



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 09:38 PM UTC

PDB ID : 3OV7 / pdb_00003ov7
Title : How the CCA-Adding Enzyme Selects Adenine over Cytosine in Position 76 of tRNA
Authors : Pan, B.C.; Xiong, Y.; Steitz, T.A.
Deposited on : 2010-09-15
Resolution : 3.00 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

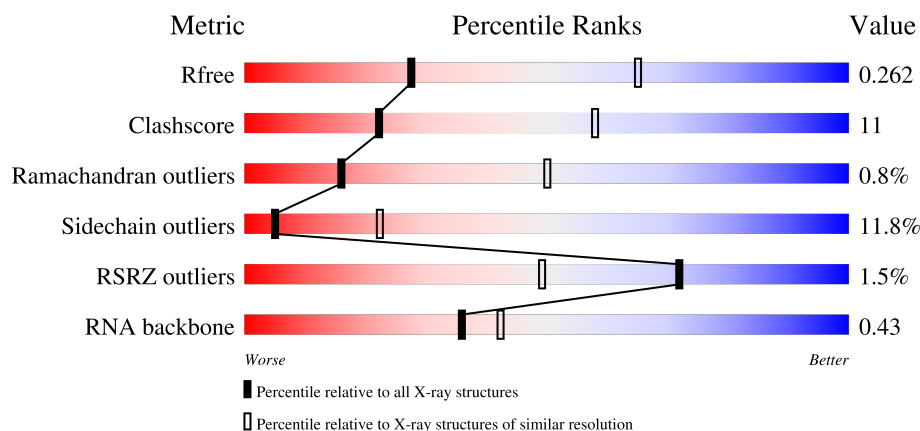
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)
RNA backbone	3983	1109 (3.20-2.80)

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	441	71% 25% •
1	B	441	2% 66% 29% 5% •
2	C	34	62% 32% 6%
2	D	34	6% 44% 29% 24% •

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 8853 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 CCA-Adding Enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	441	Total	C	N	O	S	0	0	0
			3655	2347	636	655	17			
1	B	437	Total	C	N	O	S	0	0	0
			3629	2333	632	651	13			

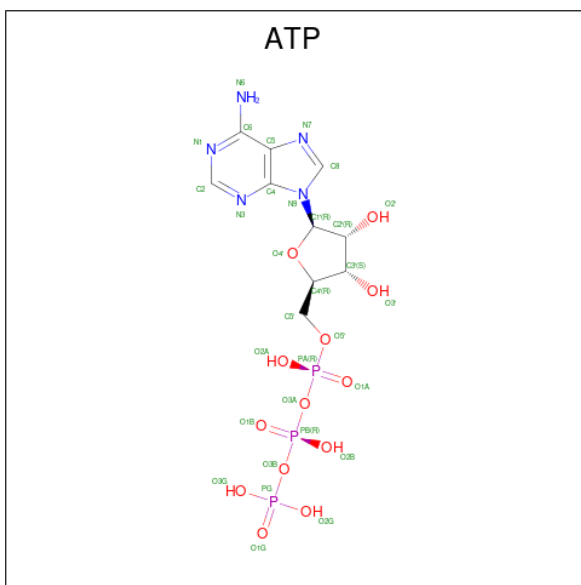
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	438	CYS	-	expression tag	UNP O28126
A	439	CYS	-	expression tag	UNP O28126
A	440	CYS	-	expression tag	UNP O28126
A	441	MET	-	expression tag	UNP O28126
B	438	CYS	-	expression tag	UNP O28126
B	439	CYS	-	expression tag	UNP O28126
B	440	CYS	-	expression tag	UNP O28126
B	441	MET	-	expression tag	UNP O28126

- Molecule 2 is a RNA chain called RNA (34-MER).

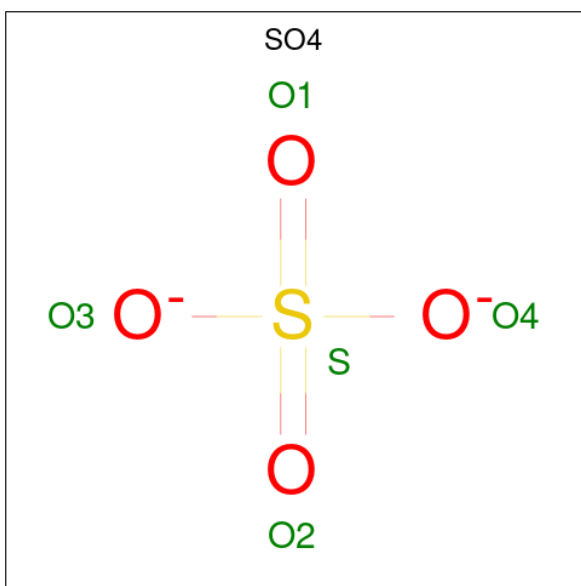
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	34	Total	C	N	O	P	3	0	0
			720	322	124	240	34			
2	D	34	Total	C	N	O	P	3	0	0
			720	322	124	240	34			

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



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

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O_4S).

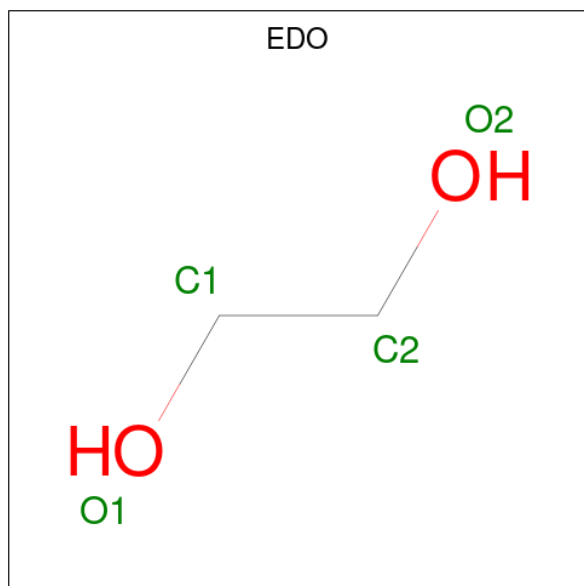


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0

- Molecule 5 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mn	0	0
			1	1		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

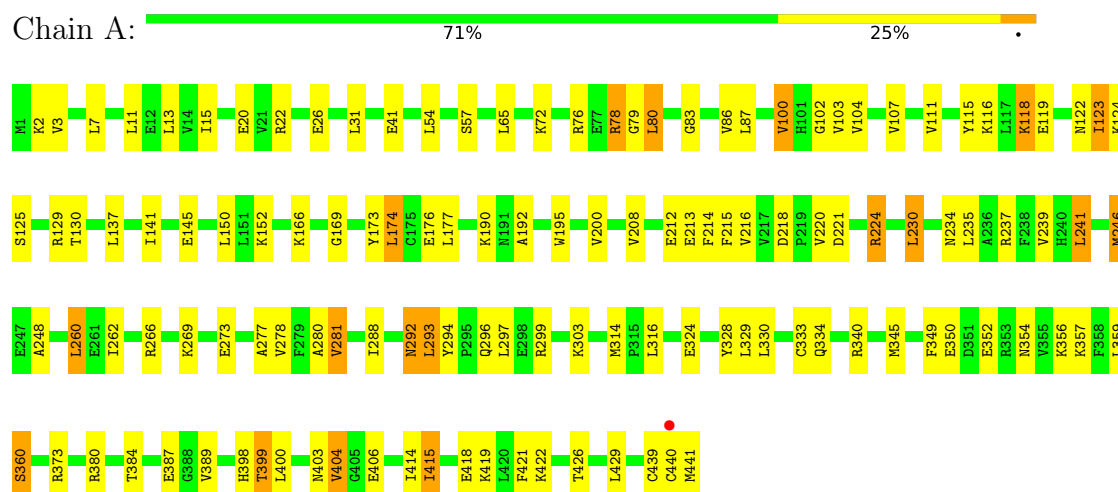
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	37	Total	O	0	0
			37	37		
7	B	8	Total	O	0	0
			8	8		
7	C	6	Total	O	0	0
			6	6		
7	D	1	Total	O	0	0
			1	1		

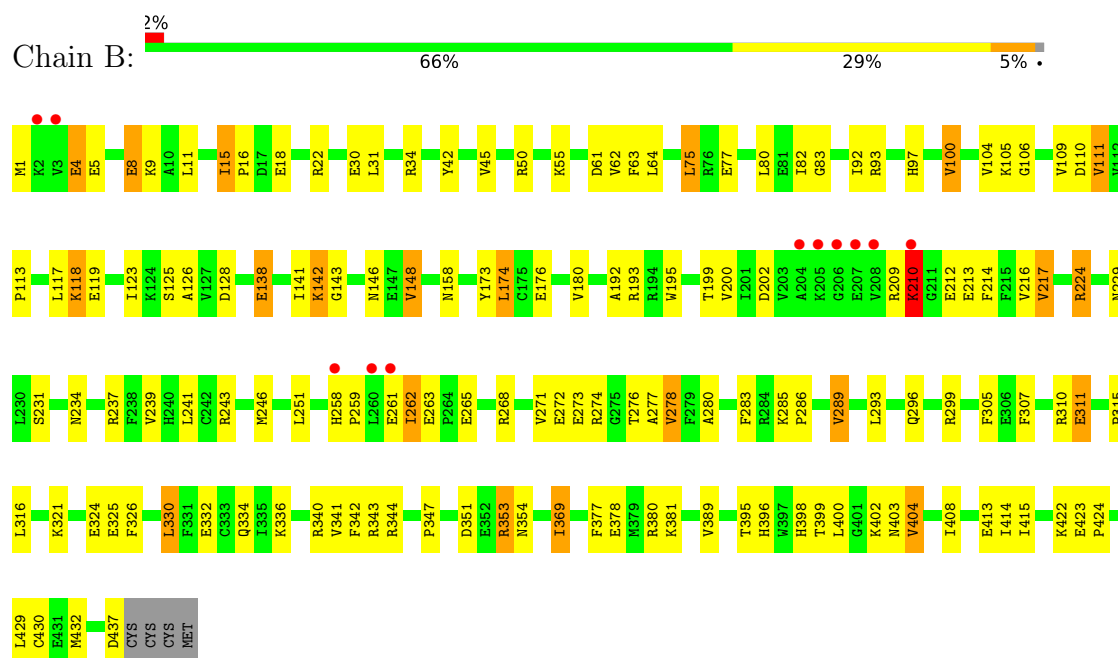
3 Residue-property plots

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.

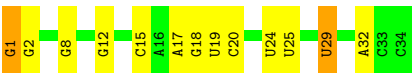
• Molecule 1: CCA-Adding Enzyme



• Molecule 1: CCA-Adding Enzyme



• Molecule 2: RNA (34-MER)



● Molecule 2: RNA (34-MER)



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	111.39Å 216.09Å 58.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.51 – 3.00 49.51 – 3.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.51-3.00) 99.9 (49.51-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.205 , 0.267 0.215 , 0.262	Depositor DCC
R_{free} test set	1602 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	50.0	Xtriage
Anisotropy	0.291	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 64.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	8853	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.34% 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: MN, SO4, ATP, 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	0.64	0/3738	0.95	5/5021 (0.1%)
1	B	0.61	3/3712 (0.1%)	0.90	3/4987 (0.1%)
2	C	2.19	2/803 (0.2%)	1.18	4/1246 (0.3%)
2	D	1.61	2/803 (0.2%)	1.10	3/1246 (0.2%)
All	All	0.99	7/9056 (0.1%)	0.97	15/12500 (0.1%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1	G	OP3-P	53.20	2.54	1.48
2	D	1	G	P-O5'	35.95	2.31	1.59
2	C	1	G	P-O5'	26.73	2.13	1.59
2	D	1	G	OP3-P	-23.06	1.02	1.48
1	B	210	LYS	CD-CE	9.07	1.79	1.52
1	B	4	GLU	CD-OE1	7.55	1.39	1.25
1	B	4	GLU	CD-OE2	5.30	1.35	1.25

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	29	U	C5'-C4'-C3'	-7.72	104.41	116.00
2	C	1	G	O5'-P-OP1	-6.87	90.11	110.70
1	A	130	THR	CA-C-N	-6.80	112.69	119.56
1	A	130	THR	C-N-CA	-6.80	112.69	119.56
2	D	17	A	C1'-O4'-C4'	-6.66	103.04	109.70
2	C	1	G	P-O5'-C5'	-6.14	102.46	120.90
2	D	17	A	C2'-C3'-O3'	5.65	117.97	109.50
1	A	248	ALA	CA-C-N	5.37	125.31	119.78
1	A	248	ALA	C-N-CA	5.37	125.31	119.78
1	B	210	LYS	CD-CE-NZ	-5.30	94.93	111.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	258	HIS	CA-C-N	5.18	126.31	119.84
1	B	258	HIS	C-N-CA	5.18	126.31	119.84
2	C	32	A	N9-C1'-C2'	5.18	119.76	112.00
2	D	32	A	N9-C1'-C2'	5.16	119.74	112.00
1	A	235	LEU	N-CA-C	-5.13	105.59	111.14

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	3655	0	3657	75	0
1	B	3629	0	3633	102	0
2	C	720	0	366	12	0
2	D	720	0	366	14	0
3	A	31	0	12	0	0
3	B	31	0	12	2	0
4	A	5	0	0	1	0
4	B	5	0	0	0	0
5	A	1	0	0	0	0
6	B	4	0	6	3	0
7	A	37	0	0	2	0
7	B	8	0	0	0	0
7	C	6	0	0	1	0
7	D	1	0	0	0	0
All	All	8853	0	8052	178	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 11.

All (178) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:LYS:CD	1:B:210:LYS:CE	1.79	1.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:GLY:HA3	1:A:100:VAL:HG11	1.37	1.06
1:B:353:ARG:HG2	1:B:353:ARG:HH11	1.24	1.00
1:B:400:LEU:O	1:B:404:VAL:HG23	1.80	0.82
1:B:5:GLU:HA	1:B:8:GLU:HG3	1.62	0.81
2:D:17:A:H4'	2:D:18:G:OP2	1.81	0.80
1:B:210:LYS:CE	1:B:210:LYS:CG	2.64	0.76
1:B:311:GLU:HA	1:B:311:GLU:OE1	1.85	0.76
1:B:210:LYS:CD	1:B:210:LYS:NZ	2.50	0.75
1:A:299:ARG:NH1	1:A:399:THR:O	2.20	0.74
1:A:296:GLN:HE22	1:A:403:ASN:HD22	1.35	0.74
1:B:123:ILE:HG23	1:B:128:ASP:HB2	1.72	0.72
1:A:83:GLY:HA3	1:A:100:VAL:CG1	2.18	0.72
1:A:212:GLU:HA	1:A:212:GLU:OE1	1.91	0.71
1:B:351:ASP:OD1	1:B:354:ASN:HB2	1.92	0.68
1:B:100:VAL:HG13	1:B:111:VAL:HG13	1.75	0.68
1:B:93:ARG:HD2	1:B:289:VAL:HG13	1.75	0.67
2:D:25:U:H2'	2:D:26:A:C8	2.30	0.67
1:A:380:ARG:HH22	1:B:334:GLN:HE22	1.41	0.67
1:A:293:LEU:HD22	1:A:297:LEU:HG	1.77	0.66
1:A:334:GLN:NE2	1:B:380:ARG:HH12	1.93	0.66
1:B:353:ARG:HH11	1:B:353:ARG:CG	2.04	0.66
1:B:104:VAL:C	1:B:106:GLY:H	2.03	0.65
1:A:418:GLU:HA	1:A:421:PHE:CD1	2.30	0.65
1:B:354:ASN:HD22	2:D:15:C:H6	1.43	0.64
1:A:292:ASN:ND2	2:C:1:G:H1'	2.13	0.63
1:A:352:GLU:HG3	1:B:243:ARG:HD3	1.81	0.62
1:B:83:GLY:HA3	1:B:100:VAL:HG11	1.82	0.62
1:A:400:LEU:O	1:A:404:VAL:HG22	2.00	0.62
1:A:83:GLY:CA	1:A:100:VAL:HG11	2.23	0.61
1:B:305:PHE:CE1	1:B:315:PRO:HB2	2.35	0.61
1:B:395:THR:OG1	1:B:396:HIS:HD2	1.84	0.61
1:A:334:GLN:HE22	1:B:380:ARG:HH22	1.47	0.60
7:A:475:HOH:O	1:B:340:ARG:HD2	2.02	0.59
1:A:314:MET:HB3	1:A:334:GLN:HB2	1.84	0.59
1:A:439:CYS:SG	6:B:801:EDO:H22	2.42	0.59
1:A:292:ASN:HD21	2:C:1:G:H1'	1.67	0.59
1:B:274:ARG:NH1	1:B:332:GLU:OE2	2.36	0.59
1:A:22:ARG:O	1:A:26:GLU:HG2	2.02	0.58
1:B:246:MET:HA	1:B:246:MET:HE2	1.85	0.58
1:A:266:ARG:HA	1:A:269:LYS:HE2	1.84	0.58
1:B:97:HIS:HE1	1:B:113:PRO:O	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:GLY:HA2	1:A:230:LEU:HD23	1.86	0.58
1:B:398:HIS:HD2	2:D:23:U:OP1	1.87	0.57
2:D:25:U:H2'	2:D:26:A:H8	1.68	0.57
1:B:402:LYS:HG2	2:D:1:G:H5''	1.86	0.57
1:A:354:ASN:HD22	2:C:15:C:H6	1.53	0.57
1:B:148:VAL:CG1	1:B:176:GLU:HA	2.35	0.56
1:B:353:ARG:HG2	1:B:353:ARG:NH1	2.05	0.56
1:B:347:PRO:HB3	2:D:15:C:C5	2.40	0.56
1:A:303:LYS:NZ	4:A:601:SO4:O2	2.32	0.56
1:A:237:ARG:HG3	1:B:369:ILE:HD12	1.86	0.56
1:B:15:ILE:O	1:B:55:LYS:HD2	2.05	0.56
1:B:138:GLU:O	1:B:142:LYS:HG2	2.06	0.55
1:A:115:TYR:CE1	1:A:124:LYS:HB3	2.41	0.55
1:B:75:LEU:HD22	1:B:113:PRO:HB2	1.87	0.55
1:A:292:ASN:HD21	2:C:1:G:C1'	2.19	0.55
1:B:262:ILE:HG21	1:B:430:CYS:SG	2.47	0.54
1:B:296:GLN:HE22	1:B:403:ASN:HD22	1.55	0.54
1:B:110:ASP:OD1	2:D:34:C:H5''	2.07	0.54
1:A:123:ILE:CG2	1:A:125:SER:O	2.56	0.54
1:A:54:LEU:HD11	1:A:152:LYS:HB3	1.90	0.54
1:A:440:CYS:SG	1:A:441:MET:N	2.81	0.54
1:B:148:VAL:HG11	1:B:176:GLU:HA	1.89	0.53
1:A:76:ARG:O	1:A:80:LEU:HD22	2.09	0.53
1:A:281:VAL:HG13	1:A:329:LEU:HB2	1.90	0.53
1:A:129:ARG:CZ	1:A:224:ARG:HG3	2.39	0.53
1:A:234:ASN:HD22	1:A:237:ARG:HD2	1.73	0.53
1:A:118:LYS:HD3	1:A:119:GLU:HG2	1.90	0.52
1:B:429:LEU:HD23	1:B:432:MET:HE1	1.90	0.52
1:A:334:GLN:HE21	1:B:380:ARG:HH12	1.56	0.52
1:B:4:GLU:O	1:B:8:GLU:HG2	2.09	0.52
1:B:30:GLU:HG3	1:B:34:ARG:HE	1.75	0.52
1:B:45:VAL:O	1:B:61:ASP:HB2	2.09	0.52
1:A:345:MET:HG3	1:A:373:ARG:HD3	1.91	0.52
1:B:429:LEU:HA	1:B:432:MET:HE2	1.91	0.52
1:A:116:LYS:HA	7:A:476:HOH:O	2.10	0.52
1:B:396:HIS:HE1	2:D:21:C:O2'	1.93	0.52
1:A:190:LYS:HG3	1:A:246:MET:SD	2.50	0.52
1:A:415:ILE:HG23	1:A:419:LYS:HD2	1.92	0.51
1:B:234:ASN:ND2	1:B:237:ARG:HH11	2.08	0.51
1:A:288:ILE:HD12	1:A:292:ASN:HB3	1.91	0.51
1:A:426:THR:HA	1:A:429:LEU:HD12	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:299:ARG:NH1	1:B:399:THR:O	2.44	0.51
1:A:218:ASP:HB3	1:A:221:ASP:O	2.11	0.50
2:D:16:A:H3'	2:D:17:A:H5''	1.93	0.50
1:A:439:CYS:HB2	6:B:801:EDO:H22	1.92	0.50
1:B:50:ARG:NH2	3:B:502:ATP:O3'	2.45	0.50
1:B:325:GLU:HG3	1:B:326:PHE:CD1	2.46	0.50
1:B:64:LEU:HB2	1:B:113:PRO:HA	1.93	0.50
1:B:64:LEU:HD21	1:B:82:ILE:HG21	1.93	0.50
1:B:18:GLU:HB3	1:B:22:ARG:HH21	1.77	0.49
1:B:104:VAL:O	1:B:106:GLY:N	2.45	0.49
1:B:199:THR:O	1:B:214:PHE:HA	2.13	0.49
1:B:123:ILE:CG2	1:B:128:ASP:HB2	2.39	0.49
1:A:292:ASN:ND2	2:C:1:G:C1'	2.76	0.49
1:B:429:LEU:HD23	1:B:432:MET:CE	2.43	0.49
1:B:104:VAL:C	1:B:106:GLY:N	2.71	0.48
2:C:24:U:H2'	2:C:25:U:H6	1.78	0.48
1:B:118:LYS:HD3	1:B:119:GLU:N	2.28	0.48
1:B:296:GLN:HE22	1:B:403:ASN:ND2	2.11	0.48
1:A:87:LEU:HD13	1:A:102:GLY:HA3	1.94	0.48
1:B:423:GLU:HB3	1:B:424:PRO:HD2	1.95	0.48
2:C:24:U:H2'	2:C:25:U:C6	2.49	0.48
1:A:200:VAL:HG22	1:A:215:PHE:HB3	1.96	0.48
1:B:237:ARG:O	1:B:241:LEU:HB2	2.13	0.48
1:B:192:ALA:O	1:B:195:TRP:HB2	2.14	0.48
1:B:5:GLU:HA	1:B:8:GLU:CG	2.40	0.47
1:B:277:ALA:HB2	1:B:336:LYS:HG2	1.95	0.47
2:C:29:U:H5''	2:C:29:U:C6	2.50	0.47
1:A:11:LEU:HD22	1:A:15:ILE:HD11	1.96	0.47
1:A:380:ARG:HH12	1:B:334:GLN:NE2	2.13	0.47
1:B:118:LYS:HD3	1:B:119:GLU:HB2	1.95	0.47
1:B:342:PHE:HB3	1:B:380:ARG:NH1	2.29	0.47
1:B:271:VAL:CG1	1:B:278:VAL:CG1	2.93	0.47
1:B:307:PHE:HA	1:B:310:ARG:NH1	2.30	0.46
1:B:296:GLN:HB3	1:B:404:VAL:HG22	1.98	0.46
1:A:418:GLU:HA	1:A:421:PHE:HD1	1.79	0.45
1:B:280:ALA:HB2	1:B:330:LEU:HD23	1.98	0.45
1:B:9:LYS:HD3	1:B:251:LEU:HD22	1.98	0.45
1:A:65:LEU:HB3	1:A:116:LYS:HB2	1.99	0.45
1:B:202:ASP:OD2	1:B:217:VAL:HG21	2.16	0.45
1:A:174:LEU:HD21	1:A:216:VAL:HG21	1.98	0.45
1:B:261:GLU:O	1:B:262:ILE:C	2.59	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:260:LEU:HD12	1:A:260:LEU:HA	1.82	0.45
1:A:328:TYR:N	1:A:328:TYR:CD1	2.84	0.45
1:B:158:ASN:CG	1:B:241:LEU:HD21	2.42	0.45
1:A:31:LEU:HD13	1:A:104:VAL:HG21	1.99	0.45
1:B:396:HIS:O	1:B:399:THR:HG22	2.17	0.45
1:B:400:LEU:O	1:B:404:VAL:CG2	2.60	0.45
1:A:78:ARG:O	1:A:79:GLY:C	2.60	0.44
1:B:224:ARG:NH1	2:D:31:C:OP1	2.48	0.44
1:A:354:ASN:ND2	2:C:15:C:H6	2.15	0.44
2:C:29:U:H5'	2:C:29:U:H6	1.82	0.44
1:B:176:GLU:O	1:B:180:VAL:HG23	2.18	0.44
1:A:357:LYS:O	1:A:360:SER:HB3	2.18	0.44
1:B:271:VAL:HG11	1:B:278:VAL:CG1	2.48	0.44
1:B:125:SER:HB3	1:B:128:ASP:OD1	2.18	0.43
1:A:7:LEU:HD22	1:A:150:LEU:HD13	2.01	0.43
1:B:174:LEU:HD21	1:B:216:VAL:HG21	2.00	0.43
1:A:316:LEU:O	1:B:343:ARG:NH1	2.51	0.43
1:B:1:MET:HE2	1:B:5:GLU:HB3	2.00	0.43
1:B:193:ARG:HG2	1:B:243:ARG:HG3	2.00	0.43
1:A:192:ALA:HA	1:A:195:TRP:CE2	2.54	0.43
1:B:285:LYS:HA	1:B:286:PRO:HD2	1.93	0.43
1:B:342:PHE:HZ	1:B:344:ARG:HH21	1.66	0.43
1:A:384:THR:OG1	1:A:387:GLU:HG3	2.18	0.42
1:B:229:ASN:HD22	2:D:32:A:H5'	1.84	0.42
1:A:277:ALA:HB3	1:A:333:CYS:HB2	2.02	0.42
1:B:341:VAL:CG2	1:B:377:PHE:HB3	2.49	0.42
1:A:123:ILE:HG23	1:A:125:SER:O	2.18	0.42
1:B:15:ILE:HD11	1:B:146:ASN:HD21	1.84	0.42
1:A:213:GLU:HG3	1:A:214:PHE:O	2.20	0.42
1:A:241:LEU:HD13	1:A:241:LEU:HA	1.89	0.42
1:A:398:HIS:HB3	7:C:45:HOH:O	2.20	0.42
1:B:342:PHE:O	1:B:377:PHE:HA	2.20	0.42
1:A:357:LYS:HB3	2:C:15:C:H4'	2.02	0.41
1:B:265:GLU:OE2	1:B:268:ARG:NH2	2.54	0.41
2:D:9:A:H2'	2:D:10:U:O4'	2.20	0.41
1:B:15:ILE:HA	1:B:16:PRO:HD3	1.78	0.41
1:B:77:GLU:HA	1:B:77:GLU:OE1	2.19	0.41
1:B:283:PHE:HE1	1:B:404:VAL:HG12	1.85	0.41
1:A:129:ARG:NE	1:A:224:ARG:HG3	2.35	0.41
1:A:173:TYR:CE2	1:A:177:LEU:HD11	2.55	0.41
1:A:280:ALA:HA	1:A:329:LEU:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:GLN:HG2	2:C:2:G:H5'	2.02	0.41
1:A:349:PHE:CE1	1:A:350:GLU:HG3	2.55	0.41
1:B:125:SER:OG	1:B:126:ALA:N	2.52	0.41
1:A:439:CYS:CB	6:B:801:EDO:H22	2.50	0.41
1:B:200:VAL:HB	1:B:209:ARG:HB3	2.03	0.41
1:A:166:LYS:HG3	1:A:294:TYR:CE1	2.56	0.41
1:A:280:ALA:HB2	1:A:330:LEU:HD23	2.03	0.41
1:B:173:TYR:HB2	3:B:502:ATP:H2'	2.03	0.41
1:B:305:PHE:CZ	1:B:315:PRO:HG2	2.56	0.41
1:B:42:TYR:HA	1:B:63:PHE:O	2.22	0.40
1:B:141:ILE:O	1:B:143:GLY:N	2.55	0.40
1:A:380:ARG:HH22	1:B:334:GLN:NE2	2.11	0.40
2:D:4:A:H61	2:D:28:U:H3	1.70	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	439/441 (100%)	415 (94%)	23 (5%)	1 (0%)	43	76
1	B	435/441 (99%)	403 (93%)	26 (6%)	6 (1%)	9	36
All	All	874/882 (99%)	818 (94%)	49 (6%)	7 (1%)	16	50

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	122	ASN
1	B	142	LYS
1	B	212	GLU
1	B	105	LYS
1	B	324	GLU

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Mol	Chain	Res	Type
1	B	259	PRO
1	B	262	ILE

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	391/391 (100%)	344 (88%)	47 (12%)	5	22
1	B	387/391 (99%)	342 (88%)	45 (12%)	5	23
All	All	778/782 (100%)	686 (88%)	92 (12%)	5	22

All (92) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	3	VAL
1	A	13	LEU
1	A	20	GLU
1	A	41	GLU
1	A	57	SER
1	A	72	LYS
1	A	78	ARG
1	A	80	LEU
1	A	86	VAL
1	A	100	VAL
1	A	103	VAL
1	A	107	VAL
1	A	111	VAL
1	A	118	LYS
1	A	123	ILE
1	A	137	LEU
1	A	141	ILE
1	A	145	GLU
1	A	174	LEU
1	A	176	GLU

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Mol	Chain	Res	Type
1	A	208	VAL
1	A	220	VAL
1	A	224	ARG
1	A	230	LEU
1	A	239	VAL
1	A	241	LEU
1	A	246	MET
1	A	260	LEU
1	A	262	ILE
1	A	273	GLU
1	A	278	VAL
1	A	281	VAL
1	A	292	ASN
1	A	293	LEU
1	A	324	GLU
1	A	340	ARG
1	A	356	LYS
1	A	359	LEU
1	A	360	SER
1	A	389	VAL
1	A	399	THR
1	A	404	VAL
1	A	406	GLU
1	A	414	ILE
1	A	415	ILE
1	A	422	LYS
1	B	8	GLU
1	B	11	LEU
1	B	15	ILE
1	B	31	LEU
1	B	62	VAL
1	B	75	LEU
1	B	80	LEU
1	B	92	ILE
1	B	100	VAL
1	B	109	VAL
1	B	111	VAL
1	B	117	LEU
1	B	118	LYS
1	B	138	GLU
1	B	148	VAL
1	B	174	LEU

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Mol	Chain	Res	Type
1	B	210	LYS
1	B	213	GLU
1	B	217	VAL
1	B	224	ARG
1	B	231	SER
1	B	239	VAL
1	B	263	GLU
1	B	272	GLU
1	B	273	GLU
1	B	276	THR
1	B	278	VAL
1	B	289	VAL
1	B	293	LEU
1	B	311	GLU
1	B	316	LEU
1	B	321	LYS
1	B	330	LEU
1	B	353	ARG
1	B	369	ILE
1	B	378	GLU
1	B	381	LYS
1	B	389	VAL
1	B	404	VAL
1	B	408	ILE
1	B	413	GLU
1	B	414	ILE
1	B	415	ILE
1	B	422	LYS
1	B	437	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	97	HIS
1	A	101	HIS
1	A	234	ASN
1	A	240	HIS
1	A	292	ASN
1	A	334	GLN
1	A	362	ASN
1	A	396	HIS
1	A	403	ASN

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Mol	Chain	Res	Type
1	B	97	HIS
1	B	158	ASN
1	B	229	ASN
1	B	234	ASN
1	B	334	GLN
1	B	354	ASN
1	B	396	HIS
1	B	398	HIS
1	B	403	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	33/34 (97%)	6 (18%)	1 (3%)
2	D	33/34 (97%)	10 (30%)	1 (3%)
All	All	66/68 (97%)	16 (24%)	2 (3%)

All (16) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	C	8	G
2	C	12	G
2	C	17	A
2	C	18	G
2	C	19	U
2	C	20	C
2	D	9	A
2	D	10	U
2	D	17	A
2	D	18	G
2	D	19	U
2	D	22	A
2	D	25	U
2	D	26	A
2	D	29	U
2	D	34	C

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	C	19	U

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Mol	Chain	Res	Type
2	D	17	A

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 1 is 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$
4	SO4	A	601	-	4,4,4	0.24	0	6,6,6	0.26	0
4	SO4	B	603	-	4,4,4	0.27	0	6,6,6	0.23	0
6	EDO	B	801	-	3,3,3	0.42	0	2,2,2	0.37	0
3	ATP	A	501	5	32,33,33	1.75	7 (21%)	48,52,52	1.78	10 (20%)
3	ATP	B	502	-	32,33,33	1.81	7 (21%)	48,52,52	1.73	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
3	ATP	B	502	-	-	6/22/38/38	0/3/3/3
6	EDO	B	801	-	-	1/1/1/1	-
3	ATP	A	501	5	-	4/22/38/38	0/3/3/3

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	502	ATP	C5-C4	4.88	1.47	1.39
3	A	501	ATP	C5-C4	4.21	1.46	1.39
3	B	502	ATP	PA-O3A	4.13	1.64	1.59
3	B	502	ATP	PB-O3A	4.09	1.63	1.59
3	A	501	ATP	PA-O3A	3.93	1.63	1.59
3	A	501	ATP	PB-O3A	3.77	1.63	1.59
3	A	501	ATP	PB-O3B	3.76	1.63	1.59
3	B	502	ATP	PB-O3B	3.63	1.63	1.59
3	A	501	ATP	C5-N7	-2.98	1.33	1.39
3	B	502	ATP	C5-C6	2.91	1.49	1.41
3	B	502	ATP	C5-N7	-2.28	1.34	1.39
3	B	502	ATP	C8-N7	2.27	1.36	1.31
3	A	501	ATP	C8-N7	2.25	1.36	1.31
3	A	501	ATP	C5-C6	2.23	1.47	1.41

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	ATP	C5-C4-N3	-5.61	119.00	126.72
3	A	501	ATP	C5-C4-N3	-5.03	119.79	126.72
3	B	502	ATP	N3-C4-N9	4.51	134.84	127.17
3	A	501	ATP	N3-C2-N1	-4.51	121.76	128.58
3	A	501	ATP	N3-C4-N9	4.19	134.29	127.17
3	B	502	ATP	C2-N3-C4	3.88	121.31	111.83
3	B	502	ATP	N3-C2-N1	-3.73	122.93	128.58
3	A	501	ATP	C2-N3-C4	3.68	120.82	111.83
3	B	502	ATP	C4-C5-N7	-3.22	106.91	110.58
3	A	501	ATP	C4-C5-N7	-3.07	107.07	110.58
3	A	501	ATP	C2-N1-C6	2.95	123.57	118.73
3	A	501	ATP	C4-N9-C8	2.82	108.70	105.74
3	A	501	ATP	C5-N7-C8	2.79	107.83	103.45
3	A	501	ATP	O2'-C2'-C1'	-2.48	101.58	110.10
3	A	501	ATP	N9-C8-N7	-2.44	110.48	113.94
3	B	502	ATP	C5-N7-C8	2.39	107.20	103.45
3	B	502	ATP	C4-N9-C8	2.38	108.23	105.74
3	B	502	ATP	O3G-PG-O2G	2.28	116.35	107.80
3	B	502	ATP	C6-C5-N7	2.27	136.46	132.09

There are no chirality outliers.

All (11) torsion outliers are listed below:

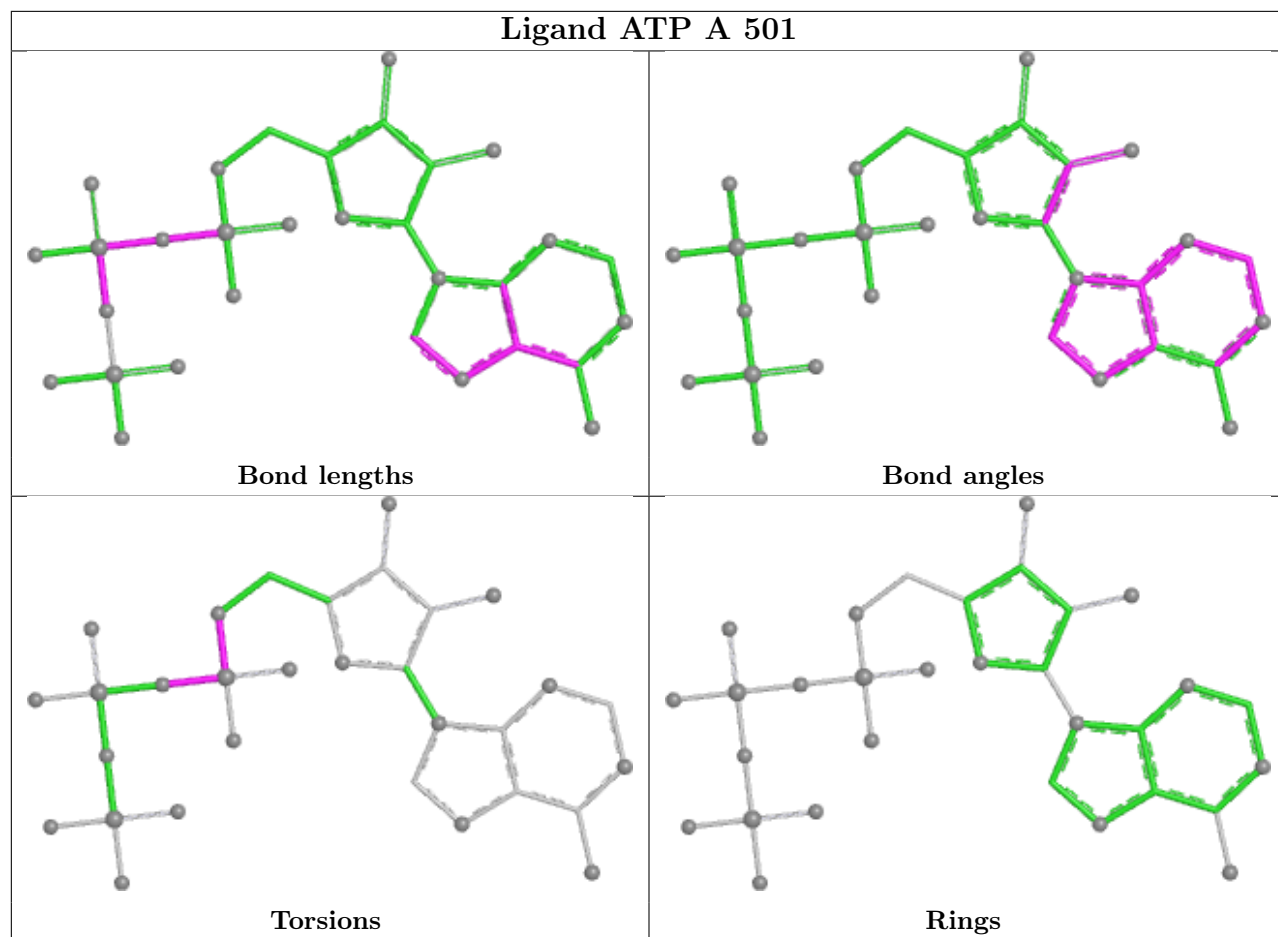
Mol	Chain	Res	Type	Atoms
3	A	501	ATP	C5'-O5'-PA-O1A
3	B	502	ATP	C5'-O5'-PA-O1A
3	B	502	ATP	C5'-O5'-PA-O3A
3	B	502	ATP	PB-O3B-PG-O1G
3	A	501	ATP	PB-O3A-PA-O5'
3	B	502	ATP	PB-O3A-PA-O5'
3	A	501	ATP	C5'-O5'-PA-O2A
3	A	501	ATP	C5'-O5'-PA-O3A
3	B	502	ATP	C5'-O5'-PA-O2A
6	B	801	EDO	O1-C1-C2-O2
3	B	502	ATP	PB-O3A-PA-O2A

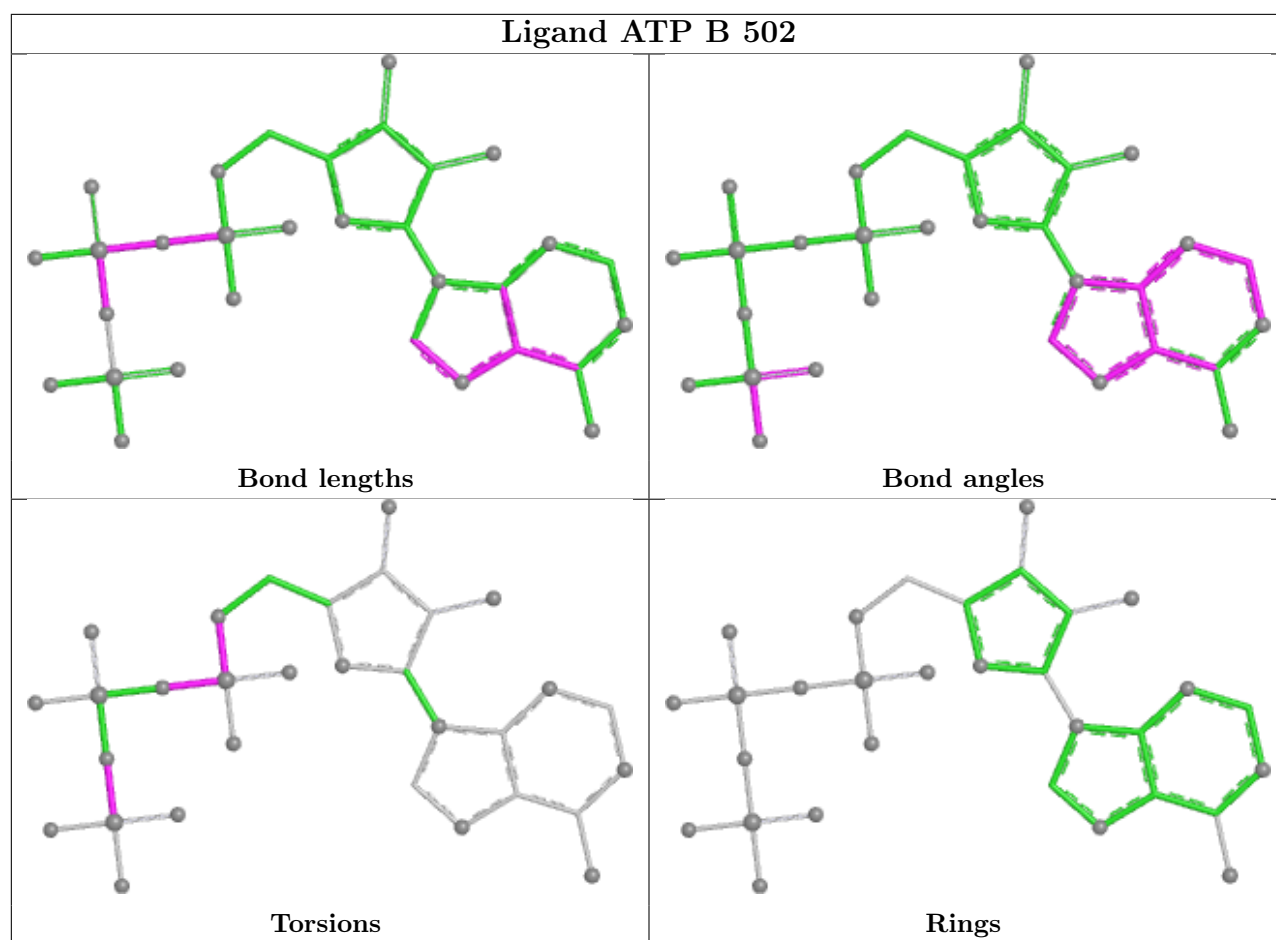
There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	601	SO4	1	0
6	B	801	EDO	3	0
3	B	502	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	441/441 (100%)	-0.41	1 (0%) 91 83	10, 33, 89, 180	1 (0%)
1	B	437/441 (99%)	0.19	11 (2%) 58 35	10, 71, 163, 241	0
2	C	34/34 (100%)	0.10	0 100 100	25, 95, 145, 158	1 (2%)
2	D	34/34 (100%)	0.22	2 (5%) 28 14	41, 102, 189, 205	1 (2%)
All	All	946/950 (99%)	-0.09	14 (1%) 72 49	10, 49, 154, 241	3 (0%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	260	LEU	3.2
1	B	205	LYS	3.1
1	B	208	VAL	2.9
1	B	204	ALA	2.7
1	B	261	GLU	2.6
1	B	206	GLY	2.6
1	B	210	LYS	2.3
2	D	17	A	2.3
1	B	258	HIS	2.3
2	D	18	G	2.2
1	A	440	CYS	2.2
1	B	207	GLU	2.2
1	B	2	LYS	2.1
1	B	3	VAL	2.1

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](#)

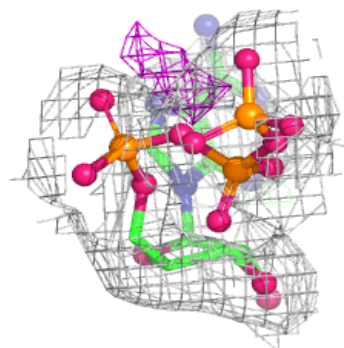
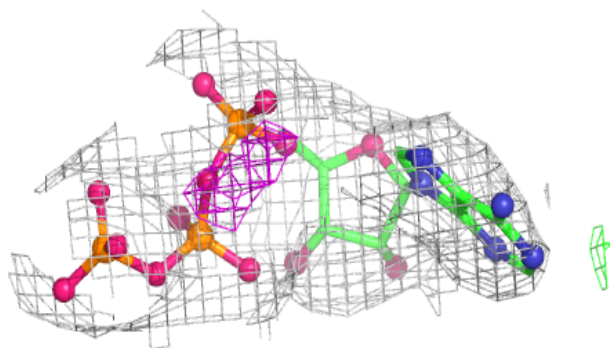
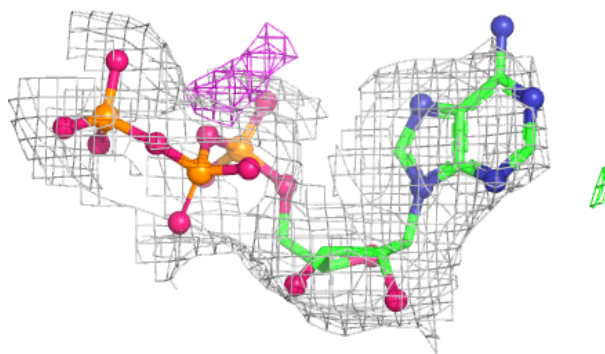
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	SO4	A	601	5/5	0.84	0.17	86,86,86,87	0
3	ATP	B	502	31/31	0.89	0.10	70,74,75,75	0
5	MN	A	701	1/1	0.89	0.24	49,49,49,49	0
6	EDO	B	801	4/4	0.89	0.09	43,43,44,44	0
4	SO4	B	603	5/5	0.93	0.10	69,70,70,70	0
3	ATP	A	501	31/31	0.97	0.06	22,26,28,28	0

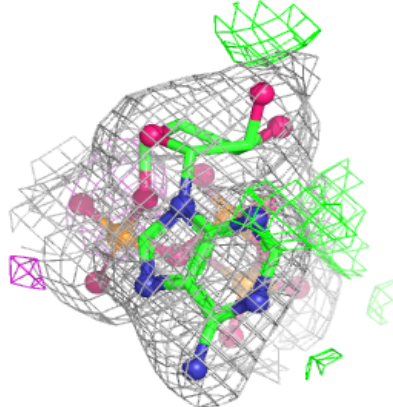
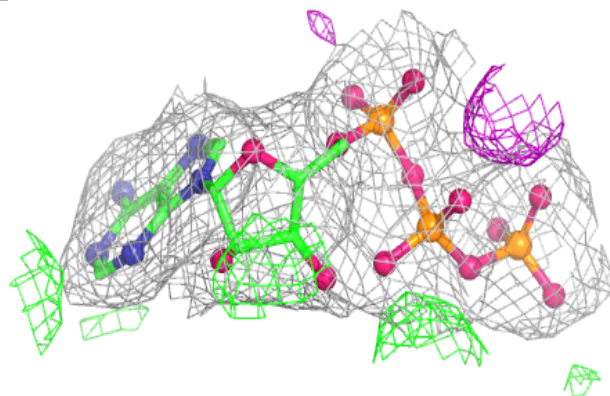
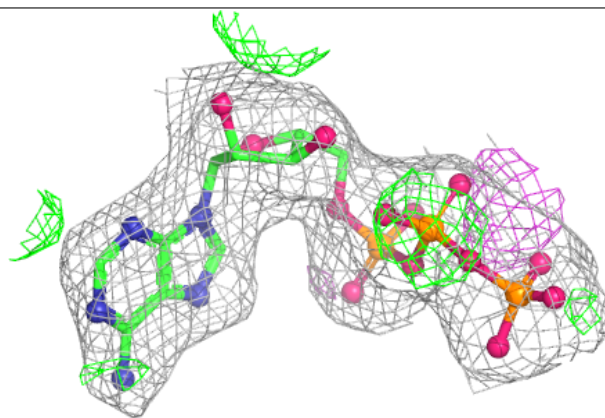
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ATP B 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 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.