



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:24 PM UTC

PDB ID : 5OJK / pdb_00005ojk
Title : Crystal structure of the human neuroligin 1 cholinesterase domain containing spliced sequence B (SSB) (NL1(-A+B))
Authors : Elegheert, J.; Aricescu, A.R.
Deposited on : 2017-07-21
Resolution : 2.55 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

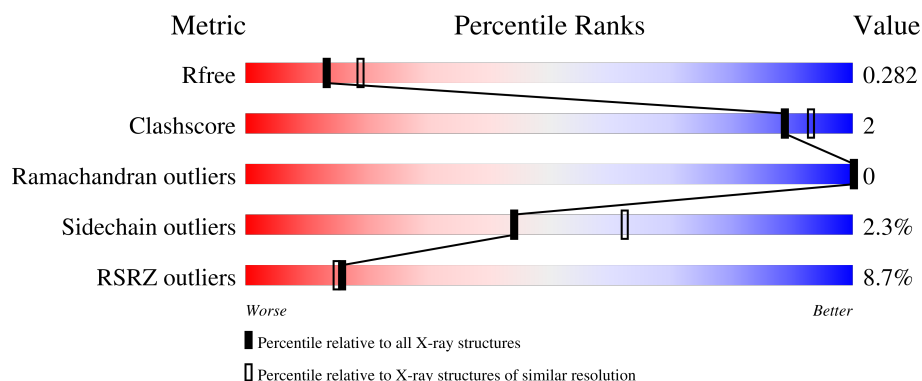
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	585	8% 87% 5% 8%
1	B	585	8% 85% 6% 8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8583 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.

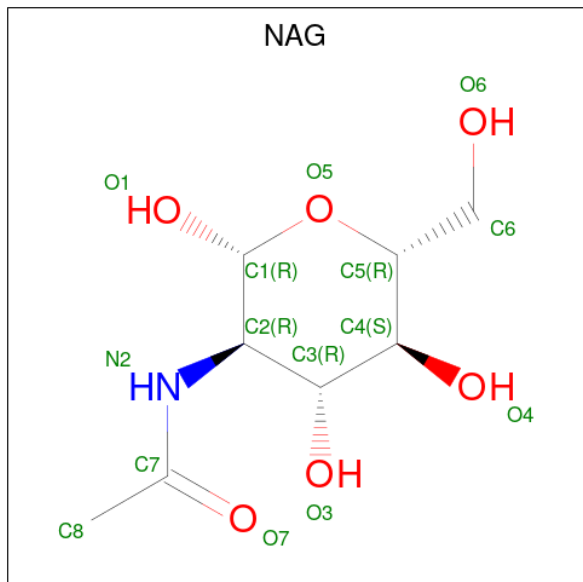
- Molecule 1 is a protein called Neuroligin-1,Neuroligin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	537	Total	C	N	O	S	0	1	0
			4199	2694	700	788	17			
1	B	537	Total	C	N	O	S	0	2	0
			4207	2704	700	786	17			

There are 24 discrepancies between the modelled and reference sequences:

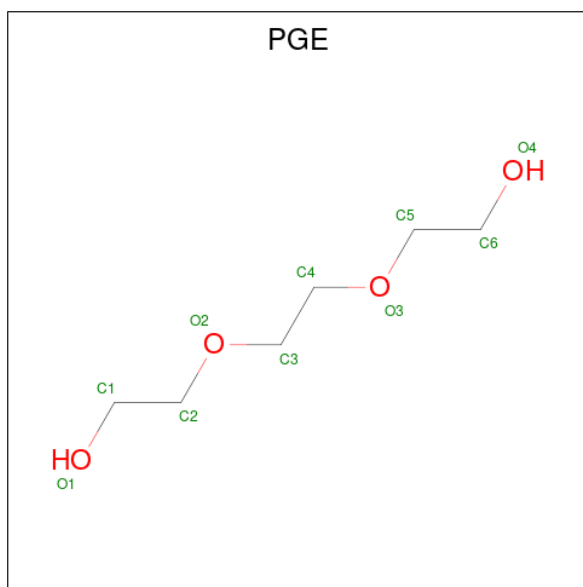
Chain	Residue	Modelled	Actual	Comment	Reference
A	43	GLU	-	expression tag	UNP Q8N2Q7
A	44	THR	-	expression tag	UNP Q8N2Q7
A	45	GLY	-	expression tag	UNP Q8N2Q7
A	636	ARG	-	expression tag	UNP Q8N2Q7
A	637	THR	-	expression tag	UNP Q8N2Q7
A	638	LYS	-	expression tag	UNP Q8N2Q7
A	639	HIS	-	expression tag	UNP Q8N2Q7
A	640	HIS	-	expression tag	UNP Q8N2Q7
A	641	HIS	-	expression tag	UNP Q8N2Q7
A	642	HIS	-	expression tag	UNP Q8N2Q7
A	643	HIS	-	expression tag	UNP Q8N2Q7
A	644	HIS	-	expression tag	UNP Q8N2Q7
B	43	GLU	-	expression tag	UNP Q8N2Q7
B	44	THR	-	expression tag	UNP Q8N2Q7
B	45	GLY	-	expression tag	UNP Q8N2Q7
B	636	ARG	-	expression tag	UNP Q8N2Q7
B	637	THR	-	expression tag	UNP Q8N2Q7
B	638	LYS	-	expression tag	UNP Q8N2Q7
B	639	HIS	-	expression tag	UNP Q8N2Q7
B	640	HIS	-	expression tag	UNP Q8N2Q7
B	641	HIS	-	expression tag	UNP Q8N2Q7
B	642	HIS	-	expression tag	UNP Q8N2Q7
B	643	HIS	-	expression tag	UNP Q8N2Q7
B	644	HIS	-	expression tag	UNP Q8N2Q7

- 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	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			10	6	4		
3	A	1	Total	C	O	0	0
			10	6	4		
3	B	1	Total	C	O	0	0
			10	6	4		
3	B	1	Total	C	O	0	0
			10	6	4		

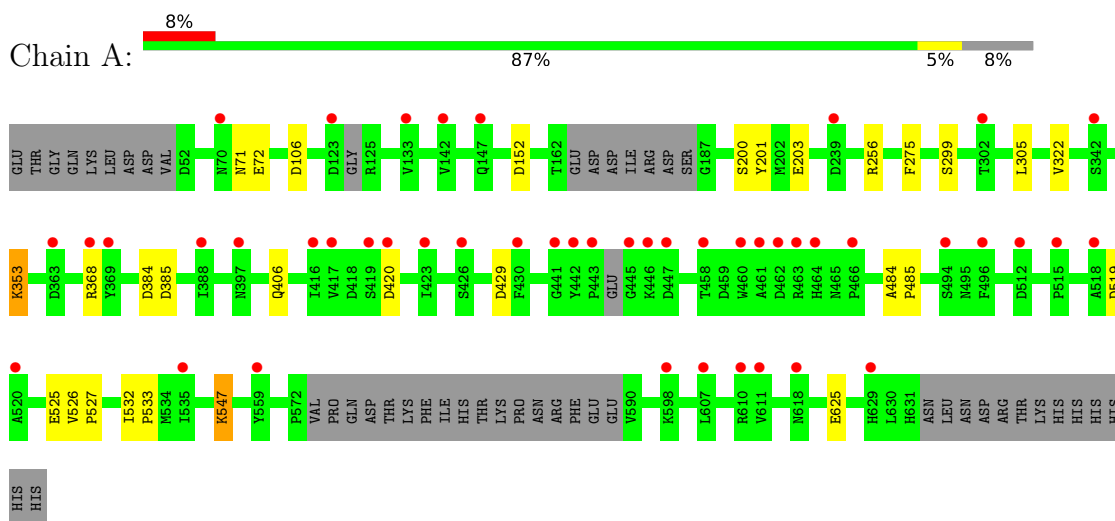
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	55	Total	O	0	0
			55	55		
4	B	54	Total	O	0	0
			54	54		

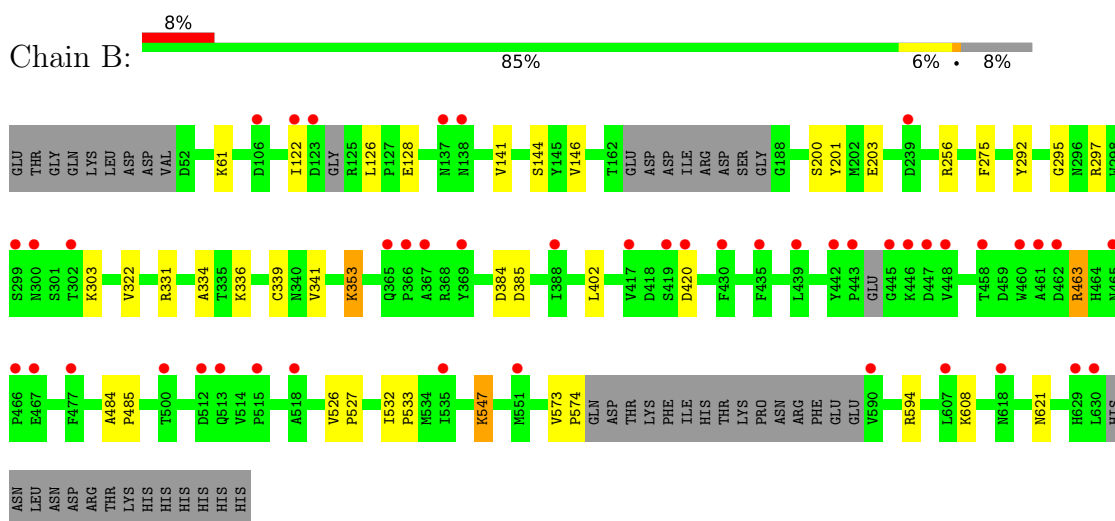
3 Residue-property plots [i](#)

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: Neuroligin-1,Neuroligin-1



• Molecule 1: Neuroligin-1,Neuroligin-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.57Å 118.05Å 214.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	51.70 – 2.55 51.70 – 2.55	Depositor EDS
% Data completeness (in resolution range)	96.7 (51.70-2.55) 87.9 (51.70-2.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.21 (at 2.55Å)	Xtriage
Refinement program	PHENIX (dev_2044)	Depositor
R, R_{free}	0.236 , 0.278 0.240 , 0.282	Depositor DCC
R_{free} test set	2008 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	40.8	Xtriage
Anisotropy	0.444	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8583	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 53.04 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.4623e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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, PGE

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	0.29	0/4316	0.61	0/5893
1	B	0.30	0/4331	0.63	0/5913
All	All	0.30	0/8647	0.62	0/11806

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	4199	0	4017	14	1
1	B	4207	0	4053	18	1
2	A	14	0	13	0	0
2	B	14	0	13	0	0
3	A	20	0	28	1	0
3	B	20	0	28	1	0
4	A	55	0	0	2	0
4	B	54	0	0	2	0
All	All	8583	0	8152	31	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 2.

All (31) 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:331:ARG:NH1	1:B:341:VAL:O	2.19	0.74
1:A:625:GLU:OE2	1:B:463:ARG:NH2	2.29	0.65
1:A:71:ASN:O	4:A:801:HOH:O	2.15	0.64
1:A:353:LYS:NZ	3:A:702:PGE:O2	2.33	0.57
1:A:547:LYS:H	1:A:547:LYS:HD2	1.71	0.55
1:A:406:GLN:NE2	1:A:519:ASP:OD2	2.40	0.54
1:A:368:ARG:N	4:A:802:HOH:O	2.42	0.51
1:B:621:ASN:ND2	4:B:804:HOH:O	2.45	0.50
1:B:141:VAL:O	1:B:144:SER:OG	2.30	0.48
1:A:384:ASP:OD1	1:A:385:ASP:N	2.44	0.47
1:B:353:LYS:HE2	3:B:702:PGE:H2	1.97	0.46
1:B:334:ALA:O	1:B:339:CYS:N	2.38	0.46
1:A:200:SER:O	1:A:201:TYR:HB2	2.16	0.45
1:B:256:ARG:NH2	1:B:292:TYR:O	2.49	0.45
1:B:200:SER:O	1:B:201:TYR:HB2	2.17	0.45
1:B:384:ASP:OD1	1:B:385:ASP:N	2.48	0.44
1:A:525:GLU:OE1	1:A:525:GLU:N	2.46	0.44
1:A:532:ILE:N	1:A:533:PRO:CD	2.81	0.43
1:B:146:VAL:HG12	1:B:146:VAL:O	2.18	0.43
1:A:484:ALA:HB3	1:A:485:PRO:HD3	2.00	0.43
1:B:126:LEU:O	1:B:128:GLU:N	2.53	0.42
1:A:71:ASN:OD1	1:A:72:GLU:N	2.52	0.42
1:B:295:GLY:N	1:B:303:LYS:O	2.32	0.42
1:B:532:ILE:N	1:B:533:PRO:CD	2.82	0.42
1:B:484:ALA:HB3	1:B:485:PRO:HD3	2.02	0.42
1:B:526:VAL:N	1:B:527:PRO:HD2	2.35	0.42
1:A:256[B]:ARG:HG2	1:A:305:LEU:HD11	2.02	0.42
1:A:526:VAL:N	1:A:527:PRO:HD2	2.35	0.41
1:B:573:VAL:HG23	1:B:574:PRO:HA	2.03	0.41
1:B:61:LYS:NZ	4:B:810:HOH:O	2.54	0.41
1:B:547:LYS:H	1:B:547:LYS:HG2	1.72	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)
1:A:152:ASP:OD2	1:B:297:ARG:NH2[4_655]	2.11	0.09

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	528/585 (90%)	500 (95%)	28 (5%)	0	100	100
1	B	529/585 (90%)	502 (95%)	27 (5%)	0	100	100
All	All	1057/1170 (90%)	1002 (95%)	55 (5%)	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	451/503 (90%)	442 (98%)	9 (2%)	48	67
1	B	455/503 (90%)	443 (97%)	12 (3%)	40	59
All	All	906/1006 (90%)	885 (98%)	21 (2%)	44	62

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	106	ASP
1	A	203	GLU
1	A	275	PHE
1	A	299	SER
1	A	322	VAL
1	A	353	LYS
1	A	420	ASP
1	A	429	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	547	LYS
1	B	122	ILE
1	B	203	GLU
1	B	275	PHE
1	B	322	VAL
1	B	336	LYS
1	B	353	LYS
1	B	402	LEU
1	B	420	ASP
1	B	463	ARG
1	B	547	LYS
1	B	594	ARG
1	B	608	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	GLN
1	A	197	HIS
1	A	226	ASN
1	A	618	ASN
1	B	79	GLN
1	B	226	ASN
1	B	285	ASN
1	B	544	ASN
1	B	621	ASN
1	B	629	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

6 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	$\# Z > 2$
3	PGE	A	703	-	9,9,9	0.35	0	8,8,8	0.21	0
2	NAG	A	701	1	14,14,15	0.20	0	17,19,21	0.53	0
3	PGE	A	702	-	9,9,9	0.35	0	8,8,8	0.20	0
3	PGE	B	702	-	9,9,9	0.35	0	8,8,8	0.45	0
3	PGE	B	703	-	9,9,9	0.33	0	8,8,8	0.31	0
2	NAG	B	701	1	14,14,15	0.25	0	17,19,21	0.34	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
3	PGE	A	703	-	-	2/7/7/7	-
2	NAG	A	701	1	-	2/6/23/26	0/1/1/1
3	PGE	A	702	-	-	2/7/7/7	-
3	PGE	B	702	-	-	4/7/7/7	-
3	PGE	B	703	-	-	2/7/7/7	-
2	NAG	B	701	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	NAG	O5-C5-C6-O6
2	A	701	NAG	C4-C5-C6-O6
3	A	703	PGE	O1-C1-C2-O2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
3	B	702	PGE	O2-C3-C4-O3
3	B	703	PGE	O3-C5-C6-O4
3	A	703	PGE	O2-C3-C4-O3
3	A	702	PGE	O2-C3-C4-O3
3	A	702	PGE	C6-C5-O3-C4
3	B	703	PGE	C6-C5-O3-C4
3	B	702	PGE	C1-C2-O2-C3
3	B	702	PGE	C4-C3-O2-C2
3	B	702	PGE	C3-C4-O3-C5

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	702	PGE	1	0
3	B	702	PGE	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	537/585 (91%)	0.69	47 (8%) 15 15	20, 50, 75, 112	1 (0%)
1	B	537/585 (91%)	0.68	46 (8%) 16 15	24, 47, 72, 107	2 (0%)
All	All	1074/1170 (91%)	0.69	93 (8%) 16 15	20, 49, 74, 112	3 (0%)

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	442	TYR	5.3
1	B	442	TYR	5.2
1	B	419	SER	4.8
1	B	443	PRO	3.8
1	B	445	GLY	3.6
1	A	445	GLY	3.4
1	A	443	PRO	3.4
1	B	629	HIS	3.3
1	A	461	ALA	3.3
1	A	629	HIS	3.1
1	B	515	PRO	3.0
1	B	535	ILE	3.0
1	A	423	ILE	2.9
1	B	430	PHE	2.9
1	A	419	SER	2.9
1	B	365	GLN	2.9
1	B	618	ASN	2.9
1	A	430	PHE	2.9
1	B	461	ALA	2.8
1	B	366	PRO	2.8
1	B	462	ASP	2.8
1	A	420	ASP	2.8
1	B	369	TYR	2.7
1	A	518	ALA	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	147	GLN	2.7
1	A	142	VAL	2.7
1	B	467	GLU	2.7
1	A	512	ASP	2.7
1	B	435	PHE	2.7
1	A	123	ASP	2.6
1	A	466	PRO	2.6
1	A	515	PRO	2.6
1	B	137	ASN	2.6
1	B	417	VAL	2.6
1	A	607	LEU	2.6
1	A	70	ASN	2.5
1	A	302	THR	2.5
1	B	466	PRO	2.5
1	B	123	ASP	2.5
1	A	446	LYS	2.5
1	A	417	VAL	2.5
1	B	500	THR	2.5
1	B	367	ALA	2.5
1	B	106	ASP	2.4
1	B	299	SER	2.4
1	B	477	PHE	2.4
1	A	133	VAL	2.4
1	B	138[A]	ASN	2.4
1	B	122	ILE	2.4
1	A	239	ASP	2.4
1	B	607	LEU	2.4
1	A	618	ASN	2.4
1	B	239	ASP	2.3
1	B	512	ASP	2.3
1	A	458	THR	2.3
1	A	494	SER	2.3
1	B	439	LEU	2.3
1	A	611	VAL	2.3
1	A	363	ASP	2.3
1	A	463	ARG	2.3
1	B	302	THR	2.2
1	A	426	SER	2.2
1	B	446	LYS	2.2
1	B	465	ASN	2.2
1	B	458	THR	2.2
1	A	441	GLY	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	342	SER	2.2
1	A	535	ILE	2.2
1	B	447	ASP	2.2
1	A	368	ARG	2.2
1	B	300	ASN	2.2
1	A	598	LYS	2.2
1	B	630	LEU	2.2
1	A	496	PHE	2.2
1	B	590	VAL	2.2
1	A	520	ALA	2.2
1	A	388	ILE	2.2
1	A	462	ASP	2.1
1	A	460	TRP	2.1
1	B	388	ILE	2.1
1	A	369	TYR	2.1
1	A	447	ASP	2.1
1	B	460	TRP	2.1
1	A	397	ASN	2.1
1	B	420	ASP	2.1
1	B	513	GLN	2.1
1	B	448	VAL	2.1
1	B	518	ALA	2.0
1	A	416	ILE	2.0
1	A	610	ARG	2.0
1	B	551	MET	2.0
1	A	464	HIS	2.0
1	A	559	TYR	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(Å ²)	Q<0.9
2	NAG	A	701	14/15	0.64	0.18	62,73,83,85	0
3	PGE	A	703	10/10	0.77	0.18	46,59,74,80	0
2	NAG	B	701	14/15	0.80	0.14	59,70,77,82	0
3	PGE	B	703	10/10	0.82	0.16	44,51,56,58	0
3	PGE	A	702	10/10	0.87	0.12	44,48,65,65	0
3	PGE	B	702	10/10	0.92	0.10	42,54,58,63	0

6.5 Other polymers [i](#)

There are no such residues in this entry.