



wwPDB X-ray Structure Validation Summary Report ⓘ

May 14, 2020 – 08:38 pm BST

PDB ID : 1ZWS
Title : Crystal structure of the catalytic domain of human DRP-1 kinase
Authors : Kursula, P.; Schunck, H.; Wilmanns, M.
Deposited on : 2005-06-06
Resolution : 2.90 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 ⓘ 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	:	2.11
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)	:	2.11

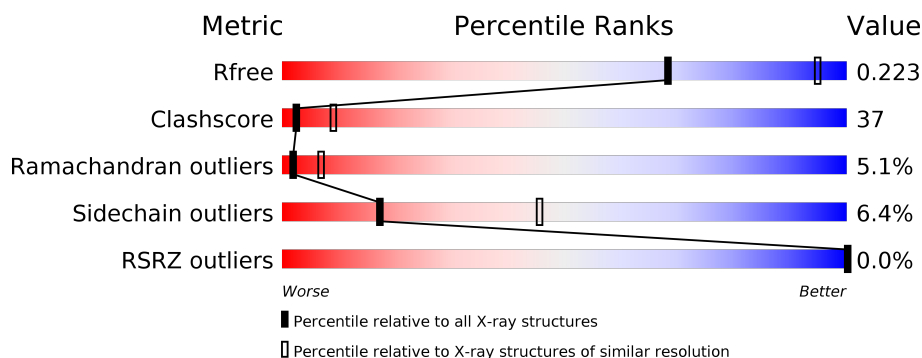
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 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	288	<div> <div></div> <div>38%50%8%.</div> </div>
1	B	288	<div> <div></div> <div>39%50%7%.</div> </div>
1	C	288	<div> <div></div> <div>37%51%8%.</div> </div>
1	D	288	<div> <div></div> <div>39%49%8%.</div> </div>
1	E	288	<div> <div></div> <div>38%51%7%.</div> </div>
1	F	288	<div> <div></div> <div>37%51%9%.</div> </div>

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Mol	Chain	Length	Quality of chain
1	G	288	<div><div></div><div>40%</div><div>49%</div><div>8%</div><div></div></div>
1	H	288	<div><div></div><div>40%</div><div>50%</div><div>7%</div><div></div></div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 18024 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 DAP-kinase related protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	278	Total	C	N	O	S	0	0	0
			2253	1448	375	427	3			
1	B	278	Total	C	N	O	S	0	0	0
			2253	1448	375	427	3			
1	C	278	Total	C	N	O	S	0	0	0
			2253	1448	375	427	3			
1	D	278	Total	C	N	O	S	0	0	0
			2253	1448	375	427	3			
1	E	278	Total	C	N	O	S	0	0	0
			2253	1448	375	427	3			
1	F	278	Total	C	N	O	S	0	0	0
			2253	1448	375	427	3			
1	G	278	Total	C	N	O	S	0	0	0
			2253	1448	375	427	3			
1	H	278	Total	C	N	O	S	0	0	0
			2253	1448	375	427	3			

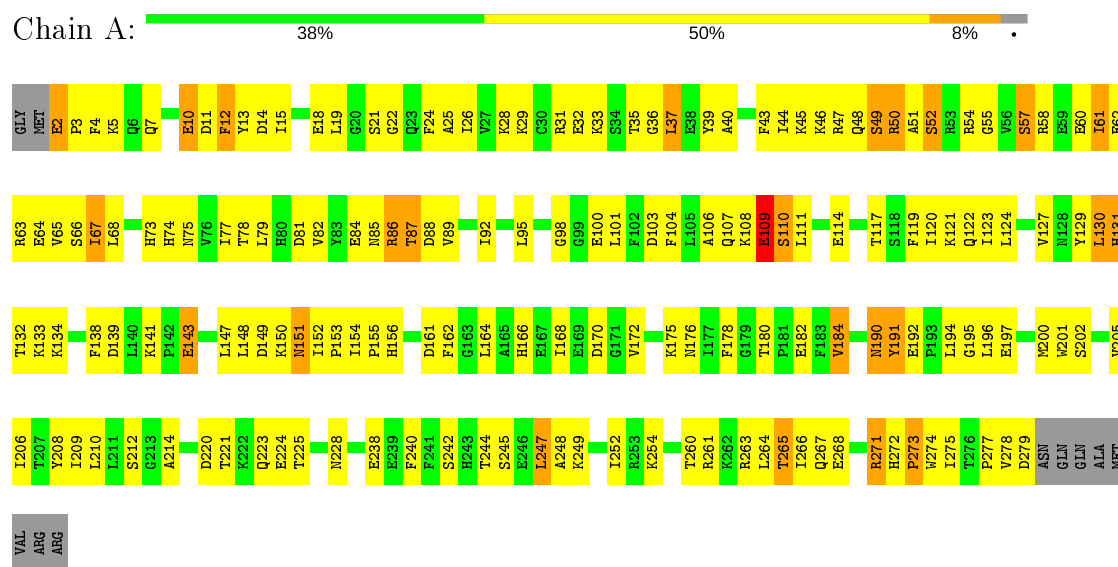
There are 8 discrepancies between the modelled and reference sequences:

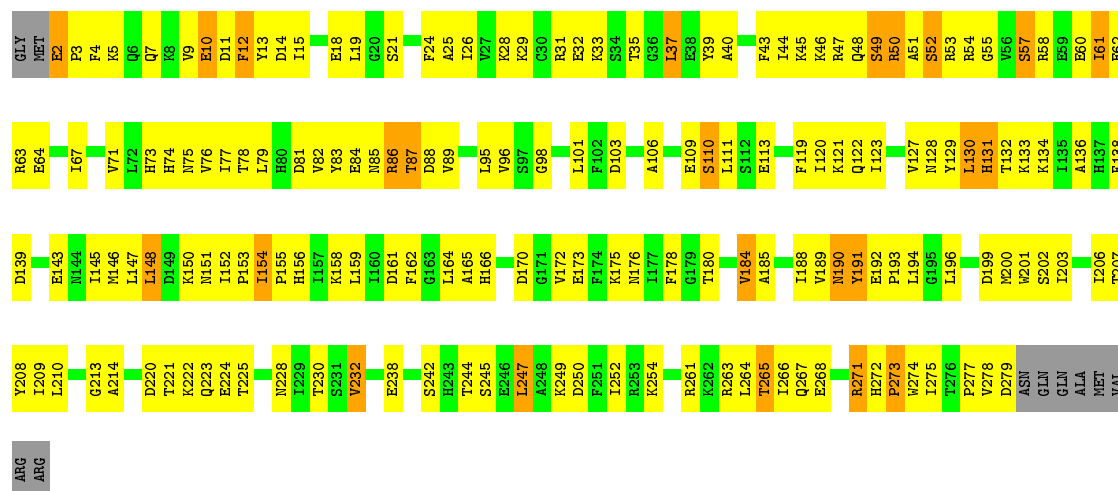
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	CLONING ARTIFACT	UNP Q9UIK4
B	0	GLY	-	CLONING ARTIFACT	UNP Q9UIK4
C	0	GLY	-	CLONING ARTIFACT	UNP Q9UIK4
D	0	GLY	-	CLONING ARTIFACT	UNP Q9UIK4
E	0	GLY	-	CLONING ARTIFACT	UNP Q9UIK4
F	0	GLY	-	CLONING ARTIFACT	UNP Q9UIK4
G	0	GLY	-	CLONING ARTIFACT	UNP Q9UIK4
H	0	GLY	-	CLONING ARTIFACT	UNP Q9UIK4

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 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: DAP-kinase related protein 1





E64	E67	V71	L72	H73	H74	N75	N76	I77	T78	D81	Y83	E84	N85	R86	T87	D88	I92	L95	V96	S97	G98	G99	D103	F104	L105	A106	Q107	K108	E109	S110	L111	E114	E115	A116	T117	S118	F119	I120	K121	Q122	I123	L124	V127	H128	Y129	L130	H131	T132	K133	K134									
GLY	MET	P3	F4	K5	Q6	Q7		E10	D11	F12	Y13	D14	I15			E18	L19	G20	S21		F24	A25	I26	V27	K28	K29	G30	R31	E32	K33	S34	T35	G36	L37	E38	F39	A40		F43	V44	K45	R46	K47	Q48	S49	R50	A51	S52	R53	R54	G55	V56	S57	R58	E59	E60	I61	P62	R63



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.62Å 211.83Å 128.42Å 90.00° 101.82° 90.00°	Depositor
Resolution (Å)	20.00 – 2.90 19.61 – 2.90	Depositor EDS
% Data completeness (in resolution range)	90.8 (20.00-2.90) 90.8 (19.61-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.28 (at 2.88Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.220 , 0.258 0.220 , 0.223	Depositor DCC
R_{free} test set	2751 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	52.3	Xtriage
Anisotropy	0.576	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 25.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.18$	Xtriage
Estimated twinning fraction	0.418 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	18024	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 4.24% 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](#)

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	0.52	0/2299	0.67	0/3104
1	B	0.52	1/2299 (0.0%)	0.67	0/3104
1	C	0.52	0/2299	0.68	0/3104
1	D	0.51	0/2299	0.68	0/3104
1	E	0.51	0/2299	0.66	0/3104
1	F	0.50	0/2299	0.68	0/3104
1	G	0.51	0/2299	0.68	0/3104
1	H	0.51	0/2299	0.67	0/3104
All	All	0.51	1/18392 (0.0%)	0.68	0/24832

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	30	CYS	CB-SG	-7.72	1.69	1.82

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	A	2253	0	2257	170	0
1	B	2253	0	2257	160	0
1	C	2253	0	2257	183	0
1	D	2253	0	2257	176	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2253	0	2257	181	1
1	F	2253	0	2257	174	0
1	G	2253	0	2257	165	0
1	H	2253	0	2257	163	0
All	All	18024	0	18056	1339	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 37.

The worst 5 of 1339 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:H:152:ILE:HD11	1:H:155:PRO:HA	1.39	1.03
1:C:152:ILE:HD11	1:C:155:PRO:HA	1.41	1.03
1:A:152:ILE:HD11	1:A:155:PRO:HA	1.39	1.03
1:F:152:ILE:HD11	1:F:155:PRO:HA	1.44	0.98
1:G:152:ILE:HD11	1:G:155:PRO:HA	1.48	0.94

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:E:38:GLU:OE1	1:E:259:GLU:OE1[1_655]	2.09	0.11

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	276/288 (96%)	213 (77%)	46 (17%)	17 (6%)	1 4
1	B	276/288 (96%)	215 (78%)	48 (17%)	13 (5%)	2 8

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	276/288 (96%)	210 (76%)	52 (19%)	14 (5%)	2	7
1	D	276/288 (96%)	213 (77%)	46 (17%)	17 (6%)	1	4
1	E	276/288 (96%)	218 (79%)	46 (17%)	12 (4%)	2	10
1	F	276/288 (96%)	214 (78%)	50 (18%)	12 (4%)	2	10
1	G	276/288 (96%)	212 (77%)	49 (18%)	15 (5%)	2	6
1	H	276/288 (96%)	210 (76%)	53 (19%)	13 (5%)	2	8
All	All	2208/2304 (96%)	1705 (77%)	390 (18%)	113 (5%)	2	7

5 of 113 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	50	ARG
1	B	50	ARG
1	B	51	ALA
1	B	52	SER
1	B	110	SER

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	249/257 (97%)	231 (93%)	18 (7%)	14	39
1	B	249/257 (97%)	235 (94%)	14 (6%)	21	52
1	C	249/257 (97%)	233 (94%)	16 (6%)	17	45
1	D	249/257 (97%)	235 (94%)	14 (6%)	21	52
1	E	249/257 (97%)	233 (94%)	16 (6%)	17	45
1	F	249/257 (97%)	231 (93%)	18 (7%)	14	39
1	G	249/257 (97%)	233 (94%)	16 (6%)	17	45
1	H	249/257 (97%)	234 (94%)	15 (6%)	19	49
All	All	1992/2056 (97%)	1865 (94%)	127 (6%)	17	45

5 of 127 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	190	ASN
1	E	168	ILE
1	H	87	THR
1	D	261	ARG
1	E	12	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 48 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	144	ASN
1	E	85	ASN
1	H	144	ASN
1	D	190	ASN
1	D	228	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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 ⓘ

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, 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	278/288 (96%)	-0.35	0 100 100	40, 40, 40, 40	0
1	B	278/288 (96%)	-0.21	0 100 100	40, 40, 40, 40	0
1	C	278/288 (96%)	-0.34	0 100 100	40, 40, 40, 40	0
1	D	278/288 (96%)	-0.33	0 100 100	40, 40, 40, 40	0
1	E	278/288 (96%)	-0.24	0 100 100	40, 40, 40, 40	0
1	F	278/288 (96%)	-0.33	1 (0%) 92 93	40, 40, 40, 40	0
1	G	278/288 (96%)	-0.38	0 100 100	40, 40, 40, 40	0
1	H	278/288 (96%)	-0.34	0 100 100	40, 40, 40, 40	0
All	All	2224/2304 (96%)	-0.31	1 (0%) 100 100	40, 40, 40, 40	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	279	ASP	2.6

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 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.