



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 16, 2026 – 10:28 AM UTC

PDB ID : 4AP5 / pdb_00004ap5
Title : Crystal structure of human POFUT2
Authors : Chen, C.; Keusch, J.J.; Klein, D.; Hess, D.; Hofsteenge, J.; Gut, H.
Deposited on : 2012-03-30
Resolution : 3.00 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

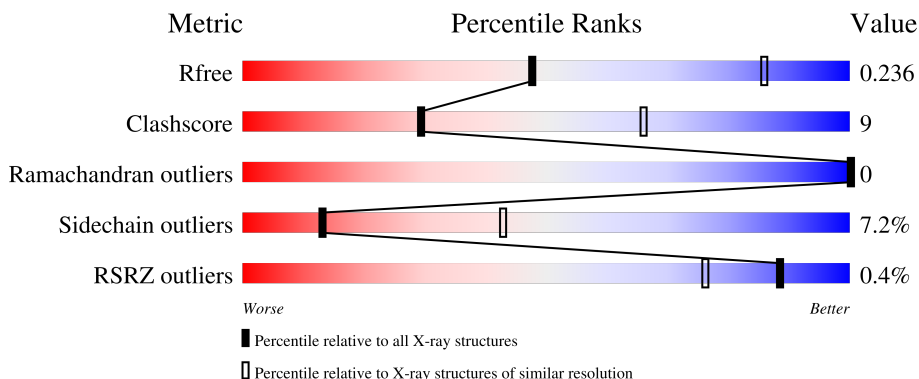
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	408	 70% 23% • 5%
1	B	408	 70% 24% • 5%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6687 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 GDP-FUCOSE PROTEIN O-FUCOSYLTRANSFERASE 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	389	Total	C	N	O	S	0	1	0
			3262	2090	575	586	11			
1	B	389	Total	C	N	O	S	0	0	0
			3251	2084	571	585	11			

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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	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 CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Mg	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	62	Total	O	0	0
			62	62		
5	B	54	Total	O	0	0
			54	54		

4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	118.63Å 118.63Å 196.23Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.66 – 3.00 29.66 – 3.00	Depositor EDS
% Data completeness (in resolution range)	96.3 (29.66-3.00) 96.1 (29.66-3.00)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.09 (at 3.00Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.173 , 0.236 0.172 , 0.236	Depositor DCC
R_{free} test set	1594 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	50.9	Xtriage
Anisotropy	0.487	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 42.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.017 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6687	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.39% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, CL, MG

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	0.54	0/3355	0.96	7/4542 (0.2%)
1	B	0.48	0/3344	0.89	7/4528 (0.2%)
All	All	0.51	0/6699	0.92	14/9070 (0.2%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	102	ILE	N-CA-C	6.10	113.92	107.76
1	A	52	PRO	N-CA-C	5.70	117.65	110.70
1	A	129	SER	N-CA-C	-5.68	103.03	110.53
1	B	52	PRO	N-CA-C	5.66	117.61	110.70
1	B	417	LYS	N-CA-C	5.64	117.23	111.14
1	B	403	ASP	CA-C-N	5.55	125.65	119.32
1	B	403	ASP	C-N-CA	5.55	125.65	119.32
1	A	363	LYS	CB-CA-C	-5.43	109.33	115.79
1	A	196	GLN	N-CA-C	5.38	119.32	112.12
1	A	365	GLY	N-CA-C	-5.36	106.19	113.37
1	A	257	HIS	N-CA-C	5.16	119.72	113.38
1	B	102	ILE	CB-CA-C	-5.15	105.20	110.71
1	B	365	GLY	N-CA-C	-5.10	107.99	114.16
1	B	129	SER	N-CA-C	-5.06	103.37	110.35

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	3262	0	3174	62	0
1	B	3251	0	3162	54	0
2	A	28	0	26	0	0
2	B	28	0	26	0	0
3	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	62	0	0	0	1
5	B	54	0	0	1	1
All	All	6687	0	6388	112	1

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 9.

All (112) 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:378:ALA:O	1:B:397:ARG:NH2	2.14	0.81
1:B:410:ARG:HG2	1:B:422:PRO:HD3	1.70	0.73
1:A:284:LEU:HG	1:A:379:ARG:HH12	1.54	0.72
1:A:88[A]:ARG:HH11	1:A:88[A]:ARG:HG3	1.56	0.70
1:A:97:ILE:HD12	1:A:99:GLN:HE21	1.57	0.69
1:B:309:LEU:HD22	1:B:342:LEU:HD13	1.73	0.69
1:A:294:ARG:HB3	1:A:384:THR:HG21	1.77	0.67
1:A:219:ARG:NH1	1:A:221:GLU:OE2	2.26	0.66
1:A:378:ALA:O	1:A:397:ARG:NH2	2.30	0.64
1:B:270:GLN:HB3	1:B:276:MET:HB2	1.81	0.62
1:B:392:ARG:NH2	5:B:2045:HOH:O	2.28	0.62
1:B:285:GLY:O	1:B:379:ARG:NH1	2.31	0.62
1:B:137:VAL:HB	1:B:190:VAL:HG22	1.82	0.61
1:A:86:TRP:HB2	1:A:102:ILE:HB	1.83	0.61
1:B:120:ILE:HB	1:B:124:GLN:HG2	1.83	0.61
1:A:104:TRP:N	1:A:121:GLU:OE2	2.35	0.60
1:B:379:ARG:HH11	1:B:379:ARG:HG3	1.65	0.60
1:A:379:ARG:HH11	1:A:379:ARG:HG3	1.67	0.59
1:A:294:ARG:NH1	1:A:297:ASP:OD2	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:HIS:HB2	1:B:160:PRO:HB3	1.82	0.59
1:A:132:PRO:HG3	1:A:184:GLU:HG2	1.85	0.59
1:A:312:ALA:O	1:A:316:ILE:HG13	2.03	0.58
1:A:143:TYR:CE2	1:A:155:LYS:HE2	2.39	0.56
1:A:106:GLU:OE1	1:A:363:LYS:NZ	2.39	0.56
1:B:198:SER:HB3	1:B:226:ASP:HA	1.87	0.55
1:A:51:ASN:HB2	1:A:221:GLU:OE2	2.08	0.54
1:A:177:GLY:HA2	1:B:92:TRP:CD1	2.43	0.53
1:A:143:TYR:CZ	1:A:155:LYS:HE2	2.44	0.53
1:B:43:ARG:NH2	1:B:128:GLU:OE1	2.40	0.53
1:B:83:LEU:HB3	1:B:104:TRP:CD1	2.44	0.53
1:B:206:LEU:HD21	1:B:215:VAL:HG11	1.90	0.53
1:A:97:ILE:HD12	1:A:99:GLN:NE2	2.24	0.52
1:B:410:ARG:HH22	1:B:416:GLU:CD	2.17	0.52
1:B:94:SER:HB3	1:B:96:ASP:HB2	1.92	0.51
1:A:88[A]:ARG:HH11	1:A:88[A]:ARG:CG	2.23	0.51
1:B:244:ARG:HA	1:B:247:ARG:HH21	1.76	0.51
1:B:399:ILE:HA	1:B:427:ILE:HD11	1.94	0.50
1:B:116:ASN:ND2	1:B:236:THR:HA	2.27	0.50
1:A:198:SER:CB	1:A:226:ASP:HA	2.41	0.49
1:B:219:ARG:NH1	1:B:221:GLU:OE2	2.45	0.49
1:B:69:LEU:HD13	1:B:199:ALA:O	2.13	0.49
1:A:126:ILE:HD13	1:A:132:PRO:HB3	1.95	0.49
1:A:362:TYR:O	1:A:366:GLY:HA3	2.13	0.49
1:A:270:GLN:HB3	1:A:276:MET:HB2	1.93	0.49
1:B:86:TRP:HB2	1:B:102:ILE:HB	1.94	0.48
1:B:304:GLN:O	1:B:413:GLY:HA3	2.13	0.48
1:A:99:GLN:OE1	1:A:364:ASP:HB2	2.13	0.48
1:B:51:ASN:OD1	1:B:52:PRO:HD2	2.13	0.48
1:B:296:LYS:HB2	1:B:333:ASP:OD2	2.13	0.47
1:B:403:ASP:HB3	1:B:406:THR:HG23	1.97	0.47
1:B:410:ARG:NH2	1:B:416:GLU:OE1	2.45	0.47
1:A:298:PHE:CD1	1:A:306:VAL:HG22	2.50	0.47
1:B:296:LYS:HB2	1:B:296:LYS:HE3	1.62	0.47
1:B:309:LEU:HD11	1:B:338:GLU:HG2	1.97	0.47
1:A:64:ILE:HG21	1:A:233:TYR:CD2	2.50	0.46
1:B:336:ARG:O	1:B:340:GLU:HG2	2.15	0.46
1:A:96:ASP:HB3	1:A:356:TRP:HH2	1.80	0.46
1:A:88[A]:ARG:HG3	1:A:88[A]:ARG:NH1	2.23	0.46
1:A:320:MET:HE1	1:A:348:GLU:HG3	1.98	0.46
1:A:395:GLU:O	1:A:399:ILE:HG13	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:392:ARG:HD3	1:A:392:ARG:HA	1.75	0.45
1:A:47:LEU:O	1:A:216:MET:HA	2.17	0.45
1:A:165:LEU:HD22	1:B:300:TRP:CD1	2.51	0.45
1:A:101:ARG:HG2	1:A:182:TYR:CE1	2.51	0.45
1:B:54:GLU:O	1:B:88:ARG:NH1	2.41	0.45
1:B:409:ASN:HA	1:B:421:GLN:OE1	2.16	0.45
1:A:109:ASP:HB3	1:A:241:VAL:HG22	1.97	0.45
1:A:176:ARG:HH12	1:B:95:PRO:HA	1.82	0.45
1:B:292:HIS:CE1	1:B:389:PHE:HB3	2.53	0.45
1:B:392:ARG:HD3	1:B:392:ARG:HA	1.38	0.45
1:A:42:ARG:HA	1:A:42:ARG:HD3	1.68	0.44
1:A:206:LEU:CD2	1:A:215:VAL:HG11	2.48	0.44
1:A:122:TYR:CE2	1:A:126:ILE:HD11	2.53	0.44
1:B:47:LEU:O	1:B:216:MET:HA	2.17	0.44
1:B:368:ALA:O	1:B:372:GLN:HG3	2.17	0.44
1:A:205:LEU:C	1:A:205:LEU:HD23	2.41	0.44
1:B:403:ASP:HA	1:B:404:PRO:HD2	1.80	0.43
1:A:298:PHE:CE1	1:A:306:VAL:HG22	2.53	0.43
1:A:256:ARG:HA	1:A:256:ARG:HD3	1.75	0.43
1:B:408:TYR:HB3	1:B:422:PRO:HG2	2.00	0.43
1:A:339:TYR:CE1	1:A:351:ARG:HD2	2.53	0.43
1:B:64:ILE:HA	1:B:236:THR:HG21	2.00	0.43
1:A:351:ARG:HG2	1:A:352:PHE:N	2.34	0.43
1:A:304:GLN:O	1:A:416:GLU:HB2	2.19	0.43
1:B:95:PRO:HD2	1:B:364:ASP:OD2	2.18	0.43
1:A:327:LYS:HE2	1:A:327:LYS:HB2	1.75	0.42
1:B:112:SER:HB3	1:B:241:VAL:HG13	2.00	0.42
1:B:206:LEU:HD23	1:B:206:LEU:HA	1.84	0.42
1:B:286:GLY:O	1:B:379:ARG:HB2	2.19	0.42
1:A:188:LEU:HD23	1:A:188:LEU:HA	1.83	0.42
1:A:379:ARG:HH11	1:A:379:ARG:CG	2.31	0.42
1:A:385:SER:HB2	1:A:410:ARG:HD2	2.01	0.42
1:B:355:THR:OG1	1:B:358:GLU:HG3	2.20	0.42
1:A:100:VAL:HG22	1:A:101:ARG:HG3	2.02	0.42
1:A:110:LEU:HD11	1:A:121:GLU:OE1	2.20	0.42
1:A:62:VAL:HA	1:A:65:ARG:HD3	2.01	0.41
1:A:399:ILE:HA	1:A:427:ILE:HD11	2.02	0.41
1:B:203:ALA:O	1:B:207:LEU:HG	2.20	0.41
1:A:403:ASP:OD1	1:A:404:PRO:HD2	2.21	0.41
1:B:274:MET:HE3	1:B:274:MET:HB2	1.92	0.41
1:A:372:GLN:HG2	1:A:393:ILE:HG12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:291:VAL:HG22	1:A:330:VAL:HG22	2.01	0.41
1:A:373:TRP:CZ2	1:A:377:HIS:HE1	2.38	0.41
1:B:64:ILE:HG21	1:B:233:TYR:CD1	2.56	0.41
1:B:339:TYR:CE1	1:B:351:ARG:HD2	2.55	0.41
1:A:284:LEU:HG	1:A:379:ARG:NH1	2.29	0.41
1:A:373:TRP:CZ2	1:A:377:HIS:CE1	3.09	0.41
1:A:73:LEU:HD13	1:A:79:TRP:CD2	2.56	0.41
1:A:62:VAL:O	1:A:65:ARG:HB2	2.22	0.40
1:B:100:VAL:HG22	1:B:101:ARG:N	2.36	0.40
1:B:110:LEU:HA	1:B:110:LEU:HD23	1.88	0.40
1:B:231:LYS:HE3	1:B:231:LYS:HB2	1.75	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:2030:HOH:O	5:B:2047:HOH:O[3_655]	2.00	0.20

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	388/408 (95%)	374 (96%)	14 (4%)	0	100	100
1	B	387/408 (95%)	369 (95%)	18 (5%)	0	100	100
All	All	775/816 (95%)	743 (96%)	32 (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	349/360 (97%)	323 (93%)	26 (7%)	13	42
1	B	348/360 (97%)	323 (93%)	25 (7%)	13	43
All	All	697/720 (97%)	646 (93%)	51 (7%)	13	42

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	LEU
1	A	65	ARG
1	A	72	THR
1	A	80	VAL
1	A	88[A]	ARG
1	A	88[B]	ARG
1	A	98	HIS
1	A	114	ASN
1	A	119	VAL
1	A	121	GLU
1	A	189	ASN
1	A	198	SER
1	A	208	ARG
1	A	211	SER
1	A	213	ARG
1	A	214	SER
1	A	219	ARG
1	A	241	VAL
1	A	275	LYS
1	A	291	VAL
1	A	306	VAL
1	A	328	VAL
1	A	364	ASP
1	A	406	THR
1	A	419	CYS
1	A	428	THR
1	B	47	LEU

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Mol	Chain	Res	Type
1	B	50	VAL
1	B	98	HIS
1	B	115	LYS
1	B	117	ILE
1	B	119	VAL
1	B	164	GLN
1	B	196	GLN
1	B	211	SER
1	B	213	ARG
1	B	241	VAL
1	B	276	MET
1	B	278	VAL
1	B	291	VAL
1	B	294	ARG
1	B	306	VAL
1	B	324	ARG
1	B	326	ASP
1	B	357	GLU
1	B	384	THR
1	B	388	THR
1	B	392	ARG
1	B	402	LEU
1	B	406	THR
1	B	428	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	HIS
1	A	114	ASN
1	A	169	GLN
1	A	209	ASN
1	A	270	GLN
1	A	377	HIS
1	B	93	GLN
1	B	222	ASN
1	B	304	GLN

5.3.3 RNA ⓘ

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 [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

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	B	1431	1	14,14,15	0.65	0	17,19,21	1.76	3 (17%)
2	NAG	A	1430	1	14,14,15	0.58	0	17,19,21	1.26	2 (11%)
2	NAG	B	1430	1	14,14,15	0.43	0	17,19,21	1.15	1 (5%)
2	NAG	A	1431	1	14,14,15	0.52	0	17,19,21	0.71	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	B	1431	1	-	3/6/23/26	0/1/1/1
2	NAG	A	1430	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1430	1	-	2/6/23/26	0/1/1/1
2	NAG	A	1431	1	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1431	NAG	C2-N2-C7	4.07	128.36	122.90
2	B	1431	NAG	C1-O5-C5	3.64	117.06	112.19
2	B	1431	NAG	C1-C2-N2	3.18	115.44	110.43
2	A	1430	NAG	O5-C1-C2	3.12	116.11	111.29
2	A	1430	NAG	C1-O5-C5	2.82	115.96	112.19
2	B	1430	NAG	O5-C1-C2	2.48	115.13	111.29

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1431	NAG	C1-C2-N2-C7
2	A	1431	NAG	O5-C5-C6-O6
2	A	1431	NAG	C4-C5-C6-O6
2	B	1430	NAG	C4-C5-C6-O6
2	B	1430	NAG	O5-C5-C6-O6
2	B	1431	NAG	C4-C5-C6-O6
2	A	1431	NAG	C8-C7-N2-C2
2	B	1431	NAG	O5-C5-C6-O6
2	A	1431	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 [i](#)

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	389/408 (95%)	-0.79	2 (0%)	87 72	16, 40, 86, 120	1 (0%)
1	B	389/408 (95%)	-0.57	1 (0%)	90 79	25, 58, 106, 136	0
All	All	778/816 (95%)	-0.68	3 (0%)	88 76	16, 49, 100, 136	1 (0%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	356	TRP	3.1
1	A	88[A]	ARG	2.9
1	A	269	PHE	2.3

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	1431	14/15	0.31	0.16	93,112,123,125	0
2	NAG	A	1430	14/15	0.36	0.23	109,129,136,139	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	A	1431	14/15	0.71	0.13	69,98,111,111	0
2	NAG	B	1430	14/15	0.84	0.17	98,118,130,130	0
4	MG	B	1432	1/1	0.92	0.25	47,47,47,47	0
3	CL	A	1432	1/1	0.96	0.07	48,48,48,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.