



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2014 – 03:42 AM GMT

PDB ID : 2D2O
Title : Structure of a complex of Thermoactinomyces vulgaris R-47 alpha-amylase 2 with maltohexaose demonstrates the important role of aromatic residues at the reducing end of the substrate binding cleft
Authors : Ohtaki, A.; Mizuno, M.; Yoshida, H.; Tonozuka, T.; Sakano, Y.; Kamitori, S.
Deposited on : 2005-09-13
Resolution : 2.10 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

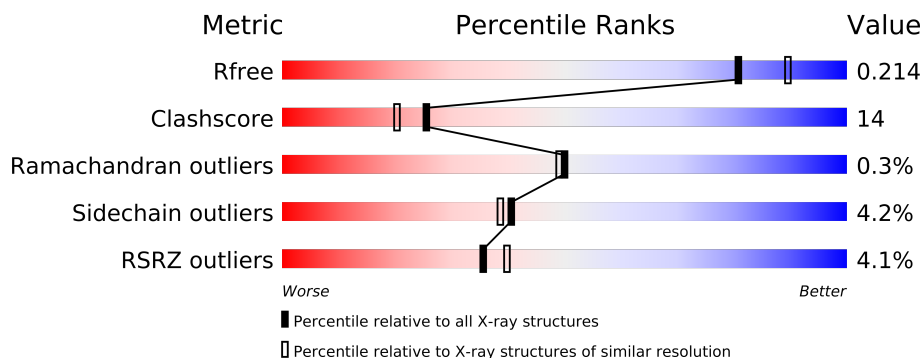
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.10 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	3012 (2.10-2.10)
Clashscore	79885	3649 (2.10-2.10)
Ramachandran outliers	78287	3610 (2.10-2.10)
Sidechain outliers	78261	3611 (2.10-2.10)
RSRZ outliers	66119	3013 (2.10-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	585	
1	B	585	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10668 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Neopullulanase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	585	Total	C	N	O	S	0	0	0
			4776	3056	832	873	15			
1	B	585	Total	C	N	O	S	0	0	0
			4776	3056	832	873	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	325	ASN	ASP	ENGINEERED	UNP Q08751
B	325	ASN	ASP	ENGINEERED	UNP Q08751

- Molecule 2 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	6	Total	C	O	0	0
			67	36	31		
2	B	6	Total	C	O	0	0
			67	36	31		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	325	ASN	ASP	ENGINEERED	UNP Q08751
B	325	ASN	ASP	ENGINEERED	UNP Q08751

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 4 is water.

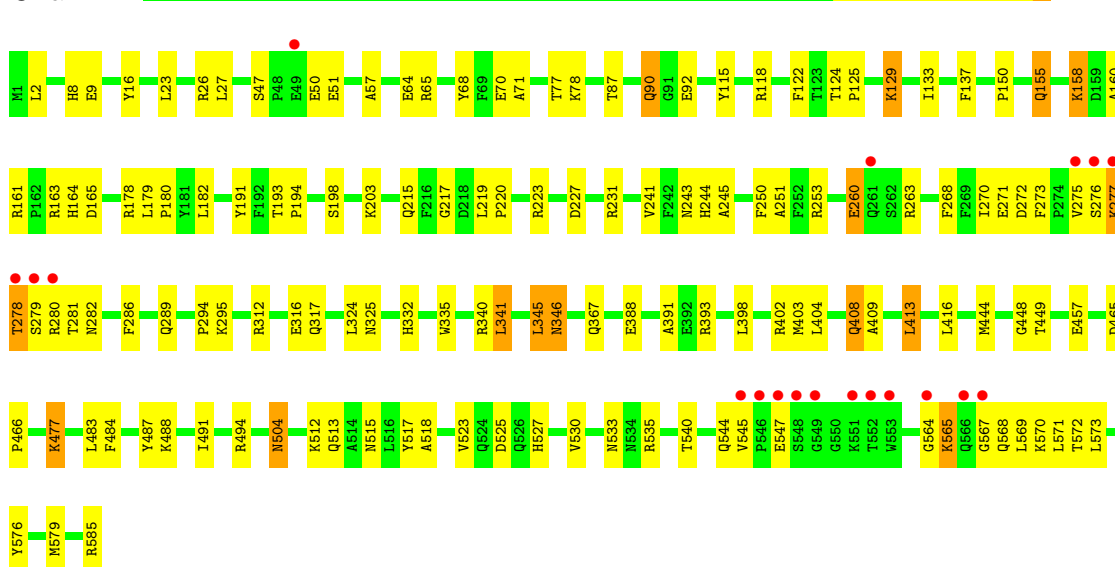
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	498	Total 498	O 498	0	0
4	B	482	Total 482	O 482	0	0

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.

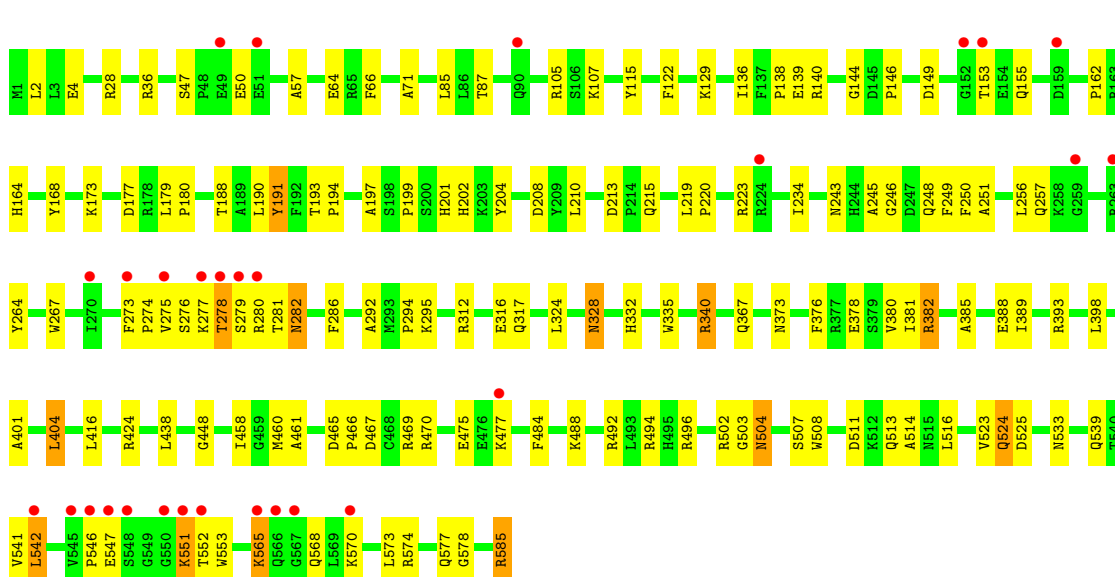
• Molecule 1: Neopullulanase 2

Chain A:



• Molecule 1: Neopullulanase 2

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	113.69Å 118.71Å 112.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.00 – 2.10 46.37 – 1.98	Depositor EDS
% Data completeness (in resolution range)	(Not available) (46.00-2.10) 93.7 (46.37-1.98)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.71 (at 1.98Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.184 , 0.217 0.179 , 0.214	Depositor DCC
R_{free} test set	6518 reflections (8.28%)	DCC
Wilson B-factor (Å ²)	26.5	Xtriage
Anisotropy	0.329	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 48.2	EDS
Estimated twinning fraction	0.016 for l,-k,h	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 98994 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10668	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	0.35	0/4906	0.59	1/6641 (0.0%)
1	B	0.34	0/4906	0.59	0/6641
All	All	0.35	0/9812	0.59	1/13282 (0.0%)

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
2	A	1	0
2	B	1	0
All	All	2	0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	137	PHE	N-CA-C	-5.21	96.92	111.00

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	706	GLC	C1
2	B	706	GLC	C1

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4776	0	4609	132	0
1	B	4776	0	4609	130	0
2	A	67	0	57	5	0
2	B	67	0	57	4	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	498	0	0	15	0
4	B	482	0	0	15	0
All	All	10668	0	9332	259	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 14.

All (259) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:2:LEU:HD21	1:B:2:LEU:HD21	1.25	1.14
1:A:129:LYS:HD2	1:A:129:LYS:H	1.24	1.02
1:A:525:ASP:HB3	1:A:585:ARG:HD3	1.49	0.94
1:A:278:THR:HG22	1:A:279:SER:H	1.35	0.91
1:A:544:GLN:HE22	1:A:568:GLN:HE21	1.20	0.90
1:A:408:GLN:HE21	1:A:408:GLN:H	1.16	0.89
1:B:511:ASP:OD2	1:B:513:GLN:HG2	1.71	0.89
1:A:129:LYS:HD2	1:A:129:LYS:N	1.91	0.83
1:B:28:ARG:HD3	1:B:66:PHE:CD2	2.12	0.83
1:B:552:THR:HB	4:B:1045:HOH:O	1.77	0.83
1:A:565:LYS:HD3	1:A:565:LYS:H	1.44	0.82
1:A:544:GLN:NE2	1:A:568:GLN:HE21	1.79	0.80
1:B:213:ASP:OD1	1:B:215:GLN:HG2	1.83	0.79
1:A:544:GLN:HE22	1:A:568:GLN:NE2	1.80	0.78
1:A:281:THR:H	1:A:289:GLN:HE22	1.32	0.76
1:B:129:LYS:HD2	1:B:502:ARG:NH2	2.00	0.76
1:B:533:ASN:HD21	1:B:539:GLN:HE21	1.31	0.75
1:A:2:LEU:CD2	1:B:2:LEU:HD21	2.13	0.75
1:A:57:ALA:HB2	1:A:71:ALA:HB2	1.68	0.75

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:271:GLU:HG2	1:A:282:ASN:O	1.86	0.75
1:B:202:HIS:HD2	1:B:204:TYR:H	1.35	0.74
1:A:223:ARG:HD3	1:A:317:GLN:HE21	1.53	0.74
1:A:408:GLN:NE2	1:A:408:GLN:H	1.87	0.73
1:A:477:LYS:HA	1:A:477:LYS:HE3	1.71	0.72
1:B:585:ARG:HH11	1:B:585:ARG:HB3	1.52	0.72
1:B:140:ARG:HH12	1:B:201:HIS:HB2	1.55	0.71
1:A:26:ARG:HD3	1:A:70:GLU:OE2	1.91	0.70
1:A:243:ASN:HD22	1:A:244:HIS:HD2	1.39	0.70
1:B:140:ARG:HG2	1:B:469:ARG:O	1.91	0.70
1:A:125:PRO:O	1:A:129:LYS:HE2	1.92	0.70
1:B:484:PHE:CE1	1:B:488:LYS:HD2	2.28	0.68
1:A:2:LEU:HD21	1:B:2:LEU:CD2	2.14	0.68
1:B:243:ASN:HD21	1:B:295:LYS:NZ	1.91	0.68
1:B:542:LEU:HD21	1:B:570:LYS:HG2	1.76	0.67
1:A:312:ARG:O	1:A:316:GLU:HG3	1.94	0.67
1:B:278:THR:HG23	1:B:280:ARG:H	1.59	0.67
1:B:202:HIS:CD2	1:B:204:TYR:H	2.12	0.67
1:A:515:ASN:HB2	1:A:535:ARG:NH2	2.10	0.67
1:B:328:ASN:HD22	1:B:328:ASN:N	1.93	0.67
1:B:188:THR:HG21	4:B:1454:HOH:O	1.95	0.66
1:B:47:SER:HB3	1:B:50:GLU:HG3	1.77	0.66
1:A:409:ALA:O	1:A:413:LEU:HD13	1.96	0.66
1:A:341:LEU:HD22	1:A:345:LEU:HD22	1.77	0.65
1:A:540:THR:OG1	1:A:572:THR:HG22	1.97	0.64
1:B:256:LEU:HD23	1:B:275:VAL:HB	1.79	0.64
1:B:475:GLU:H	1:B:475:GLU:CD	2.00	0.64
1:B:565:LYS:HE2	1:B:568:GLN:HB2	1.78	0.64
1:B:107:LYS:HG3	4:B:1399:HOH:O	1.98	0.64
1:A:133:ILE:N	1:A:133:ILE:HD12	2.12	0.64
2:A:703:GLC:H5	2:A:704:GLC:H62	1.80	0.63
1:A:525:ASP:HB2	4:A:1289:HOH:O	1.98	0.63
1:A:388:GLU:OE1	1:B:107:LYS:HE3	1.98	0.63
1:B:275:VAL:O	1:B:282:ASN:ND2	2.32	0.63
1:A:504:ASN:C	1:A:504:ASN:HD22	2.00	0.62
1:A:340:ARG:NH1	4:A:1365:HOH:O	2.31	0.62
1:B:448:GLY:O	1:B:494:ARG:NH2	2.32	0.61
1:A:465:ASP:OD2	2:A:701:GLC:O2	2.16	0.61
1:B:279:SER:HB3	4:B:1391:HOH:O	2.01	0.61
1:B:504:ASN:HD22	1:B:504:ASN:C	2.04	0.61
2:B:701:GLC:H61	4:B:1204:HOH:O	2.00	0.61
1:A:8:HIS:HD2	1:A:26:ARG:O	1.84	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:340:ARG:NH2	4:B:1415:HOH:O	2.30	0.61
1:B:328:ASN:HD22	1:B:328:ASN:H	1.49	0.60
1:B:340:ARG:HB3	1:B:340:ARG:HH21	1.66	0.60
1:B:585:ARG:HH11	1:B:585:ARG:CB	2.15	0.60
1:A:391:ALA:H	1:A:515:ASN:ND2	2.00	0.60
2:B:703:GLC:H5	2:B:704:GLC:H62	1.83	0.60
1:B:488:LYS:O	1:B:492:ARG:HG3	2.01	0.60
1:A:276:SER:O	1:A:277:LYS:HB2	2.02	0.59
1:B:173:LYS:HD3	1:B:177:ASP:OD2	2.02	0.59
1:A:408:GLN:N	1:A:408:GLN:HE21	1.96	0.58
1:B:467:ASP:OD1	1:B:470:ARG:HD3	2.03	0.58
1:A:564:GLY:HA2	1:A:568:GLN:O	2.03	0.58
1:A:260:GLU:HG3	1:A:273:PHE:CE2	2.38	0.58
1:A:444:MET:HE1	1:A:491:ILE:HG12	1.85	0.57
1:B:514:ALA:HB1	1:B:516:LEU:HD13	1.85	0.57
1:A:219:LEU:HB3	1:A:220:PRO:HD3	1.87	0.57
1:B:382:ARG:HH21	1:B:382:ARG:HB2	1.69	0.57
1:B:525:ASP:HB3	1:B:585:ARG:HD3	1.86	0.57
2:A:701:GLC:H61	4:A:1149:HOH:O	2.03	0.57
1:B:57:ALA:HB2	1:B:71:ALA:HB2	1.86	0.56
1:B:164:HIS:CE1	1:B:466:PRO:HD3	2.41	0.56
1:B:547:GLU:HB2	4:B:1369:HOH:O	2.05	0.56
1:A:124:THR:OG1	1:A:129:LYS:HE3	2.06	0.56
1:B:332:HIS:HD2	1:B:367:GLN:OE1	1.89	0.56
1:A:332:HIS:HD2	1:A:367:GLN:OE1	1.90	0.55
1:A:565:LYS:HD3	1:A:565:LYS:N	2.18	0.55
1:A:545:VAL:HA	1:A:569:LEU:HD21	1.89	0.55
1:A:241:VAL:HG13	1:A:325:ASN:HD22	1.72	0.55
1:A:133:ILE:HD13	1:A:449:THR:CG2	2.37	0.55
1:A:160:ALA:O	1:A:161:ARG:HD2	2.06	0.54
1:B:328:ASN:H	1:B:328:ASN:ND2	2.03	0.54
1:B:282:ASN:C	1:B:282:ASN:HD22	2.10	0.54
1:A:164:HIS:CE1	2:A:701:GLC:H2	2.42	0.54
1:A:278:THR:HG22	1:A:279:SER:N	2.14	0.54
1:B:504:ASN:ND2	1:B:504:ASN:C	2.61	0.53
1:B:416:LEU:H	1:B:416:LEU:HD23	1.72	0.53
1:A:115:TYR:CZ	1:B:295:LYS:HE2	2.43	0.53
1:A:90:GLN:HE21	1:A:90:GLN:HA	1.74	0.53
1:A:26:ARG:HG2	1:A:70:GLU:HG3	1.91	0.53
1:B:382:ARG:CB	1:B:382:ARG:HH21	2.22	0.52
1:A:271:GLU:HG3	1:A:272:ASP:OD2	2.10	0.52
1:B:136:ILE:O	1:B:138:PRO:HD3	2.08	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:193:THR:HB	1:A:194:PRO:CD	2.40	0.52
1:B:470:ARG:HG3	4:B:1112:HOH:O	2.09	0.52
1:A:545:VAL:HA	1:A:569:LEU:CD2	2.39	0.52
1:B:153:THR:HG23	1:B:168:TYR:HA	1.90	0.52
1:A:448:GLY:O	1:A:494:ARG:NH2	2.41	0.52
1:B:28:ARG:HD3	1:B:66:PHE:CE2	2.45	0.52
1:B:492:ARG:O	1:B:496:ARG:HB2	2.09	0.52
1:B:542:LEU:HD12	1:B:568:GLN:HB3	1.90	0.52
1:B:105:ARG:HH11	1:B:105:ARG:HG2	1.74	0.52
1:A:515:ASN:HB2	1:A:535:ARG:HH21	1.74	0.52
1:A:504:ASN:ND2	1:A:504:ASN:C	2.61	0.52
1:A:268:PHE:HB2	1:A:270:ILE:CD1	2.39	0.52
1:B:275:VAL:O	1:B:276:SER:HB3	2.10	0.51
1:A:118:ARG:HG2	4:B:1007:HOH:O	2.10	0.51
1:A:295:LYS:HE2	1:B:115:TYR:CZ	2.46	0.51
1:B:378:GLU:HG3	4:B:1362:HOH:O	2.09	0.51
1:A:158:LYS:HD2	1:A:158:LYS:O	2.11	0.51
1:B:64:GLU:OE1	1:B:393:ARG:HD3	2.11	0.51
1:B:164:HIS:CE1	2:B:701:GLC:H2	2.46	0.51
1:A:47:SER:HB3	1:A:50:GLU:HG3	1.92	0.51
1:B:328:ASN:ND2	1:B:328:ASN:N	2.59	0.50
1:B:524:GLN:HB3	4:B:1430:HOH:O	2.11	0.50
1:A:484:PHE:CE1	1:A:488:LYS:HD2	2.46	0.50
1:B:546:PRO:HG2	1:B:553:TRP:CH2	2.47	0.50
1:A:540:THR:HA	1:A:571:LEU:O	2.11	0.50
1:A:276:SER:OG	1:A:277:LYS:N	2.44	0.50
1:B:219:LEU:HB3	1:B:220:PRO:HD3	1.93	0.50
1:A:281:THR:H	1:A:289:GLN:NE2	2.06	0.50
1:B:243:ASN:ND2	1:B:295:LYS:HZ2	2.10	0.50
1:B:508:TRP:CZ2	1:B:546:PRO:HD3	2.47	0.50
1:B:542:LEU:CD2	1:B:570:LYS:HG2	2.42	0.49
1:A:227:ASP:O	1:A:231:ARG:HG2	2.12	0.49
1:B:139:GLU:OE2	1:B:140:ARG:NH1	2.45	0.49
1:B:507:SER:HB3	4:B:1120:HOH:O	2.13	0.49
1:A:133:ILE:HD13	1:A:449:THR:HG22	1.93	0.49
1:B:210:LEU:HD22	1:B:210:LEU:N	2.28	0.49
1:A:253:ARG:HH21	1:A:253:ARG:HG2	1.78	0.49
1:A:65:ARG:HG2	1:A:65:ARG:HH21	1.78	0.49
1:A:393:ARG:HG3	1:A:393:ARG:HH21	1.76	0.49
1:A:512:LYS:HG3	1:A:513:GLN:N	2.27	0.49
1:A:178:ARG:NH2	4:A:1156:HOH:O	2.45	0.49
1:B:146:PRO:HA	1:B:149:ASP:OD1	2.12	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:215:GLN:HG3	4:A:1300:HOH:O	2.12	0.48
1:B:278:THR:OG1	1:B:279:SER:N	2.46	0.48
1:A:488:LYS:HB3	4:A:1385:HOH:O	2.13	0.48
1:B:551:LYS:HE3	4:B:1045:HOH:O	2.12	0.48
1:A:517:TYR:OH	4:A:1004:HOH:O	2.16	0.48
1:B:477:LYS:NZ	1:B:477:LYS:HB3	2.28	0.48
1:A:268:PHE:HB2	1:A:270:ILE:HD11	1.95	0.47
1:B:273:PHE:HA	1:B:274:PRO:C	2.34	0.47
1:A:565:LYS:CD	1:A:565:LYS:H	2.20	0.47
1:B:282:ASN:C	1:B:282:ASN:ND2	2.67	0.47
1:A:155:GLN:HB3	1:A:155:GLN:HE21	1.52	0.47
1:A:263:ARG:HD3	4:A:1119:HOH:O	2.14	0.47
1:B:245:ALA:O	1:B:294:PRO:HD2	2.15	0.47
1:B:199:PRO:HG3	1:B:248:GLN:NE2	2.30	0.47
1:B:565:LYS:O	1:B:565:LYS:HD2	2.15	0.46
1:A:87:THR:HA	1:A:92:GLU:O	2.16	0.46
1:A:324:LEU:HD13	1:A:335:TRP:CZ3	2.50	0.46
1:B:401:ALA:O	1:B:404:LEU:HB2	2.15	0.46
1:B:312:ARG:O	1:B:316:GLU:HG3	2.14	0.46
1:A:281:THR:OG1	1:A:289:GLN:NE2	2.49	0.46
1:A:64:GLU:HB2	4:A:1313:HOH:O	2.15	0.46
1:B:340:ARG:HB3	1:B:340:ARG:NH2	2.30	0.46
1:A:279:SER:O	1:A:289:GLN:NE2	2.49	0.46
1:A:150:PRO:HG3	1:A:215:GLN:CD	2.36	0.46
1:A:51:GLU:OE1	1:A:51:GLU:HA	2.16	0.46
1:A:565:LYS:N	1:A:565:LYS:CD	2.79	0.45
1:B:376:PHE:O	1:B:380:VAL:HG23	2.17	0.45
1:A:57:ALA:CB	1:A:71:ALA:HB2	2.40	0.45
1:A:286:PHE:HZ	2:A:704:GLC:H61	1.80	0.45
1:B:140:ARG:NH1	1:B:201:HIS:HB2	2.27	0.45
1:A:272:ASP:O	1:A:282:ASN:ND2	2.50	0.45
1:A:163:ARG:NH1	1:A:165:ASP:OD1	2.49	0.45
1:B:551:LYS:HD2	1:B:551:LYS:O	2.17	0.45
1:A:244:HIS:CD2	1:A:286:PHE:HB2	2.51	0.45
1:B:276:SER:O	1:B:277:LYS:HD2	2.17	0.45
1:A:564:GLY:HA3	1:A:569:LEU:HD13	1.99	0.45
1:B:164:HIS:ND1	4:B:1167:HOH:O	2.36	0.45
1:A:565:LYS:NZ	1:A:570:LYS:HE3	2.32	0.45
1:B:467:ASP:O	1:B:470:ARG:HG3	2.17	0.45
1:B:324:LEU:HD13	1:B:335:TRP:CZ3	2.52	0.44
1:B:250:PHE:CG	1:B:251:ALA:N	2.85	0.44
1:A:402:ARG:HD2	4:A:1261:HOH:O	2.18	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:458:ILE:CD1	1:B:460:MET:HG3	2.47	0.44
1:B:524:GLN:HB3	1:B:525:ASP:H	1.48	0.44
1:A:164:HIS:CE1	1:A:466:PRO:HD3	2.52	0.44
1:A:64:GLU:HG3	4:A:1332:HOH:O	2.17	0.44
1:A:416:LEU:HD23	1:A:416:LEU:H	1.81	0.44
1:B:542:LEU:HD11	1:B:565:LYS:NZ	2.32	0.44
1:B:388:GLU:HB2	4:B:1187:HOH:O	2.17	0.44
1:A:124:THR:OG1	1:A:129:LYS:CE	2.66	0.44
1:B:193:THR:HB	1:B:194:PRO:CD	2.47	0.43
1:A:564:GLY:CA	1:A:568:GLN:O	2.67	0.43
1:A:281:THR:N	1:A:289:GLN:HE22	2.10	0.43
1:B:249:PHE:CE2	1:B:251:ALA:HB3	2.53	0.43
1:A:533:ASN:O	1:A:576:TYR:HA	2.17	0.43
1:A:527:HIS:HD2	4:A:1315:HOH:O	2.00	0.43
1:A:545:VAL:HG23	1:A:567:GLY:C	2.39	0.43
1:B:542:LEU:HD11	1:B:565:LYS:HZ3	1.83	0.43
1:B:105:ARG:NH1	1:B:105:ARG:HG2	2.34	0.43
1:A:65:ARG:HB3	1:B:4:GLU:HG3	1.99	0.43
1:B:197:ALA:HB3	1:B:208:ASP:HB3	2.01	0.43
1:A:245:ALA:O	1:A:294:PRO:HD2	2.19	0.43
1:A:133:ILE:CD1	1:A:449:THR:CG2	2.97	0.43
1:A:518:ALA:HA	1:A:530:VAL:O	2.19	0.43
1:B:275:VAL:HG12	1:B:275:VAL:O	2.18	0.42
1:A:27:LEU:C	1:A:27:LEU:HD23	2.39	0.42
1:A:275:VAL:HG12	1:A:276:SER:N	2.34	0.42
1:B:416:LEU:CD2	1:B:416:LEU:H	2.32	0.42
1:A:270:ILE:HD12	1:A:270:ILE:N	2.34	0.42
1:A:569:LEU:HD22	1:A:569:LEU:N	2.35	0.42
1:B:193:THR:HB	1:B:194:PRO:HD2	2.02	0.42
1:A:9:GLU:HA	4:A:1205:HOH:O	2.18	0.42
1:B:190:LEU:HG	1:B:234:ILE:CG2	2.50	0.42
1:B:516:LEU:HD11	1:B:539:GLN:NE2	2.33	0.42
1:B:162:PRO:HG2	1:B:470:ARG:HA	2.00	0.42
1:B:546:PRO:CG	1:B:553:TRP:CH2	3.02	0.42
1:A:444:MET:CE	1:A:491:ILE:HG12	2.49	0.42
1:A:523:VAL:HG13	1:A:523:VAL:O	2.20	0.42
1:B:424:ARG:HG3	1:B:461:ALA:C	2.39	0.42
1:B:286:PHE:HZ	2:B:704:GLC:H61	1.83	0.42
1:B:381:ILE:O	1:B:385:ALA:HB3	2.20	0.42
1:A:77:THR:O	1:A:78:LYS:HB2	2.20	0.42
1:A:179:LEU:N	1:A:180:PRO:CD	2.83	0.41
1:A:133:ILE:N	1:A:133:ILE:CD1	2.80	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:457:GLU:HA	1:A:487:TYR:CE1	2.55	0.41
1:B:577:GLN:HG2	1:B:578:GLY:N	2.36	0.41
1:A:281:THR:HG22	4:A:1419:HOH:O	2.20	0.41
1:B:36:ARG:HB3	1:B:87:THR:HB	2.03	0.41
1:B:264:TYR:O	1:B:267:TRP:HB2	2.20	0.41
1:B:465:ASP:HA	1:B:466:PRO:HA	1.88	0.41
1:B:144:GLY:C	1:B:173:LYS:HD2	2.41	0.41
1:B:546:PRO:HG2	1:B:553:TRP:CZ2	2.55	0.41
1:B:191:TYR:C	1:B:191:TYR:CD1	2.93	0.41
1:A:217:GLY:HA2	4:A:1015:HOH:O	2.19	0.41
1:B:279:SER:O	1:B:281:THR:N	2.54	0.41
1:A:68:TYR:CD2	1:A:403:MET:HG3	2.55	0.41
1:B:523:VAL:HG13	1:B:523:VAL:O	2.20	0.41
1:A:250:PHE:CG	1:A:251:ALA:N	2.88	0.41
1:B:382:ARG:HG3	1:B:388:GLU:HB3	2.02	0.41
1:B:223:ARG:CZ	1:B:317:GLN:OE1	2.68	0.41
1:B:541:VAL:HG22	1:B:542:LEU:N	2.36	0.41
1:A:345:LEU:HA	1:A:345:LEU:HD12	1.96	0.41
1:A:198:SER:OG	1:A:203:LYS:HA	2.21	0.41
1:B:574:ARG:HG3	1:B:577:GLN:OE1	2.21	0.41
1:A:16:TYR:C	1:A:23:LEU:HD12	2.40	0.41
1:B:179:LEU:N	1:B:180:PRO:CD	2.84	0.41
1:B:503:GLY:HA2	1:B:523:VAL:HB	2.03	0.41
1:A:346:ASN:HD22	1:A:346:ASN:C	2.23	0.41
1:A:573:LEU:HD11	1:A:579:MET:HG3	2.03	0.40
1:B:382:ARG:HB3	1:B:389:ILE:HG12	2.03	0.40
1:B:246:GLY:HA2	1:B:292:ALA:O	2.22	0.40
1:A:223:ARG:HA	1:A:223:ARG:HH11	1.85	0.40
1:A:477:LYS:CA	1:A:477:LYS:HE3	2.47	0.40
1:B:146:PRO:HA	1:B:149:ASP:CG	2.41	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	583/585 (100%)	555 (95%)	26 (4%)	2 (0%)	50	49
1	B	583/585 (100%)	551 (94%)	31 (5%)	1 (0%)	56	57
All	All	1166/1170 (100%)	1106 (95%)	57 (5%)	3 (0%)	50	49

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	277	LYS
1	A	278	THR
1	B	278	THR

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	493/493 (100%)	472 (96%)	21 (4%)	40	37
1	B	493/493 (100%)	473 (96%)	20 (4%)	41	40
All	All	986/986 (100%)	945 (96%)	41 (4%)	40	38

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90	GLN
1	A	122	PHE
1	A	129	LYS
1	A	155	GLN
1	A	158	LYS
1	A	182	LEU
1	A	191	TYR
1	A	260	GLU
1	A	280	ARG
1	A	341	LEU
1	A	345	LEU
1	A	346	ASN
1	A	398	LEU
1	A	404	LEU

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Mol	Chain	Res	Type
1	A	408	GLN
1	A	413	LEU
1	A	477	LYS
1	A	483	LEU
1	A	504	ASN
1	A	547	GLU
1	A	565	LYS
1	B	85	LEU
1	B	122	PHE
1	B	155	GLN
1	B	191	TYR
1	B	257	GLN
1	B	282	ASN
1	B	328	ASN
1	B	340	ARG
1	B	373	ASN
1	B	382	ARG
1	B	398	LEU
1	B	404	LEU
1	B	438	LEU
1	B	504	ASN
1	B	524	GLN
1	B	542	LEU
1	B	551	LYS
1	B	565	LYS
1	B	573	LEU
1	B	585	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (41) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	HIS
1	A	90	GLN
1	A	135	GLN
1	A	155	GLN
1	A	244	HIS
1	A	248	GLN
1	A	289	GLN
1	A	317	GLN
1	A	325	ASN
1	A	332	HIS
1	A	346	ASN

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Mol	Chain	Res	Type
1	A	367	GLN
1	A	390	HIS
1	A	408	GLN
1	A	443	GLN
1	A	504	ASN
1	A	509	HIS
1	A	515	ASN
1	A	527	HIS
1	A	539	GLN
1	A	563	HIS
1	A	568	GLN
1	B	135	GLN
1	B	164	HIS
1	B	202	HIS
1	B	243	ASN
1	B	244	HIS
1	B	248	GLN
1	B	282	ASN
1	B	328	ASN
1	B	332	HIS
1	B	357	HIS
1	B	367	GLN
1	B	373	ASN
1	B	504	ASN
1	B	509	HIS
1	B	533	ASN
1	B	539	GLN
1	B	544	GLN
1	B	563	HIS
1	B	566	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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

12 carbohydrates 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	GLC	A	701	2	10,11,12	0.39	0	11,15,17	0.41	0
2	GLC	A	702	2	10,11,12	0.57	0	11,15,17	0.19	0
2	GLC	A	703	2	10,11,12	0.55	0	11,15,17	0.80	0
2	GLC	A	704	2	10,11,12	0.51	0	11,15,17	0.42	0
2	GLC	A	705	2	10,11,12	0.55	0	11,15,17	0.31	0
2	GLC	A	706	2	12,12,12	0.41	0	17,17,17	0.73	0
2	GLC	B	701	2	10,11,12	0.41	0	11,15,17	0.36	0
2	GLC	B	702	2	10,11,12	0.49	0	11,15,17	0.24	0
2	GLC	B	703	2	10,11,12	0.49	0	11,15,17	0.72	0
2	GLC	B	704	2	10,11,12	0.54	0	11,15,17	0.38	0
2	GLC	B	705	2	10,11,12	0.47	0	11,15,17	0.28	0
2	GLC	B	706	2	12,12,12	0.40	0	17,17,17	0.45	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	GLC	A	701	2	-	0/2/19/22	0/1/1/1
2	GLC	A	702	2	-	0/2/19/22	0/1/1/1
2	GLC	A	703	2	-	0/2/19/22	0/1/1/1
2	GLC	A	704	2	-	0/2/19/22	0/1/1/1
2	GLC	A	705	2	-	0/2/19/22	0/1/1/1
2	GLC	A	706	2	1/1/5/5	0/2/22/22	0/1/1/1
2	GLC	B	701	2	-	0/2/19/22	0/1/1/1
2	GLC	B	702	2	-	0/2/19/22	0/1/1/1
2	GLC	B	703	2	-	0/2/19/22	0/1/1/1
2	GLC	B	704	2	-	0/2/19/22	0/1/1/1
2	GLC	B	705	2	-	0/2/19/22	0/1/1/1
2	GLC	B	706	2	1/1/5/5	0/2/22/22	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	706	GLC	C1
2	B	706	GLC	C1

There are no torsion outliers.

There are no ring outliers.

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.

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	585/585 (100%)	-0.04	19 (3%)	45 50	15, 25, 47, 72	0
1	B	585/585 (100%)	0.05	29 (4%)	28 30	17, 27, 50, 83	0
All	All	1170/1170 (100%)	0.01	48 (4%)	35 40	15, 26, 48, 83	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	549	GLY	5.6
1	B	277	LYS	4.8
1	A	277	LYS	4.7
1	B	279	SER	4.7
1	B	551	LYS	4.7
1	B	548	SER	4.3
1	B	565	LYS	4.2
1	A	280	ARG	4.0
1	B	566	GLN	4.0
1	B	275	VAL	3.9
1	A	551	LYS	3.8
1	A	548	SER	3.8
1	A	547	GLU	3.5
1	B	552	THR	3.3
1	B	570	LYS	3.3
1	B	547	GLU	3.1
1	B	49	GLU	3.1
1	A	276	SER	3.0
1	B	278	THR	3.0
1	A	546	PRO	3.0
1	B	273	PHE	3.0
1	A	278	THR	2.9
1	A	275	VAL	2.8
1	B	280	ARG	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	263	ARG	2.8
1	A	279	SER	2.7
1	B	152	GLY	2.6
1	A	552	THR	2.5
1	B	90	GLN	2.5
1	B	153	THR	2.4
1	B	51	GLU	2.4
1	B	545	VAL	2.4
1	A	567	GLY	2.3
1	B	270	ILE	2.3
1	B	542	LEU	2.3
1	A	545	VAL	2.2
1	B	259	GLY	2.2
1	A	553	TRP	2.2
1	A	49	GLU	2.2
1	A	564	GLY	2.2
1	A	566	GLN	2.2
1	B	546	PRO	2.1
1	B	477	LYS	2.1
1	B	159	ASP	2.1
1	A	261	GLN	2.1
1	B	224	ARG	2.0
1	B	567	GLY	2.0
1	B	550	GLY	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 ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	GLC	B	705	11/12	0.20	7.62	33,38,40,43	11
2	GLC	A	701	11/12	0.42	5.32	42,47,50,51	11

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	GLC	B	701	11/12	0.38	4.22	43,47,49,51	11
2	GLC	A	705	11/12	0.20	4.05	23,29,34,37	11
2	GLC	B	706	12/12	0.26	3.75	51,64,66,67	0
2	GLC	A	706	12/12	0.24	3.53	40,48,52,55	11
2	GLC	A	704	11/12	0.10	0.04	20,24,27,33	0
2	GLC	B	704	11/12	0.10	-0.44	24,28,31,33	0
2	GLC	B	703	11/12	0.09	-0.79	23,26,27,28	0
2	GLC	A	702	11/12	0.08	-0.80	21,23,26,33	0
2	GLC	A	703	11/12	0.09	-0.89	17,21,22,23	0
2	GLC	B	702	11/12	0.08	-1.15	27,29,31,37	0

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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	CA	A	1001	1/1	0.09	-0.21	24,24,24,24	0
3	CA	B	1002	1/1	0.05	-2.28	32,32,32,32	0

6.5 Other polymers

There are no such residues in this entry.