



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 10:07 AM GMT

PDB ID : 3EAP
Title : Crystal structure of the RhoGAP domain of ARHGAP11A
Authors : Shen, Y.; Shen, L.; Tong, Y.; Tempel, W.; MacKenzie, F.; Arrowsmith, C.H.;
Edwards, A.M.; Bountra, C.; Weigelt, J.; Bochkarev, A.; Park, H.; Structural
Genomics Consortium; Structural Genomics Consortium (SGC)
Deposited on : 2008-08-26
Resolution : 2.30 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

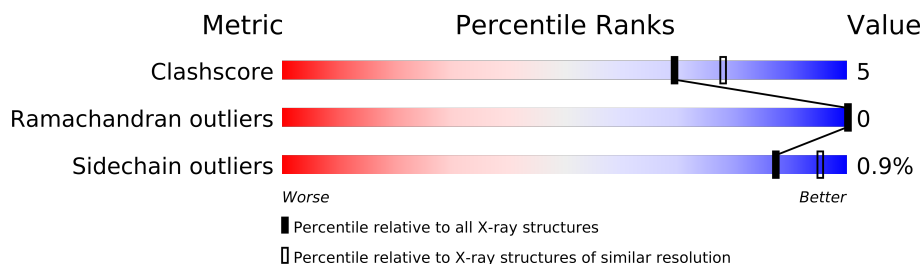
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	FAILED
Percentile statistics	:	21963
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.30 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))
Clashscore	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS failed to run properly.

Mol	Chain	Length	Quality of chain
1	A	271	
1	B	271	
1	C	271	
1	D	271	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6525 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Rho GTPase-activating protein 11A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	225	Total	C	N	O	S	0	0	0
			1665	1079	281	299	6			
1	B	213	Total	C	N	O	S	0	0	0
			1606	1041	275	286	4			
1	C	216	Total	C	N	O	S	0	0	0
			1613	1049	273	285	6			
1	D	212	Total	C	N	O	S	0	0	0
			1577	1026	266	280	5			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-17	MET	-	EXPRESSION TAG	UNP Q6P4F7
A	-16	HIS	-	EXPRESSION TAG	UNP Q6P4F7
A	-15	HIS	-	EXPRESSION TAG	UNP Q6P4F7
A	-14	HIS	-	EXPRESSION TAG	UNP Q6P4F7
A	-13	HIS	-	EXPRESSION TAG	UNP Q6P4F7
A	-12	HIS	-	EXPRESSION TAG	UNP Q6P4F7
A	-11	HIS	-	EXPRESSION TAG	UNP Q6P4F7
A	-10	SER	-	EXPRESSION TAG	UNP Q6P4F7
A	-9	SER	-	EXPRESSION TAG	UNP Q6P4F7
A	-8	GLY	-	EXPRESSION TAG	UNP Q6P4F7
A	-7	ARG	-	EXPRESSION TAG	UNP Q6P4F7
A	-6	GLU	-	EXPRESSION TAG	UNP Q6P4F7
A	-5	ASN	-	EXPRESSION TAG	UNP Q6P4F7
A	-4	LEU	-	EXPRESSION TAG	UNP Q6P4F7
A	-3	TYR	-	EXPRESSION TAG	UNP Q6P4F7
A	-2	PHE	-	EXPRESSION TAG	UNP Q6P4F7
A	-1	GLN	-	EXPRESSION TAG	UNP Q6P4F7
A	0	GLY	-	EXPRESSION TAG	UNP Q6P4F7
B	-17	MET	-	EXPRESSION TAG	UNP Q6P4F7
B	-16	HIS	-	EXPRESSION TAG	UNP Q6P4F7
B	-15	HIS	-	EXPRESSION TAG	UNP Q6P4F7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	EXPRESSION TAG	UNP Q6P4F7
B	-13	HIS	-	EXPRESSION TAG	UNP Q6P4F7
B	-12	HIS	-	EXPRESSION TAG	UNP Q6P4F7
B	-11	HIS	-	EXPRESSION TAG	UNP Q6P4F7
B	-10	SER	-	EXPRESSION TAG	UNP Q6P4F7
B	-9	SER	-	EXPRESSION TAG	UNP Q6P4F7
B	-8	GLY	-	EXPRESSION TAG	UNP Q6P4F7
B	-7	ARG	-	EXPRESSION TAG	UNP Q6P4F7
B	-6	GLU	-	EXPRESSION TAG	UNP Q6P4F7
B	-5	ASN	-	EXPRESSION TAG	UNP Q6P4F7
B	-4	LEU	-	EXPRESSION TAG	UNP Q6P4F7
B	-3	TYR	-	EXPRESSION TAG	UNP Q6P4F7
B	-2	PHE	-	EXPRESSION TAG	UNP Q6P4F7
B	-1	GLN	-	EXPRESSION TAG	UNP Q6P4F7
B	0	GLY	-	EXPRESSION TAG	UNP Q6P4F7
C	-17	MET	-	EXPRESSION TAG	UNP Q6P4F7
C	-16	HIS	-	EXPRESSION TAG	UNP Q6P4F7
C	-15	HIS	-	EXPRESSION TAG	UNP Q6P4F7
C	-14	HIS	-	EXPRESSION TAG	UNP Q6P4F7
C	-13	HIS	-	EXPRESSION TAG	UNP Q6P4F7
C	-12	HIS	-	EXPRESSION TAG	UNP Q6P4F7
C	-11	HIS	-	EXPRESSION TAG	UNP Q6P4F7
C	-10	SER	-	EXPRESSION TAG	UNP Q6P4F7
C	-9	SER	-	EXPRESSION TAG	UNP Q6P4F7
C	-8	GLY	-	EXPRESSION TAG	UNP Q6P4F7
C	-7	ARG	-	EXPRESSION TAG	UNP Q6P4F7
C	-6	GLU	-	EXPRESSION TAG	UNP Q6P4F7
C	-5	ASN	-	EXPRESSION TAG	UNP Q6P4F7
C	-4	LEU	-	EXPRESSION TAG	UNP Q6P4F7
C	-3	TYR	-	EXPRESSION TAG	UNP Q6P4F7
C	-2	PHE	-	EXPRESSION TAG	UNP Q6P4F7
C	-1	GLN	-	EXPRESSION TAG	UNP Q6P4F7
C	0	GLY	-	EXPRESSION TAG	UNP Q6P4F7
D	-17	MET	-	EXPRESSION TAG	UNP Q6P4F7
D	-16	HIS	-	EXPRESSION TAG	UNP Q6P4F7
D	-15	HIS	-	EXPRESSION TAG	UNP Q6P4F7
D	-14	HIS	-	EXPRESSION TAG	UNP Q6P4F7
D	-13	HIS	-	EXPRESSION TAG	UNP Q6P4F7
D	-12	HIS	-	EXPRESSION TAG	UNP Q6P4F7
D	-11	HIS	-	EXPRESSION TAG	UNP Q6P4F7
D	-10	SER	-	EXPRESSION TAG	UNP Q6P4F7
D	-9	SER	-	EXPRESSION TAG	UNP Q6P4F7

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-8	GLY	-	EXPRESSION TAG	UNP Q6P4F7
D	-7	ARG	-	EXPRESSION TAG	UNP Q6P4F7
D	-6	GLU	-	EXPRESSION TAG	UNP Q6P4F7
D	-5	ASN	-	EXPRESSION TAG	UNP Q6P4F7
D	-4	LEU	-	EXPRESSION TAG	UNP Q6P4F7
D	-3	TYR	-	EXPRESSION TAG	UNP Q6P4F7
D	-2	PHE	-	EXPRESSION TAG	UNP Q6P4F7
D	-1	GLN	-	EXPRESSION TAG	UNP Q6P4F7
D	0	GLY	-	EXPRESSION TAG	UNP Q6P4F7

- Molecule 2 is UNKNOWN ATOM OR ION (three-letter code: UNX) (formula: X).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	4	Total X 4 4	0	0
2	A	1	Total X 1 1	0	0
2	C	1	Total X 1 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	10	Total O 10 10	0	0
3	B	19	Total O 19 19	0	0
3	C	15	Total O 15 15	0	0
3	D	14	Total O 14 14	0	0

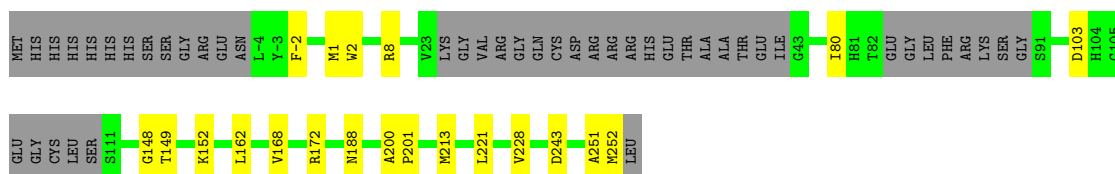
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS failed to run properly.

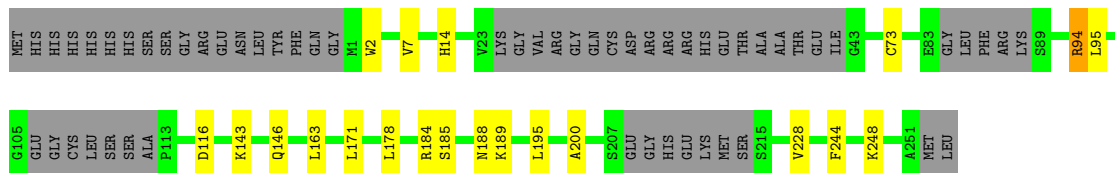
• Molecule 1: Rho GTPase-activating protein 11A

Chain A:



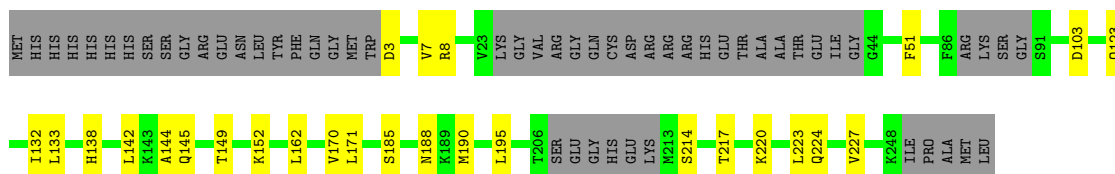
• Molecule 1: Rho GTPase-activating protein 11A

Chain B:



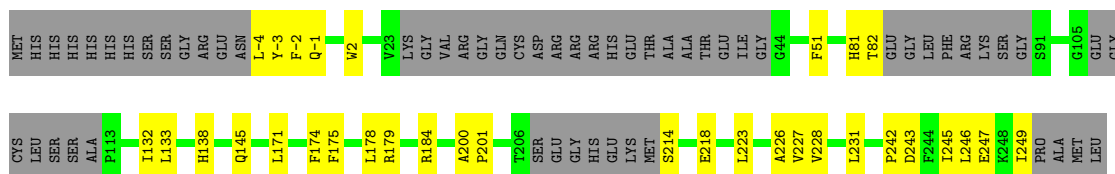
• Molecule 1: Rho GTPase-activating protein 11A

Chain C:



• Molecule 1: Rho GTPase-activating protein 11A

Chain D:



4 Data and refinement statistics

EDS failed to run properly - this section will therefore be incomplete.

Property	Value	
Space group	P 1 21 1	D
Cell constants a, b, c, α , β , γ	46.07Å 106.15Å 107.33Å 90.00° 97.16° 90.00°	D
Resolution (Å)	20.00 – 2.30	D
% Data completeness (in resolution range)	98.1 (20.00-2.30)	D
R_{merge}	0.06	D
R_{sym}	(Not available)	D
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 2.30Å)	Y
Refinement program	REFMAC 5.4.0069, resolve, tlsmd, FFAS03/SCWRL, coot, molprobity	D
R, R_{free}	0.227 , 0.277	D
Wilson B-factor (Å ²)	44.5	Y
Anisotropy	0.028	Y
Estimated twinning fraction	No twinning to report.	Y
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Y
Outliers	0 of 44568 reflections	Y
Total number of atoms	6525	ww
Average B, all atoms (Å ²)	60.0	ww

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.87% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	0.70	0/1701	0.69	0/2321
1	B	0.72	1/1639 (0.1%)	0.69	0/2230
1	C	0.80	0/1645	0.69	0/2235
1	D	0.71	0/1610	0.72	1/2192 (0.0%)
All	All	0.73	1/6595 (0.0%)	0.70	1/8978 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	73	CYS	CB-SG	-5.37	1.73	1.81

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	179	ARG	NE-CZ-NH1	-5.43	117.58	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1665	0	1583	14	0
1	B	1606	0	1567	14	0
1	C	1613	0	1569	14	0
1	D	1577	0	1511	23	0
2	A	1	0	0	0	0
2	B	4	0	0	0	0
2	C	1	0	0	0	0
3	A	10	0	0	0	0
3	B	19	0	0	1	0
3	C	15	0	0	0	0
3	D	14	0	0	0	0
All	All	6525	0	6230	60	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 5.

All (60) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:94:ARG:CG	1:B:94:ARG:HH21	1.64	1.10
1:B:94:ARG:NH2	1:B:94:ARG:HG2	1.46	1.01
1:A:1:MET:O	1:A:251:ALA:HB1	1.88	0.73
1:C:223:LEU:O	1:C:227:VAL:HG23	2.01	0.61
1:C:3:ASP:O	1:C:7:VAL:HG23	2.04	0.57
1:C:190:MET:HG3	1:C:195:LEU:HG	1.87	0.56
1:B:94:ARG:HG2	1:B:94:ARG:HH21	0.68	0.56
1:A:200:ALA:HB3	1:A:201:PRO:HD3	1.88	0.55
1:B:185:SER:HA	1:B:188:ASN:OD1	2.06	0.55
1:D:223:LEU:O	1:D:227:VAL:HG23	2.06	0.55
1:C:142:LEU:HD21	1:C:224:GLN:HG3	1.88	0.54
1:B:94:ARG:NH2	1:B:94:ARG:CG	2.35	0.54
1:B:244:PHE:O	1:B:248:LYS:HD3	2.09	0.52
1:A:243:ASP:OD1	1:D:-3:TYR:CD2	2.62	0.52
1:C:145:GLN:HB2	1:C:223:LEU:HD22	1.91	0.52
1:B:143:LYS:O	1:B:146:GLN:HG2	2.10	0.52
1:B:163:LEU:HD13	1:B:171:LEU:HD12	1.93	0.51
1:D:-4:LEU:O	1:D:-1:GLN:HG2	2.10	0.51
1:D:171:LEU:HD22	1:D:175:PHE:HE2	1.76	0.50
1:D:145:GLN:HG2	1:D:226:ALA:HB3	1.94	0.50
1:B:14:HIS:HE1	3:B:274:HOH:O	1.94	0.49
1:A:149:THR:HG21	1:D:184:ARG:NH1	2.27	0.49
1:A:-2:PHE:HB2	1:D:-2:PHE:HB2	1.95	0.49
1:D:133:LEU:O	1:D:138:HIS:NE2	2.41	0.48
1:A:213:MET:SD	1:A:221:LEU:HD12	2.54	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:81:HIS:O	1:D:82:THR:C	2.51	0.47
1:C:214:SER:OG	1:C:217:THR:HG23	2.14	0.47
1:A:243:ASP:OD1	1:D:-3:TYR:CE2	2.67	0.47
1:C:132:ILE:HD13	1:C:171:LEU:HD21	1.97	0.47
1:D:242:PRO:HD2	1:D:245:ILE:HD13	1.98	0.46
1:B:2:TRP:CD1	1:B:7:VAL:HG21	2.51	0.46
1:D:171:LEU:HD22	1:D:175:PHE:CE2	2.50	0.45
1:A:200:ALA:HA	1:A:228:VAL:HG21	1.98	0.45
1:C:185:SER:HA	1:C:188:ASN:OD1	2.16	0.45
1:A:251:ALA:O	1:A:252:MET:C	2.55	0.45
1:D:171:LEU:HA	1:D:171:LEU:HD23	1.91	0.44
1:D:145:GLN:CG	1:D:226:ALA:HB3	2.48	0.44
1:A:148:GLY:O	1:A:152:LYS:HG3	2.17	0.44
1:A:168:VAL:O	1:A:172:ARG:HG3	2.17	0.44
1:C:144:ALA:O	1:C:152:LYS:HB2	2.18	0.44
1:B:178:LEU:HD22	1:B:195:LEU:HD13	1.99	0.44
1:B:184:ARG:HH11	1:C:149:THR:CB	2.31	0.43
1:D:243:ASP:O	1:D:247:GLU:HG3	2.19	0.43
1:B:95:LEU:HD13	1:B:116:ASP:HA	2.01	0.43
1:B:200:ALA:HA	1:B:228:VAL:HG21	2.01	0.42
1:D:214:SER:O	1:D:218:GLU:N	2.45	0.42
1:C:123:GLN:HE21	1:C:123:GLN:HB2	1.69	0.42
1:D:145:GLN:HB3	1:D:223:LEU:HD22	2.01	0.42
1:C:133:LEU:O	1:C:138:HIS:NE2	2.40	0.41
1:D:246:LEU:HA	1:D:249:ILE:HD12	2.02	0.41
1:D:174:PHE:CE2	1:D:178:LEU:HD11	2.56	0.41
1:C:51:PHE:CE1	1:C:170:VAL:HG21	2.55	0.41
1:D:227:VAL:O	1:D:231:LEU:HG	2.21	0.41
1:A:2:TRP:CE2	1:A:251:ALA:HB2	2.55	0.41
1:A:80:ILE:O	1:A:188:ASN:HB3	2.20	0.41
1:D:200:ALA:HA	1:D:228:VAL:HG21	2.02	0.41
1:C:8:ARG:NH1	1:C:162:LEU:HD22	2.36	0.41
1:A:8:ARG:HD3	1:A:162:LEU:CD2	2.51	0.41
1:D:51:PHE:HZ	1:D:132:ILE:HD12	1.86	0.40
1:D:200:ALA:HB3	1:D:201:PRO:HD3	2.03	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 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	217/271 (80%)	211 (97%)	6 (3%)	0	100	100
1	B	203/271 (75%)	201 (99%)	2 (1%)	0	100	100
1	C	208/271 (77%)	204 (98%)	4 (2%)	0	100	100
1	D	202/271 (74%)	195 (96%)	7 (4%)	0	100	100
All	All	830/1084 (77%)	811 (98%)	19 (2%)	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	164/233 (70%)	163 (99%)	1 (1%)	92	97
1	B	163/233 (70%)	161 (99%)	2 (1%)	82	92
1	C	160/233 (69%)	158 (99%)	2 (1%)	80	91
1	D	156/233 (67%)	155 (99%)	1 (1%)	92	97
All	All	643/932 (69%)	637 (99%)	6 (1%)	87	95

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	103	ASP
1	B	94	ARG
1	B	189	LYS
1	C	103	ASP
1	C	220	LYS

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Mol	Chain	Res	Type
1	D	2	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	202	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

Of 6 ligands modelled in this entry, 6 are unknown - 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.

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 ⓘ

EDS failed to run properly - this section will therefore be empty.

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

EDS failed to run properly - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS failed to run properly - this section will therefore be empty.

6.4 Ligands ⓘ

EDS failed to run properly - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS failed to run properly - this section will therefore be empty.