



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:18 PM UTC

PDB ID : 3IBQ / pdb_00003ibq
Title : Crystal structure of pyridoxal kinase from Lactobacillus plantarum in complex with ATP
Authors : Bagaria, A.; Kumaran, D.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2009-07-16
Resolution : 2.00 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

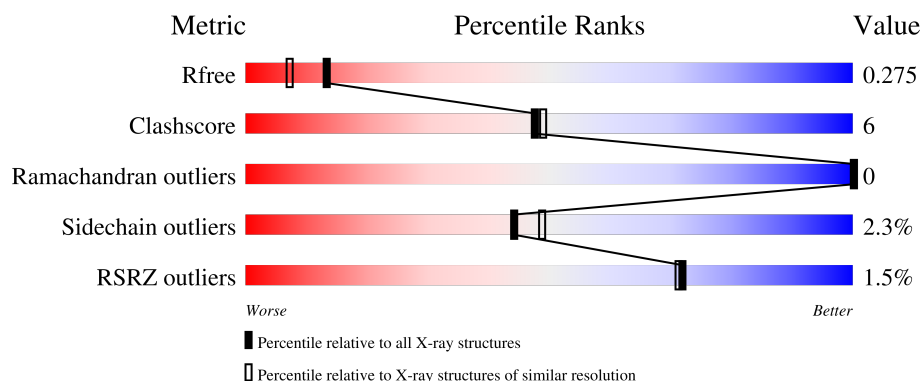
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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 $\leq 5\%$. 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	282	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2238 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.

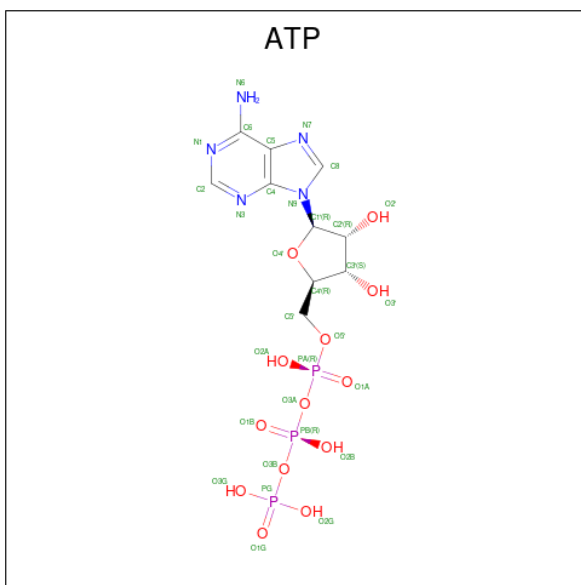
- Molecule 1 is a protein called Pyridoxal kinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	272	2037	1291	346	393	3	4	0	0	0

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	expression tag	UNP Q88YB5
A	0	SER	-	expression tag	UNP Q88YB5
A	1	LEU	-	expression tag	UNP Q88YB5
A	273	GLU	-	expression tag	UNP Q88YB5
A	274	GLY	-	expression tag	UNP Q88YB5
A	275	HIS	-	expression tag	UNP Q88YB5
A	276	HIS	-	expression tag	UNP Q88YB5
A	277	HIS	-	expression tag	UNP Q88YB5
A	278	HIS	-	expression tag	UNP Q88YB5
A	279	HIS	-	expression tag	UNP Q88YB5
A	280	HIS	-	expression tag	UNP Q88YB5

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



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

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

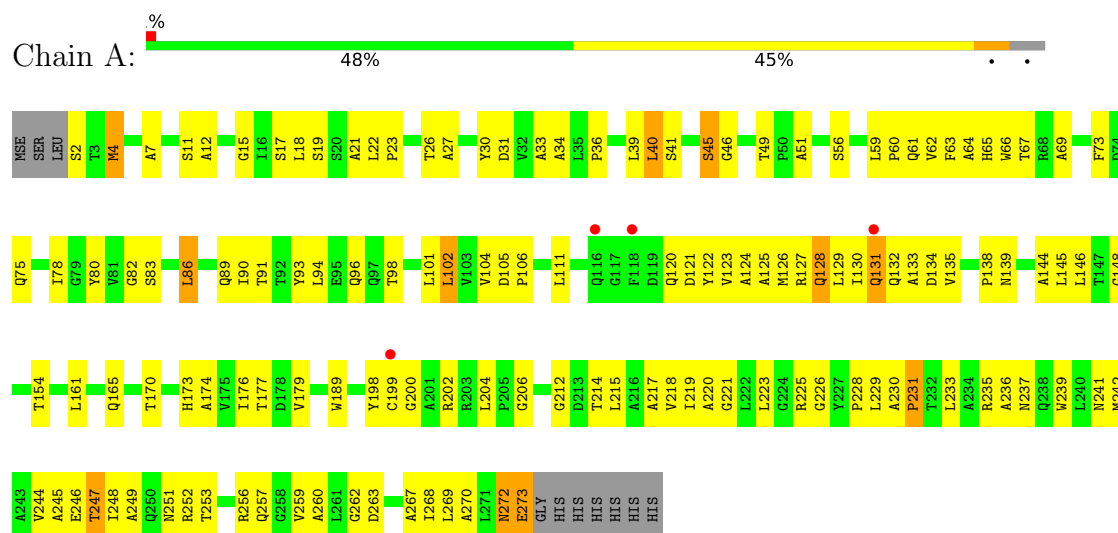
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	169	Total O 169 169	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: Pyridoxal kinase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	64.36Å 69.58Å 132.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.46 – 2.00 38.46 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (38.46-2.00) 99.8 (38.46-2.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	14.92 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.175 , 0.219 (Not available) , 0.275	Depositor DCC
R_{free} test set	1056 reflections (5.16%)	wwPDB-VP
Wilson B-factor (Å ²)	16.2	Xtriage
Anisotropy	0.011	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 48.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2238	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.02% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ATP

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	2.42	129/2071 (6.2%)	1.21	3/2834 (0.1%)

All (129) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	242	MSE	SE-CE	-12.31	1.58	1.95
1	A	59	LEU	C-O	-11.19	1.13	1.24
1	A	22	LEU	C-O	-11.07	1.14	1.24
1	A	4	MSE	SE-CE	-9.59	1.66	1.95
1	A	4	MSE	CG-SE	-8.34	1.70	1.95
1	A	244	VAL	C-O	-8.26	1.14	1.24
1	A	176	ILE	C-O	-8.20	1.15	1.24
1	A	272	ASN	C-N	8.19	1.44	1.33
1	A	221	GLY	C-O	-8.02	1.14	1.23
1	A	73	PHE	C-O	-7.99	1.14	1.23
1	A	33	ALA	CA-CB	-7.64	1.43	1.53
1	A	145	LEU	C-O	-7.55	1.15	1.24
1	A	230	ALA	C-O	-7.55	1.17	1.24
1	A	131	GLN	C-O	-7.53	1.15	1.24
1	A	123	VAL	C-O	-7.43	1.15	1.24
1	A	247	THR	C-O	-7.39	1.15	1.24
1	A	218	VAL	C-O	-7.38	1.15	1.24
1	A	129	LEU	C-O	-7.38	1.15	1.24
1	A	12	ALA	CA-CB	-7.37	1.42	1.53
1	A	267	ALA	CA-CB	-7.32	1.41	1.53
1	A	27	ALA	C-O	-7.30	1.15	1.24
1	A	26	THR	C-O	-7.28	1.15	1.24
1	A	60	PRO	C-O	-7.21	1.15	1.24
1	A	242	MSE	C-O	-7.12	1.15	1.24
1	A	66	TRP	C-O	-7.08	1.15	1.24
1	A	202	ARG	C-O	-7.08	1.15	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	215	LEU	C-O	-7.04	1.16	1.24
1	A	248	ILE	C-O	-7.03	1.16	1.24
1	A	69	ALA	C-O	-6.93	1.15	1.24
1	A	7	ALA	C-O	-6.83	1.16	1.24
1	A	96	GLN	C-O	-6.80	1.15	1.24
1	A	229	LEU	C-O	-6.77	1.16	1.24
1	A	206	GLY	C-O	-6.76	1.16	1.24
1	A	219	ILE	C-O	-6.72	1.16	1.24
1	A	130	ILE	C-O	-6.70	1.15	1.24
1	A	268	ILE	C-O	-6.69	1.16	1.24
1	A	94	LEU	C-O	-6.68	1.16	1.24
1	A	51	ALA	CA-CB	-6.68	1.44	1.53
1	A	134	ASP	C-O	-6.62	1.16	1.24
1	A	236	ALA	C-O	-6.62	1.16	1.24
1	A	223	LEU	C-O	-6.61	1.16	1.24
1	A	91	THR	C-O	-6.55	1.16	1.24
1	A	174	ALA	CA-CB	-6.51	1.44	1.53
1	A	241	ASN	C-O	-6.48	1.16	1.24
1	A	90	ILE	C-O	-6.43	1.16	1.24
1	A	262	GLY	C-O	-6.39	1.15	1.23
1	A	22	LEU	N-CA	-6.36	1.41	1.46
1	A	127	ARG	C-O	-6.35	1.16	1.24
1	A	252	ARG	C-O	-6.30	1.16	1.23
1	A	62	VAL	C-O	-6.28	1.17	1.24
1	A	18	LEU	C-O	-6.25	1.16	1.24
1	A	36	PRO	C-O	-6.25	1.16	1.23
1	A	124	ALA	C-O	-6.25	1.16	1.24
1	A	126	MSE	SE-CE	-6.19	1.76	1.95
1	A	231	PRO	C-O	-6.19	1.16	1.24
1	A	78	ILE	C-O	-6.17	1.17	1.24
1	A	220	ALA	C-O	-6.15	1.17	1.24
1	A	233	LEU	C-O	-6.13	1.17	1.24
1	A	133	ALA	C-O	-6.11	1.16	1.23
1	A	214	THR	C-O	-6.10	1.17	1.24
1	A	39	LEU	C-O	-6.08	1.17	1.24
1	A	235	ARG	C-O	-6.05	1.17	1.24
1	A	239	TRP	C-O	-6.05	1.17	1.24
1	A	121	ASP	C-O	-6.04	1.17	1.24
1	A	217	ALA	C-O	-6.00	1.17	1.24
1	A	120	GLN	C-O	-5.97	1.16	1.24
1	A	89	GLN	C-O	-5.96	1.17	1.24
1	A	80	TYR	C-O	-5.96	1.16	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	148	GLY	C-O	-5.96	1.16	1.24
1	A	257	GLN	C-O	-5.95	1.16	1.23
1	A	125	ALA	C-O	-5.93	1.17	1.24
1	A	204	LEU	C-O	-5.93	1.16	1.23
1	A	269	LEU	C-O	-5.93	1.16	1.24
1	A	260	ALA	C-O	-5.91	1.16	1.23
1	A	19	SER	C-O	-5.88	1.16	1.24
1	A	245	ALA	C-O	-5.84	1.17	1.24
1	A	21	ALA	C-O	-5.84	1.17	1.24
1	A	144	ALA	CA-CB	-5.78	1.44	1.53
1	A	61	GLN	C-O	-5.77	1.17	1.24
1	A	179	VAL	C-O	-5.72	1.18	1.24
1	A	101	LEU	C-O	-5.68	1.16	1.23
1	A	122	TYR	C-O	-5.66	1.17	1.24
1	A	237	ASN	C-O	-5.66	1.17	1.24
1	A	17	SER	C-O	-5.65	1.16	1.24
1	A	34	ALA	C-O	-5.64	1.16	1.23
1	A	111	LEU	C-O	-5.64	1.15	1.23
1	A	267	ALA	C-O	-5.64	1.17	1.24
1	A	270	ALA	C-O	-5.60	1.17	1.24
1	A	263	ASP	C-O	-5.58	1.17	1.24
1	A	256	ARG	C-O	-5.58	1.16	1.24
1	A	86	LEU	C-O	-5.55	1.17	1.24
1	A	51	ALA	C-O	-5.54	1.17	1.24
1	A	64	ALA	C-O	-5.49	1.17	1.24
1	A	40	LEU	C-O	-5.47	1.17	1.23
1	A	15	GLY	C-O	-5.45	1.18	1.24
1	A	64	ALA	CA-CB	-5.45	1.45	1.53
1	A	189	TRP	C-O	-5.44	1.17	1.23
1	A	128	GLN	C-O	-5.44	1.17	1.24
1	A	135	VAL	C-O	-5.40	1.18	1.24
1	A	98	THR	C-O	-5.40	1.16	1.23
1	A	63	PHE	C-O	-5.39	1.17	1.24
1	A	173	HIS	C-O	-5.36	1.17	1.24
1	A	217	ALA	CA-CB	-5.34	1.45	1.53
1	A	41	SER	C-O	-5.34	1.18	1.24
1	A	146	LEU	C-O	-5.33	1.17	1.24
1	A	165	GLN	C-O	-5.31	1.17	1.24
1	A	31	ASP	C-O	-5.30	1.17	1.24
1	A	105	ASP	C-O	-5.29	1.17	1.24
1	A	226	GLY	C-O	-5.29	1.17	1.24
1	A	46	GLY	C-O	-5.29	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	198	TYR	C-O	-5.28	1.17	1.23
1	A	161	LEU	C-O	-5.28	1.19	1.24
1	A	138	PRO	C-O	-5.27	1.17	1.23
1	A	253	THR	C-O	-5.24	1.17	1.24
1	A	248	ILE	CG1-CD1	-5.23	1.31	1.51
1	A	102	LEU	C-O	-5.20	1.18	1.24
1	A	249	ALA	C-O	-5.19	1.17	1.24
1	A	272	ASN	C-O	-5.17	1.18	1.24
1	A	12	ALA	C-O	-5.17	1.18	1.24
1	A	170	THR	C-O	-5.16	1.17	1.23
1	A	23	PRO	C-O	-5.15	1.17	1.24
1	A	200	GLY	C-O	-5.13	1.17	1.23
1	A	246	GLU	C-O	-5.13	1.18	1.24
1	A	45	SER	C-O	-5.11	1.17	1.23
1	A	65	HIS	C-O	-5.09	1.18	1.24
1	A	93	TYR	C-O	-5.09	1.18	1.24
1	A	30	TYR	C-O	-5.05	1.17	1.23
1	A	67	THR	C-O	-5.03	1.18	1.24
1	A	49	THR	C-O	-5.03	1.17	1.24

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	177	THR	N-CA-C	8.68	124.00	113.23
1	A	82	GLY	N-CA-C	5.43	121.21	114.37
1	A	45	SER	N-CA-C	5.19	117.89	107.41

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	2037	0	2031	22	0
2	A	31	0	11	2	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	169	0	0	4	0
All	All	2238	0	2042	23	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 6.

All (23) 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:132:GLN:HG2	4:A:342:HOH:O	1.42	1.16
1:A:272:ASN:O	1:A:273:GLU:HB2	1.69	0.91
2:A:281:ATP:O3G	2:A:281:ATP:O1B	1.94	0.81
1:A:139:ASN:HB3	4:A:358:HOH:O	1.86	0.73
1:A:251:ASN:HB3	4:A:373:HOH:O	1.89	0.70
1:A:56:SER:HA	1:A:86:LEU:HD11	1.76	0.66
1:A:4:MSE:HE2	1:A:75:GLN:HB3	1.81	0.63
1:A:4:MSE:CE	1:A:75:GLN:HB3	2.30	0.61
1:A:228:PRO:HG2	1:A:231:PRO:CG	2.30	0.61
1:A:228:PRO:HG2	1:A:231:PRO:HG3	1.88	0.56
1:A:83:SER:OG	1:A:86:LEU:HD13	2.09	0.52
1:A:104:VAL:O	1:A:106:PRO:HD3	2.11	0.51
1:A:4:MSE:CE	1:A:75:GLN:CB	2.88	0.51
1:A:225:ARG:NH2	1:A:272:ASN:OD1	2.42	0.51
1:A:2:SER:N	4:A:366:HOH:O	2.46	0.48
1:A:4:MSE:HE3	1:A:75:GLN:HB2	1.94	0.48
1:A:199:CYS:SG	1:A:199:CYS:O	2.73	0.47
1:A:212:GLY:HA3	2:A:281:ATP:O1B	2.16	0.46
1:A:4:MSE:HE3	1:A:75:GLN:CB	2.46	0.45
1:A:228:PRO:HG2	1:A:231:PRO:HG2	1.97	0.45
1:A:128:GLN:O	1:A:131:GLN:HG2	2.17	0.43
1:A:102:LEU:CD2	1:A:132:GLN:HG3	2.48	0.43
1:A:247:THR:OG1	1:A:259:VAL:HG23	2.18	0.43

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	270/282 (96%)	259 (96%)	11 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	213/217 (98%)	208 (98%)	5 (2%)	44	49

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	SER
1	A	40	LEU
1	A	45	SER
1	A	154	THR
1	A	273	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	GLN
1	A	97	GLN
1	A	131	GLN

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Mol	Chain	Res	Type
1	A	184	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\# Z > 2$
2	ATP	A	281	3	32,33,33	2.07	9 (28%)	48,52,52	2.07	16 (33%)

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	ATP	A	281	3	-	3/22/38/38	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	281	ATP	C5-N7	-5.53	1.29	1.39
2	A	281	ATP	C8-N9	-4.62	1.29	1.37
2	A	281	ATP	C4-N9	-3.62	1.30	1.37
2	A	281	ATP	C5-C4	3.24	1.44	1.39
2	A	281	ATP	C5-C6	2.81	1.48	1.41
2	A	281	ATP	O3'-C3'	-2.47	1.36	1.43
2	A	281	ATP	C2'-C3'	-2.46	1.46	1.53
2	A	281	ATP	C6-N1	-2.19	1.29	1.35
2	A	281	ATP	PG-O2G	-2.15	1.46	1.54

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	281	ATP	C5-C4-N3	-5.56	119.06	126.72
2	A	281	ATP	N3-C4-N9	4.40	134.65	127.17
2	A	281	ATP	C4-C5-N7	-4.35	105.60	110.58
2	A	281	ATP	O3G-PG-O2G	3.91	122.45	107.80
2	A	281	ATP	O2G-PG-O3B	-3.83	91.78	104.64
2	A	281	ATP	C5-N7-C8	3.35	108.72	103.45
2	A	281	ATP	C2-N3-C4	3.14	119.49	111.83
2	A	281	ATP	O3'-C3'-C2'	-2.98	102.26	111.82
2	A	281	ATP	O3A-PA-O1A	2.82	119.19	110.70
2	A	281	ATP	C2-N1-C6	2.82	123.36	118.73
2	A	281	ATP	C6-C5-N7	2.77	137.42	132.09
2	A	281	ATP	C4-N9-C8	2.70	108.58	105.74
2	A	281	ATP	O3A-PB-O1B	2.48	118.18	110.70
2	A	281	ATP	N3-C2-N1	-2.36	125.01	128.58
2	A	281	ATP	O2'-C2'-C3'	-2.33	104.34	111.82
2	A	281	ATP	N9-C8-N7	-2.16	110.87	113.94

There are no chirality outliers.

All (3) torsion outliers are listed below:

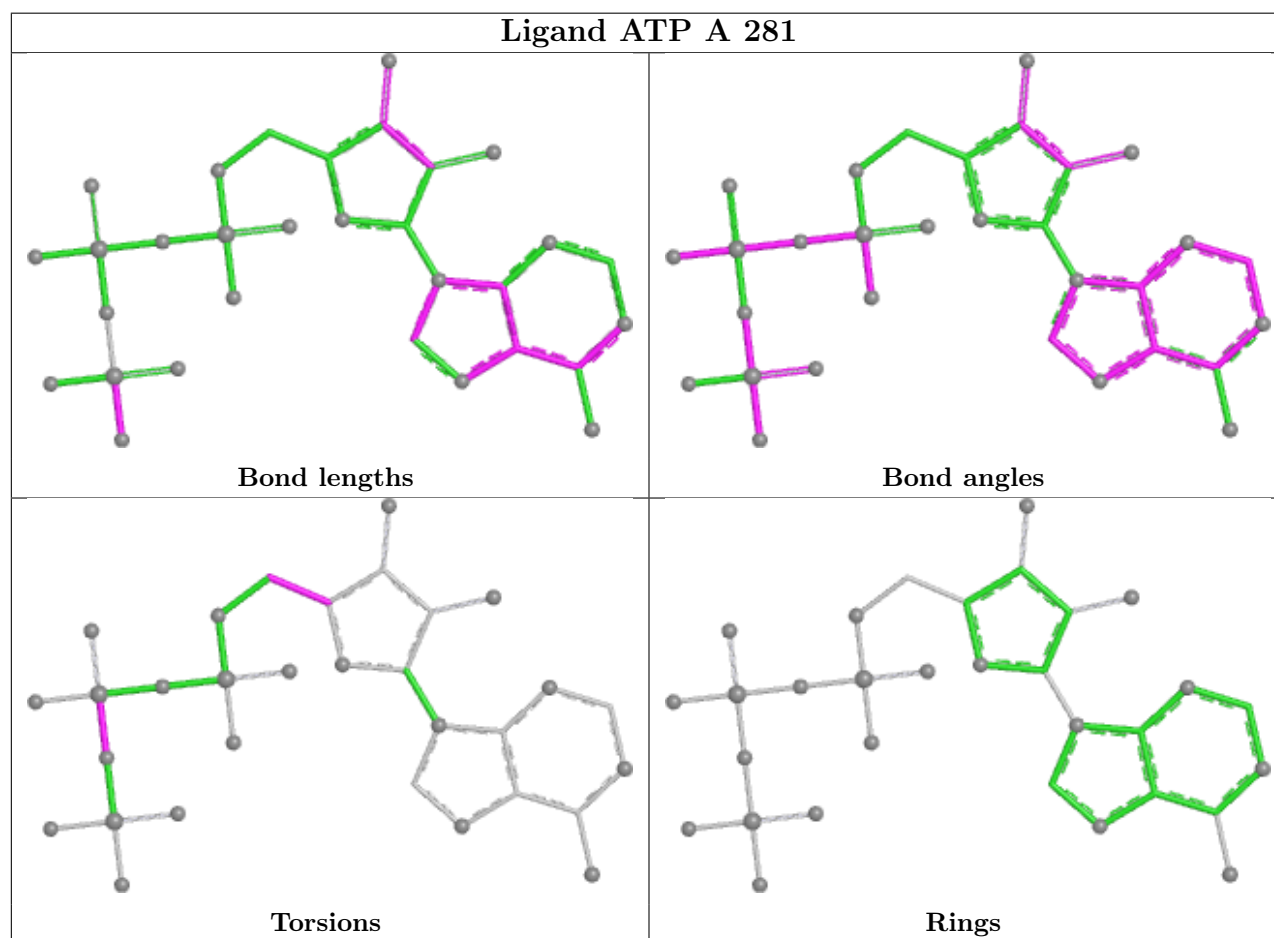
Mol	Chain	Res	Type	Atoms
2	A	281	ATP	O4'-C4'-C5'-O5'
2	A	281	ATP	C3'-C4'-C5'-O5'
2	A	281	ATP	PG-O3B-PB-O2B

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	281	ATP	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	268/282 (95%)	-0.04	4 (1%) 72 71	6, 14, 31, 40	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	199	CYS	4.2
1	A	116	GLN	3.2
1	A	131	GLN	3.1
1	A	118	PHE	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 oligosaccharides in this entry.

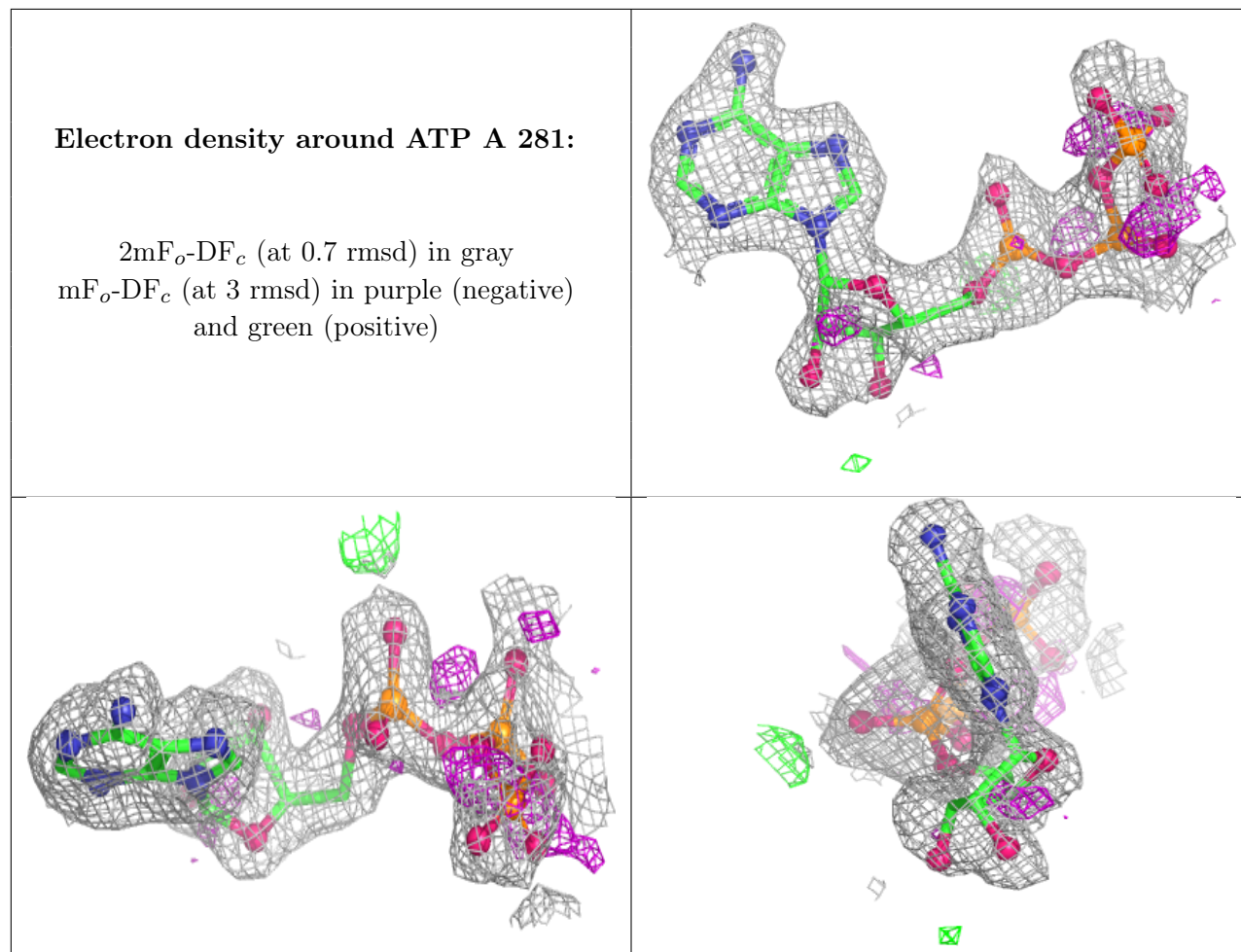
6.4 Ligands [i](#)

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, 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	B-factors(Å ²)	Q<0.9
2	ATP	A	281	31/31	0.78	0.14	26,38,68,69	0
3	MG	A	282	1/1	0.83	0.10	40,40,40,40	0

The following is a graphical depiction of the model fit to experimental electron density of all

instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers ⓘ

There are no such residues in this entry.