



wwPDB X-ray Structure Validation Summary Report ⓘ

May 15, 2020 – 02:38 pm BST

PDB ID : 6DUQ
Title : Structure of a Rho-NusG KOW domain complex
Authors : Berger, J.M.; Lawson, M.R.
Deposited on : 2018-06-21
Resolution : 3.70 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 ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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

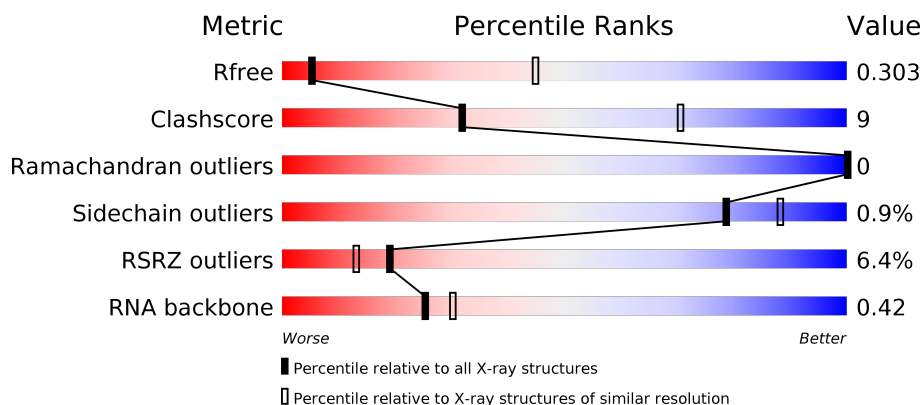
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.70 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	1049 (3.88-3.52)
Clashscore	141614	1027 (3.86-3.54)
Ramachandran outliers	138981	1069 (3.88-3.52)
Sidechain outliers	138945	1065 (3.88-3.52)
RSRZ outliers	127900	1578 (3.90-3.50)
RNA backbone	3102	1027 (4.40-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	422	<div> <div>5%</div> <div> <div></div> <div>77%</div> <div>20%</div> <div>..</div> </div> </div>
1	B	422	<div> <div>3%</div> <div> <div></div> <div>77%</div> <div>20%</div> <div>.</div> </div> </div>
1	C	422	<div> <div>5%</div> <div> <div></div> <div>77%</div> <div>21%</div> <div>..</div> </div> </div>
1	D	422	<div> <div>3%</div> <div> <div></div> <div>77%</div> <div>19%</div> <div>.</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	E	422	
1	F	422	
1	G	422	
1	H	422	
1	I	422	
1	J	422	
1	K	422	
1	L	422	
2	M	60	
2	N	60	
2	O	60	
2	P	60	
2	Q	60	
2	R	60	
2	S	60	
2	T	60	
2	U	60	
2	V	60	
2	W	60	
2	X	60	
3	Y	12	
3	Z	12	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 43612 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 Transcription termination factor Rho.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	409	Total	C	N	O	S	0	0	0
			3207	2019	566	606	16			
1	A	414	Total	C	N	O	S	0	0	0
			3232	2034	571	611	16			
1	G	414	Total	C	N	O	S	0	0	0
			3232	2034	571	611	16			
1	B	409	Total	C	N	O	S	0	0	0
			3207	2019	566	606	16			
1	H	409	Total	C	N	O	S	0	0	0
			3207	2019	566	606	16			
1	C	414	Total	C	N	O	S	0	0	0
			3232	2034	571	611	16			
1	I	414	Total	C	N	O	S	0	0	0
			3232	2034	571	611	16			
1	D	409	Total	C	N	O	S	0	0	0
			3207	2019	566	606	16			
1	K	414	Total	C	N	O	S	0	0	0
			3232	2034	571	611	16			
1	F	409	Total	C	N	O	S	0	0	0
			3207	2019	566	606	16			
1	J	409	Total	C	N	O	S	0	0	0
			3207	2019	566	606	16			
1	E	414	Total	C	N	O	S	0	0	0
			3232	2034	571	611	16			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	-2	MET	-	initiating methionine	UNP F4TMM7
L	-1	GLY	-	expression tag	UNP F4TMM7
L	0	HIS	-	expression tag	UNP F4TMM7
A	-2	MET	-	initiating methionine	UNP F4TMM7
A	-1	GLY	-	expression tag	UNP F4TMM7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	0	HIS	-	expression tag	UNP F4TMM7
G	-2	MET	-	initiating methionine	UNP F4TMM7
G	-1	GLY	-	expression tag	UNP F4TMM7
G	0	HIS	-	expression tag	UNP F4TMM7
B	-2	MET	-	initiating methionine	UNP F4TMM7
B	-1	GLY	-	expression tag	UNP F4TMM7
B	0	HIS	-	expression tag	UNP F4TMM7
H	-2	MET	-	initiating methionine	UNP F4TMM7
H	-1	GLY	-	expression tag	UNP F4TMM7
H	0	HIS	-	expression tag	UNP F4TMM7
C	-2	MET	-	initiating methionine	UNP F4TMM7
C	-1	GLY	-	expression tag	UNP F4TMM7
C	0	HIS	-	expression tag	UNP F4TMM7
I	-2	MET	-	initiating methionine	UNP F4TMM7
I	-1	GLY	-	expression tag	UNP F4TMM7
I	0	HIS	-	expression tag	UNP F4TMM7
D	-2	MET	-	initiating methionine	UNP F4TMM7
D	-1	GLY	-	expression tag	UNP F4TMM7
D	0	HIS	-	expression tag	UNP F4TMM7
K	-2	MET	-	initiating methionine	UNP F4TMM7
K	-1	GLY	-	expression tag	UNP F4TMM7
K	0	HIS	-	expression tag	UNP F4TMM7
F	-2	MET	-	initiating methionine	UNP F4TMM7
F	-1	GLY	-	expression tag	UNP F4TMM7
F	0	HIS	-	expression tag	UNP F4TMM7
J	-2	MET	-	initiating methionine	UNP F4TMM7
J	-1	GLY	-	expression tag	UNP F4TMM7
J	0	HIS	-	expression tag	UNP F4TMM7
E	-2	MET	-	initiating methionine	UNP F4TMM7
E	-1	GLY	-	expression tag	UNP F4TMM7
E	0	HIS	-	expression tag	UNP F4TMM7

- Molecule 2 is a protein called Transcription termination/antitermination protein NusG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	M	43	Total	C	N	O	S	0	0	0
			323	206	51	65	1			
2	N	43	Total	C	N	O	S	0	0	0
			323	206	51	65	1			
2	O	43	Total	C	N	O	S	0	0	0
			323	206	51	65	1			
2	P	43	Total	C	N	O	S	0	0	0
			323	206	51	65	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	Q	43	Total	C	N	O	S	0	0	0
			323	206	51	65	1			
2	R	43	Total	C	N	O	S	0	0	0
			323	206	51	65	1			
2	S	43	Total	C	N	O	S	0	0	0
			323	206	51	65	1			
2	T	43	Total	C	N	O	S	0	0	0
			323	206	51	65	1			
2	U	43	Total	C	N	O	S	0	0	0
			323	206	51	65	1			
2	V	43	Total	C	N	O	S	0	0	0
			323	206	51	65	1			
2	W	43	Total	C	N	O	S	0	0	0
			319	204	51	63	1			
2	X	43	Total	C	N	O	S	0	0	0
			319	204	51	63	1			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
M	122	SER	-	expression tag	UNP F4T6L5
M	123	ASN	-	expression tag	UNP F4T6L5
M	124	ALA	-	expression tag	UNP F4T6L5
N	122	SER	-	expression tag	UNP F4T6L5
N	123	ASN	-	expression tag	UNP F4T6L5
N	124	ALA	-	expression tag	UNP F4T6L5
O	122	SER	-	expression tag	UNP F4T6L5
O	123	ASN	-	expression tag	UNP F4T6L5
O	124	ALA	-	expression tag	UNP F4T6L5
P	122	SER	-	expression tag	UNP F4T6L5
P	123	ASN	-	expression tag	UNP F4T6L5
P	124	ALA	-	expression tag	UNP F4T6L5
Q	122	SER	-	expression tag	UNP F4T6L5
Q	123	ASN	-	expression tag	UNP F4T6L5
Q	124	ALA	-	expression tag	UNP F4T6L5
R	122	SER	-	expression tag	UNP F4T6L5
R	123	ASN	-	expression tag	UNP F4T6L5
R	124	ALA	-	expression tag	UNP F4T6L5
S	122	SER	-	expression tag	UNP F4T6L5
S	123	ASN	-	expression tag	UNP F4T6L5
S	124	ALA	-	expression tag	UNP F4T6L5
T	122	SER	-	expression tag	UNP F4T6L5
T	123	ASN	-	expression tag	UNP F4T6L5

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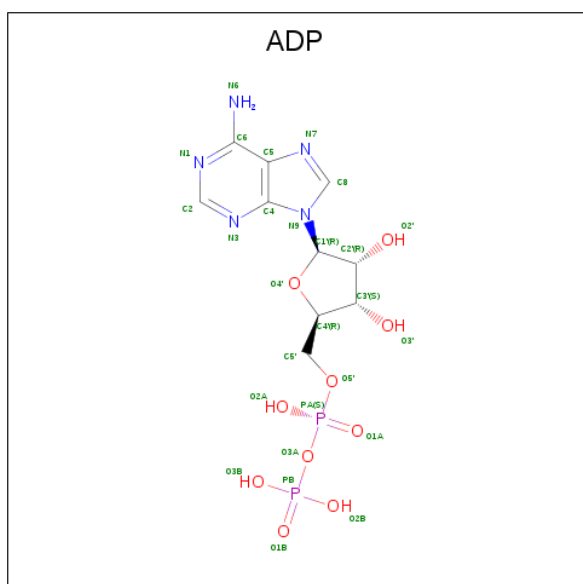
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Chain	Residue	Modelled	Actual	Comment	Reference
T	124	ALA	-	expression tag	UNP F4T6L5
U	122	SER	-	expression tag	UNP F4T6L5
U	123	ASN	-	expression tag	UNP F4T6L5
U	124	ALA	-	expression tag	UNP F4T6L5
V	122	SER	-	expression tag	UNP F4T6L5
V	123	ASN	-	expression tag	UNP F4T6L5
V	124	ALA	-	expression tag	UNP F4T6L5
W	122	SER	-	expression tag	UNP F4T6L5
W	123	ASN	-	expression tag	UNP F4T6L5
W	124	ALA	-	expression tag	UNP F4T6L5
X	122	SER	-	expression tag	UNP F4T6L5
X	123	ASN	-	expression tag	UNP F4T6L5
X	124	ALA	-	expression tag	UNP F4T6L5

- Molecule 3 is a RNA chain called rU12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	Y	6	Total	C	N	O	P	0	6	0
			363	162	36	147	18			
3	Z	6	Total	C	N	O	P	0	6	0
			363	162	36	147	18			

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	L	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	G	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	H	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	I	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	D	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	K	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		
4	J	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	E	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

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

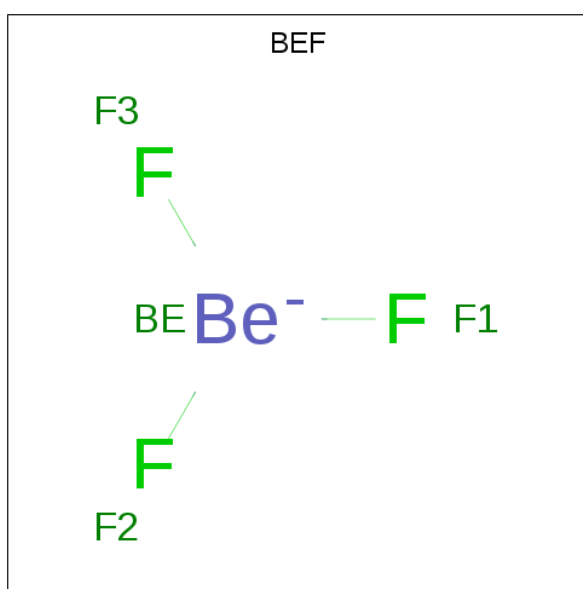
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	G	1	Total	Mg	0	0
			1	1		
5	J	1	Total	Mg	0	0
			1	1		
5	D	1	Total	Mg	0	0
			1	1		
5	K	1	Total	Mg	0	0
			1	1		
5	E	1	Total	Mg	0	0
			1	1		
5	H	1	Total	Mg	0	0
			1	1		
5	B	1	Total	Mg	0	0
			1	1		
5	I	1	Total	Mg	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	1	Total 1	Mg 1	0	0
5	A	1	Total 1	Mg 1	0	0
5	L	1	Total 1	Mg 1	0	0
5	F	1	Total 1	Mg 1	0	0

- Molecule 6 is BERYLLIUM TRIFLUORIDE ION (three-letter code: BEF) (formula: BeF₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	L	1	Total 4	Be 1	F 3	0	0
6	A	1	Total 4	Be 1	F 3	0	0
6	G	1	Total 4	Be 1	F 3	0	0
6	B	1	Total 4	Be 1	F 3	0	0
6	H	1	Total 4	Be 1	F 3	0	0
6	C	1	Total 4	Be 1	F 3	0	0
6	I	1	Total 4	Be 1	F 3	0	0

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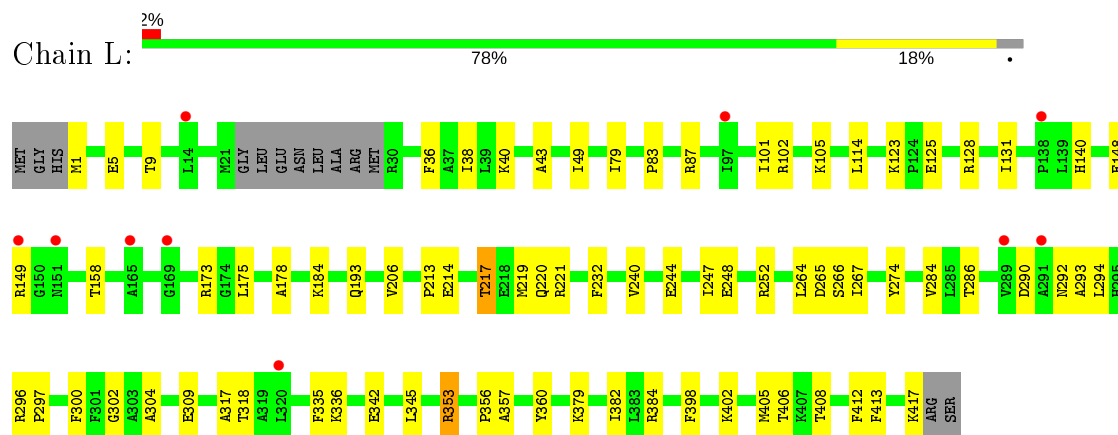
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
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6	K	1	Total 4	Be 1	F 3	0	0
6	F	1	Total 4	Be 1	F 3	0	0
6	J	1	Total 4	Be 1	F 3	0	0
6	E	1	Total 4	Be 1	F 3	0	0

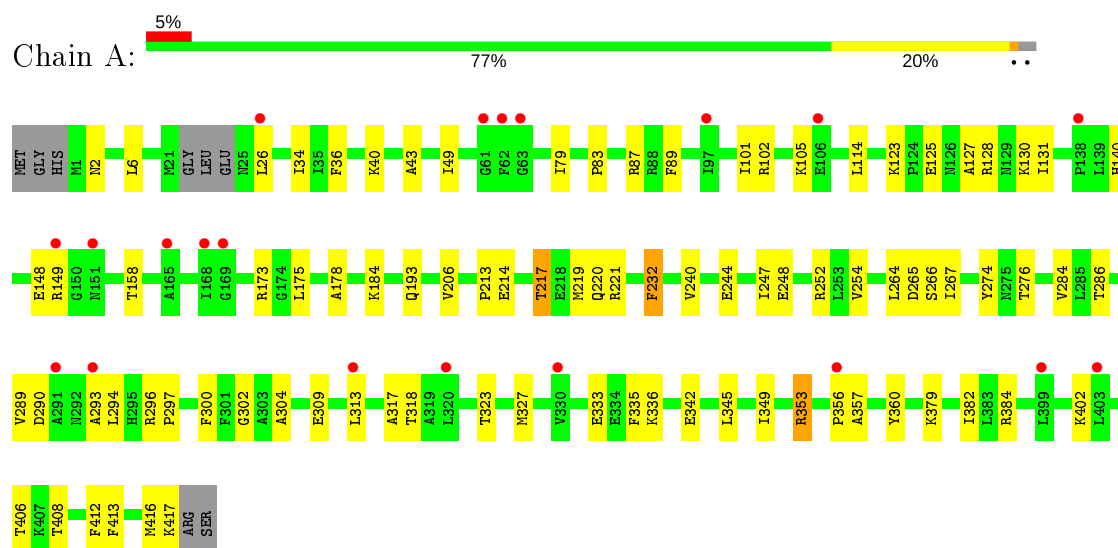
3 Residue-property plots [i](#)

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.

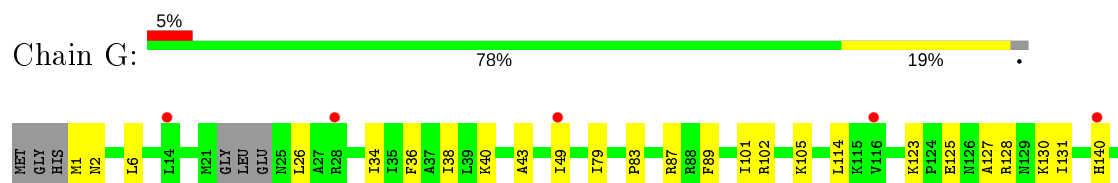
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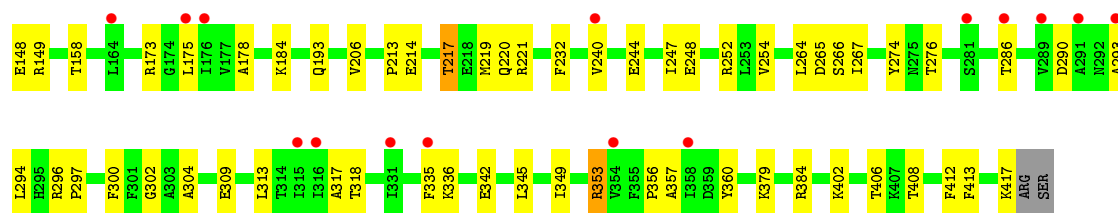


- Molecule 1: Transcription termination factor Rho

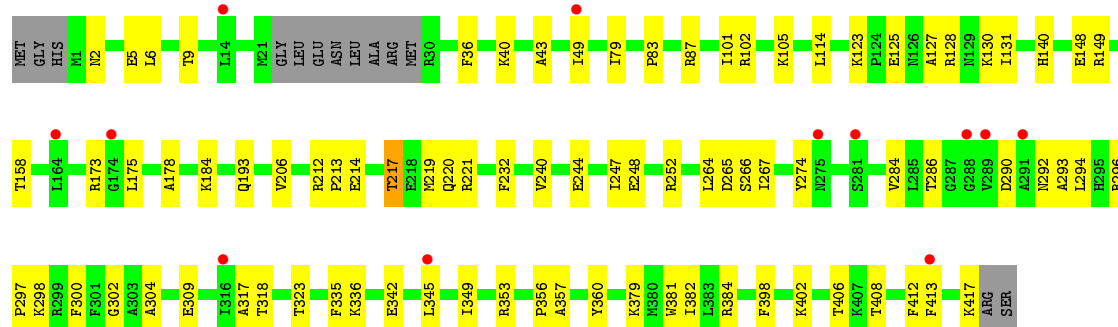
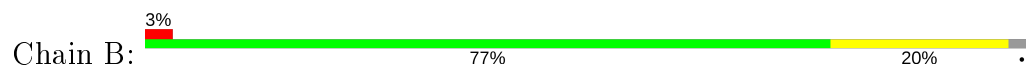


- Molecule 1: Transcription termination factor Rho

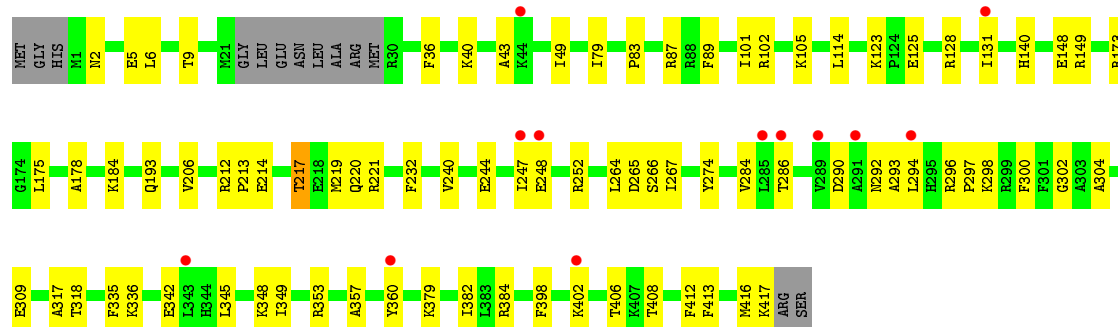
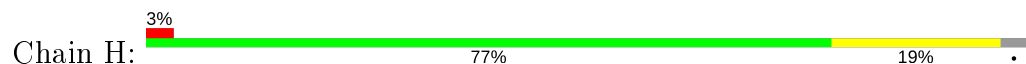




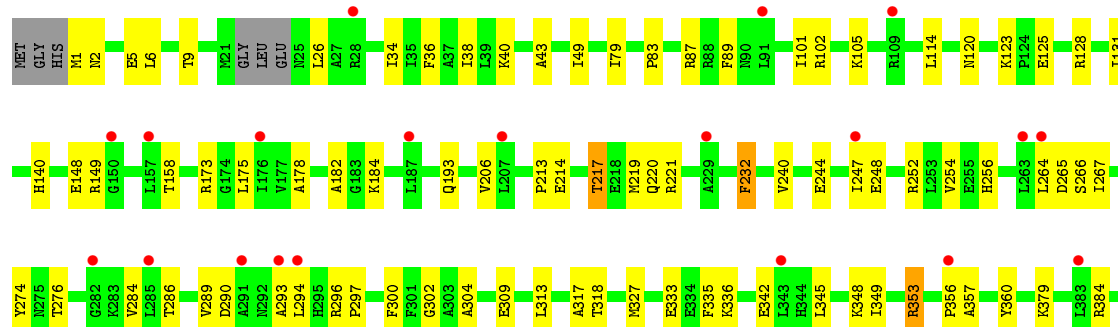
• Molecule 1: Transcription termination factor Rho



• Molecule 1: Transcription termination factor Rho

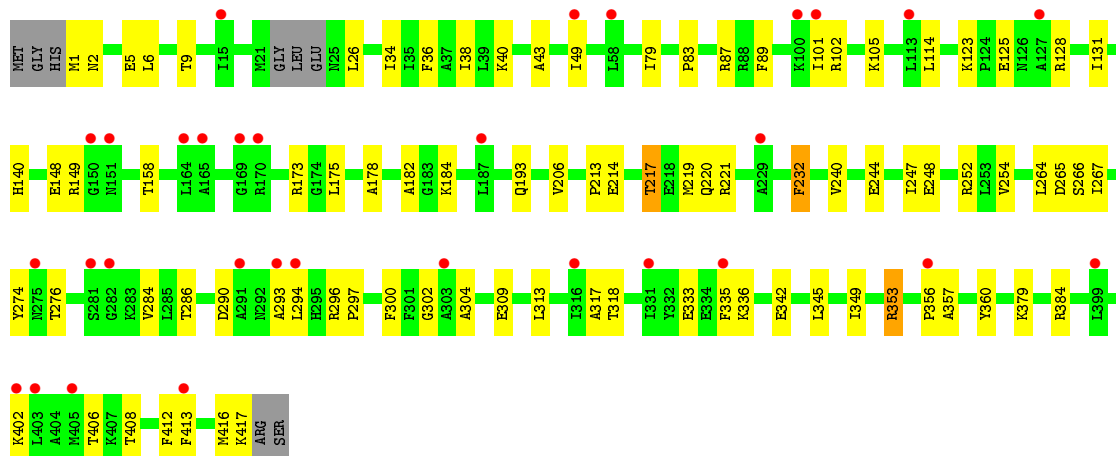
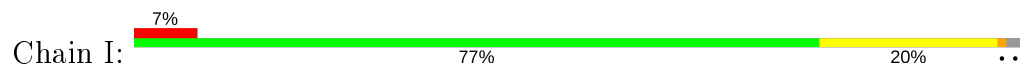


• Molecule 1: Transcription termination factor Rho

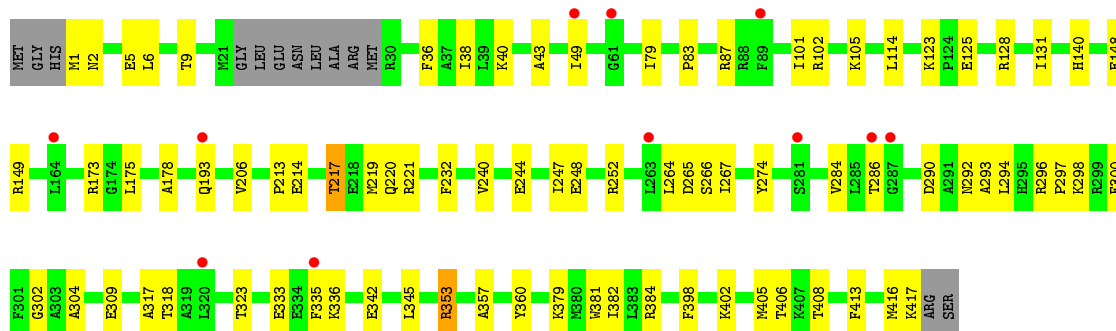
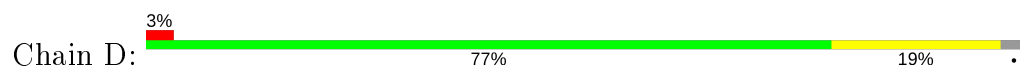




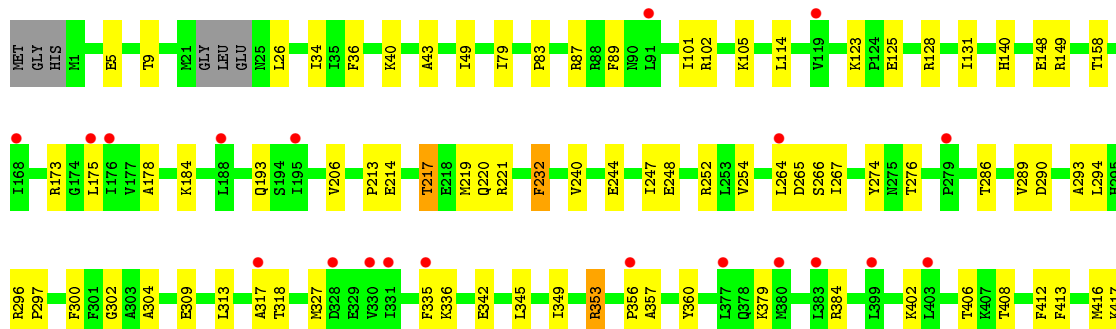
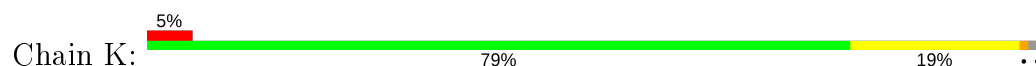
• Molecule 1: Transcription termination factor Rho



• Molecule 1: Transcription termination factor Rho




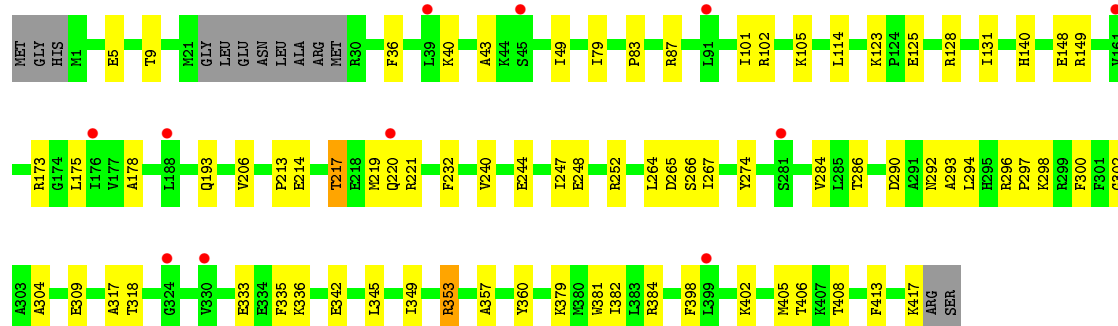
• Molecule 1: Transcription termination factor Rho




ARG
SER

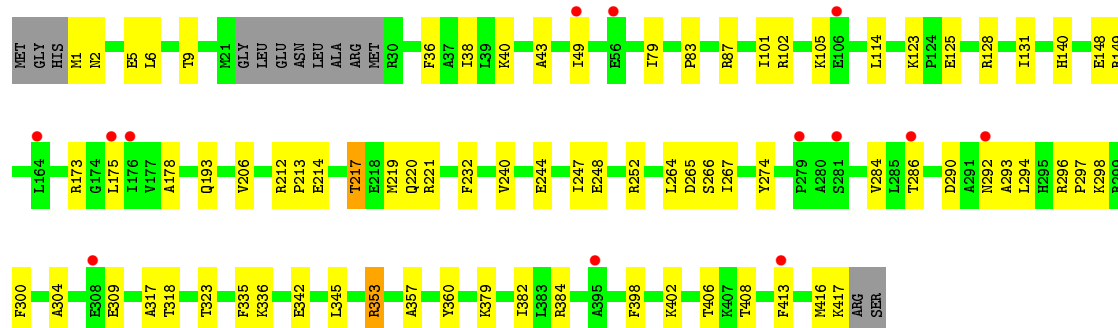
• Molecule 1: Transcription termination factor Rho

Chain F: 



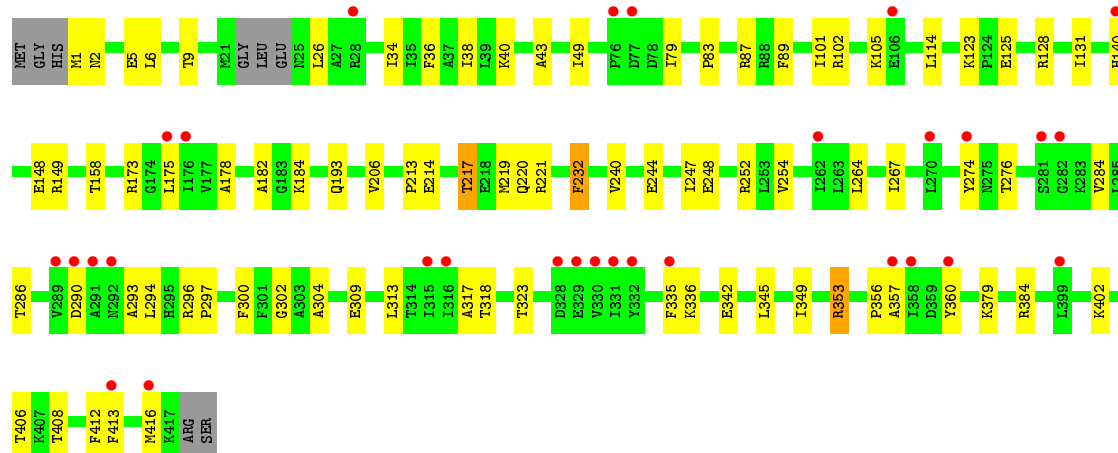
• Molecule 1: Transcription termination factor Rho

Chain J: 

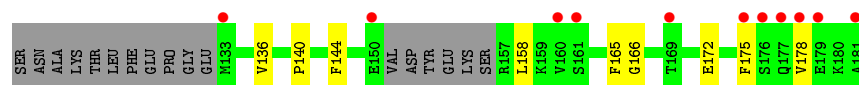


• Molecule 1: Transcription termination factor Rho

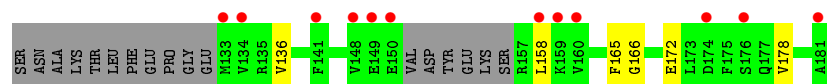
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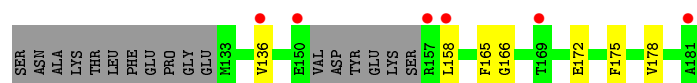
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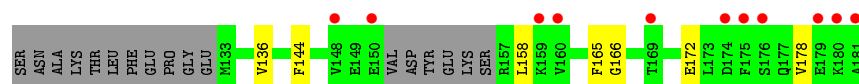
- Molecule 2: Transcription termination/antitermination protein NusG



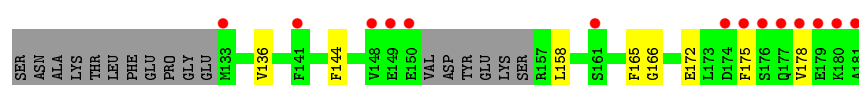
- Molecule 2: Transcription termination/antitermination protein NusG



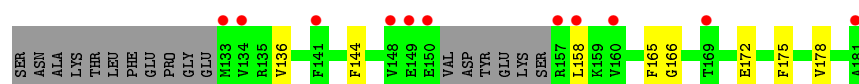
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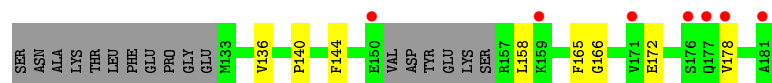
- Molecule 2: Transcription termination/antitermination protein NusG



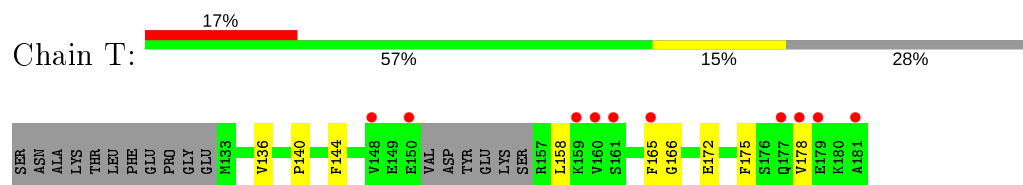
- Molecule 2: Transcription termination/antitermination protein NusG



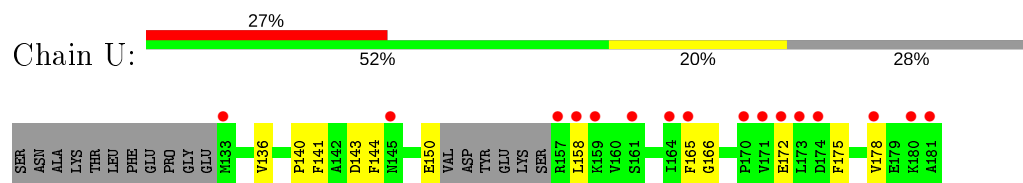
- Molecule 2: Transcription termination/antitermination protein NusG



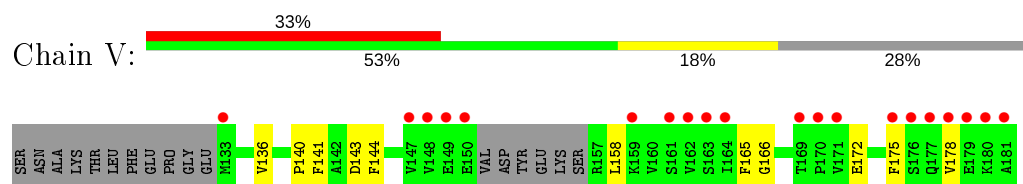
- Molecule 2: Transcription termination/antitermination protein NusG



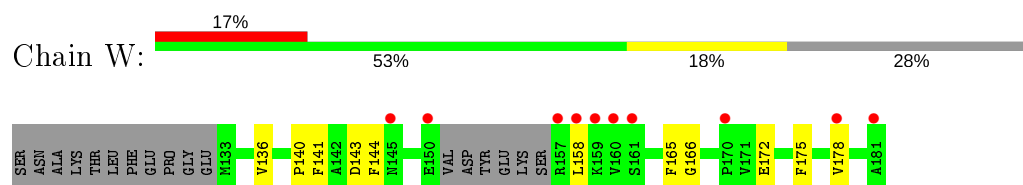
- Molecule 2: Transcription termination/antitermination protein NusG



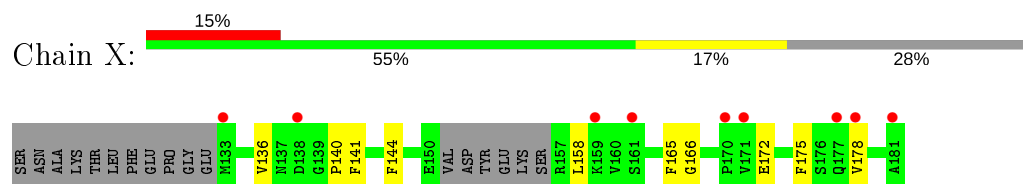
- Molecule 2: Transcription termination/antitermination protein NusG



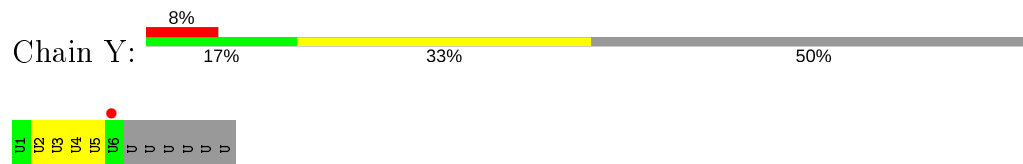
- Molecule 2: Transcription termination/antitermination protein NusG



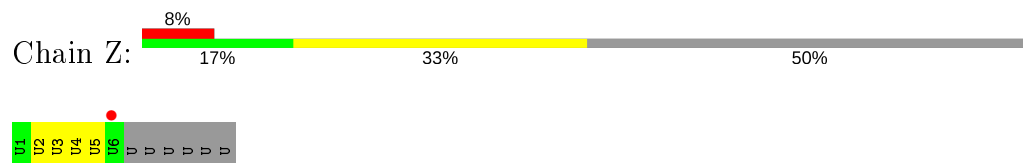
- Molecule 2: Transcription termination/antitermination protein NusG



- Molecule 3: rU12



- Molecule 3: rU12



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	126.35Å 126.32Å 166.82Å 89.94° 89.90° 60.01°	Depositor
Resolution (Å)	39.03 – 3.70 39.03 – 3.40	Depositor EDS
% Data completeness (in resolution range)	97.7 (39.03-3.70) 96.7 (39.03-3.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.31 (at 3.40Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, R_{free}	0.287 , 0.300 0.289 , 0.303	Depositor DCC
R_{free} test set	1958 reflections (1.63%)	wwPDB-VP
Wilson B-factor (Å ²)	103.0	Xtriage
Anisotropy	0.151	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 99.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	0.317 for h-k,h,l 0.317 for k,-h+k,l 0.307 for -k,h-k,l 0.307 for -h+k,-h,l 0.126 for -h+k,k,-l 0.127 for h,h-k,-l 0.316 for -h,-k,l 0.125 for k,h,-l 0.126 for -k,-h,-l 0.126 for h-k,-k,-l 0.125 for -h,-h+k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	43612	wwPDB-VP
Average B, all atoms (Å ²)	201.0	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 44.00 % 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 1.6248e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, BEF, 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	0.31	0/3280	0.48	0/4422
1	B	0.34	0/3255	0.50	0/4387
1	C	0.32	0/3280	0.48	0/4422
1	D	0.33	0/3255	0.50	0/4387
1	E	0.32	0/3280	0.48	0/4422
1	F	0.34	0/3255	0.50	0/4387
1	G	0.32	0/3280	0.48	0/4422
1	H	0.33	0/3255	0.49	0/4387
1	I	0.32	0/3280	0.48	0/4422
1	J	0.33	0/3255	0.50	0/4387
1	K	0.31	0/3280	0.48	0/4422
1	L	0.34	0/3255	0.50	0/4387
2	M	0.41	0/327	0.60	0/440
2	N	0.40	0/327	0.60	0/440
2	O	0.40	0/327	0.60	0/440
2	P	0.40	0/327	0.60	0/440
2	Q	0.39	0/327	0.60	0/440
2	R	0.40	0/327	0.60	0/440
2	S	0.55	0/327	0.69	0/440
2	T	0.55	0/327	0.69	0/440
2	U	0.58	0/327	0.71	0/440
2	V	0.47	0/327	0.64	0/440
2	W	0.44	0/323	0.61	0/435
2	X	0.54	0/323	0.69	0/435
3	Y	0.48	0/396	1.02	0/600
3	Z	0.48	0/396	1.05	0/600
All	All	0.34	0/43918	0.52	0/59324

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3232	0	3273	62	0
1	B	3207	0	3260	69	0
1	C	3232	0	3273	62	0
1	D	3207	0	3260	68	0
1	E	3232	0	3273	61	0
1	F	3207	0	3260	67	0
1	G	3232	0	3273	54	0
1	H	3207	0	3260	67	0
1	I	3232	0	3273	60	0
1	J	3207	0	3260	69	0
1	K	3232	0	3273	58	0
1	L	3207	0	3260	63	0
2	M	323	0	307	6	0
2	N	323	0	307	3	0
2	O	323	0	307	4	0
2	P	323	0	307	4	0
2	Q	323	0	307	5	0
2	R	323	0	307	5	0
2	S	323	0	307	15	0
2	T	323	0	307	16	0
2	U	323	0	307	19	0
2	V	323	0	307	14	0
2	W	319	0	303	20	0
2	X	319	0	303	15	0
3	Y	363	0	183	22	0
3	Z	363	0	183	16	0
4	A	27	0	12	0	0
4	B	27	0	12	0	0
4	C	27	0	12	0	0
4	D	27	0	12	0	0
4	E	27	0	12	0	0
4	F	27	0	12	0	0
4	G	27	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	27	0	12	0	0
4	I	27	0	12	0	0
4	J	27	0	12	0	0
4	K	27	0	12	0	0
4	L	27	0	12	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
5	E	1	0	0	0	0
5	F	1	0	0	0	0
5	G	1	0	0	0	0
5	H	1	0	0	0	0
5	I	1	0	0	0	0
5	J	1	0	0	0	0
5	K	1	0	0	0	0
5	L	1	0	0	0	0
6	A	4	0	0	0	0
6	B	4	0	0	0	0
6	C	4	0	0	0	0
6	D	4	0	0	0	0
6	E	4	0	0	0	0
6	F	4	0	0	0	0
6	G	4	0	0	0	0
6	H	4	0	0	0	0
6	I	4	0	0	0	0
6	J	4	0	0	0	0
6	K	4	0	0	0	0
6	L	4	0	0	0	0
All	All	43612	0	43384	776	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 9.

The worst 5 of 776 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:F:402:LYS:HG3	2:U:165:PHE:HE1	1.17	1.08
1:K:286:THR:HA	3:Z:5[A]:U:H4'	1.48	0.95
1:F:402:LYS:HG3	2:U:165:PHE:CE1	2.05	0.92
1:J:402:LYS:HG3	2:W:165:PHE:HE1	1.38	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:286:THR:HA	3:Z:5[B]:U:H4'	1.56	0.88

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	410/422 (97%)	403 (98%)	7 (2%)	0	100	100
1	B	405/422 (96%)	398 (98%)	7 (2%)	0	100	100
1	C	410/422 (97%)	403 (98%)	7 (2%)	0	100	100
1	D	405/422 (96%)	398 (98%)	7 (2%)	0	100	100
1	E	410/422 (97%)	403 (98%)	7 (2%)	0	100	100
1	F	405/422 (96%)	398 (98%)	7 (2%)	0	100	100
1	G	410/422 (97%)	403 (98%)	7 (2%)	0	100	100
1	H	405/422 (96%)	398 (98%)	7 (2%)	0	100	100
1	I	410/422 (97%)	403 (98%)	7 (2%)	0	100	100
1	J	405/422 (96%)	398 (98%)	7 (2%)	0	100	100
1	K	410/422 (97%)	403 (98%)	7 (2%)	0	100	100
1	L	405/422 (96%)	398 (98%)	7 (2%)	0	100	100
2	M	39/60 (65%)	38 (97%)	1 (3%)	0	100	100
2	N	39/60 (65%)	38 (97%)	1 (3%)	0	100	100
2	O	39/60 (65%)	38 (97%)	1 (3%)	0	100	100
2	P	39/60 (65%)	38 (97%)	1 (3%)	0	100	100
2	Q	39/60 (65%)	38 (97%)	1 (3%)	0	100	100
2	R	39/60 (65%)	38 (97%)	1 (3%)	0	100	100
2	S	39/60 (65%)	38 (97%)	1 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	T	39/60 (65%)	37 (95%)	2 (5%)	0	100	100
2	U	39/60 (65%)	37 (95%)	2 (5%)	0	100	100
2	V	39/60 (65%)	38 (97%)	1 (3%)	0	100	100
2	W	39/60 (65%)	37 (95%)	2 (5%)	0	100	100
2	X	39/60 (65%)	38 (97%)	1 (3%)	0	100	100
All	All	5358/5784 (93%)	5259 (98%)	99 (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	348/361 (96%)	344 (99%)	4 (1%)	73	85
1	B	348/361 (96%)	345 (99%)	3 (1%)	78	88
1	C	348/361 (96%)	344 (99%)	4 (1%)	73	85
1	D	348/361 (96%)	345 (99%)	3 (1%)	78	88
1	E	348/361 (96%)	344 (99%)	4 (1%)	73	85
1	F	348/361 (96%)	345 (99%)	3 (1%)	78	88
1	G	348/361 (96%)	345 (99%)	3 (1%)	78	88
1	H	348/361 (96%)	345 (99%)	3 (1%)	78	88
1	I	348/361 (96%)	344 (99%)	4 (1%)	73	85
1	J	348/361 (96%)	345 (99%)	3 (1%)	78	88
1	K	348/361 (96%)	344 (99%)	4 (1%)	73	85
1	L	348/361 (96%)	345 (99%)	3 (1%)	78	88
2	M	35/52 (67%)	35 (100%)	0	100	100
2	N	35/52 (67%)	35 (100%)	0	100	100
2	O	35/52 (67%)	35 (100%)	0	100	100
2	P	35/52 (67%)	35 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	Q	35/52 (67%)	35 (100%)	0	100	100
2	R	35/52 (67%)	35 (100%)	0	100	100
2	S	35/52 (67%)	35 (100%)	0	100	100
2	T	35/52 (67%)	35 (100%)	0	100	100
2	U	35/52 (67%)	35 (100%)	0	100	100
2	V	35/52 (67%)	35 (100%)	0	100	100
2	W	34/52 (65%)	34 (100%)	0	100	100
2	X	34/52 (65%)	34 (100%)	0	100	100
All	All	4594/4956 (93%)	4553 (99%)	41 (1%)	78	88

5 of 41 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	353	ARG
1	I	353	ARG
1	E	217	THR
1	C	406	THR
1	I	217	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	Y	0/12	-	-
3	Z	0/12	-	-
All	All	0/24	-	-

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 36 ligands modelled in this entry, 12 are monoatomic - leaving 24 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
4	ADP	L	1000	5,6	24,29,29	0.96	1 (4%)	29,45,45	1.44	6 (20%)
4	ADP	A	1000	5,6	24,29,29	0.94	1 (4%)	29,45,45	1.39	6 (20%)
6	BEF	H	1002	4	0,3,3	0.00	-	-		
4	ADP	E	1000	5,6	24,29,29	0.95	1 (4%)	29,45,45	1.42	5 (17%)
4	ADP	H	1000	5,6	24,29,29	0.96	1 (4%)	29,45,45	1.46	6 (20%)
4	ADP	G	1000	5,6	24,29,29	0.95	1 (4%)	29,45,45	1.40	6 (20%)
6	BEF	F	1002	4	0,3,3	0.00	-	-		
4	ADP	B	1000	5,6	24,29,29	0.96	1 (4%)	29,45,45	1.45	6 (20%)
4	ADP	J	1000	5,6	24,29,29	0.96	1 (4%)	29,45,45	1.43	6 (20%)
6	BEF	K	1002	4	0,3,3	0.00	-	-		
6	BEF	A	1002	4	0,3,3	0.00	-	-		
6	BEF	C	1002	4	0,3,3	0.00	-	-		
6	BEF	G	1002	4	0,3,3	0.00	-	-		
6	BEF	J	1002	4	0,3,3	0.00	-	-		
6	BEF	I	1002	4	0,3,3	0.00	-	-		
6	BEF	B	1002	4	0,3,3	0.00	-	-		
6	BEF	E	1002	4	0,3,3	0.00	-	-		
6	BEF	D	1002	4	0,3,3	0.00	-	-		
4	ADP	C	1000	5,6	24,29,29	0.96	1 (4%)	29,45,45	1.40	6 (20%)
6	BEF	L	1002	4	0,3,3	0.00	-	-		
4	ADP	I	1000	5,6	24,29,29	0.95	1 (4%)	29,45,45	1.44	6 (20%)
4	ADP	K	1000	5,6	24,29,29	0.94	1 (4%)	29,45,45	1.42	6 (20%)
4	ADP	D	1000	5,6	24,29,29	0.96	1 (4%)	29,45,45	1.44	6 (20%)
4	ADP	F	1000	5,6	24,29,29	0.96	1 (4%)	29,45,45	1.45	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
4	ADP	L	1000	5,6	-	0/12/32/32	0/3/3/3
4	ADP	A	1000	5,6	-	0/12/32/32	0/3/3/3
4	ADP	C	1000	5,6	-	0/12/32/32	0/3/3/3
4	ADP	E	1000	5,6	-	0/12/32/32	0/3/3/3
4	ADP	H	1000	5,6	-	0/12/32/32	0/3/3/3
4	ADP	G	1000	5,6	-	0/12/32/32	0/3/3/3
4	ADP	I	1000	5,6	-	0/12/32/32	0/3/3/3
4	ADP	K	1000	5,6	-	0/12/32/32	0/3/3/3
4	ADP	B	1000	5,6	-	0/12/32/32	0/3/3/3
4	ADP	D	1000	5,6	-	0/12/32/32	0/3/3/3
4	ADP	F	1000	5,6	-	0/12/32/32	0/3/3/3
4	ADP	J	1000	5,6	-	0/12/32/32	0/3/3/3

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	I	1000	ADP	C5-C4	2.47	1.47	1.40
4	E	1000	ADP	C5-C4	2.47	1.47	1.40
4	G	1000	ADP	C5-C4	2.45	1.47	1.40
4	C	1000	ADP	C5-C4	2.44	1.47	1.40
4	K	1000	ADP	C5-C4	2.43	1.47	1.40

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	1000	ADP	N3-C2-N1	-3.38	123.40	128.68
4	H	1000	ADP	PA-O3A-PB	-3.35	121.31	132.83
4	K	1000	ADP	N3-C2-N1	-3.34	123.46	128.68
4	G	1000	ADP	N3-C2-N1	-3.33	123.48	128.68
4	E	1000	ADP	N3-C2-N1	-3.32	123.49	128.68

There are no chirality outliers.

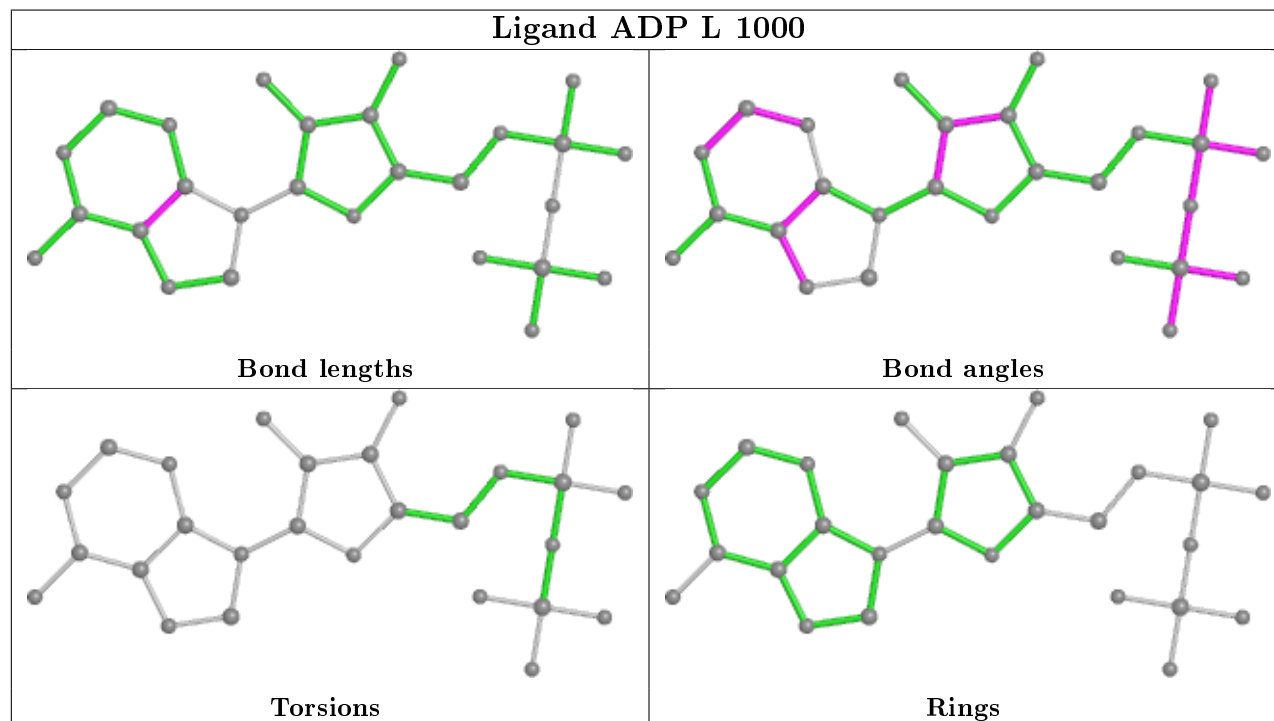
There are no torsion outliers.

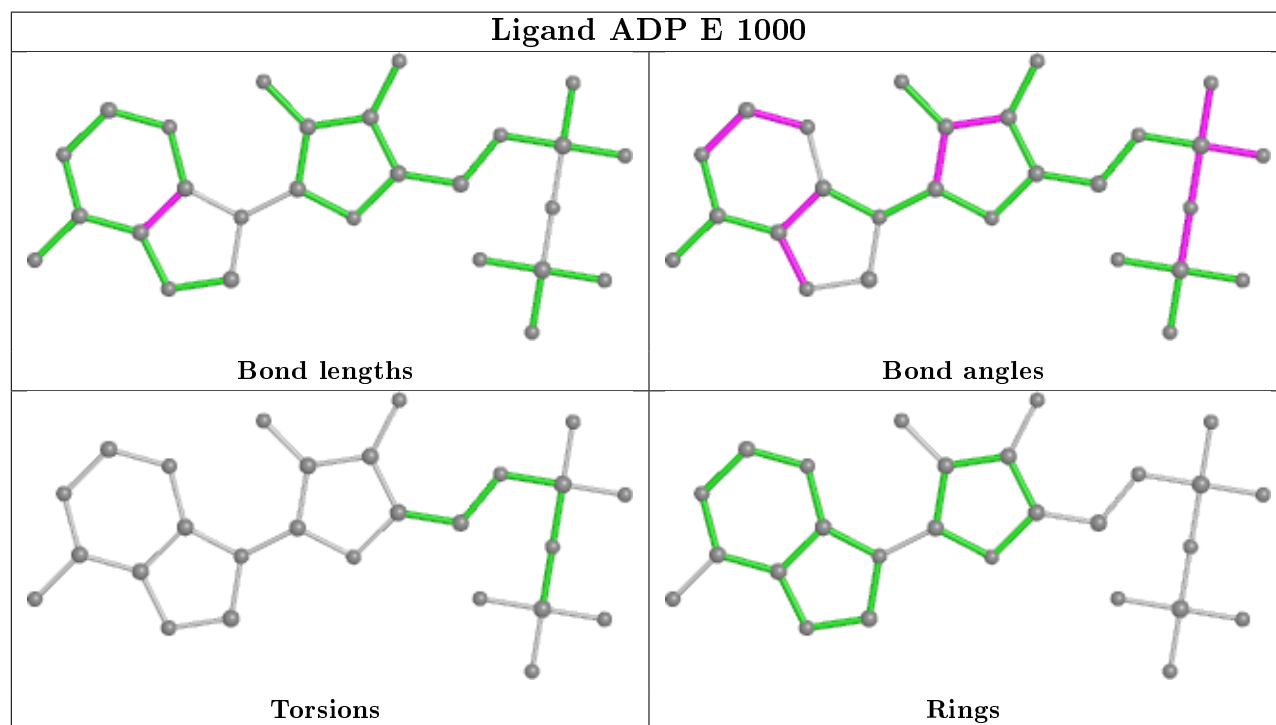
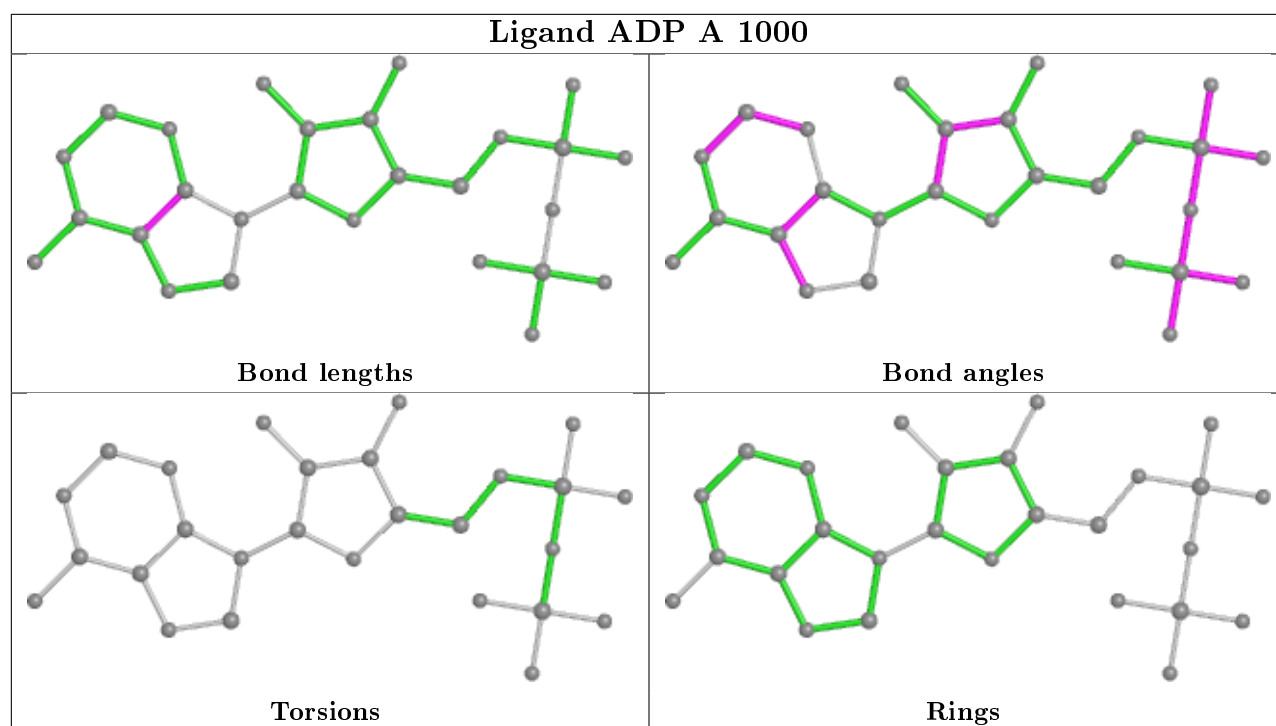
There are no ring outliers.

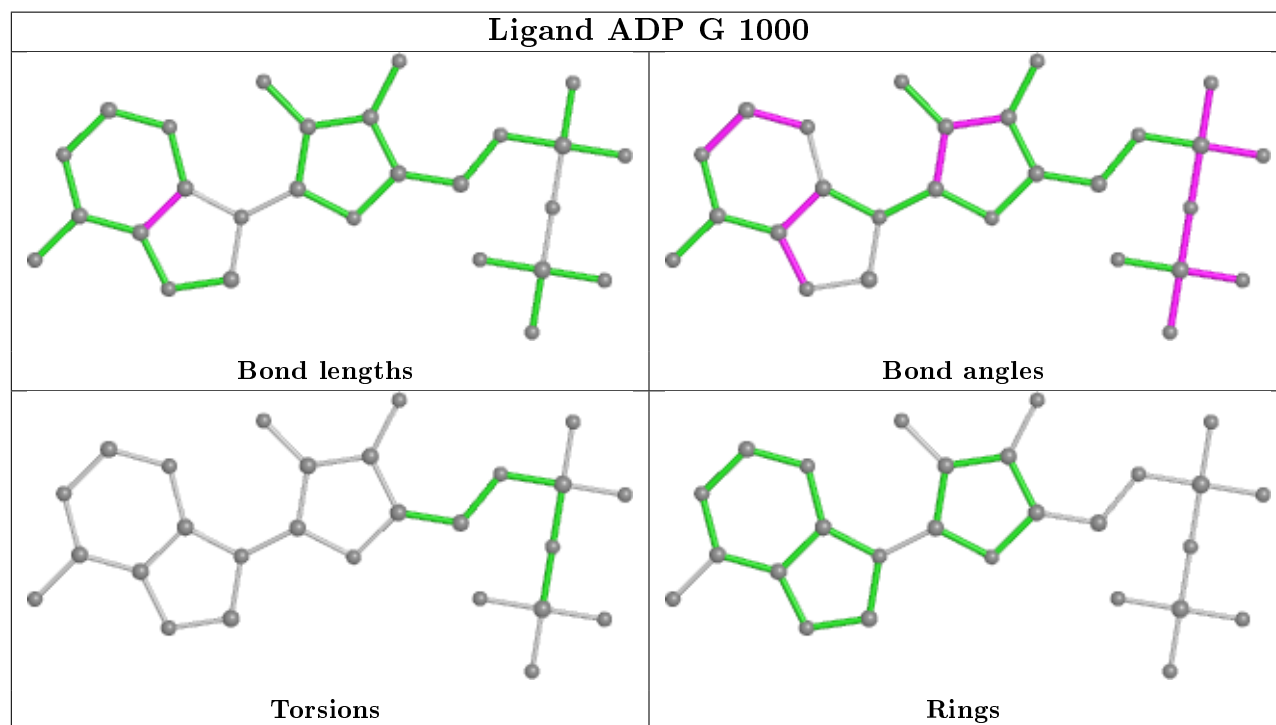
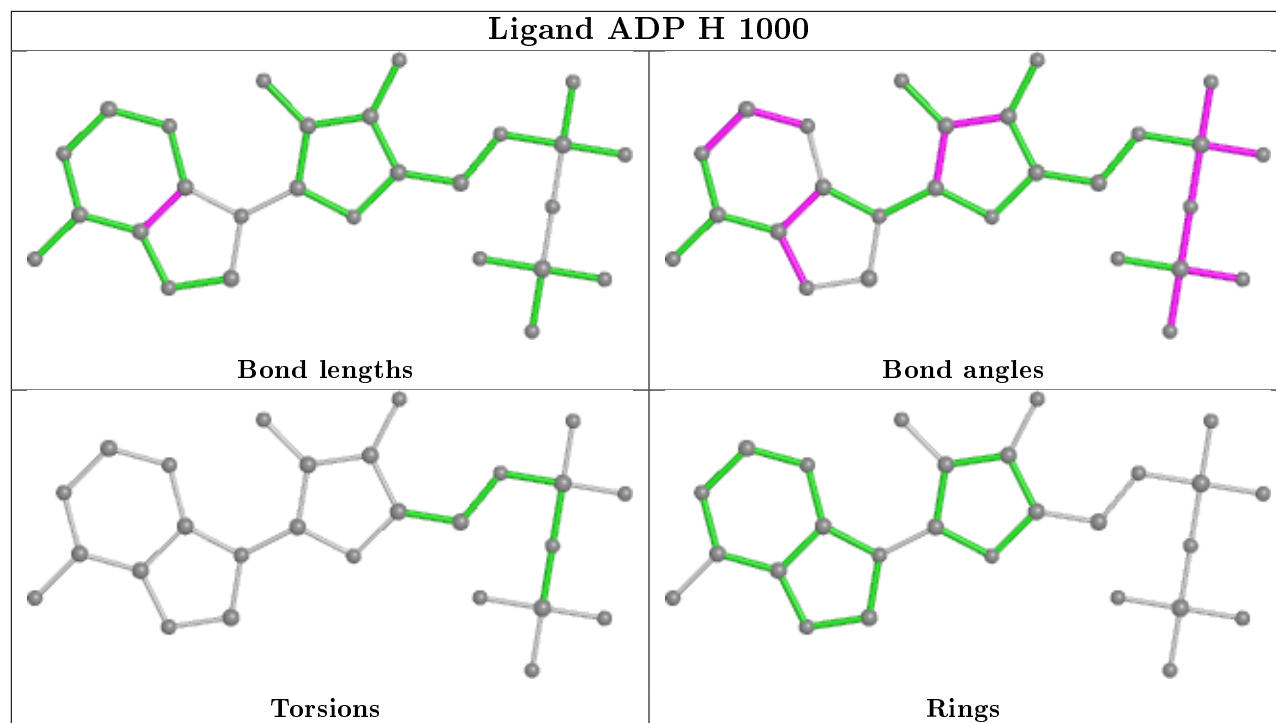
No monomer is involved in short contacts.

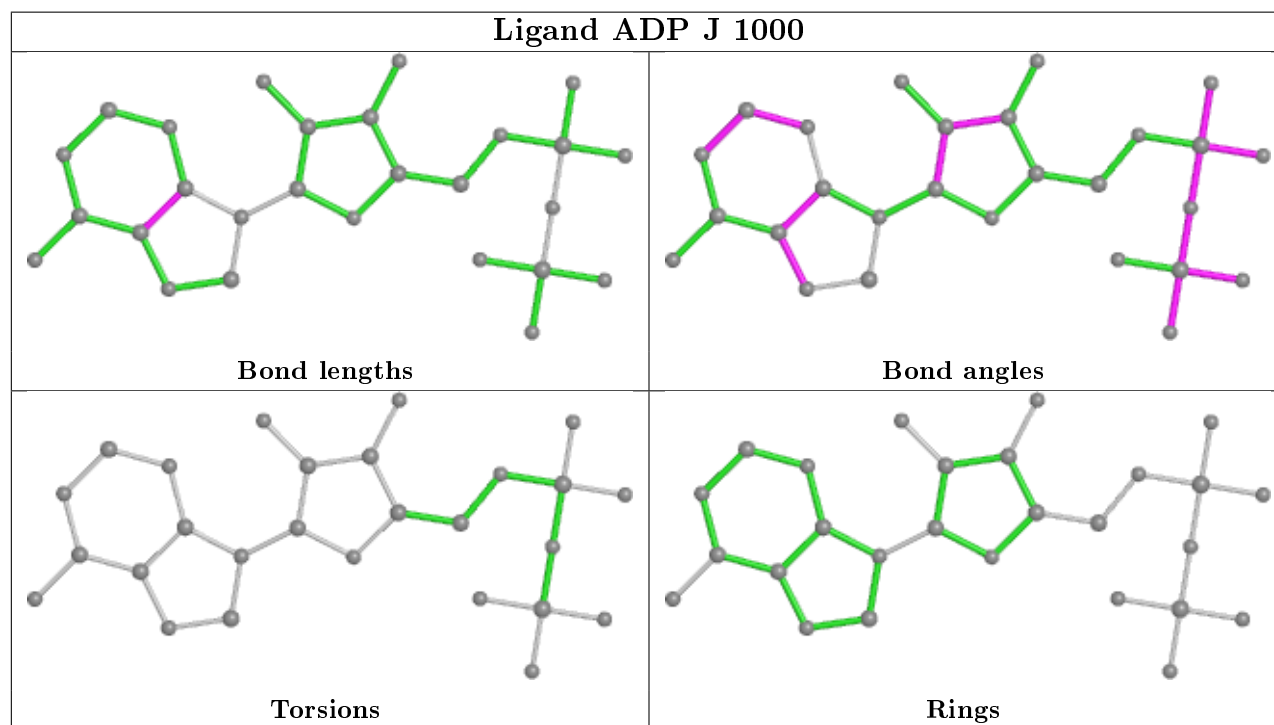
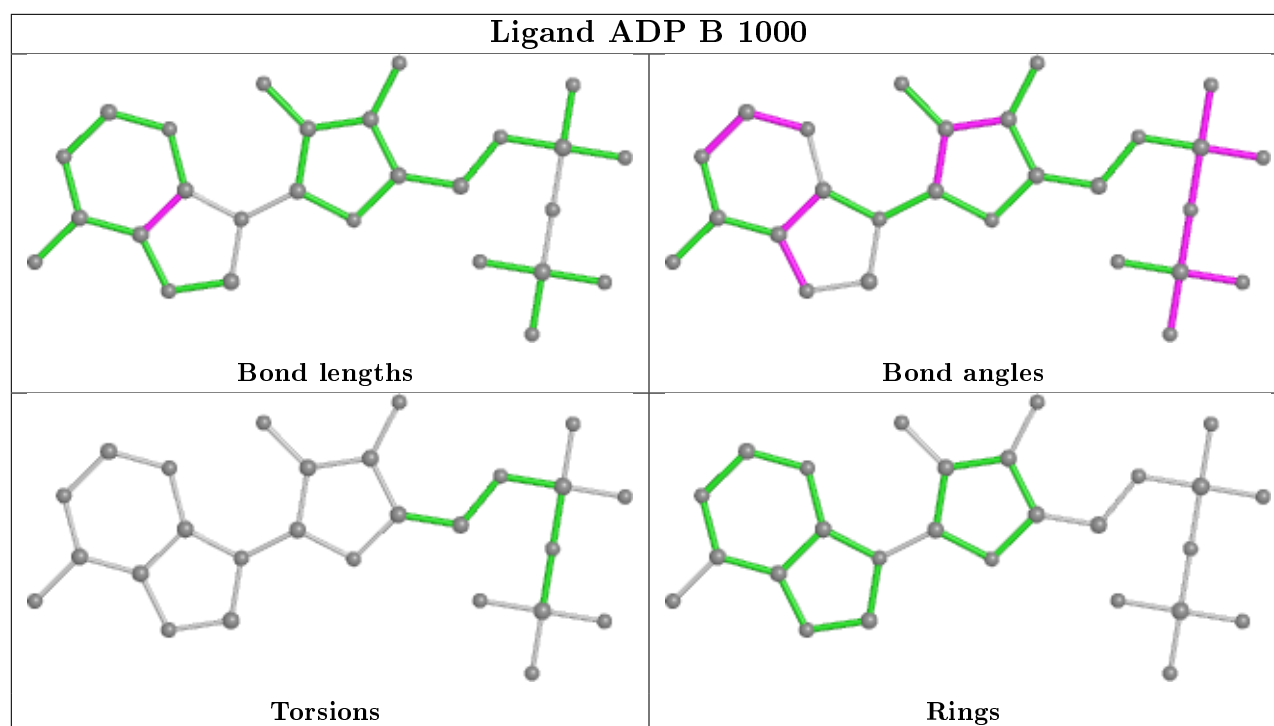
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

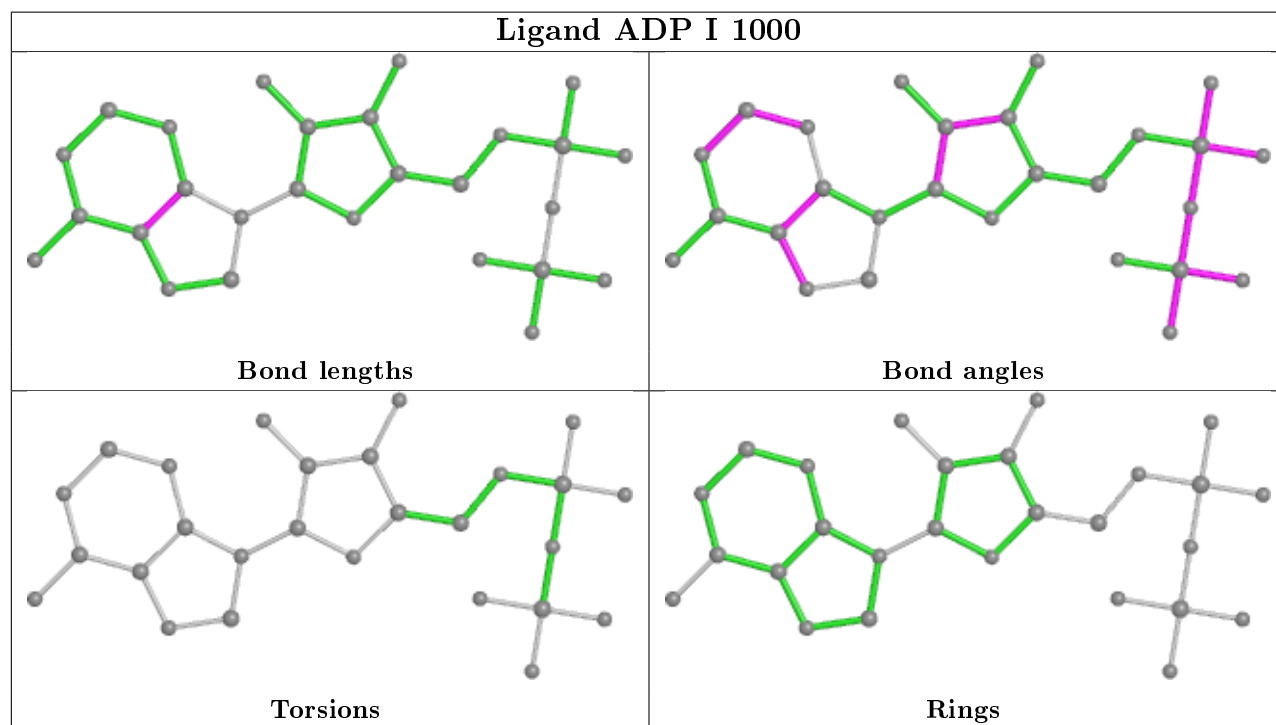
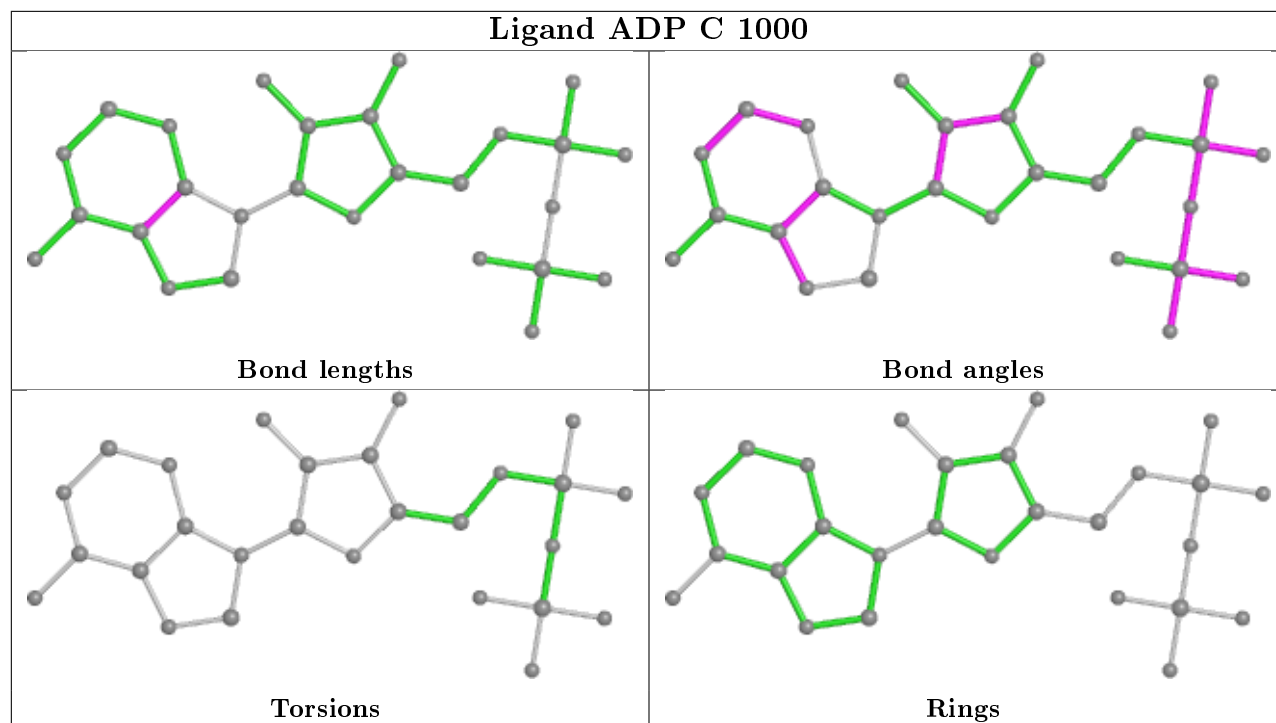
addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

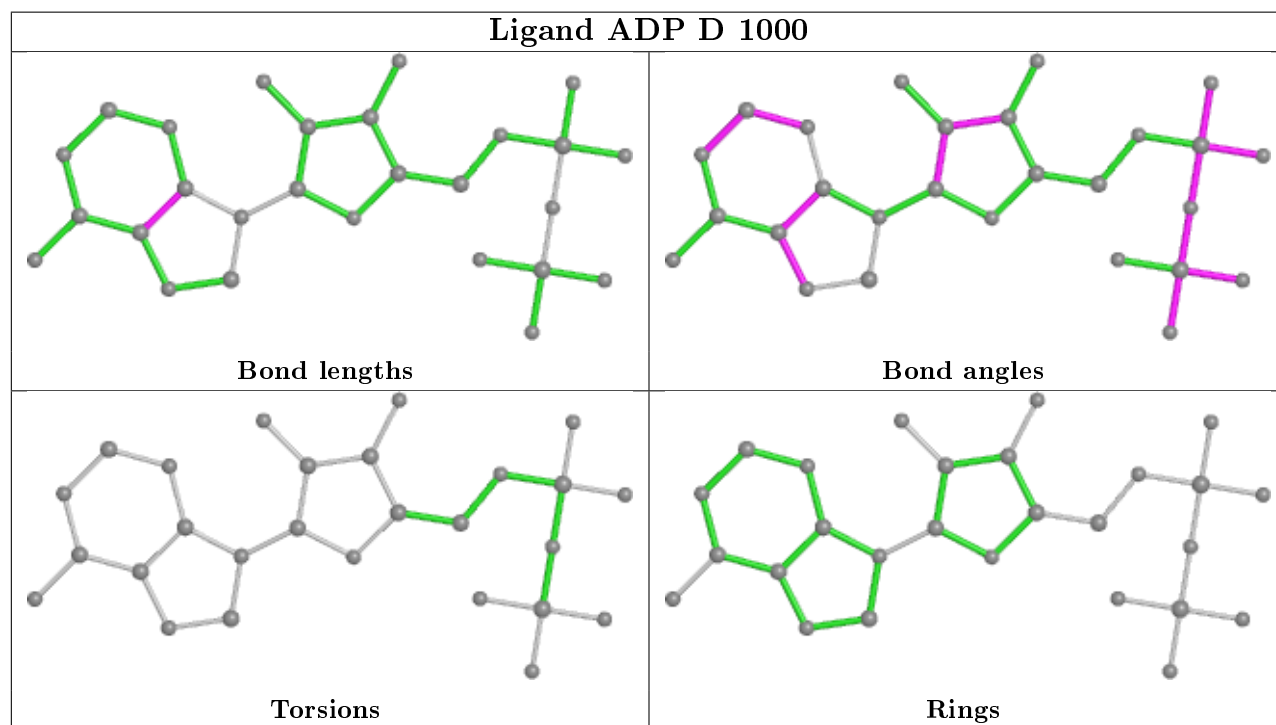
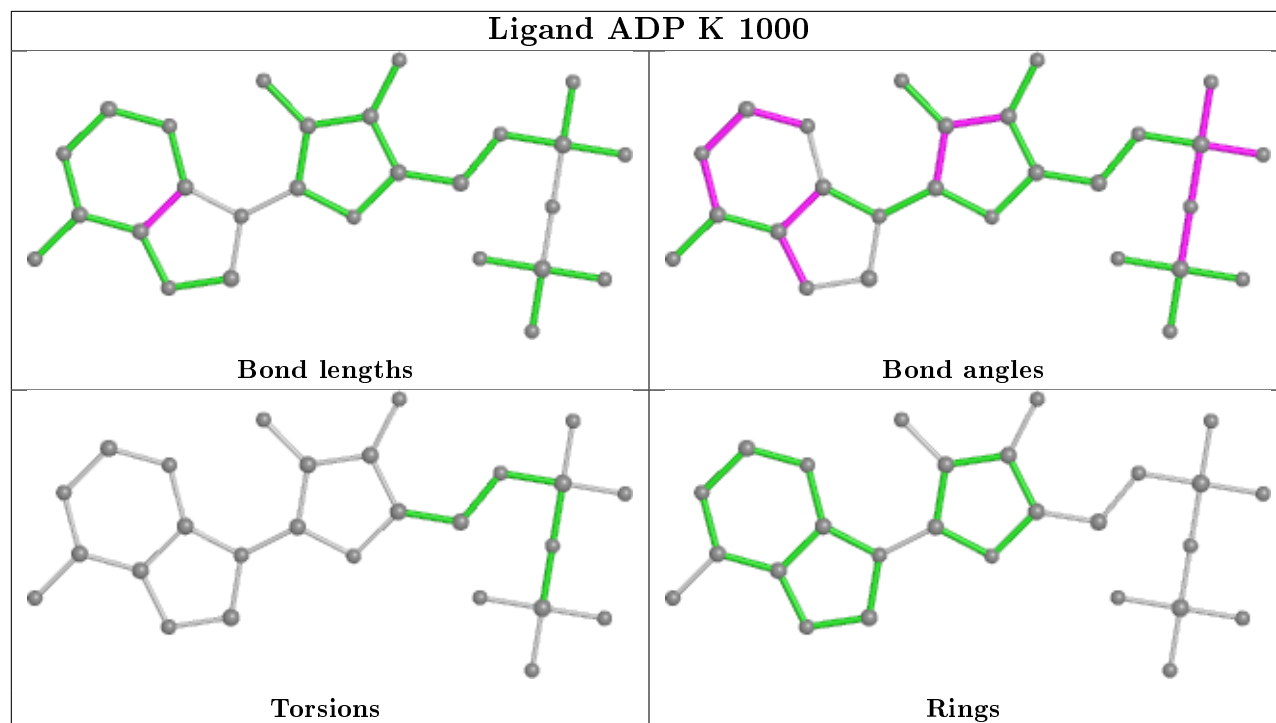


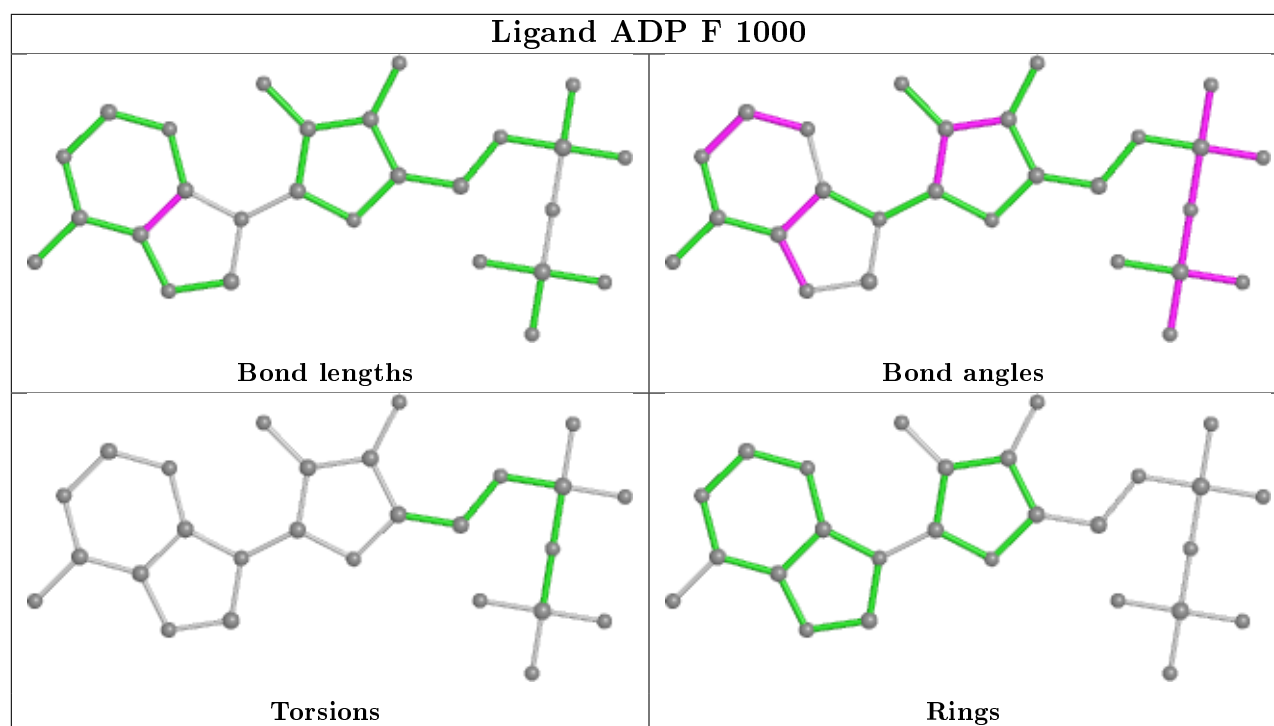












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	414/422 (98%)	0.07	20 (4%) 30 22	156, 200, 267, 301	0
1	B	409/422 (96%)	0.04	12 (2%) 51 39	110, 167, 245, 283	0
1	C	414/422 (98%)	-0.03	23 (5%) 24 17	141, 196, 260, 300	0
1	D	409/422 (96%)	-0.03	11 (2%) 54 42	119, 174, 249, 295	0
1	E	414/422 (98%)	0.16	30 (7%) 15 11	130, 219, 294, 330	0
1	F	409/422 (96%)	-0.12	11 (2%) 54 42	109, 170, 236, 273	0
1	G	414/422 (98%)	0.13	20 (4%) 30 22	158, 197, 262, 289	0
1	H	409/422 (96%)	-0.06	12 (2%) 51 39	116, 170, 252, 287	0
1	I	414/422 (98%)	0.19	31 (7%) 14 10	143, 210, 277, 312	0
1	J	409/422 (96%)	0.02	13 (3%) 47 35	114, 177, 243, 285	0
1	K	414/422 (98%)	-0.02	20 (4%) 30 22	143, 208, 271, 300	0
1	L	409/422 (96%)	0.00	10 (2%) 59 47	113, 169, 248, 293	0
2	M	43/60 (71%)	0.84	11 (25%) 0 0	219, 281, 308, 314	0
2	N	43/60 (71%)	1.11	12 (27%) 0 0	220, 291, 337, 346	0
2	O	43/60 (71%)	0.56	6 (13%) 2 3	233, 304, 350, 362	0
2	P	43/60 (71%)	0.99	11 (25%) 0 0	228, 287, 334, 351	0
2	Q	43/60 (71%)	1.36	14 (32%) 0 0	217, 305, 355, 367	0
2	R	43/60 (71%)	1.13	11 (25%) 0 0	233, 299, 356, 372	0
2	S	43/60 (71%)	0.72	7 (16%) 1 1	287, 329, 349, 357	0
2	T	43/60 (71%)	1.06	10 (23%) 0 0	303, 322, 348, 354	0
2	U	43/60 (71%)	1.43	16 (37%) 0 0	323, 350, 360, 363	0
2	V	43/60 (71%)	1.62	20 (46%) 0 0	309, 337, 364, 373	0
2	W	43/60 (71%)	1.74	10 (23%) 0 0	318, 340, 356, 363	0
2	X	43/60 (71%)	0.61	9 (20%) 1 0	335, 355, 374, 376	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å²)	Q<0.9
3	Y	6/12 (50%)	1.34	1 (16%) 1 1	50, 59, 70, 70	0
3	Z	6/12 (50%)	1.22	1 (16%) 1 1	50, 57, 60, 65	0
All	All	5466/5808 (94%)	0.13	352 (6%) 19 13	50, 193, 323, 376	0

The worst 5 of 352 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	W	161	SER	11.6
2	Q	150	GLU	9.9
2	U	181	ALA	9.5
2	W	150	GLU	9.2
2	R	150	GLU	9.1

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates 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, 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
6	BEF	E	1002	4/4	0.67	0.20	162,178,182,194	0
6	BEF	I	1002	4/4	0.79	0.26	158,158,163,176	0
6	BEF	C	1002	4/4	0.83	0.25	151,152,156,160	0
6	BEF	G	1002	4/4	0.84	0.24	157,164,172,177	0
4	ADP	E	1000	27/27	0.87	0.20	162,186,230,258	0
6	BEF	A	1002	4/4	0.89	0.24	156,172,179,191	0
4	ADP	K	1000	27/27	0.91	0.25	155,176,228,246	0
6	BEF	L	1002	4/4	0.92	0.28	142,142,142,152	0
4	ADP	J	1000	27/27	0.92	0.23	142,147,173,190	0
5	MG	D	1001	1/1	0.92	0.23	132,132,132,132	0

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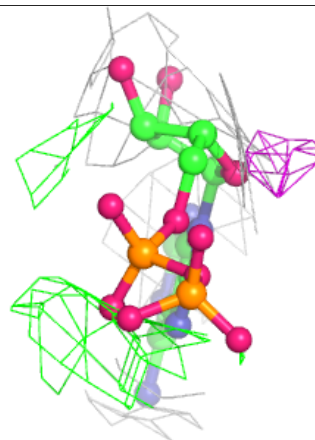
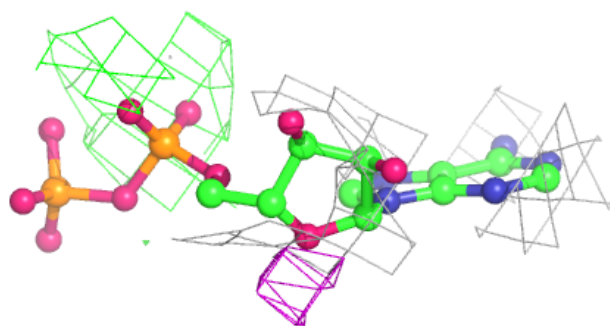
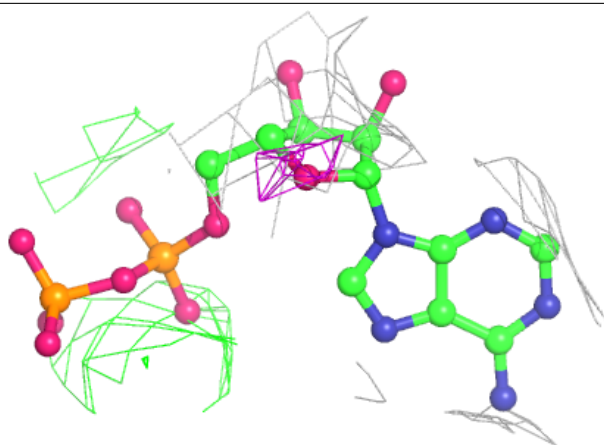
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
6	BEF	K	1002	4/4	0.93	0.17	155,172,183,192	0
4	ADP	G	1000	27/27	0.93	0.23	157,174,210,230	0
4	ADP	C	1000	27/27	0.93	0.26	151,165,204,232	0
4	ADP	I	1000	27/27	0.94	0.17	158,175,215,221	0
6	BEF	J	1002	4/4	0.94	0.16	142,142,142,142	0
4	ADP	F	1000	27/27	0.94	0.23	142,147,170,189	0
4	ADP	B	1000	27/27	0.94	0.25	133,133,154,177	0
4	ADP	L	1000	27/27	0.95	0.23	142,144,165,188	0
4	ADP	A	1000	27/27	0.95	0.20	156,171,210,235	0
6	BEF	B	1002	4/4	0.95	0.27	133,133,133,133	0
4	ADP	D	1000	27/27	0.95	0.22	132,132,157,181	0
6	BEF	H	1002	4/4	0.95	0.30	134,134,134,135	0
4	ADP	H	1000	27/27	0.95	0.25	134,135,156,193	0
5	MG	E	1001	1/1	0.97	0.29	162,162,162,162	0
5	MG	I	1001	1/1	0.97	0.28	158,158,158,158	0
6	BEF	D	1002	4/4	0.97	0.27	132,132,132,132	0
5	MG	G	1001	1/1	0.97	0.34	157,157,157,157	0
5	MG	K	1001	1/1	0.97	0.16	155,155,155,155	0
5	MG	F	1001	1/1	0.98	0.18	142,142,142,142	0
5	MG	H	1001	1/1	0.98	0.34	134,134,134,134	0
5	MG	A	1001	1/1	0.99	0.25	156,156,156,156	0
5	MG	C	1001	1/1	0.99	0.33	151,151,151,151	0
5	MG	J	1001	1/1	0.99	0.23	142,142,142,142	0
6	BEF	F	1002	4/4	0.99	0.17	142,142,142,142	0
5	MG	L	1001	1/1	0.99	0.30	142,142,142,142	0
5	MG	B	1001	1/1	0.99	0.33	133,133,133,133	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.

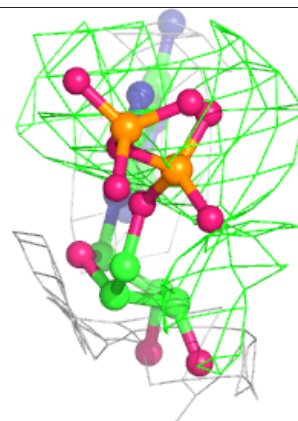
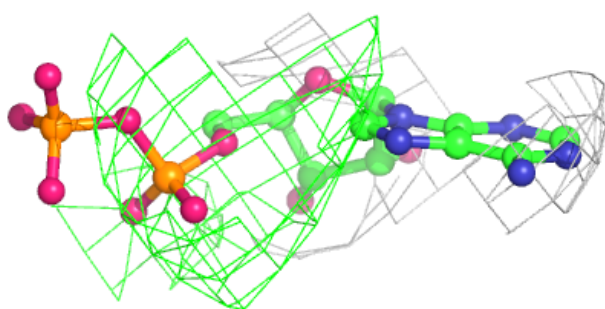
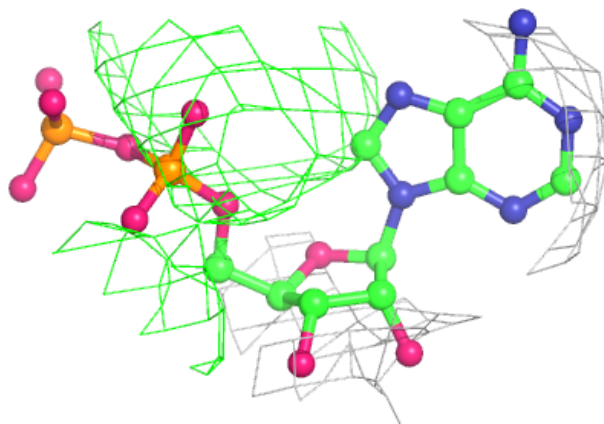
Electron density around ADP E 1000:

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



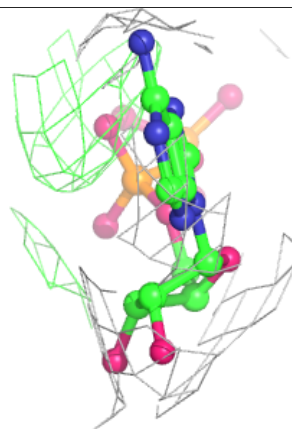
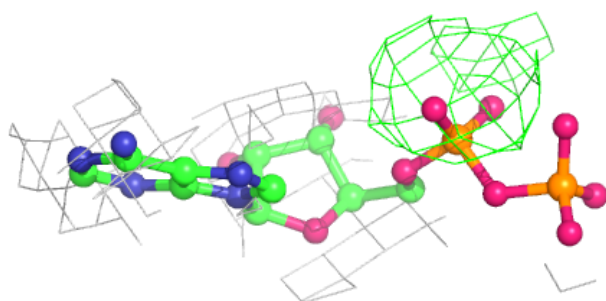
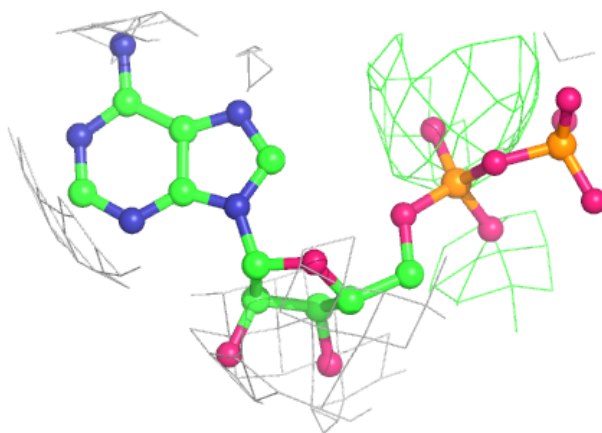
Electron density around ADP K 1000:

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

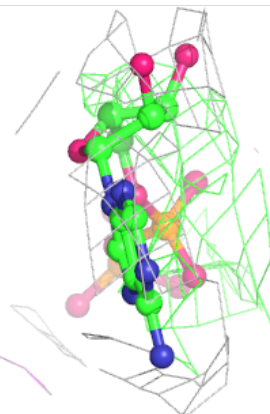
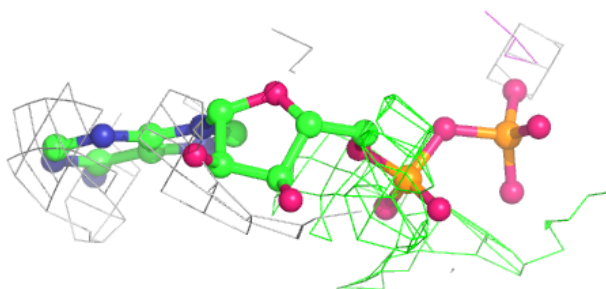
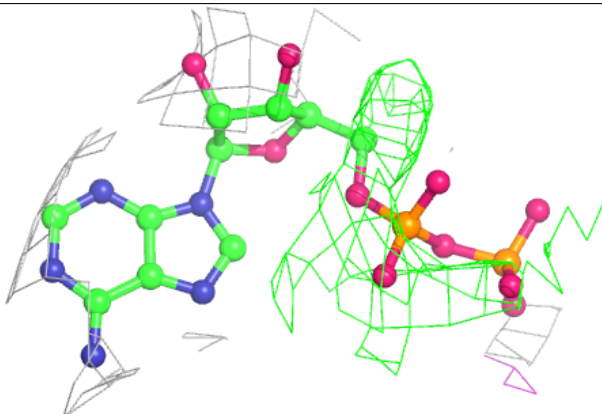


Electron density around ADP J 1000:

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

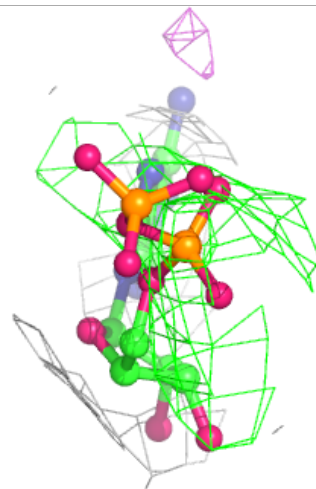
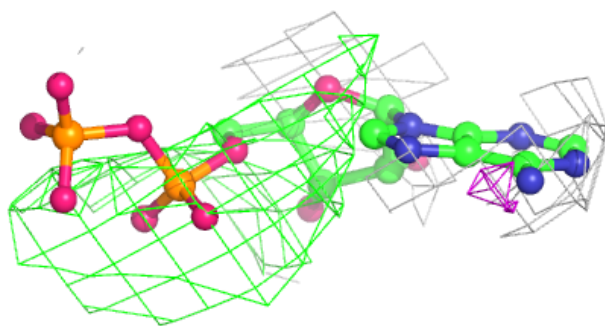
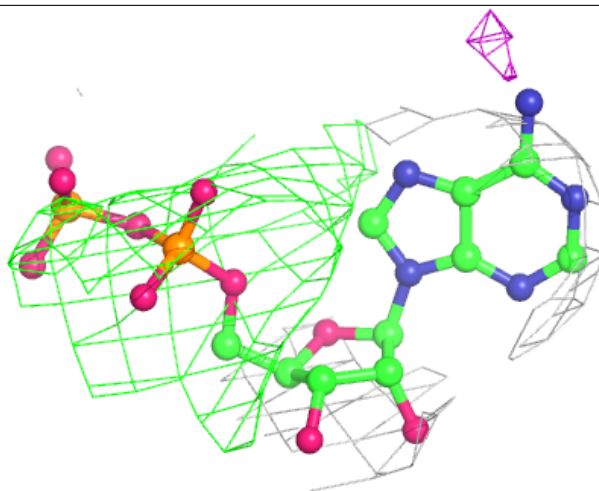
**Electron density around ADP G 1000:**

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



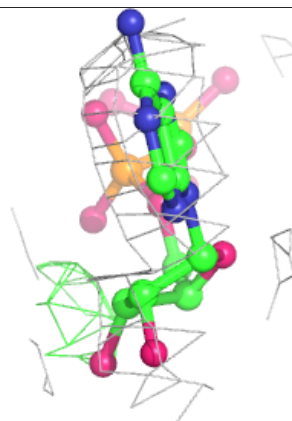
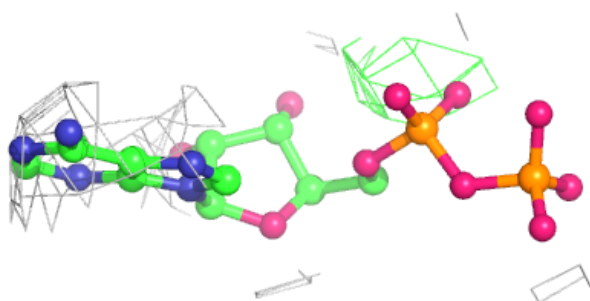
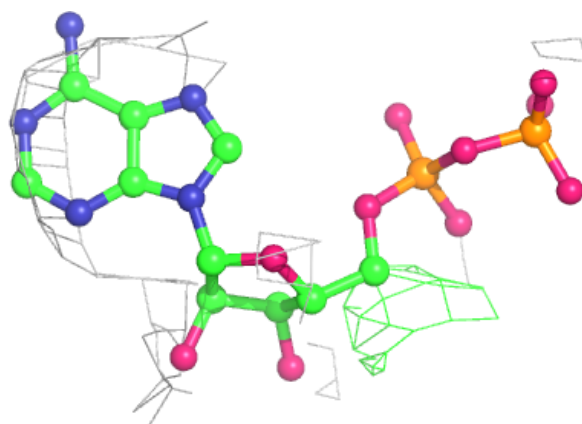
Electron density around ADP C 1000:

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



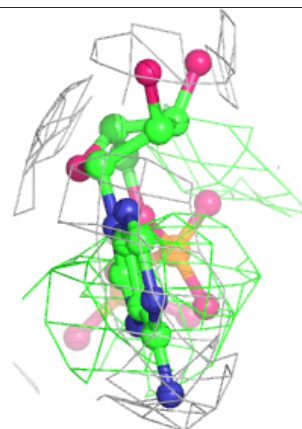
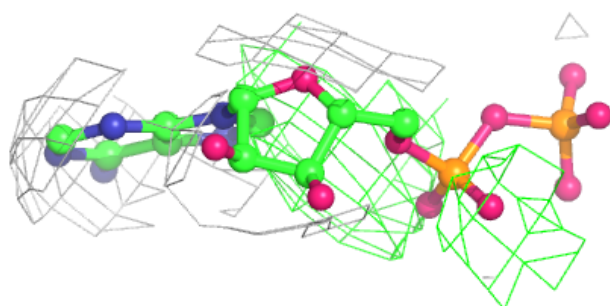
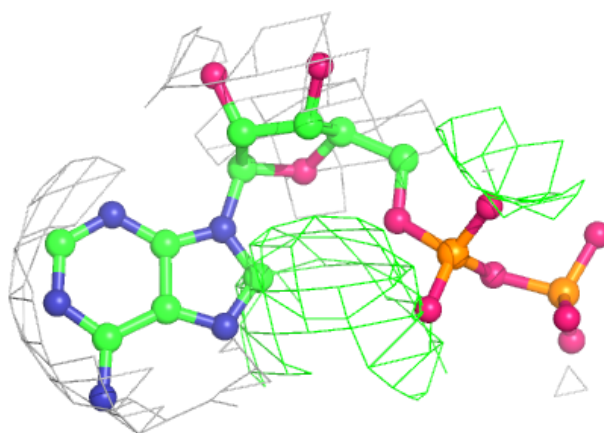
Electron density around ADP I 1000:

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

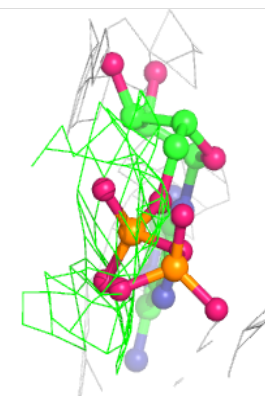
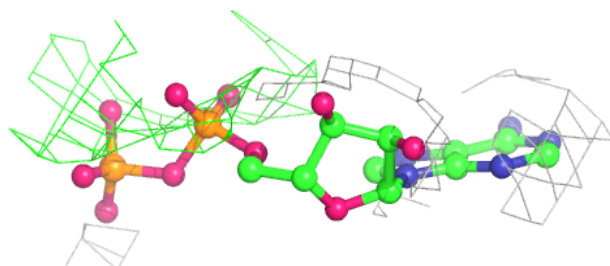
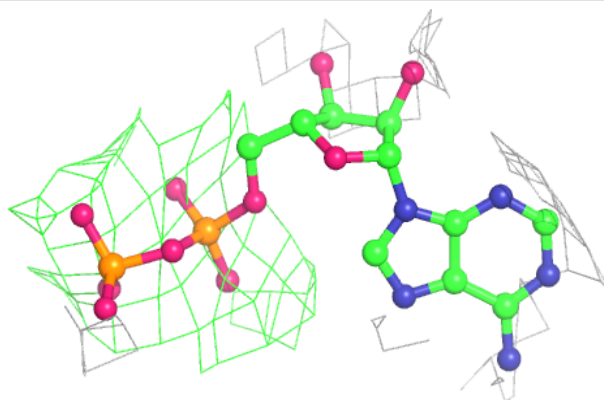


Electron density around ADP F 1000:

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

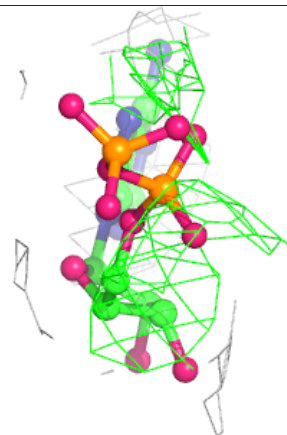
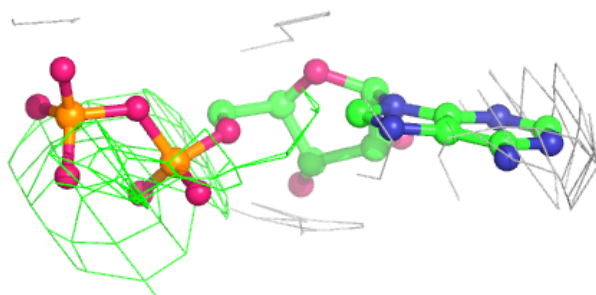
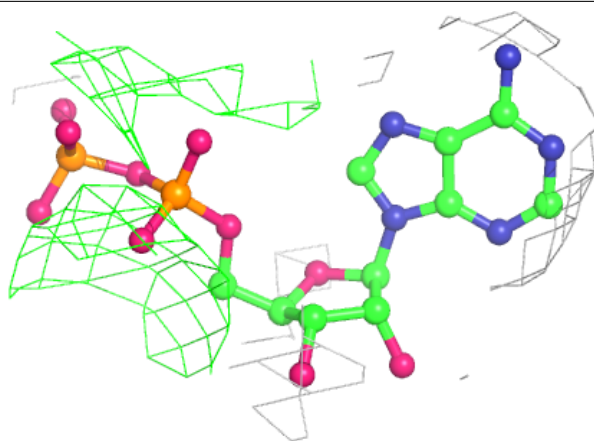
**Electron density around ADP B 1000:**

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



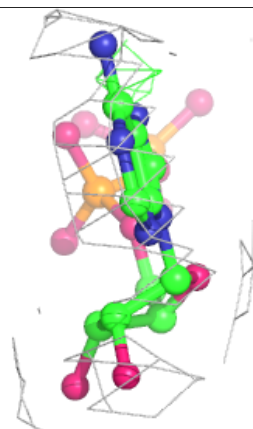
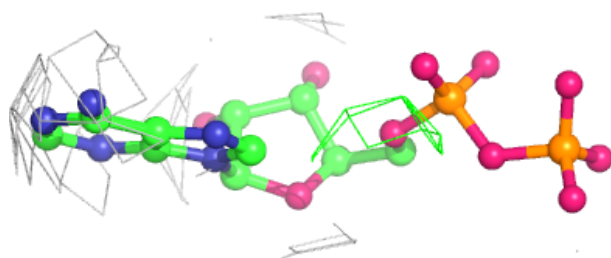
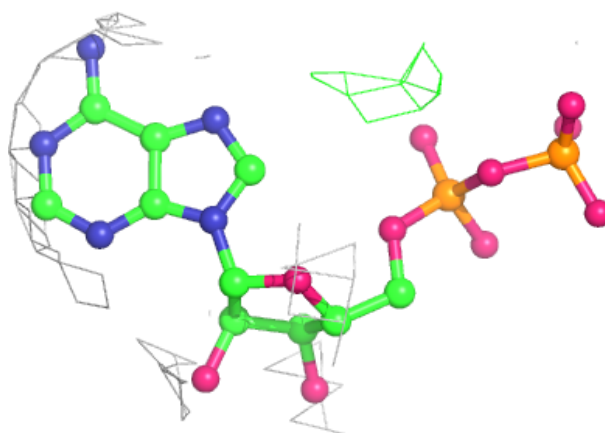
Electron density around ADP L 1000:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

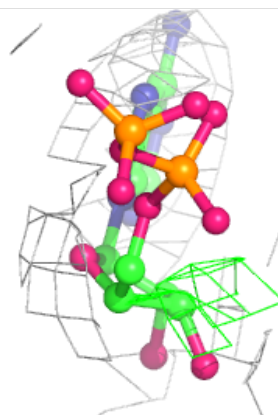
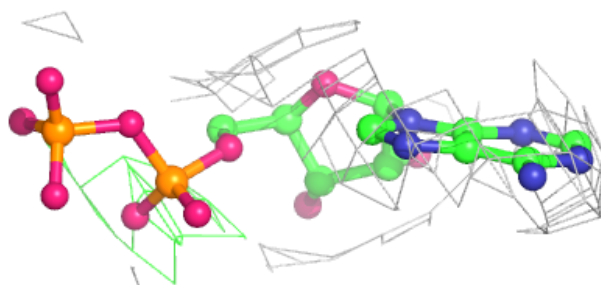
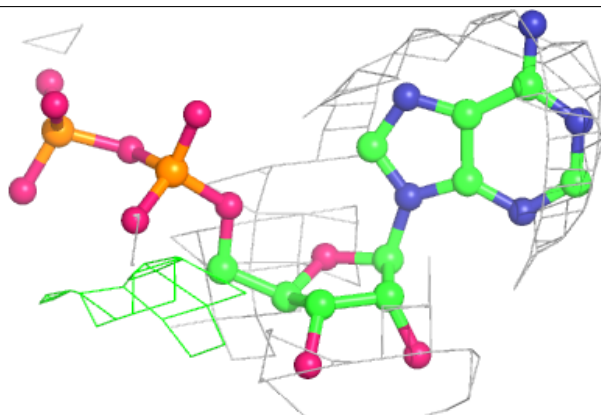


Electron density around ADP A 1000:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

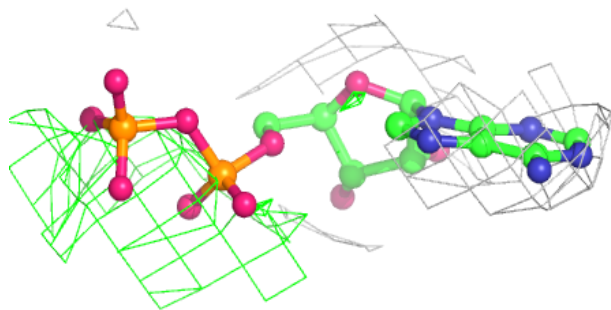
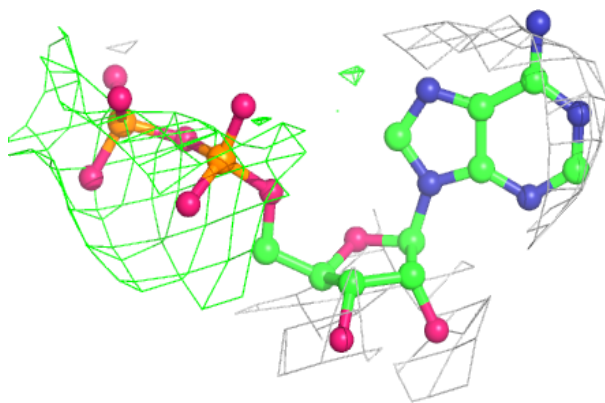
**Electron density around ADP D 1000:**

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



Electron density around ADP H 1000:

$2mF_o-DF_c$ (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.