



wwPDB EM Map/Model Validation Report ⓘ

Apr 28, 2016 – 02:58 AM EDT

PDB ID : 5IPV
EMDB ID: : EMD-8106
Title : Cryo-EM structure of GluN1/GluN2B NMDA receptor in the DCKA/D-APV-bound conformation, state 1
Authors : Zhu, S.; Stein, A.R.; Yoshioka, C.; Lee, C.H.; Goehring, A.; McHaourab, S.H.; Gouaux, E.
Deposited on : 2016-03-10
Resolution : 9.25 Å(reported)

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.
For rigid body fitted models, validation errors reported here could stem from errors in the original structure(s) used in the fitting.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/EMValidationReportHelp>

MolProbity : 4.02b-467
Mogul : unknown
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027457

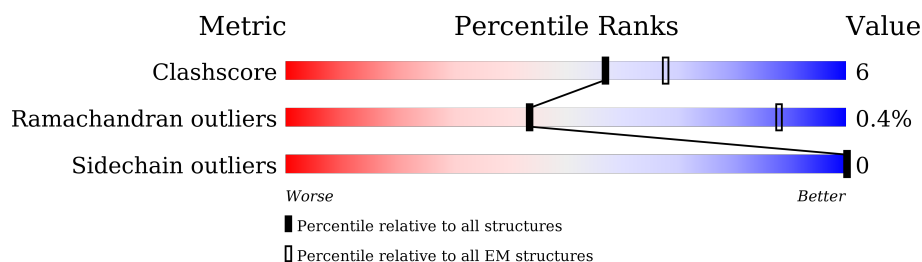
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 9.25 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)	EM structures (#Entries)
Clashscore	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	822	
1	C	822	
2	B	825	
2	D	825	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10209 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 N-methyl-D-aspartate receptor subunit NR1-8a.

Mol	Chain	Residues	Atoms				AltConf	Trace
1	A	653	Total	C	N	O	0	0
			2612	1306	653	653		
1	C	653	Total	C	N	O	0	0
			2612	1306	653	653		

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	51	PHE	LYS	engineered mutation	UNP C0KD18
A	52	PHE	ARG	engineered mutation	UNP C0KD18
A	300	GLN	ASN	engineered mutation	UNP C0KD18
A	350	GLN	ASN	engineered mutation	UNP C0KD18
A	368	ASP	ASN	engineered mutation	UNP C0KD18
A	440	ASP	ASN	engineered mutation	UNP C0KD18
A	469	ASP	ASN	engineered mutation	UNP C0KD18
A	493	ALA	LYS	engineered mutation	UNP C0KD18
A	494	ALA	LYS	engineered mutation	UNP C0KD18
A	495	ALA	GLU	engineered mutation	UNP C0KD18
A	?	-	LYS	deletion	UNP C0KD18
A	?	-	VAL	deletion	UNP C0KD18
A	?	-	ASN	deletion	UNP C0KD18
A	?	-	SER	deletion	UNP C0KD18
A	?	-	GLU	deletion	UNP C0KD18
A	?	-	GLU	deletion	UNP C0KD18
A	?	-	GLU	deletion	UNP C0KD18
A	?	-	GLU	deletion	UNP C0KD18
A	602	ARG	GLY	engineered mutation	UNP C0KD18
A	609	LEU	ILE	engineered mutation	UNP C0KD18
A	648	ARG	ASP	engineered mutation	UNP C0KD18
A	761	GLU	ASN	engineered mutation	UNP C0KD18
A	829	SER	-	expression tag	UNP C0KD18
A	830	ARG	-	expression tag	UNP C0KD18
A	831	ALA	-	expression tag	UNP C0KD18
A	832	GLU	-	expression tag	UNP C0KD18

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Chain	Residue	Modelled	Actual	Comment	Reference
A	833	ALA	-	expression tag	UNP C0KD18
A	834	LYS	-	expression tag	UNP C0KD18
A	835	ARG	-	expression tag	UNP C0KD18
A	836	MET	-	expression tag	UNP C0KD18
A	837	LYS	-	expression tag	UNP C0KD18
A	838	GLY	-	expression tag	UNP C0KD18
A	839	LEU	-	expression tag	UNP C0KD18
A	840	GLU	-	expression tag	UNP C0KD18
A	841	VAL	-	expression tag	UNP C0KD18
A	842	LEU	-	expression tag	UNP C0KD18
A	843	PHE	-	expression tag	UNP C0KD18
A	844	GLN	-	expression tag	UNP C0KD18
C	51	PHE	LYS	engineered mutation	UNP C0KD18
C	52	PHE	ARG	engineered mutation	UNP C0KD18
C	300	GLN	ASN	engineered mutation	UNP C0KD18
C	350	GLN	ASN	engineered mutation	UNP C0KD18
C	368	ASP	ASN	engineered mutation	UNP C0KD18
C	440	ASP	ASN	engineered mutation	UNP C0KD18
C	469	ASP	ASN	engineered mutation	UNP C0KD18
C	493	ALA	LYS	engineered mutation	UNP C0KD18
C	494	ALA	LYS	engineered mutation	UNP C0KD18
C	495	ALA	GLU	engineered mutation	UNP C0KD18
C	?	-	LYS	deletion	UNP C0KD18
C	?	-	VAL	deletion	UNP C0KD18
C	?	-	ASN	deletion	UNP C0KD18
C	?	-	SER	deletion	UNP C0KD18
C	?	-	GLU	deletion	UNP C0KD18
C	?	-	GLU	deletion	UNP C0KD18
C	?	-	GLU	deletion	UNP C0KD18
C	?	-	GLU	deletion	UNP C0KD18
C	602	ARG	GLY	engineered mutation	UNP C0KD18
C	609	LEU	ILE	engineered mutation	UNP C0KD18
C	648	ARG	ASP	engineered mutation	UNP C0KD18
C	761	GLU	ASN	engineered mutation	UNP C0KD18
C	829	SER	-	expression tag	UNP C0KD18
C	830	ARG	-	expression tag	UNP C0KD18
C	831	ALA	-	expression tag	UNP C0KD18
C	832	GLU	-	expression tag	UNP C0KD18
C	833	ALA	-	expression tag	UNP C0KD18
C	834	LYS	-	expression tag	UNP C0KD18
C	835	ARG	-	expression tag	UNP C0KD18
C	836	MET	-	expression tag	UNP C0KD18

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Chain	Residue	Modelled	Actual	Comment	Reference
C	837	LYS	-	expression tag	UNP C0KD18
C	838	GLY	-	expression tag	UNP C0KD18
C	839	LEU	-	expression tag	UNP C0KD18
C	840	GLU	-	expression tag	UNP C0KD18
C	841	VAL	-	expression tag	UNP C0KD18
C	842	LEU	-	expression tag	UNP C0KD18
C	843	PHE	-	expression tag	UNP C0KD18
C	844	GLN	-	expression tag	UNP C0KD18

- Molecule 2 is a protein called Ionotropic glutamate receptor subunit NR2B.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	B	623	Total	C	N	O	0	0
			2493	1247	623	623		
2	D	622	Total	C	N	O	0	0
			2492	1248	622	622		

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	20	SER	MET	engineered mutation	UNP A7XY94
B	21	ARG	GLY	engineered mutation	UNP A7XY94
B	22	ALA	CYS	engineered mutation	UNP A7XY94
B	64	GLU	ALA	engineered mutation	UNP A7XY94
B	69	GLN	ASN	engineered mutation	UNP A7XY94
B	343	ASP	ASN	engineered mutation	UNP A7XY94
B	?	-	LYS	deletion	UNP A7XY94
B	?	-	TYR	deletion	UNP A7XY94
B	?	-	TYR	deletion	UNP A7XY94
B	?	-	VAL	deletion	UNP A7XY94
B	486	VAL	THR	engineered mutation	UNP A7XY94
B	?	-	ARG	deletion	UNP A7XY94
B	?	-	CYS	deletion	UNP A7XY94
B	?	-	LEU	deletion	UNP A7XY94
B	?	-	ALA	deletion	UNP A7XY94
B	?	-	ASP	deletion	UNP A7XY94
B	?	-	GLY	deletion	UNP A7XY94
B	?	-	ARG	deletion	UNP A7XY94
B	?	-	GLU	deletion	UNP A7XY94
B	?	-	PRO	deletion	UNP A7XY94
B	?	-	GLY	deletion	UNP A7XY94
B	601	LEU	VAL	engineered mutation	UNP A7XY94

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Chain	Residue	Modelled	Actual	Comment	Reference
B	640	ARG	GLU	engineered mutation	UNP A7XY94
B	641	ARG	GLU	engineered mutation	UNP A7XY94
D	20	SER	MET	engineered mutation	UNP A7XY94
D	21	ARG	GLY	engineered mutation	UNP A7XY94
D	22	ALA	CYS	engineered mutation	UNP A7XY94
D	64	GLU	ALA	engineered mutation	UNP A7XY94
D	69	GLN	ASN	engineered mutation	UNP A7XY94
D	343	ASP	ASN	engineered mutation	UNP A7XY94
D	?	-	LYS	deletion	UNP A7XY94
D	?	-	TYR	deletion	UNP A7XY94
D	?	-	TYR	deletion	UNP A7XY94
D	?	-	VAL	deletion	UNP A7XY94
D	486	VAL	THR	engineered mutation	UNP A7XY94
D	?	-	ARG	deletion	UNP A7XY94
D	?	-	CYS	deletion	UNP A7XY94
D	?	-	LEU	deletion	UNP A7XY94
D	?	-	ALA	deletion	UNP A7XY94
D	?	-	ASP	deletion	UNP A7XY94
D	?	-	GLY	deletion	UNP A7XY94
D	?	-	ARG	deletion	UNP A7XY94
D	?	-	GLU	deletion	UNP A7XY94
D	?	-	PRO	deletion	UNP A7XY94
D	?	-	GLY	deletion	UNP A7XY94
D	601	LEU	VAL	engineered mutation	UNP A7XY94
D	640	ARG	GLU	engineered mutation	UNP A7XY94
D	641	ARG	GLU	engineered mutation	UNP A7XY94

4 Experimental information

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of images	393513	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	Not provided	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	Not provided	Depositor
Minimum defocus (nm)	Not provided	Depositor
Maximum defocus (nm)	Not provided	Depositor
Magnification	Not provided	Depositor
Image detector	Not provided	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	$\# Z > 2$	RMSZ	$\# Z > 2$
1	A	0.28	0/2609	0.70	5/3256 (0.2%)
1	C	0.30	0/2609	0.97	5/3256 (0.2%)
2	B	0.32	0/2489	0.71	4/3105 (0.1%)
2	D	0.32	0/2489	0.70	4/3107 (0.1%)
All	All	0.30	0/10196	0.78	18/12724 (0.1%)

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	C	0	1

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	C	784	ARG	O-C-N	23.06	159.59	122.70
1	C	784	ARG	C-N-CA	-20.77	69.79	121.70
1	C	487	ARG	N-CA-C	19.94	164.83	111.00
1	C	784	ARG	CA-C-N	-19.45	74.41	117.20
1	A	784	ARG	N-CA-C	11.99	143.36	111.00
2	B	534	SER	N-CA-C	-9.31	85.86	111.00
1	A	785	TYR	N-CA-C	8.63	134.31	111.00
2	D	353	MET	N-CA-C	7.96	132.50	111.00
1	A	787	GLU	N-CA-C	7.86	132.22	111.00
1	A	786	GLN	N-CA-C	-7.34	91.19	111.00
1	A	123	THR	N-CA-C	6.49	128.53	111.00
1	C	123	THR	N-CA-C	6.46	128.44	111.00
2	B	533	ARG	N-CA-C	-6.12	94.47	111.00
2	D	354	HIS	N-CA-C	-5.42	96.36	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	D	399	LEU	N-CA-C	5.34	125.43	111.00
2	B	399	LEU	N-CA-C	5.32	125.36	111.00
2	B	37	VAL	C-N-CA	5.16	133.13	122.30
2	D	37	VAL	C-N-CA	5.11	133.03	122.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	487	ARG	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	2612	0	699	30	0
1	C	2612	0	699	14	0
2	B	2493	0	685	31	0
2	D	2492	0	688	21	0
All	All	10209	0	2771	81	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 6.

All (81) 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:79:CYS:CA	2:B:74:LYS:CA	1.77	1.62
2:B:345:SER:O	2:B:353:MET:N	1.79	1.15
1:A:79:CYS:CA	2:B:74:LYS:N	2.17	1.07
2:D:409:PHE:CB	2:D:454:CYS:H	1.68	1.06
1:A:133:ILE:CA	2:B:130:ALA:CA	2.36	1.04
2:D:409:PHE:CB	2:D:454:CYS:N	2.22	1.02
1:A:308:CYS:O	2:B:71:SER:O	1.88	0.91
2:D:409:PHE:HA	2:D:453:PHE:N	1.86	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:409:PHE:HA	2:D:453:PHE:H	1.34	0.88
1:A:79:CYS:C	2:B:74:LYS:CA	2.45	0.84
1:A:308:CYS:O	2:B:72:ASP:CA	2.28	0.81
1:A:310:GLY:HA2	2:B:70:GLU:CA	2.14	0.78
2:B:50:GLU:O	2:B:53:ASP:N	2.20	0.74
2:D:50:GLU:O	2:D:53:ASP:N	2.20	0.73
1:C:660:PRO:O	1:C:664:ASN:N	2.22	0.72
1:C:464:LEU:O	1:C:468:MET:N	2.22	0.72
1:A:464:LEU:O	1:A:468:MET:N	2.23	0.72
1:C:144:TYR:O	1:C:147:GLN:N	2.21	0.71
2:B:345:SER:O	2:B:353:MET:CA	2.38	0.70
1:A:308:CYS:CA	2:B:72:ASP:CA	2.69	0.69
1:A:308:CYS:O	2:B:71:SER:C	2.30	0.69
1:C:483:GLY:HA2	1:C:497:ASN:O	1.93	0.69
1:A:483:GLY:HA2	1:A:497:ASN:O	1.93	0.69
1:A:660:PRO:O	1:A:664:ASN:N	2.22	0.68
1:A:310:GLY:H	2:B:70:GLU:C	1.97	0.67
2:D:330:LEU:O	2:D:334:LEU:N	2.30	0.64
2:B:330:LEU:O	2:B:334:LEU:N	2.30	0.64
2:D:480:GLY:HA2	2:D:488:ASN:O	1.98	0.63
1:A:308:CYS:C	2:B:72:ASP:CA	2.66	0.63
2:B:346:PHE:CA	2:B:353:MET:H	2.10	0.63
2:D:45:ILE:O	2:D:49:HIS:N	2.32	0.63
2:D:47:ASP:O	2:D:51:LYS:N	2.29	0.63
2:B:480:GLY:HA2	2:B:488:ASN:O	1.98	0.62
1:A:785:TYR:O	1:A:786:GLN:C	2.38	0.62
2:D:409:PHE:CA	2:D:454:CYS:H	2.12	0.62
2:B:345:SER:C	2:B:353:MET:H	2.03	0.61
2:B:45:ILE:O	2:B:49:HIS:N	2.32	0.61
2:D:355:PRO:O	2:D:356:LYS:C	2.39	0.61
2:B:493:GLU:O	2:B:499:ALA:N	2.31	0.60
2:D:493:GLU:O	2:D:499:ALA:N	2.31	0.60
2:B:47:ASP:O	2:B:51:LYS:N	2.29	0.60
2:D:406:GLU:O	2:D:410:VAL:O	2.20	0.59
1:C:265:GLY:HA3	1:C:381:TRP:C	2.23	0.58
1:A:150:VAL:O	1:A:154:MET:N	2.35	0.58
1:A:265:GLY:HA3	1:A:381:TRP:C	2.23	0.58
1:A:149:LEU:O	1:A:153:GLU:N	2.35	0.58
2:D:526:GLY:O	2:D:716:ALA:N	2.36	0.57
1:A:102:LEU:O	1:A:105:THR:N	2.37	0.57
2:B:526:GLY:O	2:B:716:ALA:N	2.36	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:102:LEU:O	1:C:105:THR:N	2.37	0.57
1:A:498:GLY:O	1:A:501:GLY:N	2.38	0.56
1:C:150:VAL:O	1:C:154:MET:N	2.35	0.56
1:C:498:GLY:O	1:C:501:GLY:N	2.38	0.55
1:A:535:GLY:HA2	1:A:745:ARG:CA	2.37	0.54
1:A:308:CYS:O	2:B:72:ASP:N	2.42	0.53
2:B:346:PHE:CA	2:B:353:MET:N	2.72	0.52
1:C:365:GLY:HA2	1:C:375:ASN:H	1.75	0.52
1:A:365:GLY:HA2	1:A:375:ASN:H	1.75	0.52
1:C:149:LEU:O	1:C:153:GLU:N	2.35	0.51
2:B:174:TYR:C	2:B:176:ASP:H	2.14	0.51
2:D:303:LEU:O	2:D:307:ASN:N	2.36	0.50
1:A:79:CYS:CA	2:B:74:LYS:H	2.16	0.49
2:B:303:LEU:O	2:B:307:ASN:N	2.36	0.49
2:B:351:TYR:O	2:B:352:GLN:C	2.50	0.49
1:A:310:GLY:CA	2:B:70:GLU:CA	2.87	0.47
2:D:174:TYR:C	2:D:176:ASP:H	2.17	0.47
2:B:345:SER:O	2:B:354:HIS:N	2.47	0.46
1:A:785:TYR:C	1:A:786:GLN:O	2.46	0.46
1:C:144:TYR:O	1:C:145:SER:C	2.53	0.46
2:D:409:PHE:CB	2:D:453:PHE:CA	2.94	0.46
2:D:280:TRP:C	2:D:282:TYR:H	2.20	0.45
2:D:409:PHE:HA	2:D:453:PHE:CA	2.47	0.44
2:D:409:PHE:CB	2:D:453:PHE:C	2.86	0.43
2:D:280:TRP:C	2:D:282:TYR:N	2.73	0.42
1:A:113:PHE:CA	2:B:102:ALA:CA	2.98	0.42
1:C:482:PHE:O	1:C:499:MET:N	2.53	0.41
1:A:265:GLY:HA3	1:A:381:TRP:O	2.21	0.41
1:A:482:PHE:O	1:A:499:MET:N	2.53	0.41
1:A:710:ALA:O	1:A:715:LYS:N	2.55	0.40
1:C:709:GLN:O	1:C:713:ASP:N	2.41	0.40
1:C:30:GLY:HA2	1:C:63:THR:O	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 PDB entries followed by that with respect to all EM

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	647/822 (79%)	614 (95%)	29 (4%)	4 (1%)	30	74
1	C	647/822 (79%)	617 (95%)	28 (4%)	2 (0%)	46	83
2	B	615/825 (74%)	582 (95%)	31 (5%)	2 (0%)	46	83
2	D	614/825 (74%)	578 (94%)	33 (5%)	3 (0%)	34	77
All	All	2523/3294 (77%)	2391 (95%)	121 (5%)	11 (0%)	43	80

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	141	VAL
1	A	142	PRO
1	A	143	PRO
2	B	143	PRO
2	B	408	PRO
2	D	143	PRO
2	D	410	VAL
1	A	124	ARG
1	C	124	ARG
1	C	143	PRO
2	D	407	ALA

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 PDB entries followed by that with respect to all EM 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	
2	D	1/727 (0%)	1 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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.