



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 19, 2024 – 11:04 PM EST

PDB ID : 4LNI
Title : B. subtilis glutamine synthetase structures reveal large active site conformational changes and basis for isoenzyme specific regulation: structure of the transition state complex
Authors : Schumacher, M.A.; Chinnam, N.; Tonthat, N.; Fisher, S.; Wray, L.
Deposited on : 2013-07-11
Resolution : 2.58 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

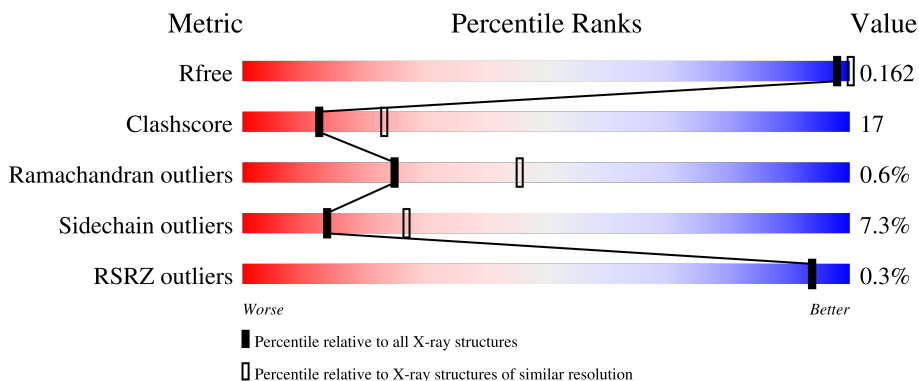
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.58 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	3676 (2.60-2.56)
Clashscore	141614	4049 (2.60-2.56)
Ramachandran outliers	138981	3979 (2.60-2.56)
Sidechain outliers	138945	3979 (2.60-2.56)
RSRZ outliers	127900	3614 (2.60-2.56)

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	443	 67% 28% 5%
1	B	443	 67% 27% 6%
1	C	443	 68% 27% . .
1	D	443	 65% 29% 5% .

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Mol	Chain	Length	Quality of chain
1	E	443	 68% 27% 5%
1	F	443	 63% 33% .
1	G	443	 % 65% 31% .
1	H	443	 67% 27% . .
1	I	443	 67% 28% 5% .
1	J	443	 67% 28% .
1	K	443	 70% 26% .
1	L	443	 67% 27% 6%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	P3S	D	501	-	-	X	-
2	P3S	F	601	-	-	X	-
2	P3S	G	601	-	-	X	-

2 Entry composition

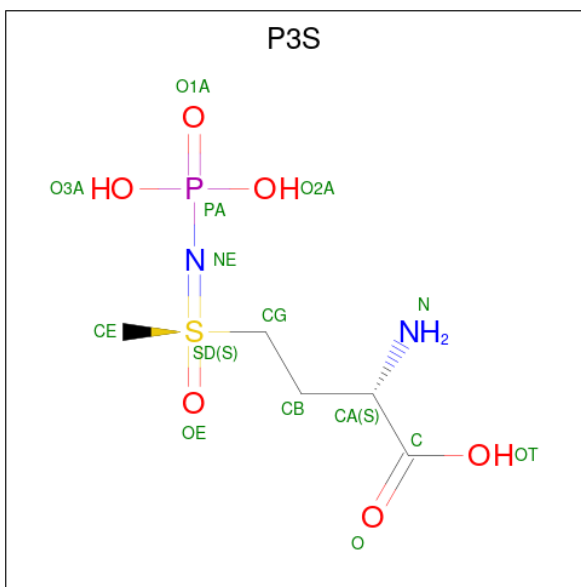
There are 5 unique types of molecules in this entry. The entry contains 86960 atoms, of which 41847 are hydrogens and 0 are deuteriums.

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 Glutamine synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
1	A	443	7012	2259	3477	590	670	16	0	0	0
1	B	443	7012	2259	3477	590	670	16	0	0	0
1	C	443	7012	2259	3477	590	670	16	0	0	0
1	D	443	7012	2259	3477	590	670	16	0	0	0
1	E	443	7012	2259	3477	590	670	16	0	0	0
1	F	443	7012	2259	3477	590	670	16	0	0	0
1	G	443	7009	2259	3474	590	670	16	0	0	0
1	H	443	7012	2259	3477	590	670	16	0	0	0
1	I	443	7012	2259	3477	590	670	16	0	0	0
1	J	441	6980	2250	3459	587	668	16	0	0	0
1	K	443	7012	2259	3477	590	670	16	0	0	0
1	L	443	7012	2259	3477	590	670	16	0	0	0

- Molecule 2 is L-METHIONINE-S-SULFOXIMINE PHOSPHATE (three-letter code: P3S) (formula: C₅H₁₃N₂O₆PS).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
			Total	C	H	N	O	P	S		
2	A	1	Total	C	H	N	O	P	S	0	0
			27	5	12	2	6	1	1		
2	B	1	Total	C	H	N	O	P	S	0	0
			27	5	12	2	6	1	1		
2	C	1	Total	C	H	N	O	P	S	0	0
			27	5	12	2	6	1	1		
2	D	1	Total	C	H	N	O	P	S	0	0
			27	5	12	2	6	1	1		
2	E	1	Total	C	H	N	O	P	S	0	0
			27	5	12	2	6	1	1		
2	F	1	Total	C	H	N	O	P	S	0	0
			27	5	12	2	6	1	1		
2	G	1	Total	C	H	N	O	P	S	0	0
			27	5	12	2	6	1	1		
2	H	1	Total	C	H	N	O	P	S	0	0
			27	5	12	2	6	1	1		
2	I	1	Total	C	H	N	O	P	S	0	0
			27	5	12	2	6	1	1		
2	J	1	Total	C	H	N	O	P	S	0	0
			27	5	12	2	6	1	1		
2	K	1	Total	C	H	N	O	P	S	0	0
			27	5	12	2	6	1	1		
2	L	1	Total	C	H	N	O	P	S	0	0
			27	5	12	2	6	1	1		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
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		
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		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	3	Total Mg 3 3	0	0
4	B	3	Total Mg 3 3	0	0
4	C	3	Total Mg 3 3	0	0
4	D	3	Total Mg 3 3	0	0
4	E	3	Total Mg 3 3	0	0
4	F	3	Total Mg 3 3	0	0
4	G	3	Total Mg 3 3	0	0
4	H	3	Total Mg 3 3	0	0
4	I	3	Total Mg 3 3	0	0
4	J	3	Total Mg 3 3	0	0
4	K	3	Total Mg 3 3	0	0
4	L	3	Total Mg 3 3	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	177	Total O 177 177	0	0
5	B	178	Total O 178 178	0	0
5	C	183	Total O 183 183	0	0
5	D	200	Total O 200 200	0	0
5	E	168	Total O 168 168	0	0
5	F	166	Total O 166 166	0	0
5	G	184	Total O 184 184	0	0
5	H	174	Total O 174 174	0	0

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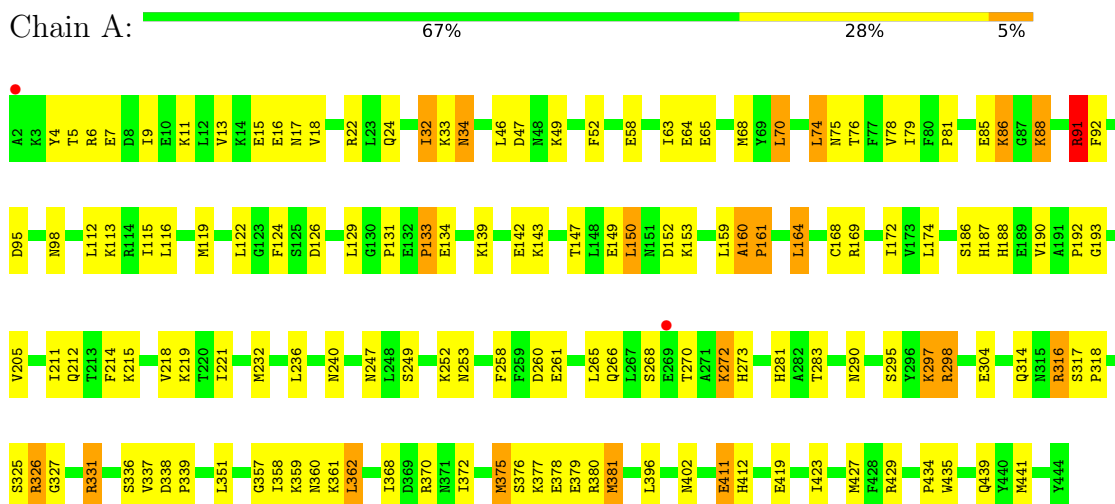
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	I	189	Total 189	O 189	0	0
5	J	194	Total 194	O 194	0	0
5	K	191	Total 191	O 191	0	0
5	L	163	Total 163	O 163	0	0

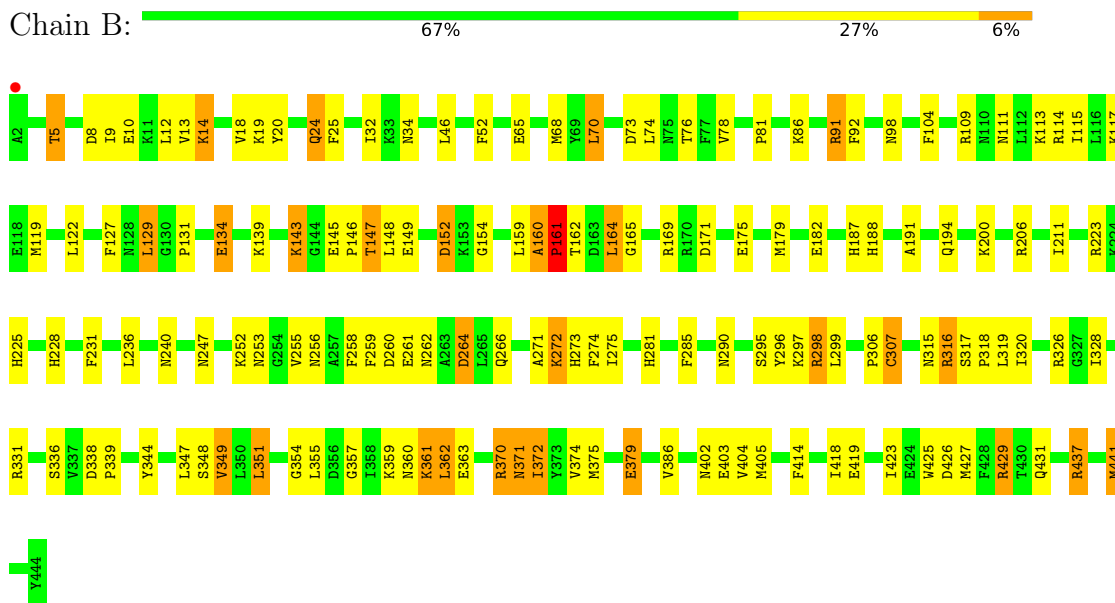
3 Residue-property plots [i](#)

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.

- Molecule 1: Glutamine synthetase

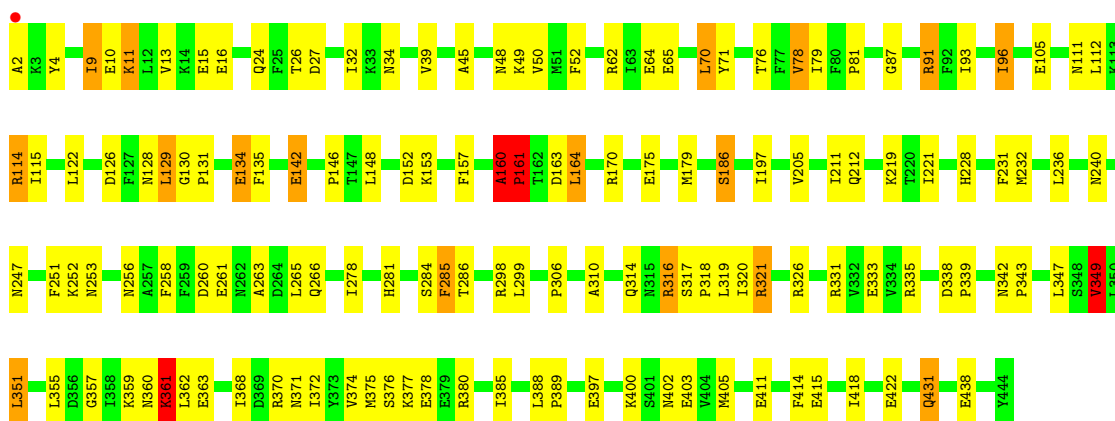


- Molecule 1: Glutamine synthetase



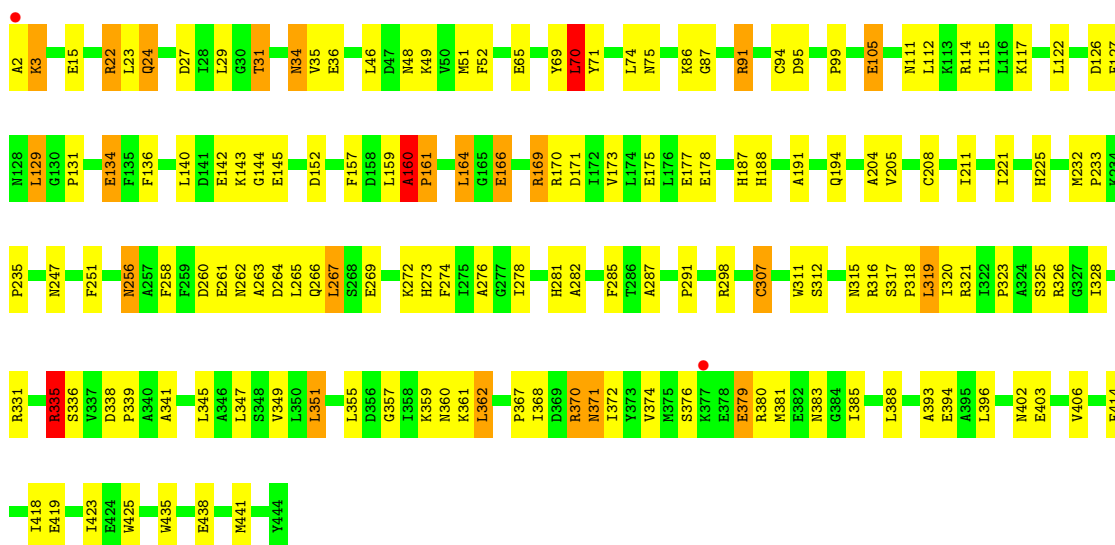
- Molecule 1: Glutamine synthetase

Chain C:  68% 27% 5% 0%



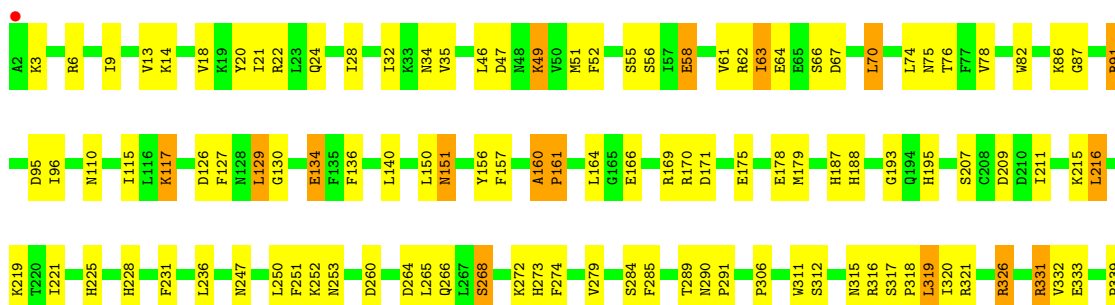
• Molecule 1: Glutamine synthetase

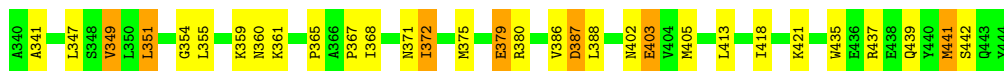
Chain D:  65% 29% 5% 1%



• Molecule 1: Glutamine synthetase

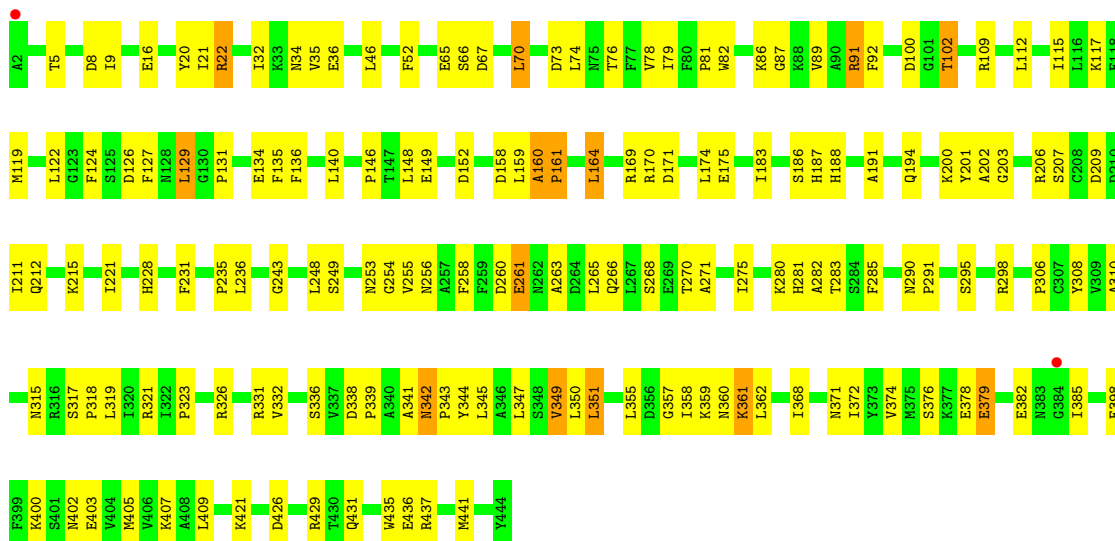
Chain E:  68% 27% 5% 0%





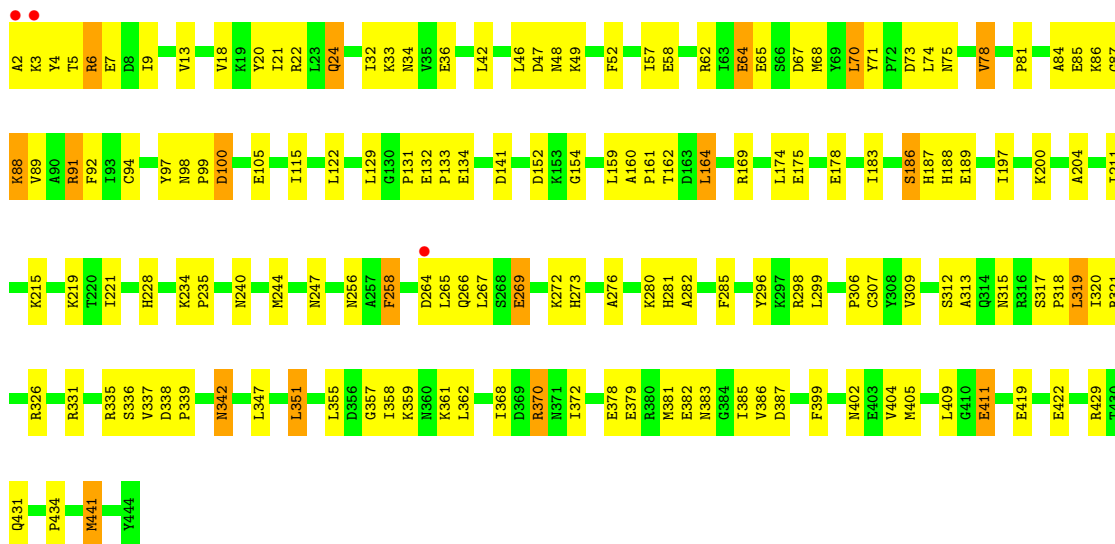
- Molecule 1: Glutamine synthetase

Chain F: 63% 33% .



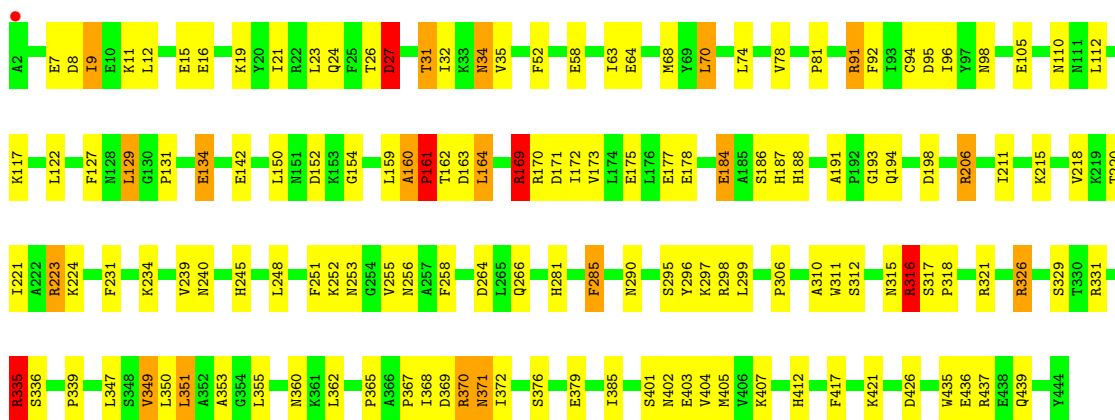
- Molecule 1: Glutamine synthetase

Chain G: % 65% 31% .



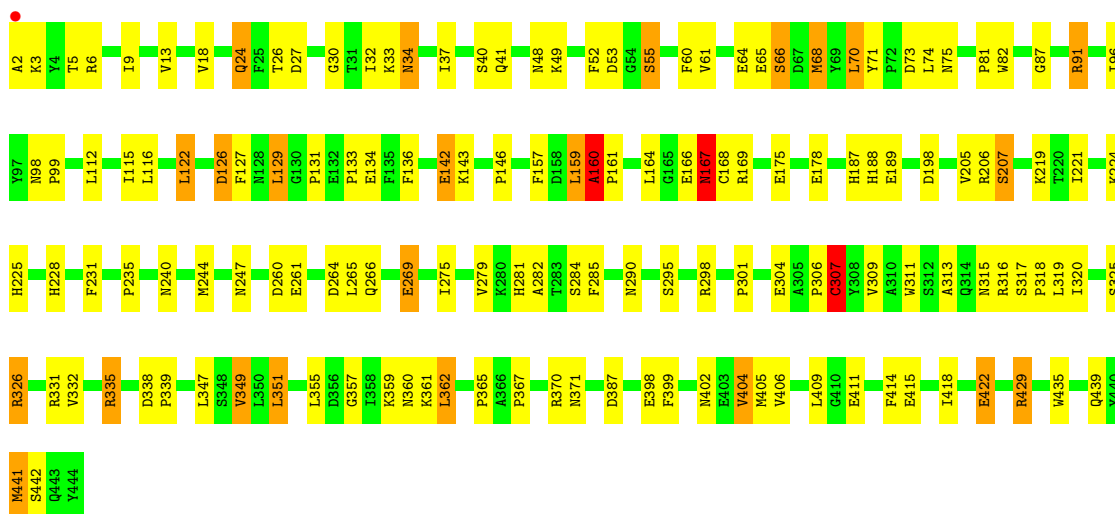
- Molecule 1: Glutamine synthetase

Chain H: 67% 27% . .



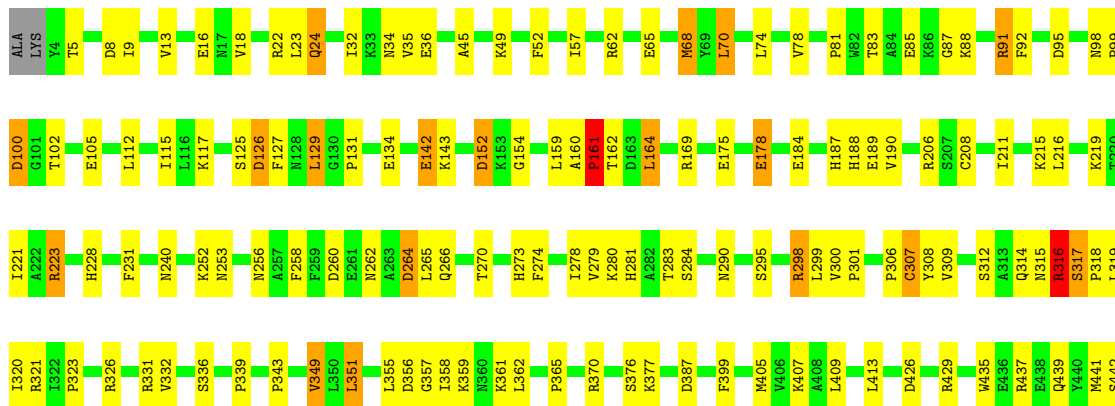
● Molecule 1: Glutamine synthetase

Chain I: 67% 28% 5%



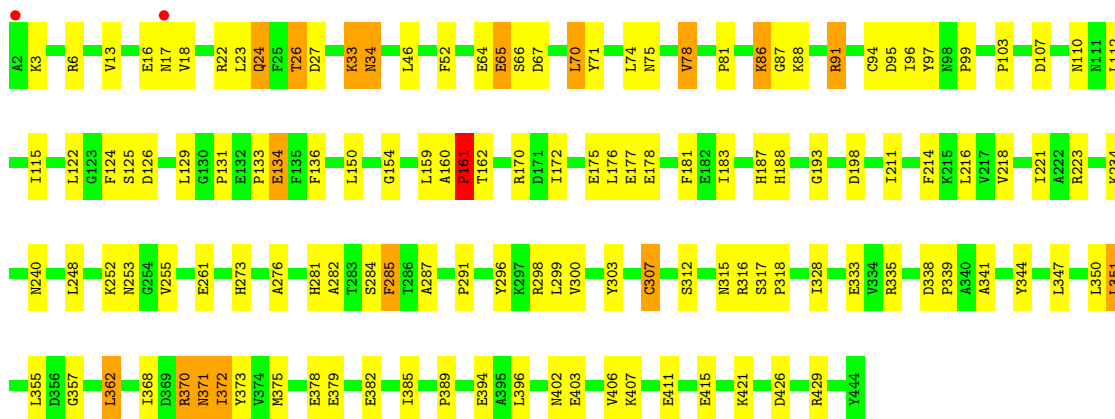
● Molecule 1: Glutamine synthetase

Chain J: 67% 28%

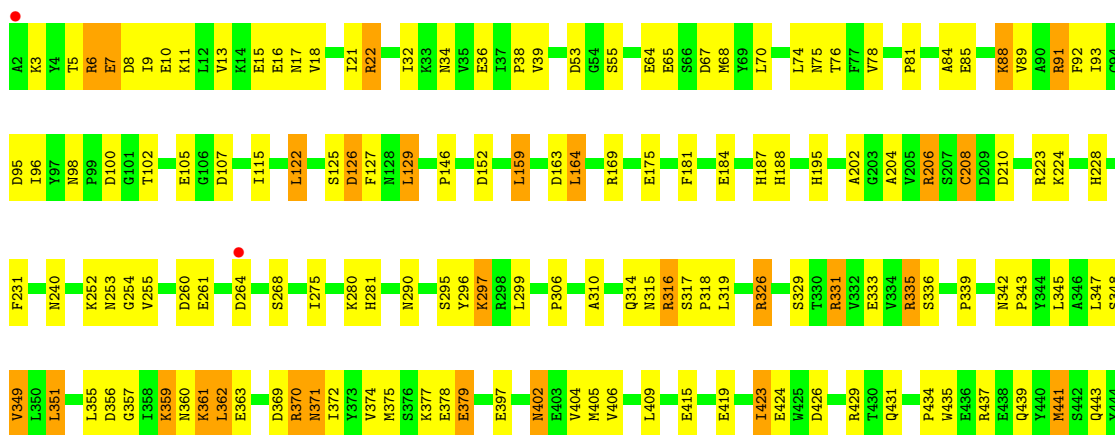


0443
Y444

- Molecule 1: Glutamine synthetase

Chain K:  70% 26%

- Molecule 1: Glutamine synthetase

Chain L:  67% 27% 6%

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	110.20Å 141.60Å 142.10Å 60.29° 67.38° 76.20°	Depositor
Resolution (Å)	117.10 – 2.58 117.10 – 2.58	Depositor EDS
% Data completeness (in resolution range)	90.8 (117.10-2.58) 90.8 (117.10-2.58)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.14 (at 2.58Å)	Xtrriage
Refinement program	PHENIX 1.6.4_486, CNS 1.2	Depositor
R, R_{free}	0.165 , 0.223 0.158 , 0.162	Depositor DCC
R_{free} test set	21985 reflections (11.18%)	wwPDB-VP
Wilson B-factor (Å ²)	19.1	Xtrriage
Anisotropy	1.014	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	86960	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.84% of the height of the origin peak. No significant pseudotranslation is detected.*

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: P3S, MG, ADP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.96	0/3618	0.95	6/4895 (0.1%)
1	B	0.98	3/3618 (0.1%)	0.97	10/4895 (0.2%)
1	C	0.98	4/3618 (0.1%)	1.02	10/4895 (0.2%)
1	D	0.97	4/3618 (0.1%)	0.96	9/4895 (0.2%)
1	E	0.97	3/3618 (0.1%)	1.03	14/4895 (0.3%)
1	F	0.95	1/3618 (0.0%)	0.96	5/4895 (0.1%)
1	G	0.95	3/3618 (0.1%)	0.92	6/4895 (0.1%)
1	H	0.97	2/3618 (0.1%)	0.99	9/4895 (0.2%)
1	I	0.96	2/3618 (0.1%)	0.99	11/4895 (0.2%)
1	J	0.98	6/3604 (0.2%)	1.01	10/4877 (0.2%)
1	K	1.01	2/3618 (0.1%)	1.00	8/4895 (0.2%)
1	L	0.96	2/3618 (0.1%)	0.89	2/4895 (0.0%)
All	All	0.97	32/43402 (0.1%)	0.98	100/58722 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	2
1	D	0	1
1	E	0	1
1	F	0	1
1	H	0	1
1	I	0	1
All	All	0	9

The worst 5 of 32 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	161	PRO	N-CD	7.91	1.58	1.47
1	D	166	GLU	CG-CD	7.69	1.63	1.51
1	C	134	GLU	CG-CD	7.57	1.63	1.51
1	I	307	CYS	CB-SG	-6.89	1.70	1.82
1	J	105	GLU	CG-CD	6.76	1.62	1.51

The worst 5 of 100 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	160	ALA	C-N-CD	-26.90	61.42	120.60
1	K	160	ALA	C-N-CD	-23.72	68.41	120.60
1	C	160	ALA	C-N-CD	-22.46	71.20	120.60
1	H	160	ALA	C-N-CD	-22.29	71.56	120.60
1	J	160	ALA	C-N-CD	-22.20	71.76	120.60

There are no chirality outliers.

5 of 9 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	160	ALA	Peptide
1	B	160	ALA	Peptide
1	C	160	ALA	Peptide,Mainchain
1	D	160	ALA	Peptide
1	E	160	ALA	Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3535	3477	3466	118	0
1	B	3535	3477	3466	113	0
1	C	3535	3477	3466	114	0
1	D	3535	3477	3466	128	0
1	E	3535	3477	3466	117	0
1	F	3535	3477	3466	140	0
1	G	3535	3474	3466	141	0
1	H	3535	3477	3466	126	0
1	I	3535	3477	3466	129	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	J	3521	3459	3448	115	0
1	K	3535	3477	3466	111	0
1	L	3535	3477	3466	129	0
2	A	15	12	10	4	0
2	B	15	12	10	1	0
2	C	15	12	10	4	0
2	D	15	12	10	6	0
2	E	15	12	10	1	0
2	F	15	12	10	6	0
2	G	15	12	10	8	0
2	H	15	12	10	5	0
2	I	15	12	10	2	0
2	J	15	12	10	3	0
2	K	15	12	10	2	0
2	L	15	12	10	1	0
3	A	27	0	12	0	0
3	B	27	0	12	0	0
3	C	27	0	12	0	0
3	D	27	0	12	0	0
3	E	27	0	12	1	0
3	F	27	0	12	1	0
3	G	27	0	12	0	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
4	A	3	0	0	0	0
4	B	3	0	0	0	0
4	C	3	0	0	0	0
4	D	3	0	0	0	0
4	E	3	0	0	0	0
4	F	3	0	0	0	0
4	G	3	0	0	0	0
4	H	3	0	0	0	0
4	I	3	0	0	0	0
4	J	3	0	0	0	0
4	K	3	0	0	0	0
4	L	3	0	0	0	0
5	A	177	0	0	11	0
5	B	178	0	0	14	0
5	C	183	0	0	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	200	0	0	13	0
5	E	168	0	0	11	0
5	F	166	0	0	15	0
5	G	184	0	0	19	0
5	H	174	0	0	17	0
5	I	189	0	0	16	0
5	J	194	0	0	20	0
5	K	191	0	0	19	0
5	L	163	0	0	11	0
All	All	45113	41847	41838	1414	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 17.

The worst 5 of 1414 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)
2:H:601:P3S:SD	2:H:601:P3S:CE	2.01	1.48
2:G:601:P3S:SD	2:G:601:P3S:CE	2.01	1.46
2:F:601:P3S:SD	2:F:601:P3S:CE	2.02	1.45
2:D:501:P3S:CE	2:D:501:P3S:SD	2.03	1.43
1:I:357:GLY:HA2	1:I:362:LEU:CD1	1.64	1.27

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	441/443 (100%)	407 (92%)	32 (7%)	2 (0%)	29 50
1	B	441/443 (100%)	417 (95%)	22 (5%)	2 (0%)	29 50

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	441/443 (100%)	414 (94%)	22 (5%)	5 (1%)	14	28
1	D	441/443 (100%)	407 (92%)	31 (7%)	3 (1%)	22	41
1	E	441/443 (100%)	417 (95%)	21 (5%)	3 (1%)	22	41
1	F	441/443 (100%)	410 (93%)	28 (6%)	3 (1%)	22	41
1	G	441/443 (100%)	403 (91%)	34 (8%)	4 (1%)	17	34
1	H	441/443 (100%)	420 (95%)	20 (4%)	1 (0%)	47	69
1	I	441/443 (100%)	414 (94%)	24 (5%)	3 (1%)	22	41
1	J	439/443 (99%)	413 (94%)	23 (5%)	3 (1%)	22	41
1	K	441/443 (100%)	416 (94%)	23 (5%)	2 (0%)	29	50
1	L	441/443 (100%)	416 (94%)	24 (5%)	1 (0%)	47	69
All	All	5290/5316 (100%)	4954 (94%)	304 (6%)	32 (1%)	25	45

5 of 32 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	161	PRO
1	B	161	PRO
1	C	161	PRO
1	C	361	LYS
1	D	161	PRO

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	382/382 (100%)	353 (92%)	29 (8%)	13	25
1	B	382/382 (100%)	354 (93%)	28 (7%)	14	27
1	C	382/382 (100%)	358 (94%)	24 (6%)	18	35
1	D	382/382 (100%)	358 (94%)	24 (6%)	18	35
1	E	382/382 (100%)	357 (94%)	25 (6%)	17	33
1	F	382/382 (100%)	363 (95%)	19 (5%)	24	45

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	382/382 (100%)	354 (93%)	28 (7%)	14	27
1	H	382/382 (100%)	350 (92%)	32 (8%)	11	20
1	I	382/382 (100%)	349 (91%)	33 (9%)	10	19
1	J	381/382 (100%)	349 (92%)	32 (8%)	11	20
1	K	382/382 (100%)	359 (94%)	23 (6%)	19	37
1	L	382/382 (100%)	346 (91%)	36 (9%)	8	16
All	All	4583/4584 (100%)	4250 (93%)	333 (7%)	14	27

5 of 333 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	335	ARG
1	K	112	LEU
1	I	411	GLU
1	J	240	ASN
1	L	17	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 148 such sidechains are listed below:

Mol	Chain	Res	Type
1	J	240	ASN
1	L	314	GLN
1	J	412	HIS
1	K	314	GLN
1	E	24	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 60 ligands modelled in this entry, 36 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
2	P3S	I	501	4	11,14,14	2.40	4 (36%)	12,21,21	4.70	7 (58%)
3	ADP	I	502	4	24,29,29	1.15	2 (8%)	29,45,45	1.81	8 (27%)
3	ADP	F	602	4	24,29,29	1.15	3 (12%)	29,45,45	2.20	10 (34%)
3	ADP	L	502	4	24,29,29	1.13	2 (8%)	29,45,45	1.61	7 (24%)
2	P3S	C	501	4	11,14,14	2.33	4 (36%)	12,21,21	8.06	7 (58%)
2	P3S	L	501	4	11,14,14	2.45	4 (36%)	12,21,21	6.22	4 (33%)
2	P3S	H	601	4	11,14,14	2.66	6 (54%)	12,21,21	5.36	5 (41%)
2	P3S	A	501	4	11,14,14	2.57	5 (45%)	12,21,21	5.85	4 (33%)
3	ADP	D	502	4	24,29,29	1.00	1 (4%)	29,45,45	1.72	5 (17%)
2	P3S	D	501	4	11,14,14	2.50	5 (45%)	12,21,21	7.32	5 (41%)
2	P3S	G	601	4	11,14,14	2.48	4 (36%)	12,21,21	6.16	6 (50%)
2	P3S	J	501	4	11,14,14	2.24	3 (27%)	12,21,21	5.83	5 (41%)
3	ADP	B	502	4	24,29,29	0.93	0	29,45,45	1.79	7 (24%)
3	ADP	J	502	4	24,29,29	0.87	1 (4%)	29,45,45	1.61	6 (20%)
2	P3S	B	501	4	11,14,14	2.43	4 (36%)	12,21,21	5.31	9 (75%)
3	ADP	H	602	4	24,29,29	1.10	3 (12%)	29,45,45	1.76	5 (17%)
2	P3S	F	601	4	11,14,14	2.38	5 (45%)	12,21,21	7.23	5 (41%)
3	ADP	C	502	4	24,29,29	0.88	0	29,45,45	2.16	9 (31%)
3	ADP	A	502	4	24,29,29	0.96	0	29,45,45	1.62	5 (17%)
3	ADP	E	502	4	24,29,29	0.92	1 (4%)	29,45,45	1.98	5 (17%)
3	ADP	K	502	4	24,29,29	1.00	2 (8%)	29,45,45	2.17	9 (31%)
2	P3S	E	501	4	11,14,14	2.30	4 (36%)	12,21,21	5.69	8 (66%)
2	P3S	K	501	4	11,14,14	2.34	4 (36%)	12,21,21	4.84	4 (33%)
3	ADP	G	602	4	24,29,29	0.92	1 (4%)	29,45,45	1.56	8 (27%)

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	P3S	I	501	4	-	4/9/16/16	-
3	ADP	I	502	4	-	2/12/32/32	0/3/3/3
3	ADP	F	602	4	-	5/12/32/32	0/3/3/3
3	ADP	L	502	4	-	5/12/32/32	0/3/3/3
2	P3S	C	501	4	-	4/9/16/16	-
2	P3S	L	501	4	-	4/9/16/16	-
2	P3S	H	601	4	-	4/9/16/16	-
2	P3S	A	501	4	-	4/9/16/16	-
3	ADP	D	502	4	-	3/12/32/32	0/3/3/3
2	P3S	D	501	4	-	4/9/16/16	-
2	P3S	G	601	4	-	4/9/16/16	-
2	P3S	J	501	4	-	4/9/16/16	-
3	ADP	B	502	4	-	4/12/32/32	0/3/3/3
3	ADP	J	502	4	-	4/12/32/32	0/3/3/3
2	P3S	B	501	4	-	4/9/16/16	-
3	ADP	H	602	4	-	6/12/32/32	0/3/3/3
2	P3S	F	601	4	-	4/9/16/16	-
3	ADP	C	502	4	-	2/12/32/32	0/3/3/3
3	ADP	A	502	4	-	2/12/32/32	0/3/3/3
3	ADP	E	502	4	-	3/12/32/32	0/3/3/3
3	ADP	K	502	4	-	4/12/32/32	0/3/3/3
2	P3S	E	501	4	-	4/9/16/16	-
2	P3S	K	501	4	-	4/9/16/16	-
3	ADP	G	602	4	-	2/12/32/32	0/3/3/3

The worst 5 of 68 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	P3S	CE-SD	5.52	2.03	1.74
2	F	601	P3S	CE-SD	5.29	2.02	1.74
2	H	601	P3S	CE-SD	5.16	2.01	1.74
2	G	601	P3S	CE-SD	5.14	2.01	1.74
2	J	501	P3S	CE-SD	5.06	2.01	1.74

The worst 5 of 153 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	P3S	OE-SD-CG	23.26	126.57	108.37
2	D	501	P3S	OE-SD-CG	20.64	124.53	108.37
2	F	601	P3S	OE-SD-CG	20.42	124.35	108.37
2	L	501	P3S	OE-SD-CG	16.18	121.03	108.37
2	A	501	P3S	OE-SD-CE	15.58	134.05	109.24

There are no chirality outliers.

5 of 90 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	P3S	O-C-CA-N
3	B	502	ADP	PA-O3A-PB-O2B
3	D	502	ADP	PA-O3A-PB-O2B
3	E	502	ADP	PA-O3A-PB-O2B
3	F	602	ADP	C5'-O5'-PA-O3A

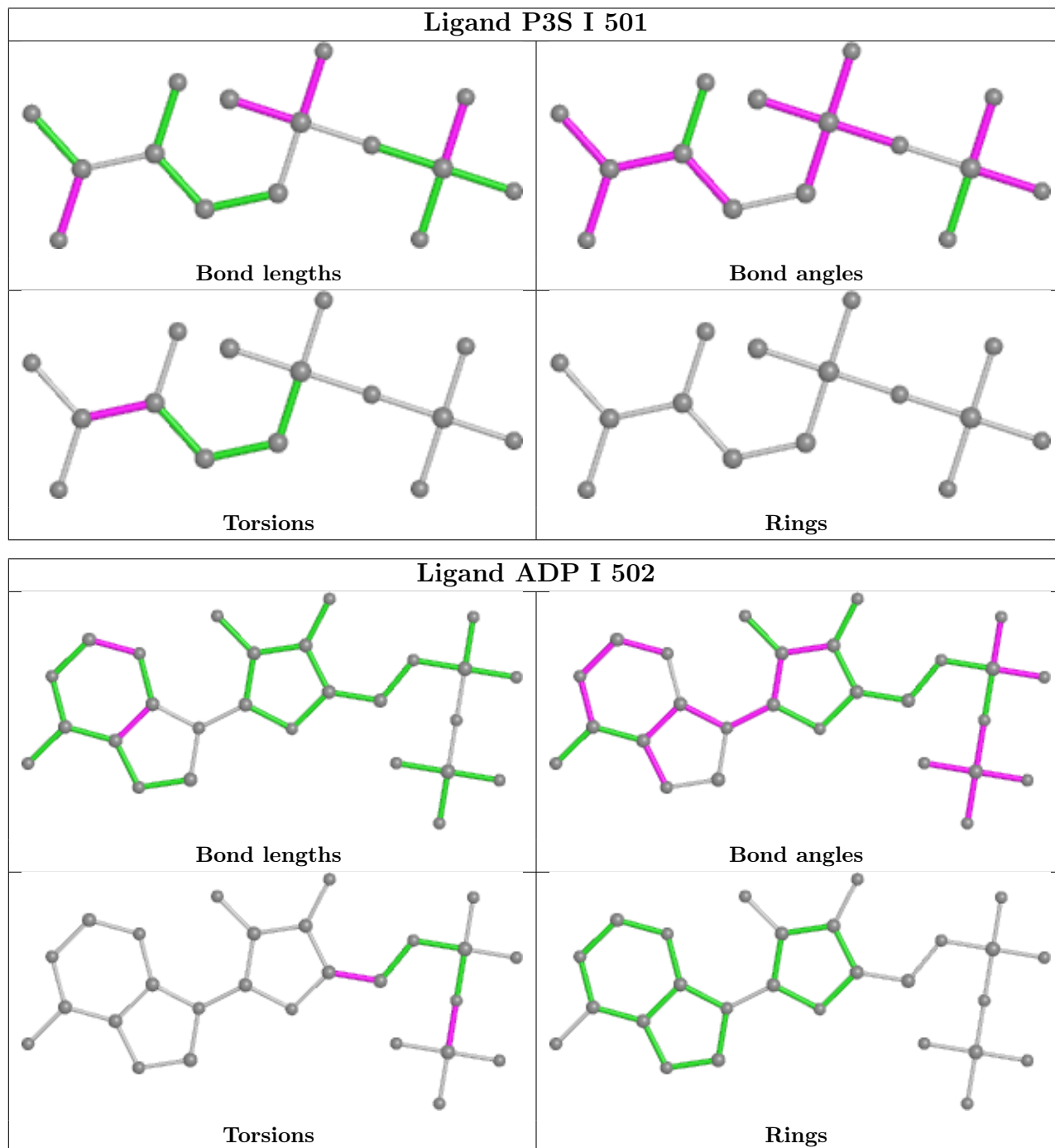
There are no ring outliers.

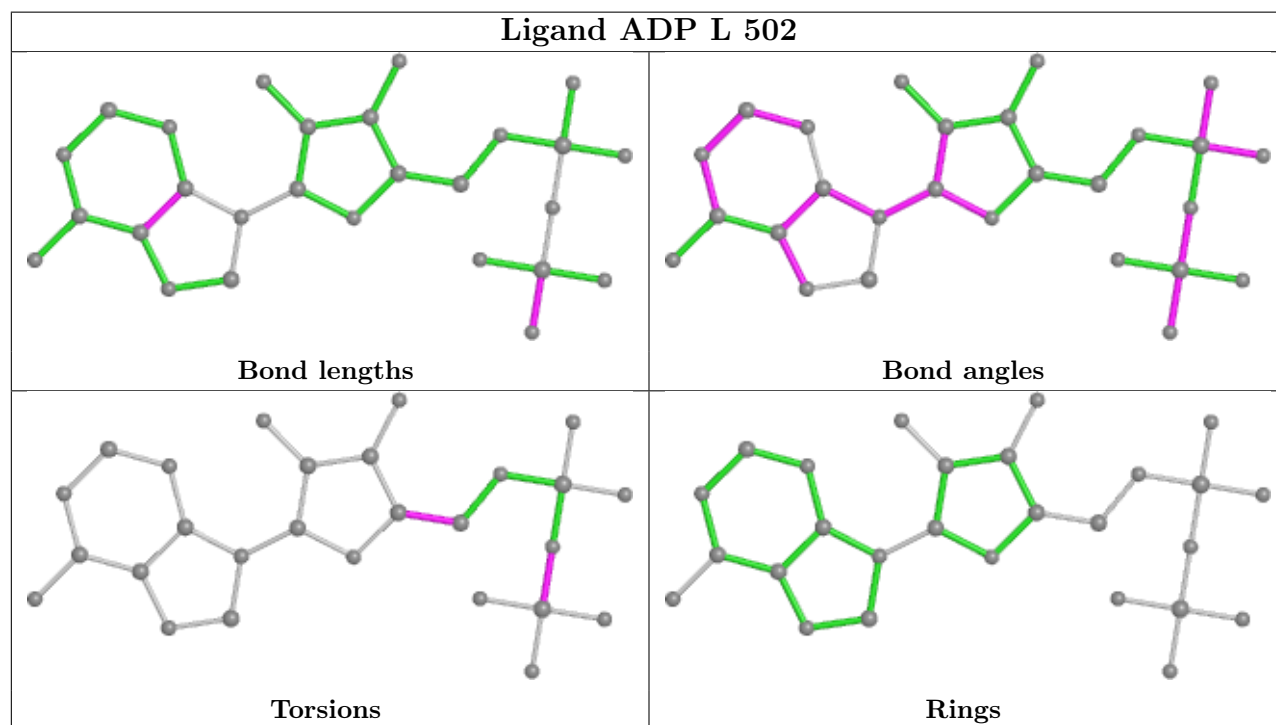
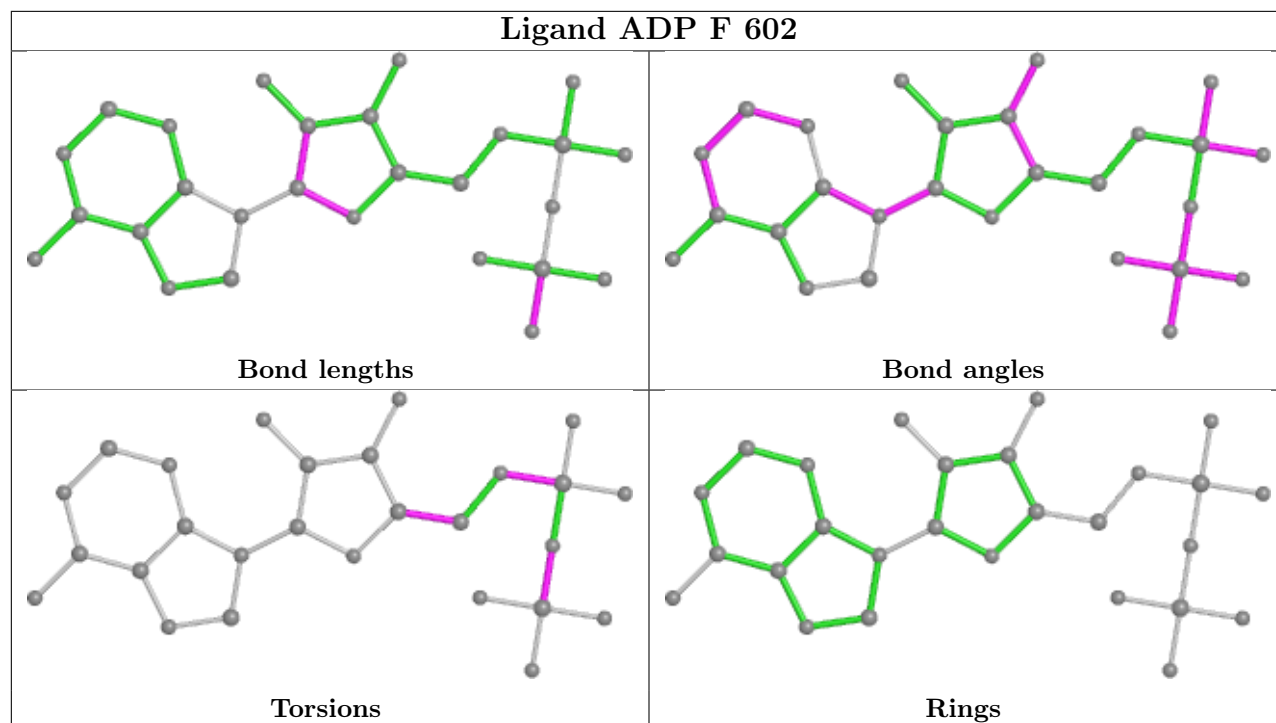
14 monomers are involved in 45 short contacts:

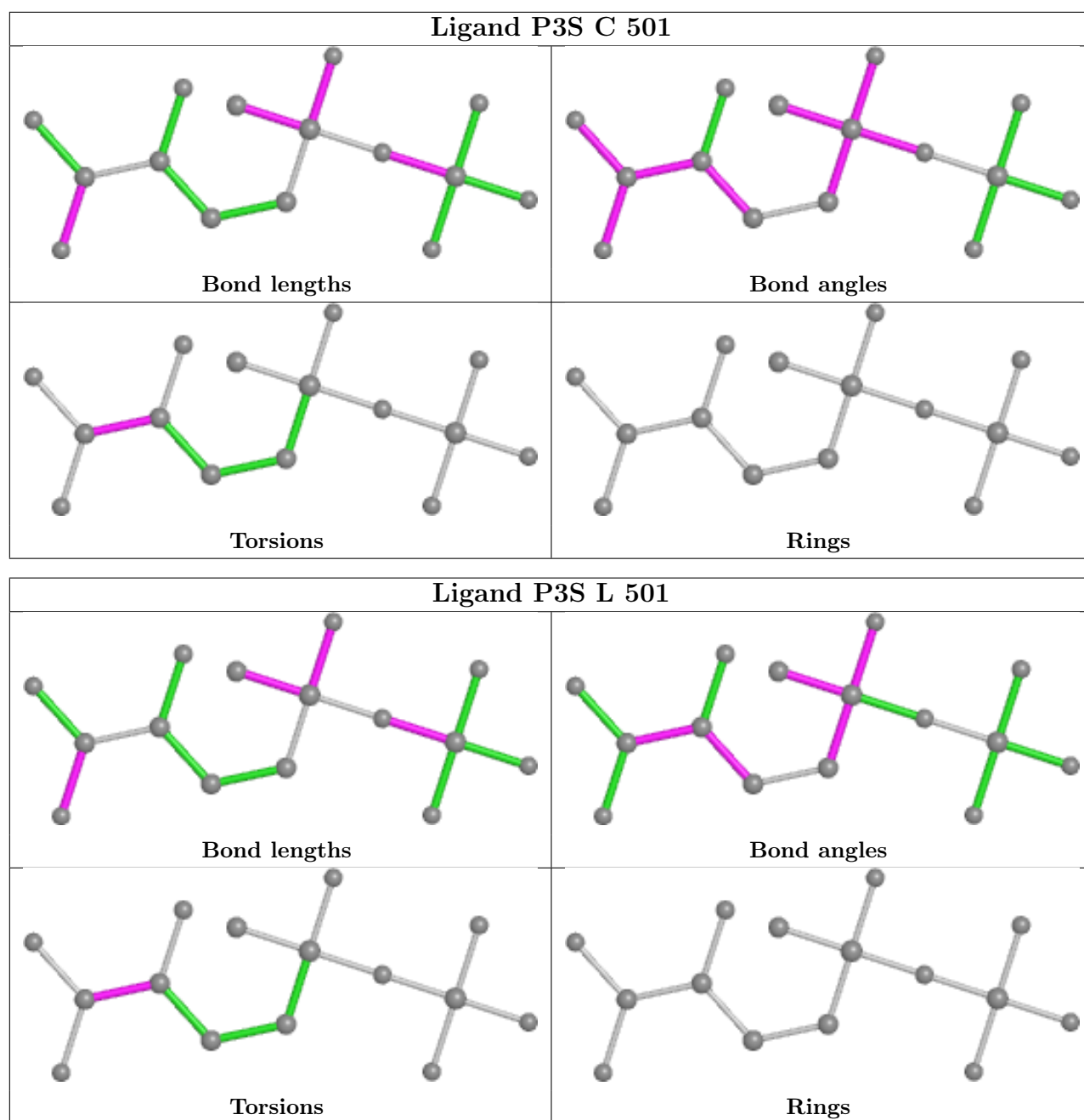
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	I	501	P3S	2	0
3	F	602	ADP	1	0
2	C	501	P3S	4	0
2	L	501	P3S	1	0
2	H	601	P3S	5	0
2	A	501	P3S	4	0
2	D	501	P3S	6	0
2	G	601	P3S	8	0
2	J	501	P3S	3	0
2	B	501	P3S	1	0
2	F	601	P3S	6	0
3	E	502	ADP	1	0
2	E	501	P3S	1	0
2	K	501	P3S	2	0

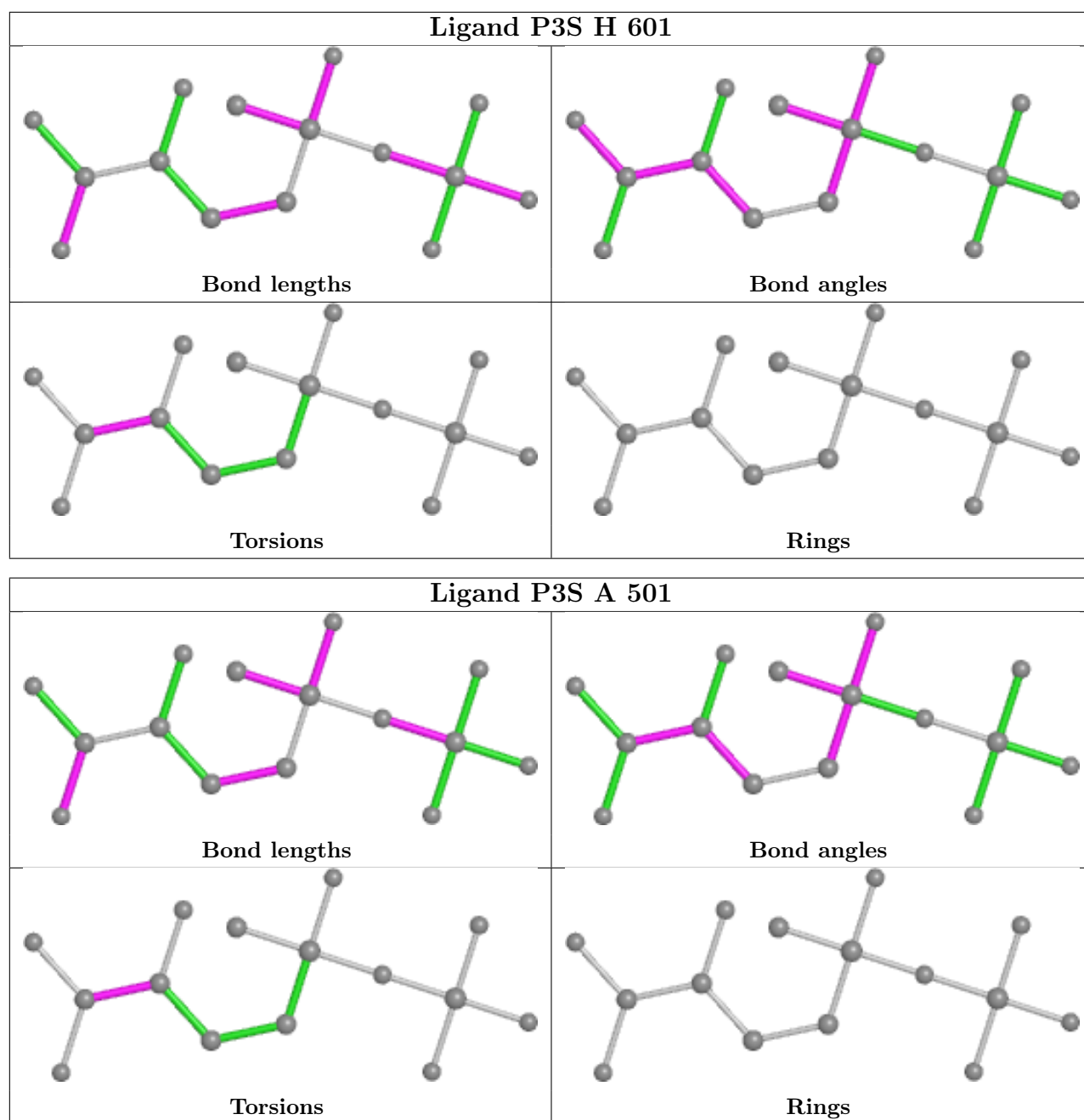
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

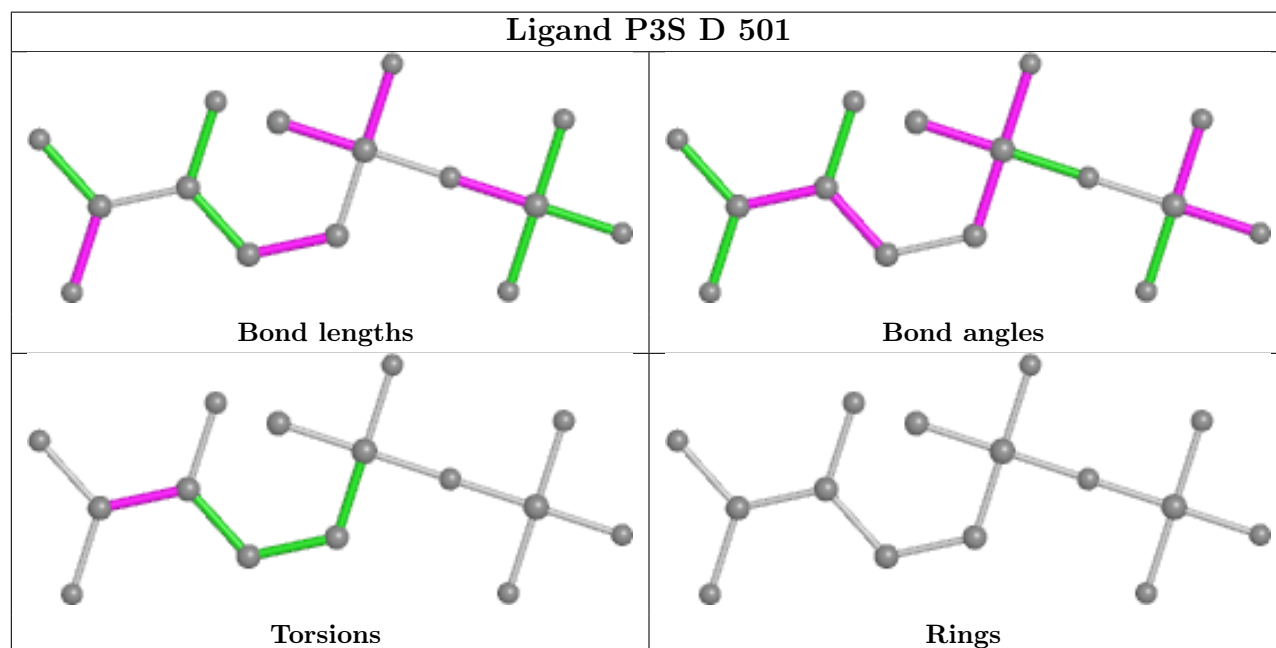
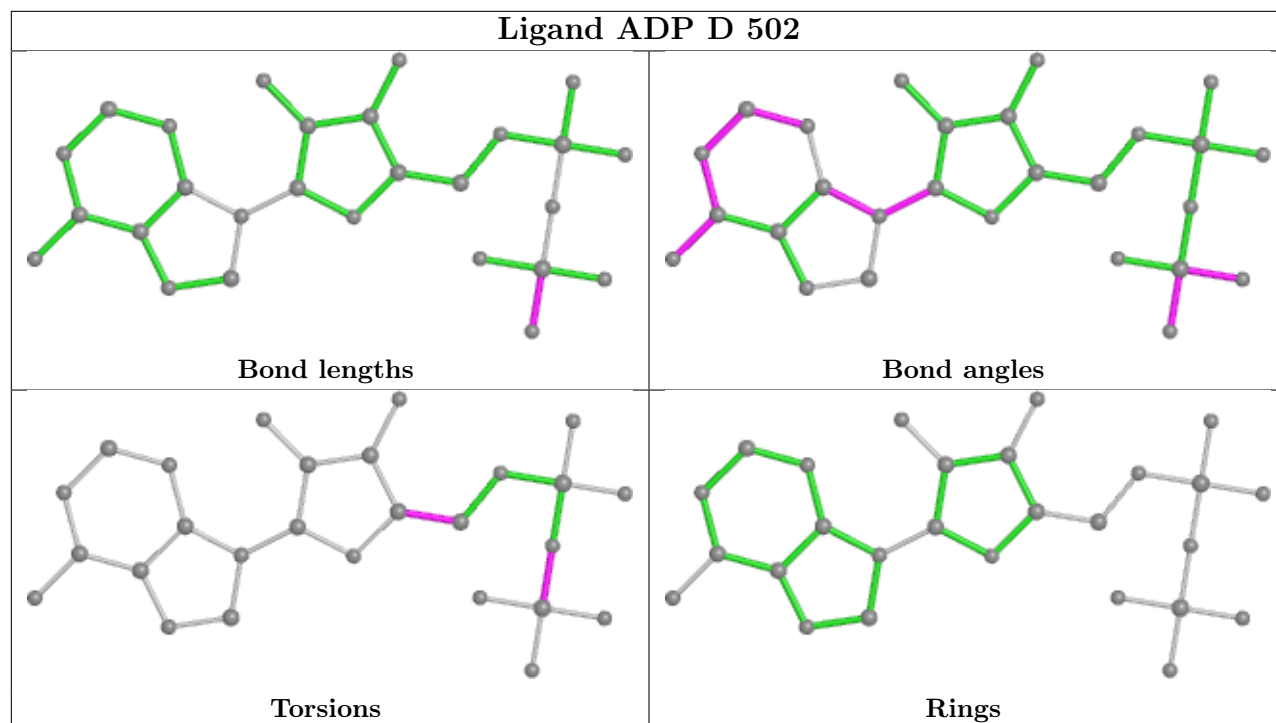
highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified 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.

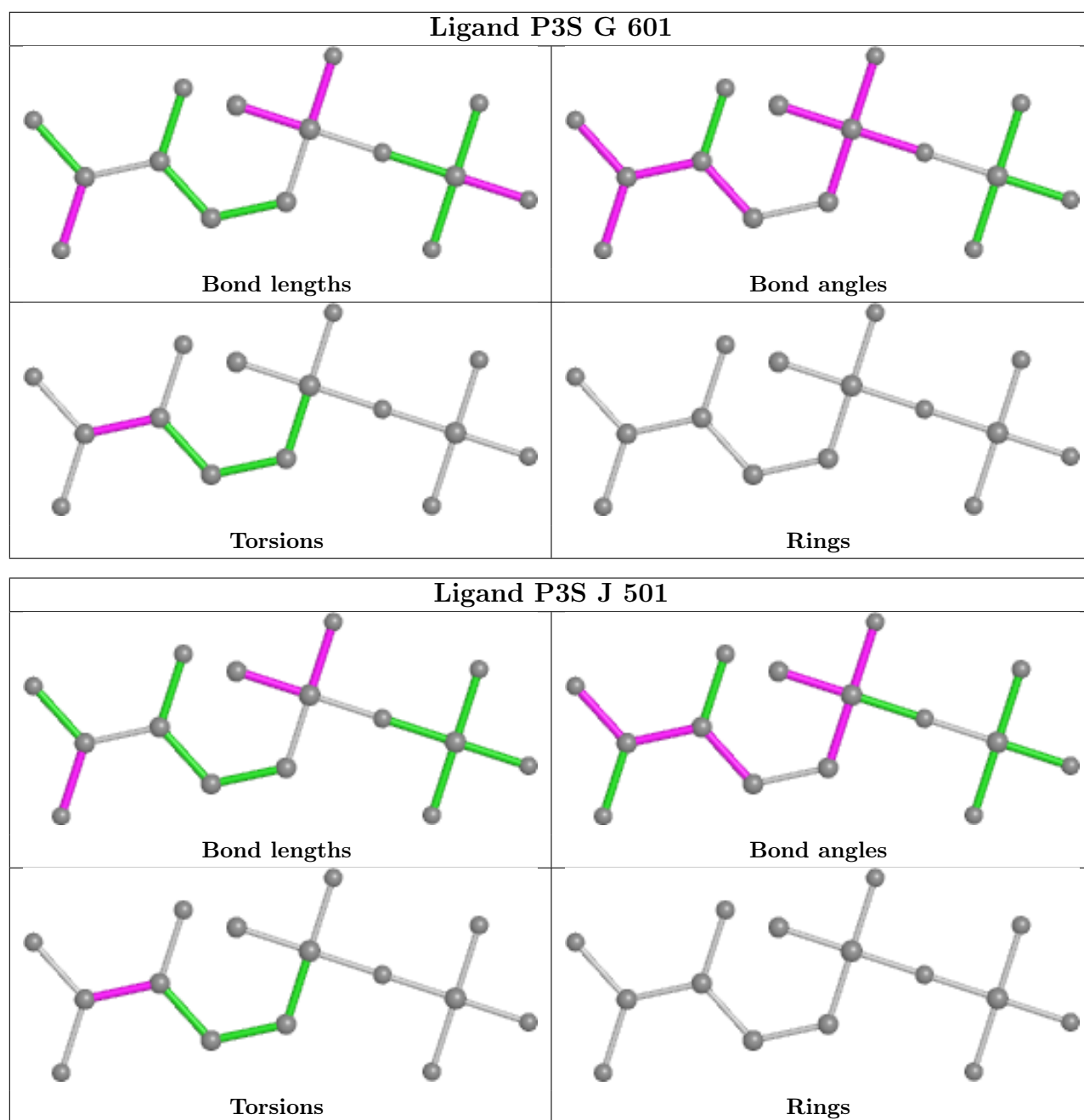


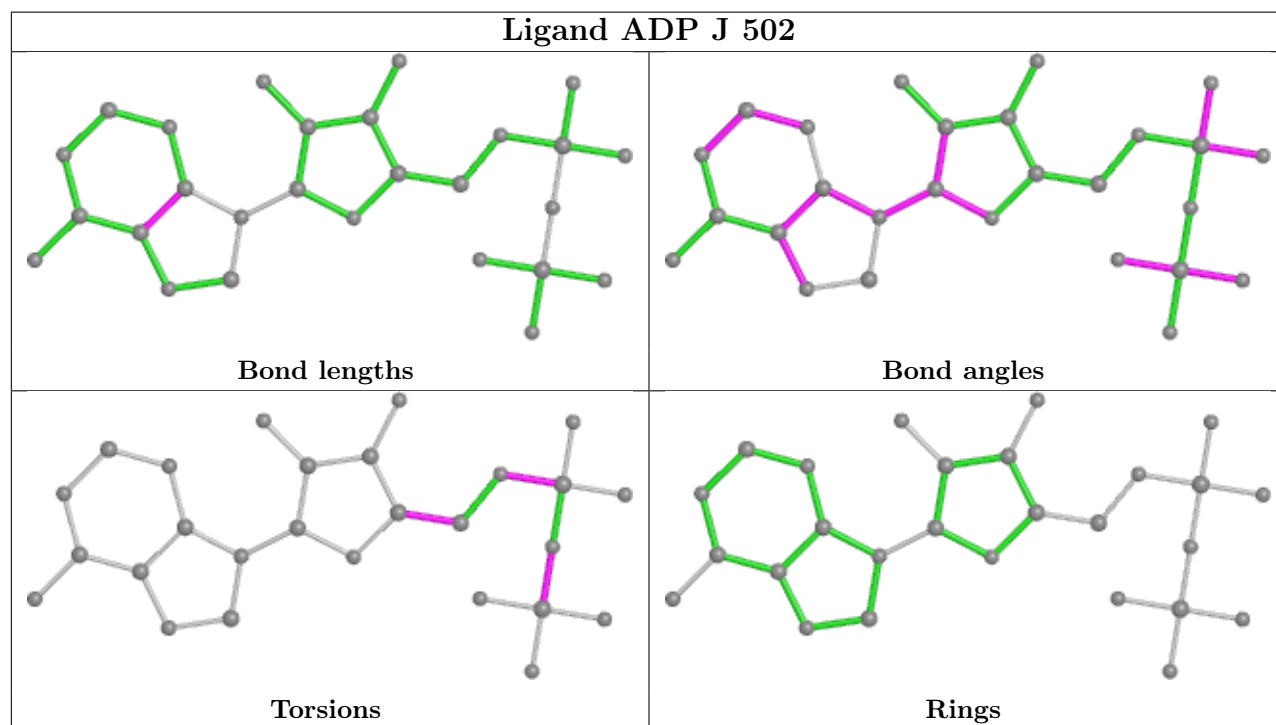
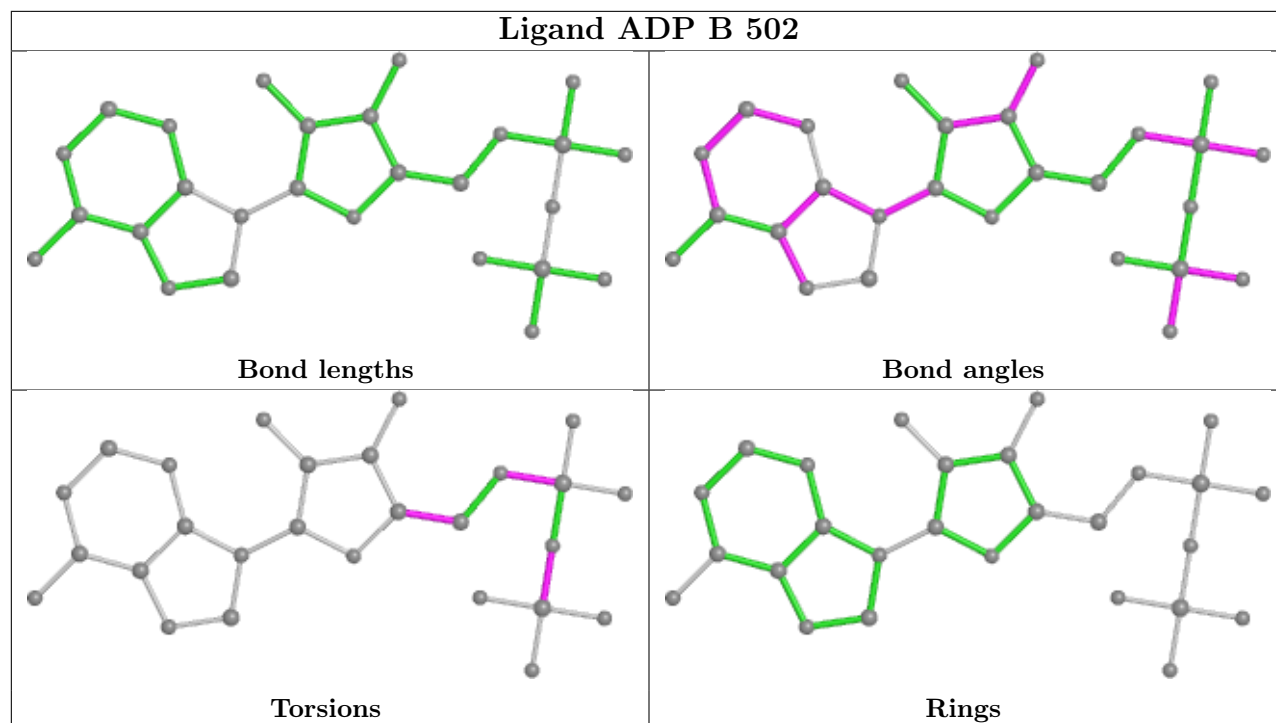


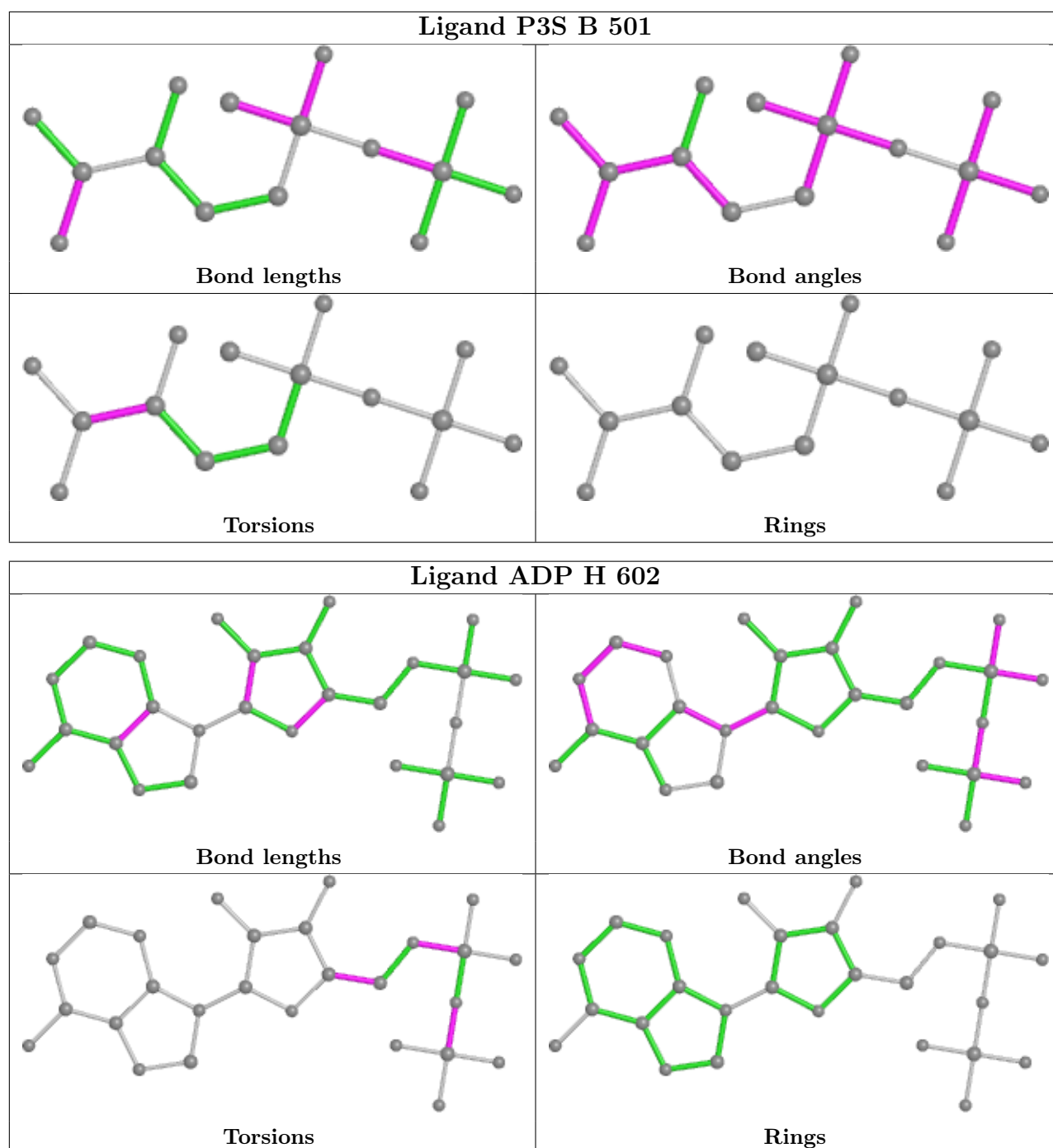


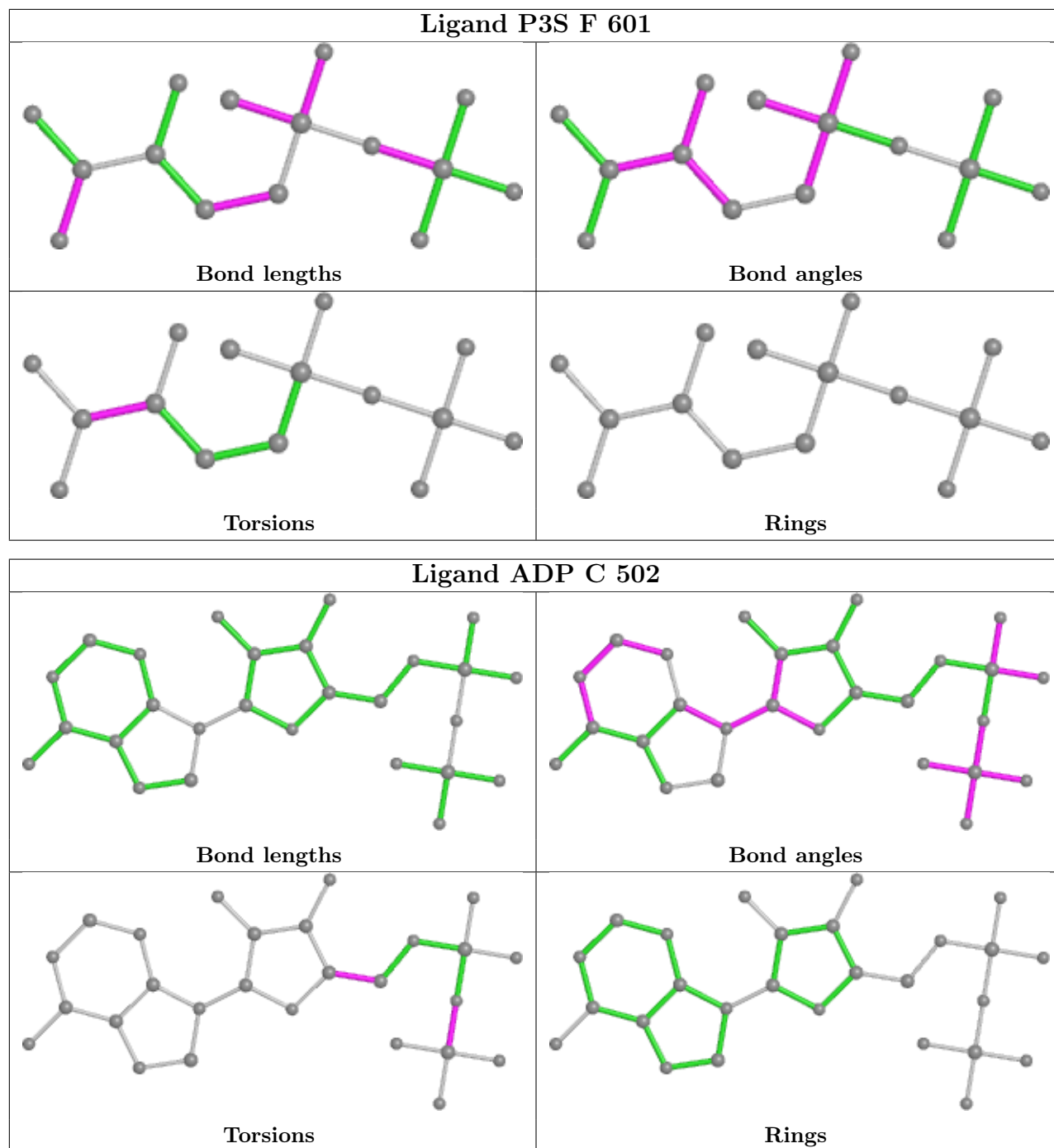


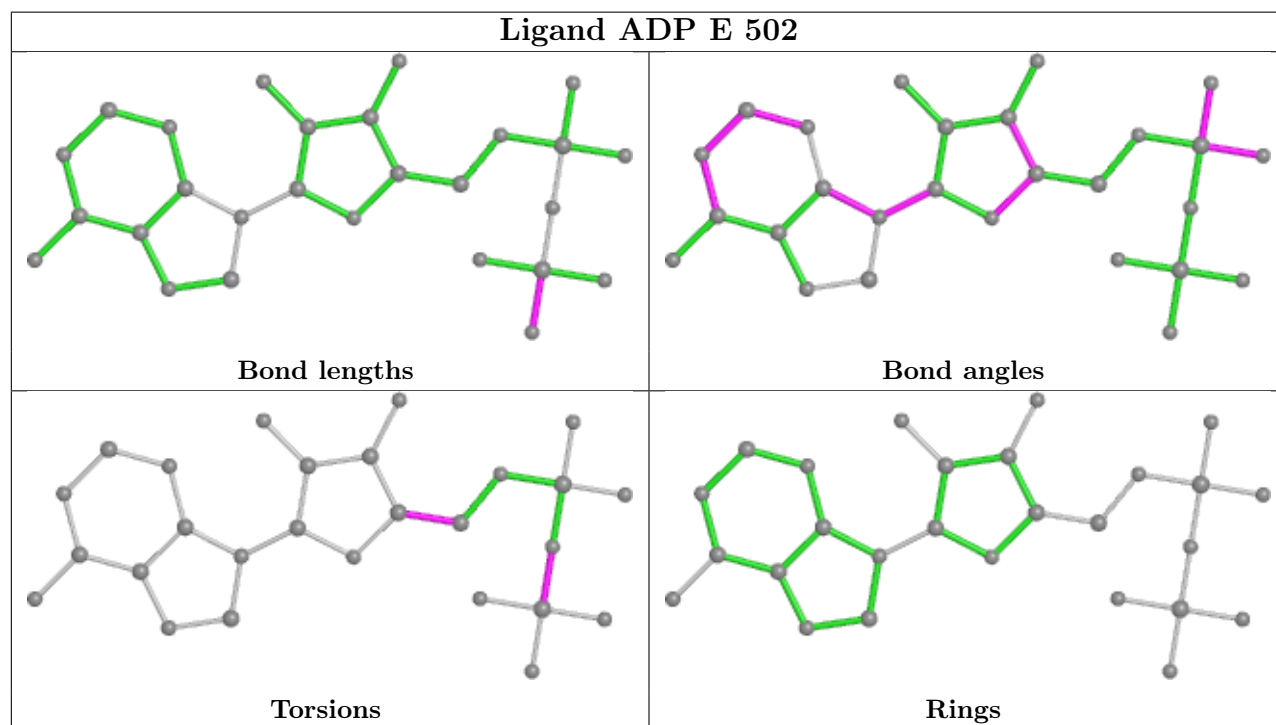
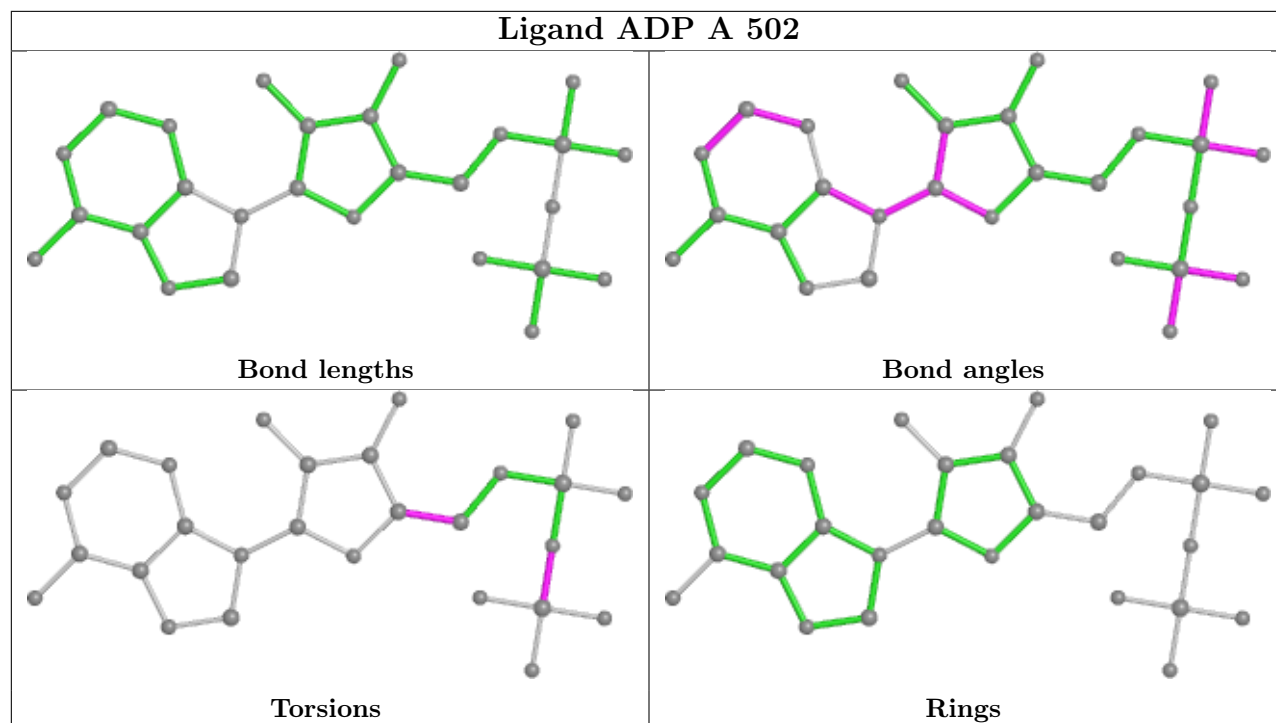


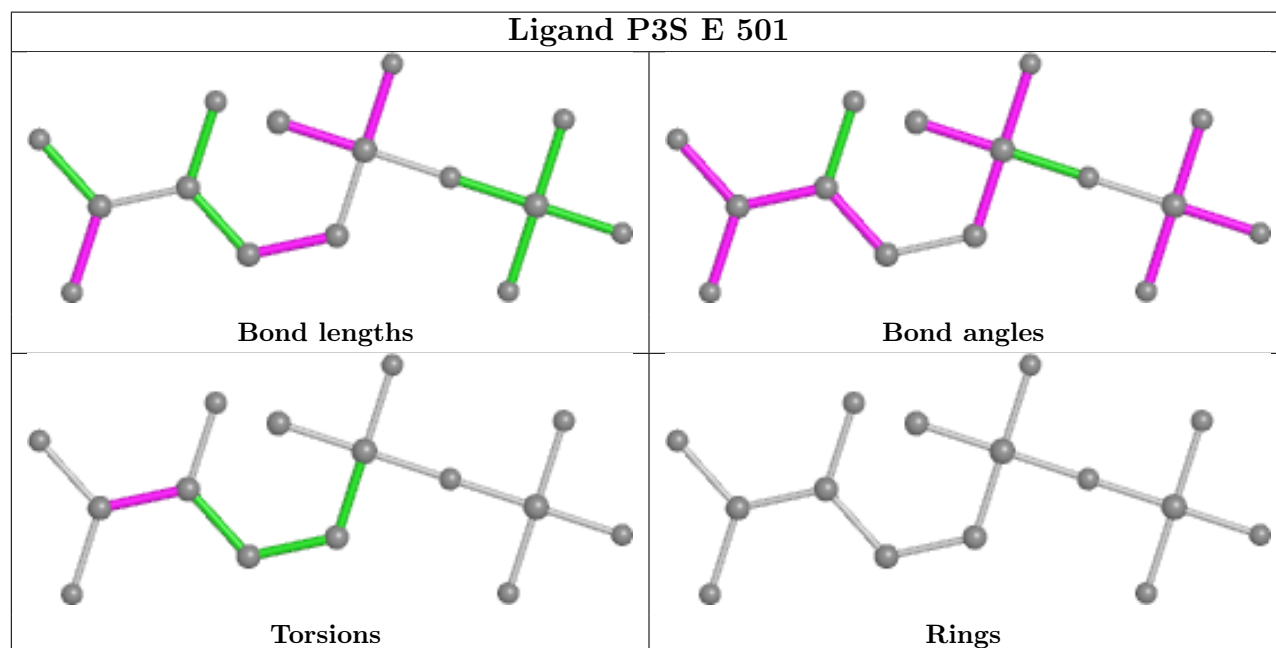
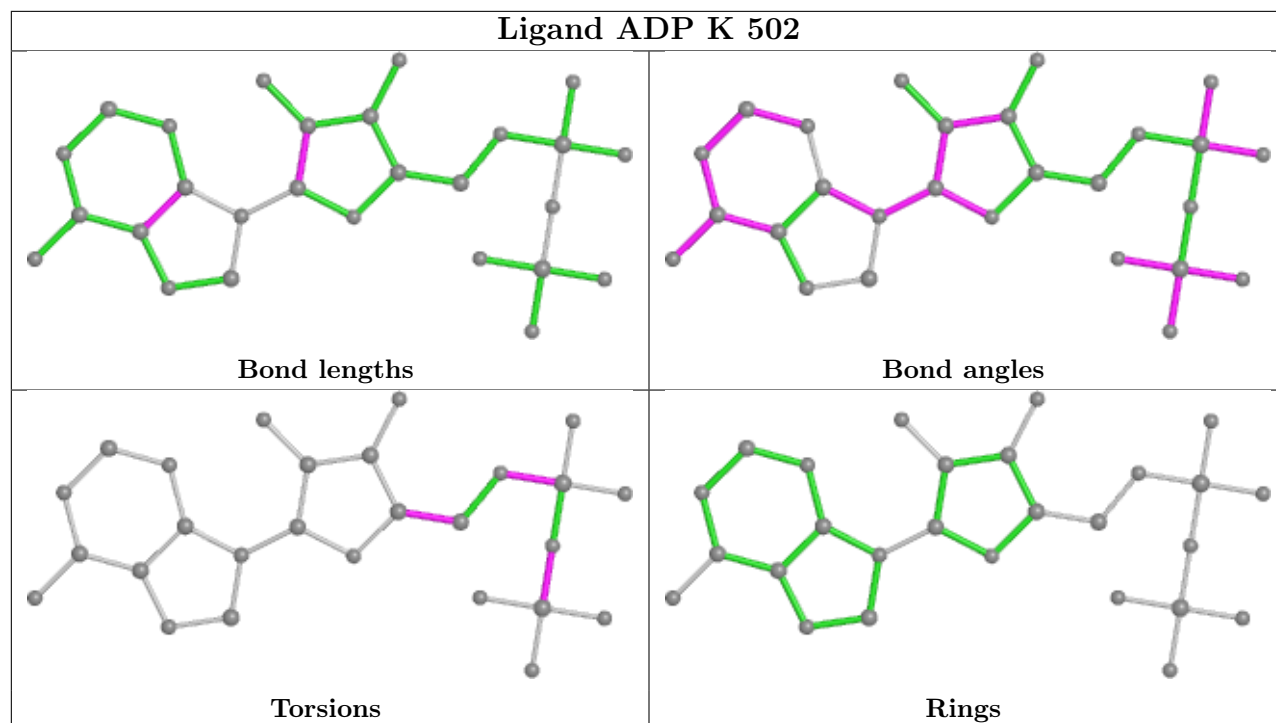


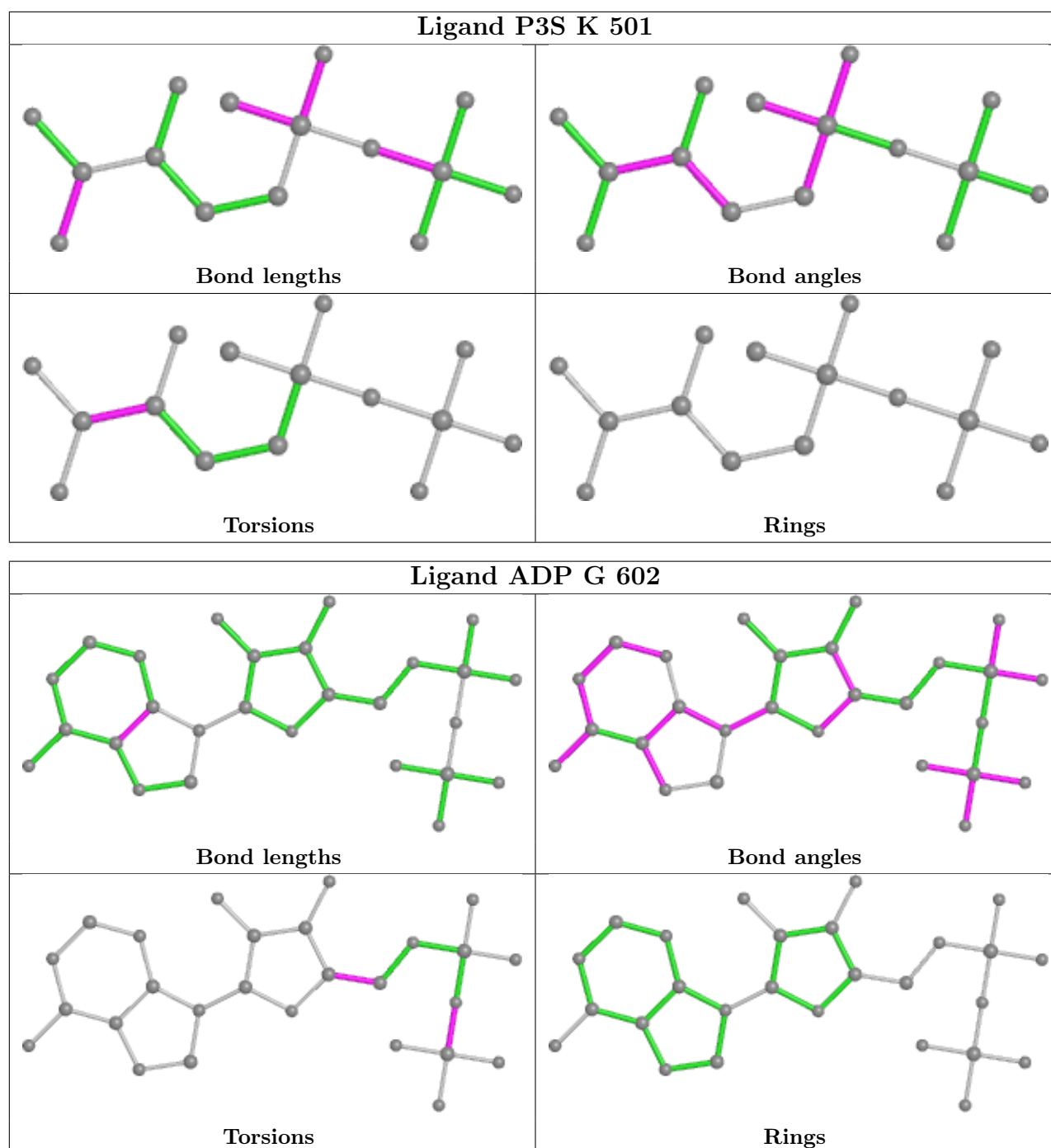












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	443/443 (100%)	-0.21	2 (0%) 91 90	8, 19, 44, 77	0
1	B	443/443 (100%)	-0.20	1 (0%) 95 95	8, 19, 45, 79	0
1	C	443/443 (100%)	-0.25	1 (0%) 95 95	8, 19, 45, 61	0
1	D	443/443 (100%)	-0.20	2 (0%) 91 90	9, 19, 42, 67	0
1	E	443/443 (100%)	-0.21	1 (0%) 95 95	7, 19, 42, 72	0
1	F	443/443 (100%)	-0.21	2 (0%) 91 90	9, 20, 45, 74	0
1	G	443/443 (100%)	-0.20	3 (0%) 87 86	8, 19, 45, 73	0
1	H	443/443 (100%)	-0.20	1 (0%) 95 95	7, 19, 44, 74	0
1	I	443/443 (100%)	-0.24	1 (0%) 95 95	7, 19, 43, 70	0
1	J	441/443 (99%)	-0.21	0 100 100	7, 19, 42, 65	0
1	K	443/443 (100%)	-0.22	2 (0%) 91 90	7, 19, 44, 67	0
1	L	443/443 (100%)	-0.23	2 (0%) 91 90	9, 21, 45, 72	0
All	All	5314/5316 (99%)	-0.22	18 (0%) 94 94	7, 19, 44, 79	0

The worst 5 of 18 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	ALA	8.6
1	E	2	ALA	6.8
1	B	2	ALA	5.4
1	K	2	ALA	4.7
1	F	2	ALA	4.0

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 monosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MG	G	605	1/1	0.94	0.14	7,7,7,7	0
4	MG	E	504	1/1	0.96	0.15	3,3,3,3	0
4	MG	B	505	1/1	0.96	0.16	7,7,7,7	0
4	MG	I	505	1/1	0.96	0.17	6,6,6,6	0
4	MG	D	505	1/1	0.97	0.17	1,1,1,1	0
4	MG	C	505	1/1	0.97	0.16	5,5,5,5	0
4	MG	J	505	1/1	0.97	0.17	0,0,0,0	0
4	MG	L	505	1/1	0.97	0.15	0,0,0,0	0
2	P3S	L	501	15/15	0.98	0.16	13,18,26,32	0
3	ADP	A	502	27/27	0.98	0.15	14,18,24,25	0
3	ADP	B	502	27/27	0.98	0.14	11,19,23,28	0
3	ADP	D	502	27/27	0.98	0.14	11,19,28,32	0
3	ADP	F	602	27/27	0.98	0.14	10,20,27,30	0
3	ADP	G	602	27/27	0.98	0.13	16,22,31,32	0
3	ADP	I	502	27/27	0.98	0.15	13,18,25,26	0
3	ADP	L	502	27/27	0.98	0.14	16,24,35,37	0
4	MG	B	504	1/1	0.98	0.16	4,4,4,4	0
2	P3S	A	501	15/15	0.98	0.18	8,16,28,34	0
2	P3S	B	501	15/15	0.98	0.16	7,15,22,31	0
2	P3S	E	501	15/15	0.98	0.17	7,15,26,33	0
4	MG	E	503	1/1	0.98	0.21	1,1,1,1	0
2	P3S	F	601	15/15	0.98	0.18	12,21,32,45	0
4	MG	F	604	1/1	0.98	0.14	6,6,6,6	0
4	MG	G	603	1/1	0.98	0.15	4,4,4,4	0
2	P3S	G	601	15/15	0.98	0.17	13,24,38,39	0
2	P3S	H	601	15/15	0.98	0.15	9,13,21,26	0
2	P3S	I	501	15/15	0.98	0.19	9,13,27,28	0
4	MG	K	504	1/1	0.98	0.12	5,5,5,5	0
2	P3S	K	501	15/15	0.98	0.18	10,15,21,23	0
2	P3S	J	501	15/15	0.99	0.17	8,16,22,24	0
4	MG	C	503	1/1	0.99	0.20	0,0,0,0	0
2	P3S	D	501	15/15	0.99	0.18	8,18,34,41	0

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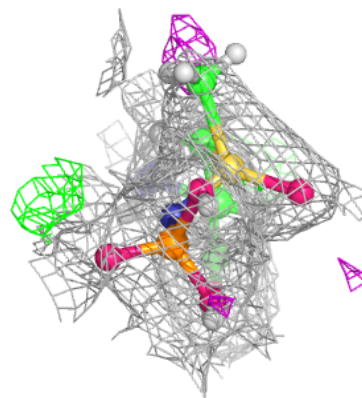
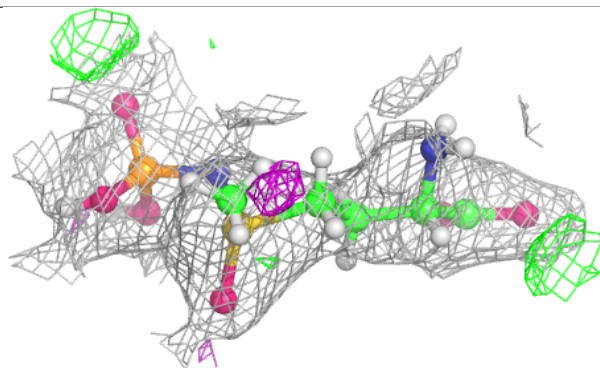
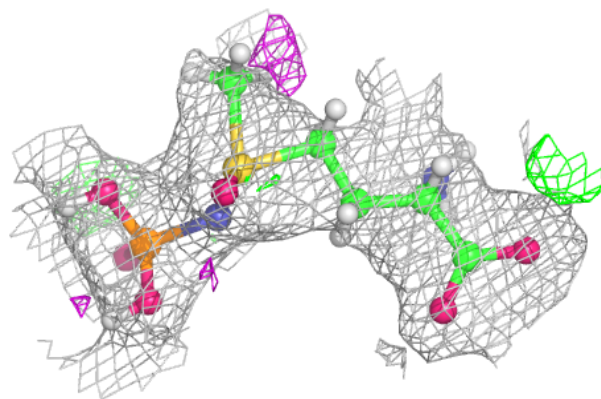
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MG	D	503	1/1	0.99	0.18	5,5,5,5	0
4	MG	D	504	1/1	0.99	0.10	2,2,2,2	0
3	ADP	H	602	27/27	0.99	0.14	11,18,23,25	0
3	ADP	C	502	27/27	0.99	0.15	11,18,21,23	0
3	ADP	J	502	27/27	0.99	0.15	13,18,24,33	0
4	MG	E	505	1/1	0.99	0.16	4,4,4,4	0
4	MG	F	603	1/1	0.99	0.15	5,5,5,5	0
3	ADP	K	502	27/27	0.99	0.14	8,15,22,25	0
4	MG	F	605	1/1	0.99	0.14	5,5,5,5	0
2	P3S	C	501	15/15	0.99	0.17	4,13,27,40	0
4	MG	G	604	1/1	0.99	0.12	10,10,10,10	0
4	MG	A	503	1/1	0.99	0.17	4,4,4,4	0
4	MG	H	603	1/1	0.99	0.17	1,1,1,1	0
4	MG	H	604	1/1	0.99	0.15	5,5,5,5	0
4	MG	H	605	1/1	0.99	0.16	1,1,1,1	0
4	MG	I	503	1/1	0.99	0.18	0,0,0,0	0
4	MG	I	504	1/1	0.99	0.15	5,5,5,5	0
4	MG	A	504	1/1	0.99	0.15	4,4,4,4	0
4	MG	J	503	1/1	0.99	0.17	5,5,5,5	0
4	MG	J	504	1/1	0.99	0.12	4,4,4,4	0
4	MG	A	505	1/1	0.99	0.17	6,6,6,6	0
4	MG	B	503	1/1	0.99	0.22	0,0,0,0	0
4	MG	K	505	1/1	0.99	0.19	1,1,1,1	0
4	MG	L	503	1/1	0.99	0.15	2,2,2,2	0
4	MG	L	504	1/1	0.99	0.12	7,7,7,7	0
3	ADP	E	502	27/27	0.99	0.12	8,15,23,31	0
4	MG	K	503	1/1	1.00	0.17	0,0,0,0	0
4	MG	C	504	1/1	1.00	0.11	6,6,6,6	0

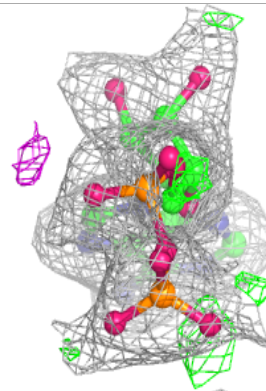
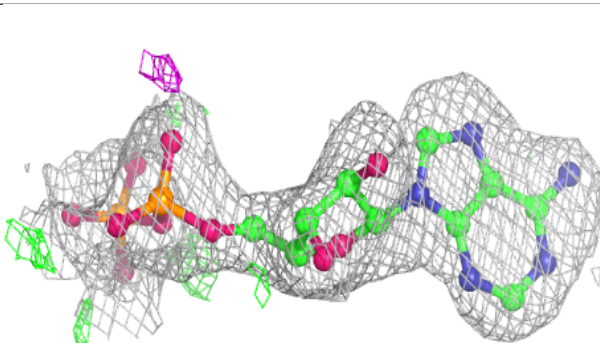
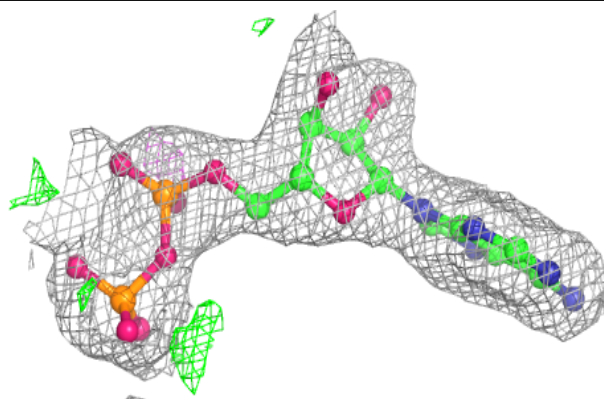
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around P3S L 501:

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

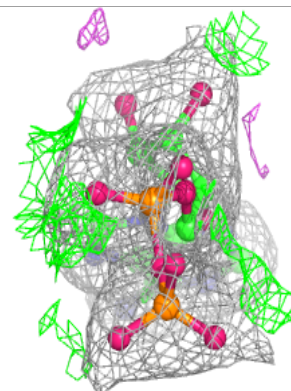
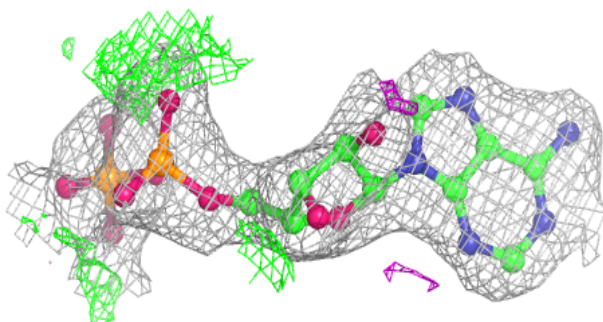
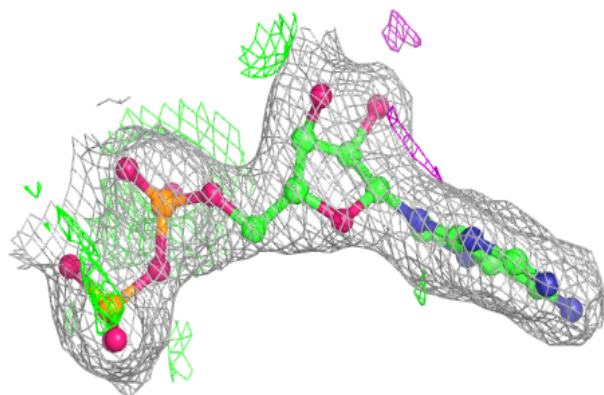
**Electron density around ADP A 502:**

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

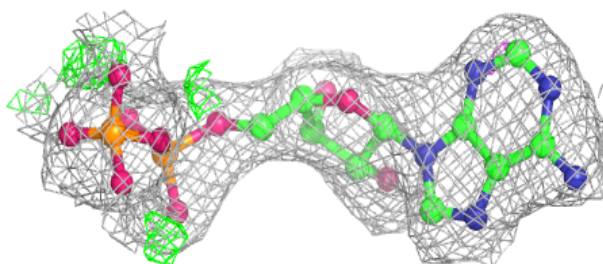
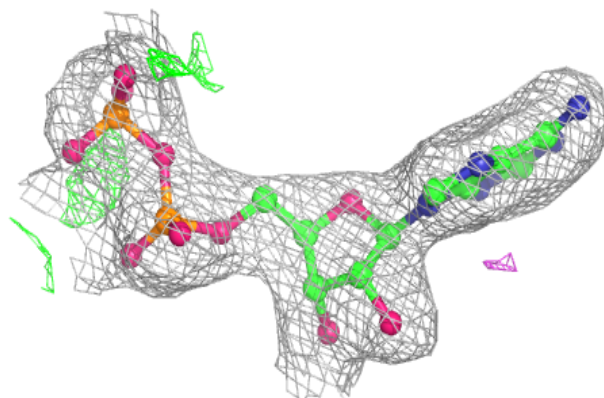


Electron density around ADP B 502:

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

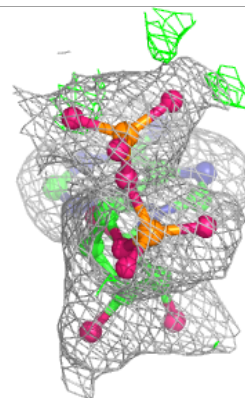
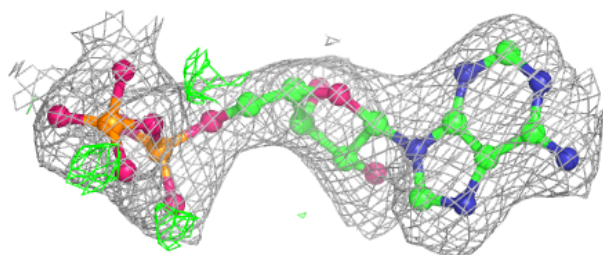
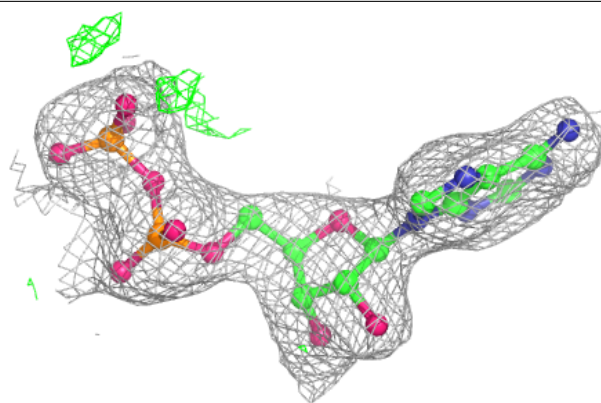
**Electron density around ADP D 502:**

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

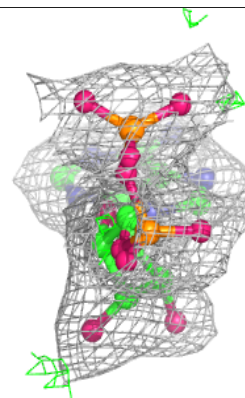
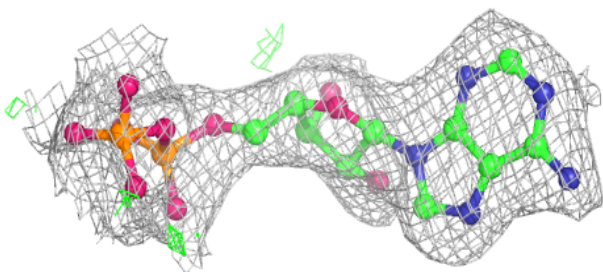
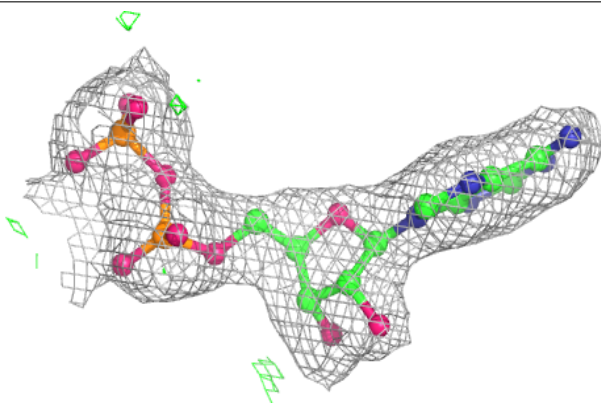


Electron density around ADP F 602:

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

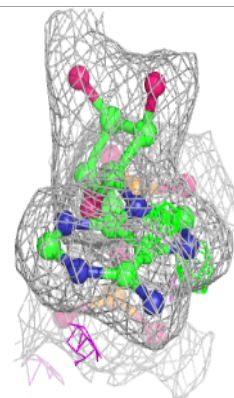
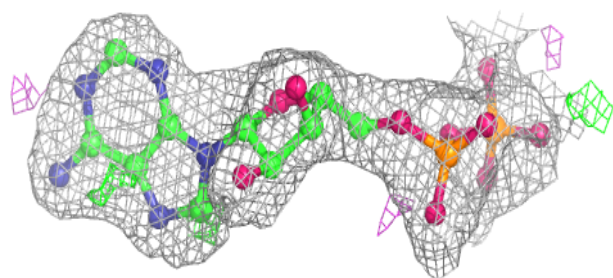
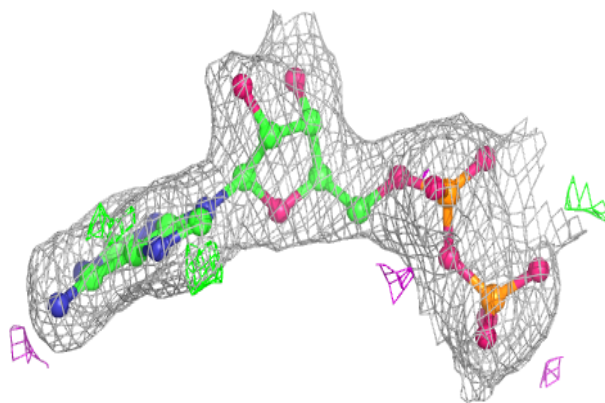
**Electron density around ADP G 602:**

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

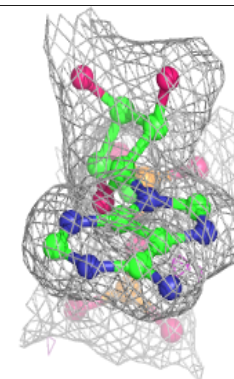
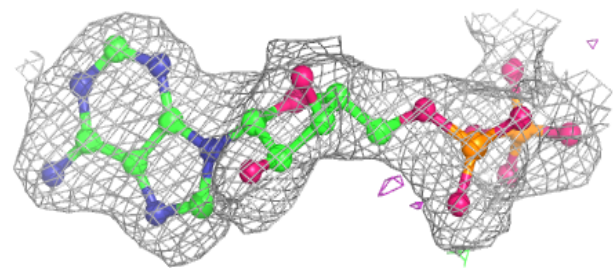
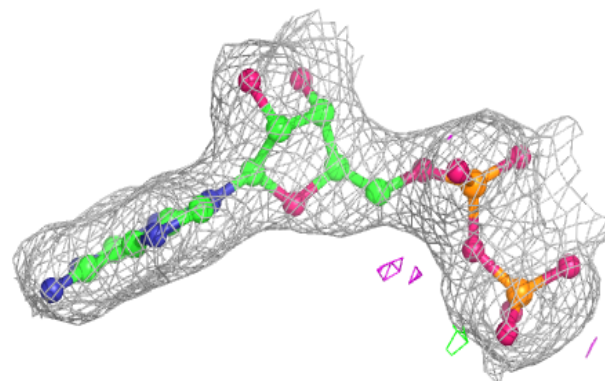


Electron density around ADP I 502:

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

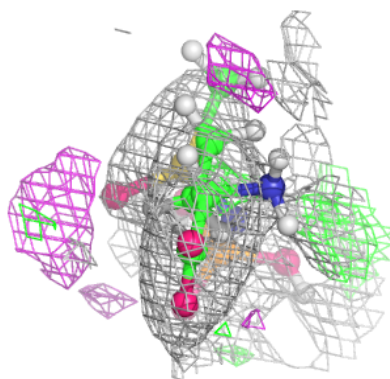
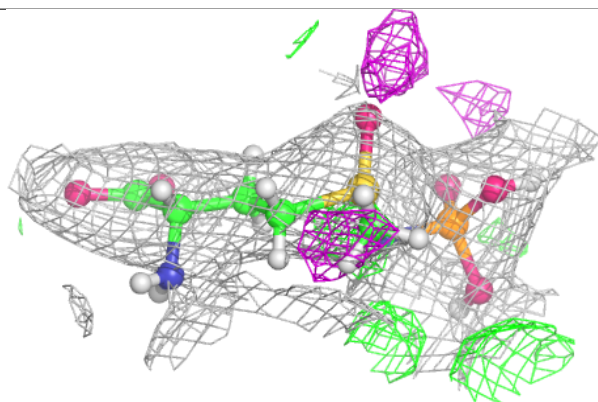
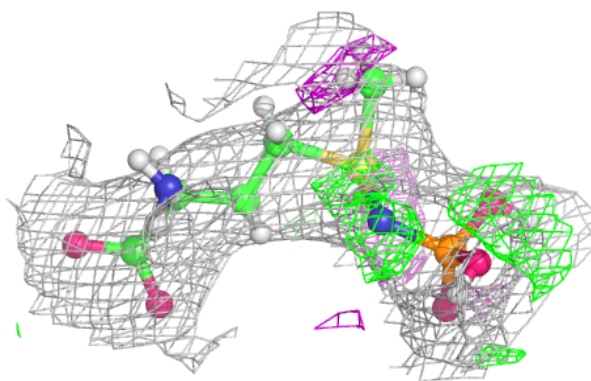
**Electron density around ADP L 502:**

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

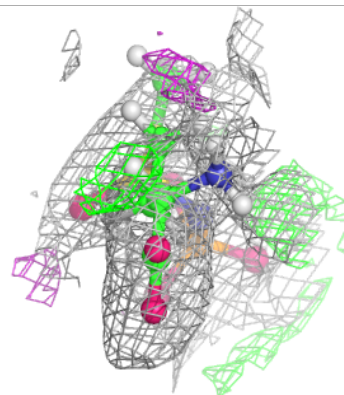
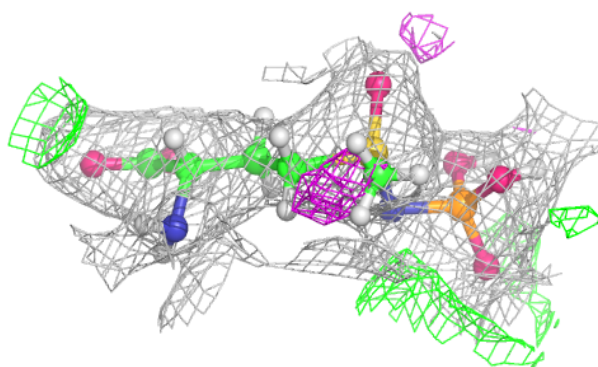
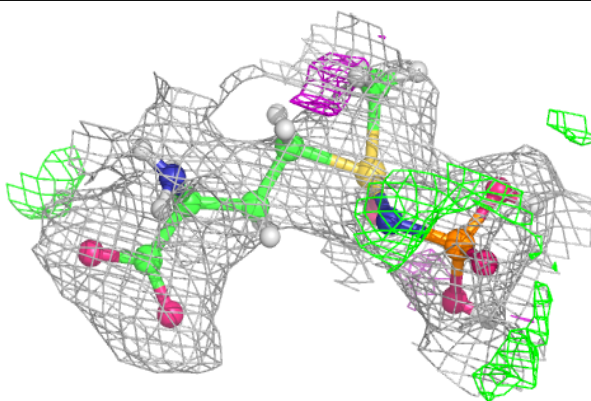


Electron density around P3S A 501:

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

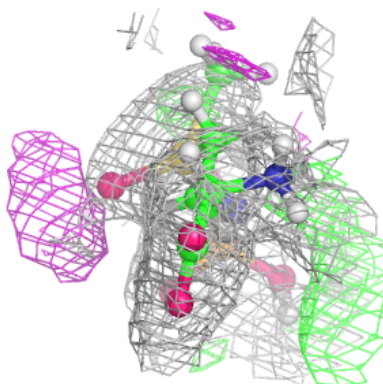
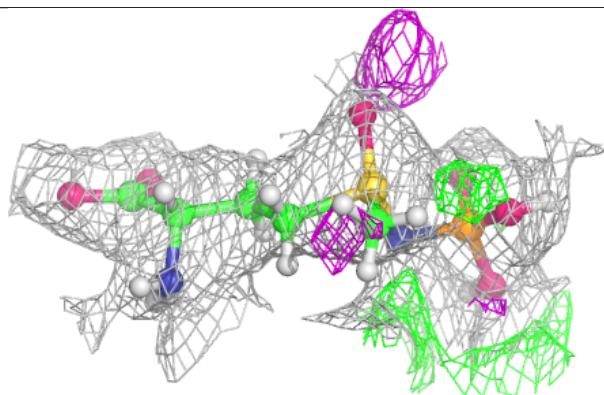
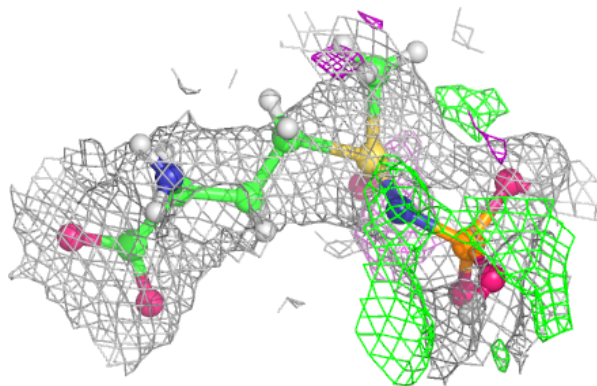
**Electron density around P3S B 501:**

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

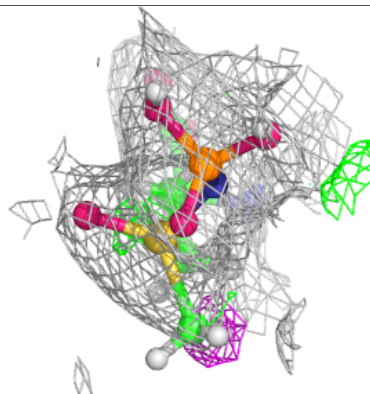
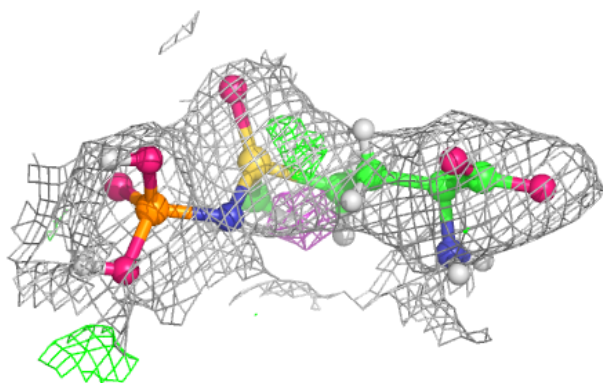
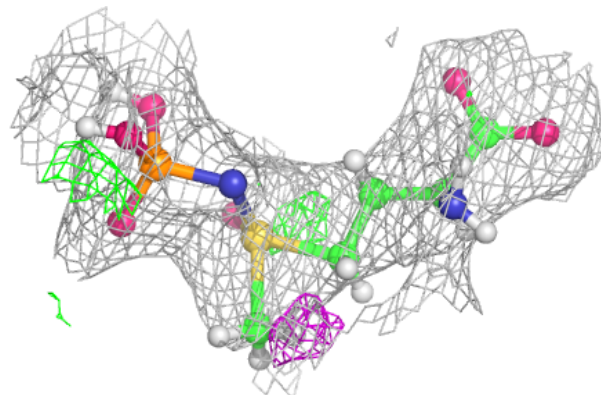


Electron density around P3S E 501:

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

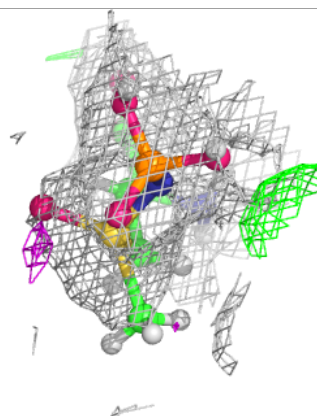
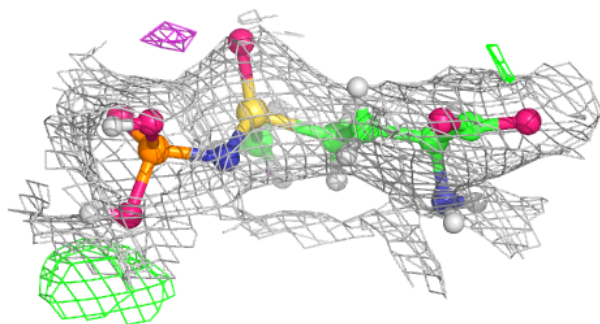
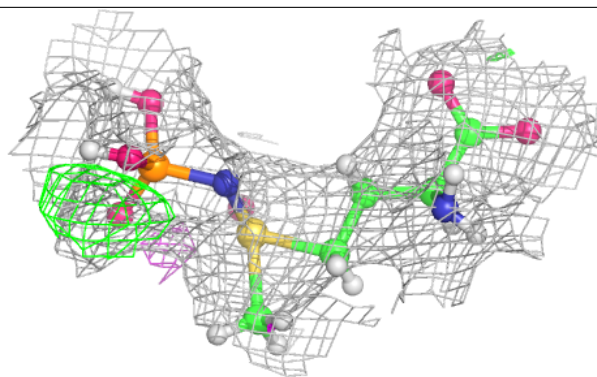
**Electron density around P3S F 601:**

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

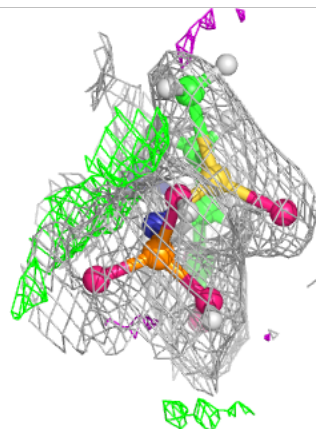
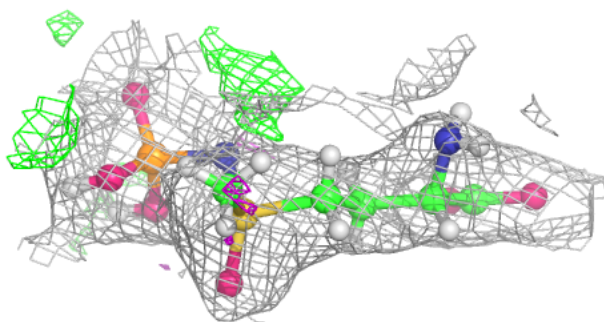
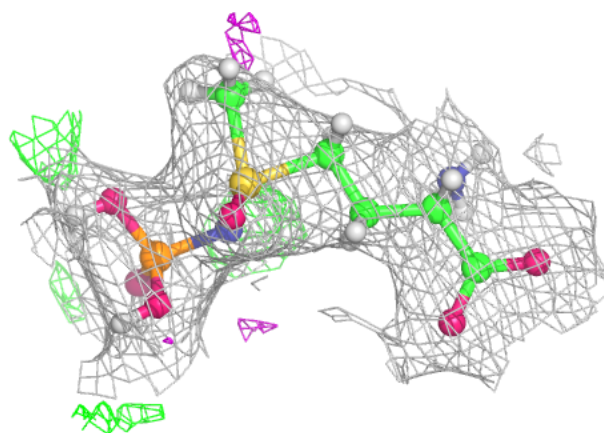


Electron density around P3S G 601:

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

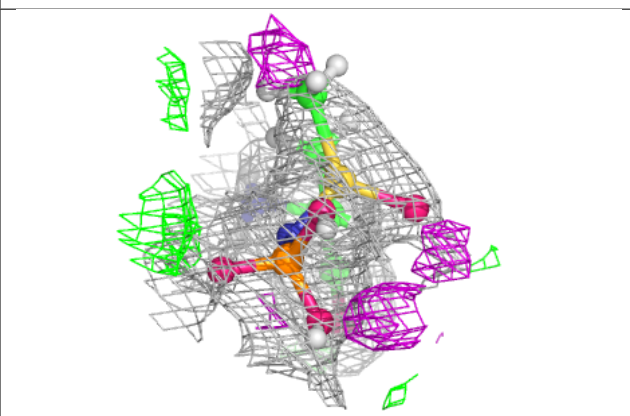
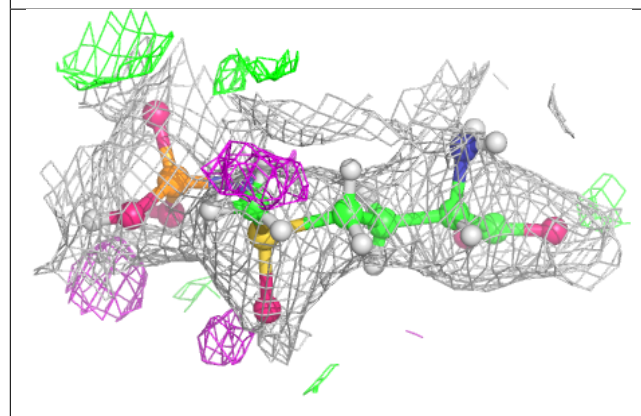
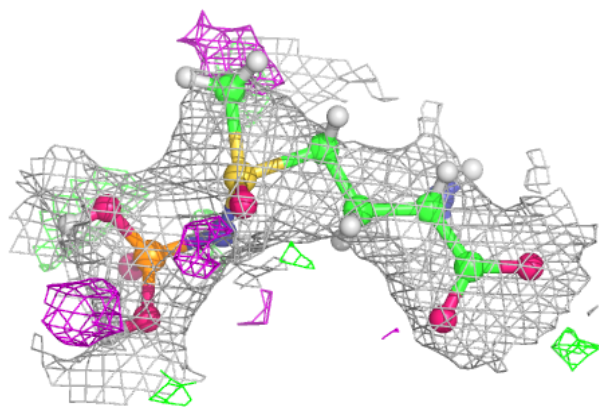
**Electron density around P3S H 601:**

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

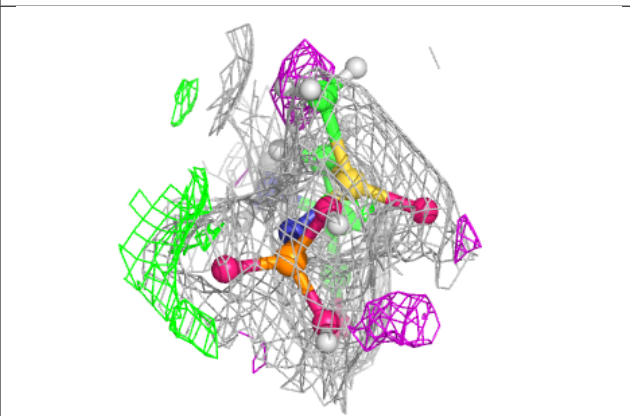
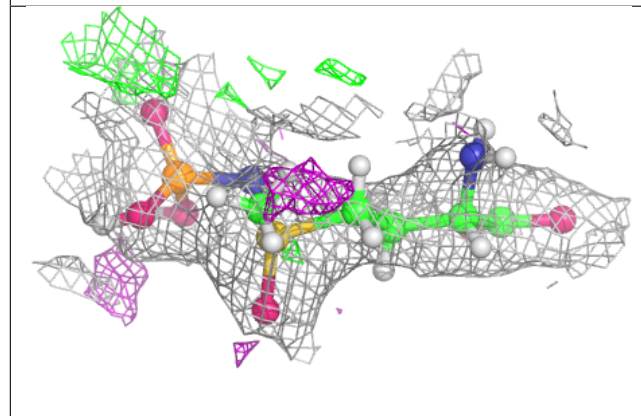
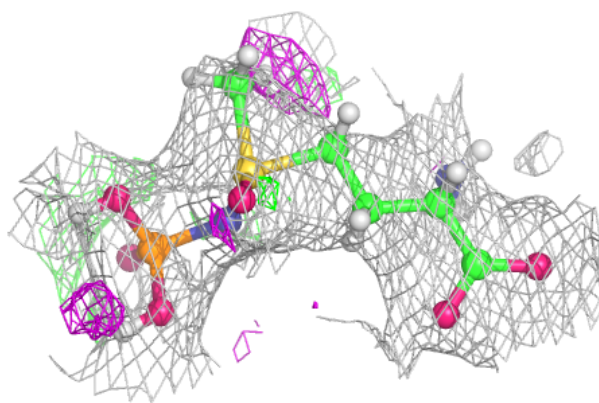


Electron density around P3S I 501:

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

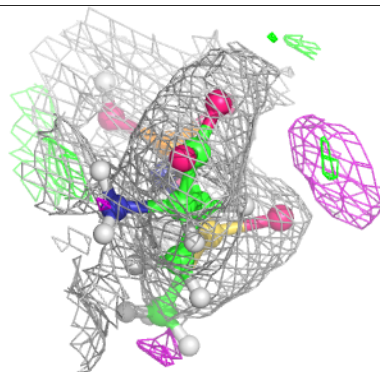
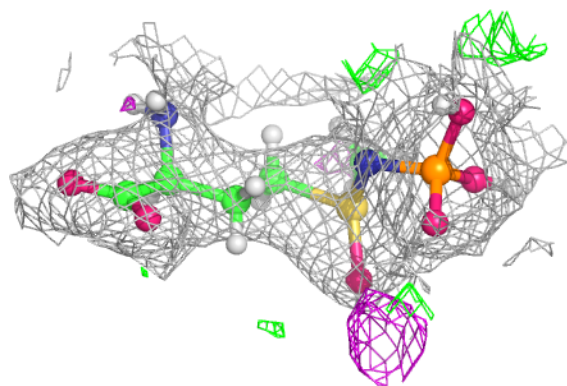
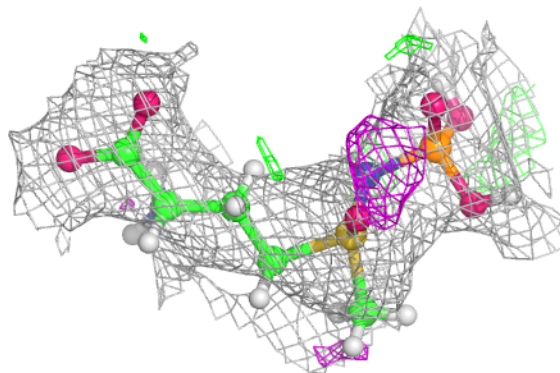
**Electron density around P3S K 501:**

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



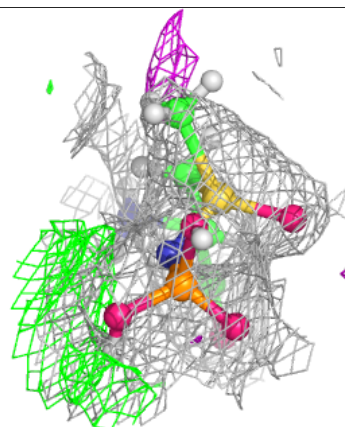
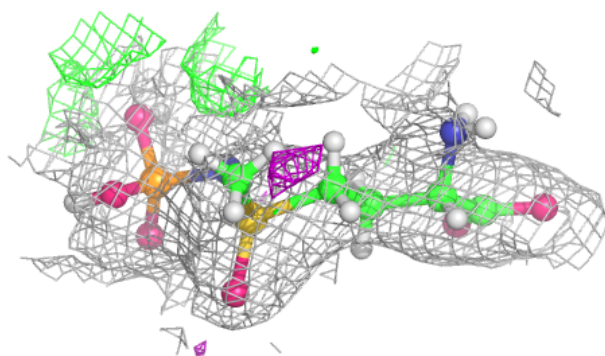
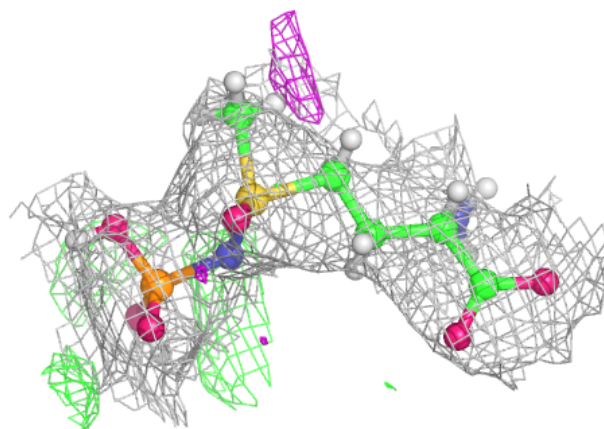
Electron density around P3S J 501:

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

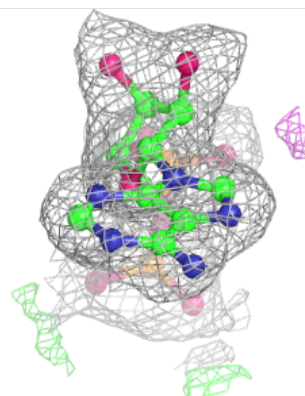
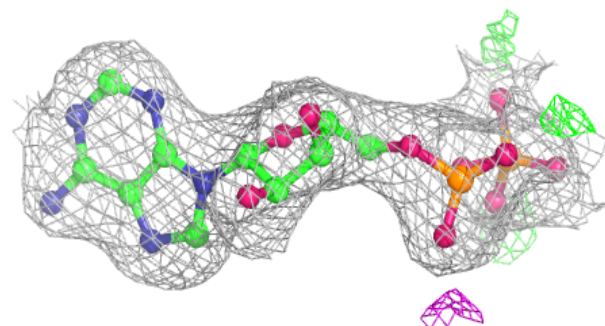
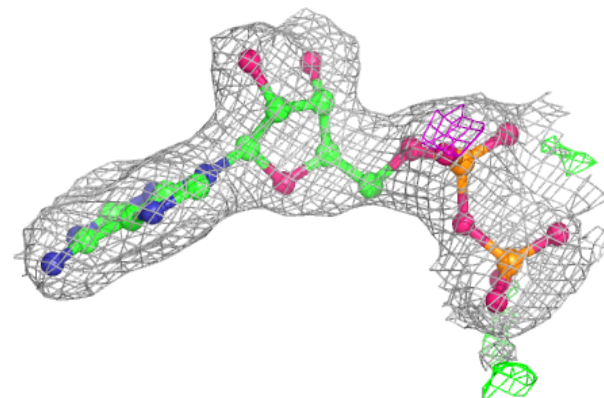


Electron density around P3S D 501:

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

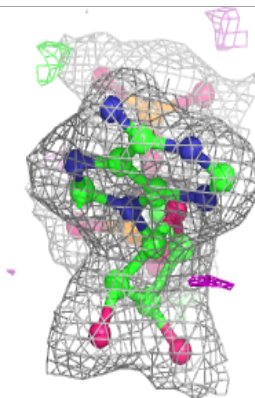
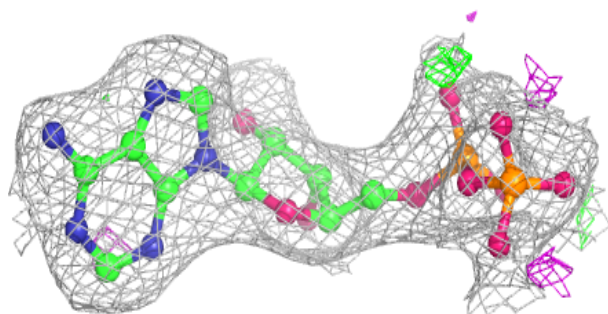
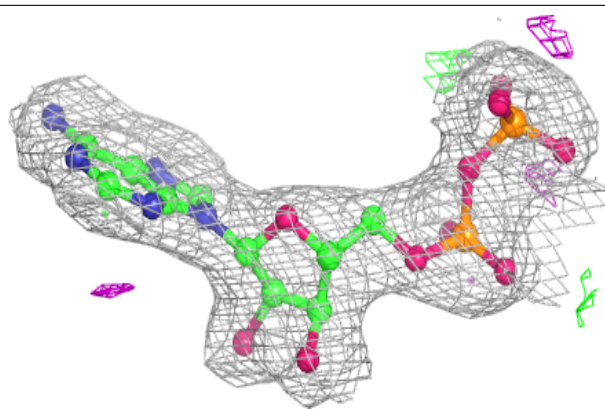
**Electron density around ADP H 602:**

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

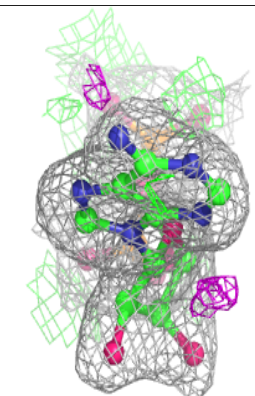
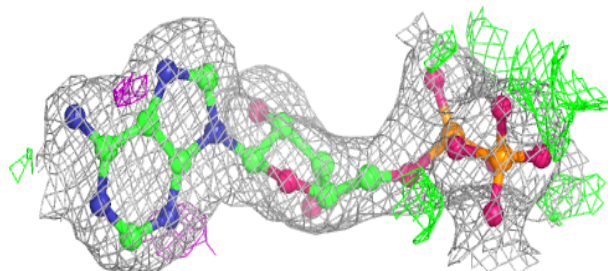
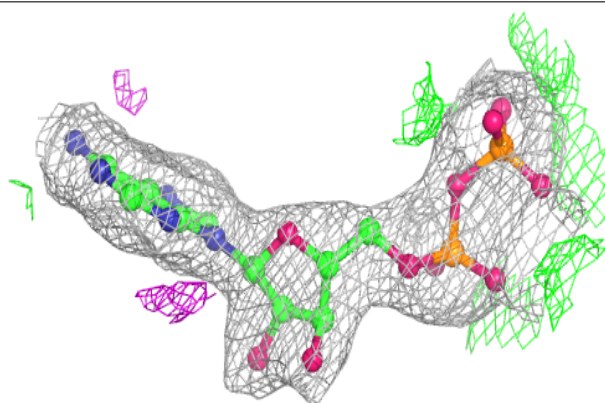


Electron density around ADP C 502:

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

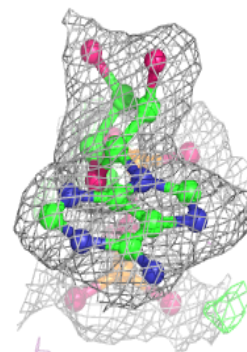
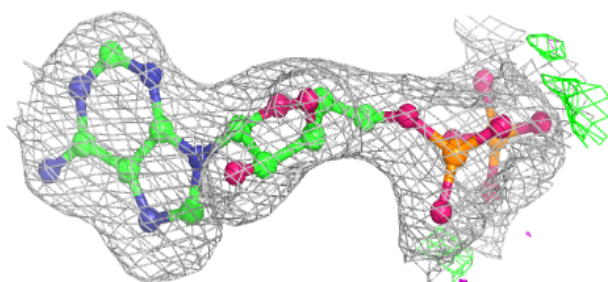
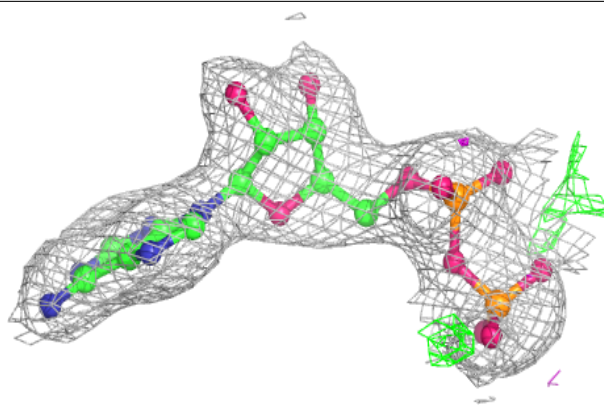
**Electron density around ADP J 502:**

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

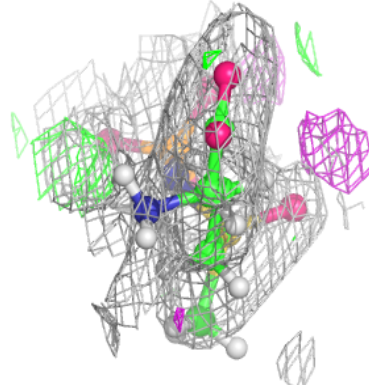
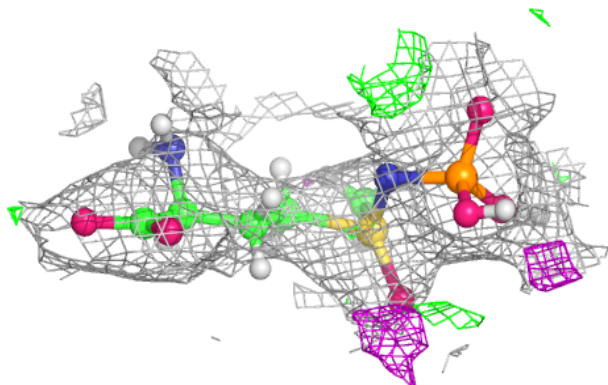
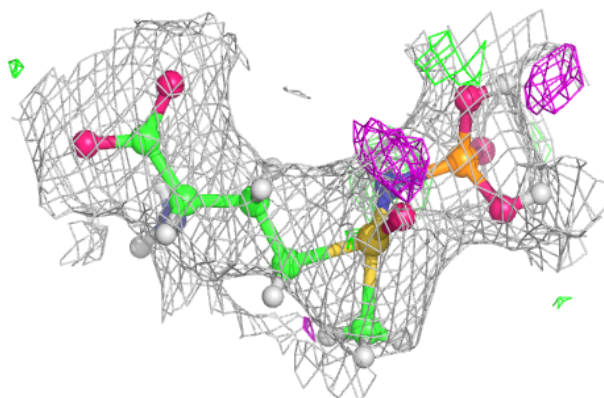


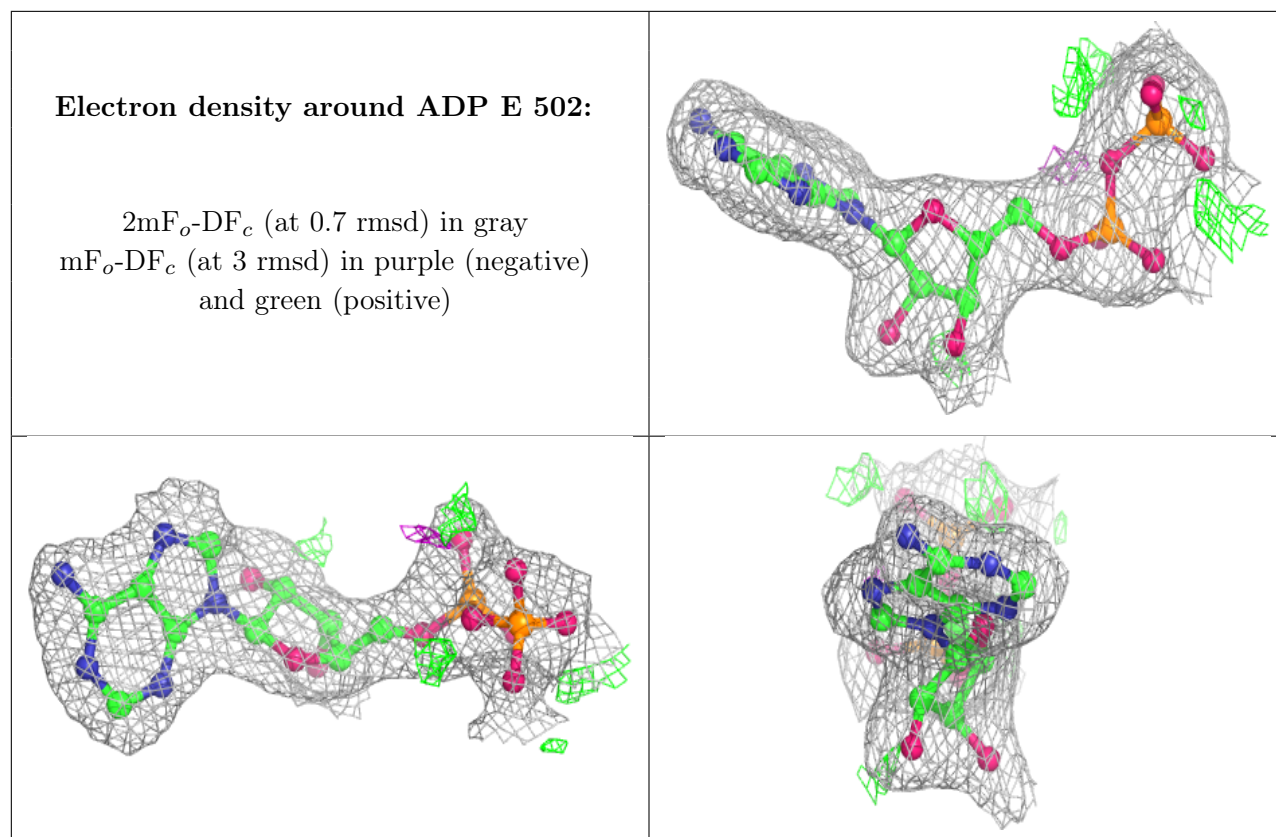
Electron density around ADP K 502:

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

**Electron density around P3S C 501:**

$2mF_o-DF_c$ (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.