Full wwPDB X-ray Structure Validation Report ⓘ

Sep 28, 2020 – 03:11 PM BST

PDB ID	:	5OQU
Title	:	The crystal structure of CK2alpha in complex with compound 5
Authors	:	Brear, P.; De Fusco, C.; Iegre, J.; Yoshida, M.; Mitchell, S.; Rossmann, M.; Carro, L.; Sore, H.; Hyvonen, M.; Spring, D.
Deposited on	:	2017-08-14
Resolution	:	2.32 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:

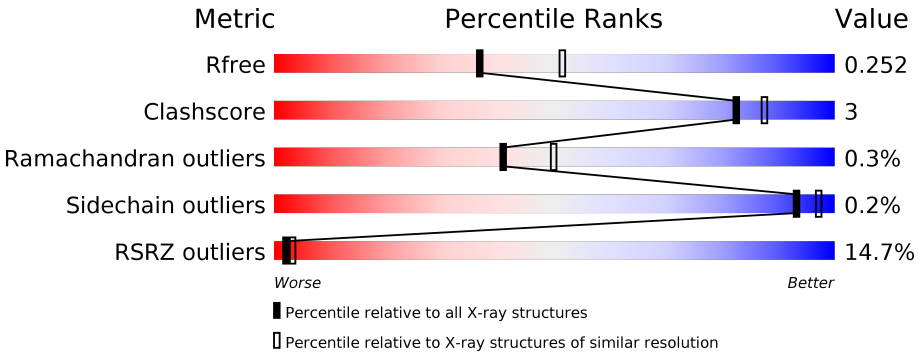
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.14.6
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & 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 2.32 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	5974 (2.34-2.30)
Clashscore	141614	6604 (2.34-2.30)
Ramachandran outliers	138981	6523 (2.34-2.30)
Sidechain outliers	138945	6523 (2.34-2.30)
RSRZ outliers	127900	5855 (2.34-2.30)

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 <=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	352	6% 87% 5% 8%
1	B	352	21% 85% 7% 8%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 5671 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 Casein kinase II subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	324	Total	C	N	O	S	0	3	0
			2763	1770	485	497	11			
1	A	324	Total	C	N	O	S	0	3	0
			2763	1770	485	497	11			

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-22	GLY	-	expression tag	UNP P68400
B	-21	SER	-	expression tag	UNP P68400
B	-20	MET	-	expression tag	UNP P68400
B	-19	ASP	-	expression tag	UNP P68400
B	-18	ILE	-	expression tag	UNP P68400
B	-17	GLU	-	expression tag	UNP P68400
B	-16	PHE	-	expression tag	UNP P68400
B	-15	ASP	-	expression tag	UNP P68400
B	-14	ASP	-	expression tag	UNP P68400
B	-13	ASP	-	expression tag	UNP P68400
B	-12	ALA	-	expression tag	UNP P68400
B	-11	ASP	-	expression tag	UNP P68400
B	-10	ASP	-	expression tag	UNP P68400
B	-9	ASP	-	expression tag	UNP P68400
B	-8	GLY	-	expression tag	UNP P68400
B	-7	SER	-	expression tag	UNP P68400
B	-6	GLY	-	expression tag	UNP P68400
B	-5	SER	-	expression tag	UNP P68400
B	-4	GLY	-	expression tag	UNP P68400
B	-3	SER	-	expression tag	UNP P68400
B	-2	GLY	-	expression tag	UNP P68400
B	-1	SER	-	expression tag	UNP P68400
B	0	GLY	-	expression tag	UNP P68400
B	1	SER	-	expression tag	UNP P68400
B	21	SER	ARG	engineered mutation	UNP P68400

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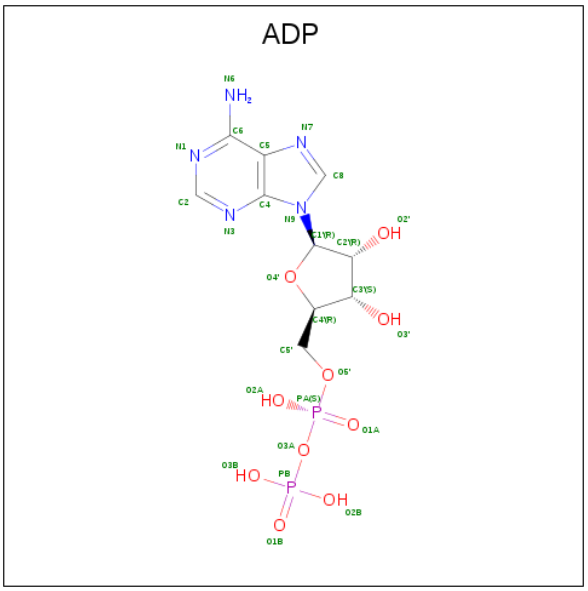
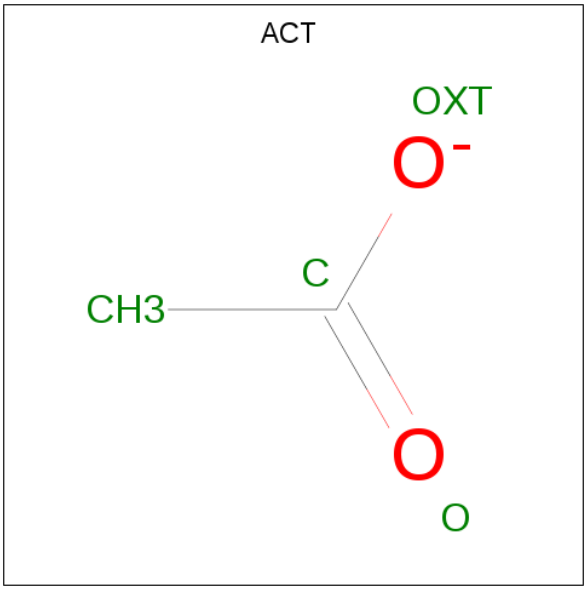
Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	GLY	-	expression tag	UNP P68400
A	-21	SER	-	expression tag	UNP P68400
A	-20	MET	-	expression tag	UNP P68400
A	-19	ASP	-	expression tag	UNP P68400
A	-18	ILE	-	expression tag	UNP P68400
A	-17	GLU	-	expression tag	UNP P68400
A	-16	PHE	-	expression tag	UNP P68400
A	-15	ASP	-	expression tag	UNP P68400
A	-14	ASP	-	expression tag	UNP P68400
A	-13	ASP	-	expression tag	UNP P68400
A	-12	ALA	-	expression tag	UNP P68400
A	-11	ASP	-	expression tag	UNP P68400
A	-10	ASP	-	expression tag	UNP P68400
A	-9	ASP	-	expression tag	UNP P68400
A	-8	GLY	-	expression tag	UNP P68400
A	-7	SER	-	expression tag	UNP P68400
A	-6	GLY	-	expression tag	UNP P68400
A	-5	SER	-	expression tag	UNP P68400
A	-4	GLY	-	expression tag	UNP P68400
A	-3	SER	-	expression tag	UNP P68400
A	-2	GLY	-	expression tag	UNP P68400
A	-1	SER	-	expression tag	UNP P68400
A	0	GLY	-	expression tag	UNP P68400
A	1	SER	-	expression tag	UNP P68400
A	21	SER	ARG	engineered mutation	UNP P68400

- 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	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

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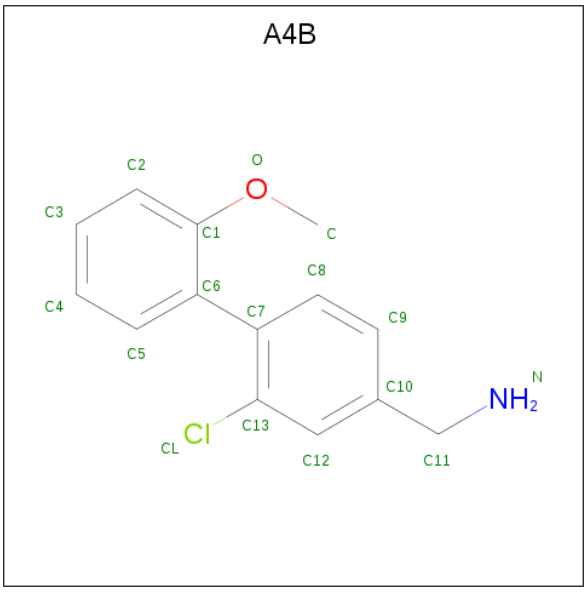
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0

- Molecule 4 is [3-chloranyl-4-(2-methoxyphenyl)phenyl]methanamine (three-letter code: A4B) (formula: $C_{14}H_{14}ClNO$).

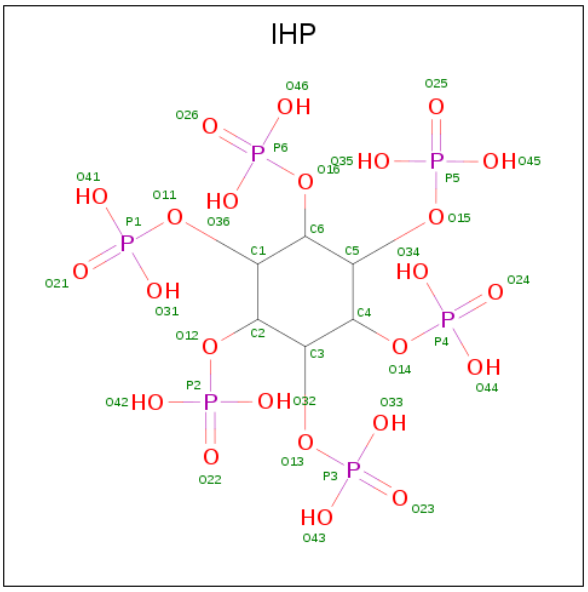


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C Cl N O 17 14 1 1 1	0	0

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	2	Total Mg 2 2	0	0

- Molecule 6 is INOSITOL HEXAKISPHOSPHATE (three-letter code: IHP) (formula: $C_6H_{18}O_{24}P_6$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O P 36 6 24 6	0	0

- Molecule 7 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total Cl 1 1	0	0

- Molecule 8 is water.

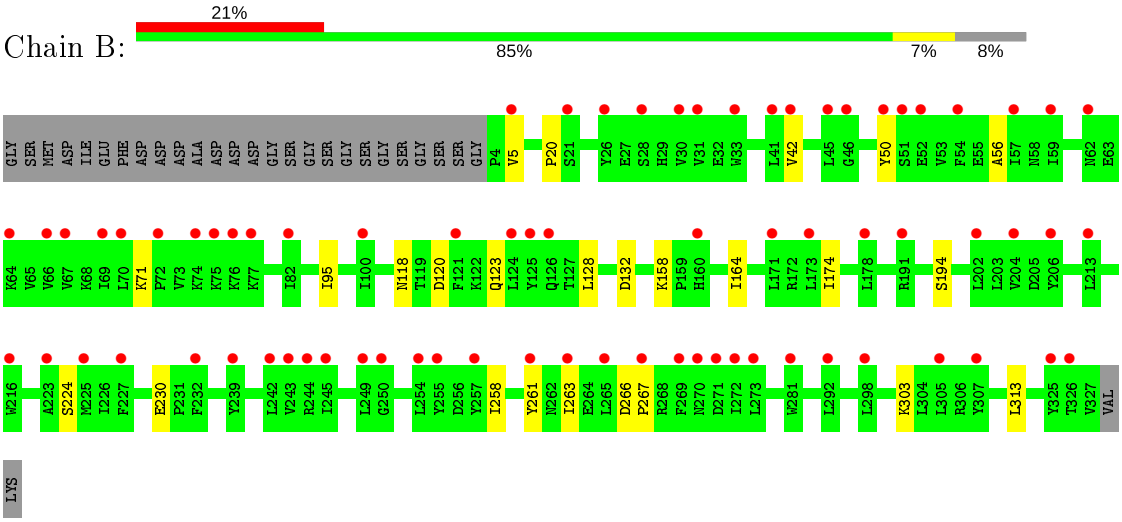
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	B	2	Total O 2 2	0	0
8	A	13	Total O 13 13	0	0

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.

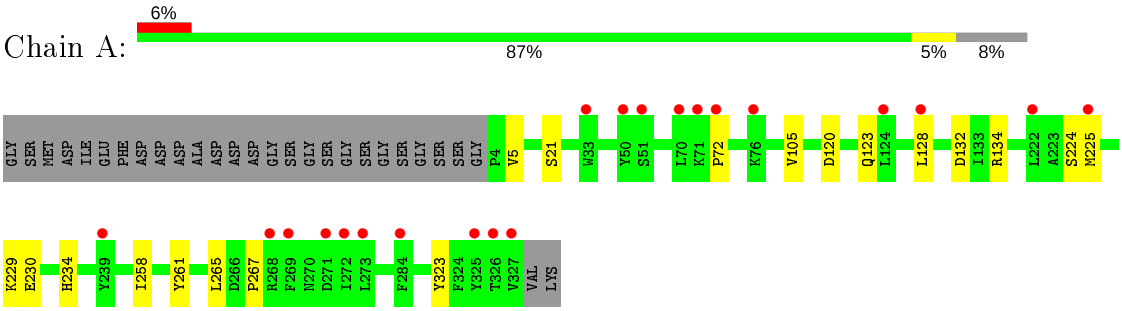
- Molecule 1: Casein kinase II subunit alpha

Chain B:



- Molecule 1: Casein kinase II subunit alpha

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	64.67Å 67.92Å 332.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	166.40 – 2.32 166.40 – 2.32	Depositor EDS
% Data completeness (in resolution range)	99.4 (166.40-2.32) 99.8 (166.40-2.32)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.50 (at 2.32Å)	Xtriage
Refinement program	BUSTER 2.10.1	Depositor
R, R_{free}	0.220 , 0.244 0.228 , 0.252	Depositor DCC
R_{free} test set	1601 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å^2)	54.1	Xtriage
Anisotropy	0.663	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 55.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.135 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5671	wwPDB-VP
Average B, all atoms (Å^2)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.28% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, A4B, ADP, CL, IHP, ACT

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	0.56	0/2838	0.65	0/3839
1	B	0.52	0/2838	0.65	0/3839
All	All	0.54	0/5676	0.65	0/7678

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

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	2763	0	2702	13	3
1	B	2763	0	2702	17	0
2	A	27	0	12	0	0
2	B	27	0	12	0	0
3	A	12	0	9	0	0
3	B	8	0	6	0	0
4	A	17	0	0	4	0
5	A	2	0	0	0	0
6	A	36	0	6	0	3
7	A	1	0	0	1	0
8	A	13	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	B	2	0	0	0	0
All	All	5671	0	5449	28	3

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 3.

All (28) 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:258:ILE:HG23	1:B:263:ILE:HB	1.66	0.77
1:B:224:SER:HB2	1:B:230:GLU:HG2	1.73	0.69
1:A:224:SER:HB2	1:A:230:GLU:HG2	1.74	0.68
4:A:402:A4B:N	7:A:409:CL:CL	2.64	0.66
1:B:266:ASP:OD1	1:B:267:PRO:HD2	2.03	0.59
1:B:50:TYR:HB3	1:B:71:LYS:HB3	1.86	0.57
1:B:50:TYR:CD1	1:B:50:TYR:N	2.72	0.55
1:B:158:LYS:HE2	1:B:194[A]:SER:OG	2.07	0.54
1:A:225:MET:CG	4:A:402:A4B:C2	2.87	0.53
1:A:5:VAL:HB	1:A:261:TYR:HA	1.92	0.52
1:B:20:PRO:HD2	1:A:105:VAL:HG21	1.91	0.51
1:B:20:PRO:CG	1:A:105:VAL:HG11	2.41	0.51
1:B:303:LYS:HB3	1:B:313:LEU:HG	1.94	0.50
1:A:225:MET:HG2	4:A:402:A4B:C2	2.41	0.50
1:A:120:ASP:HB2	1:A:123:GLN:HB2	1.96	0.48
1:A:134:ARG:HG2	1:A:323:TYR:CZ	2.48	0.47
1:B:120:ASP:HB3	1:B:123:GLN:HB2	1.96	0.47
1:B:5:VAL:HB	1:B:261:TYR:HA	1.98	0.45
1:B:20:PRO:HG2	1:A:105:VAL:HG11	1.97	0.45
1:A:224:SER:CB	1:A:230:GLU:HG2	2.45	0.45
1:A:225:MET:HG3	4:A:402:A4B:C2	2.47	0.44
1:B:224:SER:CB	1:B:230:GLU:HG2	2.44	0.42
1:A:258:ILE:HD13	1:A:265:LEU:HG	2.02	0.42
1:B:95:ILE:HB	1:B:174:ILE:HG22	2.02	0.41
1:B:128:LEU:HD22	1:B:132:ASP:HB3	2.01	0.41
1:B:118:ASN:HD22	1:B:164:ILE:H	1.69	0.41
1:B:42:VAL:HG23	1:B:56:ALA:HA	2.02	0.41
1:A:128:LEU:HD22	1:A:132:ASP:HB3	2.02	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:LYS:NZ	6:A:408:IHP:O33[4_597]	1.52	0.68
1:A:229:LYS:CE	6:A:408:IHP:O33[4_597]	2.02	0.18
1:A:234:HIS:ND1	6:A:408:IHP:O34[4_597]	2.06	0.14

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	325/352 (92%)	310 (95%)	13 (4%)	2 (1%)	25 30
1	B	325/352 (92%)	308 (95%)	17 (5%)	0	100 100
All	All	650/704 (92%)	618 (95%)	30 (5%)	2 (0%)	41 50

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	267	PRO
1	A	72	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	302/319 (95%)	301 (100%)	1 (0%)	92 96
1	B	302/319 (95%)	302 (100%)	0	100 100
All	All	604/638 (95%)	603 (100%)	1 (0%)	93 97

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	58	ASN
1	B	118	ASN
1	A	118	ASN
1	A	262	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry

Of 12 ligands modelled in this entry, 3 are monoatomic - leaving 9 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	$\#\|Z\| > 2$
3	ACT	A	406	-	1,3,3	5.26	1 (100%)	0,3,3	0.00	-
2	ADP	B	401	-	24,29,29	1.07	1 (4%)	29,45,45	1.33	3 (10%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
3	ACT	B	403	-	1,3,3	4.89	1 (100%)	0,3,3	0.00	-
3	ACT	A	407	-	1,3,3	6.59	1 (100%)	0,3,3	0.00	-
6	IHP	A	408	-	36,36,36	1.13	3 (8%)	54,60,60	0.82	1 (1%)
3	ACT	B	402	-	1,3,3	5.77	1 (100%)	0,3,3	0.00	-
4	A4B	A	402	-	18,18,18	0.08	0	24,24,24	0.27	0
3	ACT	A	405	-	1,3,3	5.23	1 (100%)	0,3,3	0.00	-
2	ADP	A	401	5	24,29,29	0.80	1 (4%)	29,45,45	0.91	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
4	A4B	A	402	-	-	0/8/8/8	0/2/2/2
6	IHP	A	408	-	-	2/30/54/54	0/1/1/1
2	ADP	A	401	5	-	0/12/32/32	0/3/3/3
2	ADP	B	401	-	-	2/12/32/32	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	407	ACT	CH3-C	6.59	1.57	1.48
3	B	402	ACT	CH3-C	5.77	1.56	1.48
3	A	406	ACT	CH3-C	5.26	1.55	1.48
3	A	405	ACT	CH3-C	5.23	1.55	1.48
3	B	403	ACT	CH3-C	4.89	1.55	1.48
6	A	408	IHP	P4-O14	3.98	1.66	1.59
6	A	408	IHP	P3-O13	2.90	1.64	1.59
2	B	401	ADP	C2'-C1'	-2.59	1.49	1.53
6	A	408	IHP	P1-O11	2.20	1.63	1.59
2	A	401	ADP	PB-O1B	2.15	1.57	1.50

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	401	ADP	N3-C2-N1	-3.50	123.20	128.68
2	B	401	ADP	C4-C5-N7	-2.99	106.28	109.40
6	A	408	IHP	C5-C4-C3	2.69	116.30	110.41
2	B	401	ADP	C2-N1-C6	2.36	122.79	118.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	401	ADP	C5-C6-N6	2.18	123.67	120.35
2	A	401	ADP	O3B-PB-O3A	2.12	111.76	104.64

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	408	IHP	C5-C4-O14-P4
6	A	408	IHP	C3-C4-O14-P4
2	B	401	ADP	PB-O3A-PA-O2A
2	B	401	ADP	PB-O3A-PA-O1A

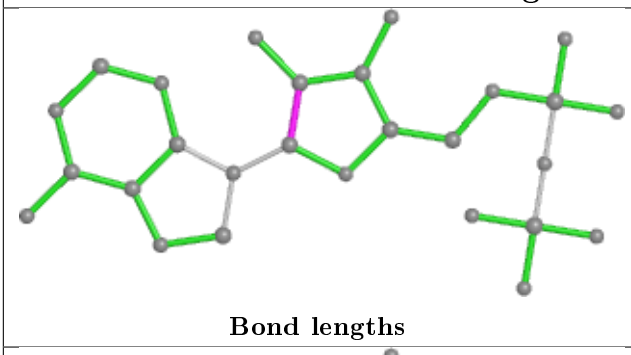
There are no ring outliers.

2 monomers are involved in 7 short contacts:

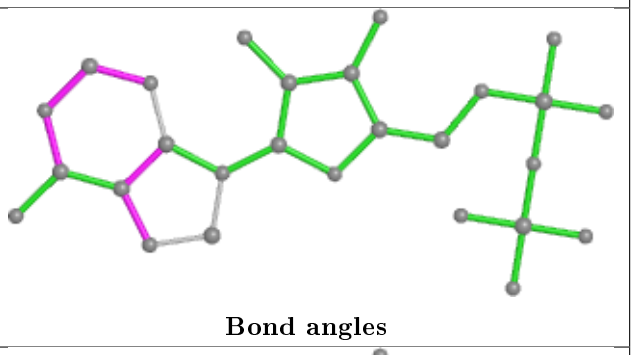
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	408	IHP	0	3
4	A	402	A4B	4	0

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 then 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.

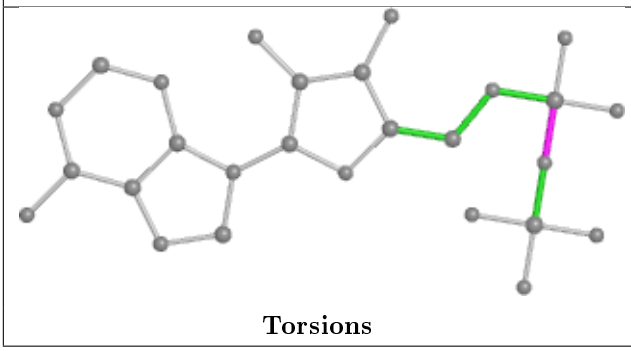
Ligand ADP B 401



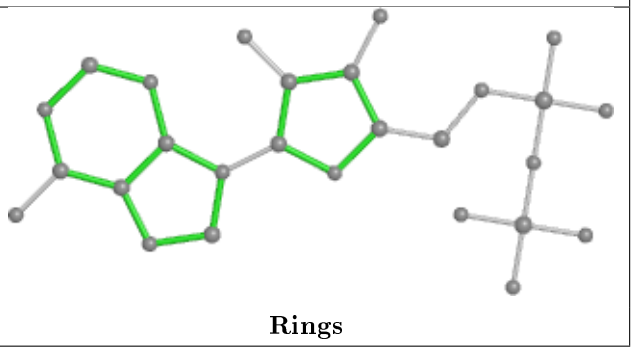
Bond lengths



Bond angles

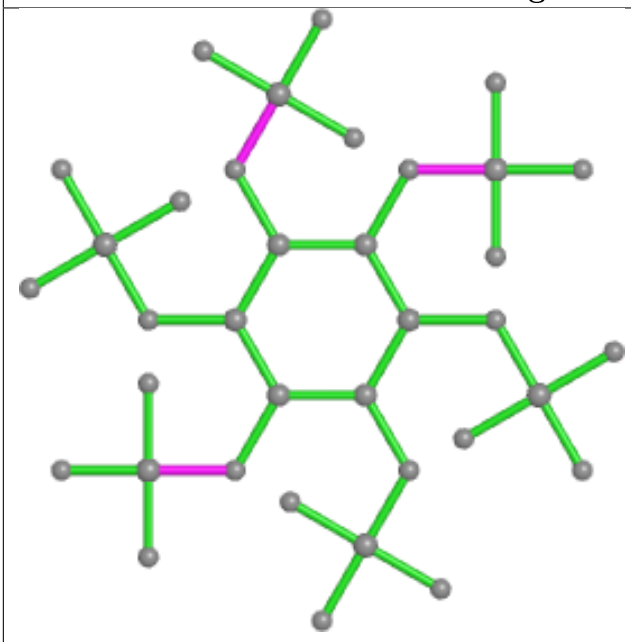


Torsions

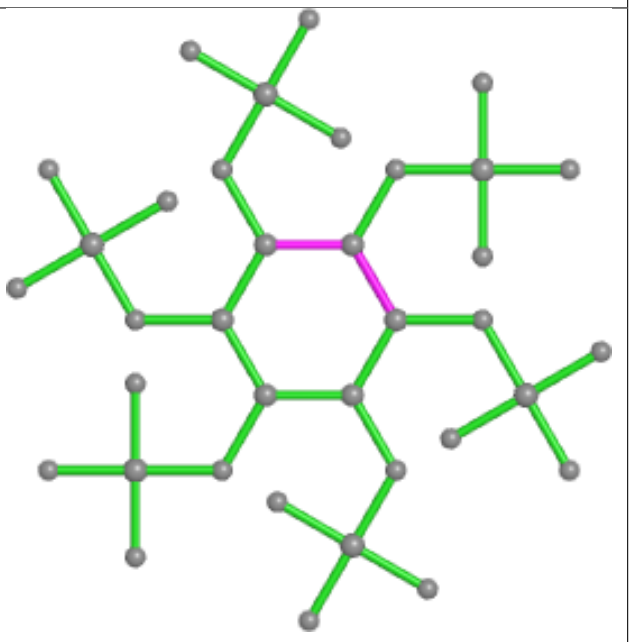


Rings

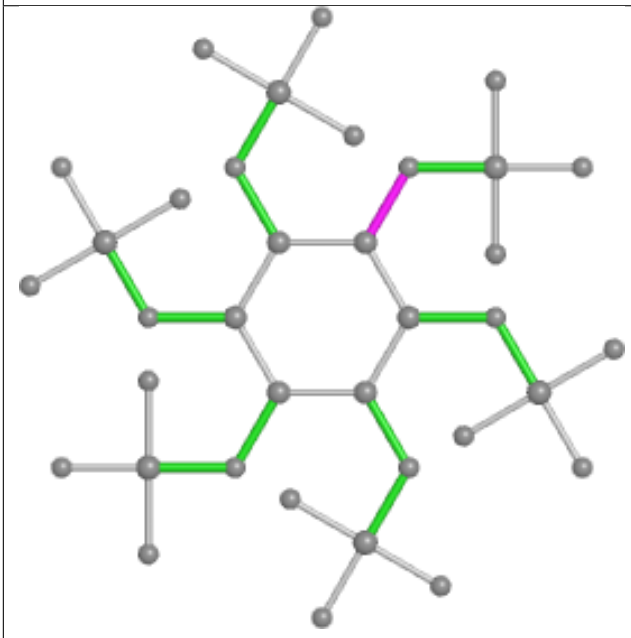
Ligand IHP A 408



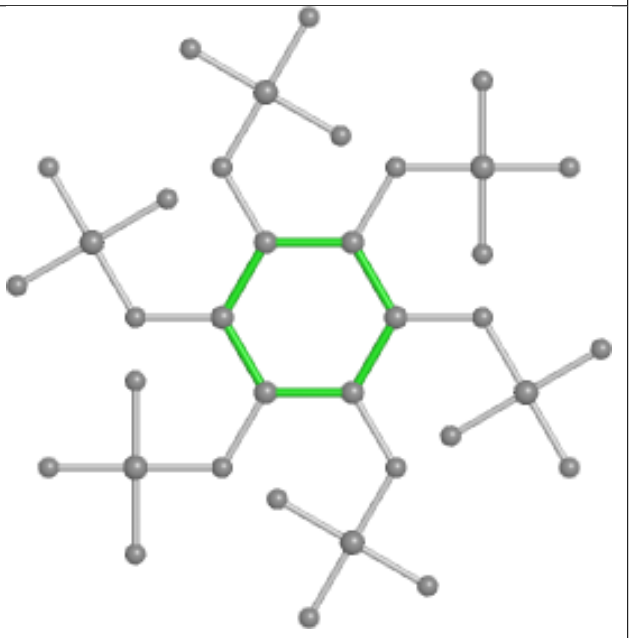
Bond lengths



Bond angles

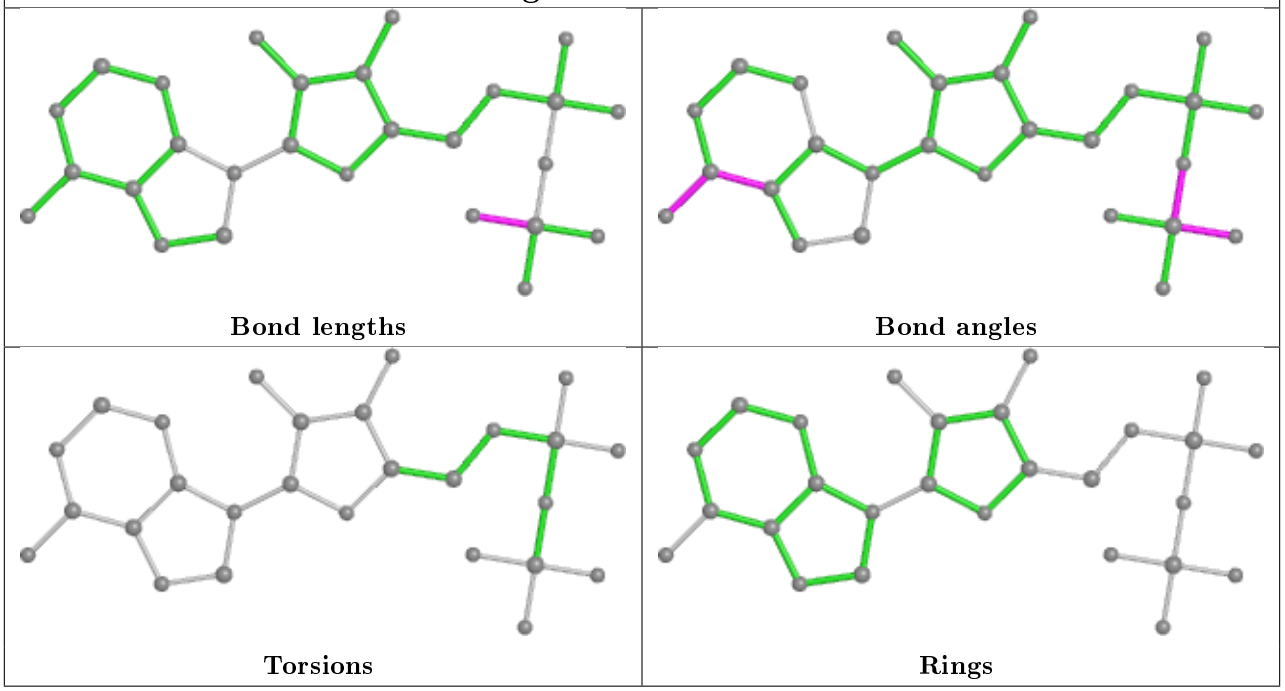


Torsions



Rings

Ligand ADP A 401



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, 95^{th} 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	324/352 (92%)	0.79	21 (6%) 18 24	34, 58, 96, 114	0
1	B	324/352 (92%)	1.20	74 (22%) 0 1	46, 81, 115, 140	0
All	All	648/704 (92%)	0.99	95 (14%) 2 3	34, 70, 111, 140	0

All (95) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	307	TYR	5.9
1	B	50	TYR	5.5
1	B	292	LEU	5.2
1	A	33	TRP	5.0
1	B	257	TYR	4.7
1	B	273	LEU	4.7
1	A	325	TYR	4.6
1	B	77	LYS	4.1
1	B	325	TYR	4.0
1	B	41	LEU	3.9
1	B	269	PHE	3.9
1	B	242	LEU	3.9
1	B	33	TRP	3.9
1	A	72	PRO	3.9
1	B	54	PHE	3.8
1	B	204	VAL	3.8
1	B	223	ALA	3.8
1	B	267	PRO	3.7
1	B	225	MET	3.7
1	B	75	LYS	3.5
1	B	254	LEU	3.5
1	B	125	TYR	3.5
1	B	62	ASN	3.4
1	B	227	PHE	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	42	VAL	3.4
1	A	71	LYS	3.4
1	B	216	TRP	3.3
1	A	50	TYR	3.3
1	B	76	LYS	3.3
1	A	271	ASP	3.3
1	B	249	LEU	3.2
1	B	265	LEU	3.2
1	B	245	ILE	3.1
1	B	206	TYR	3.1
1	B	270	ASN	3.0
1	B	250	GLY	3.0
1	B	326	THR	3.0
1	B	243	VAL	2.9
1	A	128	LEU	2.9
1	B	261	TYR	2.9
1	B	30	VAL	2.9
1	B	28	SER	2.9
1	B	298	LEU	2.9
1	B	305	LEU	2.9
1	B	51	SER	2.8
1	A	273	LEU	2.8
1	B	72	PRO	2.7
1	A	51	SER	2.7
1	B	5	VAL	2.7
1	B	202	LEU	2.7
1	B	281	TRP	2.7
1	B	171	LEU	2.7
1	B	178	LEU	2.7
1	A	124	LEU	2.6
1	B	121	PHE	2.6
1	B	173	LEU	2.6
1	A	70	LEU	2.6
1	B	255	TYR	2.6
1	B	239	TYR	2.5
1	B	52	GLU	2.5
1	B	271	ASP	2.4
1	B	66	VAL	2.4
1	B	124	LEU	2.4
1	B	57	ILE	2.4
1	A	326	THR	2.4
1	B	59	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	269	PHE	2.3
1	B	46	GLY	2.3
1	B	74	LYS	2.3
1	A	284	PHE	2.3
1	B	263	ILE	2.3
1	A	268	ARG	2.2
1	A	272	ILE	2.2
1	B	64	LYS	2.2
1	B	26	TYR	2.2
1	A	76	LYS	2.2
1	A	327	VAL	2.2
1	B	21	SER	2.2
1	B	69	ILE	2.2
1	B	82	ILE	2.2
1	B	213	LEU	2.2
1	B	100	ILE	2.2
1	B	45	LEU	2.2
1	B	244	ARG	2.2
1	B	272	ILE	2.1
1	A	222	LEU	2.1
1	B	160	HIS	2.1
1	B	232	PHE	2.1
1	A	239	TYR	2.1
1	B	191	ARG	2.0
1	A	225	MET	2.0
1	B	67	VAL	2.0
1	B	126	GLN	2.0
1	B	31	VAL	2.0
1	B	70	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)

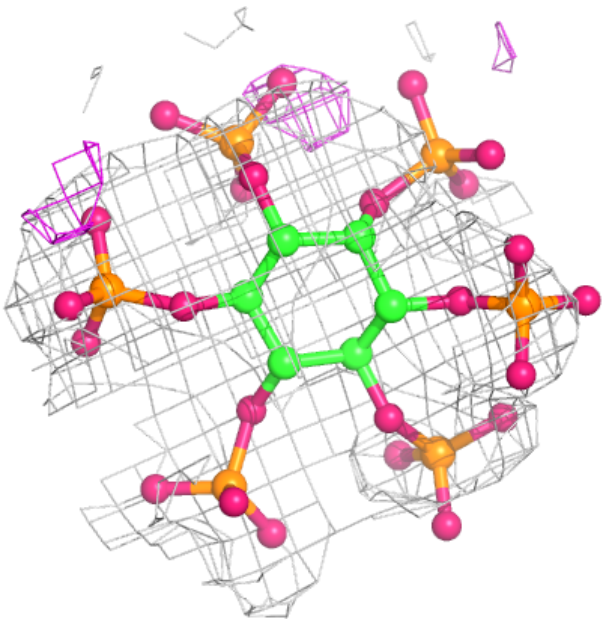
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

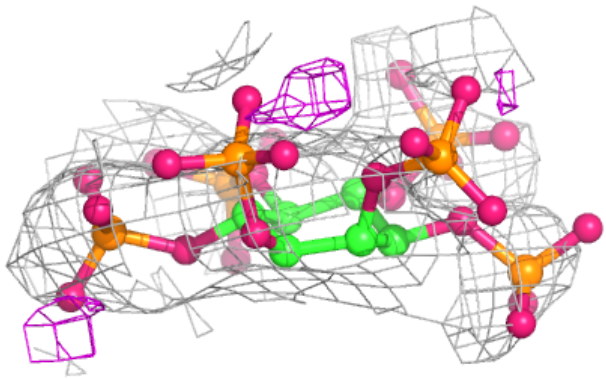
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	ACT	B	402	4/4	0.71	0.28	79,80,80,81	0
6	IHP	A	408	36/36	0.74	0.24	124,128,132,133	36
4	A4B	A	402	17/17	0.79	0.25	62,69,79,90	0
3	ACT	B	403	4/4	0.81	0.33	78,81,83,86	0
3	ACT	A	405	4/4	0.81	0.31	72,73,74,75	0
2	ADP	B	401	27/27	0.87	0.24	96,108,126,128	0
5	MG	A	403	1/1	0.90	0.12	61,61,61,61	0
7	CL	A	409	1/1	0.91	0.24	93,93,93,93	0
3	ACT	A	406	4/4	0.91	0.26	66,70,71,73	0
2	ADP	A	401	27/27	0.93	0.16	54,71,87,90	0
3	ACT	A	407	4/4	0.94	0.29	61,63,63,64	0
5	MG	A	404	1/1	0.97	0.09	64,64,64,64	0

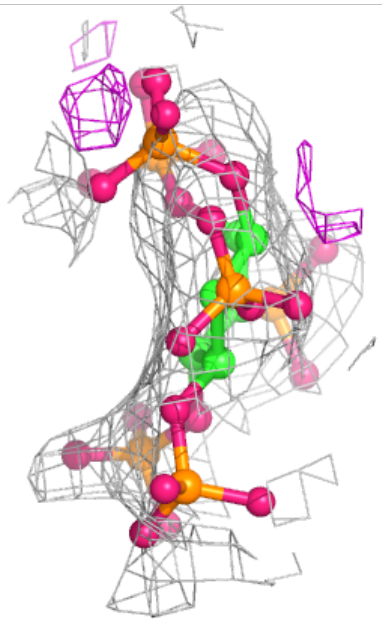
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around IHP A 408:

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

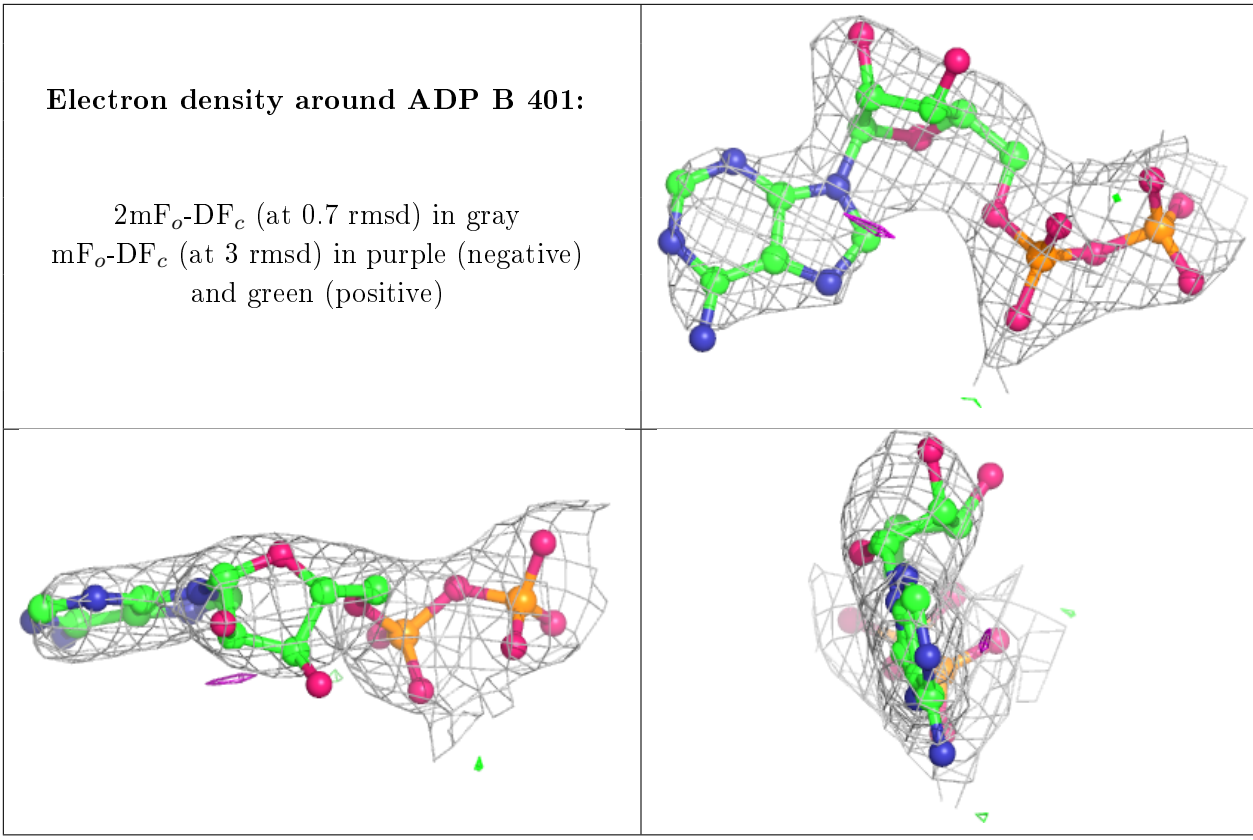






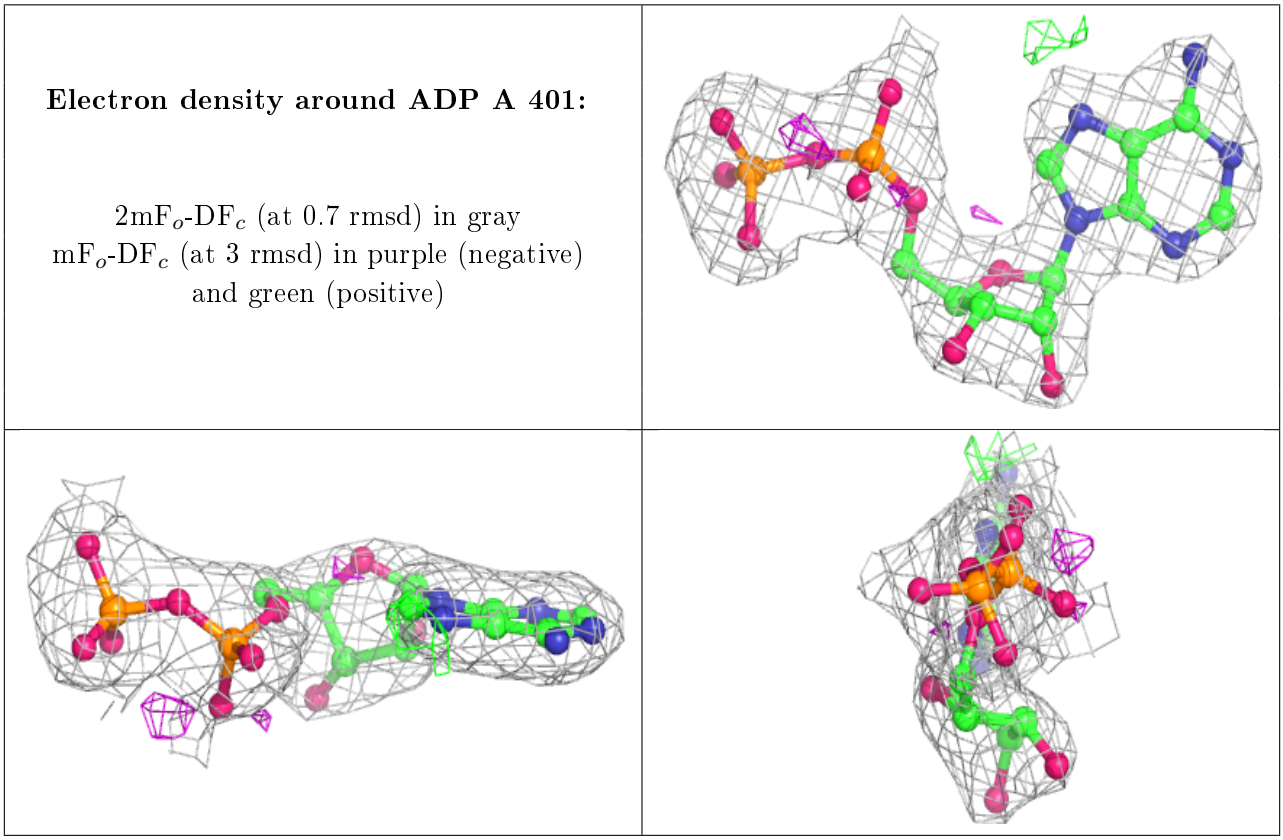
Electron density around ADP B 401:

$2mF_o$-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 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 (i)

There are no such residues in this entry.