



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2023 – 01:27 PM EDT

PDB ID : 2R83
Title : Crystal structure analysis of human synaptotagmin 1 C2A-C2B
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Deposited on : 2007-09-10
Resolution : 2.70 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.35
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35

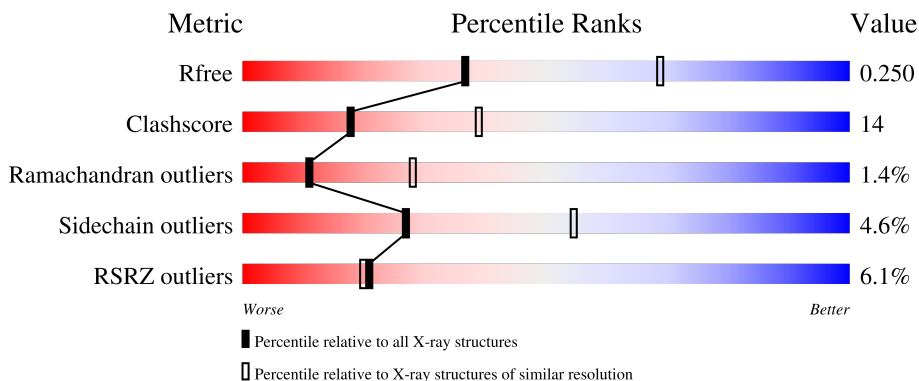
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	284	 6% 72% 23% ..
1	B	284	 6% 72% 24% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4563 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 Synaptotagmin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	279	Total	C	N	O	S	0	0	0
			2256	1455	375	418	8			
1	B	279	Total	C	N	O	S	0	0	0
			2256	1455	375	418	8			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	138	GLY	-	expression tag	UNP P21579
A	139	SER	-	expression tag	UNP P21579
A	220	ALA	GLY	conflict	UNP P21579
B	138	GLY	-	expression tag	UNP P21579
B	139	SER	-	expression tag	UNP P21579
B	220	ALA	GLY	conflict	UNP P21579

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cl		
2	A	6	Total	Cl	0	0
			6	6		
2	B	2	Total	Cl	0	0
			2	2		

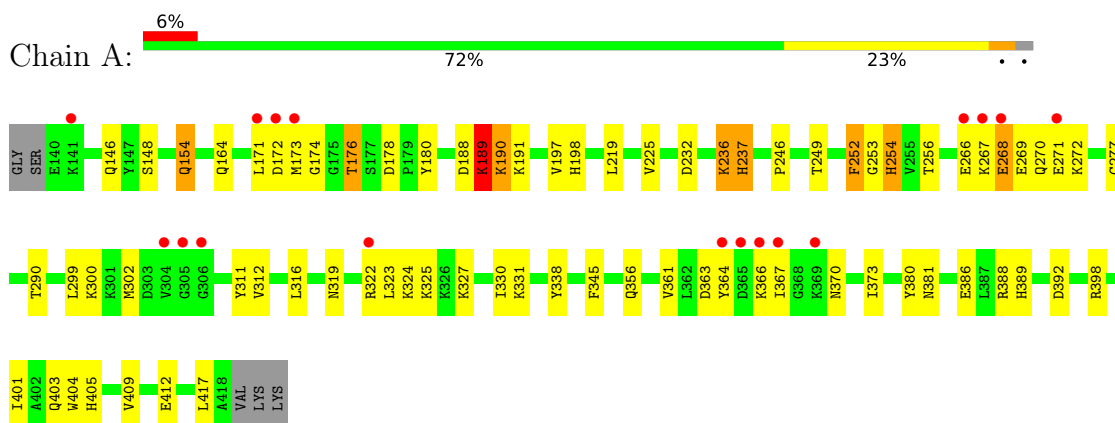
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	19	Total	O	0	0
			19	19		
3	B	24	Total	O	0	0
			24	24		

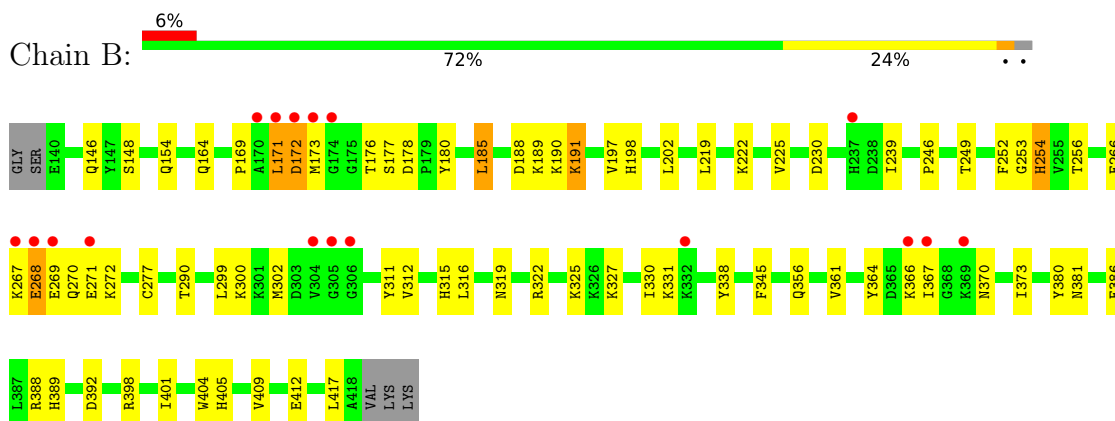
3 Residue-property plots [i](#)

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: Synaptotagmin-1



- Molecule 1: Synaptotagmin-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	82.37Å 86.31Å 147.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.31 – 2.70 46.31 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.2 (46.31-2.70) 99.2 (46.31-2.70)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 2.69Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.230 , 0.253 0.228 , 0.250	Depositor DCC
R_{free} test set	2921 reflections (9.97%)	wwPDB-VP
Wilson B-factor (Å ²)	48.3	Xtrriage
Anisotropy	0.411	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	0.033 for k,h,-l	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4563	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.35% 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: CL

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	0.39	0/2305	0.64	0/3111
1	B	0.40	0/2305	0.63	0/3111
All	All	0.40	0/4610	0.63	0/6222

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 [i](#)

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	2256	0	2279	59	0
1	B	2256	0	2279	64	0
2	A	6	0	0	0	0
2	B	2	0	0	0	0
3	A	19	0	0	2	0
3	B	24	0	0	0	0
All	All	4563	0	4558	123	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 14.

All (123) 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:299:LEU:HD22	1:A:373:ILE:HD11	1.50	0.91
1:B:299:LEU:HD22	1:B:373:ILE:HD11	1.52	0.89
1:B:169:PRO:HG2	1:B:171:LEU:HD22	1.60	0.82
1:A:266:GLU:HG3	1:A:267:LYS:H	1.42	0.82
1:B:266:GLU:HG3	1:B:267:LYS:H	1.44	0.82
1:A:191:LYS:HA	1:A:412:GLU:OE2	1.79	0.81
1:A:271:GLU:O	1:A:271:GLU:HG3	1.81	0.80
1:B:271:GLU:OE2	1:B:409:VAL:HG13	1.82	0.80
1:A:389:HIS:CD2	1:A:405:HIS:HE1	2.03	0.77
1:B:389:HIS:CD2	1:B:405:HIS:HE1	2.03	0.77
1:B:271:GLU:HG3	1:B:271:GLU:O	1.82	0.77
1:B:169:PRO:HG2	1:B:171:LEU:CD2	2.16	0.76
1:A:271:GLU:OE2	1:A:409:VAL:HG13	1.91	0.70
1:B:191:LYS:HA	1:B:412:GLU:OE2	1.91	0.70
1:B:146:GLN:HE21	1:B:164:GLN:HG2	1.57	0.70
1:B:169:PRO:O	1:B:171:LEU:HD23	1.91	0.70
1:A:299:LEU:HD22	1:A:373:ILE:CD1	2.23	0.69
1:A:146:GLN:HE21	1:A:164:GLN:HG2	1.57	0.67
1:B:409:VAL:HG23	1:B:412:GLU:HG2	1.77	0.67
1:B:239:ILE:HD13	1:B:268:GLU:HG2	1.76	0.66
1:A:268:GLU:HA	1:A:268:GLU:OE1	1.96	0.66
1:B:299:LEU:HD22	1:B:373:ILE:CD1	2.24	0.65
1:A:409:VAL:HG23	1:A:412:GLU:HG2	1.79	0.64
1:B:266:GLU:HG3	1:B:267:LYS:N	2.13	0.64
1:A:266:GLU:HG3	1:A:267:LYS:N	2.12	0.62
1:B:268:GLU:HA	1:B:268:GLU:OE1	1.98	0.62
1:B:171:LEU:HG	1:B:172:ASP:OD1	2.01	0.61
1:B:185:LEU:HD22	1:B:222:LYS:HB3	1.83	0.61
1:B:176:THR:HG22	1:B:202:LEU:HG	1.82	0.60
1:B:389:HIS:CD2	1:B:405:HIS:CE1	2.88	0.60
1:A:171:LEU:HD21	3:A:436:HOH:O	2.00	0.60
1:B:189:LYS:HE2	1:B:269:GLU:HG3	1.84	0.59
1:A:389:HIS:CD2	1:A:405:HIS:CE1	2.89	0.59
1:B:409:VAL:CG2	1:B:412:GLU:HG2	2.31	0.59
1:B:322:ARG:HD2	1:B:325:LYS:NZ	2.19	0.58
1:A:409:VAL:CG2	1:A:412:GLU:HG2	2.33	0.58
1:A:322:ARG:HD2	1:A:325:LYS:NZ	2.19	0.57
1:A:311:TYR:CD2	1:A:327:LYS:HE3	2.40	0.57
1:B:311:TYR:CD2	1:B:327:LYS:HE3	2.41	0.56
1:A:389:HIS:HD2	1:A:405:HIS:HE1	1.53	0.55
1:A:312:VAL:HG22	1:A:361:VAL:HG22	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:380:TYR:O	1:B:381:ASN:HB2	2.06	0.55
1:B:312:VAL:HG22	1:B:361:VAL:HG22	1.89	0.55
1:B:389:HIS:HD2	1:B:405:HIS:CE1	2.25	0.54
1:A:232:ASP:OD1	1:A:232:ASP:C	2.46	0.54
1:A:386:GLU:OE1	1:A:405:HIS:HD2	1.91	0.54
1:A:380:TYR:O	1:A:381:ASN:HB2	2.06	0.54
1:B:188:ASP:O	1:B:189:LYS:HB3	2.07	0.54
1:A:389:HIS:HD2	1:A:405:HIS:CE1	2.25	0.53
1:A:277:CYS:HB2	1:A:404:TRP:CD2	2.44	0.53
1:B:386:GLU:OE1	1:B:405:HIS:HD2	1.92	0.53
1:B:190:LYS:HG2	1:B:271:GLU:OE2	2.09	0.52
1:B:271:GLU:OE2	1:B:409:VAL:CG1	2.56	0.52
1:B:277:CYS:HB2	1:B:404:TRP:CD2	2.44	0.52
1:A:225:VAL:HG11	1:A:268:GLU:OE2	2.10	0.51
1:A:236:LYS:O	1:A:237:HIS:HB2	2.10	0.51
1:B:225:VAL:HG11	1:B:268:GLU:OE2	2.11	0.51
1:B:277:CYS:HB2	1:B:404:TRP:CE3	2.46	0.51
1:A:189:LYS:HG2	1:A:269:GLU:HG3	1.93	0.50
1:A:277:CYS:HB2	1:A:404:TRP:CE3	2.46	0.50
1:B:386:GLU:N	1:B:386:GLU:OE2	2.44	0.50
1:B:398:ARG:NH2	1:B:398:ARG:HB2	2.27	0.49
1:A:331:LYS:HE2	1:A:338:TYR:O	2.13	0.48
1:A:236:LYS:O	1:A:237:HIS:CB	2.62	0.48
1:A:403:GLN:NE2	3:A:423:HOH:O	2.47	0.48
1:B:271:GLU:CD	1:B:409:VAL:HG13	2.33	0.47
1:A:154:GLN:HE21	1:A:154:GLN:HB3	1.55	0.47
1:A:398:ARG:NH2	1:A:398:ARG:HB2	2.29	0.47
1:B:180:TYR:HB3	1:B:197:VAL:HA	1.96	0.47
1:A:146:GLN:NE2	1:A:164:GLN:HG2	2.29	0.47
1:B:331:LYS:HE2	1:B:338:TYR:O	2.15	0.47
1:B:389:HIS:HE1	1:B:401:ILE:O	1.98	0.47
1:A:180:TYR:HB3	1:A:197:VAL:HA	1.97	0.46
1:A:271:GLU:CD	1:A:409:VAL:HG13	2.35	0.46
1:A:311:TYR:CE2	1:A:327:LYS:HE3	2.51	0.46
1:B:172:ASP:C	1:B:173:MET:HG2	2.36	0.46
1:B:148:SER:HA	1:B:256:THR:O	2.16	0.46
1:A:367:ILE:O	1:A:367:ILE:HG12	2.16	0.46
1:B:246:PRO:HB2	1:B:249:THR:HG23	1.98	0.46
1:B:300:LYS:HE2	1:B:302:MET:HE1	1.98	0.46
1:A:148:SER:HA	1:A:256:THR:O	2.16	0.46
1:A:188:ASP:C	1:A:190:LYS:H	2.19	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:190:LYS:HE2	1:A:271:GLU:OE2	2.15	0.46
1:A:386:GLU:OE2	1:A:386:GLU:N	2.49	0.45
1:B:176:THR:HB	1:B:177:SER:H	1.56	0.45
1:A:171:LEU:HD22	1:A:176:THR:O	2.16	0.45
1:B:367:ILE:HG12	1:B:367:ILE:O	2.15	0.45
1:B:188:ASP:C	1:B:190:LYS:H	2.20	0.45
1:B:172:ASP:O	1:B:173:MET:CB	2.65	0.45
1:B:356:GLN:NE2	1:B:417:LEU:HD22	2.32	0.45
1:B:311:TYR:CE2	1:B:327:LYS:HE3	2.53	0.44
1:B:146:GLN:NE2	1:B:164:GLN:HG2	2.29	0.44
1:A:246:PRO:HB2	1:A:249:THR:HG23	2.00	0.44
1:B:185:LEU:CD2	1:B:222:LYS:HB3	2.46	0.44
1:B:252:PHE:O	1:B:253:GLY:C	2.55	0.44
1:A:330:ILE:HD13	1:A:364:TYR:CG	2.53	0.44
1:B:172:ASP:OD1	1:B:230:ASP:OD2	2.36	0.44
1:A:389:HIS:HE1	1:A:401:ILE:O	2.00	0.43
1:A:271:GLU:OE2	1:A:409:VAL:CG1	2.64	0.43
1:A:322:ARG:HD2	1:A:325:LYS:HZ1	1.82	0.43
1:A:356:GLN:NE2	1:A:417:LEU:HD22	2.34	0.43
1:B:388:ARG:HD2	1:B:392:ASP:OD2	2.19	0.43
1:A:300:LYS:HG2	1:A:302:MET:HE3	2.00	0.43
1:B:316:LEU:HD22	1:B:345:PHE:CD2	2.53	0.42
1:A:364:TYR:CE2	1:A:366:LYS:HG2	2.55	0.42
1:B:330:ILE:HD13	1:B:364:TYR:CG	2.54	0.42
1:A:370:ASN:N	1:A:370:ASN:HD22	2.17	0.42
1:B:300:LYS:HG2	1:B:302:MET:HE3	2.02	0.42
1:B:239:ILE:CD1	1:B:268:GLU:HG2	2.47	0.42
1:B:364:TYR:CE2	1:B:366:LYS:HG2	2.54	0.42
1:A:300:LYS:HE2	1:A:302:MET:HE1	2.00	0.42
1:A:302:MET:HG3	1:A:363:ASP:HB2	2.02	0.42
1:B:300:LYS:CD	1:B:302:MET:HE1	2.50	0.42
1:B:370:ASN:HD22	1:B:370:ASN:N	2.17	0.41
1:A:316:LEU:HD22	1:A:345:PHE:CD2	2.55	0.41
1:A:172:ASP:OD2	1:A:172:ASP:N	2.47	0.41
1:A:252:PHE:O	1:A:253:GLY:C	2.57	0.41
1:A:178:ASP:HA	1:A:198:HIS:O	2.20	0.41
1:A:388:ARG:HD2	1:A:392:ASP:OD2	2.19	0.41
1:B:315:HIS:CG	1:B:322:ARG:NH1	2.89	0.41
1:A:323:LEU:O	1:A:324:LYS:HG2	2.21	0.40
1:B:178:ASP:HA	1:B:198:HIS:O	2.21	0.40
1:B:364:TYR:CZ	1:B:366:LYS:HG2	2.55	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	277/284 (98%)	250 (90%)	21 (8%)	6 (2%)	6	17
1	B	277/284 (98%)	252 (91%)	23 (8%)	2 (1%)	22	46
All	All	554/568 (98%)	502 (91%)	44 (8%)	8 (1%)	11	28

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	236	LYS
1	A	237	HIS
1	B	172	ASP
1	A	254	HIS
1	B	254	HIS
1	A	189	LYS
1	A	252	PHE
1	A	174	GLY

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	249/253 (98%)	237 (95%)	12 (5%)	25	53
1	B	249/253 (98%)	238 (96%)	11 (4%)	28	56

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	498/506 (98%)	475 (95%)	23 (5%)	27	54

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	154	GLN
1	A	173	MET
1	A	176	THR
1	A	189	LYS
1	A	190	LYS
1	A	219	LEU
1	A	254	HIS
1	A	268	GLU
1	A	270	GLN
1	A	272	LYS
1	A	290	THR
1	A	319	ASN
1	B	154	GLN
1	B	171	LEU
1	B	185	LEU
1	B	191	LYS
1	B	219	LEU
1	B	254	HIS
1	B	268	GLU
1	B	270	GLN
1	B	272	LYS
1	B	290	THR
1	B	319	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	146	GLN
1	A	154	GLN
1	A	157	GLN
1	A	164	GLN
1	A	207	ASN
1	A	319	ASN
1	A	340	ASN
1	A	353	GLN
1	A	356	GLN

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Mol	Chain	Res	Type
1	A	370	ASN
1	A	381	ASN
1	A	389	HIS
1	A	396	ASN
1	A	403	GLN
1	A	405	HIS
1	A	408	GLN
1	B	146	GLN
1	B	154	GLN
1	B	157	GLN
1	B	164	GLN
1	B	207	ASN
1	B	319	ASN
1	B	340	ASN
1	B	353	GLN
1	B	356	GLN
1	B	370	ASN
1	B	381	ASN
1	B	389	HIS
1	B	403	GLN
1	B	405	HIS
1	B	408	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	279/284 (98%)	0.17	17 (6%) 21 20	22, 44, 77, 97	0
1	B	279/284 (98%)	0.21	17 (6%) 21 20	21, 43, 81, 97	0
All	All	558/568 (98%)	0.19	34 (6%) 21 20	21, 44, 81, 97	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	171	LEU	6.8
1	B	172	ASP	5.5
1	A	367	ILE	4.3
1	A	271	GLU	4.3
1	B	306	GLY	4.3
1	B	174	GLY	4.2
1	B	367	ILE	4.0
1	A	266	GLU	3.8
1	A	173	MET	3.7
1	A	172	ASP	3.6
1	A	268	GLU	3.4
1	B	271	GLU	3.2
1	A	364	TYR	2.9
1	A	304	VAL	2.8
1	A	366	LYS	2.7
1	B	268	GLU	2.7
1	A	171	LEU	2.6
1	A	306	GLY	2.5
1	B	305	GLY	2.5
1	B	366	LYS	2.3
1	A	369	LYS	2.3
1	B	267	LYS	2.3
1	B	237	HIS	2.2
1	A	322	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	365	ASP	2.2
1	A	305	GLY	2.2
1	A	267	LYS	2.1
1	B	369	LYS	2.1
1	B	170	ALA	2.1
1	B	173	MET	2.1
1	A	141	LYS	2.0
1	B	332	LYS	2.0
1	B	269	GLU	2.0
1	B	304	VAL	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 monosaccharides 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	CL	A	108	1/1	0.88	0.16	55,55,55,55	0
2	CL	A	104	1/1	0.96	0.07	36,36,36,36	0
2	CL	B	107	1/1	0.97	0.12	45,45,45,45	0
2	CL	A	101	1/1	0.98	0.10	37,37,37,37	0
2	CL	A	105	1/1	0.99	0.09	39,39,39,39	0
2	CL	A	103	1/1	0.99	0.07	36,36,36,36	0
2	CL	B	106	1/1	0.99	0.11	34,34,34,34	0
2	CL	A	102	1/1	0.99	0.09	38,38,38,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.