



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 25, 2026 – 10:39 AM UTC

PDB ID : 9JC7 / pdb_00009jc7
Title : Crystal structure of alanyl-tRNA synthetase in complex with ATP and L-alanine
Authors : Son, S.Y.; Cha, S.S.
Deposited on : 2024-08-28
Resolution : 1.90 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

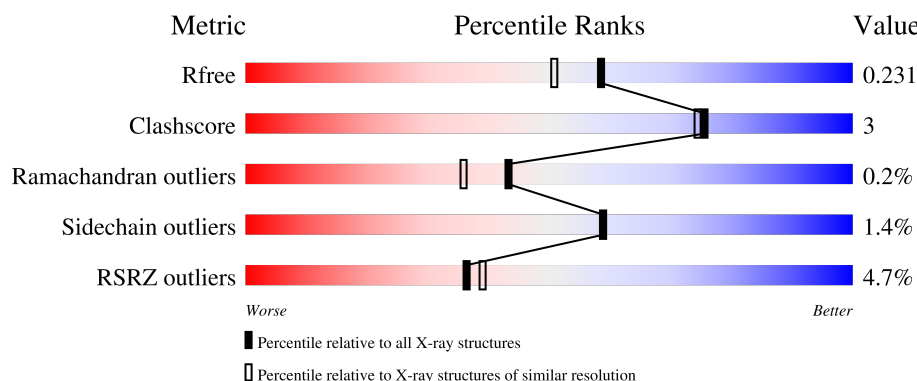
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	435	6% 92% 5%
1	B	435	4% 91% 6%

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
3	ALA	A	502	-	X	-	-
3	ALA	B	502	-	X	-	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 7055 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.

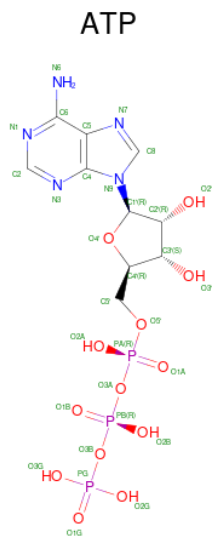
- Molecule 1 is a protein called Alanine-tRNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	425	Total	C	N	O	S	0	0	0
			3347	2122	584	625	16			
1	B	425	Total	C	N	O	S	0	0	0
			3347	2122	584	625	16			

There are 12 discrepancies between the modelled and reference sequences:

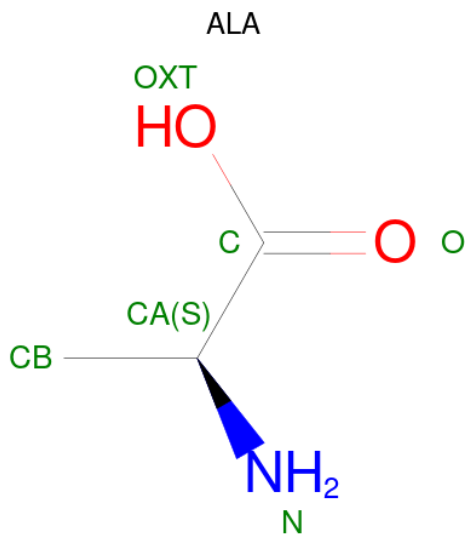
Chain	Residue	Modelled	Actual	Comment	Reference
A	430	HIS	-	expression tag	UNP A0A172UEB3
A	431	HIS	-	expression tag	UNP A0A172UEB3
A	432	HIS	-	expression tag	UNP A0A172UEB3
A	433	HIS	-	expression tag	UNP A0A172UEB3
A	434	HIS	-	expression tag	UNP A0A172UEB3
A	435	HIS	-	expression tag	UNP A0A172UEB3
B	430	HIS	-	expression tag	UNP A0A172UEB3
B	431	HIS	-	expression tag	UNP A0A172UEB3
B	432	HIS	-	expression tag	UNP A0A172UEB3
B	433	HIS	-	expression tag	UNP A0A172UEB3
B	434	HIS	-	expression tag	UNP A0A172UEB3
B	435	HIS	-	expression tag	UNP A0A172UEB3

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is ALANINE (CCD ID: ALA) (formula: $\text{C}_3\text{H}_7\text{NO}_2$) (labeled as "Ligand of Interest" by depositor).



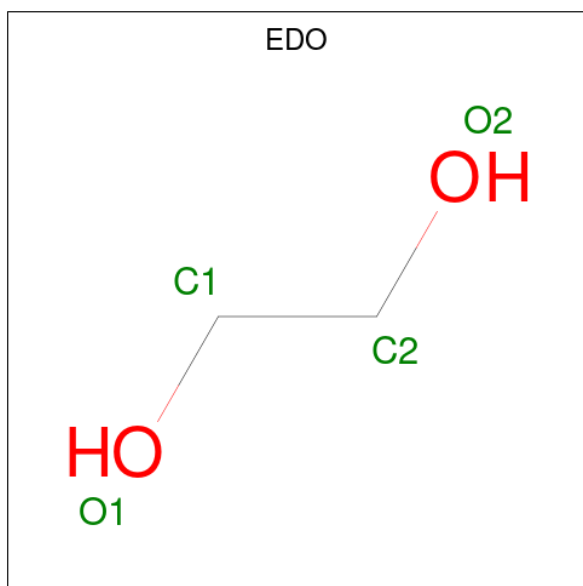
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			6	3	1	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			6	3	1	2		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



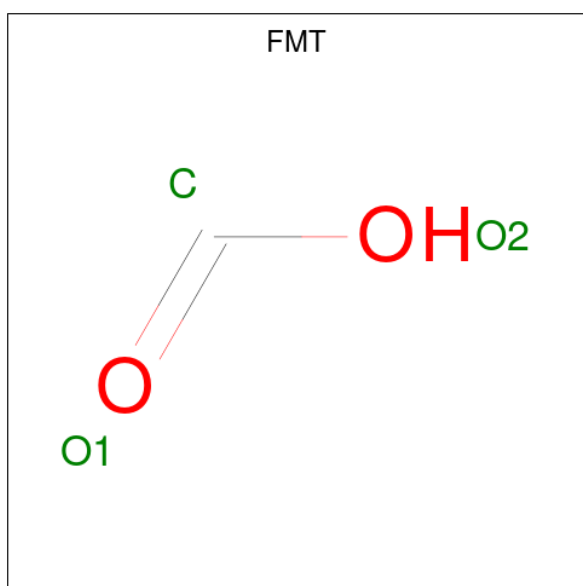
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total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

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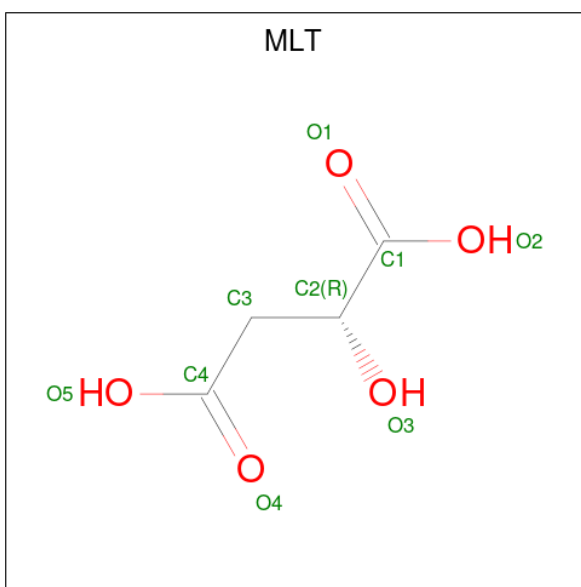
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is FORMIC ACID (CCD ID: FMT) (formula: CH_2O_2).



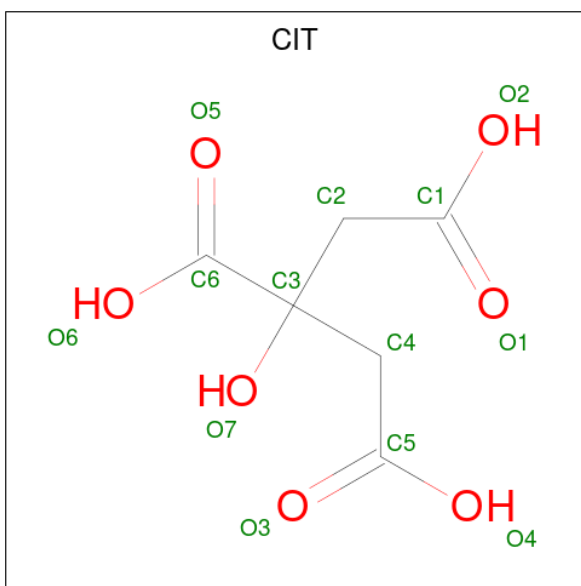
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			3	1	2		
5	B	1	Total	C	O	0	0
			3	1	2		
5	B	1	Total	C	O	0	0
			3	1	2		

- Molecule 6 is D-MALATE (CCD ID: MLT) (formula: $\text{C}_4\text{H}_6\text{O}_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			9	4	5		

- Molecule 7 is CITRIC ACID (CCD ID: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			13	6	7		

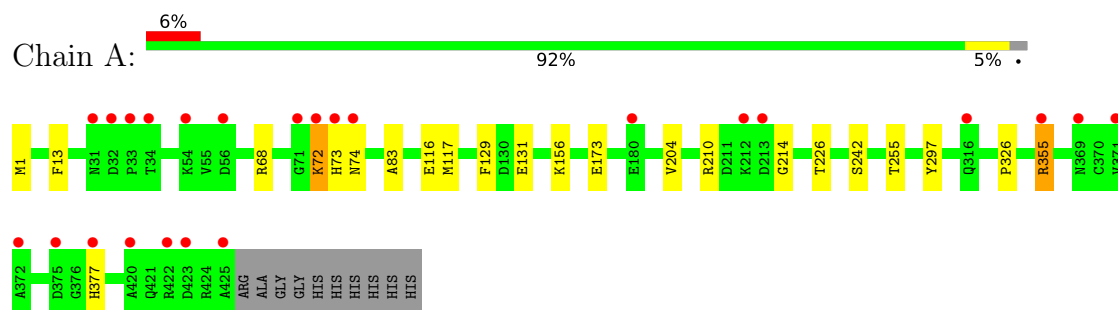
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	95	Total 95	O 95	0	0
8	B	101	Total 101	O 101	0	0

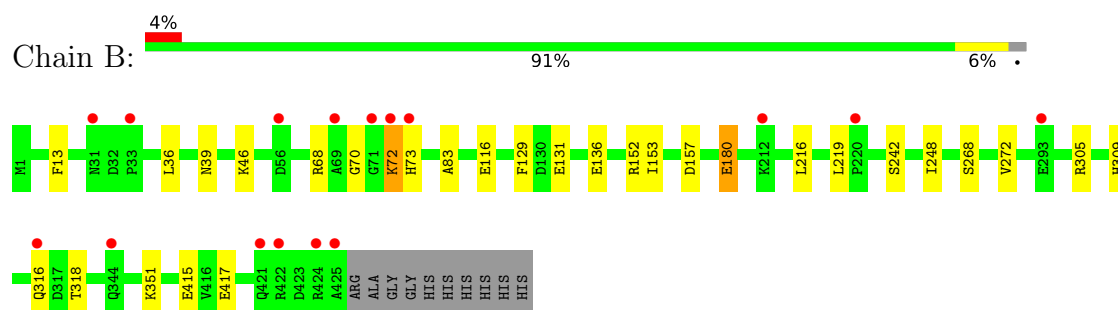
3 Residue-property plots [i](#)

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: Alanine-tRNA ligase



• Molecule 1: Alanine-tRNA ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	78.23Å 55.59Å 125.72Å 90.00° 102.07° 90.00°	Depositor
Resolution (Å)	29.80 – 1.90 29.80 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.80-1.90) 99.9 (29.80-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.65 (at 1.91Å)	Xtriage
Refinement program	PHENIX 1.11.1 _2575	Depositor
R, R_{free}	0.200 , 0.230 0.200 , 0.231	Depositor DCC
R_{free} test set	1999 reflections (2.39%)	wwPDB-VP
Wilson B-factor (Å ²)	28.0	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 33.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7055	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.31% 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: CIT, FMT, ATP, MLT, EDO

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	0.36	0/3427	0.54	0/4640
1	B	0.34	0/3427	0.52	0/4640
All	All	0.35	0/6854	0.53	0/9280

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	3347	0	3237	17	0
1	B	3347	0	3237	22	0
2	A	31	0	12	1	0
2	B	31	0	12	2	0
3	A	6	0	4	2	0
3	B	6	0	4	1	0
4	A	28	0	42	3	0
4	B	32	0	48	5	0
5	A	3	0	1	0	0
5	B	6	0	2	0	0
6	A	9	0	4	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	B	13	0	5	1	0
8	A	95	0	0	2	0
8	B	101	0	0	1	0
All	All	7055	0	6608	41	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 3.

All (41) 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:72:LYS:H	1:A:72:LYS:HD2	1.49	0.77
1:A:210:ARG:NH1	1:A:214:GLY:O	2.25	0.69
1:B:309:HIS:NE2	4:B:508:EDO:H21	2.09	0.68
1:A:72:LYS:HD2	1:A:72:LYS:N	2.10	0.66
1:A:173:GLU:OE1	8:A:601:HOH:O	2.13	0.65
8:A:606:HOH:O	1:B:415:GLU:HG3	1.97	0.65
1:A:68:ARG:NH2	2:A:501:ATP:O1A	2.29	0.63
1:A:13:PHE:CD1	1:A:116:GLU:HG3	2.38	0.58
2:B:501:ATP:O2A	3:B:502:ALA:N	2.36	0.58
1:B:39:ASN:HD22	1:B:73:HIS:CE1	2.23	0.56
1:A:297:TYR:HD2	4:A:504:EDO:H22	1.68	0.56
1:B:13:PHE:CD1	1:B:116:GLU:HG3	2.40	0.56
2:B:501:ATP:O2B	7:B:505:CIT:H41	2.07	0.55
1:B:316:GLN:HB2	4:B:510:EDO:H11	1.89	0.54
1:B:83:ALA:HB1	1:B:242:SER:HA	1.93	0.51
1:A:255:THR:HG21	1:A:326:PRO:HG3	1.93	0.50
1:A:129:PHE:CE2	1:A:131:GLU:HB2	2.48	0.49
1:A:204:VAL:HG21	3:A:502:ALA:HB1	1.94	0.49
1:B:136:GLU:OE1	1:B:152:ARG:HD3	2.12	0.49
1:B:318:THR:HB	4:B:509:EDO:H21	1.95	0.48
1:A:83:ALA:HB1	1:A:242:SER:HA	1.94	0.48
1:B:129:PHE:CE2	1:B:131:GLU:HB2	2.50	0.47
1:A:1:MET:HE1	1:A:117:MET:HE1	1.97	0.46
1:B:68:ARG:HB3	1:B:73:HIS:O	2.16	0.45
1:B:219:LEU:HD23	1:B:219:LEU:HA	1.82	0.45
1:A:355:ARG:C	1:A:355:ARG:HD3	2.43	0.44
1:B:153:ILE:HG21	1:B:157:ASP:HB2	1.99	0.44
1:A:297:TYR:HB3	4:A:504:EDO:H22	2.00	0.44
1:B:305:ARG:HH21	4:B:508:EDO:H22	1.82	0.43
1:B:180:GLU:H	1:B:180:GLU:HG2	1.42	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:297:TYR:CD2	4:A:504:EDO:H22	2.52	0.43
1:A:355:ARG:HD3	1:A:355:ARG:O	2.18	0.43
1:B:73:HIS:O	1:B:73:HIS:ND1	2.52	0.43
1:B:39:ASN:HD22	1:B:73:HIS:CD2	2.37	0.42
1:B:268:SER:O	1:B:272:VAL:HG23	2.19	0.42
1:B:36:LEU:HD12	1:B:70:GLY:HA3	2.01	0.42
1:B:73:HIS:HE1	8:B:665:HOH:O	2.03	0.42
1:B:417:GLU:OE2	4:B:511:EDO:H21	2.20	0.41
1:B:46:LYS:HB2	1:B:46:LYS:HE3	1.80	0.41
1:A:204:VAL:CG2	3:A:502:ALA:HB1	2.50	0.41
1:B:72:LYS:HD3	1:B:72:LYS:HA	1.49	0.40

There are no symmetry-related clashes.

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	423/435 (97%)	410 (97%)	11 (3%)	2 (0%)	24	16
1	B	423/435 (97%)	414 (98%)	9 (2%)	0	100	100
All	All	846/870 (97%)	824 (97%)	20 (2%)	2 (0%)	43	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	73	HIS
1	A	74	ASN

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	348/355 (98%)	343 (99%)	5 (1%)	59	59
1	B	348/355 (98%)	343 (99%)	5 (1%)	59	59
All	All	696/710 (98%)	686 (99%)	10 (1%)	59	59

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	LYS
1	A	156	LYS
1	A	226	THR
1	A	355	ARG
1	A	377	HIS
1	B	72	LYS
1	B	180	GLU
1	B	216	LEU
1	B	248	ILE
1	B	351	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	39	ASN
1	B	250	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

24 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	ATP	A	501	-	32,33,33	1.38	5 (15%)	48,52,52	1.70	10 (20%)
4	EDO	B	506	-	3,3,3	0.51	0	2,2,2	0.16	0
4	EDO	B	509	-	3,3,3	0.42	0	2,2,2	0.42	0
4	EDO	B	512	-	3,3,3	0.45	0	2,2,2	0.48	0
6	MLT	A	509	-	8,8,8	1.17	0	10,10,10	1.81	4 (40%)
4	EDO	A	511	-	3,3,3	0.51	0	2,2,2	0.28	0
2	ATP	B	501	-	32,33,33	1.36	4 (12%)	48,52,52	1.59	8 (16%)
4	EDO	A	506	-	3,3,3	0.57	0	2,2,2	0.33	0
4	EDO	B	513	-	3,3,3	0.48	0	2,2,2	0.51	0
4	EDO	A	505	-	3,3,3	0.49	0	2,2,2	0.29	0
3	ALA	A	502	-	5,5,5	0.96	0	6,6,6	1.64	2 (33%)
4	EDO	A	508	-	3,3,3	0.59	0	2,2,2	0.37	0
4	EDO	A	503	-	3,3,3	0.40	0	2,2,2	0.43	0
5	FMT	A	507	-	2,2,2	0.67	0	1,1,1	0.13	0
4	EDO	B	508	-	3,3,3	0.60	0	2,2,2	0.50	0
7	CIT	B	505	-	12,12,12	1.65	1 (8%)	17,17,17	1.86	4 (23%)
5	FMT	B	504	-	2,2,2	0.69	0	1,1,1	0.26	0
4	EDO	A	510	-	3,3,3	0.31	0	2,2,2	0.70	0
5	FMT	B	503	-	2,2,2	0.60	0	1,1,1	0.27	0
3	ALA	B	502	-	5,5,5	1.09	1 (20%)	6,6,6	1.70	2 (33%)
4	EDO	B	507	-	3,3,3	0.51	0	2,2,2	0.44	0
4	EDO	B	511	-	3,3,3	0.42	0	2,2,2	0.46	0
4	EDO	B	510	-	3,3,3	0.48	0	2,2,2	0.43	0
4	EDO	A	504	-	3,3,3	0.28	0	2,2,2	0.80	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	ATP	A	501	-	-	3/22/38/38	0/3/3/3
4	EDO	B	506	-	-	1/1/1/1	-
4	EDO	B	509	-	-	0/1/1/1	-
4	EDO	B	512	-	-	1/1/1/1	-
6	MLT	A	509	-	-	0/8/8/8	-
4	EDO	A	511	-	-	1/1/1/1	-
2	ATP	B	501	-	-	1/22/38/38	0/3/3/3
4	EDO	A	506	-	-	1/1/1/1	-
4	EDO	B	513	-	-	0/1/1/1	-
4	EDO	A	505	-	-	0/1/1/1	-
3	ALA	A	502	-	-	4/4/4/4	-
4	EDO	A	508	-	-	0/1/1/1	-
4	EDO	A	503	-	-	1/1/1/1	-
4	EDO	B	508	-	-	0/1/1/1	-
7	CIT	B	505	-	-	2/16/16/16	-
4	EDO	A	510	-	-	0/1/1/1	-
3	ALA	B	502	-	-	4/4/4/4	-
4	EDO	B	507	-	-	0/1/1/1	-
4	EDO	B	511	-	-	0/1/1/1	-
4	EDO	B	510	-	-	0/1/1/1	-
4	EDO	A	504	-	-	0/1/1/1	-

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	ATP	C5-C4	4.30	1.46	1.39
2	A	501	ATP	C5-C4	4.20	1.46	1.39
7	B	505	CIT	C3-C6	-4.12	1.49	1.53
2	B	501	ATP	PB-O3B	2.96	1.62	1.59
2	A	501	ATP	PB-O3B	2.88	1.62	1.59
2	B	501	ATP	C8-N7	2.86	1.37	1.31
2	A	501	ATP	C8-N7	2.71	1.36	1.31
2	B	501	ATP	C5-C6	2.32	1.47	1.41
2	A	501	ATP	C5-C6	2.24	1.47	1.41
3	B	502	ALA	OXT-C	-2.20	1.23	1.30
2	A	501	ATP	C4-N9	-2.03	1.33	1.37

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	505	CIT	O7-C3-C6	-5.91	100.57	108.96
2	B	501	ATP	C5-C4-N3	-5.17	119.59	126.72
2	A	501	ATP	C5-C4-N3	-4.91	119.95	126.72
2	B	501	ATP	N3-C4-N9	4.19	134.29	127.17
2	A	501	ATP	N3-C4-N9	4.18	134.28	127.17
2	A	501	ATP	N3-C2-N1	-3.80	122.84	128.58
2	A	501	ATP	C4-N9-C8	3.51	109.43	105.74
2	A	501	ATP	C2-N3-C4	3.46	120.28	111.83
2	B	501	ATP	C2-N3-C4	3.42	120.19	111.83
2	B	501	ATP	N3-C2-N1	-3.21	123.72	128.58
3	A	502	ALA	OXT-C-O	-3.11	117.03	124.08
3	B	502	ALA	OXT-C-O	-3.10	117.04	124.08
6	A	509	MLT	O2-C1-C2	2.89	118.86	112.74
2	B	501	ATP	C4-N9-C8	2.67	108.54	105.74
2	A	501	ATP	C4-C5-N7	-2.54	107.68	110.58
3	B	502	ALA	OXT-C-CA	2.53	122.50	113.79
6	A	509	MLT	O5-C4-O4	-2.46	117.01	123.33
2	B	501	ATP	C4-C5-N7	-2.45	107.78	110.58
7	B	505	CIT	O2-C1-C2	2.43	122.05	114.35
7	B	505	CIT	O7-C3-C4	2.28	114.59	109.38
3	A	502	ALA	OXT-C-CA	2.27	121.59	113.79
2	A	501	ATP	N9-C8-N7	-2.23	110.78	113.94
2	A	501	ATP	C2-N1-C6	2.22	122.38	118.73
2	A	501	ATP	O2B-PB-O1B	2.22	122.77	112.44
7	B	505	CIT	O2-C1-O1	-2.16	117.77	123.33
6	A	509	MLT	O5-C4-C3	2.15	120.69	114.00
6	A	509	MLT	O3-C2-C3	2.14	115.11	109.96
2	B	501	ATP	O3G-PG-O2G	2.04	115.47	107.80
2	A	501	ATP	C6-C5-N7	2.04	136.01	132.09
2	B	501	ATP	O2B-PB-O1B	2.01	121.81	112.44

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	ATP	PB-O3B-PG-O3G
3	A	502	ALA	O-C-CA-N
3	B	502	ALA	OXT-C-CA-N
3	B	502	ALA	OXT-C-CA-CB
3	A	502	ALA	OXT-C-CA-N
3	A	502	ALA	O-C-CA-CB
3	A	502	ALA	OXT-C-CA-CB
3	B	502	ALA	O-C-CA-CB

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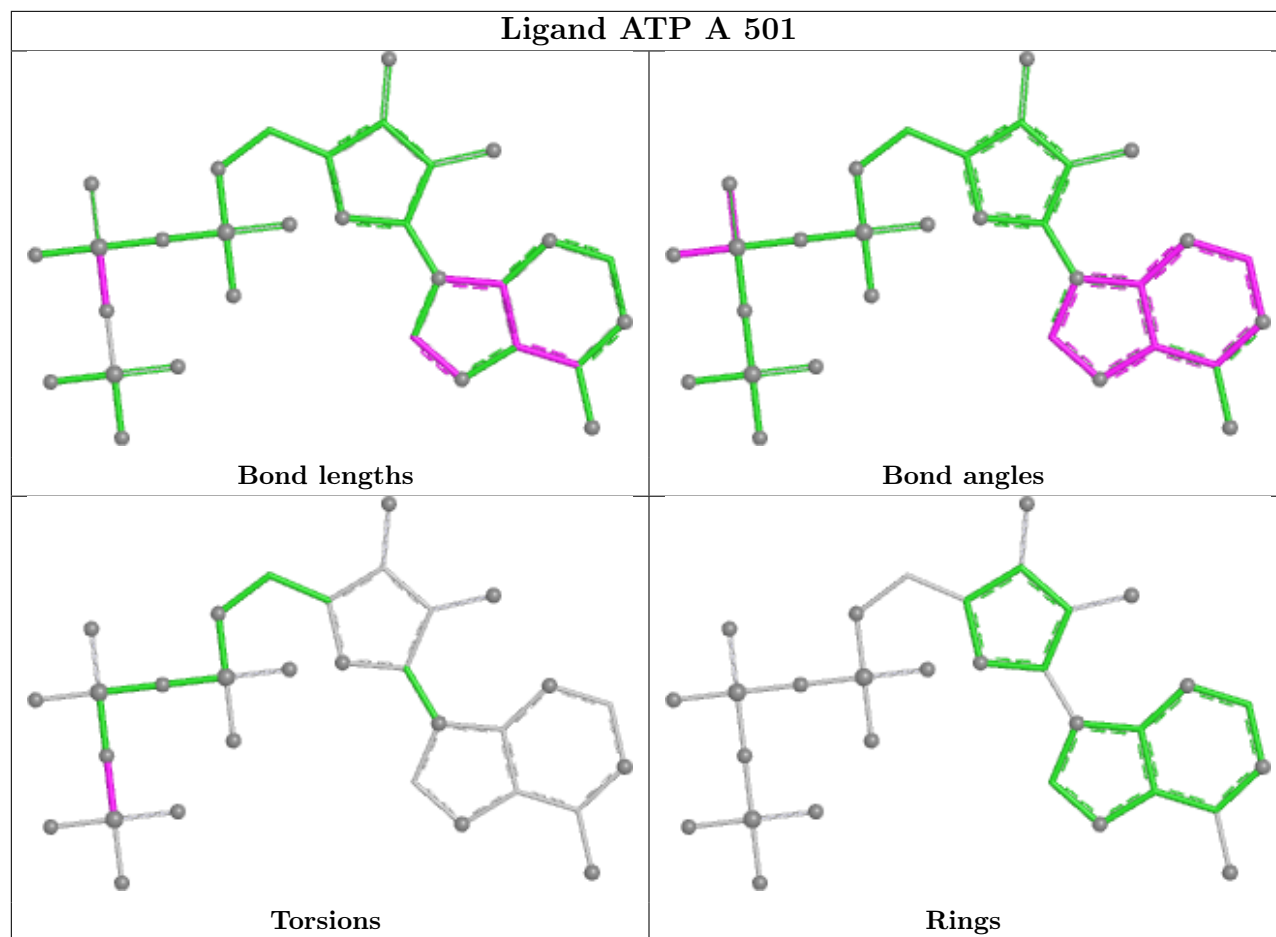
Mol	Chain	Res	Type	Atoms
2	B	501	ATP	PB-O3A-PA-O1A
7	B	505	CIT	C1-C2-C3-C6
4	A	511	EDO	O1-C1-C2-O2
4	A	503	EDO	O1-C1-C2-O2
4	B	512	EDO	O1-C1-C2-O2
2	A	501	ATP	PB-O3B-PG-O1G
2	A	501	ATP	PB-O3B-PG-O2G
4	A	506	EDO	O1-C1-C2-O2
4	B	506	EDO	O1-C1-C2-O2
3	B	502	ALA	O-C-CA-N
7	B	505	CIT	C3-C4-C5-O3

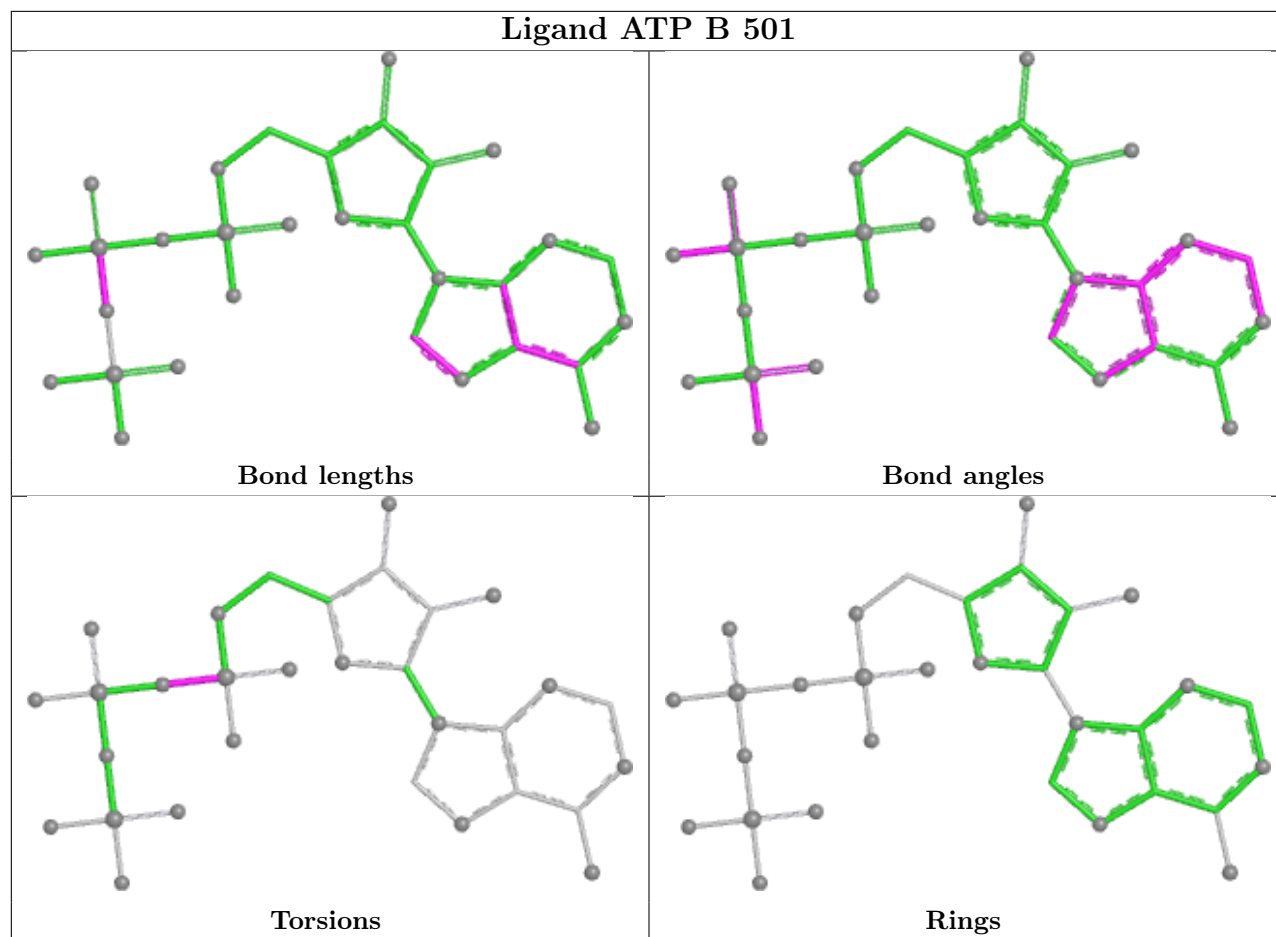
There are no ring outliers.

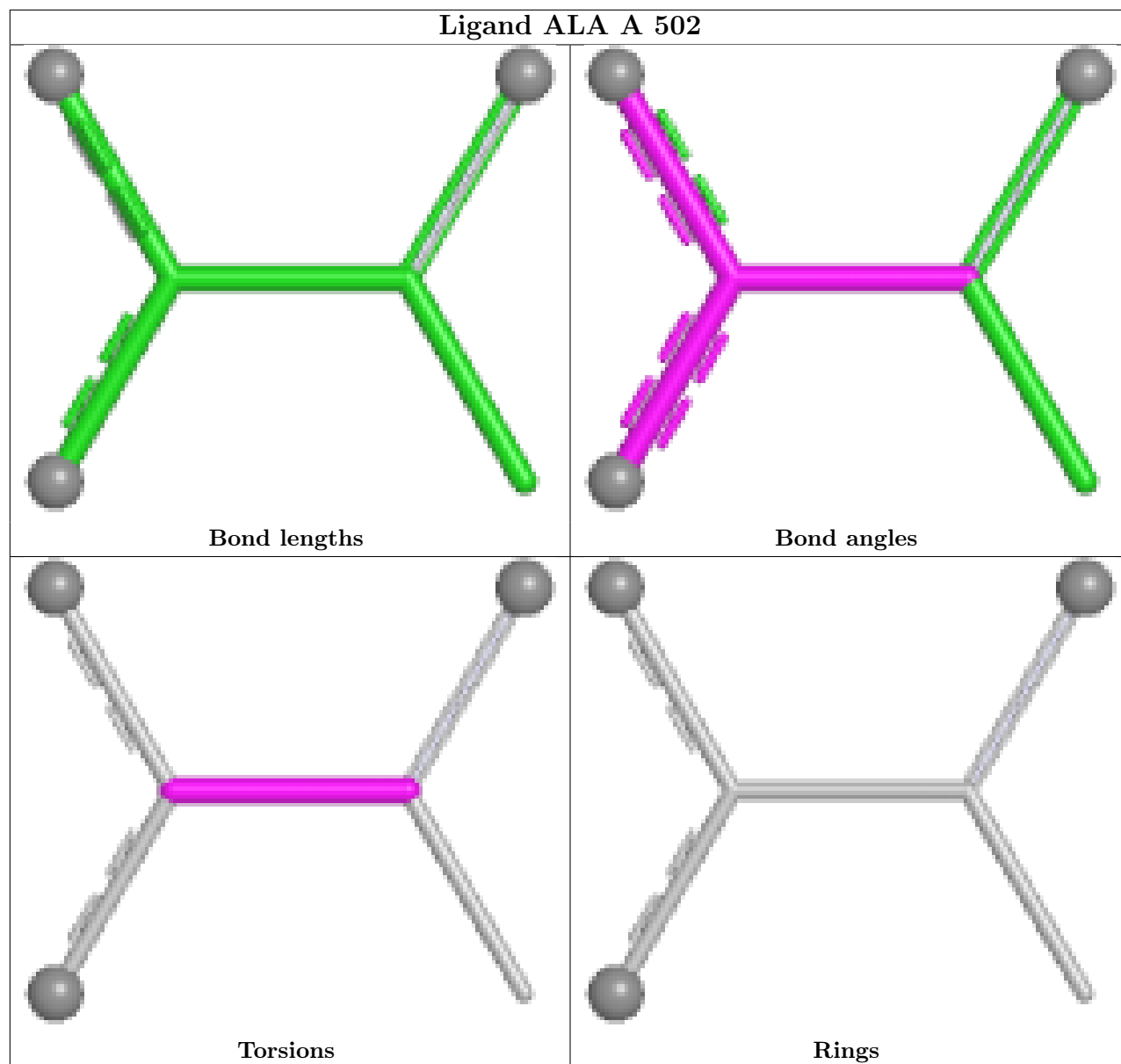
10 monomers are involved in 13 short contacts:

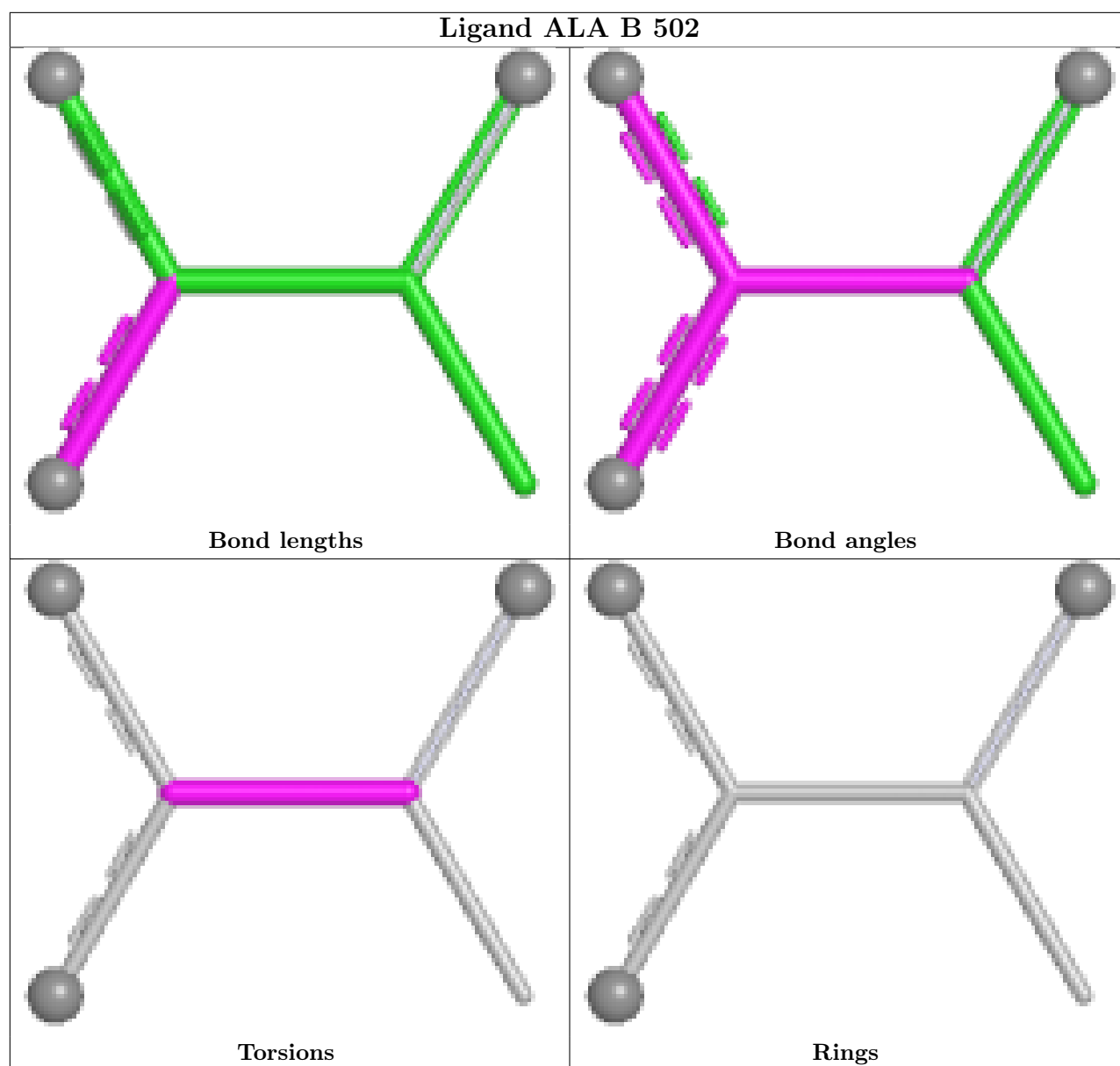
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	ATP	1	0
4	B	509	EDO	1	0
2	B	501	ATP	2	0
3	A	502	ALA	2	0
4	B	508	EDO	2	0
7	B	505	CIT	1	0
3	B	502	ALA	1	0
4	B	511	EDO	1	0
4	B	510	EDO	1	0
4	A	504	EDO	3	0

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

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	425/435 (97%)	0.24	24 (5%) 30 32	18, 27, 49, 65	0
1	B	425/435 (97%)	0.21	16 (3%) 44 47	19, 28, 47, 65	0
All	All	850/870 (97%)	0.22	40 (4%) 36 39	18, 28, 48, 65	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	425	ALA	5.1
1	B	73	HIS	5.1
1	A	73	HIS	5.0
1	A	425	ALA	4.0
1	A	72	LYS	3.9
1	A	33	PRO	3.5
1	A	74	ASN	3.4
1	A	212	LYS	3.4
1	A	31	ASN	3.3
1	A	423	ASP	3.3
1	B	72	LYS	3.1
1	B	212	LYS	3.0
1	B	56	ASP	3.0
1	A	34	THR	2.9
1	B	220	PRO	2.9
1	B	69	ALA	2.9
1	A	54	LYS	2.9
1	B	33	PRO	2.8
1	A	372	ALA	2.8
1	A	213	ASP	2.8
1	B	31	ASN	2.7
1	B	344	GLN	2.5
1	B	71	GLY	2.5
1	B	293	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	422	ARG	2.5
1	A	420	ALA	2.4
1	A	371	VAL	2.4
1	A	56	ASP	2.4
1	B	421	GLN	2.3
1	B	424	ARG	2.3
1	A	377	HIS	2.3
1	A	32	ASP	2.2
1	A	355	ARG	2.2
1	A	316	GLN	2.2
1	B	422	ARG	2.1
1	A	369	ASN	2.1
1	A	180	GLU	2.0
1	A	71	GLY	2.0
1	A	375	ASP	2.0
1	B	316	GLN	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($\text{\AA}^2$)	Q<0.9
3	ALA	A	502	6/6	0.75	0.17	31,39,49,52	0
7	CIT	B	505	13/13	0.78	0.14	34,37,45,49	0
6	MLT	A	509	9/9	0.80	0.14	32,39,43,43	0
4	EDO	A	506	4/4	0.83	0.14	33,35,36,39	0
4	EDO	B	510	4/4	0.83	0.16	39,43,45,47	0
5	FMT	B	504	3/3	0.84	0.12	44,44,47,48	0
3	ALA	B	502	6/6	0.85	0.13	34,37,42,49	0

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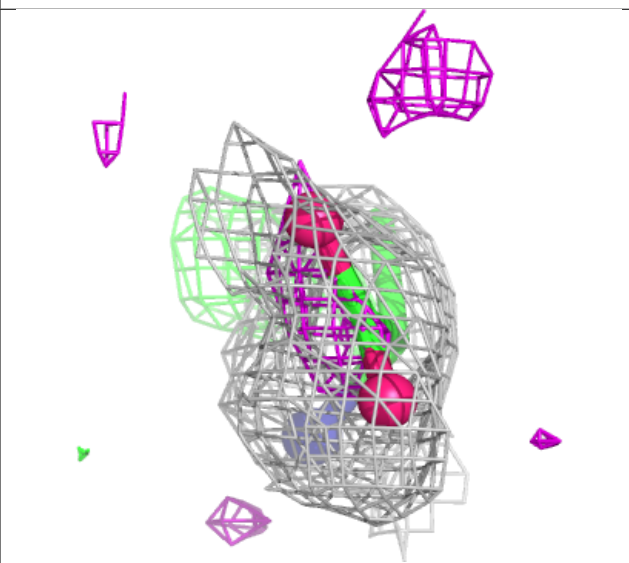
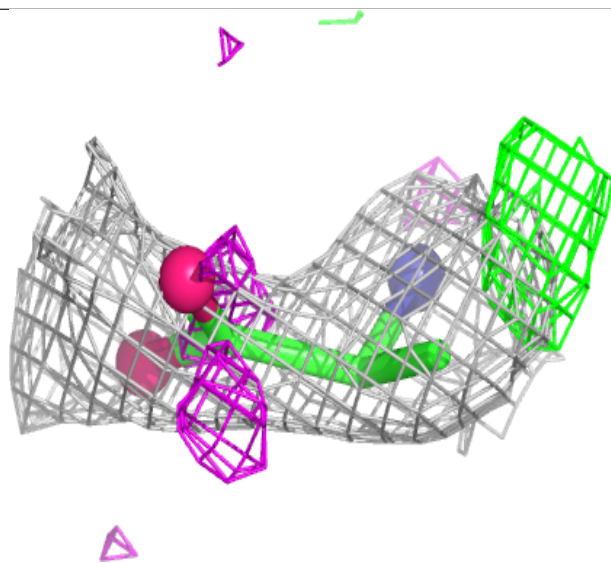
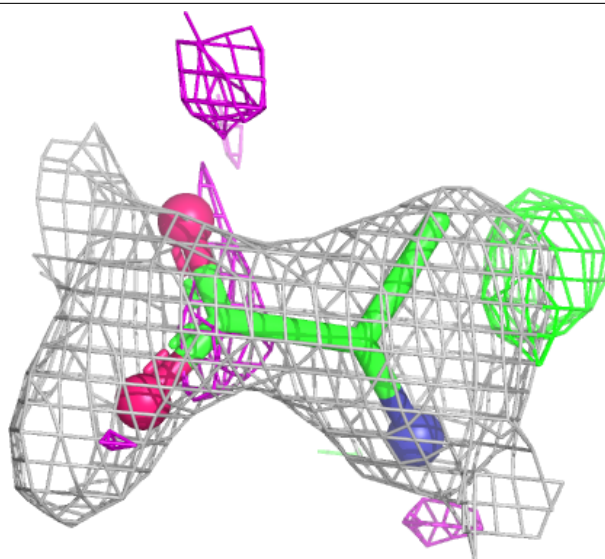
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	EDO	A	510	4/4	0.85	0.18	29,32,36,37	0
5	FMT	B	503	3/3	0.87	0.12	31,31,39,43	0
4	EDO	B	508	4/4	0.88	0.18	24,26,32,35	0
4	EDO	B	513	4/4	0.88	0.11	31,34,37,42	0
4	EDO	B	509	4/4	0.88	0.20	29,29,33,38	0
4	EDO	A	508	4/4	0.89	0.12	25,30,33,41	0
4	EDO	B	506	4/4	0.90	0.11	34,34,35,41	0
4	EDO	B	507	4/4	0.90	0.10	34,35,35,45	0
4	EDO	B	511	4/4	0.91	0.14	32,36,38,43	0
4	EDO	A	504	4/4	0.91	0.15	30,31,34,34	0
5	FMT	A	507	3/3	0.92	0.10	35,35,40,44	0
2	ATP	B	501	31/31	0.93	0.09	20,25,69,71	0
4	EDO	A	511	4/4	0.93	0.12	29,33,35,36	0
4	EDO	A	503	4/4	0.94	0.09	31,32,32,33	0
4	EDO	A	505	4/4	0.94	0.09	25,29,30,31	0
2	ATP	A	501	31/31	0.95	0.07	16,22,58,63	0
4	EDO	B	512	4/4	0.95	0.10	34,34,35,36	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.

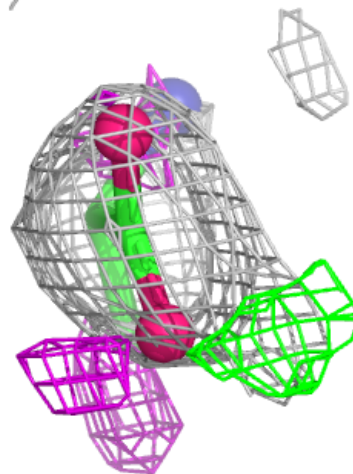
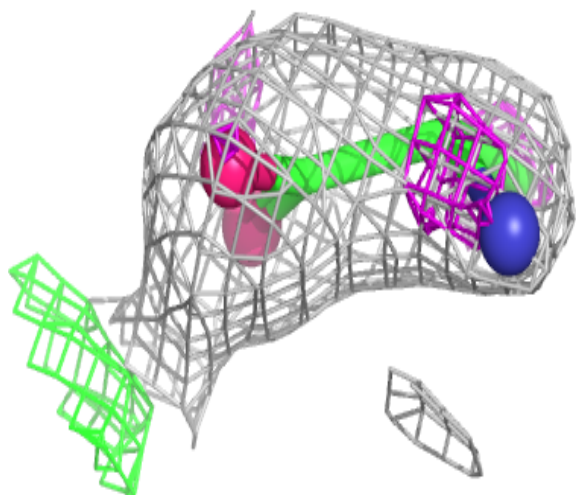
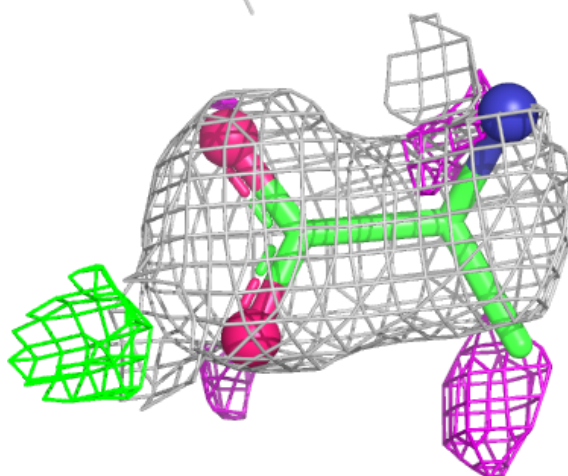
Electron density around ALA A 502:

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



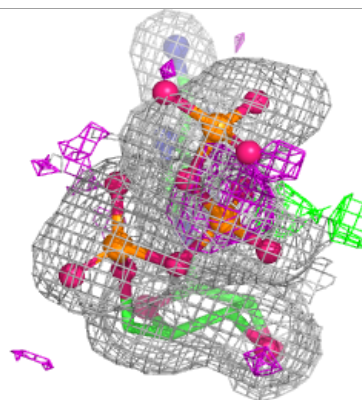
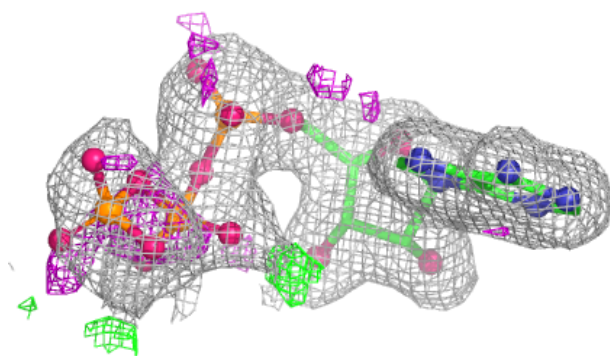
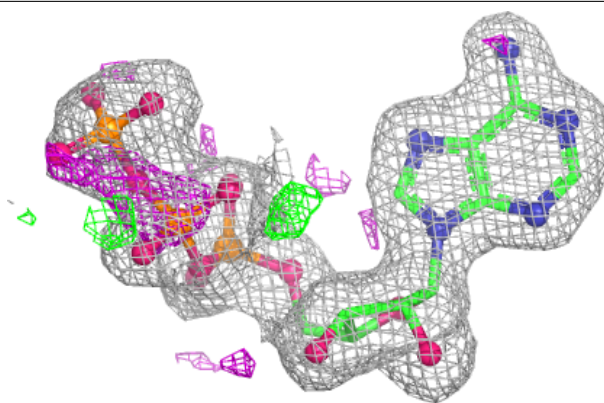
Electron density around ALA B 502:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

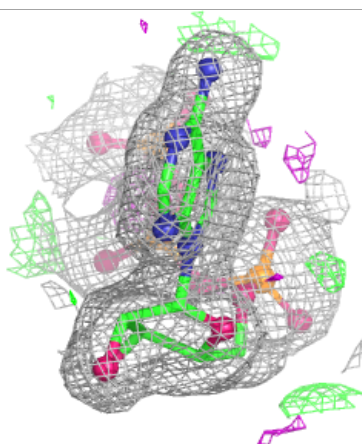
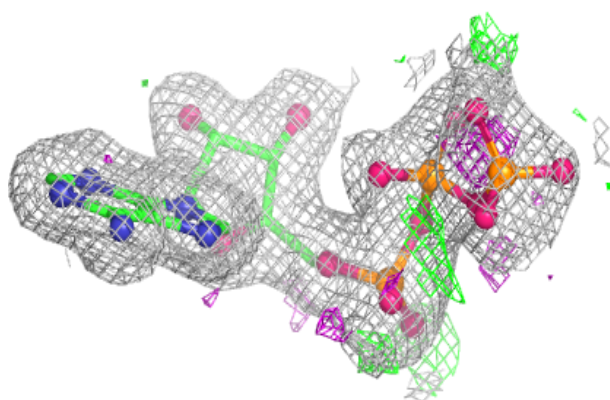
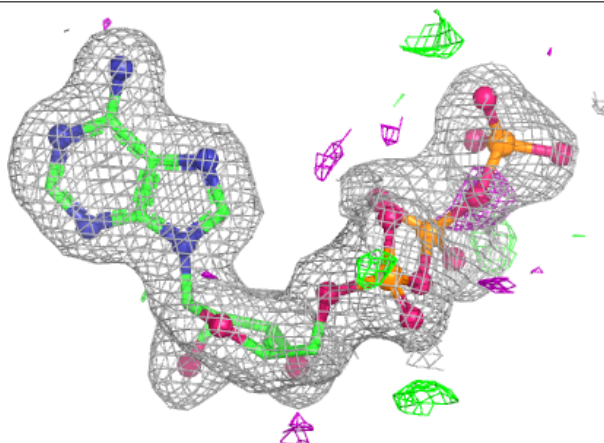


Electron density around ATP B 501:

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

**Electron density around ATP A 501:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.