



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:56 AM UTC

PDB ID : 6HVG / pdb_00006hvg
Title : Crystal Structure of Truncated Alternansucrase from *Leuconostoc mesenteroides* NRRL B-1355
Authors : Molina, M.; Cioci, G.; Moulis, C.; Remaud-Simeon, M.
Deposited on : 2018-10-11
Resolution : 2.80 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

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
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

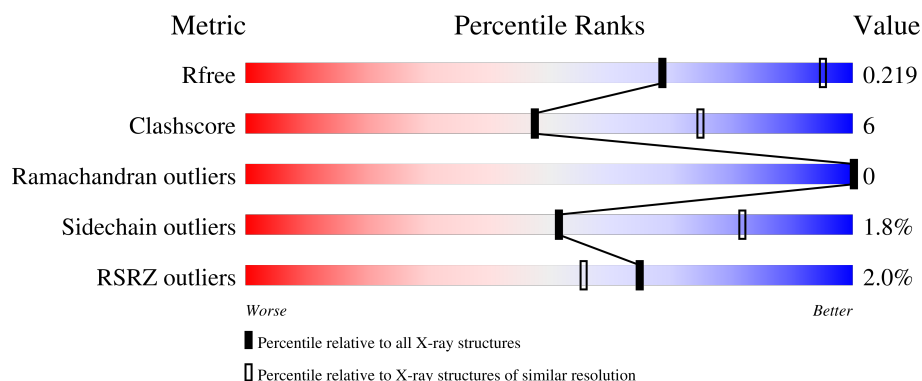
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1435	80% 9% 11%
1	B	1435	68% 13% 18%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 19473 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 Alternansucrase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1278	Total	C	N	O	S	0	0	0
			10013	6256	1706	2025	26			
1	B	1174	Total	C	N	O	S	0	0	0
			9180	5744	1563	1847	26			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	15	MET	-	initiating methionine	UNP Q9RE05
A	16	ALA	-	expression tag	UNP Q9RE05
A	17	HIS	-	expression tag	UNP Q9RE05
A	18	HIS	-	expression tag	UNP Q9RE05
A	19	HIS	-	expression tag	UNP Q9RE05
A	20	HIS	-	expression tag	UNP Q9RE05
A	21	HIS	-	expression tag	UNP Q9RE05
A	22	HIS	-	expression tag	UNP Q9RE05
A	23	VAL	-	expression tag	UNP Q9RE05
A	24	THR	-	expression tag	UNP Q9RE05
A	25	SER	-	expression tag	UNP Q9RE05
A	26	LEU	-	expression tag	UNP Q9RE05
A	27	TYR	-	expression tag	UNP Q9RE05
A	28	LYS	-	expression tag	UNP Q9RE05
A	29	LYS	-	expression tag	UNP Q9RE05
A	30	ALA	-	expression tag	UNP Q9RE05
A	31	GLY	-	expression tag	UNP Q9RE05
A	32	SER	-	expression tag	UNP Q9RE05
A	33	ALA	-	expression tag	UNP Q9RE05
A	34	ALA	-	expression tag	UNP Q9RE05
A	35	ALA	-	expression tag	UNP Q9RE05
A	36	PRO	-	expression tag	UNP Q9RE05
A	37	PHE	-	expression tag	UNP Q9RE05
A	38	THR	-	expression tag	UNP Q9RE05
A	1426	LYS	-	expression tag	UNP Q9RE05

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1427	GLY	-	expression tag	UNP Q9RE05
A	1428	GLY	-	expression tag	UNP Q9RE05
A	1429	ARG	-	expression tag	UNP Q9RE05
A	1430	ALA	-	expression tag	UNP Q9RE05
A	1431	ASP	-	expression tag	UNP Q9RE05
A	1432	PRO	-	expression tag	UNP Q9RE05
A	1433	ALA	-	expression tag	UNP Q9RE05
A	1434	PHE	-	expression tag	UNP Q9RE05
A	1435	LEU	-	expression tag	UNP Q9RE05
A	1436	TYR	-	expression tag	UNP Q9RE05
A	1437	LYS	-	expression tag	UNP Q9RE05
A	1438	VAL	-	expression tag	UNP Q9RE05
A	1439	VAL	-	expression tag	UNP Q9RE05
A	1440	SER	-	expression tag	UNP Q9RE05
A	1441	ALA	-	expression tag	UNP Q9RE05
A	1442	TRP	-	expression tag	UNP Q9RE05
A	1443	SER	-	expression tag	UNP Q9RE05
A	1444	HIS	-	expression tag	UNP Q9RE05
A	1445	PRO	-	expression tag	UNP Q9RE05
A	1446	GLN	-	expression tag	UNP Q9RE05
A	1447	PHE	-	expression tag	UNP Q9RE05
A	1448	GLU	-	expression tag	UNP Q9RE05
A	1449	LYS	-	expression tag	UNP Q9RE05
B	15	MET	-	initiating methionine	UNP Q9RE05
B	16	ALA	-	expression tag	UNP Q9RE05
B	17	HIS	-	expression tag	UNP Q9RE05
B	18	HIS	-	expression tag	UNP Q9RE05
B	19	HIS	-	expression tag	UNP Q9RE05
B	20	HIS	-	expression tag	UNP Q9RE05
B	21	HIS	-	expression tag	UNP Q9RE05
B	22	HIS	-	expression tag	UNP Q9RE05
B	23	VAL	-	expression tag	UNP Q9RE05
B	24	THR	-	expression tag	UNP Q9RE05
B	25	SER	-	expression tag	UNP Q9RE05
B	26	LEU	-	expression tag	UNP Q9RE05
B	27	TYR	-	expression tag	UNP Q9RE05
B	28	LYS	-	expression tag	UNP Q9RE05
B	29	LYS	-	expression tag	UNP Q9RE05
B	30	ALA	-	expression tag	UNP Q9RE05
B	31	GLY	-	expression tag	UNP Q9RE05
B	32	SER	-	expression tag	UNP Q9RE05
B	33	ALA	-	expression tag	UNP Q9RE05

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Chain	Residue	Modelled	Actual	Comment	Reference
B	34	ALA	-	expression tag	UNP Q9RE05
B	35	ALA	-	expression tag	UNP Q9RE05
B	36	PRO	-	expression tag	UNP Q9RE05
B	37	PHE	-	expression tag	UNP Q9RE05
B	38	THR	-	expression tag	UNP Q9RE05
B	1426	LYS	-	expression tag	UNP Q9RE05
B	1427	GLY	-	expression tag	UNP Q9RE05
B	1428	GLY	-	expression tag	UNP Q9RE05
B	1429	ARG	-	expression tag	UNP Q9RE05
B	1430	ALA	-	expression tag	UNP Q9RE05
B	1431	ASP	-	expression tag	UNP Q9RE05
B	1432	PRO	-	expression tag	UNP Q9RE05
B	1433	ALA	-	expression tag	UNP Q9RE05
B	1434	PHE	-	expression tag	UNP Q9RE05
B	1435	LEU	-	expression tag	UNP Q9RE05
B	1436	TYR	-	expression tag	UNP Q9RE05
B	1437	LYS	-	expression tag	UNP Q9RE05
B	1438	VAL	-	expression tag	UNP Q9RE05
B	1439	VAL	-	expression tag	UNP Q9RE05
B	1440	SER	-	expression tag	UNP Q9RE05
B	1441	ALA	-	expression tag	UNP Q9RE05
B	1442	TRP	-	expression tag	UNP Q9RE05
B	1443	SER	-	expression tag	UNP Q9RE05
B	1444	HIS	-	expression tag	UNP Q9RE05
B	1445	PRO	-	expression tag	UNP Q9RE05
B	1446	GLN	-	expression tag	UNP Q9RE05
B	1447	PHE	-	expression tag	UNP Q9RE05
B	1448	GLU	-	expression tag	UNP Q9RE05
B	1449	LYS	-	expression tag	UNP Q9RE05

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Ca 1 1	0	0
2	B	1	Total Ca 1 1	0	0

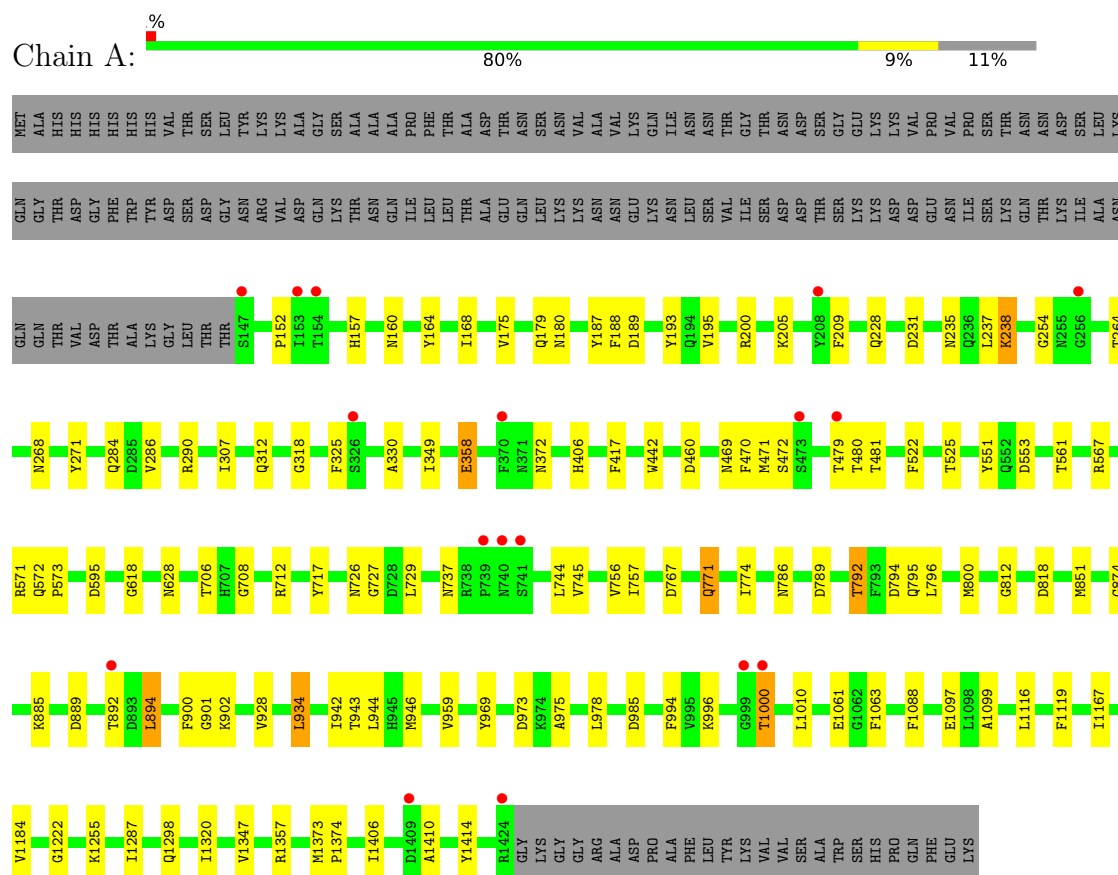
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	166	Total 166	O 166	0	0
3	B	112	Total 112	O 112	0	0

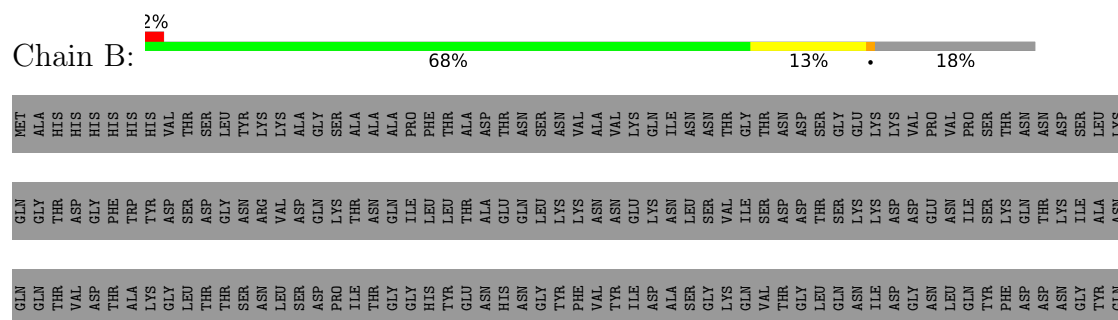
3 Residue-property plots

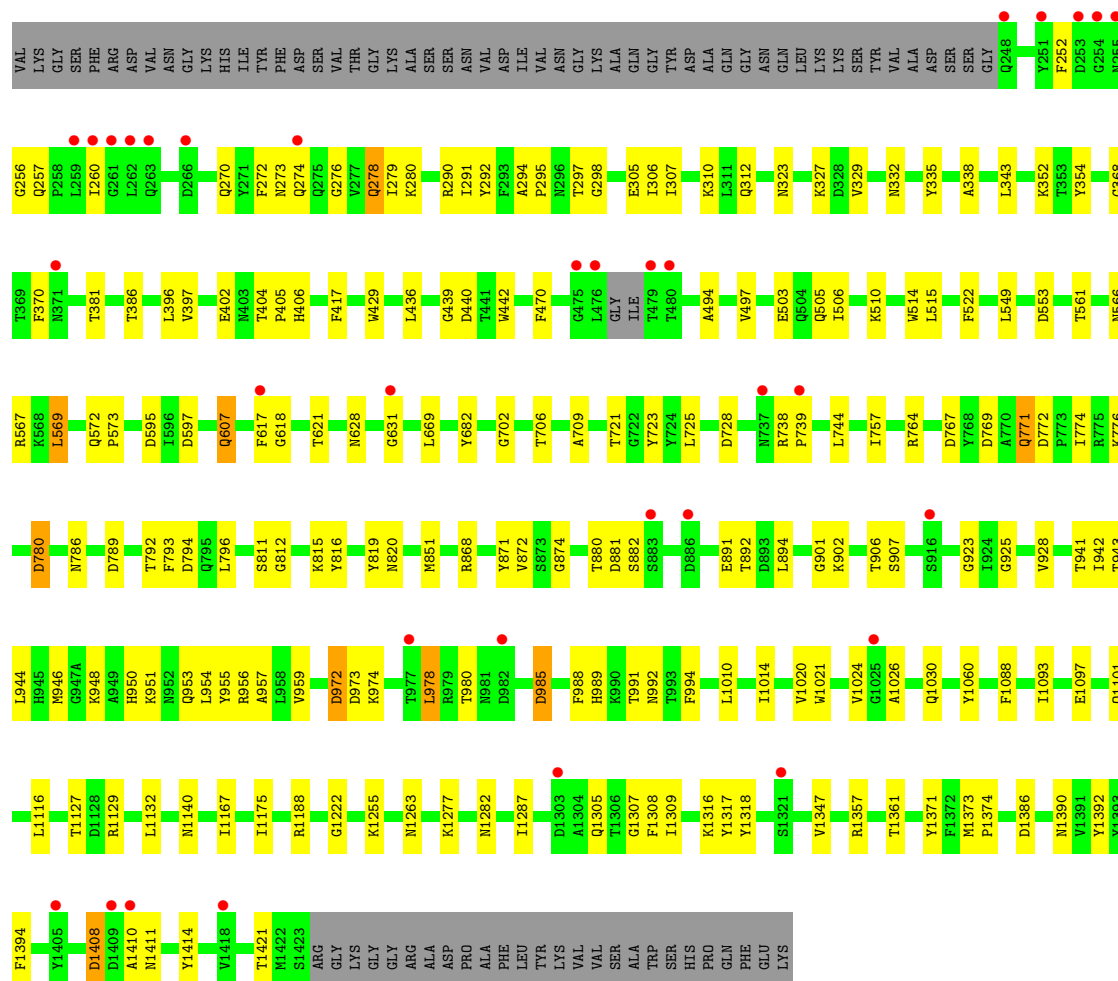
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Alternansucrase



• Molecule 1: Alternansucrase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.23Å 134.80Å 237.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.82 – 2.80 47.82 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.3 (47.82-2.80) 99.3 (47.82-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.98 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.207 , 0.237 (Not available) , 0.219	Depositor DCC
R_{free} test set	4021 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	60.3	Xtriage
Anisotropy	0.189	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 30.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	19473	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.23% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.66	0/10225	0.81	5/13863 (0.0%)
1	B	0.68	0/9374	0.83	4/12709 (0.0%)
All	All	0.67	0/19599	0.82	9/26572 (0.0%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	902	LYS	N-CA-C	7.92	119.55	111.07
1	A	889	ASP	N-CA-C	5.95	118.28	111.02
1	A	1000	THR	N-CA-C	5.73	115.97	107.88
1	A	902	LYS	N-CA-C	5.46	116.91	111.07
1	B	278	GLN	N-CA-C	5.37	118.69	110.20
1	A	900	PHE	N-CA-C	5.26	119.78	112.90
1	A	729	LEU	N-CA-C	-5.25	106.13	112.54
1	B	728	ASP	N-CA-C	5.18	116.74	111.14
1	B	1014	ILE	N-CA-C	5.02	114.81	107.88

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	10013	0	9418	79	0
1	B	9180	0	8628	128	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	166	0	0	5	0
3	B	112	0	0	1	0
All	All	19473	0	18046	207	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 6.

All (207) 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:944:LEU:HB3	1:A:946:MET:HE3	1.72	0.70
1:B:944:LEU:HB3	1:B:946:MET:HE3	1.74	0.70
1:B:290:ARG:NH2	1:B:305:GLU:OE1	2.26	0.67
1:B:948:LYS:O	1:B:951:LYS:HG3	1.95	0.67
1:A:706:THR:HG21	3:A:2212:HOH:O	1.94	0.66
1:B:617:PHE:CE1	1:B:1132:LEU:HD12	2.31	0.65
1:B:972:ASP:OD2	1:B:972:ASP:N	2.29	0.65
1:A:892:THR:HG23	1:A:934:LEU:HD21	1.79	0.64
1:B:402:GLU:O	1:B:405:PRO:HD2	1.98	0.64
1:A:1222:GLY:O	1:A:1255:LYS:HE2	1.99	0.63
1:A:152:PRO:HA	1:A:180:ASN:O	1.99	0.61
1:A:417:PHE:HB2	1:A:1287:ILE:HD13	1.81	0.61
1:B:874:GLY:HA3	1:B:901:GLY:O	2.01	0.60
1:B:386:THR:HG22	3:B:2152:HOH:O	2.00	0.60
1:A:874:GLY:HA3	1:A:901:GLY:O	2.02	0.60
1:B:767:ASP:O	1:B:771:GLN:NE2	2.34	0.60
1:B:417:PHE:HB2	1:B:1287:ILE:HD13	1.83	0.59
1:B:1222:GLY:O	1:B:1255:LYS:HE2	2.01	0.59
1:B:767:ASP:C	1:B:771:GLN:HE21	2.11	0.58
1:B:273:ASN:OD1	1:B:276:GLY:N	2.35	0.58
1:B:567:ARG:O	1:B:569:LEU:HD13	2.04	0.58
1:B:868:ARG:HA	1:B:872:VAL:CG2	2.33	0.58
1:B:270:GLN:HG2	1:B:298:GLY:O	2.04	0.58
1:B:706:THR:HA	1:B:819:TYR:O	2.03	0.58
1:B:1408:ASP:HB2	1:B:1414:TYR:HE2	1.68	0.58
1:B:429:TRP:CH2	1:B:515:LEU:HD23	2.38	0.57
1:A:786:ASN:O	1:A:789:ASP:HB2	2.05	0.57
1:B:956:ARG:HG2	1:B:1024:VAL:HG12	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1347:VAL:HG11	1:A:1357:ARG:HD2	1.86	0.56
1:A:726:ASN:O	1:A:727:GLY:C	2.48	0.56
1:A:290:ARG:O	1:A:318:GLY:HA2	2.06	0.56
1:B:1347:VAL:HG11	1:B:1357:ARG:HD2	1.87	0.56
1:A:284:GLN:HG3	1:A:286:VAL:CG2	2.35	0.56
1:A:894:LEU:HD23	1:A:942:ILE:HD13	1.88	0.56
1:A:231:ASP:HB3	1:A:237:LEU:HD21	1.89	0.55
1:A:330:ALA:O	1:A:1410:ALA:HA	2.06	0.55
1:A:745:VAL:HG22	1:A:756:VAL:HG21	1.88	0.55
1:B:506:ILE:HG13	1:B:514:TRP:HZ2	1.71	0.55
1:B:1410:ALA:O	1:B:1411:ASN:CB	2.54	0.55
1:B:506:ILE:O	1:B:510:LYS:N	2.37	0.55
1:B:1373:MET:HB3	1:B:1374:PRO:CD	2.37	0.55
1:B:1373:MET:HB3	1:B:1374:PRO:HD2	1.87	0.55
1:A:745:VAL:HG22	1:A:756:VAL:CG2	2.37	0.55
1:A:996:LYS:HB2	1:A:1000:THR:H	1.72	0.55
1:B:954:LEU:HD23	1:B:1024:VAL:HG21	1.88	0.54
1:A:284:GLN:HG3	1:A:286:VAL:HG23	1.87	0.54
1:A:708:GLY:O	1:A:712:ARG:NH2	2.40	0.54
1:B:280:LYS:HB3	1:B:295:PRO:O	2.08	0.53
1:B:950:HIS:ND1	1:B:955:TYR:OH	2.38	0.53
1:B:617:PHE:CE1	1:B:621:THR:HG21	2.44	0.53
1:B:1305:GLN:O	1:B:1318:TYR:HB3	2.08	0.53
1:B:252:PHE:HA	1:B:257:GLN:O	2.08	0.53
1:A:175:VAL:CG1	1:A:179:GLN:HG3	2.39	0.52
1:A:271:TYR:CD1	1:A:286:VAL:HG21	2.44	0.52
1:A:358:GLU:H	1:A:358:GLU:CD	2.16	0.52
1:B:1371:TYR:HB2	1:B:1394:PHE:CZ	2.45	0.52
1:A:175:VAL:HG11	1:A:179:GLN:HG3	1.92	0.52
1:A:472:SER:HB3	1:A:479:THR:HA	1.91	0.52
1:B:396:LEU:HB2	1:B:1309:ILE:HB	1.91	0.52
1:A:188:PHE:HA	1:A:193:TYR:O	2.10	0.52
1:B:631:GLY:HA3	1:B:669:LEU:O	2.09	0.52
1:B:793:PHE:O	1:B:794:ASP:C	2.53	0.52
1:B:894:LEU:HD13	1:B:942:ILE:HD13	1.92	0.51
1:B:1410:ALA:O	1:B:1411:ASN:HB2	2.10	0.51
1:A:756:VAL:HG22	3:A:2157:HOH:O	2.10	0.51
1:B:553:ASP:OD1	1:B:567:ARG:NH1	2.43	0.51
1:B:851:MET:HE1	1:B:1088:PHE:HZ	1.75	0.51
1:B:989:HIS:O	1:B:992:ASN:ND2	2.43	0.51
1:B:354:TYR:HE1	1:B:381:THR:HG22	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:738:ARG:HG3	1:B:739:PRO:HD2	1.93	0.51
1:B:973:ASP:OD1	1:B:974:LYS:HG3	2.11	0.51
1:B:335:TYR:CD2	1:B:343:LEU:HD12	2.46	0.51
1:A:706:THR:CG2	1:A:818:ASP:OD2	2.59	0.51
1:A:851:MET:HE1	1:A:1088:PHE:HZ	1.76	0.51
1:A:522:PHE:O	1:A:525:THR:HB	2.11	0.50
1:B:631:GLY:O	1:B:1167:ILE:HA	2.12	0.50
1:B:503:GLU:OE1	1:B:503:GLU:HA	2.12	0.50
1:A:1097:GLU:HA	1:A:1167:ILE:HB	1.94	0.50
1:A:595:ASP:HB2	3:A:2160:HOH:O	2.12	0.49
1:B:989:HIS:HB2	1:B:991:THR:O	2.12	0.49
1:B:617:PHE:CZ	1:B:1132:LEU:HD12	2.47	0.49
1:B:506:ILE:HG13	1:B:514:TRP:CZ2	2.48	0.49
1:B:764:ARG:NH1	1:B:769:ASP:O	2.45	0.49
1:B:819:TYR:O	1:B:820:ASN:HB2	2.12	0.49
1:A:189:ASP:OD1	1:A:200:ARG:NH2	2.46	0.49
1:A:470:PHE:CD2	1:A:522:PHE:HB2	2.48	0.49
1:B:780:ASP:OD2	1:B:815:LYS:HE2	2.13	0.49
1:A:1298:GLN:NE2	1:A:1320:ILE:O	2.46	0.48
1:B:868:ARG:HA	1:B:872:VAL:HG23	1.94	0.48
1:B:874:GLY:CA	1:B:901:GLY:O	2.61	0.48
1:B:1097:GLU:HA	1:B:1167:ILE:HB	1.95	0.48
1:A:325:PHE:CE1	1:A:349:ILE:HD11	2.49	0.48
1:B:252:PHE:CD2	1:B:256:GLY:O	2.67	0.48
1:B:406:HIS:HB3	1:B:442:TRP:CE3	2.48	0.48
1:B:549:LEU:HD23	1:B:1188:ARG:HA	1.95	0.48
1:B:279:ILE:HD13	1:B:279:ILE:N	2.28	0.48
1:B:470:PHE:CD2	1:B:522:PHE:HB2	2.49	0.48
1:A:238:LYS:HG2	1:A:254:GLY:O	2.14	0.47
1:B:307:ILE:HG13	1:B:312:GLN:NE2	2.28	0.47
1:A:205:LYS:HE3	1:A:235:ASN:OD1	2.13	0.47
1:A:264:THR:HA	1:A:268:ASN:O	2.14	0.47
1:B:894:LEU:CD1	1:B:942:ILE:HD13	2.43	0.47
1:B:1026:ALA:HB1	1:B:1030:GLN:NE2	2.28	0.47
1:A:717:TYR:OH	1:A:885:LYS:HG2	2.14	0.47
1:B:953:GLN:N	1:B:980:THR:OG1	2.48	0.47
1:A:1063:PHE:CG	1:A:1099:ALA:HB2	2.50	0.47
1:B:294:ALA:HB3	1:B:297:THR:HG23	1.97	0.46
1:B:368:GLY:HA3	1:B:370:PHE:CE2	2.51	0.46
1:B:404:THR:N	1:B:405:PRO:CD	2.78	0.46
1:B:1308:PHE:HA	1:B:1316:LYS:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:252:PHE:HB3	1:B:256:GLY:C	2.40	0.46
1:B:272:PHE:HE1	1:B:278:GLN:HB2	1.81	0.46
1:B:252:PHE:HD2	1:B:257:GLN:C	2.22	0.46
1:B:906:THR:HG22	1:B:907:SER:N	2.29	0.46
1:A:189:ASP:OD2	1:A:200:ARG:NH2	2.49	0.46
1:A:228:GLN:NE2	3:A:2110:HOH:O	2.48	0.46
1:B:291:ILE:HG22	1:B:292:TYR:N	2.30	0.46
1:A:157:HIS:CE1	1:A:168:ILE:HB	2.51	0.46
1:A:744:LEU:HD11	1:A:757:ILE:HB	1.98	0.46
1:B:291:ILE:CG2	1:B:292:TYR:N	2.79	0.46
1:B:572:GLN:HB2	1:B:573:PRO:HD2	1.98	0.46
1:A:792:THR:HG22	1:A:795:GLN:H	1.80	0.45
1:B:1390:ASN:HB3	1:B:1421:THR:HG22	1.98	0.45
1:A:470:PHE:CD2	1:A:471:MET:HE2	2.52	0.45
1:A:187:TYR:HB2	1:A:209:PHE:CZ	2.52	0.45
1:B:702:GLY:HA2	1:B:706:THR:OG1	2.17	0.45
1:B:368:GLY:C	1:B:370:PHE:CE2	2.95	0.45
1:B:505:GLN:HB3	1:B:514:TRP:CZ2	2.52	0.45
1:A:470:PHE:CE2	1:A:522:PHE:HB2	2.52	0.45
1:A:978:LEU:HD22	1:A:994:PHE:HE1	1.81	0.45
1:B:772:ASP:O	1:B:776:LYS:HG3	2.17	0.44
1:B:925:GLY:O	1:B:1020:VAL:HA	2.18	0.44
1:B:815:LYS:HG2	1:B:816:TYR:CZ	2.52	0.44
1:A:470:PHE:HD2	1:A:471:MET:HE2	1.82	0.44
1:B:709:ALA:HB2	1:B:819:TYR:OH	2.17	0.44
1:B:272:PHE:CE1	1:B:278:GLN:HB2	2.53	0.44
1:B:744:LEU:HD11	1:B:757:ILE:HB	2.00	0.44
1:A:307:ILE:HB	1:A:312:GLN:OE1	2.18	0.44
1:A:794:ASP:OD2	1:A:794:ASP:N	2.50	0.44
1:B:793:PHE:O	1:B:796:LEU:N	2.51	0.44
1:B:973:ASP:OD1	1:B:974:LYS:N	2.50	0.44
1:A:1061:GLU:HA	1:A:1097:GLU:HB3	2.00	0.44
1:B:618:GLY:N	1:B:628:ASN:OD1	2.51	0.44
1:B:1277:LYS:HE3	1:B:1282:ASN:HA	2.00	0.44
1:A:969:TYR:HB3	1:A:975:ALA:HB2	2.00	0.43
1:B:494:ALA:O	1:B:497:VAL:HG22	2.17	0.43
1:B:811:SER:O	1:B:1010:LEU:CD2	2.66	0.43
1:B:811:SER:O	1:B:1010:LEU:HD23	2.18	0.43
1:A:553:ASP:OD1	1:A:567:ARG:NH1	2.52	0.43
1:A:978:LEU:HD22	1:A:994:PHE:CE1	2.53	0.43
1:B:406:HIS:HB3	1:B:442:TRP:CZ3	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:894:LEU:CD1	1:B:942:ILE:CD1	2.97	0.43
1:A:572:GLN:HB2	1:A:573:PRO:HD2	2.00	0.43
1:B:294:ALA:HB3	1:B:297:THR:CG2	2.48	0.43
1:B:323:ASN:OD1	1:B:338:ALA:HB2	2.19	0.43
1:B:881:ASP:OD2	1:B:891:GLU:HA	2.19	0.43
1:A:943:THR:CG2	1:A:985:ASP:HB3	2.49	0.42
1:A:561:THR:O	1:A:561:THR:HG22	2.18	0.42
1:B:595:ASP:OD2	1:B:1263:ASN:ND2	2.38	0.42
1:A:708:GLY:O	1:A:712:ARG:NE	2.53	0.42
1:A:1406:ILE:CG2	1:A:1414:TYR:HB2	2.50	0.42
1:A:561:THR:O	1:A:561:THR:CG2	2.67	0.42
1:B:566:ASN:O	1:B:597:ASP:OD2	2.36	0.42
1:B:941:THR:HA	1:B:988:PHE:O	2.20	0.42
1:B:943:THR:CG2	1:B:985:ASP:HB3	2.50	0.42
1:B:1408:ASP:OD1	1:B:1410:ALA:N	2.38	0.42
1:B:812:GLY:HA3	1:B:1010:LEU:HD23	2.01	0.42
1:B:370:PHE:CD2	1:B:370:PHE:N	2.88	0.42
1:A:1116:LEU:HA	1:A:1119:PHE:CZ	2.55	0.42
1:B:956:ARG:CG	1:B:1024:VAL:HG12	2.48	0.41
1:B:682:TYR:CD1	1:B:682:TYR:C	2.98	0.41
1:B:1307:GLY:O	1:B:1317:TYR:HA	2.19	0.41
1:A:551:TYR:CE1	1:A:1184:VAL:HG21	2.55	0.41
1:A:1373:MET:HB3	1:A:1374:PRO:CD	2.50	0.41
1:B:957:ALA:N	1:B:1021:TRP:CE3	2.88	0.41
1:A:469:ASN:OD1	1:A:480:THR:O	2.38	0.41
1:A:205:LYS:CE	1:A:235:ASN:OD1	2.68	0.41
1:A:1406:ILE:HG22	1:A:1414:TYR:HB2	2.01	0.41
1:B:307:ILE:HG13	1:B:312:GLN:HE22	1.86	0.41
1:B:871:TYR:HB3	1:B:923:GLY:O	2.20	0.41
1:B:978:LEU:HG	1:B:994:PHE:CE1	2.56	0.41
1:A:767:ASP:O	1:A:771:GLN:OE1	2.38	0.41
1:B:721:THR:C	1:B:723:TYR:N	2.77	0.41
1:B:1101:GLN:HB2	1:B:1132:LEU:HD22	2.02	0.41
1:B:771:GLN:H	1:B:771:GLN:HG3	1.47	0.41
1:B:1127:THR:O	1:B:1175:ILE:HD11	2.21	0.41
1:A:160:ASN:HA	1:A:164:TYR:O	2.21	0.41
1:A:195:VAL:HG21	1:A:200:ARG:NH2	2.36	0.41
1:B:306:ILE:HD12	1:B:306:ILE:O	2.20	0.41
1:A:330:ALA:HB1	1:A:1410:ALA:O	2.20	0.41
1:A:771:GLN:H	1:A:771:GLN:HG3	1.45	0.41
1:B:1386:ASP:HB3	1:B:1392:TYR:HE1	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1408:ASP:CB	1:B:1414:TYR:HE2	2.32	0.41
1:A:406:HIS:HB3	1:A:442:TRP:CZ3	2.57	0.41
1:B:561:THR:O	1:B:561:THR:CG2	2.69	0.41
1:B:260:ILE:HD13	1:B:274:GLN:O	2.21	0.40
1:B:327:LYS:HA	1:B:332:ASN:O	2.21	0.40
1:B:1060:TYR:HB2	1:B:1093:ILE:HG13	2.03	0.40
1:A:618:GLY:N	1:A:628:ASN:OD1	2.54	0.40
1:B:436:LEU:CD2	1:B:439:GLY:HA2	2.51	0.40
1:B:786:ASN:O	1:B:789:ASP:HB2	2.21	0.40
1:B:792:THR:HG22	1:B:793:PHE:N	2.37	0.40
1:A:796:LEU:HD11	1:A:800:MET:HE2	2.04	0.40
1:A:812:GLY:HA3	1:A:1010:LEU:HD23	2.02	0.40
1:B:561:THR:O	1:B:561:THR:HG22	2.20	0.40
1:A:571:ARG:HB2	3:A:2189:HOH:O	2.22	0.40
1:B:607:GLN:OE1	1:B:1129:ARG:NH2	2.48	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	1276/1435 (89%)	1225 (96%)	51 (4%)	0	100	100
1	B	1170/1435 (82%)	1120 (96%)	50 (4%)	0	100	100
All	All	2446/2870 (85%)	2345 (96%)	101 (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	1067/1213 (88%)	1053 (99%)	14 (1%)	61	86
1	B	971/1213 (80%)	948 (98%)	23 (2%)	43	77
All	All	2038/2426 (84%)	2001 (98%)	37 (2%)	51	82

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	238	LYS
1	A	358	GLU
1	A	372	ASN
1	A	460	ASP
1	A	481	THR
1	A	737	ASN
1	A	771	GLN
1	A	774	ILE
1	A	792	THR
1	A	894	LEU
1	A	928	VAL
1	A	934	LEU
1	A	959	VAL
1	A	973	ASP
1	B	310	LYS
1	B	329	VAL
1	B	352	LYS
1	B	397	VAL
1	B	440	ASP
1	B	569	LEU
1	B	607	GLN
1	B	725	LEU
1	B	771	GLN
1	B	774	ILE
1	B	780	ASP
1	B	880	THR
1	B	882	SER
1	B	892	THR
1	B	928	VAL
1	B	959	VAL
1	B	972	ASP

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Mol	Chain	Res	Type
1	B	978	LEU
1	B	985	ASP
1	B	1116	LEU
1	B	1140	ASN
1	B	1361	THR
1	B	1408	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	270	GLN
1	A	299	ASN
1	A	332	ASN
1	A	532	GLN
1	A	688	ASN
1	A	807	GLN
1	A	938	ASN
1	A	1078	ASN
1	A	1101	GLN
1	A	1324	GLN
1	B	296	ASN
1	B	312	GLN
1	B	807	GLN
1	B	817	ASN
1	B	1078	ASN
1	B	1101	GLN
1	B	1162	ASN
1	B	1203	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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

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

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	1278/1435 (89%)	-0.07	17 (1%) 75 66	37, 60, 94, 131	0
1	B	1174/1435 (81%)	0.15	33 (2%) 55 45	40, 66, 111, 161	0
All	All	2452/2870 (85%)	0.04	50 (2%) 65 56	37, 63, 103, 161	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	147	SER	4.2
1	B	251	TYR	4.2
1	B	1409	ASP	4.0
1	B	259	LEU	3.8
1	A	740	ASN	3.8
1	B	263	GLN	3.7
1	A	370	PHE	3.6
1	B	479	THR	3.6
1	B	262	LEU	3.3
1	B	248	GLN	3.3
1	B	274	GLN	3.3
1	B	739	PRO	3.1
1	B	1025	GLY	3.0
1	B	260	ILE	3.0
1	B	255	ASN	2.9
1	B	480	THR	2.8
1	A	999	GLY	2.8
1	B	371	ASN	2.7
1	A	479	THR	2.7
1	B	1405	TYR	2.7
1	A	739	PRO	2.7
1	B	261	GLY	2.7
1	B	737	ASN	2.6
1	B	475	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	253	ASP	2.5
1	B	631	GLY	2.4
1	B	617	PHE	2.4
1	B	977	THR	2.4
1	B	476	LEU	2.4
1	A	1409	ASP	2.3
1	A	1000	THR	2.3
1	B	1418	VAL	2.3
1	A	256	GLY	2.3
1	A	326	SER	2.3
1	A	208	TYR	2.3
1	B	982	ASP	2.2
1	B	1410	ALA	2.2
1	A	153	ILE	2.2
1	B	1303	ASP	2.2
1	B	1321	SER	2.2
1	B	883	SER	2.2
1	B	266	ASP	2.2
1	A	154	THR	2.1
1	A	473	SER	2.1
1	B	886	ASP	2.1
1	A	741	SER	2.0
1	B	916	SER	2.0
1	A	1424	ARG	2.0
1	B	254	GLY	2.0
1	A	892	THR	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CA	B	2000	1/1	0.96	0.06	56,56,56,56	0
2	CA	A	2000	1/1	1.00	0.05	47,47,47,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.