



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 07:55 PM GMT

PDB ID : 4X6Q
Title : An Isoform-specific Myristylation Switch Targets RIIB PKA Holoenzymes to Membranes
Authors : Zhang, P.; Ye, F.; Bastidas, A.C.; Kornev, A.P.; Ginsberg, M.H.; Taylor, S.S.
Deposited on : 2014-12-08
Resolution : 2.52 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20026982

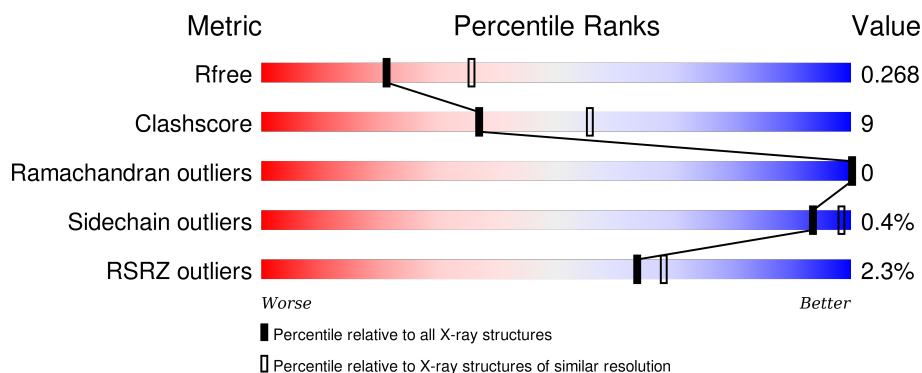
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.52 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}	91344	4241 (2.54-2.50)
Clashscore	102246	4968 (2.54-2.50)
Ramachandran outliers	100387	4873 (2.54-2.50)
Sidechain outliers	100360	4875 (2.54-2.50)
RSRZ outliers	91569	4253 (2.54-2.50)

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. 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	B	416	
2	C	350	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4995 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 cAMP-dependent protein kinase type II-beta regulatory subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	270	Total	C	N	O	S	0	0	0
			2132	1344	374	399	15			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	230	LYS	ARG	conflict	UNP P31324

- Molecule 2 is a protein called cAMP-dependent protein kinase catalytic subunit alpha.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	C	337	Total	C	N	O	P	S	0	0	0
			2787	1803	466	507	3	8			

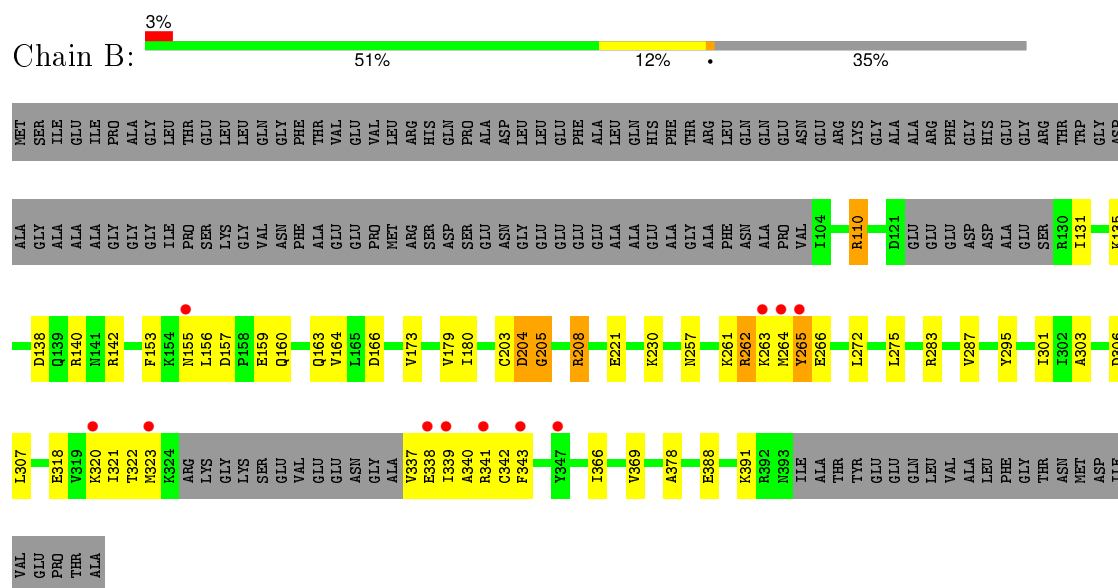
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	41	Total	O	0	0
			41	41		
3	C	35	Total	O	0	0
			35	35		

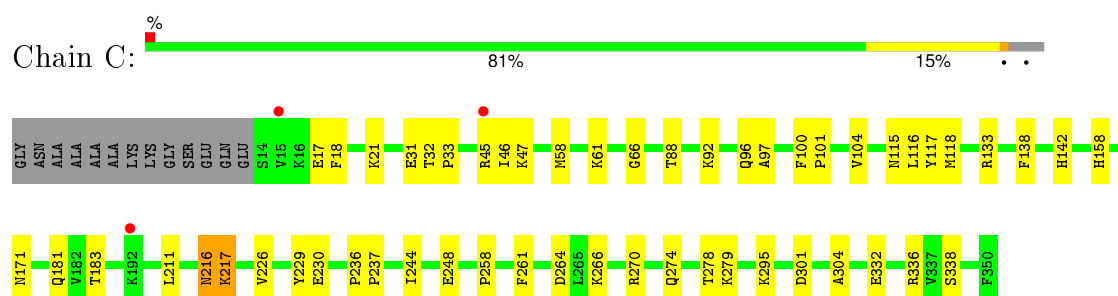
3 Residue-property plots

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.

- Molecule 1: cAMP-dependent protein kinase type II-beta regulatory subunit



- Molecule 2: cAMP-dependent protein kinase catalytic subunit alpha



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	152.71Å 214.04Å 61.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.82 – 2.52 48.07 – 2.52	Depositor EDS
% Data completeness (in resolution range)	95.3 (43.82-2.52) 95.1 (48.07-2.52)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.198 , 0.261 0.215 , 0.268	Depositor DCC
R_{free} test set	1661 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	57.1	Xtriage
Anisotropy	0.571	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 53.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 33259 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4995	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.99% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	B	0.72	3/2164 (0.1%)	0.99	12/2908 (0.4%)
2	C	0.54	0/2824	0.65	1/3801 (0.0%)
All	All	0.63	3/4988 (0.1%)	0.81	13/6709 (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	3
2	C	0	1
All	All	0	4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	342	CYS	C-N	-13.21	1.03	1.34
1	B	208	ARG	C-N	-12.73	1.04	1.34
1	B	343	PHE	C-N	10.46	1.58	1.34

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	208	ARG	O-C-N	-20.19	90.39	122.70
1	B	262	ARG	CB-CA-C	-15.24	79.92	110.40
1	B	208	ARG	CA-C-N	14.33	148.73	117.20
1	B	262	ARG	N-CA-C	12.48	144.69	111.00
1	B	156	LEU	N-CA-CB	-7.63	95.14	110.40
2	C	216	ASN	N-CA-C	-7.42	90.98	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	203	CYS	N-CA-C	6.92	129.68	111.00
1	B	208	ARG	C-N-CA	6.36	137.60	121.70
1	B	263	LYS	N-CA-C	-6.08	94.57	111.00
1	B	110	ARG	NE-CZ-NH1	5.34	122.97	120.30
1	B	342	CYS	O-C-N	5.33	131.23	122.70
1	B	342	CYS	C-N-CA	-5.32	108.40	121.70
1	B	307	LEU	CA-CB-CG	5.20	127.27	115.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	205	GLY	Peptide
1	B	208	ARG	Peptide
1	B	265	TYR	Peptide
2	C	216	ASN	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	B	2132	0	2137	53	0
2	C	2787	0	2767	52	0
3	B	41	0	0	3	0
3	C	35	0	0	4	0
All	All	4995	0	4904	92	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 9.

All (92) 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:B:110:ARG:HH21	2:C:133:ARG:HH12	1.24	0.85
2:C:31:GLU:C	2:C:33:PRO:HD3	1.97	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:110:ARG:NH2	2:C:133:ARG:HH12	1.75	0.83
1:B:221:GLU:HG2	2:C:244:ILE:HD11	1.62	0.80
1:B:110:ARG:NH2	2:C:230:GLU:OE1	2.19	0.76
1:B:221:GLU:HG2	2:C:244:ILE:CD1	2.17	0.74
1:B:323:MET:O	1:B:337:VAL:N	2.22	0.73
1:B:204:ASP:H	1:B:205:GLY:HA3	1.54	0.72
1:B:155:ASN:HD21	2:C:211:LEU:HD13	1.54	0.70
2:C:97:ALA:O	3:C:414:HOH:O	2.10	0.70
1:B:159:GLU:OE2	1:B:257:ASN:ND2	2.24	0.69
2:C:31:GLU:O	2:C:33:PRO:HD3	1.93	0.69
1:B:157:ASP:OD1	1:B:160:GLN:HB2	1.93	0.69
1:B:155:ASN:ND2	2:C:211:LEU:HD13	2.08	0.69
1:B:261:LYS:O	1:B:265:TYR:HB2	1.94	0.67
2:C:295:LYS:NZ	3:C:421:HOH:O	2.12	0.67
1:B:221:GLU:CG	2:C:244:ILE:HD13	2.25	0.66
2:C:18:PHE:HA	2:C:21:LYS:HE2	1.78	0.66
1:B:272:LEU:HD22	1:B:275:LEU:HD12	1.77	0.66
1:B:303:ALA:HB3	1:B:306:ASP:OD2	1.95	0.66
2:C:17:GLU:O	2:C:21:LYS:HG3	1.99	0.62
1:B:221:GLU:CG	2:C:244:ILE:CD1	2.78	0.61
1:B:140:ARG:NH1	1:B:166:ASP:OD1	2.33	0.61
1:B:264:MET:N	3:B:532:HOH:O	2.23	0.61
2:C:158:HIS:CE1	2:C:217:LYS:HB3	2.36	0.59
1:B:153:PHE:HE2	1:B:164:VAL:HG11	1.69	0.57
2:C:32:THR:N	2:C:33:PRO:HD3	2.14	0.57
1:B:157:ASP:OD2	1:B:261:LYS:HE3	2.03	0.57
2:C:61:LYS:HE2	2:C:66:GLY:HA2	1.87	0.56
1:B:266:GLU:HG3	1:B:283:ARG:HD2	1.88	0.56
1:B:262:ARG:O	1:B:287:VAL:HG11	2.06	0.56
1:B:173:VAL:HG11	1:B:179:VAL:HG22	1.88	0.54
2:C:92:LYS:HA	2:C:118:MET:HE1	1.90	0.54
2:C:295:LYS:HG2	3:C:421:HOH:O	2.09	0.53
2:C:46:ILE:HD11	2:C:61:LYS:HB2	1.90	0.53
1:B:318:GLU:HG2	1:B:366:ILE:HD12	1.91	0.53
1:B:341:ARG:HH21	1:B:366:ILE:HD11	1.74	0.52
1:B:110:ARG:NH2	2:C:133:ARG:NH1	2.52	0.52
2:C:33:PRO:HB3	2:C:96:GLN:HE22	1.75	0.51
1:B:378:ALA:HB1	2:C:278:THR:HG21	1.91	0.51
2:C:33:PRO:HB3	2:C:96:GLN:NE2	2.26	0.51
2:C:236:PRO:HB2	3:C:403:HOH:O	2.10	0.50
2:C:115:ASN:HB2	2:C:117:TYR:CE2	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:320:LYS:HE2	1:B:322:THR:CG2	2.42	0.50
1:B:204:ASP:HB2	1:B:205:GLY:HA3	1.93	0.50
1:B:157:ASP:OD2	1:B:261:LYS:CE	2.60	0.50
2:C:88:THR:HG21	2:C:116:LEU:HD12	1.94	0.49
1:B:159:GLU:O	1:B:163:GLN:HG3	2.13	0.49
1:B:318:GLU:H	1:B:369:VAL:HG22	1.77	0.49
2:C:88:THR:HG21	2:C:116:LEU:CD1	2.43	0.48
1:B:295:TYR:CZ	1:B:301:ILE:HG23	2.48	0.48
2:C:226:VAL:HG13	2:C:237:PRO:HD2	1.95	0.47
1:B:388:GLU:OE2	1:B:391:LYS:HE3	2.15	0.47
2:C:100:PHE:CD1	2:C:101:PRO:HD2	2.49	0.47
2:C:244:ILE:O	2:C:248:GLU:HG3	2.15	0.47
1:B:138:ASP:OD2	1:B:142:ARG:HD3	2.14	0.47
1:B:283:ARG:O	1:B:287:VAL:HG23	2.15	0.47
2:C:138:PHE:HB3	2:C:142:HIS:HB3	1.97	0.47
2:C:266:LYS:O	2:C:270:ARG:HG3	2.14	0.47
2:C:274:GLN:HG2	2:C:279:LYS:HB2	1.96	0.46
1:B:323:MET:HB2	1:B:339:ILE:HD11	1.97	0.46
2:C:45:ARG:HD2	2:C:58:MET:SD	2.55	0.46
1:B:204:ASP:N	1:B:205:GLY:HA3	2.20	0.46
2:C:47:LYS:HG2	2:C:332:GLU:OE2	2.16	0.46
1:B:320:LYS:HG2	1:B:322:THR:HG23	1.98	0.46
2:C:32:THR:OG1	2:C:32:THR:O	2.27	0.45
1:B:131:ILE:O	3:B:535:HOH:O	2.21	0.45
2:C:96:GLN:HG2	2:C:97:ALA:N	2.32	0.45
1:B:341:ARG:NH2	1:B:366:ILE:HD11	2.32	0.44
2:C:115:ASN:HB2	2:C:117:TYR:CZ	2.52	0.44
2:C:301:ASP:OD2	2:C:304:ALA:HB2	2.16	0.44
1:B:221:GLU:HG3	2:C:244:ILE:HD13	1.96	0.44
2:C:264:ASP:OD1	2:C:295:LYS:HD2	2.17	0.44
2:C:336:ARG:NH2	2:C:338:SEP:O3P	2.51	0.44
1:B:321:ILE:HB	1:B:340:ALA:HB3	2.00	0.43
1:B:320:LYS:HE2	1:B:322:THR:HG21	1.99	0.43
1:B:204:ASP:CB	1:B:205:GLY:HA3	2.47	0.43
1:B:204:ASP:H	1:B:205:GLY:CA	2.27	0.43
2:C:217:LYS:H	2:C:217:LYS:HG3	1.50	0.42
1:B:180:ILE:HD13	1:B:230:LYS:HD2	2.02	0.42
2:C:104:VAL:CG2	2:C:181:GLN:HB3	2.49	0.42
1:B:391:LYS:HB2	1:B:391:LYS:HE2	1.74	0.42
1:B:337:VAL:HG13	1:B:338:GLU:N	2.36	0.41
2:C:100:PHE:CG	2:C:101:PRO:HD2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:258:PRO:HD2	2:C:261:PHE:CD1	2.56	0.41
2:C:18:PHE:CA	2:C:21:LYS:HE2	2.48	0.41
1:B:323:MET:CB	1:B:339:ILE:HD11	2.51	0.41
2:C:171:ASN:O	2:C:183:THR:HG22	2.21	0.41
1:B:155:ASN:HD21	2:C:211:LEU:CD1	2.28	0.40
1:B:135:LYS:HE3	1:B:135:LYS:HB3	1.83	0.40
2:C:229:TYR:CE1	2:C:237:PRO:HB3	2.56	0.40
1:B:131:ILE:HG22	3:B:535:HOH:O	2.21	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	B	264/416 (64%)	251 (95%)	13 (5%)	0	100	100
2	C	332/350 (95%)	317 (96%)	15 (4%)	0	100	100
All	All	596/766 (78%)	568 (95%)	28 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	B	229/340 (67%)	228 (100%)	1 (0%)	93	98

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	C	295/302 (98%)	294 (100%)	1 (0%)	94	98
All	All	524/642 (82%)	522 (100%)	2 (0%)	93	98

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	204	ASP
2	C	217	LYS

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

Mol	Chain	Res	Type
1	B	346	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 non-standard protein/DNA/RNA residues are modelled in this entry.

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
2	SEP	C	139	2	7,9,10	1.82	3 (42%)	8,12,14	1.50	2 (25%)
2	TPO	C	197	2	7,10,11	1.48	1 (14%)	10,14,16	1.65	3 (30%)
2	SEP	C	338	2	7,9,10	1.54	1 (14%)	8,12,14	2.11	1 (12%)

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
2	SEP	C	139	2	-	0/5/8/10	0/0/0/0
2	TPO	C	197	2	-	0/8/11/13	0/0/0/0
2	SEP	C	338	2	-	0/5/8/10	0/0/0/0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	139	SEP	P-O3P	2.04	1.61	1.54
2	C	139	SEP	P-O2P	2.07	1.61	1.54
2	C	197	TPO	P-O1P	2.87	1.59	1.50
2	C	338	SEP	P-O1P	3.00	1.60	1.50
2	C	139	SEP	P-O1P	3.61	1.62	1.50

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	197	TPO	P-OG1-CB	-2.81	109.12	121.42
2	C	139	SEP	O-C-CA	-2.30	119.56	125.72
2	C	197	TPO	C-CA-N	2.03	114.43	109.95
2	C	197	TPO	OG1-P-O1P	2.62	113.73	107.48
2	C	139	SEP	OG-CB-CA	3.08	110.94	108.26
2	C	338	SEP	OG-CB-CA	5.17	112.77	108.26

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	338	SEP	1	0

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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	208:ARG	C	209:CYS	N	1.04
1	B	342:CYS	C	343:PHE	N	1.03

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	B	270/416 (64%)	0.05	11 (4%) 41 46	40, 60, 87, 105	0
2	C	334/350 (95%)	0.01	3 (0%) 85 88	38, 59, 84, 103	0
All	All	604/766 (78%)	0.03	14 (2%) 64 68	38, 60, 85, 105	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	343	PHE	5.7
1	B	263	LYS	5.3
1	B	264	MET	4.9
1	B	323	MET	3.9
1	B	339	ILE	3.6
2	C	45	ARG	3.2
1	B	320	LYS	3.1
2	C	15	VAL	2.9
1	B	265	TYR	2.7
1	B	341	ARG	2.6
1	B	338	GLU	2.3
1	B	347	TYR	2.3
2	C	192	LYS	2.2
1	B	155	ASN	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	SEP	C	338	10/11	0.95	0.13	-	73,77,84,87	0
2	TPO	C	197	11/12	0.98	0.17	-	39,44,47,49	0
2	SEP	C	139	10/11	0.95	0.11	-	60,68,78,86	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.