



Full wwPDB NMR Structure Validation Report ⓘ

Mar 9, 2026 – 04:40 PM UTC

PDB ID : 1K5W / pdb_00001k5w
Title : THREE-DIMENSIONAL STRUCTURE OF THE SYNAPTOTAGMIN 1
C2B-DOMAIN: SYNAPTOTAGMIN 1 AS A PHOSPHOLIPID BINDING
MACHINE
Authors : Fernandez, I.; Arac, D.; Ubach, J.; Gerber, S.H.; Shin, O.; Gao, Y.; Anderson,
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Deposited on : 2001-10-12

This is a Full wwPDB NMR 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/NMRValidationReportHelp>

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
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

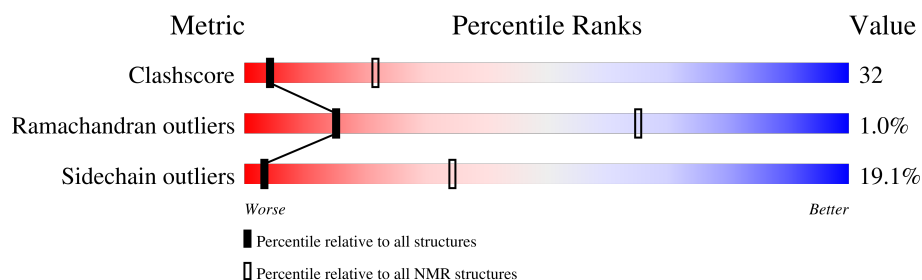
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment was not calculated.

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)	NMR archive (#Entries)
Clashscore	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and 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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	152	

2 Ensemble composition and analysis

This entry contains 20 models. Model 3 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *structure with the minimum noe energy*.

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:273-A:365, A:370-A:378, A:386-A:414 (131)	0.26	3

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

The models can be grouped into 3 clusters. No single-model clusters were found.

Cluster number	Models
1	1, 2, 3, 4, 5, 7, 9, 10, 13, 14, 17, 18, 20
2	6, 8, 11, 12, 16
3	15, 19

3 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 2410 atoms, of which 1225 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Synaptotagmin I.

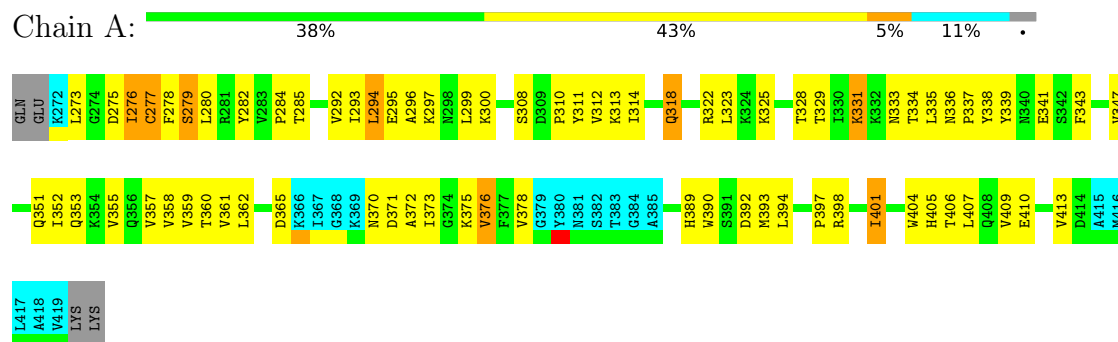
Mol	Chain	Residues	Atoms						Trace
1	A	148	Total	C	H	N	O	S	0
			2408	762	1225	202	214	5	

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	
2	A	2	Total	Ca
			2	2

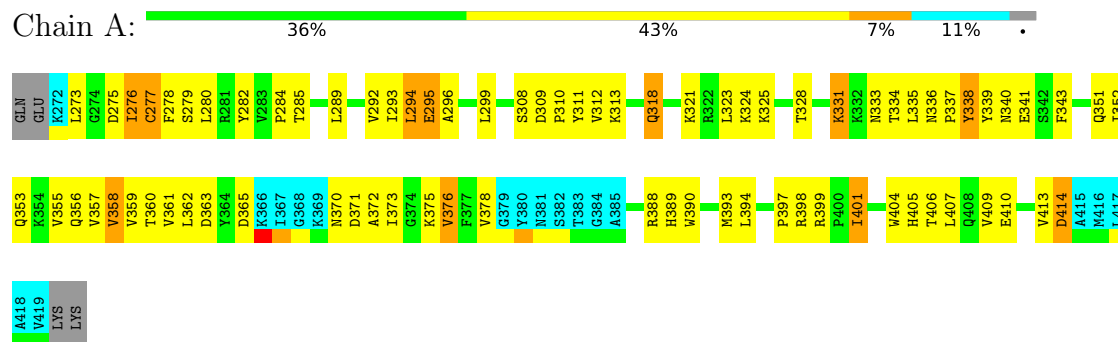
4.2.2 Score per residue for model 2

• Molecule 1: Synaptotagmin I



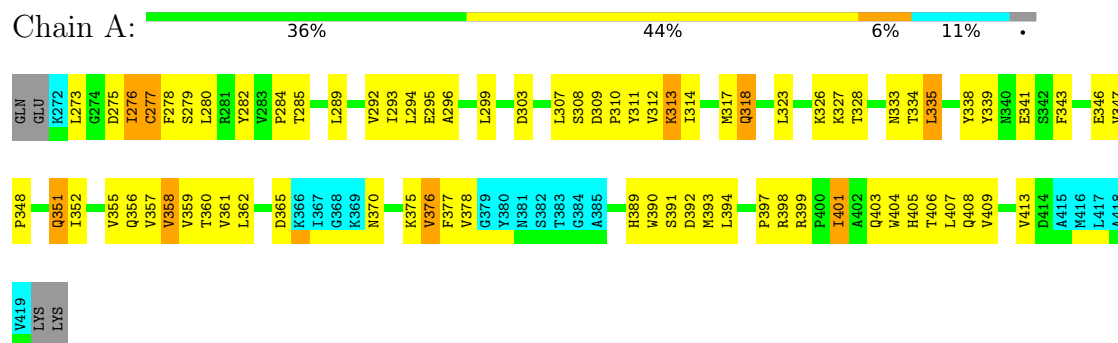
4.2.3 Score per residue for model 3 (medoid)

• Molecule 1: Synaptotagmin I



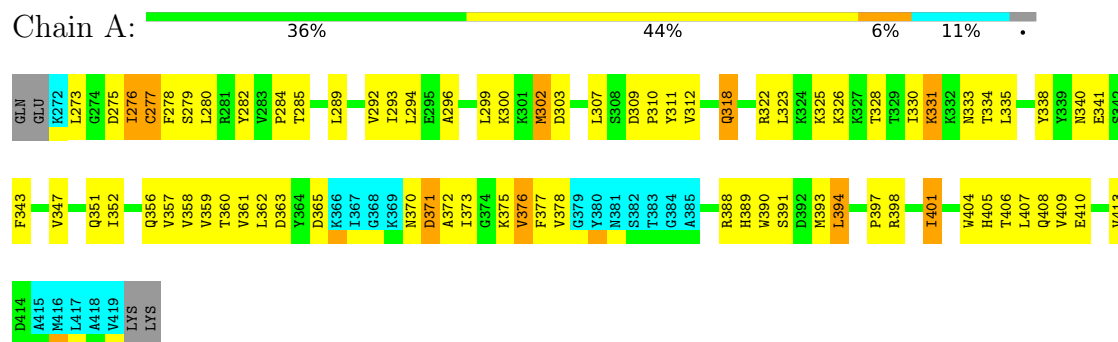
4.2.4 Score per residue for model 4

• Molecule 1: Synaptotagmin I



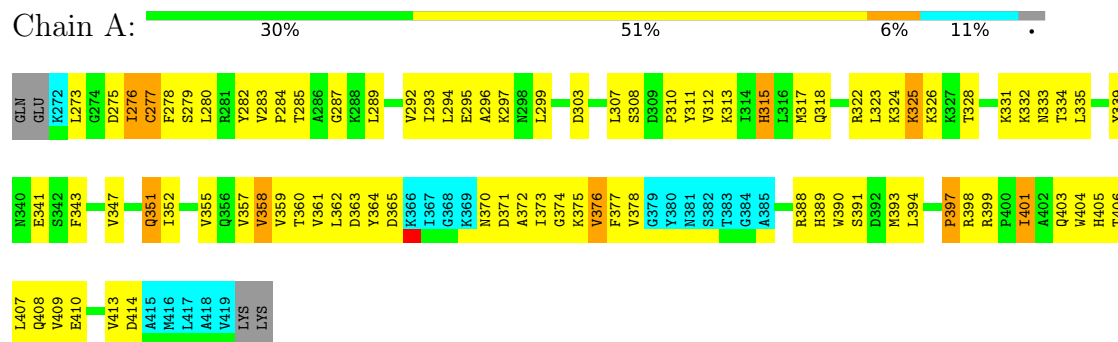
4.2.5 Score per residue for model 5

• Molecule 1: Synaptotagmin I



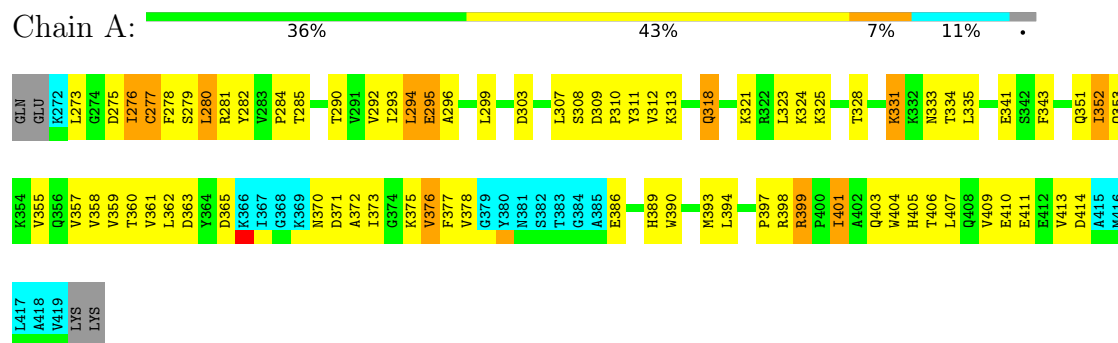
4.2.6 Score per residue for model 6

• Molecule 1: Synaptotagmin I



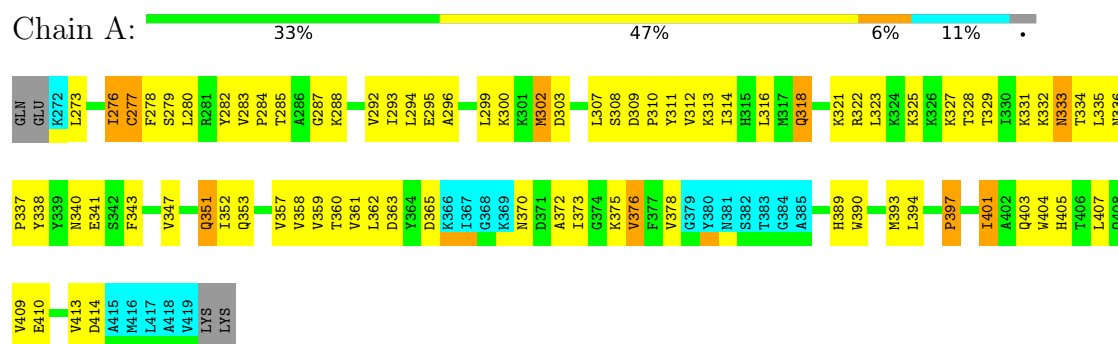
4.2.7 Score per residue for model 7

• Molecule 1: Synaptotagmin I



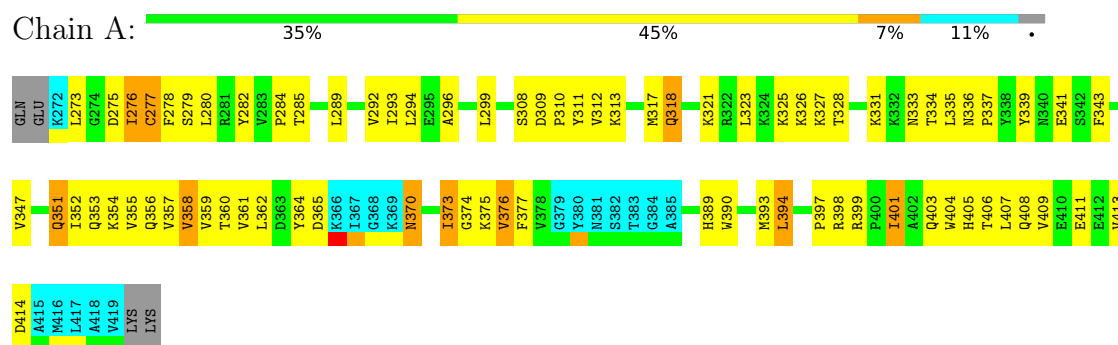
4.2.8 Score per residue for model 8

• Molecule 1: Synaptotagmin I



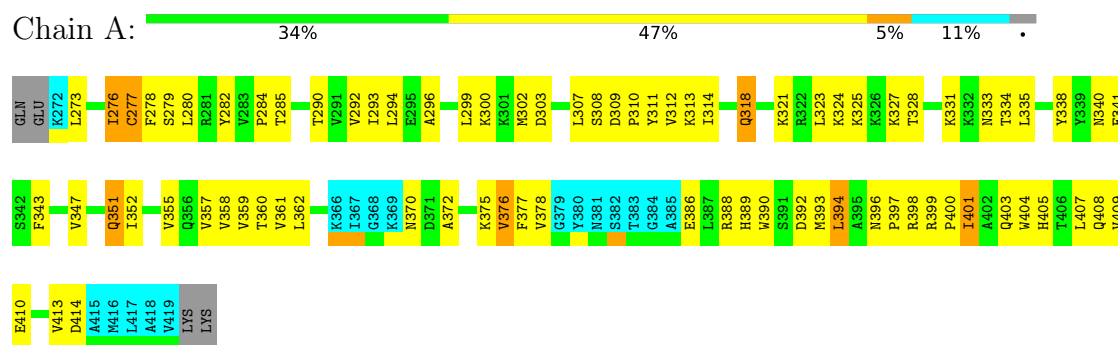
4.2.9 Score per residue for model 9

• Molecule 1: Synaptotagmin I



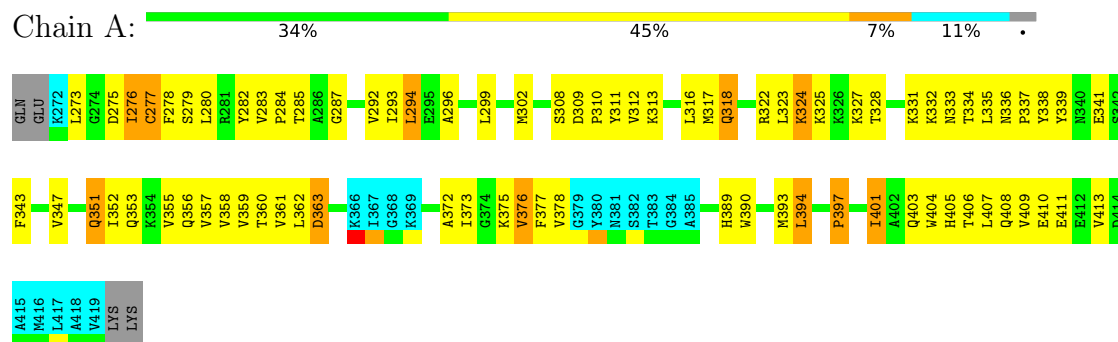
4.2.10 Score per residue for model 10

• Molecule 1: Synaptotagmin I



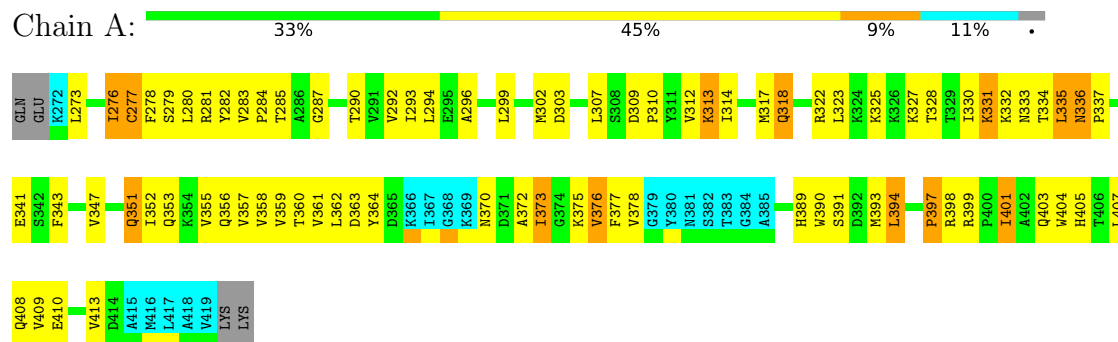
4.2.11 Score per residue for model 11

• Molecule 1: Synaptotagmin I



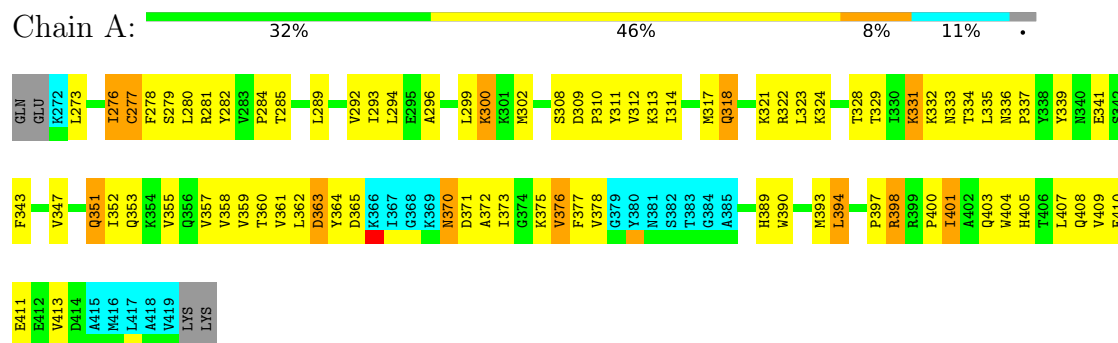
4.2.12 Score per residue for model 12

• Molecule 1: Synaptotagmin I



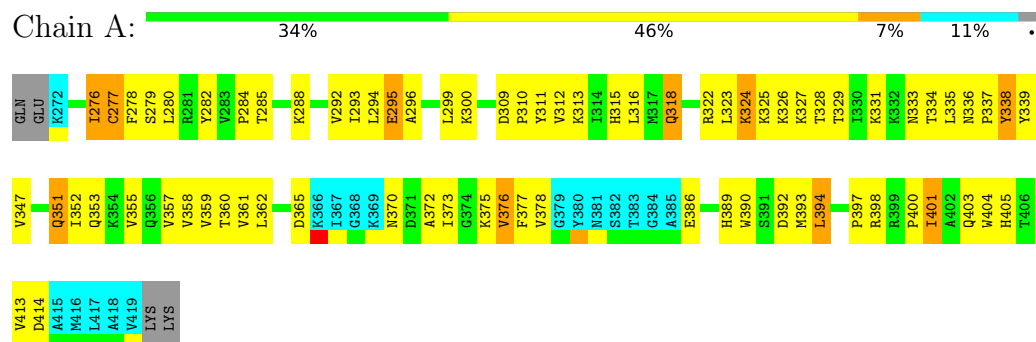
4.2.13 Score per residue for model 13

• Molecule 1: Synaptotagmin I



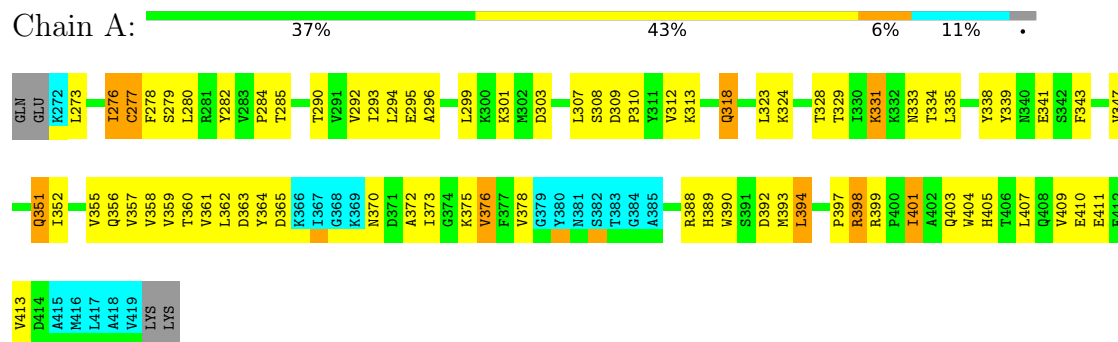
4.2.14 Score per residue for model 14

• Molecule 1: Synaptotagmin I



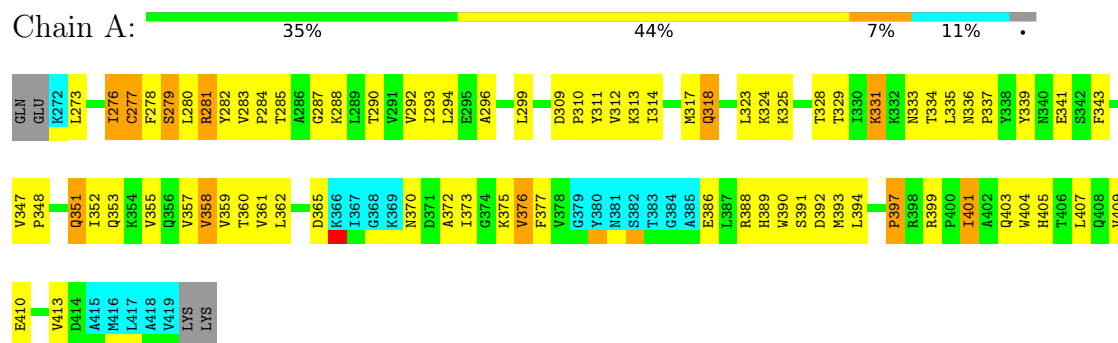
4.2.15 Score per residue for model 15

• Molecule 1: Synaptotagmin I



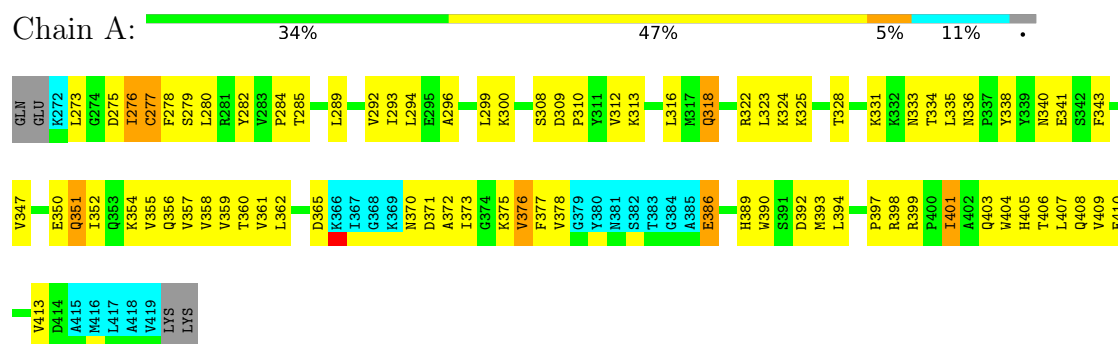
4.2.16 Score per residue for model 16

• Molecule 1: Synaptotagmin I



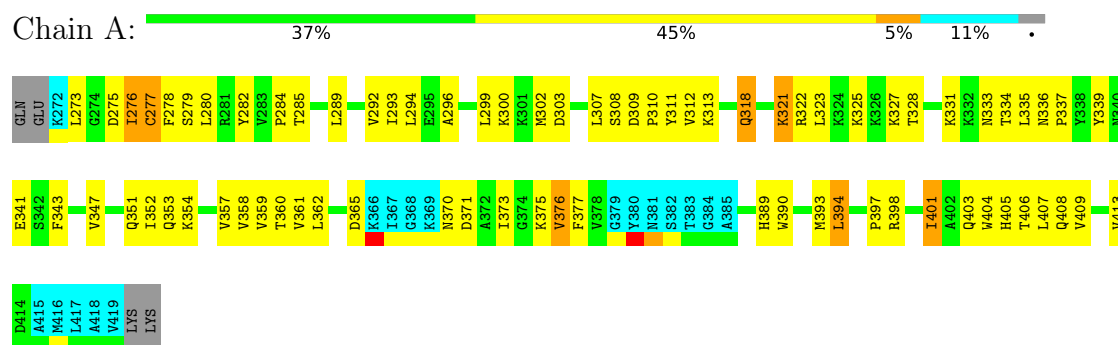
4.2.17 Score per residue for model 17

• Molecule 1: Synaptotagmin I



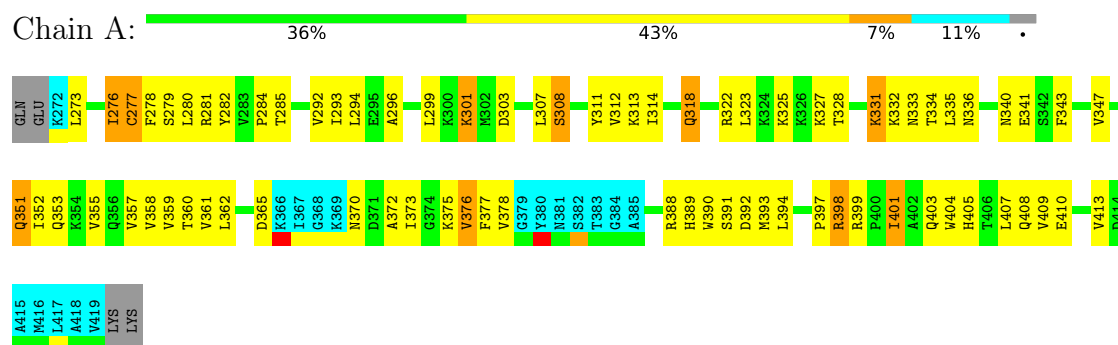
4.2.18 Score per residue for model 18

• Molecule 1: Synaptotagmin I



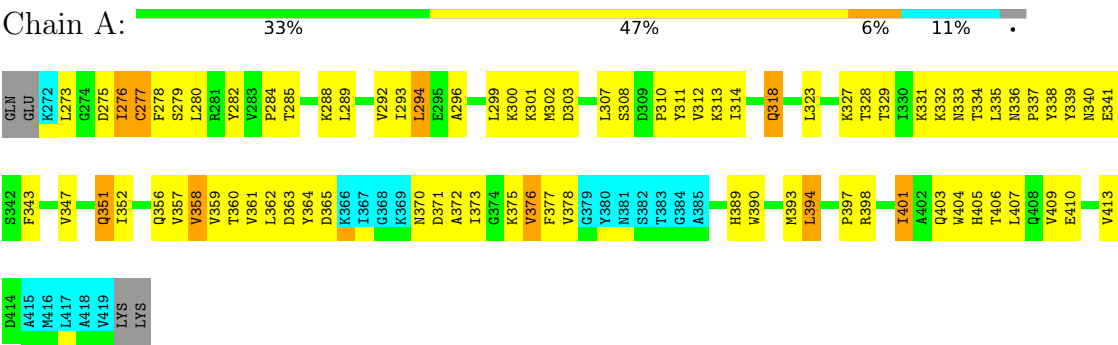
4.2.19 Score per residue for model 19

• Molecule 1: Synaptotagmin I



4.2.20 Score per residue for model 20

● Molecule 1: Synaptotagmin I



5 Refinement protocol and experimental data overview

The models were refined using the following method: *torsion angle dynamics*.

Of the 500 calculated structures, 20 were deposited, based on the following criterion: *STRUCTURES WITH THE MINIMUM NOE ENERGY*.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
CNS	refinement	0.9

No chemical shift data was provided.

6 Model quality [i](#)

6.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section:
CA

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

6.2 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 each 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 averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	1065	1095	1094	70±4
All	All	21340	21900	21880	1396

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 32.

All unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:352:ILE:HD11	1:A:393:MET:HE3	0.91	1.41	10	9
1:A:282:TYR:CD1	1:A:352:ILE:HD13	0.89	2.03	16	20
1:A:318:GLN:CB	1:A:323:LEU:HD11	0.84	2.02	20	16
1:A:312:VAL:HG22	1:A:361:VAL:HG22	0.83	1.49	14	20
1:A:360:THR:HG23	1:A:375:LYS:HB3	0.83	1.48	3	20
1:A:296:ALA:CB	1:A:299:LEU:HD11	0.81	2.06	19	16
1:A:276:ILE:HD13	1:A:277:CYS:N	0.81	1.91	18	20
1:A:280:LEU:HD12	1:A:389:HIS:CD2	0.79	2.11	16	18
1:A:409:VAL:O	1:A:413:VAL:HG23	0.78	1.77	19	20
1:A:393:MET:HE1	1:A:399:ARG:O	0.76	1.80	7	1
1:A:376:VAL:HG11	1:A:405:HIS:HB3	0.76	1.56	4	20
1:A:277:CYS:HG	1:A:404:TRP:CG	0.76	1.99	8	20

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:316:LEU:HD23	1:A:324:LYS:HD3	0.75	1.58	11	2
1:A:352:ILE:HD11	1:A:393:MET:CE	0.74	2.13	15	9
1:A:372:ALA:HB1	1:A:410:GLU:OE2	0.72	1.84	1	16
1:A:273:LEU:HD22	1:A:373:ILE:O	0.72	1.85	15	16
1:A:296:ALA:HB3	1:A:299:LEU:HD11	0.72	1.62	19	19
1:A:347:VAL:HG13	1:A:351:GLN:HB3	0.72	1.61	9	16
1:A:352:ILE:CG2	1:A:394:LEU:HD21	0.71	2.16	7	20
1:A:318:GLN:HB3	1:A:323:LEU:HD11	0.71	1.60	2	5
1:A:375:LYS:HG3	1:A:413:VAL:HG11	0.71	1.62	15	20
1:A:302:MET:HE3	1:A:373:ILE:HG12	0.69	1.63	5	1
1:A:352:ILE:HD11	1:A:393:MET:SD	0.69	2.28	3	9
1:A:376:VAL:HG22	1:A:407:LEU:HD23	0.69	1.62	20	20
1:A:299:LEU:HD13	1:A:310:PRO:CG	0.69	2.18	9	13
1:A:313:LYS:HE3	1:A:362:LEU:HD13	0.68	1.65	19	14
1:A:313:LYS:HE2	1:A:362:LEU:HD13	0.68	1.65	18	5
1:A:328:THR:HG22	1:A:343:PHE:CE1	0.68	2.24	15	19
1:A:334:THR:O	1:A:335:LEU:HD23	0.68	1.88	7	20
1:A:372:ALA:HB1	1:A:410:GLU:CD	0.67	2.14	14	7
1:A:275:ASP:OD1	1:A:406:THR:HG23	0.67	1.89	4	11
1:A:351:GLN:O	1:A:355:VAL:HG22	0.67	1.89	17	5
1:A:293:ILE:HG22	1:A:339:TYR:HB2	0.67	1.65	3	13
1:A:279:SER:OG	1:A:292:VAL:HB	0.67	1.90	8	2
1:A:296:ALA:HB1	1:A:299:LEU:HD21	0.66	1.64	18	20
1:A:318:GLN:CG	1:A:323:LEU:HD11	0.66	2.21	3	11
1:A:278:PHE:CZ	1:A:405:HIS:CD2	0.66	2.84	2	20
1:A:390:TRP:CZ3	1:A:393:MET:HE1	0.66	2.25	19	19
1:A:318:GLN:HB2	1:A:323:LEU:HD11	0.65	1.68	15	13
1:A:289:LEU:HD11	1:A:393:MET:HE1	0.64	1.69	6	9
1:A:277:CYS:HG	1:A:404:TRP:CD1	0.63	2.12	7	18
1:A:318:GLN:HG2	1:A:355:VAL:HG12	0.62	1.71	11	6
1:A:360:THR:HG23	1:A:375:LYS:CB	0.62	2.23	1	20
1:A:316:LEU:HD23	1:A:324:LYS:CD	0.62	2.25	11	2
1:A:302:MET:HE3	1:A:373:ILE:CG1	0.60	2.25	5	1
1:A:352:ILE:HG21	1:A:394:LEU:HD21	0.60	1.71	16	12
1:A:361:VAL:HG21	1:A:407:LEU:HD11	0.60	1.73	17	15
1:A:331:LYS:O	1:A:334:THR:HG22	0.60	1.96	1	16
1:A:313:LYS:CE	1:A:362:LEU:HD13	0.60	2.27	1	9
1:A:318:GLN:HG2	1:A:323:LEU:HD11	0.59	1.73	9	2
1:A:293:ILE:HG13	1:A:328:THR:HG21	0.57	1.76	20	20
1:A:278:PHE:CZ	1:A:405:HIS:CG	0.57	2.93	19	20
1:A:294:LEU:HD22	1:A:295:GLU:OE1	0.56	1.99	2	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:289:LEU:HD11	1:A:393:MET:CE	0.56	2.30	18	8
1:A:331:LYS:HB3	1:A:334:THR:HG21	0.56	1.76	2	11
1:A:328:THR:HG22	1:A:343:PHE:CE2	0.55	2.37	6	1
1:A:357:VAL:O	1:A:378:VAL:HG22	0.55	2.01	1	1
1:A:352:ILE:HG23	1:A:394:LEU:HD21	0.54	1.78	13	16
1:A:293:ILE:HG22	1:A:339:TYR:CB	0.54	2.33	9	8
1:A:314:ILE:CG2	1:A:357:VAL:HG13	0.54	2.33	4	1
1:A:321:LYS:HG3	1:A:323:LEU:HD21	0.54	1.78	18	1
1:A:318:GLN:HB3	1:A:355:VAL:HG12	0.54	1.78	9	1
1:A:276:ILE:HD13	1:A:276:ILE:C	0.53	2.28	2	20
1:A:378:VAL:HG23	1:A:378:VAL:O	0.53	2.03	8	15
1:A:284:PRO:HD3	1:A:397:PRO:HB3	0.53	1.79	16	1
1:A:397:PRO:C	1:A:398:ARG:HG2	0.52	2.28	1	15
1:A:293:ILE:HD11	1:A:314:ILE:HD11	0.52	1.80	10	1
1:A:376:VAL:HG22	1:A:407:LEU:CD2	0.52	2.35	13	5
1:A:318:GLN:HG3	1:A:323:LEU:HD11	0.52	1.80	17	3
1:A:292:VAL:HG13	1:A:341:GLU:O	0.52	2.05	4	18
1:A:289:LEU:HD21	1:A:357:VAL:CG2	0.52	2.34	13	5
1:A:375:LYS:CG	1:A:413:VAL:HG21	0.52	2.35	13	3
1:A:364:TYR:CD1	1:A:365:ASP:N	0.51	2.79	15	3
1:A:314:ILE:HG23	1:A:357:VAL:HG13	0.51	1.82	4	7
1:A:284:PRO:CG	1:A:397:PRO:CB	0.51	2.89	7	15
1:A:357:VAL:HG23	1:A:390:TRP:CZ3	0.51	2.41	20	18
1:A:377:PHE:CD1	1:A:413:VAL:HG13	0.51	2.41	6	3
1:A:347:VAL:HG11	1:A:352:ILE:HA	0.51	1.83	16	4
1:A:278:PHE:CE1	1:A:405:HIS:CD2	0.51	2.99	18	20
1:A:339:TYR:C	1:A:340:ASN:ND2	0.50	2.69	3	2
1:A:282:TYR:O	1:A:397:PRO:C	0.50	2.53	16	5
1:A:299:LEU:HD13	1:A:310:PRO:HG3	0.50	1.82	9	9
1:A:284:PRO:CG	1:A:397:PRO:HB3	0.50	2.37	17	6
1:A:273:LEU:HD12	1:A:408:GLN:C	0.50	2.32	4	11
1:A:282:TYR:CE1	1:A:352:ILE:HD13	0.50	2.42	11	7
1:A:312:VAL:CG1	1:A:359:VAL:HG13	0.49	2.36	17	17
1:A:375:LYS:HG2	1:A:413:VAL:HG21	0.49	1.82	5	14
1:A:308:SER:HB2	1:A:335:LEU:HD23	0.49	1.83	15	12
1:A:347:VAL:HG13	1:A:351:GLN:CB	0.49	2.34	6	2
1:A:352:ILE:O	1:A:390:TRP:NE1	0.49	2.45	4	14
1:A:284:PRO:CD	1:A:397:PRO:HB2	0.49	2.38	1	11
1:A:399:ARG:HG3	1:A:400:PRO:HD2	0.49	1.83	10	1
1:A:284:PRO:HD3	1:A:397:PRO:CB	0.49	2.38	16	5
1:A:329:THR:HG23	1:A:331:LYS:HE3	0.49	1.84	14	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:280:LEU:O	1:A:401:ILE:N	0.49	2.44	5	7
1:A:329:THR:HG22	1:A:341:GLU:OE2	0.48	2.08	16	1
1:A:328:THR:HG22	1:A:343:PHE:HE1	0.48	1.68	19	13
1:A:316:LEU:HD23	1:A:324:LYS:HG3	0.48	1.86	17	1
1:A:279:SER:N	1:A:292:VAL:O	0.48	2.39	2	2
1:A:390:TRP:CH2	1:A:393:MET:HE1	0.48	2.44	10	3
1:A:280:LEU:HD12	1:A:389:HIS:NE2	0.48	2.23	7	4
1:A:310:PRO:O	1:A:331:LYS:HG2	0.48	2.09	12	1
1:A:389:HIS:CD2	1:A:401:ILE:CG2	0.47	2.97	11	20
1:A:351:GLN:O	1:A:355:VAL:HG13	0.47	2.10	17	5
1:A:303:ASP:HB2	1:A:307:LEU:O	0.47	2.10	19	12
1:A:358:VAL:HG11	1:A:375:LYS:HE2	0.47	1.86	3	4
1:A:308:SER:O	1:A:334:THR:N	0.47	2.46	19	10
1:A:290:THR:HG23	1:A:343:PHE:O	0.47	2.10	15	3
1:A:378:VAL:CG1	1:A:405:HIS:CE1	0.47	2.98	13	5
1:A:293:ILE:N	1:A:341:GLU:O	0.47	2.48	8	20
1:A:311:TYR:CD1	1:A:362:LEU:O	0.47	2.68	5	16
1:A:353:GLN:HE21	1:A:394:LEU:HD22	0.47	1.69	8	2
1:A:322:ARG:O	1:A:323:LEU:HD23	0.47	2.10	14	1
1:A:328:THR:HG22	1:A:343:PHE:CZ	0.46	2.46	16	2
1:A:299:LEU:HD13	1:A:310:PRO:CB	0.46	2.40	13	7
1:A:358:VAL:HG13	1:A:377:PHE:HB3	0.46	1.88	6	4
1:A:375:LYS:CD	1:A:410:GLU:HG2	0.46	2.41	1	1
1:A:278:PHE:CZ	1:A:359:VAL:HG21	0.46	2.46	5	1
1:A:280:LEU:CD1	1:A:389:HIS:CE1	0.46	2.99	10	4
1:A:309:ASP:N	1:A:310:PRO:CD	0.45	2.79	12	16
1:A:311:TYR:O	1:A:361:VAL:HA	0.45	2.12	18	4
1:A:312:VAL:CG1	1:A:359:VAL:CG1	0.45	2.95	4	20
1:A:283:VAL:O	1:A:287:GLY:N	0.45	2.49	6	5
1:A:282:TYR:CG	1:A:352:ILE:HD13	0.45	2.43	7	1
1:A:280:LEU:HD12	1:A:389:HIS:CG	0.45	2.45	20	2
1:A:296:ALA:HB1	1:A:299:LEU:HD11	0.45	1.81	19	1
1:A:315:HIS:CD2	1:A:325:LYS:CG	0.45	3.00	6	1
1:A:300:LYS:HD2	1:A:302:MET:HE2	0.45	1.86	20	2
1:A:276:ILE:C	1:A:276:ILE:CD1	0.45	2.90	14	11
1:A:359:VAL:N	1:A:376:VAL:O	0.45	2.49	6	5
1:A:273:LEU:HD13	1:A:374:GLY:CA	0.44	2.42	6	2
1:A:364:TYR:CD2	1:A:370:ASN:ND2	0.44	2.85	9	1
1:A:313:LYS:N	1:A:360:THR:O	0.44	2.49	8	13
1:A:397:PRO:O	1:A:398:ARG:CD	0.44	2.66	2	8
1:A:338:TYR:CE2	1:A:340:ASN:HB2	0.44	2.48	5	4

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:330:ILE:HG21	1:A:364:TYR:CB	0.44	2.43	12	1
1:A:278:PHE:CD2	1:A:359:VAL:HG11	0.44	2.48	15	9
1:A:352:ILE:CG1	1:A:394:LEU:HD21	0.44	2.43	18	2
1:A:399:ARG:CG	1:A:400:PRO:HD2	0.44	2.42	10	1
1:A:301:LYS:HE3	1:A:335:LEU:HD11	0.44	1.90	20	1
1:A:318:GLN:O	1:A:318:GLN:HG3	0.44	2.11	10	1
1:A:282:TYR:CZ	1:A:352:ILE:HB	0.44	2.48	14	13
1:A:375:LYS:CG	1:A:413:VAL:HG11	0.44	2.38	15	3
1:A:278:PHE:CE2	1:A:359:VAL:HG21	0.44	2.48	5	1
1:A:284:PRO:CD	1:A:397:PRO:CB	0.43	2.96	19	6
1:A:353:GLN:NE2	1:A:394:LEU:HD13	0.43	2.29	16	1
1:A:280:LEU:O	1:A:400:PRO:HA	0.43	2.13	13	3
1:A:301:LYS:HG3	1:A:335:LEU:HD21	0.43	1.87	19	1
1:A:328:THR:HG22	1:A:343:PHE:HE2	0.43	1.69	6	1
1:A:284:PRO:HD3	1:A:397:PRO:HB2	0.43	1.91	8	4
1:A:336:ASN:N	1:A:337:PRO:HD3	0.43	2.29	8	12
1:A:302:MET:HG3	1:A:363:ASP:HB3	0.43	1.91	13	2
1:A:377:PHE:N	1:A:377:PHE:CD1	0.43	2.87	18	2
1:A:331:LYS:HB2	1:A:334:THR:CG2	0.43	2.44	12	1
1:A:375:LYS:O	1:A:413:VAL:HG21	0.42	2.14	8	1
1:A:280:LEU:HB3	1:A:393:MET:HE3	0.42	1.90	17	2
1:A:307:LEU:HD13	1:A:333:ASN:C	0.42	2.39	8	1
1:A:364:TYR:CD1	1:A:364:TYR:C	0.42	2.97	20	3
1:A:357:VAL:HG23	1:A:390:TRP:CH2	0.42	2.49	14	3
1:A:390:TRP:CZ3	1:A:393:MET:CE	0.42	3.02	2	1
1:A:347:VAL:CG1	1:A:352:ILE:N	0.42	2.82	10	2
1:A:295:GLU:CD	1:A:338:TYR:CE1	0.42	2.98	14	5
1:A:314:ILE:HG23	1:A:357:VAL:CG1	0.42	2.45	12	1
1:A:377:PHE:CE1	1:A:386:GLU:OE1	0.42	2.73	7	1
1:A:302:MET:SD	1:A:371:ASP:HB2	0.42	2.54	5	1
1:A:316:LEU:HD21	1:A:347:VAL:CG2	0.41	2.45	8	1
1:A:294:LEU:HD23	1:A:338:TYR:CE2	0.41	2.50	20	1
1:A:282:TYR:O	1:A:397:PRO:CB	0.41	2.68	11	3
1:A:358:VAL:HG11	1:A:375:LYS:NZ	0.41	2.30	1	1
1:A:355:VAL:HG23	1:A:390:TRP:CZ2	0.41	2.50	4	1
1:A:377:PHE:CD1	1:A:377:PHE:N	0.41	2.88	19	1
1:A:378:VAL:O	1:A:390:TRP:CE3	0.41	2.74	13	1
1:A:295:GLU:OE1	1:A:338:TYR:CZ	0.41	2.74	2	1
1:A:277:CYS:O	1:A:294:LEU:CB	0.41	2.69	3	1
1:A:388:ARG:O	1:A:392:ASP:N	0.41	2.51	16	1
1:A:318:GLN:NE2	1:A:354:LYS:CD	0.41	2.83	17	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:294:LEU:CD2	1:A:338:TYR:CE2	0.41	3.04	1	2
1:A:377:PHE:CZ	1:A:386:GLU:OE1	0.41	2.74	17	2
1:A:277:CYS:N	1:A:295:GLU:O	0.41	2.50	8	1
1:A:352:ILE:O	1:A:390:TRP:CZ2	0.41	2.74	10	1
1:A:282:TYR:OH	1:A:347:VAL:O	0.41	2.39	17	2
1:A:280:LEU:HD13	1:A:393:MET:HE3	0.41	1.92	9	1
1:A:364:TYR:CD2	1:A:370:ASN:OD1	0.41	2.74	13	1
1:A:338:TYR:C	1:A:339:TYR:CD1	0.41	2.99	14	1
1:A:293:ILE:O	1:A:341:GLU:N	0.41	2.54	19	1
1:A:281:ARG:O	1:A:290:THR:CB	0.41	2.69	12	2
1:A:312:VAL:O	1:A:328:THR:HG23	0.41	2.16	15	1
1:A:282:TYR:OH	1:A:348:PRO:C	0.40	2.64	16	2
1:A:311:TYR:HB3	1:A:330:ILE:HA	0.40	1.93	5	1
1:A:311:TYR:CD2	1:A:362:LEU:O	0.40	2.74	8	1
1:A:375:LYS:HD2	1:A:410:GLU:HG2	0.40	1.93	13	1
1:A:315:HIS:CD2	1:A:325:LYS:CB	0.40	3.04	6	1
1:A:314:ILE:N	1:A:314:ILE:HD12	0.40	2.31	2	1
1:A:280:LEU:HB3	1:A:393:MET:CE	0.40	2.47	6	1
1:A:280:LEU:HG	1:A:389:HIS:CE1	0.40	2.52	6	1
1:A:351:GLN:O	1:A:355:VAL:N	0.40	2.54	6	1
1:A:312:VAL:HG11	1:A:359:VAL:HG13	0.40	1.94	8	1
1:A:389:HIS:NE2	1:A:401:ILE:O	0.40	2.54	11	1
1:A:302:MET:HG3	1:A:363:ASP:CB	0.40	2.46	12	1
1:A:372:ALA:HB1	1:A:410:GLU:OE1	0.40	2.16	13	1
1:A:397:PRO:O	1:A:398:ARG:HD3	0.40	2.16	19	1
1:A:302:MET:SD	1:A:363:ASP:HB3	0.40	2.57	8	1
1:A:318:GLN:NE2	1:A:351:GLN:CG	0.40	2.84	5	1
1:A:295:GLU:HB3	1:A:338:TYR:CD1	0.40	2.52	8	1
1:A:396:ASN:C	1:A:398:ARG:H	0.40	2.25	10	1

6.3 Torsion angles ⓘ

6.3.1 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 PDB entries followed by that with respect to all NMR entries. 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	131/152 (86%)	116±2 (89±1%)	13±2 (10±2%)	1±0 (1±0%)	15	65

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	2620/3040 (86%)	2330 (89%)	264 (10%)	26 (1%)	15 65

All 3 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	333	ASN	20
1	A	397	PRO	5
1	A	398	ARG	1

6.3.2 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 PDB entries followed by that with respect to all NMR entries. 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	120/135 (89%)	97±3 (81±2%)	23±3 (19±2%)	3 34
All	All	2400/2700 (89%)	1942 (81%)	458 (19%)	3 34

All 57 unique residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	276	ILE	20
1	A	277	CYS	20
1	A	285	THR	20
1	A	294	LEU	20
1	A	318	GLN	20
1	A	358	VAL	20
1	A	376	VAL	20
1	A	401	ILE	20
1	A	370	ASN	19
1	A	279	SER	18
1	A	351	GLN	17
1	A	403	GLN	17
1	A	325	LYS	16
1	A	331	LYS	12
1	A	322	ARG	11
1	A	327	LYS	11

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Mol	Chain	Res	Type	Models (Total)
1	A	394	LEU	11
1	A	324	LYS	10
1	A	399	ARG	10
1	A	300	LYS	9
1	A	371	ASP	9
1	A	321	LYS	8
1	A	332	LYS	8
1	A	363	ASP	8
1	A	391	SER	7
1	A	392	ASP	7
1	A	414	ASP	7
1	A	281	ARG	5
1	A	329	THR	5
1	A	388	ARG	5
1	A	326	LYS	5
1	A	411	GLU	5
1	A	373	ILE	4
1	A	386	GLU	4
1	A	295	GLU	4
1	A	356	GLN	4
1	A	302	MET	4
1	A	398	ARG	4
1	A	288	LYS	4
1	A	336	ASN	3
1	A	352	ILE	2
1	A	297	LYS	2
1	A	338	TYR	2
1	A	313	LYS	2
1	A	335	LEU	2
1	A	315	HIS	2
1	A	317	MET	2
1	A	354	LYS	2
1	A	301	LYS	2
1	A	340	ASN	2
1	A	342	SER	1
1	A	375	LYS	1
1	A	346	GLU	1
1	A	323	LEU	1
1	A	280	LEU	1
1	A	350	GLU	1
1	A	308	SER	1

6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

6.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided