



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:27 PM UTC

PDB ID : 5GV3 / pdb_00005gv3
Title : Crystal structure of the membrane-distal domain of mouse lysosome-associated membrane protein 2 (LAMP-2)
Authors : Tomabechi, Y.; Ehara, H.; Kukimoto-Niino, M.; Shirouzu, M.
Deposited on : 2016-09-01
Resolution : 2.10 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

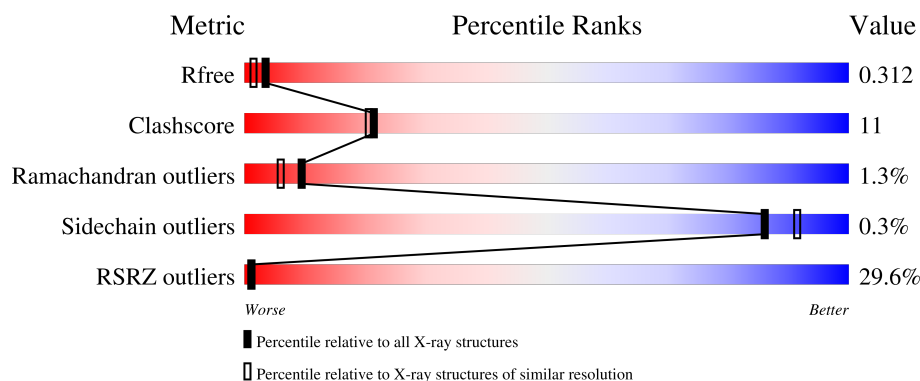
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	169	24% 76% 17% .. 5%
1	B	169	33% 72% 24% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2773 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 Lysosome-associated membrane glycoprotein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	161	Total	C	N	O	S	0	0	0
			1261	802	205	248	6			
1	B	163	Total	C	N	O	S	0	0	0
			1275	809	209	251	6			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	25	ASP	-	expression tag	UNP P17047
A	190	GLY	-	expression tag	UNP P17047
A	191	SER	-	expression tag	UNP P17047
A	192	SER	-	expression tag	UNP P17047
A	193	LEU	-	expression tag	UNP P17047
B	25	ASP	-	expression tag	UNP P17047
B	190	GLY	-	expression tag	UNP P17047
B	191	SER	-	expression tag	UNP P17047
B	192	SER	-	expression tag	UNP P17047
B	193	LEU	-	expression tag	UNP P17047

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	9	Total	Zn	3	0
			9	9		
2	B	6	Total	Zn	2	0
			6	6		

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

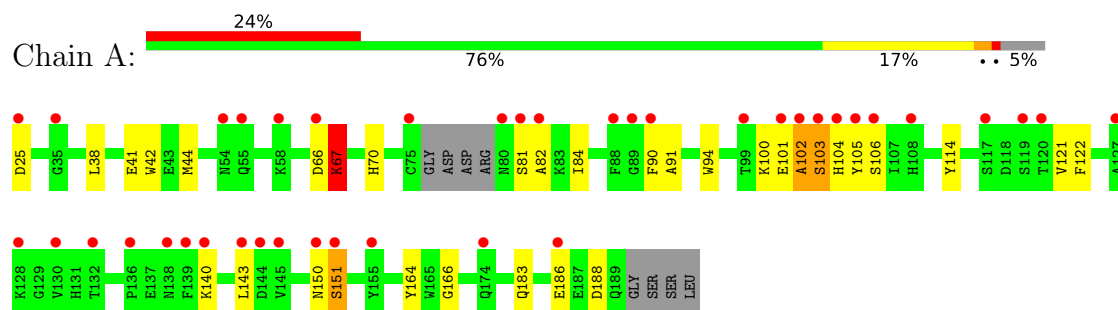
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	50	Total	O	0	0
			50	50		
4	B	32	Total	O	0	0
			32	32		

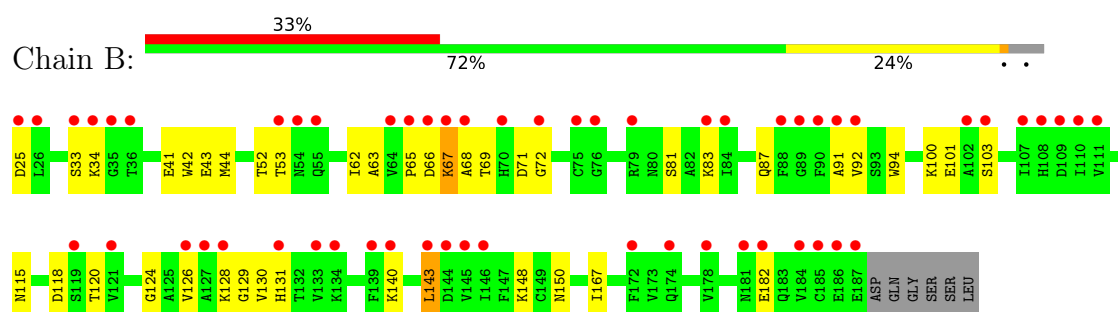
3 Residue-property plots [i](#)

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: Lysosome-associated membrane glycoprotein 2



- Molecule 1: Lysosome-associated membrane glycoprotein 2



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	59.61Å 103.93Å 129.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.49 – 2.10 40.49 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.7 (40.49-2.10) 98.7 (40.49-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.9_1690	Depositor
R, R_{free}	0.269 , 0.307 0.274 , 0.312	Depositor DCC
R_{free} test set	1982 reflections (8.27%)	wwPDB-VP
Wilson B-factor (Å ²)	45.4	Xtriage
Anisotropy	0.642	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 59.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.020 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.035 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2773	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.12% 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: ZN, 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	0.54	0/1289	0.95	4/1757 (0.2%)
1	B	0.57	1/1304 (0.1%)	0.90	3/1778 (0.2%)
All	All	0.55	1/2593 (0.0%)	0.93	7/3535 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	43	GLU	CB-CG	7.32	1.74	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	102	ALA	N-CA-C	-7.08	104.22	112.92
1	B	167	ILE	N-CA-C	5.60	116.80	108.23
1	B	68	ALA	N-CA-C	-5.56	102.66	110.50
1	A	90	PHE	N-CA-C	-5.44	106.81	113.50
1	B	43	GLU	CA-CB-CG	5.43	124.96	114.10
1	A	151	SER	N-CA-C	5.05	117.19	110.53
1	A	67	LYS	N-CA-C	-5.01	100.13	110.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	67	LYS	Peptide

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	1261	0	1210	28	3
1	B	1275	0	1224	28	3
2	A	9	0	0	0	0
2	B	6	0	0	0	1
3	A	70	0	65	1	0
3	B	70	0	65	2	0
4	A	50	0	0	3	1
4	B	32	0	0	3	1
All	All	2773	0	2564	57	5

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 11.

All (57) 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:42:TRP:HE1	1:A:44:MET:HE3	1.14	1.13
1:B:42:TRP:HE1	1:B:44:MET:HE3	1.22	1.05
1:B:91:ALA:HB1	1:B:115:ASN:HB3	1.47	0.97
1:B:69:THR:HG23	1:B:87:GLN:OE1	1.71	0.90
1:A:42:TRP:NE1	1:A:44:MET:HE3	1.90	0.87
1:B:44:MET:SD	4:B:331:HOH:O	2.36	0.82
1:B:100:LYS:HE2	1:B:143:LEU:HD21	1.63	0.79
1:B:101:GLU:OE2	1:B:140:LYS:NZ	2.14	0.79
1:A:100:LYS:HE3	1:A:143:LEU:HD21	1.68	0.75
1:A:183:GLN:NE2	4:A:303:HOH:O	2.20	0.73
1:A:104:HIS:HB3	1:A:143:LEU:HD13	1.71	0.71
1:A:151:SER:OG	1:A:188:ASP:OD1	2.07	0.71
1:A:102:ALA:HB3	1:A:140:LYS:HZ3	1.61	0.66
1:A:44:MET:HE2	1:A:164:TYR:CE1	2.33	0.64
1:B:150:ASN:OD1	4:B:304:HOH:O	2.15	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:42:TRP:HE1	1:B:44:MET:CE	2.04	0.61
1:A:44:MET:HE1	1:A:94:TRP:CZ2	2.36	0.61
1:A:44:MET:HE1	1:A:94:TRP:CE2	2.35	0.60
1:B:42:TRP:NE1	1:B:44:MET:HE3	2.06	0.59
1:B:71:ASP:O	1:B:87:GLN:NE2	2.32	0.59
1:A:150:ASN:HA	1:A:166:GLY:HA2	1.84	0.59
1:A:186:GLU:OE1	4:A:302:HOH:O	2.17	0.59
1:A:103:SER:C	1:A:140:LYS:HE3	2.30	0.57
3:A:210:NAG:H3	3:A:210:NAG:H83	1.84	0.57
1:B:66:ASP:OD1	1:B:67:LYS:HG2	2.05	0.57
1:A:106:SER:HB3	1:A:140:LYS:NZ	2.21	0.56
1:A:102:ALA:HB3	1:A:140:LYS:NZ	2.21	0.56
1:A:106:SER:HB3	1:A:140:LYS:HZ2	1.73	0.53
1:B:81:SER:OG	3:B:209:NAG:H83	2.09	0.52
1:A:44:MET:HE2	1:A:164:TYR:HE1	1.73	0.52
1:B:53:THR:N	1:B:124:GLY:O	2.42	0.51
1:B:103:SER:O	1:B:143:LEU:HB2	2.12	0.50
1:A:66:ASP:O	1:A:67:LYS:HG2	2.11	0.50
1:B:25:ASP:N	4:B:307:HOH:O	2.44	0.49
1:B:44:MET:HE1	1:B:94:TRP:CE2	2.48	0.49
1:B:148:LYS:NZ	1:B:182:GLU:CD	2.52	0.49
1:A:121:VAL:HG13	1:A:122:PHE:CD2	2.47	0.49
1:A:70:HIS:HE1	4:A:339:HOH:O	1.97	0.48
1:B:33:SER:HA	1:B:34:LYS:HA	1.58	0.47
1:B:72:GLY:O	1:B:83:LYS:HE3	2.14	0.47
1:B:62:ILE:HD13	1:B:92:VAL:HG21	1.97	0.46
1:B:118:ASP:OD1	1:B:120:THR:OG1	2.24	0.46
1:A:44:MET:HE2	1:A:164:TYR:CD1	2.50	0.45
1:B:63:ALA:O	1:B:65:PRO:HD3	2.17	0.45
1:A:100:LYS:HG2	1:A:101:GLU:HG3	1.98	0.44
1:B:130:VAL:HG21	3:B:208:NAG:H82	1.99	0.44
1:A:91:ALA:O	1:A:114:TYR:HA	2.19	0.43
1:A:38:LEU:HD21	1:A:84:ILE:HD11	2.00	0.43
1:A:140:LYS:HD2	1:A:140:LYS:HA	1.61	0.42
1:B:44:MET:HE1	1:B:94:TRP:CD2	2.54	0.42
1:B:52:THR:C	1:B:126:VAL:HG23	2.43	0.42
1:B:66:ASP:HA	1:B:67:LYS:HA	1.55	0.42
1:A:25:ASP:N	1:A:25:ASP:OD1	2.51	0.42
1:A:101:GLU:N	1:A:104:HIS:O	2.39	0.42
1:B:128:LYS:HA	1:B:129:GLY:HA2	1.73	0.41
1:A:105:TYR:HD2	1:A:143:LEU:HD12	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:148:LYS:HZ1	1:B:182:GLU:CD	2.01	0.40

All (5) 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:41:GLU:OE2	2:B:206:ZN:ZN[4_566]	1.63	0.57
1:A:150:ASN:ND2	1:B:41:GLU:OE2[4_566]	2.07	0.13
4:A:335:HOH:O	4:B:312:HOH:O[4_566]	2.10	0.10
1:A:150:ASN:OD1	1:B:25:ASP:OD2[4_566]	2.16	0.04
1:B:101:GLU:OE2	1:B:131:HIS:ND1[8_566]	2.19	0.01

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	157/169 (93%)	146 (93%)	7 (4%)	4 (2%)	4	1
1	B	161/169 (95%)	147 (91%)	14 (9%)	0	100	100
All	All	318/338 (94%)	293 (92%)	21 (7%)	4 (1%)	9	6

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	103	SER
1	A	67	LYS
1	A	82	ALA
1	A	81	SER

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	144/150 (96%)	144 (100%)	0	100	100
1	B	145/150 (97%)	144 (99%)	1 (1%)	76	83
All	All	289/300 (96%)	288 (100%)	1 (0%)	86	91

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	143	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	168	HIS
1	B	156	ASN
1	B	183	GLN

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

Of 25 ligands modelled in this entry, 15 are monoatomic - leaving 10 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
3	NAG	A	214	1	14,14,15	0.39	0	17,19,21	0.44	0
3	NAG	A	210	1	14,14,15	0.59	0	17,19,21	1.31	1 (5%)
3	NAG	B	207	1	14,14,15	0.51	0	17,19,21	0.58	0
3	NAG	A	212	1	14,14,15	0.25	0	17,19,21	0.67	0
3	NAG	B	211	1	14,14,15	0.46	0	17,19,21	0.39	0
3	NAG	A	211	1	14,14,15	0.33	0	17,19,21	0.47	0
3	NAG	B	210	1	14,14,15	0.18	0	17,19,21	0.51	0
3	NAG	B	209	1	14,14,15	0.46	0	17,19,21	0.81	1 (5%)
3	NAG	A	213	1	14,14,15	0.59	0	17,19,21	0.47	0
3	NAG	B	208	1	14,14,15	0.31	0	17,19,21	0.65	1 (5%)

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	NAG	A	214	1	-	1/6/23/26	0/1/1/1
3	NAG	A	210	1	-	6/6/23/26	0/1/1/1
3	NAG	B	207	1	-	2/6/23/26	0/1/1/1
3	NAG	A	212	1	-	2/6/23/26	0/1/1/1
3	NAG	B	211	1	-	0/6/23/26	0/1/1/1
3	NAG	A	211	1	-	0/6/23/26	0/1/1/1
3	NAG	B	210	1	-	0/6/23/26	0/1/1/1
3	NAG	B	209	1	-	3/6/23/26	0/1/1/1
3	NAG	A	213	1	-	0/6/23/26	0/1/1/1
3	NAG	B	208	1	-	2/6/23/26	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(°)
3	A	210	NAG	C2-N2-C7	4.38	128.76	122.90
3	B	209	NAG	C1-C2-N2	2.27	114.02	110.43
3	B	208	NAG	C1-O5-C5	2.22	115.17	112.19

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	209	NAG	C1-C2-N2-C7
3	B	208	NAG	C4-C5-C6-O6
3	B	207	NAG	O5-C5-C6-O6
3	B	208	NAG	O5-C5-C6-O6
3	B	209	NAG	C4-C5-C6-O6
3	B	207	NAG	C4-C5-C6-O6
3	A	210	NAG	O5-C5-C6-O6
3	B	209	NAG	O5-C5-C6-O6
3	A	210	NAG	C8-C7-N2-C2
3	A	210	NAG	O7-C7-N2-C2
3	A	210	NAG	C4-C5-C6-O6
3	A	212	NAG	C4-C5-C6-O6
3	A	212	NAG	O5-C5-C6-O6
3	A	214	NAG	O5-C5-C6-O6
3	A	210	NAG	C1-C2-N2-C7
3	A	210	NAG	C3-C2-N2-C7

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	210	NAG	1	0
3	B	209	NAG	1	0
3	B	208	NAG	1	0

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	161/169 (95%)	1.57	40 (24%)	2 2	37, 51, 72, 92	0
1	B	163/169 (96%)	1.79	56 (34%)	1 1	39, 54, 79, 98	0
All	All	324/338 (95%)	1.68	96 (29%)	1 1	37, 53, 78, 98	0

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	66	ASP	7.0
1	B	25	ASP	6.8
1	A	82	ALA	6.6
1	A	25	ASP	6.3
1	A	81	SER	6.1
1	B	90	PHE	6.1
1	B	88	PHE	5.8
1	A	90	PHE	5.7
1	A	102	ALA	5.6
1	A	103	SER	5.6
1	A	104	HIS	5.4
1	A	140	LYS	5.2
1	B	143	LEU	4.9
1	A	66	ASP	4.8
1	B	26	LEU	4.7
1	B	91	ALA	4.7
1	B	103	SER	4.6
1	B	33	SER	4.3
1	B	67	LYS	4.2
1	B	108	HIS	4.1
1	B	54	ASN	4.1
1	A	150	ASN	4.0
1	A	101	GLU	3.9
1	B	68	ALA	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	89	GLY	3.7
1	A	143	LEU	3.6
1	B	185	CYS	3.6
1	A	108	HIS	3.5
1	B	102	ALA	3.5
1	B	55	GLN	3.5
1	B	186	GLU	3.4
1	B	140	LYS	3.4
1	B	134	LYS	3.2
1	B	139	PHE	3.2
1	B	131	HIS	3.2
1	B	181	ASN	3.2
1	B	65	PRO	3.0
1	B	126	VAL	3.0
1	A	75	CYS	3.0
1	A	144	ASP	2.9
1	B	182	GLU	2.9
1	A	151	SER	2.9
1	B	92	VAL	2.8
1	B	70	HIS	2.8
1	A	35	GLY	2.8
1	A	120	THR	2.8
1	B	121	VAL	2.8
1	B	35	GLY	2.7
1	B	109	ASP	2.7
1	B	128	LYS	2.7
1	A	88	PHE	2.7
1	A	139	PHE	2.7
1	A	145	VAL	2.7
1	B	184	VAL	2.7
1	B	36	THR	2.6
1	B	144	ASP	2.6
1	B	187	GLU	2.5
1	A	136	PRO	2.5
1	A	55	GLN	2.5
1	B	111	VAL	2.5
1	A	99	THR	2.4
1	A	106	SER	2.4
1	B	127	ALA	2.4
1	B	178	VAL	2.4
1	A	128	LYS	2.4
1	A	54	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	64	VAL	2.4
1	B	133	VAL	2.4
1	A	186	GLU	2.3
1	B	174	GLN	2.3
1	A	132	THR	2.3
1	B	75	CYS	2.3
1	A	80	ASN	2.3
1	B	79	ARG	2.3
1	B	83	LYS	2.3
1	A	119	SER	2.3
1	B	76	GLY	2.2
1	A	174	GLN	2.2
1	A	58	LYS	2.2
1	B	53	THR	2.2
1	B	84	ILE	2.2
1	B	72	GLY	2.2
1	A	105	TYR	2.2
1	B	34	LYS	2.2
1	A	127	ALA	2.2
1	B	145	VAL	2.1
1	A	138	ASN	2.1
1	A	155	TYR	2.1
1	B	146	ILE	2.1
1	A	89	GLY	2.1
1	B	110	ILE	2.1
1	A	130	VAL	2.1
1	B	107	ILE	2.1
1	A	117	SER	2.1
1	B	172	PHE	2.0
1	B	119	SER	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 oligosaccharides 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. 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	ZN	A	206	1/1	0.42	0.24	64,64,64,64	1
2	ZN	B	204	1/1	0.49	0.18	93,93,93,93	1
3	NAG	A	213	14/15	0.49	0.17	70,86,92,93	0
2	ZN	A	204	1/1	-	-	68,68,68,68	1
2	ZN	A	205	1/1	-	-	80,80,80,80	1
2	ZN	A	207	1/1	0.52	0.17	102,102,102,102	1
2	ZN	A	208	1/1	0.57	0.19	69,69,69,69	1
3	NAG	A	212	14/15	0.62	0.17	73,89,104,106	0
2	ZN	A	209	1/1	-	-	79,79,79,79	1
2	ZN	A	203	1/1	0.62	0.16	90,90,90,90	1
3	NAG	B	209	14/15	0.63	0.16	60,70,82,85	0
3	NAG	B	208	14/15	0.68	0.14	68,79,84,85	0
3	NAG	B	207	14/15	0.77	0.16	57,71,82,83	0
2	ZN	B	205	1/1	-	-	90,90,90,90	1
2	ZN	B	206	1/1	-	-	60,60,60,60	1
3	NAG	B	211	14/15	0.83	0.14	43,63,68,77	0
3	NAG	A	210	14/15	0.84	0.14	46,54,64,74	0
3	NAG	A	211	14/15	0.88	0.10	40,52,67,73	0
3	NAG	A	214	14/15	0.88	0.10	43,53,66,69	0
2	ZN	B	203	1/1	0.89	0.09	52,52,52,52	1
3	NAG	B	210	14/15	0.90	0.10	42,50,58,62	0
2	ZN	A	201	1/1	0.95	0.09	45,45,45,45	0
2	ZN	A	202	1/1	0.96	0.05	49,49,49,49	1
2	ZN	B	202	1/1	0.96	0.11	48,48,48,48	0
2	ZN	B	201	1/1	0.97	0.12	38,38,38,38	0

6.5 Other polymers

There are no such residues in this entry.