



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 28, 2020 – 03:14 PM BST

PDB ID : 6FJK
Title : Inositol 1,3,4,5,6-pentakisphosphate 2-kinase from *A. thaliana* in complex with myo-IP6 and ADP
Authors : Whitfield, H.L.; Brearley, C.A.; Hemmings, A.M.
Deposited on : 2018-01-22
Resolution : 2.02 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.14.6
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) : 2.14.6

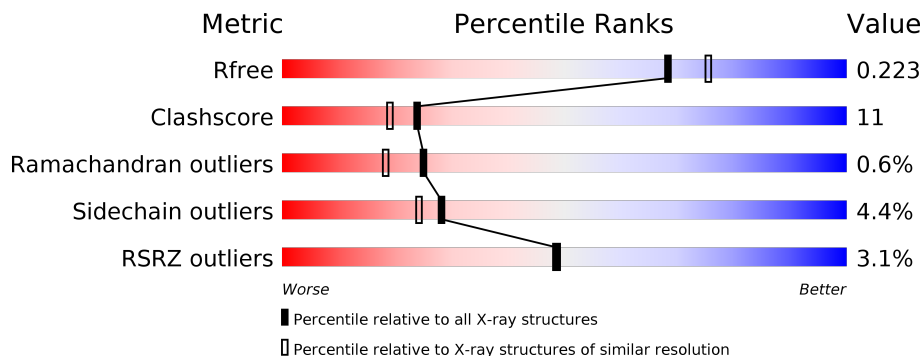
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.02 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	10434 (2.04-2.00)
Clashscore	141614	11643 (2.04-2.00)
Ramachandran outliers	138981	11493 (2.04-2.00)
Sidechain outliers	138945	11492 (2.04-2.00)
RSRZ outliers	127900	10220 (2.04-2.00)

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 on 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	470	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 75%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 13%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="font-size: small; margin-top: 5px;">3% 75% 13% • 10%</p>
1	B	470	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 65%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 21%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 11%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="font-size: small; margin-top: 5px;">3% 65% 21% • 11%</p>

2 Entry composition i

There are 7 unique types of molecules in this entry. The entry contains 7314 atoms, of which 48 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-pentakisphosphate 2-kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	424	3362	2139	567	642	14	0	0	0
1	B	416	3313	2108	559	632	14	0	0	0

There are 40 discrepancies between the modelled and reference sequences:

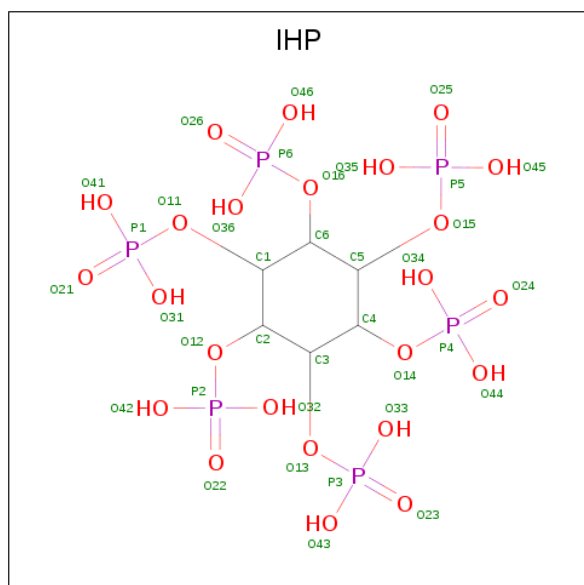
Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	MET	-	initiating methionine	UNP A0A178UAB5
A	-17	ALA	-	expression tag	UNP A0A178UAB5
A	-16	HIS	-	expression tag	UNP A0A178UAB5
A	-15	HIS	-	expression tag	UNP A0A178UAB5
A	-14	HIS	-	expression tag	UNP A0A178UAB5
A	-13	HIS	-	expression tag	UNP A0A178UAB5
A	-12	HIS	-	expression tag	UNP A0A178UAB5
A	-11	HIS	-	expression tag	UNP A0A178UAB5
A	-10	SER	-	expression tag	UNP A0A178UAB5
A	-9	SER	-	expression tag	UNP A0A178UAB5
A	-8	GLY	-	expression tag	UNP A0A178UAB5
A	-7	LEU	-	expression tag	UNP A0A178UAB5
A	-6	GLU	-	expression tag	UNP A0A178UAB5
A	-5	VAL	-	expression tag	UNP A0A178UAB5
A	-4	LEU	-	expression tag	UNP A0A178UAB5
A	-3	PHE	-	expression tag	UNP A0A178UAB5
A	-2	GLN	-	expression tag	UNP A0A178UAB5
A	-1	GLY	-	expression tag	UNP A0A178UAB5
A	0	PRO	-	expression tag	UNP A0A178UAB5
A	185	MET	ILE	conflict	UNP A0A178UAB5
B	-18	MET	-	initiating methionine	UNP A0A178UAB5
B	-17	ALA	-	expression tag	UNP A0A178UAB5
B	-16	HIS	-	expression tag	UNP A0A178UAB5
B	-15	HIS	-	expression tag	UNP A0A178UAB5
B	-14	HIS	-	expression tag	UNP A0A178UAB5

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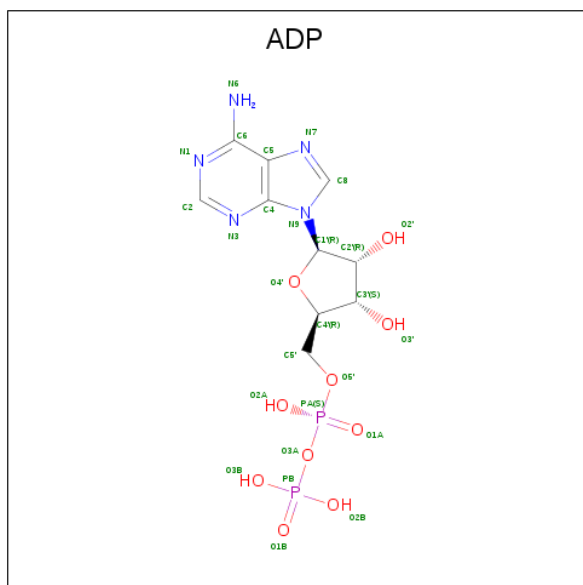
Chain	Residue	Modelled	Actual	Comment	Reference
B	-13	HIS	-	expression tag	UNP A0A178UAB5
B	-12	HIS	-	expression tag	UNP A0A178UAB5
B	-11	HIS	-	expression tag	UNP A0A178UAB5
B	-10	SER	-	expression tag	UNP A0A178UAB5
B	-9	SER	-	expression tag	UNP A0A178UAB5
B	-8	GLY	-	expression tag	UNP A0A178UAB5
B	-7	LEU	-	expression tag	UNP A0A178UAB5
B	-6	GLU	-	expression tag	UNP A0A178UAB5
B	-5	VAL	-	expression tag	UNP A0A178UAB5
B	-4	LEU	-	expression tag	UNP A0A178UAB5
B	-3	PHE	-	expression tag	UNP A0A178UAB5
B	-2	GLN	-	expression tag	UNP A0A178UAB5
B	-1	GLY	-	expression tag	UNP A0A178UAB5
B	0	PRO	-	expression tag	UNP A0A178UAB5
B	185	MET	ILE	conflict	UNP A0A178UAB5

- Molecule 2 is INOSITOL HEXAKISPHOSPHATE (three-letter code: IHP) (formula: $C_6H_{18}O_{24}P_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	A	1	Total	C	H	O	P	0	0
			42	6	6	24	6		
2	B	1	Total	C	H	O	P	0	0
			42	6	6	24	6		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	H	N	O			P
3	A	1	39	10	12	5	10	2	0	0
3	B	1	39	10	12	5	10	2	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
5	B	1	1	1	0	0
5	A	1	1	1	0	0

- Molecule 6 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: C₄H₁₂NO₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	H	N	O		
6	B	1	20	4	12	1	3	0	0

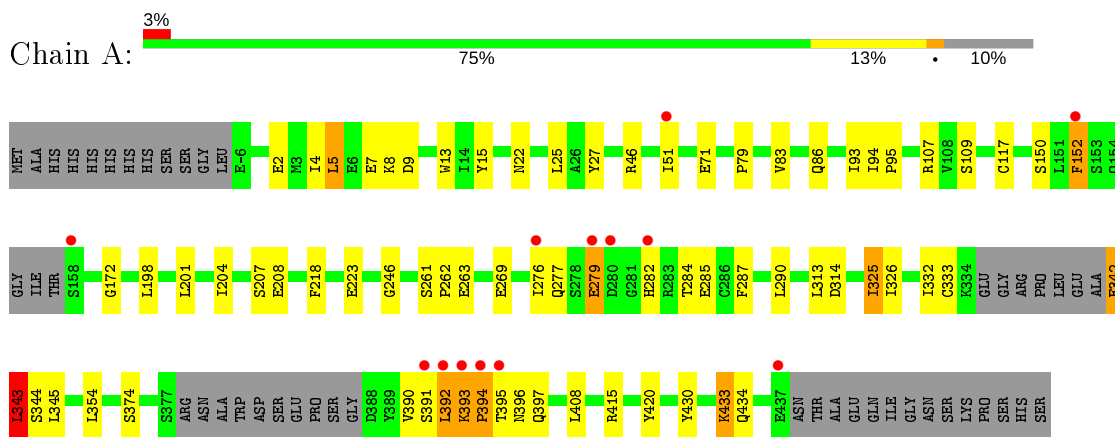
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	240	Total	O	0	0
			240	240		
7	B	212	Total	O	0	0
			212	212		

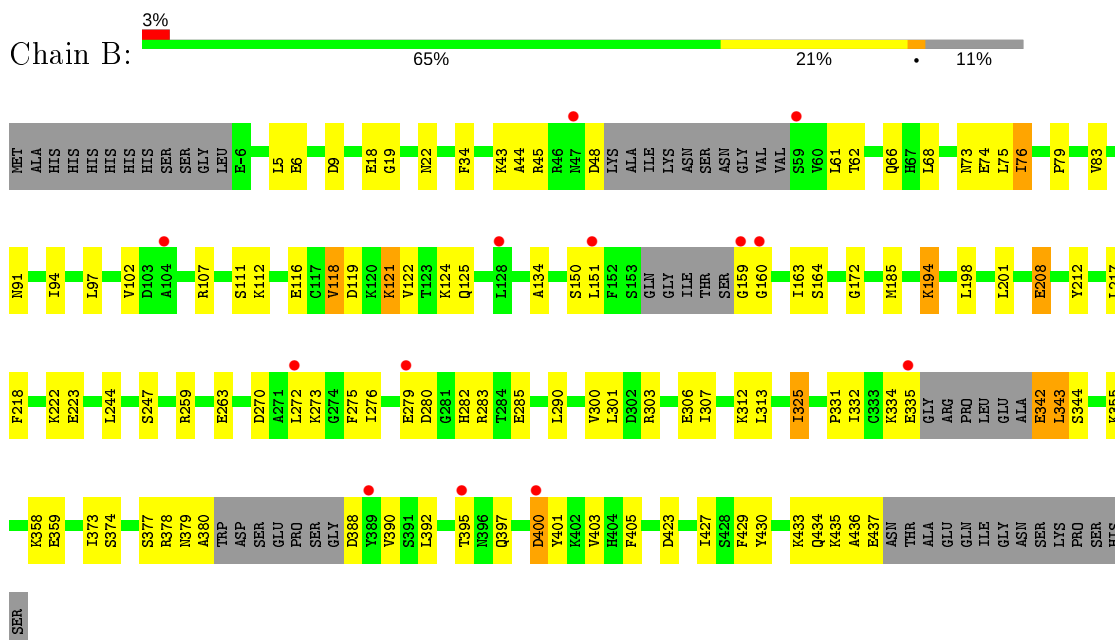
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Inositol-pentakisphosphate 2-kinase



- Molecule 1: Inositol-pentakisphosphate 2-kinase



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	59.81Å 60.34Å 83.98Å 88.53° 89.18° 63.36°	Depositor
Resolution (Å)	83.95 – 2.02 83.95 – 2.02	Depositor EDS
% Data completeness (in resolution range)	89.5 (83.95-2.02) 89.6 (83.95-2.02)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.03Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.178 , 0.225 0.179 , 0.223	Depositor DCC
R_{free} test set	3102 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.116	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 56.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.023 for -h,-k,l 0.015 for k,h,-l 0.022 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7314	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.16% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: ZN, MG, TRS, IHP, ADP

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	0.38	0/3422	0.51	0/4612
1	B	0.37	0/3372	0.51	0/4542
All	All	0.37	0/6794	0.51	0/9154

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	3362	0	3379	59	0
1	B	3313	0	3326	94	0
2	A	36	6	6	1	0
2	B	36	6	6	0	0
3	A	27	12	12	1	0
3	B	27	12	12	2	0
4	A	1	0	0	0	0
4	B	2	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	B	8	12	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	240	0	0	2	0
7	B	212	0	0	2	0
All	All	7266	48	6753	150	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 (150) 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:74:GLU:HG3	1:B:91:ASN:HD22	1.12	1.14
1:B:74:GLU:HG3	1:B:91:ASN:ND2	1.71	1.04
1:B:272:LEU:HD11	1:B:283:ARG:HB3	1.36	1.01
1:B:272:LEU:CD1	1:B:283:ARG:HB3	2.05	0.87
1:A:263:GLU:OE1	1:A:263:GLU:N	2.08	0.87
1:A:269:GLU:OE1	1:A:284:THR:HG23	1.77	0.83
1:B:272:LEU:HD11	1:B:283:ARG:CB	2.10	0.82
1:B:342:GLU:OE1	1:B:343:LEU:N	2.14	0.81
1:B:6:GLU:N	1:B:9:ASP:OD1	2.10	0.79
1:A:284:THR:HG21	7:A:601:HOH:O	1.82	0.78
1:B:74:GLU:CG	1:B:91:ASN:HD22	1.95	0.75
1:B:22:ASN:N	3:B:502:ADP:O3B	2.20	0.74
1:B:208:GLU:HG3	1:B:259:ARG:HH11	1.53	0.73
1:B:74:GLU:N	1:B:74:GLU:OE1	2.20	0.72
1:A:342:GLU:O	1:A:343:LEU:HB2	1.89	0.72
1:B:73:ASN:ND2	7:B:601:HOH:O	2.22	0.71
1:A:325:ILE:HG22	1:A:326:ILE:HG23	1.72	0.71
1:B:343:LEU:O	1:B:343:LEU:HD23	1.91	0.70
1:A:325:ILE:HG13	1:A:354:LEU:HD23	1.72	0.70
1:B:222:LYS:NZ	1:B:223:GLU:OE2	2.17	0.70
1:B:83:VAL:HG22	1:B:107:ARG:HD3	1.75	0.69
1:B:279:GLU:HG2	1:B:280:ASP:N	2.08	0.69
1:B:331:PRO:O	1:B:334:LYS:HE3	1.92	0.68
1:A:198:LEU:HD21	1:A:433:LYS:HG3	1.74	0.68
1:B:19:GLY:HA3	3:B:502:ADP:O2B	1.94	0.67
1:A:282:HIS:ND1	1:A:285:GLU:OE2	2.27	0.67
1:B:198:LEU:HD13	1:B:433:LYS:HD3	1.77	0.66
1:B:150:SER:O	1:B:151:LEU:HD23	1.95	0.66
1:B:208:GLU:HG3	1:B:259:ARG:NH1	2.09	0.66
1:B:434:GLN:O	1:B:437:GLU:HG2	1.96	0.66
1:A:223:GLU:OE1	1:A:223:GLU:N	2.27	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:121:LYS:HG3	1:B:122:VAL:HG13	1.79	0.65
1:B:379:ASN:OD1	1:B:380:ALA:N	2.29	0.64
1:B:121:LYS:HG3	1:B:122:VAL:N	2.11	0.64
1:A:430:TYR:O	1:A:434:GLN:HG2	1.97	0.64
1:B:342:GLU:OE1	1:B:344:SER:N	2.27	0.63
1:A:392:LEU:HD11	1:A:395:THR:OG1	2.00	0.62
1:B:300:VAL:HG13	1:B:301:LEU:HD12	1.82	0.62
1:B:217:LEU:HG	1:B:301:LEU:HD21	1.81	0.62
1:A:22:ASN:N	3:A:502:ADP:O2B	2.31	0.62
1:B:163:ILE:HD12	1:B:275:PHE:CG	2.35	0.62
1:A:392:LEU:HD21	1:A:397:GLN:HB2	1.81	0.61
1:A:277:GLN:O	1:A:277:GLN:HG3	1.98	0.61
1:B:61:LEU:HD12	1:B:76:ILE:HG22	1.81	0.61
1:A:392:LEU:H	1:A:392:LEU:HD23	1.65	0.61
1:B:217:LEU:HG	1:B:301:LEU:CD2	2.31	0.61
1:A:204:ILE:CD1	1:B:313:LEU:HD11	2.31	0.60
1:B:332:ILE:O	1:B:334:LYS:HG2	2.02	0.59
1:B:429:PHE:CE2	1:B:433:LYS:HD2	2.38	0.58
1:B:379:ASN:O	1:B:380:ALA:HB2	2.02	0.58
1:B:172:GLY:HA3	1:B:218:PHE:CD2	2.40	0.56
1:A:343:LEU:O	1:A:345:LEU:N	2.39	0.56
1:B:62:THR:O	1:B:66:GLN:HG3	2.06	0.56
1:A:392:LEU:H	1:A:392:LEU:CD2	2.19	0.55
1:A:392:LEU:CD2	1:A:397:GLN:H	2.20	0.55
1:B:5:LEU:HD21	1:B:34:PHE:CZ	2.42	0.55
1:A:325:ILE:HG13	1:A:354:LEU:CD2	2.35	0.55
1:A:79:PRO:HD2	1:A:83:VAL:HG11	1.88	0.55
1:B:263:GLU:N	1:B:263:GLU:OE1	2.34	0.54
1:B:272:LEU:HD12	1:B:272:LEU:O	2.07	0.54
1:B:342:GLU:O	1:B:343:LEU:HB2	2.08	0.54
1:B:163:ILE:HD12	1:B:275:PHE:CD2	2.43	0.54
1:B:270:ASP:O	1:B:273:LYS:HG2	2.09	0.53
1:B:272:LEU:HD13	1:B:276:ILE:HG13	1.91	0.53
1:B:83:VAL:CG2	1:B:107:ARG:HD3	2.40	0.52
1:B:312:LYS:NZ	1:B:359:GLU:OE1	2.34	0.52
1:A:8:LYS:NZ	1:A:9:ASP:OD1	2.43	0.52
1:B:94:ILE:HD11	1:B:102:VAL:CG2	2.39	0.52
1:A:333:CYS:SG	1:A:342:GLU:HB3	2.50	0.51
1:B:83:VAL:HG22	1:B:107:ARG:CD	2.39	0.51
1:B:303:ARG:HH11	1:B:303:ARG:HA	1.75	0.51
1:A:392:LEU:CD2	1:A:397:GLN:HB2	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:272:LEU:HD13	1:B:276:ILE:CG1	2.41	0.51
1:B:279:GLU:HG2	1:B:280:ASP:H	1.74	0.51
1:B:335:GLU:N	1:B:335:GLU:OE1	2.34	0.51
1:A:93:ILE:HG21	1:A:408:LEU:HD13	1.92	0.50
1:B:121:LYS:HG3	1:B:122:VAL:CG1	2.41	0.50
1:A:392:LEU:HB2	1:A:394:PRO:HD2	1.92	0.49
1:B:430:TYR:O	1:B:434:GLN:HG2	2.12	0.49
1:B:194:LYS:HE2	1:B:212:TYR:O	2.12	0.49
1:B:395:THR:OG1	1:B:397:GLN:HG2	2.12	0.49
1:A:313:LEU:O	1:A:314:ASP:HB3	2.13	0.49
1:A:204:ILE:HD12	1:B:313:LEU:HD11	1.95	0.49
1:B:94:ILE:CG1	1:B:102:VAL:HG21	2.43	0.48
1:A:342:GLU:HG2	1:A:343:LEU:O	2.14	0.48
1:B:164:SER:HA	1:B:373:ILE:O	2.14	0.47
1:B:44:ALA:O	1:B:134:ALA:HB1	2.15	0.47
1:A:325:ILE:HD11	1:A:354:LEU:N	2.30	0.47
1:A:325:ILE:CG1	1:A:354:LEU:HD23	2.44	0.47
1:A:332:ILE:HD12	1:A:420:TYR:CZ	2.49	0.47
1:B:97:LEU:HD13	1:B:307:ILE:HD12	1.96	0.47
1:A:290:LEU:HD12	1:A:390:VAL:HG11	1.96	0.47
1:B:159:GLY:HA2	1:B:160:GLY:HA3	1.62	0.47
1:A:392:LEU:HD23	1:A:397:GLN:H	1.78	0.47
1:B:379:ASN:CG	1:B:380:ALA:H	2.17	0.47
2:A:501:IHP:O32	2:A:501:IHP:O23	2.33	0.47
1:B:395:THR:O	1:B:397:GLN:N	2.43	0.47
1:A:172:GLY:HA3	1:A:218:PHE:CD2	2.50	0.46
1:B:121:LYS:HE3	1:B:121:LYS:HB2	1.59	0.46
1:B:121:LYS:HG3	1:B:122:VAL:H	1.80	0.46
1:B:18:GLU:OE2	1:B:43:LYS:NZ	2.48	0.46
1:A:204:ILE:HD11	1:B:313:LEU:HD11	1.96	0.46
1:A:415:ARG:HD3	7:A:637:HOH:O	2.15	0.46
1:B:282:HIS:O	1:B:285:GLU:HG2	2.16	0.46
1:B:379:ASN:O	1:B:380:ALA:CB	2.65	0.45
1:B:388:ASP:N	7:B:615:HOH:O	2.50	0.45
1:B:435:LYS:O	1:B:437:GLU:N	2.50	0.45
1:B:102:VAL:O	1:B:102:VAL:CG2	2.65	0.45
1:B:124:LYS:HE3	1:B:124:LYS:HB2	1.80	0.45
1:B:97:LEU:HD13	1:B:307:ILE:CD1	2.47	0.44
1:B:325:ILE:HA	1:B:325:ILE:HD12	1.74	0.44
1:B:79:PRO:HD2	1:B:83:VAL:HG11	1.98	0.44
1:A:279:GLU:HG3	1:A:279:GLU:H	1.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:290:LEU:CD1	1:B:390:VAL:HG21	2.47	0.44
1:A:325:ILE:HD12	1:A:325:ILE:HA	1.72	0.44
1:A:198:LEU:CD2	1:A:433:LYS:HG3	2.44	0.44
1:B:74:GLU:HG2	1:B:75:LEU:N	2.33	0.44
1:B:185:MET:CE	1:B:185:MET:HA	2.48	0.44
1:B:94:ILE:HD11	1:B:102:VAL:HG22	2.00	0.44
1:B:112:LYS:O	1:B:116:GLU:HG3	2.17	0.44
1:B:423:ASP:O	1:B:427:ILE:HG22	2.18	0.44
1:A:15:TYR:HA	1:A:25:LEU:CD2	2.48	0.43
1:A:392:LEU:N	1:A:392:LEU:HD23	2.33	0.43
1:B:392:LEU:O	1:B:395:THR:O	2.36	0.43
1:A:4:ILE:HG12	1:A:109:SER:HB2	2.01	0.43
1:A:86:GLN:OE1	1:A:107:ARG:HG3	2.19	0.43
1:B:118:VAL:HG12	1:B:119:ASP:N	2.32	0.43
1:B:68:LEU:HD11	1:B:358:LYS:HG3	2.00	0.43
1:A:393:LYS:N	1:A:394:PRO:CD	2.82	0.43
1:A:5:LEU:HA	1:A:5:LEU:HD12	1.85	0.43
1:B:355:LYS:HE2	1:B:359:GLU:OE2	2.19	0.42
1:A:276:ILE:CD1	1:A:287:PHE:HB2	2.49	0.42
1:A:46:ARG:HB2	1:B:185:MET:HG2	2.01	0.42
1:A:7:GLU:HG3	1:A:117:CYS:SG	2.59	0.42
1:B:102:VAL:O	1:B:102:VAL:HG22	2.19	0.42
1:A:277:GLN:HG2	1:A:397:GLN:OE1	2.19	0.42
1:B:374:SER:O	1:B:401:TYR:HA	2.20	0.42
1:B:83:VAL:CG2	1:B:107:ARG:CD	2.97	0.41
1:A:261:SER:HB2	1:A:262:PRO:HD2	2.03	0.41
1:A:51:ILE:HD12	1:A:51:ILE:N	2.35	0.41
1:B:378:ARG:HA	1:B:400:ASP:OD2	2.20	0.41
1:A:152:PHE:O	1:A:152:PHE:HD2	2.04	0.41
1:A:277:GLN:NE2	1:A:395:THR:O	2.53	0.41
1:A:51:ILE:H	1:A:51:ILE:HD12	1.85	0.41
1:A:13:TRP:CZ3	1:A:27:TYR:HB2	2.55	0.41
1:A:325:ILE:HG22	1:A:326:ILE:N	2.36	0.41
1:B:300:VAL:HG22	1:B:405:PHE:CZ	2.56	0.41
1:B:244:LEU:O	1:B:247:SER:OG	2.36	0.40
1:A:150:SER:OG	1:A:246:GLY:HA2	2.21	0.40
1:A:94:ILE:N	1:A:95:PRO:CD	2.84	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	416/470 (88%)	406 (98%)	7 (2%)	3 (1%)	22	15
1	B	406/470 (86%)	395 (97%)	9 (2%)	2 (0%)	29	22
All	All	822/940 (87%)	801 (97%)	16 (2%)	5 (1%)	25	18

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	344	SER
1	B	343	LEU
1	A	343	LEU
1	B	436	ALA
1	A	394	PRO

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	377/415 (91%)	360 (96%)	17 (4%)	27	23
1	B	371/415 (89%)	355 (96%)	16 (4%)	29	25
All	All	748/830 (90%)	715 (96%)	33 (4%)	28	24

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLU

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Mol	Chain	Res	Type
1	A	5	LEU
1	A	71	GLU
1	A	152	PHE
1	A	201	LEU
1	A	207	SER
1	A	208	GLU
1	A	279	GLU
1	A	325	ILE
1	A	342	GLU
1	A	343	LEU
1	A	374	SER
1	A	391	SER
1	A	392	LEU
1	A	393	LYS
1	A	396	ASN
1	A	433	LYS
1	B	45	ARG
1	B	48	ASP
1	B	76	ILE
1	B	111	SER
1	B	118	VAL
1	B	121	LYS
1	B	125	GLN
1	B	194	LYS
1	B	201	LEU
1	B	208	GLU
1	B	306	GLU
1	B	325	ILE
1	B	342	GLU
1	B	377	SER
1	B	400	ASP
1	B	403	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	91	ASN

5.3.3 RNA

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 10 ligands modelled in this entry, 5 are 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
2	IHP	A	501	4	36,36,36	1.00	2 (5%)	54,60,60	0.91	0
3	ADP	A	502	4	24,29,29	1.01	2 (8%)	29,45,45	1.45	3 (10%)
3	ADP	B	502	4	24,29,29	0.92	1 (4%)	29,45,45	1.42	5 (17%)
2	IHP	B	501	4	36,36,36	0.98	2 (5%)	54,60,60	0.99	3 (5%)
6	TRS	B	506	-	7,7,7	0.26	0	9,9,9	0.48	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	IHP	A	501	4	-	2/30/54/54	0/1/1/1
3	ADP	A	502	4	-	1/12/32/32	0/3/3/3
3	ADP	B	502	4	-	3/12/32/32	0/3/3/3
2	IHP	B	501	4	-	3/30/54/54	0/1/1/1
6	TRS	B	506	-	-	0/9/9/9	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	IHP	P3-O13	2.81	1.64	1.59
2	A	501	IHP	P3-O13	2.79	1.64	1.59
3	A	502	ADP	C5-C4	2.45	1.47	1.40
3	B	502	ADP	C5-C4	2.34	1.47	1.40
2	B	501	IHP	P2-O12	2.18	1.63	1.59
2	A	501	IHP	P2-O12	2.15	1.63	1.59
3	A	502	ADP	C2-N3	2.13	1.35	1.32

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	ADP	N3-C2-N1	-3.81	122.73	128.68
3	B	502	ADP	N3-C2-N1	-3.53	123.16	128.68
3	B	502	ADP	PA-O3A-PB	-3.45	121.00	132.83
3	A	502	ADP	PA-O3A-PB	-2.99	122.58	132.83
3	A	502	ADP	C3'-C2'-C1'	2.73	105.09	100.98
3	B	502	ADP	C3'-C2'-C1'	2.60	104.89	100.98
2	B	501	IHP	O44-P4-O34	2.28	116.34	107.64
2	B	501	IHP	O16-P6-O26	-2.05	101.50	109.39
3	B	502	ADP	N6-C6-N1	2.04	122.82	118.57
2	B	501	IHP	O12-P2-O22	-2.03	101.56	109.39
3	B	502	ADP	C4-C5-N7	-2.00	107.31	109.40

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	501	IHP	C2-O12-P2-O22
3	B	502	ADP	C5'-O5'-PA-O3A
2	A	501	IHP	C2-O12-P2-O22
2	A	501	IHP	C2-O12-P2-O32
3	B	502	ADP	C5'-O5'-PA-O1A
2	B	501	IHP	C1-O11-P1-O21
2	B	501	IHP	C2-O12-P2-O32
3	A	502	ADP	C5'-O5'-PA-O1A
3	B	502	ADP	C5'-O5'-PA-O2A

There are no ring outliers.

3 monomers are involved in 4 short contacts:

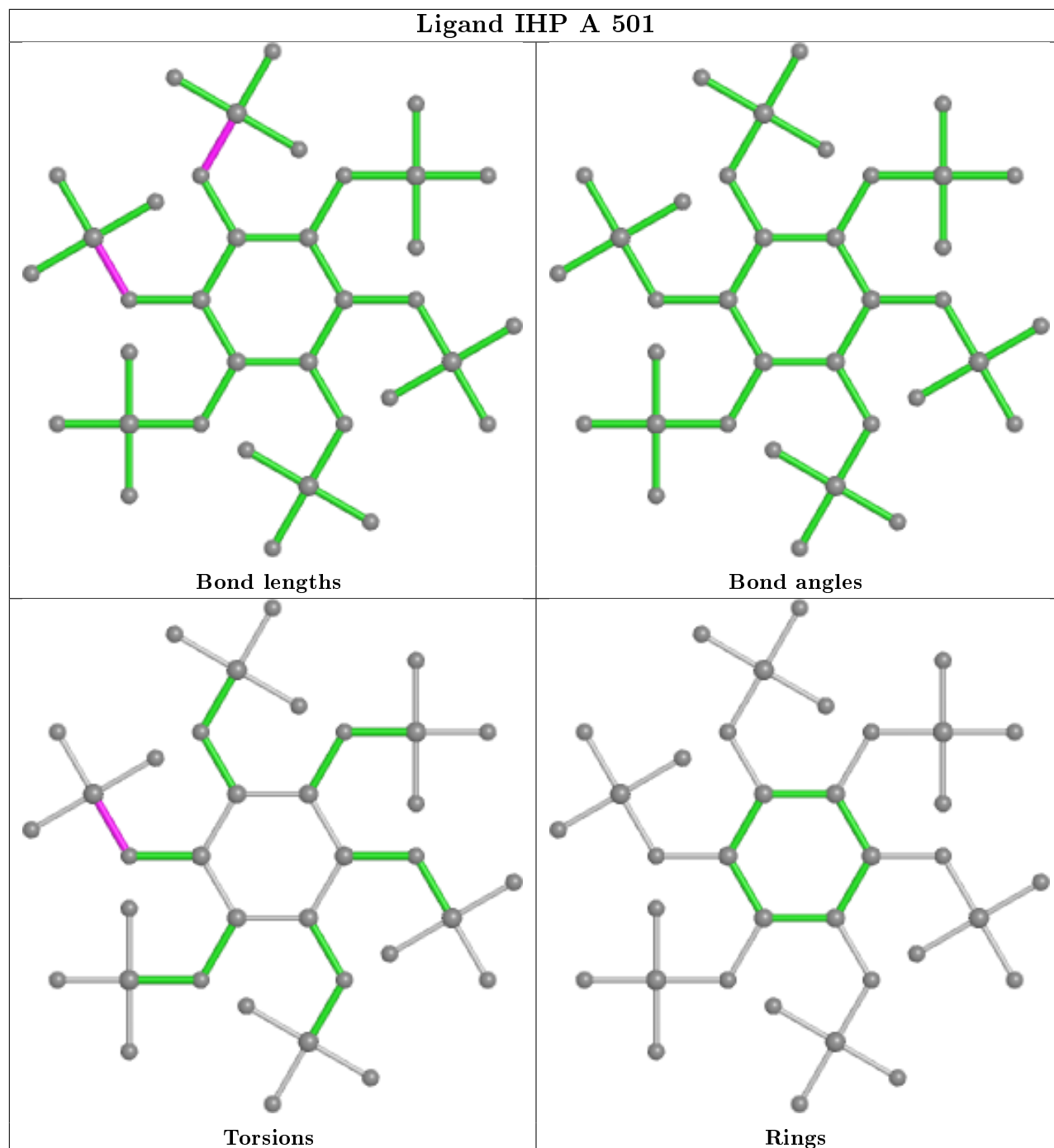
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	IHP	1	0

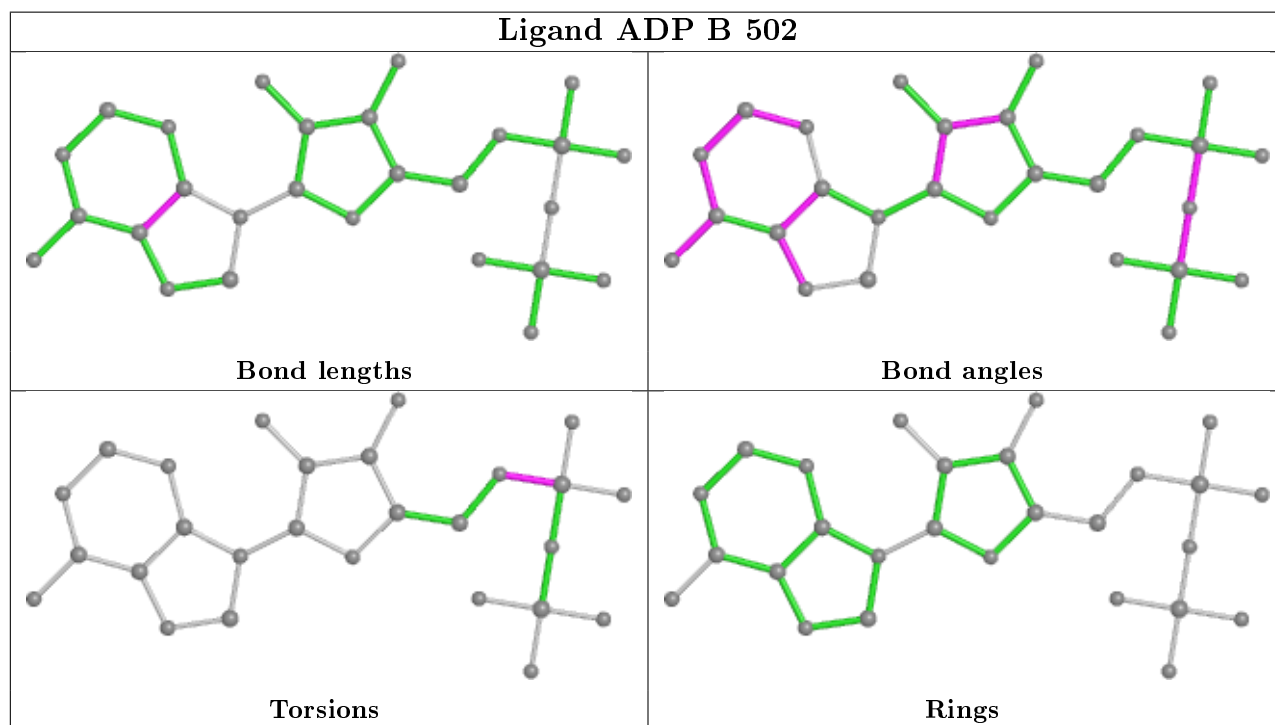
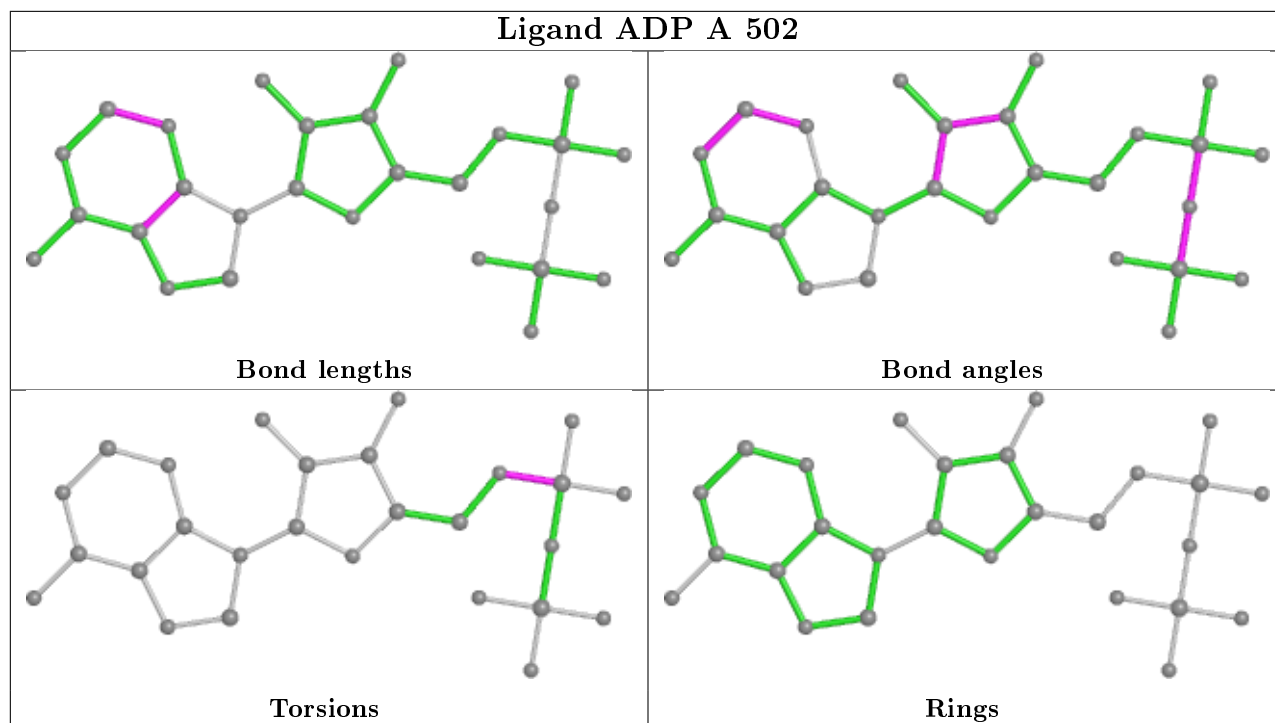
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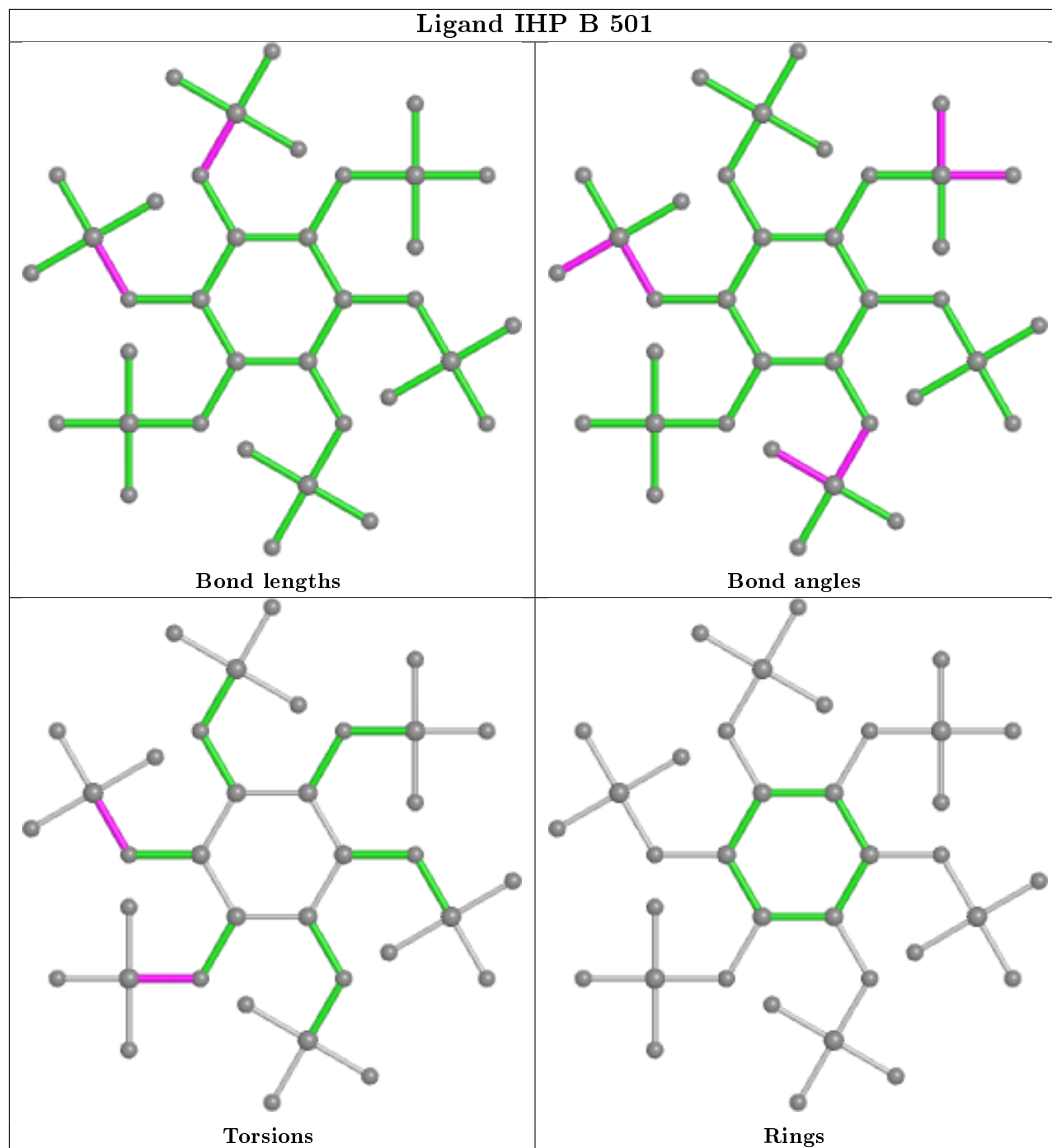
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	ADP	1	0
3	B	502	ADP	2	0

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

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	424/470 (90%)	0.02	13 (3%) 49 48	17, 35, 76, 114	0
1	B	416/470 (88%)	0.05	13 (3%) 49 48	18, 39, 73, 109	0
All	All	840/940 (89%)	0.04	26 (3%) 49 48	17, 37, 75, 114	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	158	SER	7.5
1	B	335	GLU	6.3
1	B	104	ALA	5.1
1	B	47	ASN	4.7
1	A	394	PRO	4.2
1	A	279	GLU	4.0
1	A	152	PHE	3.7
1	B	279	GLU	3.6
1	A	280	ASP	3.1
1	B	59	SER	2.8
1	A	395	THR	2.7
1	B	395	THR	2.6
1	A	392	LEU	2.6
1	A	393	LYS	2.6
1	B	272	LEU	2.6
1	B	151	LEU	2.5
1	A	391	SER	2.5
1	A	282	HIS	2.4
1	B	389	TYR	2.3
1	B	128	LEU	2.3
1	A	437	GLU	2.3
1	B	160	GLY	2.2
1	B	159	GLY	2.2
1	A	276	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	400	ASP	2.1
1	A	51	ILE	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 monosaccharides in this entry.

6.4 Ligands [i](#)

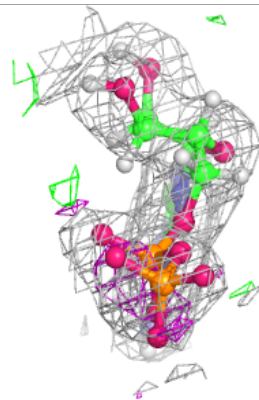
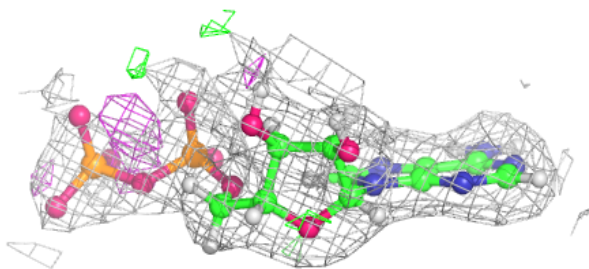
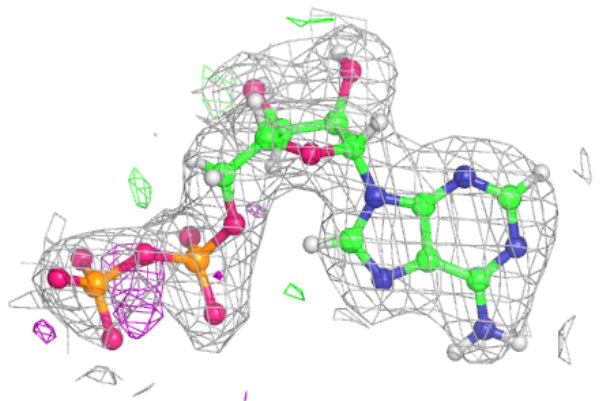
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MG	A	503	1/1	0.72	0.12	58,58,58,58	0
6	TRS	B	506	8/8	0.89	0.13	34,45,63,76	0
3	ADP	A	502	27/27	0.90	0.14	25,32,79,87	0
3	ADP	B	502	27/27	0.92	0.17	29,39,97,103	0
4	MG	B	504	1/1	0.93	0.11	46,46,46,46	0
4	MG	B	503	1/1	0.96	0.10	48,48,48,48	0
2	IHP	B	501	36/36	0.97	0.13	13,23,57,69	0
2	IHP	A	501	36/36	0.98	0.12	12,23,42,63	0
5	ZN	A	504	1/1	0.99	0.12	32,32,32,32	0
5	ZN	B	505	1/1	1.00	0.12	25,25,25,25	0

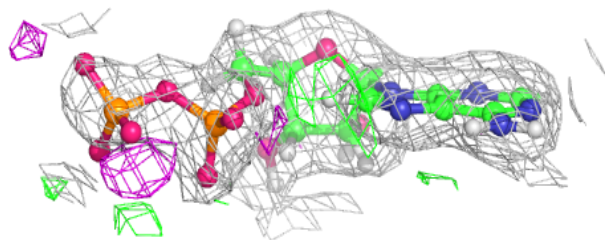
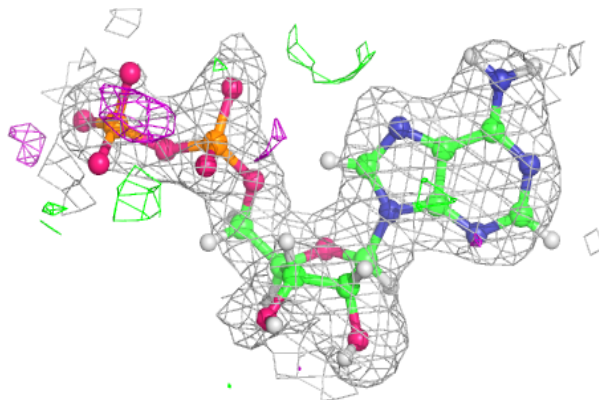
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ADP A 502:

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

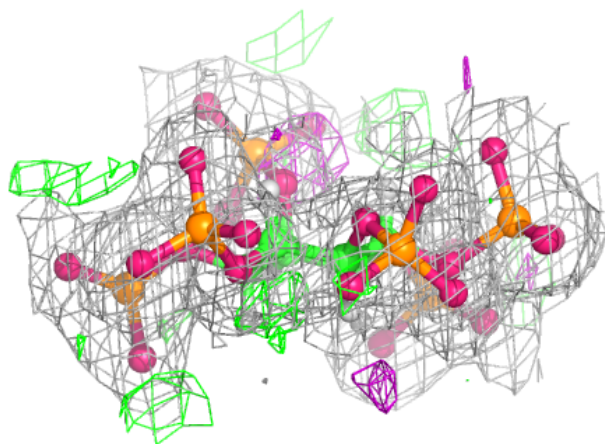
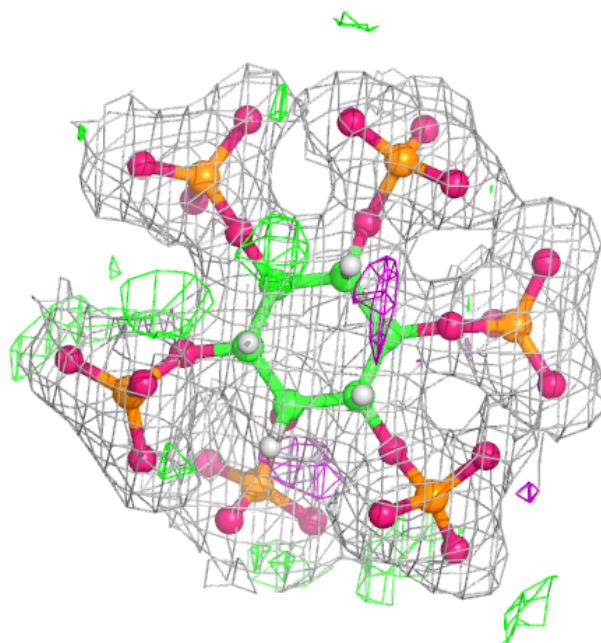
**Electron density around ADP B 502:**

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



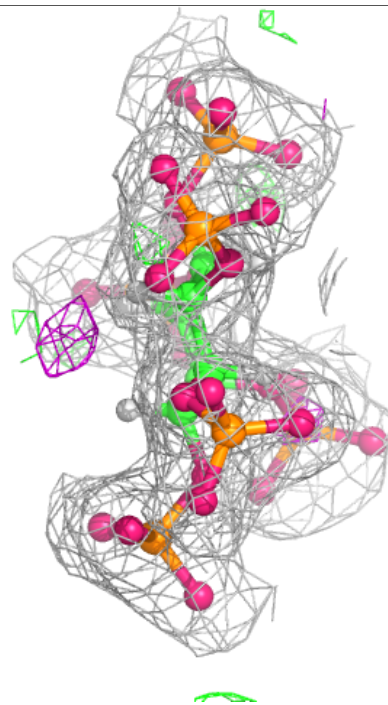
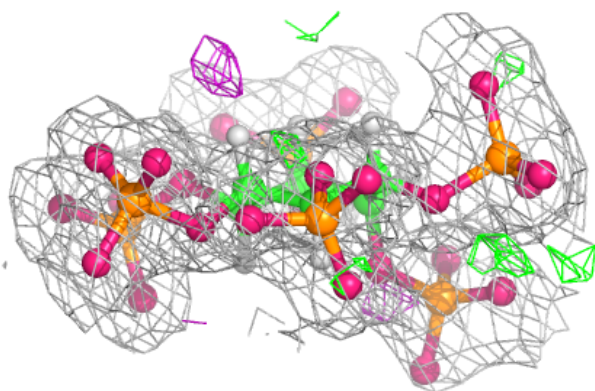
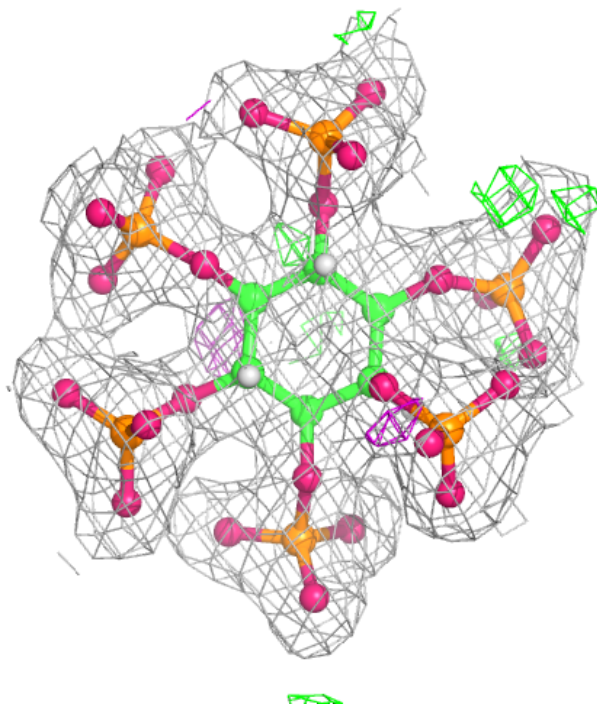
Electron density around IHP B 501:

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



Electron density around IHP 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.