



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 06:28 PM UTC

PDB ID : 2X86 / pdb_00002x86
Title : AGME bound to ADP-B-mannose
Authors : Kowatz, T.; Morrison, J.P.; Tanner, M.E.; Naismith, J.H.
Deposited on : 2010-03-06
Resolution : 2.80 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 types of validation reports are described at

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

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)
Xtriage (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

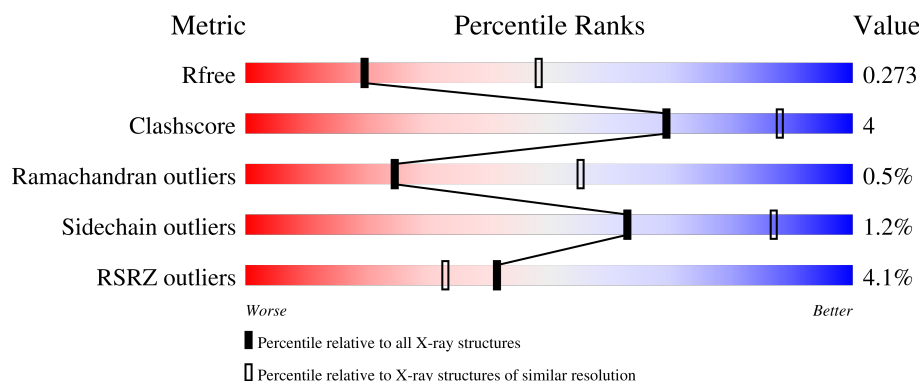
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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 ≥ 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	357	
1	B	357	
1	C	357	
1	D	357	
1	E	357	

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Mol	Chain	Length	Quality of chain
1	F	357	
1	G	357	
1	H	357	
1	I	357	
1	J	357	
1	K	357	
1	L	357	
1	M	357	
1	N	357	
1	O	357	
1	P	357	
1	Q	357	
1	R	357	
1	S	357	
1	T	357	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 51434 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 ADP-L-GLYCERO-D-MANNO-HEPTOSE-6-EPIMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	B	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	C	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	D	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	E	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	F	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	G	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	H	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	I	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	J	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	K	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	L	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	M	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	N	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	O	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	P	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			

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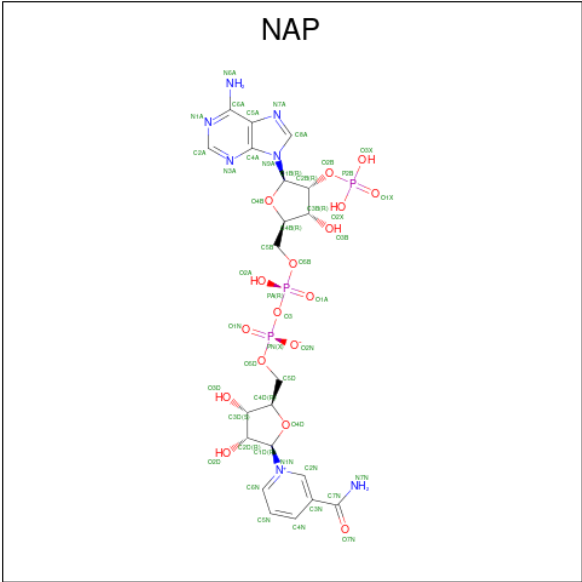
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	R	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	S	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			
1	T	307	Total	C	N	O	S	0	0	0
			2440	1566	396	469	9			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	140	PHE	TYR	engineered mutation	UNP P67911
B	140	PHE	TYR	engineered mutation	UNP P67911
C	140	PHE	TYR	engineered mutation	UNP P67911
D	140	PHE	TYR	engineered mutation	UNP P67911
E	140	PHE	TYR	engineered mutation	UNP P67911
F	140	PHE	TYR	engineered mutation	UNP P67911
G	140	PHE	TYR	engineered mutation	UNP P67911
H	140	PHE	TYR	engineered mutation	UNP P67911
I	140	PHE	TYR	engineered mutation	UNP P67911
J	140	PHE	TYR	engineered mutation	UNP P67911
K	140	PHE	TYR	engineered mutation	UNP P67911
L	140	PHE	TYR	engineered mutation	UNP P67911
M	140	PHE	TYR	engineered mutation	UNP P67911
N	140	PHE	TYR	engineered mutation	UNP P67911
O	140	PHE	TYR	engineered mutation	UNP P67911
P	140	PHE	TYR	engineered mutation	UNP P67911
Q	140	PHE	TYR	engineered mutation	UNP P67911
R	140	PHE	TYR	engineered mutation	UNP P67911
S	140	PHE	TYR	engineered mutation	UNP P67911
T	140	PHE	TYR	engineered mutation	UNP P67911

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



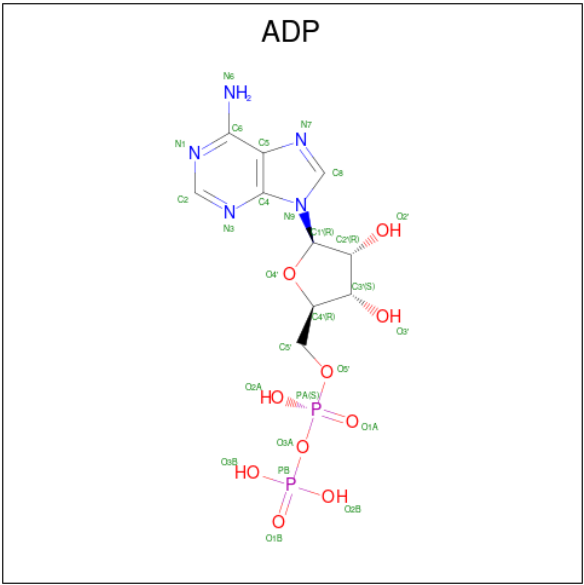
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	D	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	E	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	F	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	G	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	H	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	I	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	J	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	K	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	L	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	M	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	N	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	O	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	P	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	Q	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	R	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	S	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	T	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



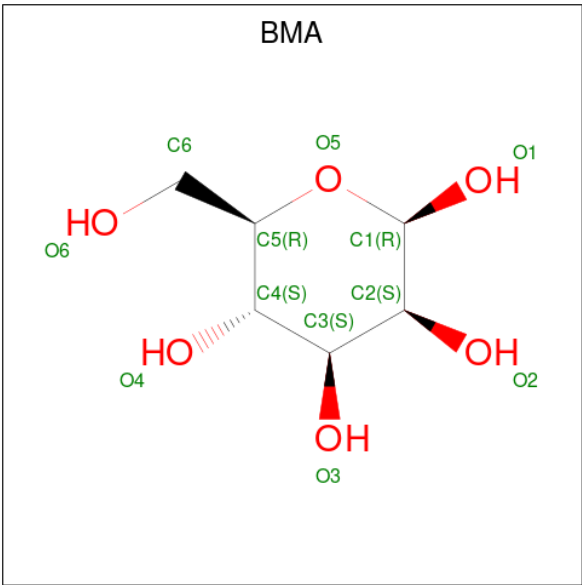
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	D	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	E	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	F	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	G	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	H	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	I	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	J	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	K	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	L	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	M	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	N	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	O	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	P	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	Q	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	R	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	S	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	T	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 4 is beta-D-mannopyranose (CCD ID: BMA) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			11	6	5		
4	C	1	Total	C	O	0	0
			11	6	5		
4	D	1	Total	C	O	0	0
			11	6	5		
4	E	1	Total	C	O	0	0
			11	6	5		
4	F	1	Total	C	O	0	0
			11	6	5		
4	G	1	Total	C	O	0	0
			11	6	5		
4	H	1	Total	C	O	0	0
			11	6	5		
4	I	1	Total	C	O	0	0
			11	6	5		
4	J	1	Total	C	O	0	0
			11	6	5		
4	K	1	Total	C	O	0	0
			11	6	5		
4	L	1	Total	C	O	0	0
			11	6	5		
4	M	1	Total	C	O	0	0
			11	6	5		
4	N	1	Total	C	O	0	0
			11	6	5		
4	O	1	Total	C	O	0	0
			11	6	5		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	P	1	Total	C	O	0	0
			11	6	5		
4	Q	1	Total	C	O	0	0
			11	6	5		
4	R	1	Total	C	O	0	0
			11	6	5		
4	S	1	Total	C	O	0	0
			11	6	5		
4	T	1	Total	C	O	0	0
			11	6	5		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	59	Total	O	0	0
			59	59		
5	B	36	Total	O	0	0
			36	36		
5	C	34	Total	O	0	0
			34	34		
5	D	60	Total	O	0	0
			60	60		
5	E	61	Total	O	0	0
			61	61		
5	F	56	Total	O	0	0
			56	56		
5	G	28	Total	O	0	0
			28	28		
5	H	49	Total	O	0	0
			49	49		
5	I	78	Total	O	0	0
			78	78		
5	J	66	Total	O	0	0
			66	66		
5	K	59	Total	O	0	0
			59	59		
5	L	51	Total	O	0	0
			51	51		
5	M	45	Total	O	0	0
			45	45		
5	N	39	Total	O	0	0
			39	39		

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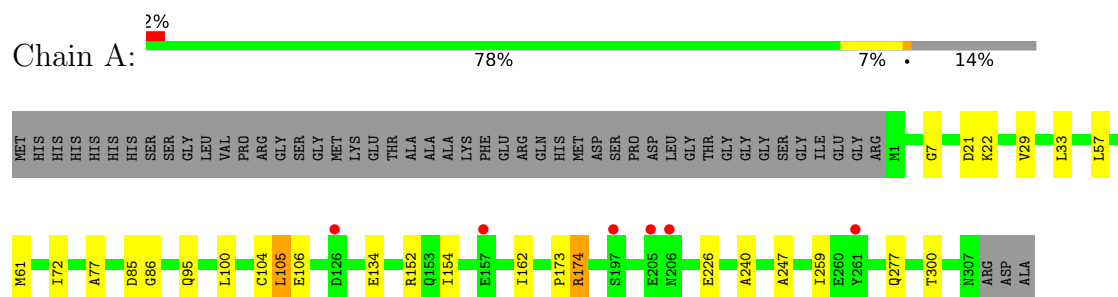
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	O	33	Total 33	O 33	0	0
5	P	49	Total 49	O 49	0	0
5	Q	50	Total 50	O 50	0	0
5	R	8	Total 8	O 8	0	0
5	S	20	Total 20	O 20	0	0
5	T	44	Total 44	O 44	0	0

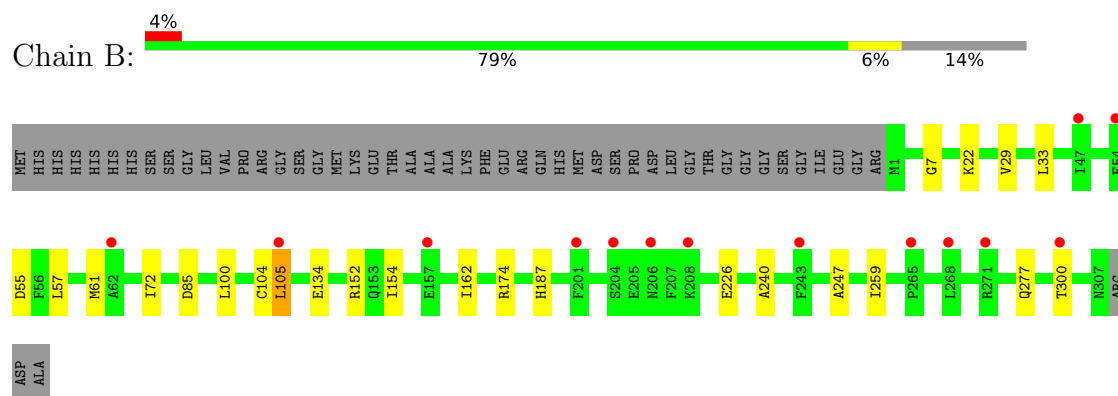
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

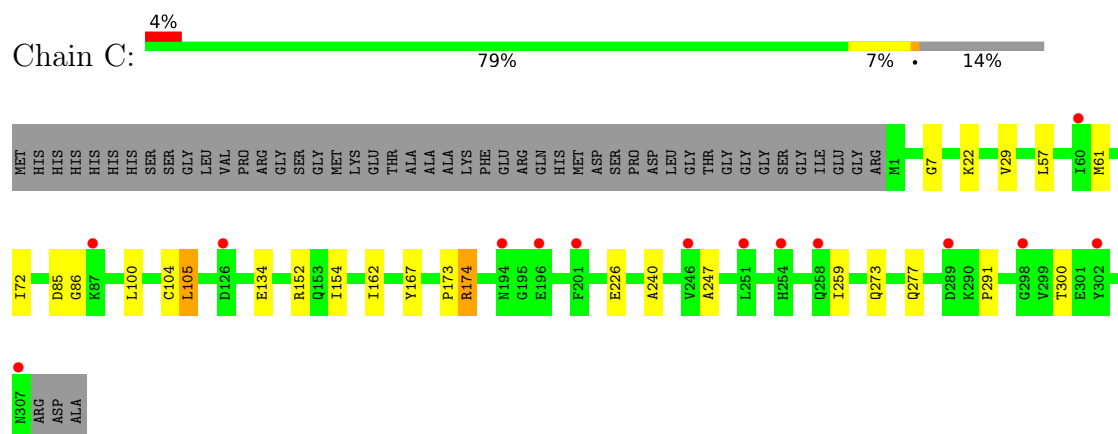
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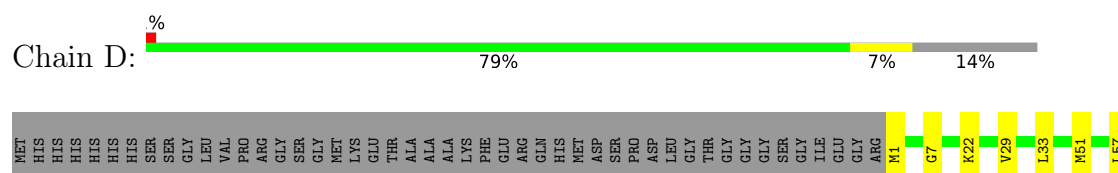
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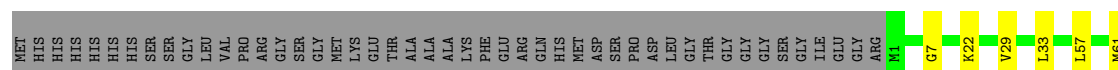
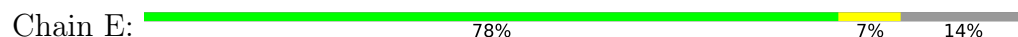
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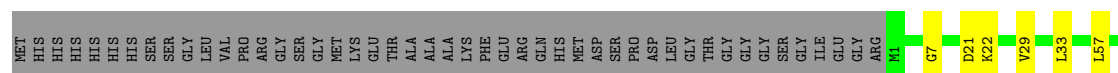
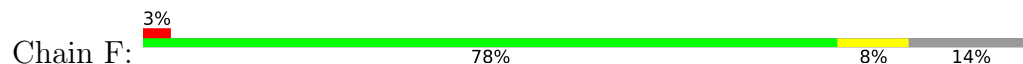
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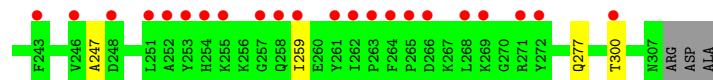
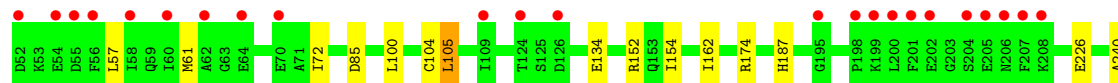
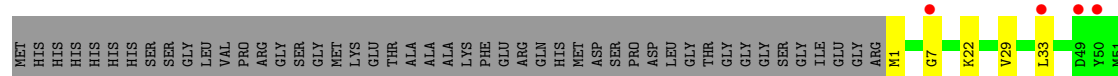
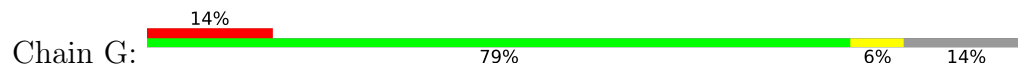
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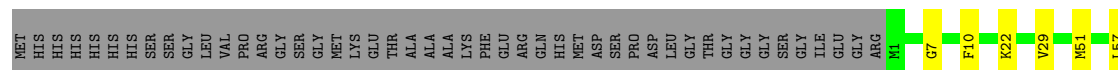
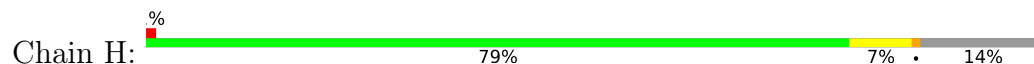
• Molecule 1: ADP-L-GLYCERO-D-MANNO-HEPTOSE-6-EPIMERASE



• Molecule 1: ADP-L-GLYCERO-D-MANNO-HEPTOSE-6-EPIMERASE



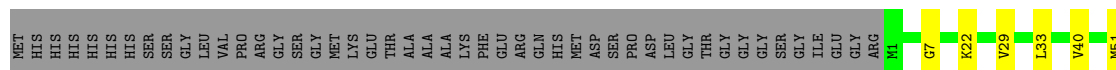
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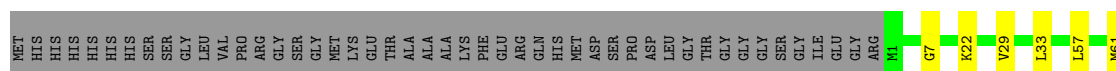
- Molecule 1: ADP-L-GLYCERO-D-MANNO-HEPTOSE-6-EPIMERASE

Chain I: 78% 7% 14%



- Molecule 1: ADP-L-GLYCERO-D-MANNO-HEPTOSE-6-EPIMERASE

Chain J: 79% 6% 14%



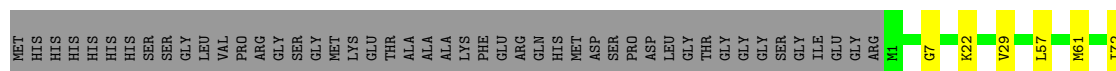
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Chain K: 2% 77% 8% 14%



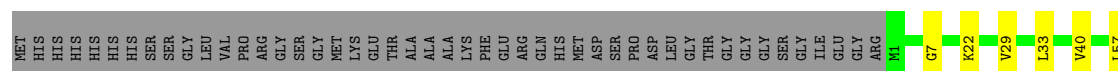
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Chain L: % 78% 7% 14%

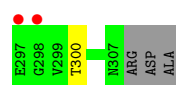
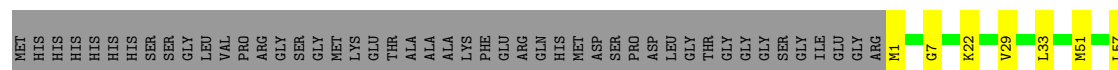
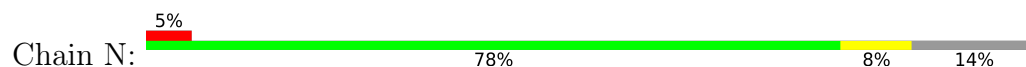


- Molecule 1: ADP-L-GLYCERO-D-MANNO-HEPTOSE-6-EPIMERASE

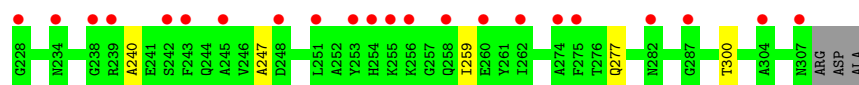
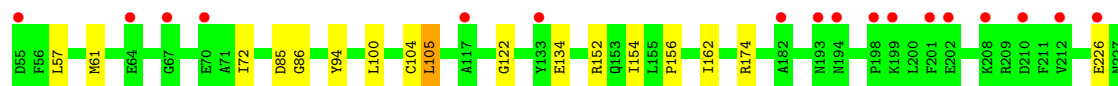
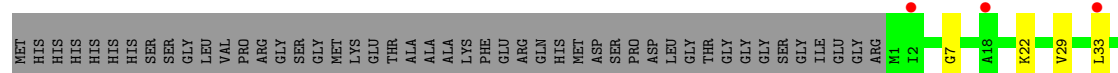
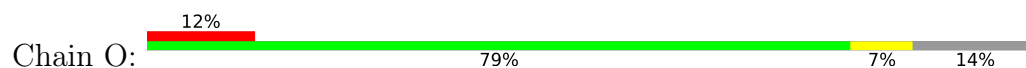
Chain M: 77% 9% 14%



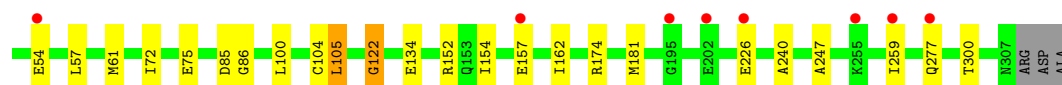
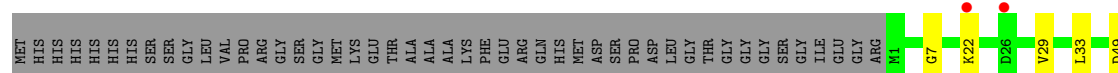
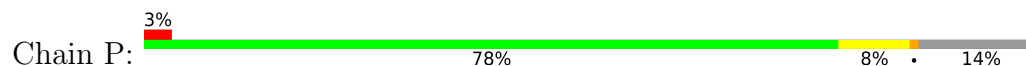
• Molecule 1: ADP-L-GLYCERO-D-MANNO-HEPTOSE-6-EPIMERASE



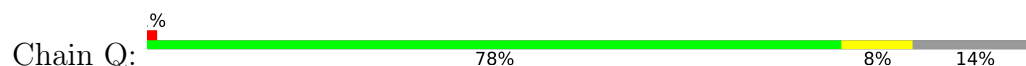
• Molecule 1: ADP-L-GLYCERO-D-MANNO-HEPTOSE-6-EPIMERASE

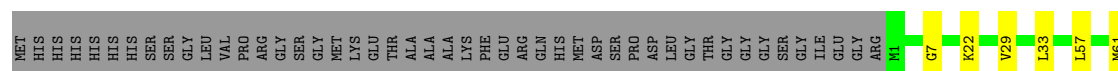


• Molecule 1: ADP-L-GLYCERO-D-MANNO-HEPTOSE-6-EPIMERASE

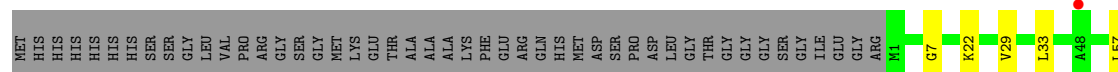
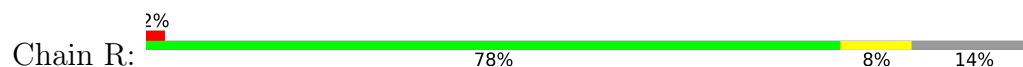


• Molecule 1: ADP-L-GLYCERO-D-MANNO-HEPTOSE-6-EPIMERASE

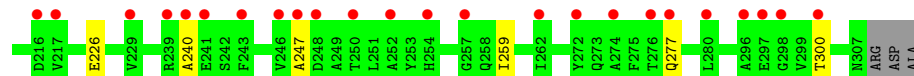
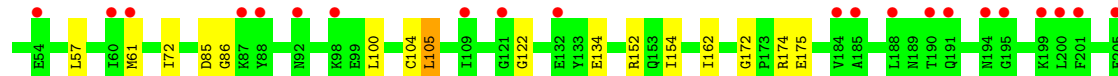
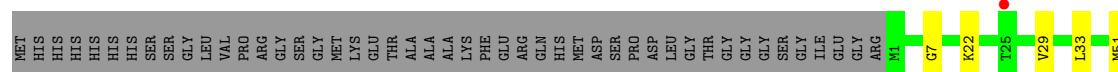
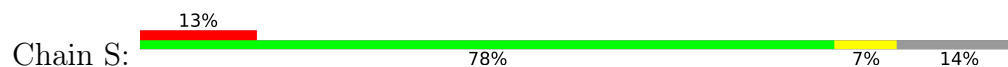




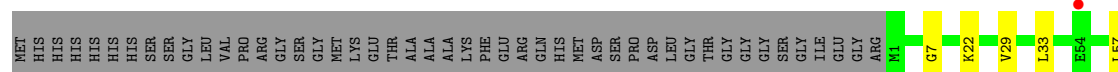
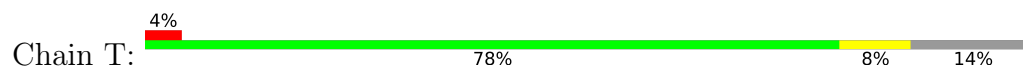
• Molecule 1: ADP-L-GLYCERO-D-MANNO-HEPTOSE-6-EPIMERASE



• Molecule 1: ADP-L-GLYCERO-D-MANNO-HEPTOSE-6-EPIMERASE



• Molecule 1: ADP-L-GLYCERO-D-MANNO-HEPTOSE-6-EPIMERASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	138.07Å 162.46Å 185.03Å 90.00° 101.45° 90.00°	Depositor
Resolution (Å)	35.05 – 2.80 35.05 – 2.80	Depositor EDS
% Data completeness (in resolution range)	96.8 (35.05-2.80) 97.0 (35.05-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.54 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.6.0060	Depositor
R, R_{free}	0.252 , 0.272 0.253 , 0.273	Depositor DCC
R_{free} test set	9658 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	41.5	Xtriage
Anisotropy	0.149	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 35.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.55$, $\langle L^2 \rangle = 0.40$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	51434	wwPDB-VP
Average B, all atoms (Å ²)	12.0	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 48.14 % 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 8.9719e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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: BMA, NAP, 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	0.87	0/2498	0.87	0/3377
1	B	0.86	0/2498	0.87	0/3377
1	C	0.85	0/2498	0.87	1/3377 (0.0%)
1	D	0.86	0/2498	0.87	0/3377
1	E	0.86	0/2498	0.88	1/3377 (0.0%)
1	F	0.86	1/2498 (0.0%)	0.87	0/3377
1	G	0.87	0/2498	0.88	1/3377 (0.0%)
1	H	0.86	1/2498 (0.0%)	0.87	0/3377
1	I	0.86	0/2498	0.87	0/3377
1	J	0.85	0/2498	0.87	0/3377
1	K	0.86	0/2498	0.87	2/3377 (0.1%)
1	L	0.86	0/2498	0.87	1/3377 (0.0%)
1	M	0.86	0/2498	0.87	1/3377 (0.0%)
1	N	0.85	0/2498	0.87	0/3377
1	O	0.85	0/2498	0.87	0/3377
1	P	0.86	0/2498	0.87	0/3377
1	Q	0.86	0/2498	0.87	1/3377 (0.0%)
1	R	0.85	0/2498	0.86	0/3377
1	S	0.84	0/2498	0.86	0/3377
1	T	0.85	0/2498	0.87	0/3377
All	All	0.86	2/49960 (0.0%)	0.87	8/67540 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	258	GLN	CD-NE2	-5.37	1.22	1.33
1	F	132	GLU	CG-CD	5.31	1.65	1.52

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	M	273	GLN	CG-CD-NE2	5.76	125.04	116.40
1	C	273	GLN	CG-CD-NE2	5.73	124.99	116.40
1	E	273	GLN	CG-CD-NE2	5.65	124.88	116.40
1	K	273	GLN	CG-CD-NE2	5.57	124.75	116.40
1	L	273	GLN	CG-CD-NE2	5.53	124.70	116.40

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	2440	0	2337	17	1
1	B	2440	0	2337	15	0
1	C	2440	0	2337	16	0
1	D	2440	0	2337	16	1
1	E	2440	0	2337	15	4
1	F	2440	0	2337	16	2
1	G	2440	0	2337	13	0
1	H	2440	0	2337	16	0
1	I	2440	0	2337	18	1
1	J	2440	0	2337	13	0
1	K	2440	0	2337	19	0
1	L	2440	0	2337	14	0
1	M	2440	0	2337	21	4
1	N	2440	0	2337	21	0
1	O	2440	0	2337	18	0
1	P	2440	0	2337	20	0
1	Q	2440	0	2337	17	0
1	R	2440	0	2337	20	0
1	S	2440	0	2337	16	0
1	T	2440	0	2337	16	1
2	A	48	0	25	2	0
2	B	48	0	25	0	0
2	C	48	0	25	3	0
2	D	48	0	25	3	0
2	E	48	0	25	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	48	0	23	3	0
2	G	48	0	23	1	0
2	H	48	0	25	3	0
2	I	48	0	25	3	0
2	J	48	0	25	3	0
2	K	48	0	25	3	0
2	L	48	0	25	3	0
2	M	48	0	25	6	0
2	N	48	0	25	1	0
2	O	48	0	25	1	0
2	P	48	0	25	3	0
2	Q	48	0	25	4	0
2	R	48	0	25	3	0
2	S	48	0	25	1	0
2	T	48	0	25	2	0
3	A	27	0	12	1	0
3	B	27	0	12	1	0
3	C	27	0	12	0	0
3	D	27	0	12	0	0
3	E	27	0	12	0	0
3	F	27	0	12	0	0
3	G	27	0	12	1	0
3	H	27	0	12	0	0
3	I	27	0	12	0	0
3	J	27	0	12	0	0
3	K	27	0	12	0	0
3	L	27	0	12	0	0
3	M	27	0	12	0	0
3	N	27	0	12	0	0
3	O	27	0	12	0	0
3	P	27	0	12	0	0
3	Q	27	0	12	0	0
3	R	27	0	12	0	0
3	S	27	0	12	0	0
3	T	27	0	12	0	0
4	A	11	0	10	1	0
4	C	11	0	10	2	0
4	D	11	0	10	2	0
4	E	11	0	10	2	0
4	F	11	0	10	2	0
4	G	11	0	10	1	0
4	H	11	0	10	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	I	11	0	10	3	0
4	J	11	0	10	2	0
4	K	11	0	10	2	0
4	L	11	0	10	2	0
4	M	11	0	10	2	0
4	N	11	0	10	1	0
4	O	11	0	10	1	0
4	P	11	0	10	2	0
4	Q	11	0	10	1	0
4	R	11	0	10	2	0
4	S	11	0	10	1	0
4	T	11	0	10	2	0
5	A	59	0	0	1	0
5	B	36	0	0	1	0
5	C	34	0	0	0	0
5	D	60	0	0	3	0
5	E	61	0	0	1	0
5	F	56	0	0	1	0
5	G	28	0	0	1	0
5	H	49	0	0	1	0
5	I	78	0	0	2	0
5	J	66	0	0	0	0
5	K	59	0	0	3	0
5	L	51	0	0	2	0
5	M	45	0	0	1	0
5	N	39	0	0	2	0
5	O	33	0	0	0	0
5	P	49	0	0	3	0
5	Q	50	0	0	1	0
5	R	8	0	0	0	0
5	S	20	0	0	0	0
5	T	44	0	0	0	0
All	All	51434	0	47666	345	7

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.

The worst 5 of 345 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:N:258:GLN:HB3	1:R:157:GLU:OE1	1.58	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:291:PRO:HB3	1:O:156:PRO:O	1.58	1.02
1:N:258:GLN:CB	1:R:157:GLU:OE1	2.10	0.99
2:H:400:NAP:C5N	4:H:402:BMA:H61	2.05	0.86
2:I:400:NAP:C5N	4:I:402:BMA:H61	2.07	0.85

The worst 5 of 7 symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:195:GLY:O	1:M:159:ASN:OD1[2_646]	1.33	0.87
1:A:21:ASP:O	1:I:277:GLN:OE1[2_546]	1.58	0.62
1:D:277:GLN:OE1	1:F:21:ASP:O[2_556]	1.79	0.41
1:E:195:GLY:O	1:M:159:ASN:CG[2_646]	1.88	0.32
1:E:195:GLY:C	1:M:159:ASN:OD1[2_646]	1.99	0.21

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	305/357 (85%)	298 (98%)	6 (2%)	1 (0%)	36	66
1	B	305/357 (85%)	297 (97%)	7 (2%)	1 (0%)	36	66
1	C	305/357 (85%)	296 (97%)	8 (3%)	1 (0%)	36	66
1	D	305/357 (85%)	297 (97%)	7 (2%)	1 (0%)	36	66
1	E	305/357 (85%)	297 (97%)	6 (2%)	2 (1%)	18	47
1	F	305/357 (85%)	298 (98%)	6 (2%)	1 (0%)	36	66
1	G	305/357 (85%)	297 (97%)	7 (2%)	1 (0%)	36	66
1	H	305/357 (85%)	296 (97%)	8 (3%)	1 (0%)	36	66
1	I	305/357 (85%)	297 (97%)	6 (2%)	2 (1%)	18	47
1	J	305/357 (85%)	297 (97%)	6 (2%)	2 (1%)	18	47

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	K	305/357 (85%)	298 (98%)	5 (2%)	2 (1%)	18	47
1	L	305/357 (85%)	296 (97%)	7 (2%)	2 (1%)	18	47
1	M	305/357 (85%)	297 (97%)	7 (2%)	1 (0%)	36	66
1	N	305/357 (85%)	297 (97%)	6 (2%)	2 (1%)	18	47
1	O	305/357 (85%)	297 (97%)	6 (2%)	2 (1%)	18	47
1	P	305/357 (85%)	296 (97%)	7 (2%)	2 (1%)	18	47
1	Q	305/357 (85%)	297 (97%)	6 (2%)	2 (1%)	18	47
1	R	305/357 (85%)	297 (97%)	6 (2%)	2 (1%)	18	47
1	S	305/357 (85%)	296 (97%)	7 (2%)	2 (1%)	18	47
1	T	305/357 (85%)	296 (97%)	7 (2%)	2 (1%)	18	47
All	All	6100/7140 (85%)	5937 (97%)	131 (2%)	32 (0%)	24	55

5 of 32 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	174	ARG
1	B	174	ARG
1	C	174	ARG
1	D	174	ARG
1	E	174	ARG

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	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	B	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	C	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	D	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	E	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	F	255/292 (87%)	252 (99%)	3 (1%)	63	87

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	H	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	I	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	J	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	K	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	L	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	M	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	N	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	O	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	P	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	Q	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	R	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	S	255/292 (87%)	252 (99%)	3 (1%)	63	87
1	T	255/292 (87%)	252 (99%)	3 (1%)	63	87
All	All	5100/5840 (87%)	5040 (99%)	60 (1%)	63	87

5 of 60 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	J	105	LEU
1	S	105	LEU
1	L	300	THR
1	S	100	LEU
1	T	300	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 78 such sidechains are listed below:

Mol	Chain	Res	Type
1	O	41	ASN
1	S	41	ASN
1	O	234	ASN
1	Q	234	ASN
1	T	221	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

59 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAP	F	400	-	50,52,52	3.74	8 (16%)	71,80,80	2.95	20 (28%)
4	BMA	H	402	3	11,11,12	0.77	0	15,15,17	1.07	1 (6%)
2	NAP	D	400	-	50,52,52	1.77	5 (10%)	71,80,80	1.45	9 (12%)
2	NAP	J	400	-	50,52,52	1.76	6 (12%)	71,80,80	1.68	13 (18%)
4	BMA	I	402	3	11,11,12	0.33	0	15,15,17	0.91	1 (6%)
3	ADP	G	401	4	28,29,29	1.12	2 (7%)	43,45,45	1.96	11 (25%)
2	NAP	M	400	-	50,52,52	1.72	6 (12%)	71,80,80	1.63	11 (15%)
4	BMA	N	402	3	11,11,12	0.52	0	15,15,17	0.76	0
3	ADP	P	401	4	28,29,29	1.11	2 (7%)	43,45,45	1.85	9 (20%)
2	NAP	G	400	-	50,52,52	3.26	12 (24%)	71,80,80	2.24	20 (28%)
2	NAP	P	400	-	50,52,52	2.34	11 (22%)	71,80,80	1.63	14 (19%)
3	ADP	H	401	4	28,29,29	1.15	2 (7%)	43,45,45	2.03	13 (30%)
4	BMA	E	402	3	11,11,12	0.62	0	15,15,17	1.08	1 (6%)
3	ADP	I	401	4	28,29,29	1.12	2 (7%)	43,45,45	2.01	11 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ADP	T	401	4	28,29,29	1.21	2 (7%)	43,45,45	1.96	10 (23%)
2	NAP	B	400	-	50,52,52	1.75	6 (12%)	71,80,80	1.59	14 (19%)
4	BMA	D	402	3	11,11,12	0.57	0	15,15,17	1.07	2 (13%)
4	BMA	Q	402	3	11,11,12	0.58	0	15,15,17	0.81	0
4	BMA	S	402	3	11,11,12	0.47	0	15,15,17	0.88	0
2	NAP	N	400	-	50,52,52	1.75	7 (14%)	71,80,80	1.49	12 (16%)
3	ADP	O	401	4	28,29,29	1.24	2 (7%)	43,45,45	1.94	8 (18%)
4	BMA	A	402	3	11,11,12	0.68	0	15,15,17	0.77	0
4	BMA	M	402	3	11,11,12	0.48	0	15,15,17	0.99	1 (6%)
2	NAP	R	400	-	50,52,52	1.75	7 (14%)	71,80,80	1.69	13 (18%)
2	NAP	H	400	-	50,52,52	1.97	7 (14%)	71,80,80	1.49	11 (15%)
3	ADP	D	401	4	28,29,29	1.19	2 (7%)	43,45,45	2.07	11 (25%)
4	BMA	G	402	3	11,11,12	0.60	0	15,15,17	0.58	0
2	NAP	S	400	-	50,52,52	1.75	5 (10%)	71,80,80	1.51	10 (14%)
3	ADP	A	401	4	28,29,29	1.10	2 (7%)	43,45,45	1.95	11 (25%)
4	BMA	O	402	3	11,11,12	0.55	0	15,15,17	0.86	1 (6%)
2	NAP	Q	400	-	50,52,52	1.87	6 (12%)	71,80,80	1.67	11 (15%)
4	BMA	K	402	3	11,11,12	0.69	0	15,15,17	0.83	0
4	BMA	C	402	3	11,11,12	0.74	0	15,15,17	0.88	0
3	ADP	C	401	4	28,29,29	1.17	2 (7%)	43,45,45	1.84	11 (25%)
3	ADP	B	401	-	28,29,29	1.50	6 (21%)	43,45,45	1.81	10 (23%)
4	BMA	T	402	3	11,11,12	0.65	0	15,15,17	0.75	0
3	ADP	K	401	4	28,29,29	1.11	1 (3%)	43,45,45	2.02	12 (27%)
4	BMA	R	402	3	11,11,12	0.68	0	15,15,17	2.10	4 (26%)
2	NAP	C	400	-	50,52,52	2.31	5 (10%)	71,80,80	1.90	17 (23%)
4	BMA	J	402	3	11,11,12	0.58	0	15,15,17	0.92	1 (6%)
4	BMA	F	402	3	11,11,12	0.52	0	15,15,17	0.82	0
3	ADP	N	401	4	28,29,29	1.15	2 (7%)	43,45,45	2.01	11 (25%)
2	NAP	I	400	-	50,52,52	1.82	6 (12%)	71,80,80	1.55	10 (14%)
2	NAP	L	400	-	50,52,52	1.83	6 (12%)	71,80,80	1.48	9 (12%)
2	NAP	O	400	-	50,52,52	1.73	9 (18%)	71,80,80	1.76	17 (23%)
2	NAP	A	400	-	50,52,52	1.71	7 (14%)	71,80,80	1.95	18 (25%)
2	NAP	K	400	-	50,52,52	2.20	9 (18%)	71,80,80	1.88	19 (26%)
3	ADP	E	401	4	28,29,29	1.09	1 (3%)	43,45,45	1.88	11 (25%)
2	NAP	T	400	-	50,52,52	2.03	7 (14%)	71,80,80	1.66	14 (19%)
3	ADP	M	401	4	28,29,29	1.18	2 (7%)	43,45,45	1.94	10 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ADP	J	401	4	28,29,29	1.12	1 (3%)	43,45,45	2.17	13 (30%)
2	NAP	E	400	-	50,52,52	1.89	6 (12%)	71,80,80	1.62	13 (18%)
3	ADP	L	401	4	28,29,29	1.14	2 (7%)	43,45,45	1.92	8 (18%)
4	BMA	L	402	3	11,11,12	0.65	0	15,15,17	0.69	0
3	ADP	Q	401	4	28,29,29	1.10	2 (7%)	43,45,45	1.95	12 (27%)
3	ADP	S	401	4	28,29,29	1.10	1 (3%)	43,45,45	1.97	10 (23%)
3	ADP	F	401	4	28,29,29	1.16	3 (10%)	43,45,45	2.06	9 (20%)
4	BMA	P	402	3	11,11,12	0.70	0	15,15,17	1.21	2 (13%)
3	ADP	R	401	4	28,29,29	1.32	3 (10%)	43,45,45	1.87	9 (20%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAP	F	400	-	-	7/35/67/67	0/5/5/5
4	BMA	H	402	3	-	2/2/19/22	0/1/1/1
2	NAP	D	400	-	-	3/35/67/67	0/5/5/5
2	NAP	J	400	-	-	14/35/67/67	0/5/5/5
4	BMA	I	402	3	-	1/2/19/22	0/1/1/1
3	ADP	G	401	4	-	0/16/32/32	0/3/3/3
2	NAP	M	400	-	-	7/35/67/67	0/5/5/5
4	BMA	N	402	3	-	2/2/19/22	0/1/1/1
3	ADP	P	401	4	-	0/16/32/32	0/3/3/3
2	NAP	G	400	-	-	7/35/67/67	0/5/5/5
2	NAP	P	400	-	-	9/35/67/67	0/5/5/5
3	ADP	H	401	4	-	1/16/32/32	0/3/3/3
4	BMA	E	402	3	-	2/2/19/22	0/1/1/1
3	ADP	I	401	4	-	1/16/32/32	0/3/3/3
3	ADP	T	401	4	-	1/16/32/32	0/3/3/3
2	NAP	B	400	-	-	16/35/67/67	0/5/5/5
4	BMA	D	402	3	-	2/2/19/22	0/1/1/1
4	BMA	Q	402	3	-	2/2/19/22	0/1/1/1
4	BMA	S	402	3	-	2/2/19/22	0/1/1/1
2	NAP	N	400	-	-	8/35/67/67	0/5/5/5
3	ADP	O	401	4	-	2/16/32/32	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	BMA	A	402	3	-	2/2/19/22	0/1/1/1
4	BMA	M	402	3	-	2/2/19/22	0/1/1/1
2	NAP	R	400	-	-	10/35/67/67	0/5/5/5
2	NAP	H	400	-	-	14/35/67/67	0/5/5/5
3	ADP	D	401	4	-	1/16/32/32	0/3/3/3
4	BMA	G	402	3	-	0/2/19/22	0/1/1/1
2	NAP	S	400	-	-	6/35/67/67	0/5/5/5
3	ADP	A	401	4	-	1/16/32/32	0/3/3/3
4	BMA	O	402	3	-	2/2/19/22	0/1/1/1
2	NAP	Q	400	-	-	13/35/67/67	0/5/5/5
4	BMA	K	402	3	-	2/2/19/22	0/1/1/1
4	BMA	C	402	3	-	1/2/19/22	0/1/1/1
3	ADP	C	401	4	-	1/16/32/32	0/3/3/3
3	ADP	B	401	-	-	0/16/32/32	0/3/3/3
4	BMA	T	402	3	-	2/2/19/22	0/1/1/1
3	ADP	K	401	4	-	3/16/32/32	0/3/3/3
4	BMA	R	402	3	-	2/2/19/22	0/1/1/1
2	NAP	C	400	-	-	12/35/67/67	0/5/5/5
4	BMA	J	402	3	-	2/2/19/22	0/1/1/1
4	BMA	F	402	3	-	0/2/19/22	0/1/1/1
3	ADP	N	401	4	-	1/16/32/32	0/3/3/3
2	NAP	I	400	-	-	13/35/67/67	0/5/5/5
2	NAP	L	400	-	-	7/35/67/67	0/5/5/5
2	NAP	O	400	-	-	10/35/67/67	0/5/5/5
2	NAP	A	400	-	-	10/35/67/67	0/5/5/5
2	NAP	K	400	-	-	11/35/67/67	0/5/5/5
3	ADP	E	401	4	-	3/16/32/32	0/3/3/3
2	NAP	T	400	-	-	11/35/67/67	0/5/5/5
3	ADP	M	401	4	-	1/16/32/32	0/3/3/3
3	ADP	J	401	4	-	2/16/32/32	0/3/3/3
2	NAP	E	400	-	-	12/35/67/67	0/5/5/5
3	ADP	L	401	4	-	1/16/32/32	0/3/3/3
4	BMA	L	402	3	-	2/2/19/22	0/1/1/1
3	ADP	Q	401	4	-	0/16/32/32	0/3/3/3
3	ADP	S	401	4	-	3/16/32/32	0/3/3/3
3	ADP	F	401	4	-	0/16/32/32	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	BMA	P	402	3	-	2/2/19/22	0/1/1/1
3	ADP	R	401	4	-	2/16/32/32	0/3/3/3

The worst 5 of 183 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	400	NAP	O7N-C7N	15.99	1.53	1.24
2	G	400	NAP	C2N-N1N	12.18	1.48	1.35
2	F	400	NAP	C7N-N7N	-11.82	1.11	1.33
2	G	400	NAP	O7N-C7N	11.58	1.45	1.24
2	H	400	NAP	O7N-C7N	11.11	1.44	1.24

The worst 5 of 499 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	400	NAP	C5N-C4N-C3N	14.08	134.21	120.36
2	F	400	NAP	C6N-N1N-C2N	7.92	128.61	121.88
2	F	400	NAP	C3N-C7N-N7N	7.50	126.98	117.74
2	F	400	NAP	O7N-C7N-C3N	-7.37	110.58	119.60
2	G	400	NAP	C2N-C3N-C4N	7.20	126.63	118.26

There are no chirality outliers.

5 of 256 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	400	NAP	C2B-C1B-N9A-C8A
2	A	400	NAP	C5D-O5D-PN-O2N
2	A	400	NAP	O4D-C1D-N1N-C2N
2	B	400	NAP	C5B-O5B-PA-O2A
2	B	400	NAP	C5B-O5B-PA-O3

There are no ring outliers.

41 monomers are involved in 53 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	400	NAP	3	0
4	H	402	BMA	2	0
2	D	400	NAP	3	0
2	J	400	NAP	3	0
4	I	402	BMA	3	0

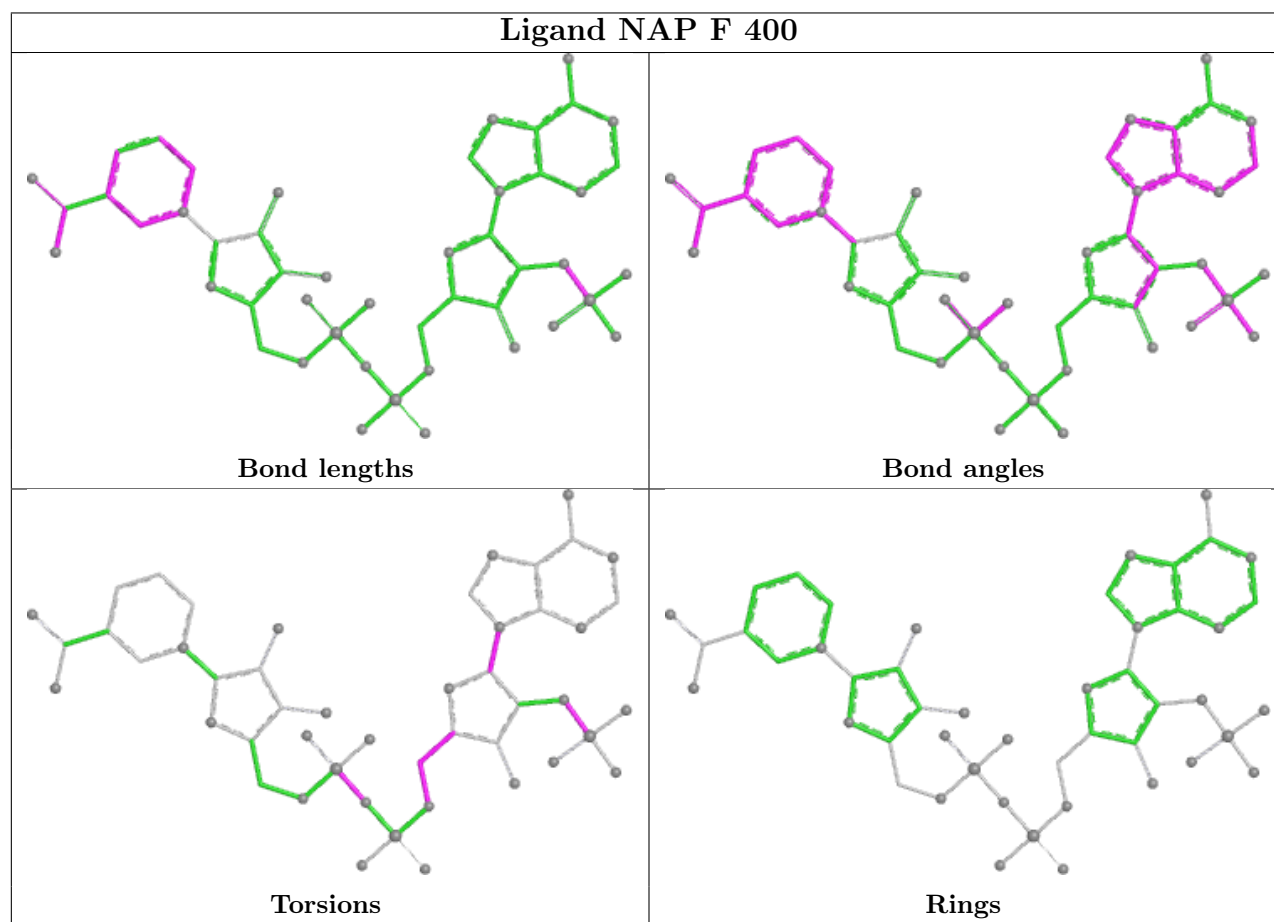
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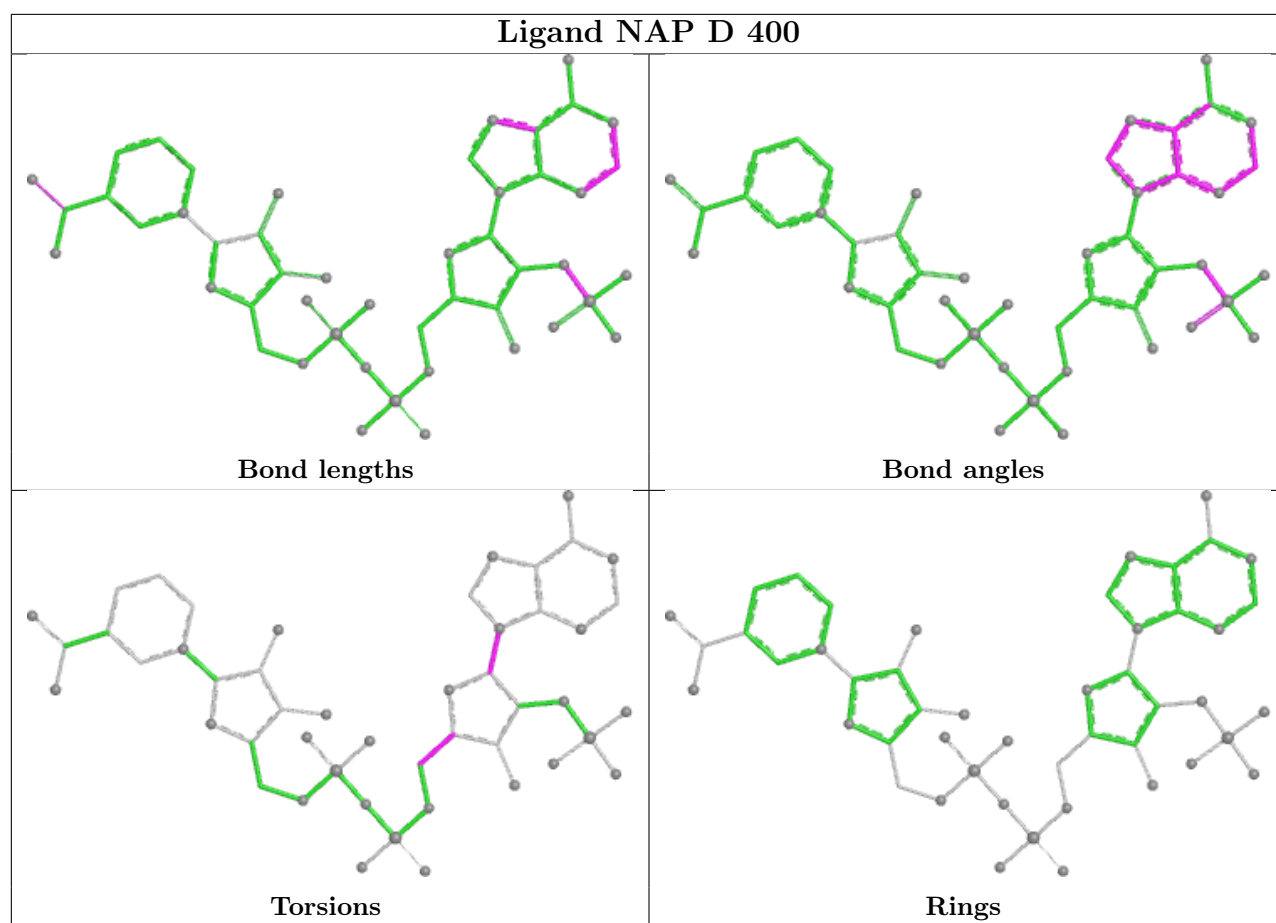
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	401	ADP	1	0
2	M	400	NAP	6	0
4	N	402	BMA	1	0
2	G	400	NAP	1	0
2	P	400	NAP	3	0
4	E	402	BMA	2	0
4	D	402	BMA	2	0
4	Q	402	BMA	1	0
4	S	402	BMA	1	0
2	N	400	NAP	1	0
4	A	402	BMA	1	0
4	M	402	BMA	2	0
2	R	400	NAP	3	0
2	H	400	NAP	3	0
4	G	402	BMA	1	0
2	S	400	NAP	1	0
3	A	401	ADP	1	0
4	O	402	BMA	1	0
2	Q	400	NAP	4	0
4	K	402	BMA	2	0
4	C	402	BMA	2	0
3	B	401	ADP	1	0
4	T	402	BMA	2	0
4	R	402	BMA	2	0
2	C	400	NAP	3	0
4	J	402	BMA	2	0
4	F	402	BMA	2	0
2	I	400	NAP	3	0
2	L	400	NAP	3	0
2	O	400	NAP	1	0
2	A	400	NAP	2	0
2	K	400	NAP	3	0
2	T	400	NAP	2	0
2	E	400	NAP	2	0
4	L	402	BMA	2	0
4	P	402	BMA	2	0

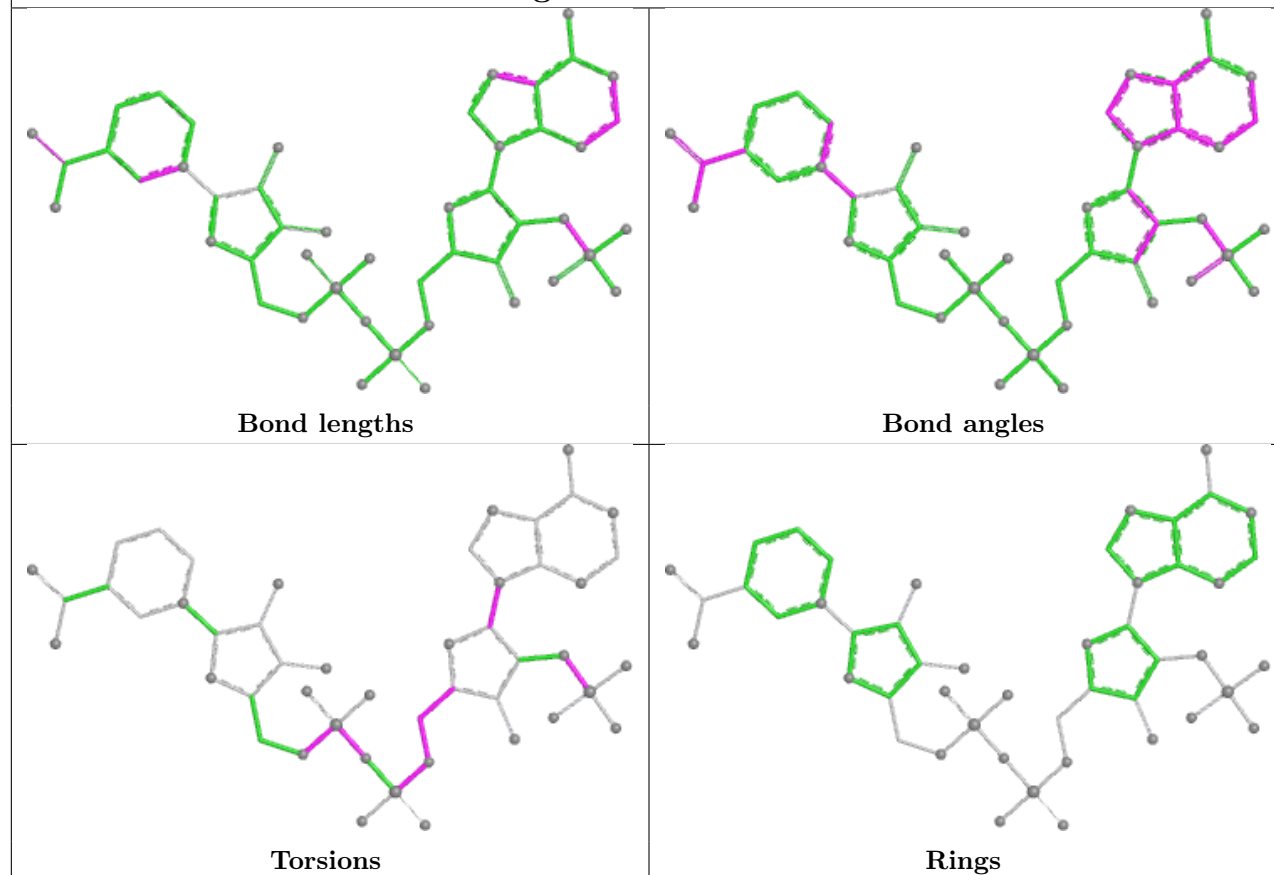
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

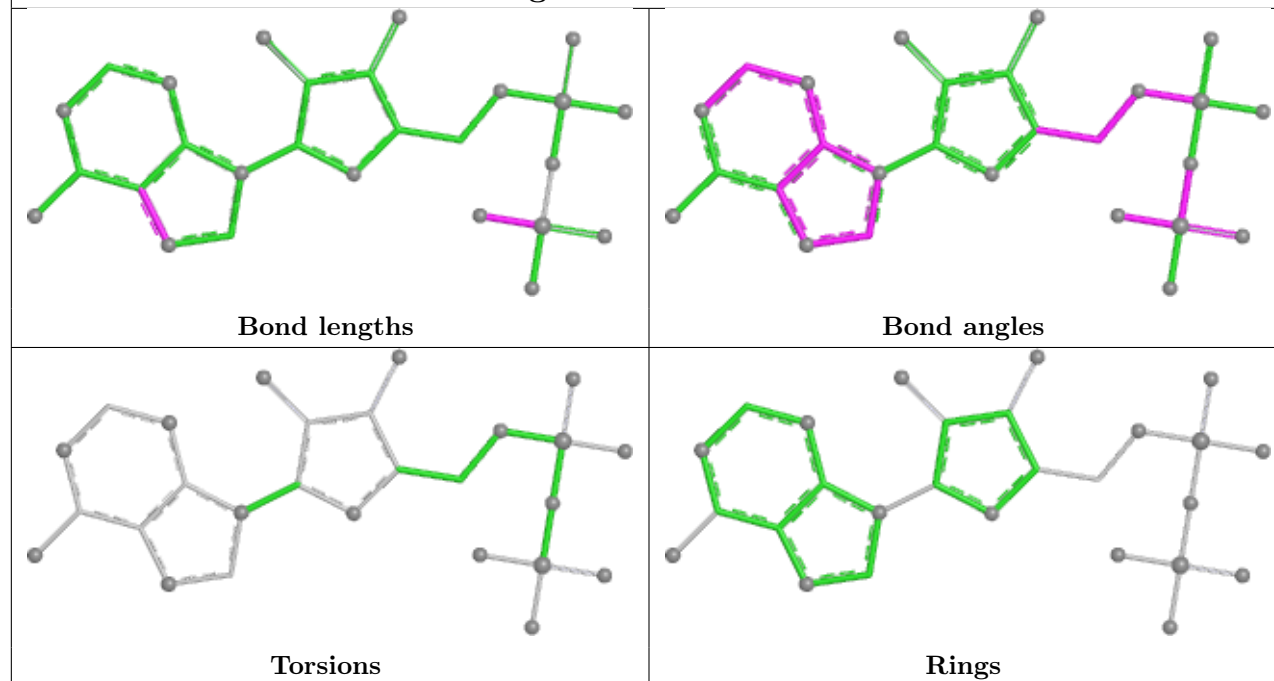


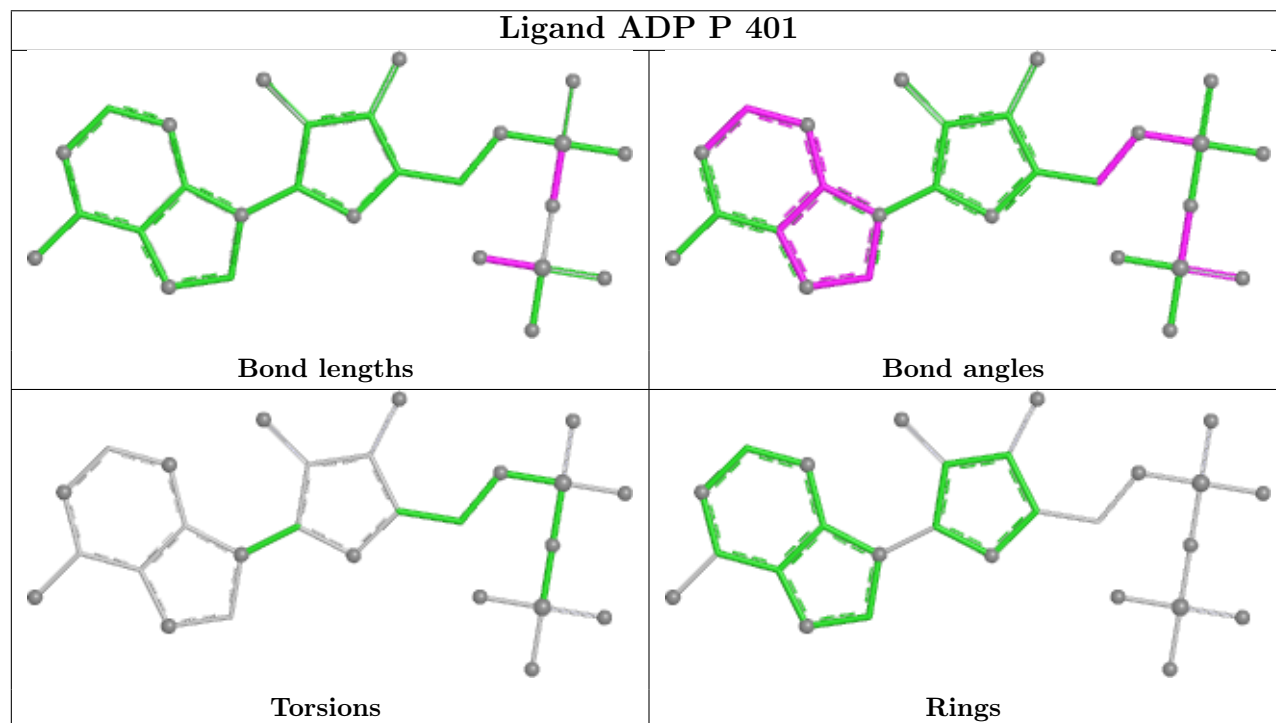
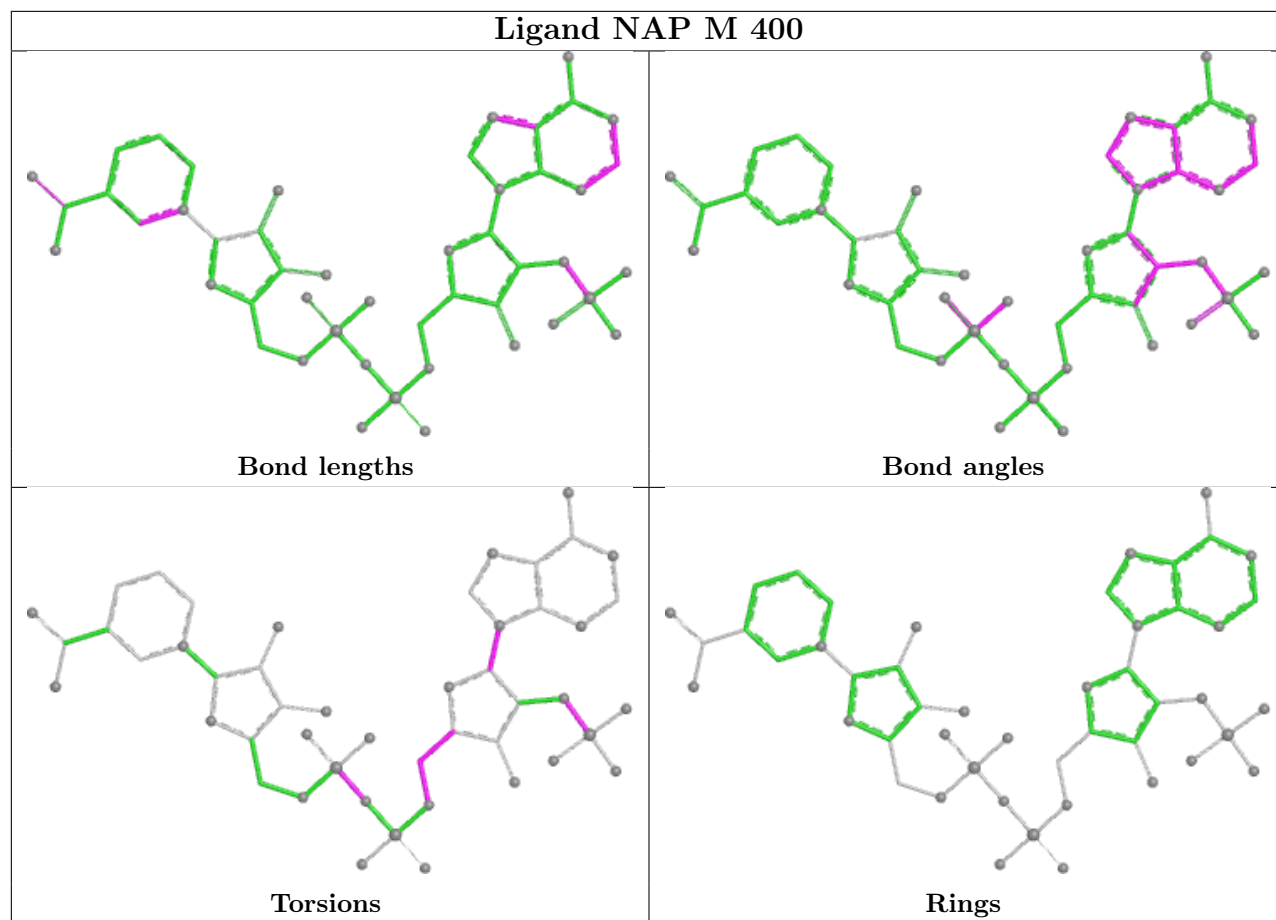


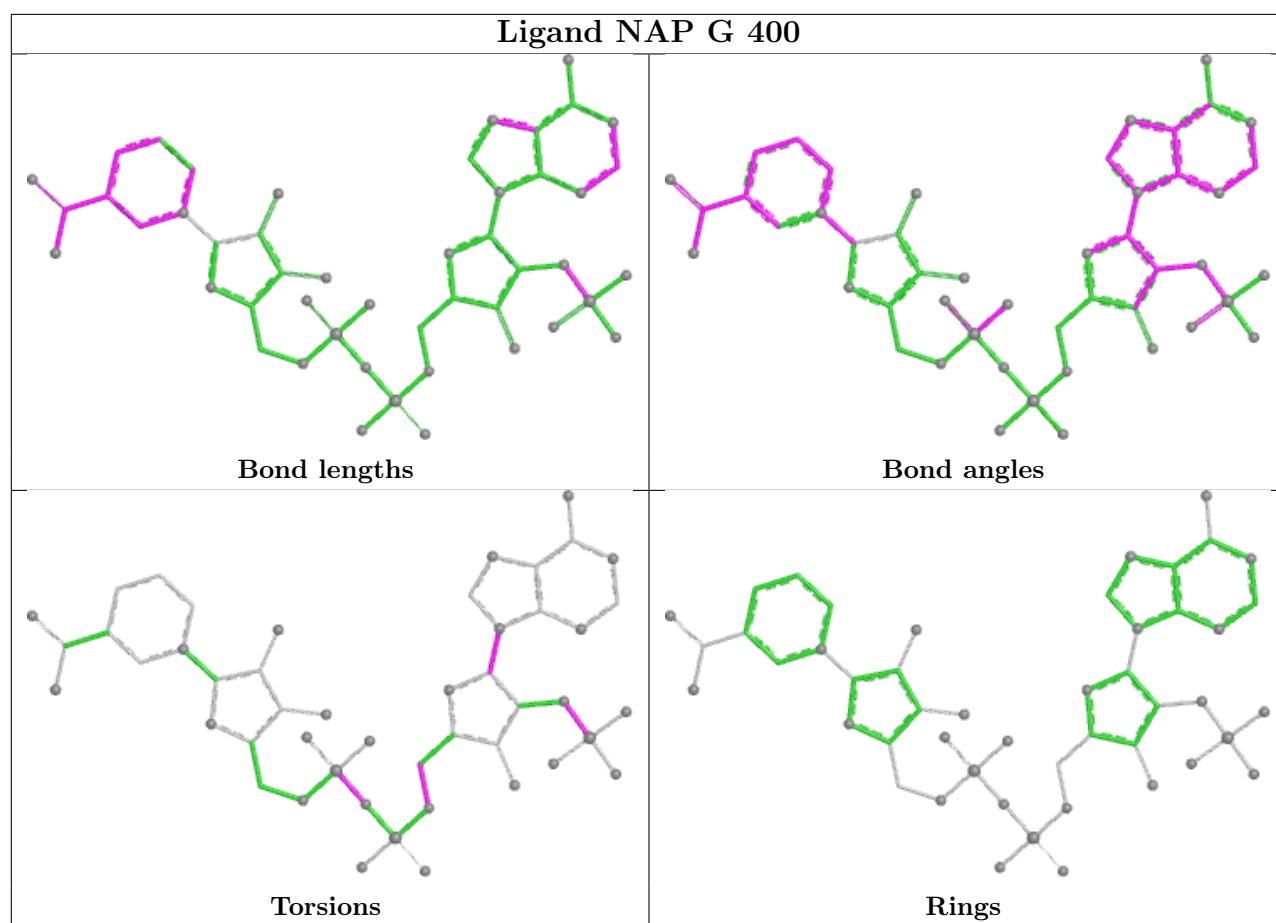
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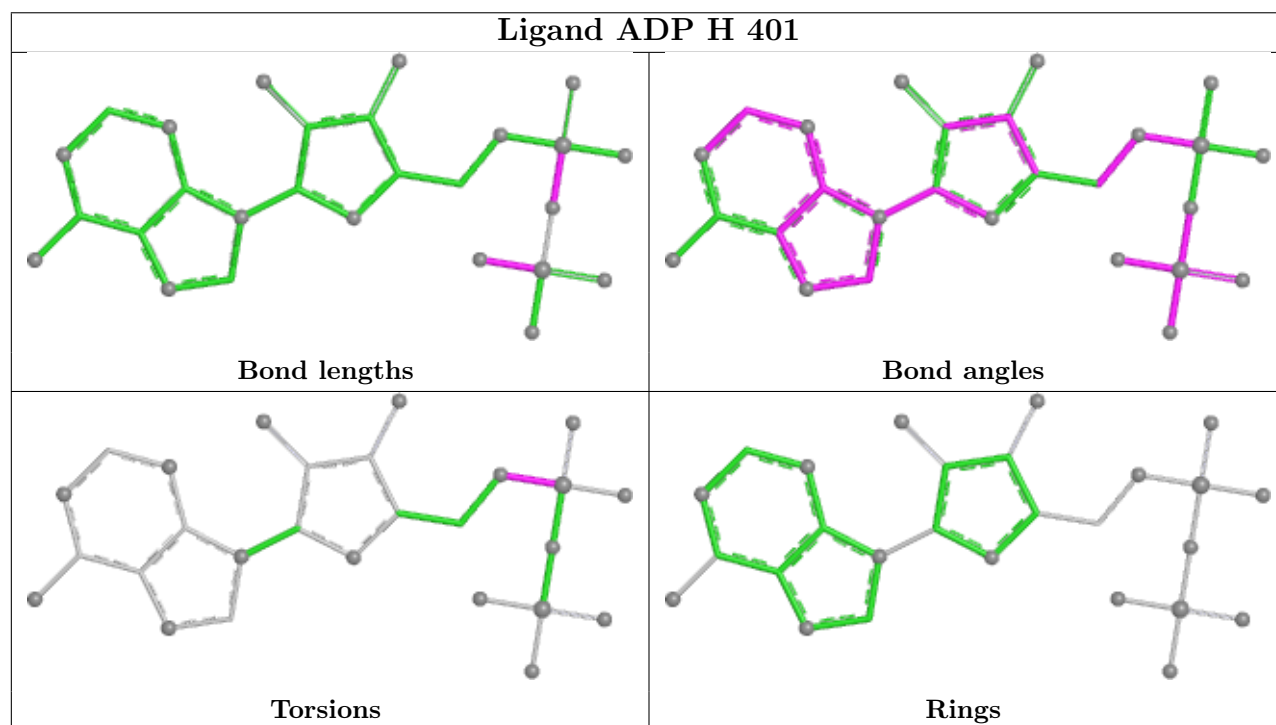
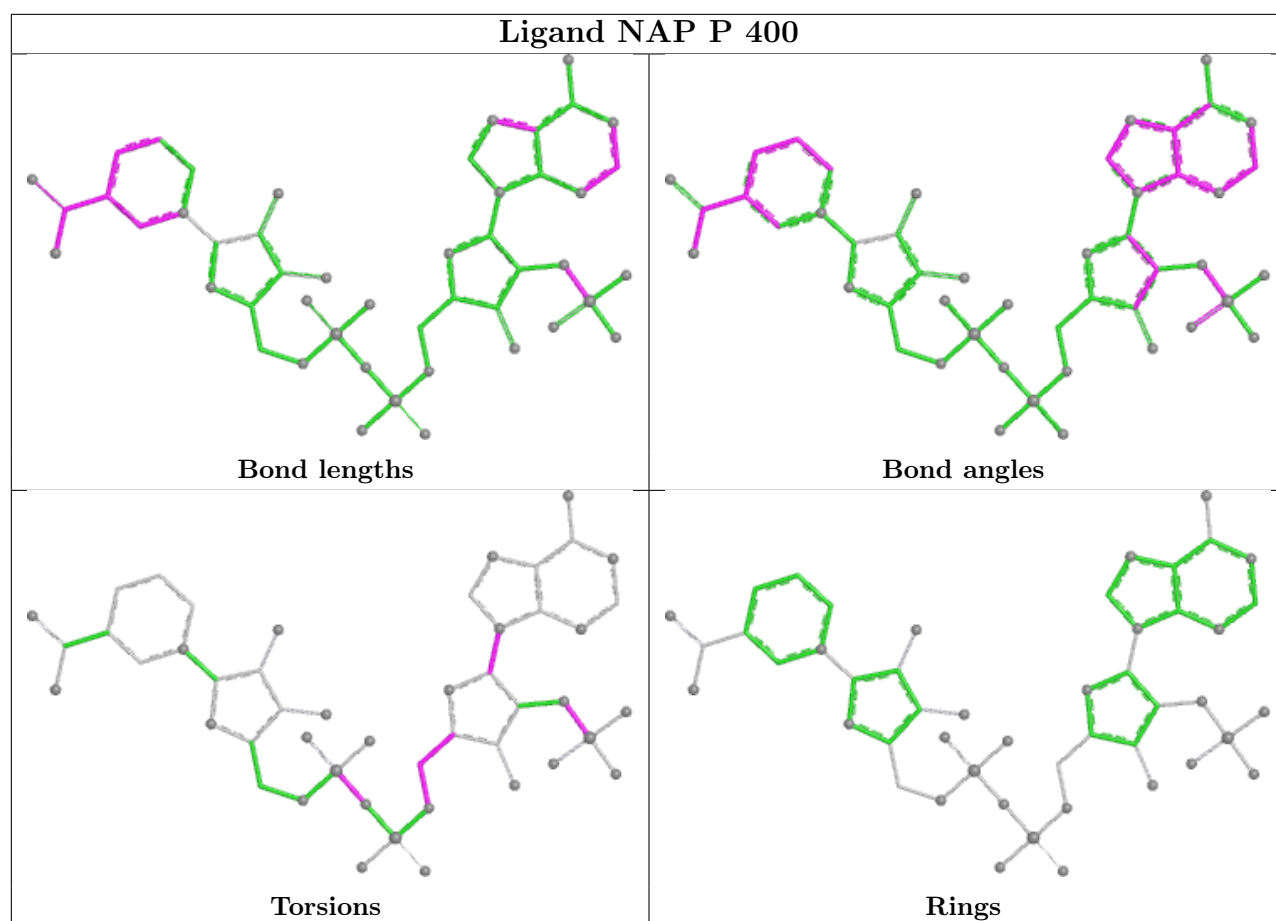


Ligand ADP G 401

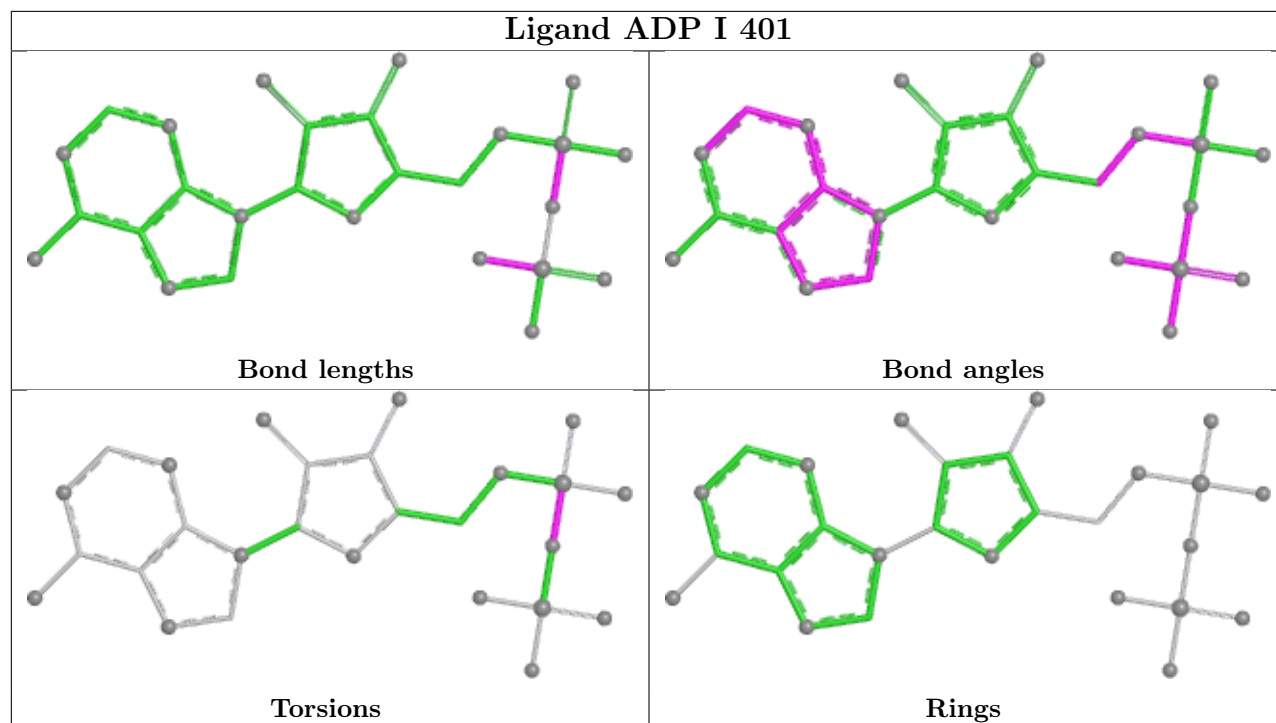




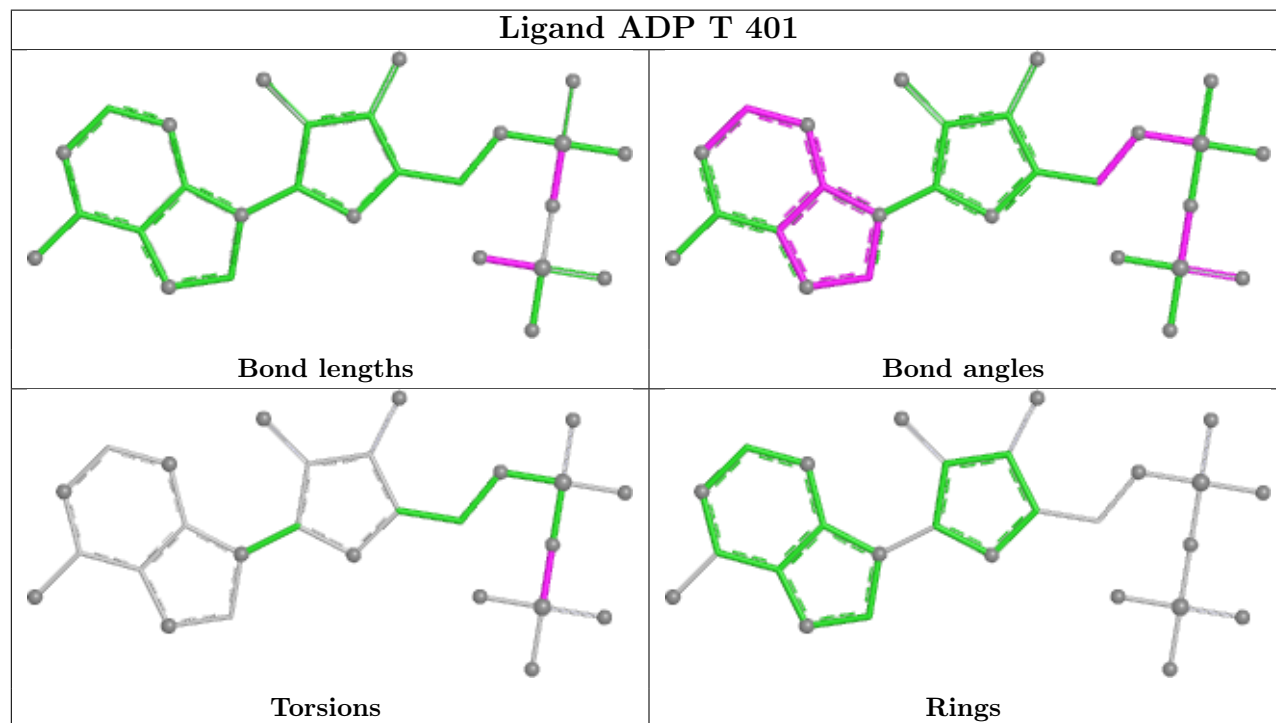


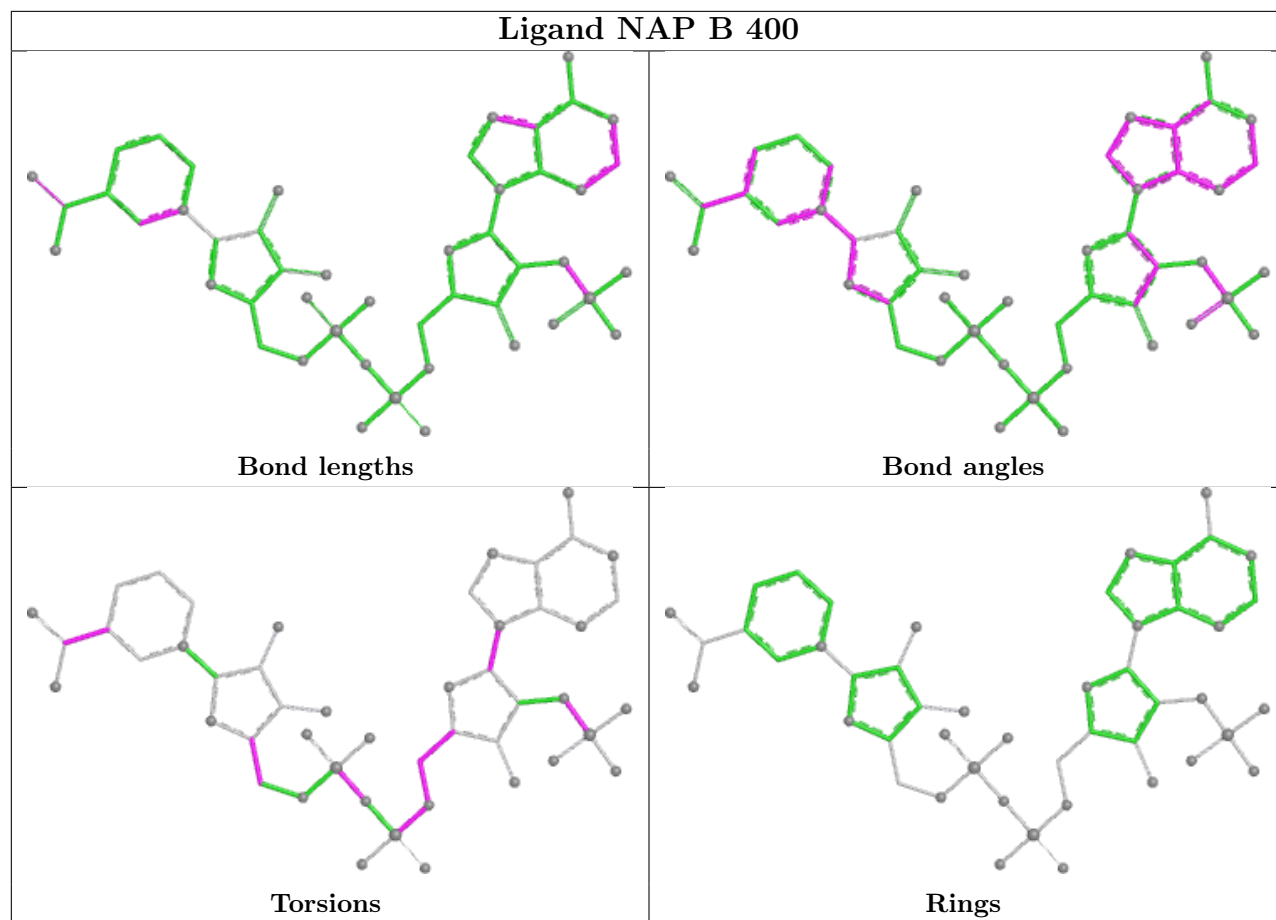


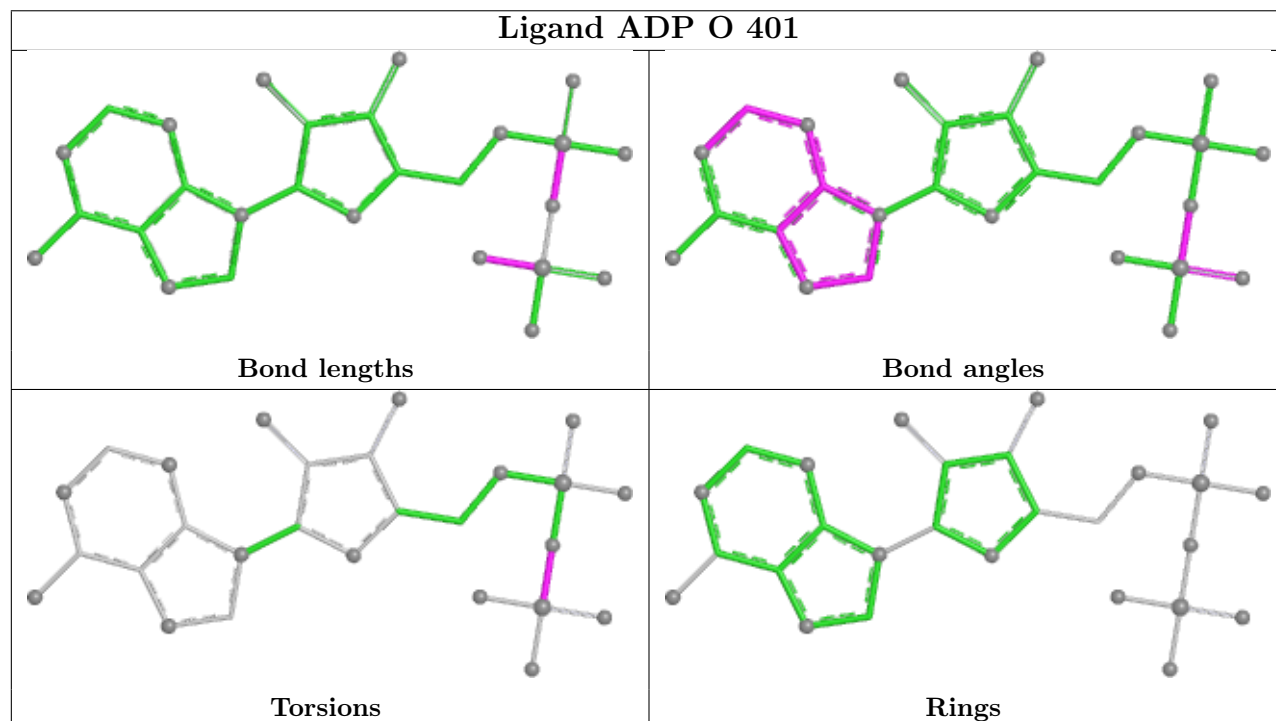
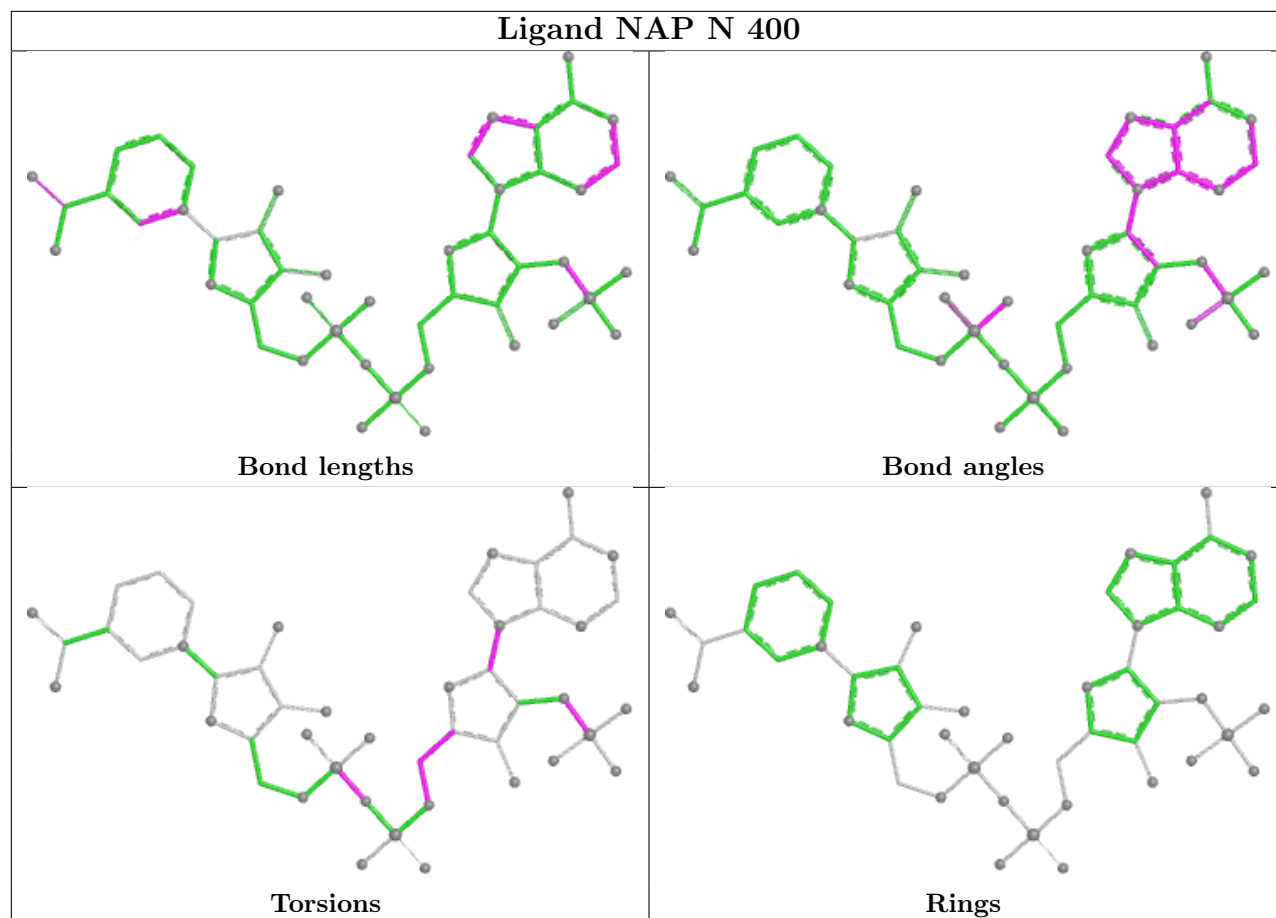
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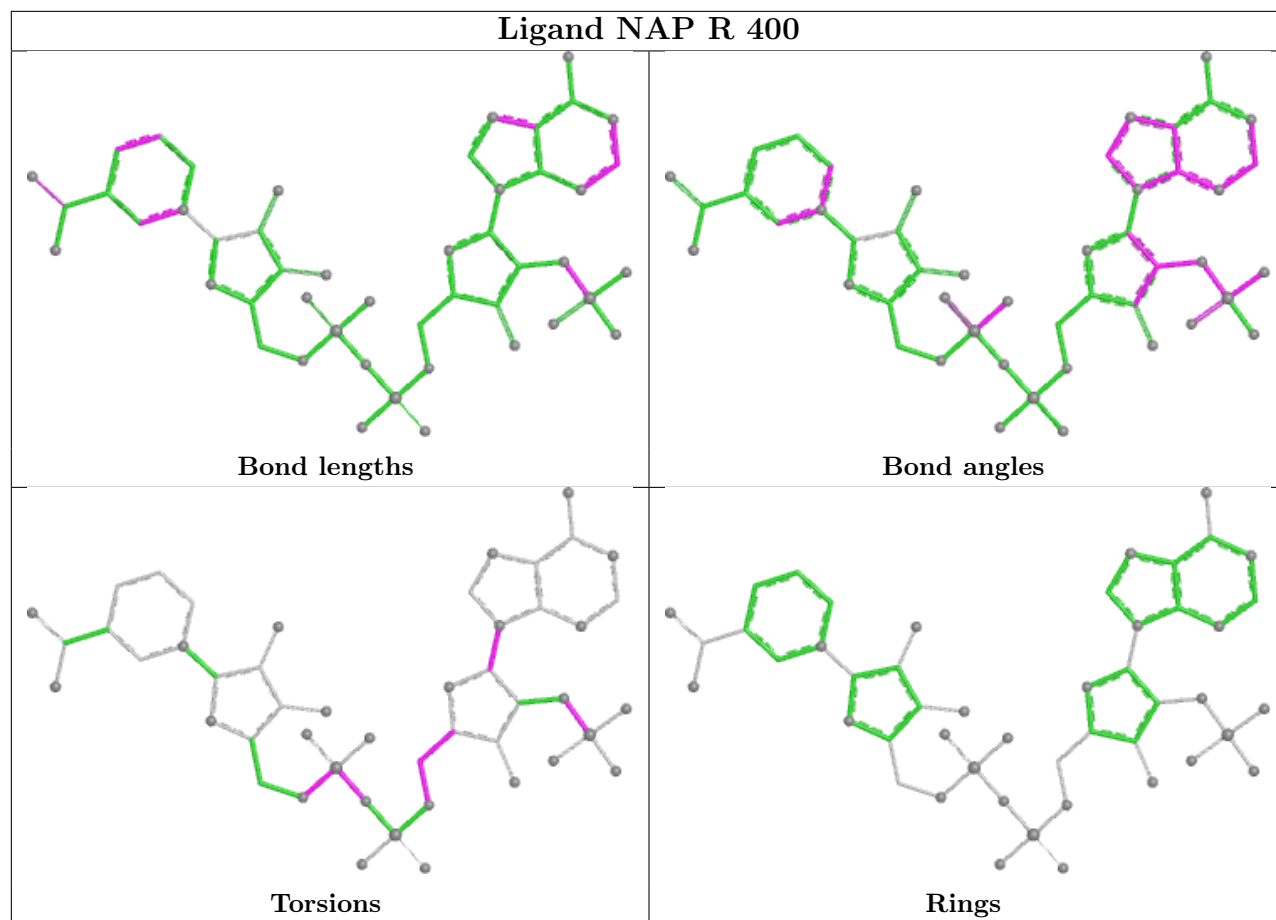


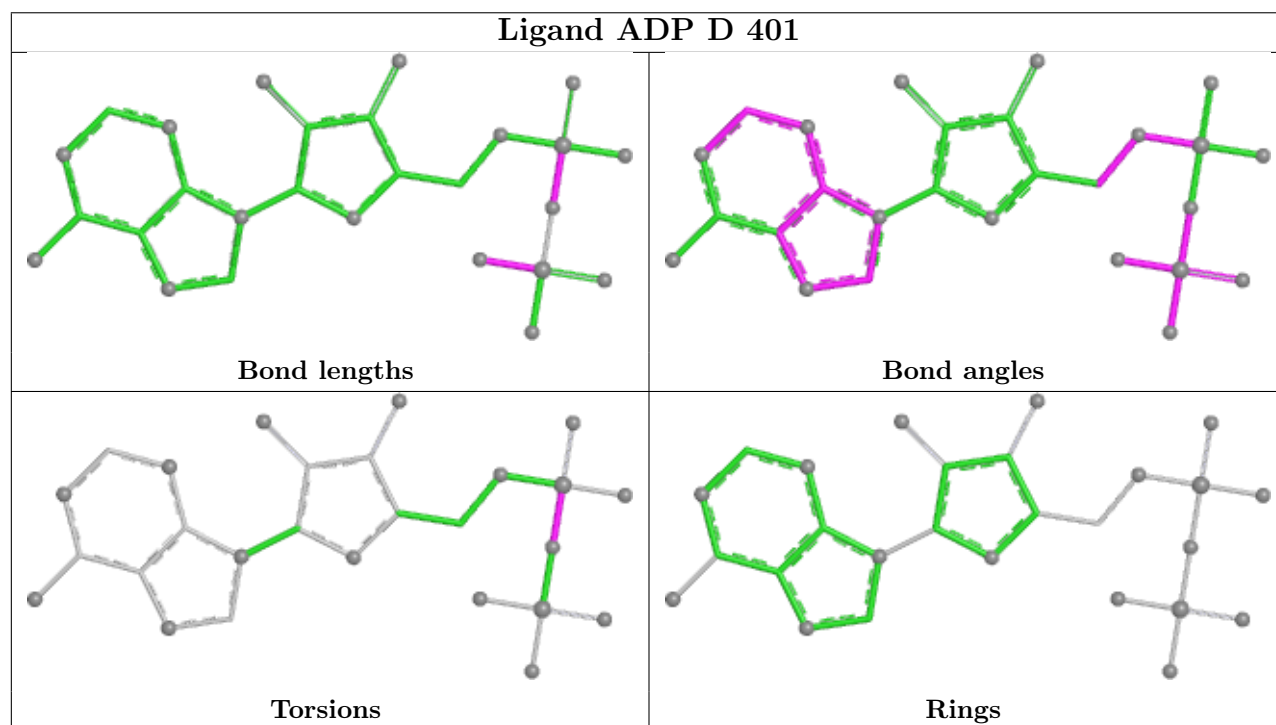
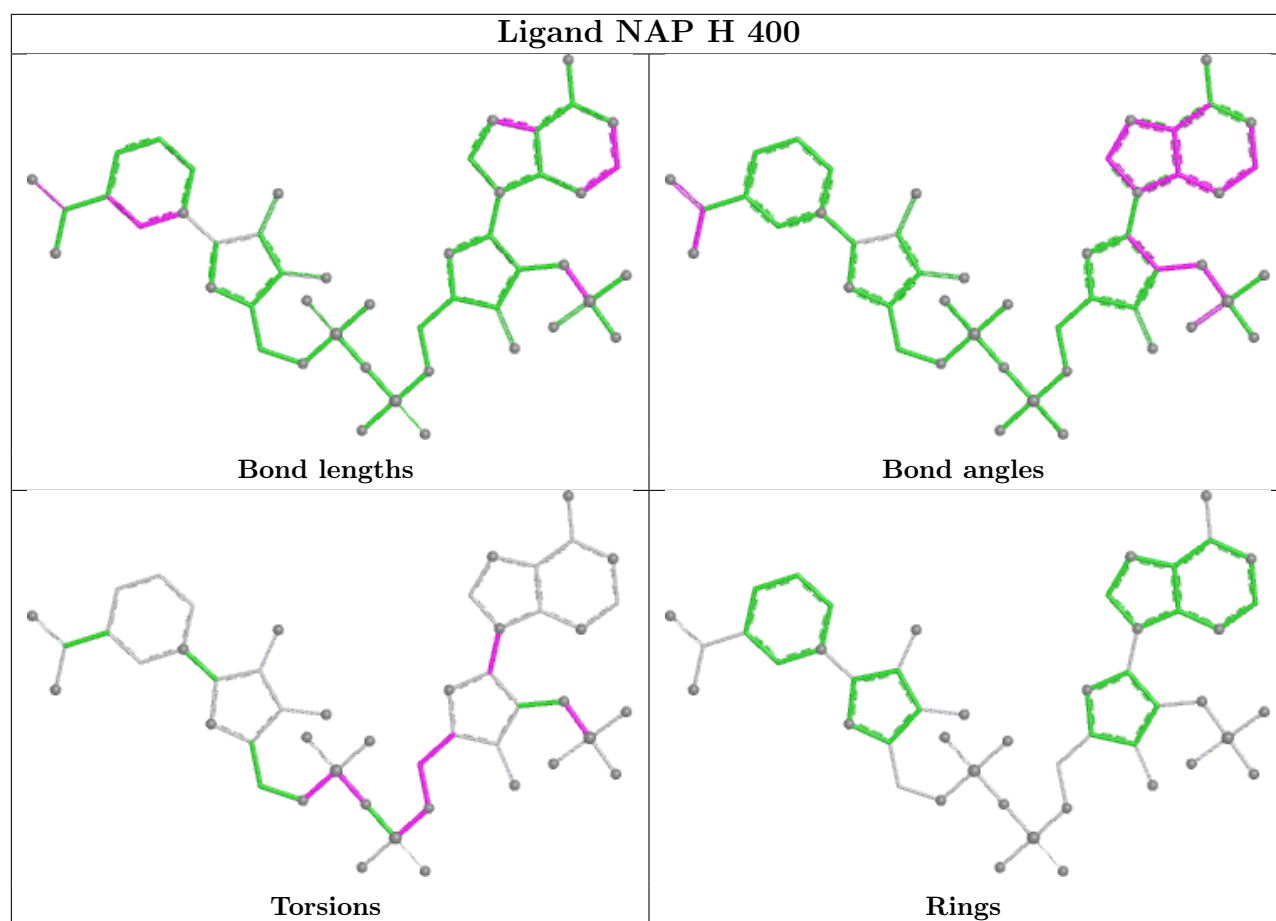
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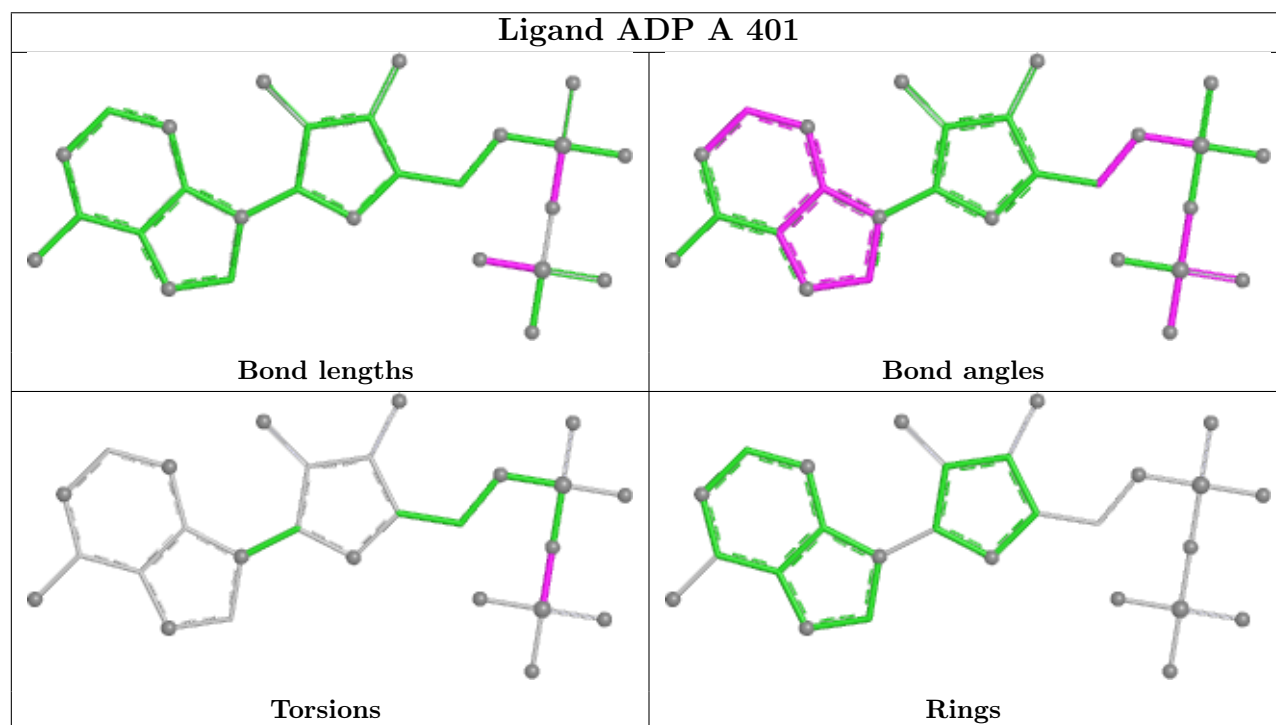
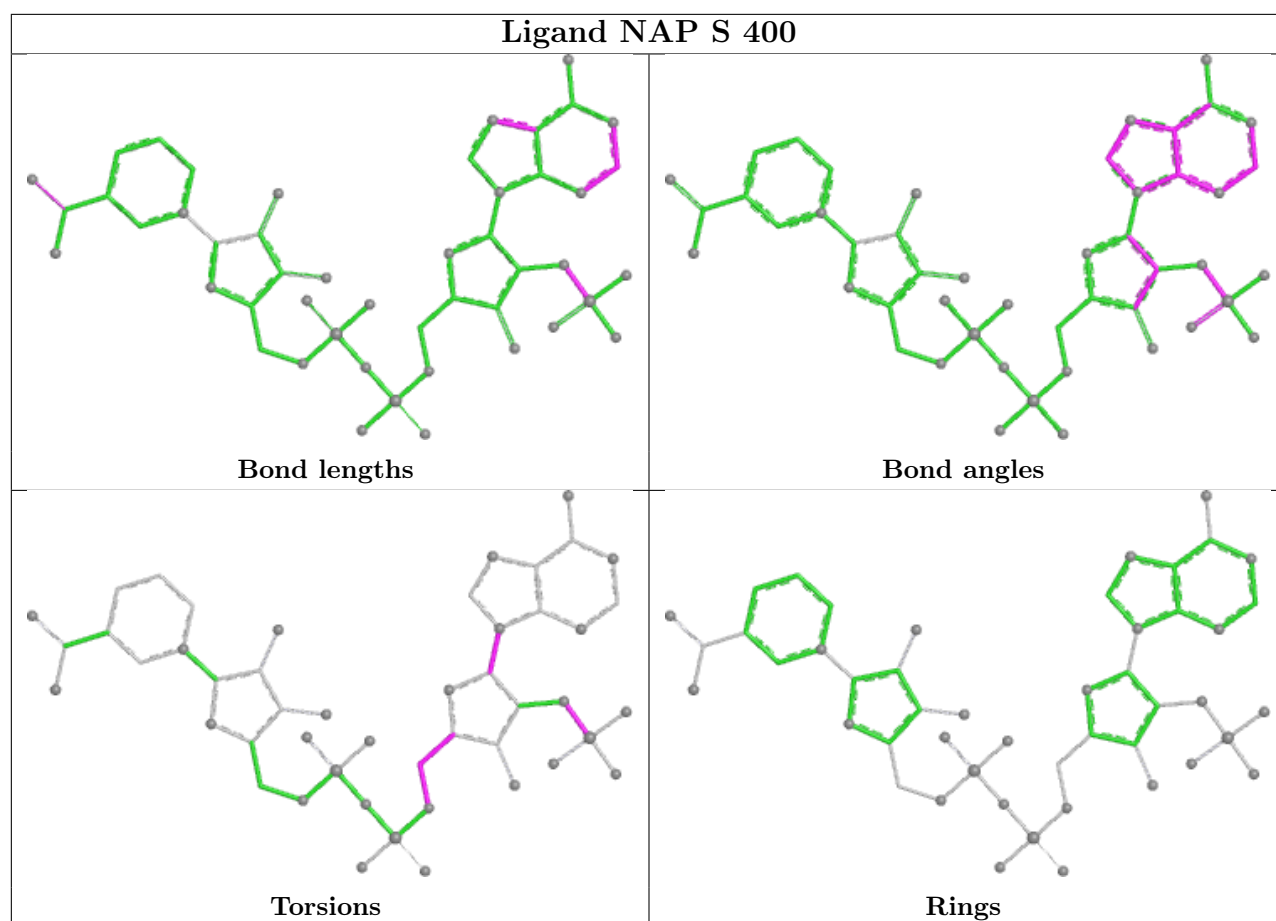


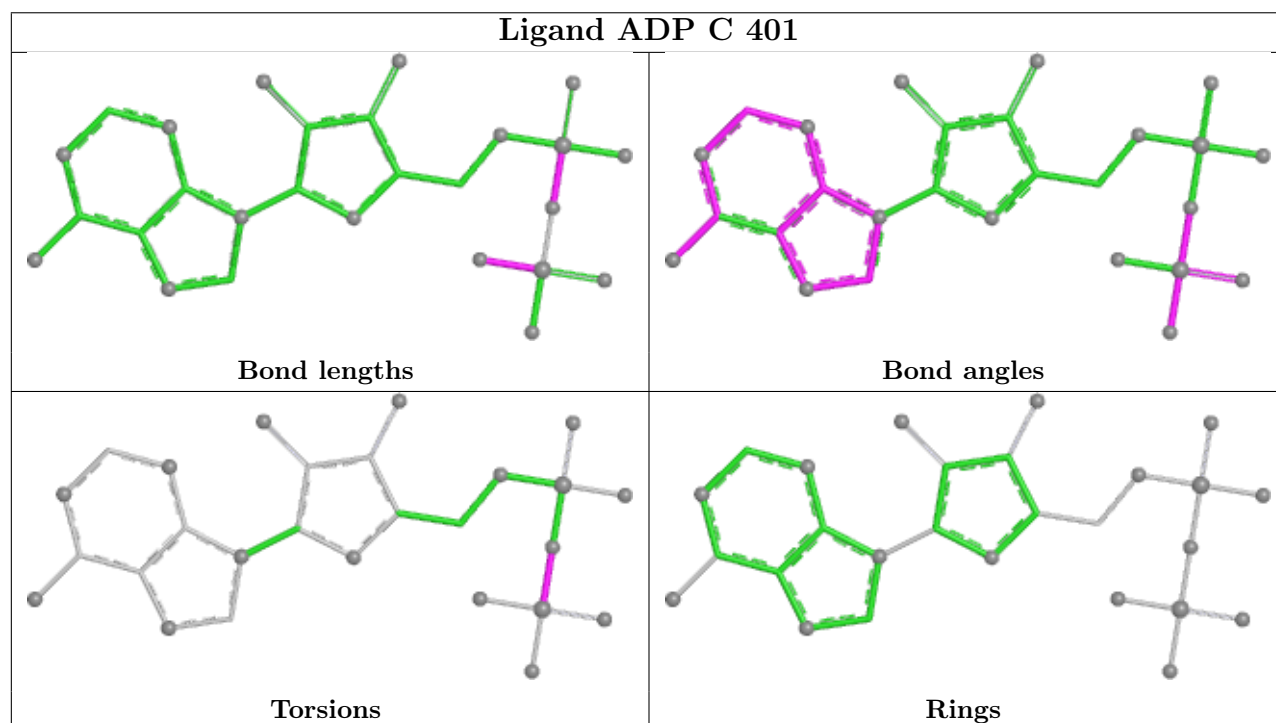
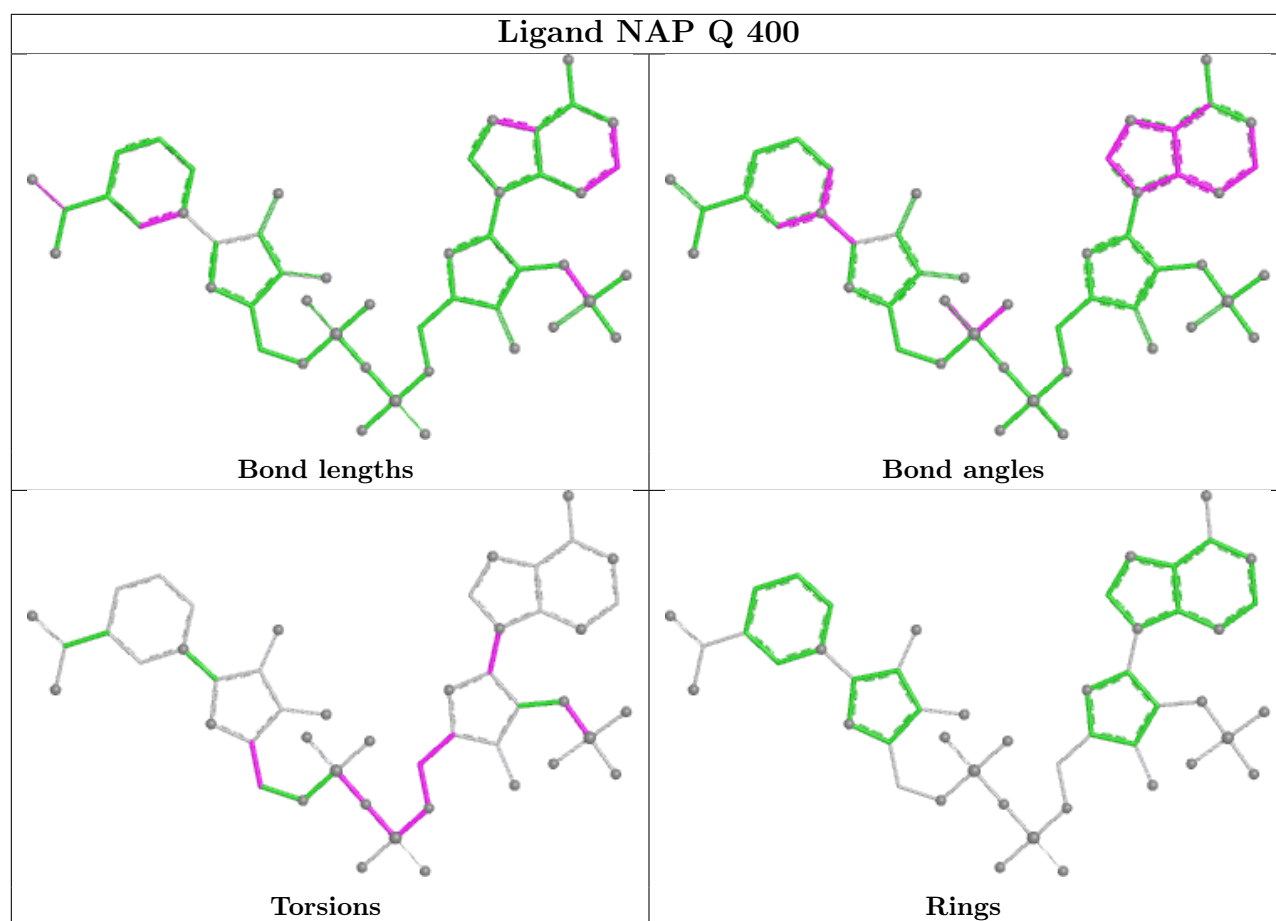


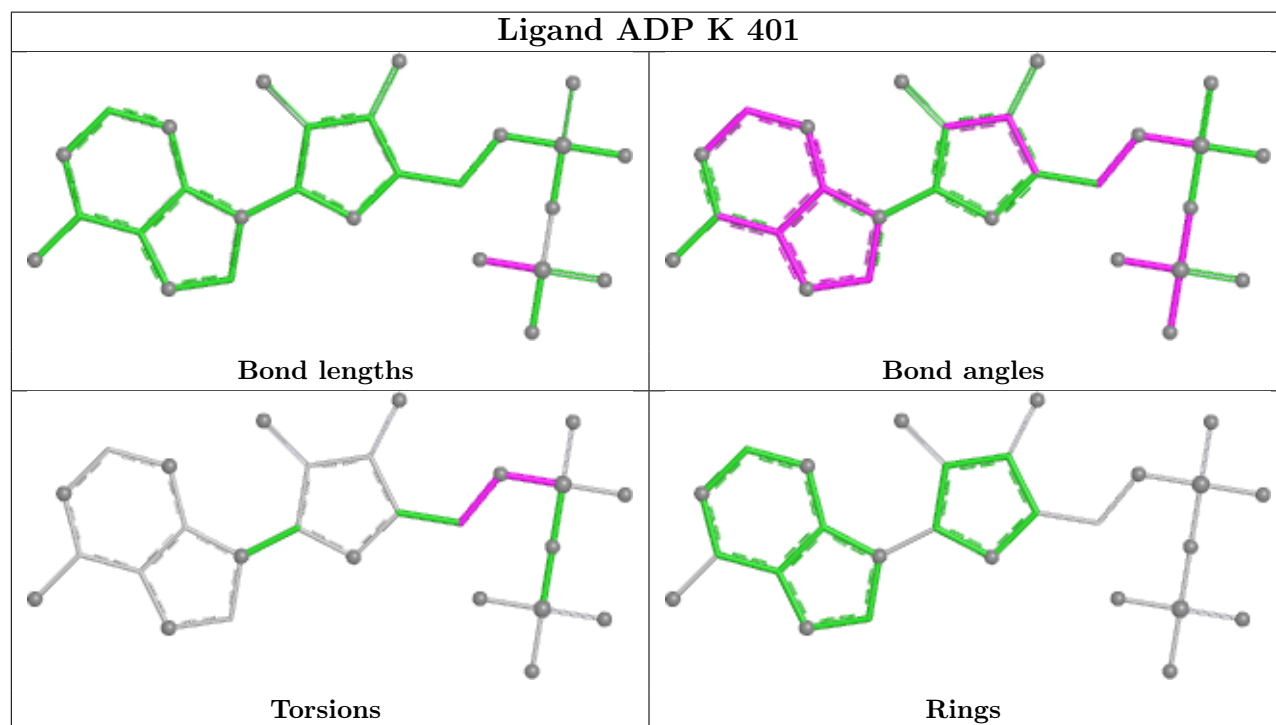
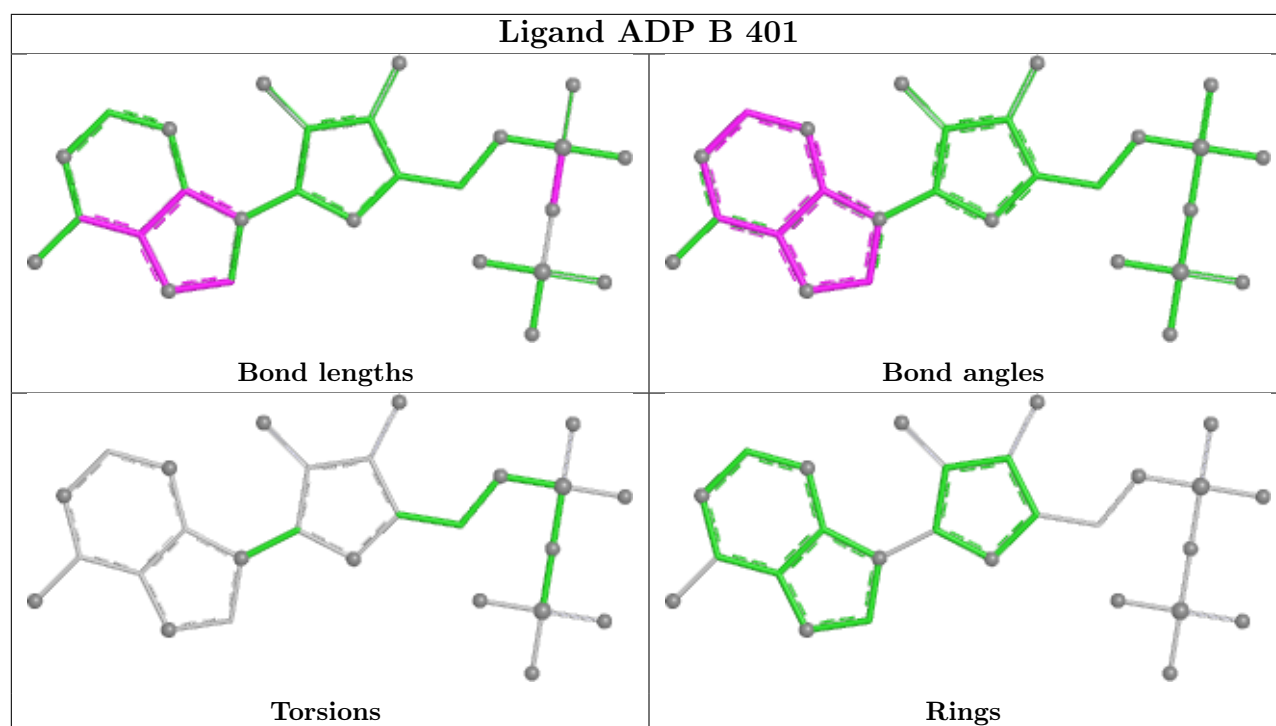


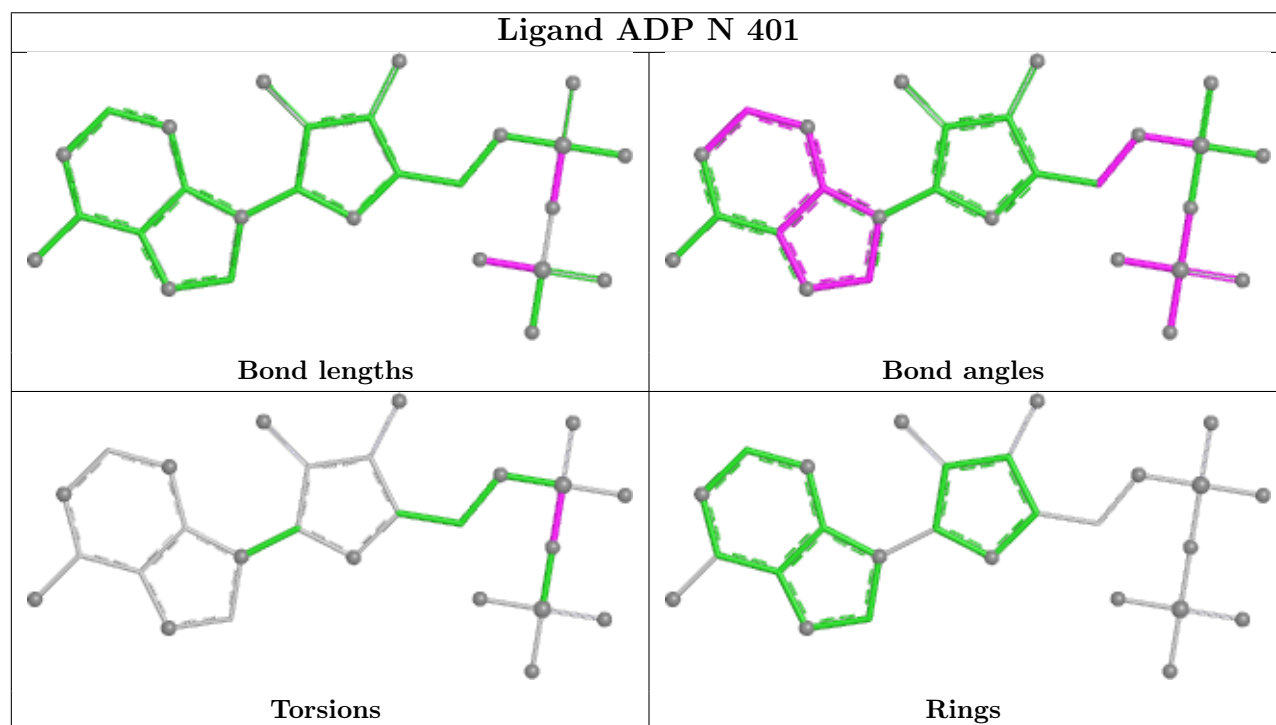
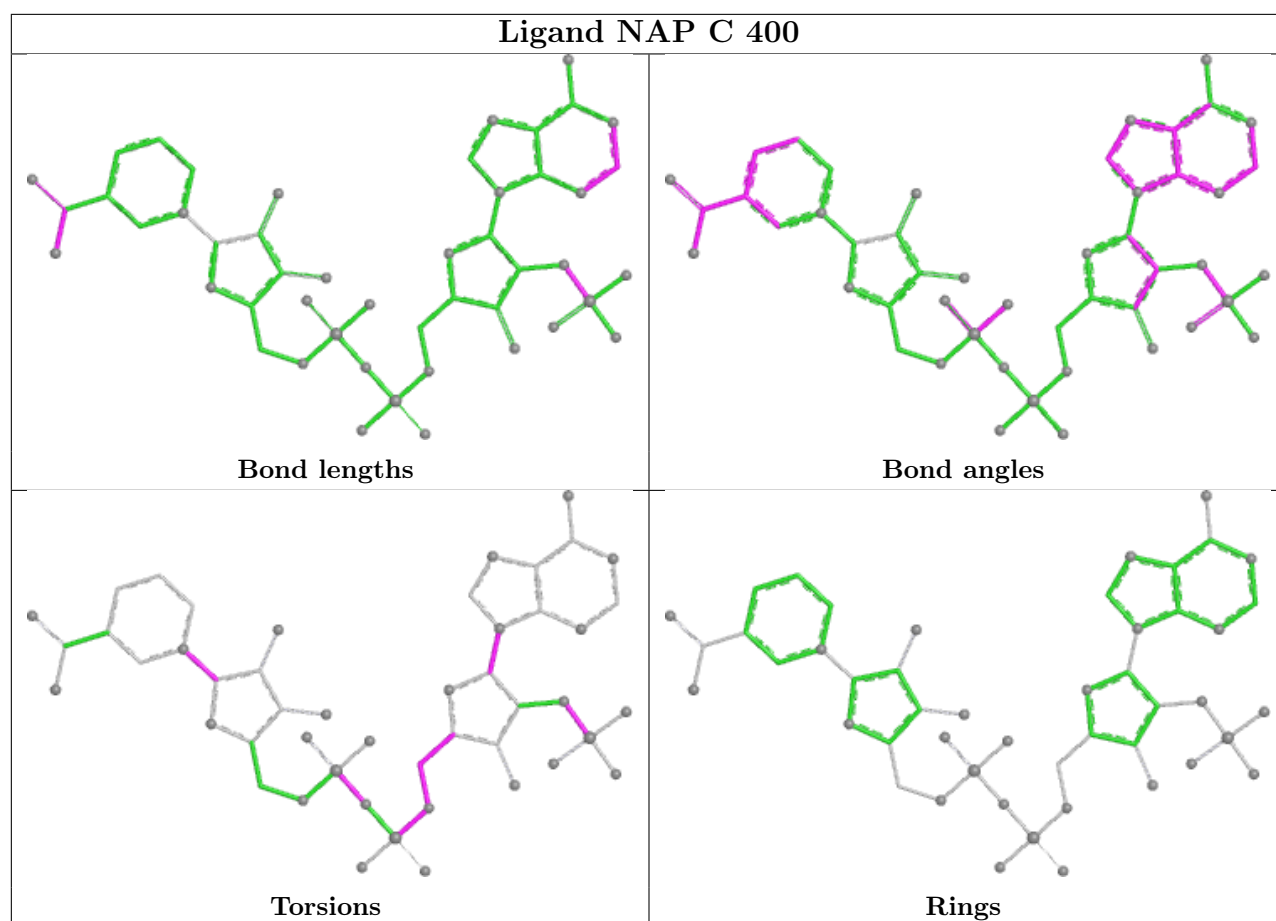


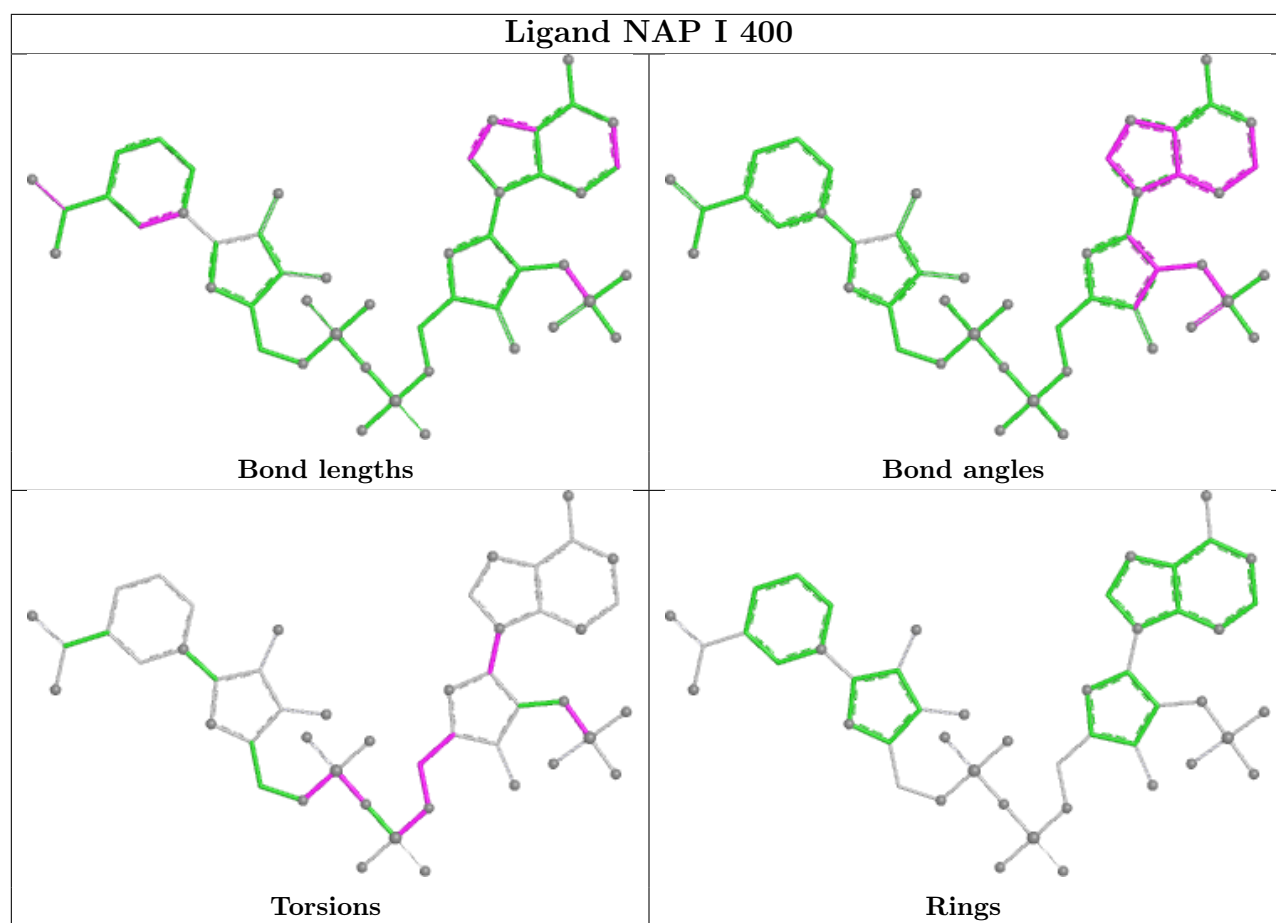


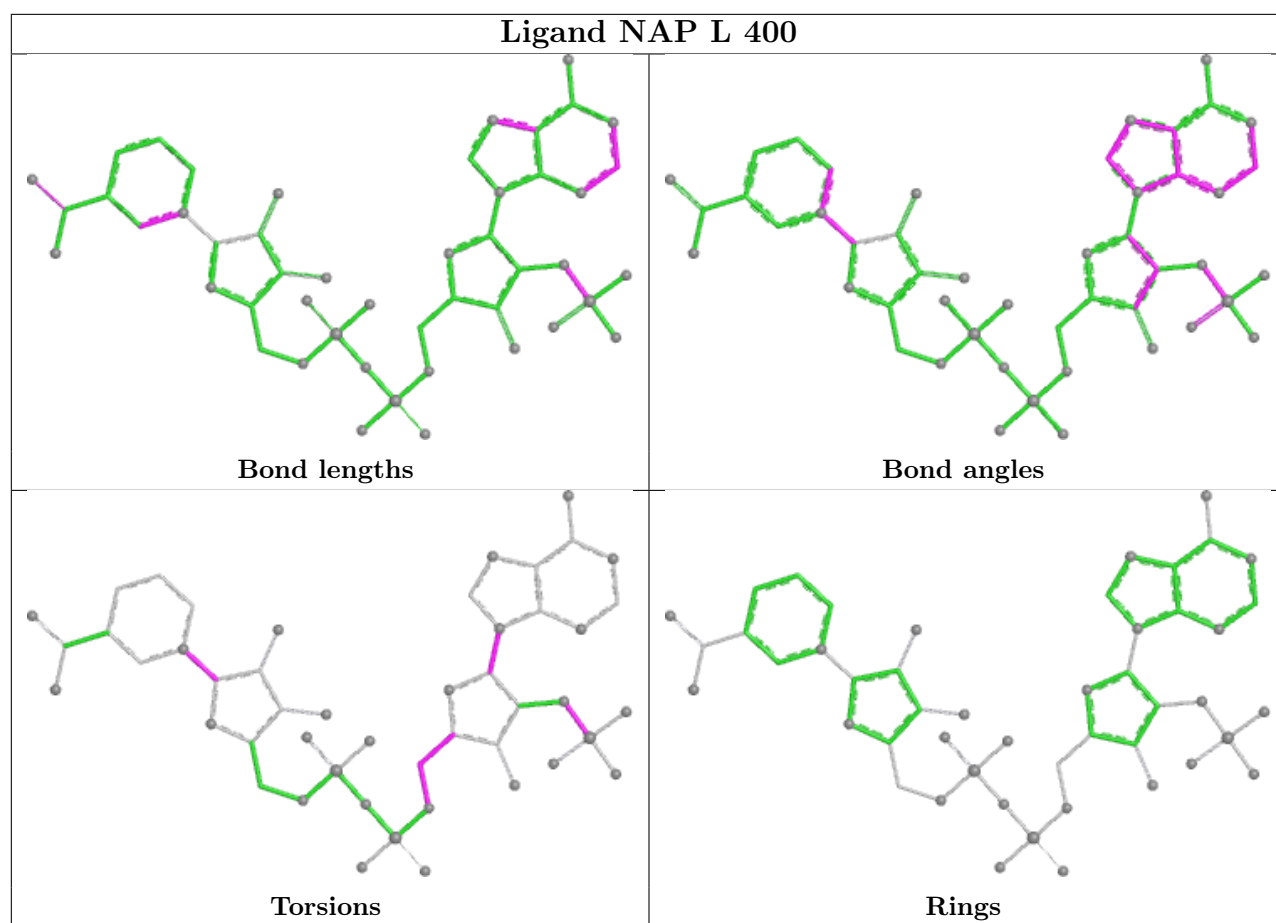


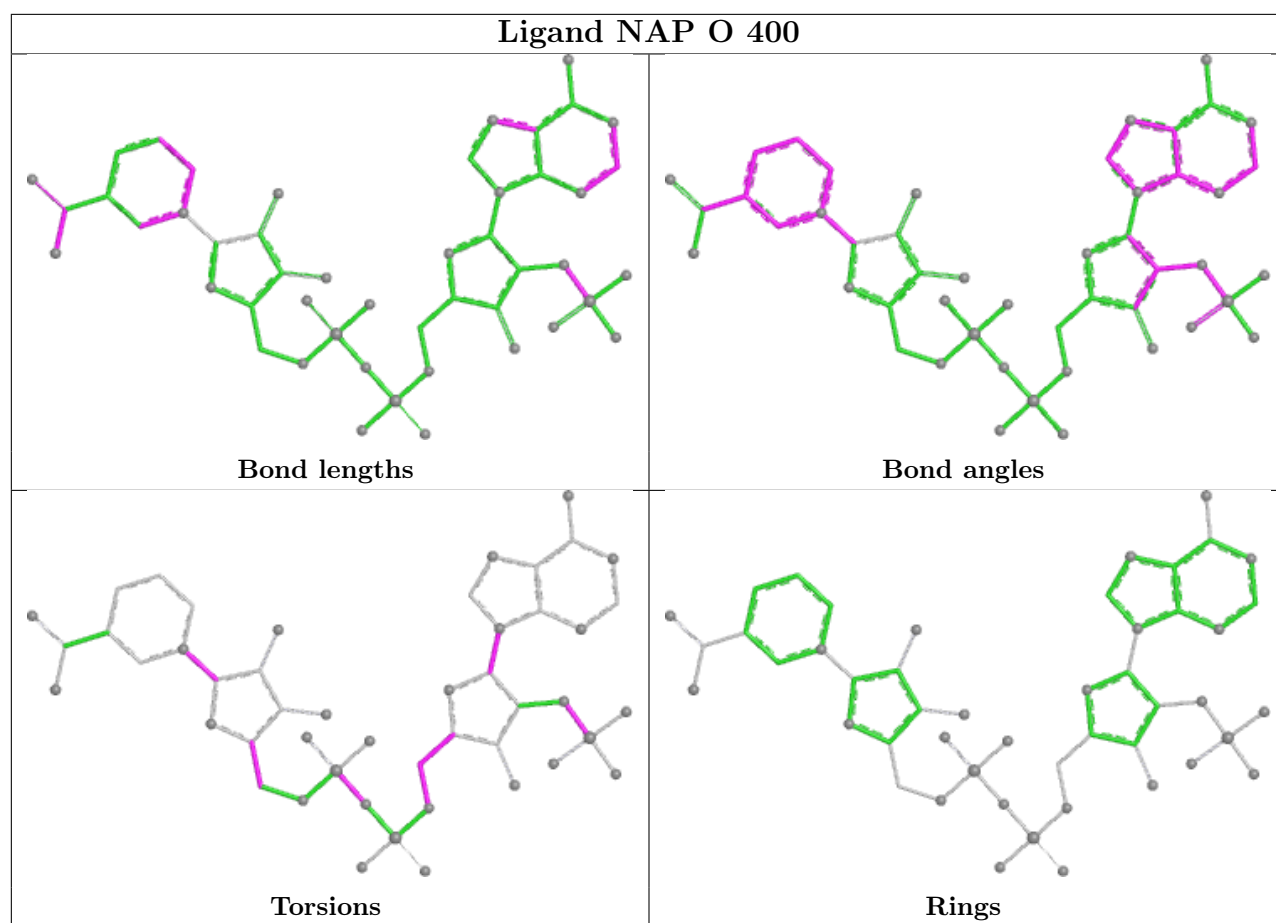


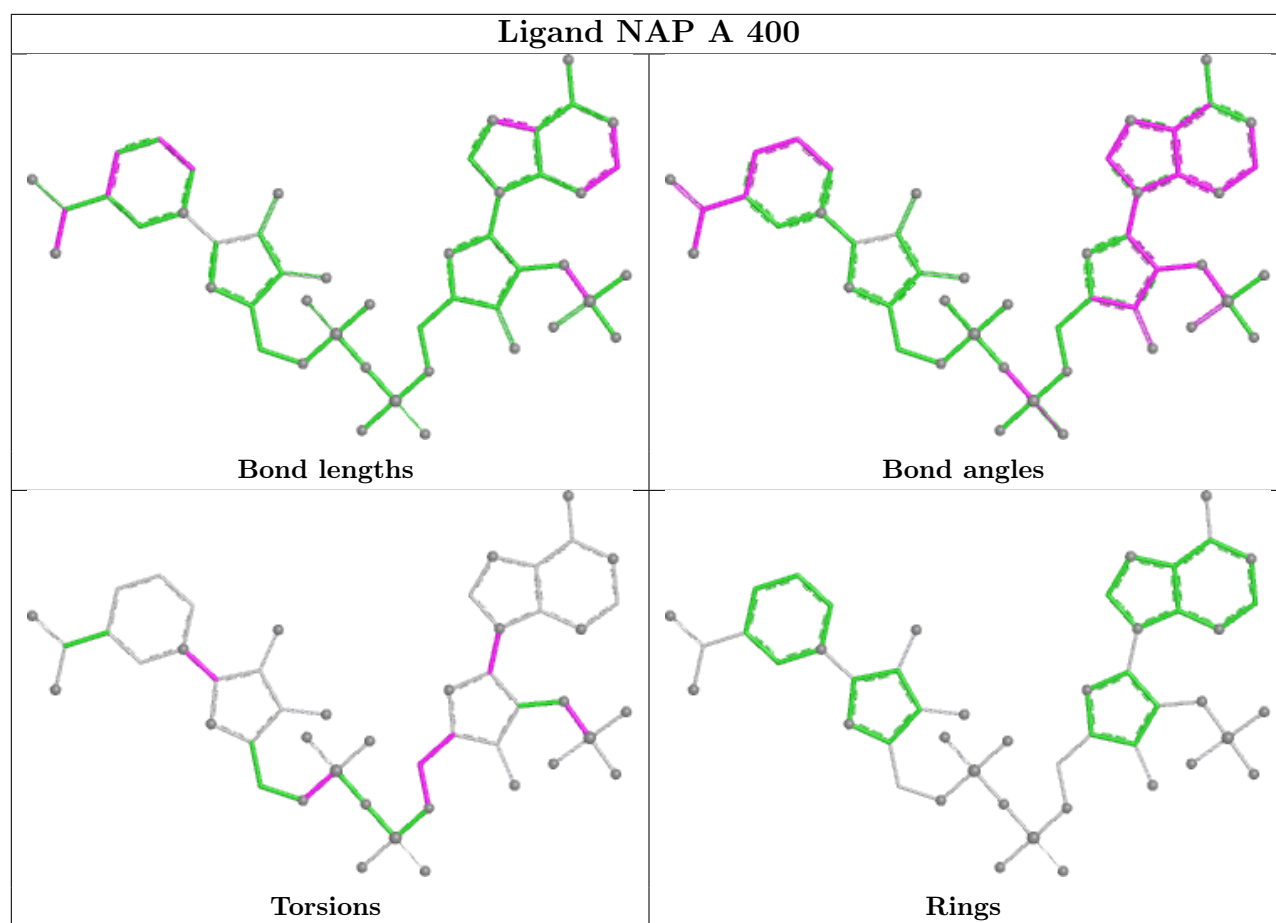


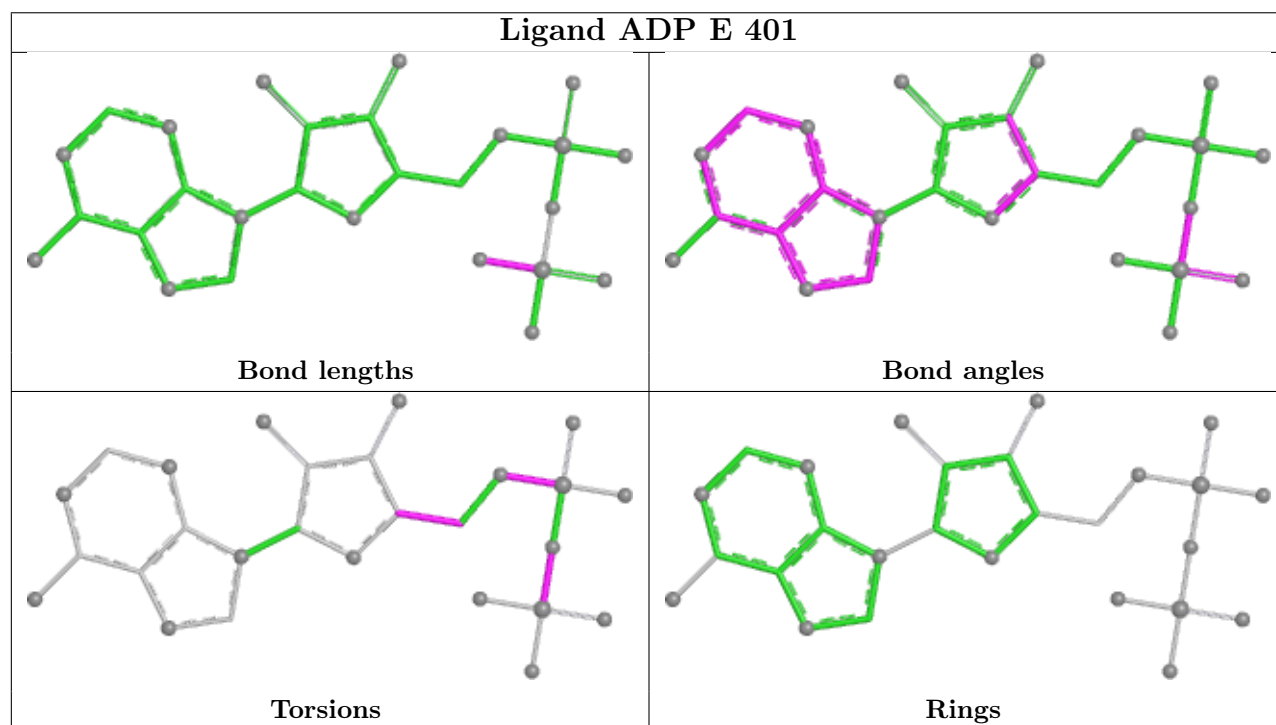
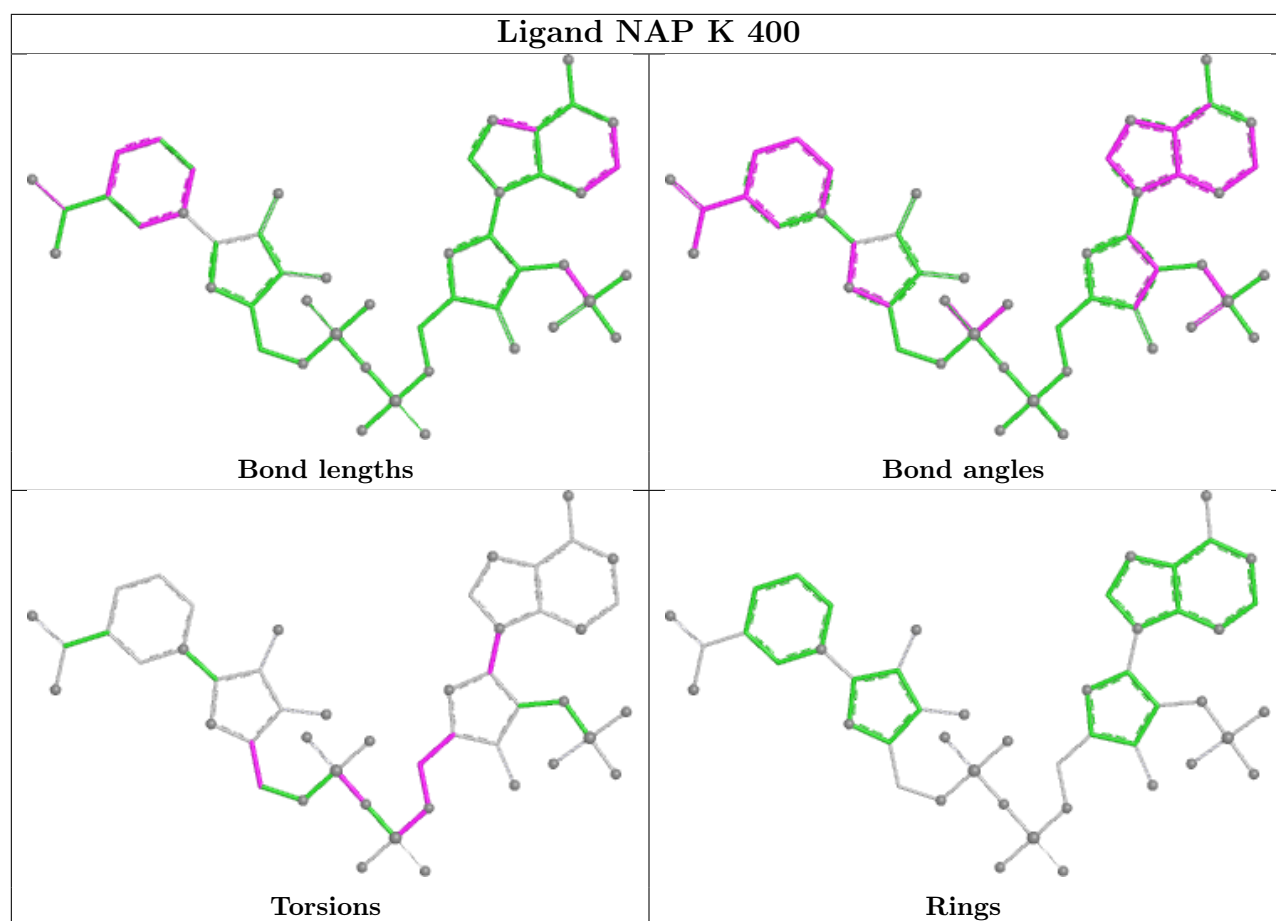


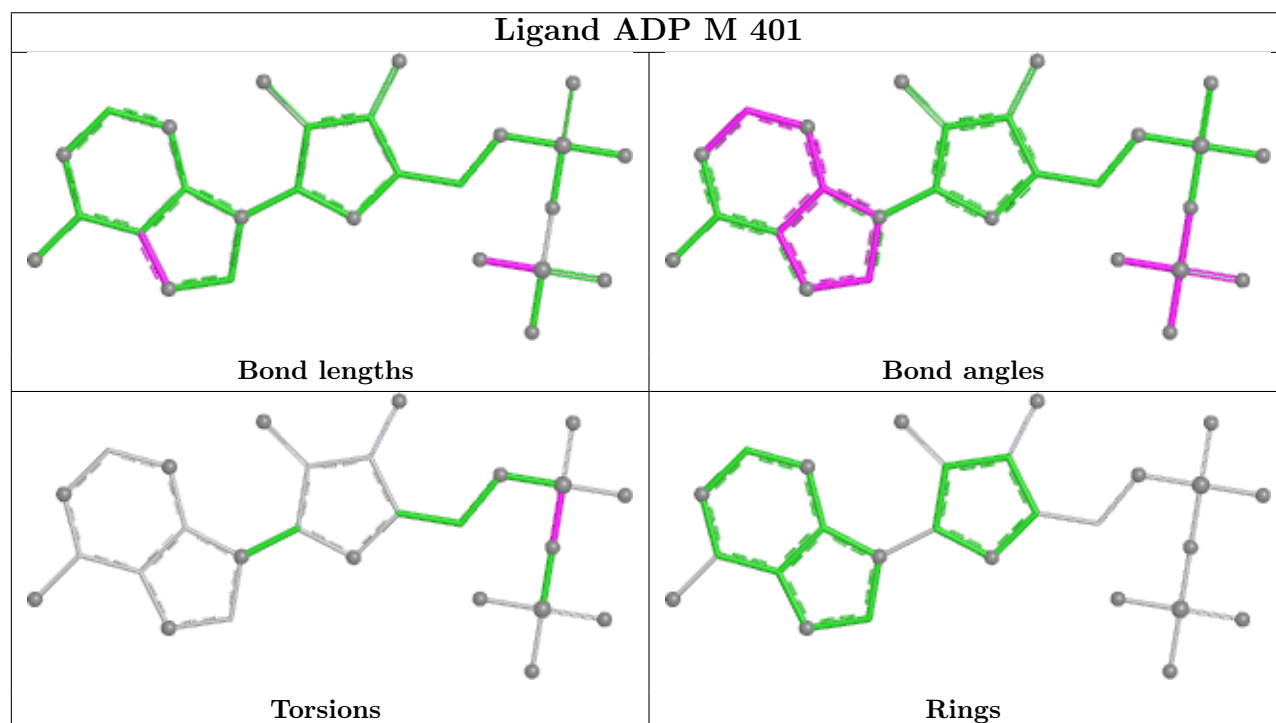
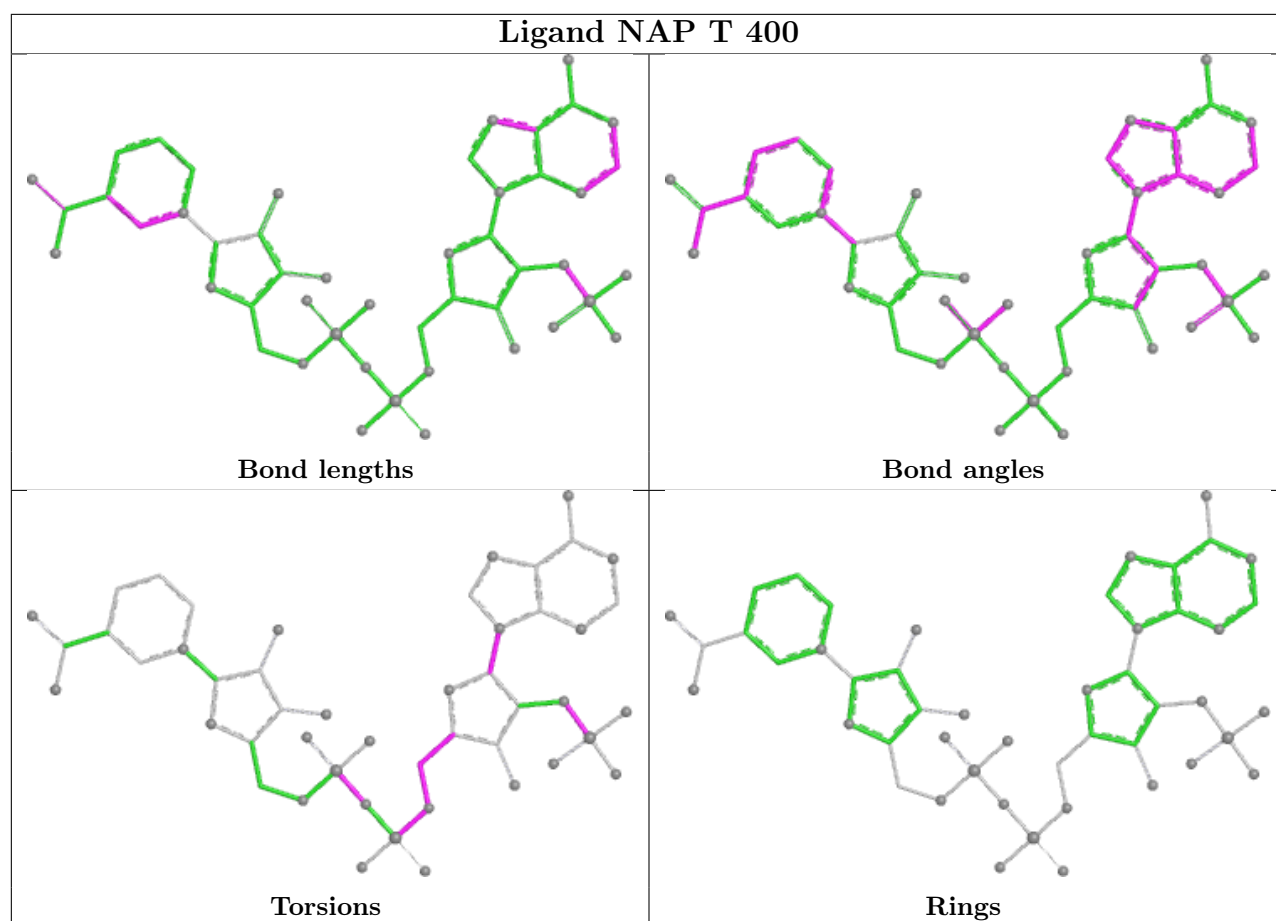




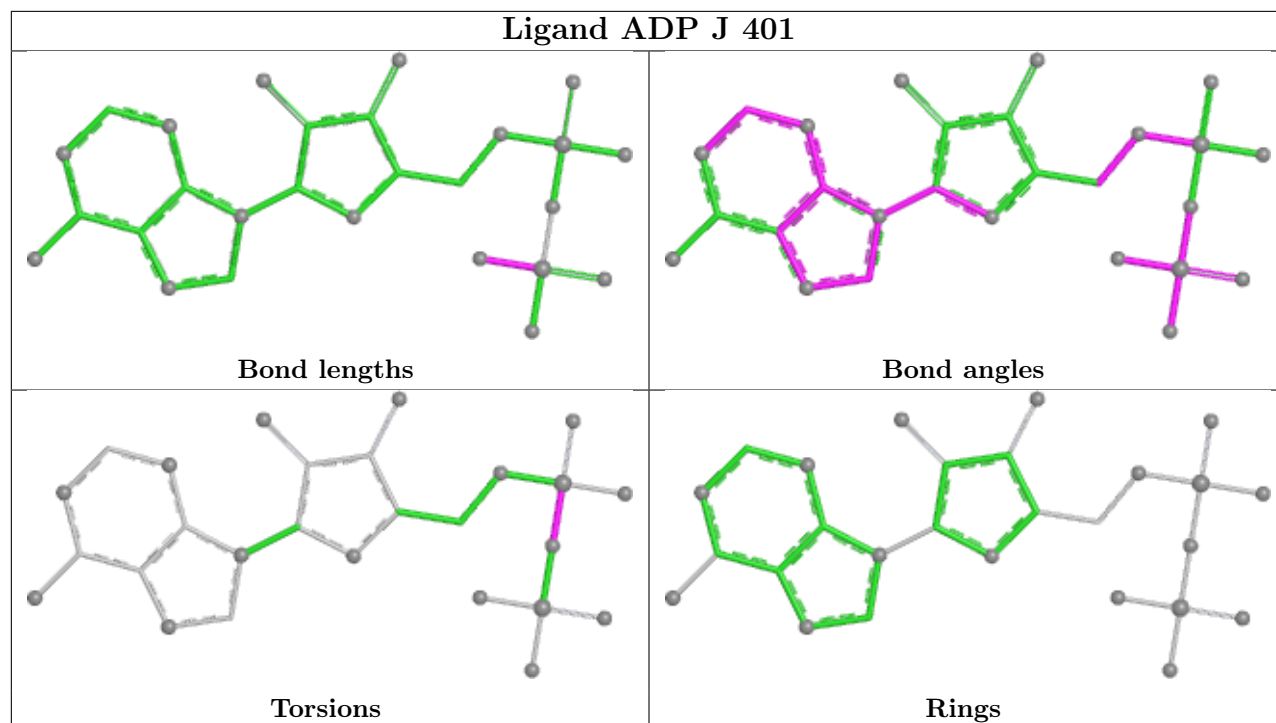




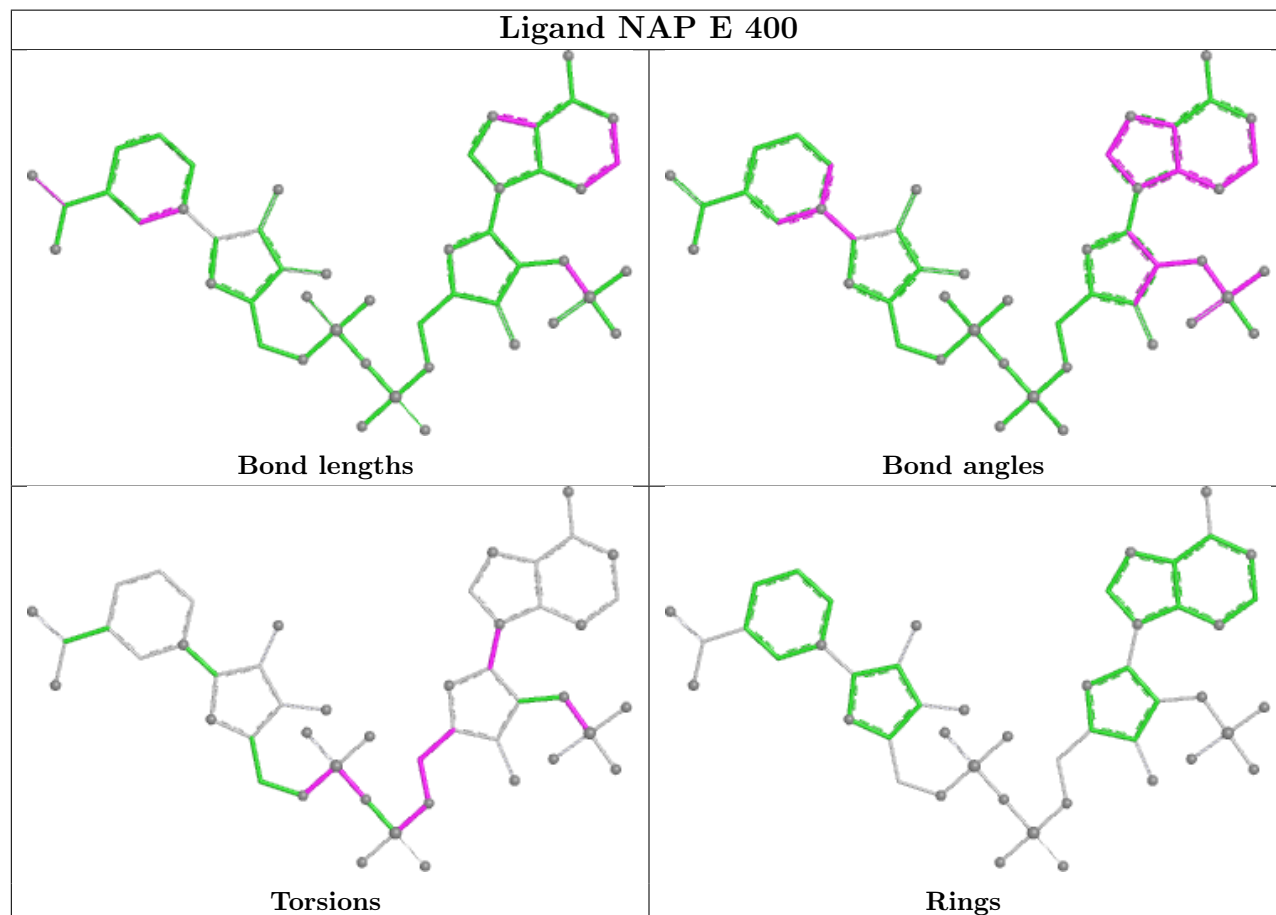


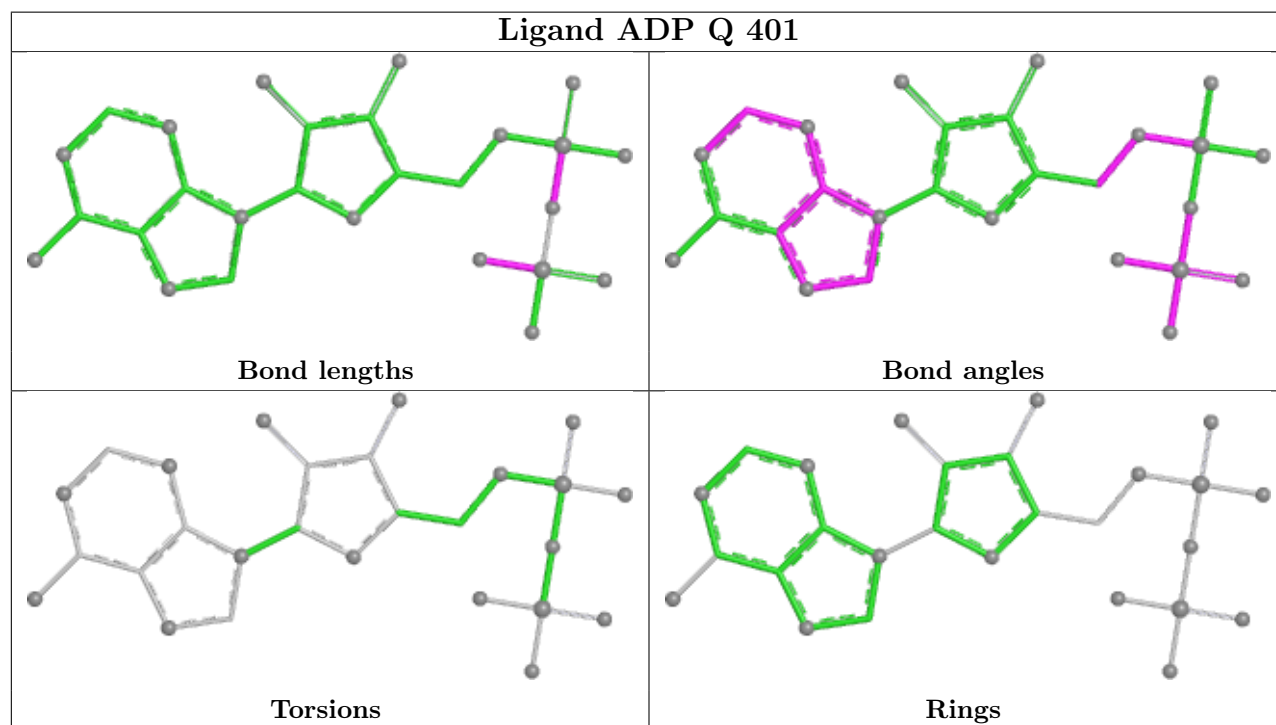
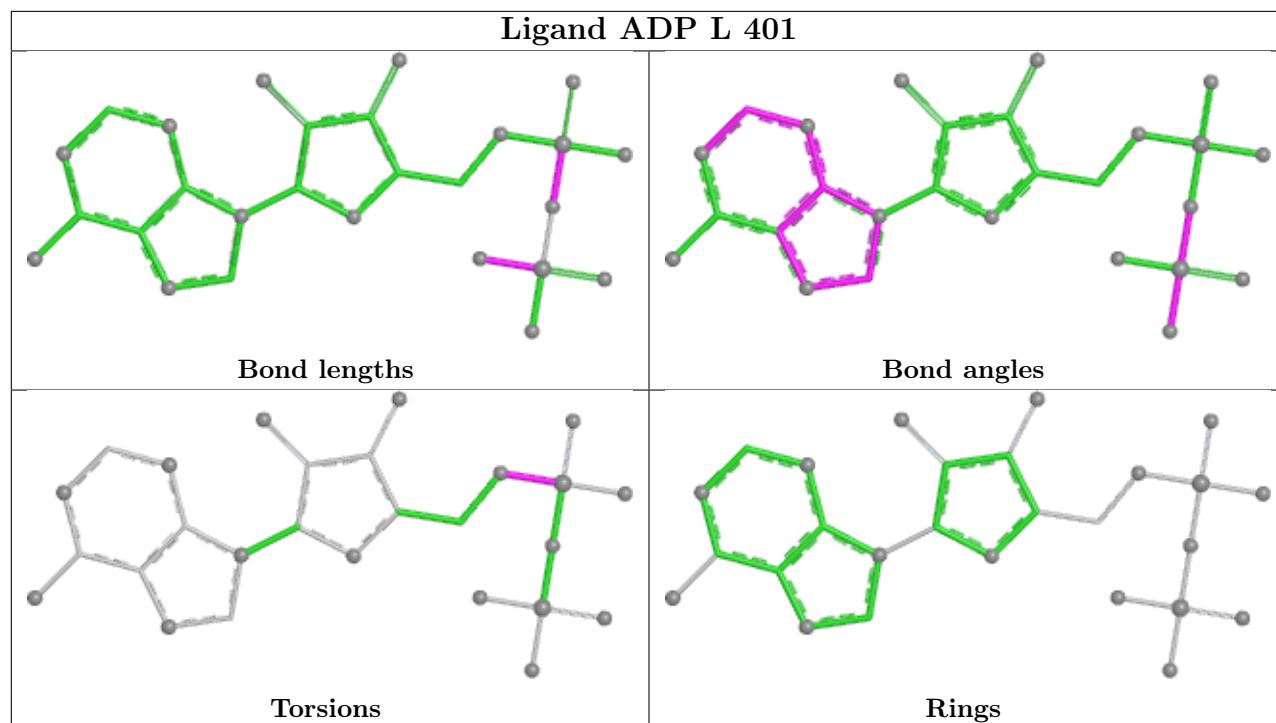


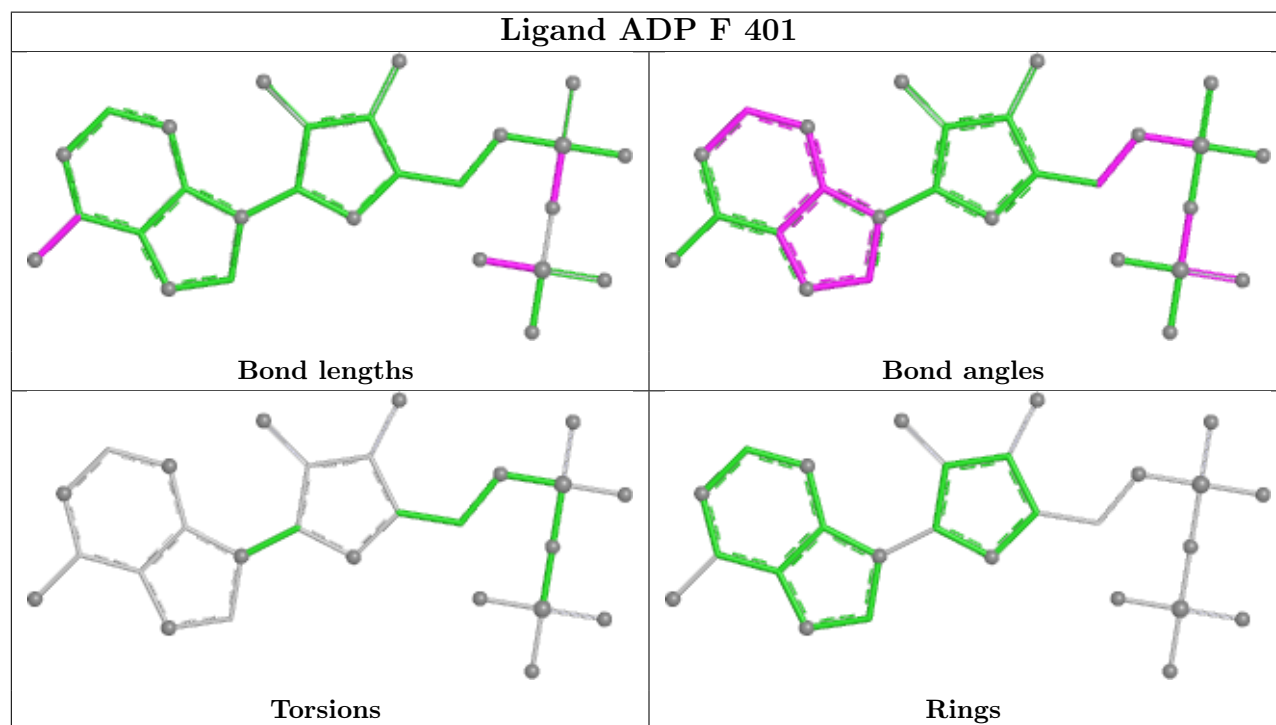
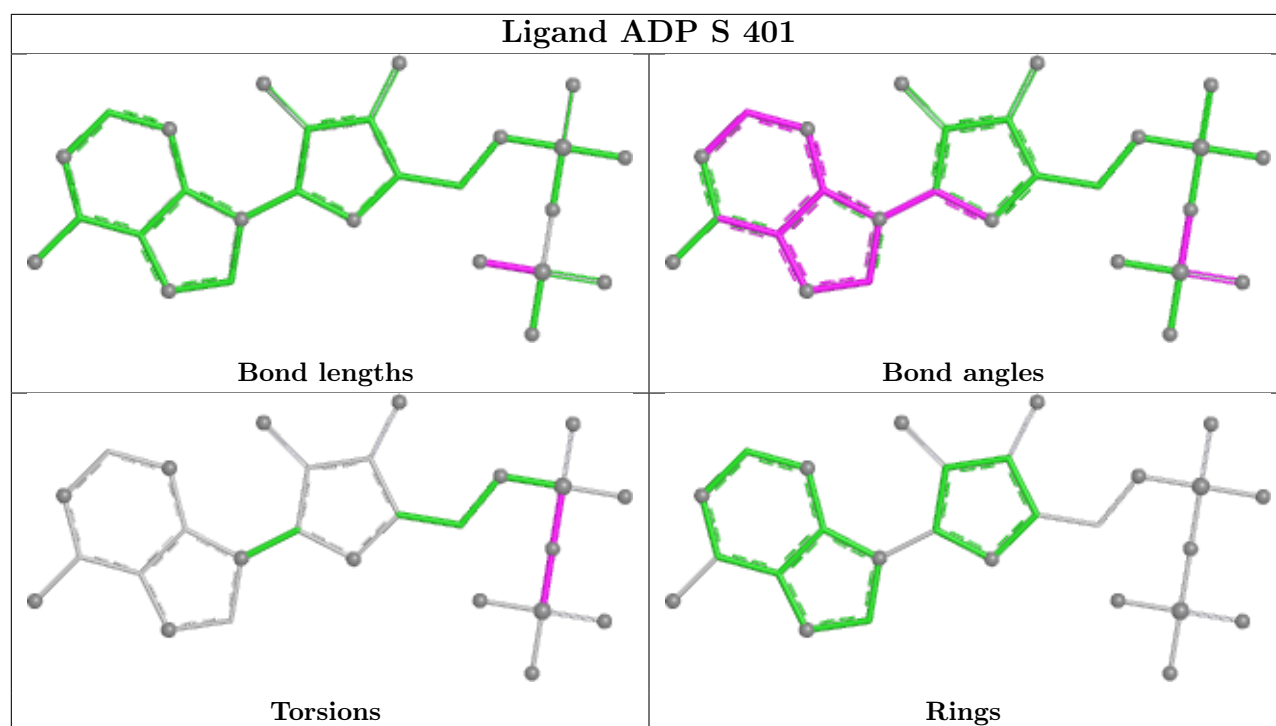
Ligand ADP J 401

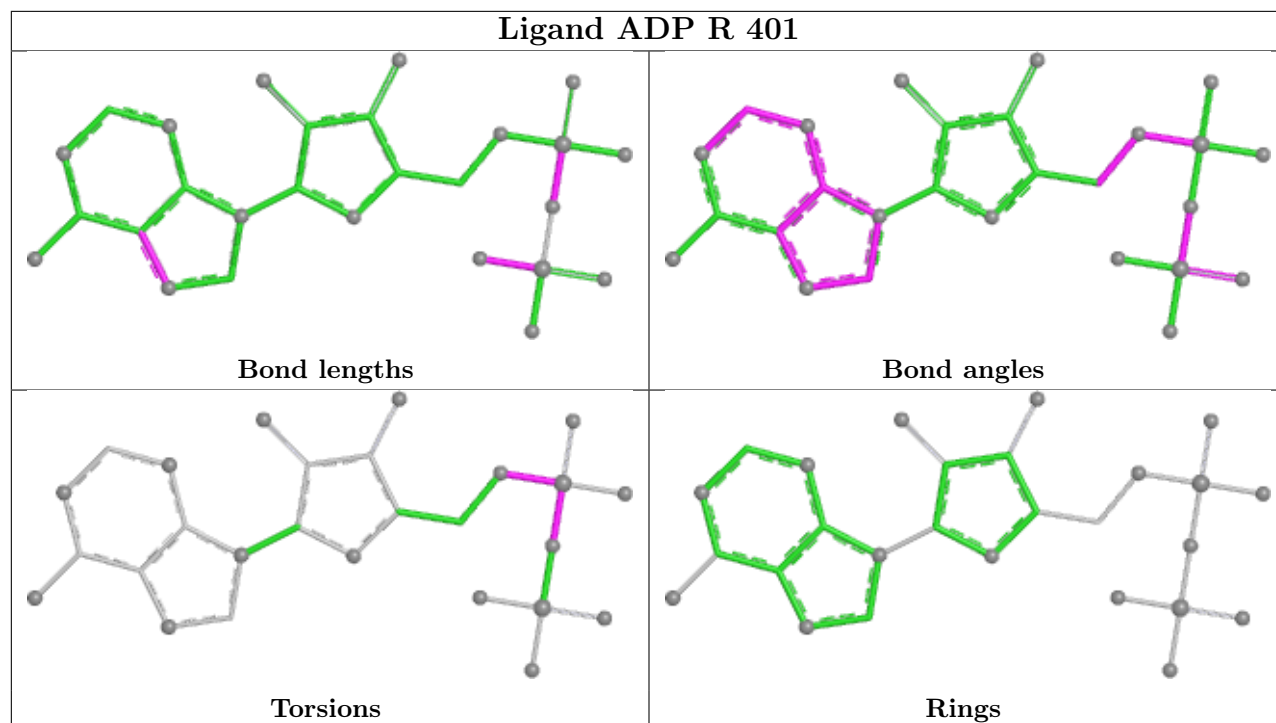


Ligand NAP E 400









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	307/357 (85%)	0.39	6 (1%) 65 56	6, 11, 26, 36	0
1	B	307/357 (85%)	0.74	14 (4%) 37 29	4, 12, 24, 36	0
1	C	307/357 (85%)	0.69	14 (4%) 37 29	6, 12, 20, 30	0
1	D	307/357 (85%)	0.19	2 (0%) 84 77	6, 11, 20, 30	0
1	E	307/357 (85%)	0.11	1 (0%) 90 86	4, 10, 19, 26	0
1	F	307/357 (85%)	0.30	9 (2%) 53 43	5, 11, 25, 36	0
1	G	307/357 (85%)	1.15	49 (15%) 5 4	4, 11, 32, 48	0
1	H	307/357 (85%)	0.38	4 (1%) 75 66	6, 11, 20, 28	0
1	I	307/357 (85%)	0.13	1 (0%) 90 86	6, 11, 19, 27	0
1	J	307/357 (85%)	0.15	1 (0%) 90 86	5, 11, 20, 27	0
1	K	307/357 (85%)	0.38	6 (1%) 65 56	6, 12, 19, 31	0
1	L	307/357 (85%)	0.10	3 (0%) 79 72	5, 11, 20, 27	0
1	M	307/357 (85%)	0.16	1 (0%) 90 86	5, 11, 22, 30	0
1	N	307/357 (85%)	0.70	19 (6%) 26 20	5, 11, 26, 32	0
1	O	307/357 (85%)	1.09	42 (13%) 6 5	6, 12, 22, 34	0
1	P	307/357 (85%)	0.41	10 (3%) 49 39	5, 11, 20, 33	0
1	Q	307/357 (85%)	0.20	3 (0%) 79 72	5, 11, 19, 30	0
1	R	307/357 (85%)	0.47	6 (1%) 65 56	6, 11, 18, 32	0
1	S	307/357 (85%)	1.21	46 (14%) 5 4	4, 12, 22, 36	0
1	T	307/357 (85%)	0.51	13 (4%) 40 32	5, 11, 20, 27	0
All	All	6140/7140 (85%)	0.47	250 (4%) 41 33	4, 11, 22, 48	0

The worst 5 of 250 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	S	243	PHE	6.2
1	O	201	PHE	5.1
1	B	265	PRO	5.0
1	S	257	GLY	4.7
1	S	240	ALA	4.7

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 [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ADP	G	401	27/27	0.73	0.15	30,32,34,35	0
4	BMA	G	402	11/12	0.74	0.24	29,31,32,32	0
4	BMA	C	402	11/12	0.76	0.20	20,20,21,22	0
4	BMA	A	402	11/12	0.76	0.20	33,35,41,44	0
4	BMA	S	402	11/12	0.77	0.20	22,23,24,24	0
4	BMA	H	402	11/12	0.78	0.18	22,23,23,23	0
3	ADP	S	401	27/27	0.78	0.16	17,19,19,21	0
3	ADP	O	401	27/27	0.79	0.15	14,16,17,18	0
3	ADP	B	401	27/27	0.81	0.14	15,19,20,20	0
4	BMA	R	402	11/12	0.82	0.18	21,23,26,26	0
4	BMA	P	402	11/12	0.82	0.16	24,25,28,29	0
3	ADP	N	401	27/27	0.83	0.14	10,12,20,25	0
4	BMA	N	402	11/12	0.83	0.17	27,28,31,34	0
4	BMA	O	402	11/12	0.83	0.20	18,18,19,19	0
4	BMA	J	402	11/12	0.84	0.15	22,23,26,27	0
4	BMA	Q	402	11/12	0.85	0.16	18,18,19,21	0
3	ADP	C	401	27/27	0.85	0.14	14,15,17,19	0
3	ADP	T	401	27/27	0.85	0.13	10,11,18,21	0
3	ADP	A	401	27/27	0.86	0.12	14,17,23,28	0

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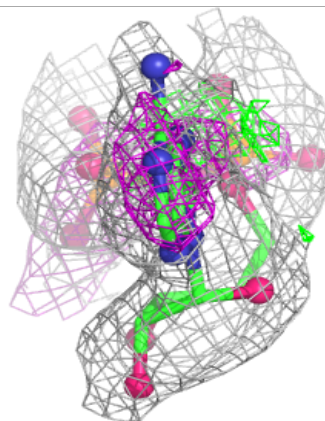
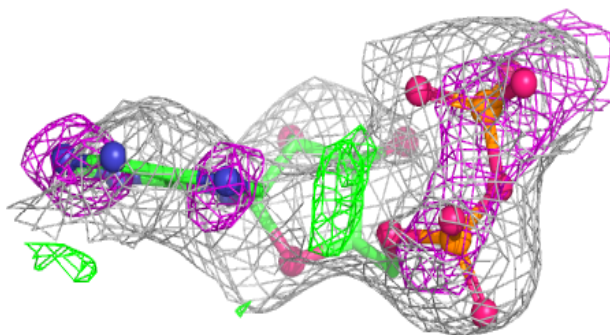
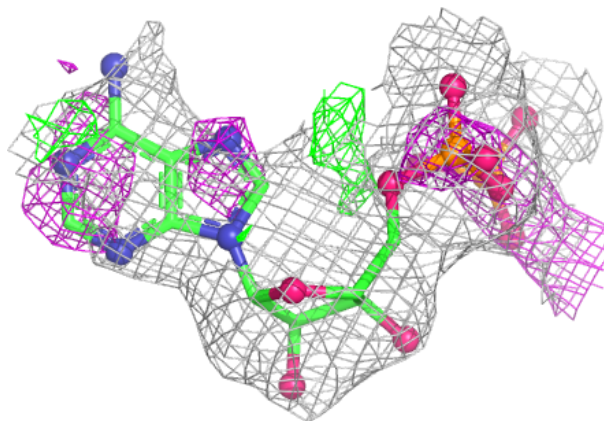
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ADP	F	401	27/27	0.86	0.13	13,14,19,21	0
4	BMA	E	402	11/12	0.87	0.13	19,20,23,23	0
4	BMA	T	402	11/12	0.87	0.14	24,27,29,30	0
4	BMA	D	402	11/12	0.88	0.12	20,21,22,24	0
4	BMA	F	402	11/12	0.88	0.17	24,25,28,30	0
4	BMA	K	402	11/12	0.89	0.12	25,27,30,32	0
3	ADP	H	401	27/27	0.89	0.12	15,16,20,22	0
4	BMA	I	402	11/12	0.90	0.10	16,17,17,18	0
2	NAP	B	400	48/48	0.91	0.12	10,14,27,28	0
2	NAP	O	400	48/48	0.91	0.11	10,12,26,27	0
4	BMA	L	402	11/12	0.91	0.13	15,17,19,20	0
2	NAP	S	400	48/48	0.91	0.11	12,14,23,23	0
3	ADP	K	401	27/27	0.91	0.11	14,14,21,23	0
2	NAP	G	400	48/48	0.92	0.11	10,13,29,30	0
3	ADP	D	401	27/27	0.92	0.10	12,13,17,19	0
2	NAP	T	400	48/48	0.92	0.11	9,12,25,26	0
3	ADP	Q	401	27/27	0.92	0.10	12,12,14,16	0
2	NAP	C	400	48/48	0.92	0.11	10,14,36,37	0
2	NAP	R	400	48/48	0.92	0.11	9,14,37,37	0
3	ADP	R	401	27/27	0.93	0.10	14,14,16,18	0
2	NAP	N	400	48/48	0.93	0.11	9,12,26,27	0
3	ADP	L	401	27/27	0.93	0.10	10,12,14,16	0
3	ADP	P	401	27/27	0.93	0.09	8,10,19,22	0
4	BMA	M	402	11/12	0.93	0.14	17,19,19,20	0
3	ADP	M	401	27/27	0.93	0.10	12,12,16,16	0
2	NAP	K	400	48/48	0.94	0.10	9,13,21,22	0
2	NAP	L	400	48/48	0.94	0.09	8,16,29,30	0
2	NAP	M	400	48/48	0.94	0.09	10,15,26,26	0
2	NAP	D	400	48/48	0.94	0.10	13,18,34,35	0
2	NAP	E	400	48/48	0.94	0.10	8,11,28,30	0
2	NAP	P	400	48/48	0.94	0.09	9,12,22,24	0
3	ADP	J	401	27/27	0.94	0.09	10,11,16,20	0
2	NAP	Q	400	48/48	0.94	0.09	11,14,28,29	0
2	NAP	A	400	48/48	0.94	0.10	9,15,27,29	0
2	NAP	H	400	48/48	0.94	0.09	10,12,25,26	0
2	NAP	I	400	48/48	0.94	0.09	10,13,26,27	0
2	NAP	J	400	48/48	0.94	0.10	9,13,21,23	0
3	ADP	E	401	27/27	0.95	0.08	8,10,15,18	0
2	NAP	F	400	48/48	0.95	0.09	8,13,24,25	0
3	ADP	I	401	27/27	0.95	0.08	12,12,15,16	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.

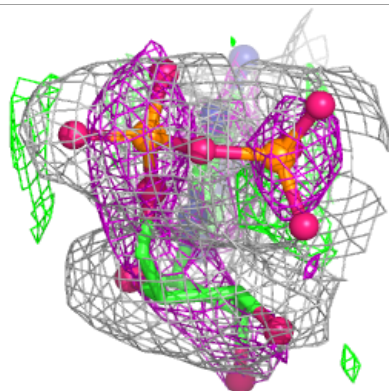
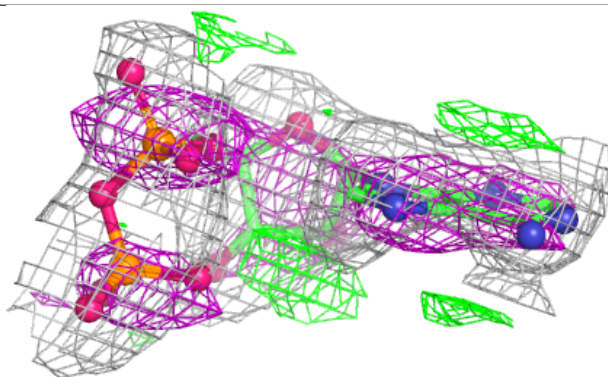
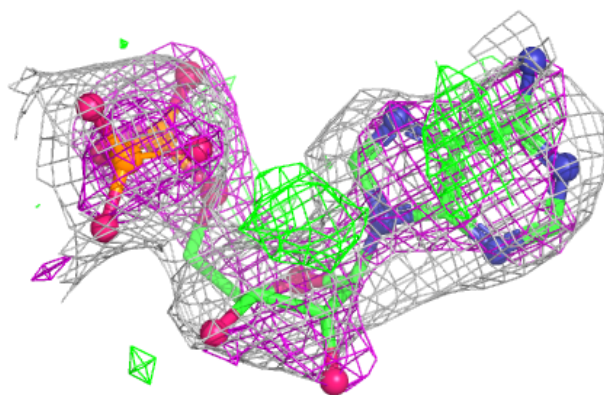
Electron density around ADP G 401:

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and green (positive)

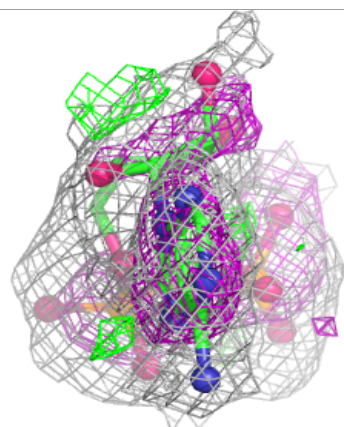
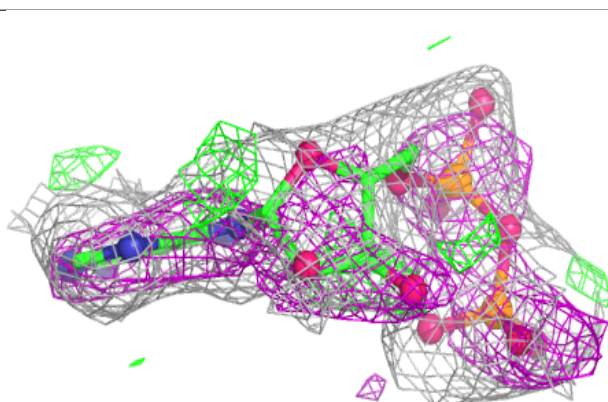
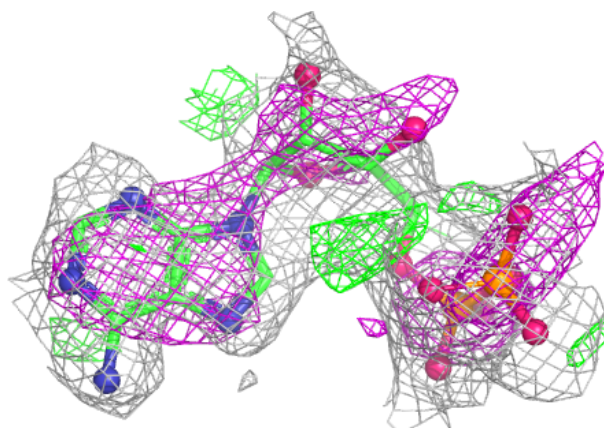


Electron density around ADP S 401:

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and green (positive)

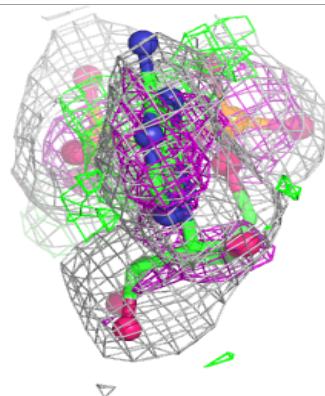
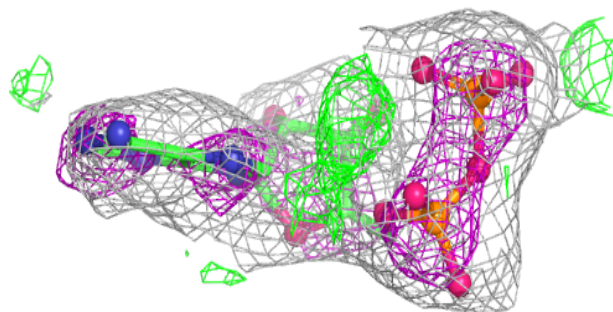
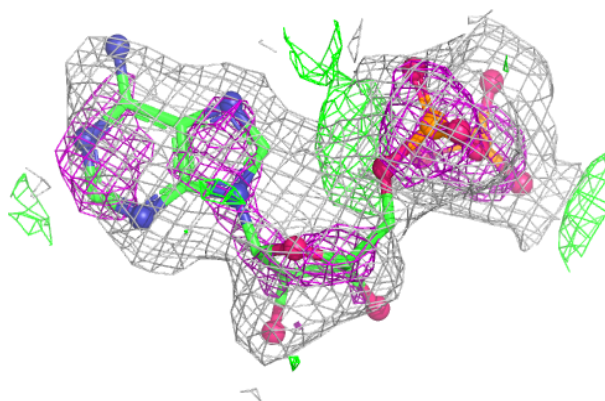
**Electron density around ADP O 401:**

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and green (positive)

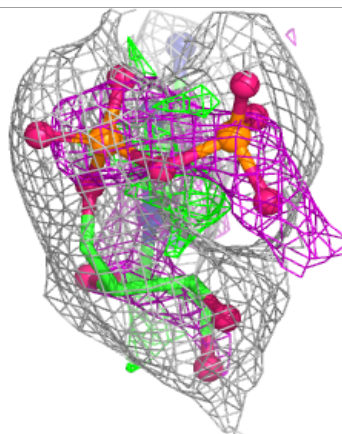
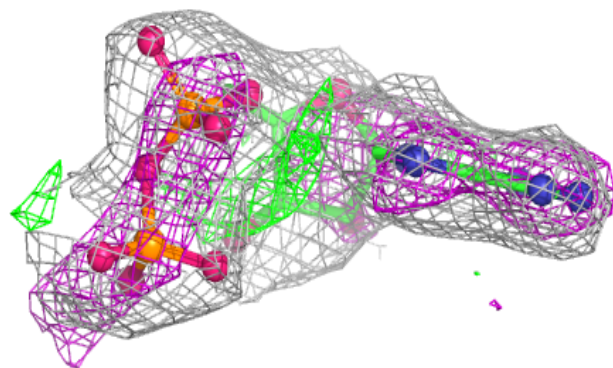
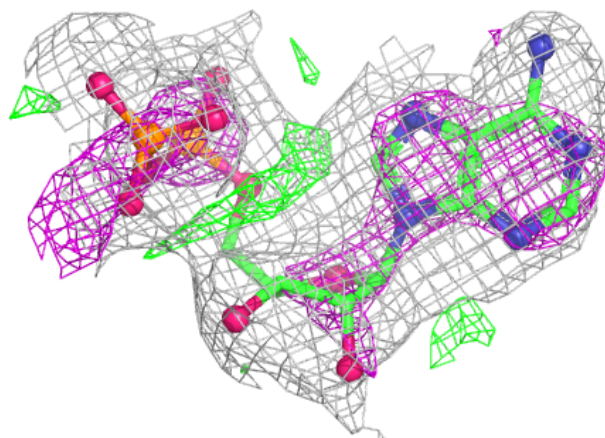


Electron density around ADP B 401:

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and green (positive)

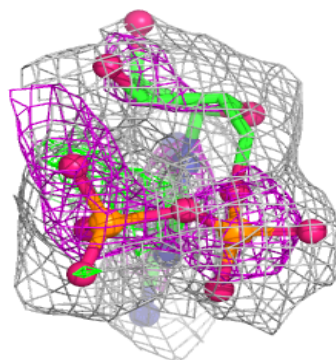
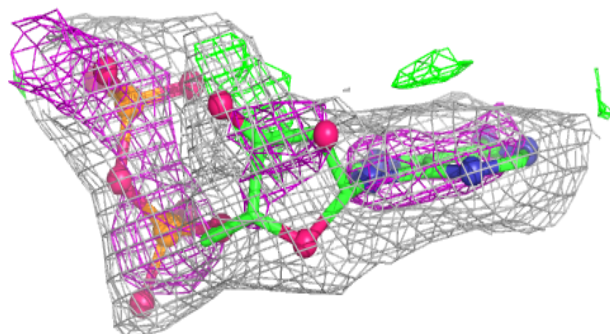
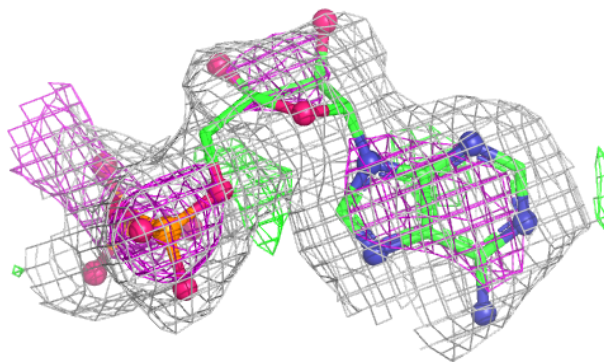
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and green (positive)

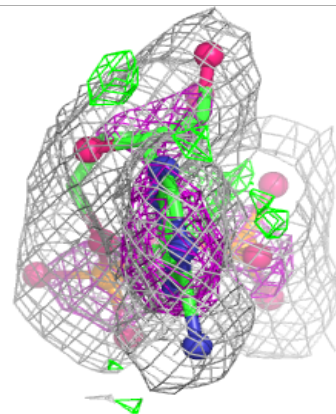
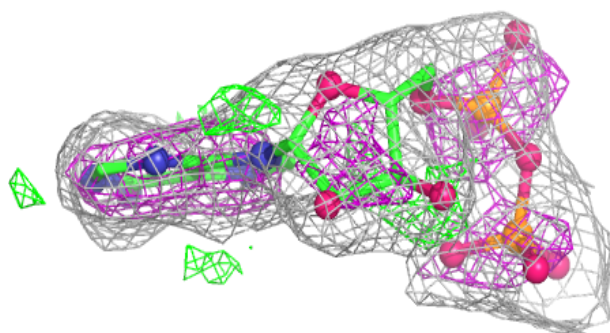
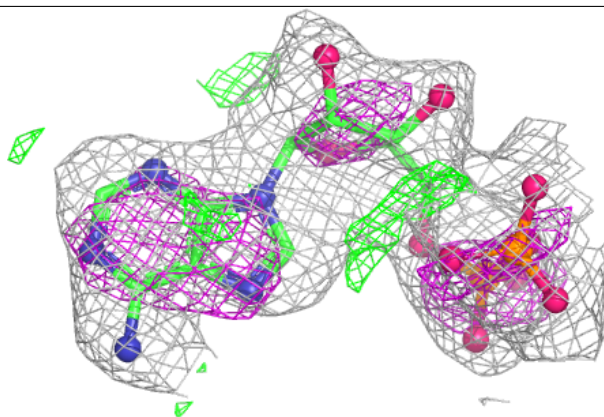


Electron density around ADP C 401:

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and green (positive)

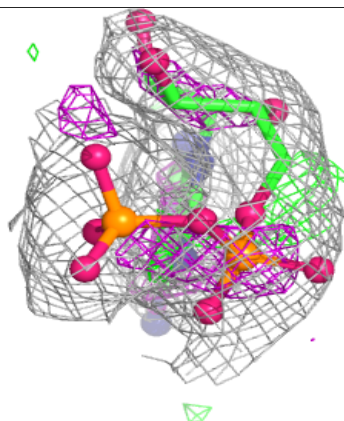
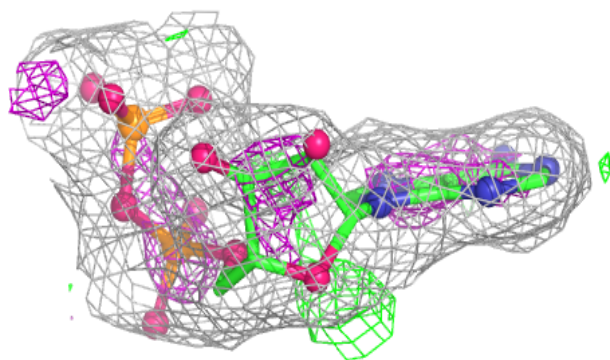
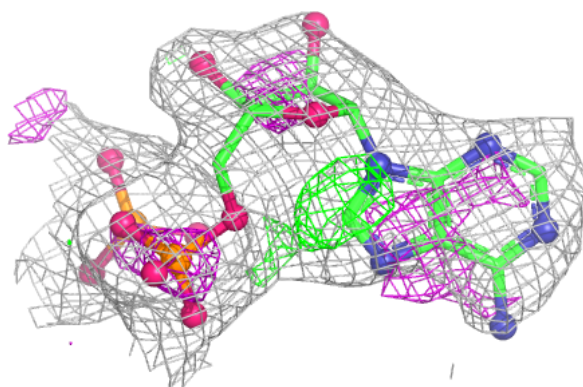
**Electron density around ADP T 401:**

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and green (positive)

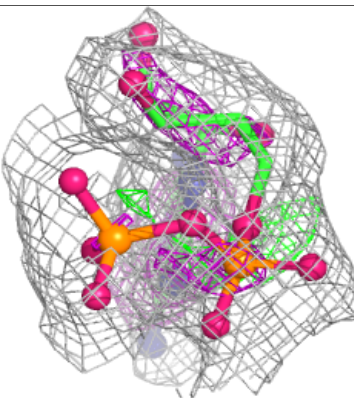
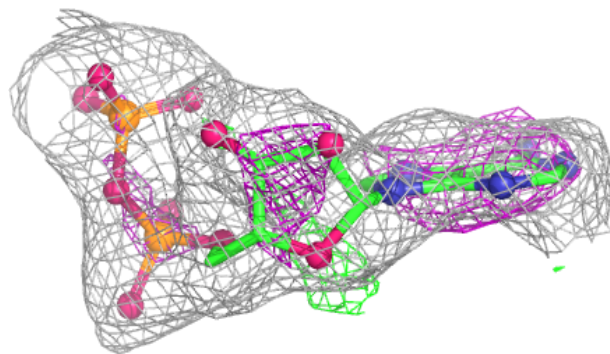
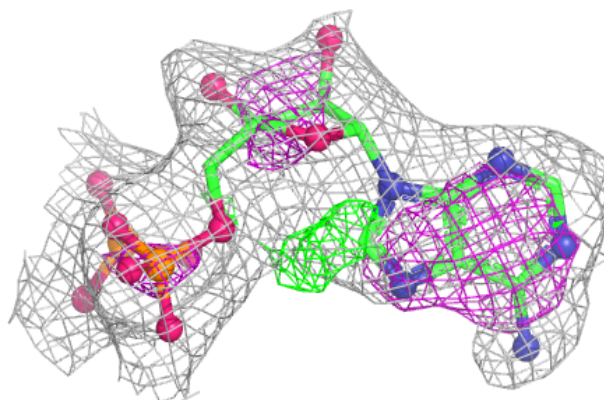


Electron density around ADP A 401:

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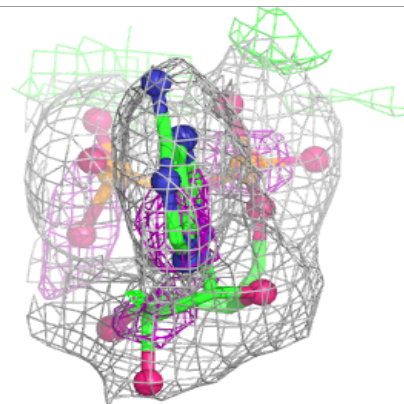
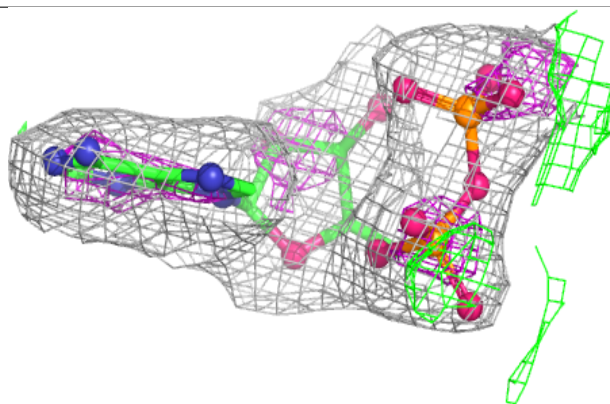
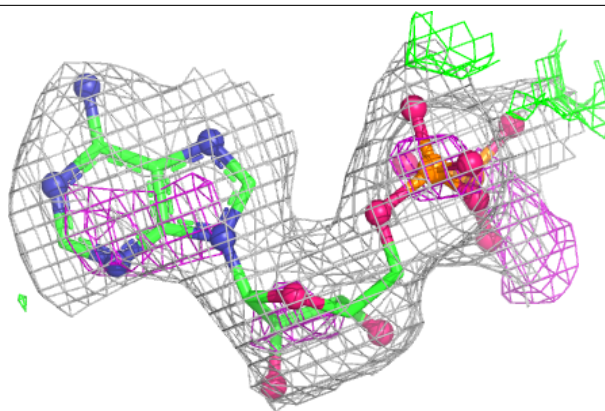
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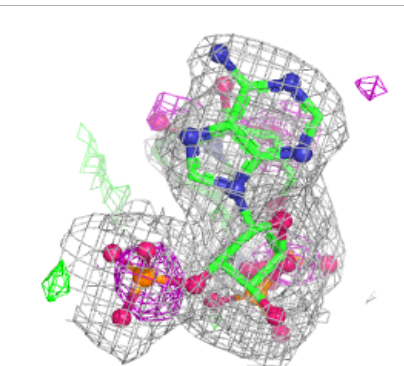
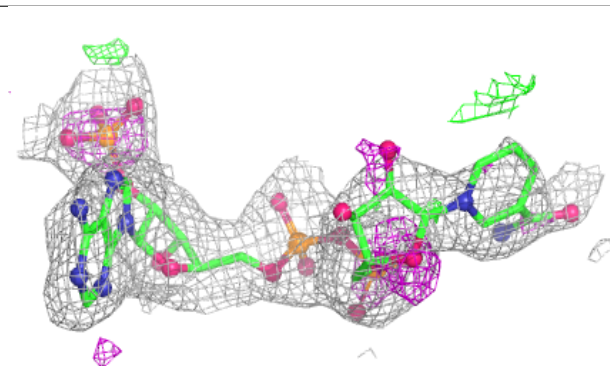
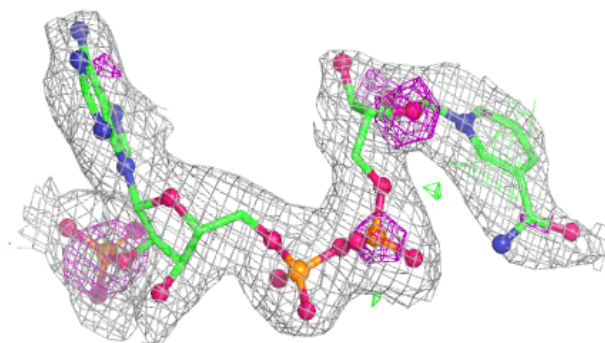


Electron density around ADP H 401:

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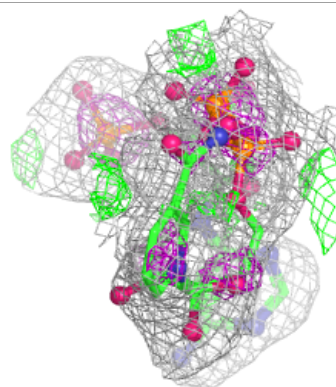
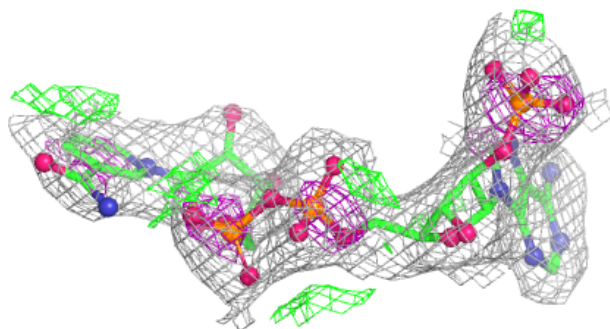
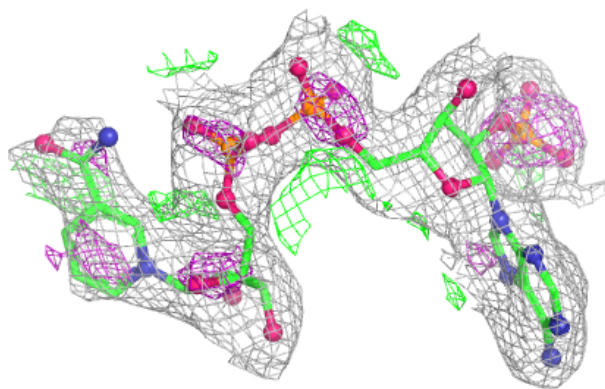
**Electron density around NAP B 400:**

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and green (positive)

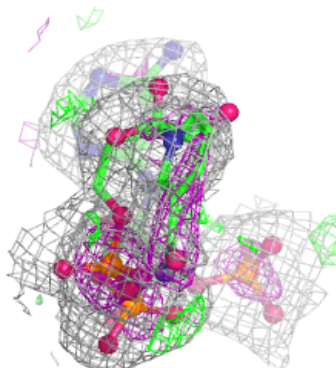
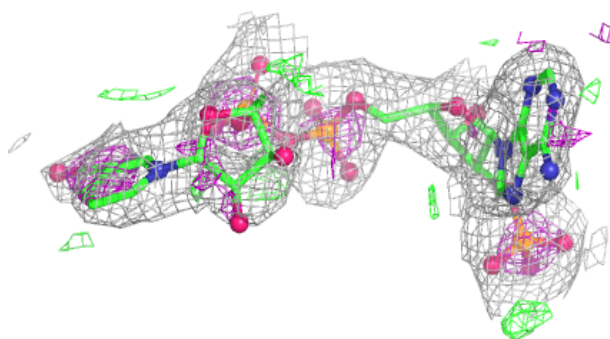
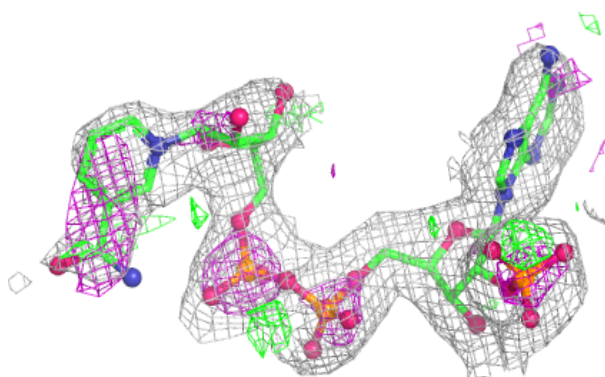


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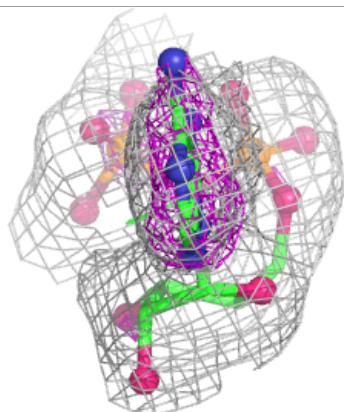
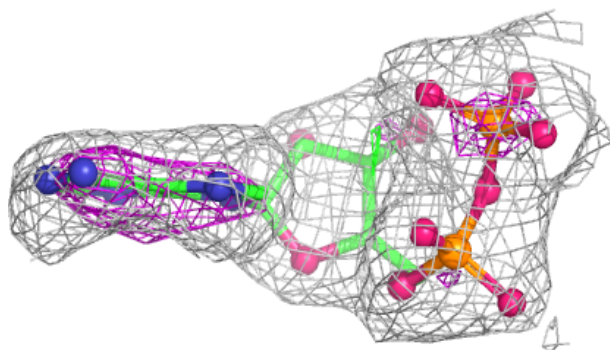
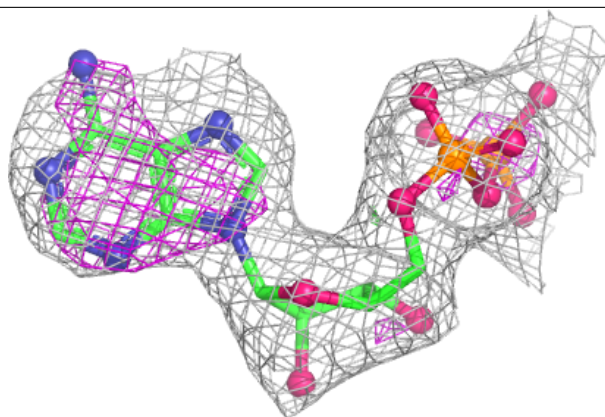
**Electron density around NAP S 400:**

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and green (positive)

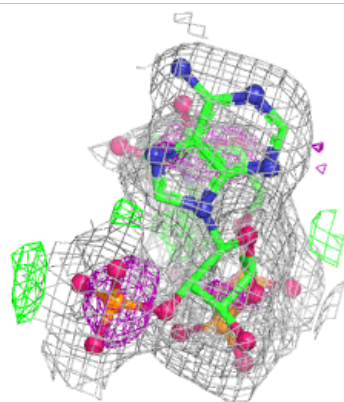
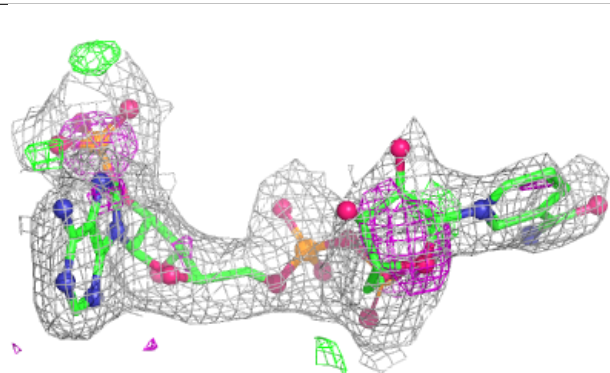
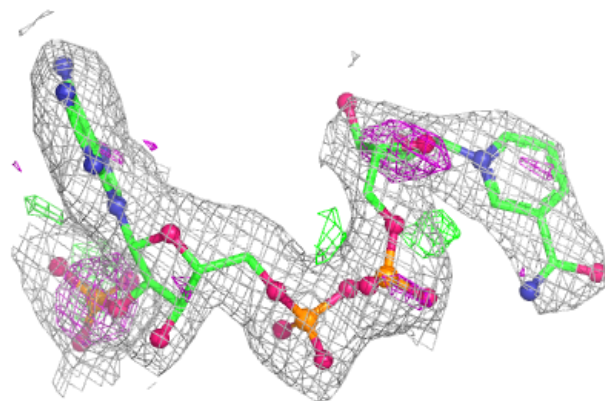


Electron density around ADP K 401:

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

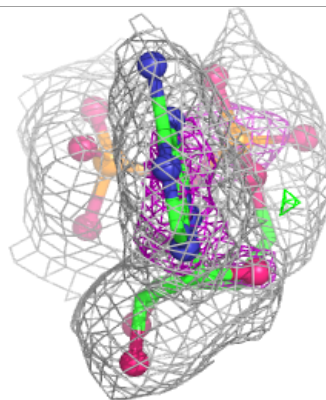
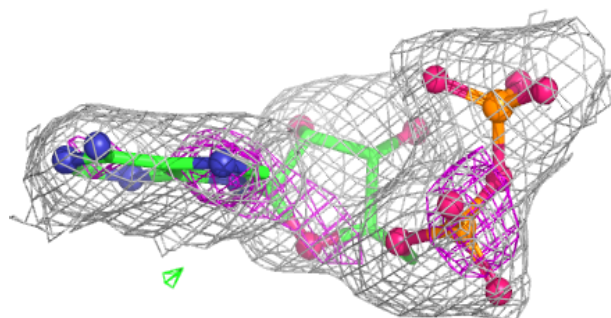
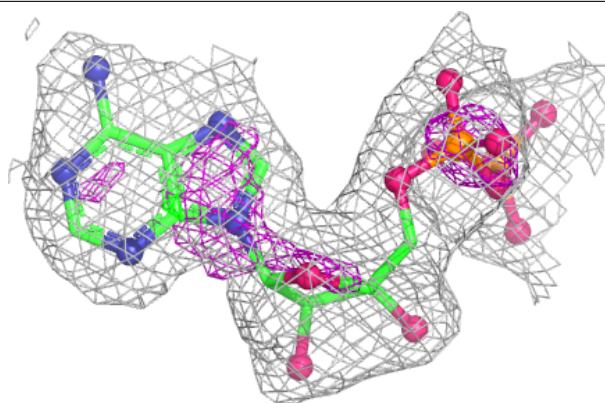
**Electron density around NAP G 400:**

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

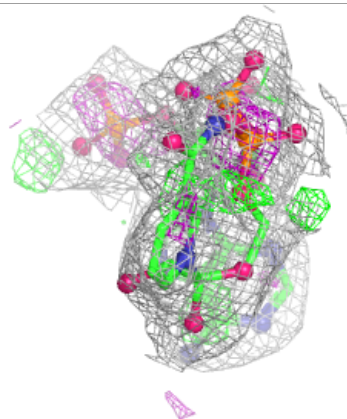
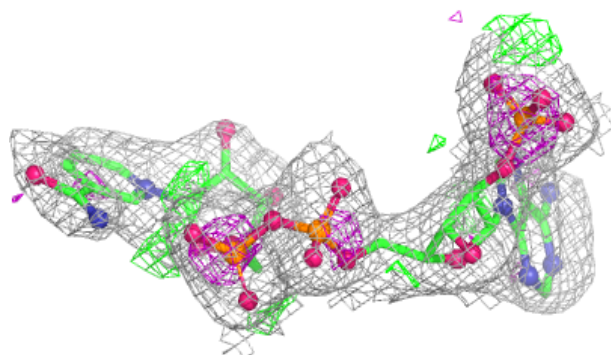
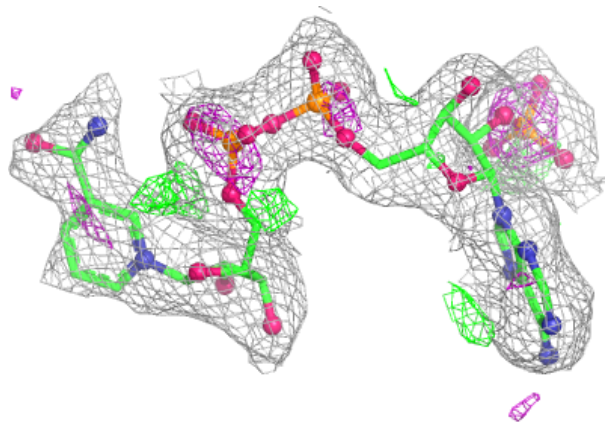


Electron density around ADP D 401:

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

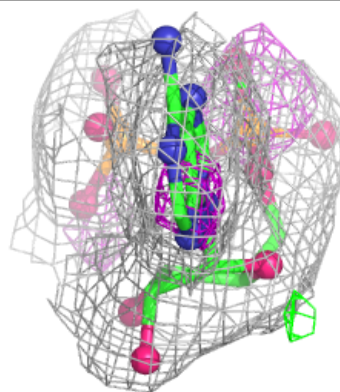
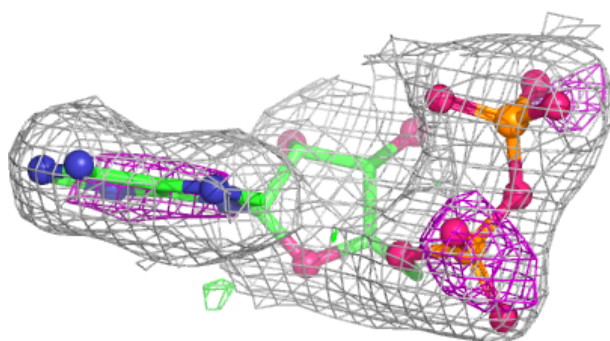
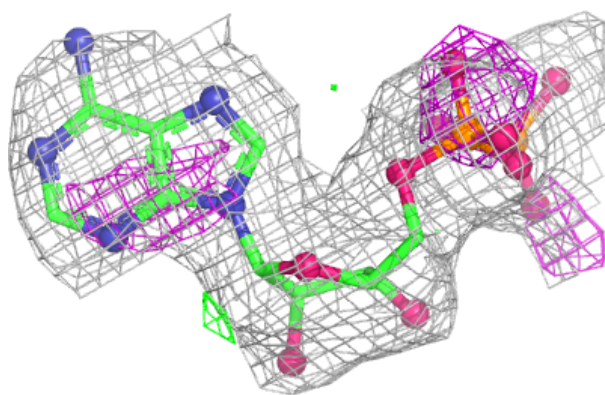
**Electron density around NAP T 400:**

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

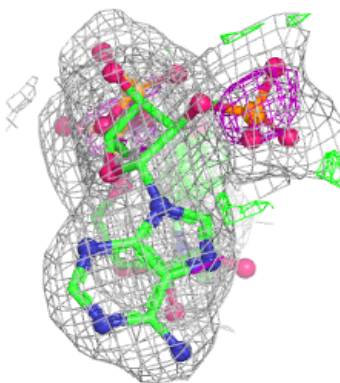
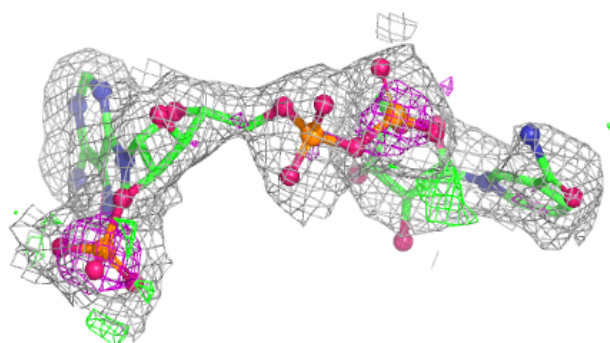
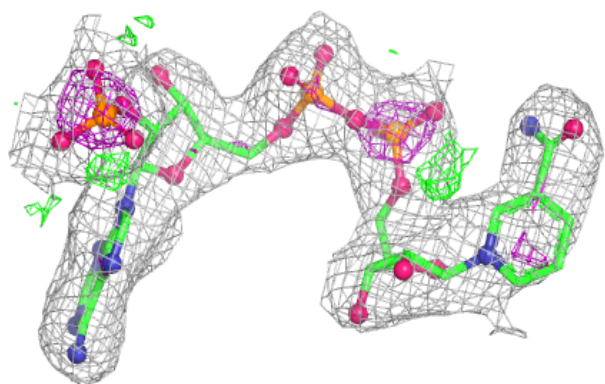


Electron density around ADP Q 401:

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

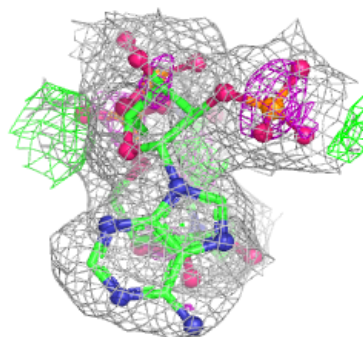
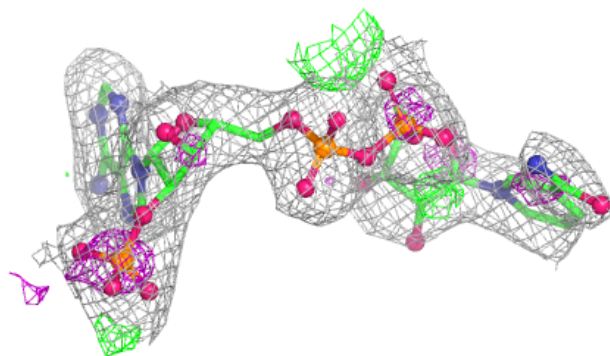
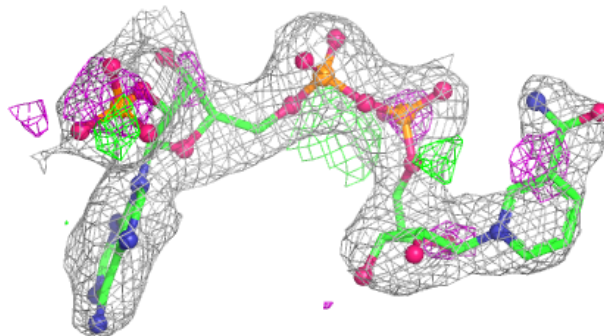
**Electron density around NAP C 400:**

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

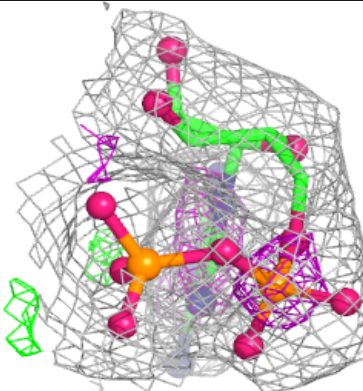
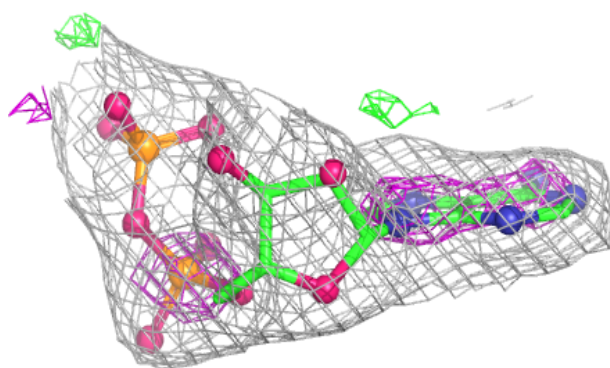
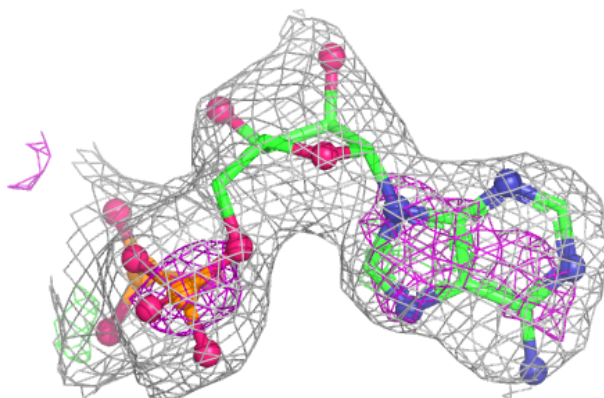


Electron density around NAP R 400:

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

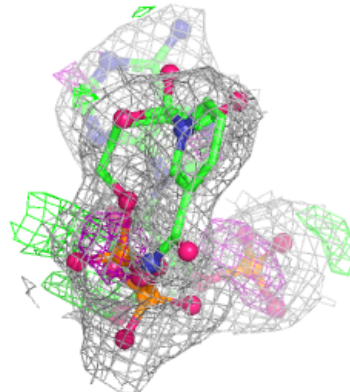
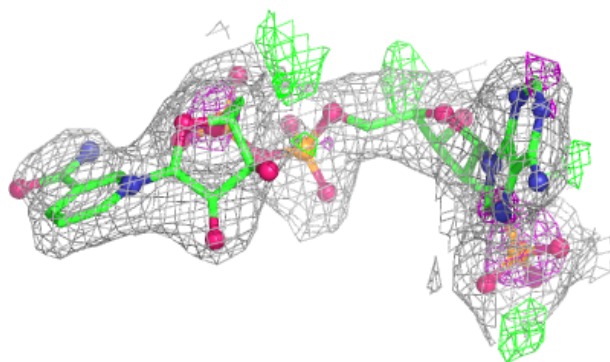
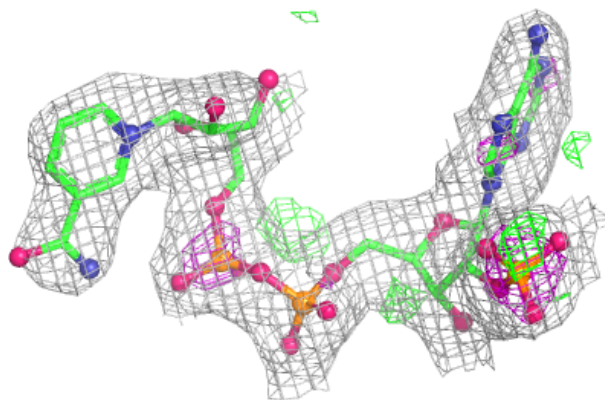
**Electron density around ADP R 401:**

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

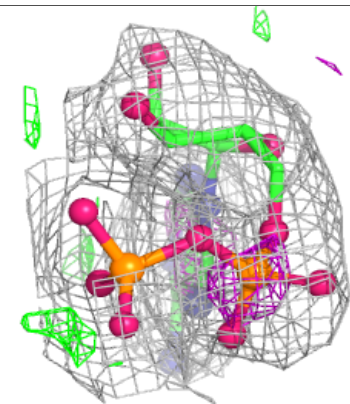
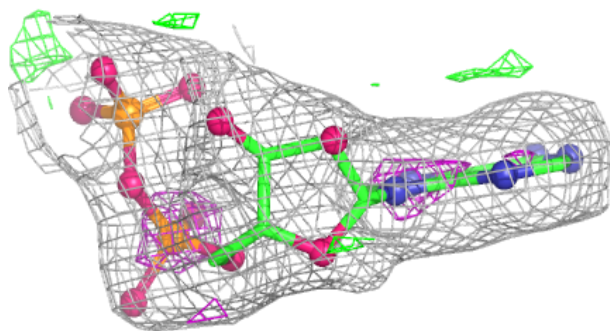
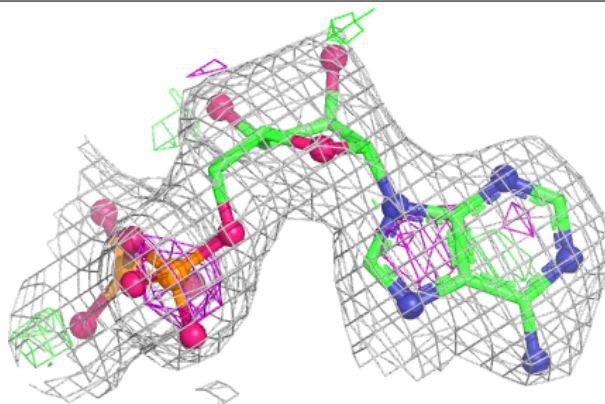


Electron density around NAP N 400:

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

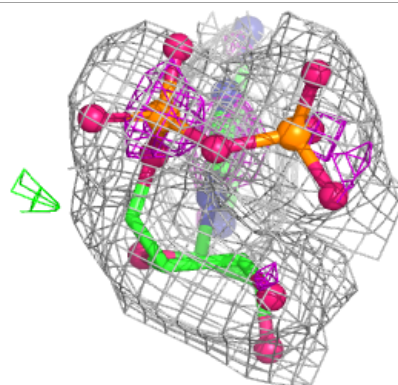
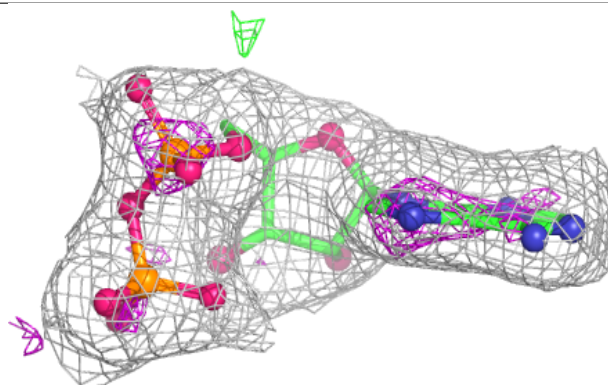
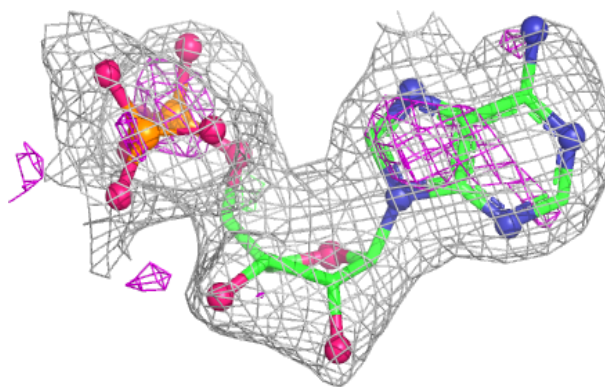
**Electron density around ADP L 401:**

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

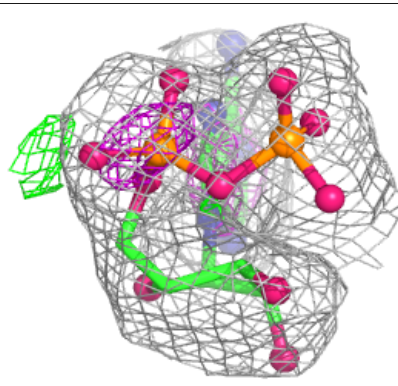
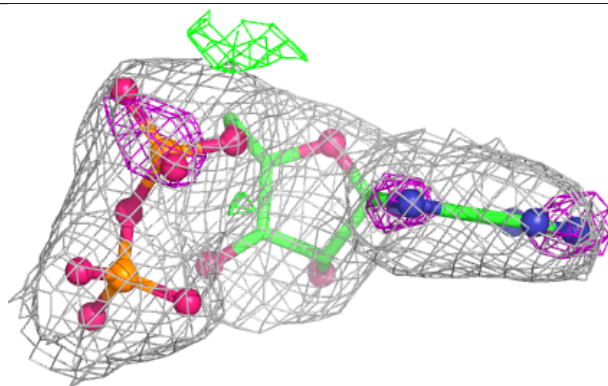
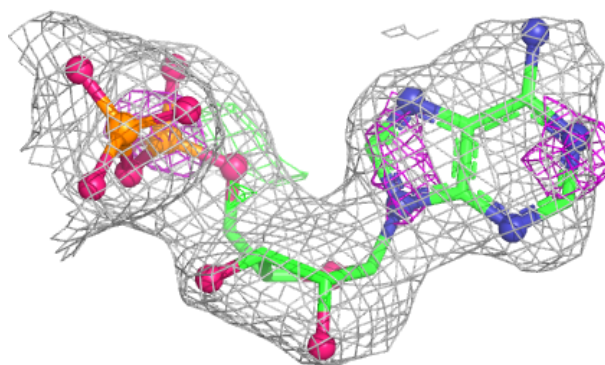


Electron density around ADP P 401:

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

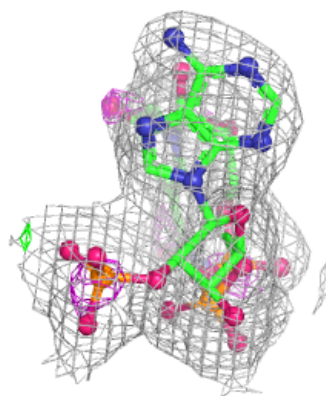
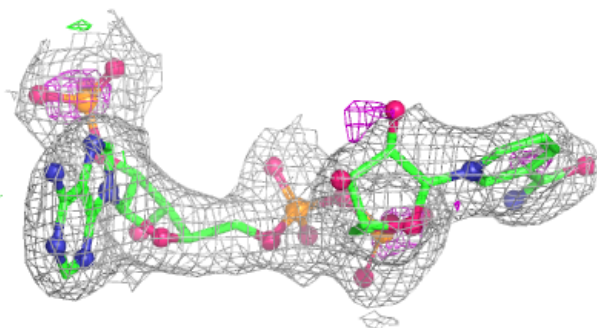
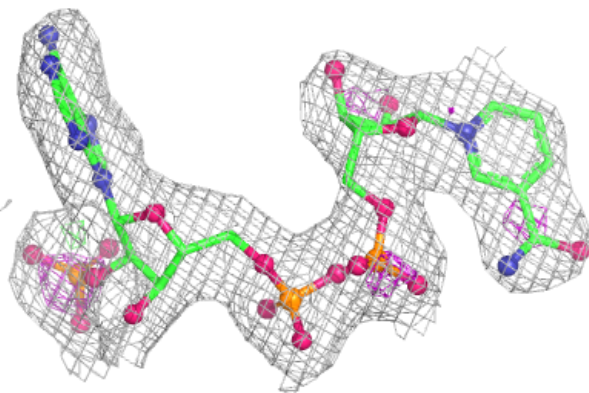
**Electron density around ADP M 401:**

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

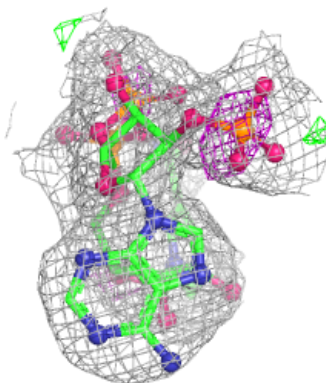
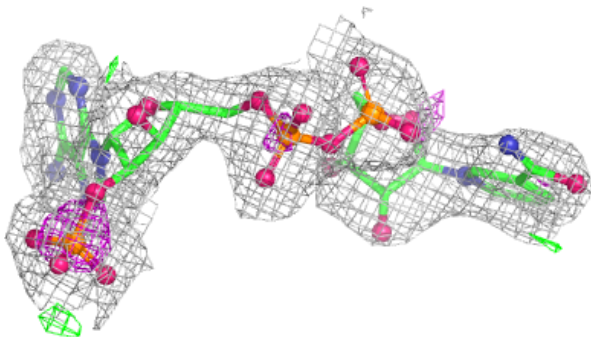
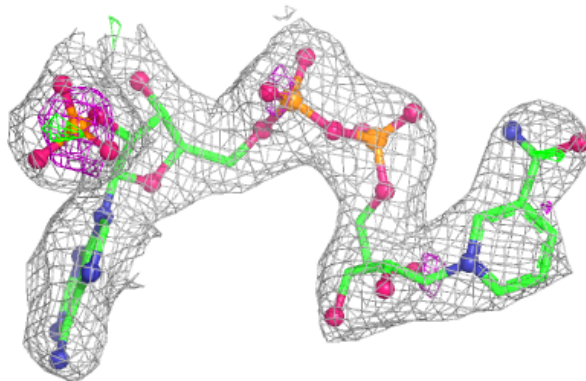


Electron density around NAP K 400:

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

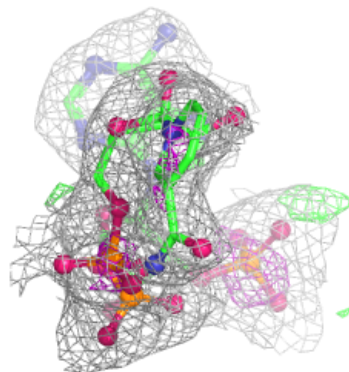
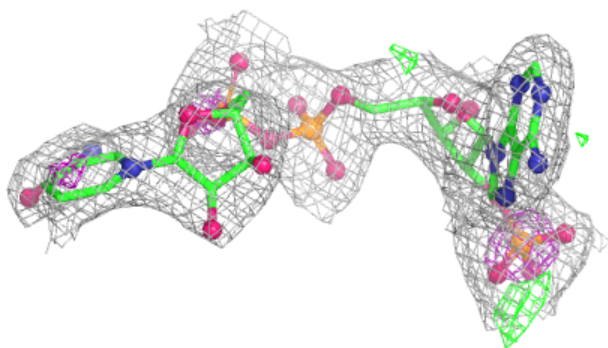
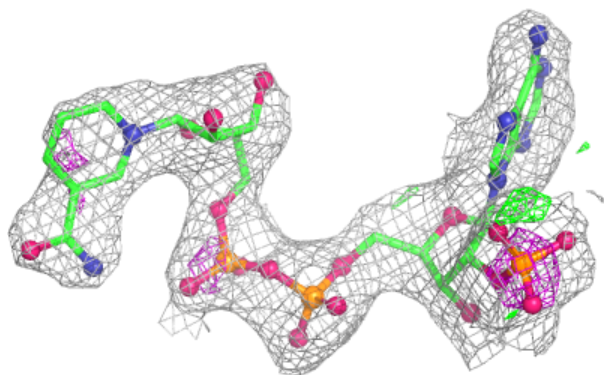
**Electron density around NAP L 400:**

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

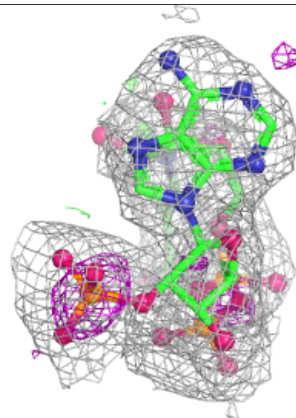
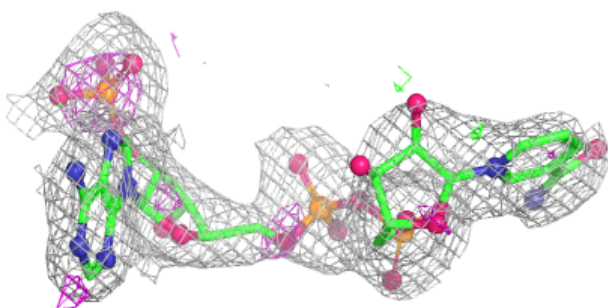
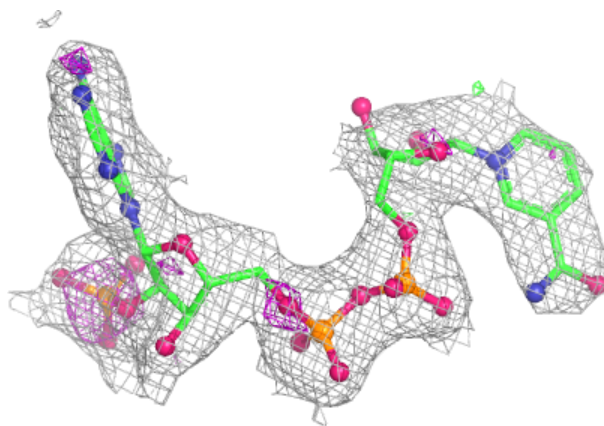


Electron density around NAP M 400:

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

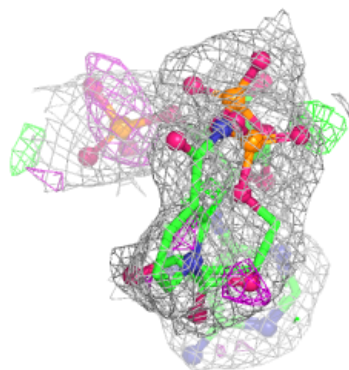
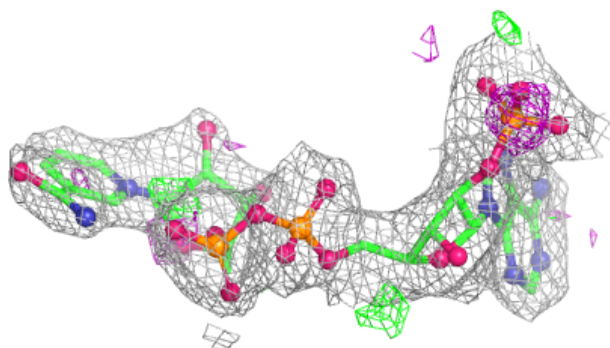
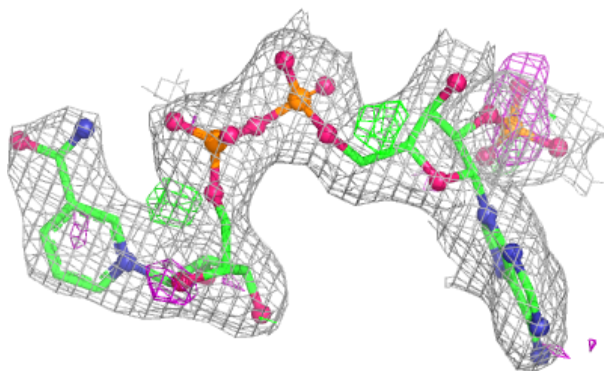
**Electron density around NAP D 400:**

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

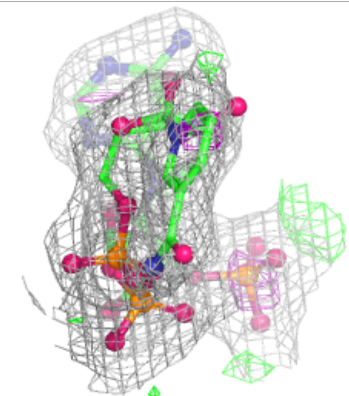
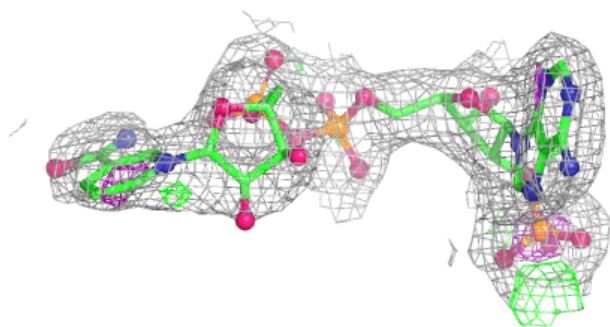
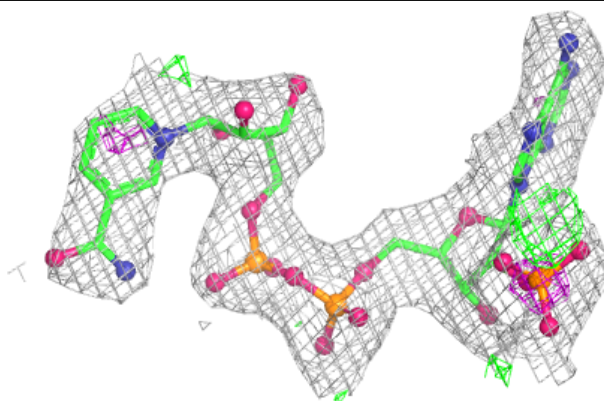


Electron density around NAP E 400:

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

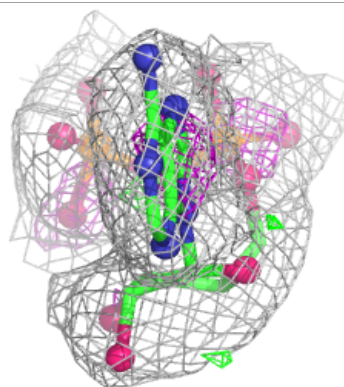
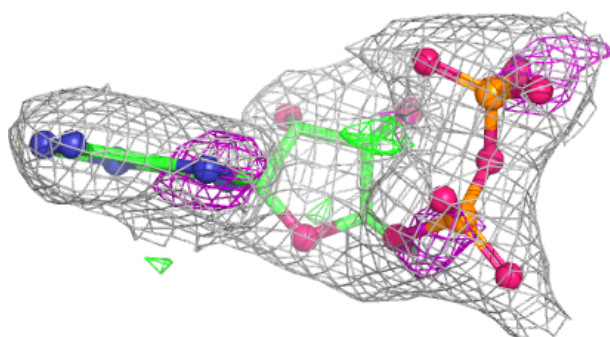
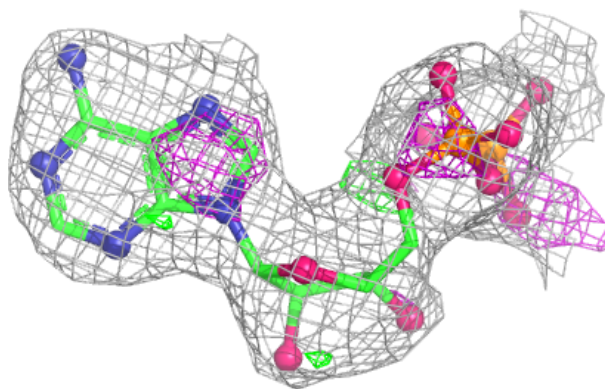
**Electron density around NAP P 400:**

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

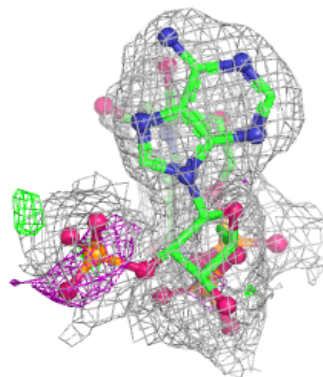
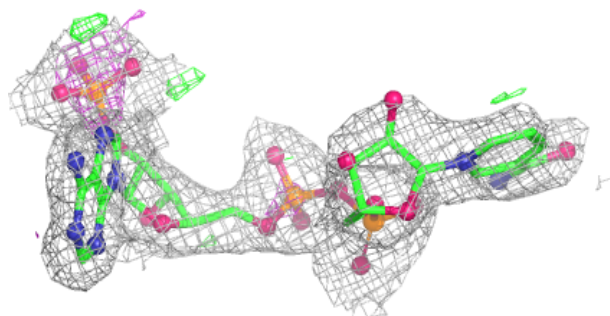
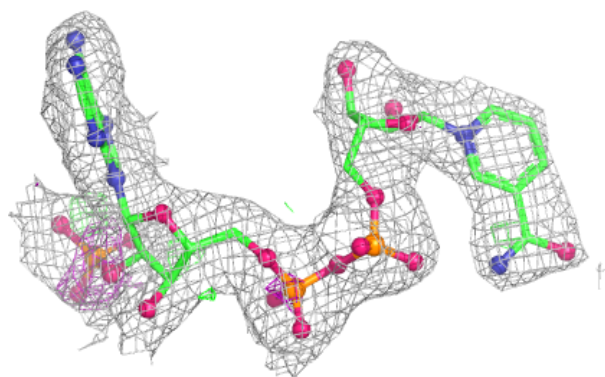


Electron density around ADP J 401:

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

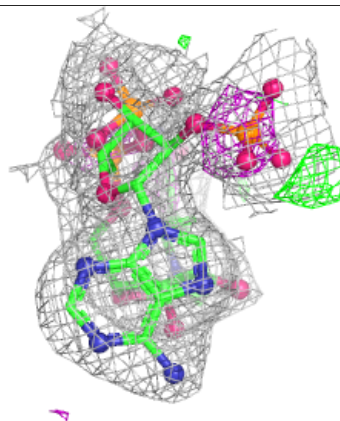
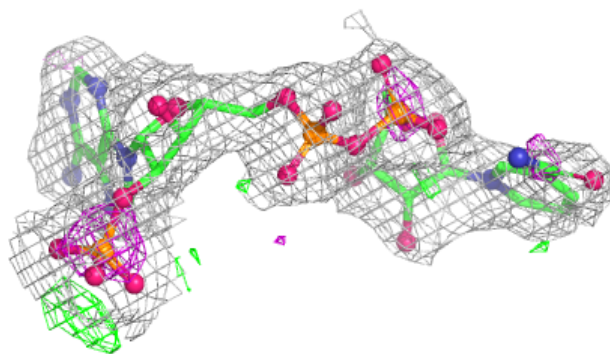
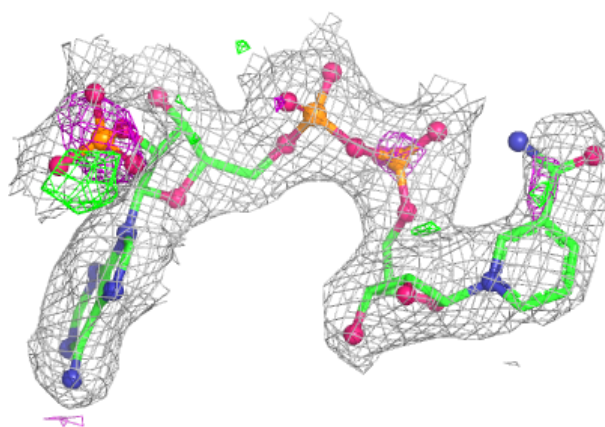
**Electron density around NAP Q 400:**

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

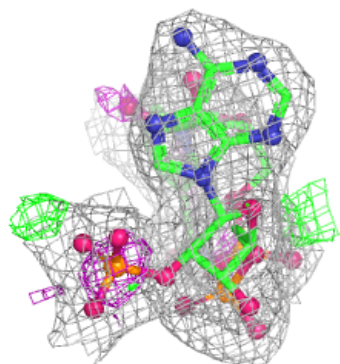
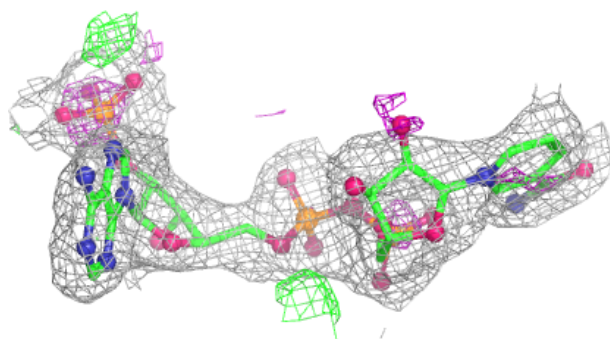
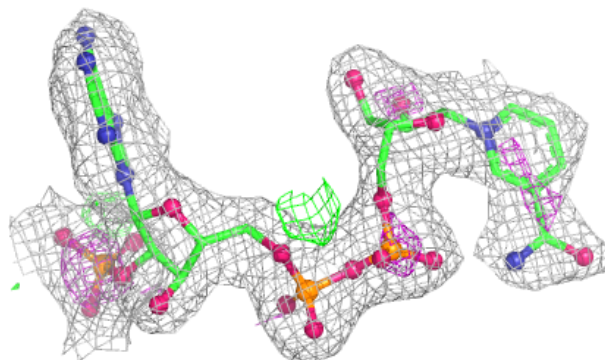


Electron density around NAP A 400:

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

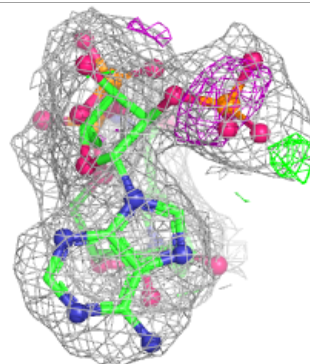
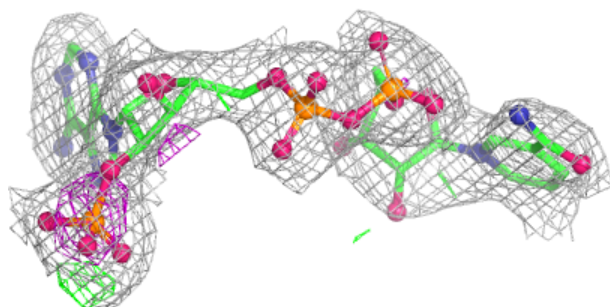
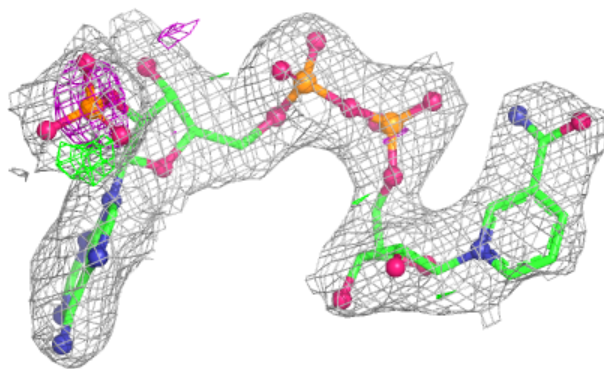
**Electron density around NAP H 400:**

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

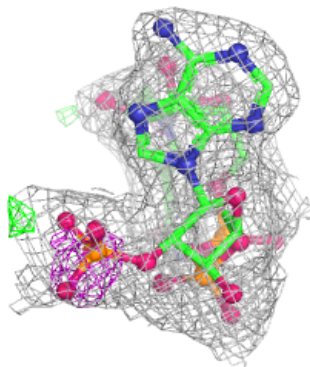
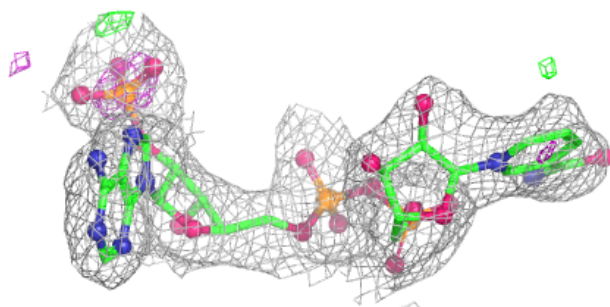
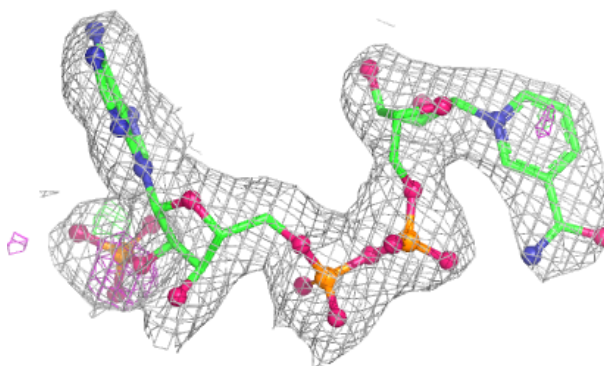


Electron density around NAP I 400:

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

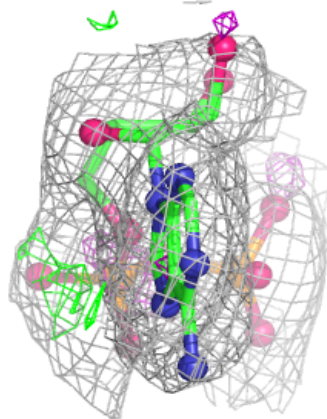
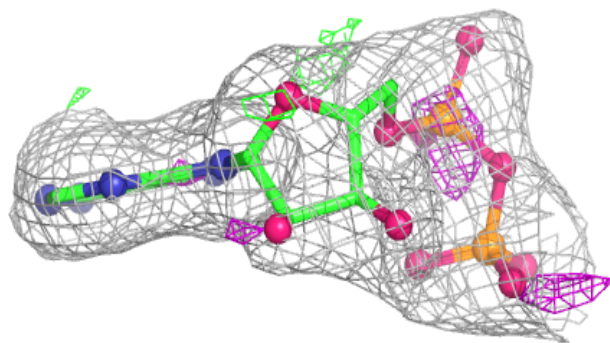
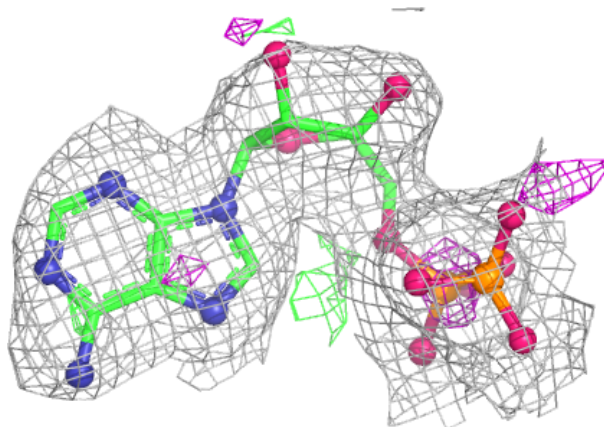
**Electron density around NAP J 400:**

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

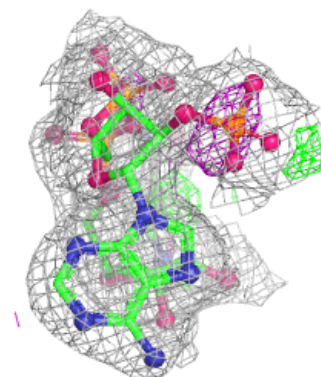
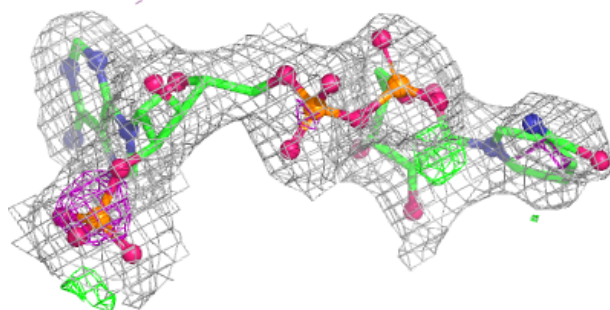
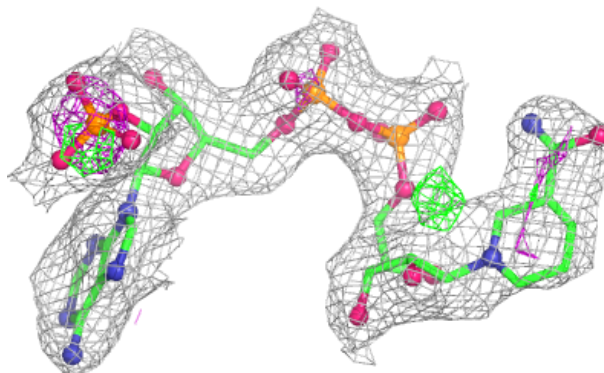


Electron density around ADP E 401:

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

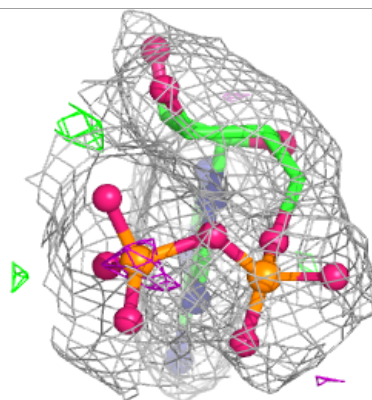
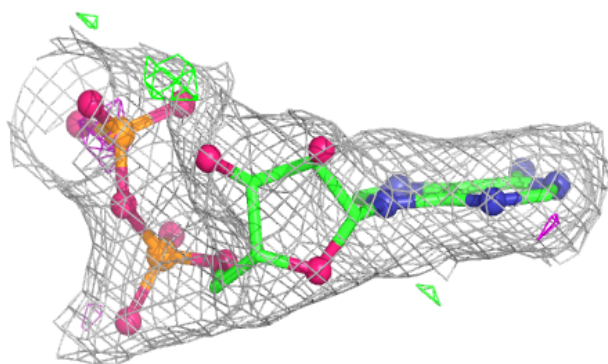
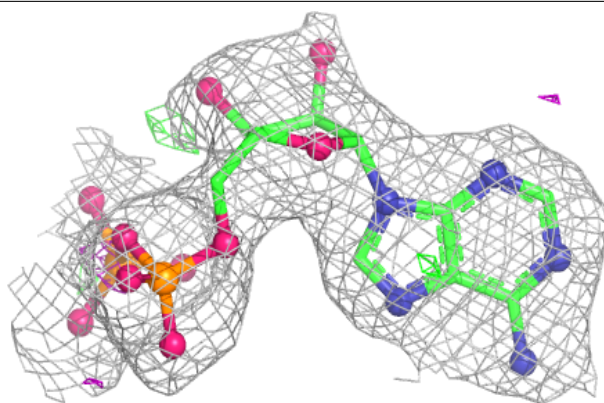
**Electron density around NAP F 400:**

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



Electron density around ADP I 401:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.