



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 21, 2018 – 11:18 AM EST

PDB ID : 5ORH
Title : The crystal structure of CK2alpha in complex with compound 2
Authors : Brear, P.; De Fusco, C.; Iegre, J.; Yoshida, M.; Mitchell, S.; Rossmann, M.; Carro, L.; Sore, H.; Hyvonen, M.; Spring, D.
Deposited on : 2017-08-16
Resolution : 1.75 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030736
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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-20030736

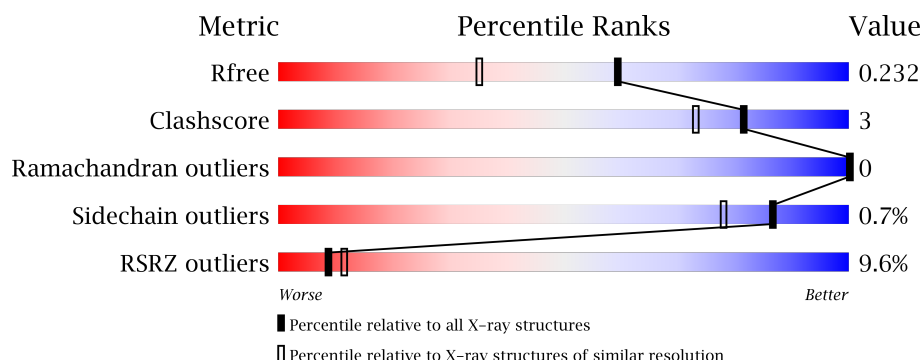
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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}	100719	1762 (1.76-1.76)
Clashscore	112137	1889 (1.76-1.76)
Ramachandran outliers	110173	1868 (1.76-1.76)
Sidechain outliers	110143	1868 (1.76-1.76)
RSRZ outliers	101464	1770 (1.76-1.76)

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	A	352	
1	B	352	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	A4N	B	401	-	-	-	X
2	A4N	B	402	-	-	X	-
3	ACT	B	405	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5902 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 Casein kinase II subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	324	Total	C	N	O	S	0	2	0
			2754	1766	481	495	12			
1	B	325	Total	C	N	O	S	0	2	0
			2758	1768	482	496	12			

There are 50 discrepancies between the modelled and reference sequences:

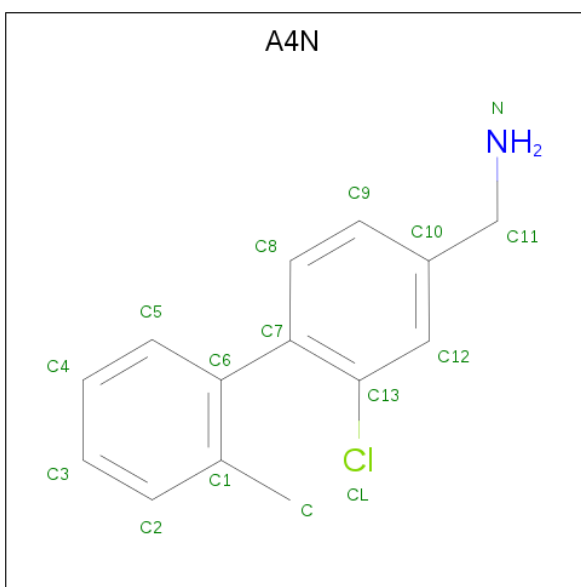
Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	GLY	-	expression tag	UNP P68400
A	-21	SER	-	expression tag	UNP P68400
A	-20	MET	-	expression tag	UNP P68400
A	-19	ASP	-	expression tag	UNP P68400
A	-18	ILE	-	expression tag	UNP P68400
A	-17	GLU	-	expression tag	UNP P68400
A	-16	PHE	-	expression tag	UNP P68400
A	-15	ASP	-	expression tag	UNP P68400
A	-14	ASP	-	expression tag	UNP P68400
A	-13	ASP	-	expression tag	UNP P68400
A	-12	ALA	-	expression tag	UNP P68400
A	-11	ASP	-	expression tag	UNP P68400
A	-10	ASP	-	expression tag	UNP P68400
A	-9	ASP	-	expression tag	UNP P68400
A	-8	GLY	-	expression tag	UNP P68400
A	-7	SER	-	expression tag	UNP P68400
A	-6	GLY	-	expression tag	UNP P68400
A	-5	SER	-	expression tag	UNP P68400
A	-4	GLY	-	expression tag	UNP P68400
A	-3	SER	-	expression tag	UNP P68400
A	-2	GLY	-	expression tag	UNP P68400
A	-1	SER	-	expression tag	UNP P68400
A	0	GLY	-	expression tag	UNP P68400
A	1	SER	-	expression tag	UNP P68400
A	21	SER	ARG	engineered mutation	UNP P68400

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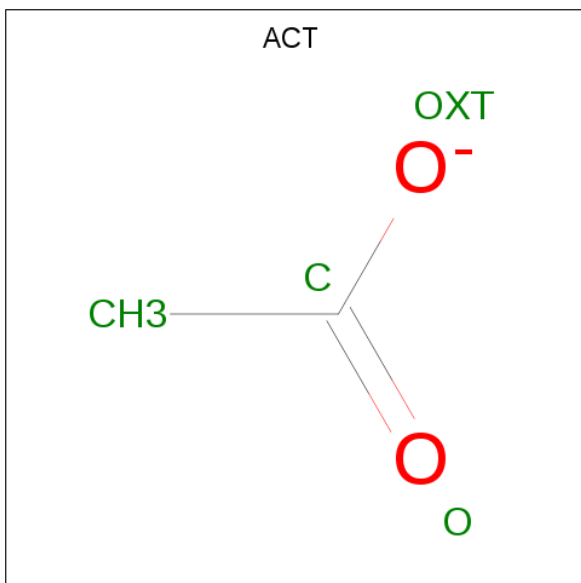
Chain	Residue	Modelled	Actual	Comment	Reference
B	-22	GLY	-	expression tag	UNP P68400
B	-21	SER	-	expression tag	UNP P68400
B	-20	MET	-	expression tag	UNP P68400
B	-19	ASP	-	expression tag	UNP P68400
B	-18	ILE	-	expression tag	UNP P68400
B	-17	GLU	-	expression tag	UNP P68400
B	-16	PHE	-	expression tag	UNP P68400
B	-15	ASP	-	expression tag	UNP P68400
B	-14	ASP	-	expression tag	UNP P68400
B	-13	ASP	-	expression tag	UNP P68400
B	-12	ALA	-	expression tag	UNP P68400
B	-11	ASP	-	expression tag	UNP P68400
B	-10	ASP	-	expression tag	UNP P68400
B	-9	ASP	-	expression tag	UNP P68400
B	-8	GLY	-	expression tag	UNP P68400
B	-7	SER	-	expression tag	UNP P68400
B	-6	GLY	-	expression tag	UNP P68400
B	-5	SER	-	expression tag	UNP P68400
B	-4	GLY	-	expression tag	UNP P68400
B	-3	SER	-	expression tag	UNP P68400
B	-2	GLY	-	expression tag	UNP P68400
B	-1	SER	-	expression tag	UNP P68400
B	0	GLY	-	expression tag	UNP P68400
B	1	SER	-	expression tag	UNP P68400
B	21	SER	ARG	engineered mutation	UNP P68400

- Molecule 2 is [3-chloranyl-4-(2-methylphenyl)phenyl]methanamine (three-letter code: A4N) (formula: C₁₄H₁₄ClN).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	0	0
			16	14	1	1		
2	A	1	Total	C	Cl	N	0	0
			16	14	1	1		
2	B	1	Total	C	Cl	N	0	0
			16	14	1	1		
2	B	1	Total	C	Cl	N	0	0
			16	14	1	1		

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).

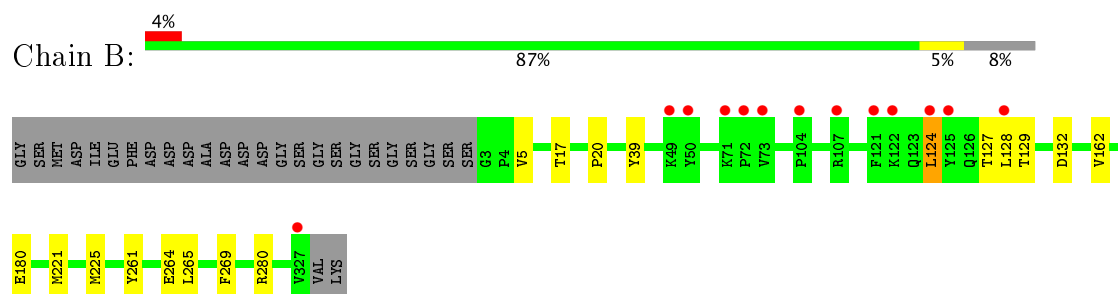
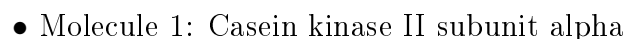


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	105	Total O 105 105	0	0
4	B	209	Total O 209 209	0	0

- Molecule 1: Casein kinase II subunit alpha



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	64.52Å 67.97Å 333.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	83.33 – 1.75 83.33 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.9 (83.33-1.75) 99.9 (83.33-1.75)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 1.75Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.208 , 0.227 0.210 , 0.232	Depositor DCC
R_{free} test set	3778 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	28.0	Xtriage
Anisotropy	0.460	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 50.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.054 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5902	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.60% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: A4N, ACT

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	0.50	0/2829	0.61	0/3826
1	B	0.64	0/2833	0.64	0/3832
All	All	0.58	0/5662	0.63	0/7658

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

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	2754	0	2693	18	0
1	B	2758	0	2695	15	1
2	A	32	0	0	2	0
2	B	32	0	0	8	0
3	B	12	0	9	0	0
4	A	105	0	0	0	0
4	B	209	0	0	2	0
All	All	5902	0	5397	33	1

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 3.

All (33) 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:225:MET:HG3	2:B:402:A4N:C2	2.02	0.90
1:A:116:VAL:HB	1:A:163[B]:MET:CE	2.25	0.67
1:B:124:LEU:HA	1:B:127:THR:HG22	1.79	0.63
1:B:221:MET:HE1	2:B:402:A4N:CL	2.36	0.63
1:B:225:MET:HG3	2:B:402:A4N:C1	2.29	0.62
1:A:116:VAL:HB	1:A:163[B]:MET:HE1	1.82	0.61
1:A:128:LEU:HD22	1:A:132:ASP:HB3	1.84	0.59
1:A:133:ILE:HG12	1:A:225:MET:HE3	1.85	0.58
1:B:39:TYR:O	2:B:401:A4N:N	2.39	0.56
1:A:105:VAL:HG21	1:B:20:PRO:HG2	1.89	0.54
1:B:5:VAL:HB	1:B:261:TYR:HA	1.90	0.53
1:B:225:MET:CG	2:B:402:A4N:C2	2.84	0.53
1:A:5:VAL:HB	1:A:261:TYR:HA	1.91	0.52
1:B:17:THR:HG22	4:B:602:HOH:O	2.13	0.49
1:A:49:LYS:HD3	1:A:50:TYR:CE1	2.49	0.48
2:B:402:A4N:N	4:B:502:HOH:O	2.36	0.47
1:B:129:THR:HG22	1:B:132:ASP:CG	2.35	0.47
1:A:116:VAL:HB	1:A:163[B]:MET:HE2	1.98	0.46
1:B:129:THR:HG23	1:B:132:ASP:H	1.81	0.46
1:B:162:VAL:HG11	2:B:402:A4N:CL	2.53	0.45
1:A:47:ARG:HG3	1:A:52:GLU:HG2	1.99	0.45
1:A:33:TRP:CZ3	1:A:100:ILE:HG22	2.52	0.45
1:A:26:TYR:HE1	1:A:76:LYS:HG2	1.83	0.43
1:A:128:LEU:HD22	1:A:132:ASP:CB	2.48	0.43
1:B:124:LEU:O	1:B:128:LEU:HG	2.19	0.42
1:A:121:PHE:HZ	1:A:230:GLU:HG3	1.85	0.42
1:B:265:LEU:HD22	1:B:269:PHE:CD2	2.56	0.41
1:A:140:ILE:HD11	2:A:402:A4N:CL	2.58	0.41
1:A:303:LYS:HB3	1:A:313:LEU:HG	2.02	0.41
1:A:225:MET:HG3	2:A:402:A4N:C1	2.51	0.41
1:A:128:LEU:HD12	1:A:225:MET:CE	2.50	0.41
1:A:103:ASP:HA	1:A:104:PRO:HD2	1.91	0.40
1:B:225:MET:CG	2:B:402:A4N:C1	2.98	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:264:GLU:OE2	1:B:280:ARG:NH2[8_597]	1.74	0.46

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	324/352 (92%)	314 (97%)	10 (3%)	0	100	100
1	B	325/352 (92%)	319 (98%)	6 (2%)	0	100	100
All	All	649/704 (92%)	633 (98%)	16 (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	301/319 (94%)	299 (99%)	2 (1%)	87	78
1	B	301/319 (94%)	299 (99%)	2 (1%)	87	78
All	All	602/638 (94%)	598 (99%)	4 (1%)	87	78

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	ILE
1	A	205	ASP
1	B	124	LEU
1	B	180	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	118	ASN
1	A	262	ASN
1	B	118	ASN
1	B	168	HIS
1	B	186	GLN
1	B	262	ASN
1	B	270	ASN

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 ⓘ

7 ligands 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	A4N	A	401	-	17,17,17	0.19	0	23,23,23	0.20	0
2	A4N	A	402	-	17,17,17	0.23	0	23,23,23	0.41	0
2	A4N	B	401	-	17,17,17	0.19	0	23,23,23	0.20	0
2	A4N	B	402	-	17,17,17	0.21	0	23,23,23	0.20	0
3	ACT	B	403	-	1,3,3	5.54	1 (100%)	0,3,3	0.00	-
3	ACT	B	404	-	1,3,3	6.13	1 (100%)	0,3,3	0.00	-
3	ACT	B	405	-	1,3,3	5.60	1 (100%)	0,3,3	0.00	-

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	A4N	A	401	-	-	0/6/6/6	0/2/2/2
2	A4N	A	402	-	-	0/6/6/6	0/2/2/2
2	A4N	B	401	-	-	0/6/6/6	0/2/2/2
2	A4N	B	402	-	-	0/6/6/6	0/2/2/2
3	ACT	B	403	-	-	0/0/0/0	0/0/0/0
3	ACT	B	404	-	-	0/0/0/0	0/0/0/0
3	ACT	B	405	-	-	0/0/0/0	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	403	ACT	CH3-C	5.54	1.55	1.48
3	B	405	ACT	CH3-C	5.60	1.56	1.48
3	B	404	ACT	CH3-C	6.13	1.56	1.48

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	402	A4N	2	0
2	B	401	A4N	1	0
2	B	402	A4N	7	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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	324/352 (92%)	0.85	49 (15%) 3 4	25, 42, 72, 91	0
1	B	325/352 (92%)	0.36	13 (4%) 39 46	18, 28, 58, 90	0
All	All	649/704 (92%)	0.61	62 (9%) 9 11	18, 35, 69, 91	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	50	TYR	9.2
1	B	124	LEU	9.0
1	A	50	TYR	8.4
1	A	125	TYR	8.2
1	B	121	PHE	7.6
1	A	47	ARG	5.8
1	A	278	ARG	5.7
1	A	270	ASN	5.7
1	A	73	VAL	5.1
1	A	126	GLN	5.1
1	A	326	THR	5.0
1	A	49	LYS	5.0
1	B	125	TYR	4.9
1	A	48	GLY	4.2
1	A	75	LYS	4.0
1	A	286	HIS	3.9
1	A	288	GLU	3.9
1	A	121	PHE	3.8
1	A	104	PRO	3.8
1	A	107	ARG	3.7
1	A	307	TYR	3.7
1	B	107	ARG	3.7
1	A	325	TYR	3.6
1	A	40	GLN	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	49	LYS	3.5
1	A	105	VAL	3.4
1	A	257	TYR	3.3
1	A	255	TYR	3.3
1	B	72	PRO	3.3
1	A	283	ARG	3.1
1	A	267	PRO	3.1
1	A	292	LEU	2.9
1	B	73	VAL	2.9
1	A	117	ASN	2.9
1	A	227	PHE	2.8
1	A	131	TYR	2.8
1	A	225	MET	2.7
1	A	251	THR	2.7
1	A	239	TYR	2.7
1	A	252	GLU	2.7
1	B	71	LYS	2.6
1	A	254	LEU	2.6
1	B	128	LEU	2.6
1	A	290	GLN	2.6
1	A	271	ASP	2.5
1	A	237	ASP	2.5
1	B	327	VAL	2.5
1	B	104	PRO	2.4
1	A	71	LYS	2.4
1	A	45	LEU	2.3
1	A	72	PRO	2.3
1	A	33	TRP	2.2
1	A	124	LEU	2.2
1	A	281	TRP	2.2
1	A	284	PHE	2.2
1	A	273	LEU	2.2
1	B	122	LYS	2.2
1	A	122	LYS	2.2
1	A	74	LYS	2.1
1	A	272	ILE	2.1
1	A	106	SER	2.0
1	A	57	ILE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. 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
3	ACT	B	405	4/4	0.88	0.15	3.00	59,61,61,62	0
2	A4N	B	401	16/16	0.85	0.18	2.16	34,37,46,47	0
2	A4N	A	402	16/16	0.76	0.19	1.62	32,47,49,56	0
2	A4N	A	401	16/16	0.76	0.24	1.36	45,53,59,61	0
3	ACT	B	404	4/4	0.94	0.10	0.76	29,30,31,38	0
2	A4N	B	402	16/16	0.71	0.18	0.46	33,38,42,54	0
3	ACT	B	403	4/4	0.94	0.10	-0.38	36,41,42,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.