



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 12:55 am GMT

PDB ID : 1I0E
Title : CRYSTAL STRUCTURE OF CREATINE KINASE FROM HUMAN MUSCLE
Authors : Shen, Y.-Q.; Tang, L.; Zhou, H.-M.; Lin, Z.-J.
Deposited on : 2001-01-29
Resolution : 3.50 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

<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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
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)	:	recalc28949

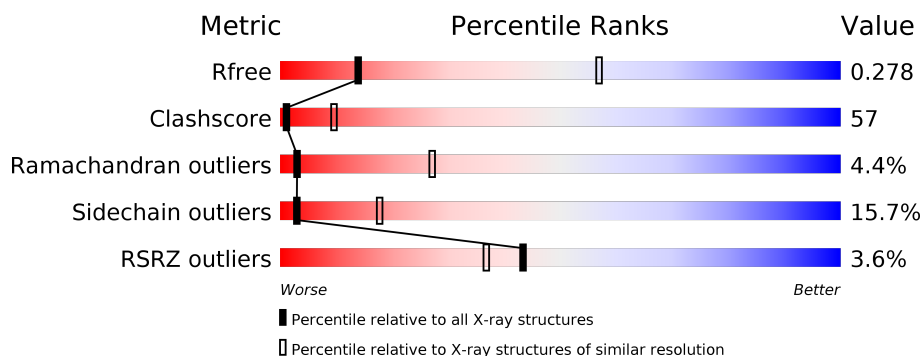
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1195 (3.60-3.40)
Clashscore	112137	1322 (3.60-3.40)
Ramachandran outliers	110173	1283 (3.60-3.40)
Sidechain outliers	110143	1284 (3.60-3.40)
RSRZ outliers	101464	1226 (3.60-3.40)

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	381	<div> <div>4%</div> <div> <div></div> <div>38%</div> <div>44%</div> <div>13%</div> <div>• •</div> </div> </div>
1	B	381	<div> <div>3%</div> <div> <div></div> <div>38%</div> <div>44%</div> <div>13%</div> <div>• •</div> </div> </div>
1	C	381	<div> <div>4%</div> <div> <div></div> <div>38%</div> <div>44%</div> <div>13%</div> <div>• •</div> </div> </div>
1	D	381	<div> <div>3%</div> <div> <div></div> <div>38%</div> <div>45%</div> <div>12%</div> <div>• •</div> </div> </div>

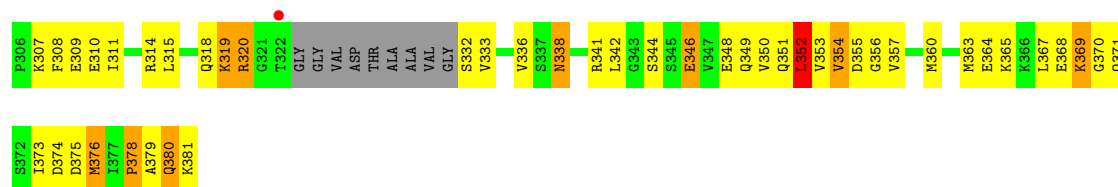
2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 11608 atoms, of which 0 are hydrogens and 0 are deuteriums.

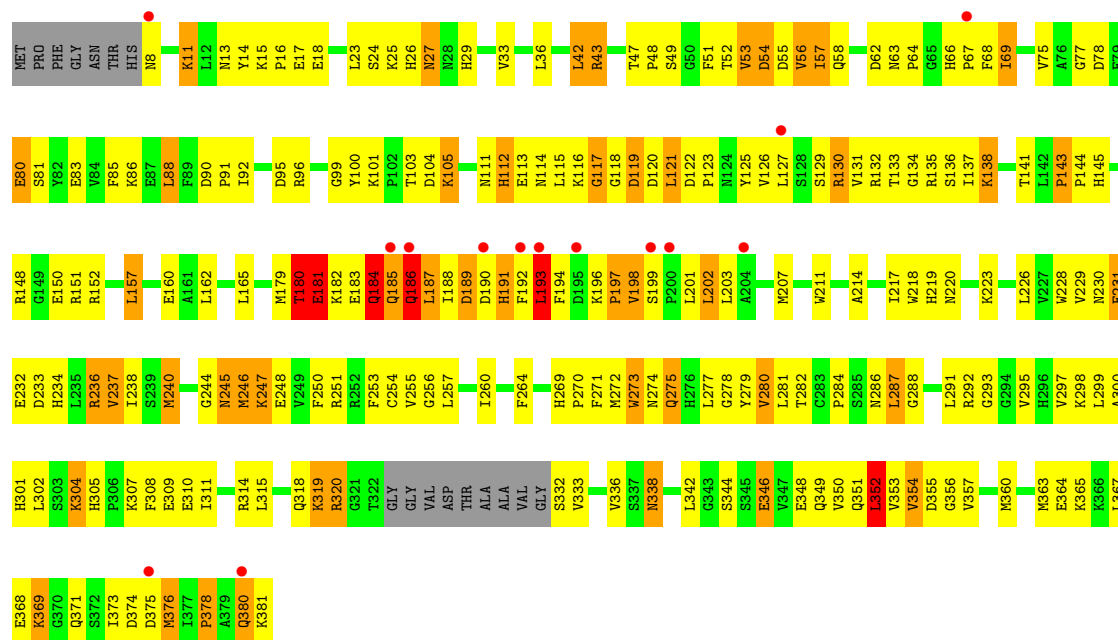
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called CREATINE KINASE,M CHAIN.

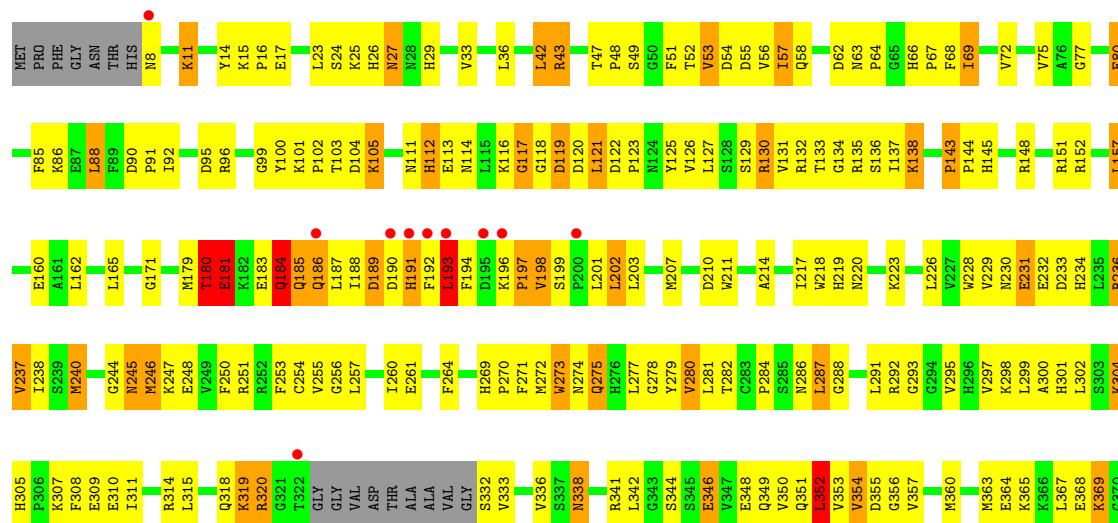
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	365	Total	C	N	O	S	0	0	0
			2902	1834	513	541	14			
1	B	365	Total	C	N	O	S	0	0	0
			2902	1834	513	541	14			
1	C	365	Total	C	N	O	S	0	0	0
			2902	1834	513	541	14			
1	D	365	Total	C	N	O	S	0	0	0
			2902	1834	513	541	14			



• Molecule 1: CREATINE KINASE,M CHAIN



• Molecule 1: CREATINE KINASE,M CHAIN



Q371	S372	I373	D374	D375	M376	I377	F378	A379	Q380	K381
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4 Data and refinement statistics

Property	Value	Source
Space group	P 31 1 2	Depositor
Cell constants a, b, c, α , β , γ	89.33Å 89.33Å 402.96Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 – 3.50 29.27 – 3.50	Depositor EDS
% Data completeness (in resolution range)	89.7 (8.00-3.50) 88.3 (29.27-3.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 3.47Å)	Xtriage
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.211 , 0.287 0.210 , 0.278	Depositor DCC
R_{free} test set	1915 reflections (9.83%)	DCC
Wilson B-factor (Å ²)	61.5	Xtriage
Anisotropy	0.214	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 84.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.052 for -h,-k,l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	11608	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.52% 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	0.40	0/2965	0.63	0/3992
1	B	0.46	0/2965	0.65	0/3992
1	C	0.42	0/2965	0.64	0/3992
1	D	0.43	0/2965	0.65	0/3992
All	All	0.43	0/11860	0.64	0/15968

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	A	2902	0	2869	325	0
1	B	2902	0	2869	333	0
1	C	2902	0	2869	327	1
1	D	2902	0	2869	326	1
All	All	11608	0	11476	1311	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 57.

The worst 5 of 1311 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:D:344:SER:HB2	1:D:349:GLN:HE21	1.12	1.14
1:B:105:LYS:HZ3	1:B:105:LYS:HA	1.09	1.14
1:C:344:SER:HB2	1:C:349:GLN:HE21	1.10	1.11
1:B:344:SER:HB2	1:B:349:GLN:HE21	1.11	1.10
1:B:181:GLU:HA	1:B:185:GLN:H	1.19	1.08

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:C:18:GLU:O	1:D:152:ARG:NH1[6_566]	1.99	0.21

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	361/381 (95%)	281 (78%)	64 (18%)	16 (4%)	3	27
1	B	361/381 (95%)	284 (79%)	61 (17%)	16 (4%)	3	27
1	C	361/381 (95%)	279 (77%)	66 (18%)	16 (4%)	3	27
1	D	361/381 (95%)	281 (78%)	64 (18%)	16 (4%)	3	27
All	All	1444/1524 (95%)	1125 (78%)	255 (18%)	64 (4%)	3	27

5 of 64 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	ASP
1	A	181	GLU
1	A	198	VAL
1	B	119	ASP
1	B	181	GLU

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	318/335 (95%)	268 (84%)	50 (16%)	3	18
1	B	318/335 (95%)	268 (84%)	50 (16%)	3	18
1	C	318/335 (95%)	267 (84%)	51 (16%)	3	16
1	D	318/335 (95%)	269 (85%)	49 (15%)	3	18
All	All	1272/1340 (95%)	1072 (84%)	200 (16%)	3	18

5 of 200 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	319	LYS
1	C	105	LYS
1	D	275	GLN
1	B	346	GLU
1	C	27	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 71 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	349	GLN
1	C	114	ASN
1	D	275	GLN
1	B	351	GLN
1	C	26	HIS

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 [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, 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	365/381 (95%)	-0.19	17 (4%) 32 26	6, 46, 93, 100	0
1	B	365/381 (95%)	-0.51	11 (3%) 51 42	2, 23, 85, 100	0
1	C	365/381 (95%)	-0.28	14 (3%) 41 35	5, 42, 93, 100	0
1	D	365/381 (95%)	-0.40	10 (2%) 55 46	2, 31, 83, 100	0
All	All	1460/1524 (95%)	-0.34	52 (3%) 43 37	2, 36, 91, 100	0

The worst 5 of 52 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	193	LEU	6.5
1	D	200	PRO	6.1
1	A	195	ASP	5.9
1	A	322	THR	5.8
1	A	200	PRO	5.8

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.