



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:39 PM GMT

PDB ID : 3UG5
Title : Crystal structure of alpha-L-arabinofuranosidase from *Thermotoga maritima* xylose complex
Authors : Im, D.-H.; Miyazaki, K.; Wakagi, T.; Fushinobu, S.
Deposited on : 2011-11-02
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (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

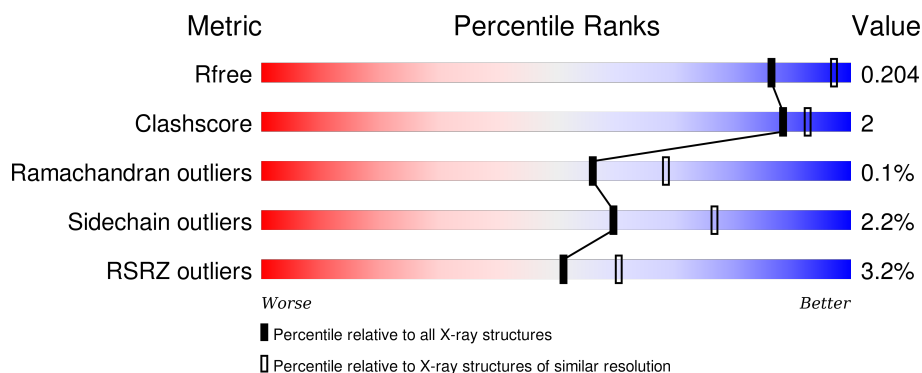
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



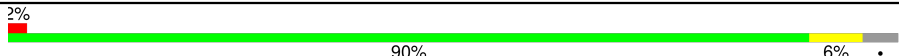
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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>3%</div> <div>88%</div> <div>7%</div> <div>•</div> </div>
1	B	504	<div> <div>4%</div> <div>87%</div> <div>8%</div> <div>• •</div> </div>
1	C	504	<div> <div>3%</div> <div>89%</div> <div>7%</div> <div>•</div> </div>
1	D	504	<div> <div>3%</div> <div>85%</div> <div>10%</div> <div>• 5%</div> </div>
1	E	504	<div> <div>3%</div> <div>88%</div> <div>7%</div> <div>• •</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	504	

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	XYP	A	488	-	-	X	X
3	XYP	A	489	-	-	-	X
3	XYP	B	488	-	-	-	X
3	XYP	C	488	-	-	-	X
3	XYP	D	488	-	-	-	X
3	XYP	D	489	-	-	-	X
3	XYP	E	488	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 24750 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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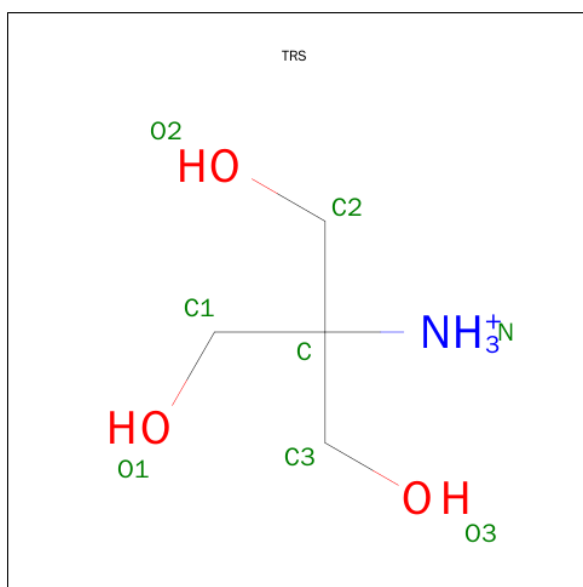
Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	GLY	-	EXPRESSION TAG	UNP Q9WYB7
C	-1	SER	-	EXPRESSION TAG	UNP Q9WYB7
C	0	HIS	-	EXPRESSION TAG	UNP Q9WYB7
C	4	GLY	ARG	ENGINEERED MUTATION	UNP Q9WYB7
D	-19	MET	-	EXPRESSION TAG	UNP Q9WYB7
D	-18	GLY	-	EXPRESSION TAG	UNP Q9WYB7
D	-17	SER	-	EXPRESSION TAG	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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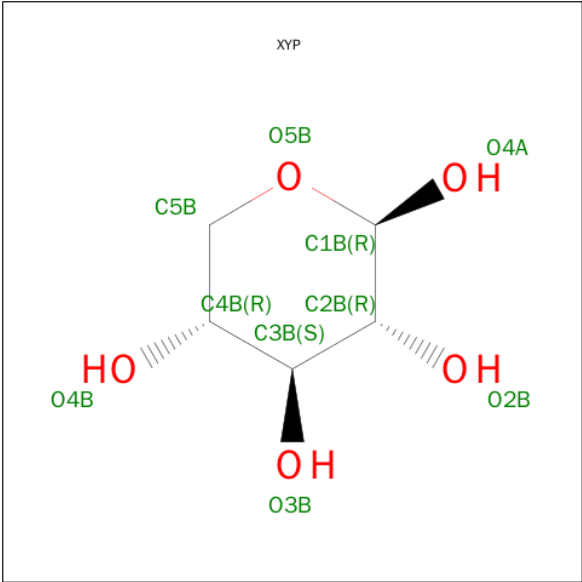
Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	GLY	-	EXPRESSION TAG	UNP Q9WYB7
E	-1	SER	-	EXPRESSION TAG	UNP Q9WYB7
E	0	HIS	-	EXPRESSION TAG	UNP Q9WYB7
E	4	GLY	ARG	ENGINEERED MUTATION	UNP Q9WYB7
F	-19	MET	-	EXPRESSION TAG	UNP Q9WYB7
F	-18	GLY	-	EXPRESSION TAG	UNP Q9WYB7
F	-17	SER	-	EXPRESSION TAG	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: C₄H₁₂NO₃).



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

- Molecule 3 is SUGAR (BETA-D-XYLOPYRANOSE) (three-letter code: XYP) (formula: C₅H₁₀O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			10	5	5		
3	A	1	Total	C	O	0	0
			10	5	5		
3	A	1	Total	C	O	0	0
			10	5	5		
3	B	1	Total	C	O	0	0
			10	5	5		
3	B	1	Total	C	O	0	0
			10	5	5		
3	B	1	Total	C	O	0	0
			10	5	5		
3	C	1	Total	C	O	0	0
			10	5	5		
3	C	1	Total	C	O	0	0
			10	5	5		
3	D	1	Total	C	O	0	0
			10	5	5		
3	D	1	Total	C	O	0	0
			10	5	5		
3	D	1	Total	C	O	0	0
			10	5	5		
3	D	1	Total	C	O	0	0
			10	5	5		
3	E	1	Total	C	O	0	0
			10	5	5		
3	E	1	Total	C	O	0	0
			10	5	5		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	F	1	Total	C	O	0	0
			10	5	5		

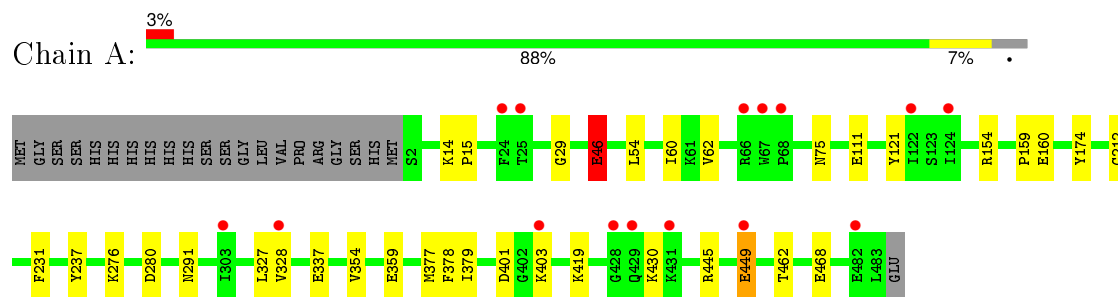
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	213	Total	O	0	0
			213	213		
4	B	175	Total	O	0	0
			175	175		
4	C	239	Total	O	0	0
			239	239		
4	D	205	Total	O	0	0
			205	205		
4	E	191	Total	O	0	0
			191	191		
4	F	224	Total	O	0	0
			224	224		

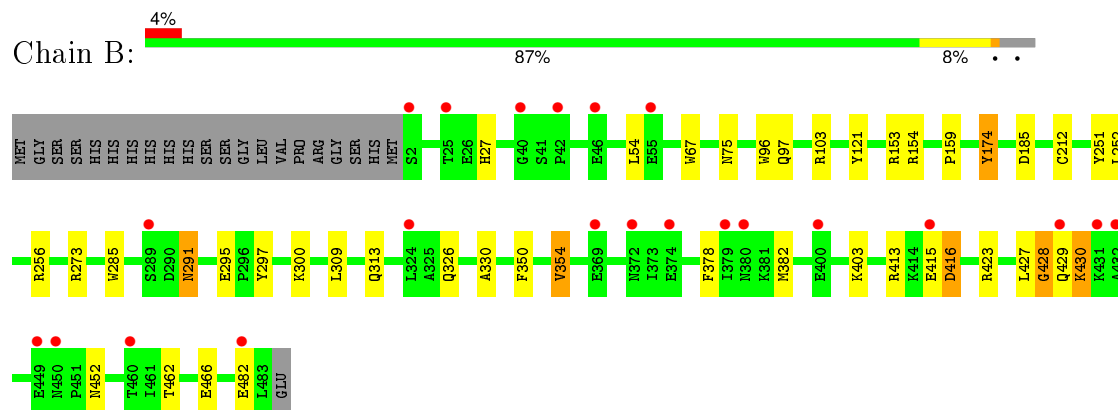
3 Residue-property plots [i](#)

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

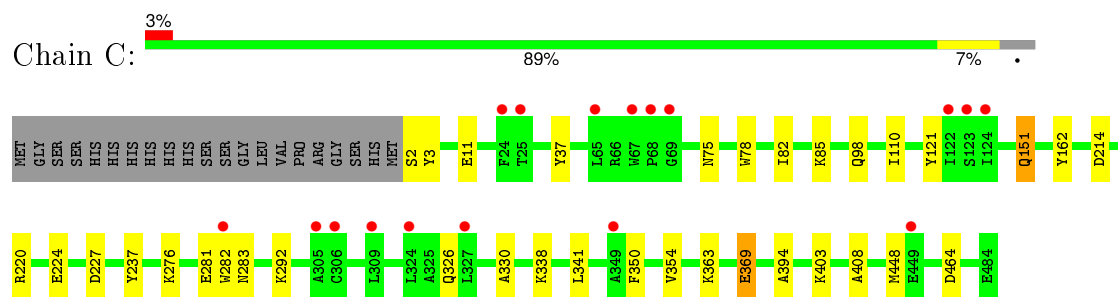
- Molecule 1: Alpha-L-arabinofuranosidase



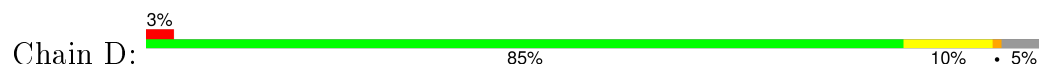
- Molecule 1: Alpha-L-arabinofuranosidase

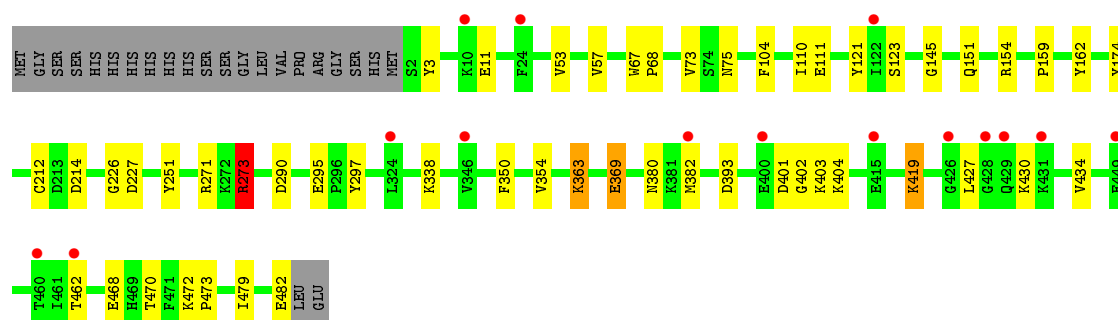


- Molecule 1: Alpha-L-arabinofuranosidase

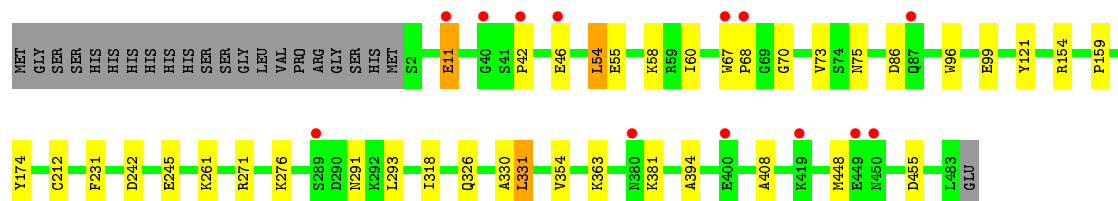
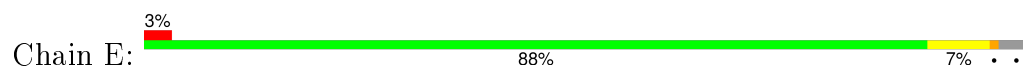


- Molecule 1: Alpha-L-arabinofuranosidase

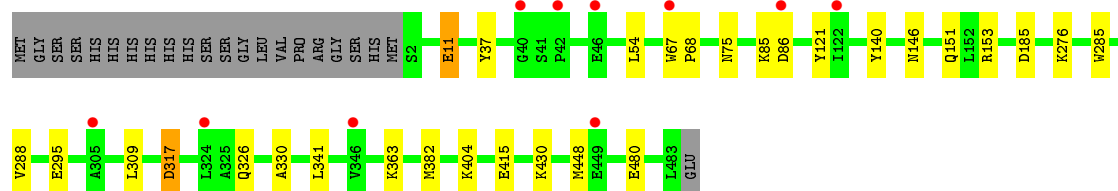
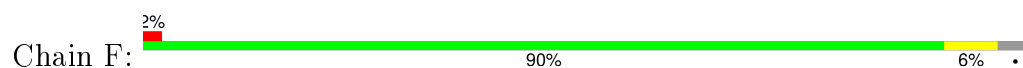




- Molecule 1: Alpha-L-arabinofuranosidase



- 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, α , β , γ	102.20Å 160.93Å 156.17Å 90.00° 91.84° 90.00°	Depositor
Resolution (Å)	29.59 – 2.30 29.59 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.3 (29.59-2.30) 99.3 (29.59-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.97 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.168 , 0.209 0.165 , 0.204	Depositor DCC
R_{free} test set	11084 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	34.9	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 41.0	EDS
Estimated twinning fraction	0.004 for -h,-l,-k 0.000 for -h,l,k 0.012 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 220848 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	24750	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: XYP, TRS

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	1.04	7/3973 (0.2%)	0.86	3/5392 (0.1%)
1	B	1.05	1/3973 (0.0%)	0.89	7/5392 (0.1%)
1	C	1.03	3/3982 (0.1%)	0.85	4/5404 (0.1%)
1	D	1.04	3/3965 (0.1%)	0.85	3/5381 (0.1%)
1	E	0.98	3/3973 (0.1%)	0.84	4/5392 (0.1%)
1	F	0.98	0/3973	0.85	2/5392 (0.0%)
All	All	1.02	17/23839 (0.1%)	0.86	23/32353 (0.1%)

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	46	GLU	CG-CD	8.16	1.64	1.51
1	A	46	GLU	CG-CD	7.68	1.63	1.51
1	C	98	GLN	CG-CD	6.80	1.66	1.51
1	A	160	GLU	CG-CD	6.32	1.61	1.51
1	A	449	GLU	CG-CD	6.29	1.61	1.51
1	E	46	GLU	CB-CG	6.02	1.63	1.52
1	E	55	GLU	CG-CD	5.82	1.60	1.51
1	A	62	VAL	CB-CG1	5.69	1.64	1.52
1	D	104	PHE	CE2-CZ	5.56	1.48	1.37
1	B	174	TYR	CD1-CE1	5.53	1.47	1.39
1	A	337	GLU	CG-CD	5.53	1.60	1.51
1	C	224	GLU	CB-CG	-5.29	1.42	1.52
1	D	111	GLU	CG-CD	5.24	1.59	1.51
1	A	46	GLU	CB-CG	5.18	1.61	1.52
1	C	151	GLN	CG-CD	5.07	1.62	1.51
1	A	359	GLU	CG-CD	5.03	1.59	1.51
1	D	151	GLN	CG-CD	5.00	1.62	1.51

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	273	ARG	NE-CZ-NH1	-6.98	116.81	120.30
1	B	185	ASP	CB-CG-OD1	6.91	124.52	118.30
1	F	317	ASP	CB-CG-OD1	6.51	124.16	118.30
1	D	290	ASP	CB-CG-OD1	6.39	124.06	118.30
1	B	54	LEU	CB-CG-CD1	-6.31	100.28	111.00
1	E	455	ASP	CB-CG-OD1	6.27	123.94	118.30
1	B	103	ARG	NE-CZ-NH2	-6.05	117.28	120.30
1	C	214	ASP	CB-CG-OD1	6.01	123.71	118.30
1	A	280	ASP	CB-CG-OD1	5.61	123.35	118.30
1	B	382	MET	CG-SD-CE	5.61	109.18	100.20
1	A	291	ASN	CB-CA-C	-5.46	99.47	110.40
1	C	220	ARG	NE-CZ-NH2	-5.44	117.58	120.30
1	C	227	ASP	CB-CG-OD1	5.40	123.16	118.30
1	C	464	ASP	CB-CG-OD2	-5.40	113.44	118.30
1	B	67	TRP	C-N-CD	-5.33	108.87	120.60
1	B	153	ARG	NE-CZ-NH1	5.33	122.97	120.30
1	A	54	LEU	CA-CB-CG	5.31	127.52	115.30
1	D	214	ASP	CB-CG-OD2	-5.26	113.56	118.30
1	E	86	ASP	CB-CG-OD1	5.17	122.95	118.30
1	F	86	ASP	CB-CG-OD1	5.11	122.90	118.30
1	E	291	ASN	CB-CA-C	-5.10	100.21	110.40
1	B	251	TYR	CA-CB-CG	-5.04	103.82	113.40
1	E	271	ARG	NE-CZ-NH1	5.03	122.81	120.30

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	3876	0	3793	18	0
1	B	3876	0	3793	15	0
1	C	3885	0	3799	19	0
1	D	3868	0	3782	23	0
1	E	3876	0	3793	18	0
1	F	3876	0	3793	17	0
2	A	16	0	24	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	16	0	24	0	0
2	C	16	0	24	5	0
2	D	16	0	24	0	0
2	E	16	0	24	2	0
2	F	16	0	24	3	0
3	A	30	0	30	8	0
3	B	30	0	30	2	0
3	C	20	0	20	0	0
3	D	40	0	40	0	0
3	E	20	0	20	0	0
3	F	10	0	10	1	0
4	A	213	0	0	2	0
4	B	175	0	0	0	0
4	C	239	0	0	3	0
4	D	205	0	0	0	0
4	E	191	0	0	0	0
4	F	224	0	0	5	0
All	All	24750	0	23047	114	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 (114) 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)
2:C:486:TRS:H31	4:C:1244:HOH:O	1.69	0.92
1:C:237:TYR:OH	2:C:486:TRS:H21	1.75	0.86
1:F:11:GLU:HG2	1:F:363:LYS:HB2	1.60	0.84
3:A:488:XYP:H4B	4:A:690:HOH:O	1.77	0.83
1:F:185:ASP:HB3	4:F:1239:HOH:O	1.81	0.79
1:A:29:GLY:O	3:A:488:XYP:H5B2	1.84	0.77
1:A:29:GLY:CA	3:A:488:XYP:H5B2	2.14	0.77
1:D:11:GLU:HG2	1:D:363:LYS:HD3	1.66	0.75
1:E:11:GLU:HG2	1:E:363:LYS:HB2	1.69	0.74
2:F:486:TRS:H21	4:F:1247:HOH:O	1.88	0.73
1:A:29:GLY:O	3:A:488:XYP:C5B	2.38	0.71
1:A:29:GLY:C	3:A:488:XYP:H5B2	2.11	0.69
1:A:29:GLY:HA2	3:A:488:XYP:H5B2	1.75	0.69
1:D:419:LYS:HE2	1:D:468:GLU:OE1	1.93	0.68
1:C:237:TYR:HH	2:C:486:TRS:H21	1.60	0.66
1:F:151:GLN:HG2	4:F:497:HOH:O	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:350:PHE:O	1:B:354:VAL:HB	1.96	0.66
1:C:237:TYR:OH	2:C:486:TRS:C2	2.44	0.66
1:C:11:GLU:HG2	1:C:363:LYS:HE3	1.77	0.66
1:E:293:LEU:HB3	1:E:331:LEU:HD13	1.77	0.65
1:A:154:ARG:HG2	1:A:159:PRO:HA	1.78	0.65
1:B:256:ARG:HE	3:B:489:XYP:H5B1	1.63	0.64
1:A:237:TYR:OH	2:A:486:TRS:H21	1.98	0.63
1:D:3:TYR:CZ	1:D:369:GLU:HG3	2.35	0.61
1:C:281:GLU:OE2	2:C:486:TRS:H21	2.01	0.59
1:D:295:GLU:HG2	1:D:297:TYR:CE1	2.38	0.59
1:B:96:TRP:HE3	1:B:291:ASN:HB3	1.68	0.58
1:B:326:GLN:HB2	1:B:330:ALA:O	2.03	0.58
2:A:486:TRS:H31	4:A:1245:HOH:O	2.04	0.58
1:E:96:TRP:HZ2	2:E:486:TRS:H21	1.70	0.57
1:E:11:GLU:HG2	1:E:363:LYS:CB	2.34	0.57
1:D:434:VAL:HG22	1:D:479:ILE:HG12	1.87	0.57
1:D:403:LYS:O	1:D:482:GLU:HG2	2.06	0.56
1:E:326:GLN:HB2	1:E:330:ALA:O	2.05	0.56
1:B:252:LEU:HD11	3:B:489:XYP:H4B	1.90	0.54
1:F:341:LEU:HD12	1:F:341:LEU:C	2.29	0.53
1:A:29:GLY:O	3:A:488:XYP:H5B1	2.09	0.52
1:C:3:TYR:OH	1:C:369:GLU:HG3	2.09	0.52
1:A:46:GLU:CD	1:A:46:GLU:H	2.13	0.52
3:A:489:XYP:H5B2	1:D:271:ARG:HD2	1.91	0.52
1:F:285:TRP:CZ3	2:F:486:TRS:O3	2.64	0.51
1:C:3:TYR:CZ	1:C:369:GLU:HG3	2.45	0.51
1:B:423:ARG:NH1	1:B:466:GLU:OE2	2.44	0.51
1:D:154:ARG:HG2	1:D:159:PRO:HA	1.92	0.49
1:B:96:TRP:CE3	1:B:291:ASN:HB3	2.47	0.49
1:F:37:TYR:OH	1:F:85:LYS:NZ	2.42	0.49
1:E:261:LYS:HG3	1:E:318:ILE:HG12	1.95	0.48
1:C:350:PHE:O	1:C:354:VAL:HB	2.13	0.48
1:D:226:GLY:O	1:D:273:ARG:HD3	2.14	0.48
1:F:382:MET:CE	4:F:1205:HOH:O	2.62	0.48
1:C:338:LYS:HG2	4:C:516:HOH:O	2.13	0.47
1:F:288:VAL:HG22	1:F:295:GLU:OE1	2.13	0.47
1:C:151:GLN:HG2	4:C:908:HOH:O	2.13	0.47
1:C:341:LEU:C	1:C:341:LEU:HD12	2.35	0.47
1:D:174:TYR:CE2	1:D:212:CYS:HB3	2.50	0.46
1:E:60:ILE:O	1:E:354:VAL:HG11	2.15	0.46
1:E:96:TRP:CZ2	2:E:486:TRS:H21	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:285:TRP:CE3	1:B:378:PHE:HE1	2.33	0.46
1:D:67:TRP:CE2	1:D:68:PRO:HB3	2.51	0.46
1:E:154:ARG:HG2	1:E:159:PRO:HA	1.97	0.46
1:D:472:LYS:HB3	1:D:473:PRO:HD2	1.98	0.45
1:E:394:ALA:HA	1:E:408:ALA:O	2.16	0.45
1:A:14:LYS:HB2	1:A:15:PRO:HD2	1.99	0.45
1:B:427:LEU:O	1:B:428:GLY:O	2.34	0.45
1:C:276:LYS:HD3	1:C:276:LYS:HA	1.80	0.45
1:A:419:LYS:HD3	1:A:468:GLU:OE1	2.17	0.44
1:C:394:ALA:HA	1:C:408:ALA:O	2.17	0.44
1:D:73:VAL:HG21	1:D:123:SER:O	2.18	0.44
1:C:37:TYR:OH	1:C:85:LYS:NZ	2.46	0.44
1:D:11:GLU:HG2	1:D:363:LYS:HB2	2.00	0.44
1:C:282:TRP:O	1:C:283:ASN:HB2	2.18	0.43
1:B:309:LEU:O	1:B:313:GLN:HG3	2.17	0.43
1:D:380:ASN:OD1	1:D:382:MET:HG3	2.17	0.43
1:F:404:LYS:NZ	1:F:480:GLU:OE2	2.52	0.43
1:E:11:GLU:CG	1:E:363:LYS:HB2	2.42	0.43
1:D:419:LYS:HG2	1:D:470:THR:OG1	2.19	0.43
1:E:231:PHE:HA	1:E:276:LYS:O	2.17	0.43
1:A:60:ILE:O	1:A:354:VAL:HG21	2.18	0.43
1:F:276:LYS:HE3	1:F:317:ASP:O	2.18	0.43
1:A:174:TYR:CE2	1:A:212:CYS:HB3	2.54	0.43
1:A:378:PHE:CG	1:A:379:ILE:N	2.87	0.43
1:B:300:LYS:HE3	1:B:300:LYS:HB3	1.73	0.43
1:F:341:LEU:HD12	1:F:341:LEU:O	2.19	0.43
1:D:350:PHE:O	1:D:354:VAL:HB	2.19	0.43
1:A:327:LEU:HG	1:A:328:VAL:HG23	2.00	0.43
1:B:154:ARG:HG2	1:B:159:PRO:HA	1.99	0.43
1:B:295:GLU:HG2	1:B:297:TYR:CE1	2.54	0.43
1:D:227:ASP:O	1:D:273:ARG:HD2	2.19	0.42
1:D:251:TYR:OH	1:D:393:ASP:OD2	2.33	0.42
2:F:486:TRS:H31	3:F:487:XYP:O2B	2.19	0.42
1:E:70:GLY:O	1:E:73:VAL:HG12	2.19	0.42
1:A:401:ASP:OD1	1:A:403:LYS:HB2	2.20	0.42
1:F:382:MET:HE1	4:F:1205:HOH:O	2.19	0.42
1:F:67:TRP:CE2	1:F:68:PRO:HB3	2.54	0.42
1:E:242:ASP:HB3	1:E:245:GLU:HB2	2.02	0.42
1:C:403:LYS:HD3	1:C:403:LYS:HA	1.92	0.42
1:F:309:LEU:HA	1:F:309:LEU:HD23	1.92	0.42
1:E:67:TRP:CE2	1:E:68:PRO:HB3	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:401:ASP:OD1	1:D:402:GLY:N	2.53	0.42
1:F:326:GLN:HB2	1:F:330:ALA:O	2.20	0.42
1:E:174:TYR:CE2	1:E:212:CYS:HB3	2.55	0.41
1:A:231:PHE:HA	1:A:276:LYS:O	2.19	0.41
1:C:326:GLN:HB2	1:C:330:ALA:O	2.20	0.41
1:E:54:LEU:O	1:E:58:LYS:HG3	2.21	0.41
1:D:53:VAL:O	1:D:57:VAL:HG23	2.20	0.41
1:D:110:ILE:HG21	1:D:162:TYR:CD1	2.56	0.41
1:F:67:TRP:CD1	1:F:68:PRO:HA	2.56	0.41
1:B:415:GLU:HB2	1:B:416:ASP:OD2	2.21	0.41
1:B:174:TYR:CE2	1:B:212:CYS:HB3	2.56	0.41
1:C:78:TRP:CZ3	1:C:82:ILE:HD13	2.56	0.40
1:A:377:MET:O	1:A:378:PHE:HB3	2.21	0.40
1:C:110:ILE:HG21	1:C:162:TYR:CD1	2.56	0.40
1:D:145:GLY:HA3	1:E:99:GLU:OE2	2.21	0.40
1:F:140:TYR:O	1:F:153:ARG:HD3	2.20	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	480/504 (95%)	463 (96%)	17 (4%)	0	100	100
1	B	480/504 (95%)	453 (94%)	24 (5%)	3 (1%)	30	36
1	C	481/504 (95%)	464 (96%)	17 (4%)	0	100	100
1	D	479/504 (95%)	452 (94%)	27 (6%)	0	100	100
1	E	480/504 (95%)	463 (96%)	17 (4%)	0	100	100
1	F	480/504 (95%)	452 (94%)	28 (6%)	0	100	100
All	All	2880/3024 (95%)	2747 (95%)	130 (4%)	3 (0%)	56	68

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	428	GLY
1	B	27	HIS
1	B	430	LYS

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%)	409 (98%)	8 (2%)	65	81
1	B	417/436 (96%)	403 (97%)	14 (3%)	44	59
1	C	418/436 (96%)	412 (99%)	6 (1%)	74	86
1	D	416/436 (95%)	405 (97%)	11 (3%)	54	71
1	E	417/436 (96%)	409 (98%)	8 (2%)	65	81
1	F	417/436 (96%)	409 (98%)	8 (2%)	65	81
All	All	2502/2616 (96%)	2447 (98%)	55 (2%)	60	77

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	GLU
1	A	75	ASN
1	A	111	GLU
1	A	121	TYR
1	A	430	LYS
1	A	445	ARG
1	A	449	GLU
1	A	462	THR
1	B	75	ASN
1	B	97	GLN
1	B	121	TYR
1	B	273	ARG
1	B	291	ASN
1	B	354	VAL

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Mol	Chain	Res	Type
1	B	403	LYS
1	B	413	ARG
1	B	416	ASP
1	B	429	GLN
1	B	430	LYS
1	B	452	ASN
1	B	462	THR
1	B	482	GLU
1	C	2	SER
1	C	75	ASN
1	C	121	TYR
1	C	292	LYS
1	C	369	GLU
1	C	448	MET
1	D	75	ASN
1	D	121	TYR
1	D	273	ARG
1	D	338	LYS
1	D	363	LYS
1	D	369	GLU
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	42	PRO
1	E	54	LEU
1	E	75	ASN
1	E	121	TYR
1	E	331	LEU
1	E	381	LYS
1	E	448	MET
1	F	11	GLU
1	F	54	LEU
1	F	75	ASN
1	F	121	TYR
1	F	146	ASN
1	F	415	GLU
1	F	430	LYS
1	F	448	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no

such sidechains identified.

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 ⓘ

27 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	TRS	A	485	-	7,7,7	1.00	1 (14%)	9,9,9	1.02	1 (11%)
2	TRS	A	486	-	7,7,7	0.72	0	9,9,9	2.12	3 (33%)
3	XYP	A	487	-	10,10,10	1.25	1 (10%)	12,14,14	1.39	2 (16%)
3	XYP	A	488	-	10,10,10	4.25	5 (50%)	12,14,14	3.73	5 (41%)
3	XYP	A	489	-	10,10,10	2.98	3 (30%)	12,14,14	2.22	3 (25%)
2	TRS	B	485	-	7,7,7	1.08	1 (14%)	9,9,9	1.70	1 (11%)
2	TRS	B	486	-	7,7,7	0.47	0	9,9,9	2.35	5 (55%)
3	XYP	B	487	-	10,10,10	1.08	1 (10%)	12,14,14	0.42	0
3	XYP	B	488	-	10,10,10	2.50	3 (30%)	12,14,14	3.64	9 (75%)
3	XYP	B	489	-	10,10,10	2.67	5 (50%)	12,14,14	1.36	1 (8%)
2	TRS	C	485	-	7,7,7	1.37	1 (14%)	9,9,9	2.28	4 (44%)
2	TRS	C	486	-	7,7,7	1.13	0	9,9,9	1.67	3 (33%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	XYP	C	487	-	10,10,10	1.43	1 (10%)	12,14,14	1.04	1 (8%)
3	XYP	C	488	-	10,10,10	2.54	3 (30%)	12,14,14	2.28	4 (33%)
2	TRS	D	485	-	7,7,7	0.58	0	9,9,9	1.51	3 (33%)
2	TRS	D	486	-	7,7,7	0.89	0	9,9,9	1.89	4 (44%)
3	XYP	D	487	-	10,10,10	2.18	4 (40%)	12,14,14	1.02	2 (16%)
3	XYP	D	488	-	10,10,10	2.54	3 (30%)	12,14,14	1.32	1 (8%)
3	XYP	D	489	-	10,10,10	2.26	4 (40%)	12,14,14	1.60	1 (8%)
3	XYP	D	490	-	10,10,10	1.89	4 (40%)	12,14,14	2.03	3 (25%)
2	TRS	E	485	-	7,7,7	1.06	0	9,9,9	1.26	1 (11%)
2	TRS	E	486	-	7,7,7	0.43	0	9,9,9	1.57	2 (22%)
3	XYP	E	487	-	10,10,10	1.87	1 (10%)	12,14,14	1.68	2 (16%)
3	XYP	E	488	-	10,10,10	2.63	4 (40%)	12,14,14	2.85	6 (50%)
2	TRS	F	485	-	7,7,7	0.85	0	9,9,9	1.07	1 (11%)
2	TRS	F	486	-	7,7,7	0.90	0	9,9,9	2.35	5 (55%)
3	XYP	F	487	-	10,10,10	1.92	2 (20%)	12,14,14	1.40	2 (16%)

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
2	TRS	A	486	-	-	0/9/9/9	0/0/0/0
3	XYP	A	487	-	-	0/0/17/17	0/1/1/1
3	XYP	A	488	-	-	0/0/17/17	0/1/1/1
3	XYP	A	489	-	-	0/0/17/17	0/1/1/1
2	TRS	B	485	-	-	0/9/9/9	0/0/0/0
2	TRS	B	486	-	-	0/9/9/9	0/0/0/0
3	XYP	B	487	-	-	0/0/17/17	0/1/1/1
3	XYP	B	488	-	-	0/0/17/17	0/1/1/1
3	XYP	B	489	-	-	0/0/17/17	0/1/1/1
2	TRS	C	485	-	-	0/9/9/9	0/0/0/0
2	TRS	C	486	-	-	0/9/9/9	0/0/0/0
3	XYP	C	487	-	-	0/0/17/17	0/1/1/1
3	XYP	C	488	-	-	0/0/17/17	0/1/1/1
2	TRS	D	485	-	-	0/9/9/9	0/0/0/0
2	TRS	D	486	-	-	0/9/9/9	0/0/0/0
3	XYP	D	487	-	-	0/0/17/17	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	XYP	D	488	-	-	0/0/17/17	0/1/1/1
3	XYP	D	489	-	-	0/0/17/17	0/1/1/1
3	XYP	D	490	-	-	0/0/17/17	0/1/1/1
2	TRS	E	485	-	-	0/9/9/9	0/0/0/0
2	TRS	E	486	-	-	0/9/9/9	0/0/0/0
3	XYP	E	487	-	-	0/0/17/17	0/1/1/1
3	XYP	E	488	-	-	0/0/17/17	0/1/1/1
2	TRS	F	485	-	-	0/9/9/9	0/0/0/0
2	TRS	F	486	-	-	0/9/9/9	0/0/0/0
3	XYP	F	487	-	-	0/0/17/17	0/1/1/1

All (47) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	485	TRS	C-N	-3.12	1.46	1.50
2	B	485	TRS	C-N	-2.59	1.46	1.50
2	A	485	TRS	C-N	-2.57	1.46	1.50
3	B	489	XYP	O5B-C5B	2.00	1.47	1.43
3	A	488	XYP	C3B-C2B	2.01	1.57	1.52
3	E	488	XYP	C5B-C4B	2.02	1.57	1.52
3	A	487	XYP	O5B-C1B	2.07	1.46	1.43
3	B	489	XYP	O4A-C1B	2.19	1.47	1.39
3	B	487	XYP	O5B-C5B	2.20	1.47	1.43
3	D	489	XYP	C1B-C2B	2.24	1.57	1.52
3	E	488	XYP	O5B-C5B	2.28	1.47	1.43
3	D	490	XYP	C1B-C2B	2.30	1.57	1.52
3	D	487	XYP	C4B-C3B	2.35	1.55	1.52
3	D	487	XYP	O4A-C1B	2.39	1.47	1.39
3	A	489	XYP	C1B-C2B	2.40	1.57	1.52
3	D	488	XYP	O4A-C1B	2.41	1.48	1.39
3	D	490	XYP	O5B-C1B	2.46	1.46	1.43
3	B	489	XYP	C1B-C2B	2.64	1.58	1.52
3	D	490	XYP	O3B-C3B	2.65	1.49	1.43
3	D	489	XYP	C4B-C3B	2.73	1.56	1.52
3	D	489	XYP	O5B-C5B	2.73	1.48	1.43
3	B	488	XYP	O5B-C1B	2.80	1.47	1.43
3	F	487	XYP	O5B-C5B	2.80	1.48	1.43
3	B	489	XYP	C4B-C3B	2.81	1.56	1.52
3	C	487	XYP	O5B-C1B	2.84	1.47	1.43
3	D	488	XYP	C4B-C3B	3.03	1.56	1.52
3	D	490	XYP	O4A-C1B	3.11	1.50	1.39
3	C	488	XYP	O5B-C5B	3.30	1.49	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	488	XYP	C4B-C3B	3.44	1.57	1.52
3	D	487	XYP	O5B-C1B	3.88	1.48	1.43
3	D	487	XYP	O5B-C5B	4.06	1.50	1.43
3	A	488	XYP	O5B-C1B	4.10	1.49	1.43
3	E	488	XYP	O5B-C1B	4.19	1.49	1.43
3	B	488	XYP	O2B-C2B	4.31	1.53	1.43
3	E	487	XYP	O5B-C1B	4.71	1.50	1.43
3	F	487	XYP	O5B-C1B	4.75	1.50	1.43
3	D	489	XYP	O5B-C1B	5.19	1.50	1.43
3	B	488	XYP	C4B-C3B	5.23	1.59	1.52
3	C	488	XYP	O5B-C1B	5.71	1.51	1.43
3	A	489	XYP	C4B-C3B	5.73	1.60	1.52
3	D	488	XYP	O5B-C1B	5.87	1.51	1.43
3	E	488	XYP	C4B-C3B	5.89	1.60	1.52
3	A	488	XYP	O5B-C5B	6.19	1.54	1.43
3	A	489	XYP	O5B-C1B	6.21	1.52	1.43
3	B	489	XYP	O5B-C1B	6.24	1.52	1.43
3	A	488	XYP	O4B-C4B	6.78	1.58	1.43
3	A	488	XYP	C4B-C3B	8.30	1.63	1.52

All (75) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	488	XYP	C1B-C2B-C3B	-7.38	99.45	110.43
3	E	488	XYP	C1B-C2B-C3B	-6.35	100.99	110.43
2	A	486	TRS	O1-C1-C	-4.48	102.12	111.18
3	C	488	XYP	C1B-C2B-C3B	-4.29	104.05	110.43
2	C	485	TRS	C3-C-C2	-4.24	101.60	110.78
2	F	486	TRS	C3-C-C1	-4.21	101.67	110.78
2	B	486	TRS	O1-C1-C	-3.84	103.42	111.18
2	A	486	TRS	C2-C-C1	-3.32	103.59	110.78
2	B	486	TRS	C3-C-C1	-3.16	103.94	110.78
2	C	485	TRS	C1-C-N	-3.08	102.48	108.09
3	B	488	XYP	O4A-C1B-O5B	-2.83	101.83	109.90
3	A	489	XYP	C4B-C3B-C2B	-2.72	106.42	111.04
2	E	485	TRS	C2-C-C1	-2.69	104.96	110.78
2	F	485	TRS	O1-C1-C	-2.54	106.04	111.18
3	B	488	XYP	O5B-C5B-C4B	-2.54	106.74	110.86
2	C	486	TRS	O1-C1-C	-2.50	106.12	111.18
3	A	487	XYP	O4B-C4B-C3B	-2.50	105.09	110.12
2	D	485	TRS	O1-C1-C	-2.45	106.22	111.18
2	C	486	TRS	C2-C-C1	-2.43	105.52	110.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	486	TRS	O1-C1-C	-2.33	106.46	111.18
2	E	486	TRS	C3-C-C1	-2.29	105.82	110.78
3	C	487	XYP	O4B-C4B-C3B	-2.27	105.56	110.12
2	D	486	TRS	C3-C-C2	-2.24	105.94	110.78
3	B	488	XYP	C4B-C3B-C2B	-2.18	107.33	111.04
2	F	486	TRS	O3-C3-C	-2.18	106.78	111.18
3	A	488	XYP	C5B-C4B-C3B	-2.13	107.02	109.54
2	B	486	TRS	O2-C2-C	-2.11	106.92	111.18
2	A	485	TRS	O3-C3-C	-2.08	106.97	111.18
2	D	485	TRS	C3-C-C1	-2.05	106.33	110.78
3	D	487	XYP	C5B-C4B-C3B	2.00	111.91	109.54
2	D	486	TRS	C3-C-N	2.01	111.75	108.09
3	F	487	XYP	O4B-C4B-C5B	2.02	113.25	109.21
2	F	486	TRS	C2-C-C1	2.08	115.29	110.78
3	D	487	XYP	O4A-C1B-C2B	2.16	114.99	109.21
3	E	487	XYP	O4A-C1B-C2B	2.19	115.09	109.21
3	E	488	XYP	O2B-C2B-C1B	2.21	114.68	109.82
3	E	488	XYP	C5B-C4B-C3B	2.22	112.17	109.54
3	C	488	XYP	O3B-C3B-C4B	2.25	114.07	110.00
2	F	486	TRS	C1-C-N	2.25	112.19	108.09
2	E	486	TRS	C1-C-N	2.33	112.32	108.09
3	E	488	XYP	O2B-C2B-C3B	2.37	115.68	110.34
2	A	486	TRS	C2-C-N	2.38	112.42	108.09
2	D	485	TRS	C3-C-N	2.40	112.45	108.09
3	D	488	XYP	C5B-C4B-C3B	2.54	112.55	109.54
2	C	486	TRS	C2-C-N	2.59	112.79	108.09
3	A	488	XYP	O4A-C1B-O5B	2.66	117.48	109.90
3	D	490	XYP	C5B-C4B-C3B	2.71	112.75	109.54
2	B	486	TRS	C1-C-N	2.72	113.03	108.09
2	C	485	TRS	C2-C-N	2.74	113.08	108.09
2	C	485	TRS	C3-C-C1	2.81	116.87	110.78
2	B	486	TRS	C3-C-N	2.86	113.30	108.09
3	F	487	XYP	O4A-C1B-O5B	2.88	118.10	109.90
3	A	487	XYP	C5B-C4B-C3B	2.93	113.01	109.54
2	B	485	TRS	C2-C-N	2.98	113.52	108.09
3	B	488	XYP	O4B-C4B-C3B	3.12	116.39	110.12
3	E	488	XYP	O4B-C4B-C3B	3.23	116.61	110.12
2	D	486	TRS	C2-C-N	3.38	114.24	108.09
3	B	488	XYP	O3B-C3B-C4B	3.38	116.11	110.00
3	C	488	XYP	C5B-C4B-C3B	3.41	113.58	109.54
3	D	490	XYP	O4A-C1B-C2B	3.43	118.40	109.21
3	B	489	XYP	C5B-C4B-C3B	3.66	113.87	109.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	489	XYP	O4B-C4B-C3B	3.93	118.02	110.12
3	A	488	XYP	O3B-C3B-C4B	3.97	117.17	110.00
3	B	488	XYP	O2B-C2B-C3B	4.12	119.61	110.34
2	F	486	TRS	O1-C1-C	4.13	119.54	111.18
3	C	488	XYP	O5B-C5B-C4B	4.18	117.64	110.86
3	E	487	XYP	C5B-C4B-C3B	4.22	114.54	109.54
3	D	489	XYP	O5B-C5B-C4B	4.29	117.81	110.86
3	D	490	XYP	O5B-C5B-C4B	4.31	117.85	110.86
3	E	488	XYP	O3B-C3B-C4B	4.40	117.94	110.00
3	B	488	XYP	O2B-C2B-C1B	4.40	119.51	109.82
3	A	489	XYP	O3B-C3B-C4B	4.68	118.45	110.00
3	B	488	XYP	O4A-C1B-C2B	4.88	122.29	109.21
3	A	488	XYP	O4B-C4B-C3B	5.28	120.74	110.12
3	A	488	XYP	O5B-C5B-C4B	10.19	127.39	110.86

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	486	TRS	2	0
3	A	488	XYP	7	0
3	A	489	XYP	1	0
3	B	489	XYP	2	0
2	C	486	TRS	5	0
2	E	486	TRS	2	0
2	F	486	TRS	3	0
3	F	487	XYP	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	482/504 (95%)	-0.15	15 (3%)	52	62	21, 30, 50, 63	0
1	B	482/504 (95%)	-0.04	22 (4%)	36	45	23, 33, 52, 66	0
1	C	483/504 (95%)	-0.13	17 (3%)	48	56	21, 30, 48, 64	0
1	D	481/504 (95%)	-0.15	15 (3%)	52	62	22, 32, 50, 65	0
1	E	482/504 (95%)	-0.20	13 (2%)	58	67	23, 32, 50, 64	0
1	F	482/504 (95%)	-0.19	10 (2%)	67	74	23, 32, 49, 64	0
All	All	2892/3024 (95%)	-0.14	92 (3%)	51	60	21, 32, 50, 66	0

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	449	GLU	4.8
1	D	431	LYS	4.4
1	D	429	GLN	4.3
1	D	449	GLU	4.1
1	B	431	LYS	4.0
1	F	449	GLU	3.9
1	A	428	GLY	3.8
1	C	449	GLU	3.6
1	E	449	GLU	3.6
1	C	324	LEU	3.5
1	E	46	GLU	3.4
1	B	460	THR	3.3
1	B	2	SER	3.2
1	B	429	GLN	3.2
1	A	449	GLU	3.2
1	F	46	GLU	3.2
1	C	68	PRO	3.1
1	C	327	LEU	3.1
1	D	382	MET	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	25	THR	3.0
1	C	67	TRP	3.0
1	D	460	THR	3.0
1	C	305	ALA	3.0
1	D	462	THR	2.9
1	A	429	GLN	2.9
1	B	415	GLU	2.8
1	A	68	PRO	2.8
1	B	42	PRO	2.7
1	F	42	PRO	2.7
1	E	68	PRO	2.7
1	B	289	SER	2.7
1	C	24	PHE	2.6
1	A	25	THR	2.6
1	C	349	ALA	2.6
1	C	122	ILE	2.6
1	A	24	PHE	2.6
1	D	10	LYS	2.6
1	B	324	LEU	2.6
1	C	123	SER	2.6
1	E	40	GLY	2.5
1	C	124	ILE	2.5
1	F	40	GLY	2.5
1	D	24	PHE	2.5
1	D	426	GLY	2.5
1	A	67	TRP	2.5
1	C	309	LEU	2.5
1	E	87	GLN	2.4
1	B	380	ASN	2.4
1	B	432	ALA	2.4
1	E	380	ASN	2.4
1	E	42	PRO	2.4
1	E	11	GLU	2.4
1	E	400	GLU	2.4
1	B	379	ILE	2.3
1	B	25	THR	2.3
1	F	305	ALA	2.3
1	A	303	ILE	2.3
1	D	415	GLU	2.3
1	A	328	VAL	2.3
1	B	46	GLU	2.3
1	B	40	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	122	ILE	2.2
1	C	65	LEU	2.2
1	B	369	GLU	2.2
1	A	403	LYS	2.2
1	A	431	LYS	2.2
1	F	122	ILE	2.2
1	F	346	VAL	2.2
1	B	374	GLU	2.2
1	B	450	ASN	2.2
1	E	450	ASN	2.2
1	D	346	VAL	2.1
1	D	428	GLY	2.1
1	E	419	LYS	2.1
1	C	306	CYS	2.1
1	F	86	ASP	2.1
1	C	69	GLY	2.1
1	A	122	ILE	2.1
1	E	289	SER	2.1
1	C	282	TRP	2.1
1	D	400	GLU	2.1
1	F	67	TRP	2.1
1	B	55	GLU	2.1
1	B	372	ASN	2.1
1	A	66	ARG	2.0
1	D	324	LEU	2.0
1	E	67	TRP	2.0
1	B	482	GLU	2.0
1	B	400	GLU	2.0
1	A	124	ILE	2.0
1	A	482	GLU	2.0
1	F	324	LEU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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, 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	LLDF	B-factors(Å ²)	Q<0.9
3	XYP	E	488	10/10	0.72	0.29	7.46	64,75,77,79	0
3	XYP	D	488	10/10	0.78	0.28	5.78	51,61,67,69	0
3	XYP	C	488	10/10	0.77	0.28	4.92	68,73,76,77	0
3	XYP	B	488	10/10	0.87	0.20	3.83	43,56,59,64	0
3	XYP	A	489	10/10	0.78	0.21	3.71	52,59,60,63	0
3	XYP	A	488	10/10	0.75	0.21	3.36	44,55,58,59	0
3	XYP	D	489	10/10	0.91	0.23	2.69	55,60,62,65	0
2	TRS	F	486	8/8	0.91	0.18	1.81	51,53,54,55	0
3	XYP	D	490	10/10	0.89	0.22	1.44	54,60,61,63	0
3	XYP	B	489	10/10	0.86	0.18	1.33	62,68,70,71	0
2	TRS	C	486	8/8	0.85	0.20	0.98	45,50,51,52	0
2	TRS	E	486	8/8	0.92	0.18	0.80	47,48,51,51	0
2	TRS	C	485	8/8	0.98	0.09	0.61	22,25,27,28	0
3	XYP	F	487	10/10	0.95	0.13	0.51	49,50,53,53	0
2	TRS	E	485	8/8	0.98	0.09	0.42	31,32,33,34	0
3	XYP	C	487	10/10	0.94	0.12	0.41	44,48,49,52	0
2	TRS	D	486	8/8	0.94	0.17	0.41	34,39,42,43	0
2	TRS	B	485	8/8	0.98	0.09	0.08	27,29,29,31	0
2	TRS	A	486	8/8	0.93	0.14	0.02	41,45,46,48	0
3	XYP	A	487	10/10	0.95	0.11	-0.03	38,40,41,41	0
2	TRS	B	486	8/8	0.95	0.14	-0.09	42,43,45,46	0
3	XYP	E	487	10/10	0.94	0.10	-0.29	43,46,49,49	0
3	XYP	D	487	10/10	0.95	0.09	-0.48	43,47,49,51	0
3	XYP	B	487	10/10	0.95	0.08	-0.88	44,47,48,48	0
2	TRS	F	485	8/8	0.98	0.08	-1.01	32,32,33,33	0
2	TRS	A	485	8/8	0.98	0.08	-1.02	28,30,32,32	0
2	TRS	D	485	8/8	0.98	0.07	-1.35	28,30,31,31	0

6.5 Other polymers ⓘ

There are no such residues in this entry.