

wwPDB X-ray Structure Validation Summary Report (i)

Dec 19, 2023 – 06:16 PM EST

PDB ID	:	1UKI
Title	:	Structural basis for the selective inhibition of JNK1 by the scaffolding protein
		JIP1 and SP600125
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		Kim, J.H.; Hyun, YL.; Jeon, Y.H.; Ro, S.; Lee, T.G.; Cho, J.M.; Hwang,
		K.Y.; Yang, CH.
Deposited on	:	2003-08-23
Resolution	:	2.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

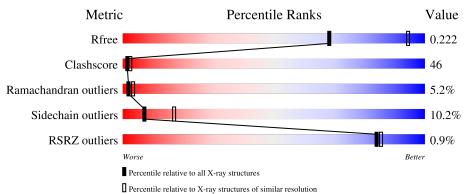
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.70 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 <=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	Q	uality of chain		
1	А	369	% 30%	51%	7% •	11%
2	В	11	45%	27%	18%	9%



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2763 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 mitogen-activated protein kinase 8 isoform 4.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	А	328	Total 2647	C 1705	N 446	0 477	S 19	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

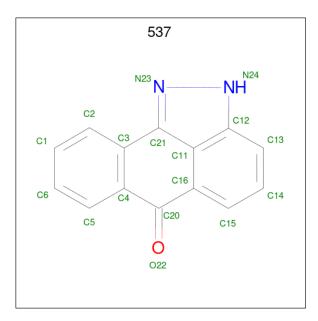
Chain	Residue	Modelled	Actual	Comment	Reference
А	364	HIS	-	expression tag	UNP P45983
А	365	HIS	-	expression tag	UNP P45983
А	366	HIS	-	expression tag	UNP P45983
A	367	HIS	-	expression tag	UNP P45983
А	368	HIS	-	expression tag	UNP P45983
А	369	HIS	-	expression tag	UNP P45983

• Molecule 2 is a protein called 11-mer peptide from C-jun-amino-terminal kinase interacting protein 1.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace		
2	В	10	Total 84	C 55	N 15	0 14	0	0	0

• Molecule 3 is 2,6-DIHYDROANTHRA/1,9-CD/PYRAZOL-6-ONE (three-letter code: 537) (formula: $C_{14}H_8N_2O$).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	Δ	1	Total	С	Ν	0	0	0
5	Л	1	17	14	2	1	0	0

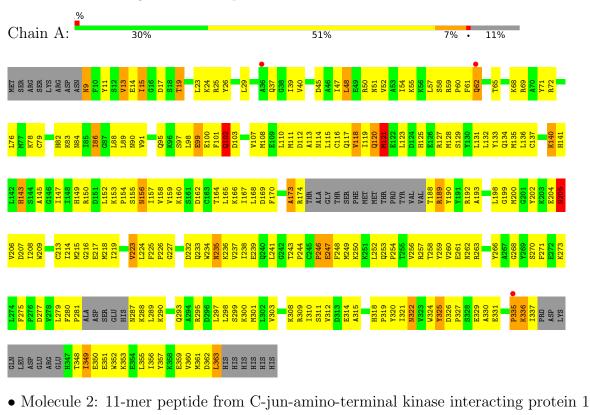
• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	15	Total O 15 15	0	0



3 Residue-property plots (i)

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: mitogen-activated protein kinase 8 isoform 4





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	61.70Å 79.46Å 82.86 Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.96 - 2.70	Depositor
Resolution (A)	19.96 - 2.70	EDS
% Data completeness	(Not available) $(19.96-2.70)$	Depositor
(in resolution range)	90.4(19.96-2.70)	EDS
R _{merge}	(Not available)	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$2.25 (at 2.71 \text{\AA})$	Xtriage
Refinement program	CNS 1.1	Depositor
P. P.	0.216 , 0.241	Depositor
R, R_{free}	0.219 , 0.222	DCC
R_{free} test set	552 reflections (5.23%)	wwPDB-VP
Wilson B-factor $(Å^2)$	48.0	Xtriage
Anisotropy	0.303	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.28 , 54.7	EDS
L-test for twinning ²	$< L > = 0.43, < L^2 > = 0.26$	Xtriage
Estimated twinning fraction	0.045 for -h,l,k	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2763	wwPDB-VP
Average B, all atoms $(Å^2)$	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.65% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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: $537\,$

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	Mol Chain		lengths	Bond angles		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.47	0/2704	0.65	0/3652	
2	В	0.52	0/86	0.80	0/114	
All	All	0.47	0/2790	0.66	0/3766	

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 (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	А	2647	0	2683	252	0
2	В	84	0	91	7	0
3	А	17	0	8	1	0
4	А	15	0	0	1	0
All	All	2763	0	2782	255	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 46.

The worst 5 of 255 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:258:THR:HG22	1:A:262:ASN:HD21	1.15	1.06	
1:A:128:MET:HG3	1:A:218:MET:HE3	1.41	1.02	
1:A:190:TYR:HA	1:A:233:GLN:HE22	1.24	0.99	
1:A:51:ASN:HB3	1:A:110:LEU:HD23	1.48	0.95	
1:A:91:VAL:HG22	1:A:360:VAL:HG13	1.49	0.95	

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	А	320/369~(87%)	245~(77%)	58 (18%)	17 (5%)	2	3
2	В	8/11 (73%)	8 (100%)	0	0	100	100
All	All	328/380~(86%)	253 (77%)	58 (18%)	17 (5%)	2	3

5 of 17 Ramachandran outliers are listed below:

Mol	Chain	\mathbf{Res}	Type
1	А	62	GLN
1	А	121	MET
1	А	189	ARG
1	А	226	PRO
1	А	173	ALA

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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	А	293/331~(88%)	266~(91%)	27~(9%)	9 21		
2	В	10/11 (91%)	6 (60%)	4 (40%)	0 0		
All	All	303/342~(89%)	272~(90%)	31 (10%)	7 17		

5 of 31 residues with a non-rotameric side chain are listed below:

Mol	Chain	Res	Type
1	А	156	ASN
2	В	158	THR
1	А	235	ASN
2	В	161	ASN
1	А	336	LYS

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 19 such side chains are listed below:

Mol	Chain	Res	Type
1	А	262	ASN
1	А	322	ASN
2	В	161	ASN
1	А	318	HIS
1	А	120	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

1 ligand is 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	Chain	Res	Tiple	Bond lengths			Bond angles		
				nes	Link	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2	
	3	537	А	0	-	16,20,20	2.59	9 (56%)	23,30,30	2.08	5 (21%)	

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
3	537	А	0	-	-	-	0/4/4/4

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	А	0	537	O22-C20	4.18	1.30	1.22
3	А	0	537	C14-C13	4.08	1.46	1.36
3	А	0	537	C15-C16	3.54	1.44	1.38
3	А	0	537	C5-C4	3.45	1.44	1.38
3	А	0	537	C14-C15	3.37	1.46	1.38

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	А	0	537	C15-C16-C20	7.39	125.91	119.61
3	А	0	537	C11-C16-C20	-3.48	117.05	119.92
3	А	0	537	C3-C21-C11	-2.42	120.15	121.92
3	А	0	537	C16-C11-C12	2.25	125.40	120.00
3	А	0	537	C5-C4-C20	2.05	121.36	119.61

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
3	А	0	537	1	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 (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, 95^{th} 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	$\mathbf{OWAB}(\mathbf{\AA}^2)$	Q < 0.9
1	А	328/369~(88%)	-0.27	3 (0%) 84 85	26, 47, 71, 83	0
2	В	10/11~(90%)	-0.26	0 100 100	42, 47, 55, 57	0
All	All	338/380~(88%)	-0.27	3 (0%) 84 85	26, 47, 71, 83	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	36	ALA	2.9
1	А	335	PRO	2.3
1	А	62	GLN	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, 95^{th} 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	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q < 0.9
3	537	А	0	17/17	0.94	0.14	37,41,43,44	0



6.5 Other polymers (i)

There are no such residues in this entry.

