



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 28, 2024 – 02:40 pm BST

PDB ID : 1W2D
Title : Human Inositol (1,4,5)-trisphosphate 3-kinase complexed with Mn²⁺/ADP/Ins(1,3,4,5)P₄
Authors : Gonzalez, B.; Schell, M.J.; Irvine, R.F.; Williams, R.L.
Deposited on : 2004-07-01
Resolution : 1.94 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	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

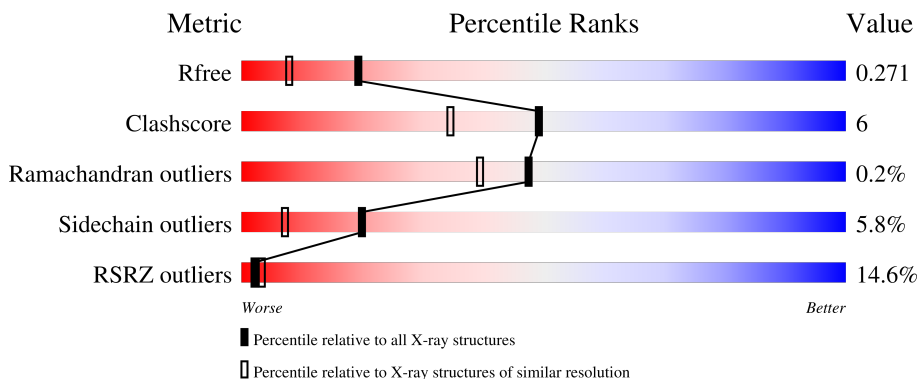
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.94 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}	130704	4310 (1.96-1.92)
Clashscore	141614	1023 (1.94-1.94)
Ramachandran outliers	138981	1007 (1.94-1.94)
Sidechain outliers	138945	1007 (1.94-1.94)
RSRZ outliers	127900	4250 (1.96-1.92)

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	265	<div> <div>9%</div> <div>82%</div> <div>15%</div> <div>.</div> </div>
1	B	265	<div> <div>19%</div> <div>78%</div> <div>14%</div> <div>6%</div> </div>

2 Entry composition [i](#)

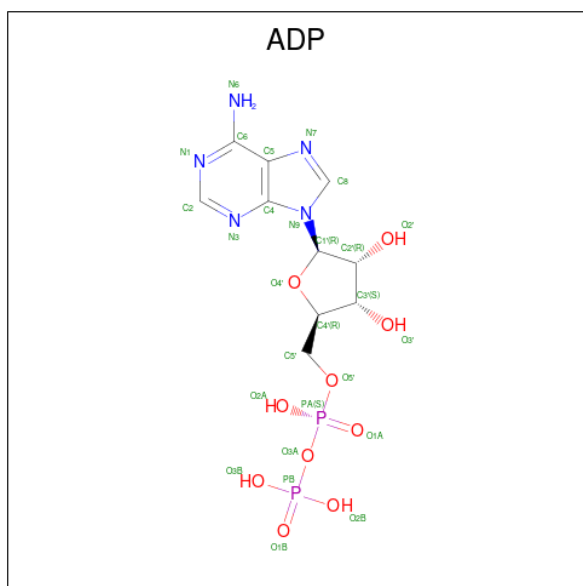
There are 6 unique types of molecules in this entry. The entry contains 4524 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 INOSITOL-TRISPHOSPHATE 3-KINASE A.

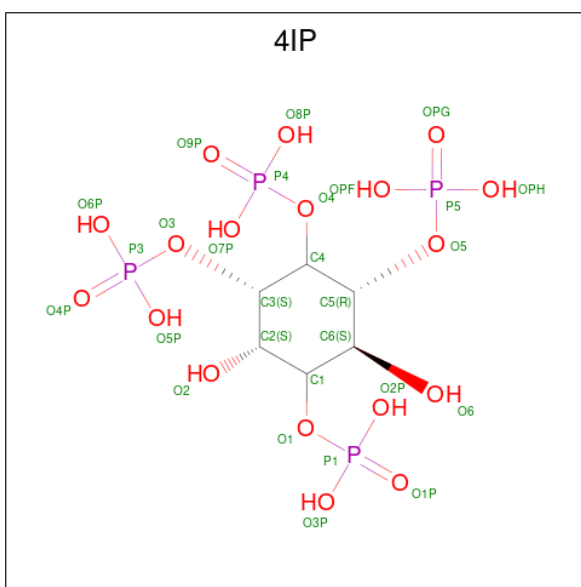
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	0	0
			2133	1334	388	399	12			
1	B	250	Total	C	N	O	S	0	0	0
			2005	1260	361	372	12			

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 3 is INOSITOL-(1,3,4,5)-TETRAKISPHOSPHATE (three-letter code: 4IP) (formula: $C_6H_{16}O_{18}P_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			28	6	18	4		
3	B	1	Total	C	O	P	0	0
			28	6	18	4		

- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mn	0	0
			1	1		
4	B	1	Total	Mn	0	0
			1	1		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

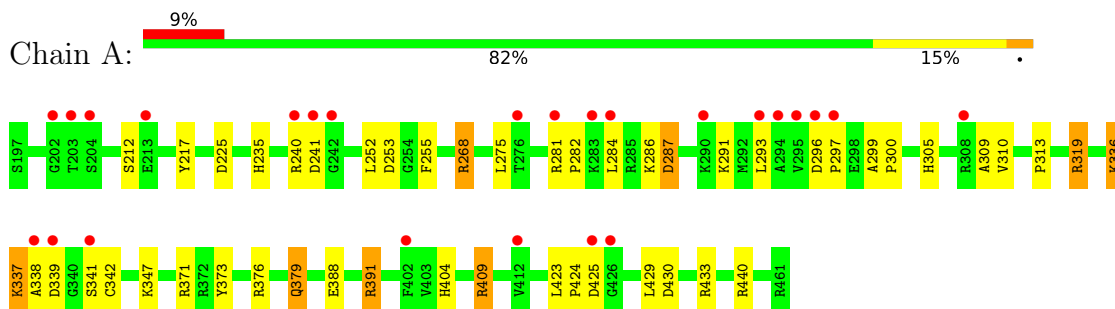
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	129	Total	O	0	0
			129	129		
6	B	130	Total	O	0	0
			130	130		

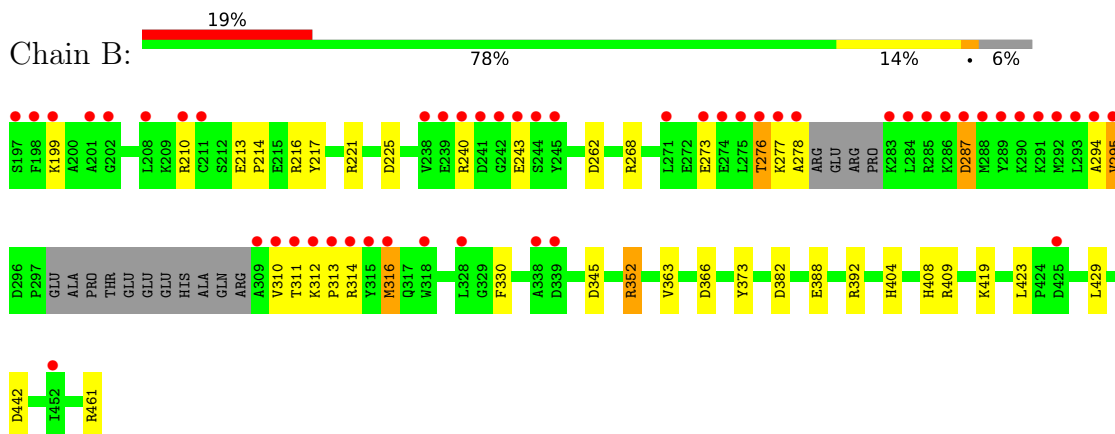
3 Residue-property plots

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: INOSITOL-TRISPHOSPHATE 3-KINASE A



• Molecule 1: INOSITOL-TRISPHOSPHATE 3-KINASE A



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	72.05Å 97.41Å 190.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.80 – 1.94 47.71 – 1.94	Depositor EDS
% Data completeness (in resolution range)	96.9 (48.80-1.94) 95.0 (47.71-1.94)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.05 (at 1.94Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.226 , 0.267 0.229 , 0.271	Depositor DCC
R_{free} test set	2409 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	27.5	Xtriage
Anisotropy	0.307	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 49.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4524	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% 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: SO4, 4IP, ADP, MN

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	0.92	0/2173	1.03	11/2924 (0.4%)
1	B	0.91	2/2040 (0.1%)	1.00	7/2741 (0.3%)
All	All	0.92	2/4213 (0.0%)	1.01	18/5665 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	363	VAL	CB-CG1	-6.73	1.38	1.52
1	B	330	PHE	CD2-CE2	-5.77	1.27	1.39

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	287	ASP	CB-CG-OD2	6.62	124.26	118.30
1	A	253	ASP	CB-CG-OD2	6.49	124.14	118.30
1	A	268	ARG	NE-CZ-NH2	-6.25	117.18	120.30
1	B	287	ASP	CB-CG-OD2	6.18	123.86	118.30
1	A	241	ASP	CB-CG-OD2	5.91	123.62	118.30
1	B	262	ASP	CB-CG-OD2	5.81	123.53	118.30
1	B	382	ASP	CB-CG-OD2	5.78	123.50	118.30
1	B	352	ARG	NE-CZ-NH1	5.62	123.11	120.30
1	B	225	ASP	CB-CG-OD2	5.41	123.16	118.30
1	A	319	ARG	NE-CZ-NH1	5.31	122.95	120.30
1	A	268	ARG	NE-CZ-NH1	5.30	122.95	120.30
1	A	371	ARG	NE-CZ-NH1	5.19	122.90	120.30
1	B	442	ASP	CB-CG-OD2	5.16	122.94	118.30
1	A	409	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	B	268	ARG	NE-CZ-NH2	-5.09	117.75	120.30
1	A	425	ASP	CB-CG-OD2	5.07	122.86	118.30
1	A	430	ASP	CB-CG-OD2	5.07	122.86	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	225	ASP	CB-CG-OD2	5.04	122.83	118.30

There are no chirality outliers.

There are no planarity outliers.

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	2133	0	2113	33	0
1	B	2005	0	1996	29	0
2	A	27	0	12	0	0
2	B	27	0	12	0	0
3	A	28	0	8	0	0
3	B	28	0	8	3	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	10	0	0	0	0
5	B	5	0	0	0	0
6	A	129	0	0	4	0
6	B	130	0	0	3	0
All	All	4524	0	4149	50	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (50) 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:338:ALA:C	1:B:294:ALA:HB1	1.98	0.82
1:A:338:ALA:HB1	1:B:294:ALA:HB1	1.71	0.72
1:A:339:ASP:HB3	1:B:295:VAL:HA	1.71	0.72
1:A:338:ALA:O	1:B:294:ALA:CB	2.38	0.72
1:A:338:ALA:CA	1:B:294:ALA:HB1	2.20	0.72
1:A:338:ALA:CB	1:B:294:ALA:HB1	2.25	0.66
1:A:376:ARG:HH11	1:A:379:GLN:HE22	1.44	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:297:PRO:HB2	6:A:2036:HOH:O	1.99	0.63
1:A:338:ALA:O	1:B:294:ALA:HB3	1.99	0.62
1:A:338:ALA:O	1:B:294:ALA:HB1	1.99	0.62
1:B:221:ARG:NH2	6:B:2009:HOH:O	2.35	0.58
1:B:419:LYS:NZ	3:B:1463:4IP:O9P	2.32	0.58
1:A:341:SER:HB2	1:B:345:ASP:HA	1.85	0.57
1:A:339:ASP:HB2	1:B:295:VAL:HG22	1.86	0.57
1:A:293:LEU:HD13	1:A:305:HIS:HE1	1.69	0.56
1:A:284:LEU:HB3	1:A:309:ALA:HB1	1.88	0.55
1:B:316:MET:HA	1:B:316:MET:HE3	1.88	0.55
1:A:338:ALA:C	1:B:294:ALA:CB	2.71	0.55
1:B:423:LEU:HD11	1:B:429:LEU:HG	1.88	0.54
1:B:373:TYR:OH	1:B:404:HIS:HD2	1.91	0.53
1:B:388:GLU:O	1:B:392:ARG:HG2	2.08	0.53
1:A:299:ALA:HB3	1:A:300:PRO:HD3	1.91	0.52
1:A:338:ALA:HB1	1:B:294:ALA:CB	2.39	0.52
1:A:293:LEU:CD1	1:A:305:HIS:HE1	2.22	0.52
1:B:366:ASP:OD2	1:B:408:HIS:HD2	1.92	0.51
1:A:255:PHE:CE1	1:A:409:ARG:HG2	2.45	0.51
1:B:419:LYS:CE	3:B:1463:4IP:O9P	2.59	0.51
1:A:337:LYS:HB2	1:A:341:SER:O	2.13	0.48
1:B:409:ARG:NE	6:B:2088:HOH:O	2.45	0.48
1:A:287:ASP:O	1:A:291:LYS:HD2	2.13	0.48
1:A:376:ARG:NH1	1:A:379:GLN:HE22	2.12	0.47
1:A:336:LYS:HG3	1:A:342:CYS:SG	2.55	0.47
1:A:275:LEU:CD1	1:A:313:PRO:HB3	2.45	0.46
1:A:388:GLU:OE2	1:A:391:ARG:HD3	2.17	0.45
1:A:235:HIS:HE1	6:A:2017:HOH:O	1.99	0.45
1:B:419:LYS:HE3	3:B:1463:4IP:O9P	2.17	0.45
1:A:282:PRO:HD2	6:A:2033:HOH:O	2.16	0.44
1:B:404:HIS:HE1	6:B:2068:HOH:O	2.00	0.44
1:B:373:TYR:OH	1:B:404:HIS:CD2	2.71	0.43
1:B:311:THR:OG1	1:B:314:ARG:HB2	2.17	0.43
1:A:373:TYR:OH	1:A:404:HIS:HD2	2.01	0.43
1:B:312:LYS:HB3	1:B:313:PRO:HD3	1.99	0.43
1:A:424:PRO:HD3	1:A:440:ARG:NH1	2.34	0.42
1:B:273:GLU:O	1:B:276:THR:HB	2.20	0.42
1:A:296:ASP:HB3	1:A:299:ALA:HB2	2.02	0.42
1:B:213:GLU:N	1:B:214:PRO:CD	2.83	0.41
1:B:277:LYS:O	1:B:278:ALA:HB2	2.20	0.41
1:A:275:LEU:HD12	1:A:313:PRO:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:423:LEU:HD11	1:A:429:LEU:HG	2.01	0.41
1:A:319:ARG:HD2	6:A:2066:HOH:O	2.20	0.41

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	263/265 (99%)	254 (97%)	9 (3%)	0	100	100
1	B	244/265 (92%)	238 (98%)	5 (2%)	1 (0%)	34	24
All	All	507/530 (96%)	492 (97%)	14 (3%)	1 (0%)	47	39

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	276	THR

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	229/229 (100%)	215 (94%)	14 (6%)	18	6
1	B	216/229 (94%)	204 (94%)	12 (6%)	21	8
All	All	445/458 (97%)	419 (94%)	26 (6%)	20	7

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	212	SER
1	A	217	TYR
1	A	240	ARG
1	A	252	LEU
1	A	268	ARG
1	A	281	ARG
1	A	286	LYS
1	A	310	VAL
1	A	336	LYS
1	A	337	LYS
1	A	347	LYS
1	A	379	GLN
1	A	391	ARG
1	A	433	ARG
1	B	199	LYS
1	B	210	ARG
1	B	216	ARG
1	B	217	TYR
1	B	240	ARG
1	B	243	GLU
1	B	287	ASP
1	B	295	VAL
1	B	310	VAL
1	B	316	MET
1	B	352	ARG
1	B	461	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	235	HIS
1	A	379	GLN
1	A	404	HIS
1	A	406	HIS
1	A	427	GLN
1	B	404	HIS
1	B	408	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 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	$\# Z > 2$
5	SO4	A	1465	-	4,4,4	0.18	0	6,6,6	0.33	0
2	ADP	B	1462	4	24,29,29	1.16	2 (8%)	29,45,45	1.77	5 (17%)
3	4IP	A	1463	4	28,28,28	1.31	5 (17%)	42,46,46	1.16	3 (7%)
3	4IP	B	1463	-	28,28,28	1.39	5 (17%)	42,46,46	1.01	1 (2%)
5	SO4	A	1466	-	4,4,4	0.28	0	6,6,6	0.22	0
5	SO4	B	1465	-	4,4,4	0.13	0	6,6,6	0.45	0
2	ADP	A	1462	-	24,29,29	1.28	2 (8%)	29,45,45	1.37	1 (3%)

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	4IP	A	1463	4	-	0/20/44/44	0/1/1/1
2	ADP	B	1462	4	-	1/12/32/32	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	4IP	B	1463	-	-	6/20/44/44	0/1/1/1
2	ADP	A	1462	-	-	1/12/32/32	0/3/3/3

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1462	ADP	C2-N3	3.64	1.38	1.32
2	A	1462	ADP	C2-N1	3.34	1.40	1.33
3	B	1463	4IP	P1-O1	2.83	1.64	1.59
2	A	1462	ADP	C2-N3	2.69	1.36	1.32
3	A	1463	4IP	P4-O4	2.60	1.64	1.59
3	B	1463	4IP	P1-O1P	2.47	1.58	1.50
3	A	1463	4IP	P4-O9P	2.42	1.58	1.50
3	A	1463	4IP	P1-O1	2.40	1.63	1.59
3	B	1463	4IP	P5-OPG	2.34	1.58	1.50
2	B	1462	ADP	C2-N1	2.22	1.38	1.33
3	B	1463	4IP	P5-O5	2.08	1.63	1.59
3	A	1463	4IP	P5-OPG	2.07	1.57	1.50
3	B	1463	4IP	P5-OPH	-2.05	1.46	1.54
3	A	1463	4IP	P5-OPH	-2.00	1.47	1.54

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1462	ADP	N3-C2-N1	-6.40	118.67	128.68
2	A	1462	ADP	N3-C2-N1	-5.19	120.57	128.68
2	B	1462	ADP	C2-N1-C6	2.73	123.43	118.75
2	B	1462	ADP	C5-C6-N6	2.71	124.47	120.35
3	A	1463	4IP	O3P-P1-O2P	2.53	117.29	107.64
3	A	1463	4IP	O6P-P3-O4P	-2.50	100.91	110.68
2	B	1462	ADP	O4'-C4'-C5'	2.46	117.47	109.37
3	B	1463	4IP	O5P-P3-O4P	2.21	119.32	110.68
2	B	1462	ADP	C1'-N9-C4	-2.20	122.77	126.64
3	A	1463	4IP	O5P-P3-O4P	2.16	119.15	110.68

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1463	4IP	C5-O5-P5-OPG
3	B	1463	4IP	C3-O3-P3-O6P

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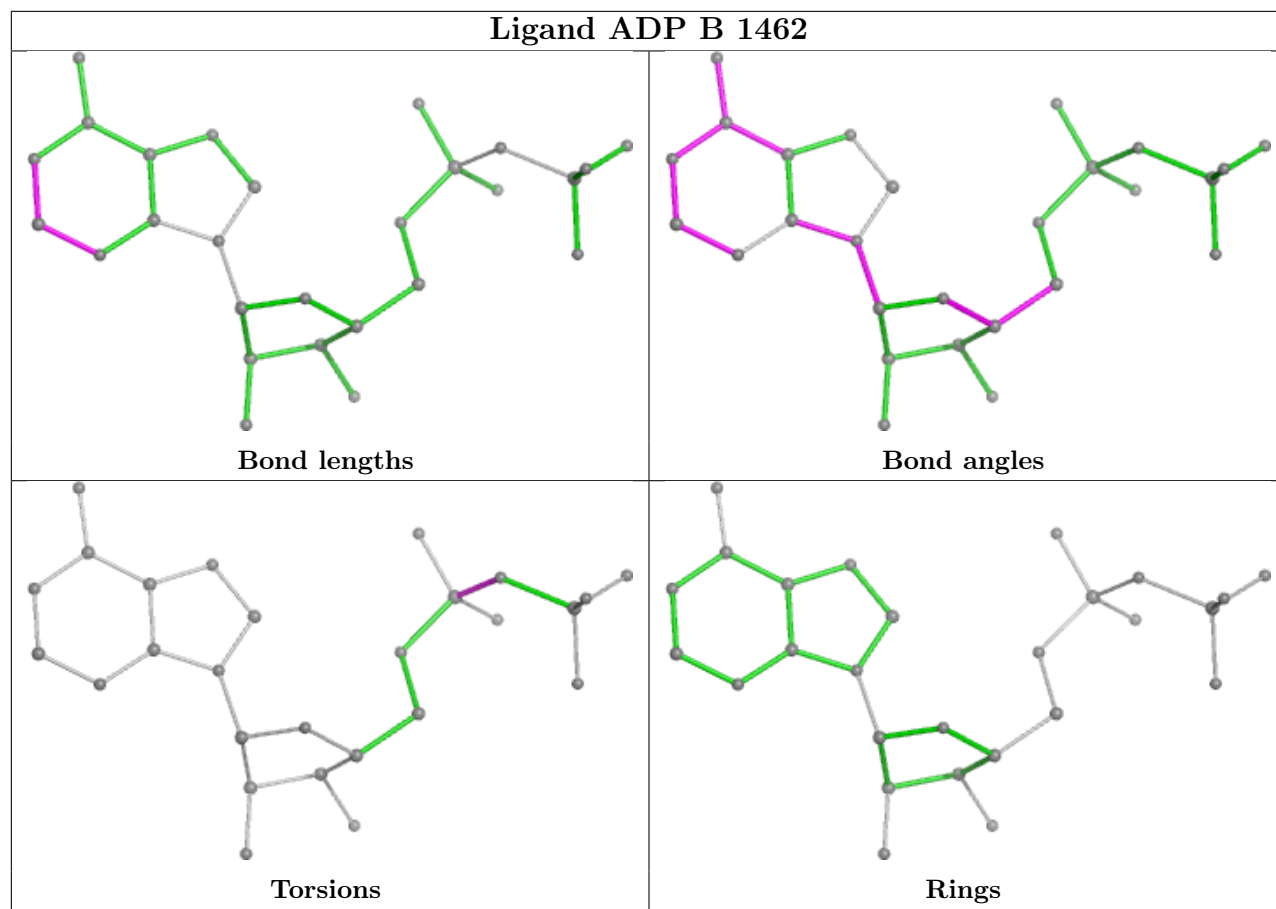
Mol	Chain	Res	Type	Atoms
3	B	1463	4IP	C5-O5-P5-OPF
2	A	1462	ADP	PB-O3A-PA-O2A
3	B	1463	4IP	C4-C5-O5-P5
2	B	1462	ADP	PB-O3A-PA-O1A
3	B	1463	4IP	C3-O3-P3-O5P
3	B	1463	4IP	C5-O5-P5-OPH

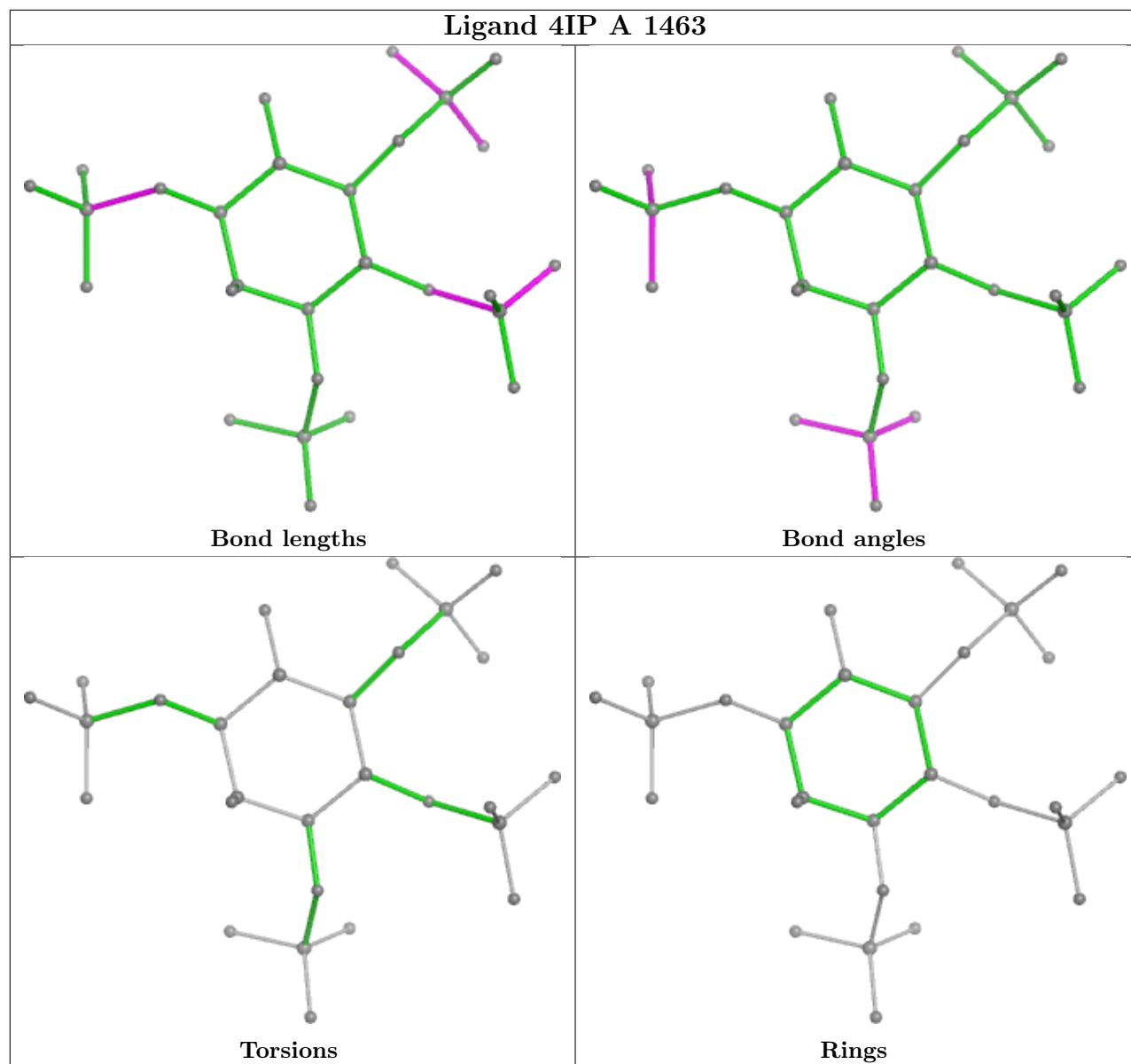
There are no ring outliers.

1 monomer is involved in 3 short contacts:

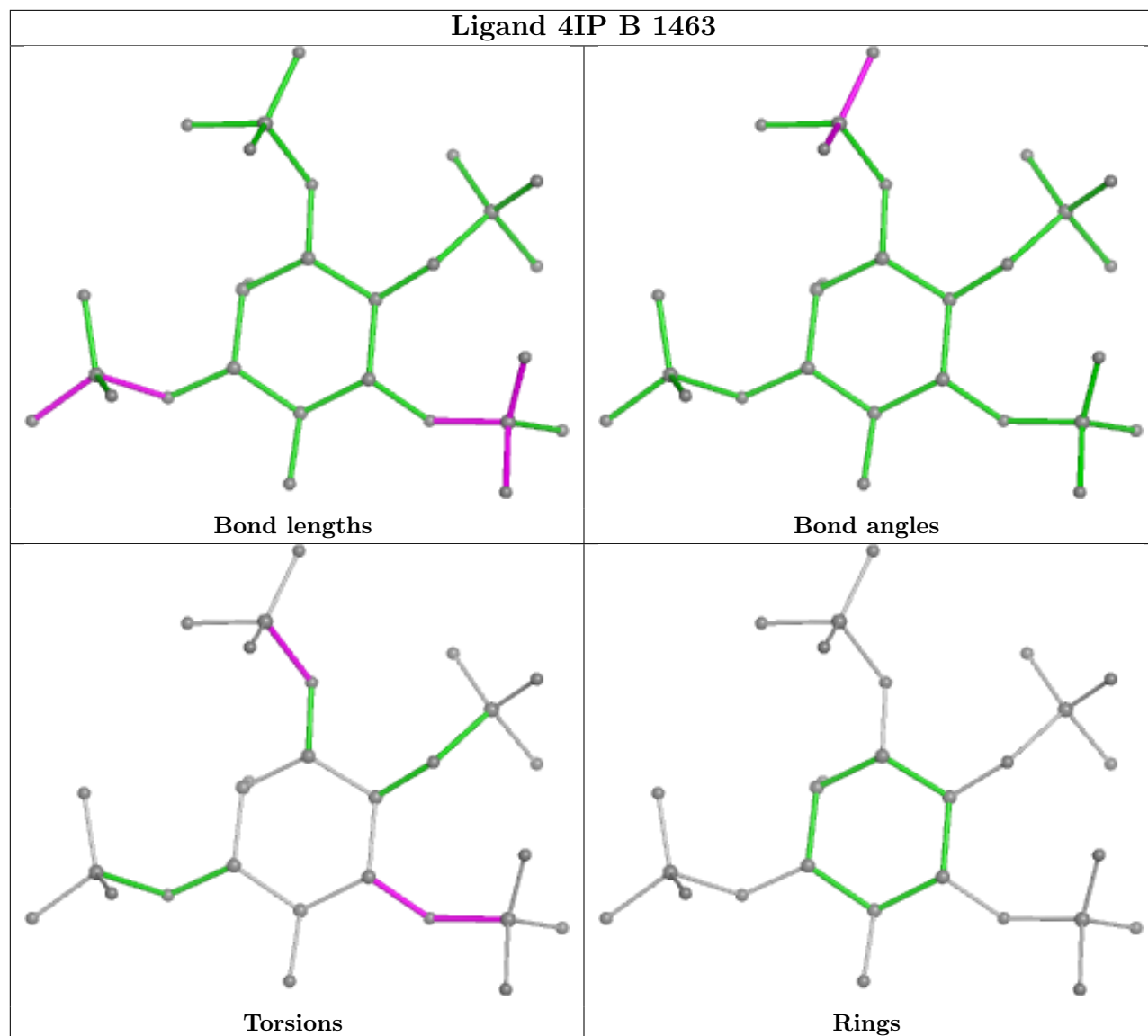
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1463	4IP	3	0

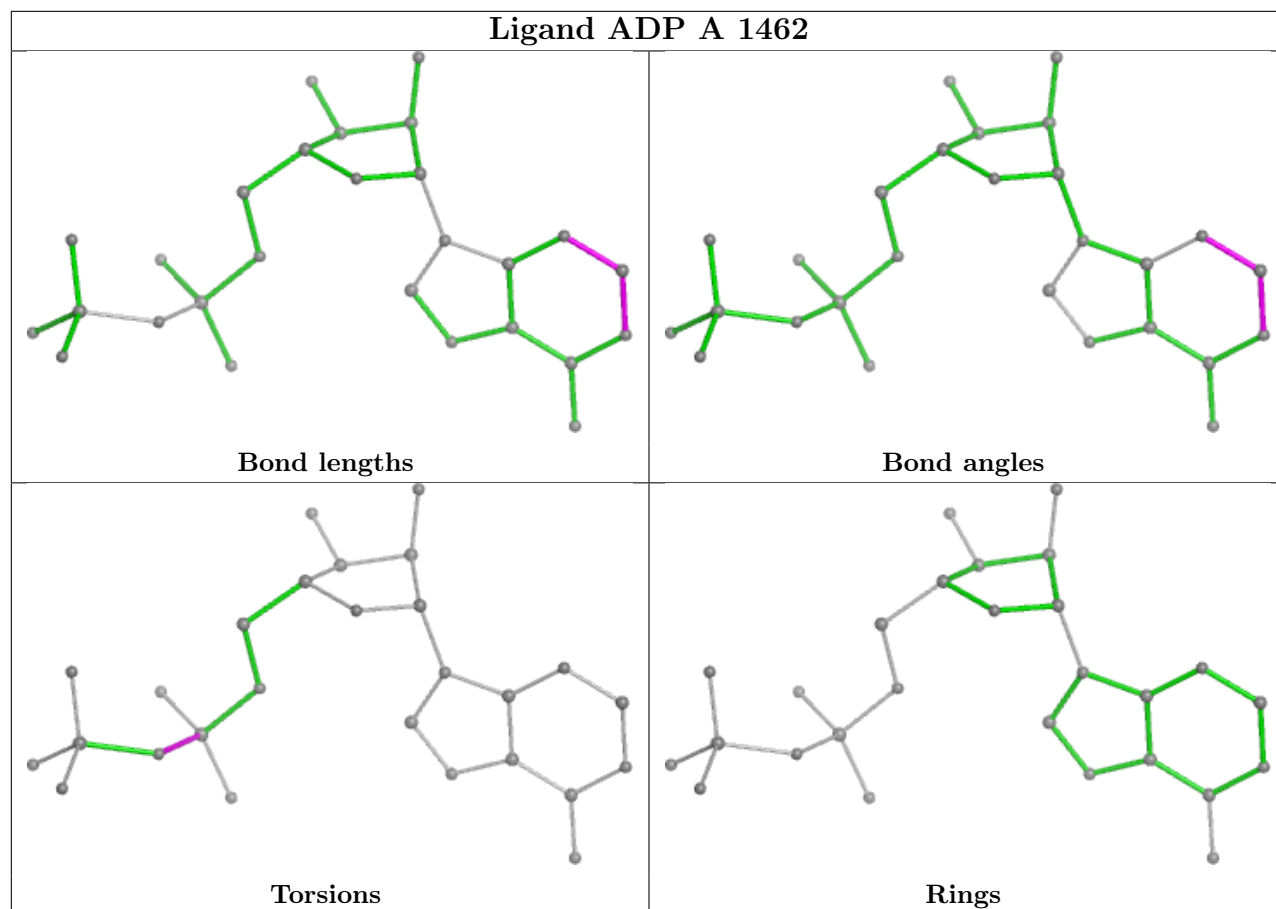
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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





Ligand 4IP B 1463





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	265/265 (100%)	0.76	25 (9%) 8 12	15, 29, 48, 53	0
1	B	250/265 (94%)	1.25	50 (20%) 1 1	14, 30, 52, 59	0
All	All	515/530 (97%)	1.00	75 (14%) 2 3	14, 30, 51, 59	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	284	LEU	9.1
1	B	310	VAL	7.9
1	B	311	THR	7.4
1	B	313	PRO	7.1
1	B	309	ALA	6.8
1	B	338	ALA	6.7
1	B	278	ALA	6.6
1	B	283	LYS	6.5
1	A	294	ALA	6.1
1	B	288	MET	6.1
1	B	286	LYS	5.9
1	B	295	VAL	5.8
1	B	287	ASP	5.7
1	B	276	THR	5.4
1	A	338	ALA	5.3
1	B	312	LYS	4.7
1	B	289	TYR	4.7
1	B	242	GLY	4.6
1	B	285	ARG	4.6
1	B	238	VAL	4.4
1	B	277	LYS	4.4
1	B	315	TYR	4.3
1	B	290	LYS	4.2
1	A	241	ASP	4.2

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Mol	Chain	Res	Type	RSRZ
1	B	275	LEU	4.1
1	A	293	LEU	4.0
1	B	294	ALA	4.0
1	B	293	LEU	4.0
1	B	198	PHE	3.9
1	A	295	VAL	3.8
1	A	213	GLU	3.7
1	B	245	TYR	3.6
1	B	202	GLY	3.5
1	A	204	SER	3.5
1	B	292	MET	3.4
1	B	316	MET	3.4
1	A	202	GLY	3.3
1	B	239	GLU	3.3
1	B	243	GLU	3.2
1	B	244	SER	3.2
1	B	291	LYS	3.2
1	B	273	GLU	3.1
1	B	425	ASP	3.1
1	A	290	LYS	3.0
1	B	211	CYS	3.0
1	B	241	ASP	3.0
1	A	339	ASP	2.9
1	A	240	ARG	2.8
1	A	296	ASP	2.8
1	B	314	ARG	2.7
1	A	281	ARG	2.7
1	A	425	ASP	2.6
1	A	283	LYS	2.6
1	A	203	THR	2.6
1	A	426	GLY	2.6
1	B	274	GLU	2.5
1	B	210	ARG	2.4
1	B	201	ALA	2.3
1	B	339	ASP	2.3
1	A	276	THR	2.3
1	A	297	PRO	2.3
1	A	284	LEU	2.2
1	B	208	LEU	2.2
1	B	199	LYS	2.2
1	B	452	ILE	2.2
1	B	240	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	402	PHE	2.1
1	B	271	LEU	2.1
1	A	308	ARG	2.1
1	B	318	TRP	2.1
1	A	341	SER	2.1
1	A	412	VAL	2.0
1	A	242	GLY	2.0
1	B	197	SER	2.0
1	B	328	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 monosaccharides in this entry.

6.4 Ligands [i](#)

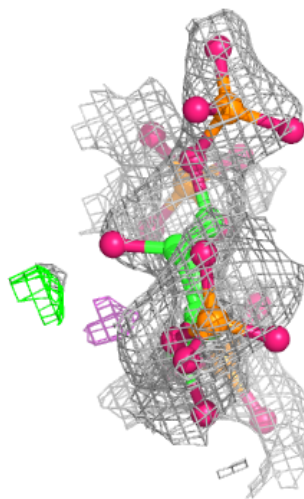
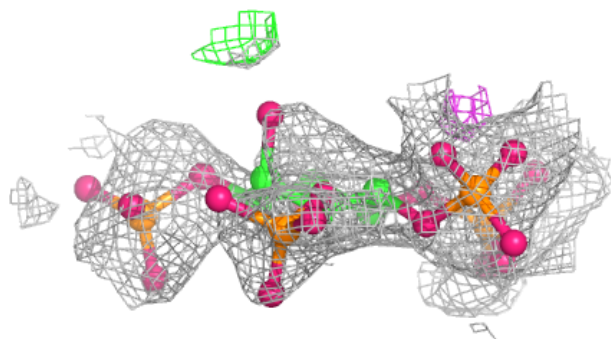
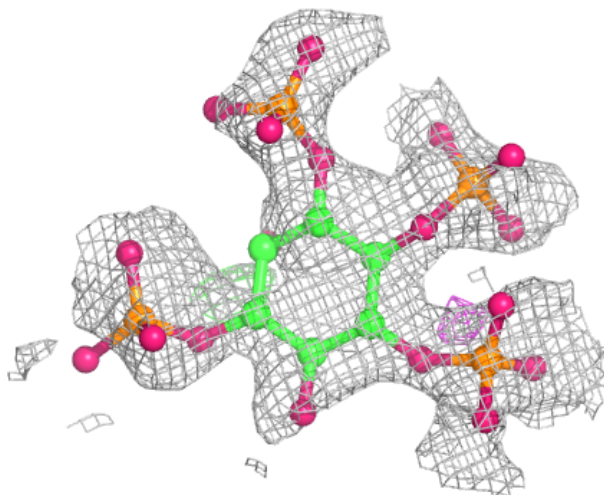
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	4IP	B	1463	28/28	0.73	0.28	55,62,65,66	28
4	MN	A	1464	1/1	0.89	0.12	51,51,51,51	1
5	SO4	B	1465	5/5	0.89	0.19	35,35,37,38	5
3	4IP	A	1463	28/28	0.90	0.11	51,55,63,65	0
5	SO4	A	1465	5/5	0.93	0.19	35,37,37,37	5
5	SO4	A	1466	5/5	0.94	0.17	34,34,36,36	5
2	ADP	B	1462	27/27	0.94	0.14	28,36,50,52	0
4	MN	B	1464	1/1	0.95	0.21	42,42,42,42	1
2	ADP	A	1462	27/27	0.95	0.10	25,31,46,47	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.

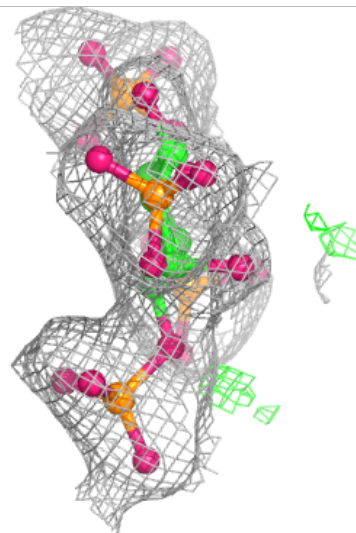
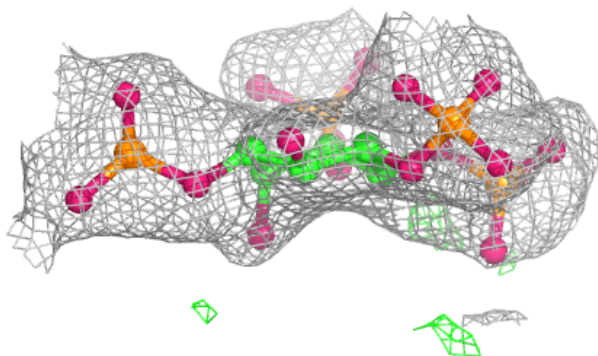
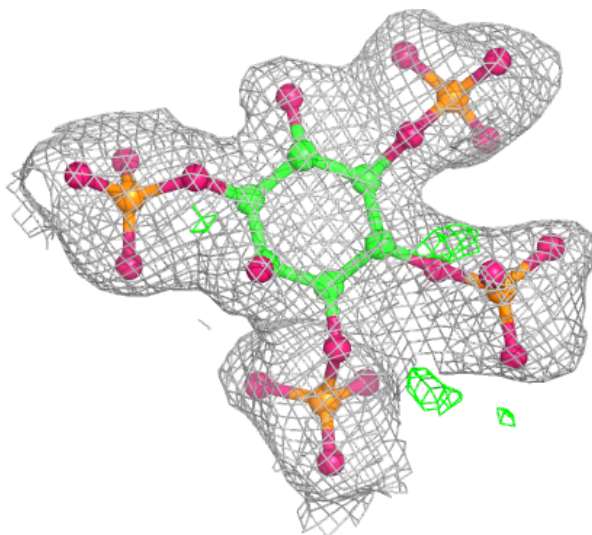
Electron density around 4IP B 1463:

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



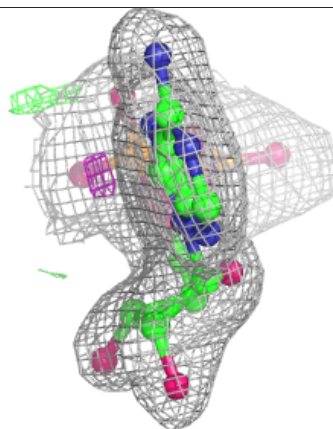
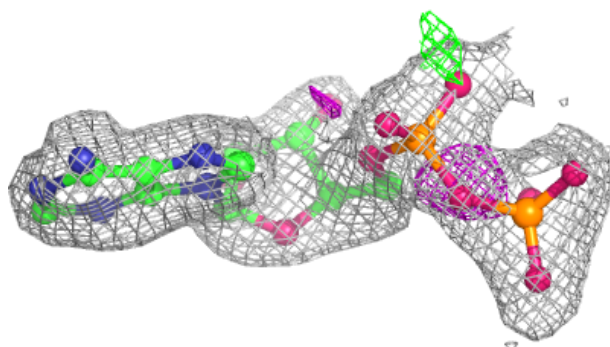
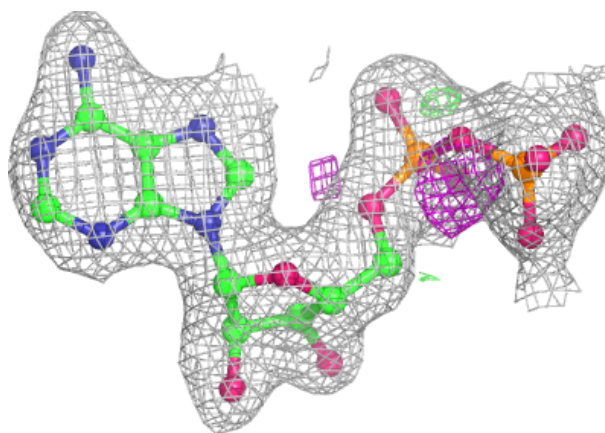
Electron density around 4IP A 1463:

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

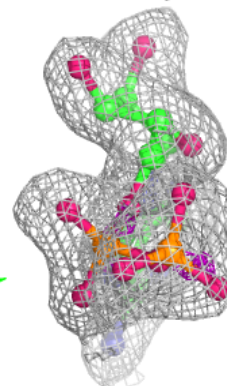
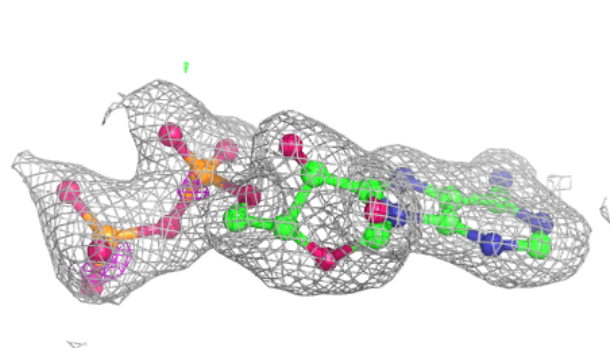
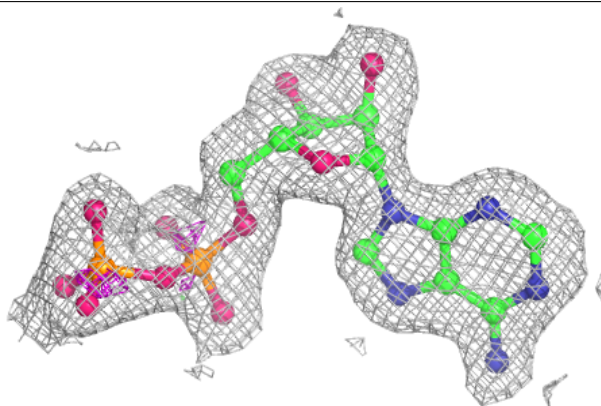


Electron density around ADP B 1462:

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

**Electron density around ADP A 1462:**

$2mF_o-DF_c$ (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.