



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 12:45 PM UTC

PDB ID : 2IDX / pdb_00002idx
Title : Structure of Human ATP:Cobalamin adenosyltransferase bound to ATP.
Authors : Schubert, H.L.; Hill, C.P.
Deposited on : 2006-09-15
Resolution : 2.50 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

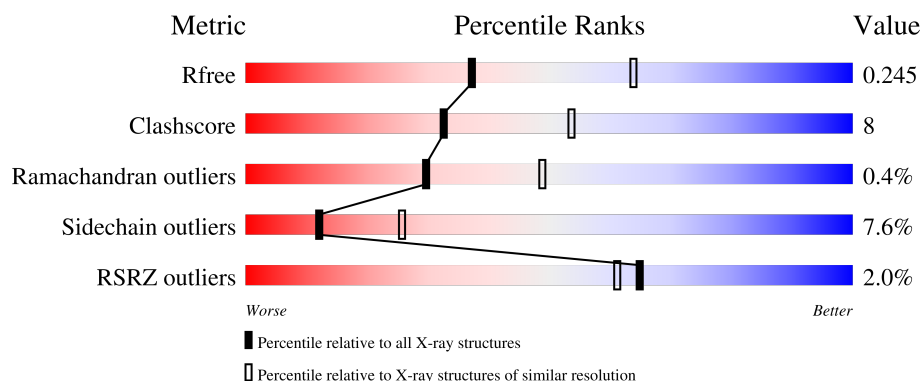
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

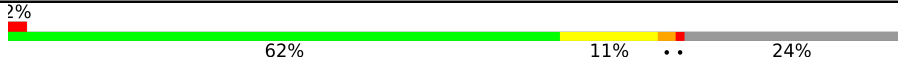
The reported resolution of this entry is 2.50 Å.

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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	196	
1	B	196	
1	C	196	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 4243 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 Cob(I)yrinic acid a,c-diamide adenosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	148	Total	C	N	O	S	0	0	0
			1154	735	194	219	6			
1	B	178	Total	C	N	O	S	0	3	0
			1412	894	238	273	7			
1	C	180	Total	C	N	O	S	0	2	0
			1424	902	242	273	7			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	55	MET	-	cloning artifact	UNP Q96EY8
B	55	MET	-	cloning artifact	UNP Q96EY8
C	55	MET	-	cloning artifact	UNP Q96EY8

- Molecule 2 is SULFATE ION (CCD ID: 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
2	C	1	Total O S 5 4 1	0	0

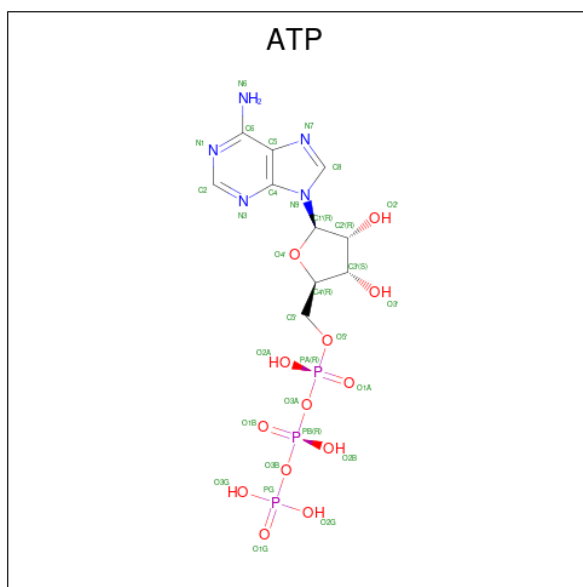
- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	2	Total Mg 2 2	0	0
4	B	1	Total Mg 1 1	0	0
4	C	1	Total Mg 1 1	0	0

- Molecule 5 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



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

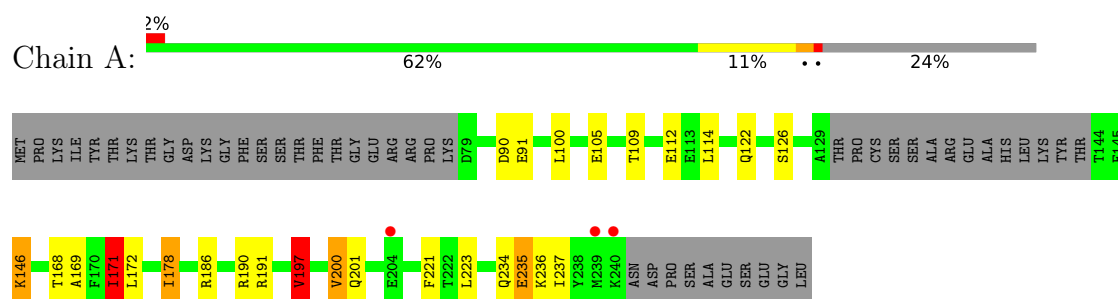
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	32	Total	O	0	0
			32	32		
6	B	69	Total	O	0	0
			69	69		
6	C	70	Total	O	0	0
			70	70		

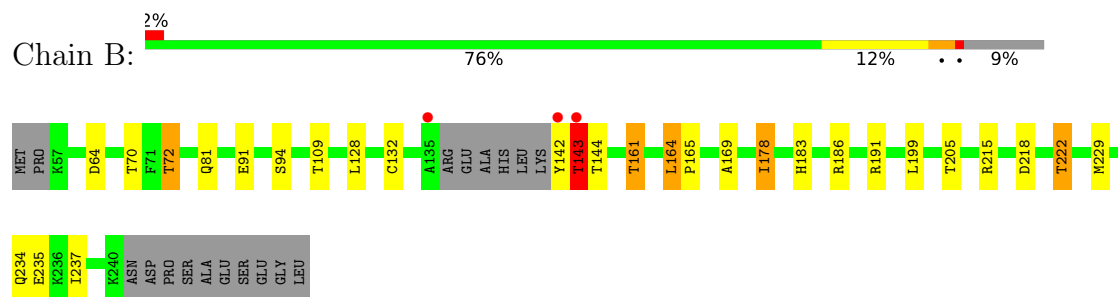
3 Residue-property plots [i](#)

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.

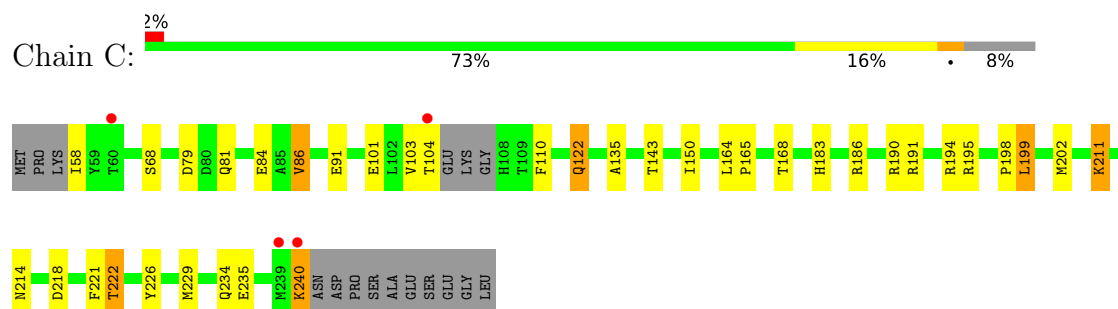
- Molecule 1: Cob(I)yrinic acid a,c-diamide adenosyltransferase



- Molecule 1: Cob(I)yrinic acid a,c-diamide adenosyltransferase



- Molecule 1: Cob(I)yrinic acid a,c-diamide adenosyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	111.24Å 111.24Å 115.53Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.50 30.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.0 (30.00-2.50) 98.9 (30.00-2.50)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.50 (at 2.51Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.197 , 0.251 0.193 , 0.245	Depositor DCC
R_{free} test set	1433 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	35.5	Xtriage
Anisotropy	0.176	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 31.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.036 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4243	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.08% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, ATP, MG, CL

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	0.68	1/1173 (0.1%)	0.89	2/1583 (0.1%)
1	B	0.77	0/1437	0.98	4/1938 (0.2%)
1	C	0.75	1/1450 (0.1%)	0.97	3/1957 (0.2%)
All	All	0.74	2/4060 (0.0%)	0.95	9/5478 (0.2%)

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	B	0	1
1	C	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	197	VAL	CA-CB	5.59	1.56	1.54
1	C	86	VAL	CA-CB	5.04	1.60	1.54

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	132	CYS	N-CA-C	6.19	118.10	111.36
1	C	202	MET	N-CA-C	-6.18	104.34	113.61
1	A	200	VAL	N-CA-C	-5.78	103.68	110.21
1	C	84	GLU	N-CA-C	-5.76	105.08	111.36
1	B	143	THR	N-CA-C	5.63	122.80	110.80
1	C	79	ASP	N-CA-CB	-5.61	101.00	110.49

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	164	LEU	CA-C-N	-5.37	115.74	119.66
1	B	164	LEU	C-N-CA	-5.37	115.74	119.66
1	A	171	ILE	CB-CA-C	-5.02	104.20	111.68

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	142	TYR	Peptide
1	C	103	VAL	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	1154	0	1159	21	0
1	B	1412	0	1404	21	0
1	C	1424	0	1416	24	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
3	A	1	0	0	0	0
4	A	2	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
5	A	31	0	12	2	0
5	B	31	0	12	0	0
6	A	32	0	0	2	0
6	B	69	0	0	6	0
6	C	70	0	0	6	1
All	All	4243	0	4003	64	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 8.

All (64) 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:146:LYS:HD3	1:A:146:LYS:H	1.23	1.04
1:B:165:PRO:HD3	1:B:229:MET:CE	1.87	1.04
1:B:164:LEU:HA	1:B:229:MET:HE1	1.48	0.95
1:B:165:PRO:HD3	1:B:229:MET:HE2	1.47	0.94
1:C:91:GLU:CD	1:C:191:ARG:HH22	1.85	0.84
1:C:240:LYS:HE3	1:C:240:LYS:C	2.02	0.84
1:B:91:GLU:CD	1:B:191:ARG:HH22	1.86	0.84
1:A:200:VAL:O	1:A:201:GLN:HB2	1.80	0.80
1:B:218:ASP:O	1:B:222:THR:HG23	1.84	0.77
1:C:58:ILE:N	6:C:661:HOH:O	2.15	0.77
1:A:146:LYS:HD3	1:A:146:LYS:N	2.00	0.76
1:A:91:GLU:CD	1:A:191:ARG:HH22	1.93	0.75
1:C:186:ARG:HD2	1:C:221:PHE:HB2	1.71	0.71
1:B:169:ALA:HB1	1:B:237:ILE:HG23	1.76	0.67
1:B:81[B]:GLN:HG3	6:B:612:HOH:O	1.94	0.66
1:C:218:ASP:O	1:C:222:THR:HG23	1.98	0.64
1:A:90:ASP:OD1	6:A:636:HOH:O	2.16	0.62
1:C:91:GLU:OE1	1:C:191:ARG:NH2	2.34	0.60
1:A:91:GLU:CD	1:A:191:ARG:NH2	2.61	0.59
1:C:150:ILE:HD13	1:C:211:LYS:HB3	1.85	0.58
1:A:178:ILE:HG12	6:A:622:HOH:O	2.03	0.58
1:B:91:GLU:OE1	1:B:191:ARG:NH2	2.36	0.58
1:B:178:ILE:HG12	6:B:655:HOH:O	2.03	0.57
1:A:186:ARG:HD2	1:A:221:PHE:HB2	1.87	0.55
1:B:183:HIS:HD2	6:B:623:HOH:O	1.90	0.54
1:B:164:LEU:HD23	1:B:229:MET:HE3	1.88	0.53
1:C:214:ASN:ND2	6:C:652:HOH:O	2.40	0.53
1:A:190:ARG:NH2	5:A:607:ATP:C4	2.79	0.51
1:C:165:PRO:HD3	1:C:229:MET:HE3	1.93	0.50
1:A:171:ILE:CG2	1:A:235:GLU:HB2	2.42	0.49
1:C:81:GLN:HB3	1:C:199:LEU:HD11	1.95	0.48
1:A:91:GLU:OE1	1:A:191:ARG:NH2	2.46	0.48
1:C:191:ARG:HG3	1:C:194:ARG:NH2	2.29	0.48
1:B:169:ALA:HB1	1:B:237:ILE:CG2	2.42	0.48
1:A:235:GLU:H	1:A:235:GLU:HG2	1.30	0.48
1:C:122:GLN:NE2	6:C:630:HOH:O	2.46	0.48
1:A:197:VAL:O	1:A:200:VAL:O	2.31	0.47
1:C:190:ARG:O	1:C:194:ARG:HG3	2.13	0.47
1:B:164:LEU:HD23	1:B:229:MET:CE	2.44	0.47
1:A:122:GLN:NE2	1:A:122:GLN:HA	2.30	0.47
1:B:164:LEU:HA	1:B:229:MET:CE	2.33	0.47
1:C:164:LEU:HD21	1:C:226:TYR:HA	1.97	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:607:ATP:H2'	1:C:68:SER:OG	2.15	0.46
1:C:183:HIS:HD2	6:C:633:HOH:O	1.98	0.46
1:C:81:GLN:HG2	6:C:638:HOH:O	2.15	0.46
1:C:195:ARG:O	1:C:198:PRO:HD2	2.16	0.46
1:C:165:PRO:HD3	1:C:229:MET:CE	2.46	0.46
1:A:191:ARG:NH1	1:C:91:GLU:OE2	2.49	0.45
1:C:110:PHE:C	1:C:110:PHE:CD1	2.95	0.44
1:B:235:GLU:HG2	6:B:631:HOH:O	2.17	0.44
1:C:135:ALA:CB	1:C:143:THR:HG21	2.46	0.44
1:A:172:LEU:HD12	1:A:236:LYS:HB3	2.00	0.43
1:A:109:THR:HG22	1:A:112:GLU:OE1	2.18	0.43
1:B:128:LEU:HD23	1:B:205:THR:HG21	2.02	0.42
1:A:100:LEU:HD21	1:A:114:LEU:HB2	2.02	0.42
1:A:146:LYS:H	1:A:146:LYS:CD	2.09	0.42
1:B:183:HIS:HE1	6:B:613:HOH:O	2.02	0.42
1:C:183:HIS:HE1	6:C:639:HOH:O	2.02	0.42
1:B:161:THR:HG21	6:B:651:HOH:O	2.19	0.41
1:B:165:PRO:HD3	1:B:229:MET:HE1	1.93	0.41
1:A:169:ALA:HB1	1:A:237:ILE:HG23	2.02	0.41
1:B:70:THR:C	1:B:72:THR:H	2.29	0.41
1:B:64:ASP:CG	1:C:211:LYS:HG2	2.45	0.41
1:A:200:VAL:O	1:A:201:GLN:CB	2.54	0.40

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)
6:C:646:HOH:O	6:C:646:HOH:O[4_555]	2.07	0.13

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	144/196 (74%)	138 (96%)	5 (4%)	1 (1%)	18	34
1	B	177/196 (90%)	168 (95%)	8 (4%)	1 (1%)	21	38
1	C	178/196 (91%)	173 (97%)	5 (3%)	0	100	100
All	All	499/588 (85%)	479 (96%)	18 (4%)	2 (0%)	30	49

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	143	THR
1	A	234	GLN

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	123/164 (75%)	114 (93%)	9 (7%)	13	27
1	B	152/164 (93%)	140 (92%)	12 (8%)	11	24
1	C	153/164 (93%)	142 (93%)	11 (7%)	13	28
All	All	428/492 (87%)	396 (92%)	32 (8%)	12	26

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	105	GLU
1	A	126	SER
1	A	146	LYS
1	A	168	THR
1	A	171	ILE
1	A	178	ILE
1	A	197	VAL
1	A	223	LEU
1	A	235	GLU
1	B	72	THR
1	B	94	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	109	THR
1	B	143	THR
1	B	144	THR
1	B	161	THR
1	B	178	ILE
1	B	186	ARG
1	B	199	LEU
1	B	215	ARG
1	B	222	THR
1	B	234	GLN
1	C	86	VAL
1	C	101	GLU
1	C	104	THR
1	C	122	GLN
1	C	168	THR
1	C	199	LEU
1	C	211	LYS
1	C	222	THR
1	C	234	GLN
1	C	235	GLU
1	C	240	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	155	GLN
1	A	201	GLN
1	A	208	ASN
1	B	183	HIS
1	C	183	HIS
1	C	208	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 5 are monoatomic - leaving 5 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	SO4	C	402	-	4,4,4	0.20	0	6,6,6	0.28	0
5	ATP	A	607	4	32,33,33	1.35	4 (12%)	48,52,52	1.81	12 (25%)
2	SO4	B	403	-	4,4,4	0.33	0	6,6,6	0.32	0
2	SO4	A	401	-	4,4,4	0.26	0	6,6,6	0.42	0
5	ATP	B	606	4	32,33,33	1.34	3 (9%)	48,52,52	1.93	9 (18%)

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
5	ATP	B	606	4	-	4/22/38/38	0/3/3/3
5	ATP	A	607	4	-	2/22/38/38	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	606	ATP	C5-C4	4.79	1.47	1.39
5	A	607	ATP	C5-C4	4.27	1.46	1.39
5	B	606	ATP	C5-C6	2.71	1.48	1.41
5	A	607	ATP	PB-O3B	2.49	1.62	1.59
5	B	606	ATP	C5-N7	-2.40	1.34	1.39
5	A	607	ATP	C5-C6	2.32	1.47	1.41
5	A	607	ATP	C4-N9	-2.22	1.33	1.37

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	606	ATP	C5-C4-N3	-6.03	118.42	126.72
5	B	606	ATP	N3-C4-N9	4.96	135.59	127.17
5	A	607	ATP	N3-C2-N1	-4.43	121.88	128.58
5	A	607	ATP	C5-C4-N3	-4.22	120.90	126.72
5	A	607	ATP	N3-C4-N9	4.15	134.22	127.17
5	B	606	ATP	C2-N3-C4	3.81	121.13	111.83
5	A	607	ATP	C4-N9-C8	3.74	109.66	105.74
5	A	607	ATP	C2-N1-C6	3.67	124.75	118.73
5	B	606	ATP	C4-C5-N7	-3.48	106.60	110.58
5	B	606	ATP	N3-C2-N1	-3.29	123.59	128.58
5	A	607	ATP	C2-N3-C4	3.24	119.73	111.83
5	B	606	ATP	C5-N7-C8	2.67	107.65	103.45
5	A	607	ATP	O2B-PB-O3B	2.67	114.48	107.27
5	B	606	ATP	C4-N9-C8	2.58	108.45	105.74
5	B	606	ATP	O3A-PB-O1B	-2.53	103.08	110.70
5	A	607	ATP	O2G-PG-O1G	2.32	119.86	110.83
5	A	607	ATP	N6-C6-N1	2.26	123.42	118.38
5	A	607	ATP	C4-C5-N7	-2.26	108.00	110.58
5	B	606	ATP	O3B-PB-O1B	-2.14	104.26	110.70
5	A	607	ATP	N9-C8-N7	-2.08	110.98	113.94
5	A	607	ATP	O3B-PB-O1B	-2.01	104.64	110.70

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	606	ATP	C5'-O5'-PA-O1A
5	A	607	ATP	PG-O3B-PB-O2B
5	B	606	ATP	PG-O3B-PB-O1B
5	B	606	ATP	PG-O3B-PB-O2B
5	B	606	ATP	C2'-C1'-N9-C8
5	A	607	ATP	PG-O3B-PB-O1B

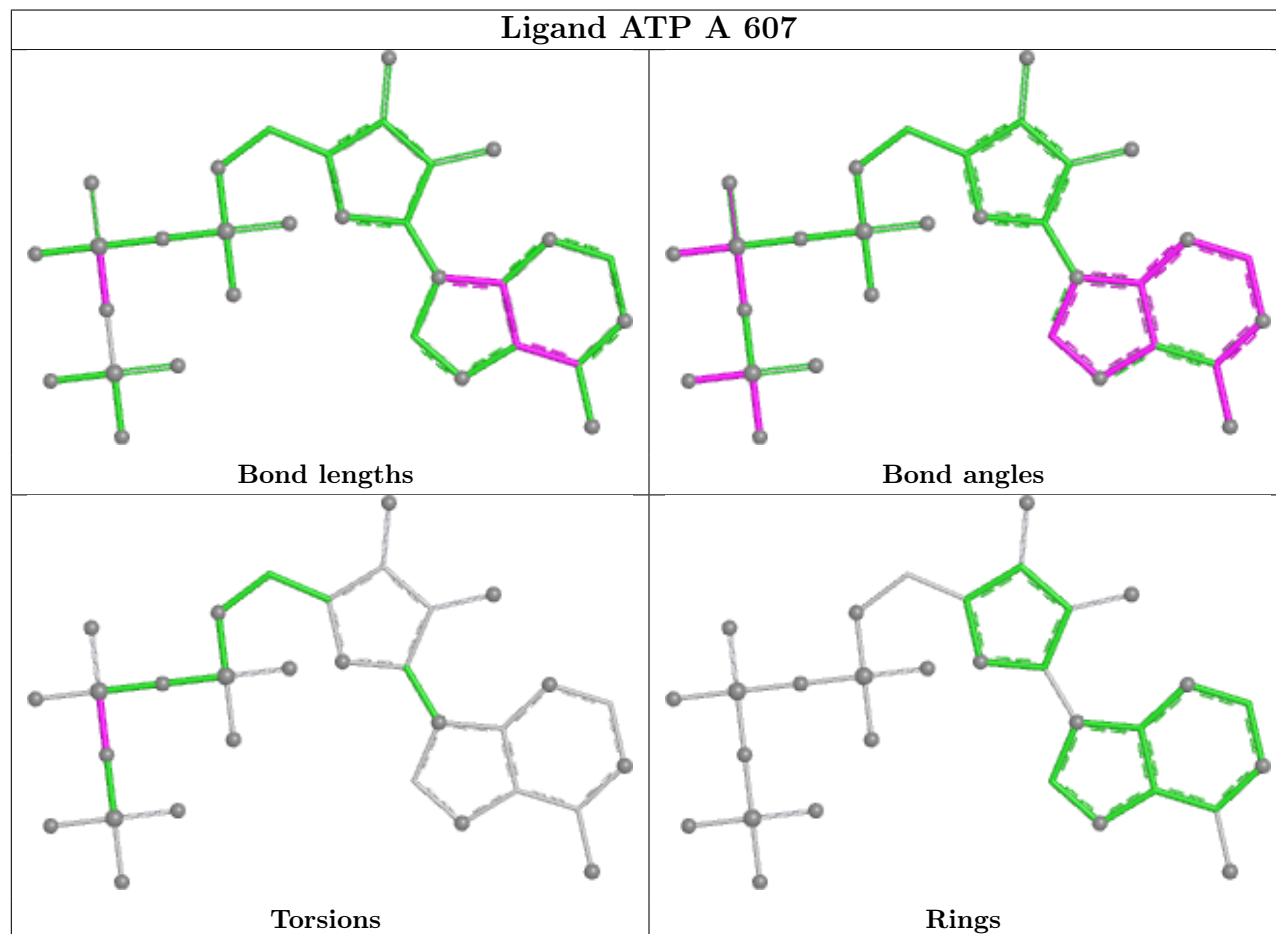
There are no ring outliers.

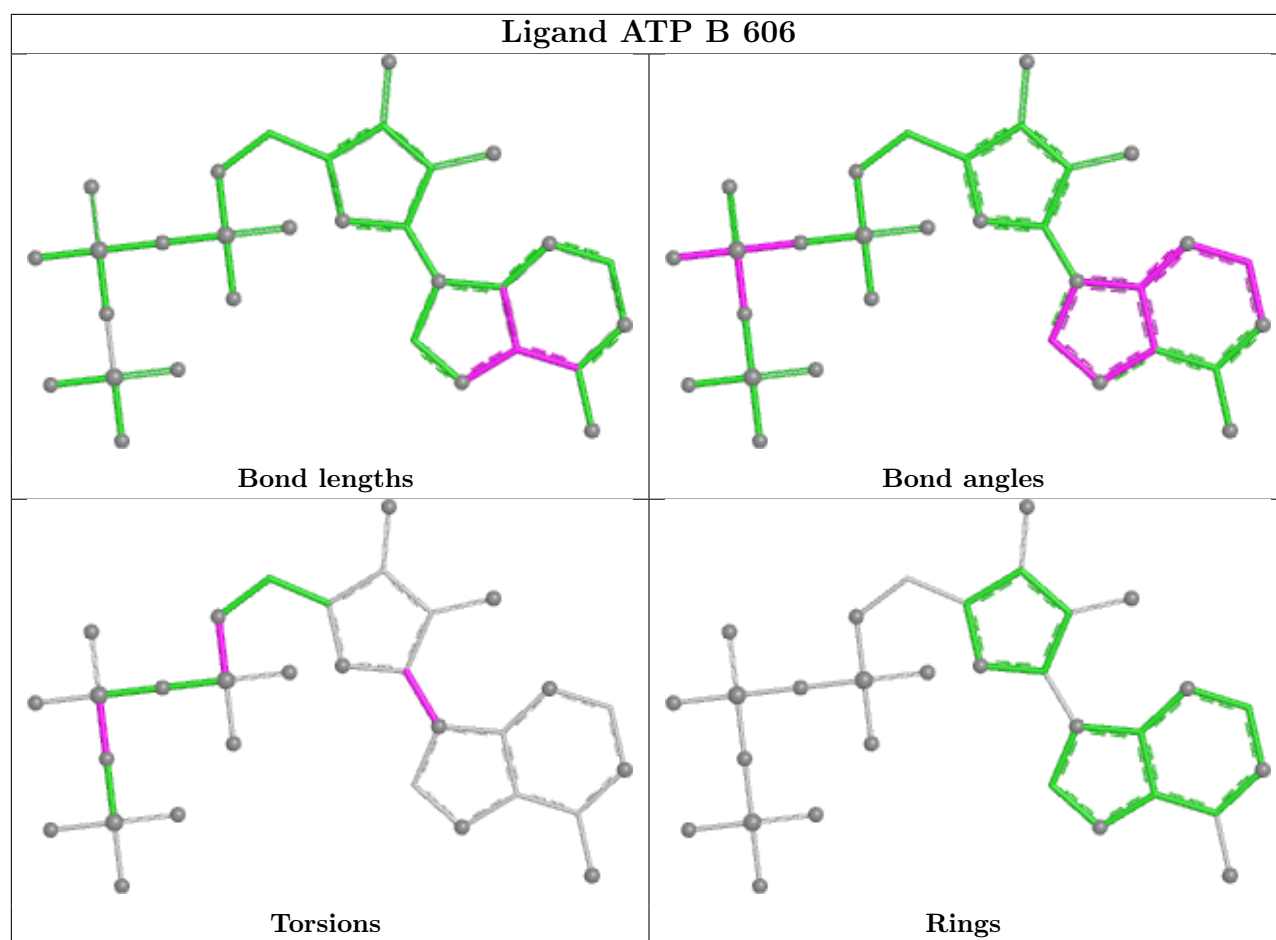
1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	607	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 [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	148/196 (75%)	-0.03	3 (2%) 65 61	22, 39, 61, 71	0
1	B	178/196 (90%)	-0.35	3 (1%) 69 65	10, 28, 50, 70	3 (1%)
1	C	180/196 (91%)	-0.35	4 (2%) 62 58	11, 30, 48, 63	2 (1%)
All	All	506/588 (86%)	-0.26	10 (1%) 65 61	10, 32, 56, 71	5 (0%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	142	TYR	4.5
1	B	143	THR	3.0
1	A	239	MET	3.0
1	C	239	MET	2.7
1	B	135	ALA	2.6
1	A	204	GLU	2.6
1	C	104	THR	2.6
1	A	240	LYS	2.6
1	C	240	LYS	2.4
1	C	60	THR	2.4

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 ⓘ

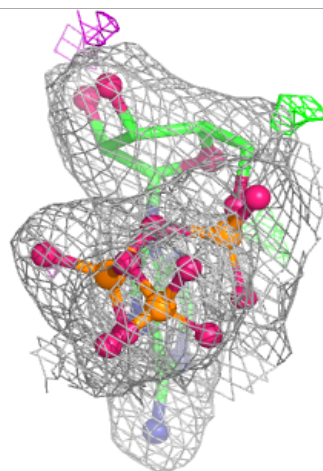
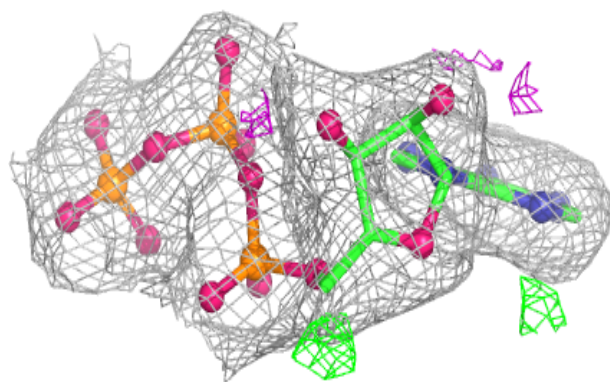
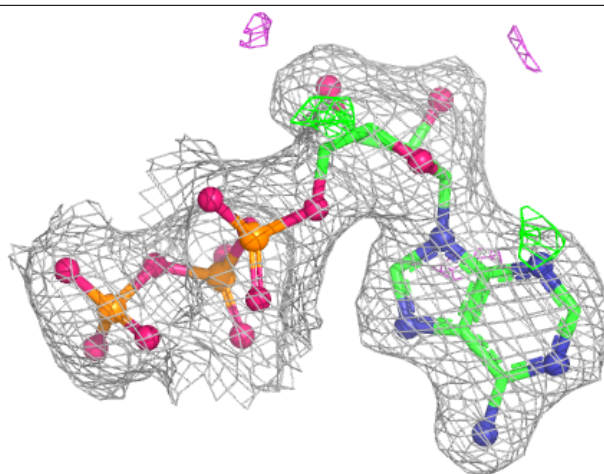
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	SO4	C	402	5/5	0.76	0.16	87,87,88,88	0
2	SO4	A	401	5/5	0.88	0.11	69,70,71,71	0
4	MG	A	606	1/1	0.95	0.05	43,43,43,43	0
4	MG	B	605	1/1	0.95	0.08	46,46,46,46	0
2	SO4	B	403	5/5	0.98	0.07	40,43,43,43	0
5	ATP	A	607	31/31	0.98	0.06	25,29,31,34	0
4	MG	A	603	1/1	0.99	0.03	33,33,33,33	0
3	CL	A	601	1/1	0.99	0.03	33,33,33,33	0
5	ATP	B	606	31/31	0.99	0.04	19,22,24,24	0
4	MG	C	604	1/1	1.00	0.02	19,19,19,19	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.

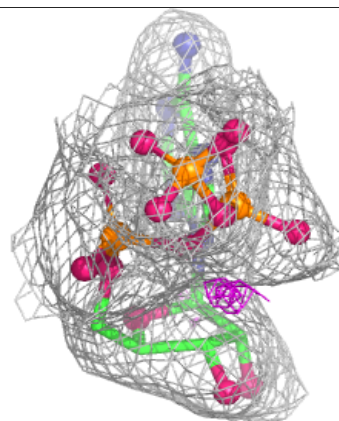
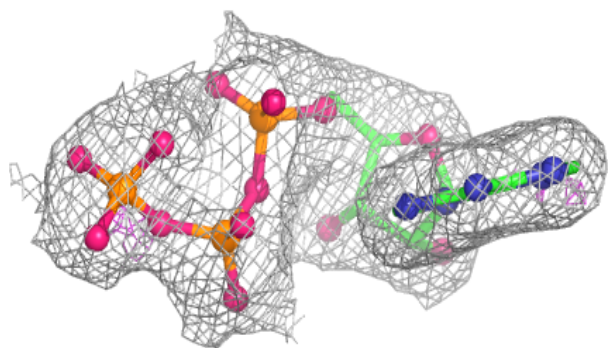
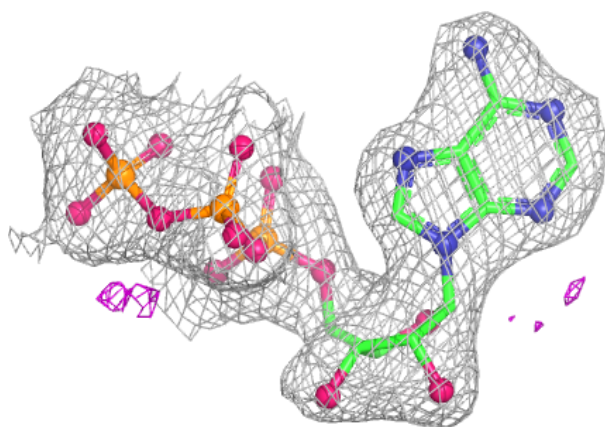
Electron density around ATP A 607:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ATP B 606:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers ⓘ

There are no such residues in this entry.