



# Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 02:24 am BST

PDB ID : 6OKV  
Title : PilT4 from *Geobacter metallireducens* bound to AMP-PNP: C2ccocco conformation  
Authors : McCallum, M.; Howell, P.L.  
Deposited on : 2019-04-15  
Resolution : 4.01 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](mailto: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.

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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	:	2.11
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

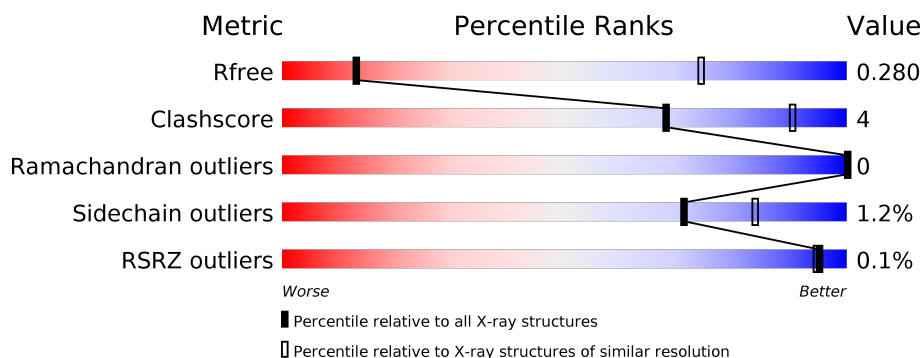
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 4.01 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	1087 (4.30-3.70)
Clashscore	141614	1148 (4.30-3.70)
Ramachandran outliers	138981	1108 (4.30-3.70)
Sidechain outliers	138945	1099 (4.30-3.70)
RSRZ outliers	127900	1028 (4.34-3.66)

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  $\geq 3$ , 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	385	 81% 10% 9%
1	B	385	 85% 6% 9%
1	C	385	 83% 8% 9%
1	D	385	 83% 10% 7%
1	E	385	 80% 11% 8%
1	F	385	 81% 11% 9%

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 16189 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 Twitching motility pilus retraction ATPase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	352	Total	C	N	O	S	0	0	0
			2656	1679	466	498	13			
1	B	351	Total	C	N	O	S	0	0	0
			2668	1688	467	499	14			
1	C	351	Total	C	N	O	S	0	0	0
			2665	1687	464	501	13			
1	D	358	Total	C	N	O	S	0	0	0
			2734	1724	484	512	14			
1	E	353	Total	C	N	O	S	0	0	0
			2661	1686	464	497	14			
1	F	352	Total	C	N	O	S	0	0	0
			2623	1663	456	490	14			

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q39VU6
A	-18	GLY	-	expression tag	UNP Q39VU6
A	-17	SER	-	expression tag	UNP Q39VU6
A	-16	SER	-	expression tag	UNP Q39VU6
A	-15	HIS	-	expression tag	UNP Q39VU6
A	-14	HIS	-	expression tag	UNP Q39VU6
A	-13	HIS	-	expression tag	UNP Q39VU6
A	-12	HIS	-	expression tag	UNP Q39VU6
A	-11	HIS	-	expression tag	UNP Q39VU6
A	-10	HIS	-	expression tag	UNP Q39VU6
A	-9	SER	-	expression tag	UNP Q39VU6
A	-8	SER	-	expression tag	UNP Q39VU6
A	-7	GLY	-	expression tag	UNP Q39VU6
A	-6	LEU	-	expression tag	UNP Q39VU6
A	-5	VAL	-	expression tag	UNP Q39VU6
A	-4	PRO	-	expression tag	UNP Q39VU6
A	-3	ARG	-	expression tag	UNP Q39VU6

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q39VU6
A	-1	SER	-	expression tag	UNP Q39VU6
A	0	HIS	-	expression tag	UNP Q39VU6
B	-19	MET	-	expression tag	UNP Q39VU6
B	-18	GLY	-	expression tag	UNP Q39VU6
B	-17	SER	-	expression tag	UNP Q39VU6
B	-16	SER	-	expression tag	UNP Q39VU6
B	-15	HIS	-	expression tag	UNP Q39VU6
B	-14	HIS	-	expression tag	UNP Q39VU6
B	-13	HIS	-	expression tag	UNP Q39VU6
B	-12	HIS	-	expression tag	UNP Q39VU6
B	-11	HIS	-	expression tag	UNP Q39VU6
B	-10	HIS	-	expression tag	UNP Q39VU6
B	-9	SER	-	expression tag	UNP Q39VU6
B	-8	SER	-	expression tag	UNP Q39VU6
B	-7	GLY	-	expression tag	UNP Q39VU6
B	-6	LEU	-	expression tag	UNP Q39VU6
B	-5	VAL	-	expression tag	UNP Q39VU6
B	-4	PRO	-	expression tag	UNP Q39VU6
B	-3	ARG	-	expression tag	UNP Q39VU6
B	-2	GLY	-	expression tag	UNP Q39VU6
B	-1	SER	-	expression tag	UNP Q39VU6
B	0	HIS	-	expression tag	UNP Q39VU6
C	-19	MET	-	expression tag	UNP Q39VU6
C	-18	GLY	-	expression tag	UNP Q39VU6
C	-17	SER	-	expression tag	UNP Q39VU6
C	-16	SER	-	expression tag	UNP Q39VU6
C	-15	HIS	-	expression tag	UNP Q39VU6
C	-14	HIS	-	expression tag	UNP Q39VU6
C	-13	HIS	-	expression tag	UNP Q39VU6
C	-12	HIS	-	expression tag	UNP Q39VU6
C	-11	HIS	-	expression tag	UNP Q39VU6
C	-10	HIS	-	expression tag	UNP Q39VU6
C	-9	SER	-	expression tag	UNP Q39VU6
C	-8	SER	-	expression tag	UNP Q39VU6
C	-7	GLY	-	expression tag	UNP Q39VU6
C	-6	LEU	-	expression tag	UNP Q39VU6
C	-5	VAL	-	expression tag	UNP Q39VU6
C	-4	PRO	-	expression tag	UNP Q39VU6
C	-3	ARG	-	expression tag	UNP Q39VU6
C	-2	GLY	-	expression tag	UNP Q39VU6
C	-1	SER	-	expression tag	UNP Q39VU6

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Chain	Residue	Modelled	Actual	Comment	Reference
C	0	HIS	-	expression tag	UNP Q39VU6
D	-19	MET	-	expression tag	UNP Q39VU6
D	-18	GLY	-	expression tag	UNP Q39VU6
D	-17	SER	-	expression tag	UNP Q39VU6
D	-16	SER	-	expression tag	UNP Q39VU6
D	-15	HIS	-	expression tag	UNP Q39VU6
D	-14	HIS	-	expression tag	UNP Q39VU6
D	-13	HIS	-	expression tag	UNP Q39VU6
D	-12	HIS	-	expression tag	UNP Q39VU6
D	-11	HIS	-	expression tag	UNP Q39VU6
D	-10	HIS	-	expression tag	UNP Q39VU6
D	-9	SER	-	expression tag	UNP Q39VU6
D	-8	SER	-	expression tag	UNP Q39VU6
D	-7	GLY	-	expression tag	UNP Q39VU6
D	-6	LEU	-	expression tag	UNP Q39VU6
D	-5	VAL	-	expression tag	UNP Q39VU6
D	-4	PRO	-	expression tag	UNP Q39VU6
D	-3	ARG	-	expression tag	UNP Q39VU6
D	-2	GLY	-	expression tag	UNP Q39VU6
D	-1	SER	-	expression tag	UNP Q39VU6
D	0	HIS	-	expression tag	UNP Q39VU6
E	-19	MET	-	expression tag	UNP Q39VU6
E	-18	GLY	-	expression tag	UNP Q39VU6
E	-17	SER	-	expression tag	UNP Q39VU6
E	-16	SER	-	expression tag	UNP Q39VU6
E	-15	HIS	-	expression tag	UNP Q39VU6
E	-14	HIS	-	expression tag	UNP Q39VU6
E	-13	HIS	-	expression tag	UNP Q39VU6
E	-12	HIS	-	expression tag	UNP Q39VU6
E	-11	HIS	-	expression tag	UNP Q39VU6
E	-10	HIS	-	expression tag	UNP Q39VU6
E	-9	SER	-	expression tag	UNP Q39VU6
E	-8	SER	-	expression tag	UNP Q39VU6
E	-7	GLY	-	expression tag	UNP Q39VU6
E	-6	LEU	-	expression tag	UNP Q39VU6
E	-5	VAL	-	expression tag	UNP Q39VU6
E	-4	PRO	-	expression tag	UNP Q39VU6
E	-3	ARG	-	expression tag	UNP Q39VU6
E	-2	GLY	-	expression tag	UNP Q39VU6
E	-1	SER	-	expression tag	UNP Q39VU6
E	0	HIS	-	expression tag	UNP Q39VU6
F	-19	MET	-	expression tag	UNP Q39VU6

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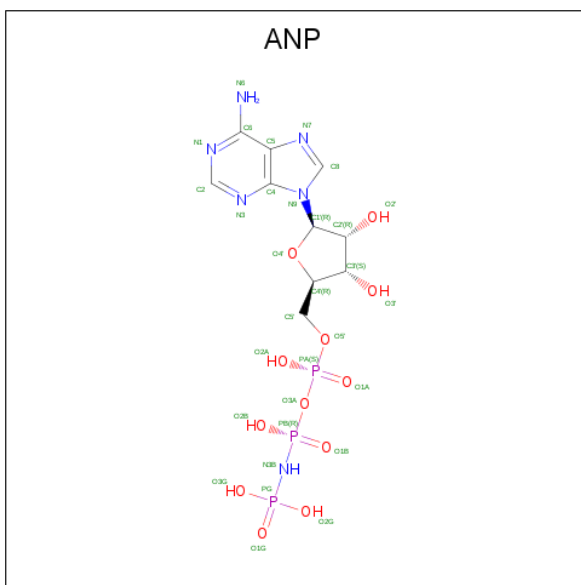
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Chain	Residue	Modelled	Actual	Comment	Reference
F	-18	GLY	-	expression tag	UNP Q39VU6
F	-17	SER	-	expression tag	UNP Q39VU6
F	-16	SER	-	expression tag	UNP Q39VU6
F	-15	HIS	-	expression tag	UNP Q39VU6
F	-14	HIS	-	expression tag	UNP Q39VU6
F	-13	HIS	-	expression tag	UNP Q39VU6
F	-12	HIS	-	expression tag	UNP Q39VU6
F	-11	HIS	-	expression tag	UNP Q39VU6
F	-10	HIS	-	expression tag	UNP Q39VU6
F	-9	SER	-	expression tag	UNP Q39VU6
F	-8	SER	-	expression tag	UNP Q39VU6
F	-7	GLY	-	expression tag	UNP Q39VU6
F	-6	LEU	-	expression tag	UNP Q39VU6
F	-5	VAL	-	expression tag	UNP Q39VU6
F	-4	PRO	-	expression tag	UNP Q39VU6
F	-3	ARG	-	expression tag	UNP Q39VU6
F	-2	GLY	-	expression tag	UNP Q39VU6
F	-1	SER	-	expression tag	UNP Q39VU6
F	0	HIS	-	expression tag	UNP Q39VU6

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

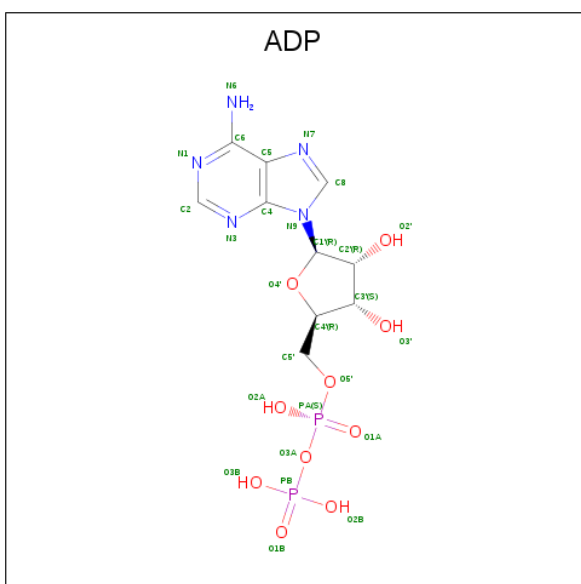
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Mg 1 1	0	0
2	A	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0
2	E	1	Total Mg 1 1	0	0

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: C<sub>10</sub>H<sub>17</sub>N<sub>6</sub>O<sub>12</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 31	C 10	N 6	O 12	P 3	0	0
3	B	1	Total 31	C 10	N 6	O 12	P 3	0	0
3	D	1	Total 31	C 10	N 6	O 12	P 3	0	0
3	E	1	Total 31	C 10	N 6	O 12	P 3	0	0

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $\text{C}_{10}\text{H}_{15}\text{N}_5\text{O}_{10}\text{P}_2$ ).




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

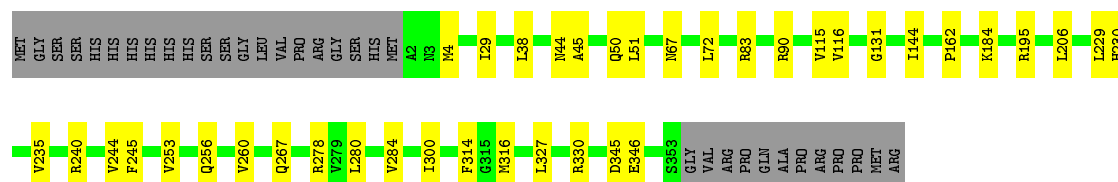


### 3 Residue-property plots [i](#)


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.

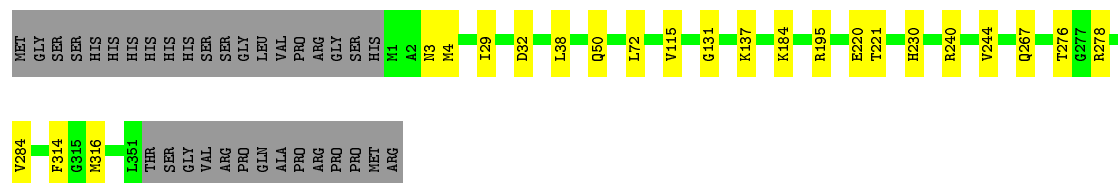
- Molecule 1: Twitching motility pilus retraction ATPase

Chain A: 




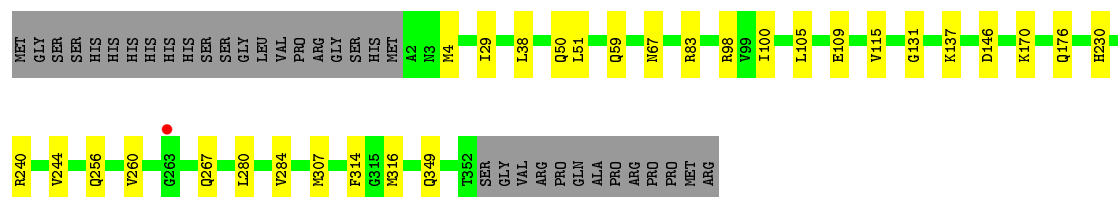
- Molecule 1: Twitching motility pilus retraction ATPase

Chain B: 




- Molecule 1: Twitching motility pilus retraction ATPase

Chain C: 



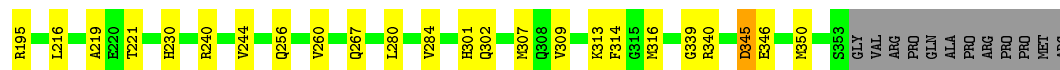
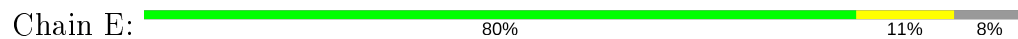
- Molecule 1: Twitching motility pilus retraction ATPase

Chain D: 

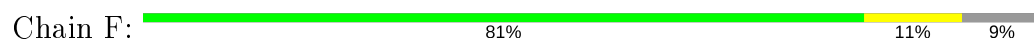




- Molecule 1: Twitching motility pilus retraction ATPase



- Molecule 1: Twitching motility pilus retraction ATPase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	98.51Å 127.13Å 187.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.26 – 4.01 49.26 – 4.01	Depositor EDS
% Data completeness (in resolution range)	60.0 (49.26-4.01) 60.1 (49.26-4.01)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.12 (at 4.00Å)	Xtriage
Refinement program	PHENIX (1.14_3260: ???)	Depositor
R, $R_{free}$	0.243 , 0.280 0.243 , 0.280	Depositor DCC
$R_{free}$ test set	593 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	84.0	Xtriage
Anisotropy	0.297	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 37.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	16189	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 38.57 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.6118e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ANP, 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	0.28	0/2701	0.55	0/3676
1	B	0.29	0/2715	0.54	0/3692
1	C	0.31	0/2712	0.61	0/3692
1	D	0.27	0/2783	0.54	0/3784
1	E	0.28	0/2708	0.56	0/3687
1	F	0.29	0/2667	0.60	1/3632 (0.0%)
All	All	0.29	0/16286	0.57	1/22163 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	195	ARG	CB-CG-CD	5.00	124.61	111.60

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	2656	0	2642	26	0
1	B	2668	0	2666	16	0
1	C	2665	0	2653	16	0
1	D	2734	0	2729	25	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2661	0	2642	26	0
1	F	2623	0	2593	27	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
3	A	31	0	13	1	0
3	B	31	0	13	2	0
3	D	31	0	13	0	0
3	E	31	0	13	1	0
4	C	27	0	11	1	0
4	F	27	0	11	1	0
All	All	16189	0	15999	122	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 4.

All (122) 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:327:LEU:O	1:A:330:ARG:NH1	2.16	0.78
1:D:161:ASP:O	1:E:195:ARG:NH2	2.19	0.75
1:A:44:ASN:OD1	1:A:45:ALA:N	2.20	0.73
1:B:29:ILE:HD12	1:B:38:LEU:HD21	1.71	0.73
1:C:98:ARG:HH22	1:D:195:ARG:C	1.92	0.72
1:A:162:PRO:HD3	1:B:195:ARG:HH12	1.55	0.70
1:A:162:PRO:HD3	1:B:195:ARG:NH1	2.07	0.70
1:C:83:ARG:HB2	1:C:100:ILE:HD11	1.74	0.70
1:C:267:GLN:HB3	1:C:280:LEU:HD11	1.74	0.70
1:F:83:ARG:HB2	1:F:100:ILE:HD11	1.74	0.68
1:F:212:ILE:HD11	1:F:241:ILE:HG23	1.75	0.67
1:E:117:ARG:O	1:E:147:LYS:NZ	2.28	0.67
1:E:309:VAL:O	1:E:313:LYS:NZ	2.28	0.67
1:C:29:ILE:HD12	1:C:38:LEU:HD21	1.76	0.66
1:A:230:HIS:CG	1:B:221:THR:HG22	2.31	0.66
1:E:219:ALA:HB1	1:E:260:VAL:HB	1.79	0.65
1:F:195:ARG:HG2	1:F:195:ARG:HH11	1.63	0.63
1:F:209:LEU:O	1:F:212:ILE:HG22	2.00	0.62
1:D:115:VAL:HG11	1:D:284:VAL:HG11	1.81	0.62
1:A:278:ARG:NH2	3:A:402:ANP:O4'	2.32	0.62
1:A:345:ASP:OD1	1:A:346:GLU:N	2.33	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:112:LEU:HB3	1:D:116:VAL:HG11	1.82	0.61
1:E:267:GLN:HB3	1:E:280:LEU:HD11	1.82	0.61
1:E:345:ASP:OD1	1:E:346:GLU:N	2.33	0.61
1:A:267:GLN:HB3	1:A:280:LEU:HD11	1.83	0.61
1:A:195:ARG:HE	1:F:83:ARG:HD3	1.65	0.60
1:A:240:ARG:NH1	1:B:220:GLU:OE1	2.33	0.60
1:B:137:LYS:NZ	3:B:402:ANP:O1B	2.22	0.60
1:E:339:GLY:O	1:E:340:ARG:NH1	2.35	0.60
1:D:67:ASN:HA	1:E:184:LYS:HE3	1.83	0.59
1:F:129:VAL:HB	1:F:141:LEU:HD11	1.85	0.59
1:D:314:PHE:HB2	1:D:316:MET:HG3	1.84	0.59
1:D:157:VAL:HG21	1:D:193:ILE:HD11	1.85	0.59
1:F:195:ARG:HG2	1:F:195:ARG:NH1	2.19	0.58
1:D:113:PRO:O	1:D:116:VAL:HG12	2.05	0.57
1:F:267:GLN:HB3	1:F:280:LEU:HD11	1.86	0.57
1:C:314:PHE:HB2	1:C:316:MET:HG3	1.85	0.57
1:A:115:VAL:HG11	1:A:284:VAL:HG11	1.85	0.57
1:B:115:VAL:HG11	1:B:284:VAL:HG11	1.87	0.56
1:D:267:GLN:HB3	1:D:280:LEU:HD11	1.86	0.56
1:F:314:PHE:HB2	1:F:316:MET:HG3	1.88	0.56
1:B:314:PHE:HB2	1:B:316:MET:HG3	1.88	0.55
1:F:115:VAL:HG11	1:F:284:VAL:HG11	1.88	0.55
1:A:67:ASN:HA	1:B:184:LYS:HE3	1.89	0.53
1:E:4:MET:CE	1:E:51:LEU:HB3	2.38	0.53
1:F:29:ILE:HD12	1:F:38:LEU:HD21	1.91	0.53
1:E:346:GLU:O	1:E:350:MET:HG3	2.08	0.53
1:C:115:VAL:HG11	1:C:284:VAL:HG11	1.92	0.52
1:D:4:MET:CE	1:D:51:LEU:HB3	2.39	0.52
1:D:29:ILE:HD12	1:D:38:LEU:HD21	1.91	0.52
1:E:29:ILE:HD12	1:E:38:LEU:HD21	1.92	0.52
1:F:134:GLY:HA2	4:F:401:ADP:O3A	2.10	0.52
1:F:324:TYR:O	1:F:328:GLN:HG2	2.09	0.52
1:C:4:MET:CE	1:C:51:LEU:HB3	2.41	0.51
1:C:240:ARG:O	1:C:244:VAL:HG13	2.11	0.51
1:E:216:LEU:HD22	1:E:256:GLN:HB3	1.92	0.51
1:A:4:MET:CE	1:A:51:LEU:HB3	2.40	0.50
1:E:240:ARG:O	1:E:244:VAL:HG13	2.11	0.50
1:E:179:VAL:HG21	1:E:186:PHE:CD1	2.46	0.50
1:E:115:VAL:HG11	1:E:284:VAL:HG11	1.93	0.50
1:D:206:LEU:HD12	1:D:229:LEU:HD11	1.94	0.50
1:A:116:VAL:HG12	1:A:144:ILE:HD11	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:ILE:HD12	1:A:38:LEU:HD21	1.92	0.50
1:E:256:GLN:O	1:E:260:VAL:HG22	2.12	0.49
1:F:4:MET:CE	1:F:51:LEU:HB3	2.42	0.49
1:A:206:LEU:HD12	1:A:229:LEU:HD11	1.93	0.49
1:A:240:ARG:O	1:A:244:VAL:HG13	2.14	0.48
1:E:314:PHE:HB2	1:E:316:MET:HG3	1.94	0.48
1:F:83:ARG:NH2	1:F:164:GLU:OE1	2.46	0.48
1:B:3:ASN:OD1	1:B:4:MET:N	2.46	0.48
1:B:131:GLY:O	1:B:230:HIS:HA	2.13	0.47
1:D:346:GLU:O	1:D:350:MET:HG3	2.13	0.47
1:E:10:GLU:OE1	1:E:38:LEU:HD13	2.14	0.47
1:D:235:VAL:HG13	1:D:300:ILE:HG22	1.96	0.47
1:C:137:LYS:HB2	4:C:401:ADP:O2B	2.15	0.47
1:F:73:SER:OG	1:F:163:ILE:O	2.28	0.46
1:C:307:MET:HG2	1:C:316:MET:HE1	1.97	0.46
1:D:307:MET:HG2	1:D:316:MET:HE1	1.97	0.46
1:C:131:GLY:O	1:C:230:HIS:HA	2.15	0.46
1:C:67:ASN:HA	1:D:184:LYS:HE3	1.98	0.46
1:D:157:VAL:CG2	1:D:193:ILE:HD11	2.46	0.46
1:F:137:LYS:O	1:F:141:LEU:HD13	2.17	0.45
1:E:131:GLY:O	1:E:230:HIS:HA	2.16	0.45
1:A:184:LYS:HE3	1:F:67:ASN:HA	1.99	0.45
1:D:240:ARG:O	1:D:244:VAL:HG13	2.16	0.45
1:A:131:GLY:O	1:A:230:HIS:HA	2.17	0.45
1:F:206:LEU:HD12	1:F:229:LEU:HD11	1.99	0.44
1:D:256:GLN:O	1:D:260:VAL:HG22	2.18	0.44
1:F:307:MET:HG2	1:F:316:MET:HE1	2.00	0.44
1:F:240:ARG:O	1:F:244:VAL:HG13	2.17	0.44
1:C:146:ASP:OD1	1:C:170:LYS:NZ	2.44	0.44
1:C:256:GLN:O	1:C:260:VAL:HG22	2.17	0.44
1:F:235:VAL:HG13	1:F:300:ILE:HG22	1.99	0.44
1:B:195:ARG:HA	1:B:195:ARG:HD3	1.73	0.43
1:D:230:HIS:CG	1:E:221:THR:HG22	2.53	0.43
1:F:256:GLN:O	1:F:260:VAL:HG22	2.17	0.43
1:A:256:GLN:O	1:A:260:VAL:HG22	2.17	0.43
1:C:105:LEU:HD22	1:C:109:GLU:OE1	2.19	0.43
1:E:137:LYS:NZ	3:E:402:ANP:O1B	2.30	0.42
1:A:116:VAL:CG1	1:A:144:ILE:HD11	2.49	0.42
1:A:314:PHE:HB2	1:A:316:MET:HG3	2.00	0.42
1:D:230:HIS:HB3	1:E:221:THR:HG22	2.02	0.42
1:F:74:PHE:CE2	1:F:82:PHE:HD2	2.37	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:122:LYS:HD3	1:E:122:LYS:HA	1.84	0.42
1:B:240:ARG:O	1:B:244:VAL:HG13	2.19	0.42
1:A:83:ARG:HD3	1:B:195:ARG:HD2	2.01	0.41
1:B:32:ASP:OD2	1:B:276:THR:O	2.38	0.41
1:E:301:HIS:NE2	1:E:302:GLN:HG3	2.35	0.41
1:A:245:PHE:CD1	1:A:253:VAL:HG21	2.56	0.41
1:F:141:LEU:N	1:F:141:LEU:HD12	2.35	0.41
1:F:131:GLY:O	1:F:230:HIS:HA	2.21	0.41
1:B:278:ARG:NH2	3:B:402:ANP:O4'	2.54	0.41
1:E:3:ASN:O	1:E:7:LEU:HG	2.21	0.41
1:C:176:GLN:OE1	1:C:176:GLN:N	2.54	0.41
1:D:131:GLY:O	1:D:230:HIS:HA	2.20	0.41
1:D:4:MET:HE1	1:D:51:LEU:HB3	2.03	0.41
1:E:307:MET:HG2	1:E:316:MET:HE1	2.02	0.41
1:A:4:MET:HE1	1:A:51:LEU:HB3	2.03	0.40
1:D:324:TYR:O	1:D:328:GLN:HG2	2.21	0.40
1:F:141:LEU:HD23	1:F:202:LEU:HD13	2.03	0.40
1:A:235:VAL:HG13	1:A:300:ILE:HG22	2.03	0.40
1:D:321:GLN:HG2	1:D:350:MET:HE1	2.02	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	350/385 (91%)	342 (98%)	8 (2%)	0	100	100
1	B	349/385 (91%)	340 (97%)	9 (3%)	0	100	100
1	C	349/385 (91%)	342 (98%)	7 (2%)	0	100	100
1	D	356/385 (92%)	347 (98%)	9 (2%)	0	100	100
1	E	351/385 (91%)	342 (97%)	9 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	350/385 (91%)	342 (98%)	8 (2%)	0	100	100
All	All	2105/2310 (91%)	2055 (98%)	50 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	286/338 (85%)	283 (99%)	3 (1%)	76	86
1	B	289/338 (86%)	286 (99%)	3 (1%)	76	86
1	C	289/338 (86%)	286 (99%)	3 (1%)	76	86
1	D	298/338 (88%)	294 (99%)	4 (1%)	69	82
1	E	285/338 (84%)	280 (98%)	5 (2%)	59	77
1	F	278/338 (82%)	276 (99%)	2 (1%)	84	90
All	All	1725/2028 (85%)	1705 (99%)	20 (1%)	71	84

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	GLN
1	A	72	LEU
1	A	90	ARG
1	B	50	GLN
1	B	72	LEU
1	B	267	GLN
1	C	50	GLN
1	C	59	GLN
1	C	349	GLN
1	D	-1	SER
1	D	50	GLN
1	D	72	LEU
1	D	301	HIS

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Mol	Chain	Res	Type
1	E	50	GLN
1	E	59	GLN
1	E	63	LYS
1	E	72	LEU
1	E	345	ASP
1	F	50	GLN
1	F	72	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	20	HIS
1	E	20	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 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	ANP	E	402	2	29,33,33	3.51	16 (55%)	31,52,52	2.14	8 (25%)
3	ANP	B	402	2	29,33,33	3.50	16 (55%)	31,52,52	2.11	7 (22%)
3	ANP	A	402	2	29,33,33	3.50	17 (58%)	31,52,52	2.16	8 (25%)
4	ADP	C	401	-	24,29,29	4.60	6 (25%)	29,45,45	2.22	4 (13%)
3	ANP	D	402	2	29,33,33	3.50	16 (55%)	31,52,52	2.12	7 (22%)
4	ADP	F	401	-	24,29,29	4.57	7 (29%)	29,45,45	2.21	4 (13%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ANP	E	402	2	-	4/14/38/38	0/3/3/3
3	ANP	B	402	2	-	4/14/38/38	0/3/3/3
3	ANP	A	402	2	-	4/14/38/38	0/3/3/3
4	ADP	C	401	-	-	2/12/32/32	0/3/3/3
3	ANP	D	402	2	-	4/14/38/38	0/3/3/3
4	ADP	F	401	-	-	3/12/32/32	0/3/3/3

All (78) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	401	ADP	O4'-C1'	15.88	1.63	1.41
4	F	401	ADP	O4'-C1'	15.75	1.63	1.41
4	F	401	ADP	C2'-C1'	-13.33	1.33	1.53
4	C	401	ADP	C2'-C1'	-13.32	1.33	1.53
3	B	402	ANP	C3'-C4'	-7.49	1.33	1.53
3	A	402	ANP	C3'-C4'	-7.49	1.33	1.53
3	E	402	ANP	C3'-C4'	-7.47	1.33	1.53
3	D	402	ANP	C3'-C4'	-7.47	1.33	1.53
3	E	402	ANP	O4'-C1'	-7.15	1.31	1.41
3	A	402	ANP	O4'-C1'	-7.14	1.31	1.41
3	B	402	ANP	O4'-C1'	-7.11	1.31	1.41
3	D	402	ANP	O4'-C1'	-7.10	1.31	1.41
3	B	402	ANP	O4'-C4'	6.97	1.60	1.45
3	A	402	ANP	O4'-C4'	6.97	1.60	1.45
3	D	402	ANP	O4'-C4'	6.93	1.60	1.45
3	E	402	ANP	O4'-C4'	6.90	1.60	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	402	ANP	PB-O3A	6.41	1.67	1.59
3	A	402	ANP	PB-O3A	6.26	1.67	1.59
3	D	402	ANP	PB-O3A	6.24	1.66	1.59
3	B	402	ANP	PB-O3A	6.12	1.66	1.59
4	F	401	ADP	O4'-C4'	-5.45	1.32	1.45
4	C	401	ADP	O4'-C4'	-5.44	1.32	1.45
3	E	402	ANP	PG-N3B	5.07	1.76	1.63
3	B	402	ANP	PG-N3B	5.03	1.76	1.63
3	A	402	ANP	PG-N3B	4.99	1.76	1.63
3	D	402	ANP	PG-N3B	4.91	1.76	1.63
3	D	402	ANP	PB-O1B	4.67	1.53	1.46
3	A	402	ANP	PB-O1B	4.65	1.53	1.46
3	E	402	ANP	PB-O1B	4.64	1.53	1.46
3	B	402	ANP	PB-O1B	4.57	1.53	1.46
3	D	402	ANP	C6-N6	4.39	1.50	1.34
3	B	402	ANP	C6-N6	4.37	1.50	1.34
3	E	402	ANP	C6-N6	4.35	1.50	1.34
3	A	402	ANP	C6-N6	4.33	1.49	1.34
3	B	402	ANP	C2-N3	4.29	1.39	1.32
3	A	402	ANP	C2-N3	4.28	1.39	1.32
3	D	402	ANP	C2-N3	4.25	1.39	1.32
3	E	402	ANP	C2-N3	4.25	1.38	1.32
4	C	401	ADP	C6-N6	3.78	1.47	1.34
4	F	401	ADP	C6-N6	3.75	1.47	1.34
3	E	402	ANP	PG-O1G	3.40	1.51	1.46
3	A	402	ANP	PG-O1G	3.37	1.51	1.46
3	D	402	ANP	PG-O1G	3.35	1.51	1.46
3	B	402	ANP	PG-O1G	3.33	1.51	1.46
3	B	402	ANP	PB-N3B	3.25	1.71	1.63
3	E	402	ANP	PB-N3B	3.24	1.71	1.63
3	B	402	ANP	PA-O5'	3.24	1.72	1.59
3	A	402	ANP	PA-O5'	3.24	1.72	1.59
3	D	402	ANP	PA-O5'	3.23	1.72	1.59
3	E	402	ANP	PA-O5'	3.21	1.72	1.59
3	D	402	ANP	PB-N3B	3.20	1.71	1.63
3	A	402	ANP	PB-N3B	3.18	1.71	1.63
3	D	402	ANP	O3'-C3'	3.08	1.50	1.43
3	B	402	ANP	O3'-C3'	3.06	1.50	1.43
3	A	402	ANP	O3'-C3'	3.02	1.50	1.43
3	E	402	ANP	O3'-C3'	3.01	1.50	1.43
4	F	401	ADP	O3'-C3'	-2.79	1.36	1.43
4	C	401	ADP	O3'-C3'	-2.73	1.36	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	ANP	PG-O2G	-2.59	1.49	1.56
4	C	401	ADP	C5-C4	-2.56	1.34	1.40
4	F	401	ADP	C5-C4	-2.55	1.34	1.40
3	A	402	ANP	PG-O2G	-2.55	1.49	1.56
3	D	402	ANP	PG-O2G	-2.53	1.49	1.56
3	D	402	ANP	PG-O3G	-2.51	1.50	1.56
3	B	402	ANP	PG-O3G	-2.51	1.50	1.56
3	E	402	ANP	PG-O2G	-2.51	1.50	1.56
3	A	402	ANP	PG-O3G	-2.49	1.50	1.56
3	E	402	ANP	PG-O3G	-2.46	1.50	1.56
3	E	402	ANP	PA-O2A	-2.12	1.45	1.55
3	D	402	ANP	O2'-C2'	-2.11	1.38	1.43
3	B	402	ANP	PA-O2A	-2.09	1.45	1.55
3	D	402	ANP	PA-O2A	-2.08	1.45	1.55
3	A	402	ANP	PA-O2A	-2.07	1.45	1.55
3	A	402	ANP	O2'-C2'	-2.06	1.38	1.43
3	B	402	ANP	O2'-C2'	-2.06	1.38	1.43
4	F	401	ADP	C2-N3	2.04	1.35	1.32
3	E	402	ANP	O2'-C2'	-2.04	1.38	1.43
3	A	402	ANP	C5-C4	-2.02	1.35	1.40

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	401	ADP	C5-C6-N6	8.05	132.59	120.35
4	F	401	ADP	C5-C6-N6	7.78	132.18	120.35
3	A	402	ANP	C1'-N9-C4	-7.42	113.60	126.64
3	E	402	ANP	C1'-N9-C4	-7.37	113.70	126.64
3	D	402	ANP	C1'-N9-C4	-7.29	113.83	126.64
3	B	402	ANP	C1'-N9-C4	-7.12	114.14	126.64
4	F	401	ADP	N3-C2-N1	-5.56	119.99	128.68
3	A	402	ANP	N3-C2-N1	-5.42	120.21	128.68
3	B	402	ANP	N3-C2-N1	-5.37	120.29	128.68
4	C	401	ADP	N3-C2-N1	-5.36	120.30	128.68
3	E	402	ANP	N3-C2-N1	-5.32	120.36	128.68
4	F	401	ADP	N6-C6-N1	-5.26	107.66	118.57
3	D	402	ANP	N3-C2-N1	-5.23	120.50	128.68
4	C	401	ADP	N6-C6-N1	-5.16	107.86	118.57
3	D	402	ANP	C3'-C2'-C1'	3.66	106.48	100.98
3	B	402	ANP	C3'-C2'-C1'	3.64	106.46	100.98
3	A	402	ANP	C3'-C2'-C1'	3.62	106.42	100.98
3	E	402	ANP	C3'-C2'-C1'	3.45	106.17	100.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	402	ANP	C5-C6-N6	2.93	124.81	120.35
3	B	402	ANP	C5-C6-N6	2.73	124.50	120.35
3	E	402	ANP	C5-C6-N6	2.73	124.50	120.35
3	A	402	ANP	C5-C6-N6	2.71	124.47	120.35
3	E	402	ANP	O2G-PG-O1G	-2.69	106.69	113.45
3	B	402	ANP	O2G-PG-O1G	-2.58	106.98	113.45
4	C	401	ADP	C3'-C2'-C1'	2.40	104.58	100.98
3	A	402	ANP	O1G-PG-N3B	-2.36	108.29	111.77
3	A	402	ANP	O2G-PG-O1G	-2.33	107.60	113.45
3	B	402	ANP	PA-O3A-PB	-2.25	124.69	132.62
3	E	402	ANP	O1G-PG-N3B	-2.23	108.48	111.77
3	A	402	ANP	C2'-C3'-C4'	2.13	106.78	102.64
4	F	401	ADP	C3'-C2'-C1'	2.13	104.18	100.98
3	D	402	ANP	O2G-PG-O1G	-2.12	108.12	113.45
3	B	402	ANP	C2'-C3'-C4'	2.11	106.75	102.64
3	E	402	ANP	PA-O3A-PB	-2.10	125.21	132.62
3	E	402	ANP	C2'-C3'-C4'	2.08	106.68	102.64
3	D	402	ANP	O1G-PG-N3B	-2.06	108.73	111.77
3	D	402	ANP	PA-O3A-PB	-2.06	125.36	132.62
3	A	402	ANP	PA-O3A-PB	-2.00	125.57	132.62

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	402	ANP	PG-N3B-PB-O1B
3	D	402	ANP	PA-O3A-PB-O1B
3	E	402	ANP	PG-N3B-PB-O1B
3	E	402	ANP	PA-O3A-PB-O1B
3	E	402	ANP	PA-O3A-PB-O2B
3	B	402	ANP	PG-N3B-PB-O1B
3	B	402	ANP	PA-O3A-PB-O1B
3	A	402	ANP	PG-N3B-PB-O1B
3	A	402	ANP	PA-O3A-PB-O1B
4	F	401	ADP	PA-O3A-PB-O2B
3	D	402	ANP	O4'-C4'-C5'-O5'
3	E	402	ANP	O4'-C4'-C5'-O5'
4	F	401	ADP	O4'-C4'-C5'-O5'
3	B	402	ANP	O4'-C4'-C5'-O5'
3	A	402	ANP	O4'-C4'-C5'-O5'
4	C	401	ADP	PA-O3A-PB-O3B
4	C	401	ADP	O4'-C4'-C5'-O5'

*Continued on next page...*

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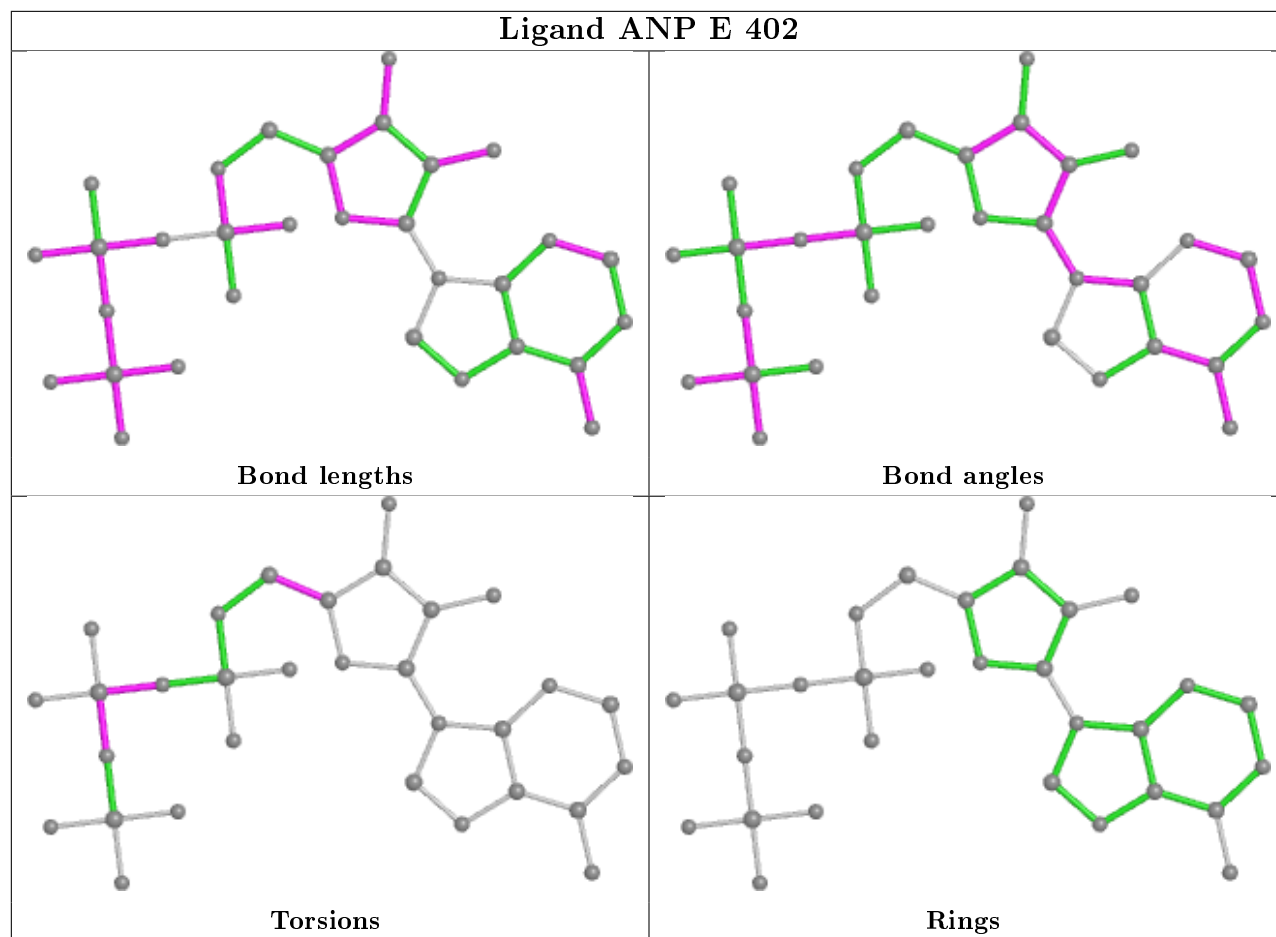
Mol	Chain	Res	Type	Atoms
3	D	402	ANP	PA-O3A-PB-O2B
3	B	402	ANP	PA-O3A-PB-O2B
3	A	402	ANP	PA-O3A-PB-O2B
4	F	401	ADP	PA-O3A-PB-O1B

There are no ring outliers.

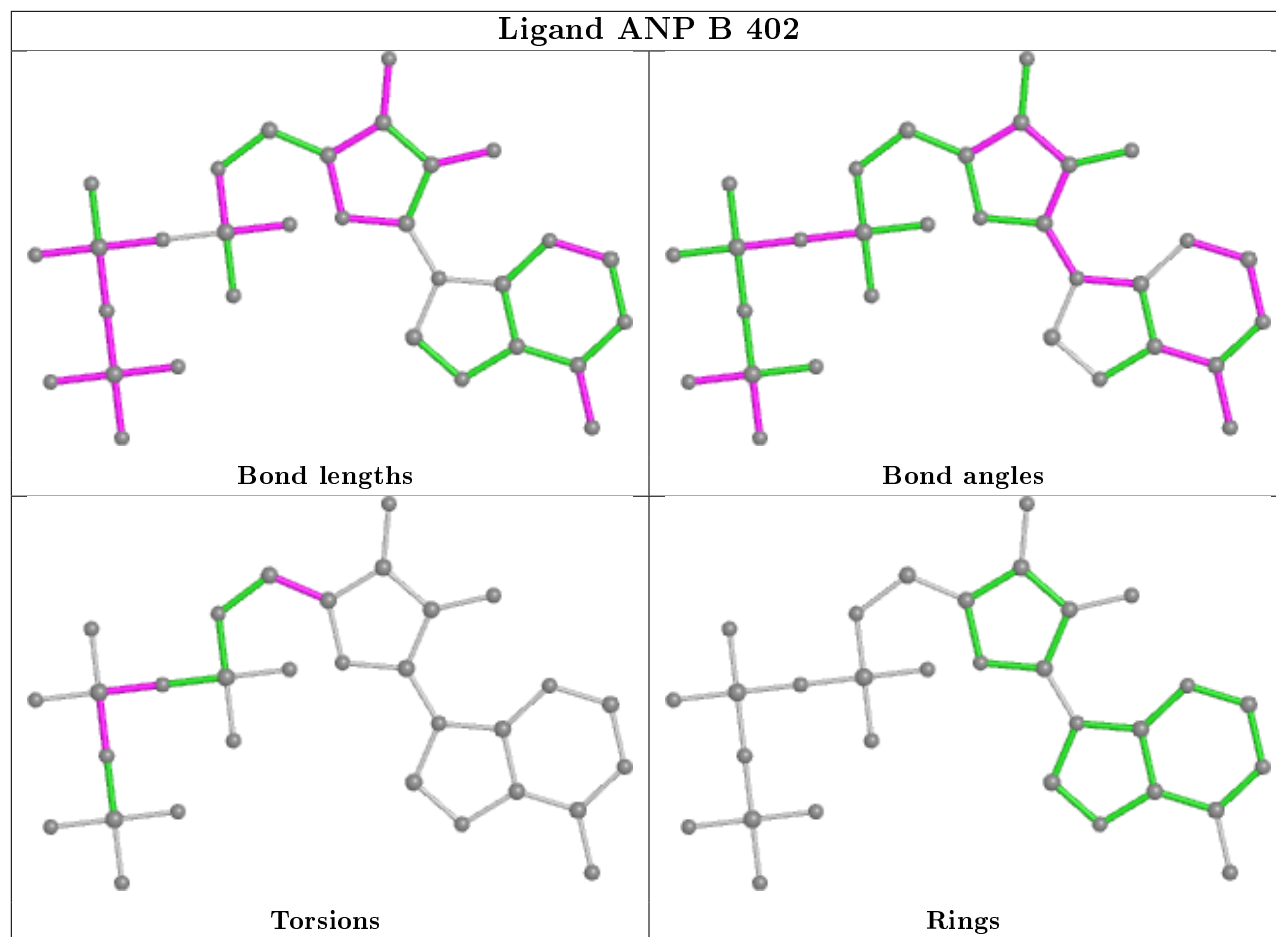
5 monomers are involved in 6 short contacts:

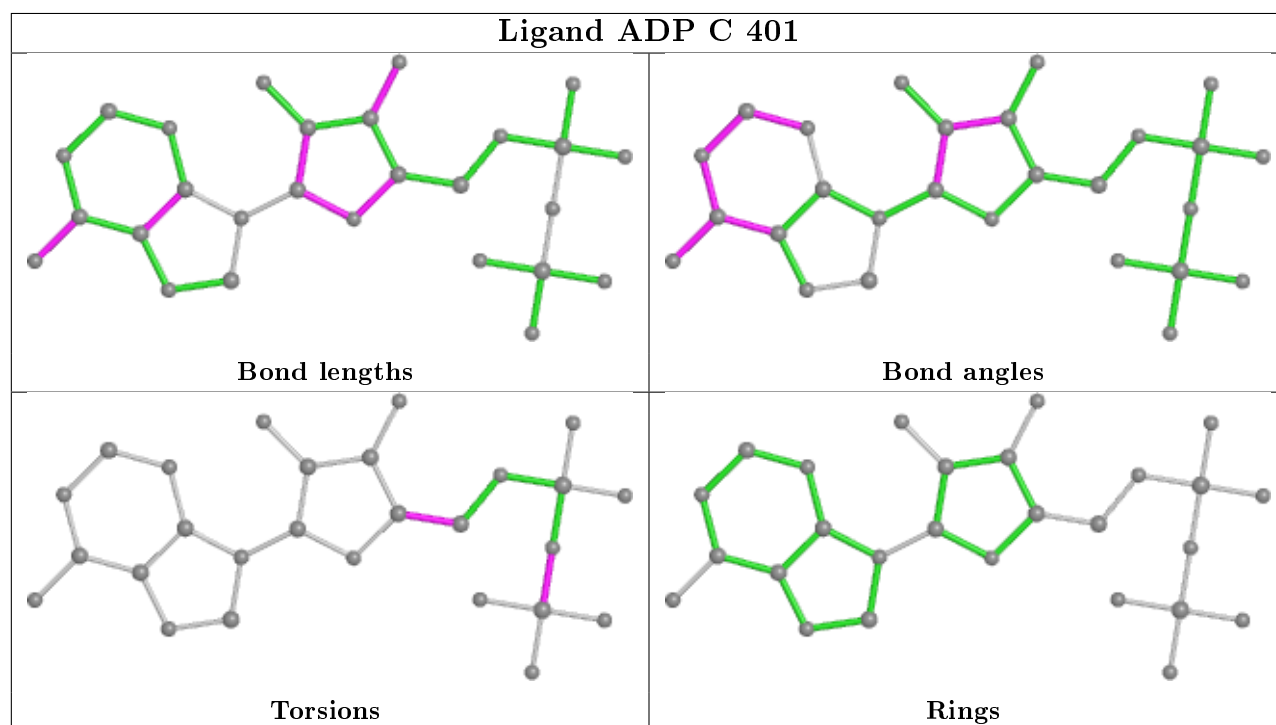
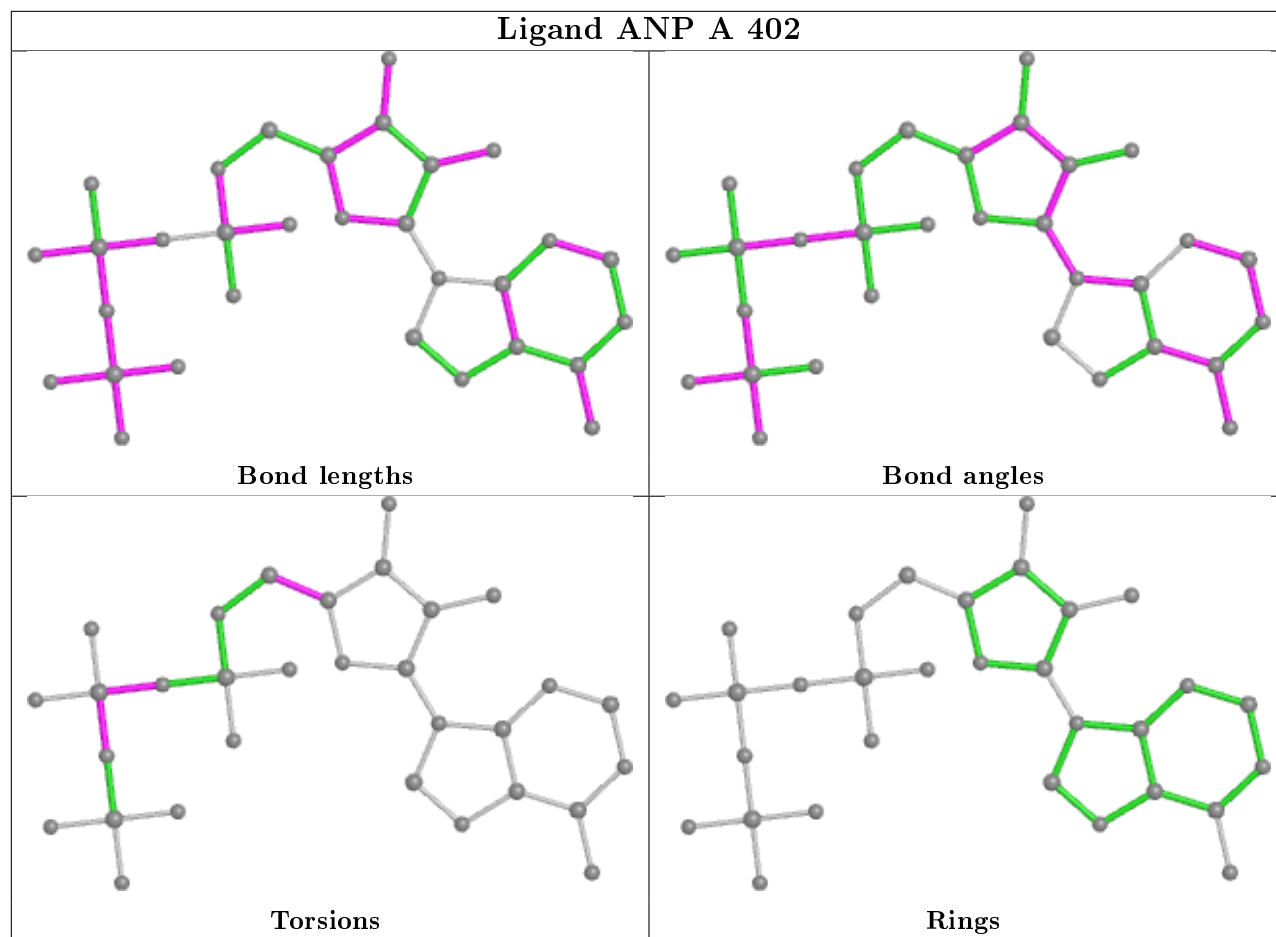
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	402	ANP	1	0
3	B	402	ANP	2	0
3	A	402	ANP	1	0
4	C	401	ADP	1	0
4	F	401	ADP	1	0

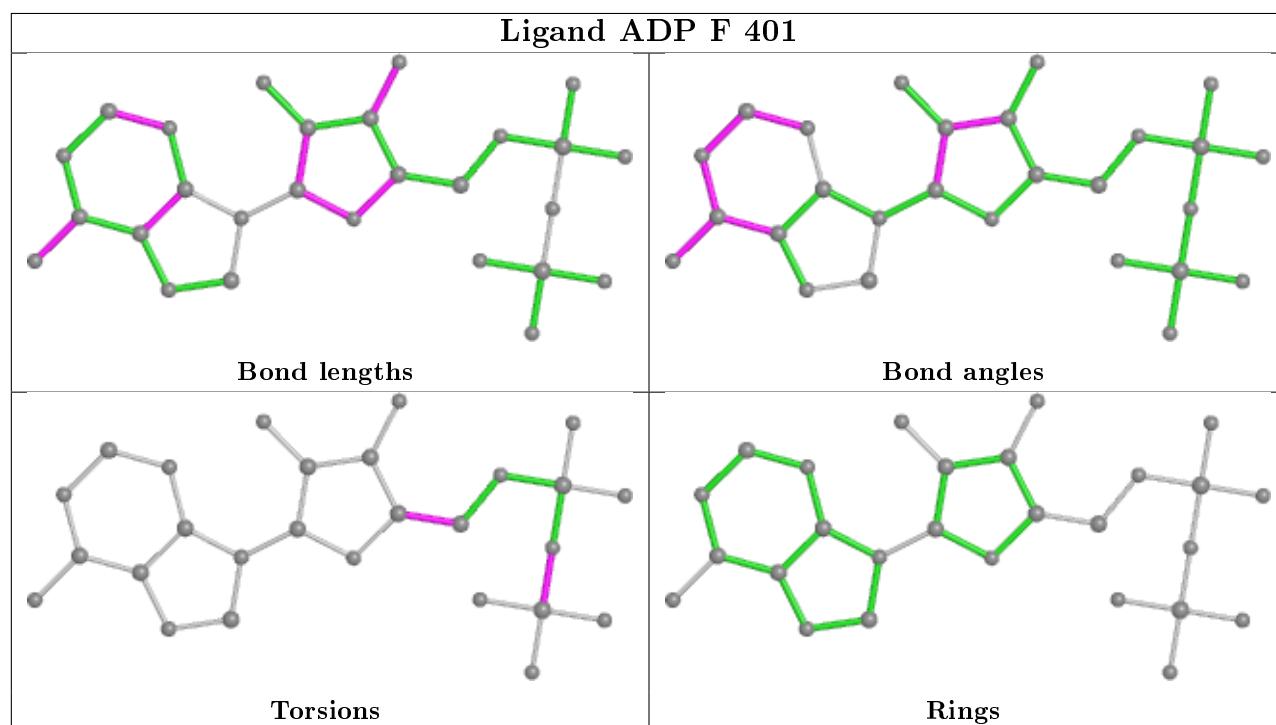
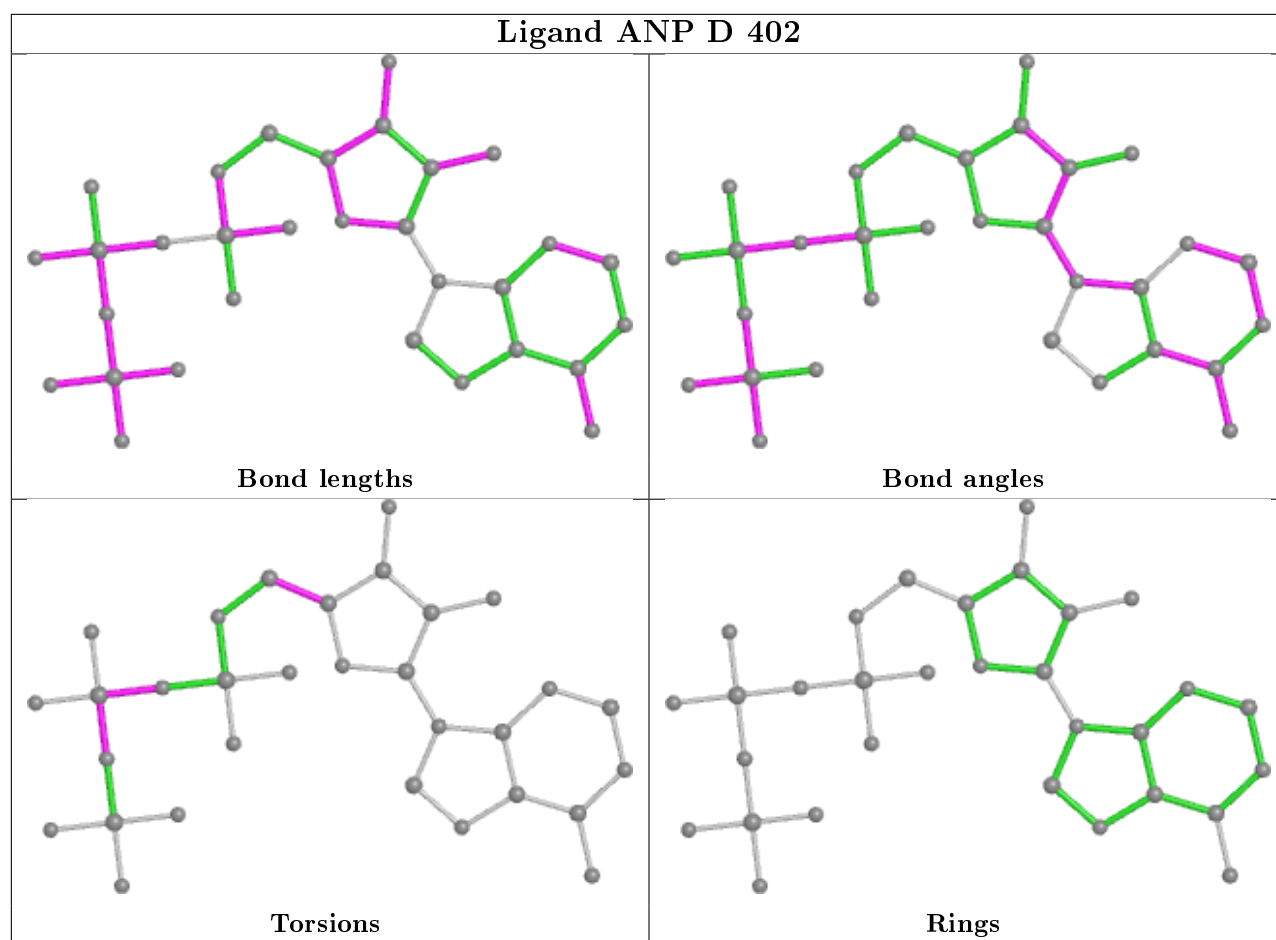
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	352/385 (91%)	-0.32	0 100 100	48, 90, 126, 141	0
1	B	351/385 (91%)	-0.34	0 100 100	73, 95, 128, 136	0
1	C	351/385 (91%)	-0.22	1 (0%) 94 90	64, 86, 118, 132	0
1	D	358/385 (92%)	-0.43	0 100 100	54, 76, 118, 157	0
1	E	353/385 (91%)	-0.35	0 100 100	57, 77, 115, 147	0
1	F	352/385 (91%)	-0.27	1 (0%) 94 90	54, 77, 144, 175	0
All	All	2117/2310 (91%)	-0.32	2 (0%) 95 94	48, 85, 126, 175	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	263	GLY	2.6
1	F	263	GLY	2.4

### 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](#)

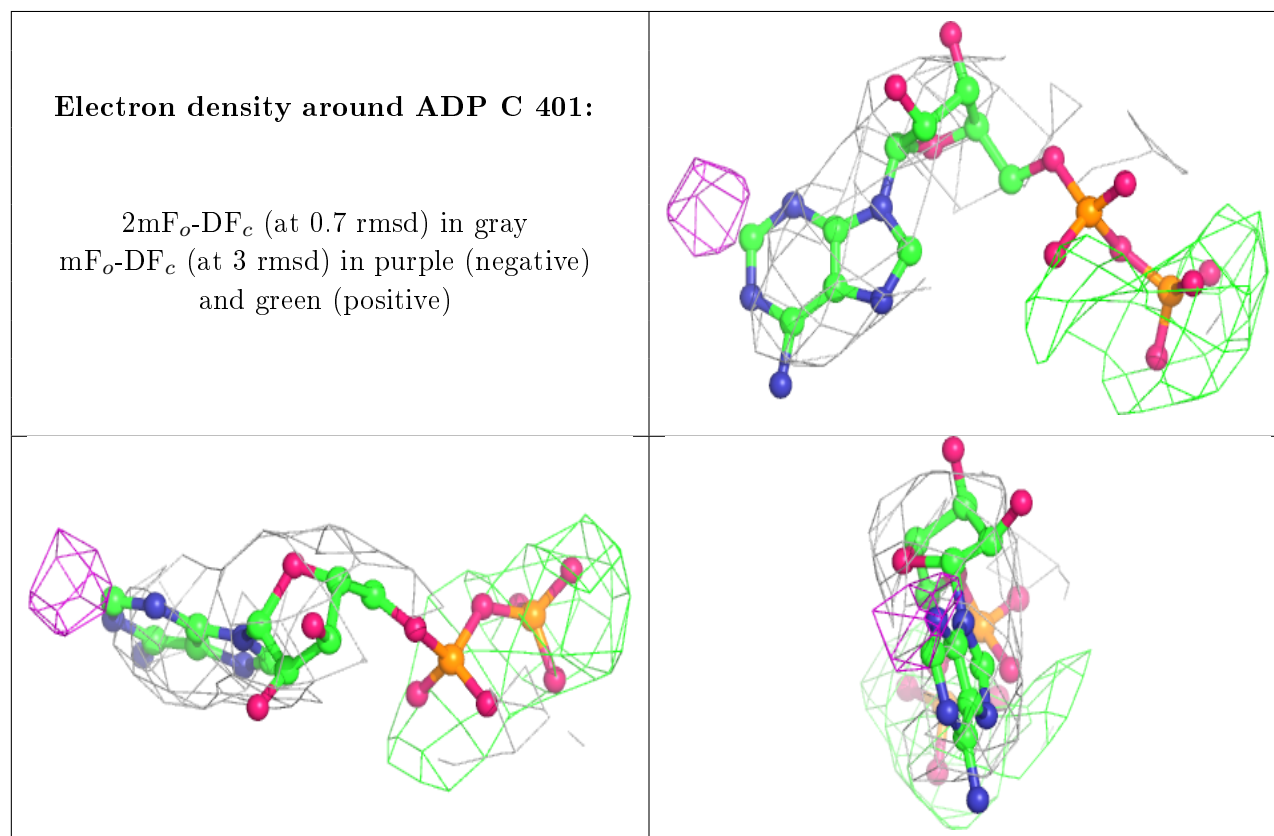
There are no carbohydrates in this entry.

### 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, 95<sup>th</sup> 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.

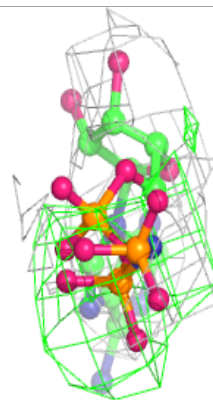
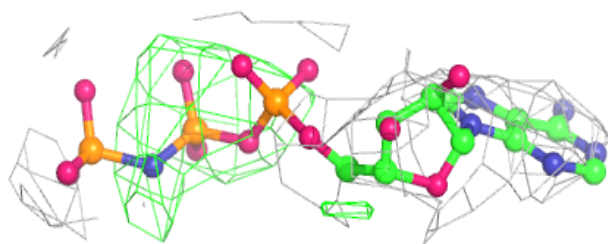
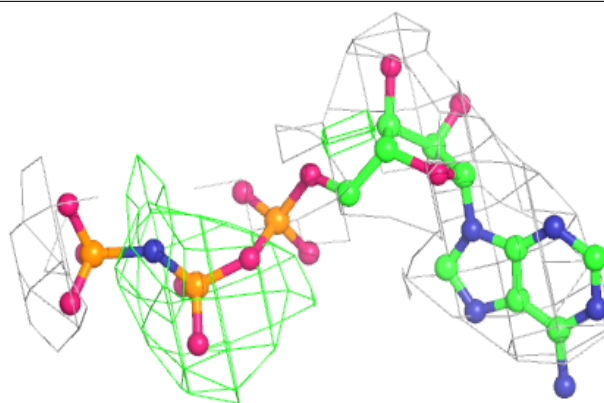
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	ADP	C	401	27/27	0.81	0.30	93,97,99,100	0
3	ANP	D	402	31/31	0.84	0.31	92,95,98,99	11
4	ADP	F	401	27/27	0.86	0.24	75,77,78,78	0
3	ANP	E	402	31/31	0.90	0.25	60,61,62,63	11
3	ANP	B	402	31/31	0.90	0.26	61,63,64,65	11
3	ANP	A	402	31/31	0.91	0.23	48,51,55,55	11
2	MG	E	401	1/1	0.94	0.17	60,60,60,60	0
2	MG	D	401	1/1	0.95	0.27	96,96,96,96	0
2	MG	B	401	1/1	0.95	0.25	60,60,60,60	0
2	MG	A	401	1/1	0.97	0.27	53,53,53,53	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.

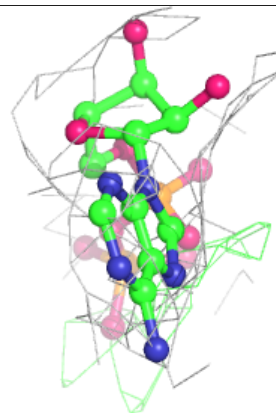
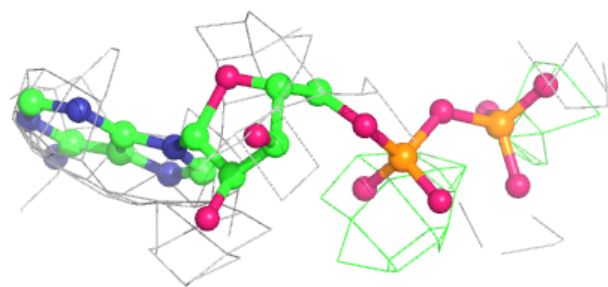
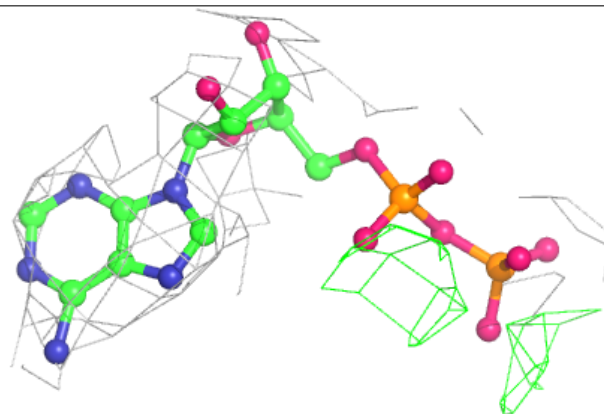


**Electron density around ANP D 402:**

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

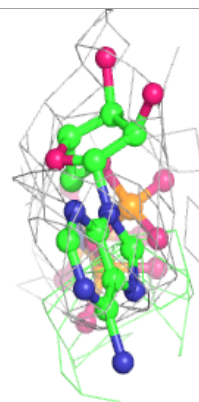
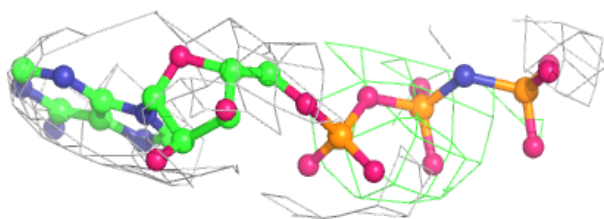
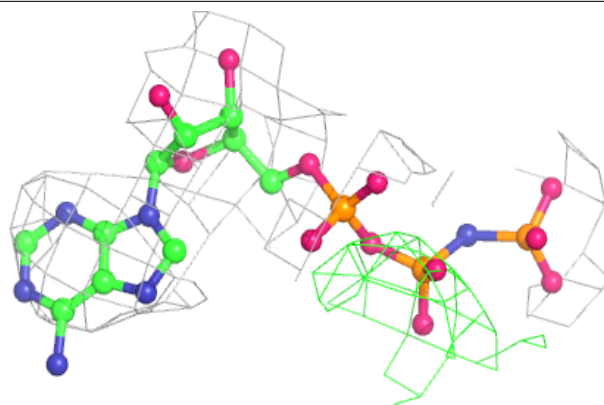
**Electron density around ADP F 401:**

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

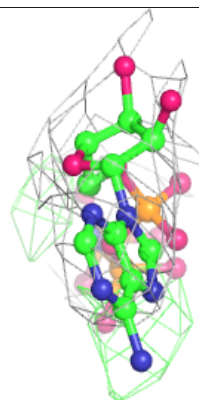
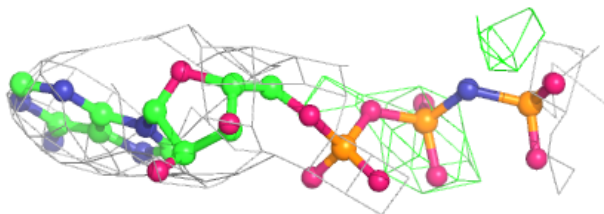
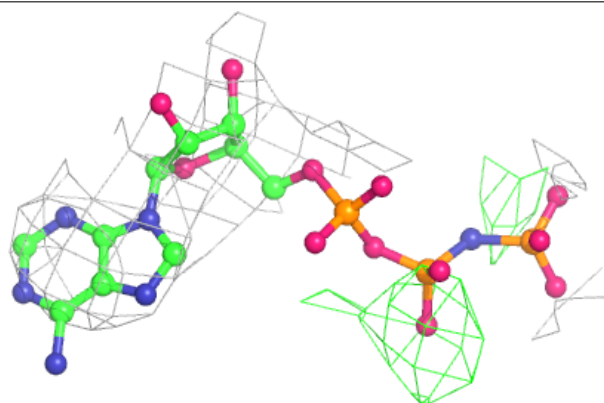


**Electron density around ANP E 402:**

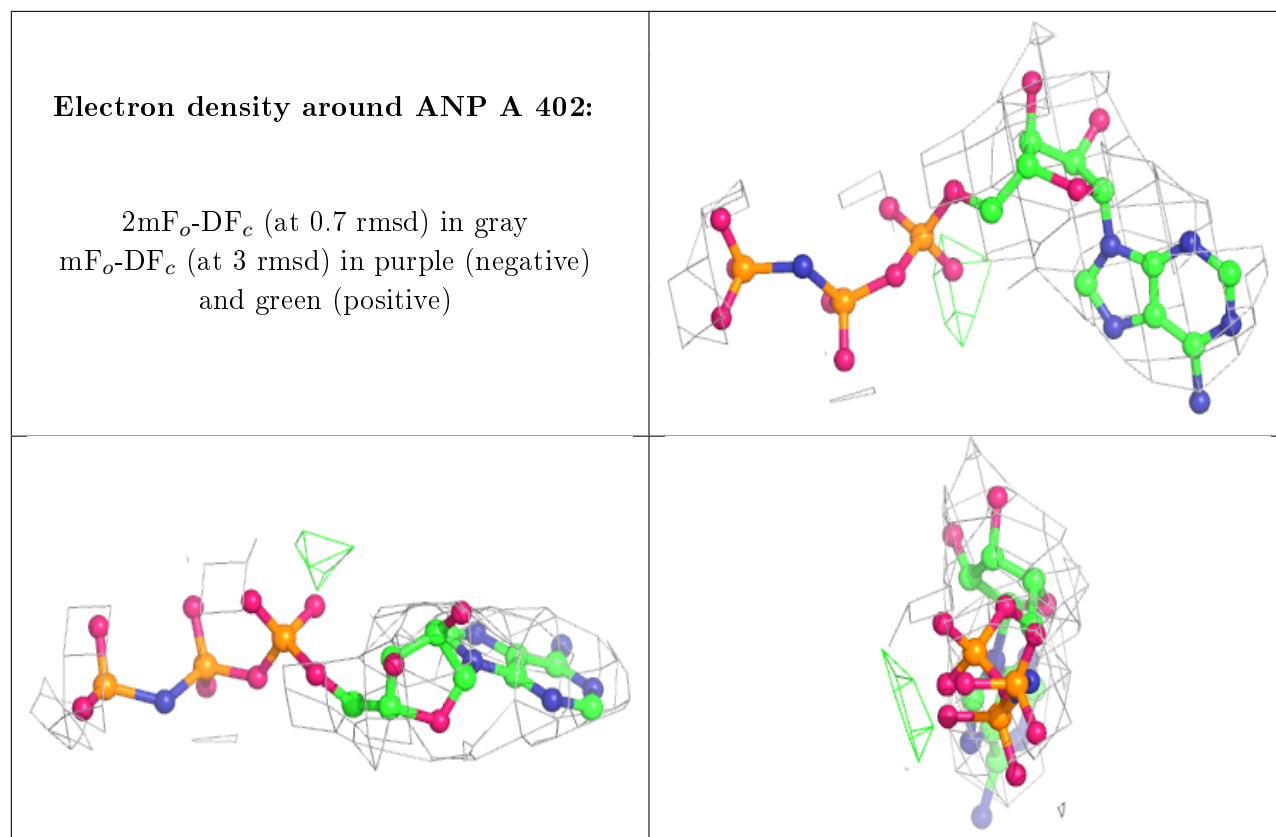
$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around ANP B 402:**

$2mF_o - DF_c$  (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.