



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:40 PM GMT

PDB ID : 1Y8P
Title : Crystal structure of the PDK3-L2 complex
Authors : Kato, M.; Chuang, J.L.; Wynn, R.M.; Chuang, D.T.
Deposited on : 2004-12-13
Resolution : 2.63 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
<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

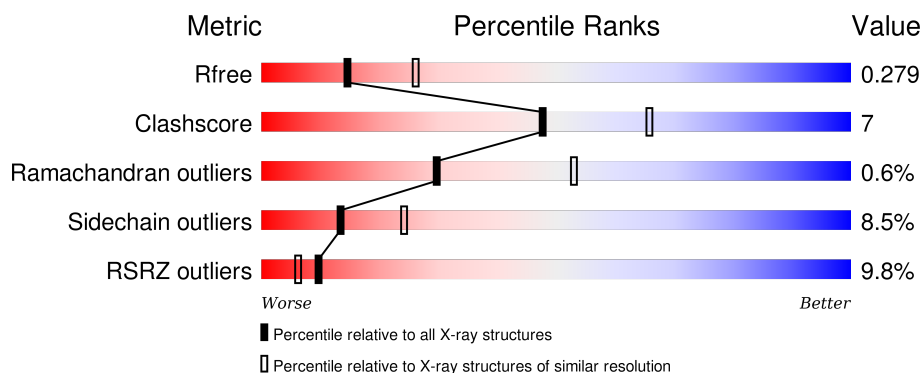
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.63 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}	91344	3377 (2.68-2.60)
Clashscore	102246	3781 (2.68-2.60)
Ramachandran outliers	100387	3722 (2.68-2.60)
Sidechain outliers	100360	3722 (2.68-2.60)
RSRZ outliers	91569	3388 (2.68-2.60)

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	419	
2	B	128	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	K	A	503	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 3949 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 [Pyruvate dehydrogenase [lipoamide]] kinase isozyme 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	381	3112	2005	520	574	13	0	1	0

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-12	GLY	-	CLONING ARTIFACT	UNP Q15120
A	-11	GLY	-	CLONING ARTIFACT	UNP Q15120
A	-10	SER	-	CLONING ARTIFACT	UNP Q15120
A	-9	HIS	-	EXPRESSION TAG	UNP Q15120
A	-8	HIS	-	EXPRESSION TAG	UNP Q15120
A	-7	HIS	-	EXPRESSION TAG	UNP Q15120
A	-6	HIS	-	EXPRESSION TAG	UNP Q15120
A	-5	HIS	-	EXPRESSION TAG	UNP Q15120
A	-4	HIS	-	EXPRESSION TAG	UNP Q15120
A	-3	GLY	-	CLONING ARTIFACT	UNP Q15120
A	-2	MET	-	CLONING ARTIFACT	UNP Q15120
A	-1	ALA	-	CLONING ARTIFACT	UNP Q15120
A	0	ARG	-	CLONING ARTIFACT	UNP Q15120
A	1	LEU	-	CLONING ARTIFACT	UNP Q15120
A	2	GLU	-	CLONING ARTIFACT	UNP Q15120
A	3	ASN	-	CLONING ARTIFACT	UNP Q15120
A	4	LEU	-	CLONING ARTIFACT	UNP Q15120
A	5	TYR	-	CLONING ARTIFACT	UNP Q15120
A	6	PHE	-	CLONING ARTIFACT	UNP Q15120
A	7	GLN	-	CLONING ARTIFACT	UNP Q15120
A	8	GLY	-	CLONING ARTIFACT	UNP Q15120

- Molecule 2 is a protein called Dihydrolipoyllysine-residue acetyltransferase component of pyruvate dehydrogenase complex.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	97	Total	C	N	O	S	0	0	0
			736	471	115	146	4			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	106	GLY	-	CLONING ARTIFACT	UNP P10515
B	107	GLY	-	CLONING ARTIFACT	UNP P10515
B	108	SER	-	CLONING ARTIFACT	UNP P10515
B	109	HIS	-	EXPRESSION TAG	UNP P10515
B	110	HIS	-	EXPRESSION TAG	UNP P10515
B	111	HIS	-	EXPRESSION TAG	UNP P10515
B	112	HIS	-	EXPRESSION TAG	UNP P10515
B	113	HIS	-	EXPRESSION TAG	UNP P10515
B	114	HIS	-	EXPRESSION TAG	UNP P10515
B	115	GLY	-	CLONING ARTIFACT	UNP P10515
B	116	MET	-	CLONING ARTIFACT	UNP P10515
B	117	ALA	-	CLONING ARTIFACT	UNP P10515
B	118	ARG	-	CLONING ARTIFACT	UNP P10515
B	119	LEU	-	CLONING ARTIFACT	UNP P10515
B	120	GLU	-	CLONING ARTIFACT	UNP P10515
B	121	ASN	-	CLONING ARTIFACT	UNP P10515
B	122	LEU	-	CLONING ARTIFACT	UNP P10515
B	123	TYR	-	CLONING ARTIFACT	UNP P10515
B	124	PHE	-	CLONING ARTIFACT	UNP P10515
B	125	GLN	-	CLONING ARTIFACT	UNP P10515

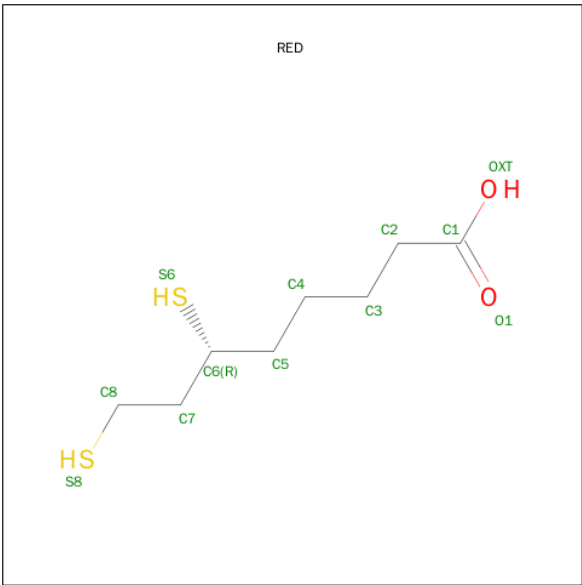
- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

- Molecule 4 is POTASSIUM ION (three-letter code: K) (formula: K).

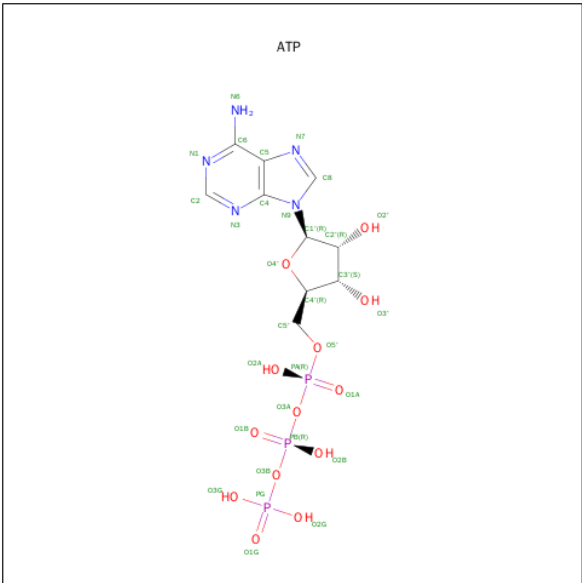
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	K	0	0
			2	2		

- Molecule 5 is DIHYDROLIPOIC ACID (three-letter code: RED) (formula: C₈H₁₆O₂S₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	O	S	0	0
			11	8	1	2		

- Molecule 6 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

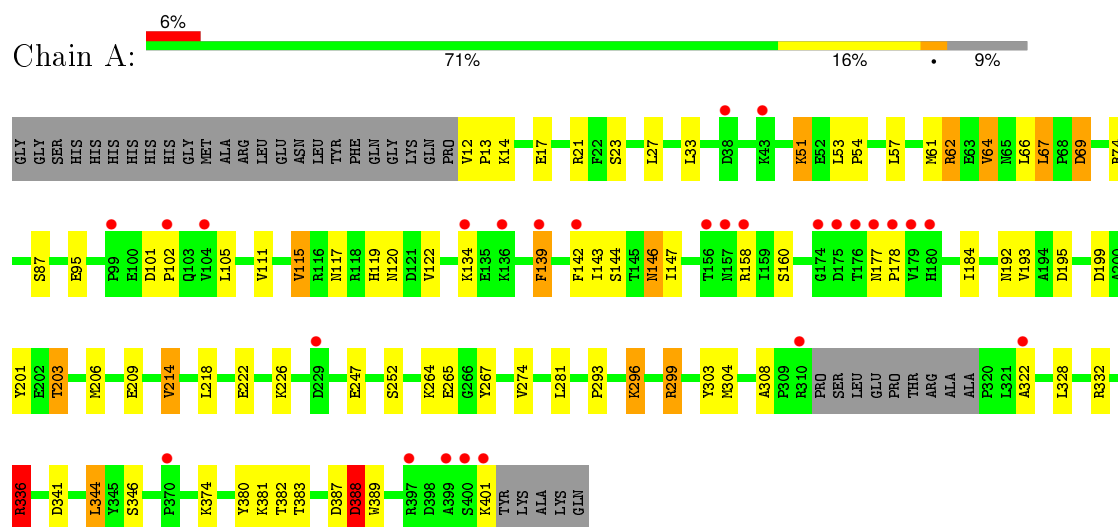
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	53	Total 53	O 53	0	0
7	B	3	Total 3	O 3	0	0

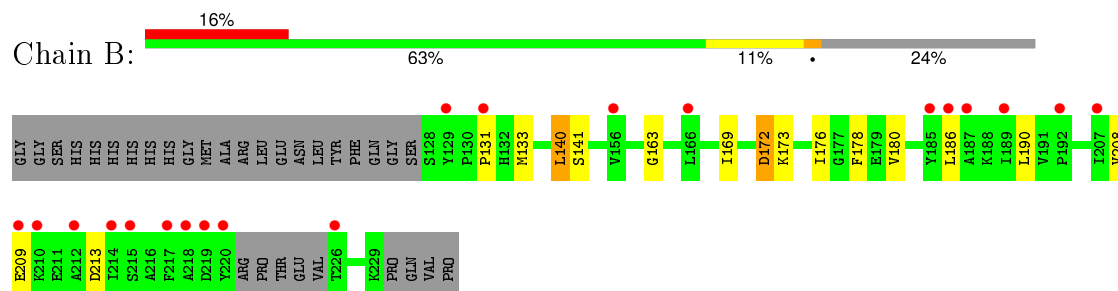
3 Residue-property plots [i](#)

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: [Pyruvate dehydrogenase [lipoamide]] kinase isozyme 3



- Molecule 2: Dihydrolipoyllysine-residue acetyltransferase component of pyruvate dehydrogenase complex



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	120.90 Å 120.90 Å 240.09 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.63 48.24 – 2.63	Depositor EDS
% Data completeness (in resolution range)	98.3 (50.00-2.63) 98.4 (48.24-2.63)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 2.65 Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.204 , 0.230 0.258 , 0.279	Depositor DCC
R_{free} test set	1566 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	68.5	Xtriage
Anisotropy	0.135	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 42.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	2 of 31043 reflections (0.006%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3949	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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, 0.375 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: MG, K, ATP, RED

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	0.74	0/3193	0.81	4/4326 (0.1%)
2	B	0.68	0/749	0.76	2/1017 (0.2%)
All	All	0.73	0/3942	0.80	6/5343 (0.1%)

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

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	336	ARG	NE-CZ-NH1	7.41	124.00	120.30
1	A	388	ASP	N-CA-CB	-6.26	99.33	110.60
1	A	344	LEU	CA-CB-CG	5.93	128.93	115.30
2	B	172	ASP	N-CA-CB	-5.49	100.73	110.60
2	B	180	VAL	CB-CA-C	-5.31	101.31	111.40
1	A	336	ARG	NE-CZ-NH2	-5.31	117.65	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	387	ASP	Peptide

5.2 Too-close contacts ⓘ

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	3112	0	3078	47	0
2	B	736	0	727	11	0
3	A	1	0	0	0	0
4	A	2	0	0	0	0
5	B	11	0	15	4	0
6	A	31	0	12	0	0
7	A	53	0	0	4	0
7	B	3	0	0	0	0
All	All	3949	0	3832	57	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 7.

All (57) 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)
2:B:173:LYS:NZ	5:B:373:RED:C1	1.82	1.43
2:B:173:LYS:HZ1	5:B:373:RED:C1	1.40	1.31
2:B:173:LYS:HZ3	5:B:373:RED:C1	1.63	0.89
1:A:303:TYR:O	1:A:304:MET:HG2	1.76	0.85
1:A:87:SER:OG	1:A:119:HIS:HE1	1.60	0.84
1:A:134:LYS:HG3	1:A:139:PHE:CD1	2.15	0.81
1:A:199:ASP:O	1:A:203:THR:HG23	1.86	0.75
1:A:122:VAL:HG11	1:A:158:ARG:HD2	1.68	0.74
2:B:173:LYS:CE	5:B:373:RED:C1	2.69	0.70
1:A:134:LYS:HG3	1:A:139:PHE:HD1	1.61	0.66
1:A:69:ASP:OD2	1:A:69:ASP:N	2.28	0.65
1:A:308:ALA:HB2	1:A:322:ALA:HB3	1.79	0.65
1:A:122:VAL:CG1	1:A:158:ARG:HD2	2.32	0.59
1:A:111:VAL:O	1:A:115:VAL:HG13	2.03	0.59
2:B:131:PRO:O	2:B:208:VAL:HG23	2.03	0.58
1:A:61:MET:HA	1:A:64:VAL:HG13	1.86	0.58
1:A:332:ARG:O	1:A:336:ARG:HG2	2.04	0.57
2:B:208:VAL:HG21	2:B:213:ASP:HB3	1.87	0.57
1:A:299:ARG:HD3	7:A:538:HOH:O	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:ASN:N	1:A:178:PRO:HD3	2.23	0.54
1:A:146:ASN:HD21	1:A:380:TYR:HA	1.71	0.54
1:A:139:PHE:HE2	1:A:143:ILE:HG22	1.72	0.54
1:A:51:LYS:HE2	1:A:95:GLU:OE2	2.07	0.53
1:A:192:ASN:HD22	1:A:195:ASP:H	1.57	0.53
1:A:199:ASP:O	1:A:203:THR:CG2	2.56	0.52
1:A:193:VAL:HG21	1:A:281:LEU:HD11	1.93	0.51
1:A:303:TYR:C	1:A:304:MET:HG2	2.31	0.50
1:A:281:LEU:HD23	1:A:281:LEU:C	2.34	0.48
2:B:140:LEU:HB3	2:B:176:ILE:HD13	1.97	0.47
1:A:57:LEU:O	1:A:61:MET:HG3	2.14	0.47
1:A:17:GLU:O	1:A:21:ARG:HG3	2.15	0.46
1:A:201:TYR:CE2	1:A:218:LEU:HB2	2.50	0.46
1:A:62:ARG:HA	1:A:62:ARG:HD2	1.77	0.46
1:A:299:ARG:CD	7:A:538:HOH:O	2.62	0.46
1:A:139:PHE:CZ	1:A:147:ILE:HD12	2.51	0.46
1:A:64:VAL:HA	1:A:67:LEU:HD22	1.98	0.46
1:A:293:PRO:HG2	1:A:296:LYS:HD3	1.97	0.45
1:A:296:LYS:HA	1:A:299:ARG:HD3	1.98	0.45
1:A:267:TYR:CD2	1:A:267:TYR:N	2.84	0.45
1:A:336:ARG:HD3	1:A:341:ASP:OD1	2.17	0.45
1:A:33:LEU:HD12	1:A:184:ILE:HG13	1.97	0.45
2:B:172:ASP:HB3	2:B:173:LYS:HG2	1.99	0.45
1:A:222:GLU:HG2	1:A:274:VAL:HB	1.99	0.44
1:A:374:LYS:HG3	2:B:163:GLY:HA3	1.99	0.43
1:A:117:ASN:O	1:A:120:ASN:HB2	2.18	0.43
1:A:247:GLU:HG3	7:A:535:HOH:O	2.17	0.43
1:A:264:LYS:HG3	1:A:265:GLU:OE1	2.19	0.43
1:A:101:ASP:OD1	1:A:102:PRO:HD2	2.18	0.43
1:A:87:SER:OG	1:A:119:HIS:CE1	2.53	0.42
1:A:53:LEU:HB2	1:A:54:PRO:HD3	2.01	0.42
1:A:308:ALA:CB	1:A:322:ALA:HB3	2.47	0.42
1:A:12:VAL:N	1:A:13:PRO:HD3	2.35	0.42
1:A:388:ASP:HB3	1:A:389:TRP:CD1	2.55	0.42
1:A:142:PHE:CD1	1:A:143:ILE:HG13	2.55	0.42
2:B:208:VAL:HG22	2:B:209:GLU:H	1.85	0.42
1:A:14:LYS:HB2	7:A:532:HOH:O	2.18	0.41
2:B:169:ILE:HD13	2:B:178:PHE:HB2	2.03	0.40

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	378/419 (90%)	356 (94%)	19 (5%)	3 (1%)	24	44
2	B	93/128 (73%)	88 (95%)	5 (5%)	0	100	100
All	All	471/547 (86%)	444 (94%)	24 (5%)	3 (1%)	30	53

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	382	THR
1	A	388	ASP
1	A	214	VAL

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	345/374 (92%)	314 (91%)	31 (9%)	12	21
2	B	79/109 (72%)	74 (94%)	5 (6%)	22	42
All	All	424/483 (88%)	388 (92%)	36 (8%)	13	24

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	SER
1	A	27	LEU
1	A	51	LYS

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Mol	Chain	Res	Type
1	A	62	ARG
1	A	64	VAL
1	A	66	LEU
1	A	67	LEU
1	A	69	ASP
1	A	74	ARG
1	A	105	LEU
1	A	115	VAL
1	A	139	PHE
1	A	144	SER
1	A	146	ASN
1	A	160	SER
1	A	203	THR
1	A	206	MET
1	A	209	GLU
1	A	214	VAL
1	A	226	LYS
1	A	252	SER
1	A	296	LYS
1	A	299	ARG
1	A	328	LEU
1	A	336	ARG
1	A	344	LEU
1	A	346	SER
1	A	381	LYS
1	A	383	THR
1	A	388	ASP
1	A	401	LYS
2	B	133	MET
2	B	140	LEU
2	B	141	SER
2	B	186	LEU
2	B	190	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	70	ASN
1	A	73	ASN
1	A	81	GLN
1	A	86	GLN

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Mol	Chain	Res	Type
1	A	119	HIS
1	A	128	GLN
1	A	146	ASN
1	A	192	ASN
1	A	240	HIS
1	A	302	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
6	ATP	A	504	3,4	24,33,33	1.12	1 (4%)	31,52,52	2.13	8 (25%)
5	RED	B	373	-	8,10,11	0.34	0	4,10,12	2.32	2 (50%)

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
6	ATP	A	504	3,4	-	0/18/38/38	0/3/3/3
5	RED	B	373	-	-	0/6/9/10	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	504	ATP	O4'-C1'	3.74	1.45	1.41

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	504	ATP	N3-C2-N1	-7.27	123.33	128.89
6	A	504	ATP	C2'-C1'-N9	-4.39	107.58	114.29
6	A	504	ATP	C4-C5-N7	-3.39	106.36	109.48
6	A	504	ATP	PA-O3A-PB	-3.05	124.18	132.73
5	B	373	RED	C7-C6-C5	-2.87	106.54	113.91
5	B	373	RED	C3-C2-C1	-2.59	105.91	113.36
6	A	504	ATP	PB-O3B-PG	-2.41	124.58	132.67
6	A	504	ATP	O3G-PG-O2G	-2.28	98.68	107.38
6	A	504	ATP	C2-N1-C6	2.15	122.62	118.77
6	A	504	ATP	O3G-PG-O3B	3.99	123.19	105.09

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	373	RED	4	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 ⓘ

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	381/419 (90%)	0.61	27 (7%) 19 13	48, 64, 86, 96	0
2	B	97/128 (75%)	1.12	20 (20%) 1 1	59, 67, 79, 125	0
All	All	478/547 (87%)	0.71	47 (9%) 10 6	48, 65, 86, 125	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	219	ASP	5.4
1	A	176	THR	4.4
1	A	310	ARG	4.2
2	B	187	ALA	4.1
2	B	226	THR	3.7
2	B	212	ALA	3.7
2	B	218	ALA	3.5
1	A	401	LYS	3.4
2	B	217	PHE	3.4
1	A	142	PHE	3.3
1	A	177	ASN	3.3
1	A	178	PRO	3.2
1	A	399	ALA	3.1
1	A	102	PRO	3.0
2	B	215	SER	3.0
1	A	174	GLY	2.9
1	A	322	ALA	2.9
1	A	400	SER	2.9
1	A	139	PHE	2.8
1	A	158	ARG	2.8
1	A	179	VAL	2.7
2	B	209	GLU	2.7
2	B	210	LYS	2.7
2	B	185	TYR	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	104	VAL	2.6
2	B	186	LEU	2.6
1	A	180	HIS	2.5
1	A	175	ASP	2.5
2	B	207	ILE	2.4
2	B	214	ILE	2.4
2	B	131	PRO	2.4
1	A	229	ASP	2.4
1	A	99	PRO	2.3
1	A	397	ARG	2.3
1	A	38	ASP	2.3
2	B	189	ILE	2.3
2	B	129	TYR	2.2
1	A	134	LYS	2.2
2	B	166	LEU	2.1
2	B	220	TYR	2.1
2	B	156	VAL	2.1
1	A	156	THR	2.1
2	B	192	PRO	2.1
1	A	370	PRO	2.0
1	A	136	LYS	2.0
1	A	43	LYS	2.0
1	A	157	ASN	2.0

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](#)

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(\AA^2)	Q<0.9
4	K	A	503	1/1	0.84	0.34	4.24	99,99,99,99	0
5	RED	B	373	11/12	0.93	0.23	0.05	54,58,69,73	0
6	ATP	A	504	31/31	0.98	0.16	-0.92	36,45,55,59	0
3	MG	A	501	1/1	0.88	0.16	-1.73	50,50,50,50	0
4	K	A	502	1/1	0.80	0.17	-2.23	103,103,103,103	0

6.5 Other polymers [i](#)

There are no such residues in this entry.