



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 05:33 PM UTC

PDB ID : 4MPS / pdb_00004mps
Title : Crystal structure of rat Beta-galactoside alpha-2,6-sialyltransferase 1 (ST6GAL1), Northeast Structural Genomics Consortium Target RnR367A
Authors : Forouhar, F.; Meng, L.; Milaninia, S.; Seetharaman, J.; Su, M.; Kornhaber, G.; Montelione, G.T.; Hunt, J.F.; Moremen, K.W.; Tong, L.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2013-09-13
Resolution : 2.40 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) : 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

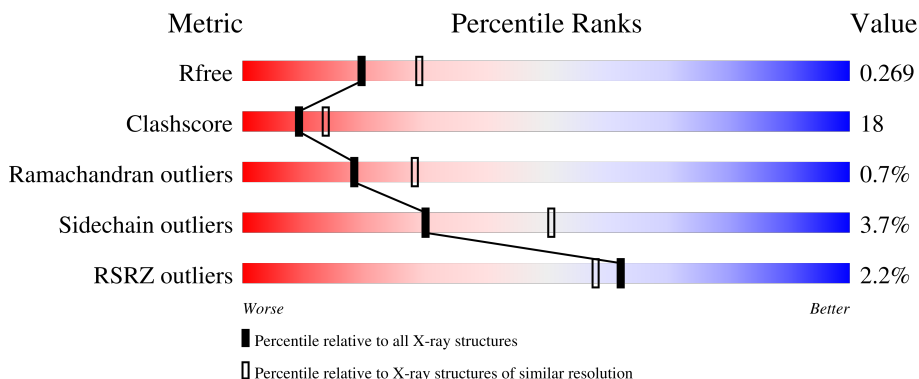
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	310	
1	B	310	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4910 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.

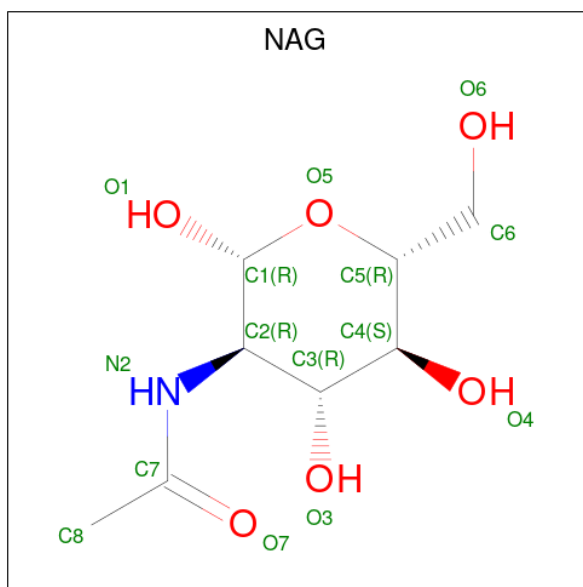
- Molecule 1 is a protein called Beta-galactoside alpha-2,6-sialyltransferase 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	293	Total 2407	C 1549	N 408	O 436	S 5	Se 9	0	0	0
1	B	273	Total 2238	C 1440	N 376	O 410	S 5	Se 7	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	94	MSE	-	expression tag	UNP G3V680
B	94	MSE	-	expression tag	UNP G3V680

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
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		

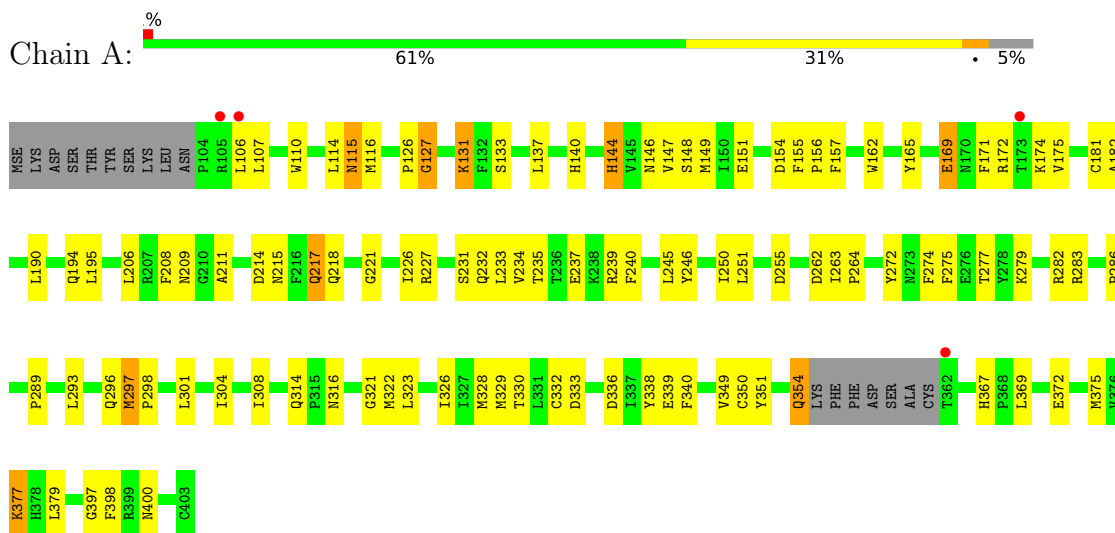
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	108	Total	O	0	0
			108	108		
3	B	101	Total	O	0	0
			101	101		

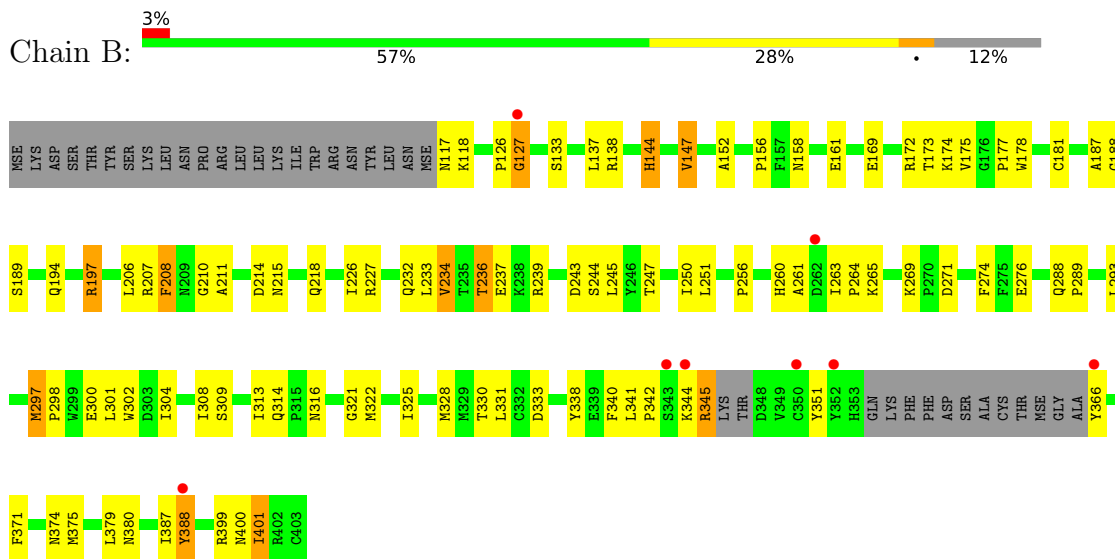
3 Residue-property plots i

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: Beta-galactoside alpha-2,6-sialyltransferase 1



- Molecule 1: Beta-galactoside alpha-2,6-sialyltransferase 1



4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	134.72Å 49.75Å 86.06Å 90.00° 92.30° 90.00°	Depositor
Resolution (Å)	46.67 – 2.40 46.67 – 2.40	Depositor EDS
% Data completeness (in resolution range)	93.4 (46.67-2.40) 93.1 (46.67-2.40)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.77 (at 2.39Å)	Xtrriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.200 , 0.259 0.209 , 0.269	Depositor DCC
R_{free} test set	4236 reflections (9.66%)	wwPDB-VP
Wilson B-factor (Å ²)	21.9	Xtrriage
Anisotropy	0.166	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 46.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.028 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4910	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.98% 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: 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	0.42	0/2466	0.94	10/3328 (0.3%)
1	B	0.42	0/2294	0.89	8/3099 (0.3%)
All	All	0.42	0/4760	0.92	18/6427 (0.3%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	340	PHE	N-CA-C	-10.33	97.16	112.04
1	B	340	PHE	N-CA-C	-7.78	100.83	112.04
1	A	133	SER	N-CA-C	-6.81	102.03	110.41
1	B	236	THR	N-CA-C	6.61	121.55	112.90
1	B	300	GLU	N-CA-C	-6.05	104.76	111.36
1	B	133	SER	N-CA-C	-5.86	102.64	110.55
1	A	400	ASN	N-CA-C	5.70	119.94	112.92
1	A	144	HIS	N-CA-C	5.56	120.06	112.88
1	A	372	GLU	N-CA-C	-5.48	105.31	111.28
1	B	172	ARG	N-CA-C	5.45	117.98	111.71
1	B	338	TYR	N-CA-C	5.36	117.25	108.52
1	A	275	PHE	N-CA-C	5.23	117.38	111.11
1	B	208	PHE	N-CA-C	5.20	117.88	109.40
1	A	333	ASP	N-CA-C	-5.17	105.34	111.69
1	B	144	HIS	N-CA-C	5.11	120.15	113.30
1	A	296	GLN	N-CA-C	5.07	118.61	112.23
1	A	107	LEU	N-CA-C	-5.04	106.23	112.38
1	A	217	GLN	N-CA-C	5.02	117.13	111.11

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2407	0	2360	84	0
1	B	2238	0	2175	84	0
2	A	28	0	26	0	0
2	B	28	0	26	0	0
3	A	108	0	0	5	0
3	B	101	0	0	8	0
All	All	4910	0	4587	168	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 18.

All (168) 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:B:197:ARG:HB2	1:B:197:ARG:HH11	1.21	1.05
1:A:297:MSE:HE3	1:A:298:PRO:HA	1.33	1.03
1:B:147:VAL:HG11	1:B:330:THR:HG21	1.49	0.95
1:A:354:GLN:HE21	1:A:354:GLN:HA	1.29	0.95
1:B:344:LYS:HD3	1:B:388:TYR:HE1	1.29	0.95
1:A:297:MSE:HE2	1:A:323:LEU:HG	1.50	0.94
1:A:297:MSE:HE3	1:A:298:PRO:CA	2.02	0.88
1:A:171:PHE:O	1:A:175:VAL:HG12	1.78	0.83
1:B:344:LYS:HD3	1:B:388:TYR:CE1	2.17	0.80
1:B:197:ARG:HB2	1:B:197:ARG:NH1	1.95	0.80
1:B:308:ILE:HG12	1:B:375:MSE:HE3	1.66	0.78
1:A:137:LEU:HD11	1:A:375:MSE:HE2	1.65	0.77
1:A:147:VAL:HG11	1:A:330:THR:HG21	1.67	0.77
1:A:147:VAL:HG13	1:A:172:ARG:NH1	2.00	0.76
1:A:137:LEU:HD23	1:A:379:LEU:HD21	1.67	0.76
1:A:304:ILE:HG22	1:A:375:MSE:HE1	1.65	0.76
1:B:345:ARG:HH21	1:B:345:ARG:HB3	1.50	0.75
3:A:668:HOH:O	1:B:177:PRO:HD3	1.88	0.73
1:A:211:ALA:O	1:A:227:ARG:NH2	2.16	0.73
1:B:276:GLU:HG3	3:B:696:HOH:O	1.86	0.73
1:A:354:GLN:HE21	1:A:354:GLN:CA	2.01	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:354:GLN:HA	1:A:354:GLN:NE2	2.05	0.71
1:B:175:VAL:HG22	1:B:250:ILE:HD13	1.73	0.71
1:B:147:VAL:CG1	1:B:330:THR:HG21	2.20	0.70
1:B:297:MSE:HG3	1:B:298:PRO:HD3	1.74	0.69
1:B:261:ALA:HB1	1:B:265:LYS:HG3	1.75	0.69
1:B:269:LYS:O	1:B:269:LYS:HD3	1.92	0.69
1:B:211:ALA:O	1:B:227:ARG:NH2	2.22	0.69
1:B:227:ARG:HB3	1:B:251:LEU:HD23	1.75	0.68
1:B:137:LEU:HD23	1:B:379:LEU:HD11	1.74	0.67
1:A:308:ILE:CD1	1:A:375:MSE:HE3	2.24	0.67
1:B:400:ASN:O	1:B:401:ILE:HD12	1.94	0.67
1:B:137:LEU:HD11	1:B:375:MSE:HE2	1.77	0.67
1:B:297:MSE:HG3	1:B:298:PRO:CD	2.25	0.67
1:B:237:GLU:OE2	1:B:239:ARG:HG3	1.95	0.67
1:B:293:LEU:HD21	1:B:297:MSE:HG2	1.77	0.66
1:A:165:TYR:O	1:A:282:ARG:HD3	1.96	0.66
1:B:152:ALA:O	1:B:158:ASN:HB3	1.96	0.66
1:A:227:ARG:HB3	1:A:251:LEU:HD23	1.79	0.65
1:A:263:ILE:HB	1:A:264:PRO:HD3	1.78	0.65
1:A:147:VAL:CG1	1:A:330:THR:HG21	2.26	0.64
1:A:293:LEU:HD11	1:A:297:MSE:HG2	1.80	0.64
1:B:206:LEU:HD12	1:B:207:ARG:N	2.14	0.63
1:A:169:GLU:H	1:A:169:GLU:CD	2.05	0.63
1:A:115:ASN:C	1:A:115:ASN:HD22	2.06	0.62
1:A:195:LEU:HD13	1:A:338:TYR:HD2	1.64	0.62
1:B:137:LEU:HD23	1:B:379:LEU:CD1	2.29	0.61
1:B:380:ASN:ND2	1:B:387:ILE:HD11	2.15	0.61
1:A:195:LEU:HD13	1:A:338:TYR:CD2	2.35	0.61
1:B:308:ILE:HG12	1:B:375:MSE:CE	2.30	0.60
1:A:297:MSE:CE	1:A:323:LEU:HG	2.30	0.60
1:A:232:GLN:HG3	1:A:272:TYR:CZ	2.37	0.59
1:B:214:ASP:O	1:B:215:ASN:HB2	2.00	0.59
1:A:301:LEU:HD22	1:A:326:ILE:HG13	1.86	0.58
1:A:114:LEU:HD12	1:A:115:ASN:N	2.20	0.57
1:B:301:LEU:HD21	1:B:322:MSE:HG3	1.87	0.57
1:B:197:ARG:HH11	1:B:197:ARG:CB	2.04	0.57
1:A:263:ILE:HG13	3:A:704:HOH:O	2.04	0.56
1:B:245:LEU:C	1:B:245:LEU:HD12	2.31	0.56
1:A:349:VAL:HG22	1:A:350:CYS:N	2.20	0.56
1:A:106:LEU:HD12	1:A:106:LEU:C	2.31	0.55
1:B:233:LEU:C	1:B:233:LEU:HD13	2.31	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:VAL:O	1:A:277:THR:HG21	2.06	0.55
1:B:308:ILE:HD12	1:B:309:SER:N	2.22	0.54
1:A:297:MSE:HG3	1:A:298:PRO:HD3	1.89	0.54
1:A:147:VAL:HG13	1:A:172:ARG:HH12	1.72	0.54
1:A:329:MSE:HE2	1:A:398:PHE:CE1	2.42	0.54
1:A:181:CYS:HB3	1:A:328:MSE:HE3	1.90	0.54
1:B:293:LEU:HD21	1:B:297:MSE:CG	2.37	0.54
1:A:308:ILE:HD13	1:A:375:MSE:HE3	1.91	0.53
1:A:262:ASP:HB2	3:A:681:HOH:O	2.09	0.52
1:A:206:LEU:C	1:A:206:LEU:HD23	2.35	0.51
1:B:345:ARG:HB3	1:B:345:ARG:NH2	2.23	0.51
1:B:313:ILE:HG22	1:B:371:PHE:CD2	2.45	0.51
1:A:237:GLU:OE1	1:A:239:ARG:NH2	2.44	0.51
1:A:240:PHE:HA	1:A:246:TYR:CE2	2.45	0.51
1:A:308:ILE:HD11	1:A:375:MSE:HE3	1.93	0.50
1:B:304:ILE:HG22	1:B:375:MSE:HE1	1.91	0.50
1:A:194:GLN:OE1	1:A:218:GLN:NE2	2.45	0.50
1:A:330:THR:HG22	1:A:398:PHE:HD2	1.76	0.50
1:A:234:VAL:HG23	1:A:274:PHE:HB2	1.94	0.49
1:A:250:ILE:HG23	1:A:289:PRO:HB2	1.95	0.49
1:B:344:LYS:HB2	1:B:388:TYR:HD1	1.77	0.49
1:A:332:CYS:O	1:A:397:GLY:HA3	2.12	0.49
1:A:283:ARG:O	1:A:286:PRO:HD3	2.13	0.49
1:B:263:ILE:HB	1:B:264:PRO:HD3	1.94	0.48
1:B:137:LEU:HD21	1:B:375:MSE:HE2	1.95	0.48
1:A:377:LYS:HD3	1:A:377:LYS:C	2.38	0.48
1:B:126:PRO:O	1:B:127:GLY:C	2.56	0.48
1:B:304:ILE:O	1:B:308:ILE:HG13	2.14	0.48
1:A:126:PRO:O	1:A:127:GLY:C	2.57	0.47
1:B:175:VAL:HG11	1:B:226:ILE:HD11	1.96	0.47
1:A:297:MSE:CE	1:A:298:PRO:HA	2.24	0.47
1:A:235:THR:OG1	1:A:274:PHE:HB3	2.14	0.47
1:B:117:ASN:HA	3:B:664:HOH:O	2.15	0.47
1:A:175:VAL:HG21	1:A:226:ILE:HD11	1.96	0.47
1:B:181:CYS:HB3	1:B:328:MSE:HE3	1.97	0.47
1:B:244:SER:O	1:B:247:THR:HB	2.14	0.47
1:A:301:LEU:HD21	1:A:322:MSE:HG3	1.95	0.47
1:B:297:MSE:HG3	1:B:298:PRO:N	2.29	0.47
1:A:330:THR:HG22	1:A:398:PHE:CD2	2.51	0.46
1:B:333:ASP:HB3	1:B:400:ASN:HD21	1.80	0.46
1:B:237:GLU:HG2	3:B:692:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:401:ILE:HG23	3:B:683:HOH:O	2.15	0.46
1:B:288:GLN:HA	1:B:289:PRO:HD3	1.86	0.46
1:A:155:PHE:CD2	1:A:156:PRO:HA	2.51	0.45
1:A:214:ASP:O	1:A:215:ASN:HB2	2.14	0.45
1:B:175:VAL:HG12	1:B:178:TRP:CH2	2.51	0.45
1:B:256:PRO:O	1:B:271:ASP:HB2	2.17	0.45
1:A:377:LYS:HD3	1:A:377:LYS:O	2.17	0.45
1:B:314:GLN:OE1	1:B:314:GLN:N	2.45	0.45
1:A:227:ARG:HD2	3:A:618:HOH:O	2.16	0.45
1:A:147:VAL:HG13	1:A:172:ARG:HH11	1.77	0.45
1:B:247:THR:HG23	3:B:693:HOH:O	2.15	0.45
1:A:297:MSE:HE3	1:A:298:PRO:N	2.31	0.44
1:B:206:LEU:HD12	1:B:207:ARG:H	1.82	0.44
1:B:302:TRP:CE2	1:B:313:ILE:HD11	2.52	0.44
1:A:140:HIS:ND1	1:A:144:HIS:ND1	2.65	0.44
1:B:144:HIS:HE1	3:B:665:HOH:O	2.00	0.44
1:B:265:LYS:HD2	1:B:265:LYS:O	2.17	0.44
1:A:131:LYS:HE2	1:A:131:LYS:HB3	1.84	0.44
1:B:399:ARG:HB2	3:B:603:HOH:O	2.18	0.44
1:A:162:TRP:CE3	1:A:279:LYS:NZ	2.86	0.44
1:A:367:HIS:C	1:A:369:LEU:H	2.26	0.44
1:B:374:ASN:N	1:B:374:ASN:HD22	2.16	0.44
1:A:171:PHE:CE1	1:A:175:VAL:HG11	2.52	0.44
1:B:194:GLN:HA	1:B:218:GLN:O	2.17	0.44
1:A:115:ASN:ND2	1:A:116:MSE:HG3	2.32	0.43
1:A:190:LEU:HG	1:A:339:GLU:HG3	1.98	0.43
1:A:146:ASN:OD1	1:A:146:ASN:C	2.61	0.43
1:A:239:ARG:HD2	3:A:626:HOH:O	2.17	0.43
1:B:117:ASN:O	1:B:118:LYS:C	2.62	0.43
1:B:234:VAL:CG2	1:B:274:PHE:HB2	2.48	0.43
1:B:210:GLY:O	1:B:211:ALA:C	2.60	0.43
1:A:231:SER:HB3	1:A:255:ASP:HB2	2.00	0.43
1:A:245:LEU:HD12	1:A:245:LEU:C	2.43	0.43
1:B:169:GLU:CD	1:B:174:LYS:HD3	2.43	0.43
1:B:232:GLN:O	1:B:236:THR:HB	2.19	0.43
1:B:234:VAL:HG23	1:B:274:PHE:HB2	1.99	0.42
1:A:209:ASN:HB3	1:A:233:LEU:HD22	2.01	0.42
1:B:187:ALA:C	1:B:189:SER:N	2.77	0.42
1:A:182:ALA:HA	1:A:336:ASP:O	2.20	0.42
1:B:208:PHE:CZ	1:B:321:GLY:HA2	2.55	0.42
1:B:263:ILE:N	1:B:264:PRO:CD	2.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:ASP:O	1:A:157:PHE:HB2	2.19	0.42
1:B:187:ALA:C	1:B:189:SER:H	2.27	0.42
1:A:115:ASN:C	1:A:115:ASN:ND2	2.74	0.42
1:B:169:GLU:HB2	1:B:173:THR:HG23	2.01	0.42
1:B:188:GLY:HA2	1:B:351:TYR:N	2.35	0.42
1:B:316:ASN:ND2	1:B:366:TYR:HD2	2.18	0.42
1:A:148:SER:C	1:A:149:MSE:HE2	2.45	0.41
1:B:138:ARG:HD3	1:B:401:ILE:CG1	2.50	0.41
1:B:331:LEU:C	1:B:399:ARG:HH21	2.28	0.41
1:B:345:ARG:HH21	1:B:345:ARG:CB	2.26	0.41
1:A:349:VAL:CG2	1:A:350:CYS:N	2.82	0.41
1:A:351:TYR:CD1	1:A:351:TYR:C	2.98	0.41
1:B:314:GLN:C	1:B:316:ASN:H	2.29	0.41
1:A:354:GLN:CA	1:A:354:GLN:NE2	2.74	0.41
1:A:208:PHE:CZ	1:A:321:GLY:HA2	2.56	0.41
1:A:217:GLN:HA	1:A:221:GLY:O	2.21	0.40
1:A:314:GLN:C	1:A:316:ASN:H	2.30	0.40
1:B:344:LYS:HB2	1:B:388:TYR:CD1	2.55	0.40
1:A:169:GLU:CD	1:A:169:GLU:N	2.76	0.40
1:A:175:VAL:HG13	1:A:175:VAL:O	2.21	0.40
1:B:138:ARG:HD3	1:B:401:ILE:HG12	2.03	0.40
1:B:210:GLY:C	1:B:227:ARG:HH21	2.29	0.40
1:B:260:HIS:CD2	3:B:604:HOH:O	2.75	0.40
1:B:322:MSE:HE3	1:B:325:ILE:HB	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	289/310 (93%)	277 (96%)	10 (4%)	2 (1%)	18 28

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	267/310 (86%)	246 (92%)	19 (7%)	2 (1%)	18	28
All	All	556/620 (90%)	523 (94%)	29 (5%)	4 (1%)	18	28

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	127	GLY
1	B	127	GLY
1	A	174	LYS
1	B	342	PRO

5.3.2 Protein sidechains [i](#)

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	267/273 (98%)	259 (97%)	8 (3%)	36	58
1	B	249/273 (91%)	238 (96%)	11 (4%)	25	43
All	All	516/546 (94%)	497 (96%)	19 (4%)	30	51

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	110	TRP
1	A	115	ASN
1	A	131	LYS
1	A	151	GLU
1	A	169	GLU
1	A	297	MSE
1	A	354	GLN
1	A	377	LYS
1	B	147	VAL
1	B	156	PRO
1	B	161	GLU
1	B	197	ARG

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Mol	Chain	Res	Type
1	B	234	VAL
1	B	243	ASP
1	B	297	MSE
1	B	341	LEU
1	B	345	ARG
1	B	388	TYR
1	B	401	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	112	ASN
1	A	115	ASN
1	A	179	GLN
1	A	192	ASN
1	A	194	GLN
1	A	215	ASN
1	A	218	GLN
1	A	354	GLN
1	A	378	HIS
1	B	192	ASN
1	B	194	GLN
1	B	218	GLN
1	B	268	GLN
1	B	316	ASN
1	B	374	ASN
1	B	400	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

4 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	502	1	14,14,15	1.14	0	17,19,21	1.58	5 (29%)
2	NAG	B	501	1	14,14,15	1.27	1 (7%)	17,19,21	1.54	5 (29%)
2	NAG	A	501	1	14,14,15	1.32	2 (14%)	17,19,21	1.61	5 (29%)
2	NAG	B	502	1	14,14,15	1.18	0	17,19,21	1.53	5 (29%)

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	502	1	-	4/6/23/26	0/1/1/1
2	NAG	B	501	1	-	0/6/23/26	0/1/1/1
2	NAG	A	501	1	-	0/6/23/26	0/1/1/1
2	NAG	B	502	1	-	4/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	NAG	C4-C5	2.62	1.58	1.53
2	B	501	NAG	C4-C5	2.12	1.57	1.53
2	A	501	NAG	C4-C3	2.07	1.57	1.52

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	502	NAG	C1-O5-C5	3.10	116.34	112.19
2	B	501	NAG	C1-O5-C5	3.07	116.30	112.19
2	A	501	NAG	C4-C3-C2	-3.04	106.56	111.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	502	NAG	C1-O5-C5	2.96	116.16	112.19
2	A	502	NAG	C4-C3-C2	-2.87	106.81	111.02
2	B	502	NAG	C4-C3-C2	-2.61	107.19	111.02
2	B	501	NAG	C4-C3-C2	-2.60	107.20	111.02
2	A	501	NAG	C1-O5-C5	2.51	115.55	112.19
2	A	501	NAG	O3-C3-C4	2.31	115.82	110.38
2	A	502	NAG	O5-C5-C6	2.26	112.07	107.66
2	B	501	NAG	O5-C5-C6	2.10	111.76	107.66
2	A	501	NAG	O5-C5-C4	-2.10	105.72	110.83
2	A	502	NAG	O5-C5-C4	-2.10	105.73	110.83
2	B	501	NAG	O3-C3-C4	2.06	115.24	110.38
2	B	502	NAG	O3-C3-C4	2.06	115.23	110.38
2	B	502	NAG	O5-C5-C4	-2.05	105.83	110.83
2	A	501	NAG	O5-C5-C6	2.05	111.65	107.66
2	A	502	NAG	O3-C3-C4	2.04	115.19	110.38
2	B	501	NAG	O5-C5-C4	-2.04	105.86	110.83
2	B	502	NAG	O5-C5-C6	2.00	111.56	107.66

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	502	NAG	O5-C5-C6-O6
2	A	502	NAG	C4-C5-C6-O6
2	B	502	NAG	O5-C5-C6-O6
2	A	502	NAG	C8-C7-N2-C2
2	B	502	NAG	C8-C7-N2-C2
2	B	502	NAG	O7-C7-N2-C2
2	B	502	NAG	C4-C5-C6-O6
2	A	502	NAG	O7-C7-N2-C2

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

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	284/310 (91%)	-0.03	4 (1%) 73 69	7, 21, 38, 65	0
1	B	266/310 (85%)	0.06	8 (3%) 52 48	8, 24, 40, 48	0
All	All	550/620 (88%)	0.02	12 (2%) 62 58	7, 23, 40, 65	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	352	TYR	2.8
1	B	350	CYS	2.7
1	B	388	TYR	2.7
1	B	262	ASP	2.6
1	A	362	THR	2.6
1	A	173	THR	2.5
1	B	127	GLY	2.5
1	A	105	ARG	2.5
1	A	106	LEU	2.3
1	B	344	LYS	2.3
1	B	343	SER	2.1
1	B	366	TYR	2.1

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(Å ²)	Q<0.9
2	NAG	B	502	14/15	0.62	0.26	89,89,89,89	0
2	NAG	B	501	14/15	0.69	0.23	67,69,70,71	0
2	NAG	A	502	14/15	0.75	0.18	58,63,67,69	0
2	NAG	A	501	14/15	0.78	0.11	29,33,35,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.