



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:38 PM GMT

PDB ID : 3UG3  
Title : Crystal structure of alpha-L-arabinofuranosidase from *Thermotoga maritima*  
ligand free form  
Authors : Im, D.-H.; Miyazaki, K.; Wakagi, T.; Fushinobu, S.  
Deposited on : 2011-11-02  
Resolution : 1.80 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

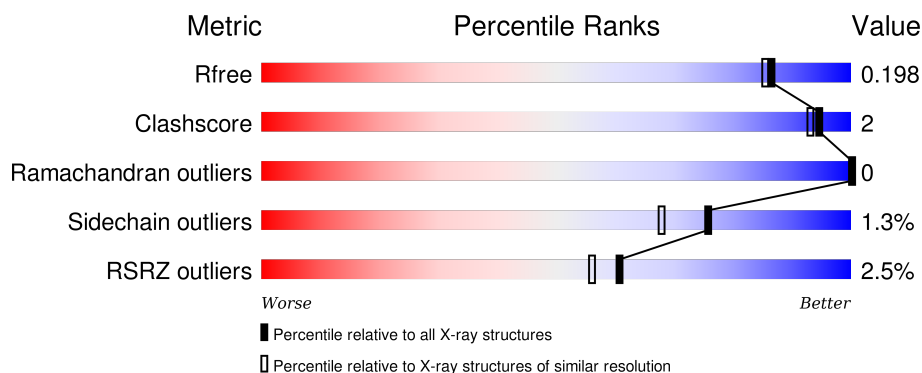
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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))
$R_{free}$	91344	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	504	<div> <div>7%</div> <div>91% 5% .</div> </div>
1	B	504	<div> <div>7%</div> <div>91% 5% .</div> </div>
1	C	504	<div> <div>0%</div> <div>92% . .</div> </div>
1	D	504	<div> <div>3%</div> <div>88% 6% . 5%</div> </div>
1	E	504	<div> <div>2%</div> <div>91% . .</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	504	 % 91% 5%

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	EDO	A	494	-	-	-	X
3	EDO	B	493	-	-	-	X
3	EDO	C	493	-	-	-	X
3	EDO	D	493	-	-	-	X
4	NA	A	496	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 25934 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 Alpha-L-arabinofuranosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	482	Total	C	N	O	S	0	0	0
			3876	2488	644	729	15			
1	B	482	Total	C	N	O	S	0	0	0
			3876	2488	644	729	15			
1	C	483	Total	C	N	O	S	0	0	0
			3885	2493	645	732	15			
1	D	481	Total	C	N	O	S	0	0	0
			3868	2482	643	728	15			
1	E	482	Total	C	N	O	S	0	0	0
			3876	2488	644	729	15			
1	F	482	Total	C	N	O	S	0	0	0
			3876	2488	644	729	15			

There are 126 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP Q9WYB7
A	-18	GLY	-	EXPRESSION TAG	UNP Q9WYB7
A	-17	SER	-	EXPRESSION TAG	UNP Q9WYB7
A	-16	SER	-	EXPRESSION TAG	UNP Q9WYB7
A	-15	HIS	-	EXPRESSION TAG	UNP Q9WYB7
A	-14	HIS	-	EXPRESSION TAG	UNP Q9WYB7
A	-13	HIS	-	EXPRESSION TAG	UNP Q9WYB7
A	-12	HIS	-	EXPRESSION TAG	UNP Q9WYB7
A	-11	HIS	-	EXPRESSION TAG	UNP Q9WYB7
A	-10	HIS	-	EXPRESSION TAG	UNP Q9WYB7
A	-9	SER	-	EXPRESSION TAG	UNP Q9WYB7
A	-8	SER	-	EXPRESSION TAG	UNP Q9WYB7
A	-7	GLY	-	EXPRESSION TAG	UNP Q9WYB7
A	-6	LEU	-	EXPRESSION TAG	UNP Q9WYB7
A	-5	VAL	-	EXPRESSION TAG	UNP Q9WYB7
A	-4	PRO	-	EXPRESSION TAG	UNP Q9WYB7
A	-3	ARG	-	EXPRESSION TAG	UNP Q9WYB7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP Q9WYB7
A	-1	SER	-	EXPRESSION TAG	UNP Q9WYB7
A	0	HIS	-	EXPRESSION TAG	UNP Q9WYB7
A	4	GLY	ARG	ENGINEERED MUTATION	UNP Q9WYB7
B	-19	MET	-	EXPRESSION TAG	UNP Q9WYB7
B	-18	GLY	-	EXPRESSION TAG	UNP Q9WYB7
B	-17	SER	-	EXPRESSION TAG	UNP Q9WYB7
B	-16	SER	-	EXPRESSION TAG	UNP Q9WYB7
B	-15	HIS	-	EXPRESSION TAG	UNP Q9WYB7
B	-14	HIS	-	EXPRESSION TAG	UNP Q9WYB7
B	-13	HIS	-	EXPRESSION TAG	UNP Q9WYB7
B	-12	HIS	-	EXPRESSION TAG	UNP Q9WYB7
B	-11	HIS	-	EXPRESSION TAG	UNP Q9WYB7
B	-10	HIS	-	EXPRESSION TAG	UNP Q9WYB7
B	-9	SER	-	EXPRESSION TAG	UNP Q9WYB7
B	-8	SER	-	EXPRESSION TAG	UNP Q9WYB7
B	-7	GLY	-	EXPRESSION TAG	UNP Q9WYB7
B	-6	LEU	-	EXPRESSION TAG	UNP Q9WYB7
B	-5	VAL	-	EXPRESSION TAG	UNP Q9WYB7
B	-4	PRO	-	EXPRESSION TAG	UNP Q9WYB7
B	-3	ARG	-	EXPRESSION TAG	UNP Q9WYB7
B	-2	GLY	-	EXPRESSION TAG	UNP Q9WYB7
B	-1	SER	-	EXPRESSION TAG	UNP Q9WYB7
B	0	HIS	-	EXPRESSION TAG	UNP Q9WYB7
B	4	GLY	ARG	ENGINEERED MUTATION	UNP Q9WYB7
C	-19	MET	-	EXPRESSION TAG	UNP Q9WYB7
C	-18	GLY	-	EXPRESSION TAG	UNP Q9WYB7
C	-17	SER	-	EXPRESSION TAG	UNP Q9WYB7
C	-16	SER	-	EXPRESSION TAG	UNP Q9WYB7
C	-15	HIS	-	EXPRESSION TAG	UNP Q9WYB7
C	-14	HIS	-	EXPRESSION TAG	UNP Q9WYB7
C	-13	HIS	-	EXPRESSION TAG	UNP Q9WYB7
C	-12	HIS	-	EXPRESSION TAG	UNP Q9WYB7
C	-11	HIS	-	EXPRESSION TAG	UNP Q9WYB7
C	-10	HIS	-	EXPRESSION TAG	UNP Q9WYB7
C	-9	SER	-	EXPRESSION TAG	UNP Q9WYB7
C	-8	SER	-	EXPRESSION TAG	UNP Q9WYB7
C	-7	GLY	-	EXPRESSION TAG	UNP Q9WYB7
C	-6	LEU	-	EXPRESSION TAG	UNP Q9WYB7
C	-5	VAL	-	EXPRESSION TAG	UNP Q9WYB7
C	-4	PRO	-	EXPRESSION TAG	UNP Q9WYB7
C	-3	ARG	-	EXPRESSION TAG	UNP Q9WYB7

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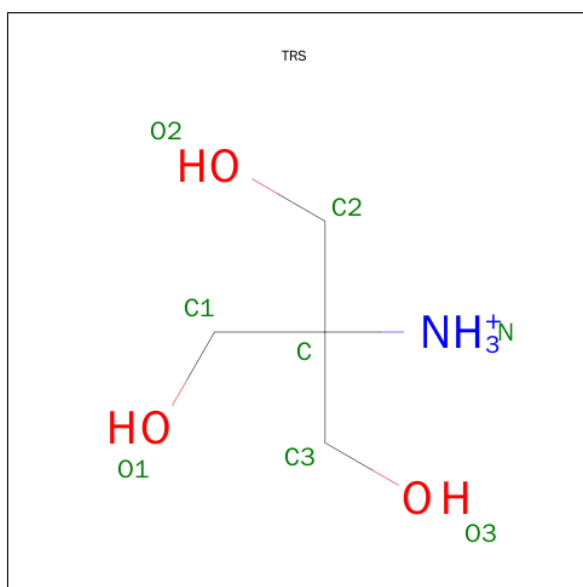
Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	GLY	-	EXPRESSION TAG	UNP Q9WYB7
C	-1	SER	-	EXPRESSION TAG	UNP Q9WYB7
C	0	HIS	-	EXPRESSION TAG	UNP Q9WYB7
C	4	GLY	ARG	ENGINEERED MUTATION	UNP Q9WYB7
D	-19	MET	-	EXPRESSION TAG	UNP Q9WYB7
D	-18	GLY	-	EXPRESSION TAG	UNP Q9WYB7
D	-17	SER	-	EXPRESSION TAG	UNP Q9WYB7
D	-16	SER	-	EXPRESSION TAG	UNP Q9WYB7
D	-15	HIS	-	EXPRESSION TAG	UNP Q9WYB7
D	-14	HIS	-	EXPRESSION TAG	UNP Q9WYB7
D	-13	HIS	-	EXPRESSION TAG	UNP Q9WYB7
D	-12	HIS	-	EXPRESSION TAG	UNP Q9WYB7
D	-11	HIS	-	EXPRESSION TAG	UNP Q9WYB7
D	-10	HIS	-	EXPRESSION TAG	UNP Q9WYB7
D	-9	SER	-	EXPRESSION TAG	UNP Q9WYB7
D	-8	SER	-	EXPRESSION TAG	UNP Q9WYB7
D	-7	GLY	-	EXPRESSION TAG	UNP Q9WYB7
D	-6	LEU	-	EXPRESSION TAG	UNP Q9WYB7
D	-5	VAL	-	EXPRESSION TAG	UNP Q9WYB7
D	-4	PRO	-	EXPRESSION TAG	UNP Q9WYB7
D	-3	ARG	-	EXPRESSION TAG	UNP Q9WYB7
D	-2	GLY	-	EXPRESSION TAG	UNP Q9WYB7
D	-1	SER	-	EXPRESSION TAG	UNP Q9WYB7
D	0	HIS	-	EXPRESSION TAG	UNP Q9WYB7
D	4	GLY	ARG	ENGINEERED MUTATION	UNP Q9WYB7
E	-19	MET	-	EXPRESSION TAG	UNP Q9WYB7
E	-18	GLY	-	EXPRESSION TAG	UNP Q9WYB7
E	-17	SER	-	EXPRESSION TAG	UNP Q9WYB7
E	-16	SER	-	EXPRESSION TAG	UNP Q9WYB7
E	-15	HIS	-	EXPRESSION TAG	UNP Q9WYB7
E	-14	HIS	-	EXPRESSION TAG	UNP Q9WYB7
E	-13	HIS	-	EXPRESSION TAG	UNP Q9WYB7
E	-12	HIS	-	EXPRESSION TAG	UNP Q9WYB7
E	-11	HIS	-	EXPRESSION TAG	UNP Q9WYB7
E	-10	HIS	-	EXPRESSION TAG	UNP Q9WYB7
E	-9	SER	-	EXPRESSION TAG	UNP Q9WYB7
E	-8	SER	-	EXPRESSION TAG	UNP Q9WYB7
E	-7	GLY	-	EXPRESSION TAG	UNP Q9WYB7
E	-6	LEU	-	EXPRESSION TAG	UNP Q9WYB7
E	-5	VAL	-	EXPRESSION TAG	UNP Q9WYB7
E	-4	PRO	-	EXPRESSION TAG	UNP Q9WYB7
E	-3	ARG	-	EXPRESSION TAG	UNP Q9WYB7

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	GLY	-	EXPRESSION TAG	UNP Q9WYB7
E	-1	SER	-	EXPRESSION TAG	UNP Q9WYB7
E	0	HIS	-	EXPRESSION TAG	UNP Q9WYB7
E	4	GLY	ARG	ENGINEERED MUTATION	UNP Q9WYB7
F	-19	MET	-	EXPRESSION TAG	UNP Q9WYB7
F	-18	GLY	-	EXPRESSION TAG	UNP Q9WYB7
F	-17	SER	-	EXPRESSION TAG	UNP Q9WYB7
F	-16	SER	-	EXPRESSION TAG	UNP Q9WYB7
F	-15	HIS	-	EXPRESSION TAG	UNP Q9WYB7
F	-14	HIS	-	EXPRESSION TAG	UNP Q9WYB7
F	-13	HIS	-	EXPRESSION TAG	UNP Q9WYB7
F	-12	HIS	-	EXPRESSION TAG	UNP Q9WYB7
F	-11	HIS	-	EXPRESSION TAG	UNP Q9WYB7
F	-10	HIS	-	EXPRESSION TAG	UNP Q9WYB7
F	-9	SER	-	EXPRESSION TAG	UNP Q9WYB7
F	-8	SER	-	EXPRESSION TAG	UNP Q9WYB7
F	-7	GLY	-	EXPRESSION TAG	UNP Q9WYB7
F	-6	LEU	-	EXPRESSION TAG	UNP Q9WYB7
F	-5	VAL	-	EXPRESSION TAG	UNP Q9WYB7
F	-4	PRO	-	EXPRESSION TAG	UNP Q9WYB7
F	-3	ARG	-	EXPRESSION TAG	UNP Q9WYB7
F	-2	GLY	-	EXPRESSION TAG	UNP Q9WYB7
F	-1	SER	-	EXPRESSION TAG	UNP Q9WYB7
F	0	HIS	-	EXPRESSION TAG	UNP Q9WYB7
F	4	GLY	ARG	ENGINEERED MUTATION	UNP Q9WYB7

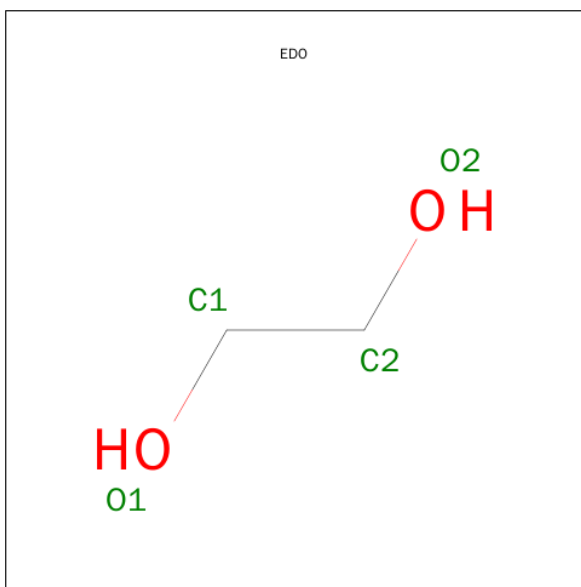
- Molecule 2 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: C<sub>4</sub>H<sub>12</sub>NO<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			8	4	1	3		
2	B	1	Total	C	N	O	0	0
			8	4	1	3		
2	C	1	Total	C	N	O	0	0
			8	4	1	3		
2	D	1	Total	C	N	O	0	0
			8	4	1	3		
2	E	1	Total	C	N	O	0	0
			8	4	1	3		
2	F	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).





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

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

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

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Na 1 1	0	0

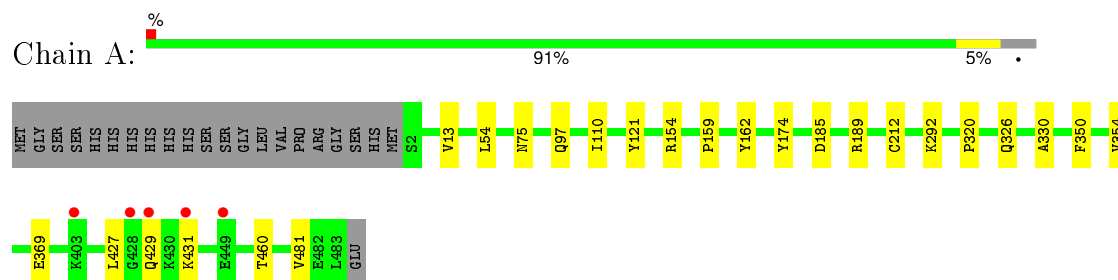
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	449	Total 449	O 449	0	0
5	B	327	Total 327	O 327	0	0
5	C	454	Total 454	O 454	0	0
5	D	391	Total 391	O 391	0	0
5	E	386	Total 386	O 386	0	0
5	F	417	Total 417	O 417	0	0

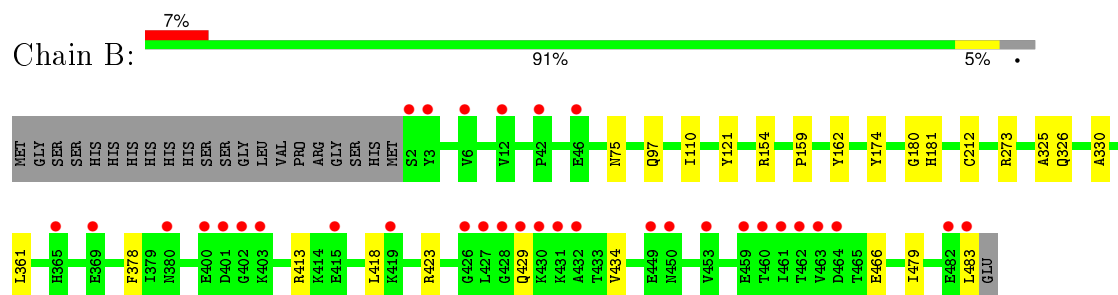
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

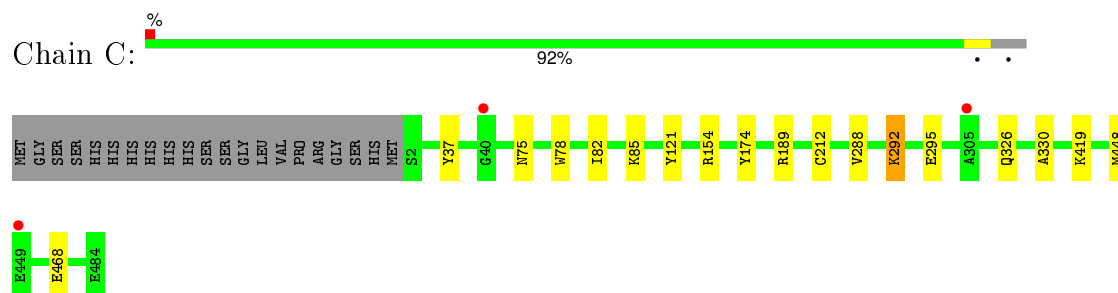
- Molecule 1: Alpha-L-arabinofuranosidase



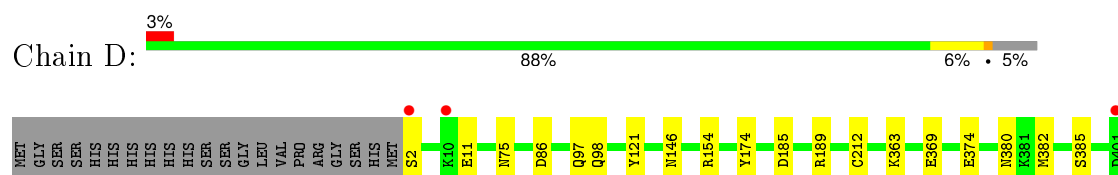
- Molecule 1: Alpha-L-arabinofuranosidase



- Molecule 1: Alpha-L-arabinofuranosidase

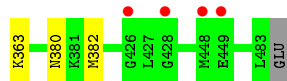
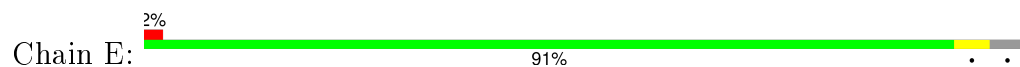


- Molecule 1: Alpha-L-arabinofuranosidase

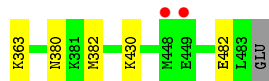
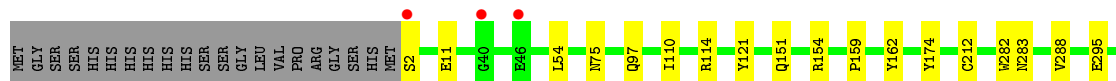
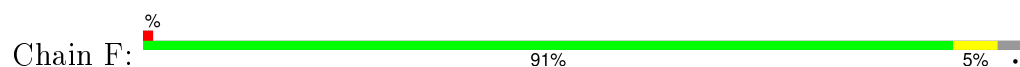




● Molecule 1: Alpha-L-arabinofuranosidase



● Molecule 1: Alpha-L-arabinofuranosidase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	101.52Å 160.65Å 155.74Å 90.00° 91.75° 90.00°	Depositor
Resolution (Å)	30.14 – 1.80 30.13 – 1.80	Depositor EDS
% Data completeness (in resolution range)	97.0 (30.14-1.80) 97.0 (30.13-1.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.02 (at 1.80Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.175 , 0.197 0.175 , 0.198	Depositor DCC
$R_{free}$ test set	22344 reflections (5.28%)	DCC
Wilson B-factor (Å <sup>2</sup> )	22.8	Xtriage
Anisotropy	0.048	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 47.8	EDS
Estimated twinning fraction	0.008 for -h,-l,-k 0.005 for -h,l,k 0.019 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 445228 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	25934	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: NA, TRS, 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	0.64	0/3973	0.67	1/5392 (0.0%)
1	B	0.62	0/3973	0.66	0/5392
1	C	0.64	0/3982	0.67	0/5404
1	D	0.64	0/3965	0.67	2/5381 (0.0%)
1	E	0.63	0/3973	0.65	0/5392
1	F	0.62	0/3973	0.67	0/5392
All	All	0.63	0/23839	0.66	3/32353 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	154	ARG	NE-CZ-NH2	6.36	123.48	120.30
1	A	185	ASP	CB-CG-OD1	5.24	123.02	118.30
1	D	86	ASP	CB-CG-OD1	5.02	122.82	118.30

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	3876	0	3793	12	0
1	B	3876	0	3793	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	3885	0	3799	9	0
1	D	3868	0	3782	17	0
1	E	3876	0	3793	13	0
1	F	3876	0	3793	10	0
2	A	8	0	12	0	0
2	B	8	0	12	0	0
2	C	8	0	12	0	0
2	D	8	0	12	0	0
2	E	8	0	12	0	0
2	F	8	0	12	0	0
3	A	40	0	60	2	0
3	B	32	0	48	2	0
3	C	32	0	48	3	0
3	D	36	0	54	1	0
3	E	32	0	48	0	0
3	F	32	0	48	1	0
4	A	1	0	0	0	0
5	A	449	0	0	5	0
5	B	327	0	0	0	0
5	C	454	0	0	2	0
5	D	391	0	0	1	0
5	E	386	0	0	3	0
5	F	417	0	0	2	0
All	All	25934	0	23131	77	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 2.

All (77) 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:D:185:ASP:HB3	5:D:1888:HOH:O	1.58	1.03
1:E:11:GLU:HG2	1:E:363:LYS:HB2	1.57	0.85
1:E:185:ASP:HB2	5:E:504:HOH:O	1.78	0.83
1:A:154:ARG:HG2	1:A:159:PRO:HA	1.71	0.73
1:C:292:LYS:HE2	5:C:586:HOH:O	1.89	0.72
1:D:11:GLU:HG2	1:D:363:LYS:HB2	1.71	0.72
1:E:154:ARG:HG2	1:E:159:PRO:HA	1.73	0.70
1:D:380:ASN:OD1	1:D:382:MET:HG3	1.94	0.68
1:F:380:ASN:OD1	1:F:382:MET:HG3	1.95	0.67
1:D:413:ARG:NH2	1:D:416:ASP:OD1	2.30	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:350:PHE:O	1:A:354:VAL:HG22	2.00	0.62
1:C:419:LYS:HE2	1:C:468:GLU:OE1	2.01	0.60
1:E:146:ASN:HB3	5:E:524:HOH:O	2.01	0.60
1:E:151:GLN:OE1	1:E:154:ARG:NH1	2.36	0.58
1:A:97:GLN:NE2	5:A:945:HOH:O	2.34	0.58
1:D:419:LYS:HE3	1:D:468:GLU:OE1	2.07	0.55
1:D:419:LYS:HG2	1:D:470:THR:OG1	2.08	0.53
1:E:380:ASN:HD21	1:E:382:MET:CE	2.23	0.52
1:D:411:ASN:OD1	1:D:413:ARG:HB2	2.10	0.52
1:F:430:LYS:HD2	1:F:482:GLU:O	2.09	0.52
1:A:292:LYS:HE2	5:A:2367:HOH:O	2.10	0.50
1:B:181:HIS:O	3:C:489:EDO:H22	2.10	0.50
1:B:180:GLY:HA2	3:C:489:EDO:H21	1.93	0.50
1:A:369:GLU:HG3	5:A:600:HOH:O	2.11	0.50
1:E:380:ASN:OD1	1:E:382:MET:HG3	2.12	0.50
1:B:434:VAL:HG22	1:B:479:ILE:HG12	1.93	0.50
1:F:110:ILE:HG21	1:F:162:TYR:CD1	2.48	0.48
1:E:341:LEU:HD12	1:E:341:LEU:C	2.34	0.48
1:F:288:VAL:HG22	1:F:295:GLU:OE1	2.14	0.48
1:A:431:LYS:HD2	1:A:460:THR:HG21	1.96	0.48
1:B:174:TYR:CE2	1:B:212:CYS:HB3	2.49	0.48
1:B:326:GLN:HB2	1:B:330:ALA:O	2.14	0.47
1:C:326:GLN:HB2	1:C:330:ALA:O	2.15	0.47
1:B:423:ARG:NH1	1:B:466:GLU:OE2	2.46	0.47
1:D:189:ARG:HH21	3:D:488:EDO:H11	1.78	0.47
1:F:154:ARG:HG2	1:F:159:PRO:HA	1.97	0.47
1:D:174:TYR:CE2	1:D:212:CYS:HB3	2.50	0.46
1:F:174:TYR:CE2	1:F:212:CYS:HB3	2.51	0.46
1:A:189:ARG:HG3	5:A:2124:HOH:O	2.15	0.46
1:E:174:TYR:CE2	1:E:212:CYS:HB3	2.50	0.46
1:C:189:ARG:NE	3:C:489:EDO:O1	2.49	0.45
1:F:11:GLU:HG2	1:F:363:LYS:HB2	1.97	0.45
3:F:489:EDO:H12	5:F:591:HOH:O	2.16	0.45
1:A:427:LEU:HD11	1:A:481:VAL:HG11	1.99	0.45
1:D:481:VAL:HG12	1:D:482:GLU:HB2	1.98	0.45
1:D:11:GLU:HG2	1:D:363:LYS:CB	2.45	0.45
1:E:110:ILE:HG21	1:E:162:TYR:CD1	2.53	0.44
1:B:413:ARG:HD3	1:B:418:LEU:HD21	2.00	0.44
1:D:374:GLU:HG2	1:D:385:SER:HB3	1.99	0.44
1:A:110:ILE:HG21	1:A:162:TYR:CD1	2.53	0.44
1:E:46:GLU:H	1:E:46:GLU:CD	2.22	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:288:VAL:HG22	1:C:295:GLU:OE2	2.18	0.43
1:D:2:SER:HB2	1:D:369:GLU:OE2	2.18	0.43
1:E:295:GLU:HG2	1:E:297:TYR:CE1	2.54	0.43
1:F:282:TRP:O	1:F:283:ASN:HB2	2.17	0.43
1:D:434:VAL:HG22	1:D:479:ILE:HG12	2.00	0.43
1:B:413:ARG:CD	1:B:418:LEU:HD21	2.49	0.43
1:F:110:ILE:O	1:F:114:ARG:HG2	2.19	0.42
1:D:435:TYR:O	1:D:477:SER:HA	2.18	0.42
1:B:154:ARG:HG2	1:B:159:PRO:HA	2.00	0.42
1:B:361:LEU:HD21	3:B:489:EDO:H22	2.01	0.42
1:F:151:GLN:HG2	5:F:501:HOH:O	2.19	0.42
1:B:378:PHE:CD2	3:B:488:EDO:H11	2.54	0.42
1:C:154:ARG:HB3	5:C:2171:HOH:O	2.19	0.42
1:D:146:ASN:HB2	5:E:1152:HOH:O	2.20	0.42
1:A:13:VAL:HG21	3:A:490:EDO:H11	2.02	0.42
1:C:174:TYR:CE2	1:C:212:CYS:HB3	2.55	0.42
1:C:37:TYR:OH	1:C:85:LYS:NZ	2.50	0.41
1:C:78:TRP:CZ3	1:C:82:ILE:HD13	2.56	0.41
1:B:110:ILE:HG21	1:B:162:TYR:CD1	2.55	0.41
3:A:489:EDO:H12	5:A:575:HOH:O	2.19	0.41
1:A:326:GLN:HB2	1:A:330:ALA:O	2.21	0.41
1:B:273:ARG:HH11	1:B:273:ARG:HG3	1.85	0.41
1:A:174:TYR:CE2	1:A:212:CYS:HB3	2.54	0.41
1:D:427:LEU:CD2	1:D:430:LYS:HD2	2.51	0.41
1:B:325:ALA:HA	1:B:326:GLN:HA	1.89	0.41
1:E:11:GLU:CG	1:E:363:LYS:HB2	2.38	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	480/504 (95%)	464 (97%)	16 (3%)	0	100	100
1	B	480/504 (95%)	462 (96%)	18 (4%)	0	100	100
1	C	481/504 (95%)	463 (96%)	18 (4%)	0	100	100
1	D	479/504 (95%)	459 (96%)	20 (4%)	0	100	100
1	E	480/504 (95%)	460 (96%)	20 (4%)	0	100	100
1	F	480/504 (95%)	461 (96%)	19 (4%)	0	100	100
All	All	2880/3024 (95%)	2769 (96%)	111 (4%)	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	417/436 (96%)	412 (99%)	5 (1%)	78	71
1	B	417/436 (96%)	412 (99%)	5 (1%)	78	71
1	C	418/436 (96%)	414 (99%)	4 (1%)	82	77
1	D	416/436 (95%)	407 (98%)	9 (2%)	60	45
1	E	417/436 (96%)	412 (99%)	5 (1%)	78	71
1	F	417/436 (96%)	412 (99%)	5 (1%)	78	71
All	All	2502/2616 (96%)	2469 (99%)	33 (1%)	76	68

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	LEU
1	A	75	ASN
1	A	121	TYR
1	A	320	PRO
1	A	429	GLN
1	B	75	ASN
1	B	97	GLN

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Mol	Chain	Res	Type
1	B	121	TYR
1	B	429	GLN
1	B	483	LEU
1	C	75	ASN
1	C	121	TYR
1	C	292	LYS
1	C	448	MET
1	D	75	ASN
1	D	97	GLN
1	D	98	GLN
1	D	121	TYR
1	D	404	LYS
1	D	419	LYS
1	D	427	LEU
1	D	430	LYS
1	D	462	THR
1	E	11	GLU
1	E	54	LEU
1	E	75	ASN
1	E	98	GLN
1	E	121	TYR
1	F	2	SER
1	F	54	LEU
1	F	75	ASN
1	F	97	GLN
1	F	121	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	98	GLN
1	E	380	ASN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 58 ligands modelled in this entry, 1 is monoatomic - leaving 57 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	TRS	A	485	-	7,7,7	0.72	0	9,9,9	1.02	0
3	EDO	A	486	-	3,3,3	0.39	0	2,2,2	0.25	0
3	EDO	A	487	-	3,3,3	0.56	0	2,2,2	0.58	0
3	EDO	A	488	-	3,3,3	0.45	0	2,2,2	0.86	0
3	EDO	A	489	-	3,3,3	0.48	0	2,2,2	0.31	0
3	EDO	A	490	-	3,3,3	0.39	0	2,2,2	0.71	0
3	EDO	A	491	-	3,3,3	0.55	0	2,2,2	0.33	0
3	EDO	A	492	-	3,3,3	0.39	0	2,2,2	0.68	0
3	EDO	A	493	-	3,3,3	0.45	0	2,2,2	0.61	0
3	EDO	A	494	-	3,3,3	0.58	0	2,2,2	0.22	0
3	EDO	A	495	-	3,3,3	0.50	0	2,2,2	0.47	0
2	TRS	B	485	-	7,7,7	0.95	1 (14%)	9,9,9	0.86	0
3	EDO	B	486	-	3,3,3	0.30	0	2,2,2	0.19	0
3	EDO	B	487	-	3,3,3	0.57	0	2,2,2	0.55	0
3	EDO	B	488	-	3,3,3	0.46	0	2,2,2	0.43	0
3	EDO	B	489	-	3,3,3	0.44	0	2,2,2	0.45	0
3	EDO	B	490	-	3,3,3	0.47	0	2,2,2	0.70	0
3	EDO	B	491	-	3,3,3	0.48	0	2,2,2	0.50	0
3	EDO	B	492	-	3,3,3	0.50	0	2,2,2	0.37	0
3	EDO	B	493	-	3,3,3	0.51	0	2,2,2	0.29	0
2	TRS	C	485	-	7,7,7	0.47	0	9,9,9	1.22	1 (11%)
3	EDO	C	486	-	3,3,3	0.50	0	2,2,2	0.25	0
3	EDO	C	487	-	3,3,3	0.54	0	2,2,2	0.53	0
3	EDO	C	488	-	3,3,3	0.40	0	2,2,2	0.68	0
3	EDO	C	489	-	3,3,3	0.44	0	2,2,2	0.55	0
3	EDO	C	490	-	3,3,3	0.44	0	2,2,2	0.53	0
3	EDO	C	491	-	3,3,3	0.48	0	2,2,2	0.37	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	EDO	C	492	-	3,3,3	0.35	0	2,2,2	0.75	0
3	EDO	C	493	-	3,3,3	0.50	0	2,2,2	0.51	0
3	EDO	D	485	-	3,3,3	0.35	0	2,2,2	0.38	0
3	EDO	D	486	-	3,3,3	0.60	0	2,2,2	0.52	0
3	EDO	D	487	-	3,3,3	0.58	0	2,2,2	0.49	0
3	EDO	D	488	-	3,3,3	0.46	0	2,2,2	0.52	0
3	EDO	D	489	-	3,3,3	0.48	0	2,2,2	0.37	0
3	EDO	D	490	-	3,3,3	0.41	0	2,2,2	0.72	0
3	EDO	D	491	-	3,3,3	0.37	0	2,2,2	0.62	0
3	EDO	D	492	-	3,3,3	0.42	0	2,2,2	0.67	0
3	EDO	D	493	-	3,3,3	0.53	0	2,2,2	0.42	0
2	TRS	D	494	-	7,7,7	0.68	0	9,9,9	1.12	1 (11%)
2	TRS	E	485	-	7,7,7	0.77	0	9,9,9	1.09	1 (11%)
3	EDO	E	486	-	3,3,3	0.50	0	2,2,2	0.25	0
3	EDO	E	487	-	3,3,3	0.47	0	2,2,2	0.64	0
3	EDO	E	488	-	3,3,3	0.45	0	2,2,2	0.43	0
3	EDO	E	489	-	3,3,3	0.48	0	2,2,2	0.41	0
3	EDO	E	490	-	3,3,3	0.52	0	2,2,2	0.31	0
3	EDO	E	491	-	3,3,3	0.38	0	2,2,2	0.68	0
3	EDO	E	492	-	3,3,3	0.41	0	2,2,2	0.64	0
3	EDO	E	493	-	3,3,3	0.47	0	2,2,2	0.36	0
2	TRS	F	485	-	7,7,7	0.71	0	9,9,9	0.96	1 (11%)
3	EDO	F	486	-	3,3,3	0.57	0	2,2,2	0.31	0
3	EDO	F	487	-	3,3,3	0.44	0	2,2,2	0.74	0
3	EDO	F	488	-	3,3,3	0.50	0	2,2,2	0.50	0
3	EDO	F	489	-	3,3,3	0.50	0	2,2,2	0.32	0
3	EDO	F	490	-	3,3,3	0.53	0	2,2,2	0.26	0
3	EDO	F	491	-	3,3,3	0.42	0	2,2,2	0.50	0
3	EDO	F	492	-	3,3,3	0.38	0	2,2,2	0.69	0
3	EDO	F	493	-	3,3,3	0.49	0	2,2,2	0.35	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	TRS	A	485	-	-	0/9/9/9	0/0/0/0
3	EDO	A	486	-	-	0/1/1/1	0/0/0/0
3	EDO	A	487	-	-	0/1/1/1	0/0/0/0
3	EDO	A	488	-	-	0/1/1/1	0/0/0/0
3	EDO	A	489	-	-	0/1/1/1	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	A	490	-	-	0/1/1/1	0/0/0/0
3	EDO	A	491	-	-	0/1/1/1	0/0/0/0
3	EDO	A	492	-	-	0/1/1/1	0/0/0/0
3	EDO	A	493	-	-	0/1/1/1	0/0/0/0
3	EDO	A	494	-	-	0/1/1/1	0/0/0/0
3	EDO	A	495	-	-	0/1/1/1	0/0/0/0
2	TRS	B	485	-	-	0/9/9/9	0/0/0/0
3	EDO	B	486	-	-	0/1/1/1	0/0/0/0
3	EDO	B	487	-	-	0/1/1/1	0/0/0/0
3	EDO	B	488	-	-	0/1/1/1	0/0/0/0
3	EDO	B	489	-	-	0/1/1/1	0/0/0/0
3	EDO	B	490	-	-	0/1/1/1	0/0/0/0
3	EDO	B	491	-	-	0/1/1/1	0/0/0/0
3	EDO	B	492	-	-	0/1/1/1	0/0/0/0
3	EDO	B	493	-	-	0/1/1/1	0/0/0/0
2	TRS	C	485	-	-	0/9/9/9	0/0/0/0
3	EDO	C	486	-	-	0/1/1/1	0/0/0/0
3	EDO	C	487	-	-	0/1/1/1	0/0/0/0
3	EDO	C	488	-	-	0/1/1/1	0/0/0/0
3	EDO	C	489	-	-	0/1/1/1	0/0/0/0
3	EDO	C	490	-	-	0/1/1/1	0/0/0/0
3	EDO	C	491	-	-	0/1/1/1	0/0/0/0
3	EDO	C	492	-	-	0/1/1/1	0/0/0/0
3	EDO	C	493	-	-	0/1/1/1	0/0/0/0
3	EDO	D	485	-	-	0/1/1/1	0/0/0/0
3	EDO	D	486	-	-	0/1/1/1	0/0/0/0
3	EDO	D	487	-	-	0/1/1/1	0/0/0/0
3	EDO	D	488	-	-	0/1/1/1	0/0/0/0
3	EDO	D	489	-	-	0/1/1/1	0/0/0/0
3	EDO	D	490	-	-	0/1/1/1	0/0/0/0
3	EDO	D	491	-	-	0/1/1/1	0/0/0/0
3	EDO	D	492	-	-	0/1/1/1	0/0/0/0
3	EDO	D	493	-	-	0/1/1/1	0/0/0/0
2	TRS	D	494	-	-	0/9/9/9	0/0/0/0
2	TRS	E	485	-	-	0/9/9/9	0/0/0/0
3	EDO	E	486	-	-	0/1/1/1	0/0/0/0
3	EDO	E	487	-	-	0/1/1/1	0/0/0/0
3	EDO	E	488	-	-	0/1/1/1	0/0/0/0
3	EDO	E	489	-	-	0/1/1/1	0/0/0/0
3	EDO	E	490	-	-	0/1/1/1	0/0/0/0
3	EDO	E	491	-	-	0/1/1/1	0/0/0/0
3	EDO	E	492	-	-	0/1/1/1	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	E	493	-	-	0/1/1/1	0/0/0/0
2	TRS	F	485	-	-	0/9/9/9	0/0/0/0
3	EDO	F	486	-	-	0/1/1/1	0/0/0/0
3	EDO	F	487	-	-	0/1/1/1	0/0/0/0
3	EDO	F	488	-	-	0/1/1/1	0/0/0/0
3	EDO	F	489	-	-	0/1/1/1	0/0/0/0
3	EDO	F	490	-	-	0/1/1/1	0/0/0/0
3	EDO	F	491	-	-	0/1/1/1	0/0/0/0
3	EDO	F	492	-	-	0/1/1/1	0/0/0/0
3	EDO	F	493	-	-	0/1/1/1	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	485	TRS	C-N	-2.38	1.47	1.50

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	485	TRS	C2-C-C1	-2.39	105.61	110.78
2	D	494	TRS	O1-C1-C	-2.35	106.43	111.18
2	F	485	TRS	O1-C1-C	-2.27	106.59	111.18
2	C	485	TRS	O3-C3-C	-2.26	106.61	111.18

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	489	EDO	1	0
3	A	490	EDO	1	0
3	B	488	EDO	1	0
3	B	489	EDO	1	0
3	C	489	EDO	3	0
3	D	488	EDO	1	0
3	F	489	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	482/504 (95%)	-0.15	5 (1%) 84 82	13, 21, 37, 46	0
1	B	482/504 (95%)	0.12	33 (6%) 20 16	14, 24, 39, 48	0
1	C	483/504 (95%)	-0.22	3 (0%) 90 88	13, 21, 34, 46	0
1	D	481/504 (95%)	-0.06	17 (3%) 48 42	14, 23, 38, 47	0
1	E	482/504 (95%)	-0.14	8 (1%) 73 69	15, 23, 36, 47	0
1	F	482/504 (95%)	-0.18	5 (1%) 84 82	15, 22, 36, 47	0
All	All	2892/3024 (95%)	-0.10	71 (2%) 61 56	13, 22, 37, 48	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	431	LYS	5.3
1	B	429	GLN	5.1
1	D	429	GLN	4.6
1	D	449	GLU	4.6
1	D	428	GLY	4.6
1	B	431	LYS	4.6
1	E	449	GLU	4.4
1	B	428	GLY	4.4
1	B	483	LEU	4.3
1	B	449	GLU	4.3
1	B	462	THR	4.1
1	F	449	GLU	4.0
1	D	462	THR	3.7
1	C	449	GLU	3.7
1	B	403	LYS	3.4
1	B	2	SER	3.4
1	B	460	THR	3.4
1	B	415	GLU	3.4
1	D	463	VAL	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	426	GLY	3.4
1	A	449	GLU	3.3
1	F	46	GLU	3.3
1	B	463	VAL	3.3
1	B	402	GLY	3.2
1	B	427	LEU	3.1
1	A	429	GLN	3.1
1	E	428	GLY	3.1
1	D	403	LYS	3.0
1	D	460	THR	2.9
1	B	464	ASP	2.9
1	A	403	LYS	2.9
1	D	419	LYS	2.9
1	E	426	GLY	2.8
1	B	450	ASN	2.8
1	B	401	ASP	2.7
1	D	10	LYS	2.7
1	D	402	GLY	2.7
1	E	10	LYS	2.6
1	E	46	GLU	2.6
1	D	464	ASP	2.6
1	B	46	GLU	2.5
1	B	369	GLU	2.5
1	B	430	LYS	2.5
1	B	453	VAL	2.5
1	B	365	HIS	2.5
1	D	426	GLY	2.4
1	F	40	GLY	2.4
1	C	305	ALA	2.4
1	B	419	LYS	2.4
1	B	6	VAL	2.4
1	A	431	LYS	2.4
1	F	2	SER	2.4
1	B	380	ASN	2.3
1	E	11	GLU	2.3
1	B	3	TYR	2.3
1	E	448	MET	2.3
1	B	461	ILE	2.2
1	D	427	LEU	2.2
1	B	12	VAL	2.2
1	B	400	GLU	2.2
1	D	2	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	432	ALA	2.2
1	C	40	GLY	2.2
1	F	448	MET	2.2
1	B	42	PRO	2.1
1	B	482	GLU	2.1
1	D	401	ASP	2.1
1	A	428	GLY	2.1
1	E	2	SER	2.1
1	B	459	GLU	2.1
1	D	467	PHE	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	EDO	C	493	4/4	0.68	0.29	16.48	53,56,57,57	0
3	EDO	B	493	4/4	0.66	0.32	11.27	63,63,64,64	0
4	NA	A	496	1/1	0.98	0.26	5.49	40,40,40,40	0
3	EDO	D	493	4/4	0.84	0.15	4.76	41,44,47,49	0
3	EDO	A	494	4/4	0.87	0.11	2.89	38,43,43,44	0
3	EDO	B	488	4/4	0.92	0.14	1.51	34,36,38,39	0
3	EDO	E	491	4/4	0.95	0.21	1.39	50,50,50,51	0
3	EDO	C	488	4/4	0.95	0.11	1.38	39,40,40,41	0
3	EDO	B	486	4/4	0.97	0.13	0.90	26,28,28,31	0
3	EDO	E	487	4/4	0.98	0.10	0.86	25,28,28,31	0
3	EDO	F	488	4/4	0.94	0.10	0.73	29,34,35,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	EDO	D	485	4/4	0.98	0.14	0.69	23,26,28,28	0
2	TRS	C	485	8/8	0.97	0.07	0.59	16,17,17,18	0
3	EDO	A	490	4/4	0.94	0.12	0.47	40,42,44,44	0
2	TRS	E	485	8/8	0.96	0.08	0.44	23,23,24,24	0
3	EDO	C	486	4/4	0.96	0.10	0.37	20,24,26,27	0
3	EDO	F	490	4/4	0.93	0.12	0.36	34,36,37,39	0
2	TRS	D	494	8/8	0.97	0.08	0.35	22,23,23,23	0
3	EDO	F	487	4/4	0.97	0.08	0.30	25,27,29,29	0
3	EDO	B	487	4/4	0.95	0.08	0.27	28,31,31,32	0
3	EDO	D	487	4/4	0.94	0.11	0.27	34,35,37,37	0
3	EDO	D	486	4/4	0.97	0.09	0.25	25,25,27,28	0
3	EDO	A	486	4/4	0.98	0.11	0.24	18,23,24,24	0
3	EDO	C	490	4/4	0.94	0.10	0.14	35,38,39,43	0
3	EDO	D	489	4/4	0.94	0.10	0.14	41,44,44,46	0
3	EDO	E	490	4/4	0.94	0.13	0.03	42,44,44,45	0
3	EDO	C	487	4/4	0.97	0.07	-0.06	28,31,31,32	0
2	TRS	F	485	8/8	0.98	0.07	-0.08	23,24,24,25	0
3	EDO	E	486	4/4	0.98	0.09	-0.23	23,23,25,26	0
2	TRS	B	485	8/8	0.96	0.07	-0.27	20,21,21,22	0
2	TRS	A	485	8/8	0.96	0.08	-0.35	22,22,22,23	0
3	EDO	F	486	4/4	0.97	0.09	-0.36	22,24,26,26	0
3	EDO	B	489	4/4	0.95	0.11	-0.56	50,50,50,51	0
3	EDO	E	488	4/4	0.96	0.09	-0.77	30,33,34,35	0
3	EDO	A	487	4/4	0.97	0.06	-1.02	22,26,26,29	0
3	EDO	A	488	4/4	0.96	0.07	-1.17	26,30,31,31	0
3	EDO	F	491	4/4	0.92	0.17	-	42,43,43,46	0
3	EDO	B	490	4/4	0.82	0.23	-	49,51,52,53	0
3	EDO	D	491	4/4	0.91	0.14	-	57,57,57,57	0
3	EDO	D	490	4/4	0.88	0.17	-	45,46,47,51	0
3	EDO	C	489	4/4	0.93	0.19	-	37,38,40,42	0
3	EDO	E	489	4/4	0.90	0.20	-	33,34,35,37	0
3	EDO	D	488	4/4	0.94	0.20	-	34,35,36,37	0
3	EDO	E	492	4/4	0.89	0.15	-	49,50,50,53	0
3	EDO	B	492	4/4	0.77	0.20	-	59,59,60,60	0
3	EDO	A	492	4/4	0.91	0.21	-	51,52,52,52	0
3	EDO	A	495	4/4	0.91	0.20	-	33,34,36,39	0
3	EDO	F	489	4/4	0.83	0.20	-	42,42,43,44	0
3	EDO	A	489	4/4	0.84	0.20	-	40,41,41,42	0
3	EDO	D	492	4/4	0.91	0.17	-	52,52,52,53	0
3	EDO	E	493	4/4	0.95	0.20	-	52,52,52,52	0
3	EDO	F	492	4/4	0.90	0.16	-	52,52,52,54	0
3	EDO	C	491	4/4	0.89	0.22	-	46,47,47,48	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	EDO	F	493	4/4	0.61	0.21	-	63,64,64,64	0
3	EDO	A	491	4/4	0.60	0.27	-	57,58,58,58	0
3	EDO	A	493	4/4	0.76	0.33	-	63,63,64,65	0
3	EDO	B	491	4/4	0.61	0.31	-	73,74,74,74	0
3	EDO	C	492	4/4	0.87	0.17	-	50,50,51,53	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.