wwPDB X-ray Structure Validation Summary Report (i)

Jun 13, 2020 – 04:46 pm BST

PDB ID	:	1W46
Title	:	P4 protein from Bacteriophage PHI12 in complex with ADP and MG
Authors	:	Mancini, E.J.; Kainov, D.E.; Grimes, J.M.; Tuma, R.; Bamford, D.H.; Stuart, D.I.
Deposited on	:	2004-07-22
Resolution	:	2.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) symbol.

The following versions of software and data (see references (i)) were used in the production of this report:

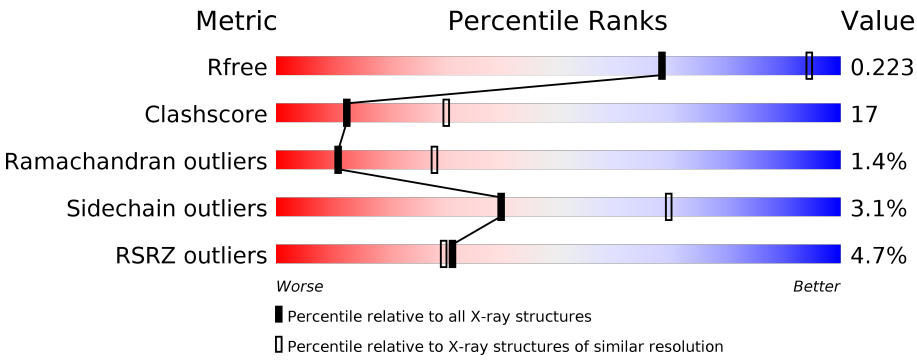
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 2.70 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	331	4% 63% 23% • 13%
1	B	331	4% 60% 25% • 13%
1	C	331	4% 62% 23% • 13%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	MG	C	701	-	-	-	X

2 Entry composition (i)

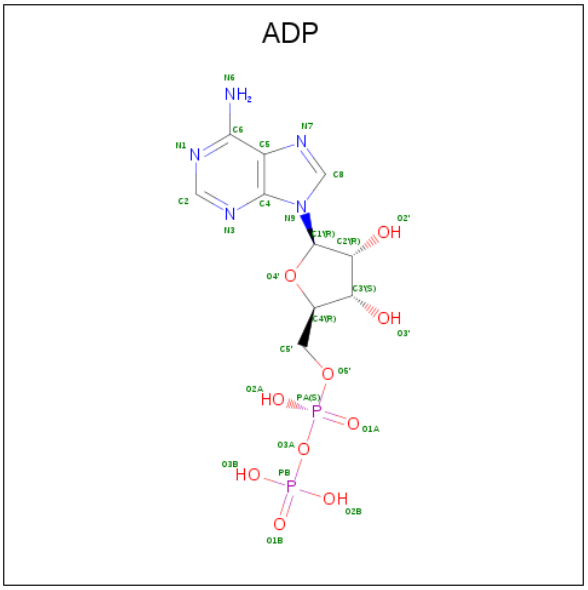
There are 4 unique types of molecules in this entry. The entry contains 6932 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 NTPASE P4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	288	2170	1361	378	424	7	0	0	0
1	B	288	2170	1361	378	424	7	0	0	0
1	C	288	2170	1361	378	424	7	0	0	0

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

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Mg 1 1	0	0
3	A	1	Total Mg 1 1	0	0
3	C	1	Total Mg 1 1	0	0

- Molecule 4 is water.

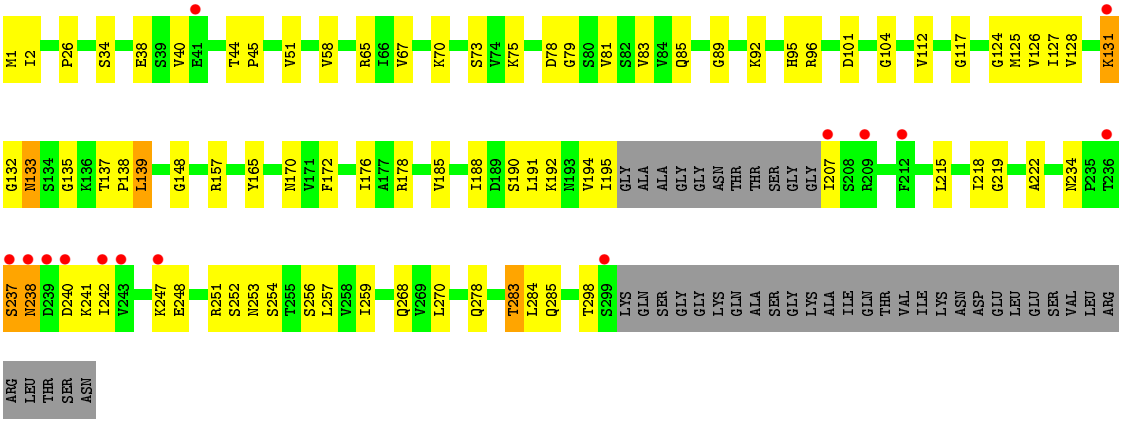
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	111	Total O 111 111	0	0
4	B	118	Total O 118 118	0	0
4	C	109	Total O 109 109	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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.

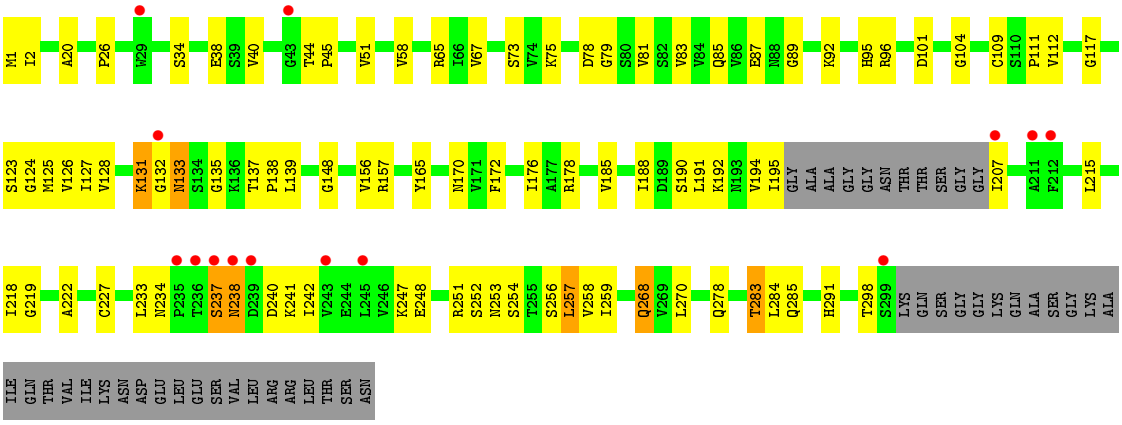
- Molecule 1: NTPASE P4

Chain A:



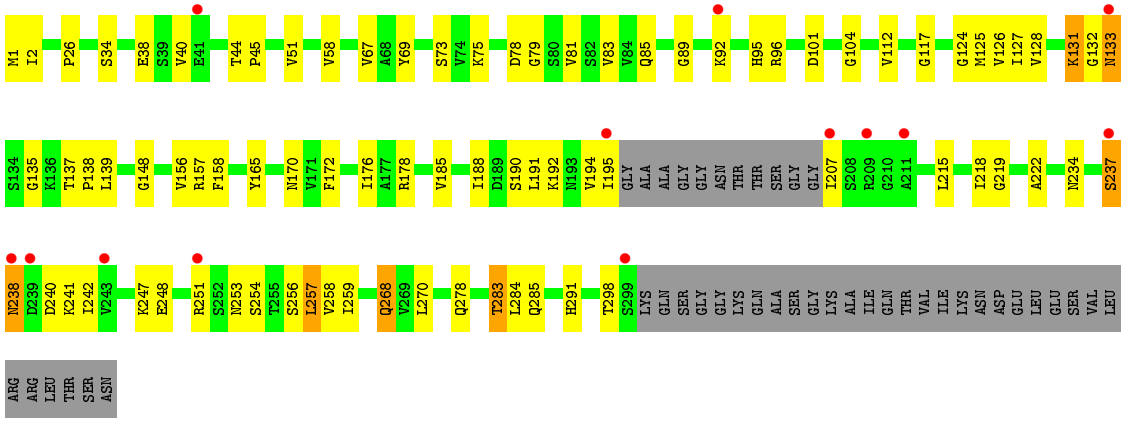
- Molecule 1: NTPASE P4

Chain B:



- Molecule 1: NTPASE P4

Chain C:


M1
I2
P26
S34
E38
S39
V40
E41
T44
P45
V51
V58
V67
A68
Y69
S73
V74
K75
D78
G79
S80
V81
S82
V83
V84
Q85
G89
K92
H95
R96
D101
G104
V112
G117
G124
M125
V126
I127
V128
K131
G132
N133
S134
G135
K136
T137
P138
L139
G148
V156
R157
F158
Y165
N170
V171
F172
I176
A177
R178
V185
I188
D189
S190
L191
K192
N193
V194
I195
GLY
ALA
ALA
GLY
GLY
ASN
THR
THR
SER
GLY
GLY
I207
S208
R209
G210
A211
L215
I218
G219
A222
N234
S237
N238
D239
D240
K241
I242
V243
K247
E248
R251
S252
N253
S254
T255
S256
L257
V258
I259
Q268
V269
L270
Q278
T283
L284
Q285
H291
T298
S299
LYS
GLN
SER
GLY
GLY
LYS
GLN
ALA
SER
GLY
LYS
ALA
ILE
GLN
THR
VAL
ILE
LYS
ASN
ASP
GLU
LEU
GLU
SER
VAL
LEU
ARG
ARG
LEU
THR
SER
ASN

4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	105.00Å 128.40Å 158.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.70 29.75 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.6 (20.00-2.70) 99.9 (29.75-2.70)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	7.05 (at 2.72Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.209 , 0.253 0.208 , 0.223	Depositor DCC
R_{free} test set	1461 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å^2)	37.5	Xtriage
Anisotropy	0.398	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 59.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	6932	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 5.78% 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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, 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	0.39	0/2207	0.64	0/2989
1	B	0.38	0/2207	0.64	0/2989
1	C	0.37	0/2207	0.64	0/2989
All	All	0.38	0/6621	0.64	0/8967

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	2170	0	2147	68	0
1	B	2170	0	2147	79	0
1	C	2170	0	2147	75	0
2	A	27	0	12	1	0
2	B	27	0	12	1	0
2	C	27	0	12	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	111	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	118	0	0	10	0
4	C	109	0	0	7	0
All	All	6932	0	6477	220	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 17.

The worst 5 of 220 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:C:117:GLY:HA2	1:C:298:THR:HG23	1.41	1.03
1:B:117:GLY:HA2	1:B:298:THR:HG23	1.41	1.02
1:A:117:GLY:HA2	1:A:298:THR:HG23	1.41	1.01
1:B:111:PRO:HG3	4:B:2050:HOH:O	1.68	0.92
1:B:240:ASP:HB3	1:B:242:ILE:HD13	1.52	0.92

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	284/331 (86%)	267 (94%)	13 (5%)	4 (1%)	11	28
1	B	284/331 (86%)	268 (94%)	12 (4%)	4 (1%)	11	28
1	C	284/331 (86%)	268 (94%)	12 (4%)	4 (1%)	11	28
All	All	852/993 (86%)	803 (94%)	37 (4%)	12 (1%)	11	28

5 of 12 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	237	SER

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Mol	Chain	Res	Type
1	A	238	ASN
1	B	237	SER
1	B	238	ASN
1	C	237	SER

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	234/265 (88%)	226 (97%)	8 (3%)	37	66
1	B	234/265 (88%)	227 (97%)	7 (3%)	41	70
1	C	234/265 (88%)	227 (97%)	7 (3%)	41	70
All	All	702/795 (88%)	680 (97%)	22 (3%)	40	69

5 of 22 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	248	GLU
1	B	278	GLN
1	C	283	THR
1	B	257	LEU
1	B	268	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 28 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	238	ASN
1	B	285	GLN
1	C	285	GLN
1	B	268	GLN
1	B	278	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 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	ADP	B	700	3	24,29,29	1.73	5 (20%)	29,45,45	1.79	7 (24%)
2	ADP	A	700	3	24,29,29	1.64	6 (25%)	29,45,45	1.81	6 (20%)
2	ADP	C	700	3	24,29,29	1.68	6 (25%)	29,45,45	1.80	6 (20%)

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	B	700	3	-	4/12/32/32	0/3/3/3
2	ADP	A	700	3	-	4/12/32/32	0/3/3/3
2	ADP	C	700	3	-	4/12/32/32	0/3/3/3

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	700	ADP	O4'-C1'	4.33	1.47	1.41
2	C	700	ADP	O4'-C1'	4.01	1.46	1.41
2	A	700	ADP	O4'-C1'	3.67	1.46	1.41
2	B	700	ADP	C2-N3	3.42	1.37	1.32
2	B	700	ADP	C5-N7	-3.22	1.28	1.39

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	700	ADP	N3-C2-N1	-6.50	118.52	128.68
2	B	700	ADP	N3-C2-N1	-6.45	118.60	128.68
2	A	700	ADP	N3-C2-N1	-6.44	118.61	128.68
2	A	700	ADP	O4'-C1'-C2'	-3.10	102.39	106.93
2	C	700	ADP	O4'-C1'-C2'	-2.96	102.60	106.93

There are no chirality outliers.

5 of 12 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	700	ADP	C3'-C4'-C5'-O5'
2	B	700	ADP	C3'-C4'-C5'-O5'
2	A	700	ADP	C3'-C4'-C5'-O5'
2	C	700	ADP	O4'-C4'-C5'-O5'
2	B	700	ADP	O4'-C4'-C5'-O5'

There are no ring outliers.

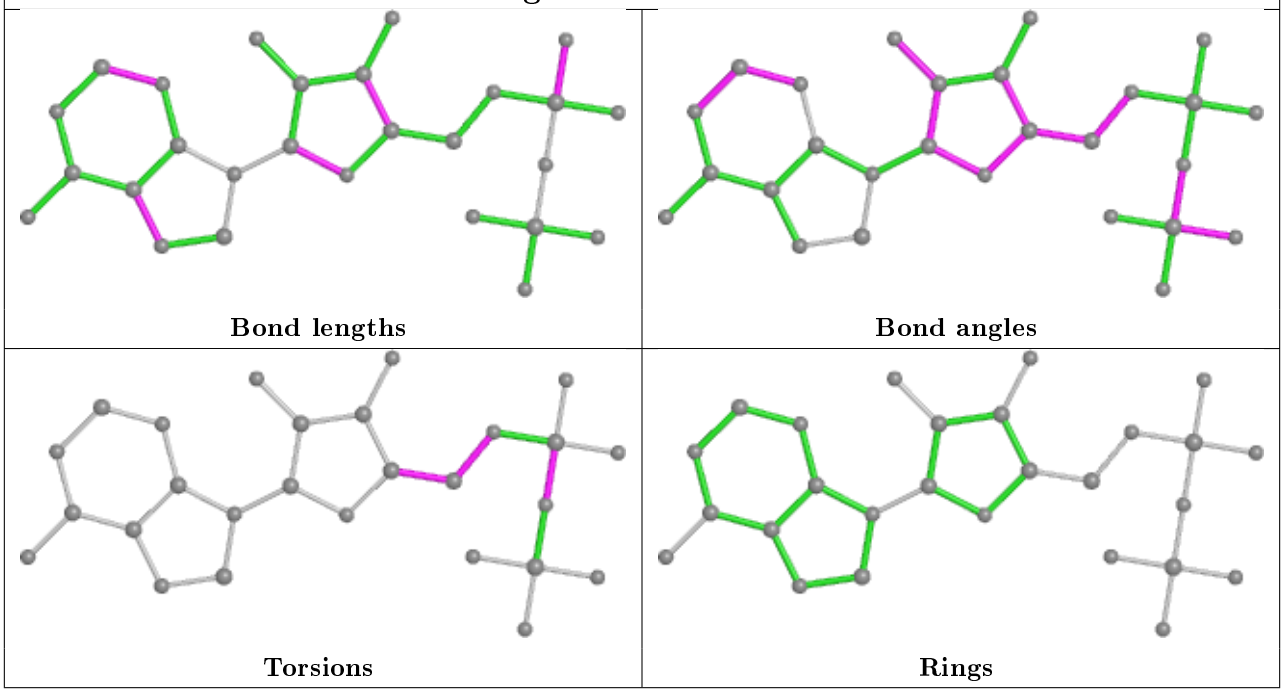
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	700	ADP	1	0
2	A	700	ADP	1	0
2	C	700	ADP	1	0

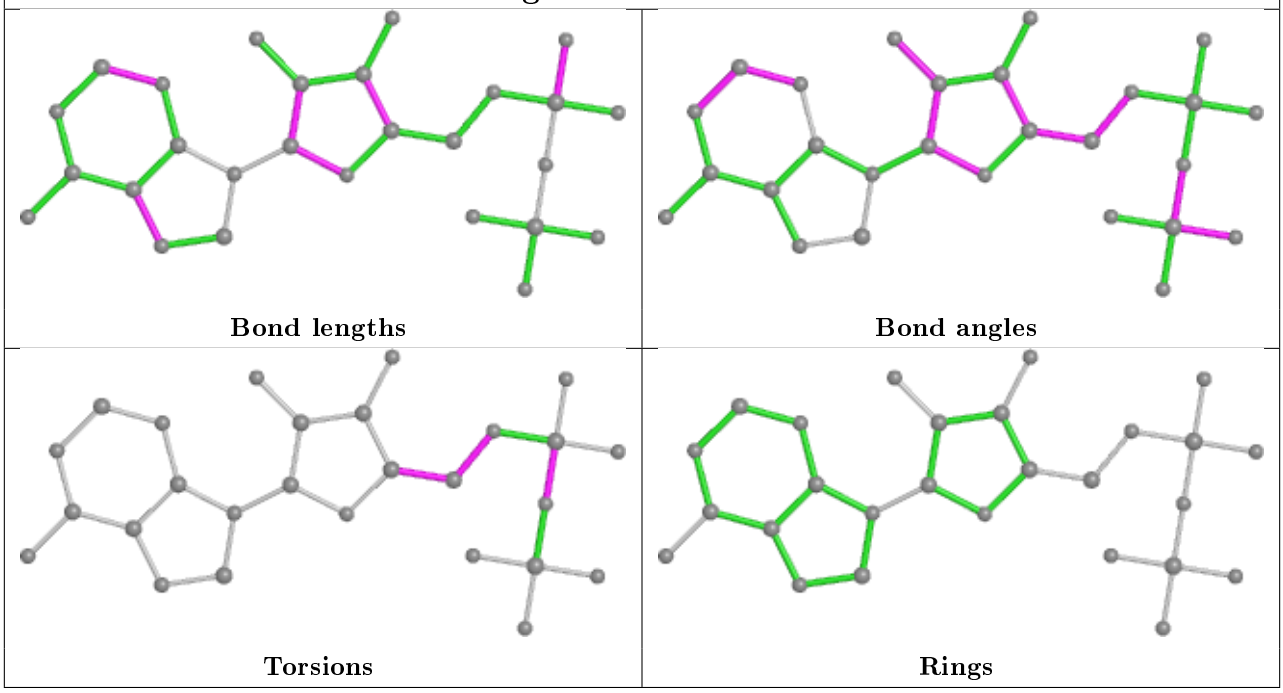
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

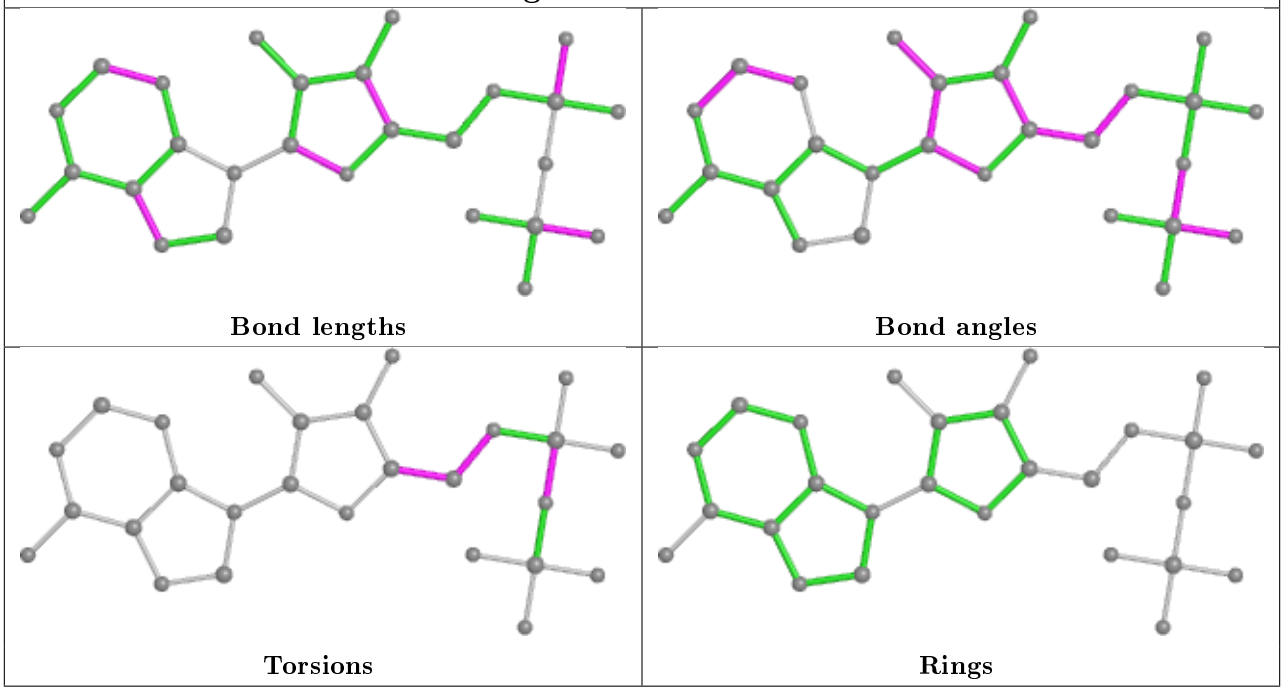
Ligand ADP B 700



Ligand ADP A 700



Ligand ADP C 700



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 (i)

6.1 Protein, DNA and RNA chains (i)

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	288/331 (87%)	-0.16	14 (4%)	29	28	14, 30, 82, 126	0
1	B	288/331 (87%)	-0.14	14 (4%)	29	28	14, 30, 82, 126	0
1	C	288/331 (87%)	-0.12	13 (4%)	33	31	15, 31, 83, 126	0
All	All	864/993 (87%)	-0.14	41 (4%)	31	30	14, 30, 83, 126	0

The worst 5 of 41 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	238	ASN	5.9
1	B	237	SER	4.7
1	A	238	ASN	4.6
1	A	239	ASP	4.5
1	C	239	ASP	4.1

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 carbohydrates in this entry.

6.4 Ligands (i)

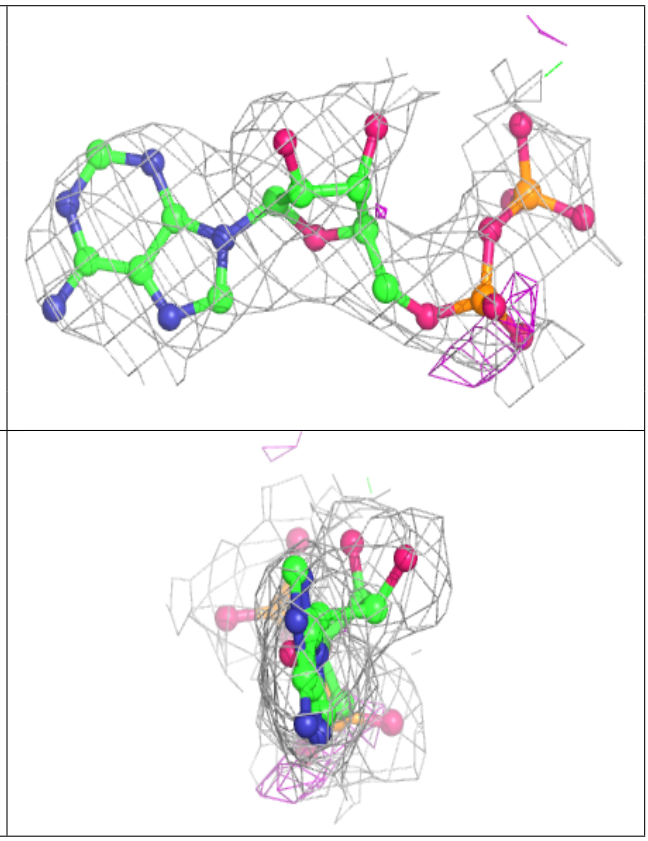
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	MG	C	701	1/1	0.75	0.42	74,74,74,74	0
3	MG	A	701	1/1	0.80	0.66	79,79,79,79	0
2	ADP	B	700	27/27	0.87	0.19	52,67,94,96	0
2	ADP	A	700	27/27	0.90	0.17	46,65,91,95	0
3	MG	B	701	1/1	0.90	0.58	74,74,74,74	0
2	ADP	C	700	27/27	0.92	0.19	50,65,90,93	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.

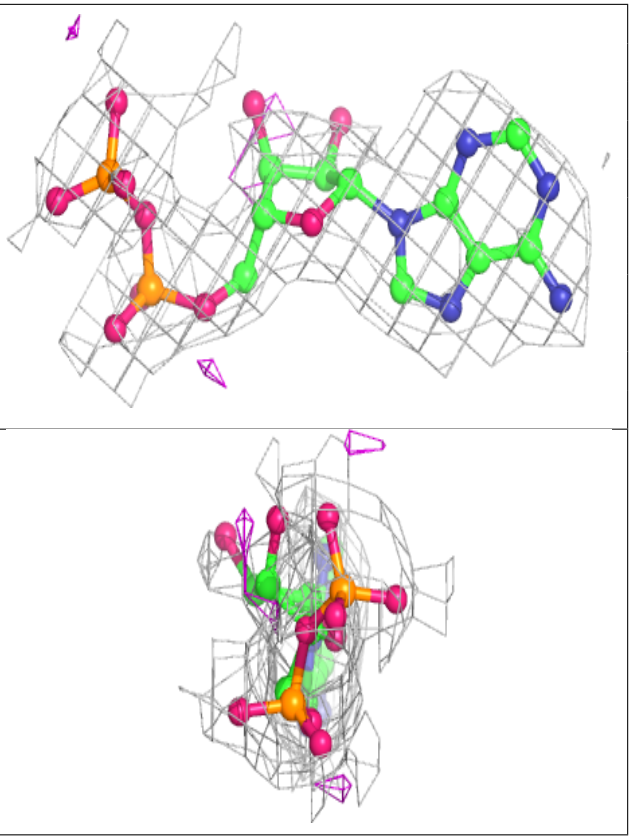
Electron density around ADP B 700:

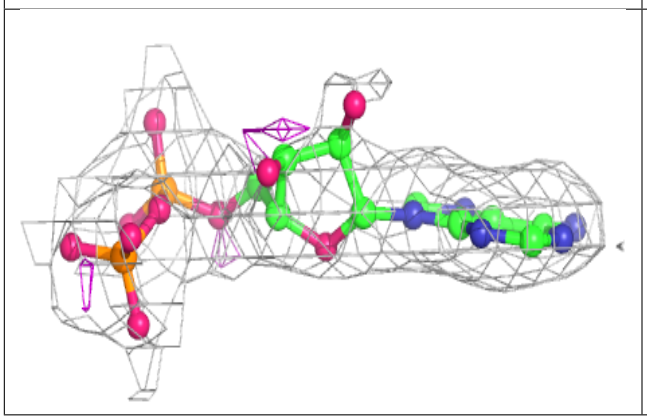
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)



Electron density around ADP A 700:

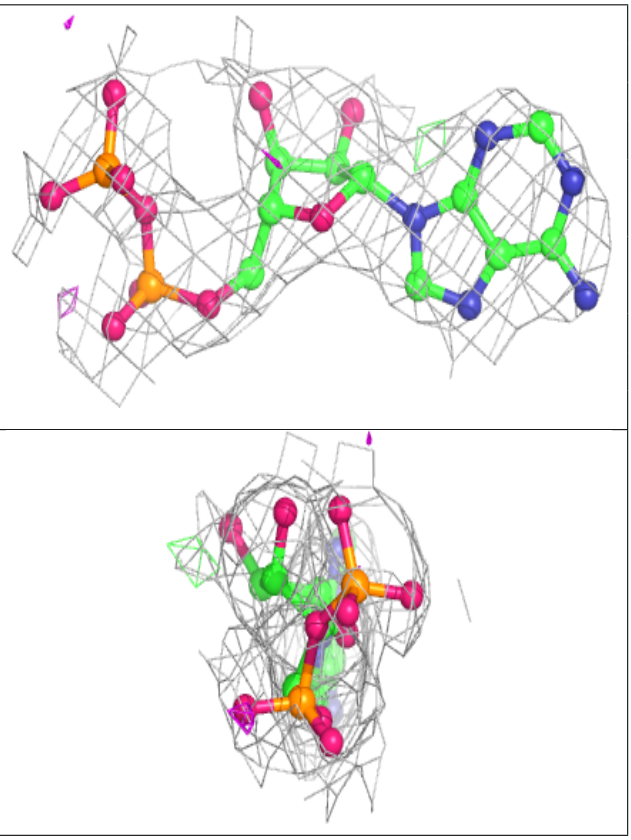
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

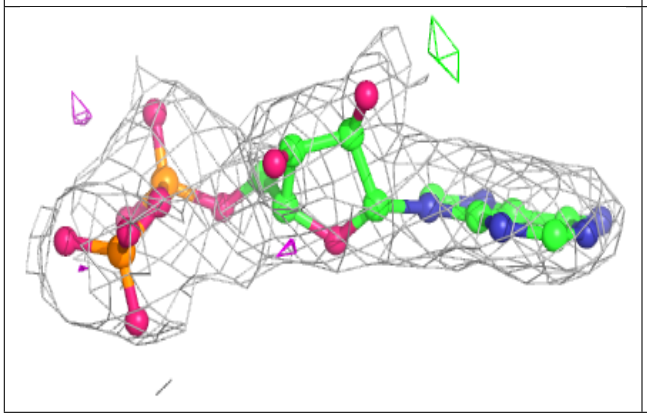




Electron density around ADP C 700:

$2mF_o$-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.