



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:30 PM UTC

PDB ID : 6ABP / pdb_00006abp
Title : SUGAR-BINDING AND CRYSTALLOGRAPHIC STUDIES OF AN ARABINOSE-BINDING PROTEIN MUTANT (MET108LEU) WHICH EXHIBITS ENHANCED AFFINITY AND ALTERED SPECIFICITY
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Deposited on : 1991-04-25
Resolution : 1.67 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

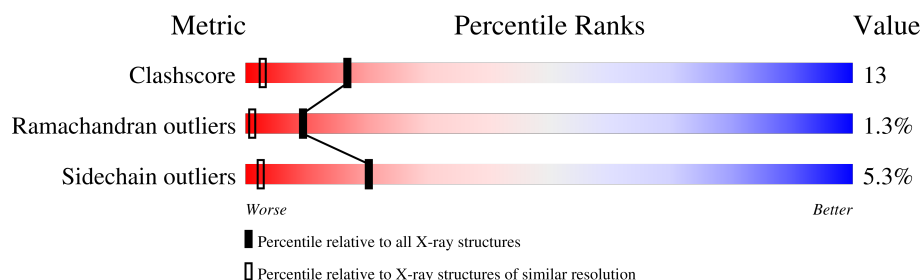
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.67 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))
Clashscore	190562	1078 (1.68-1.68)
Ramachandran outliers	187476	1068 (1.68-1.68)
Sidechain outliers	187428	1067 (1.68-1.68)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	306	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2541 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.

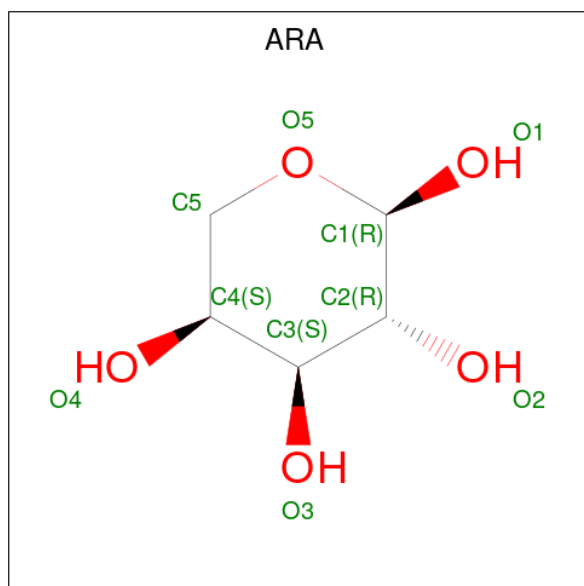
- Molecule 1 is a protein called L-ARABINOSE-BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	305	2316	1474	389	443	10	0	0	0

There is a discrepancy between the modelled and reference sequences:

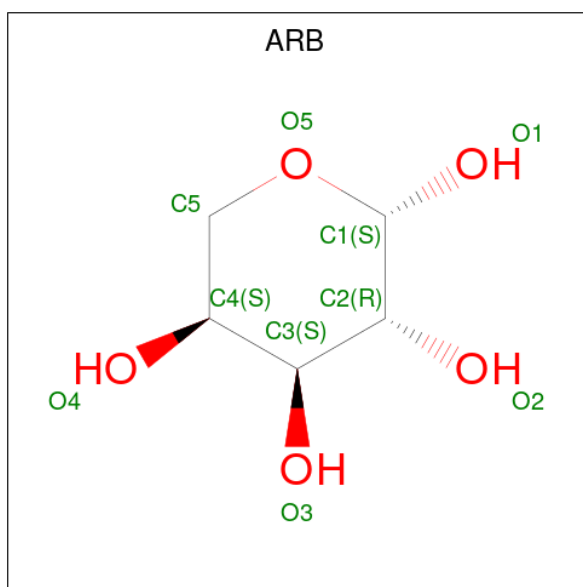
Chain	Residue	Modelled	Actual	Comment	Reference
A	108	LEU	MET	conflict	UNP P02924

- Molecule 2 is alpha-L-arabinopyranose (CCD ID: ARA) (formula: C₅H₁₀O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	10	5	5	0	1

- Molecule 3 is beta-L-arabinopyranose (CCD ID: ARB) (formula: C₅H₁₀O₅).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	205	Total	O	0	0
			205	205		

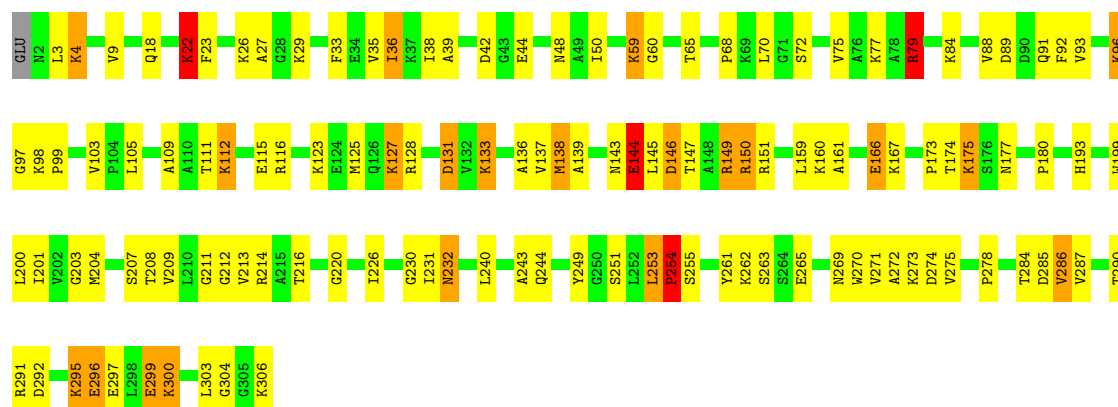
3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

• Molecule 1: L-ARABINOSE-BINDING PROTEIN

Chain A: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	55.54Å 72.12Å 78.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 1.67	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-1.67)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.160 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2541	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ARB, ARA

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	1.72	25/2361 (1.1%)	2.36	130/3190 (4.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	203	GLY	N-CA	7.47	1.53	1.45
1	A	263	SER	N-CA	6.77	1.54	1.46
1	A	249	TYR	C-O	6.17	1.31	1.24
1	A	151	ARG	CZ-NH2	6.07	1.41	1.33
1	A	136	ALA	N-CA	6.06	1.53	1.45
1	A	146	ASP	C-O	-6.00	1.17	1.24
1	A	92	PHE	C-O	5.97	1.31	1.23
1	A	240	LEU	C-O	5.94	1.31	1.24
1	A	139	ALA	N-CA	5.93	1.53	1.46
1	A	193	HIS	CA-CB	5.70	1.59	1.53
1	A	88	VAL	C-O	5.66	1.29	1.24
1	A	284	THR	CA-CB	5.47	1.63	1.53
1	A	60	GLY	N-CA	5.39	1.50	1.45
1	A	297	GLU	C-O	5.39	1.30	1.24
1	A	230	GLY	N-CA	5.38	1.51	1.45
1	A	220	GLY	CA-C	-5.28	1.44	1.51
1	A	9	VAL	C-O	5.28	1.29	1.24
1	A	220	GLY	C-O	5.26	1.30	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	116	ARG	N-CA	5.26	1.52	1.46
1	A	251	SER	C-O	5.21	1.30	1.23
1	A	42	ASP	N-CA	5.19	1.52	1.45
1	A	254	PRO	C-O	5.13	1.30	1.24
1	A	200	LEU	CA-C	5.04	1.58	1.52
1	A	177	ASN	C-N	-5.03	1.26	1.33
1	A	147	THR	CB-OG1	5.01	1.51	1.43

All (130) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	123	LYS	CA-CB-CG	11.78	137.66	114.10
1	A	193	HIS	CA-CB-CG	-11.18	102.62	113.80
1	A	79	ARG	CA-CB-CG	10.73	135.56	114.10
1	A	138	MET	CA-C-O	-9.26	110.46	120.36
1	A	151	ARG	NE-CZ-NH1	9.17	130.67	121.50
1	A	151	ARG	CA-C-O	-8.42	112.03	120.70
1	A	23	PHE	CA-CB-CG	-7.84	105.96	113.80
1	A	116	ARG	CD-NE-CZ	7.76	135.27	124.40
1	A	262	LYS	CA-C-N	-7.72	109.94	120.28
1	A	262	LYS	C-N-CA	-7.72	109.94	120.28
1	A	144	GLU	CB-CG-CD	7.67	125.64	112.60
1	A	131	ASP	CA-CB-CG	7.60	120.20	112.60
1	A	299	GLU	CB-CG-CD	7.58	125.48	112.60
1	A	244	GLN	OE1-CD-NE2	7.57	130.17	122.60
1	A	150	ARG	CD-NE-CZ	7.54	134.96	124.40
1	A	296	GLU	N-CA-CB	7.52	121.14	109.94
1	A	103	VAL	CA-C-O	-7.39	114.79	119.15
1	A	208	THR	CA-C-O	7.28	128.14	120.42
1	A	145	LEU	CA-C-O	7.24	128.21	120.46
1	A	255	SER	O-C-N	7.16	127.71	121.34
1	A	137	VAL	CA-C-O	-7.10	112.96	120.48
1	A	296	GLU	CB-CA-C	-7.05	99.53	110.81
1	A	263	SER	CA-C-N	6.94	129.47	120.44
1	A	263	SER	C-N-CA	6.94	129.47	120.44
1	A	303	LEU	CA-C-N	6.86	134.85	121.41
1	A	303	LEU	C-N-CA	6.86	134.85	121.41
1	A	149	ARG	NE-CZ-NH1	-6.84	114.66	121.50
1	A	98	LYS	CA-C-O	-6.79	113.31	120.23
1	A	295	LYS	CA-C-O	-6.72	113.77	120.70
1	A	72	SER	O-C-N	6.71	128.98	122.07
1	A	166	GLU	CA-C-O	6.71	127.68	120.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	131	ASP	O-C-N	6.70	131.26	123.02
1	A	265	GLU	CA-C-O	6.69	127.64	120.55
1	A	231	ILE	CA-C-O	-6.52	113.45	120.36
1	A	146	ASP	O-C-N	6.48	128.83	122.09
1	A	199	TRP	O-C-N	6.43	131.59	123.23
1	A	133	LYS	CA-C-N	-6.40	112.49	122.67
1	A	133	LYS	C-N-CA	-6.40	112.49	122.67
1	A	79	ARG	NE-CZ-NH1	6.38	127.89	121.50
1	A	99	PRO	CA-C-O	-6.38	114.06	121.34
1	A	291	ARG	O-C-N	6.35	128.85	122.12
1	A	296	GLU	CA-CB-CG	6.34	126.78	114.10
1	A	27	ALA	O-C-N	6.32	128.82	122.12
1	A	18	GLN	CA-C-O	6.26	127.18	120.55
1	A	65	THR	O-C-N	6.24	128.39	121.53
1	A	22	LYS	CG-CD-CE	-6.19	97.05	111.30
1	A	145	LEU	CA-C-N	6.18	128.89	120.54
1	A	145	LEU	C-N-CA	6.18	128.89	120.54
1	A	180	PRO	N-CD-CG	-6.18	93.93	103.20
1	A	200	LEU	CA-C-N	-6.17	115.10	123.12
1	A	200	LEU	C-N-CA	-6.17	115.10	123.12
1	A	137	VAL	O-C-N	6.02	129.50	123.18
1	A	211	GLY	CA-C-N	6.01	126.62	119.94
1	A	211	GLY	C-N-CA	6.01	126.62	119.94
1	A	145	LEU	O-C-N	-6.00	115.57	123.13
1	A	151	ARG	NH1-CZ-NH2	-6.00	111.51	119.30
1	A	213	VAL	CA-C-O	-5.98	114.83	121.17
1	A	174	THR	CA-C-O	-5.93	114.42	120.71
1	A	262	LYS	O-C-N	5.93	128.91	122.15
1	A	290	THR	CA-CB-OG1	-5.88	100.78	109.60
1	A	161	ALA	CA-C-O	-5.85	114.35	120.55
1	A	274	ASP	N-CA-CB	5.84	120.90	112.08
1	A	253	LEU	CA-C-N	5.77	127.06	119.84
1	A	253	LEU	C-N-CA	5.77	127.06	119.84
1	A	99	PRO	CB-CA-C	-5.75	104.05	111.46
1	A	200	LEU	O-C-N	5.70	129.71	123.27
1	A	103	VAL	O-C-N	5.69	126.52	121.57
1	A	299	GLU	CG-CD-OE1	5.68	131.46	118.40
1	A	138	MET	O-C-N	5.65	129.66	123.22
1	A	93	VAL	O-C-N	5.62	129.40	123.00
1	A	159	LEU	O-C-N	5.59	127.83	122.07
1	A	207	SER	CA-CB-OG	-5.57	99.96	111.10
1	A	193	HIS	CB-CA-C	-5.56	103.92	111.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	50	ILE	O-C-N	5.55	127.25	121.87
1	A	220	GLY	N-CA-C	5.52	122.98	115.30
1	A	36	ILE	CA-C-O	-5.48	114.24	120.65
1	A	149	ARG	NH1-CZ-NH2	5.46	126.41	119.30
1	A	38	ILE	CA-C-N	5.46	129.89	122.19
1	A	38	ILE	C-N-CA	5.46	129.89	122.19
1	A	150	ARG	CA-CB-CG	-5.44	103.22	114.10
1	A	77	LYS	CA-CB-CG	5.43	124.95	114.10
1	A	232	ASN	CA-CB-CG	-5.43	107.17	112.60
1	A	278	PRO	CA-C-O	-5.42	115.36	121.43
1	A	269	ASN	CA-CB-CG	-5.41	107.19	112.60
1	A	70	LEU	CA-C-N	5.40	126.09	119.99
1	A	70	LEU	C-N-CA	5.40	126.09	119.99
1	A	39	ALA	CA-C-O	-5.40	114.44	120.54
1	A	243	ALA	CA-C-O	5.40	126.26	120.70
1	A	286	VAL	O-C-N	5.39	128.37	123.19
1	A	111	THR	O-C-N	5.38	127.61	122.07
1	A	125	MET	O-C-N	5.36	127.80	122.12
1	A	243	ALA	O-C-N	-5.32	116.35	122.09
1	A	147	THR	O-C-N	5.31	127.75	122.12
1	A	209	VAL	N-CA-C	-5.31	105.35	110.72
1	A	97	GLY	N-CA-C	-5.30	107.90	115.64
1	A	177	ASN	CA-C-N	5.30	130.53	122.11
1	A	177	ASN	C-N-CA	5.30	130.53	122.11
1	A	226	ILE	O-C-N	5.29	128.91	123.20
1	A	145	LEU	CB-CA-C	5.27	119.39	110.79
1	A	151	ARG	O-C-N	5.26	127.78	122.09
1	A	147	THR	CA-C-N	5.24	127.57	120.44
1	A	147	THR	C-N-CA	5.24	127.57	120.44
1	A	144	GLU	CG-CD-OE1	5.24	130.44	118.40
1	A	35	VAL	CA-C-O	-5.23	114.89	120.59
1	A	173	PRO	N-CA-CB	5.22	108.15	103.19
1	A	292	ASP	CA-C-N	-5.22	113.85	121.99
1	A	292	ASP	C-N-CA	-5.22	113.85	121.99
1	A	112	LYS	CA-CB-CG	-5.21	103.67	114.10
1	A	105	LEU	CA-C-O	-5.21	114.67	120.66
1	A	44	GLU	CB-CG-CD	5.20	121.44	112.60
1	A	240	LEU	CA-C-N	5.20	130.24	121.14
1	A	240	LEU	C-N-CA	5.20	130.24	121.14
1	A	147	THR	CA-CB-CG2	5.20	119.33	110.50
1	A	105	LEU	O-C-N	5.18	129.59	123.27
1	A	48	ASN	O-C-N	5.18	127.61	122.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	167	LYS	CA-C-N	-5.17	114.86	122.83
1	A	167	LYS	C-N-CA	-5.17	114.86	122.83
1	A	249	TYR	O-C-N	-5.17	115.51	122.23
1	A	33	PHE	CA-C-O	5.16	126.91	121.33
1	A	150	ARG	NE-CZ-NH2	-5.14	114.58	119.20
1	A	249	TYR	N-CA-CB	-5.14	102.55	110.20
1	A	231	ILE	CA-C-N	5.13	131.34	121.54
1	A	231	ILE	C-N-CA	5.13	131.34	121.54
1	A	261	TYR	N-CA-CB	-5.12	102.65	110.07
1	A	291	ARG	NE-CZ-NH2	-5.11	114.61	119.20
1	A	296	GLU	O-C-N	5.10	127.39	122.09
1	A	214	ARG	CA-C-O	-5.07	115.18	120.55
1	A	109	ALA	CA-C-N	5.01	126.95	120.44
1	A	109	ALA	C-N-CA	5.01	126.95	120.44
1	A	199	TRP	CA-C-O	-5.00	115.30	120.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	128	ARG	Sidechain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2316	0	2328	60	0
2	A	10	0	8	0	0
3	A	10	0	10	0	0
4	A	205	0	0	13	0
All	All	2541	0	2346	60	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 13.

All (60) 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:146:ASP:HB2	4:A:464:HOH:O	1.23	1.38
1:A:295:LYS:HE3	1:A:306:LYS:O	1.46	1.15
1:A:295:LYS:CE	1:A:306:LYS:O	1.96	1.13
1:A:127:LYS:HB2	1:A:127:LYS:NZ	1.56	1.12
1:A:22:LYS:NZ	1:A:22:LYS:HB2	1.51	1.11
1:A:127:LYS:HZ3	1:A:127:LYS:CB	1.68	1.05
1:A:127:LYS:HB2	1:A:127:LYS:HZ3	0.80	0.96
1:A:22:LYS:HB2	1:A:22:LYS:HZ3	1.17	0.96
1:A:3:LEU:HD11	1:A:59:LYS:CG	1.96	0.95
1:A:3:LEU:HD11	1:A:59:LYS:HG3	1.52	0.92
1:A:160:LYS:HD3	1:A:166:GLU:HG2	1.53	0.91
1:A:127:LYS:NZ	1:A:127:LYS:CB	2.23	0.90
1:A:3:LEU:CD1	1:A:59:LYS:CG	2.50	0.88
1:A:3:LEU:CD1	1:A:59:LYS:HG3	2.04	0.86
1:A:146:ASP:CB	4:A:464:HOH:O	1.93	0.85
1:A:3:LEU:CD1	1:A:59:LYS:HG2	2.13	0.77
1:A:295:LYS:HE3	1:A:306:LYS:C	2.09	0.77
1:A:295:LYS:HE2	1:A:306:LYS:O	1.83	0.76
1:A:22:LYS:NZ	1:A:22:LYS:CB	2.36	0.75
1:A:26:LYS:HA	1:A:29:LYS:HD2	1.70	0.73
1:A:115:GLU:OE1	4:A:434:HOH:O	2.13	0.65
1:A:3:LEU:HD11	1:A:59:LYS:HG2	1.71	0.63
1:A:273:LYS:HB2	1:A:275:VAL:HG23	1.81	0.63
1:A:160:LYS:HD3	1:A:166:GLU:CG	2.28	0.61
1:A:75:VAL:O	1:A:79:ARG:HG2	1.99	0.61
1:A:22:LYS:HE3	4:A:475:HOH:O	2.02	0.60
1:A:3:LEU:HD12	1:A:59:LYS:HG2	1.82	0.59
1:A:59:LYS:HB3	1:A:59:LYS:HZ2	1.67	0.58
1:A:144:GLU:H	1:A:144:GLU:CD	2.13	0.56
1:A:4:LYS:HD2	1:A:36:ILE:HD11	1.88	0.56
1:A:84:LYS:HE2	4:A:487:HOH:O	2.05	0.55
1:A:306:LYS:C	4:A:435:HOH:O	2.49	0.54
1:A:146:ASP:CG	4:A:464:HOH:O	2.34	0.54
1:A:127:LYS:HE2	4:A:360:HOH:O	2.07	0.54
1:A:127:LYS:CB	1:A:127:LYS:HZ2	2.17	0.52
1:A:59:LYS:HD3	4:A:422:HOH:O	2.11	0.51
1:A:131:ASP:OD2	1:A:133:LYS:HE3	2.11	0.50
1:A:296:GLU:HG3	4:A:337:HOH:O	2.11	0.50
1:A:84:LYS:HD2	1:A:270:TRP:CE2	2.47	0.49
1:A:138:MET:CE	1:A:201:ILE:HG12	2.41	0.49
1:A:272:ALA:C	1:A:273:LYS:HG3	2.39	0.48
1:A:300:LYS:CB	1:A:300:LYS:NZ	2.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:LYS:HZ2	1:A:300:LYS:HB2	1.80	0.46
1:A:212:GLY:O	1:A:216:THR:HG23	2.15	0.46
1:A:175:LYS:NZ	4:A:342:HOH:O	2.21	0.46
1:A:112:LYS:HD3	1:A:112:LYS:HA	1.25	0.45
1:A:68:PRO:HB3	1:A:91:GLN:O	2.16	0.45
1:A:143:ASN:OD1	1:A:149:ARG:HG3	2.18	0.44
1:A:146:ASP:OD2	1:A:150:ARG:HD3	2.17	0.44
1:A:112:LYS:HE2	1:A:115:GLU:OE1	2.18	0.44
1:A:201:ILE:HD11	1:A:216:THR:HG21	1.99	0.44
1:A:286:VAL:HG22	1:A:287:VAL:N	2.33	0.43
1:A:96:LYS:HB2	1:A:96:LYS:HE3	1.29	0.43
1:A:22:LYS:HB3	1:A:22:LYS:HE2	1.49	0.42
1:A:150:ARG:CG	4:A:465:HOH:O	2.68	0.41
1:A:253:LEU:HA	1:A:254:PRO:HD3	1.92	0.41
1:A:84:LYS:HD2	1:A:270:TRP:CZ2	2.56	0.41
1:A:22:LYS:CE	4:A:475:HOH:O	2.65	0.40
1:A:270:TRP:HA	1:A:275:VAL:O	2.21	0.40
1:A:59:LYS:HB3	1:A:59:LYS:NZ	2.34	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	303/306 (99%)	296 (98%)	3 (1%)	4 (1%)	9 1

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	232	ASN
1	A	254	PRO
1	A	89	ASP

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Mol	Chain	Res	Type
1	A	304	GLY

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	243/246 (99%)	230 (95%)	13 (5%)	20 2

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LYS
1	A	22	LYS
1	A	59	LYS
1	A	79	ARG
1	A	96	LYS
1	A	127	LYS
1	A	144	GLU
1	A	175	LYS
1	A	204	MET
1	A	271	VAL
1	A	285	ASP
1	A	299	GLU
1	A	300	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	GLN
1	A	117	GLN
1	A	119	GLN
1	A	126	GLN
1	A	244	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

2 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	ARA	A	307[A]	-	10,10,10	2.71	3 (30%)	14,14,14	1.54	3 (21%)
3	ARB	A	308[B]	-	10,10,10	1.89	4 (40%)	14,14,14	1.86	2 (14%)

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	ARA	A	307[A]	-	-	-	0/1/1/1
3	ARB	A	308[B]	-	-	-	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	307[A]	ARA	C3-C2	-5.39	1.38	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	307[A]	ARA	C4-C3	4.20	1.58	1.52
2	A	307[A]	ARA	O5-C5	4.03	1.50	1.43
3	A	308[B]	ARB	C4-C3	2.79	1.56	1.52
3	A	308[B]	ARB	O4-C4	2.57	1.48	1.43
3	A	308[B]	ARB	C5-C4	2.48	1.58	1.52
3	A	308[B]	ARB	C3-C2	-2.15	1.46	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	308[B]	ARB	C5-O5-C1	4.88	122.60	112.46
3	A	308[B]	ARB	O2-C2-C3	-2.81	103.75	110.38
2	A	307[A]	ARA	O3-C3-C4	-2.56	104.84	110.05
2	A	307[A]	ARA	O3-C3-C2	2.37	115.97	110.38
2	A	307[A]	ARA	C4-C3-C2	2.08	114.51	110.86

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.