



wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 28, 2022 – 02:10 PM EDT

PDB ID : 7S01
Title : X-ray structure of the phage AR9 non-virion RNA polymerase holoenzyme in complex with a forked oligonucleotide containing the P077 promoter
Authors : Leiman, P.G.; Sokolova, M.L.; Gordeeva, J.; Fraser, A.; Severinov, K.V.
Deposited on : 2021-08-28
Resolution : 3.40 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 : 2.29
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) : 2.29

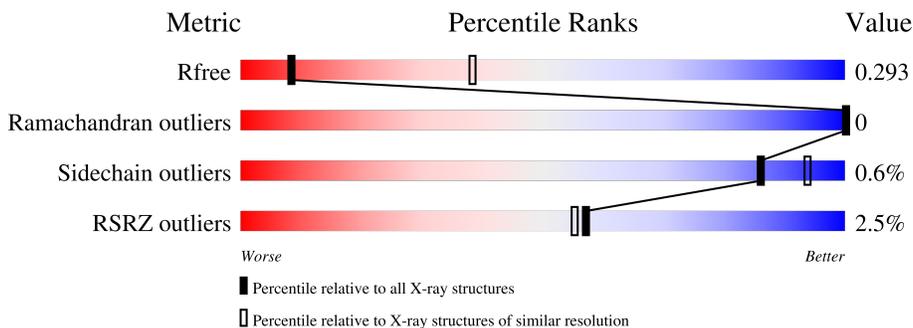
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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.



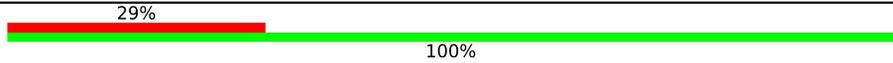
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1026 (3.48-3.32)
Ramachandran outliers	138981	1038 (3.48-3.32)
Sidechain outliers	138945	1038 (3.48-3.32)
RSRZ outliers	127900	2173 (3.50-3.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	464	5% (poor fit), 99% (0-1 outliers), 0% (2-3 outliers), 0% (not modelled)
2	d	426	0% (poor fit), 99% (0-1 outliers), 0% (2-3 outliers), 0% (not modelled)
3	c	496	4% (poor fit), 97% (0-1 outliers), 0% (2-3 outliers), 0% (not modelled)
4	D	631	0% (poor fit), 99% (0-1 outliers), 0% (2-3 outliers), 0% (not modelled)
5	C	665	0% (poor fit), 97% (0-1 outliers), 0% (2-3 outliers), 0% (not modelled)
6	N	14	0% (poor fit), 93% (0-1 outliers), 7% (2-3 outliers), 0% (not modelled)

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Mol	Chain	Length	Quality of chain
6	t	14	 29% 100%
7	T	32	 12% 66% 34%
7	n	32	 3% 69% 31%

2 Entry composition [i](#)

There are 12 unique types of molecules in this entry. The entry contains 23345 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 DNA-directed RNA polymerase subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	463	3847	2478	620	729	20	0	0	0

- Molecule 2 is a protein called DNA-directed RNA polymerase beta' subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	d	426	3488	2256	561	663	8	0	0	0

- Molecule 3 is a protein called DNA-directed RNA polymerase beta subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	c	484	4003	2580	658	754	11	0	0	0

- Molecule 4 is a protein called DNA-directed RNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	D	631	5140	3276	837	1009	18	0	0	0

- Molecule 5 is a protein called DNA-directed RNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
5	C	649	5271	3337	868	1041	25	0	0	0

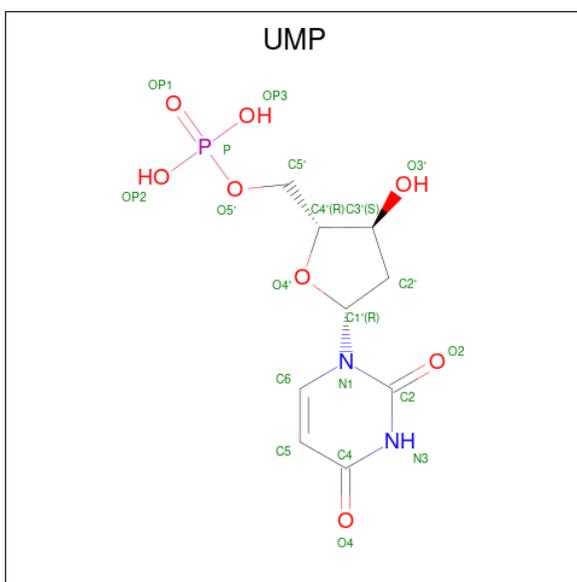
- Molecule 6 is a DNA chain called Non-template strand of the forked DNA oligonucleotide (downstream copy).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	N	13	Total	C	N	O	P	0	0	0
			267	128	49	77	13			
6	t	14	Total	C	N	O	P	0	0	0
			289	138	54	83	14			

- Molecule 7 is a DNA chain called Template strand of the forked DNA oligonucleotide (downstream copy) containing the P077 AR9 promoter motif.

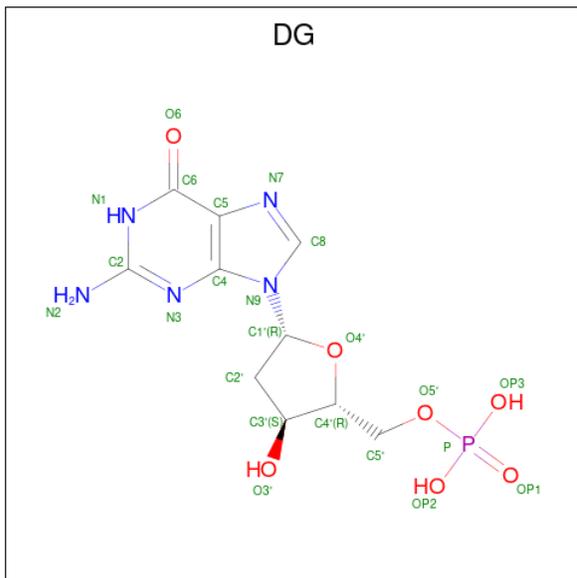
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	T	21	Total	C	N	O	P	0	0	0
			431	208	77	125	21			
7	n	22	Total	C	N	O	P	0	0	0
			450	217	80	131	22			

- Molecule 8 is 2'-DEOXYURIDINE 5'-MONOPHOSPHATE (three-letter code: UMP) (formula: $C_9H_{13}N_2O_8P$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	A	1	Total	C	N	O	P	0	0
			19	9	2	7	1		
8	A	1	Total	C	N	O	P	0	0
			19	9	2	7	1		
8	c	1	Total	C	N	O	P	0	0
			19	9	2	7	1		
8	c	1	Total	C	N	O	P	0	0
			19	9	2	7	1		
8	T	1	Total	C	N	O	P	0	0
			19	9	2	7	1		

- Molecule 9 is 2'-DEOXYGUANOSINE-5'-MONOPHOSPHATE (three-letter code: DG) (formula: $C_{10}H_{14}N_5O_7P$) (labeled as "Ligand of Interest" by depositor).

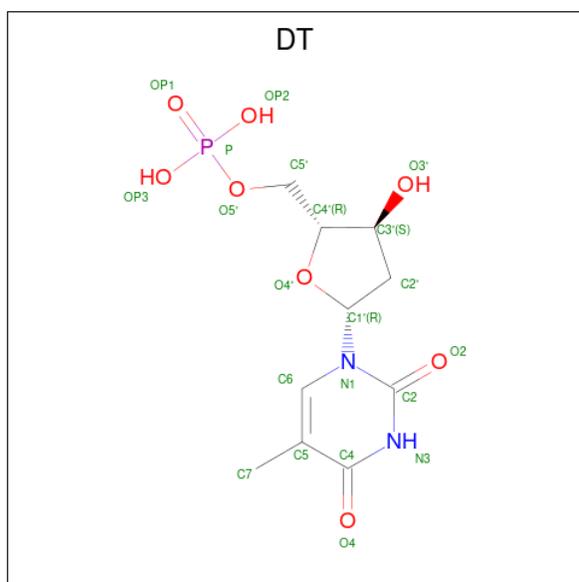


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
9	A	1	22	10	5	6	1	0	0

- Molecule 10 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

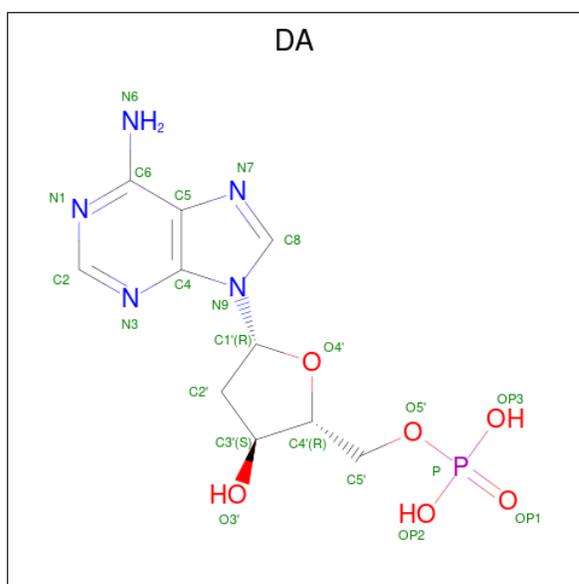
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
10	D	1	1	1	0	0

- Molecule 11 is THYMIDINE-5'-MONOPHOSPHATE (three-letter code: DT) (formula: $C_{10}H_{15}N_2O_8P$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
11	t	1	20	10	2	7	1	0	0

- Molecule 12 is 2'-DEOXYADENOSINE-5'-MONOPHOSPHATE (three-letter code: DA) (formula: $C_{10}H_{14}N_5O_6P$) (labeled as "Ligand of Interest" by depositor).

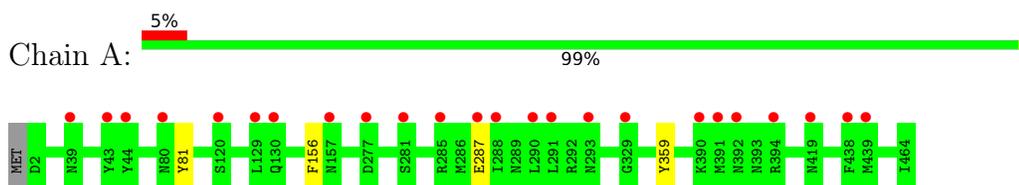


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
12	n	1	21	10	5	5	1	0	0

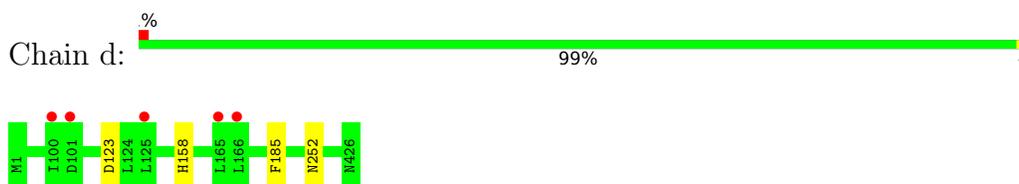
3 Residue-property plots [i](#)

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.

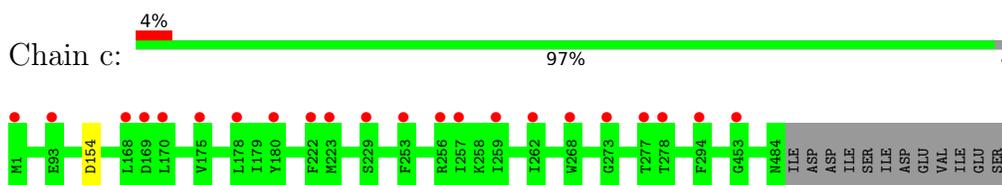
- Molecule 1: DNA-directed RNA polymerase subunit



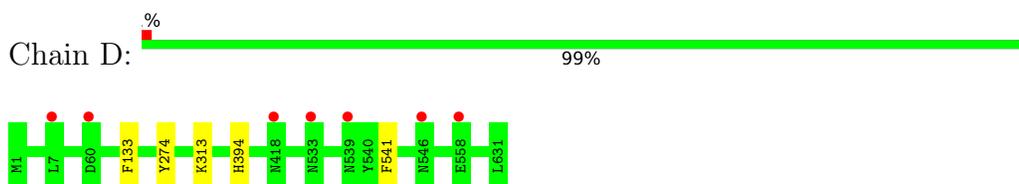
- Molecule 2: DNA-directed RNA polymerase beta' subunit



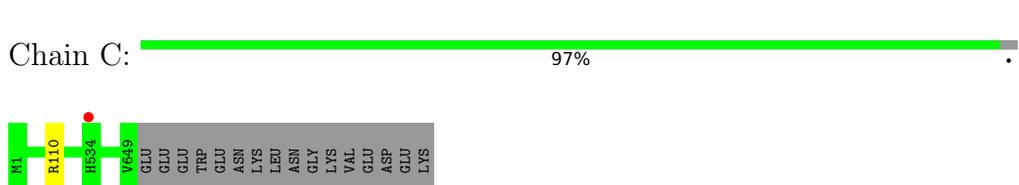
- Molecule 3: DNA-directed RNA polymerase beta subunit



- Molecule 4: DNA-directed RNA polymerase



- Molecule 5: DNA-directed RNA polymerase



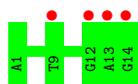
- Molecule 6: Non-template strand of the forked DNA oligonucleotide (downstream copy)

Chain N:  93% 7%



- Molecule 6: Non-template strand of the forked DNA oligonucleotide (downstream copy)

Chain t:  29% 100%



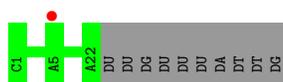
- Molecule 7: Template strand of the forked DNA oligonucleotide (downstream copy) containing the P077 AR9 promoter motif

Chain T:  12% 66% 34%



- Molecule 7: Template strand of the forked DNA oligonucleotide (downstream copy) containing the P077 AR9 promoter motif

Chain n:  3% 69% 31%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	176.93Å 110.45Å 222.38Å 90.00° 98.52° 90.00°	Depositor
Resolution (Å)	49.35 – 3.40 49.35 – 3.40	Depositor EDS
% Data completeness (in resolution range)	96.9 (49.35-3.40) 97.1 (49.35-3.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.13 (at 3.40Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.239 , 0.293 0.240 , 0.293	Depositor DCC
R_{free} test set	2837 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	109.7	Xtrriage
Anisotropy	0.316	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 91.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	23345	wwPDB-VP
Average B, all atoms (Å ²)	133.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.32% 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: UMP, ZN

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	0.24	0/3909	0.40	0/5259
2	d	0.24	0/3558	0.41	0/4804
3	c	0.24	0/4082	0.44	0/5502
4	D	0.25	0/5222	0.44	0/7038
5	C	0.25	0/5370	0.44	0/7241
6	N	0.53	0/299	0.93	0/459
6	t	0.57	0/324	0.91	0/498
7	T	0.54	0/483	0.96	0/743
7	n	0.54	0/504	0.95	0/775
All	All	0.28	0/23751	0.49	0/32319

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

Due to software issues we are unable to calculate clashes - this section is therefore empty.

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	461/464 (99%)	452 (98%)	9 (2%)	0	100	100
2	d	424/426 (100%)	418 (99%)	6 (1%)	0	100	100
3	c	482/496 (97%)	473 (98%)	9 (2%)	0	100	100
4	D	629/631 (100%)	623 (99%)	6 (1%)	0	100	100
5	C	647/665 (97%)	625 (97%)	22 (3%)	0	100	100
All	All	2643/2682 (98%)	2591 (98%)	52 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	444/445 (100%)	440 (99%)	4 (1%)	78	90
2	d	390/390 (100%)	386 (99%)	4 (1%)	76	88
3	c	457/469 (97%)	456 (100%)	1 (0%)	93	98
4	D	590/590 (100%)	585 (99%)	5 (1%)	81	91
5	C	593/608 (98%)	592 (100%)	1 (0%)	93	98
All	All	2474/2502 (99%)	2459 (99%)	15 (1%)	86	94

5 of 15 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	d	252	ASN
4	D	541	PHE
3	c	154	ASP
5	C	110	ARG
4	D	313	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
2	d	95	GLN
2	d	252	ASN
3	c	165	ASN
4	D	494	ASN

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 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 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
9	DG	A	502	8	18,24,25	1.09	2 (11%)	19,35,38	0.63	0
12	DA	n	101	8	18,23,24	0.66	0	17,33,36	0.69	1 (5%)
8	UMP	T	101	7,8	17,20,21	0.18	0	24,28,31	0.47	0
8	UMP	A	503	8,12	17,20,21	0.17	0	24,28,31	0.41	0
8	UMP	A	501	8,9	17,20,21	0.17	0	24,28,31	0.39	0
8	UMP	c	502	8	17,20,21	0.19	0	24,28,31	0.48	0
11	DT	t	101	-	18,21,22	0.33	0	26,30,33	0.65	1 (3%)
8	UMP	c	501	8,9	17,20,21	0.20	0	24,28,31	0.39	0

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
9	DG	A	502	8	-	0/3/21/22	0/3/3/3
12	DA	n	101	8	-	1/3/21/22	0/3/3/3
8	UMP	T	101	7,8	-	4/7/21/22	0/2/2/2
8	UMP	A	503	8,12	-	6/7/21/22	0/2/2/2
8	UMP	A	501	8,9	-	1/7/21/22	0/2/2/2
8	UMP	c	502	8	-	0/7/21/22	0/2/2/2
11	DT	t	101	-	-	1/7/21/22	0/2/2/2
8	UMP	c	501	8,9	-	6/7/21/22	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	502	DG	C5-C6	-2.72	1.41	1.47
9	A	502	DG	C8-N7	-2.45	1.30	1.35

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	t	101	DT	O4-C4-C5	-2.19	122.36	124.90
12	n	101	DA	C5-C6-N6	2.02	123.42	120.35

There are no chirality outliers.

5 of 19 torsion outliers are listed below:

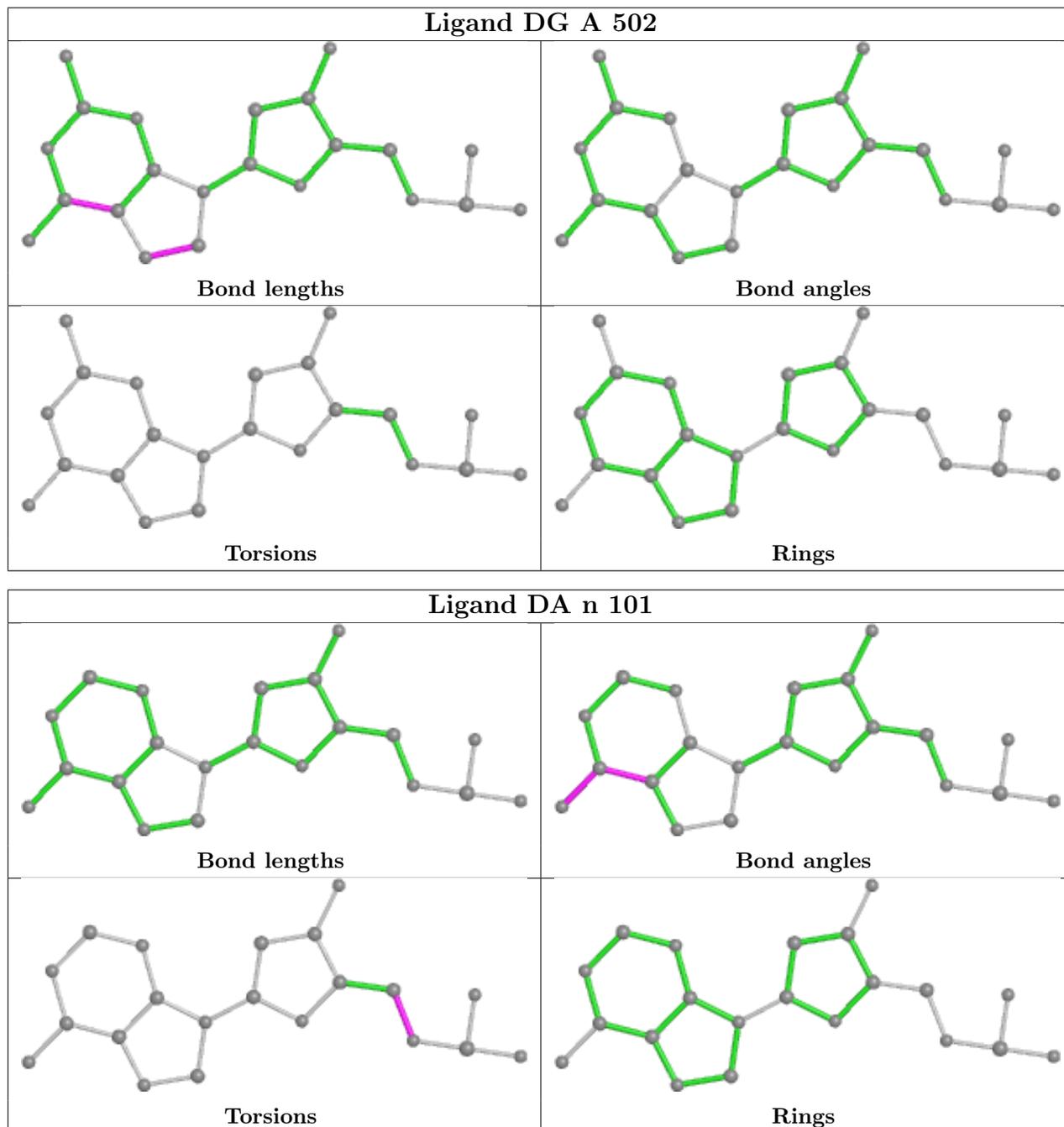
Mol	Chain	Res	Type	Atoms
8	c	501	UMP	O4'-C4'-C5'-O5'
8	A	503	UMP	O4'-C4'-C5'-O5'
8	c	501	UMP	C3'-C4'-C5'-O5'
11	t	101	DT	C3'-C4'-C5'-O5'
8	A	503	UMP	C2'-C1'-N1-C6

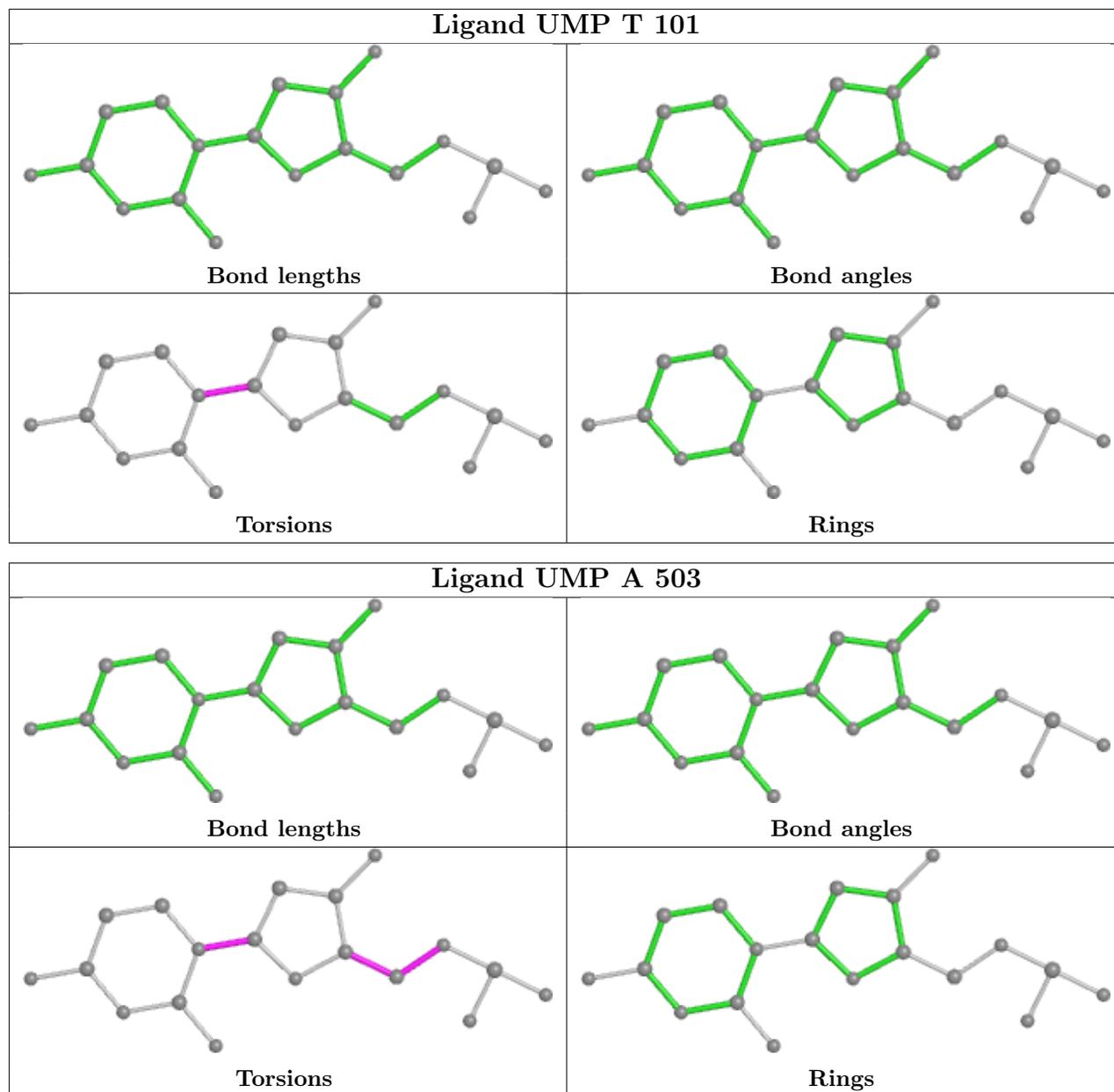
There are no ring outliers.

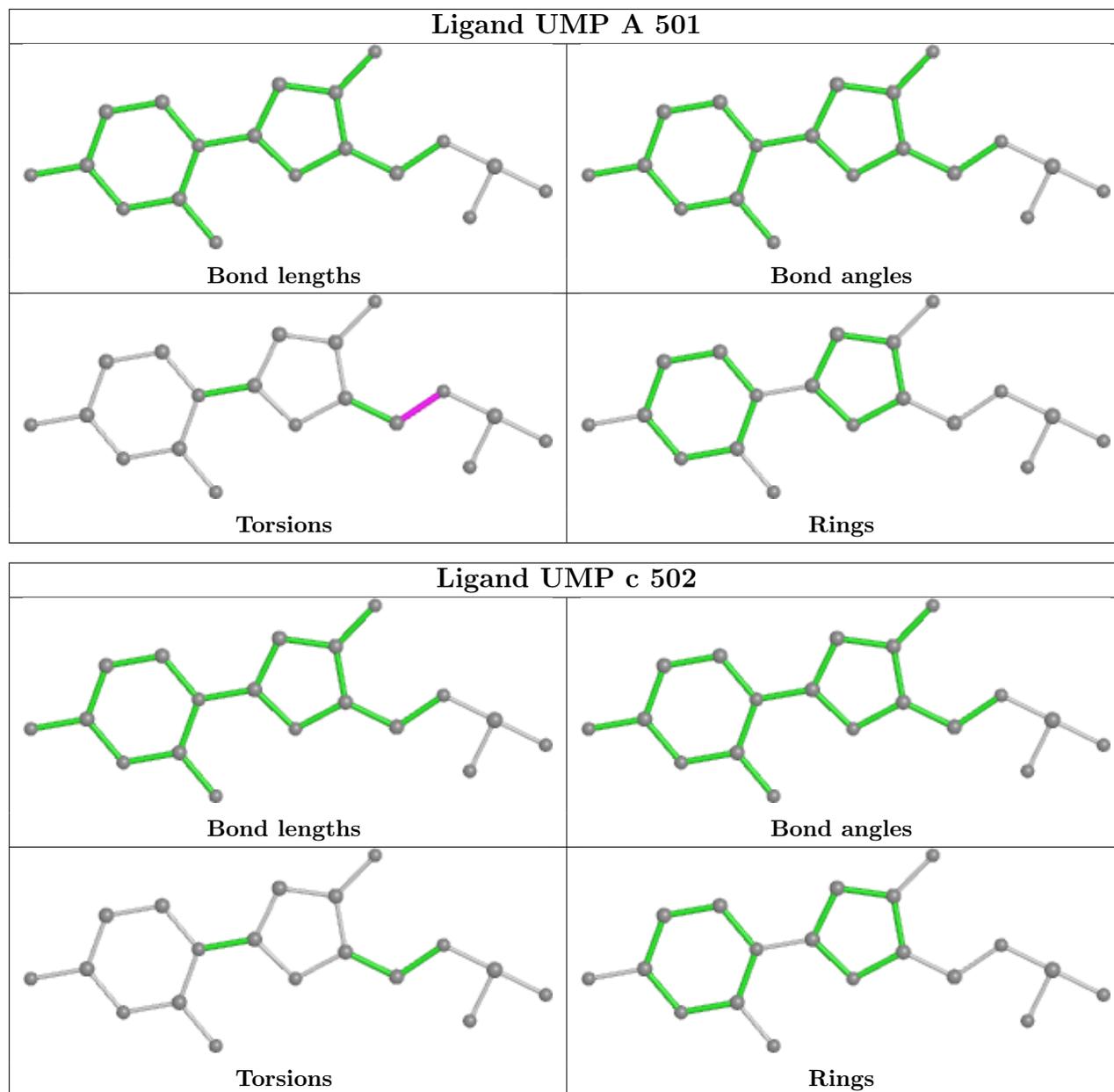
No monomer is involved in short contacts.

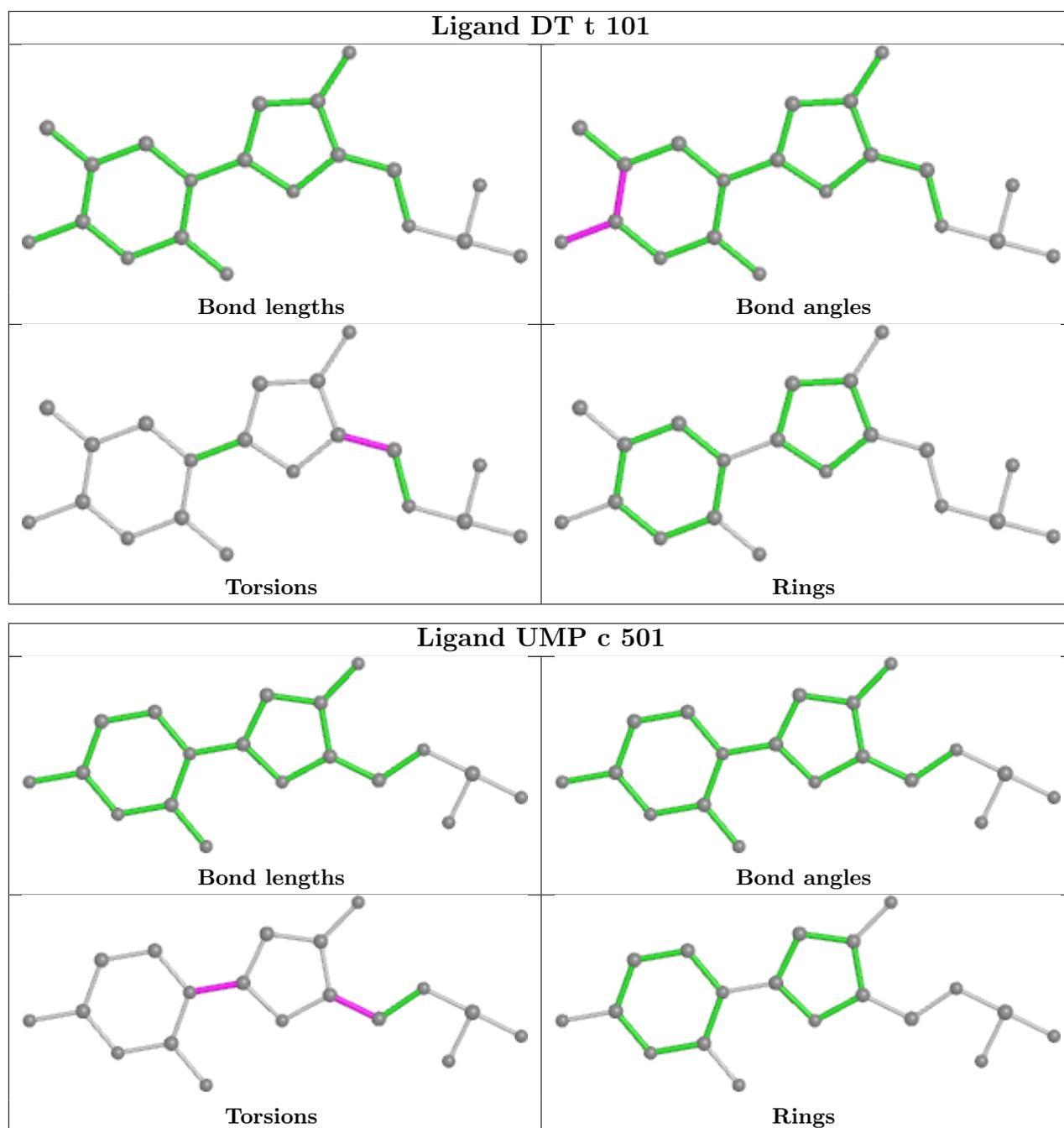
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is

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









5.7 Other polymers [\(i\)](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [\(i\)](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	463/464 (99%)	0.38	24 (5%) 27 27	92, 151, 211, 270	0
2	d	426/426 (100%)	0.15	5 (1%) 79 77	77, 116, 183, 248	0
3	c	484/496 (97%)	0.35	22 (4%) 33 33	77, 125, 193, 259	0
4	D	631/631 (100%)	0.19	7 (1%) 80 79	78, 122, 170, 208	0
5	C	649/665 (97%)	0.16	1 (0%) 95 95	66, 101, 144, 190	0
6	N	13/14 (92%)	0.52	0 100 100	210, 230, 349, 352	0
6	t	14/14 (100%)	1.50	4 (28%) 0 0	185, 246, 330, 351	0
7	T	21/32 (65%)	1.25	4 (19%) 1 1	200, 233, 285, 363	0
7	n	22/32 (68%)	0.57	1 (4%) 33 33	142, 205, 261, 275	0
All	All	2723/2774 (98%)	0.26	68 (2%) 57 55	66, 121, 197, 363	0

The worst 5 of 68 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	392	ASN	5.8
7	T	18	DA	5.6
3	c	1	MET	4.5
3	c	93	GLU	4.1
1	A	391	MET	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

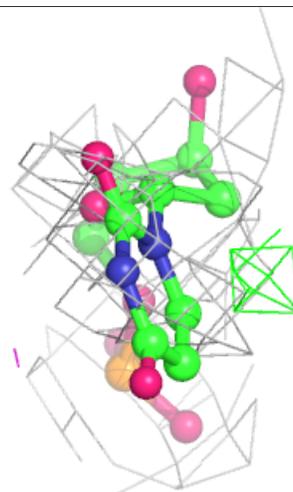
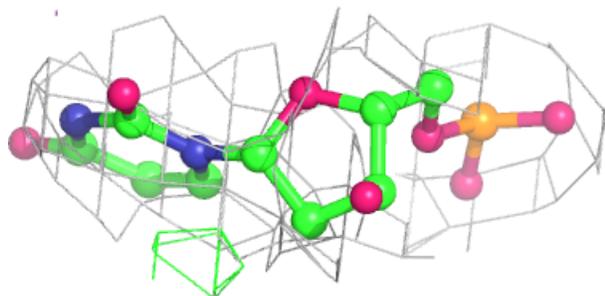
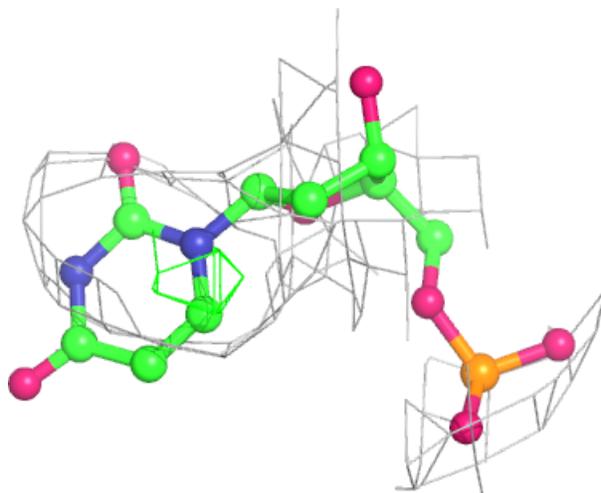
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

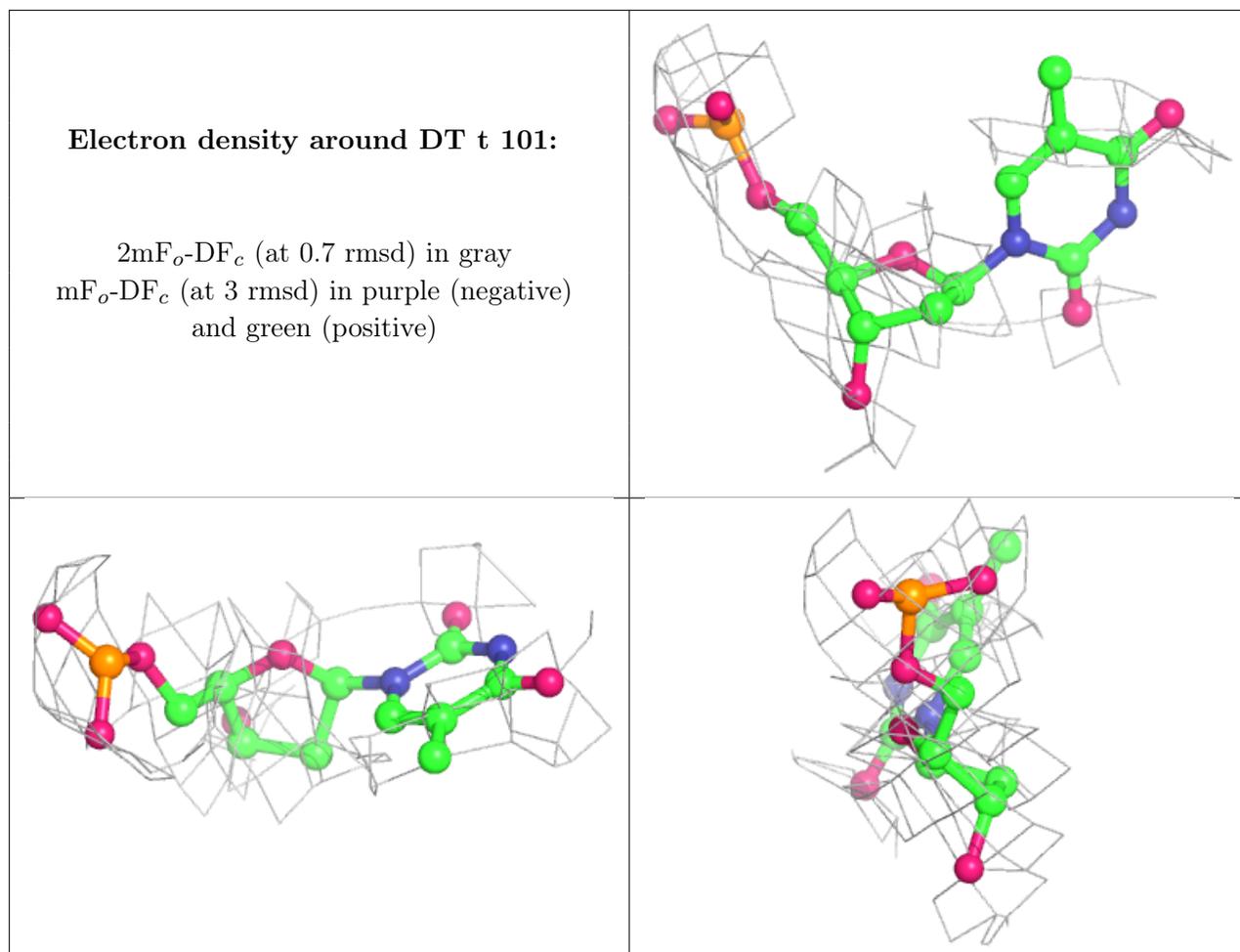
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
8	UMP	T	101	19/20	0.68	0.26	182,201,215,216	0
11	DT	t	101	20/21	0.73	0.22	240,244,248,250	0
12	DA	n	101	21/22	0.77	0.24	191,205,225,232	0
8	UMP	A	503	19/20	0.85	0.38	135,199,221,221	0
8	UMP	A	501	19/20	0.89	0.22	125,156,177,190	0
8	UMP	c	502	19/20	0.91	0.16	118,124,136,140	0
9	DG	A	502	22/23	0.92	0.20	115,124,136,141	0
8	UMP	c	501	19/20	0.94	0.22	117,123,126,128	0
10	ZN	D	801	1/1	0.96	0.08	116,116,116,116	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 UMP T 101:

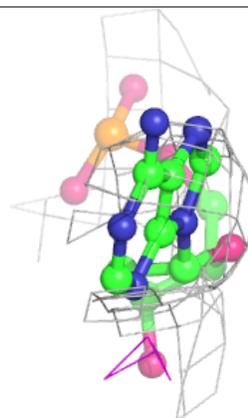
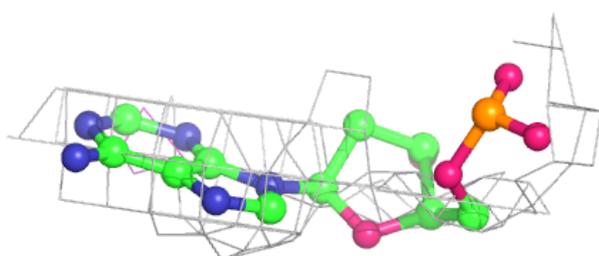
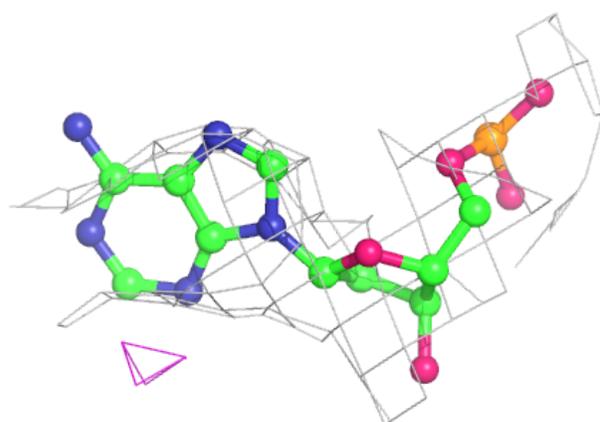
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



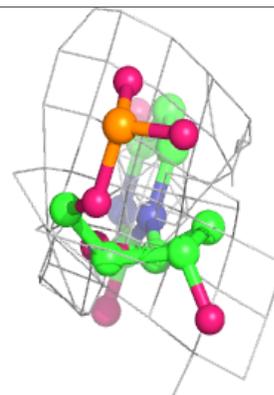
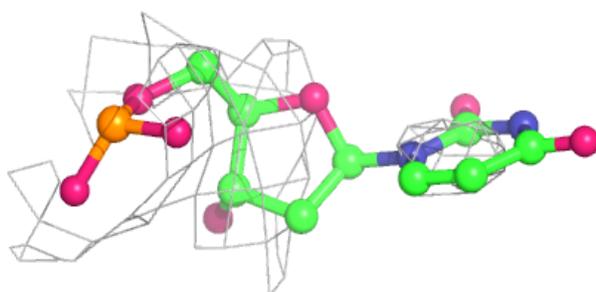
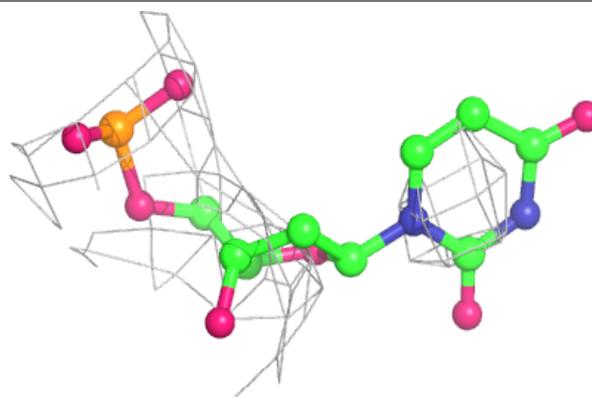


Electron density around DA n 101:

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

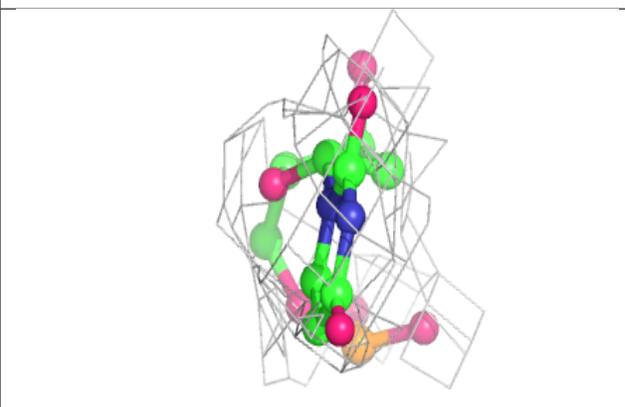
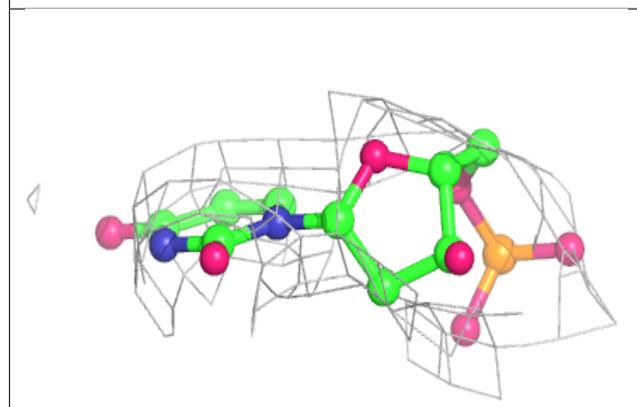
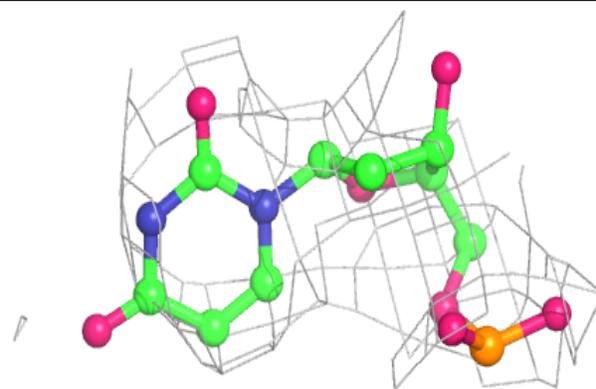
**Electron density around UMP A 503:**

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

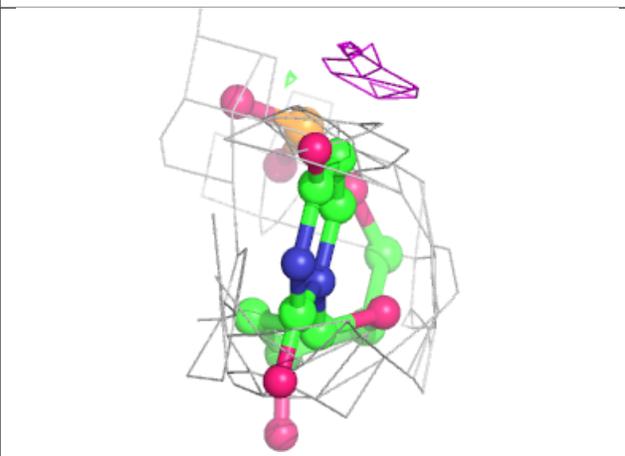
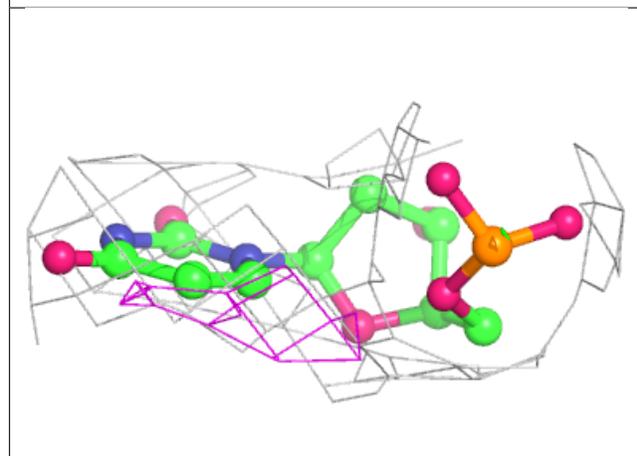
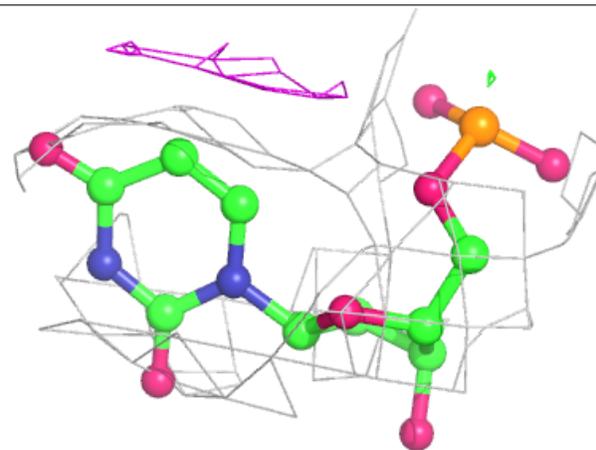


Electron density around UMP A 501:

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

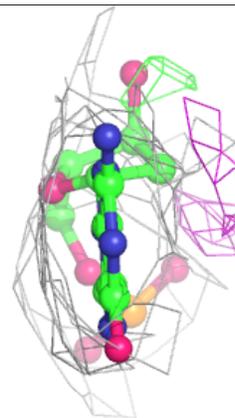
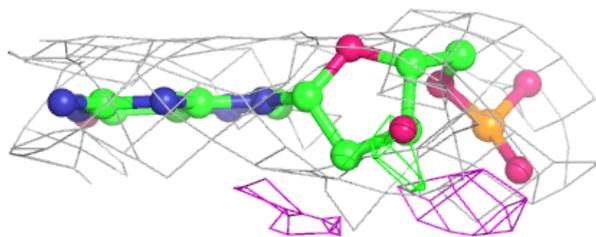
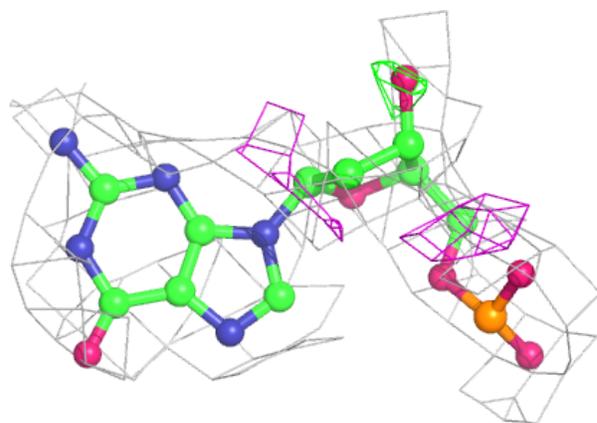
**Electron density around UMP c 502:**

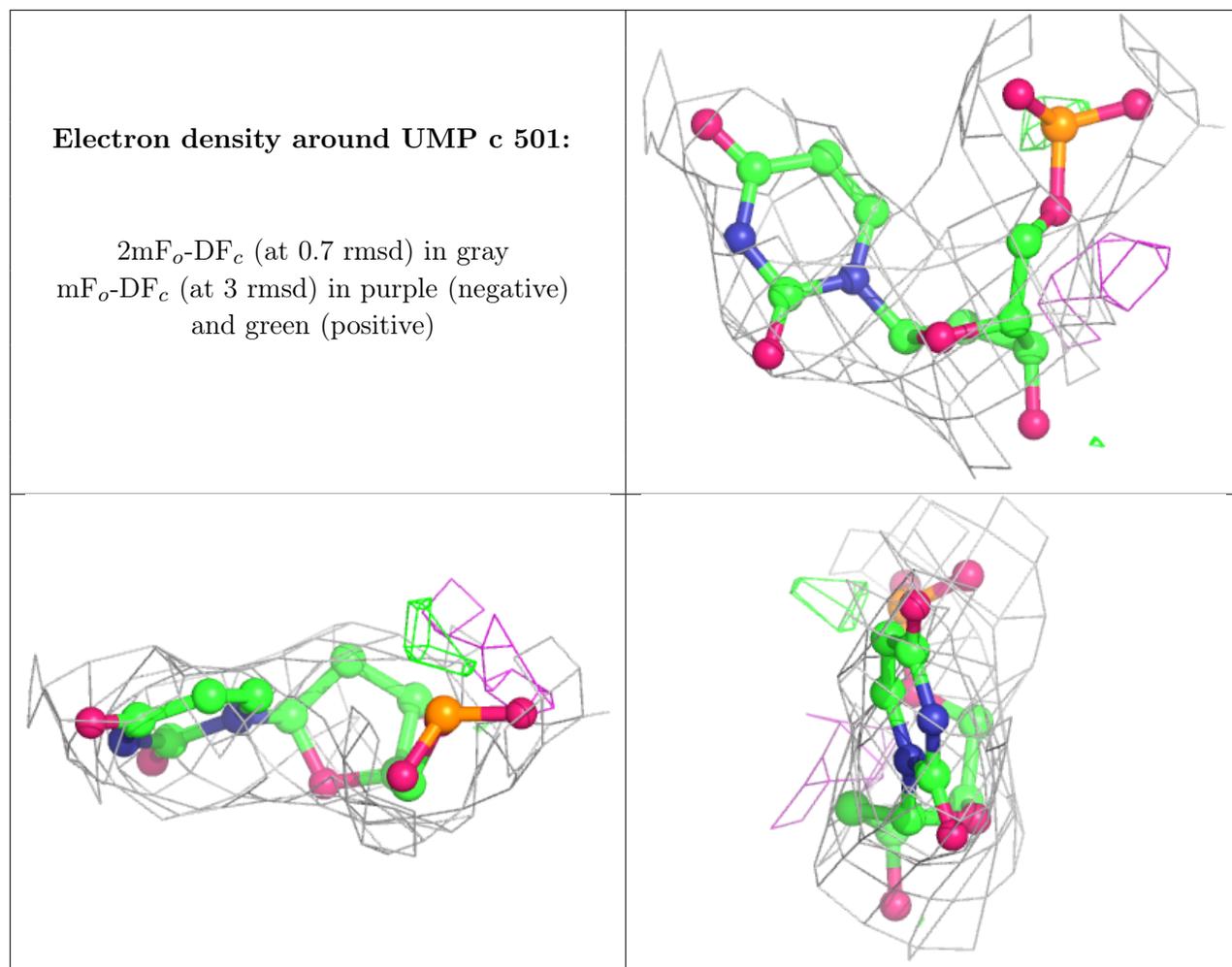
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

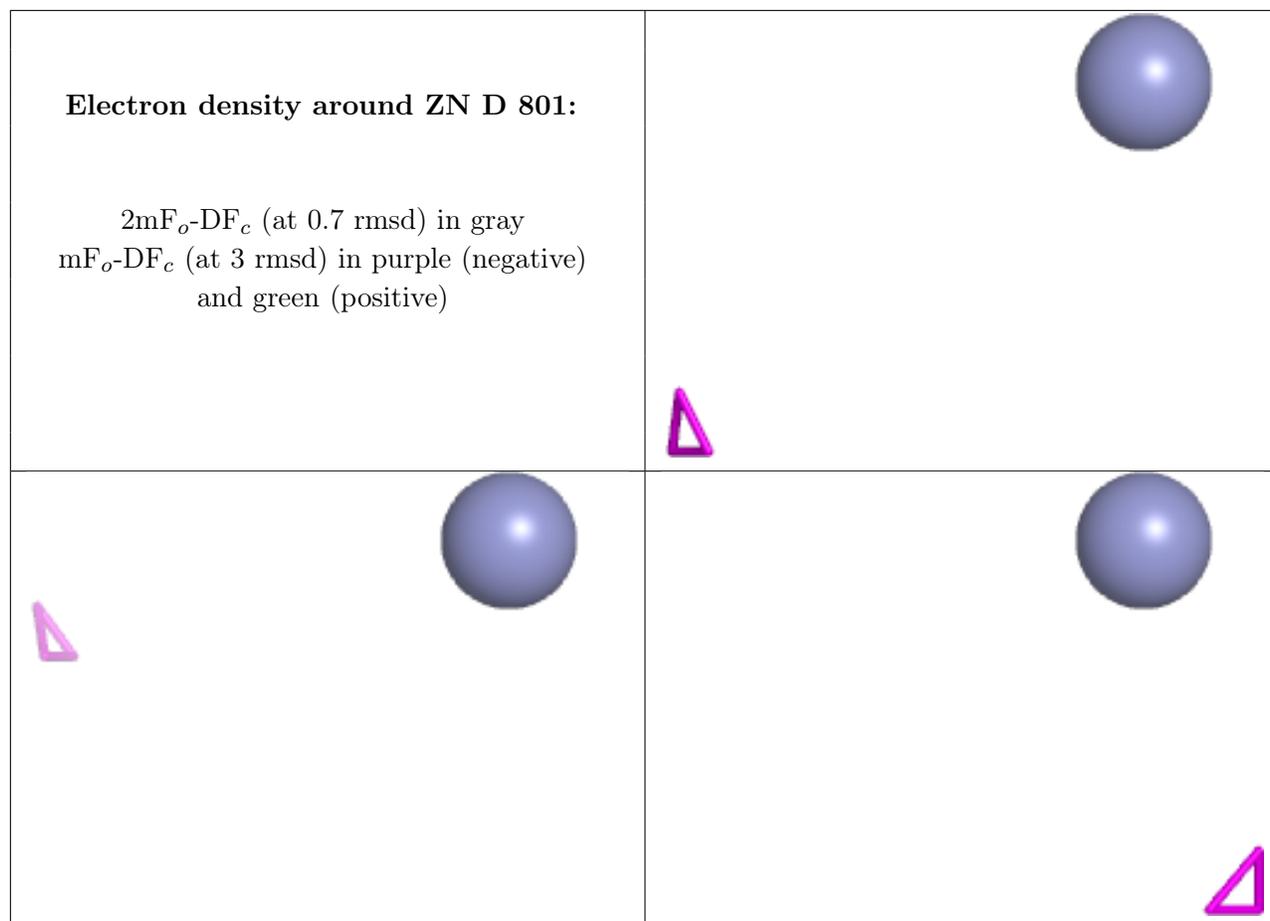


Electron density around DG A 502:

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







6.5 Other polymers [i](#)

There are no such residues in this entry.