



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:14 AM UTC

PDB ID : 1W2D / pdb_00001w2d
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	:	4-5-2 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

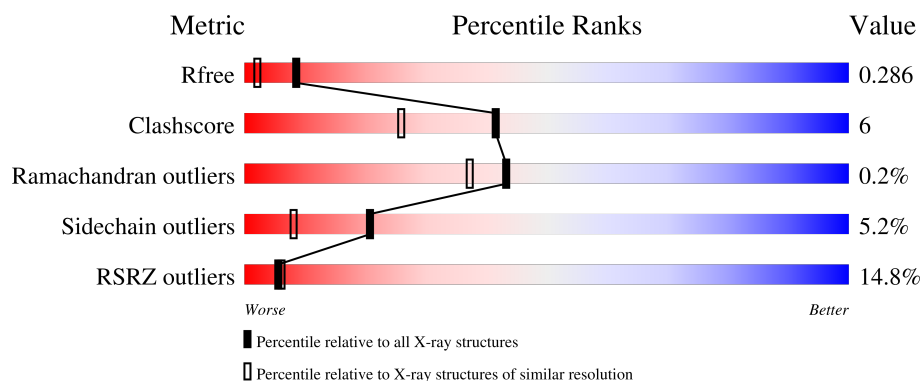
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.94 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}	180053	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	12% 82% 17% .
1	B	265	17% 79% 14% 6%

2 Entry composition [i](#)

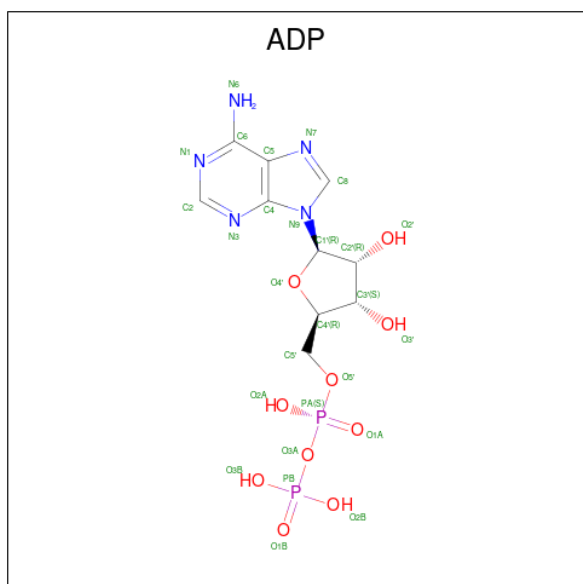
There are 6 unique types of molecules in this entry. The entry contains 4524 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

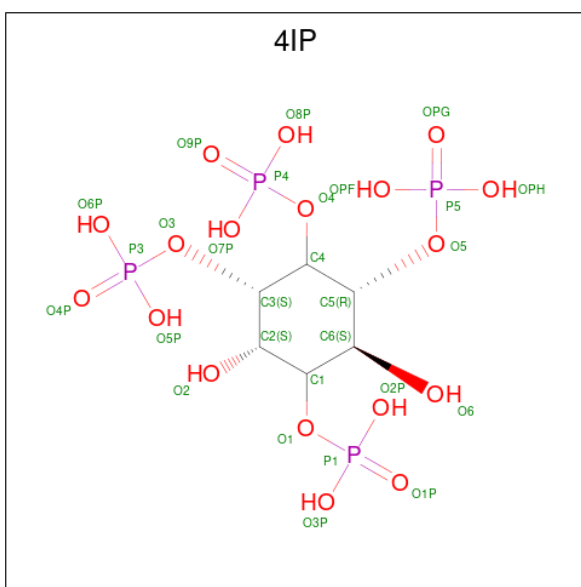
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	0	0
			2133	1334	388	399	12			
1	B	250	Total	C	N	O	S	0	0	0
			2005	1260	361	372	12			

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: 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 (CCD ID: 4IP) (formula: $C_6H_{16}O_{18}P_4$).

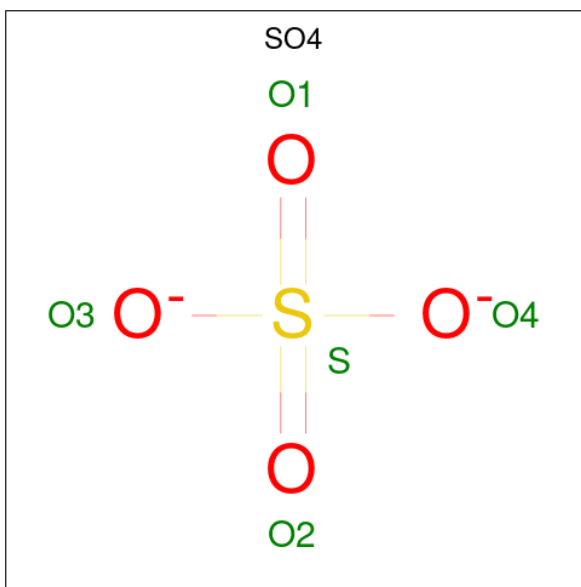


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			28	6	18	4		
3	B	1	Total	C	O	P	0	0
			28	6	18	4		

- Molecule 4 is MANGANESE (II) ION (CCD ID: 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 (CCD ID: 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		

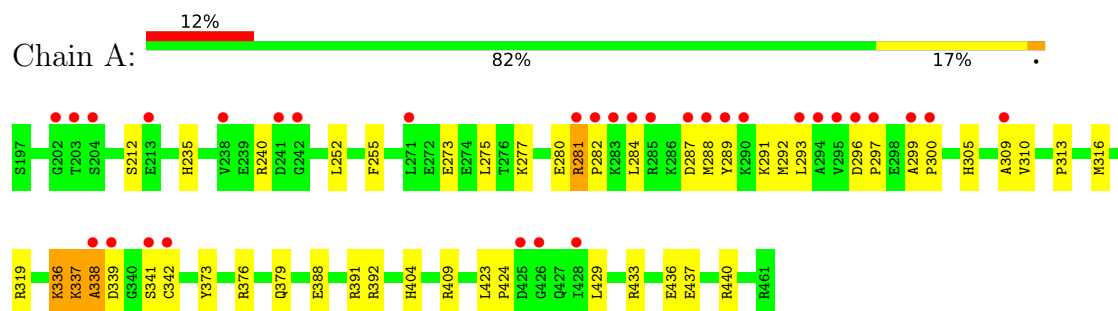
- Molecule 6 is water.

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

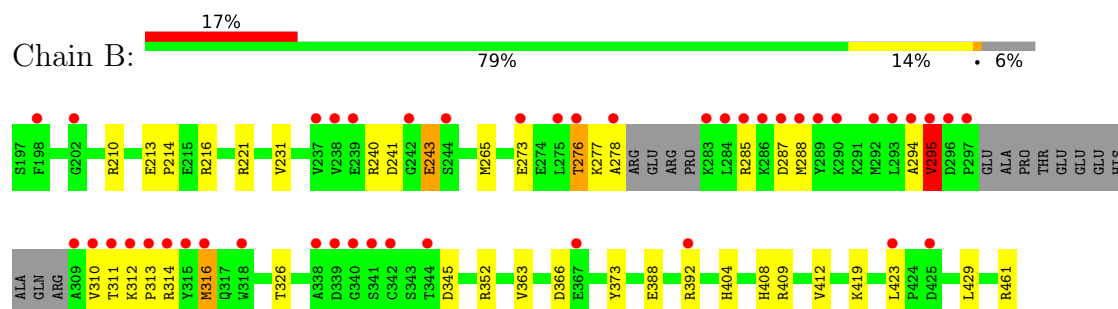
3 Residue-property plots

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 48.80 – 1.94	Depositor EDS
% Data completeness (in resolution range)	96.9 (48.80-1.94) 95.0 (48.80-1.94)	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 (Not available) , 0.286	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 , 40.1	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: MN, SO4, ADP, 4IP

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	1.14	2/2173 (0.1%)	1.11	3/2924 (0.1%)
1	B	1.13	4/2040 (0.2%)	1.07	1/2741 (0.0%)
All	All	1.14	6/4213 (0.1%)	1.09	4/5665 (0.1%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	326	THR	CA-CB	6.12	1.63	1.53
1	B	265	MET	SD-CE	5.71	1.93	1.79
1	A	316	MET	SD-CE	5.57	1.93	1.79
1	B	412	VAL	C-O	5.53	1.29	1.23
1	B	295	VAL	CA-CB	5.26	1.60	1.54
1	A	337	LYS	CA-C	-5.10	1.46	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	289	TYR	N-CA-C	6.52	118.18	111.14
1	A	288	MET	N-CA-C	-6.03	105.19	112.54
1	B	243	GLU	N-CA-C	5.88	117.90	109.14
1	A	338	ALA	N-CA-C	-5.12	105.78	111.36

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	2133	0	2113	36	0
1	B	2005	0	1996	30	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	54	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 (54) 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	0.87
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
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:C	1:B:294:ALA:CB	2.71	0.61
1:A:338:ALA:O	1:B:294:ALA:HB1	1.99	0.61
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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:285:ARG:NH2	1:B:288:MET:HE3	2.32	0.44
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:A:292:MET:HE2	1:A:292:MET:HB3	1.86	0.43
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:280:GLU:C	1:A:281:ARG:CG	2.92	0.43
1:A:424:PRO:HD3	1:A:440:ARG:NH1	2.34	0.42
1:A:436:GLU:O	1:A:437:GLU:C	2.62	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
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
1:B:316:MET:CE	1:B:316:MET:HA	2.50	0.41

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	263/265 (99%)	254 (97%)	9 (3%)	0	100	100
1	B	244/265 (92%)	238 (98%)	5 (2%)	1 (0%)	30	22
All	All	507/530 (96%)	492 (97%)	14 (3%)	1 (0%)	43	37

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	276	THR

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	229/229 (100%)	219 (96%)	10 (4%)	25	12
1	B	216/229 (94%)	203 (94%)	13 (6%)	17	5
All	All	445/458 (97%)	422 (95%)	23 (5%)	21	8

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	212	SER
1	A	240	ARG
1	A	252	LEU
1	A	273	GLU
1	A	277	LYS

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Mol	Chain	Res	Type
1	A	281	ARG
1	A	310	VAL
1	A	336	LYS
1	A	392	ARG
1	A	433	ARG
1	B	210	ARG
1	B	216	ARG
1	B	231	VAL
1	B	240	ARG
1	B	241	ASP
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	363	VAL
1	B	461	ARG

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

Mol	Chain	Res	Type
1	A	235	HIS
1	A	379	GLN
1	A	404	HIS
1	A	406	HIS
1	B	404	HIS
1	B	408	HIS

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

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	B	1465	-	4,4,4	0.18	0	6,6,6	0.44	0
3	4IP	B	1463	-	28,28,28	1.43	5 (17%)	46,46,46	1.19	5 (10%)
5	SO4	A	1465	-	4,4,4	0.18	0	6,6,6	0.30	0
2	ADP	A	1462	-	28,29,29	1.45	4 (14%)	43,45,45	1.75	9 (20%)
5	SO4	A	1466	-	4,4,4	0.28	0	6,6,6	0.15	0
3	4IP	A	1463	4	28,28,28	1.36	6 (21%)	46,46,46	1.33	6 (13%)
2	ADP	B	1462	4	28,29,29	1.31	4 (14%)	43,45,45	1.84	10 (23%)

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	ADP	A	1462	-	-	1/16/32/32	0/3/3/3
3	4IP	A	1463	4	-	0/20/44/44	0/1/1/1
3	4IP	B	1463	-	-	5/20/44/44	0/1/1/1
2	ADP	B	1462	4	-	1/16/32/32	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1462	ADP	PA-O3A	4.33	1.64	1.59
2	A	1462	ADP	PA-O3A	3.54	1.63	1.59
2	A	1462	ADP	C2-N1	3.48	1.40	1.33
3	B	1463	4IP	P1-O1	2.93	1.64	1.59
3	A	1463	4IP	P4-O4	2.68	1.64	1.59
3	B	1463	4IP	P1-O1P	2.58	1.58	1.50
3	A	1463	4IP	P4-O9P	2.54	1.58	1.50
3	A	1463	4IP	P1-O1	2.46	1.63	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1463	4IP	P5-OPG	2.45	1.58	1.50
2	A	1462	ADP	C5-N7	-2.40	1.34	1.39
2	B	1462	ADP	C2-N1	2.31	1.38	1.33
2	A	1462	ADP	C1'-N9	2.28	1.52	1.46
2	B	1462	ADP	C2-N3	2.26	1.38	1.33
2	B	1462	ADP	C8-N7	2.21	1.35	1.31
3	A	1463	4IP	P5-OPG	2.17	1.57	1.50
3	B	1463	4IP	P5-O5	2.12	1.63	1.59
3	B	1463	4IP	P5-OPH	-2.11	1.46	1.54
3	A	1463	4IP	P5-OPH	-2.07	1.47	1.54
3	A	1463	4IP	P1-O1P	2.05	1.56	1.50

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1462	ADP	N3-C2-N1	-6.54	118.67	128.58
2	A	1462	ADP	N3-C2-N1	-5.29	120.57	128.58
2	A	1462	ADP	C5-C4-N3	-4.81	120.10	126.72
3	B	1463	4IP	P1-O1-C1	-3.91	112.98	123.43
2	B	1462	ADP	N9-C8-N7	-3.38	109.14	113.94
2	B	1462	ADP	C5-N7-C8	3.25	108.55	103.45
2	A	1462	ADP	C2-N3-C4	3.19	119.63	111.83
3	A	1463	4IP	P1-O1-C1	-3.08	115.22	123.43
2	B	1462	ADP	C5-C4-N3	-3.07	122.49	126.72
2	A	1462	ADP	C5-N7-C8	3.00	108.17	103.45
2	B	1462	ADP	C2-N3-C4	2.98	119.10	111.83
2	B	1462	ADP	C2-N1-C6	2.86	123.43	118.73
2	A	1462	ADP	N3-C4-N9	2.79	131.91	127.17
2	B	1462	ADP	C4-C5-N7	-2.75	107.44	110.58
2	A	1462	ADP	C4-C5-N7	-2.68	107.51	110.58
3	A	1463	4IP	O6P-P3-O4P	-2.55	100.91	110.83
2	B	1462	ADP	O4'-C4'-C5'	2.54	117.47	109.33
3	A	1463	4IP	O3P-P1-O2P	2.53	117.29	107.80
2	A	1462	ADP	C6-C5-C4	2.41	120.46	117.18
3	A	1463	4IP	P3-O3-C3	-2.39	117.06	123.43
3	A	1463	4IP	P4-O4-C4	2.38	129.78	123.43
2	B	1462	ADP	C5-C4-N9	2.33	108.35	105.81
2	A	1462	ADP	N9-C8-N7	-2.30	110.67	113.94
2	B	1462	ADP	O3A-PA-O1A	-2.29	103.82	110.70
3	B	1463	4IP	O5P-P3-O4P	2.18	119.32	110.83
3	B	1463	4IP	OPH-P5-O5	2.17	114.31	105.85
3	A	1463	4IP	O5P-P3-O4P	2.13	119.15	110.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1463	4IP	O4-P4-O9P	-2.12	101.77	109.33
2	A	1462	ADP	O4'-C1'-N9	2.06	112.05	108.09
3	B	1463	4IP	O6P-P3-O4P	-2.01	103.02	110.83

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1463	4IP	C5-O5-P5-OPG
2	A	1462	ADP	PB-O3A-PA-O2A
3	B	1463	4IP	C3-O3-P3-O6P
3	B	1463	4IP	C5-O5-P5-OPF
3	B	1463	4IP	C4-C5-O5-P5
2	B	1462	ADP	PB-O3A-PA-O1A
3	B	1463	4IP	C5-O5-P5-OPH

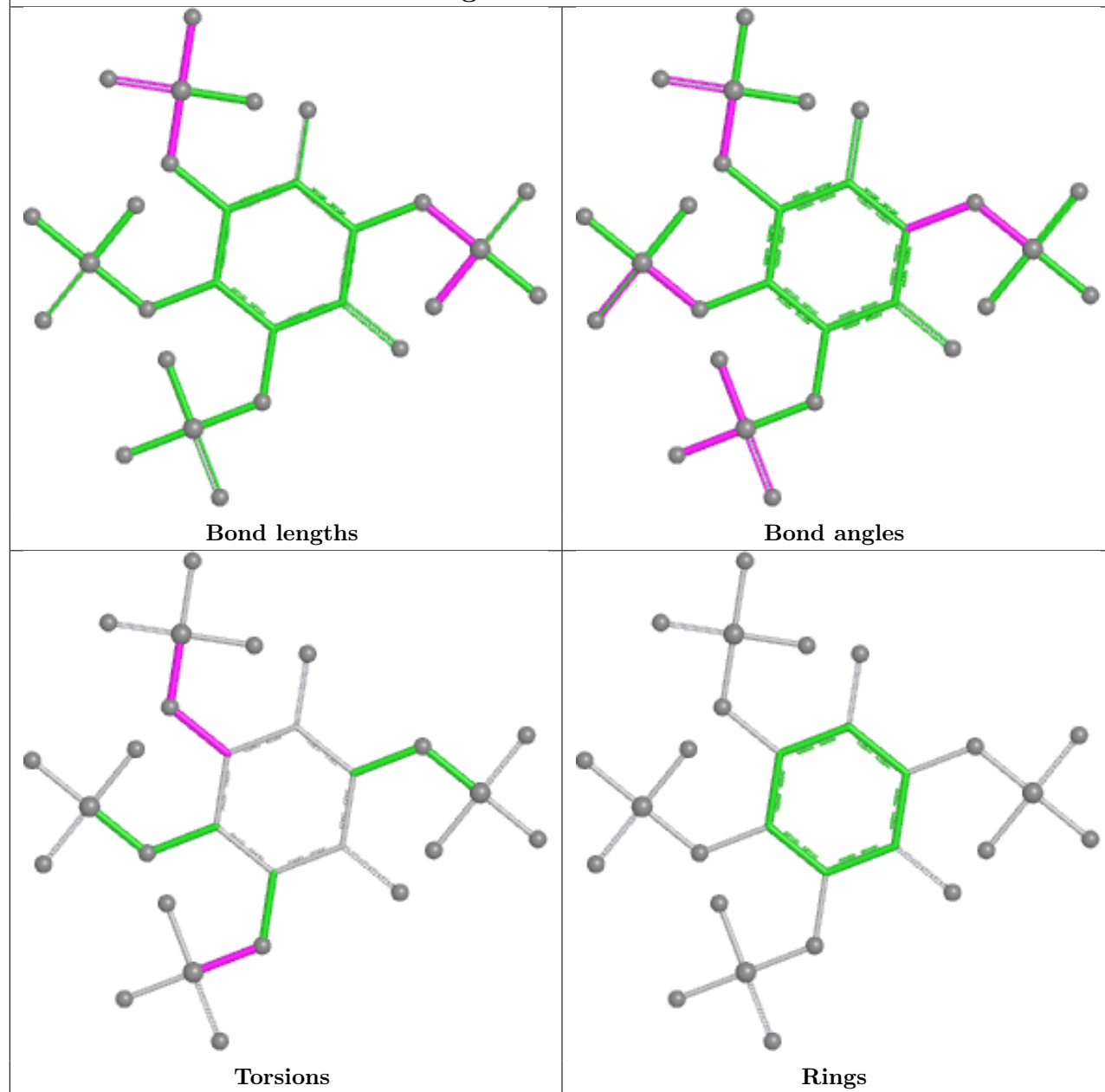
There are no ring outliers.

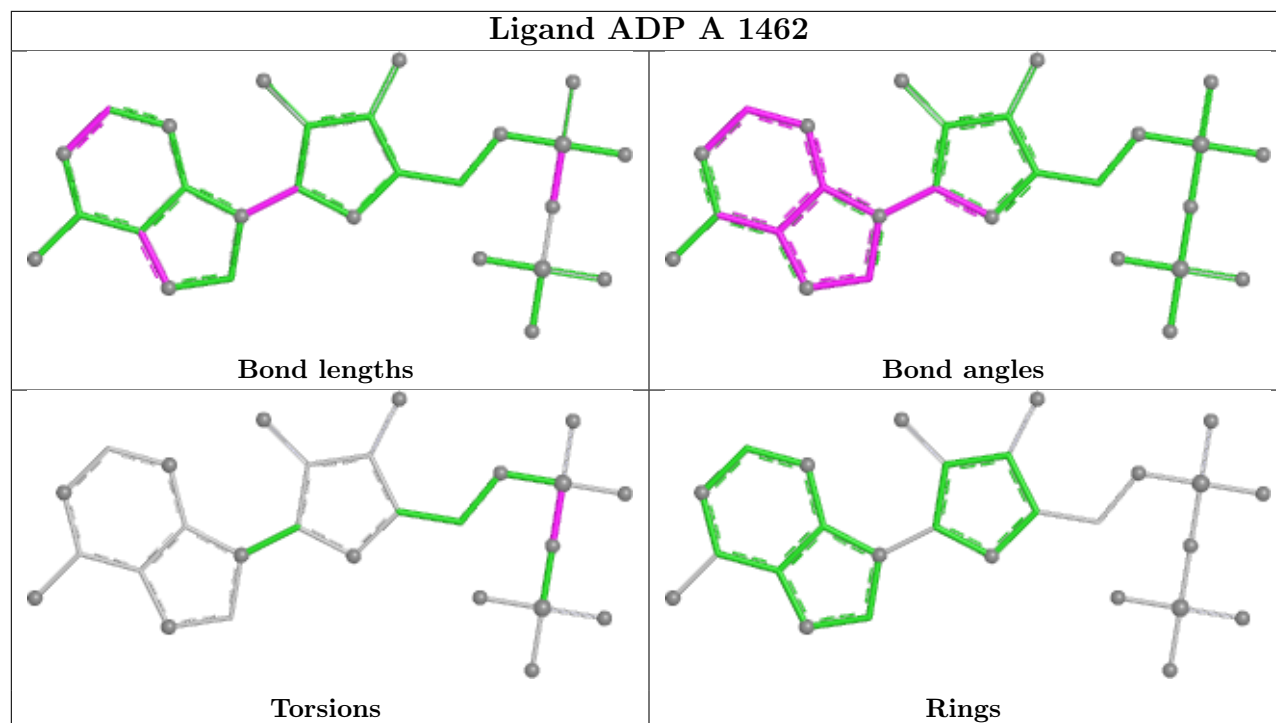
1 monomer is involved in 3 short contacts:

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

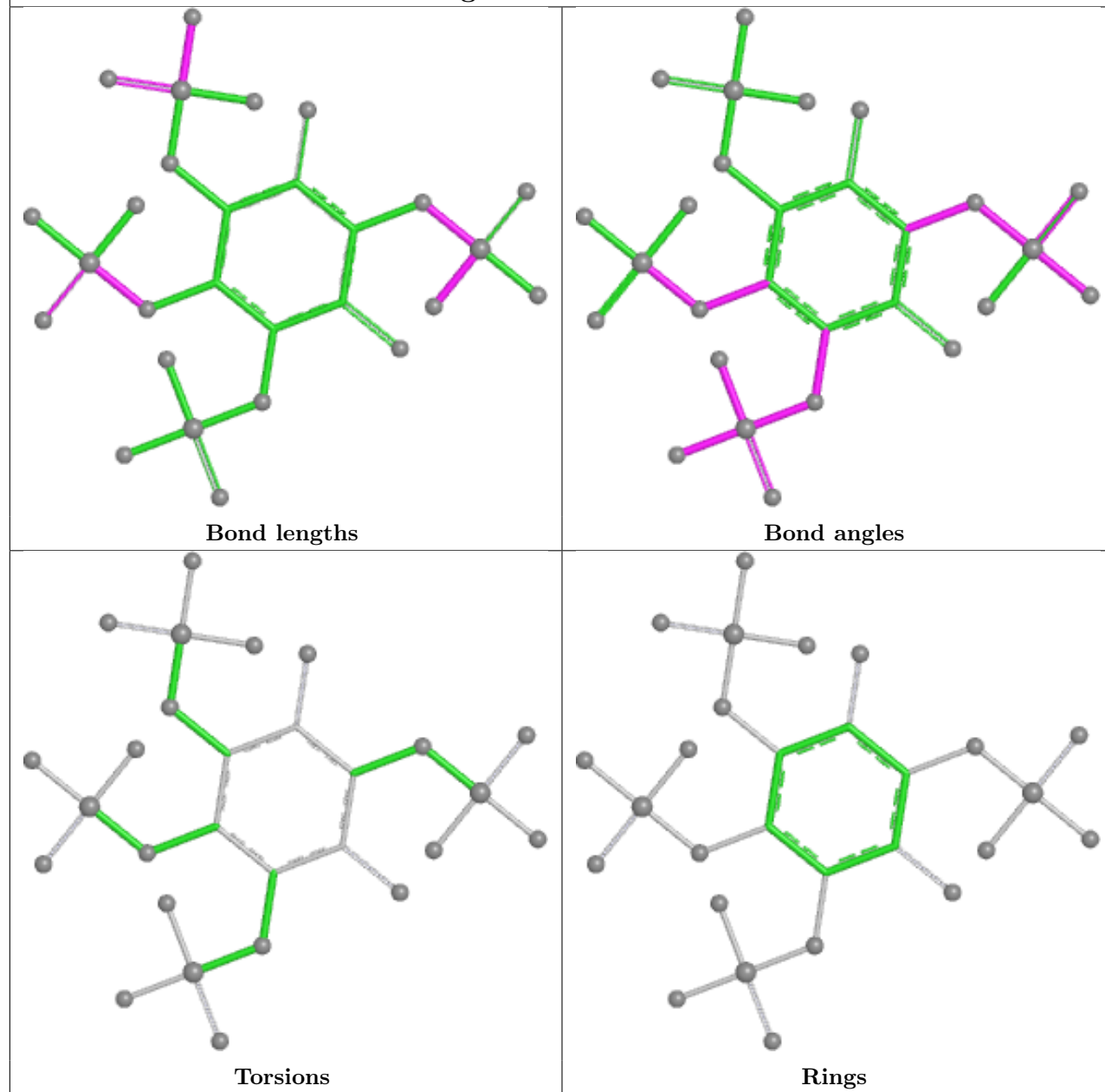
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

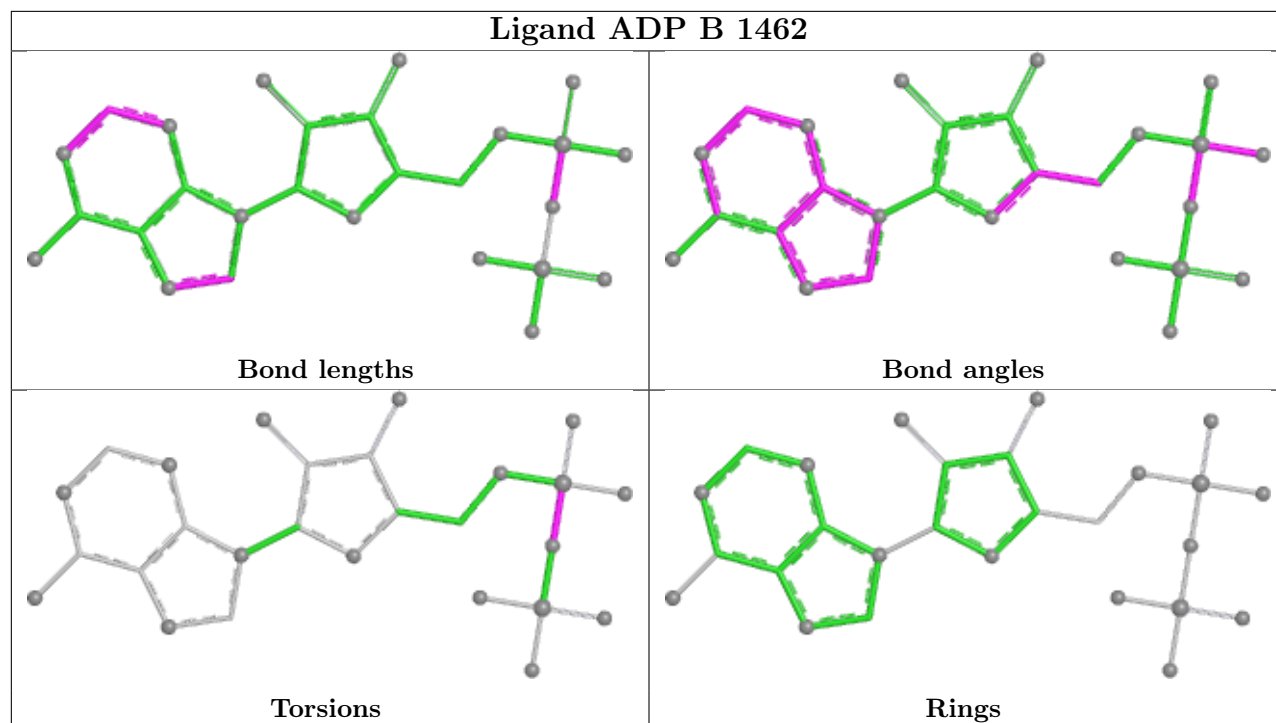
Ligand 4IP B 1463





Ligand 4IP A 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%)	0.79	32 (12%) 8 10	15, 29, 48, 53	0
1	B	250/265 (94%)	1.01	44 (17%) 4 4	14, 30, 52, 59	0
All	All	515/530 (97%)	0.89	76 (14%) 5 6	14, 30, 51, 59	0

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	310	VAL	7.9
1	B	295	VAL	7.2
1	B	338	ALA	7.1
1	A	338	ALA	6.8
1	A	294	ALA	5.0
1	B	311	THR	5.0
1	B	294	ALA	4.8
1	B	284	LEU	4.6
1	A	241	ASP	4.5
1	B	309	ALA	4.5
1	B	293	LEU	4.4
1	B	313	PRO	4.4
1	A	339	ASP	4.4
1	A	295	VAL	4.2
1	A	296	ASP	4.1
1	B	288	MET	4.1
1	A	293	LEU	4.1
1	B	242	GLY	4.0
1	B	292	MET	3.9
1	B	289	TYR	3.8
1	B	318	TRP	3.8
1	A	204	SER	3.7
1	A	341	SER	3.7
1	A	213	GLU	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	290	LYS	3.6
1	B	287	ASP	3.6
1	B	290	LYS	3.5
1	B	278	ALA	3.5
1	B	285	ARG	3.4
1	A	288	MET	3.3
1	B	316	MET	3.3
1	A	287	ASP	3.1
1	A	284	LEU	3.1
1	A	342	CYS	3.1
1	A	426	GLY	3.1
1	B	202	GLY	3.1
1	B	425	ASP	3.0
1	A	202	GLY	3.0
1	B	286	LYS	3.0
1	B	315	TYR	2.9
1	B	297	PRO	2.9
1	B	275	LEU	2.9
1	B	237	VAL	2.8
1	B	276	THR	2.8
1	A	289	TYR	2.7
1	A	297	PRO	2.7
1	A	425	ASP	2.7
1	B	244	SER	2.6
1	B	341	SER	2.6
1	B	238	VAL	2.5
1	B	314	ARG	2.4
1	B	340	GLY	2.4
1	A	238	VAL	2.4
1	B	296	ASP	2.4
1	A	203	THR	2.4
1	B	392	ARG	2.4
1	B	239	GLU	2.4
1	A	428	ILE	2.3
1	B	342	CYS	2.3
1	B	339	ASP	2.3
1	A	242	GLY	2.3
1	B	367	GLU	2.3
1	A	283	LYS	2.3
1	A	282	PRO	2.2
1	B	273	GLU	2.2
1	A	300	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	198	PHE	2.1
1	B	283	LYS	2.1
1	A	299	ALA	2.1
1	A	281	ARG	2.1
1	A	271	LEU	2.1
1	B	423	LEU	2.1
1	A	309	ALA	2.1
1	B	344	THR	2.0
1	A	285	ARG	2.0
1	B	312	LYS	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](#)

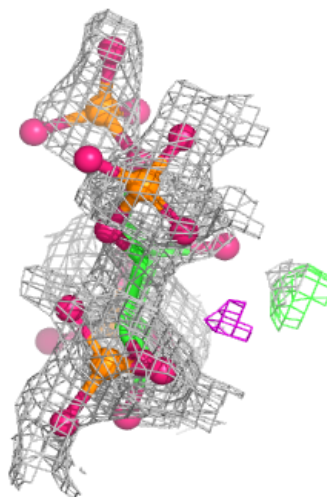
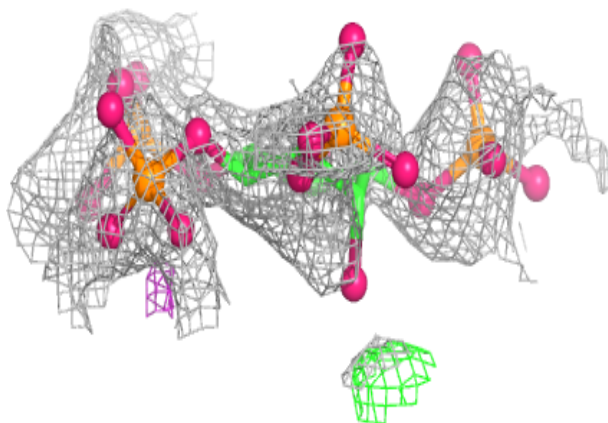
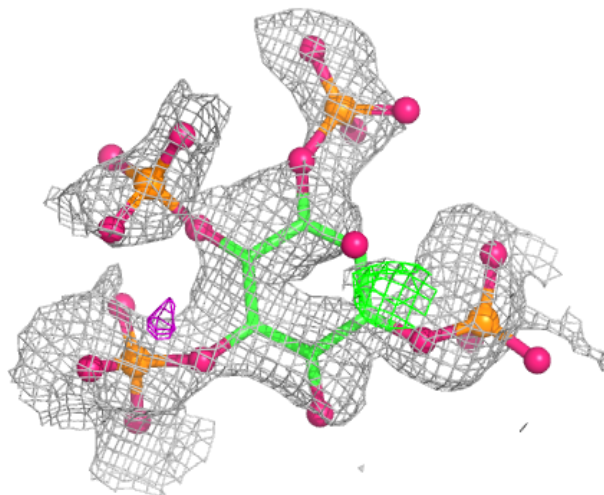
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	4IP	B	1463	28/28	0.63	0.20	55,62,65,66	28
5	SO4	B	1465	5/5	0.81	0.15	35,35,37,38	5
5	SO4	A	1465	5/5	0.86	0.21	35,37,37,37	5
5	SO4	A	1466	5/5	0.86	0.18	34,34,36,36	5
3	4IP	A	1463	28/28	0.86	0.12	51,55,63,65	0
4	MN	A	1464	1/1	0.87	0.11	51,51,51,51	1
2	ADP	B	1462	27/27	0.91	0.10	28,36,50,52	0
4	MN	B	1464	1/1	0.93	0.12	42,42,42,42	1
2	ADP	A	1462	27/27	0.94	0.08	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.

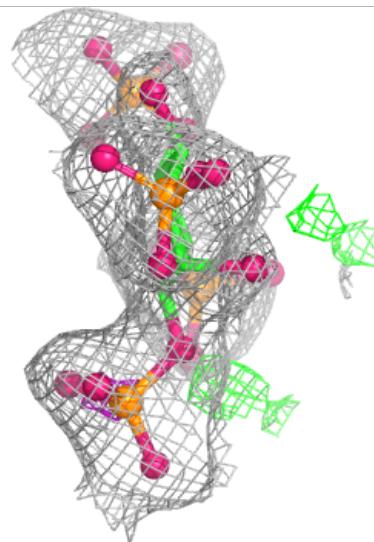
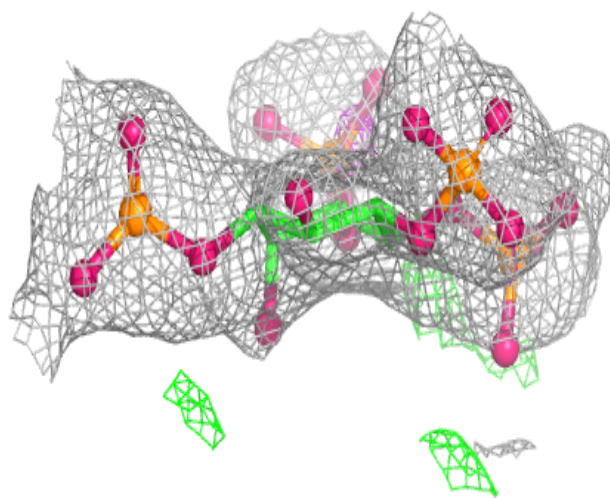
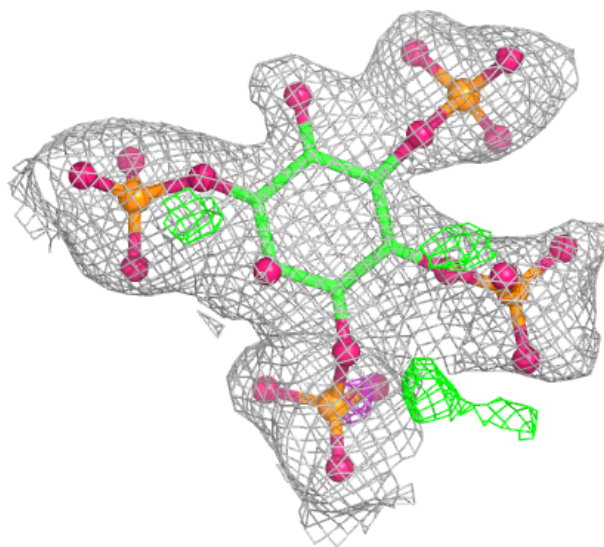
Electron density around 4IP B 1463:

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



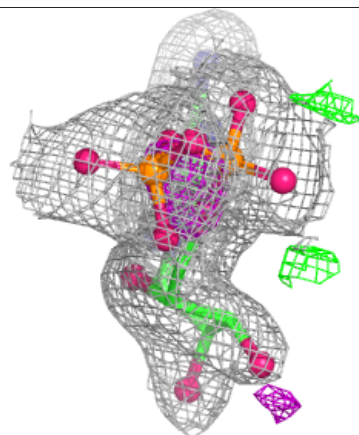
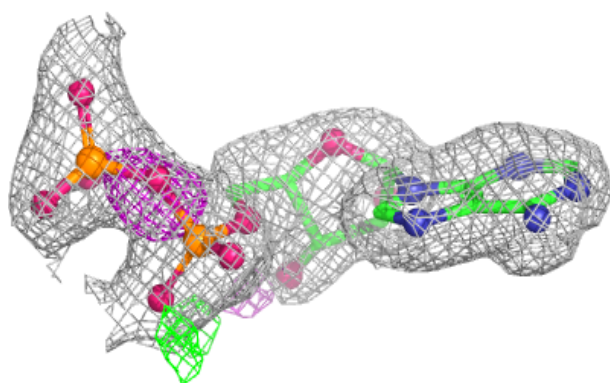
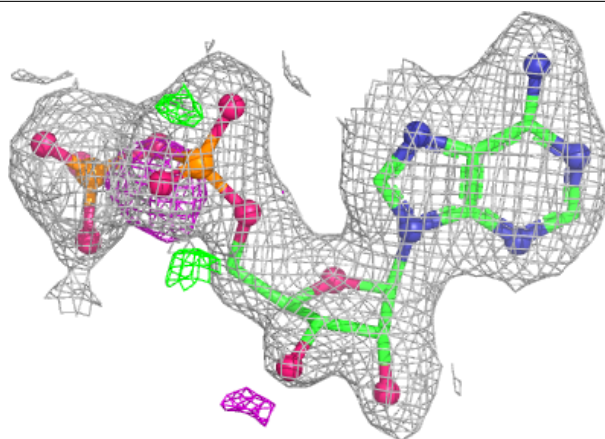
Electron density around 4IP A 1463:

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

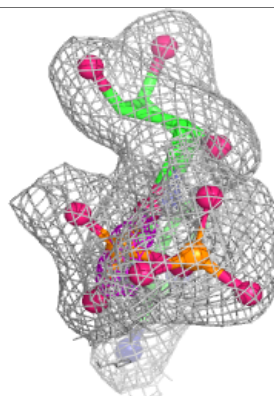
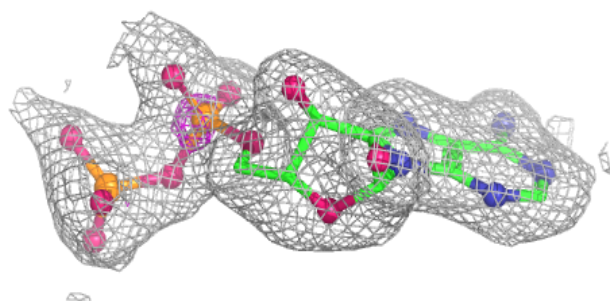
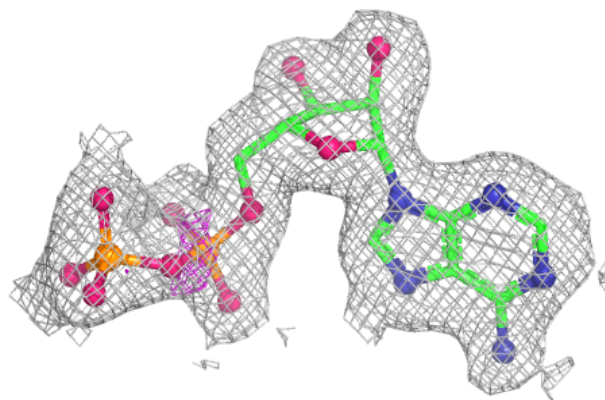


Electron density around ADP B 1462:

$2mF_o-DF_c$ (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 [i](#)

There are no such residues in this entry.