



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 09:35 PM GMT

PDB ID : 1PKL
Title : THE STRUCTURE OF LEISHMANIA PYRUVATE KINASE
Authors : Rigden, D.J.; Phillips, S.E.V.; Michels, P.A.M.; Fothergill-Gilmore, L.A.
Deposited on : 1998-09-15
Resolution : 2.35 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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

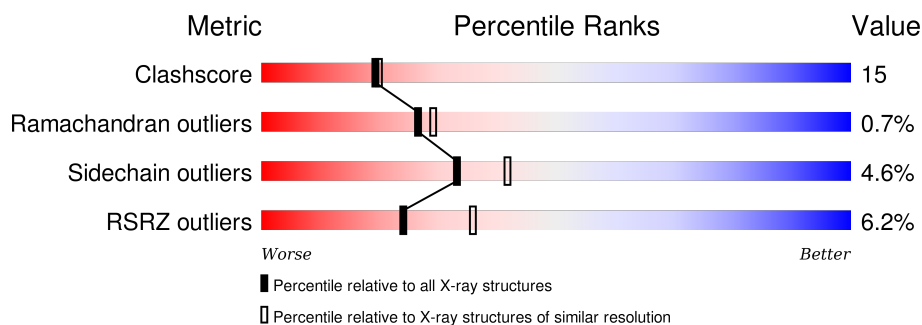
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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))
Clashscore	102246	1456 (2.38-2.34)
Ramachandran outliers	100387	1435 (2.38-2.34)
Sidechain outliers	100360	1436 (2.38-2.34)
RSRZ outliers	91569	1358 (2.38-2.34)

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	499	<div> <div>12%</div> <div>69%</div> <div>27%</div> <div>..</div> </div>
1	B	499	<div> <div>3%</div> <div>72%</div> <div>23%</div> <div>..</div> </div>
1	C	499	<div> <div>2%</div> <div>68%</div> <div>28%</div> <div>..</div> </div>
1	D	499	<div> <div>16%</div> <div>69%</div> <div>28%</div> <div>..</div> </div>
1	E	499	<div> <div>7%</div> <div>70%</div> <div>27%</div> <div>..</div> </div>
1	F	499	<div> <div>2%</div> <div>71%</div> <div>25%</div> <div>..</div> </div>
1	G	499	<div> <div>3%</div> <div>76%</div> <div>22%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
1	H	499	<div><div></div><div>3%</div><div>72%</div><div>25%</div><div>..</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 32986 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 PROTEIN (PYRUVATE KINASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	491	Total	C	N	O	S	0	0	0
			3740	2330	659	725	26			
1	B	489	Total	C	N	O	S	0	0	0
			3725	2321	655	723	26			
1	C	492	Total	C	N	O	S	0	0	0
			3748	2334	660	728	26			
1	D	493	Total	C	N	O	S	0	0	0
			3758	2340	663	729	26			
1	E	493	Total	C	N	O	S	0	0	0
			3758	2340	663	729	26			
1	F	491	Total	C	N	O	S	0	0	0
			3740	2330	659	725	26			
1	H	492	Total	C	N	O	S	0	0	0
			3748	2334	660	728	26			
1	G	498	Total	C	N	O	S	0	0	0
			3777	2351	666	734	26			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	382	SER	GLY	CONFLICT	UNP Q27686
A	389	TYR	SER	CONFLICT	UNP Q27686
A	404	ARG	ALA	CONFLICT	UNP Q27686
A	405	SER	GLY	CONFLICT	UNP Q27686
B	382	SER	GLY	CONFLICT	UNP Q27686
B	389	TYR	SER	CONFLICT	UNP Q27686
B	404	ARG	ALA	CONFLICT	UNP Q27686
B	405	SER	GLY	CONFLICT	UNP Q27686
C	382	SER	GLY	CONFLICT	UNP Q27686
C	389	TYR	SER	CONFLICT	UNP Q27686
C	404	ARG	ALA	CONFLICT	UNP Q27686
C	405	SER	GLY	CONFLICT	UNP Q27686
D	382	SER	GLY	CONFLICT	UNP Q27686

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Chain	Residue	Modelled	Actual	Comment	Reference
D	389	TYR	SER	CONFLICT	UNP Q27686
D	404	ARG	ALA	CONFLICT	UNP Q27686
D	405	SER	GLY	CONFLICT	UNP Q27686
E	382	SER	GLY	CONFLICT	UNP Q27686
E	389	TYR	SER	CONFLICT	UNP Q27686
E	404	ARG	ALA	CONFLICT	UNP Q27686
E	405	SER	GLY	CONFLICT	UNP Q27686
F	382	SER	GLY	CONFLICT	UNP Q27686
F	389	TYR	SER	CONFLICT	UNP Q27686
F	404	ARG	ALA	CONFLICT	UNP Q27686
F	405	SER	GLY	CONFLICT	UNP Q27686
H	382	SER	GLY	CONFLICT	UNP Q27686
H	389	TYR	SER	CONFLICT	UNP Q27686
H	404	ARG	ALA	CONFLICT	UNP Q27686
H	405	SER	GLY	CONFLICT	UNP Q27686
G	382	SER	GLY	CONFLICT	UNP Q27686
G	389	TYR	SER	CONFLICT	UNP Q27686
G	404	ARG	ALA	CONFLICT	UNP Q27686
G	405	SER	GLY	CONFLICT	UNP Q27686

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		

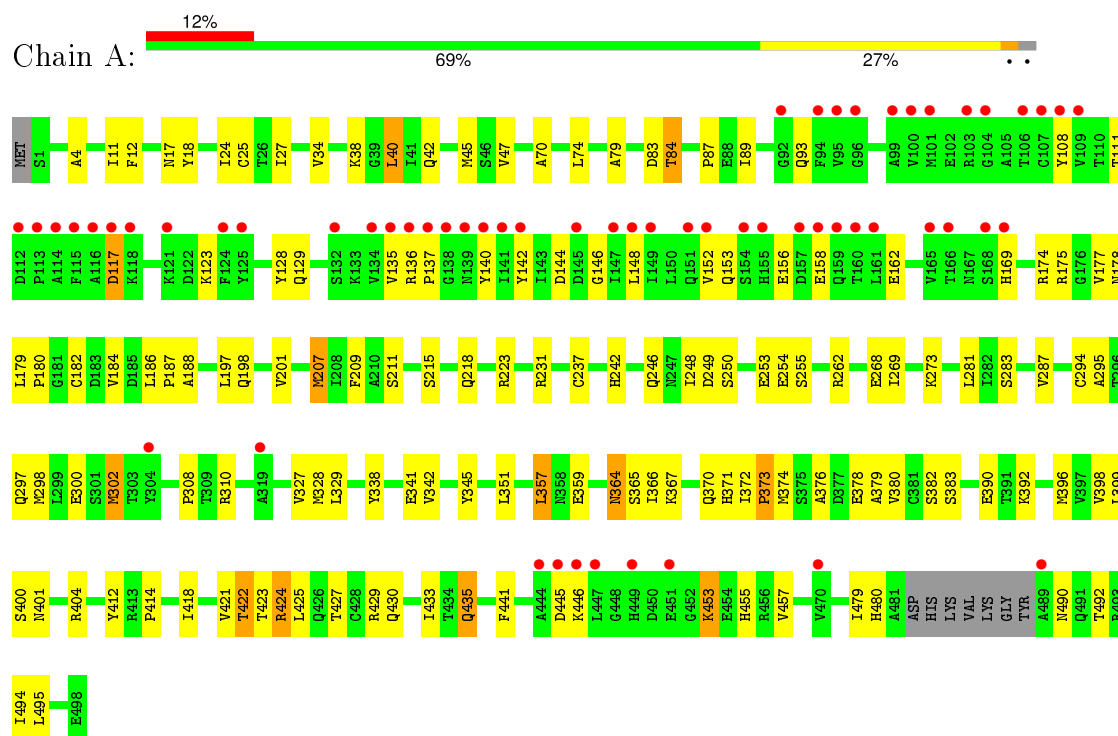
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	315	Total	O	0	0
			315	315		
3	B	366	Total	O	0	0
			366	366		
3	C	453	Total	O	0	0
			453	453		
3	D	244	Total	O	0	0
			244	244		
3	E	386	Total	O	0	0
			386	386		
3	F	399	Total	O	0	0
			399	399		
3	G	389	Total	O	0	0
			389	389		
3	H	400	Total	O	0	0
			400	400		

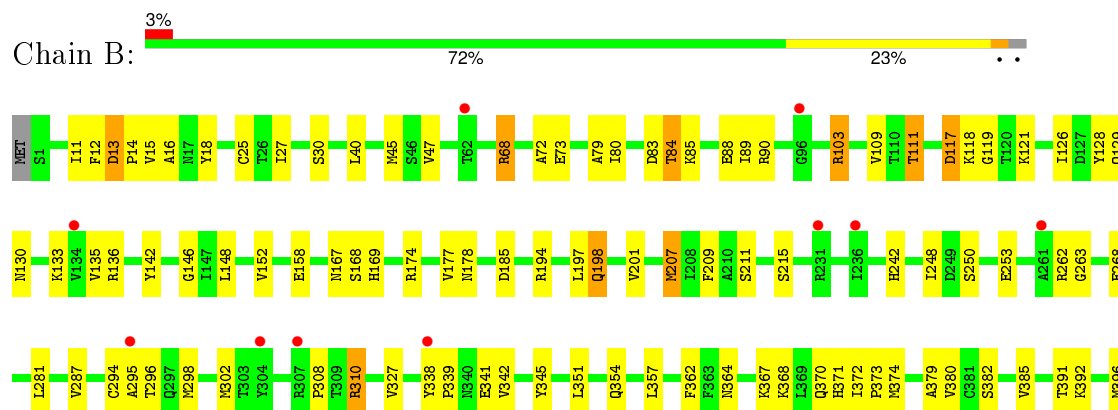
3 Residue-property plots

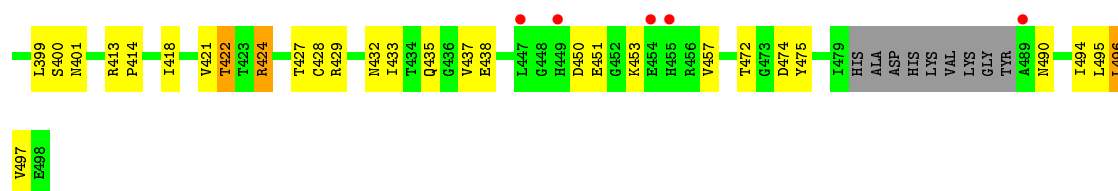
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: PROTEIN (PYRUVATE KINASE)

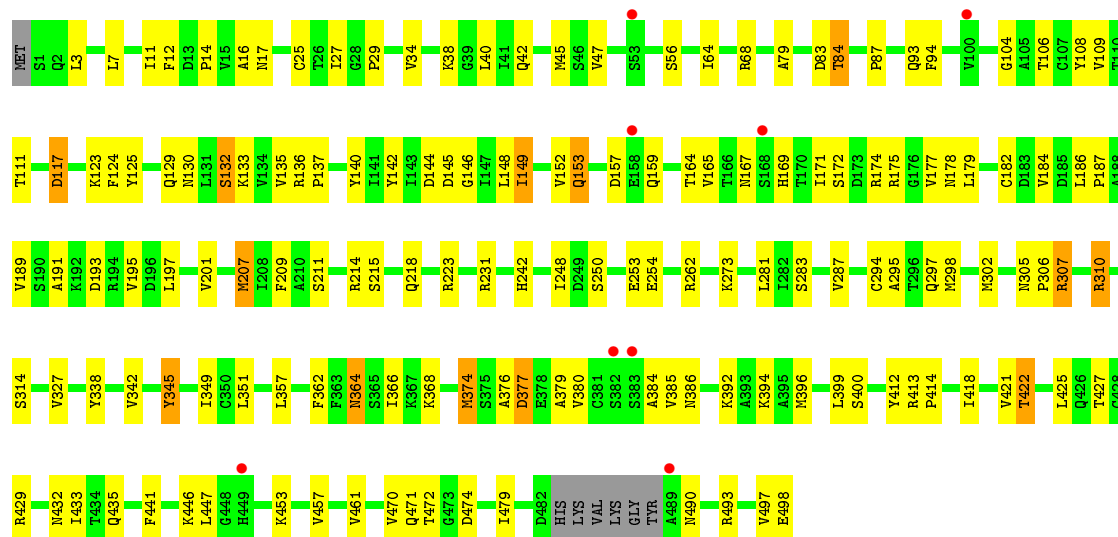


• Molecule 1: PROTEIN (PYRUVATE KINASE)

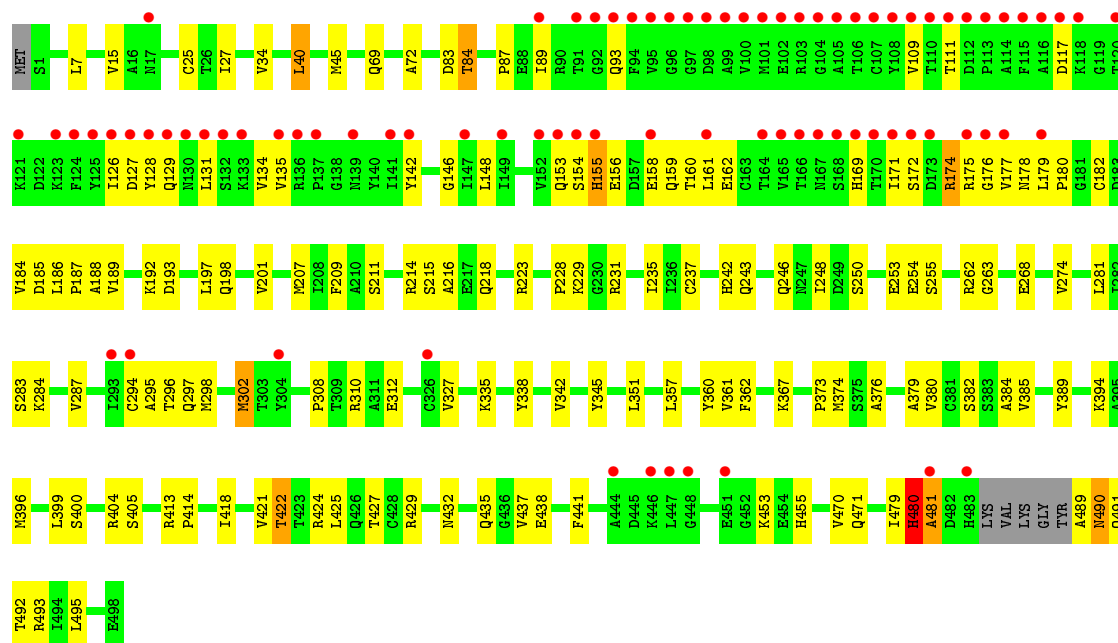




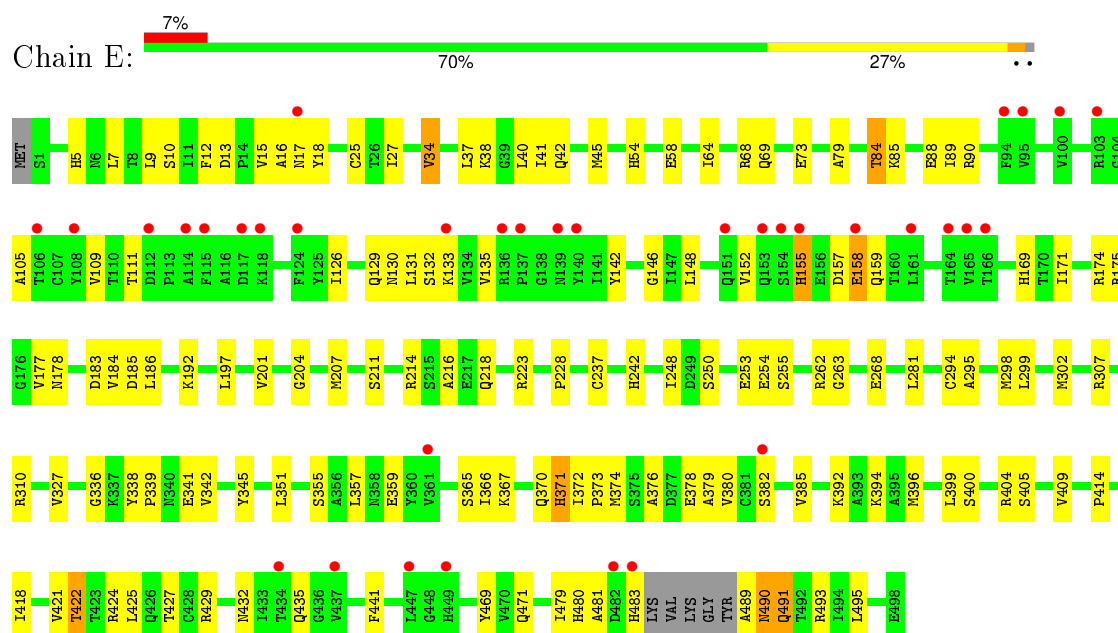
• Molecule 1: PROTEIN (PYRUVATE KINASE)



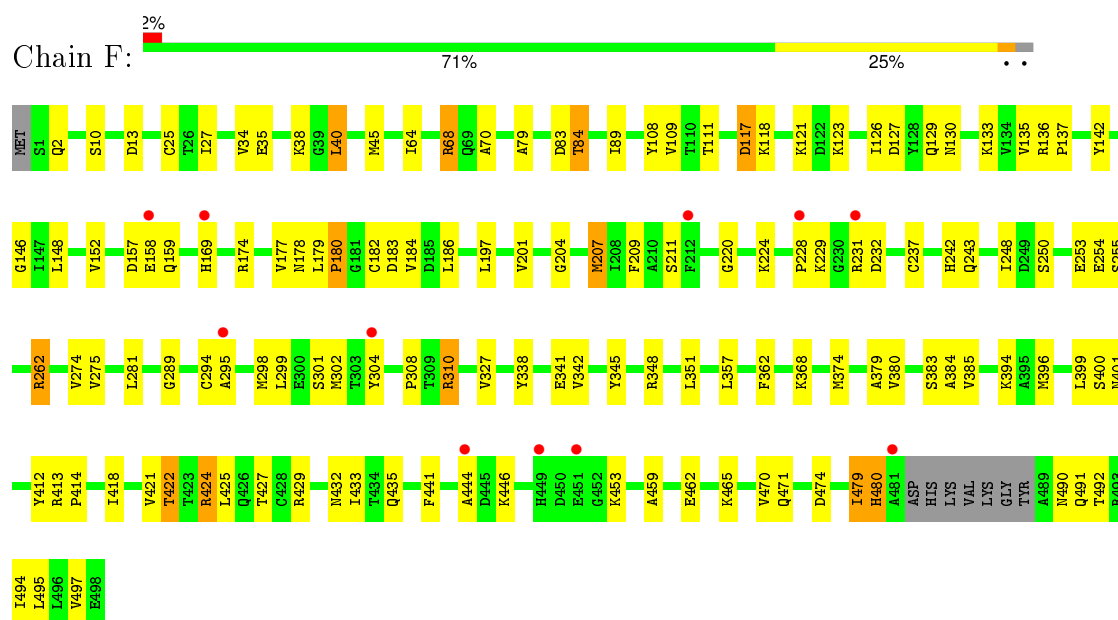
• Molecule 1: PROTEIN (PYRUVATE KINASE)



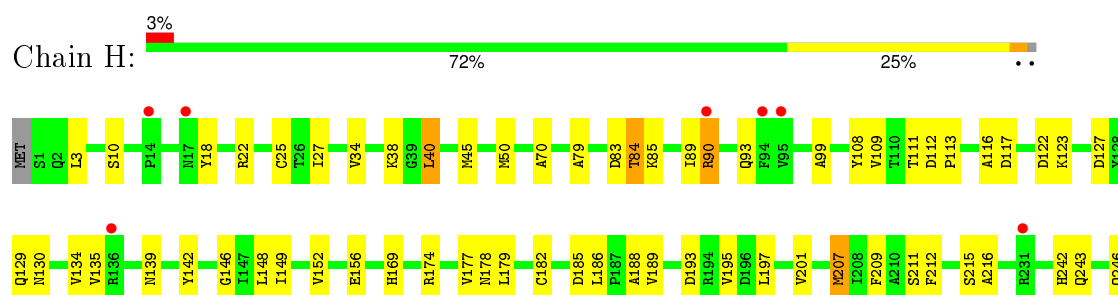
• Molecule 1: PROTEIN (PYRUVATE KINASE)

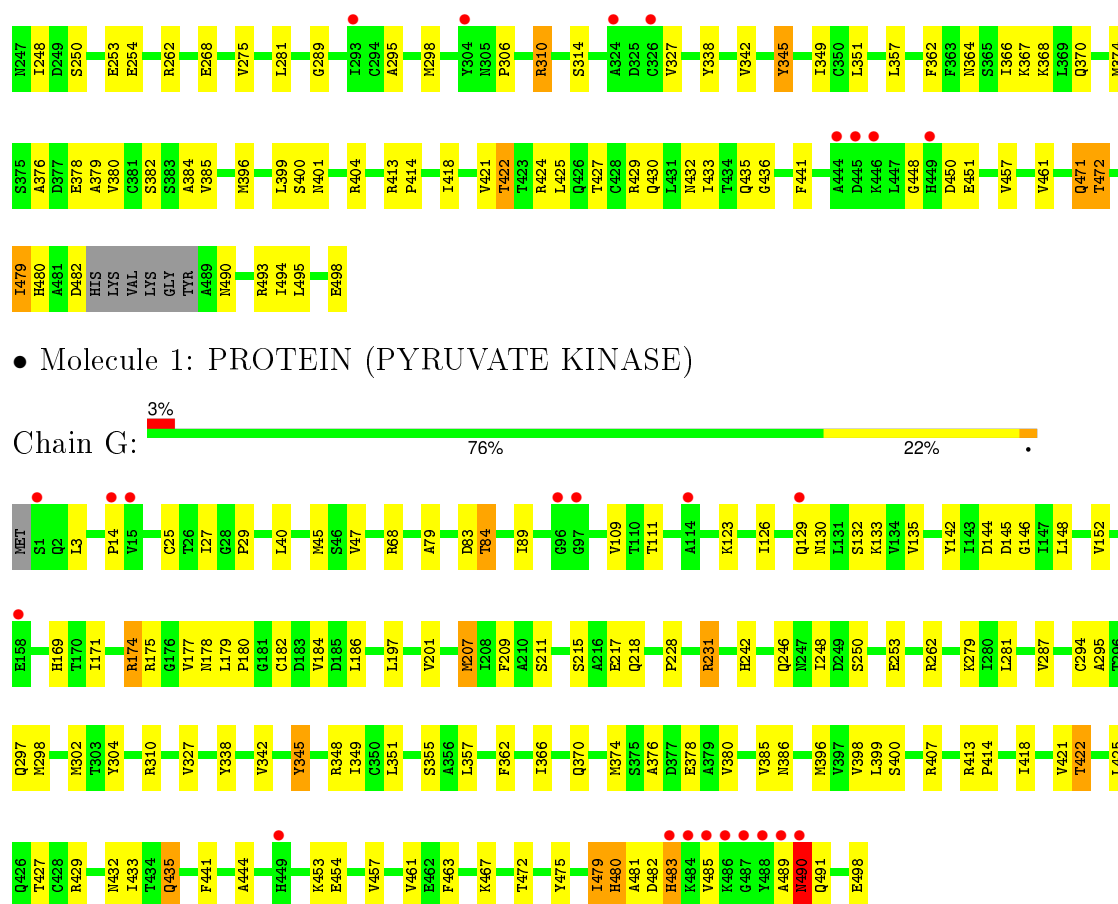


• Molecule 1: PROTEIN (PYRUVATE KINASE)

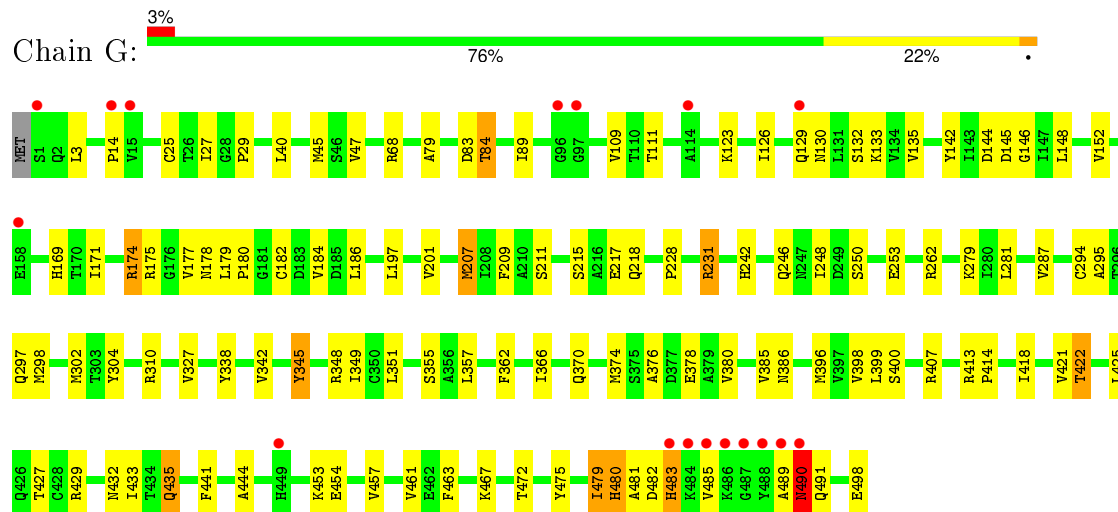


• Molecule 1: PROTEIN (PYRUVATE KINASE)





• Molecule 1: PROTEIN (PYRUVATE KINASE)



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	121.65Å 132.64Å 181.00Å 90.00° 97.11° 90.00°	Depositor
Resolution (Å)	30.00 – 2.35 29.82 – 2.35	Depositor EDS
% Data completeness (in resolution range)	67.9 (30.00-2.35) 67.9 (29.82-2.35)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.209 , 0.256 0.224 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	36.9	Xtriage
Anisotropy	0.576	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 45.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ¹	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 160774 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	32986	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 34.05 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 7.2506e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: SO4

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	0.54	2/3794 (0.1%)	0.70	4/5136 (0.1%)
1	B	0.43	0/3778	0.67	2/5114 (0.0%)
1	C	0.47	0/3802	0.68	1/5147 (0.0%)
1	D	0.45	0/3813	0.68	3/5162 (0.1%)
1	E	0.46	0/3813	0.68	1/5162 (0.0%)
1	F	0.46	0/3794	0.67	1/5136 (0.0%)
1	G	0.47	0/3832	0.68	1/5190 (0.0%)
1	H	0.45	0/3802	0.66	1/5147 (0.0%)
All	All	0.47	2/30428 (0.0%)	0.68	14/41194 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	E	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	373	PRO	N-CD	13.84	1.67	1.47
1	A	372	ILE	C-N	-7.04	1.20	1.34

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	373	PRO	CA-N-CD	-7.86	100.49	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	373	PRO	N-CA-CB	7.77	112.62	103.30
1	C	295	ALA	N-CA-C	7.47	131.17	111.00
1	D	295	ALA	N-CA-C	7.36	130.87	111.00
1	E	295	ALA	N-CA-C	7.34	130.83	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	18	TYR	Sidechain
1	E	18	TYR	Sidechain

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	3740	0	3743	121	0
1	B	3725	0	3731	110	0
1	C	3748	0	3747	132	0
1	D	3758	0	3754	115	0
1	E	3758	0	3754	126	0
1	F	3740	0	3743	113	0
1	G	3777	0	3761	106	0
1	H	3748	0	3747	116	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
2	E	5	0	0	0	0
2	F	5	0	0	0	0
2	G	5	0	0	0	0
2	H	5	0	0	0	0
3	A	315	0	0	30	0
3	B	366	0	0	29	0
3	C	453	0	0	38	0
3	D	244	0	0	18	0
3	E	386	0	0	38	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	399	0	0	31	0
3	G	389	0	0	24	0
3	H	400	0	0	37	0
All	All	32986	0	29980	903	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 15.

The worst 5 of 903 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:424:ARG:HB3	1:B:424:ARG:HH11	1.27	0.95
1:D:27:ILE:HD11	1:D:45:MET:HE1	1.51	0.91
1:A:490:ASN:HD22	1:B:494:ILE:HB	1.34	0.91
1:G:380:VAL:HG12	1:G:479:ILE:HD11	1.52	0.91
1:E:310:ARG:HH11	1:G:297:GLN:HE22	0.96	0.90

There are no symmetry-related clashes.

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	487/499 (98%)	458 (94%)	26 (5%)	3 (1%)	30	34
1	B	485/499 (97%)	457 (94%)	26 (5%)	2 (0%)	39	46
1	C	488/499 (98%)	463 (95%)	23 (5%)	2 (0%)	39	46
1	D	489/499 (98%)	455 (93%)	30 (6%)	4 (1%)	24	26
1	E	489/499 (98%)	460 (94%)	23 (5%)	6 (1%)	16	15
1	F	487/499 (98%)	462 (95%)	23 (5%)	2 (0%)	39	46
1	G	496/499 (99%)	466 (94%)	26 (5%)	4 (1%)	24	26

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	488/499 (98%)	466 (96%)	19 (4%)	3 (1%)	30	34
All	All	3909/3992 (98%)	3687 (94%)	196 (5%)	26 (1%)	26	29

5 of 26 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	483	HIS
1	G	485	VAL
1	G	490	ASN
1	A	174	ARG
1	C	174	ARG

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	410/417 (98%)	391 (95%)	19 (5%)	33	42
1	B	409/417 (98%)	386 (94%)	23 (6%)	26	31
1	C	411/417 (99%)	390 (95%)	21 (5%)	29	36
1	D	412/417 (99%)	397 (96%)	15 (4%)	42	55
1	E	412/417 (99%)	393 (95%)	19 (5%)	33	42
1	F	410/417 (98%)	393 (96%)	17 (4%)	37	48
1	G	411/417 (99%)	393 (96%)	18 (4%)	35	44
1	H	411/417 (99%)	393 (96%)	18 (4%)	35	44
All	All	3286/3336 (98%)	3136 (95%)	150 (5%)	33	42

5 of 150 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	158	GLU
1	E	155	HIS
1	G	262	ARG
1	D	262	ARG

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Mol	Chain	Res	Type
1	D	435	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 67 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	401	ASN
1	E	286	ASN
1	G	151	GLN
1	D	491	GLN
1	E	198	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

8 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	SO4	A	3000	-	4,4,4	0.22	0	6,6,6	0.24	0
2	SO4	B	3001	-	4,4,4	0.13	0	6,6,6	0.08	0
2	SO4	C	3002	-	4,4,4	0.35	0	6,6,6	0.17	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	D	3003	-	4,4,4	0.29	0	6,6,6	0.41	0
2	SO4	E	3004	-	4,4,4	0.44	0	6,6,6	0.16	0
2	SO4	F	3005	-	4,4,4	0.25	0	6,6,6	0.26	0
2	SO4	G	3007	-	4,4,4	0.21	0	6,6,6	0.10	0
2	SO4	H	3006	-	4,4,4	0.25	0	6,6,6	0.34	0

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	SO4	A	3000	-	-	0/0/0/0	0/0/0/0
2	SO4	B	3001	-	-	0/0/0/0	0/0/0/0
2	SO4	C	3002	-	-	0/0/0/0	0/0/0/0
2	SO4	D	3003	-	-	0/0/0/0	0/0/0/0
2	SO4	E	3004	-	-	0/0/0/0	0/0/0/0
2	SO4	F	3005	-	-	0/0/0/0	0/0/0/0
2	SO4	G	3007	-	-	0/0/0/0	0/0/0/0
2	SO4	H	3006	-	-	0/0/0/0	0/0/0/0

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 [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	491/499 (98%)	0.37	60 (12%) 5 10	15, 39, 100, 100	0
1	B	489/499 (97%)	0.10	15 (3%) 52 65	21, 48, 86, 100	0
1	C	492/499 (98%)	-0.07	8 (1%) 74 84	14, 38, 77, 100	0
1	D	493/499 (98%)	0.75	82 (16%) 2 4	19, 48, 100, 100	0
1	E	493/499 (98%)	0.17	35 (7%) 19 28	15, 36, 100, 100	0
1	F	491/499 (98%)	0.06	11 (2%) 65 77	19, 45, 87, 100	0
1	G	498/499 (99%)	-0.04	17 (3%) 49 62	12, 40, 91, 100	0
1	H	492/499 (98%)	0.00	15 (3%) 54 66	16, 40, 81, 100	0
All	All	3939/3992 (98%)	0.17	243 (6%) 24 36	12, 42, 100, 100	0

The worst 5 of 243 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	96	GLY	12.6
1	D	137	PRO	12.1
1	D	95	VAL	10.9
1	G	488	TYR	9.4
1	D	100	VAL	7.9

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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

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(Å ²)	Q<0.9
2	SO4	E	3004	5/5	0.98	0.08	-0.99	39,39,39,39	0
2	SO4	D	3003	5/5	0.98	0.08	-1.24	44,44,44,44	0
2	SO4	G	3007	5/5	0.99	0.06	-1.28	33,33,33,33	0
2	SO4	F	3005	5/5	0.99	0.08	-1.48	42,42,42,42	0
2	SO4	H	3006	5/5	0.99	0.07	-1.53	41,41,41,41	0
2	SO4	A	3000	5/5	0.98	0.08	-1.84	36,36,36,36	0
2	SO4	C	3002	5/5	0.99	0.07	-1.87	36,36,36,36	0
2	SO4	B	3001	5/5	0.99	0.07	-2.07	45,45,45,45	0

6.5 Other polymers

There are no such residues in this entry.