



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 11, 2021 – 04:19 AM EDT

PDB ID : 2QR8
Title : 2.0A X-ray structure of C-terminal kinase domain of p90 ribosomal S6 kinase 2 (RSK2)
Authors : Tereshko, V.; Malakhova, M.; Dong, Z.
Deposited on : 2007-07-27
Resolution : 2.00 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

<https://www.wwpdb.org/validation/2017/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
Xtriage (Phenix) : 1.13
EDS : 2.23.2
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

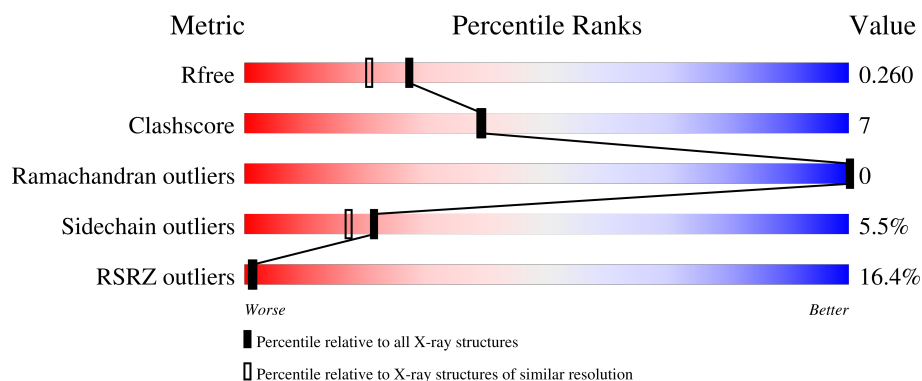
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	342	<div> <div>14%</div> <div>70%</div> <div>16%</div> <div>13%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2452 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 Ribosomal protein S6 kinase alpha-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	298	Total	C	N	O	S	0	3	0
			2374	1513	402	446	13			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	591	GLU	LYS	engineered mutation	UNP P18654

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Na	0	0
			1	1		

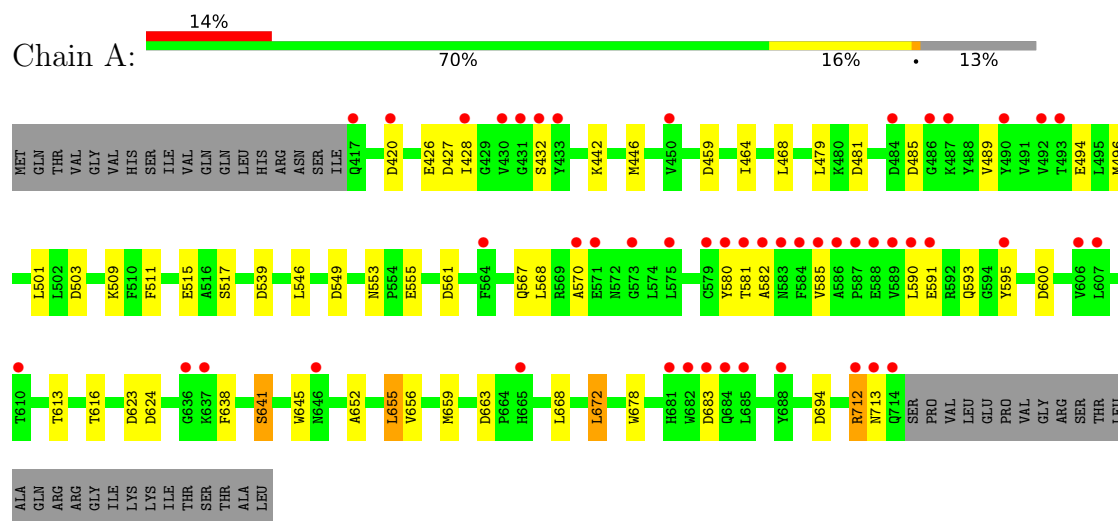
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	77	Total	O	0	0
			77	77		

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: Ribosomal protein S6 kinase alpha-3



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	46.91Å 46.91Å 291.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 19.74 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.5 (20.00-2.00) 99.5 (19.74-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.215 , 0.259 0.218 , 0.260	Depositor DCC
R_{free} test set	1197 reflections (5.17%)	wwPDB-VP
Wilson B-factor (Å ²)	42.3	Xtriage
Anisotropy	0.109	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 69.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2452	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.08% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.65	0/2440	0.97	15/3310 (0.5%)

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	549	ASP	CB-CG-OD2	6.97	124.57	118.30
1	A	623	ASP	CB-CG-OD2	6.38	124.04	118.30
1	A	683	ASP	CB-CG-OD2	5.99	123.69	118.30
1	A	481	ASP	CB-CG-OD2	5.97	123.68	118.30
1	A	459	ASP	CB-CG-OD2	5.89	123.60	118.30
1	A	694	ASP	CB-CG-OD2	5.83	123.54	118.30
1	A	539	ASP	CB-CG-OD2	5.61	123.35	118.30
1	A	427	ASP	CB-CG-OD2	5.43	123.19	118.30
1	A	663	ASP	CB-CG-OD2	5.40	123.16	118.30
1	A	485	ASP	CB-CG-OD2	5.34	123.11	118.30
1	A	561	ASP	CB-CG-OD2	5.31	123.08	118.30
1	A	420	ASP	CB-CG-OD2	5.15	122.94	118.30
1	A	600	ASP	CB-CG-OD2	5.10	122.89	118.30
1	A	624	ASP	CB-CG-OD2	5.10	122.89	118.30
1	A	503	ASP	CB-CG-OD2	5.09	122.88	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 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	2374	0	2339	35	1
2	A	1	0	0	0	0
3	A	77	0	0	4	0
All	All	2452	0	2339	35	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 7.

All (35) 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:712:ARG:HH11	1:A:713:ASN:ND2	1.54	1.04
1:A:616:THR:HG23	3:A:1020:HOH:O	1.78	0.81
1:A:616:THR:CG2	3:A:1020:HOH:O	2.31	0.78
1:A:712:ARG:HD2	1:A:713:ASN:N	2.01	0.74
1:A:428:ILE:HG22	1:A:428:ILE:O	1.87	0.74
1:A:568:LEU:HD13	1:A:593:GLN:HG3	1.70	0.72
1:A:464:ILE:HD11	1:A:489:VAL:HG11	1.72	0.71
1:A:712:ARG:HH11	1:A:713:ASN:HD22	1.37	0.70
1:A:494:GLU:HG3	1:A:496:MET:CE	2.22	0.69
1:A:580:TYR:CZ	1:A:582:ALA:HB3	2.31	0.64
1:A:712:ARG:HH11	1:A:713:ASN:HD21	1.46	0.64
1:A:655:LEU:HD22	1:A:659:MET:SD	2.40	0.62
1:A:712:ARG:NH1	1:A:713:ASN:ND2	2.39	0.62
1:A:712:ARG:HD3	1:A:713:ASN:ND2	2.18	0.59
1:A:668:LEU:HD22	1:A:672:LEU:HB3	1.84	0.59
1:A:428:ILE:O	1:A:428:ILE:CG2	2.52	0.58
1:A:570:ALA:HA	1:A:585:VAL:HG22	1.86	0.58
1:A:581:THR:HG22	1:A:581:THR:O	2.05	0.56
1:A:652:ALA:HB2	1:A:678:TRP:NE1	2.21	0.55
1:A:553:ASN:HB2	3:A:1063:HOH:O	2.10	0.51
1:A:494:GLU:HG3	1:A:496:MET:HE2	1.91	0.50
1:A:591:GLU:OE1	1:A:595:TYR:OH	2.22	0.48
1:A:517:SER:HA	1:A:678:TRP:CZ3	2.49	0.47
1:A:494:GLU:HG3	1:A:496:MET:HE1	1.94	0.47
1:A:555:GLU:HG3	3:A:1063:HOH:O	2.15	0.47
1:A:613:THR:HG21	1:A:645:TRP:CE2	2.50	0.46
1:A:712:ARG:NH1	1:A:713:ASN:HD22	2.06	0.45
1:A:652:ALA:O	1:A:656:VAL:HG23	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:712:ARG:HD2	1:A:712:ARG:C	2.39	0.42
1:A:672:LEU:HD12	1:A:672:LEU:HA	1.93	0.41
1:A:468:LEU:HD23	1:A:479:LEU:HD23	2.01	0.41
1:A:511:PHE:HA	1:A:515[A]:GLU:OE2	2.20	0.41
1:A:593:GLN:HE21	1:A:593:GLN:HB3	1.68	0.41
1:A:655:LEU:HD22	1:A:659:MET:HG3	2.03	0.40
1:A:655:LEU:CD2	1:A:659:MET:HG3	2.51	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:A:515[A]:GLU:OE2	1:A:641:SER:OG[6_555]	2.07	0.13

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	A	299/342 (87%)	290 (97%)	9 (3%)	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	A	259/296 (88%)	245 (95%)	14 (5%)	22	18

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	426	GLU
1	A	432	SER
1	A	442	LYS
1	A	446	MET
1	A	501	LEU
1	A	509	LYS
1	A	546	LEU
1	A	567	GLN
1	A	590	LEU
1	A	638	PHE
1	A	641	SER
1	A	655	LEU
1	A	672	LEU
1	A	712	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	567	GLN
1	A	593	GLN
1	A	713	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 monosaccharides in this entry.

5.6 Ligand geometry

Of 1 ligands modelled in this entry, 1 is monoatomic - 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.

No monomer is involved in short contacts.

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 ⓘ

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	298/342 (87%)	1.03	49 (16%) ⓘ ⓘ	40, 52, 66, 71	13 (4%)

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	431	GLY	9.0
1	A	433	TYR	8.3
1	A	586	ALA	5.5
1	A	573	GLY	5.3
1	A	486	GLY	5.2
1	A	484	ASP	5.0
1	A	682	TRP	4.9
1	A	582	ALA	4.7
1	A	713	ASN	4.7
1	A	430	VAL	4.7
1	A	712	ARG	4.3
1	A	584	PHE	3.9
1	A	575	LEU	3.9
1	A	585	VAL	3.7
1	A	683	ASP	3.7
1	A	681	HIS	3.5
1	A	588	GLU	3.5
1	A	714	GLN	3.5
1	A	581	THR	3.4
1	A	636	GLY	3.4
1	A	492	VAL	3.4
1	A	591	GLU	3.2
1	A	420	ASP	3.2
1	A	571	GLU	3.2
1	A	432	SER	3.2
1	A	646	ASN	3.1
1	A	579[A]	CYS	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	450	VAL	2.9
1	A	487	LYS	2.9
1	A	587	PRO	2.8
1	A	606	VAL	2.8
1	A	417	GLN	2.8
1	A	665	HIS	2.7
1	A	590	LEU	2.7
1	A	684	GLN	2.7
1	A	564	PHE	2.7
1	A	688	TYR	2.7
1	A	595	TYR	2.5
1	A	490	TYR	2.5
1	A	607	LEU	2.4
1	A	580	TYR	2.4
1	A	610	THR	2.4
1	A	428	ILE	2.3
1	A	583	ASN	2.3
1	A	493	THR	2.2
1	A	589	VAL	2.1
1	A	637	LYS	2.1
1	A	570	ALA	2.1
1	A	685	LEU	2.1

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides 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. 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	B-factors(\AA^2)	Q<0.9
2	NA	A	1000	1/1	0.98	0.10	48,48,48,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.