



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:49 PM UTC

PDB ID : 7Z6R / pdb_00007z6r
Title : Psychrobacter arcticus ATPPRT (HisGZ) R56A mutant bound to ATP and PRPP
Authors : Alphey, M.S.; Fisher, G.; da Silva, R.G.
Deposited on : 2022-03-14
Resolution : 2.55 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

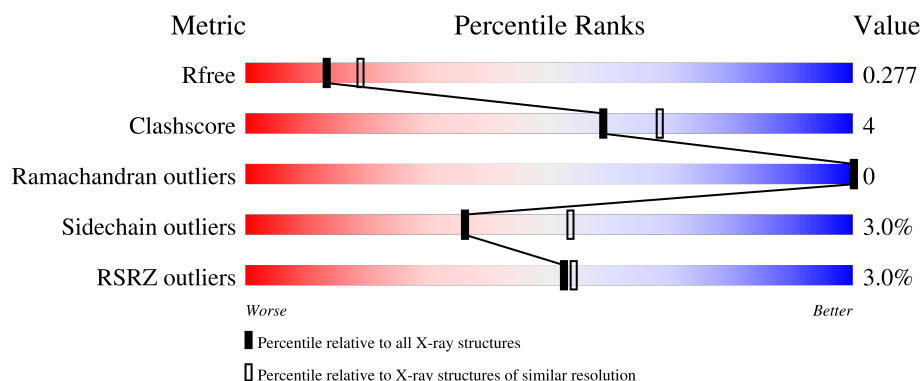
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	388	4% 85% 9% • 5%
1	B	388	% 87% 9% • •
2	C	232	5% 78% 9% • 11%
2	D	232	% 81% 7% • 11%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8522 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 ATP phosphoribosyltransferase regulatory subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	368	Total	C	N	O	S	0	0	0
			2633	1652	473	498	10			
1	B	377	Total	C	N	O	S	0	1	0
			2809	1786	499	513	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q4FTX3
B	0	GLY	-	expression tag	UNP Q4FTX3

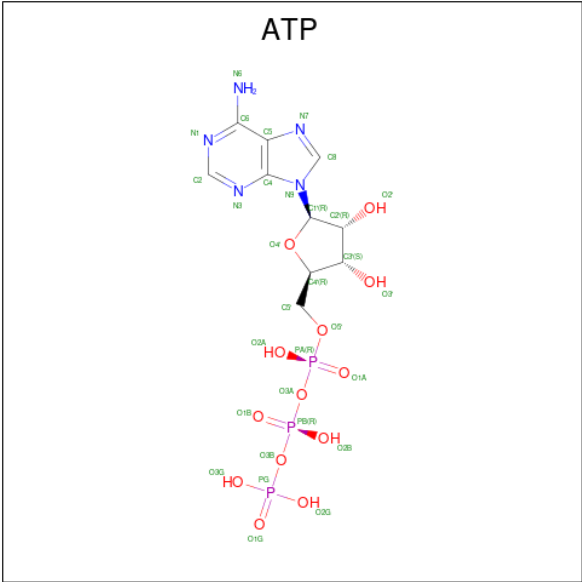
- Molecule 2 is a protein called ATP phosphoribosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	206	Total	C	N	O	S	0	0	0
			1478	938	260	276	4			
2	D	206	Total	C	N	O	S	0	0	0
			1480	946	253	277	4			

There are 4 discrepancies between the modelled and reference sequences:

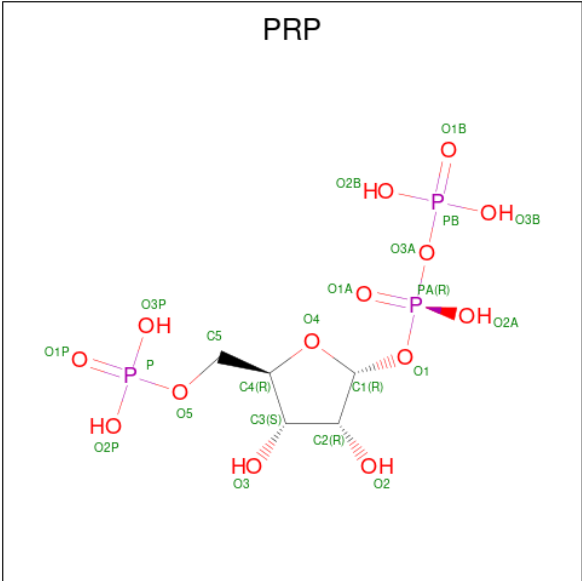
Chain	Residue	Modelled	Actual	Comment	Reference
C	0	GLY	-	expression tag	UNP Q4FQF7
C	56	ALA	ARG	engineered mutation	UNP Q4FQF7
D	0	GLY	-	expression tag	UNP Q4FQF7
D	56	ALA	ARG	engineered mutation	UNP Q4FQF7

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is 1-O-pyrophosphono-5-O-phosphono-alpha-D-ribofuranose (CCD ID: PRP) (formula: C₅H₁₃O₁₄P₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	O	P	0	0
			22	5	14	3		

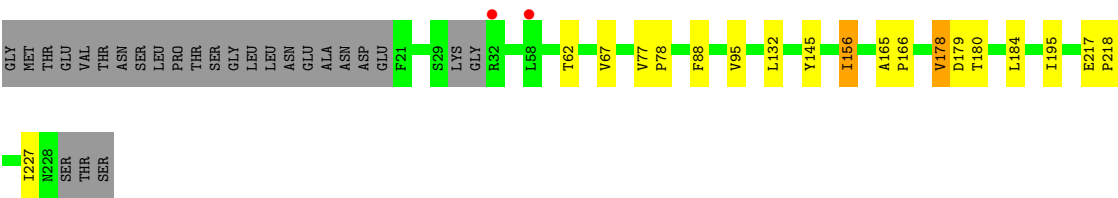
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	D	1	Total	C	O	P	0	0
			22	5	14	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	8	Total	O	0	0
			8	8		
5	B	4	Total	O	0	0
			4	4		
5	C	3	Total	O	0	0
			3	3		
5	D	1	Total	O	0	0
			1	1		



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	102.16Å 145.53Å 93.81Å 90.00° 102.48° 90.00°	Depositor
Resolution (Å)	56.97 – 2.55 56.97 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.8 (56.97-2.55) 99.8 (56.97-2.55)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.19 (at 2.55Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.233 , 0.275 0.236 , 0.277	Depositor DCC
R_{free} test set	2159 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	64.7	Xtriage
Anisotropy	0.305	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 39.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8522	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.06% 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: ATP, PRP

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	1.03	0/2682	1.51	0/3631
1	B	1.04	0/2867	1.49	0/3894
2	C	1.07	0/1500	1.46	0/2040
2	D	1.06	0/1503	1.45	0/2047
All	All	1.05	0/8552	1.48	0/11612

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	2633	0	2406	22	0
1	B	2809	0	2726	26	0
2	C	1478	0	1441	19	0
2	D	1480	0	1446	10	0
3	C	31	0	12	1	0
3	D	31	0	12	0	0
4	C	22	0	8	1	0
4	D	22	0	8	2	0
5	A	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	4	0	0	0	0
5	C	3	0	0	0	0
5	D	1	0	0	0	0
All	All	8522	0	8059	73	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 4.

All (73) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:117:LEU:CD1	2:C:177:VAL:HG12	2.26	0.66
1:B:275:ILE:HD11	1:B:282:LEU:HD22	1.78	0.65
1:B:82:LEU:HD21	1:B:290:GLY:HA2	1.80	0.64
1:A:40:MET:HE2	1:B:40:MET:HE2	1.84	0.60
2:C:78:PRO:HG2	2:C:95:VAL:HG12	1.83	0.60
1:A:63:ILE:HD12	1:A:63:ILE:N	2.19	0.58
2:C:62:THR:HG22	2:C:67:VAL:O	2.03	0.58
2:C:116:LYS:HE2	2:C:118:MET:HE2	1.86	0.57
2:C:46:VAL:HG13	2:C:62:THR:HG21	1.85	0.57
2:D:78:PRO:HG2	2:D:95:VAL:HG12	1.87	0.56
2:D:62:THR:HG22	2:D:67:VAL:O	2.06	0.55
1:A:62:ILE:HD13	1:A:73:ILE:HD11	1.88	0.54
2:C:117:LEU:HD11	2:C:177:VAL:HG12	1.90	0.53
1:B:155:ASP:OD1	1:B:254:SER:OG	2.26	0.52
2:C:165:ALA:HB3	2:C:166:PRO:HD3	1.92	0.52
2:D:165:ALA:HB3	2:D:166:PRO:HD3	1.92	0.52
1:A:102:VAL:HG23	1:A:116:PRO:HG2	1.91	0.52
1:A:30:THR:HB	1:B:352:ARG:HG2	1.93	0.51
1:B:153:HIS:CE1	1:B:276:ASN:O	2.65	0.50
2:C:107:LEU:HD12	2:C:204:VAL:HG21	1.94	0.50
2:C:130:ARG:NH2	2:C:172:ASP:OD1	2.43	0.50
1:B:275:ILE:CD1	1:B:282:LEU:HD22	2.43	0.49
1:B:21:ARG:HD3	1:B:117:LEU:HD21	1.93	0.49
1:A:287:ARG:NH1	1:A:303:GLN:HG2	2.28	0.48
2:C:179:ASP:N	4:C:402:PRP:O1P	2.46	0.48
1:B:136:ILE:HD12	1:B:245:LEU:HD21	1.95	0.48
1:A:42:GLU:HG2	1:B:2:LEU:CD2	2.43	0.48
1:B:60:PHE:CZ	1:B:105:THR:HA	2.48	0.48
2:C:177:VAL:HG21	3:C:401:ATP:C8	2.49	0.48
1:A:224:ASN:OD1	1:A:224:ASN:N	2.46	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:217:GLU:HB3	2:D:218:PRO:HD3	1.96	0.47
1:A:60:PHE:CZ	1:A:105:THR:HA	2.48	0.47
1:B:102:VAL:HG23	1:B:116:PRO:HG2	1.95	0.47
1:A:290:GLY:C	1:A:302:ARG:HB3	2.39	0.47
1:A:102:VAL:CG2	1:A:116:PRO:HG2	2.43	0.47
1:B:28:LEU:HD13	1:B:97:CYS:HB2	1.96	0.47
1:A:156:LEU:HB2	1:A:255:ILE:HD13	1.97	0.47
2:C:48:LEU:CD2	2:C:60:PHE:HB3	2.45	0.47
2:C:46:VAL:HG12	2:C:46:VAL:O	2.15	0.46
1:A:129:ILE:HD12	1:A:237:GLU:HG2	1.97	0.46
1:B:124:PHE:HB3	1:B:302:ARG:HE	1.80	0.46
1:B:275:ILE:CG1	1:B:282:LEU:HD22	2.46	0.46
2:D:178:VAL:HG11	2:D:184:LEU:HD11	1.98	0.46
1:B:171:LEU:HD11	1:B:204:PHE:CZ	2.51	0.46
2:D:179:ASP:HB3	4:D:501:PRP:O2P	2.16	0.45
2:C:118:MET:HE3	2:C:178:VAL:HG21	1.99	0.45
1:B:18:GLU:OE1	1:B:21:ARG:NH1	2.50	0.45
1:B:102:VAL:CG2	1:B:116:PRO:HG2	2.47	0.44
1:B:241:LEU:HD23	1:B:241:LEU:HA	1.83	0.44
1:B:333:PHE:HD1	1:B:344:LEU:HD23	1.83	0.44
1:A:165:LEU:CB	1:A:204:PHE:HE1	2.31	0.43
1:A:19:VAL:O	1:A:23:GLN:HG2	2.18	0.43
1:A:333:PHE:HD1	1:A:344:LEU:HD23	1.82	0.43
1:A:242:LYS:O	1:A:246:GLN:HG3	2.18	0.43
2:C:117:LEU:HD13	2:C:198:VAL:HG21	2.01	0.43
1:A:129:ILE:O	1:A:133:ILE:HD12	2.18	0.43
1:A:155:ASP:OD1	1:A:254:SER:OG	2.34	0.43
1:A:42:GLU:HB2	1:A:47:LEU:HB2	2.00	0.42
1:B:19:VAL:O	1:B:23:GLN:HG2	2.20	0.42
1:B:214:ILE:HB	1:B:255:ILE:HD12	2.02	0.42
1:B:387:VAL:HG23	1:B:387:VAL:O	2.19	0.42
2:C:48:LEU:HD21	2:C:60:PHE:HB3	2.00	0.42
2:D:180:THR:N	4:D:501:PRP:O2P	2.41	0.42
1:B:58:GLN:OE1	1:B:104:HIS:HB3	2.19	0.41
1:B:129:ILE:HD12	1:B:237:GLU:HA	2.03	0.41
2:D:145:TYR:CG	2:D:195:ILE:HG12	2.55	0.41
2:C:83:HIS:CE1	2:D:156:ILE:CD1	3.03	0.41
2:C:145:TYR:CG	2:C:195:ILE:HG12	2.56	0.41
2:C:117:LEU:HD23	2:C:141:VAL:HG11	2.03	0.40
2:D:77:VAL:HB	2:D:78:PRO:HD3	2.03	0.40
1:A:326:VAL:HA	1:A:373:HIS:O	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:GLN:HA	1:A:23:GLN:OE1	2.21	0.40
1:B:341:ARG:O	1:B:345:LEU:HG	2.21	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	358/388 (92%)	346 (97%)	12 (3%)	0	100	100
1	B	374/388 (96%)	363 (97%)	11 (3%)	0	100	100
2	C	202/232 (87%)	189 (94%)	13 (6%)	0	100	100
2	D	202/232 (87%)	190 (94%)	12 (6%)	0	100	100
All	All	1136/1240 (92%)	1088 (96%)	48 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	250/329 (76%)	242 (97%)	8 (3%)	34	51
1	B	283/329 (86%)	277 (98%)	6 (2%)	47	66
2	C	144/194 (74%)	138 (96%)	6 (4%)	26	40

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	146/194 (75%)	141 (97%)	5 (3%)	32	48
All	All	823/1046 (79%)	798 (97%)	25 (3%)	36	53

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	SER
1	A	113	SER
1	A	143	ILE
1	A	164	ARG
1	A	196	GLN
1	A	224	ASN
1	A	255	ILE
1	A	302	ARG
1	B	113	SER
1	B	129	ILE
1	B	153	HIS
1	B	241	LEU
1	B	255	ILE
1	B	275	ILE
2	C	48	LEU
2	C	88	PHE
2	C	117	LEU
2	C	118	MET
2	C	178	VAL
2	C	222	SER
2	D	88	PHE
2	D	132	LEU
2	D	156	ILE
2	D	178	VAL
2	D	227	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	HIS
1	A	26	GLN
1	A	196	GLN
1	A	212	HIS
1	A	334	ASN
1	B	15	HIS

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Mol	Chain	Res	Type
1	B	26	GLN
1	B	87	HIS
1	B	88	HIS
1	B	212	HIS
1	B	334	ASN
1	B	385	HIS
2	C	83	HIS
2	D	103	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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](#)

4 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	$\# Z > 2$
3	ATP	C	401	-	32,33,33	0.61	0	48,52,52	0.60	0
4	PRP	C	402	-	20,22,22	0.84	1 (5%)	32,35,35	0.99	2 (6%)
3	ATP	D	502	-	32,33,33	0.55	0	48,52,52	0.57	0
4	PRP	D	501	-	20,22,22	0.80	1 (5%)	32,35,35	1.09	2 (6%)

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	C	401	-	-	6/22/38/38	0/3/3/3
4	PRP	C	402	-	-	1/16/33/33	0/1/1/1
3	ATP	D	502	-	-	1/22/38/38	0/3/3/3
4	PRP	D	501	-	-	8/16/33/33	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	402	PRP	PA-O3A	2.67	1.62	1.59
4	D	501	PRP	PA-O3A	2.18	1.61	1.59

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	501	PRP	O1-C1-C2	2.78	111.17	106.65
4	C	402	PRP	O1-C1-C2	2.20	110.22	106.65
4	C	402	PRP	C1-C2-C3	2.05	104.84	102.29
4	D	501	PRP	C1-C2-C3	2.03	104.81	102.29

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	501	PRP	C1-O1-PA-O3A
4	D	501	PRP	C5-O5-P-O1P
4	D	501	PRP	C5-O5-P-O2P
4	D	501	PRP	C5-O5-P-O3P
3	C	401	ATP	C3'-C4'-C5'-O5'
3	C	401	ATP	O4'-C4'-C5'-O5'
4	D	501	PRP	O4-C4-C5-O5
4	C	402	PRP	PA-O3A-PB-O3B
4	D	501	PRP	PB-O3A-PA-O2A
3	C	401	ATP	PB-O3B-PG-O1G
4	D	501	PRP	PB-O3A-PA-O1A
3	C	401	ATP	PB-O3B-PG-O2G
3	C	401	ATP	PB-O3B-PG-O3G
3	C	401	ATP	PG-O3B-PB-O2B

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Mol	Chain	Res	Type	Atoms
3	D	502	ATP	PB-O3B-PG-O1G
4	D	501	PRP	C2-C1-O1-PA

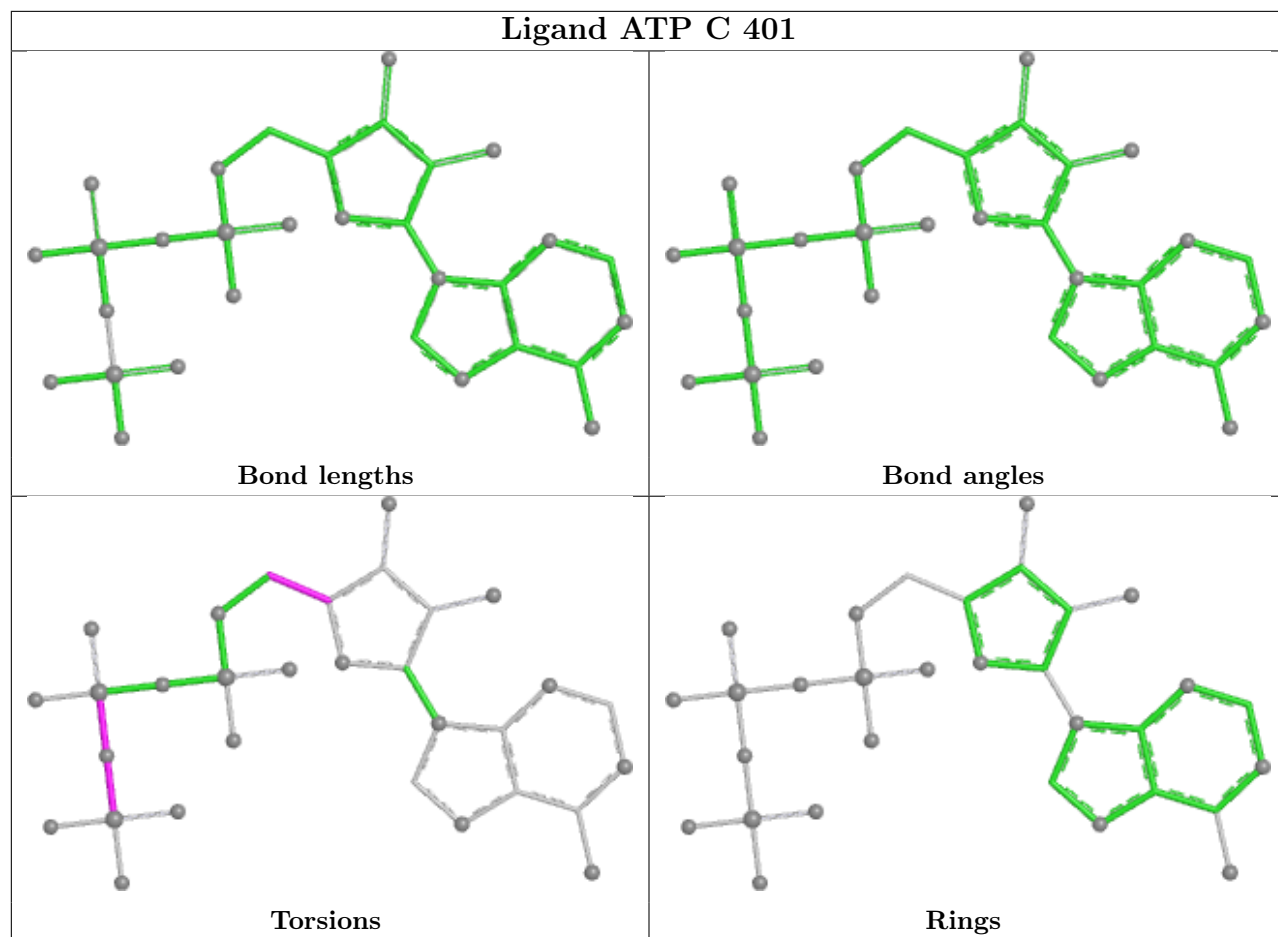
There are no ring outliers.

3 monomers are involved in 4 short contacts:

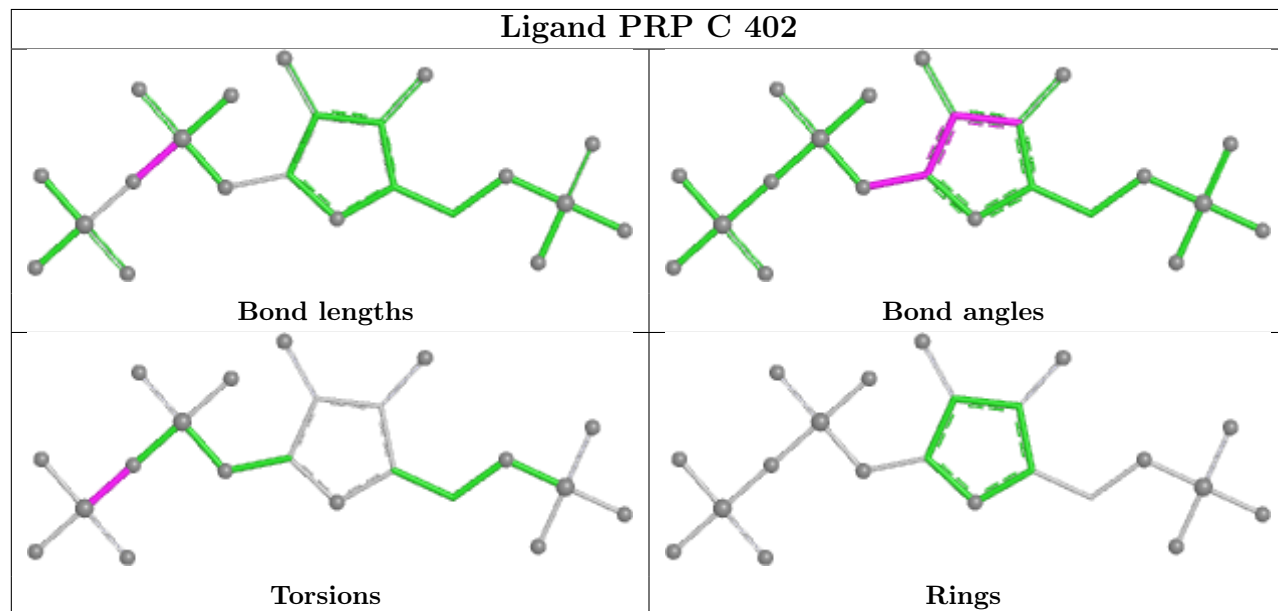
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	401	ATP	1	0
4	C	402	PRP	1	0
4	D	501	PRP	2	0

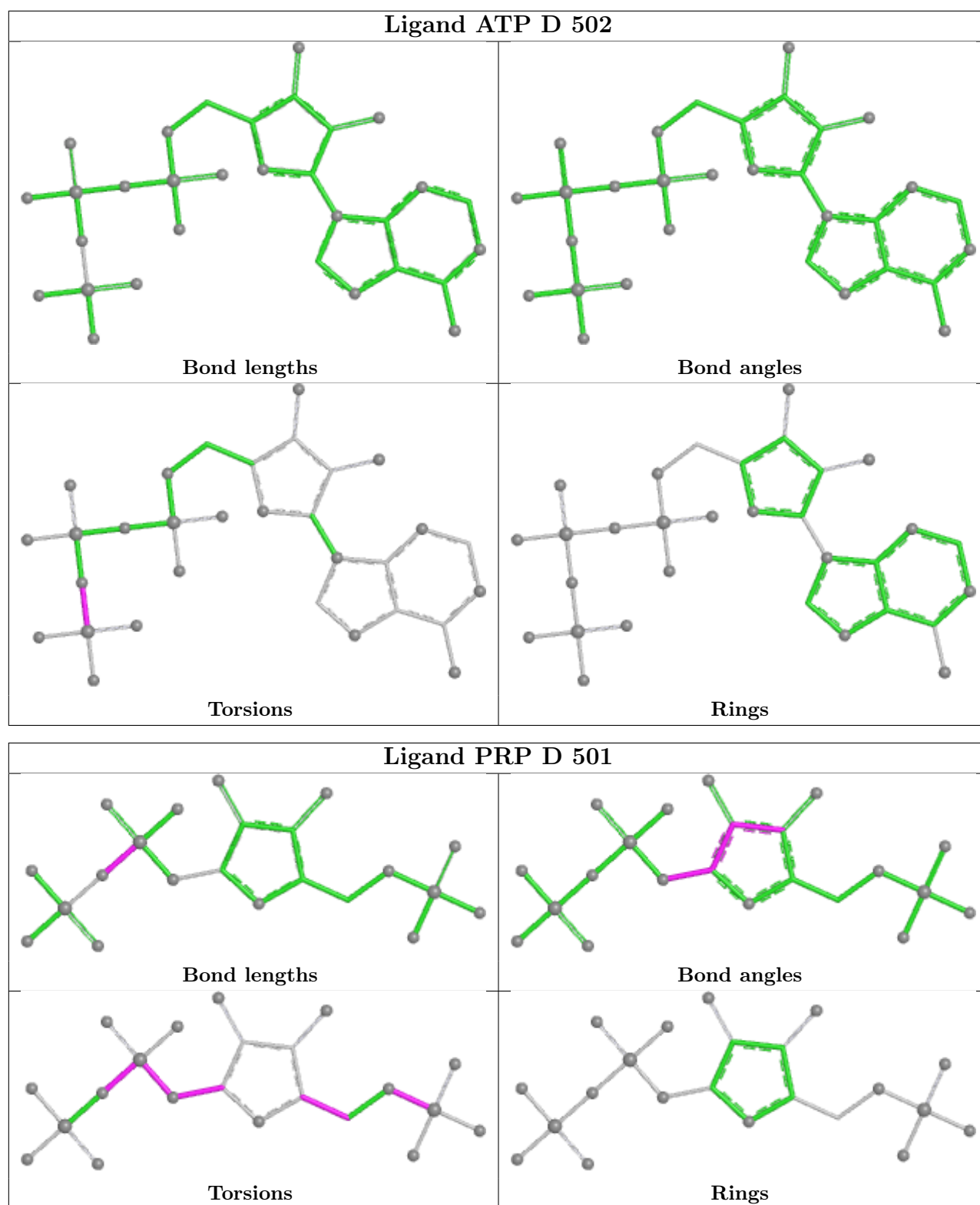
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

Ligand ATP C 401



Ligand PRP C 402





5.7 Other polymers ⓘ

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 ⓘ

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	368/388 (94%)	0.55	17 (4%) 37 38	52, 76, 123, 142	0
1	B	377/388 (97%)	0.15	4 (1%) 78 79	40, 64, 82, 98	1 (0%)
2	C	206/232 (88%)	0.58	12 (5%) 29 28	70, 87, 107, 123	0
2	D	206/232 (88%)	0.53	2 (0%) 79 80	65, 86, 113, 142	0
All	All	1157/1240 (93%)	0.42	35 (3%) 52 54	40, 76, 113, 142	1 (0%)

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	175	ASP	4.2
1	A	169	ALA	3.7
1	A	176	THR	3.7
1	A	300	GLN	3.4
1	B	290	GLY	3.1
2	C	21	PHE	3.0
1	A	165	LEU	3.0
1	A	230	LYS	3.0
2	D	32	ARG	2.9
2	C	86	ALA	2.8
1	A	217	LEU	2.8
1	A	195	CYS	2.8
2	C	58	LEU	2.6
1	B	301	PRO	2.6
2	C	69	VAL	2.4
2	C	56	ALA	2.3
1	A	222	SER	2.2
1	A	241	LEU	2.2
1	B	249	TRP	2.2
1	A	204	PHE	2.2
1	B	1	MET	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	213	ASP	2.2
1	A	178	GLN	2.1
2	C	28	LEU	2.1
2	C	81	VAL	2.1
1	A	199	PRO	2.1
1	A	225	ALA	2.1
2	C	49	LEU	2.1
2	C	46	VAL	2.0
2	C	67	VAL	2.0
1	A	191	LEU	2.0
2	C	23	GLY	2.0
1	A	231	ILE	2.0
2	C	50	GLU	2.0
2	D	58	LEU	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](#)

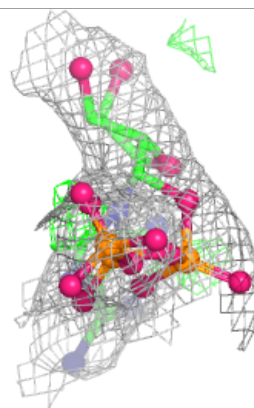
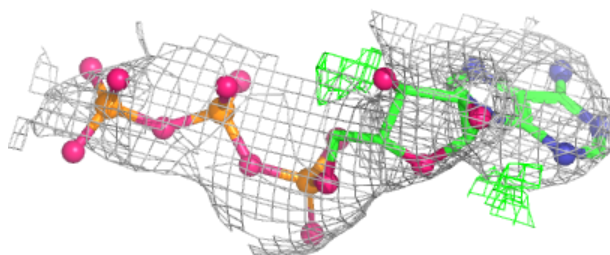
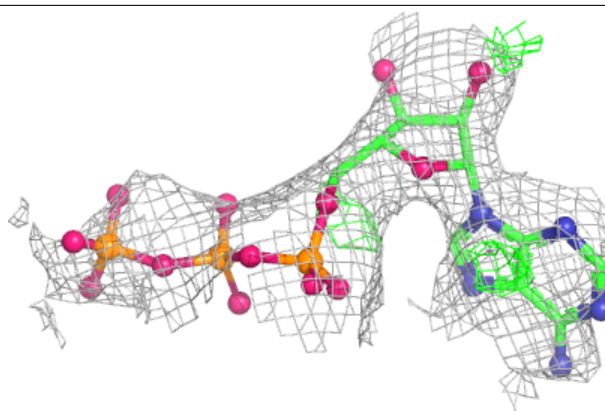
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	ATP	D	502	31/31	0.79	0.15	93,95,121,122	31
4	PRP	D	501	22/22	0.79	0.09	93,113,139,140	0
4	PRP	C	402	22/22	0.83	0.09	94,106,111,115	0
3	ATP	C	401	31/31	0.83	0.11	89,92,108,111	31

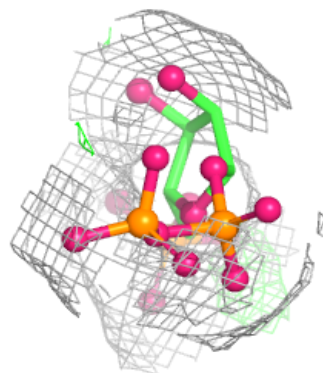
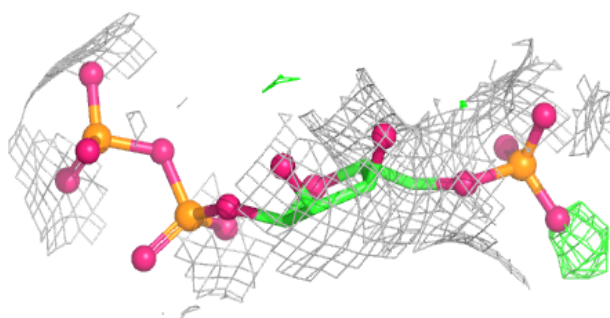
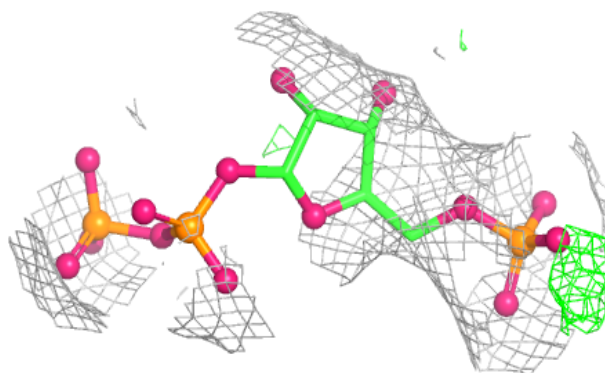
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ATP D 502:

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

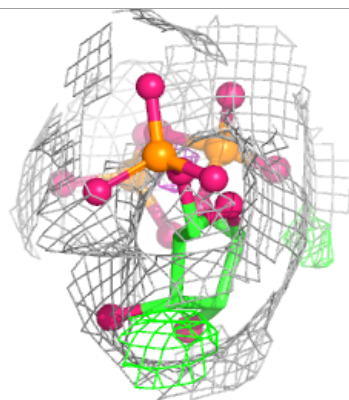
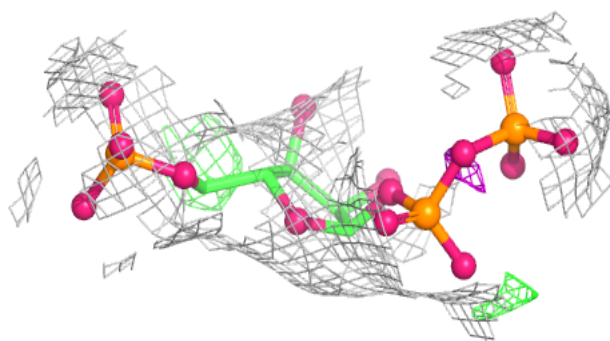
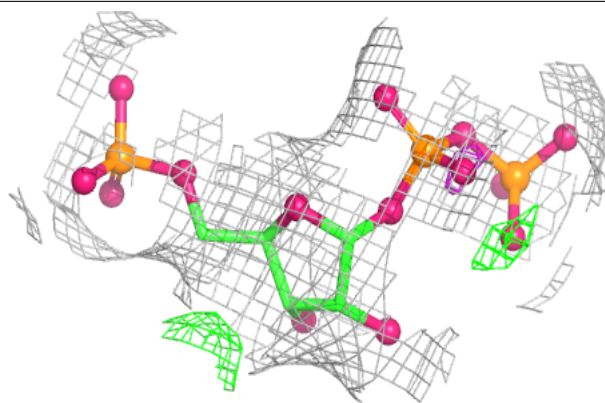
**Electron density around PRP D 501:**

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

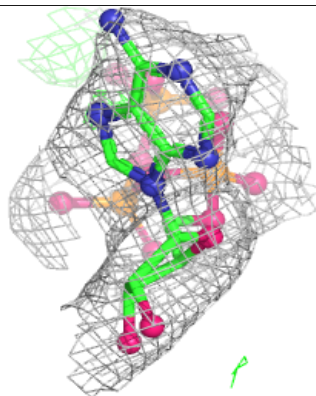
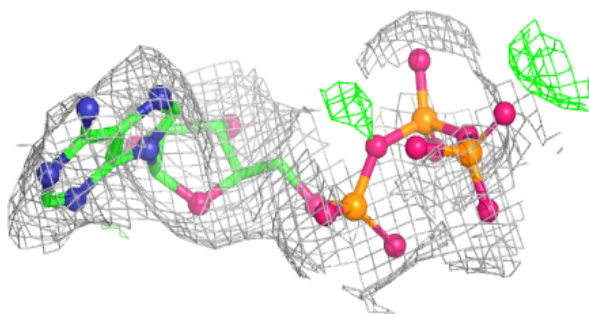
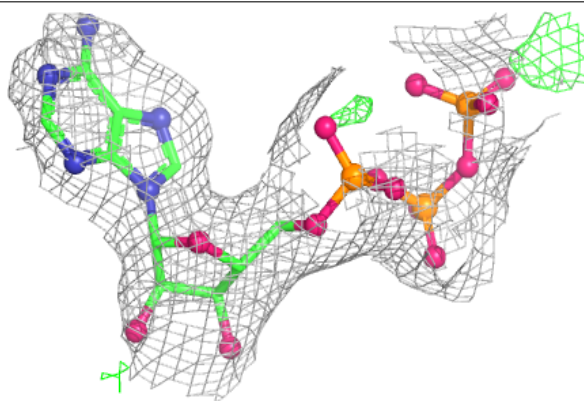


Electron density around PRP C 402:

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

**Electron density around ATP C 401:**

$2mF_o-DF_c$ (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.