



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 6S9F / pdb_00006s9f
Title : Drosophila OTK, extracellular domains 3-5
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Deposited on : 2019-07-12
Resolution : 1.97 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

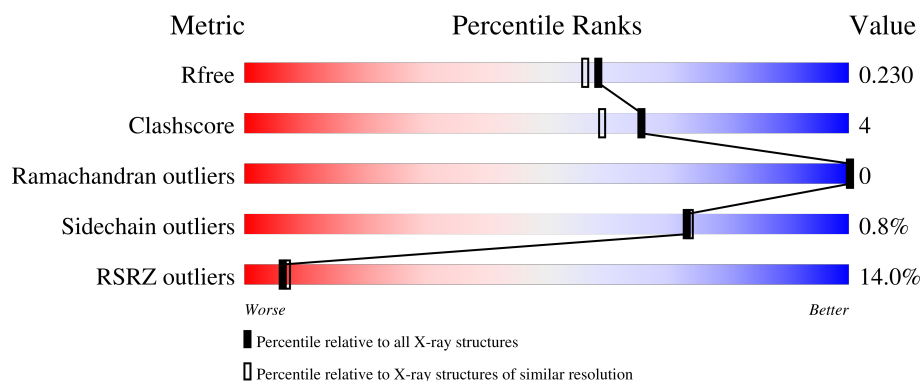
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.97 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	569	5% 49% 48%
1	B	569	10% 45% 7% 47%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5253 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.

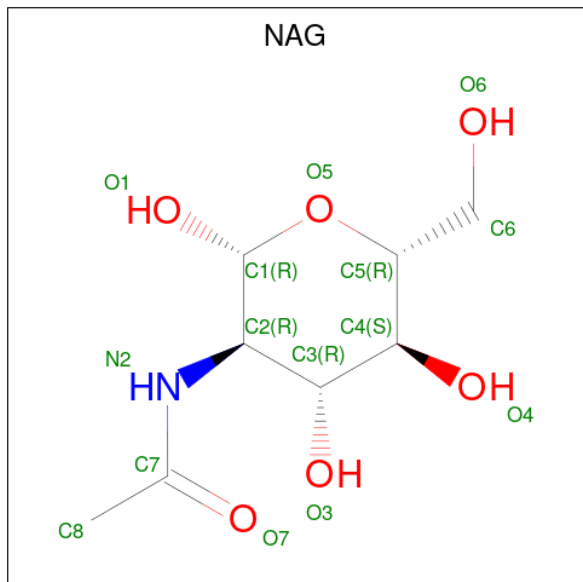
- Molecule 1 is a protein called Tyrosine-protein kinase-like otk.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	298	Total	C	N	O	S	0	0	0
			2302	1435	409	451	7			
1	B	300	Total	C	N	O	S	0	0	0
			2318	1444	413	454	7			

There are 24 discrepancies between the modelled and reference sequences:

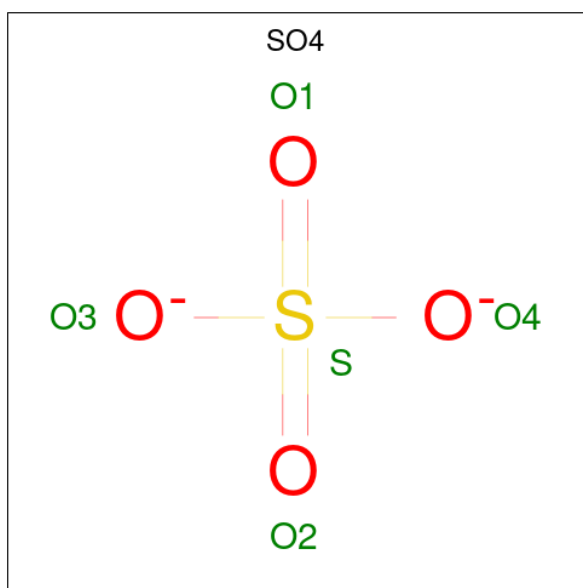
Chain	Residue	Modelled	Actual	Comment	Reference
A	21	GLU	-	expression tag	UNP Q6AWJ9
A	22	THR	-	expression tag	UNP Q6AWJ9
A	23	GLY	-	expression tag	UNP Q6AWJ9
A	581	GLY	-	expression tag	UNP Q6AWJ9
A	582	THR	-	expression tag	UNP Q6AWJ9
A	583	LYS	-	expression tag	UNP Q6AWJ9
A	584	HIS	-	expression tag	UNP Q6AWJ9
A	585	HIS	-	expression tag	UNP Q6AWJ9
A	586	HIS	-	expression tag	UNP Q6AWJ9
A	587	HIS	-	expression tag	UNP Q6AWJ9
A	588	HIS	-	expression tag	UNP Q6AWJ9
A	589	HIS	-	expression tag	UNP Q6AWJ9
B	21	GLU	-	expression tag	UNP Q6AWJ9
B	22	THR	-	expression tag	UNP Q6AWJ9
B	23	GLY	-	expression tag	UNP Q6AWJ9
B	581	GLY	-	expression tag	UNP Q6AWJ9
B	582	THR	-	expression tag	UNP Q6AWJ9
B	583	LYS	-	expression tag	UNP Q6AWJ9
B	584	HIS	-	expression tag	UNP Q6AWJ9
B	585	HIS	-	expression tag	UNP Q6AWJ9
B	586	HIS	-	expression tag	UNP Q6AWJ9
B	587	HIS	-	expression tag	UNP Q6AWJ9
B	588	HIS	-	expression tag	UNP Q6AWJ9
B	589	HIS	-	expression tag	UNP Q6AWJ9

- 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	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		
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 SULFATE ION (CCD ID: SO4) (formula: O₄S).



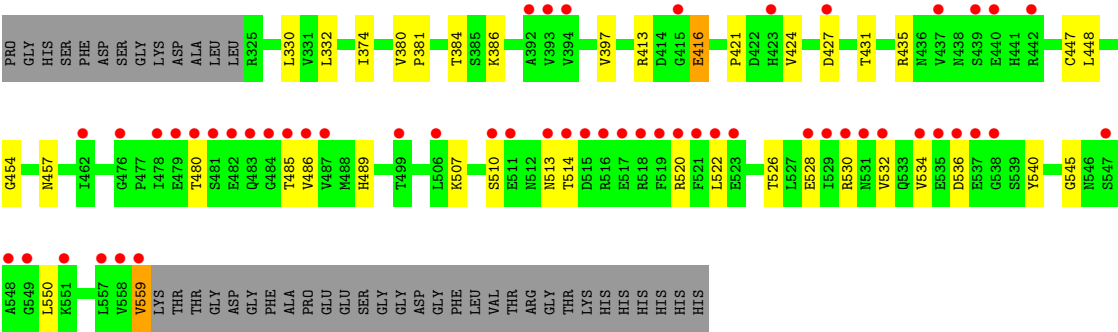
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	239	Total	O	0	0
			239	239		
4	B	191	Total	O	0	0
			191	191		

- Molecule 1: Tyrosine-protein kinase-like otk





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	80.97Å 189.91Å 131.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.48 – 1.97 47.48 – 1.97	Depositor EDS
% Data completeness (in resolution range)	97.3 (47.48-1.97) 97.3 (47.48-1.97)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 1.97Å)	Xtriage
Refinement program	PHENIX dev_3488	Depositor
R, R_{free}	0.207 , 0.235 (Not available) , 0.230	Depositor DCC
R_{free} test set	3596 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	32.2	Xtriage
Anisotropy	0.641	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 49.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5253	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% 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: SO4, 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	0.17	0/2341	0.39	0/3183
1	B	0.15	0/2358	0.37	0/3207
All	All	0.16	0/4699	0.38	0/6390

There are no bond length outliers.

There are no bond angle outliers.

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	2302	0	2285	12	0
1	B	2318	0	2296	25	0
2	A	84	0	78	0	0
2	B	84	0	78	1	0
3	A	20	0	0	0	0
3	B	15	0	0	1	0
4	A	239	0	0	2	0
4	B	191	0	0	0	0
All	All	5253	0	4737	36	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 4.

All (36) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:483:GLN:O	4:A:701:HOH:O	2.12	0.68
1:B:507:LYS:HE3	1:B:513:ASN:HD22	1.57	0.67
1:B:534:VAL:HG12	1:B:559:VAL:HG11	1.76	0.66
1:A:518:ARG:NH2	1:A:536:ASP:OD2	2.33	0.62
1:B:386:LYS:HG3	1:B:397:VAL:HG22	1.85	0.58
1:A:416:GLU:CD	1:A:416:GLU:H	2.13	0.55
1:B:435:ARG:NH2	3:B:608:SO4:O1	2.40	0.55
1:B:421:PRO:HG2	1:B:424:VAL:HG21	1.91	0.53
1:B:510:SER:O	1:B:514:THR:OG1	2.23	0.51
1:B:267:LYS:HA	1:B:374:ILE:HB	1.91	0.51
1:B:413:ARG:HE	1:B:421:PRO:HD3	1.75	0.51
1:B:522:LEU:HD12	1:B:526:THR:HB	1.93	0.51
1:A:282:ASP:HA	1:A:285:LYS:HG2	1.94	0.50
1:B:520:ARG:HB2	1:B:528:GLU:HB3	1.94	0.49
1:B:448:LEU:HD23	1:B:457:ASN:HB3	1.93	0.49
1:B:293:TRP:CD1	1:B:332:LEU:HD13	2.49	0.47
1:B:300:LEU:HD21	1:B:330:LEU:HD21	1.97	0.47
1:B:384:THR:OG1	1:B:386:LYS:NZ	2.49	0.46
1:B:276:CYS:HB2	1:B:293:TRP:CZ2	2.51	0.46
1:A:517:GLU:HB2	4:A:727:HOH:O	2.16	0.46
1:A:253:ASP:OD1	1:A:253:ASP:N	2.41	0.46
1:B:536:ASP:O	1:B:540:TYR:OH	2.33	0.45
1:B:427:ASP:HB3	1:B:431:THR:H	1.82	0.45
1:A:276:CYS:HB2	1:A:293:TRP:CZ2	2.53	0.44
1:B:416:GLU:CD	1:B:416:GLU:H	2.26	0.44
1:A:300:LEU:HD21	1:A:330:LEU:HD21	1.99	0.43
1:A:380:VAL:HA	1:A:381:PRO:HA	1.83	0.43
1:A:267:LYS:O	1:A:270:GLU:HB2	2.20	0.42
1:A:301:ARG:CZ	1:B:454:GLY:HA2	2.50	0.42
1:A:247:LEU:HD12	1:A:247:LEU:H	1.84	0.42
1:B:545:GLY:HA3	1:B:550:LEU:HD23	2.02	0.42
1:B:485:THR:HG22	1:B:530:ARG:HA	2.01	0.42
1:B:489:HIS:HB3	2:B:601:NAG:H83	2.02	0.41
1:B:480:THR:HG21	1:B:486:VAL:HG11	2.02	0.41
1:B:380:VAL:HA	1:B:381:PRO:HA	1.87	0.40
1:B:447:CYS:O	1:B:457:ASN:HA	2.22	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	294/569 (52%)	287 (98%)	7 (2%)	0	100	100
1	B	296/569 (52%)	291 (98%)	5 (2%)	0	100	100
All	All	590/1138 (52%)	578 (98%)	12 (2%)	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	257/487 (53%)	256 (100%)	1 (0%)	84	85
1	B	258/487 (53%)	255 (99%)	3 (1%)	63	61
All	All	515/974 (53%)	511 (99%)	4 (1%)	73	74

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	500	ILE
1	B	416	GLU
1	B	532	VAL
1	B	559	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	263	GLN
1	A	423	HIS
1	A	455	GLN
1	A	463	ASN
1	B	423	HIS

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 [i](#)

19 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	B	604	1	14,14,15	0.24	0	17,19,21	0.47	0
2	NAG	A	606	1	14,14,15	0.27	0	17,19,21	0.48	0
2	NAG	B	603	1	14,14,15	0.29	0	17,19,21	0.44	0
2	NAG	A	604	1	14,14,15	0.28	0	17,19,21	0.49	0
3	SO4	B	609	-	4,4,4	0.24	0	6,6,6	0.13	0
2	NAG	A	601	1	14,14,15	0.30	0	17,19,21	0.60	0
2	NAG	A	602	1	14,14,15	0.32	0	17,19,21	0.44	0
3	SO4	A	609	-	4,4,4	0.23	0	6,6,6	0.07	0
3	SO4	A	608	-	4,4,4	0.24	0	6,6,6	0.08	0
3	SO4	B	607	-	4,4,4	0.23	0	6,6,6	0.09	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	A	605	1	14,14,15	0.44	0	17,19,21	0.74	1 (5%)
2	NAG	B	601	1	14,14,15	0.27	0	17,19,21	0.47	0
2	NAG	A	603	1	14,14,15	0.22	0	17,19,21	0.47	0
3	SO4	B	608	-	4,4,4	0.23	0	6,6,6	0.08	0
3	SO4	A	607	-	4,4,4	0.23	0	6,6,6	0.08	0
2	NAG	B	606	1	14,14,15	0.29	0	17,19,21	0.56	0
2	NAG	B	605	1	14,14,15	0.46	0	17,19,21	0.46	0
3	SO4	A	610	-	4,4,4	0.24	0	6,6,6	0.13	0
2	NAG	B	602	1	14,14,15	0.33	0	17,19,21	0.58	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	604	1	-	2/6/23/26	0/1/1/1
2	NAG	A	601	1	-	2/6/23/26	0/1/1/1
2	NAG	A	603	1	-	0/6/23/26	0/1/1/1
2	NAG	A	606	1	-	0/6/23/26	0/1/1/1
2	NAG	B	606	1	-	4/6/23/26	0/1/1/1
2	NAG	B	603	1	-	0/6/23/26	0/1/1/1
2	NAG	B	605	1	-	2/6/23/26	0/1/1/1
2	NAG	A	602	1	-	0/6/23/26	0/1/1/1
2	NAG	A	605	1	-	0/6/23/26	0/1/1/1
2	NAG	B	601	1	-	2/6/23/26	0/1/1/1
2	NAG	B	602	1	-	4/6/23/26	0/1/1/1
2	NAG	A	604	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	605	NAG	C1-O5-C5	2.55	115.61	112.19

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	601	NAG	O5-C5-C6-O6
2	B	604	NAG	O5-C5-C6-O6
2	B	604	NAG	C4-C5-C6-O6
2	B	605	NAG	C4-C5-C6-O6
2	B	606	NAG	C4-C5-C6-O6
2	B	605	NAG	O5-C5-C6-O6
2	B	606	NAG	C1-C2-N2-C7
2	B	601	NAG	C4-C5-C6-O6
2	B	602	NAG	C4-C5-C6-O6
2	A	601	NAG	C3-C2-N2-C7
2	B	602	NAG	C3-C2-N2-C7
2	B	602	NAG	O5-C5-C6-O6
2	B	606	NAG	O5-C5-C6-O6
2	A	601	NAG	C1-C2-N2-C7
2	B	602	NAG	C1-C2-N2-C7
2	B	606	NAG	C3-C2-N2-C7

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	601	NAG	1	0
3	B	608	SO4	1	0

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 ⓘ

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	298/569 (52%)	0.55	26 (8%)	16 18	23, 40, 73, 101	0
1	B	300/569 (52%)	1.00	58 (19%)	3 3	24, 47, 106, 145	0
All	All	598/1138 (52%)	0.77	84 (14%)	6 7	23, 43, 89, 145	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	559	VAL	6.5
1	B	485	THR	5.9
1	B	480	THR	4.6
1	B	558	VAL	4.3
1	A	394	VAL	4.3
1	B	481	SER	4.1
1	A	534	VAL	4.1
1	B	522	LEU	4.1
1	B	521	PHE	4.1
1	A	246	ALA	4.1
1	A	485	THR	4.1
1	A	308	SER	4.0
1	B	486	VAL	3.9
1	B	529	ILE	3.9
1	B	534	VAL	3.8
1	B	246	ALA	3.8
1	A	558	VAL	3.7
1	B	393	VAL	3.6
1	B	484	GLY	3.6
1	A	561	THR	3.6
1	B	427	ASP	3.5
1	B	519	PHE	3.5
1	B	394	VAL	3.4
1	B	530	ARG	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	559	VAL	3.4
1	B	511	GLU	3.3
1	A	517	GLU	3.3
1	B	547	SER	3.3
1	B	310	PRO	3.2
1	B	523	GLU	3.2
1	A	307	GLY	3.1
1	B	557	LEU	3.1
1	A	257	VAL	3.1
1	A	511	GLU	3.1
1	B	536	ASP	3.0
1	B	514	THR	3.0
1	B	515	ASP	3.0
1	A	499	THR	3.0
1	B	516	ARG	3.0
1	A	247	LEU	2.9
1	B	478	ILE	2.8
1	B	535	GLU	2.8
1	B	510	SER	2.8
1	B	506	LEU	2.8
1	A	550	LEU	2.7
1	B	531	ASN	2.7
1	B	548	ALA	2.7
1	B	423	HIS	2.7
1	B	440	GLU	2.7
1	A	327	ASP	2.6
1	B	517	GLU	2.6
1	B	518	ARG	2.6
1	B	482	GLU	2.6
1	B	282	ASP	2.6
1	A	560	LYS	2.5
1	B	520	ARG	2.5
1	B	487	VAL	2.5
1	B	528	GLU	2.5
1	B	532	VAL	2.5
1	B	483	GLN	2.5
1	A	530	ARG	2.4
1	B	462	ILE	2.4
1	B	499	THR	2.4
1	B	392	ALA	2.4
1	A	531	ASN	2.3
1	B	476	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	484	GLY	2.2
1	B	309	ALA	2.2
1	B	442	ARG	2.2
1	A	532	VAL	2.2
1	B	437	VAL	2.2
1	B	538	GLY	2.2
1	B	479	GLU	2.2
1	B	439	SER	2.2
1	A	360	ALA	2.1
1	A	533	GLN	2.1
1	B	549	GLY	2.1
1	A	423	HIS	2.1
1	A	282	ASP	2.1
1	A	287	GLN	2.1
1	B	551	LYS	2.0
1	B	415	GLY	2.0
1	B	537	GLU	2.0
1	B	513	ASN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	B	606	14/15	0.49	0.17	93,109,117,119	0
2	NAG	B	603	14/15	0.56	0.21	75,88,101,102	0
2	NAG	A	606	14/15	0.65	0.18	99,109,120,123	0
2	NAG	B	602	14/15	0.70	0.19	87,94,103,105	0
2	NAG	B	604	14/15	0.75	0.14	80,87,103,103	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	B	605	14/15	0.78	0.23	83,95,109,113	0
3	SO4	B	608	5/5	0.79	0.11	111,111,112,112	0
2	NAG	A	601	14/15	0.81	0.14	48,60,69,72	0
3	SO4	A	609	5/5	0.82	0.12	94,97,99,100	0
2	NAG	B	601	14/15	0.86	0.18	78,86,94,99	0
3	SO4	A	607	5/5	0.86	0.17	129,131,132,133	0
3	SO4	B	609	5/5	0.89	0.11	50,57,59,68	0
3	SO4	A	608	5/5	0.90	0.10	60,65,70,75	0
2	NAG	A	604	14/15	0.90	0.11	35,46,56,58	0
2	NAG	A	603	14/15	0.91	0.11	38,46,56,64	0
3	SO4	B	607	5/5	0.92	0.10	74,75,76,77	0
3	SO4	A	610	5/5	0.94	0.12	41,46,52,54	0
2	NAG	A	602	14/15	0.94	0.08	33,38,44,47	0
2	NAG	A	605	14/15	0.95	0.07	23,30,36,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.