



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 12, 2026 – 02:10 PM UTC

PDB ID : 1MX0 / pdb\_00001mx0  
Title : Structure of topoisomerase subunit  
Authors : Corbett, K.D.; Berger, J.M.  
Deposited on : 2002-10-01  
Resolution : 2.30 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](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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtrriage (Phenix) : 2.0  
EDS : 3.0  
Buster-report : wwPDB partial adaption of 1.1.7 (2018)  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

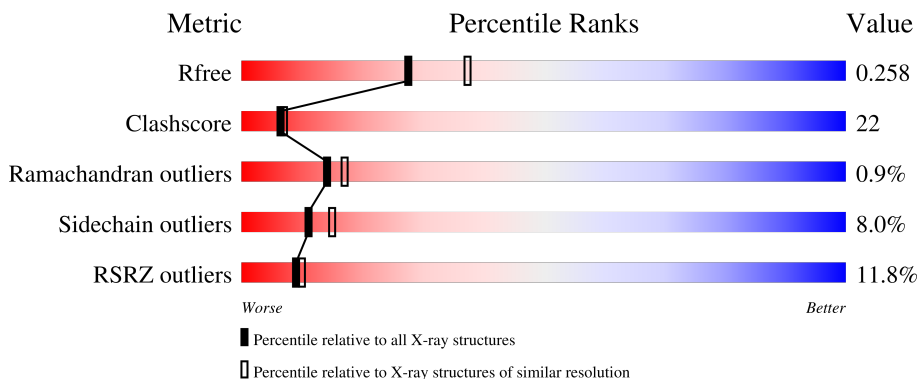
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.




Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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 of 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	472	 5% 62% 33% . .
1	B	472	 2% 66% 28% . .
1	C	472	 2% 68% 28% . .
1	D	472	 2% 71% 22% . .
1	E	472	 3% 63% 29% 5% .

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Mol	Chain	Length	Quality of chain
1	F	472	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into four segments: a red segment (53%), a green segment (39%), a yellow segment (45%), and an orange segment (12%). The segments are stacked horizontally, with the red segment on the left, followed by green, yellow, and orange. The percentages are labeled below each segment. The total length of the bar is 472 residues.</p>

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 23057 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 Type II DNA topoisomerase VI subunit B.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	461	3704	2384	620	694	1	5	0	0	0
1	B	455	3651	2352	608	685	1	5	0	0	0
1	C	466	3738	2404	626	702	1	5	0	0	0
1	D	461	3687	2374	615	692	1	5	0	0	0
1	E	456	3677	2368	614	689	1	5	0	0	0
1	F	454	3596	2316	598	676	1	5	0	0	0

There are 54 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP O05207
A	0	ALA	-	expression tag	UNP O05207
A	107	MSE	MET	modified residue	UNP O05207
A	121	MSE	MET	modified residue	UNP O05207
A	303	TYR	ASP	SEE REMARK 999	UNP O05207
A	409	MSE	MET	modified residue	UNP O05207
A	412	MSE	MET	modified residue	UNP O05207
A	435	ASP	ASN	SEE REMARK 999	UNP O05207
A	445	MSE	MET	modified residue	UNP O05207
B	-1	GLY	-	expression tag	UNP O05207
B	0	ALA	-	expression tag	UNP O05207
B	107	MSE	MET	modified residue	UNP O05207
B	121	MSE	MET	modified residue	UNP O05207
B	303	TYR	ASP	SEE REMARK 999	UNP O05207
B	409	MSE	MET	modified residue	UNP O05207
B	412	MSE	MET	modified residue	UNP O05207
B	435	ASP	ASN	SEE REMARK 999	UNP O05207

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Chain	Residue	Modelled	Actual	Comment	Reference
B	445	MSE	MET	modified residue	UNP O05207
C	-1	GLY	-	expression tag	UNP O05207
C	0	ALA	-	expression tag	UNP O05207
C	107	MSE	MET	modified residue	UNP O05207
C	121	MSE	MET	modified residue	UNP O05207
C	303	TYR	ASP	SEE REMARK 999	UNP O05207
C	409	MSE	MET	modified residue	UNP O05207
C	412	MSE	MET	modified residue	UNP O05207
C	435	ASP	ASN	SEE REMARK 999	UNP O05207
C	445	MSE	MET	modified residue	UNP O05207
D	-1	GLY	-	expression tag	UNP O05207
D	0	ALA	-	expression tag	UNP O05207
D	107	MSE	MET	modified residue	UNP O05207
D	121	MSE	MET	modified residue	UNP O05207
D	303	TYR	ASP	SEE REMARK 999	UNP O05207
D	409	MSE	MET	modified residue	UNP O05207
D	412	MSE	MET	modified residue	UNP O05207
D	435	ASP	ASN	SEE REMARK 999	UNP O05207
D	445	MSE	MET	modified residue	UNP O05207
E	-1	GLY	-	expression tag	UNP O05207
E	0	ALA	-	expression tag	UNP O05207
E	107	MSE	MET	modified residue	UNP O05207
E	121	MSE	MET	modified residue	UNP O05207
E	303	TYR	ASP	SEE REMARK 999	UNP O05207
E	409	MSE	MET	modified residue	UNP O05207
E	412	MSE	MET	modified residue	UNP O05207
E	435	ASP	ASN	SEE REMARK 999	UNP O05207
E	445	MSE	MET	modified residue	UNP O05207
F	-1	GLY	-	expression tag	UNP O05207
F	0	ALA	-	expression tag	UNP O05207
F	107	MSE	MET	modified residue	UNP O05207
F	121	MSE	MET	modified residue	UNP O05207
F	303	TYR	ASP	SEE REMARK 999	UNP O05207
F	409	MSE	MET	modified residue	UNP O05207
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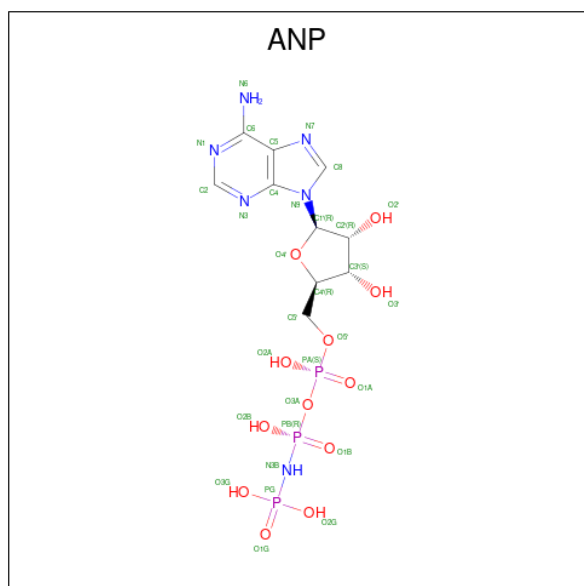
- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula:  $C_{10}H_{17}N_6O_{12}P_3$ ).



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

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

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	1	Total	Na	0	0
			1	1		

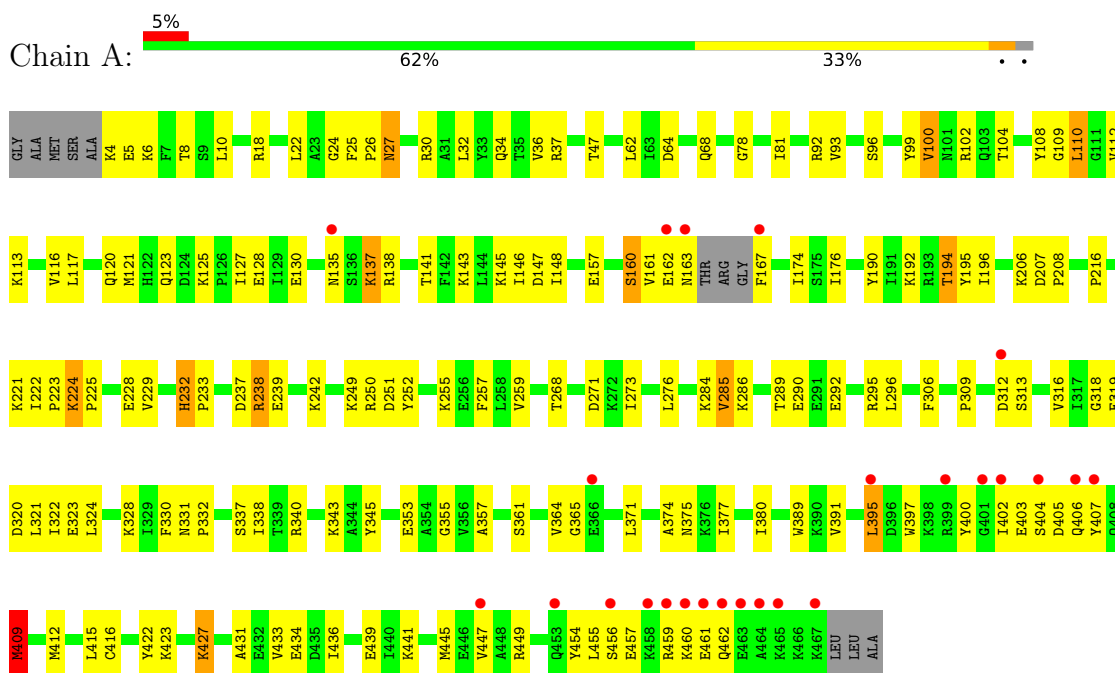
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	160	Total	O	0	0
			160	160		
5	B	145	Total	O	0	0
			145	145		
5	C	140	Total	O	0	0
			140	140		
5	D	160	Total	O	0	0
			160	160		
5	E	125	Total	O	0	0
			125	125		
5	F	81	Total	O	0	0
			81	81		

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

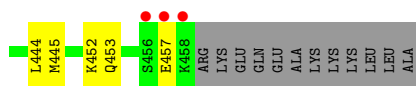
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Type II DNA topoisomerase VI subunit B

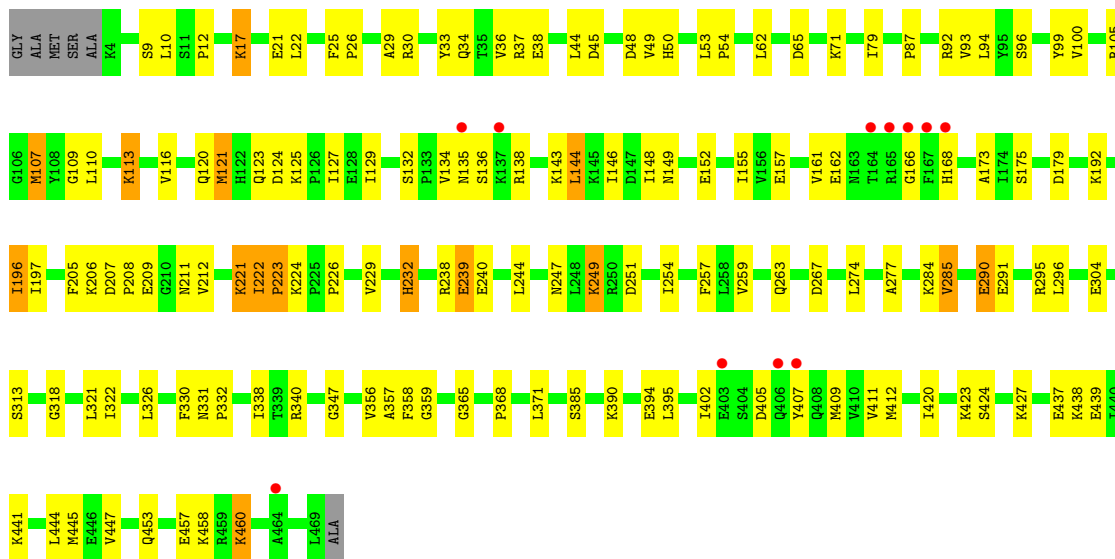


- Molecule 1: Type II DNA topoisomerase VI subunit B

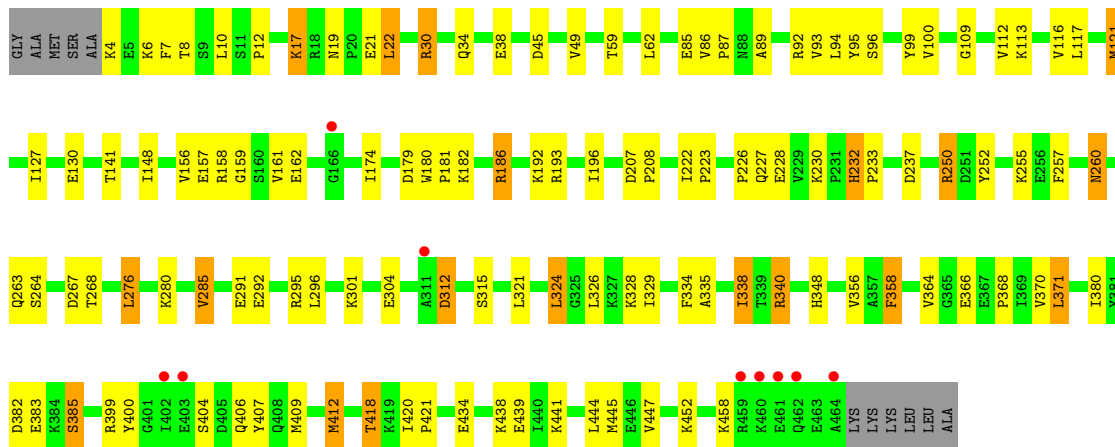




- Molecule 1: Type II DNA topoisomerase VI subunit B

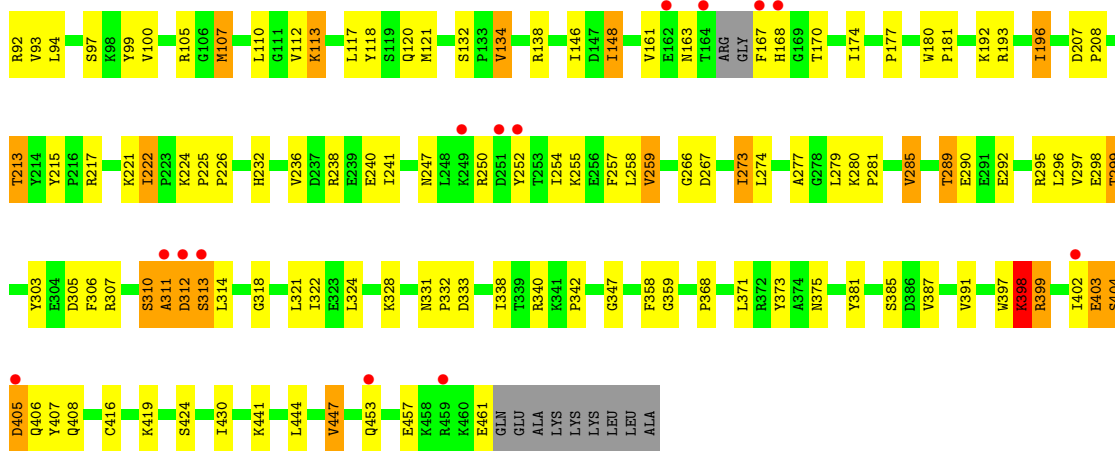


- Molecule 1: Type II DNA topoisomerase VI subunit B

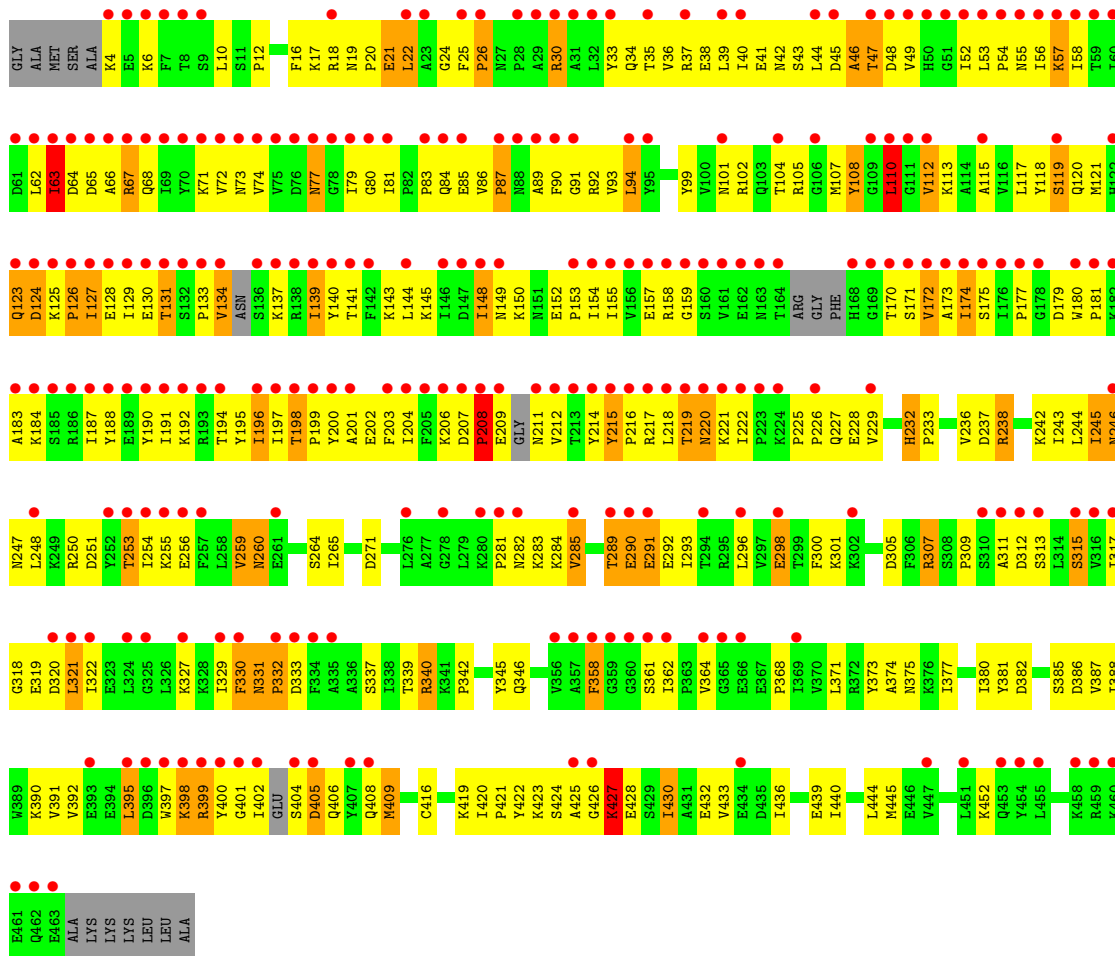


- Molecule 1: Type II DNA topoisomerase VI subunit B





• Molecule 1: Type II DNA topoisomerase VI subunit B



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	146.66Å 219.19Å 106.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 20.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.1 (20.00-2.30) 97.9 (20.00-2.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.45 (at 2.31Å)	Xtrriage
Refinement program	REFMAC 5.0	Depositor
R, $R_{free}$	0.214 , 0.263 0.216 , 0.258	Depositor DCC
$R_{free}$ test set	12540 reflections (8.36%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	26.4	Xtrriage
Anisotropy	0.171	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 49.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	23057	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.24% of the height of the origin peak. No significant pseudotranslation is detected.*

<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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NA, MG, ANP

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	0.78	0/3779	1.01	6/5104 (0.1%)
1	B	0.81	2/3727 (0.1%)	1.02	3/5038 (0.1%)
1	C	0.76	2/3814 (0.1%)	0.99	9/5152 (0.2%)
1	D	0.78	2/3763 (0.1%)	1.00	4/5087 (0.1%)
1	E	0.76	1/3752 (0.0%)	0.99	10/5067 (0.2%)
1	F	0.59	0/3666	1.06	22/4955 (0.4%)
All	All	0.75	7/22501 (0.0%)	1.01	54/30403 (0.2%)

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	412	MSE	SE-CE	-9.74	1.66	1.95
1	C	107	MSE	SE-CE	7.12	2.16	1.95
1	D	121	MSE	SE-CE	-6.80	1.75	1.95
1	E	107	MSE	SE-CE	-6.54	1.75	1.95
1	D	412	MSE	SE-CE	5.72	2.12	1.95

The worst 5 of 54 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	330	PHE	N-CA-C	-10.95	92.92	110.32
1	F	427	LYS	N-CA-C	10.10	123.64	111.02
1	F	108	TYR	N-CA-C	9.37	124.84	113.41
1	D	358	PHE	N-CA-C	9.26	123.30	110.06
1	F	110	LEU	N-CA-C	7.86	122.24	111.71

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	3704	0	3760	160	0
1	B	3651	0	3704	124	0
1	C	3738	0	3792	140	0
1	D	3687	0	3735	104	0
1	E	3677	0	3746	142	0
1	F	3596	0	3605	341	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	31	0	13	0	0
3	B	31	0	13	2	0
3	C	31	0	13	0	0
3	D	31	0	13	1	0
3	E	31	0	13	1	0
3	F	31	0	13	4	0
4	D	1	0	0	0	0
5	A	160	0	0	21	0
5	B	145	0	0	19	0
5	C	140	0	0	18	0
5	D	160	0	0	23	0
5	E	125	0	0	19	0
5	F	81	0	0	34	0
All	All	23057	0	22420	977	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 22.

The worst 5 of 977 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:107:MSE:SE	1:C:107:MSE:CE	2.16	1.43
1:C:127:ILE:HD11	1:C:144:LEU:CD1	1.56	1.34

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:LYS:HD2	1:A:4:LYS:O	1.16	1.33
1:C:127:ILE:CD1	1:C:144:LEU:HD11	1.64	1.27
1:E:402:ILE:HG22	1:E:403:GLU:O	1.30	1.25

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	457/472 (97%)	440 (96%)	17 (4%)	0	100	100
1	B	453/472 (96%)	426 (94%)	24 (5%)	3 (1%)	18	23
1	C	464/472 (98%)	448 (97%)	15 (3%)	1 (0%)	43	55
1	D	459/472 (97%)	440 (96%)	19 (4%)	0	100	100
1	E	452/472 (96%)	427 (94%)	22 (5%)	3 (1%)	18	23
1	F	444/472 (94%)	379 (85%)	48 (11%)	17 (4%)	2	1
All	All	2729/2832 (96%)	2560 (94%)	145 (5%)	24 (1%)	14	17

5 of 24 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	208	PRO
1	F	245	ILE
1	F	246	ASN
1	F	332	PRO
1	F	399	ARG

### 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	407/412 (99%)	377 (93%)	30 (7%)	13	17
1	B	402/412 (98%)	378 (94%)	24 (6%)	17	25
1	C	410/412 (100%)	384 (94%)	26 (6%)	16	23
1	D	404/412 (98%)	379 (94%)	25 (6%)	16	24
1	E	407/412 (99%)	370 (91%)	37 (9%)	9	11
1	F	389/412 (94%)	337 (87%)	52 (13%)	4	4
All	All	2419/2472 (98%)	2225 (92%)	194 (8%)	11	15

5 of 194 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	138	ARG
1	F	21	GLU
1	E	213	THR
1	E	299	THR
1	F	77	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 34 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	77	ASN
1	F	120	GLN
1	F	331	ASN
1	C	101	ASN
1	C	68	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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 13 ligands modelled in this entry, 7 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	C	920	2	33,33,33	2.42	12 (36%)	45,52,52	1.86	10 (22%)
3	ANP	F	950	2	33,33,33	2.45	13 (39%)	45,52,52	2.07	12 (26%)
3	ANP	E	940	2	33,33,33	2.35	13 (39%)	45,52,52	1.63	10 (22%)
3	ANP	A	900	2	33,33,33	2.25	10 (30%)	45,52,52	1.54	10 (22%)
3	ANP	B	910	2	33,33,33	2.30	11 (33%)	45,52,52	2.06	11 (24%)
3	ANP	D	930	2	33,33,33	2.42	12 (36%)	45,52,52	1.82	11 (24%)

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	C	920	2	-	4/18/38/38	0/3/3/3
3	ANP	F	950	2	-	4/18/38/38	0/3/3/3
3	ANP	E	940	2	-	3/18/38/38	0/3/3/3
3	ANP	A	900	2	-	4/18/38/38	0/3/3/3
3	ANP	B	910	2	-	4/18/38/38	0/3/3/3
3	ANP	D	930	2	-	4/18/38/38	0/3/3/3

The worst 5 of 71 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	920	ANP	PB-O3A	-7.83	1.49	1.59
3	F	950	ANP	PA-O3A	-7.75	1.51	1.59
3	D	930	ANP	PA-O3A	-7.06	1.51	1.59
3	C	920	ANP	PA-O3A	-6.70	1.52	1.59
3	A	900	ANP	PA-O3A	-6.67	1.52	1.59

The worst 5 of 64 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	910	ANP	C5-C4-N3	-6.16	118.23	126.72
3	F	950	ANP	C5-C4-N3	-6.02	118.43	126.72
3	C	920	ANP	C5-C4-N3	-5.71	118.86	126.72
3	D	930	ANP	C5-C4-N3	-5.38	119.31	126.72
3	E	940	ANP	C5-C4-N3	-5.02	119.81	126.72

There are no chirality outliers.

5 of 23 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	900	ANP	PB-N3B-PG-O1G
3	A	900	ANP	PG-N3B-PB-O1B
3	A	900	ANP	PA-O3A-PB-O1B
3	B	910	ANP	PB-N3B-PG-O1G
3	B	910	ANP	PG-N3B-PB-O1B

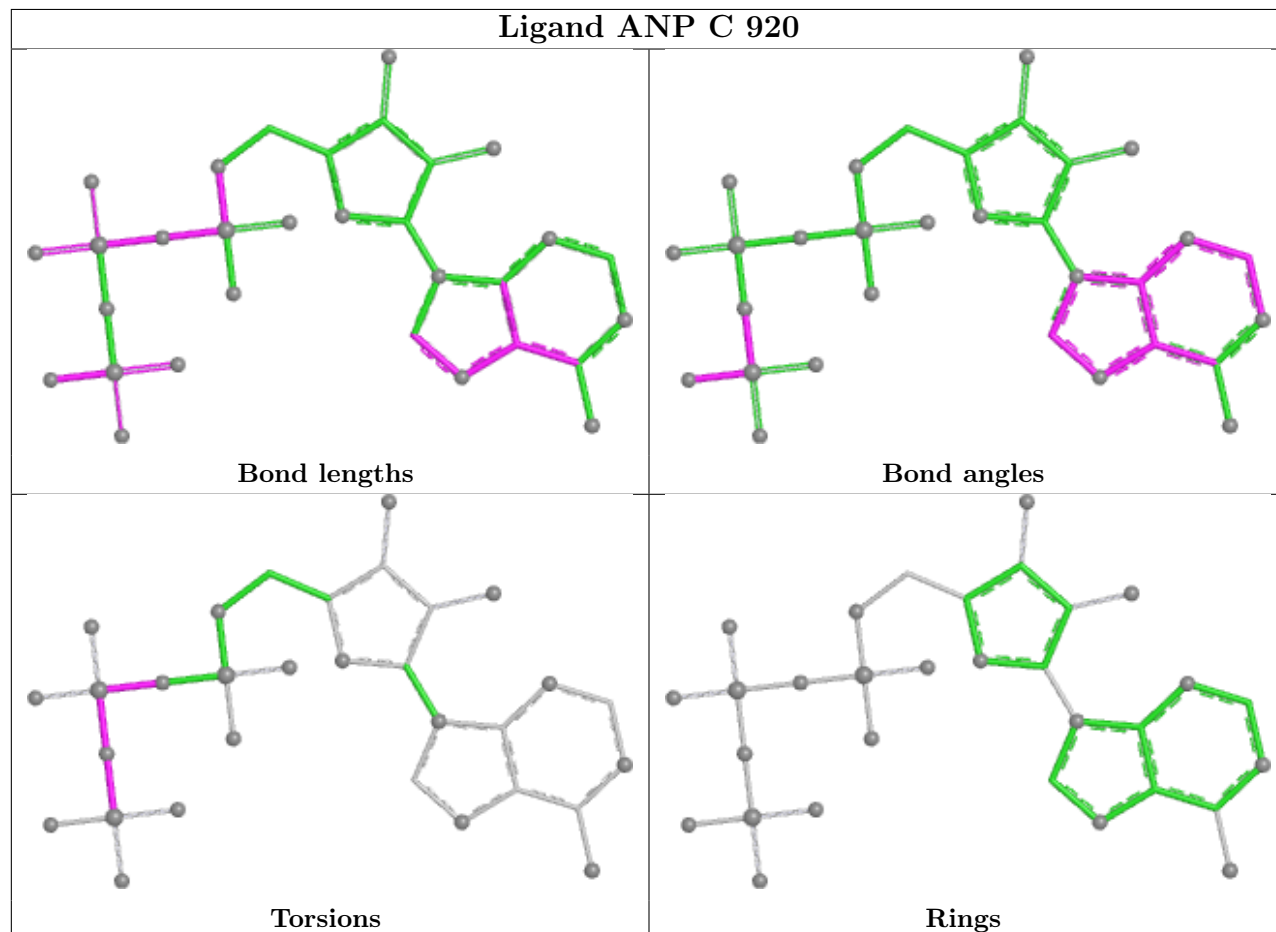
There are no ring outliers.

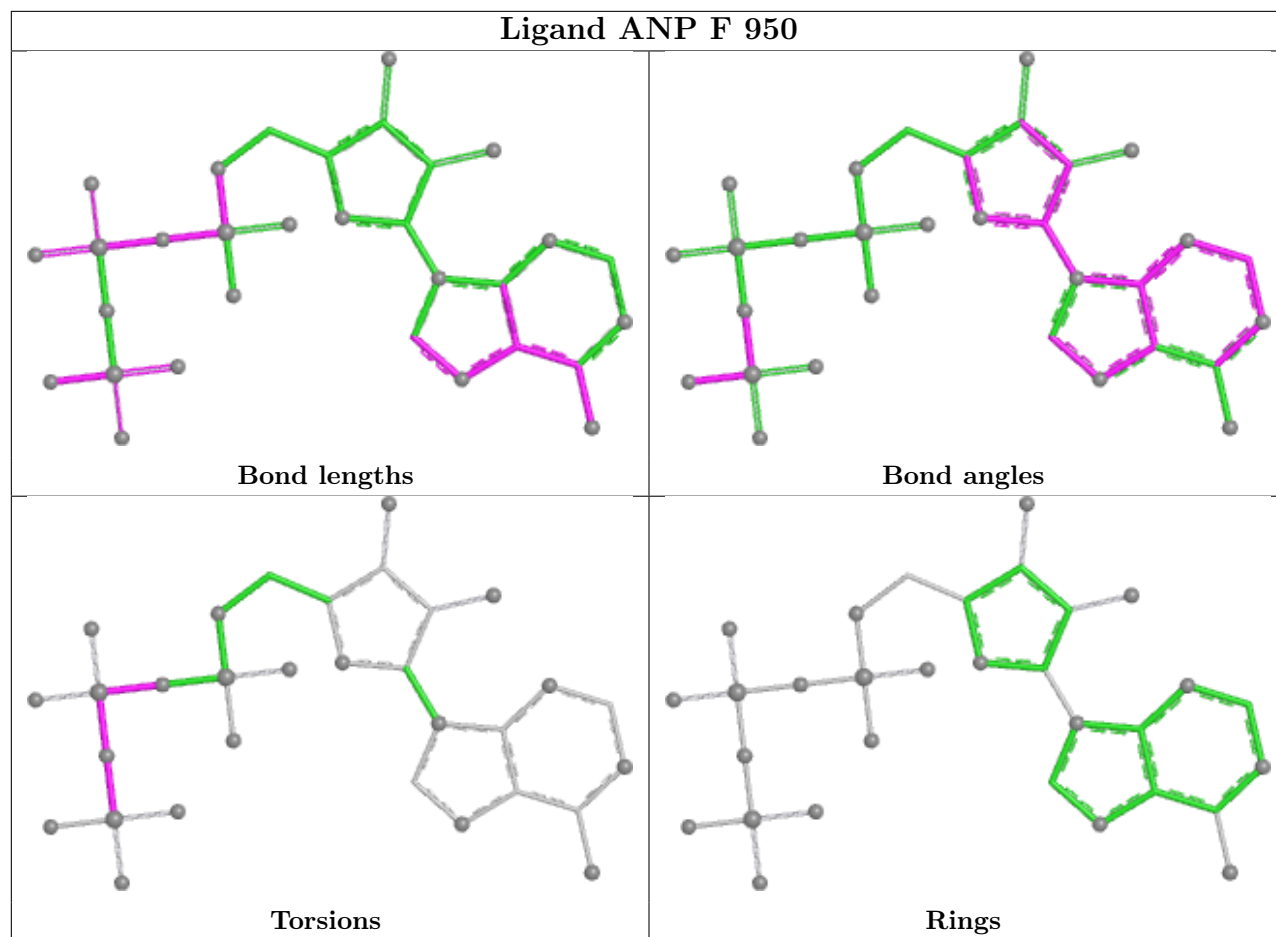
4 monomers are involved in 8 short contacts:

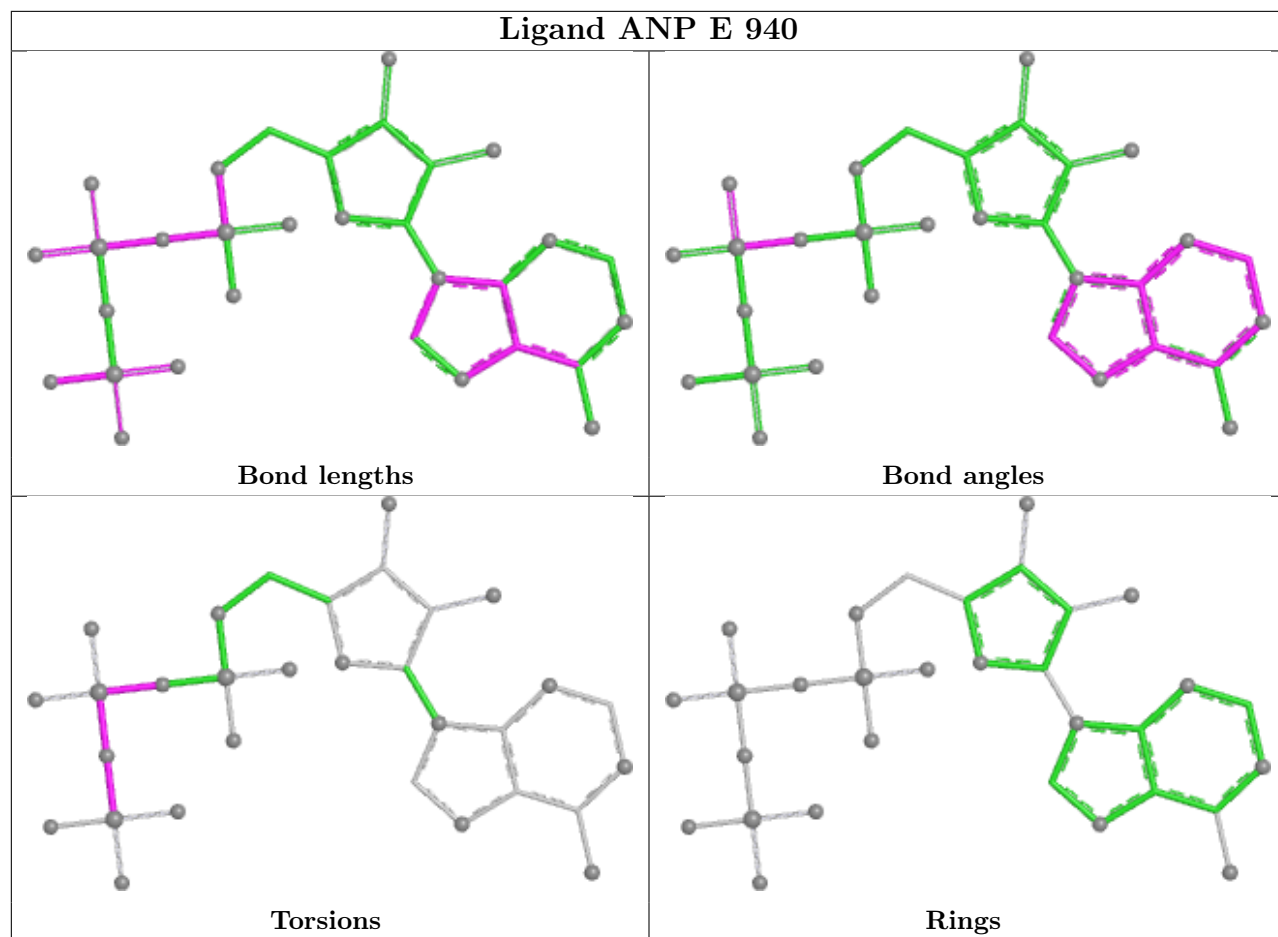
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	950	ANP	4	0
3	E	940	ANP	1	0
3	B	910	ANP	2	0
3	D	930	ANP	1	0

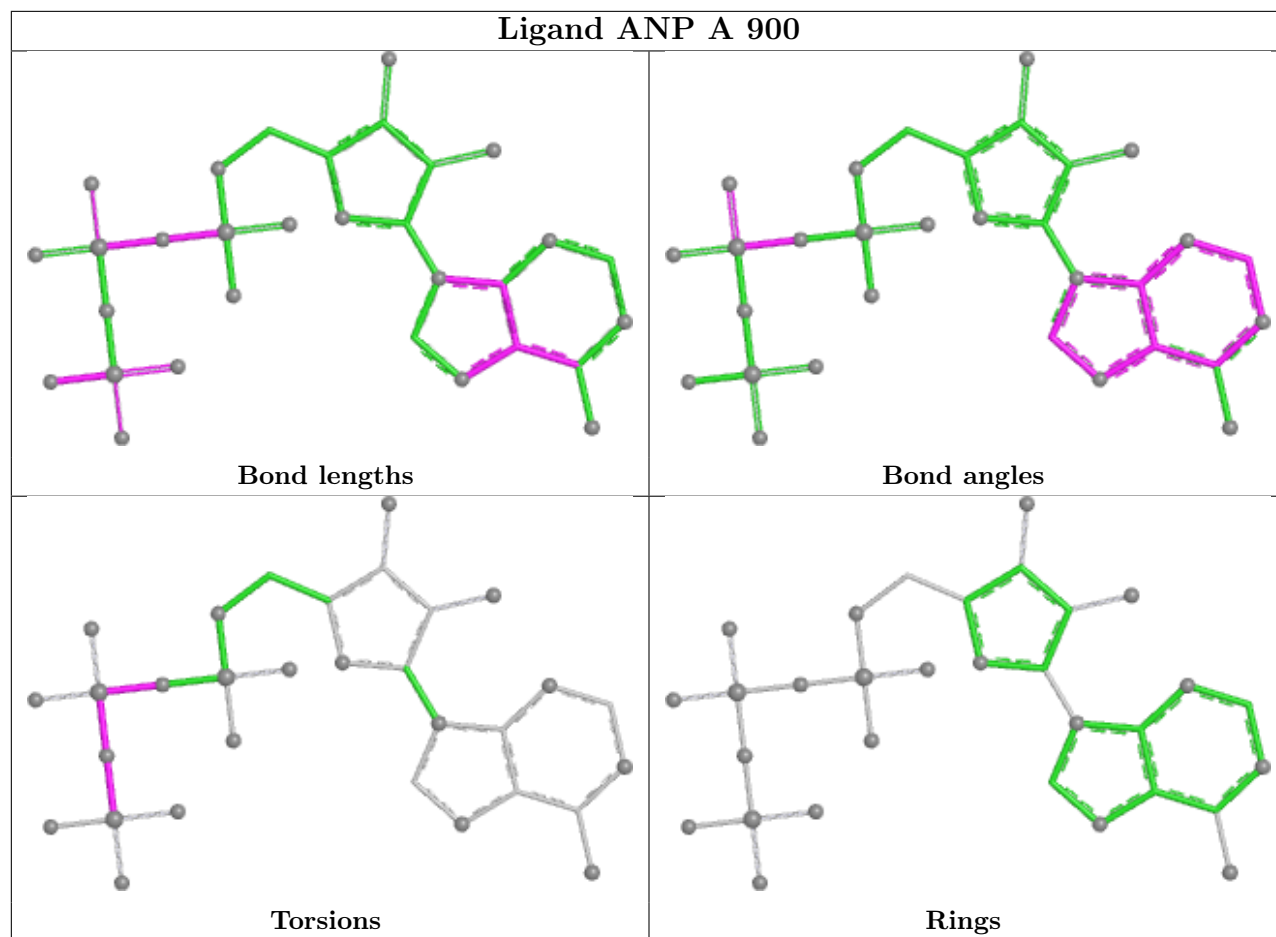
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

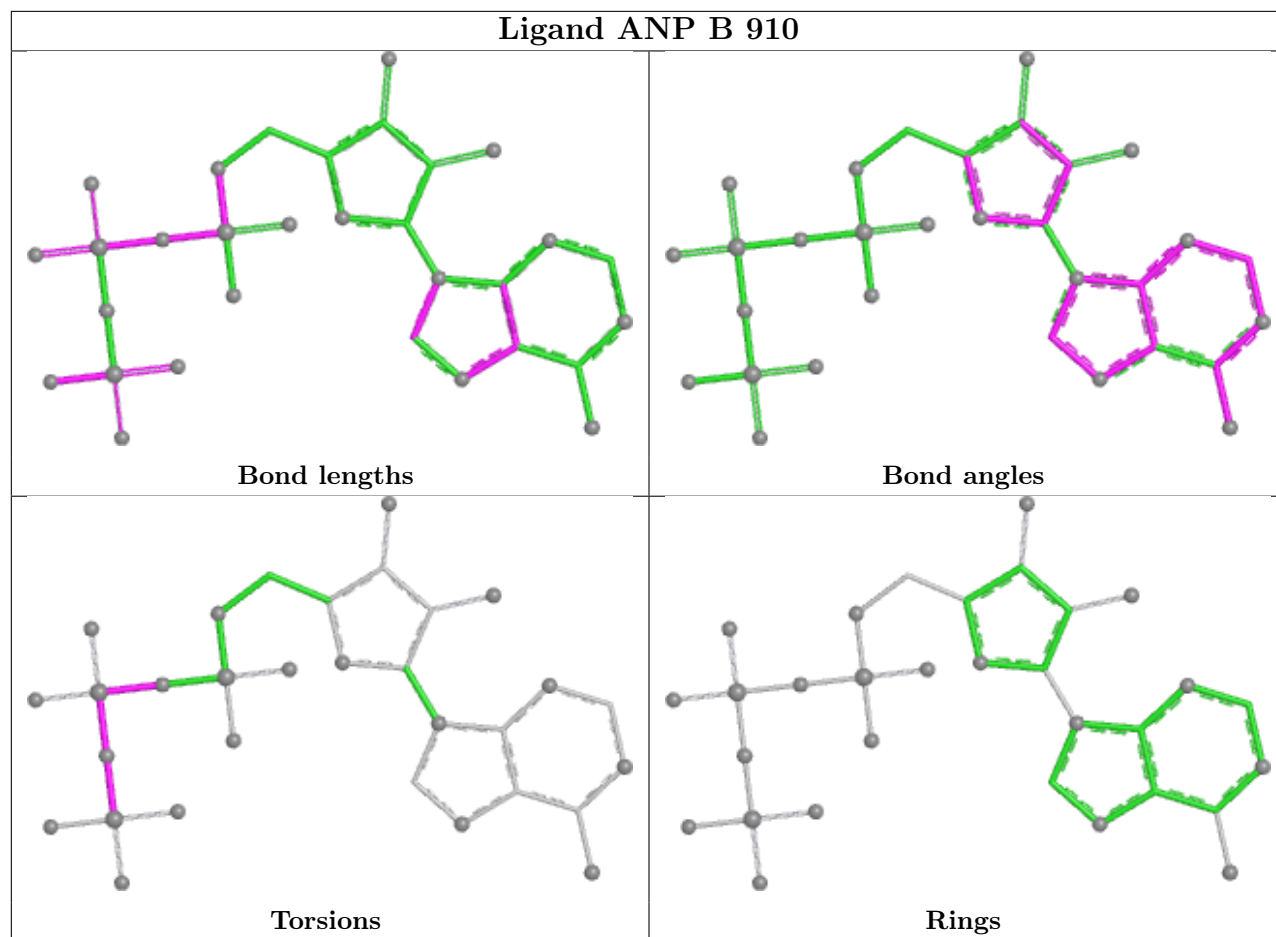
in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

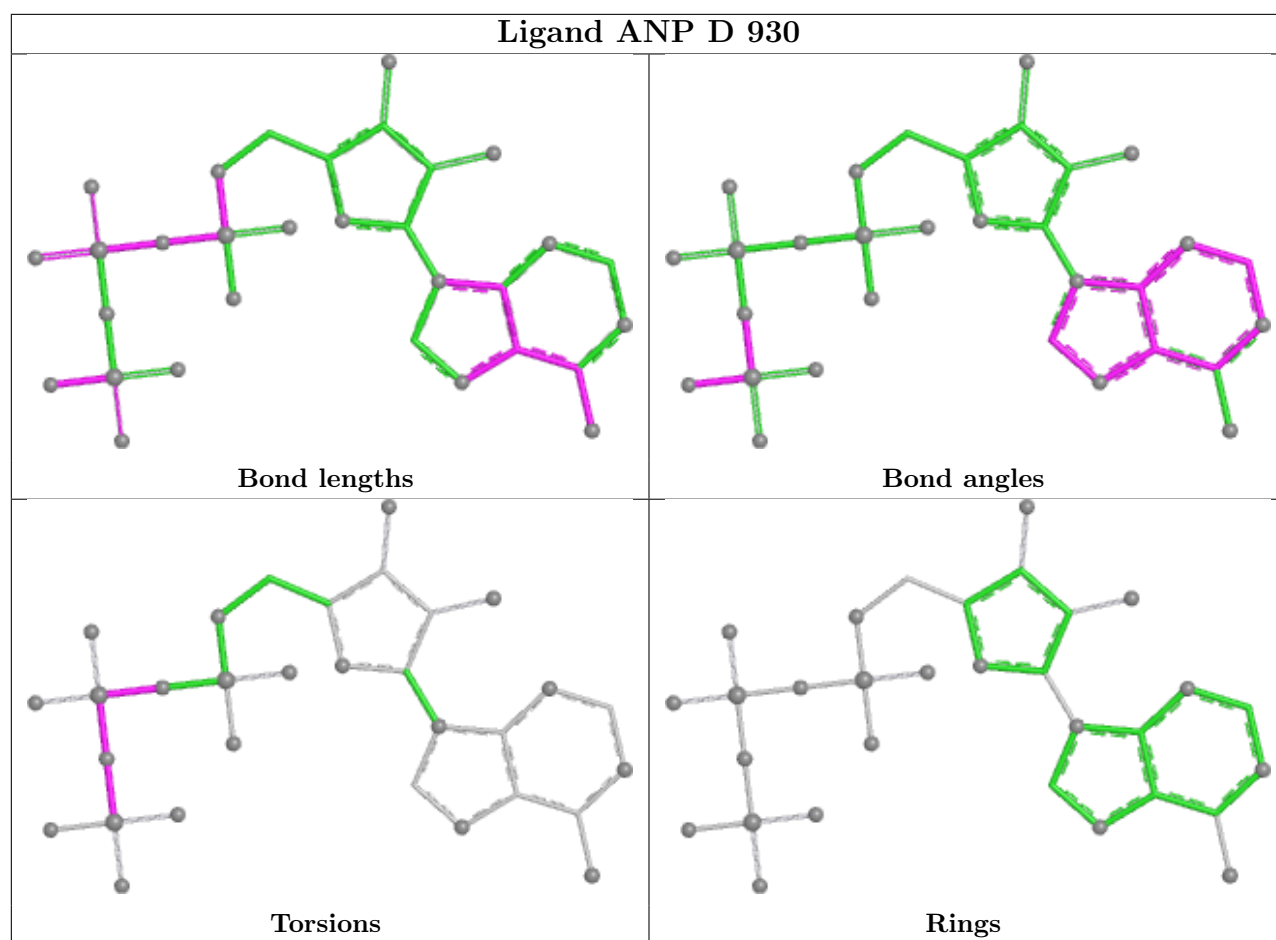












## 5.7 Other polymers [i](#)

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, 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	456/472 (96%)	-0.00	25 (5%) 30 32	4, 16, 48, 94	0
1	B	450/472 (95%)	-0.15	11 (2%) 59 62	4, 16, 40, 56	0
1	C	461/472 (97%)	-0.10	11 (2%) 59 62	7, 18, 39, 52	0
1	D	456/472 (96%)	-0.26	9 (1%) 65 66	4, 15, 37, 71	0
1	E	451/472 (95%)	0.08	16 (3%) 47 49	4, 18, 50, 72	0
1	F	449/472 (95%)	2.36	248 (55%) 0 0	18, 46, 72, 98	0
All	All	2723/2832 (96%)	0.31	320 (11%) 9 10	4, 19, 59, 98	0

The worst 5 of 320 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	212	VAL	7.9
1	F	211	ASN	7.2
1	F	217	ARG	7.1
1	F	162	GLU	6.9
1	F	159	GLY	6.7

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

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

### 6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 6.4 Ligands

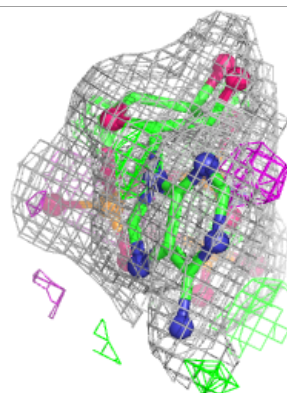
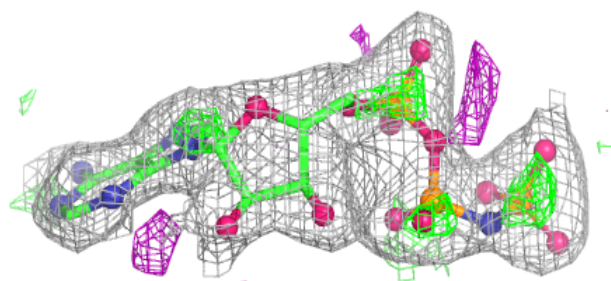
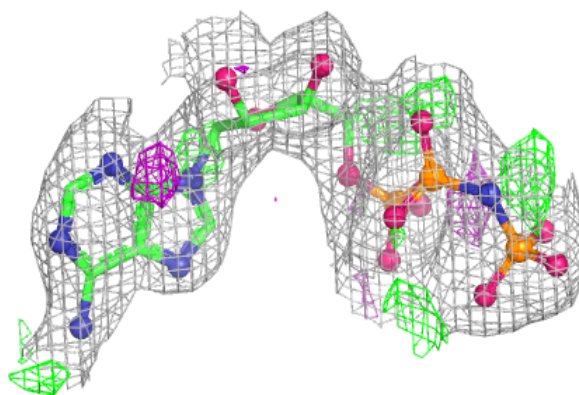
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	MG	F	501	1/1	0.73	0.16	80,80,80,80	0
4	NA	D	901	1/1	0.82	0.21	78,78,78,78	0
3	ANP	F	950	31/31	0.89	0.17	59,76,78,78	0
2	MG	E	501	1/1	0.91	0.10	51,51,51,51	0
3	ANP	B	910	31/31	0.97	0.09	38,43,46,47	0
3	ANP	C	920	31/31	0.97	0.09	42,48,52,52	0
3	ANP	E	940	31/31	0.97	0.10	43,47,49,49	0
2	MG	B	501	1/1	0.97	0.11	42,42,42,42	0
2	MG	C	501	1/1	0.97	0.05	49,49,49,49	0
2	MG	D	501	1/1	0.98	0.06	45,45,45,45	0
3	ANP	A	900	31/31	0.98	0.09	40,45,47,47	0
3	ANP	D	930	31/31	0.98	0.09	41,45,48,48	0
2	MG	A	501	1/1	0.99	0.08	43,43,43,43	0

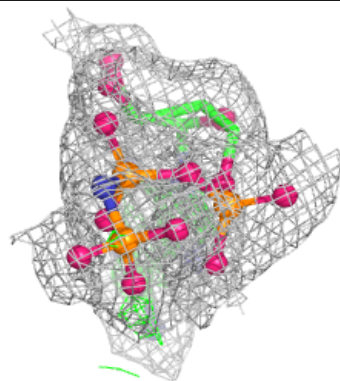
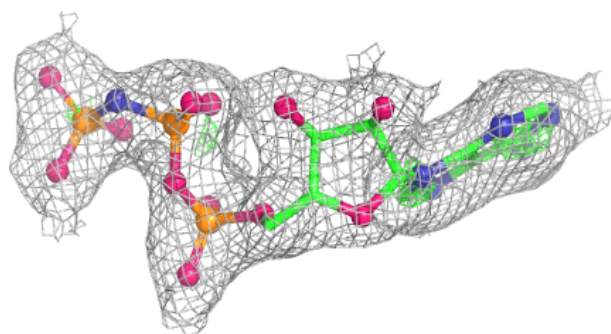
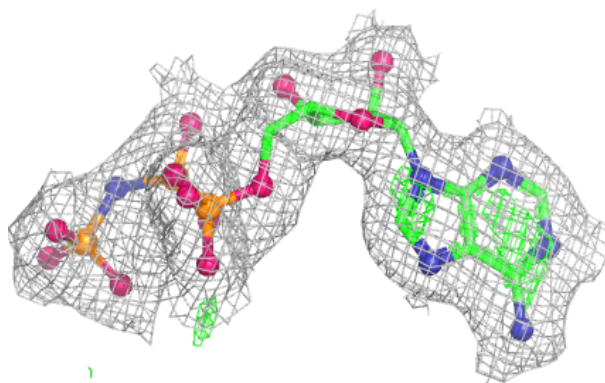
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around ANP F 950:**

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

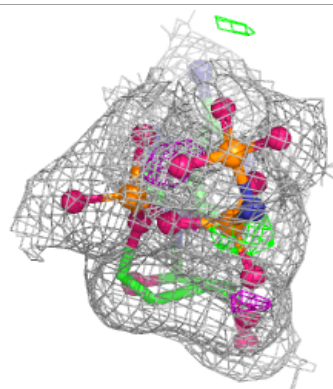
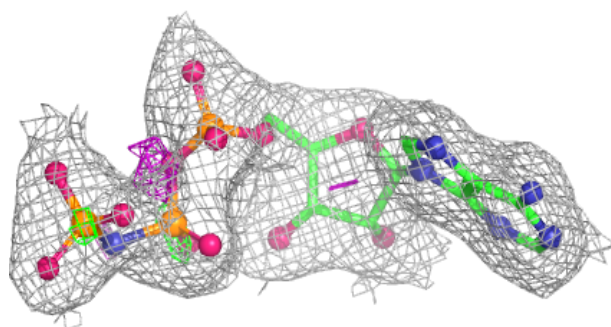
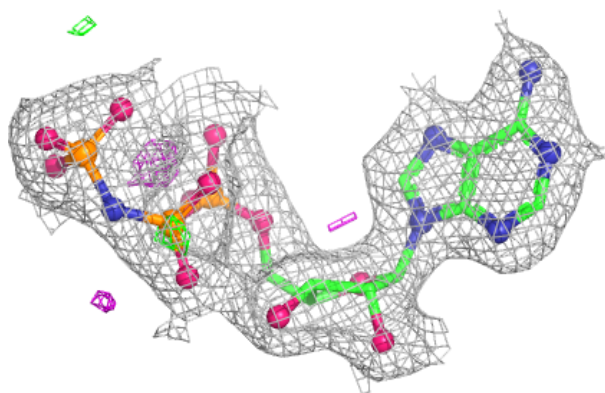
**Electron density around ANP B 910:**

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

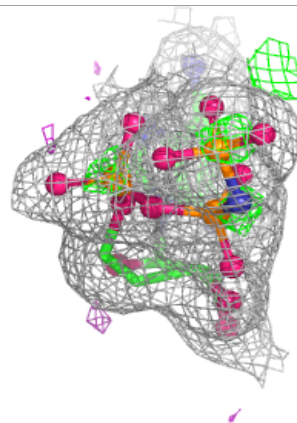
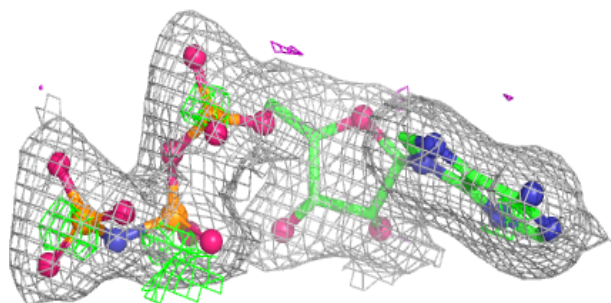
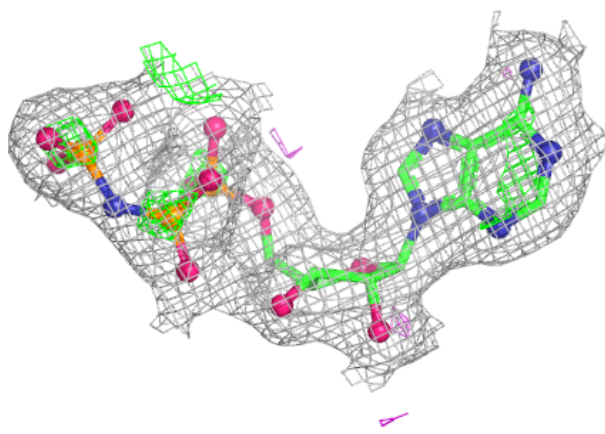


**Electron density around ANP C 920:**

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

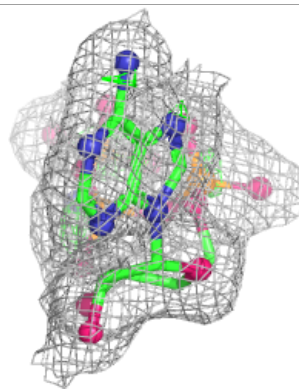
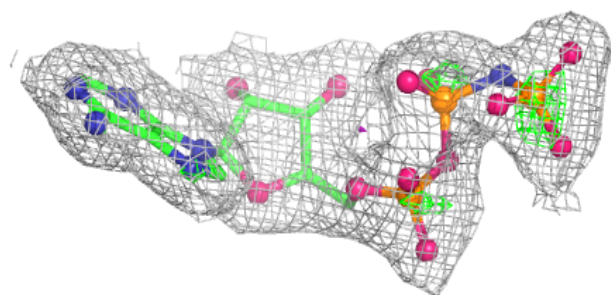
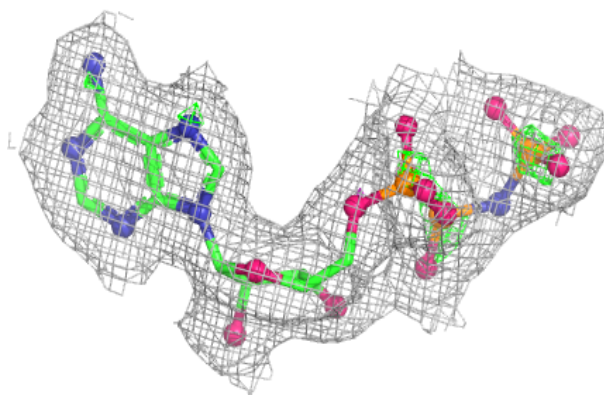
**Electron density around ANP E 940:**

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

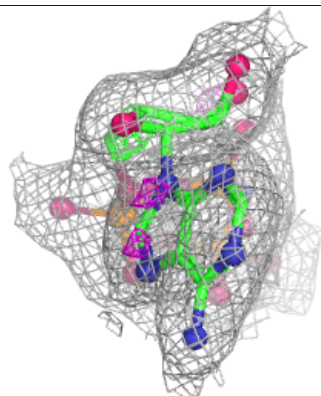
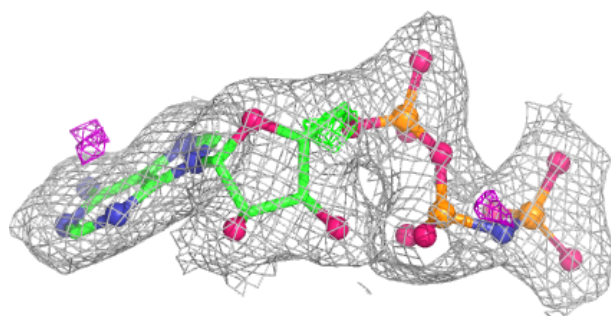
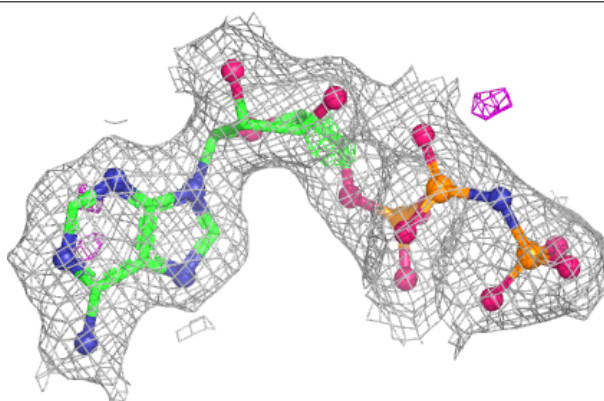


**Electron density around ANP A 900:**

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

**Electron density around ANP D 930:**

$2mF_o-DF_c$  (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.