



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 03:35 AM GMT

PDB ID : 1X0C
Title : Improved Crystal Structure of Isopullulanase from *Aspergillus niger* ATCC 9642
Authors : Mizuno, M.; Tonozone, T.; Yamamura, A.; Miyasaka, Y.; Akeboshi, H.; Kami-
tori, S.; Nishikawa, A.; Sakano, Y.
Deposited on : 2005-03-17
Resolution : 1.70 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>

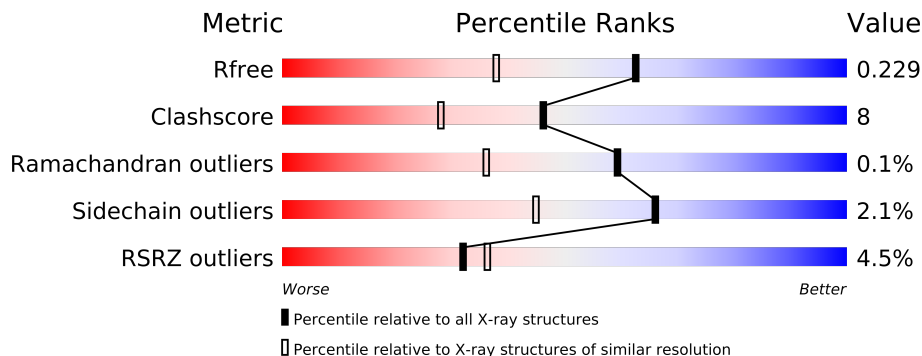
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.70 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	2456 (1.70-1.70)
Clashscore	79885	2929 (1.70-1.70)
Ramachandran outliers	78287	2878 (1.70-1.70)
Sidechain outliers	78261	2878 (1.70-1.70)
RSRZ outliers	66119	2456 (1.70-1.70)

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	549	
1	B	549	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	NAG	A	1000	-	X
2	NAG	A	1005	-	X
2	NAG	A	1007	-	X
2	NAG	A	1008	-	X
2	NAG	A	1009	-	X
2	NAG	B	1000	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
2	NAG	B	1002	-	X
2	NAG	B	1003	-	X
2	NAG	B	1006	-	X
2	NAG	B	1007	-	X
2	NAG	B	1008	-	X
2	NAG	B	1009	-	X
2	NAG	B	1010	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10079 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 Isopullulanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	549	Total	C	N	O	S	0	0	0
			4242	2676	710	845	11			
1	B	549	Total	C	N	O	S	0	0	0
			4242	2676	710	845	11			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	ARG	-	EXPRESSION TAG	UNP O00105
A	17	GLU	-	EXPRESSION TAG	UNP O00105
A	18	PHE	-	EXPRESSION TAG	UNP O00105
A	19	MET	-	EXPRESSION TAG	UNP O00105
B	16	ARG	-	EXPRESSION TAG	UNP O00105
B	17	GLU	-	EXPRESSION TAG	UNP O00105
B	18	PHE	-	EXPRESSION TAG	UNP O00105
B	19	MET	-	EXPRESSION TAG	UNP O00105

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is water.

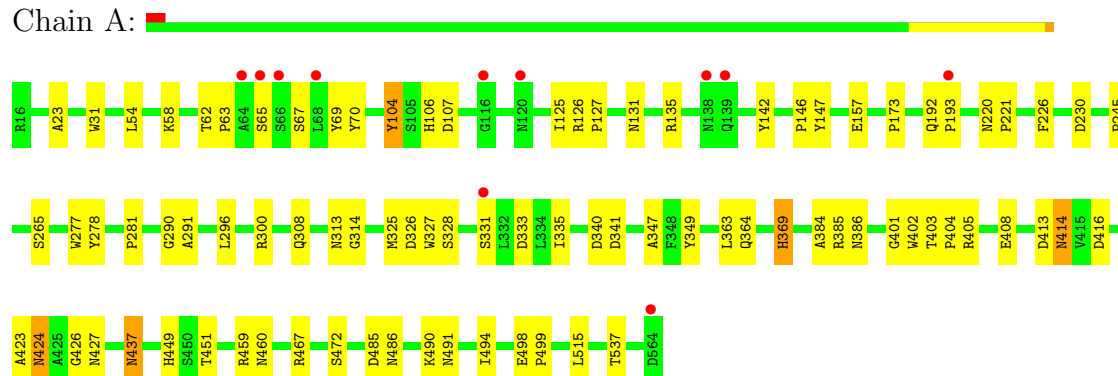
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	731	Total	O	0	0
			731	731		
3	B	542	Total	O	0	0
			542	542		

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 ($\text{RSRZ} > 2$). 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.

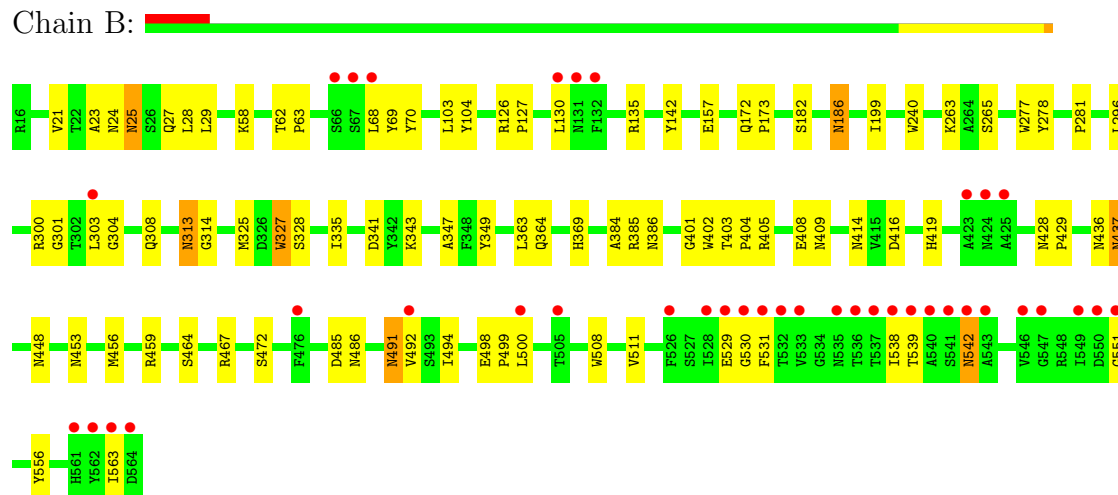
• Molecule 1: Isopullulanase

Chain A:



• Molecule 1: Isopullulanase

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	99.72Å 136.16Å 84.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.86 – 1.70 49.86 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.0 (49.86-1.70) 99.1 (49.86-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.13 (at 1.70Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.195 , 0.229 0.195 , 0.229	Depositor DCC
R_{free} test set	12531 reflections (11.08%)	DCC
Wilson B-factor (Å ²)	15.0	Xtriage
Anisotropy	0.486	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 125682 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10079	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.32% 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: NAG

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	0.32	0/4357	0.64	1/5962 (0.0%)
1	B	0.30	0/4357	0.61	0/5962
All	All	0.31	0/8714	0.63	1/11924 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	369	HIS	N-CA-C	-5.39	96.44	111.00

There are no chirality outliers.

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	4242	0	3979	58	0
1	B	4242	0	3978	78	0
2	A	154	0	143	5	0
2	B	168	0	156	9	0
3	A	731	0	0	8	0
3	B	542	0	0	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	10079	0	8256	140	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 8.

All (140) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:539:THR:H	1:B:542:ASN:HD21	1.16	0.91
1:A:54:LEU:HD11	2:A:1002:NAG:H61	1.53	0.91
1:A:414:ASN:ND2	1:A:467:ARG:HH22	1.72	0.87
1:B:539:THR:H	1:B:542:ASN:ND2	1.76	0.82
1:B:25:ASN:HD22	1:B:27:GLN:H	1.27	0.79
1:B:347:ALA:H	1:B:369:HIS:HD2	1.29	0.79
1:B:157:GLU:OE2	1:B:419:HIS:HE1	1.66	0.78
1:A:347:ALA:H	1:A:369:HIS:HD2	1.29	0.77
1:B:127:PRO:HG2	1:B:130:LEU:CD1	2.17	0.75
1:B:529:GLU:HB2	1:B:563:ILE:HB	1.73	0.70
1:B:127:PRO:HG2	1:B:130:LEU:HD13	1.74	0.69
1:B:25:ASN:ND2	1:B:28:LEU:H	1.91	0.68
1:B:303:LEU:HD23	1:B:304:GLY:O	1.93	0.68
1:B:308:GLN:NE2	1:B:327:TRP:HE1	1.93	0.65
1:A:414:ASN:HD21	1:A:467:ARG:HH22	1.42	0.65
1:A:192:GLN:HB3	1:A:193:PRO:HD2	1.79	0.64
1:A:135:ARG:HG2	1:A:142:TYR:HB2	1.78	0.63
1:A:386:ASN:HB3	3:A:1284:HOH:O	1.98	0.63
1:B:308:GLN:HE21	1:B:327:TRP:HE1	1.47	0.63
1:B:539:THR:N	1:B:542:ASN:HD21	1.94	0.60
1:B:25:ASN:ND2	1:B:27:GLN:H	1.97	0.60
1:B:157:GLU:HB3	1:B:173:PRO:HG3	1.83	0.60
1:B:436:ASN:ND2	1:B:437:ASN:H	1.99	0.60
1:B:453:ASN:HA	2:B:1008:NAG:H82	1.84	0.60
1:A:146:PRO:HG3	3:A:1498:HOH:O	2.01	0.59
1:B:186:ASN:HD22	1:B:186:ASN:C	2.05	0.59
1:B:472:SER:HB3	1:B:494:ILE:HG23	1.84	0.58
1:A:58:LYS:HB3	1:A:69:TYR:HB3	1.85	0.58
1:B:416:ASP:HB2	1:B:467:ARG:NH2	2.18	0.58
1:B:127:PRO:HG2	1:B:130:LEU:HD11	1.86	0.57
1:B:135:ARG:HG2	1:B:142:TYR:HB2	1.85	0.57
1:A:54:LEU:HD12	3:A:1364:HOH:O	2.04	0.57
1:B:436:ASN:HD22	1:B:437:ASN:H	1.53	0.57
2:A:1007:NAG:H81	3:A:1362:HOH:O	2.03	0.57
1:B:511:VAL:HA	1:B:556:TYR:CE1	2.40	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:1000:NAG:H81	3:A:1669:HOH:O	2.05	0.56
2:B:1003:NAG:H83	3:B:1392:HOH:O	2.05	0.56
1:A:127:PRO:HB3	1:A:416:ASP:OD2	2.06	0.56
1:B:542:ASN:HD22	1:B:542:ASN:C	2.09	0.56
1:A:423:ALA:H	1:A:427:ASN:ND2	2.03	0.56
1:B:23:ALA:HB2	1:B:70:TYR:CG	2.41	0.56
1:A:404:PRO:HG2	3:A:1398:HOH:O	2.06	0.56
1:A:537:THR:HG23	3:A:1320:HOH:O	2.06	0.56
1:A:424:ASN:H	1:A:427:ASN:ND2	2.04	0.55
1:A:125:ILE:HD12	1:A:125:ILE:N	2.21	0.55
1:B:498:GLU:HB3	1:B:499:PRO:HD2	1.88	0.54
1:B:281:PRO:HG3	1:B:296:LEU:HD13	1.89	0.54
1:B:491:ASN:HD22	2:B:1010:NAG:H82	1.71	0.54
1:A:347:ALA:H	1:A:369:HIS:CD2	2.18	0.54
1:A:331:SER:HB3	1:A:333:ASP:OD1	2.07	0.54
1:B:404:PRO:HG2	3:B:1207:HOH:O	2.07	0.54
1:A:449:HIS:HD2	3:A:1692:HOH:O	1.91	0.53
1:B:404:PRO:HG2	1:B:456:MET:HE1	1.91	0.53
1:A:157:GLU:HB3	1:A:173:PRO:HG3	1.90	0.53
1:A:23:ALA:HB2	1:A:70:TYR:CG	2.44	0.53
1:B:172:GLN:HG2	1:B:419:HIS:CD2	2.45	0.52
1:B:401:GLY:O	1:B:402:TRP:HB2	2.09	0.52
1:A:460:ASN:ND2	2:A:1007:NAG:C7	2.72	0.52
1:A:308:GLN:O	1:A:335:ILE:HD12	2.09	0.52
1:A:472:SER:HB3	1:A:494:ILE:HG23	1.91	0.51
1:B:419:HIS:HD2	3:B:1223:HOH:O	1.93	0.51
2:B:1002:NAG:H81	3:B:1395:HOH:O	2.11	0.51
1:B:21:VAL:HG11	1:B:68:LEU:HD21	1.93	0.50
1:A:423:ALA:H	1:A:427:ASN:HD22	1.58	0.50
1:A:62:THR:HB	1:A:63:PRO:HA	1.94	0.49
1:B:459:ARG:HA	1:B:485:ASP:O	2.13	0.49
1:B:492:VAL:HG13	1:B:531:PHE:CD1	2.48	0.49
1:B:21:VAL:O	1:B:21:VAL:HG13	2.12	0.49
1:B:62:THR:HB	1:B:63:PRO:HA	1.95	0.48
1:B:29:LEU:HD23	1:B:103:LEU:HD12	1.96	0.48
1:B:103:LEU:HA	1:B:182:SER:O	2.14	0.48
1:B:437:ASN:HD22	1:B:437:ASN:C	2.16	0.48
1:B:347:ALA:N	1:B:369:HIS:HD2	2.03	0.47
1:A:106:HIS:HD2	1:A:107:ASP:O	1.97	0.47
1:B:25:ASN:HD22	1:B:27:GLN:N	2.06	0.47
1:B:126:ARG:HA	1:B:127:PRO:C	2.34	0.47
1:A:340:ASP:OD1	1:A:364:GLN:HG3	2.14	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:58:LYS:HB3	1:B:69:TYR:HB3	1.96	0.47
1:B:25:ASN:ND2	1:B:28:LEU:N	2.63	0.47
1:A:490:LYS:O	1:A:491:ASN:HB2	2.15	0.47
1:A:126:ARG:HA	1:A:127:PRO:C	2.35	0.47
1:B:464:SER:HB2	2:B:1010:NAG:H81	1.95	0.47
1:A:408:GLU:HA	1:A:459:ARG:O	2.15	0.46
1:B:24:ASN:HD22	2:B:1000:NAG:C8	2.29	0.46
1:B:277:TRP:CD1	1:B:349:TYR:HB3	2.50	0.46
1:B:492:VAL:O	1:B:492:VAL:HG13	2.16	0.46
1:B:542:ASN:ND2	1:B:542:ASN:C	2.70	0.45
1:A:424:ASN:HD22	1:A:426:GLY:N	2.13	0.45
1:B:265:SER:HA	1:B:313:ASN:O	2.16	0.45
1:A:265:SER:HA	1:A:313:ASN:O	2.17	0.45
1:A:363:LEU:O	1:A:384:ALA:HA	2.16	0.45
1:B:467:ARG:HH11	1:B:467:ARG:HG3	1.82	0.45
1:B:508:TRP:HB3	1:B:551:GLY:HA3	1.99	0.44
1:A:401:GLY:O	1:A:402:TRP:HB2	2.17	0.44
1:B:301:GLY:HA3	1:B:327:TRP:CE2	2.53	0.44
1:A:281:PRO:HG3	1:A:296:LEU:HD13	1.99	0.44
1:B:186:ASN:ND2	1:B:186:ASN:C	2.70	0.44
1:A:300:ARG:NH1	1:A:328:SER:OG	2.49	0.44
1:B:448:ASN:ND2	2:B:1007:NAG:H82	2.33	0.44
1:A:104:TYR:CE2	1:A:147:TYR:HA	2.53	0.44
1:A:300:ARG:NE	1:A:328:SER:OG	2.50	0.44
1:A:65:SER:C	1:A:67:SER:H	2.19	0.44
1:B:303:LEU:HD13	1:B:335:ILE:HG13	1.98	0.43
1:A:424:ASN:HD22	1:A:426:GLY:H	1.65	0.43
1:B:363:LEU:O	1:B:384:ALA:HA	2.18	0.43
1:B:24:ASN:HD22	2:B:1000:NAG:H83	1.84	0.43
1:B:408:GLU:HG3	1:B:409:ASN:ND2	2.32	0.43
1:A:437:ASN:C	1:A:437:ASN:HD22	2.21	0.43
1:B:403:THR:O	1:B:405:ARG:HG3	2.18	0.43
1:B:491:ASN:ND2	2:B:1010:NAG:H82	2.33	0.43
1:B:343:LYS:HD2	1:B:343:LYS:N	2.34	0.43
1:A:498:GLU:HB3	1:A:499:PRO:HD2	2.01	0.43
1:A:300:ARG:HD2	1:A:326:ASP:O	2.19	0.43
1:A:414:ASN:HD21	1:A:467:ARG:NH2	2.14	0.42
1:A:451:THR:HA	1:A:515:LEU:HB2	2.00	0.42
1:A:403:THR:O	1:A:405:ARG:HG3	2.18	0.42
1:A:414:ASN:HD22	1:A:467:ARG:HH22	1.58	0.42
1:A:459:ARG:HA	1:A:485:ASP:O	2.20	0.42
1:A:226:PHE:O	1:A:230:ASP:HB3	2.20	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:31:TRP:CZ3	1:A:245:PRO:HB3	2.54	0.42
1:A:314:GLY:HA2	1:A:341:ASP:O	2.19	0.42
1:A:467:ARG:HG3	1:A:467:ARG:HH11	1.83	0.42
1:A:220:ASN:HB3	1:A:221:PRO:HD2	2.02	0.42
1:B:281:PRO:CG	1:B:296:LEU:HD13	2.50	0.41
1:B:314:GLY:HA2	1:B:341:ASP:O	2.20	0.41
1:B:240:TRP:CE2	1:B:263:LYS:HE3	2.54	0.41
1:B:364:GLN:HA	1:B:385:ARG:O	2.20	0.41
1:B:428:ASN:N	1:B:429:PRO:CD	2.84	0.41
1:B:500:LEU:HD21	3:B:1303:HOH:O	2.21	0.41
1:A:385:ARG:HA	1:A:413:ASP:O	2.21	0.40
1:B:511:VAL:HA	1:B:556:TYR:CZ	2.56	0.40
1:B:300:ARG:NH2	1:B:328:SER:HB2	2.36	0.40
1:B:199:ILE:HD12	3:B:1287:HOH:O	2.20	0.40
1:B:491:ASN:HA	1:B:530:GLY:O	2.21	0.40
1:B:58:LYS:HA	1:B:70:TYR:O	2.22	0.40
1:A:277:TRP:CD1	1:A:349:TYR:HB3	2.56	0.40
1:B:386:ASN:HB3	3:B:1192:HOH:O	2.21	0.40
1:B:538:ILE:HA	1:B:542:ASN:HD21	1.86	0.40
1:A:221:PRO:HG2	2:A:1000:NAG:H83	2.04	0.40
1:A:290:GLY:O	1:A:291:ALA:C	2.60	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	547/549 (100%)	517 (94%)	30 (6%)	0	100	100
1	B	547/549 (100%)	514 (94%)	32 (6%)	1 (0%)	56	33
All	All	1094/1098 (100%)	1031 (94%)	62 (6%)	1 (0%)	59	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	491	ASN

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	466/466 (100%)	457 (98%)	9 (2%)	69	50
1	B	466/466 (100%)	455 (98%)	11 (2%)	61	39
All	All	932/932 (100%)	912 (98%)	20 (2%)	66	45

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	104	TYR
1	A	131	ASN
1	A	278	TYR
1	A	325	MET
1	A	327	TRP
1	A	414	ASN
1	A	424	ASN
1	A	437	ASN
1	A	486	ASN
1	B	25	ASN
1	B	104	TYR
1	B	186	ASN
1	B	278	TYR
1	B	313	ASN
1	B	325	MET
1	B	327	TRP
1	B	414	ASN
1	B	437	ASN
1	B	486	ASN
1	B	542	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	41	GLN
1	A	51	GLN
1	A	106	HIS
1	A	120	ASN
1	A	122	ASN
1	A	131	ASN
1	A	139	GLN
1	A	369	HIS
1	A	386	ASN
1	A	414	ASN
1	A	424	ASN
1	A	427	ASN
1	A	437	ASN
1	A	449	HIS
1	A	486	ASN
1	A	516	ASN
1	B	25	ASN
1	B	41	GLN
1	B	186	ASN
1	B	192	GLN
1	B	286	GLN
1	B	308	GLN
1	B	330	ASN
1	B	369	HIS
1	B	386	ASN
1	B	414	ASN
1	B	419	HIS
1	B	436	ASN
1	B	437	ASN
1	B	486	ASN
1	B	542	ASN

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

There are no carbohydrates in this entry.

5.6 Ligand geometry

23 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	NAG	A	1000	1	12,14,15	0.46	0	15,19,21	0.73	1 (6%)
2	NAG	A	1001	1	12,14,15	0.41	0	15,19,21	0.67	0
2	NAG	A	1002	1	12,14,15	0.45	0	15,19,21	0.68	0
2	NAG	A	1003	1	12,14,15	0.43	0	15,19,21	0.77	0
2	NAG	A	1004	1	12,14,15	0.41	0	15,19,21	0.69	0
2	NAG	A	1005	1	12,14,15	0.43	0	15,19,21	0.69	0
2	NAG	A	1006	1	12,14,15	0.39	0	15,19,21	0.75	0
2	NAG	A	1007	1	12,14,15	0.39	0	15,19,21	0.76	1 (6%)
2	NAG	A	1008	1	12,14,15	0.42	0	15,19,21	0.67	0
2	NAG	A	1009	1	12,14,15	0.42	0	15,19,21	0.72	0
2	NAG	A	1010	1	12,14,15	0.41	0	15,19,21	0.71	0
2	NAG	B	1000	1	12,14,15	0.43	0	15,19,21	0.68	0
2	NAG	B	1001	1	12,14,15	0.44	0	15,19,21	0.67	0
2	NAG	B	1002	1	12,14,15	0.41	0	15,19,21	0.76	0
2	NAG	B	1003	1	12,14,15	0.43	0	15,19,21	0.68	0
2	NAG	B	1004	1	12,14,15	0.45	0	15,19,21	0.72	1 (6%)
2	NAG	B	1005	1	12,14,15	0.43	0	15,19,21	0.70	0
2	NAG	B	1006	1	12,14,15	0.45	0	15,19,21	0.68	0
2	NAG	B	1007	1	12,14,15	0.41	0	15,19,21	0.75	0
2	NAG	B	1008	1	12,14,15	0.40	0	15,19,21	0.73	0
2	NAG	B	1009	1	12,14,15	0.41	0	15,19,21	0.71	0
2	NAG	B	1010	1	12,14,15	0.42	0	15,19,21	0.71	0
2	NAG	B	1011	1	12,14,15	0.44	0	15,19,21	0.68	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	NAG	A	1000	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1001	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1002	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1003	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1004	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1005	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1006	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1007	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1008	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1009	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1010	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1000	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1001	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1002	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1003	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1004	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1005	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1006	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1007	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1008	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1009	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1010	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1011	1	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1007	NAG	C3-C2-N2	-2.12	108.53	111.76
2	B	1004	NAG	C2-N2-C7	-2.10	119.56	123.09
2	A	1000	NAG	C2-N2-C7	-2.05	119.65	123.09

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1011	NAG	O7-C7-N2-C2

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	549/549 (100%)	-0.20	11 (2%) 62 67	7, 13, 27, 48	0
1	B	549/549 (100%)	0.38	39 (7%) 16 19	9, 20, 34, 62	0
All	All	1098/1098 (100%)	0.09	50 (4%) 32 35	7, 16, 32, 62	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	564	ASP	7.4
1	B	538	ILE	6.4
1	B	563	ILE	5.5
1	B	500	LEU	5.5
1	B	535	ASN	4.6
1	B	530	GLY	4.5
1	B	66	SER	4.3
1	B	540	ALA	4.1
1	B	68	LEU	4.1
1	B	528	ILE	4.1
1	B	546	VAL	4.0
1	B	549	ILE	3.6
1	B	539	THR	3.6
1	B	531	PHE	3.4
1	A	64	ALA	3.3
1	B	543	ALA	3.3
1	A	65	SER	3.3
1	B	541	SER	3.3
1	B	492	VAL	3.2
1	B	533	VAL	3.2
1	B	424	ASN	3.2
1	B	537	THR	3.2
1	B	425	ALA	3.1
1	A	116	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	66	SER	2.9
1	B	561	HIS	2.9
1	B	505	THR	2.8
1	A	564	ASP	2.7
1	B	562	TYR	2.7
1	B	536	THR	2.7
1	A	138	ASN	2.7
1	B	423	ALA	2.7
1	B	67	SER	2.6
1	B	532	THR	2.5
1	A	139	GLN	2.5
1	B	303	LEU	2.5
1	A	68	LEU	2.4
1	B	130	LEU	2.4
1	B	550	ASP	2.4
1	A	120	ASN	2.4
1	B	131	ASN	2.4
1	B	526	PHE	2.3
1	A	193	PRO	2.2
1	A	331	SER	2.2
1	B	476	PHE	2.2
1	B	529	GLU	2.1
1	B	547	GLY	2.1
1	B	132	PHE	2.1
1	B	542	ASN	2.0
1	B	551	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 ⓘ

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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	NAG	A	1007	14/15	0.15	9.20	30,36,41,43	0
2	NAG	A	1000	14/15	0.19	8.20	32,34,39,42	0
2	NAG	B	1003	14/15	0.23	7.61	31,33,40,40	0
2	NAG	B	1000	14/15	0.15	7.37	29,33,37,39	0
2	NAG	A	1009	14/15	0.21	7.36	21,36,40,40	0
2	NAG	B	1008	14/15	0.17	6.93	38,41,43,45	0
2	NAG	B	1002	14/15	0.25	4.10	31,40,47,47	0
2	NAG	A	1008	14/15	0.10	3.39	17,22,28,39	0
2	NAG	B	1006	14/15	0.21	3.20	33,38,43,47	0
2	NAG	B	1010	14/15	0.23	3.09	42,46,48,48	0
2	NAG	B	1009	14/15	0.15	3.06	37,41,47,49	0
2	NAG	A	1005	14/15	0.12	2.91	22,28,33,37	0
2	NAG	B	1007	14/15	0.12	2.64	33,35,42,43	0
2	NAG	A	1010	14/15	0.15	1.34	24,27,31,34	0
2	NAG	B	1005	14/15	0.14	1.31	27,34,44,45	0
2	NAG	A	1002	14/15	0.17	1.22	27,29,32,36	0
2	NAG	B	1011	14/15	0.18	0.93	34,38,49,51	0
2	NAG	B	1004	14/15	0.20	0.71	32,35,40,41	0
2	NAG	B	1001	14/15	0.10	0.63	22,26,31,33	0
2	NAG	A	1004	14/15	0.14	0.62	24,29,34,34	0
2	NAG	A	1006	14/15	0.10	0.58	22,29,33,35	0
2	NAG	A	1001	14/15	0.08	0.09	13,15,17,17	0
2	NAG	A	1003	14/15	0.09	-0.12	16,18,21,22	0

6.5 Other polymers ⓘ

There are no such residues in this entry.