	183 (100%)	1 (0%)	81	88
1	H	190/217 (88%)	187 (98%)	3 (2%)	55	64
All	All	1532/1736 (88%)	1520 (99%)	12 (1%)	73	81

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	181	GLN
1	B	219	ARG
1	C	34	ILE
1	D	207	ASP
1	E	54	GLU
1	E	230	GLU
1	F	45	VAL
1	F	66	GLN
1	G	217	SER
1	H	73	ILE
1	H	105	VAL
1	H	214	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	HIS
1	A	103	GLN
1	B	12	HIS
1	C	12	HIS
1	C	103	GLN
1	D	12	HIS
1	D	21	GLN
1	E	195	HIS
1	F	12	HIS
1	F	80	ASN
1	F	103	GLN
1	F	124	GLN
1	G	12	HIS
1	H	12	HIS
1	H	82	HIS
1	H	103	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

22 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	D	303	-	3,3,3	0.58	0	2,2,2	0.21	0
2	EDO	B	303	-	3,3,3	0.34	0	2,2,2	0.74	0
2	EDO	E	301	-	3,3,3	0.47	0	2,2,2	0.56	0
2	EDO	C	303	-	3,3,3	0.40	0	2,2,2	0.42	0
2	EDO	D	301	-	3,3,3	0.60	0	2,2,2	0.45	0
2	EDO	D	304	-	3,3,3	0.56	0	2,2,2	0.26	0
2	EDO	F	302	-	3,3,3	0.48	0	2,2,2	0.56	0
2	EDO	H	301	-	3,3,3	0.49	0	2,2,2	0.37	0
2	EDO	C	301	-	3,3,3	0.38	0	2,2,2	0.63	0
2	EDO	A	302	-	3,3,3	0.36	0	2,2,2	0.38	0
2	EDO	C	302	-	3,3,3	0.43	0	2,2,2	0.28	0
2	EDO	F	303	-	3,3,3	0.56	0	2,2,2	0.26	0
2	EDO	E	302	-	3,3,3	0.47	0	2,2,2	0.47	0
2	EDO	F	301	-	3,3,3	0.40	0	2,2,2	0.67	0
2	EDO	D	302	-	3,3,3	0.65	0	2,2,2	0.23	0
2	EDO	A	301	-	3,3,3	0.46	0	2,2,2	0.51	0
2	EDO	A	303	-	3,3,3	0.51	0	2,2,2	0.38	0
2	EDO	G	302	-	3,3,3	0.44	0	2,2,2	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	EDO	B	302	-	3,3,3	0.52	0	2,2,2	0.43	0
2	EDO	B	301	-	3,3,3	0.40	0	2,2,2	0.23	0
2	EDO	G	301	-	3,3,3	0.87	0	2,2,2	0.32	0
2	EDO	E	303	-	3,3,3	0.33	0	2,2,2	0.73	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	D	303	-	-	1/1/1/1	-
2	EDO	B	303	-	-	0/1/1/1	-
2	EDO	E	301	-	-	0/1/1/1	-
2	EDO	C	303	-	-	0/1/1/1	-
2	EDO	D	301	-	-	0/1/1/1	-
2	EDO	D	304	-	-	1/1/1/1	-
2	EDO	F	302	-	-	0/1/1/1	-
2	EDO	H	301	-	-	0/1/1/1	-
2	EDO	C	301	-	-	1/1/1/1	-
2	EDO	A	302	-	-	1/1/1/1	-
2	EDO	C	302	-	-	0/1/1/1	-
2	EDO	F	303	-	-	1/1/1/1	-
2	EDO	E	302	-	-	0/1/1/1	-
2	EDO	F	301	-	-	0/1/1/1	-
2	EDO	D	302	-	-	1/1/1/1	-
2	EDO	A	301	-	-	1/1/1/1	-
2	EDO	A	303	-	-	0/1/1/1	-
2	EDO	G	302	-	-	1/1/1/1	-
2	EDO	B	302	-	-	0/1/1/1	-
2	EDO	B	301	-	-	1/1/1/1	-
2	EDO	G	301	-	-	0/1/1/1	-
2	EDO	E	303	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	G	302	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	D	304	EDO	O1-C1-C2-O2
2	D	303	EDO	O1-C1-C2-O2
2	A	302	EDO	O1-C1-C2-O2
2	D	302	EDO	O1-C1-C2-O2
2	A	301	EDO	O1-C1-C2-O2
2	F	303	EDO	O1-C1-C2-O2
2	C	301	EDO	O1-C1-C2-O2
2	B	301	EDO	O1-C1-C2-O2

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	303	EDO	1	0
2	D	304	EDO	1	0
2	A	302	EDO	1	0
2	D	302	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 ⓘ

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	235/256 (91%)	-0.48	1 (0%) 88 90	22, 33, 53, 64	0
1	B	232/256 (90%)	-0.54	0 100 100	18, 32, 48, 68	1 (0%)
1	C	235/256 (91%)	-0.49	3 (1%) 75 77	17, 33, 53, 90	1 (0%)
1	D	235/256 (91%)	-0.43	1 (0%) 88 90	22, 35, 58, 72	1 (0%)
1	E	232/256 (90%)	-0.50	1 (0%) 88 90	23, 32, 49, 71	1 (0%)
1	F	235/256 (91%)	-0.24	1 (0%) 88 90	24, 43, 67, 78	0
1	G	232/256 (90%)	-0.17	10 (4%) 40 41	23, 41, 80, 113	0
1	H	231/256 (90%)	0.20	5 (2%) 62 65	34, 54, 80, 97	0
All	All	1867/2048 (91%)	-0.33	22 (1%) 76 78	17, 37, 66, 113	4 (0%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	214	VAL	4.0
1	G	216	ALA	3.7
1	H	168	ALA	3.3
1	H	193	PHE	3.3
1	G	186	ASP	2.9
1	C	234	GLU	2.8
1	C	233	ALA	2.7
1	D	235	ALA	2.7
1	G	188	ASN	2.7
1	F	1	MET	2.7
1	H	196	PHE	2.7
1	G	212	TRP	2.6
1	A	196	PHE	2.6
1	G	232	PRO	2.5
1	G	196	PHE	2.4
1	E	196	PHE	2.3

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Mol	Chain	Res	Type	RSRZ
1	G	193	PHE	2.3
1	G	218	GLY	2.2
1	H	118	HIS	2.1
1	C	196	PHE	2.1
1	H	195	HIS	2.1
1	G	190	ARG	2.1

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
2	EDO	G	301	4/4	0.74	0.17	40,49,49,53	0
2	EDO	F	303	4/4	0.81	0.12	57,57,58,60	0
2	EDO	H	301	4/4	0.86	0.15	58,60,60,64	0
2	EDO	D	303	4/4	0.88	0.12	48,49,54,58	0
2	EDO	G	302	4/4	0.90	0.12	55,58,58,60	0
2	EDO	F	302	4/4	0.90	0.10	46,48,49,51	0
2	EDO	B	302	4/4	0.92	0.17	37,48,49,54	0
2	EDO	A	302	4/4	0.93	0.10	38,46,47,51	0
2	EDO	D	304	4/4	0.93	0.11	48,49,50,52	0
2	EDO	E	303	4/4	0.93	0.09	42,45,50,51	0
2	EDO	D	302	4/4	0.93	0.08	38,43,43,47	0
2	EDO	E	302	4/4	0.95	0.07	31,35,35,37	0
2	EDO	B	303	4/4	0.95	0.08	45,47,47,49	0
2	EDO	F	301	4/4	0.95	0.07	35,39,42,42	0
2	EDO	A	303	4/4	0.95	0.07	52,54,55,59	0
2	EDO	C	302	4/4	0.96	0.07	38,47,48,48	0
2	EDO	C	303	4/4	0.96	0.07	51,53,54,56	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	EDO	D	301	4/4	0.96	0.06	32,32,33,38	0
2	EDO	E	301	4/4	0.96	0.06	32,35,36,37	0
2	EDO	A	301	4/4	0.97	0.07	31,32,33,33	0
2	EDO	C	301	4/4	0.98	0.06	32,33,34,35	0
2	EDO	B	301	4/4	0.98	0.05	31,32,33,35	0

6.5 Other polymers [i](#)

There are no such residues in this entry.