



wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 18, 2026 – 12:04 PM UTC

PDB ID : 1Z7E / pdb_00001z7e
Title : Crystal structure of full length ArnA
Authors : Gatzeva-Topalova, P.Z.; May, A.P.; Sousa, M.C.
Deposited on : 2005-03-24
Resolution : 3.00 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

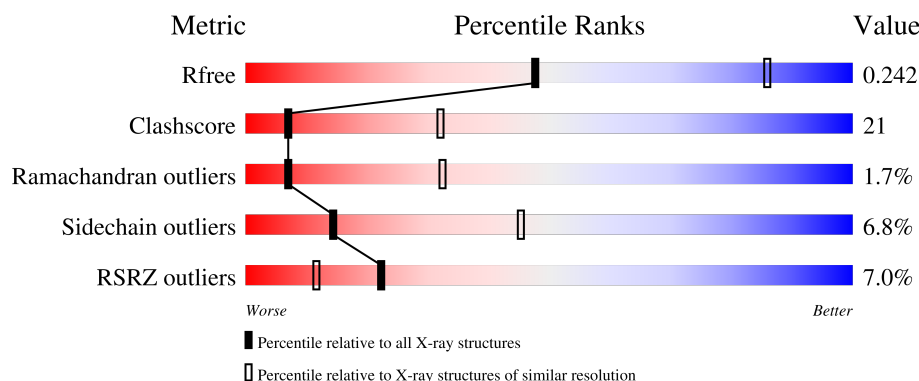
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	660	5% 56% 36% 5%
1	B	660	7% 57% 34% 5%
1	C	660	6% 58% 34% 5%
1	D	660	6% 57% 35% 5%
1	E	660	4% 58% 34% 5%

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Mol	Chain	Length	Quality of chain
1	F	660	14% 55% 37%

2 Entry composition [i](#)

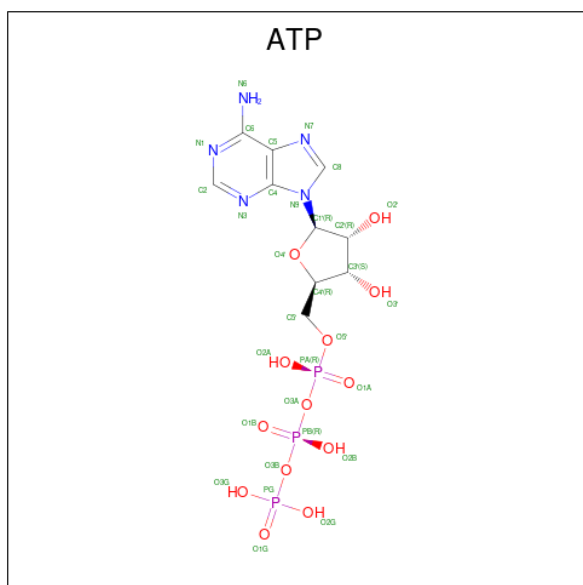
There are 3 unique types of molecules in this entry. The entry contains 30556 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 protein ArnA.

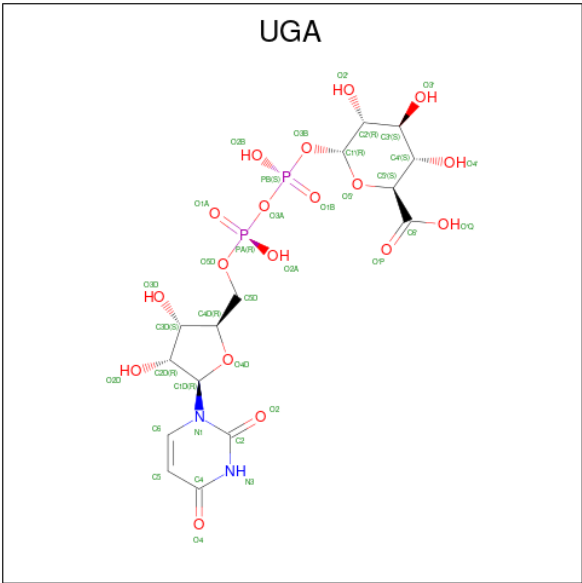
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	639	Total	C	N	O	S	0	0	0
			5020	3210	883	908	19			
1	B	639	Total	C	N	O	S	0	0	0
			5020	3210	883	908	19			
1	C	639	Total	C	N	O	S	0	0	0
			5020	3210	883	908	19			
1	D	644	Total	C	N	O	S	0	0	0
			5048	3227	888	914	19			
1	E	639	Total	C	N	O	S	0	0	0
			5020	3210	883	908	19			
1	F	639	Total	C	N	O	S	0	0	0
			5020	3210	883	908	19			

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	C	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	D	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	E	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	F	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 3 is URIDINE-5'-DIPHOSPHATE-GLUCURONIC ACID (CCD ID: UGA) (formula: C₁₅H₂₂N₂O₁₈P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			37	15	2	18	2		
3	B	1	Total	C	N	O	P	0	0
			37	15	2	18	2		
3	C	1	Total	C	N	O	P	0	0
			37	15	2	18	2		
3	D	1	Total	C	N	O	P	0	0
			37	15	2	18	2		
3	E	1	Total	C	N	O	P	0	0
			37	15	2	18	2		

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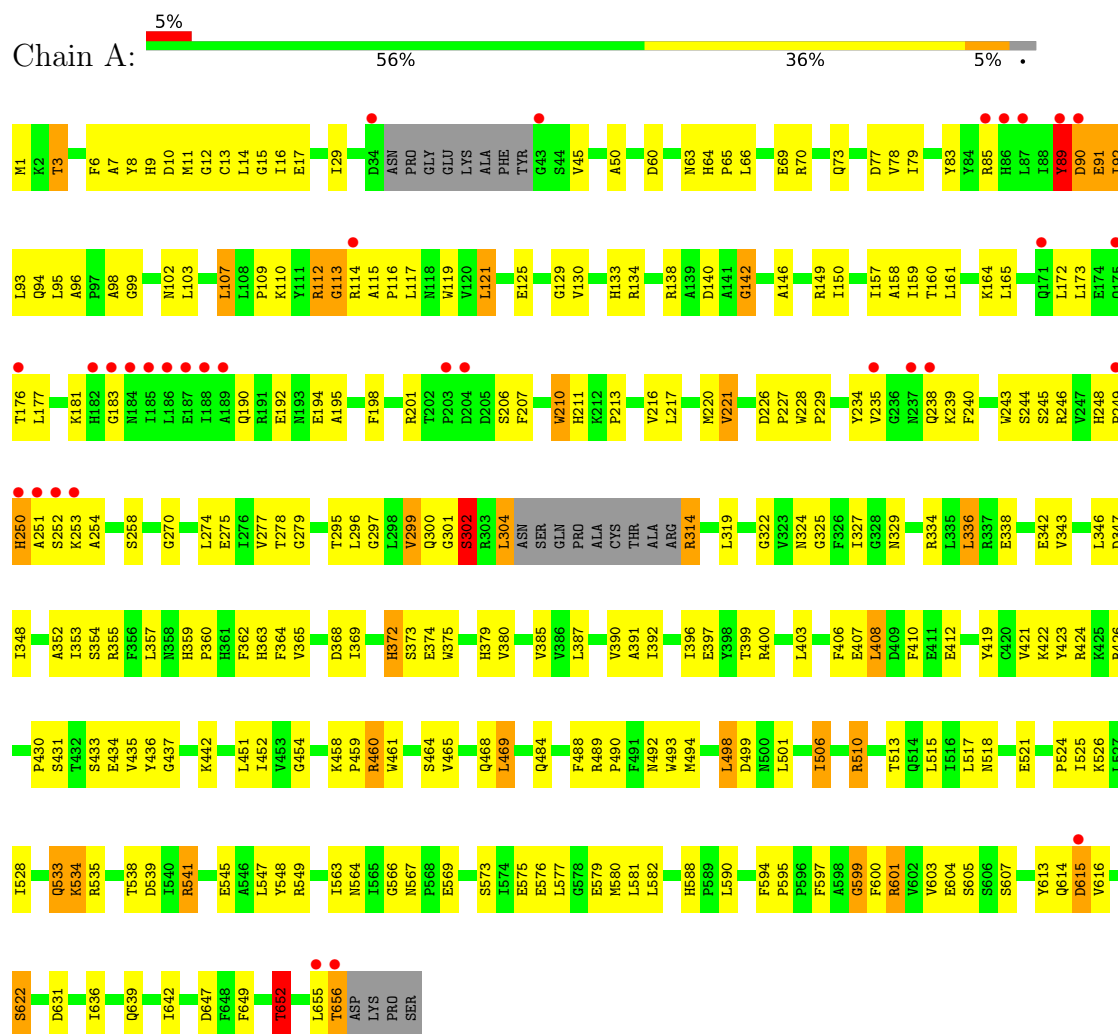
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	F	1	Total	C	N	O	P	0	0
			37	15	2	18	2		

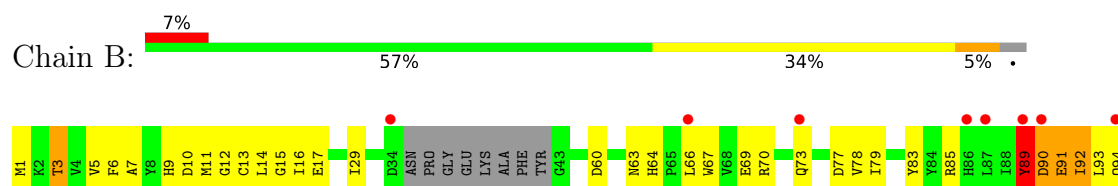
3 Residue-property plots

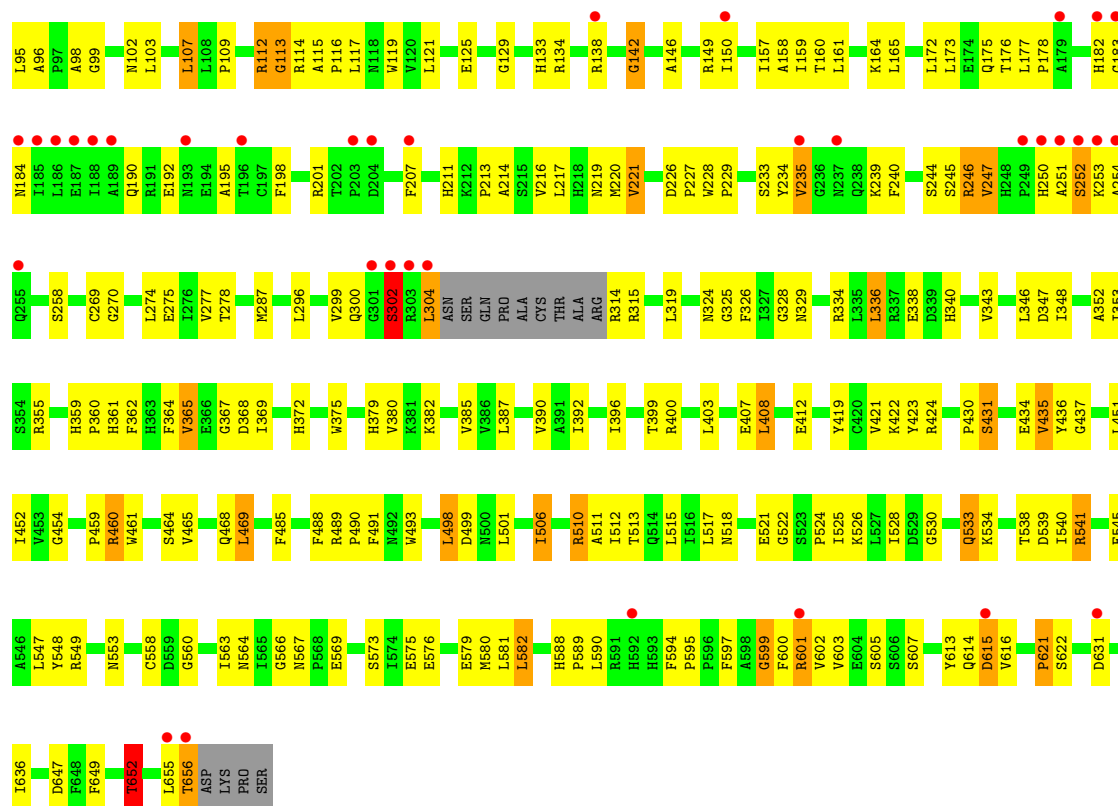
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

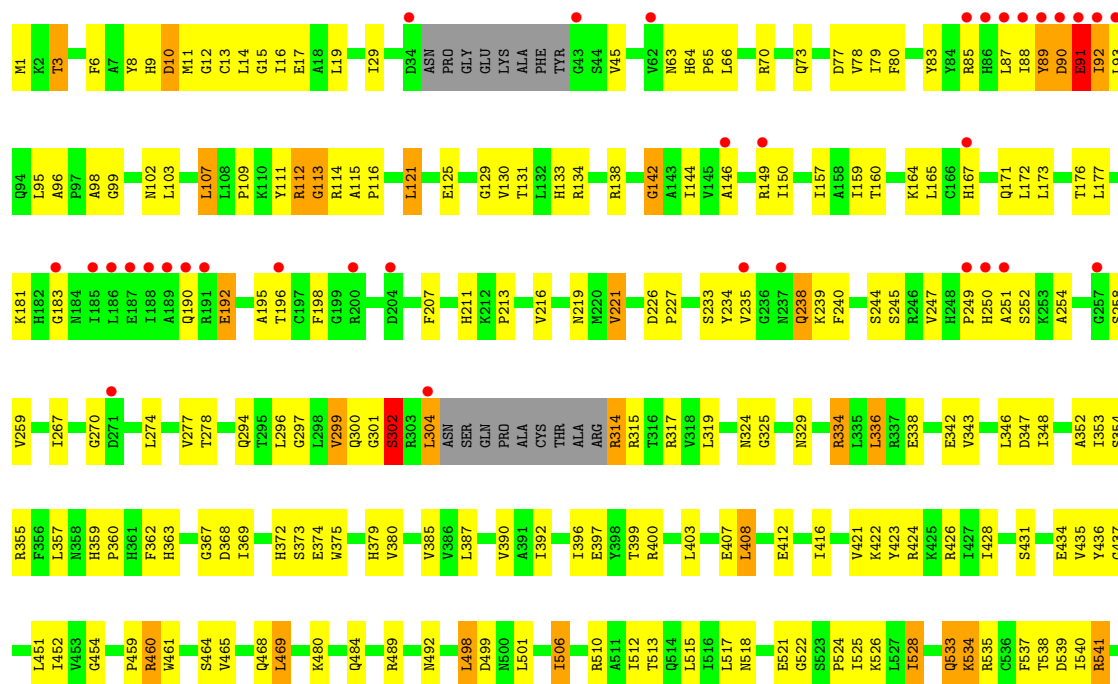


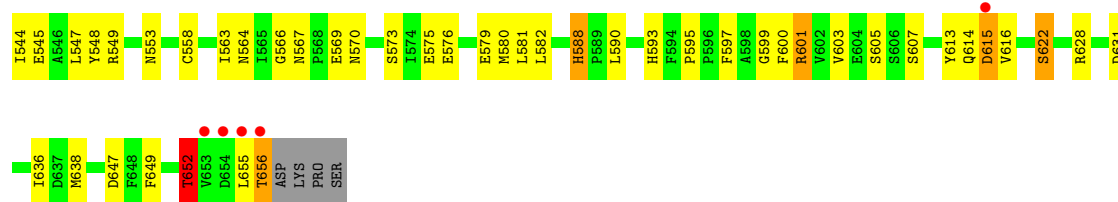
• Molecule 1: protein ArnA



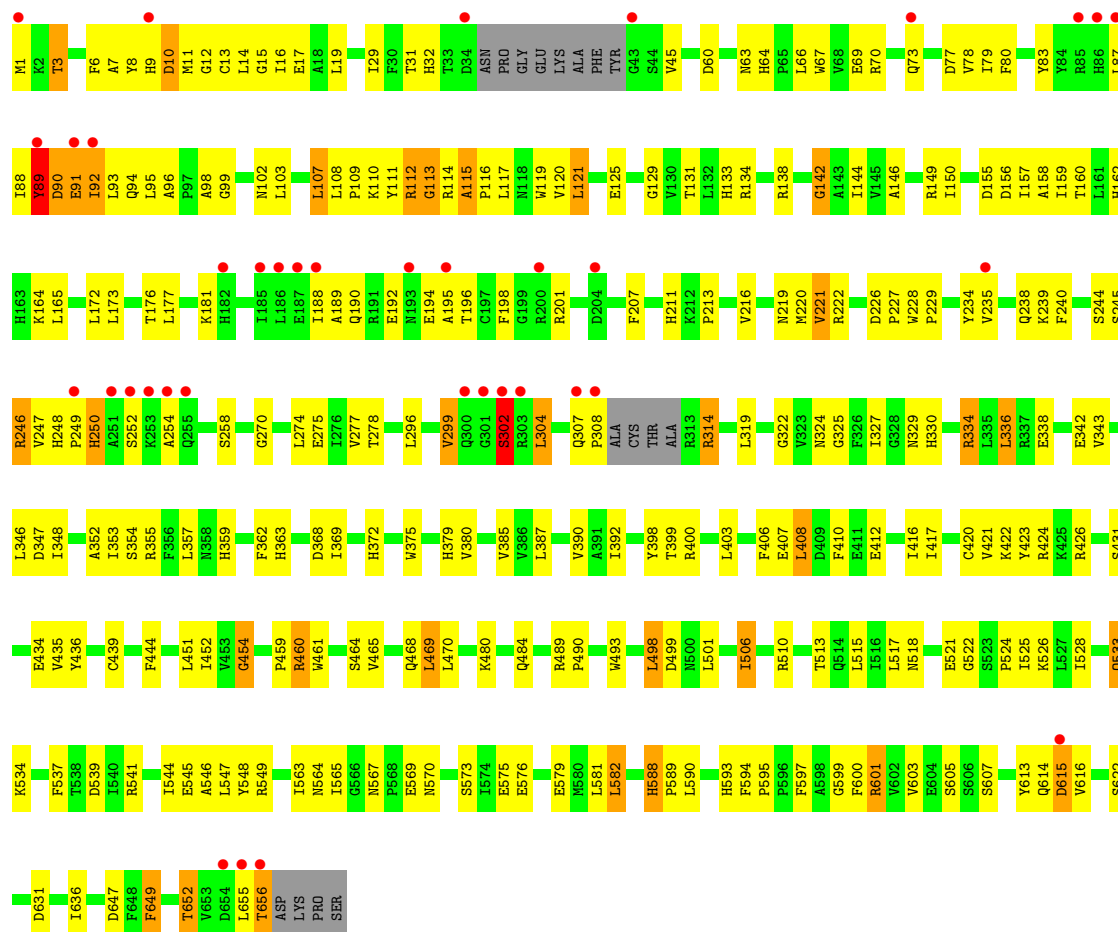


• Molecule 1: protein ArnA

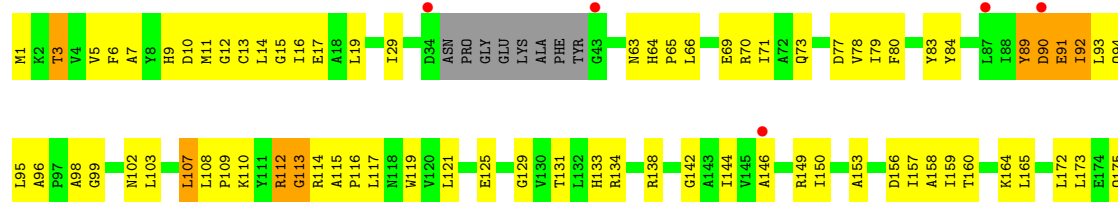


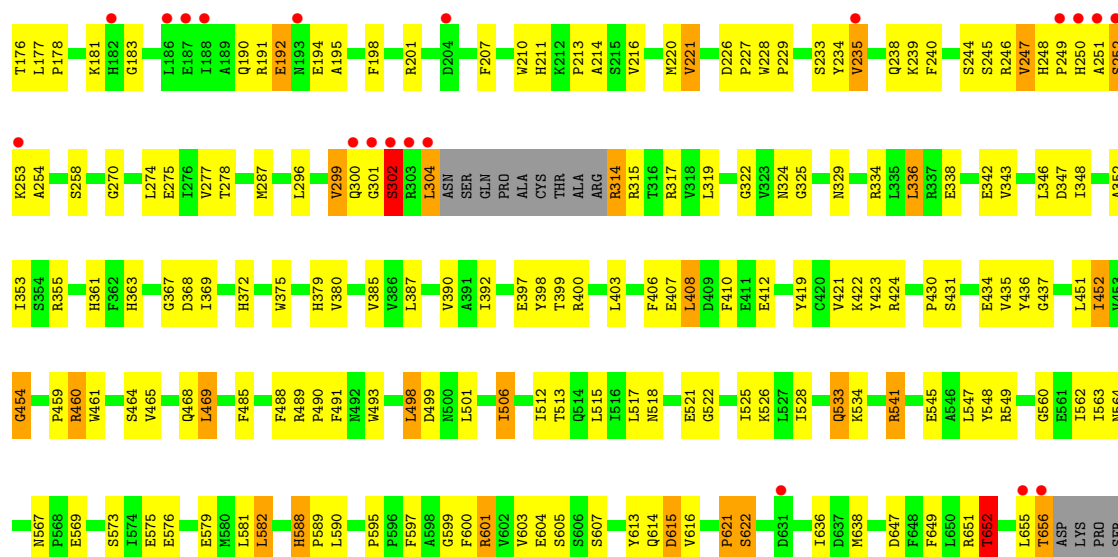


• Molecule 1: protein ArnA

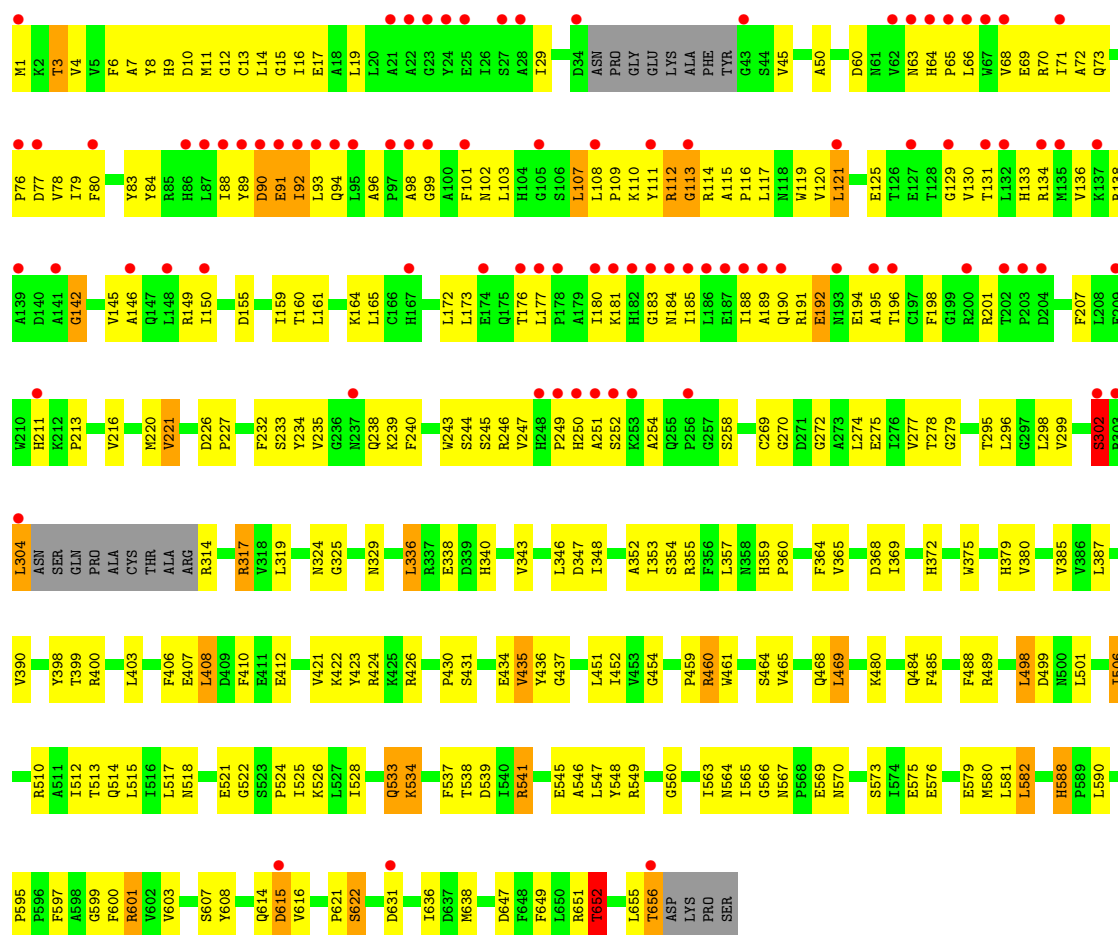


• Molecule 1: protein ArnA





• Molecule 1: protein ArnA



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	151.69Å 166.23Å 261.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.00 50.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	95.2 (50.00-3.00) 95.2 (50.00-3.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.77 (at 3.01Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.225 , 0.247 0.222 , 0.242	Depositor DCC
R_{free} test set	12667 reflections (9.56%)	wwPDB-VP
Wilson B-factor (Å ²)	47.0	Xtriage
Anisotropy	0.502	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	30556	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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: ATP, UGA

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	0.50	0/5143	0.97	21/6992 (0.3%)
1	B	0.52	0/5143	0.97	18/6992 (0.3%)
1	C	0.53	0/5143	0.98	19/6992 (0.3%)
1	D	0.50	0/5172	0.98	21/7033 (0.3%)
1	E	0.49	0/5143	0.97	21/6992 (0.3%)
1	F	0.52	0/5143	0.96	15/6992 (0.2%)
All	All	0.51	0/30887	0.97	115/41993 (0.3%)

There are no bond length outliers.

The worst 5 of 115 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	380	VAL	N-CA-C	-8.57	102.07	110.72
1	A	380	VAL	N-CA-C	-8.53	102.11	110.72
1	F	380	VAL	N-CA-C	-8.21	102.42	110.72
1	C	380	VAL	N-CA-C	-8.19	102.45	110.72
1	B	380	VAL	N-CA-C	-7.75	102.89	110.72

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5020	0	4924	218	0
1	B	5020	0	4924	214	2
1	C	5020	0	4924	210	2
1	D	5048	0	4942	209	0
1	E	5020	0	4924	215	0
1	F	5020	0	4924	226	0
2	A	31	0	12	5	0
2	B	31	0	12	4	0
2	C	31	0	12	4	0
2	D	31	0	12	4	0
2	E	31	0	12	2	0
2	F	31	0	12	3	0
3	A	37	0	19	1	0
3	B	37	0	19	1	0
3	C	37	0	19	1	0
3	D	37	0	19	1	0
3	E	37	0	19	1	0
3	F	37	0	19	1	0
All	All	30556	0	29748	1266	2

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 21.

The worst 5 of 1266 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:E:533:GLN:H	1:E:533:GLN:NE2	1.56	1.03
1:A:533:GLN:H	1:A:533:GLN:NE2	1.58	1.01
1:C:533:GLN:H	1:C:533:GLN:NE2	1.59	1.00
1:F:533:GLN:H	1:F:533:GLN:NE2	1.61	0.97
1:D:533:GLN:H	1:D:533:GLN:NE2	1.61	0.97

All (2) 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:B:184:ASN:ND2	1:C:171:GLN:OE1[1_655]	2.03	0.17
1:B:182:HIS:O	1:C:167:HIS:NE2[1_655]	2.13	0.07

5.3 Torsion angles

5.3.1 Protein backbone

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	633/660 (96%)	581 (92%)	42 (7%)	10 (2%)	7	34
1	B	633/660 (96%)	578 (91%)	43 (7%)	12 (2%)	6	30
1	C	633/660 (96%)	573 (90%)	50 (8%)	10 (2%)	7	34
1	D	638/660 (97%)	584 (92%)	44 (7%)	10 (2%)	7	34
1	E	633/660 (96%)	580 (92%)	42 (7%)	11 (2%)	7	32
1	F	633/660 (96%)	577 (91%)	45 (7%)	11 (2%)	7	32
All	All	3803/3960 (96%)	3473 (91%)	266 (7%)	64 (2%)	7	32

5 of 64 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	90	ASP
1	A	300	GLN
1	A	302	SER
1	B	90	ASP
1	B	251	ALA

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	526/559 (94%)	491 (93%)	35 (7%)	15	46
1	B	526/559 (94%)	490 (93%)	36 (7%)	14	45
1	C	526/559 (94%)	489 (93%)	37 (7%)	14	44

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	528/559 (94%)	493 (93%)	35 (7%)	15	46
1	E	526/559 (94%)	491 (93%)	35 (7%)	15	46
1	F	526/559 (94%)	488 (93%)	38 (7%)	13	43
All	All	3158/3354 (94%)	2942 (93%)	216 (7%)	14	45

5 of 216 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	121	LEU
1	E	91	GLU
1	F	435	VAL
1	D	247	VAL
1	D	460	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 137 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	147	GLN
1	F	238	GLN
1	F	518	ASN
1	C	147	GLN
1	C	133	HIS

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

12 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
3	UGA	A	1101	-	38,39,39	0.89	2 (5%)	56,60,60	0.95	3 (5%)
2	ATP	C	1003	-	32,33,33	0.90	2 (6%)	48,52,52	0.98	3 (6%)
3	UGA	E	1105	-	38,39,39	0.88	3 (7%)	56,60,60	0.97	2 (3%)
2	ATP	D	1004	-	32,33,33	1.08	3 (9%)	48,52,52	0.98	2 (4%)
3	UGA	F	1106	-	38,39,39	0.97	3 (7%)	56,60,60	0.99	2 (3%)
3	UGA	C	1103	-	38,39,39	0.85	2 (5%)	56,60,60	0.98	2 (3%)
2	ATP	E	1005	-	32,33,33	0.74	1 (3%)	48,52,52	0.99	2 (4%)
3	UGA	B	1102	-	38,39,39	0.91	3 (7%)	56,60,60	1.00	3 (5%)
2	ATP	F	1006	-	32,33,33	0.97	2 (6%)	48,52,52	0.99	3 (6%)
2	ATP	A	1001	-	32,33,33	0.84	1 (3%)	48,52,52	1.03	2 (4%)
3	UGA	D	1104	-	38,39,39	0.87	2 (5%)	56,60,60	0.93	3 (5%)
2	ATP	B	1002	-	32,33,33	0.83	1 (3%)	48,52,52	0.99	2 (4%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	UGA	A	1101	-	-	4/25/61/61	0/3/3/3
2	ATP	C	1003	-	-	4/22/38/38	0/3/3/3
3	UGA	E	1105	-	-	4/25/61/61	0/3/3/3
2	ATP	D	1004	-	-	4/22/38/38	0/3/3/3
3	UGA	F	1106	-	-	4/25/61/61	0/3/3/3
3	UGA	C	1103	-	-	4/25/61/61	0/3/3/3
2	ATP	E	1005	-	-	4/22/38/38	0/3/3/3
3	UGA	B	1102	-	-	4/25/61/61	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ATP	F	1006	-	-	4/22/38/38	0/3/3/3
2	ATP	A	1001	-	-	4/22/38/38	0/3/3/3
3	UGA	D	1104	-	-	4/25/61/61	0/3/3/3
2	ATP	B	1002	-	-	4/22/38/38	0/3/3/3

The worst 5 of 25 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	1006	ATP	PB-O3B	3.91	1.63	1.59
2	D	1004	ATP	PB-O3B	3.84	1.63	1.59
2	C	1003	ATP	PB-O3B	3.65	1.63	1.59
2	A	1001	ATP	PB-O3B	3.43	1.63	1.59
2	B	1002	ATP	PB-O3B	3.37	1.63	1.59

The worst 5 of 29 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1102	UGA	O5'-C1'-O3B	4.86	117.71	111.36
3	F	1106	UGA	O5'-C1'-O3B	4.83	117.68	111.36
3	A	1101	UGA	O5'-C1'-O3B	4.80	117.64	111.36
3	D	1104	UGA	O5'-C1'-O3B	4.61	117.39	111.36
3	C	1103	UGA	O5'-C1'-O3B	4.60	117.38	111.36

There are no chirality outliers.

5 of 48 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1001	ATP	C5'-O5'-PA-O2A
2	A	1001	ATP	C5'-O5'-PA-O3A
2	B	1002	ATP	C5'-O5'-PA-O2A
2	C	1003	ATP	C5'-O5'-PA-O2A
2	D	1004	ATP	C5'-O5'-PA-O2A

There are no ring outliers.

12 monomers are involved in 28 short contacts:

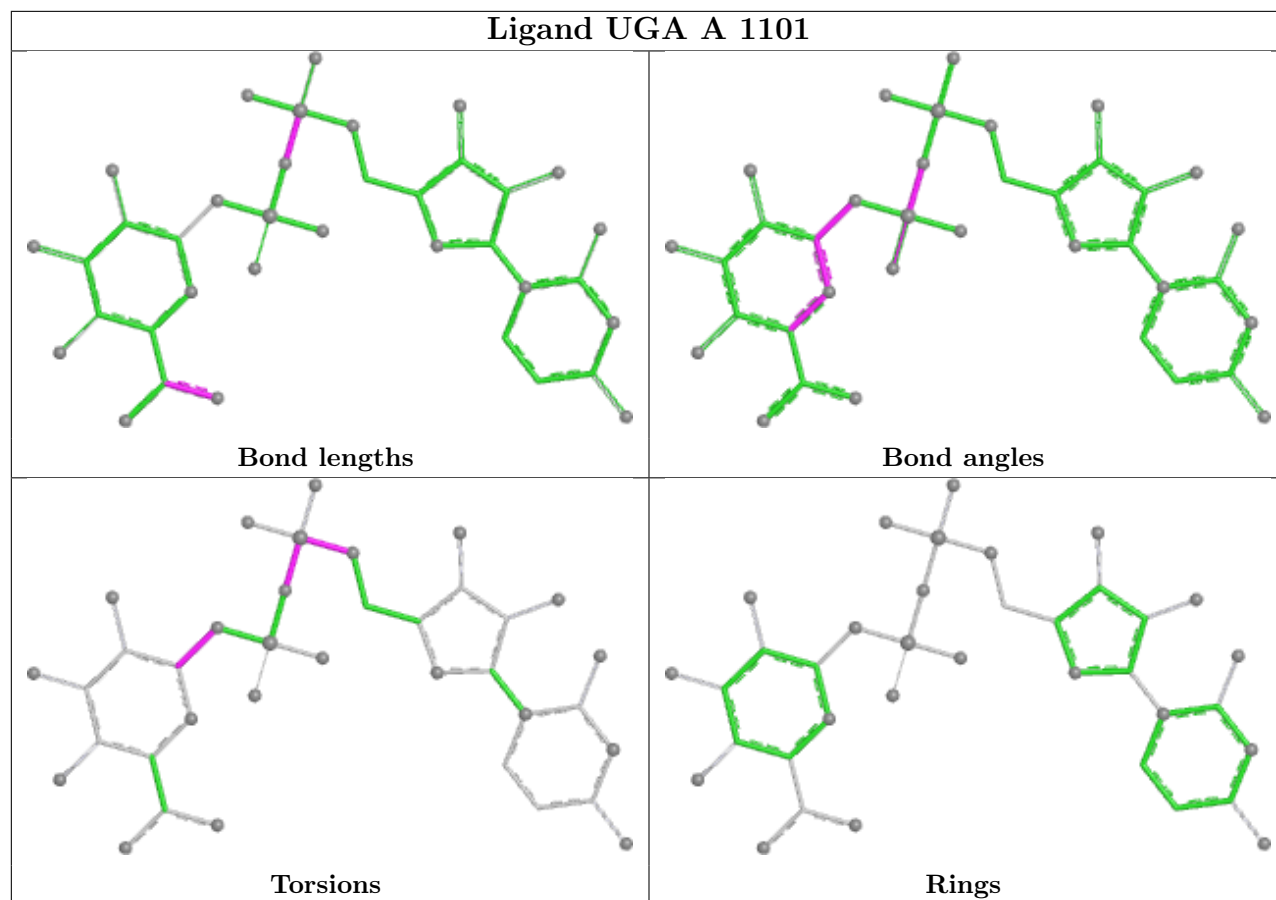
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1101	UGA	1	0
2	C	1003	ATP	4	0
3	E	1105	UGA	1	0

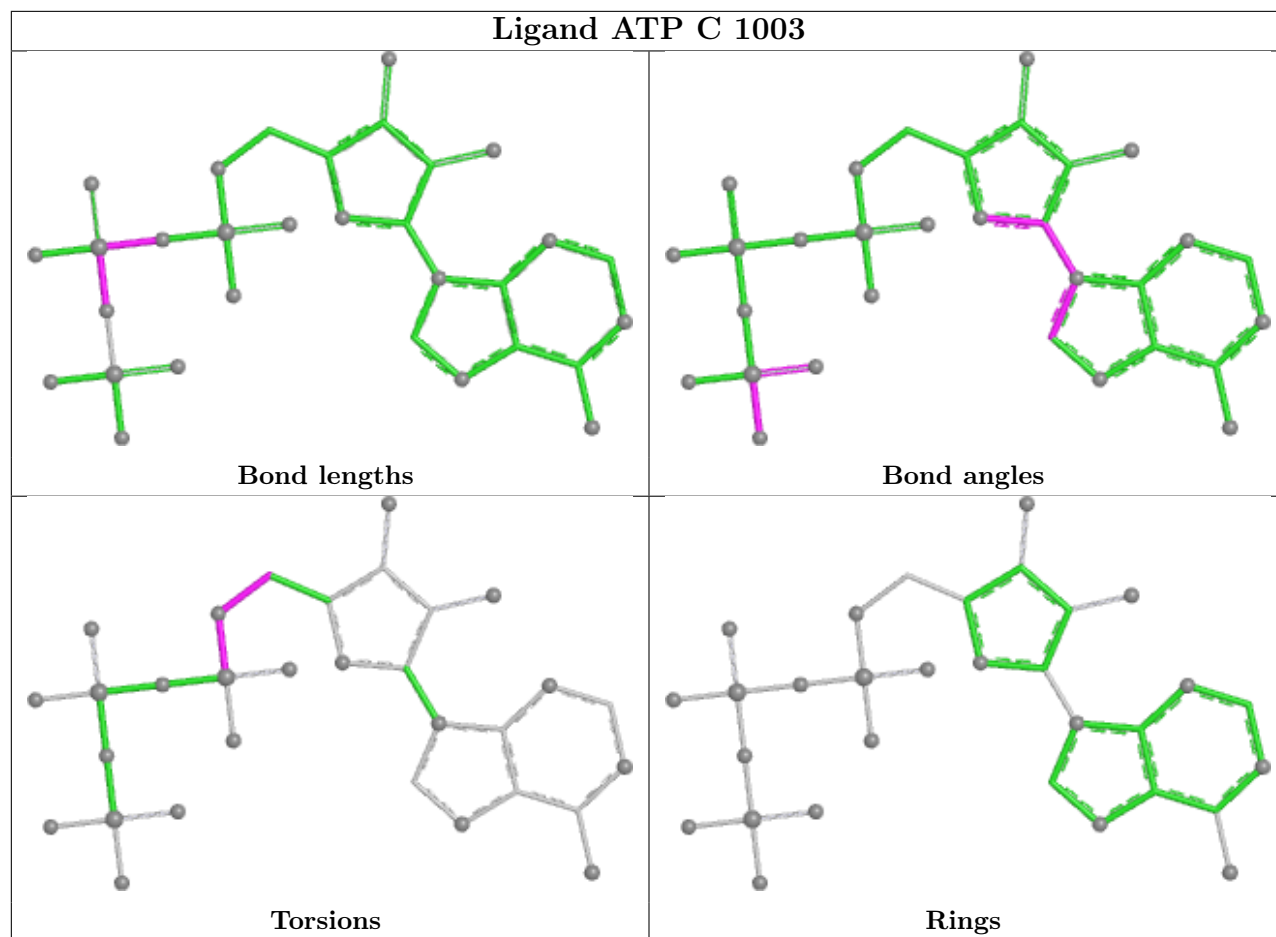
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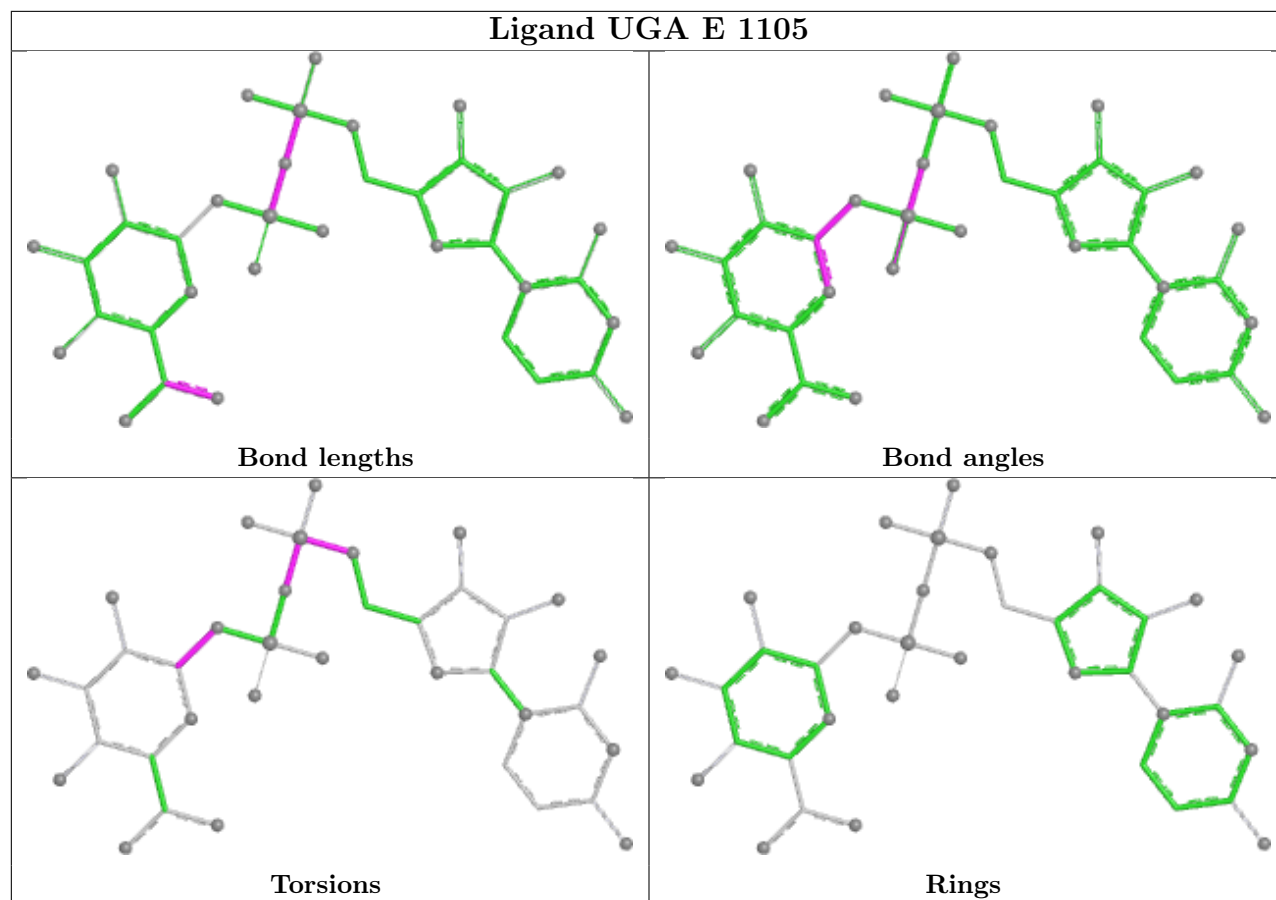
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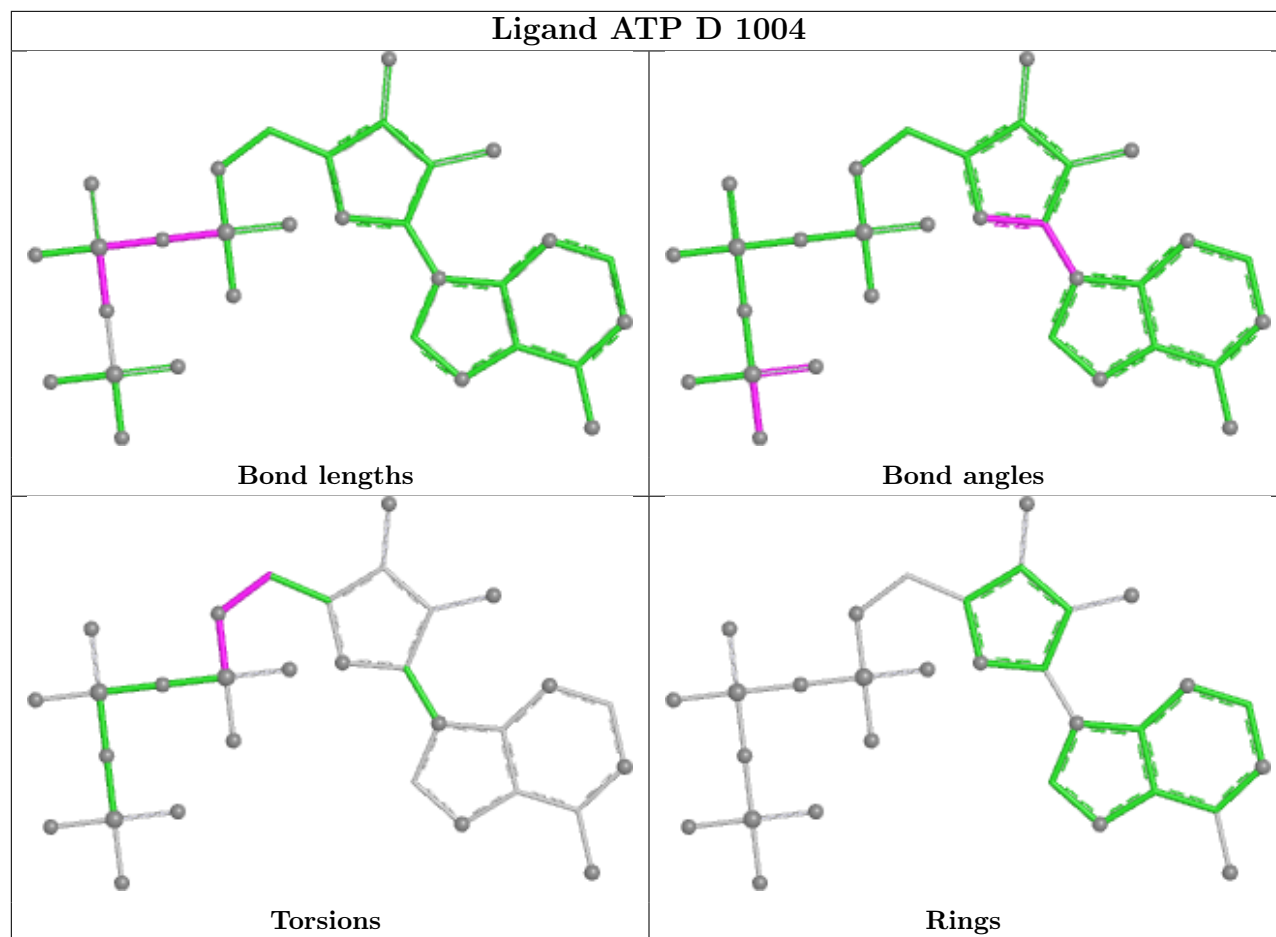
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1004	ATP	4	0
3	F	1106	UGA	1	0
3	C	1103	UGA	1	0
2	E	1005	ATP	2	0
3	B	1102	UGA	1	0
2	F	1006	ATP	3	0
2	A	1001	ATP	5	0
3	D	1104	UGA	1	0
2	B	1002	ATP	4	0

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.

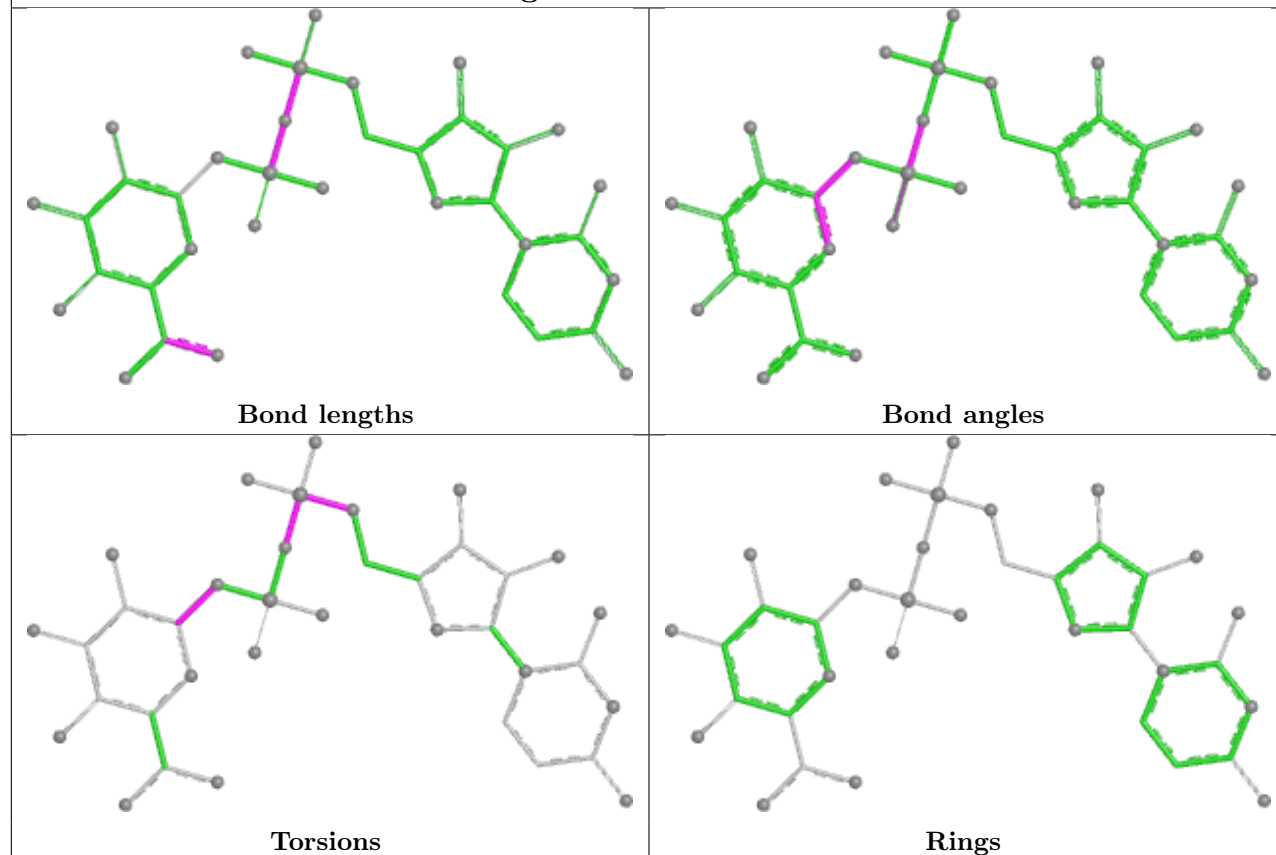




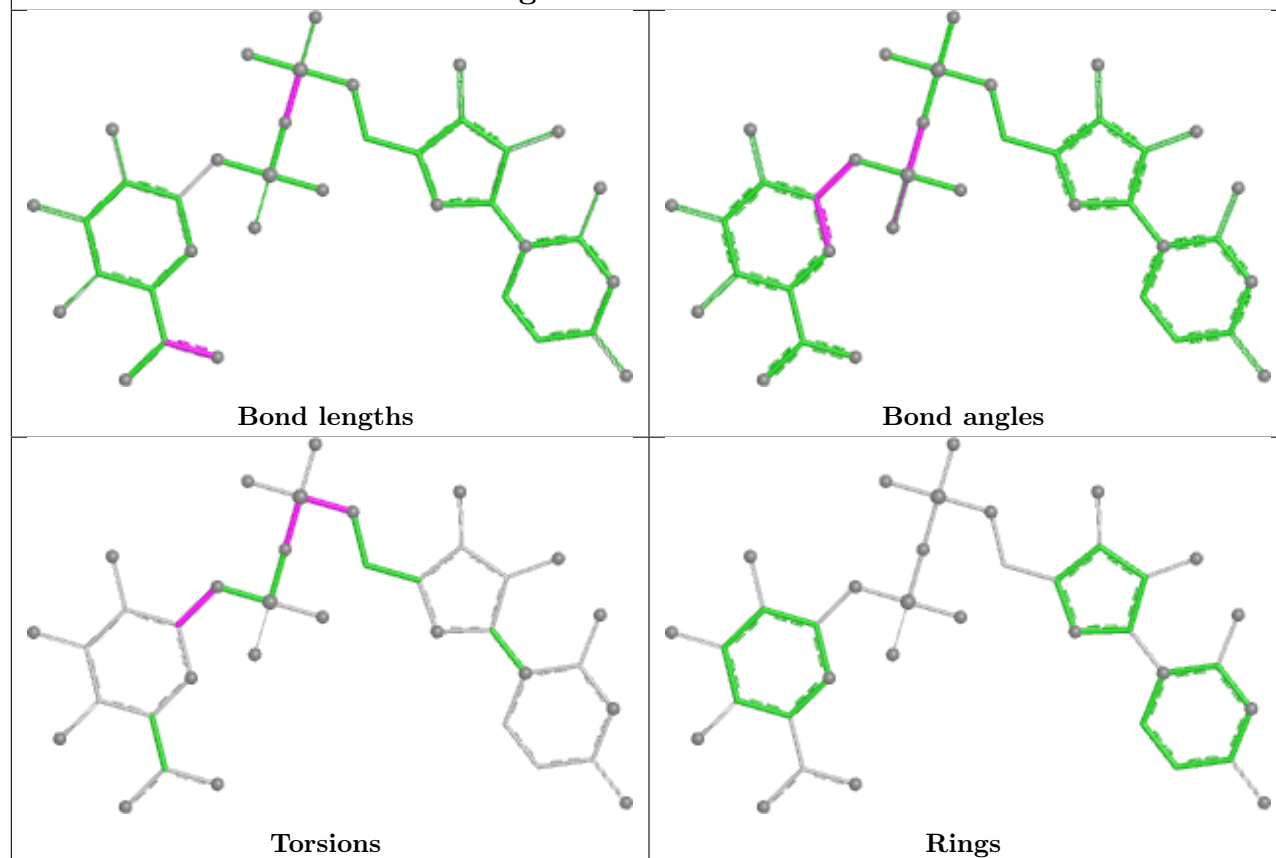


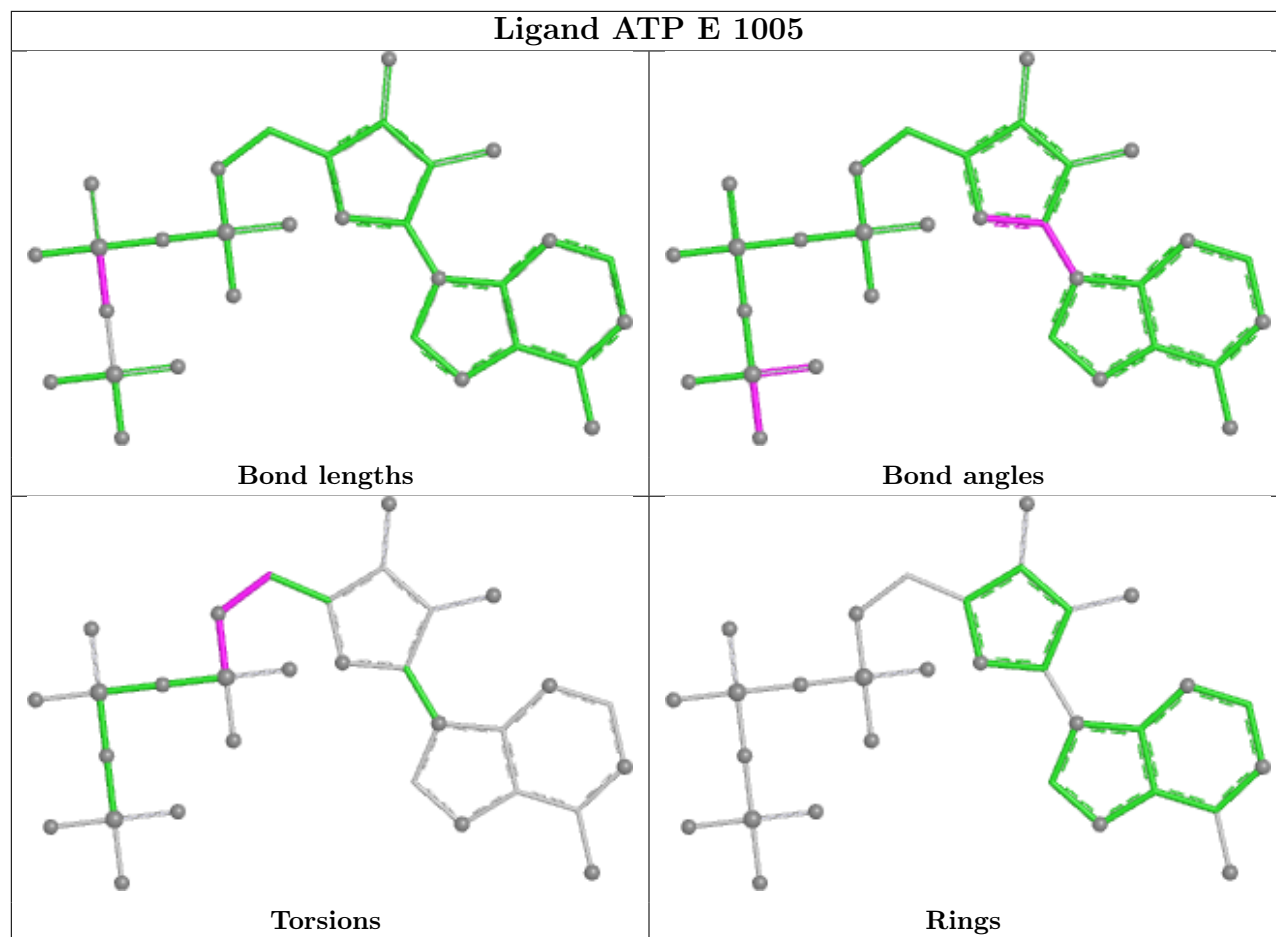


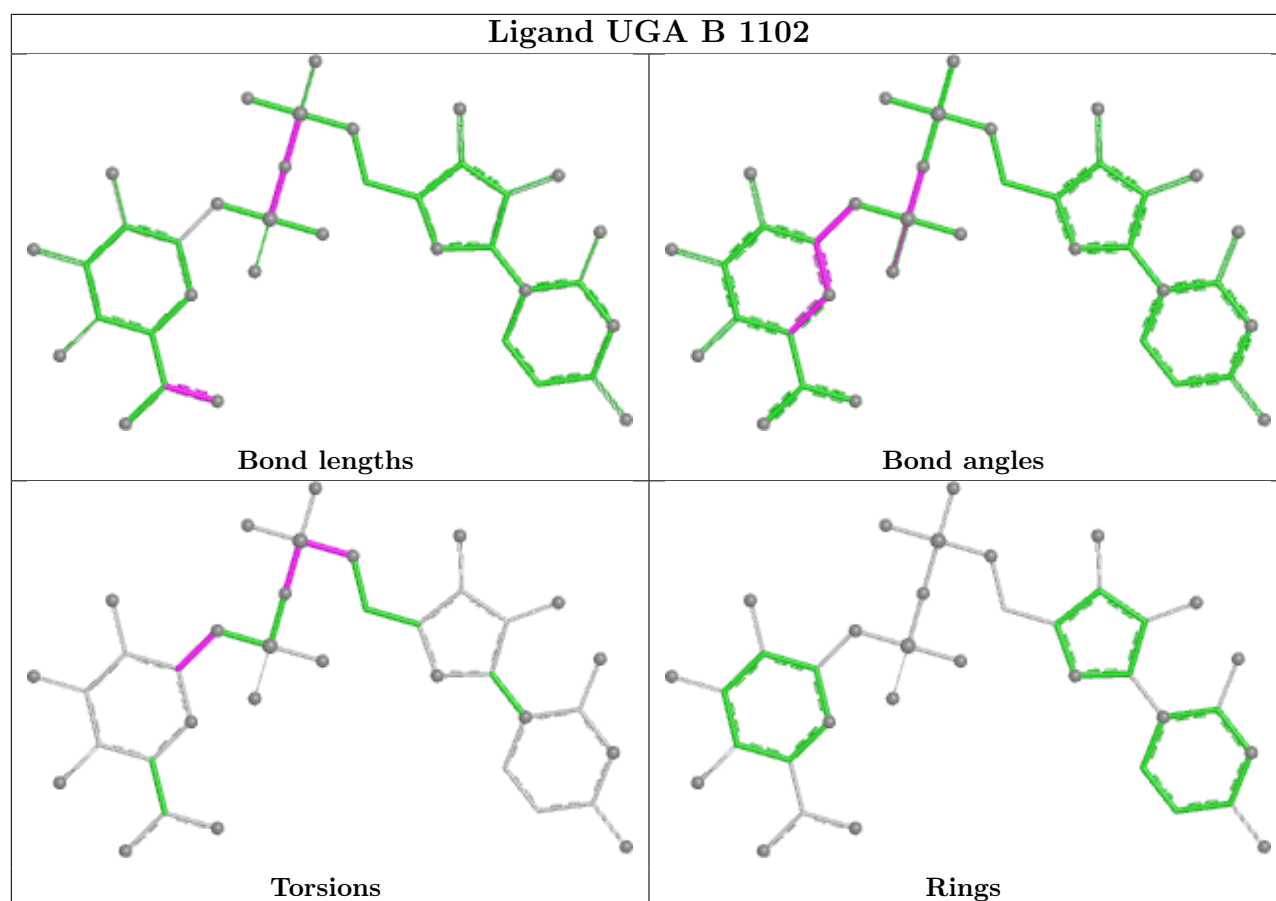
Ligand UGA F 1106

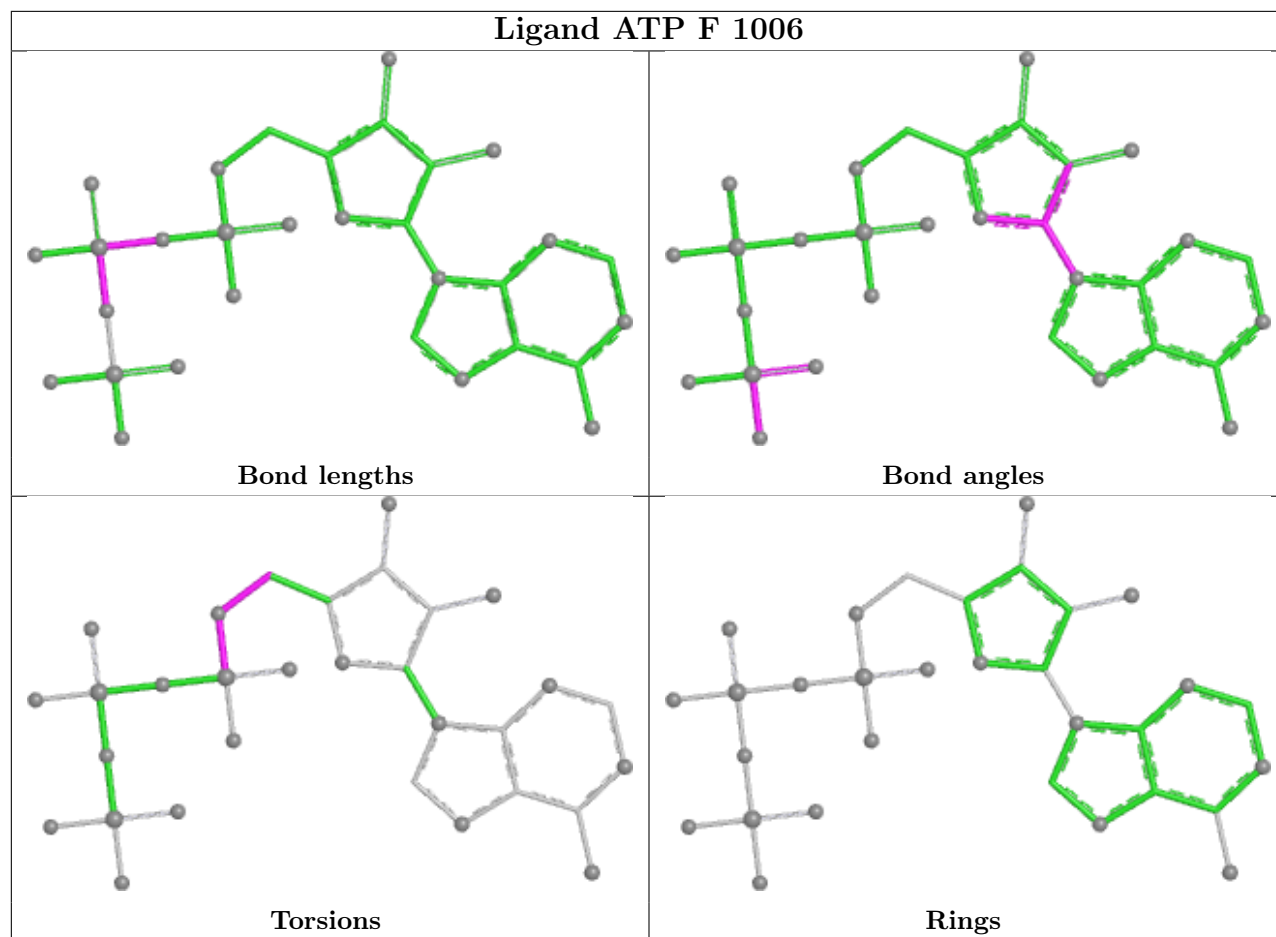


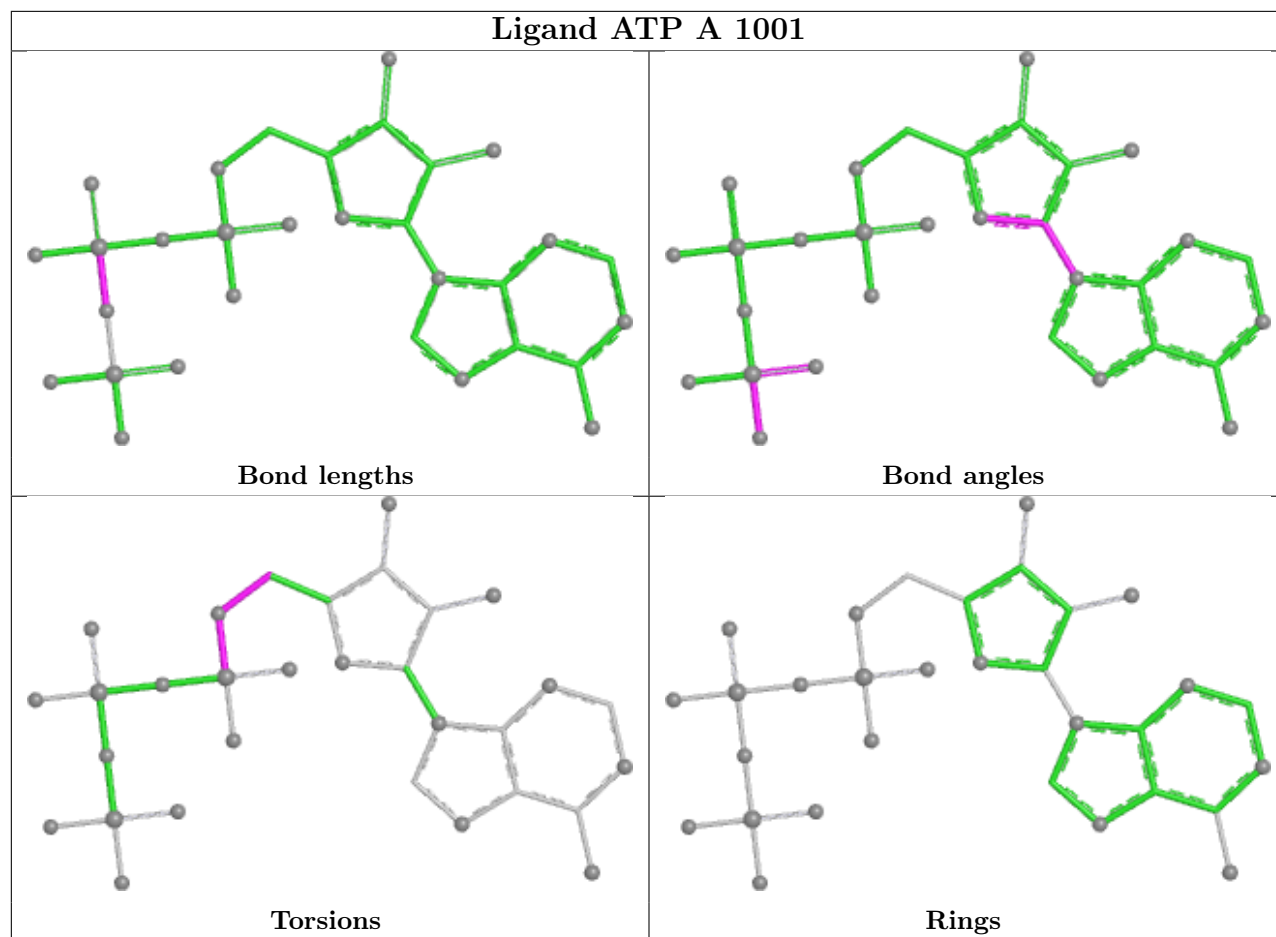
Ligand UGA C 1103



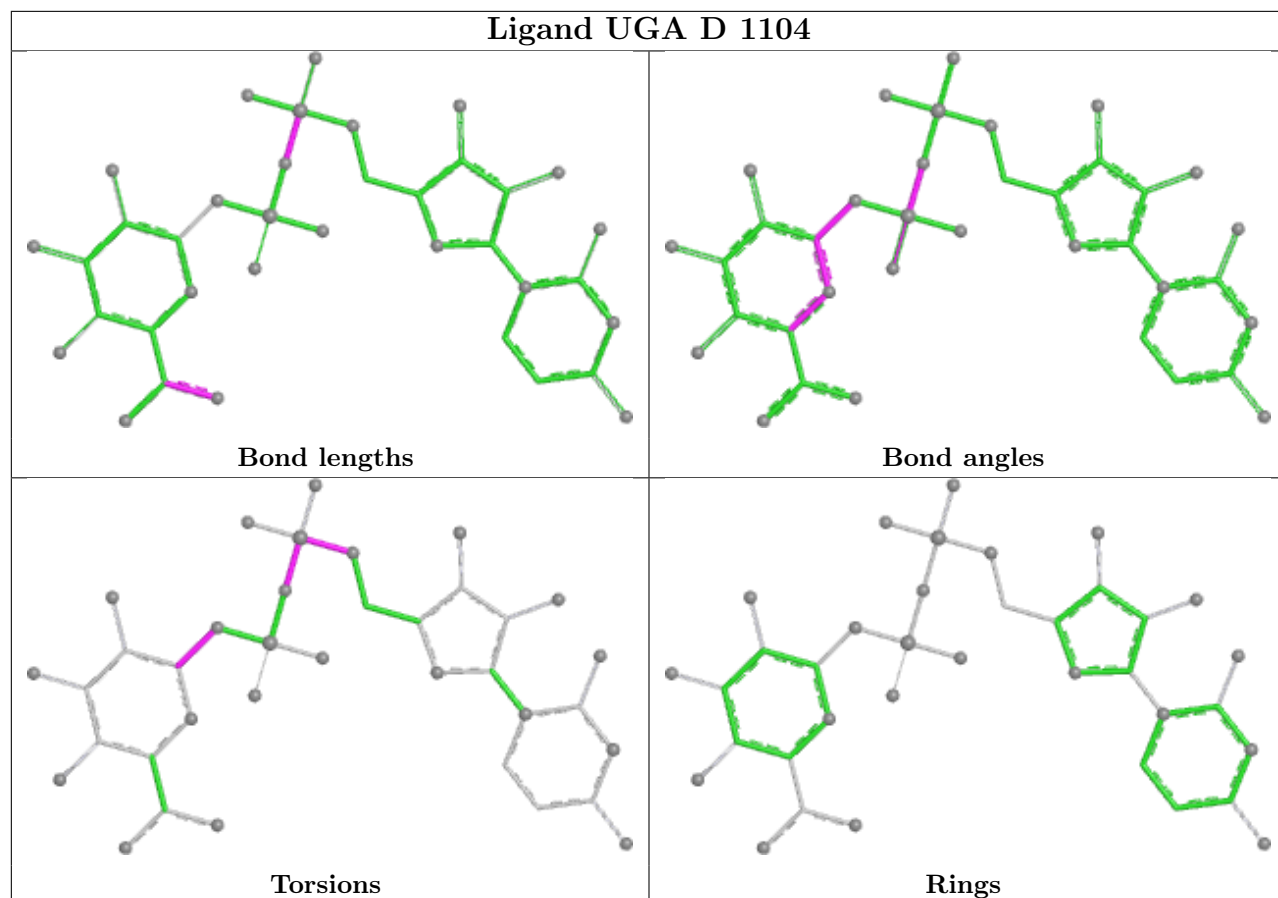


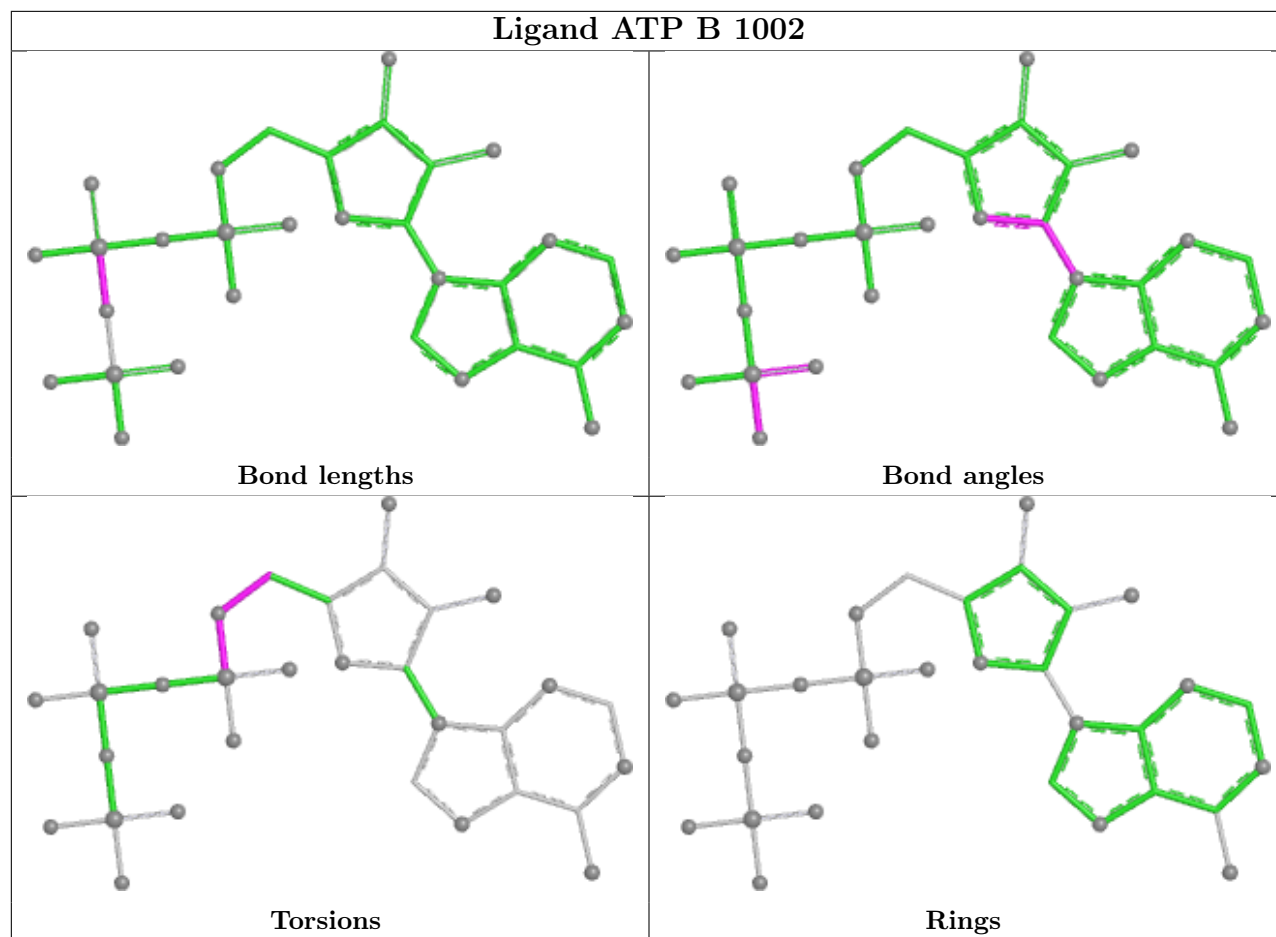






Ligand UGA D 1104





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	639/660 (96%)	0.23	32 (5%)	34	18	20, 43, 74, 90	0
1	B	639/660 (96%)	0.27	43 (6%)	24	12	19, 43, 73, 90	0
1	C	639/660 (96%)	0.33	39 (6%)	27	14	20, 43, 73, 90	0
1	D	644/660 (97%)	0.22	37 (5%)	29	15	20, 44, 74, 90	0
1	E	639/660 (96%)	0.16	25 (3%)	43	24	20, 44, 73, 90	0
1	F	639/660 (96%)	0.51	91 (14%)	6	4	19, 43, 75, 91	0
All	All	3839/3960 (96%)	0.29	267 (6%)	22	12	19, 43, 74, 91	0

The worst 5 of 267 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	656	THR	6.7
1	B	184	ASN	5.8
1	C	187	GLU	5.7
1	E	253	LYS	5.4
1	D	656	THR	5.3

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands

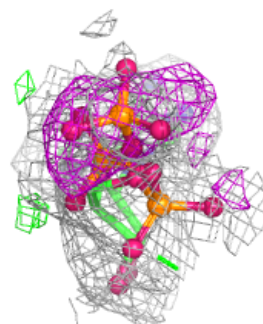
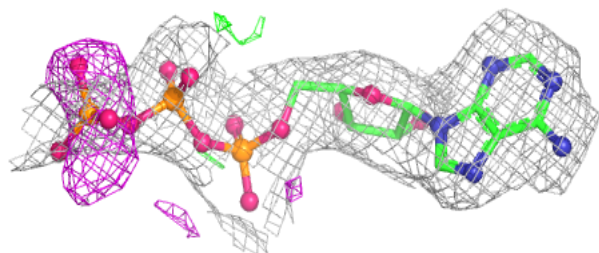
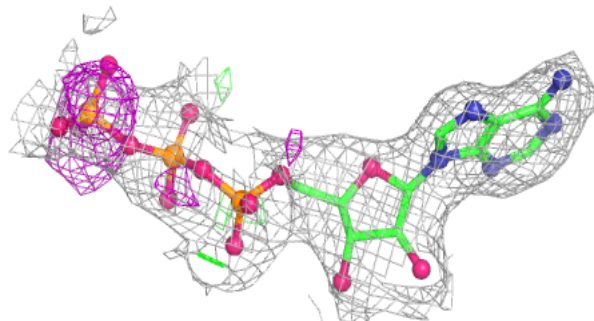
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ATP	D	1004	31/31	0.89	0.13	30,40,81,82	0
2	ATP	C	1003	31/31	0.90	0.13	32,41,82,82	0
3	UGA	D	1104	37/37	0.90	0.15	31,36,61,62	0
2	ATP	B	1002	31/31	0.91	0.13	30,39,81,82	0
2	ATP	F	1006	31/31	0.91	0.14	27,40,81,82	0
3	UGA	C	1103	37/37	0.91	0.13	31,38,61,62	0
2	ATP	A	1001	31/31	0.91	0.13	28,39,82,83	0
3	UGA	F	1106	37/37	0.91	0.14	30,36,61,61	0
3	UGA	E	1105	37/37	0.92	0.14	30,36,60,61	0
2	ATP	E	1005	31/31	0.92	0.13	30,42,83,83	0
3	UGA	A	1101	37/37	0.93	0.15	27,36,61,61	0
3	UGA	B	1102	37/37	0.93	0.14	27,36,59,60	0

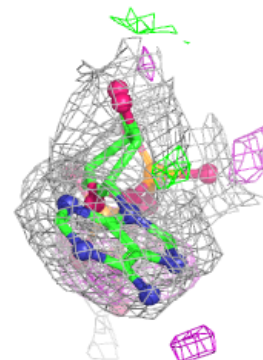
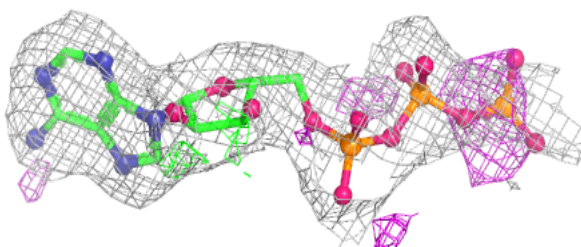
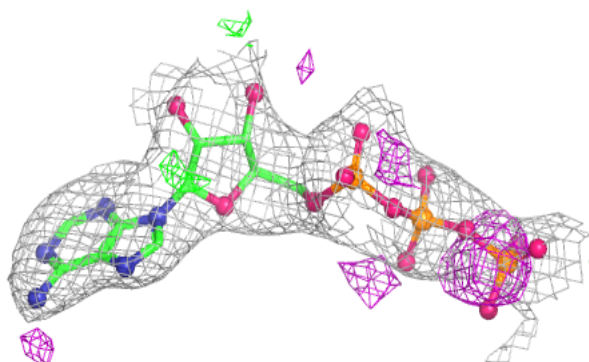
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ATP D 1004:

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

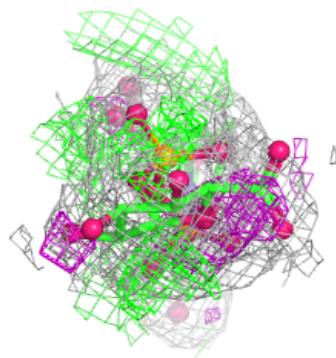
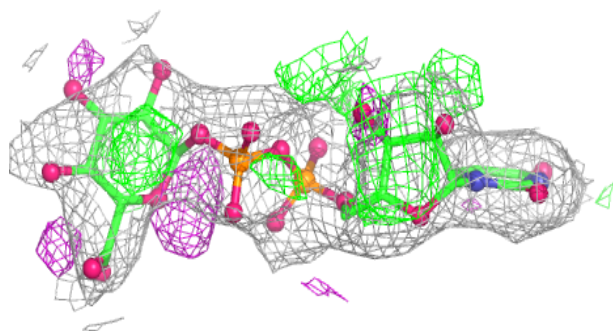
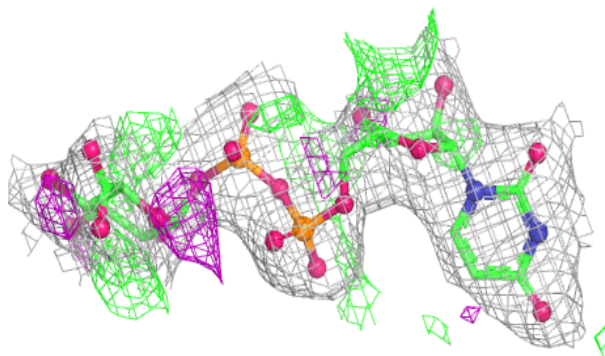
**Electron density around ATP C 1003:**

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

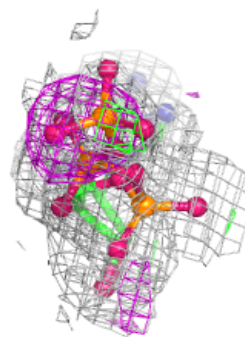
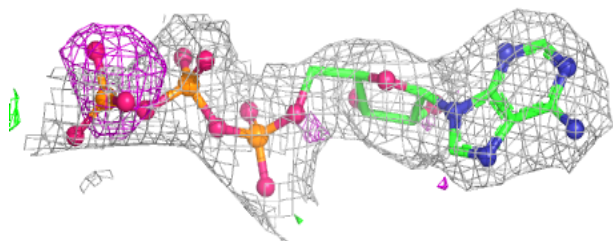
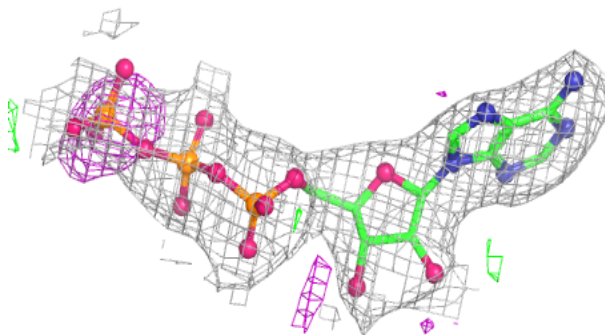


Electron density around UGA D 1104:

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

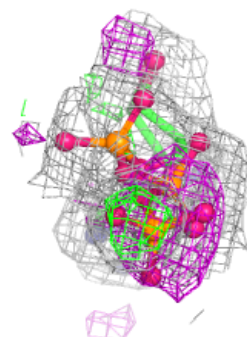
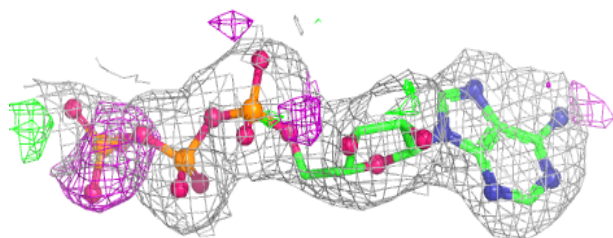
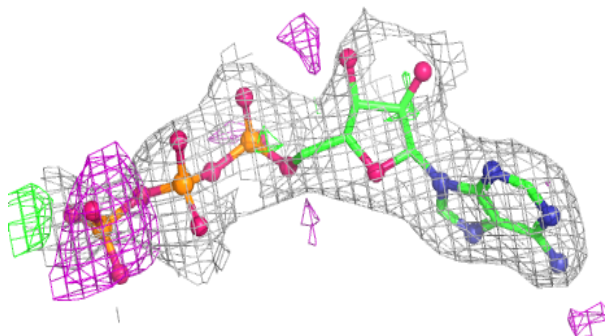
**Electron density around ATP B 1002:**

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

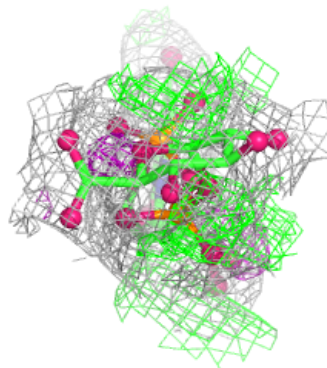
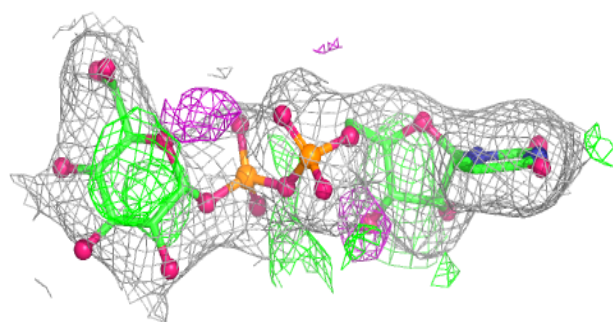
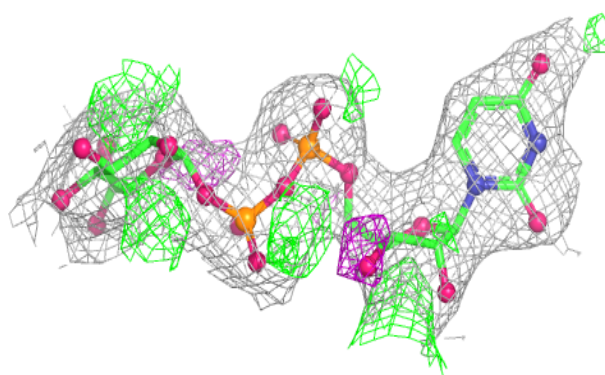


Electron density around ATP F 1006:

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

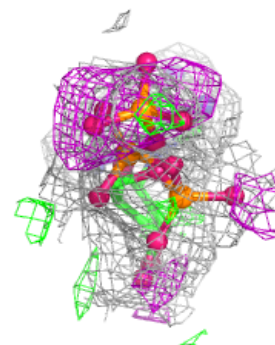
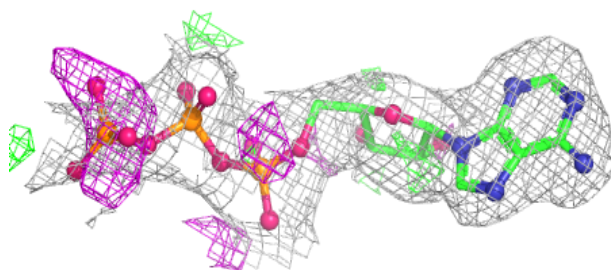
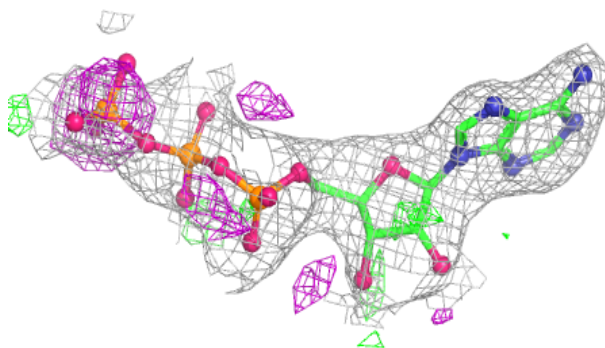
**Electron density around UGA C 1103:**

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

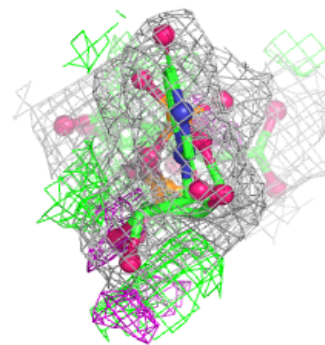
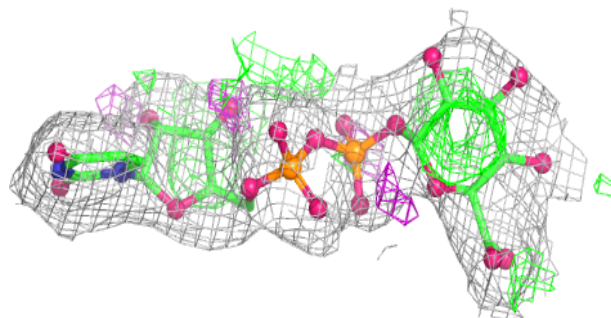
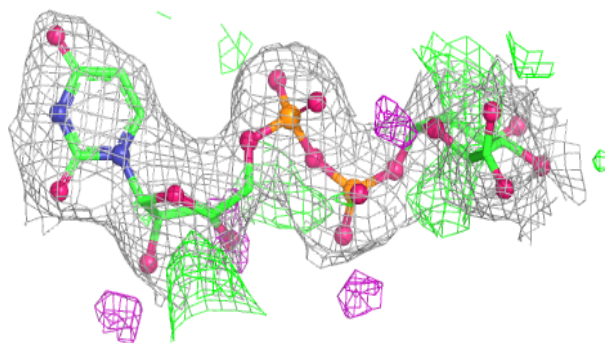


Electron density around ATP A 1001:

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

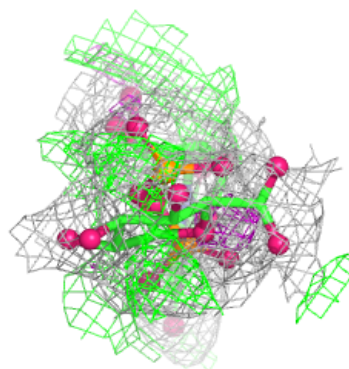
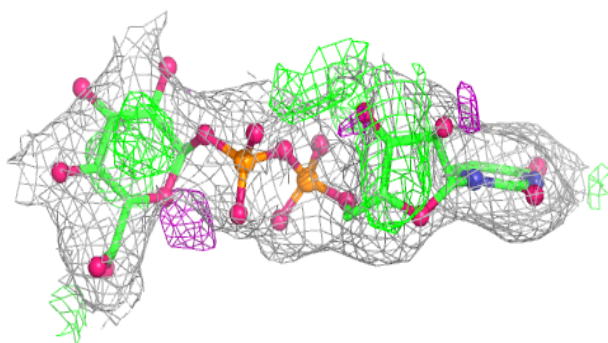
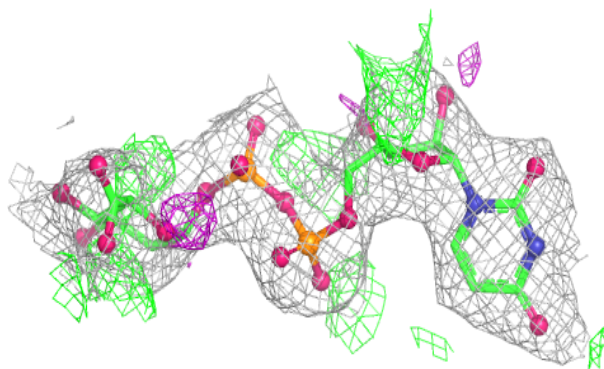
**Electron density around UGA F 1106:**

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

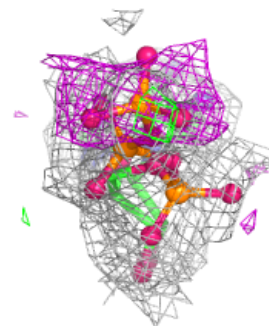
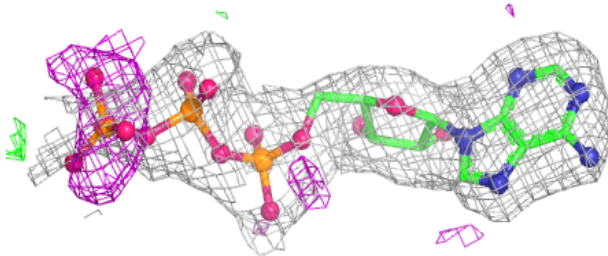
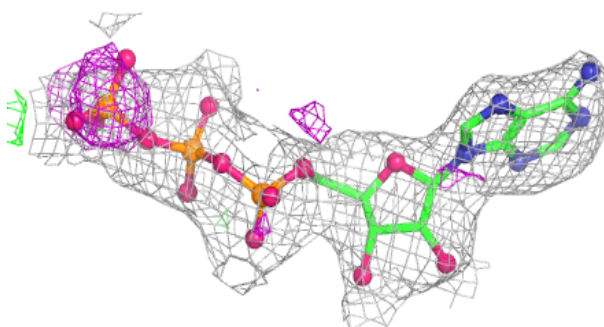


Electron density around UGA E 1105:

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

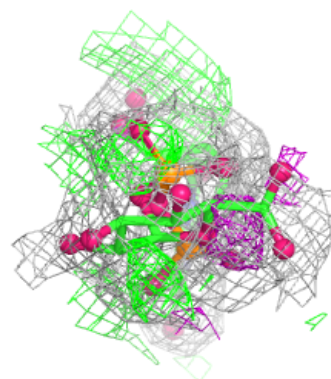
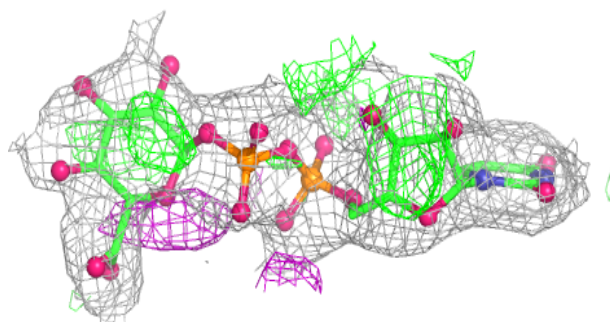
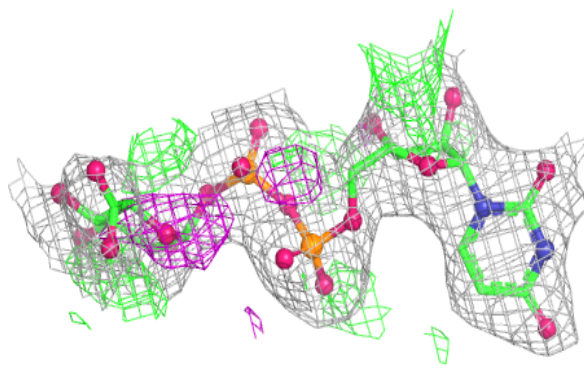
**Electron density around ATP E 1005:**

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

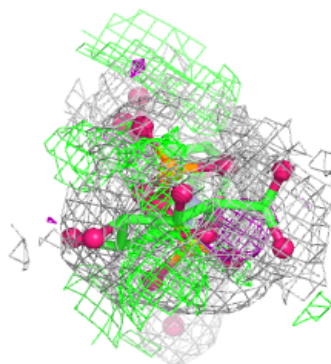
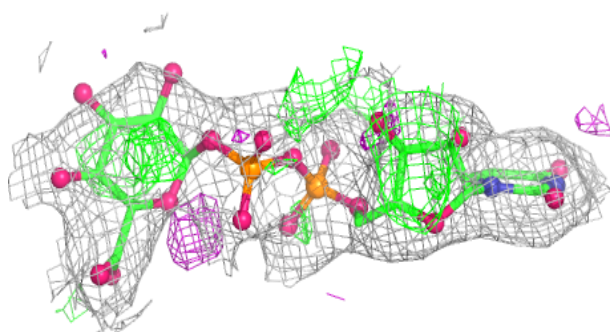
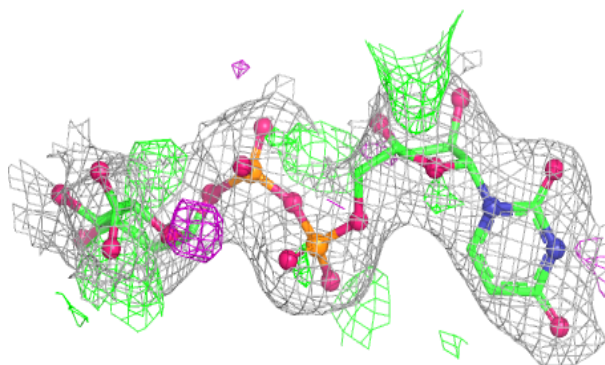


Electron density around UGA A 1101:

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

**Electron density around UGA B 1102:**

$2mF_o-DF_c$ (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.