



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2023 – 12:48 AM EDT

PDB ID : 7M4K  
Title : DNA Polymerase Lambda, TTPaS:At Ca<sup>2+</sup> Ground State Ternary Complex  
Authors : Jamsen, J.A.; Wilson, S.H.  
Deposited on : 2021-03-21  
Resolution : 1.72 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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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 : 2.35.1  
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) : 2.35.1

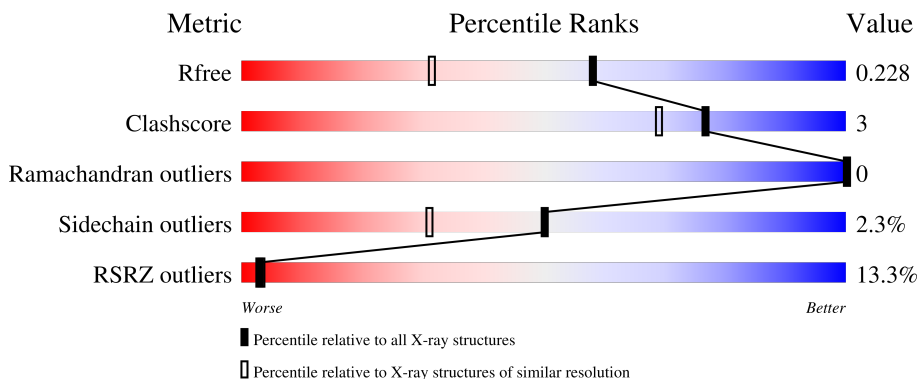
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.72 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	5722 (1.74-1.70)
Clashscore	141614	6152 (1.74-1.70)
Ramachandran outliers	138981	6051 (1.74-1.70)
Sidechain outliers	138945	6051 (1.74-1.70)
RSRZ outliers	127900	5629 (1.74-1.70)

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  $\geq 3$ , 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	329	 13% 85% 9% 6%
2	T	11	 55% 45%
3	P	6	 100%
4	D	4	 75% 25%

## 2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 3044 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 polymerase lambda.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	310	2361	1490	421	436	14	0	10	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	463	LYS	SER	conflict	UNP Q9UGP5
A	464	GLY	GLN	conflict	UNP Q9UGP5
A	?	-	GLU	deletion	UNP Q9UGP5
A	?	-	ASN	deletion	UNP Q9UGP5
A	?	-	GLY	deletion	UNP Q9UGP5
A	?	-	GLN	deletion	UNP Q9UGP5
A	?	-	GLN	deletion	UNP Q9UGP5
A	471	THR	GLN	conflict	UNP Q9UGP5
A	543	ALA	CYS	engineered mutation	UNP Q9UGP5

- Molecule 2 is a DNA chain called DNA (5'-D(\*CP\*GP\*GP\*CP\*AP\*GP\*TP\*AP\*CP\*TP\*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	T	11	224	107	43	64	10	0	0	0

- Molecule 3 is a DNA chain called DNA (5'-D(\*CP\*AP\*GP\*TP\*AP\*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	P	6	119	58	23	33	5	0	0	0

- Molecule 4 is a DNA chain called DNA (5'-D(P\*GP\*CP\*CP\*G)-3').

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
			Total	C	N	O				P
4	D	4	83	38	16	25	4	0	0	0

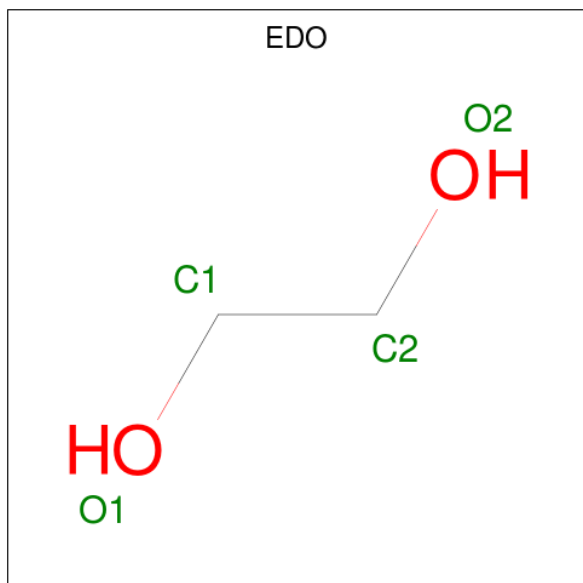
- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Ca	0	0
			1	1		

- Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	4	Total	Na	0	0
			4	4		

- Molecule 7 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



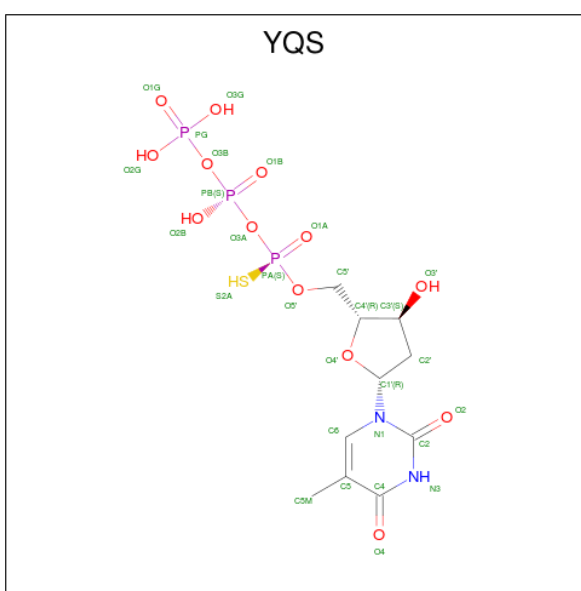
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			4	2	2		
7	A	1	Total	C	O	0	0
			4	2	2		
7	A	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			4	2	2		
7	A	1	Total	C	O	0	0
			3	2	1		
7	A	1	Total	C	O	0	0
			3	2	1		

- Molecule 8 is  $[[[2 \{R\}, 3 \{S\}, 5 \{R\}]-5-[5\text{-methyl-2,4-bis(oxidanylidene)pyrimidin-1-yl]-3\text{-oxidanyl-oxolan-2-yl}]methoxy-sulfanyl-phosphoryl] phosphono hydrogen phosphate$  (three-letter code: YQS) (formula:  $C_{10}H_{17}N_2O_{13}P_3S$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
8	A	1	Total	C	N	O	P	S	0	0
			29	10	2	13	3	1		

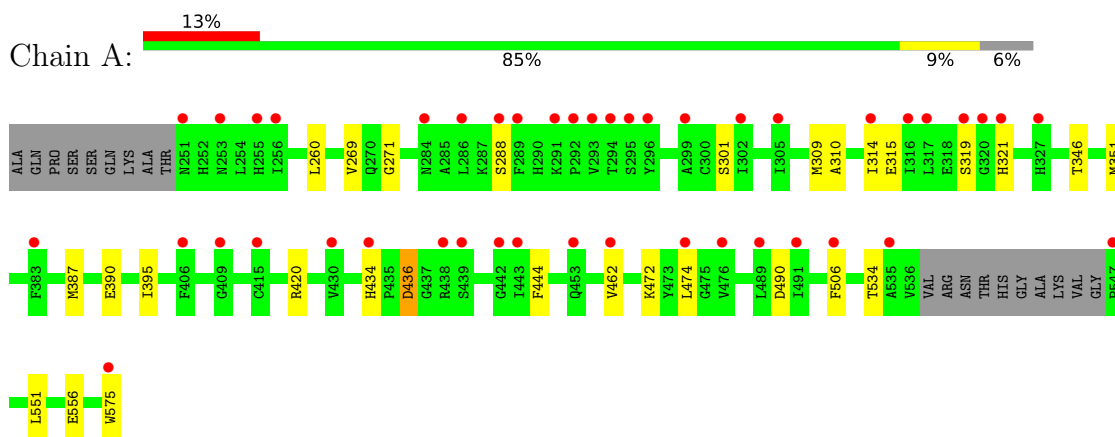
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	133	Total	O	0	5
			135	135		
9	T	36	Total	O	0	0
			36	36		
9	P	28	Total	O	0	0
			28	28		
9	D	2	Total	O	0	0
			2	2		

### 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 polymerase lambda



- Molecule 2: DNA (5'-D(\*CP\*GP\*GP\*CP\*AP\*GP\*TP\*AP\*CP\*TP\*G)-3')

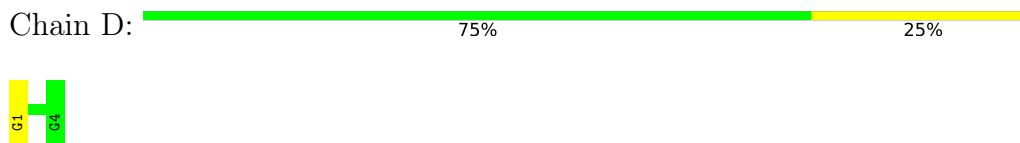


- Molecule 3: DNA (5'-D(\*CP\*AP\*GP\*TP\*AP\*C)-3')



There are no outlier residues recorded for this chain.

- Molecule 4: DNA (5'-D(P\*GP\*CP\*CP\*G)-3')



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	56.00Å 62.51Å 140.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.12 – 1.72 31.12 – 1.72	Depositor EDS
% Data completeness (in resolution range)	98.6 (31.12-1.72) 98.6 (31.12-1.72)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.59 (at 1.72Å)	Xtrriage
Refinement program	PHENIX v1.15.2-3472	Depositor
R, $R_{free}$	0.196 , 0.227 0.199 , 0.228	Depositor DCC
$R_{free}$ test set	2569 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	32.6	Xtrriage
Anisotropy	0.640	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 58.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	3044	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.84% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

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

## 5 Model quality i

### 5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: CA, YQS, NA, EDO

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	0.49	0/2414	0.65	2/3268 (0.1%)
2	T	1.56	0/251	1.36	5/386 (1.3%)
3	P	1.52	0/133	1.18	0/203
4	D	1.39	1/92 (1.1%)	0.83	0/138
All	All	0.76	1/2890 (0.0%)	0.79	7/3995 (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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1	DG	OP3-P	-9.94	1.49	1.61

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	2	DG	O4'-C1'-N9	6.37	112.46	108.00
2	T	7	DT	N3-C4-O4	5.80	123.38	119.90
1	A	351	MET	CG-SD-CE	-5.75	91.00	100.20
2	T	7	DT	C5-C4-O4	-5.30	121.19	124.90
2	T	8	DA	OP1-P-OP2	5.21	127.42	119.60
2	T	5	DA	OP1-P-OP2	5.16	127.34	119.60
1	A	420	ARG	NE-CZ-NH1	5.06	122.83	120.30



There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	288	SER	Mainchain

## 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	2361	0	2211	16	0
2	T	224	0	125	2	0
3	P	119	0	69	0	0
4	D	83	0	45	0	0
5	A	1	0	0	0	0
6	A	4	0	0	0	0
7	A	22	0	30	1	0
8	A	29	0	0	0	0
9	A	135	0	0	0	0
9	D	2	0	0	0	0
9	P	28	0	0	0	0
9	T	36	0	0	0	0
All	All	3044	0	2480	16	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 3.

All (16) 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:A:462:VAL:HG13	2:T:9:DC:H4'	1.81	0.63
1:A:269:VAL:O	1:A:346[B]:THR:HG23	2.05	0.56
1:A:462:VAL:CG1	2:T:9:DC:H4'	2.36	0.56
1:A:271:GLY:CA	1:A:346[B]:THR:HG21	2.39	0.53
1:A:319:SER:OG	1:A:321:HIS:O	2.28	0.51
1:A:310:ALA:O	1:A:314:ILE:HG23	2.11	0.50
1:A:434:HIS:CG	1:A:436:ASP:OD1	2.65	0.49
1:A:534[B]:THR:HB	1:A:551:LEU:HD11	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:575:TRP:CD1	1:A:575:TRP:N	2.82	0.46
1:A:556:GLU:H	7:A:608:EDO:H12	1.82	0.44
1:A:474:LEU:HD22	1:A:490:ASP:CG	2.39	0.43
1:A:387[B]:MET:CE	1:A:395:ILE:HD12	2.49	0.42
1:A:387[B]:MET:HE1	1:A:395:ILE:HD12	2.01	0.42
1:A:390:GLU:H	1:A:390:GLU:CD	2.23	0.41
1:A:260:LEU:HD13	1:A:309[B]:MET:HE3	2.03	0.41
1:A:315:GLU:O	1:A:319:SER:OG	2.38	0.41

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	315/329 (96%)	310 (98%)	5 (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	223/274 (81%)	218 (98%)	5 (2%)	52 33

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	301	SER
1	A	436	ASP
1	A	444	PHE
1	A	472	LYS
1	A	506	PHE

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

### 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 12 ligands modelled in this entry, 5 are monoatomic - leaving 7 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
7	EDO	A	610	-	2,2,3	0.47	0	1,1,2	0.34	0
8	YQS	A	612	5,6	24,30,30	1.06	2 (8%)	37,47,47	2.23	7 (18%)
7	EDO	A	606	-	3,3,3	0.51	0	2,2,2	0.15	0
7	EDO	A	609	-	3,3,3	0.45	0	2,2,2	0.50	0
7	EDO	A	608	-	3,3,3	0.47	0	2,2,2	0.14	0
7	EDO	A	611	-	2,2,3	0.51	0	1,1,2	0.31	0
7	EDO	A	607	-	3,3,3	0.50	0	2,2,2	0.35	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
7	EDO	A	606	-	-	0/1/1/1	-
7	EDO	A	609	-	-	1/1/1/1	-
7	EDO	A	608	-	-	0/1/1/1	-
8	YQS	A	612	5,6	-	1/18/34/34	0/2/2/2
7	EDO	A	607	-	-	0/1/1/1	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	612	YQS	C4-N3	3.12	1.44	1.38
8	A	612	YQS	PA-O5'	3.08	1.62	1.57

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	612	YQS	C4-N3-C2	-8.17	116.78	127.35
8	A	612	YQS	C5-C4-N3	6.56	120.91	115.31
8	A	612	YQS	N3-C2-N1	4.54	120.92	114.89
8	A	612	YQS	PB-O3B-PG	-3.12	122.14	132.83
8	A	612	YQS	O4-C4-C5	-3.06	121.36	124.90
8	A	612	YQS	O2-C2-N1	-2.56	119.38	122.79
8	A	612	YQS	C2'-C1'-N1	-2.03	109.09	113.77

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	612	YQS	PB-O3B-PG-O3G
7	A	609	EDO	O1-C1-C2-O2

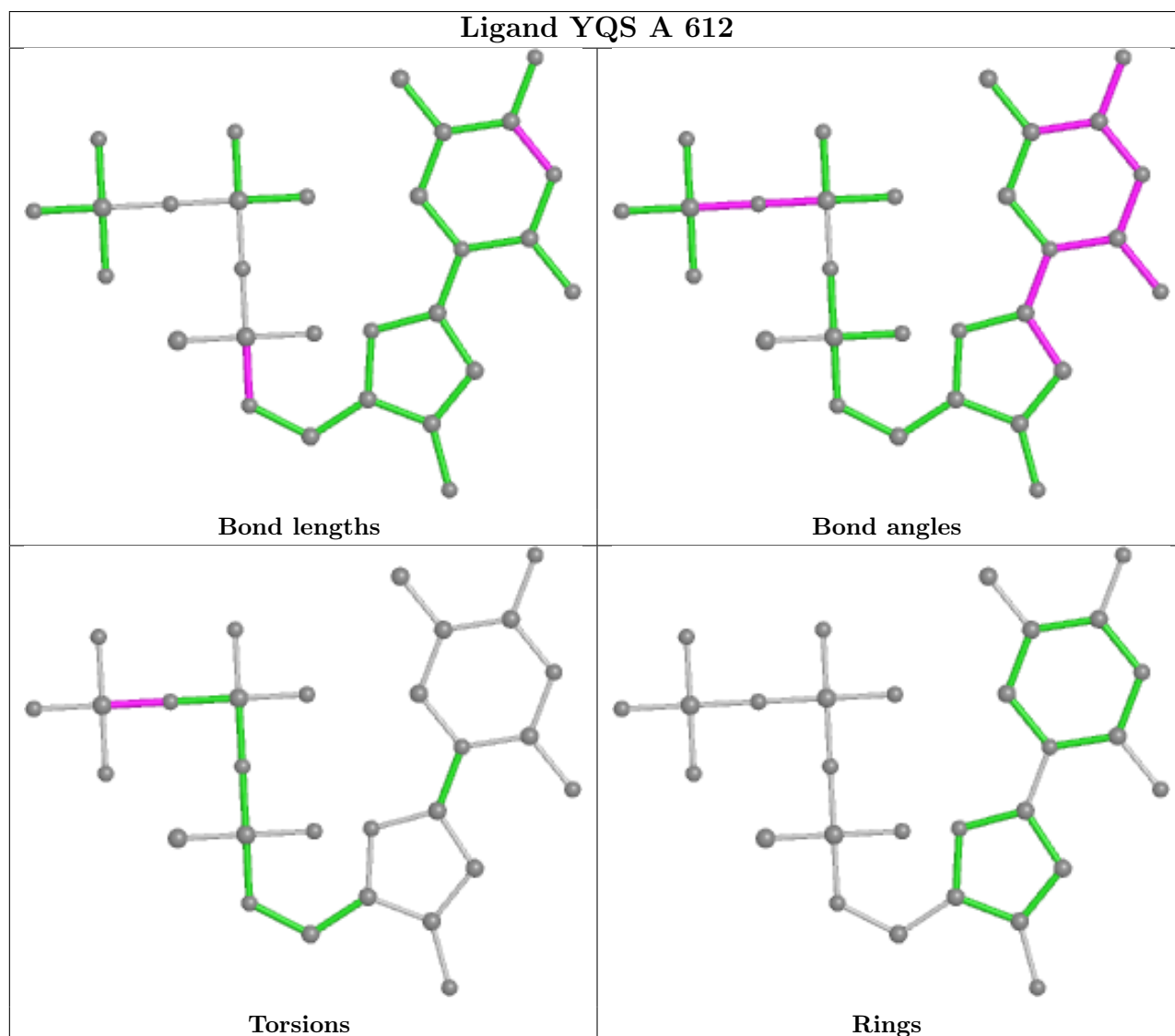
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	608	EDO	1	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

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	310/329 (94%)	0.63	44 (14%) <b>2</b>   <b>2</b>	23, 47, 87, 120	1 (0%)
2	T	11/11 (100%)	-0.45	0 <b>100</b>   <b>100</b>	33, 35, 52, 57	0
3	P	6/6 (100%)	-0.13	0 <b>100</b>   <b>100</b>	25, 27, 30, 40	0
4	D	4/4 (100%)	-0.62	0 <b>100</b>   <b>100</b>	45, 46, 51, 57	0
All	All	331/350 (94%)	0.56	44 (13%) <b>3</b>   <b>3</b>	23, 47, 87, 120	1 (0%)

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	319	SER	6.0
1	A	293	VAL	5.7
1	A	289	PHE	5.5
1	A	320	GLY	4.4
1	A	439	SER	4.1
1	A	292	PRO	3.7
1	A	299	ALA	3.6
1	A	575	TRP	3.5
1	A	434	HIS	3.5
1	A	286	LEU	3.4
1	A	295	SER	3.4
1	A	321	HIS	3.4
1	A	317	LEU	3.3
1	A	474	LEU	3.2
1	A	255	HIS	3.1
1	A	438	ARG	3.1
1	A	535[A]	ALA	3.1
1	A	489	LEU	3.1
1	A	409	GLY	3.0
1	A	296	TYR	3.0
1	A	506	PHE	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	305	ILE	2.8
1	A	491	ILE	2.8
1	A	442	GLY	2.8
1	A	314	ILE	2.7
1	A	294	THR	2.6
1	A	316	ILE	2.6
1	A	443	ILE	2.6
1	A	291	LYS	2.5
1	A	476	VAL	2.5
1	A	256	ILE	2.4
1	A	288	SER	2.4
1	A	462	VAL	2.4
1	A	430	VAL	2.3
1	A	415	CYS	2.3
1	A	453	GLN	2.3
1	A	547	PRO	2.3
1	A	406	PHE	2.3
1	A	327	HIS	2.2
1	A	251	ASN	2.2
1	A	302	ILE	2.2
1	A	383	PHE	2.1
1	A	284	ASN	2.1
1	A	253	ASN	2.0

## 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, 95<sup>th</sup> 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.

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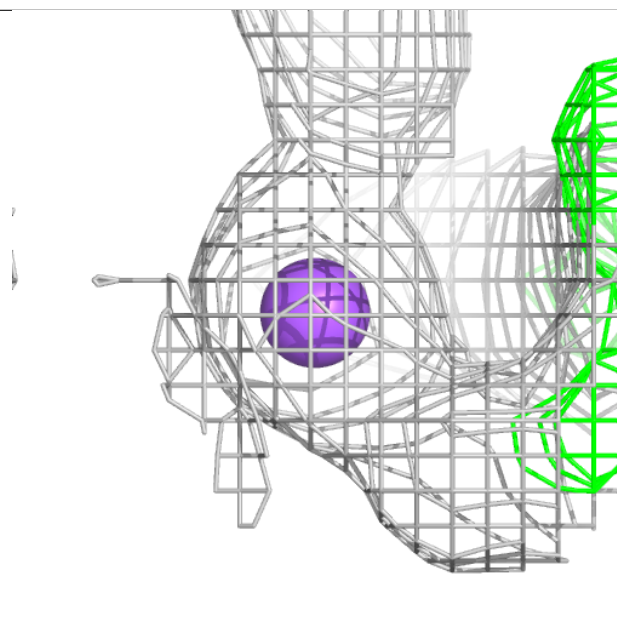
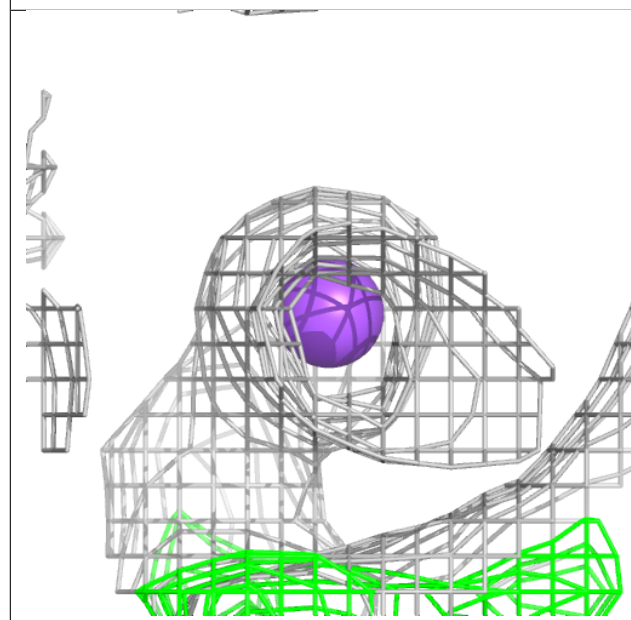
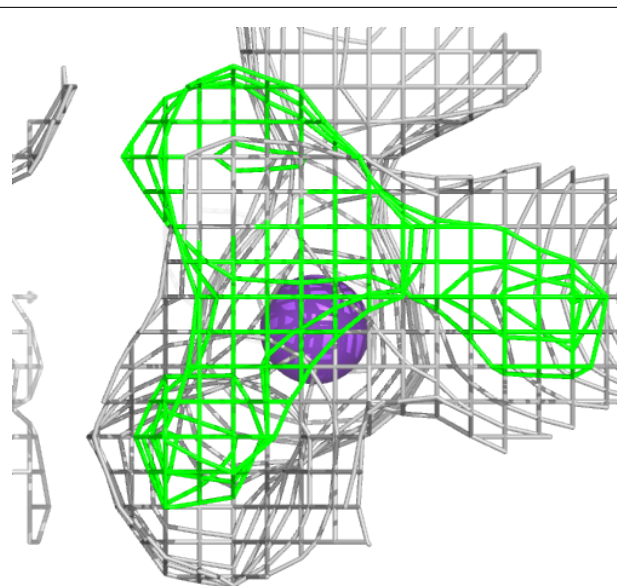
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
7	EDO	A	606	4/4	0.59	0.28	50,52,60,63	0
7	EDO	A	608	4/4	0.71	0.17	67,71,72,78	0
6	NA	A	605	1/1	0.81	0.36	59,59,59,59	0
6	NA	A	604	1/1	0.82	0.07	58,58,58,58	0
7	EDO	A	609	4/4	0.82	0.26	40,44,45,46	4
7	EDO	A	611	3/4	0.83	0.13	49,49,55,67	0
7	EDO	A	610	3/4	0.89	0.14	46,46,54,56	0
8	YQS	A	612	29/29	0.94	0.10	30,51,88,106	0
7	EDO	A	607	4/4	0.96	0.14	37,41,41,49	4
5	CA	A	601	1/1	0.97	0.03	45,45,45,45	0
6	NA	A	602	1/1	0.98	0.07	48,48,48,48	0
6	NA	A	603	1/1	0.99	0.09	24,24,24,24	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.

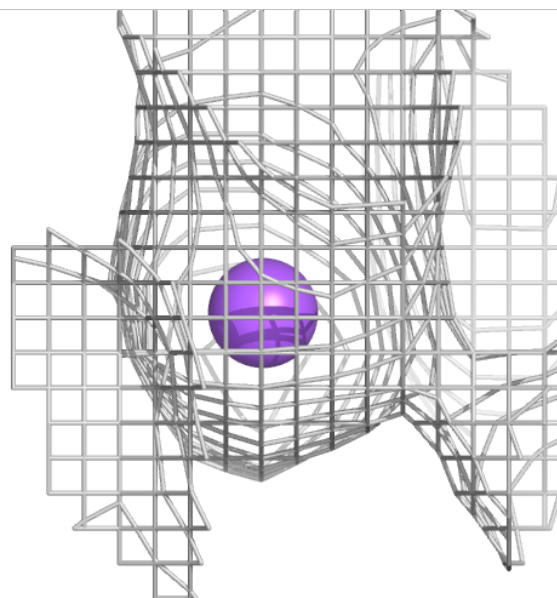
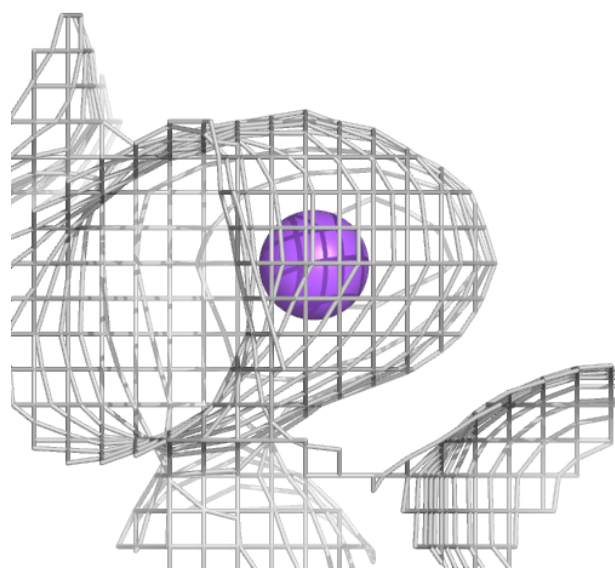
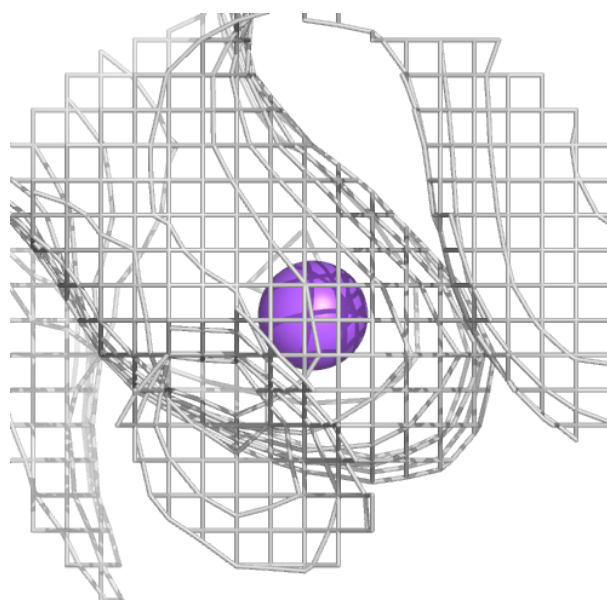
**Electron density around NA A 605:**

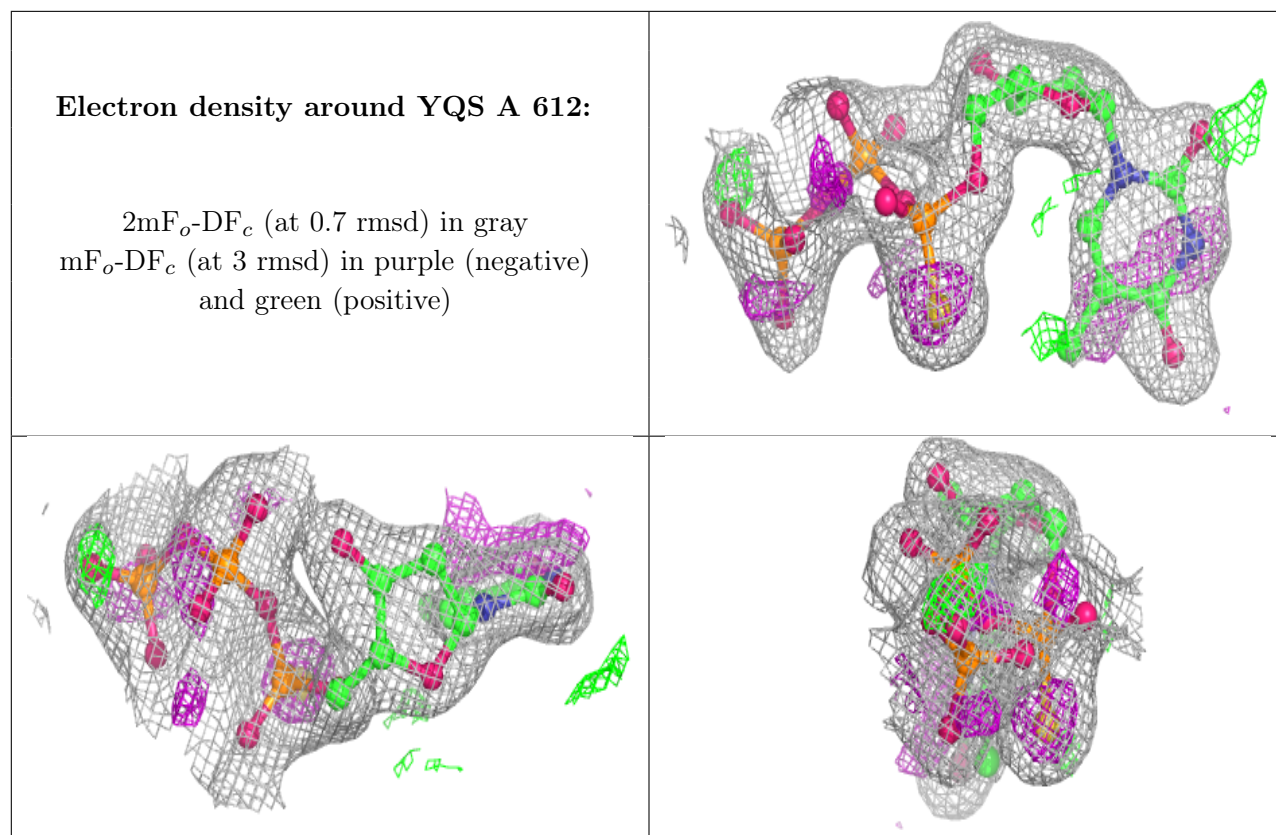
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around NA A 604:**

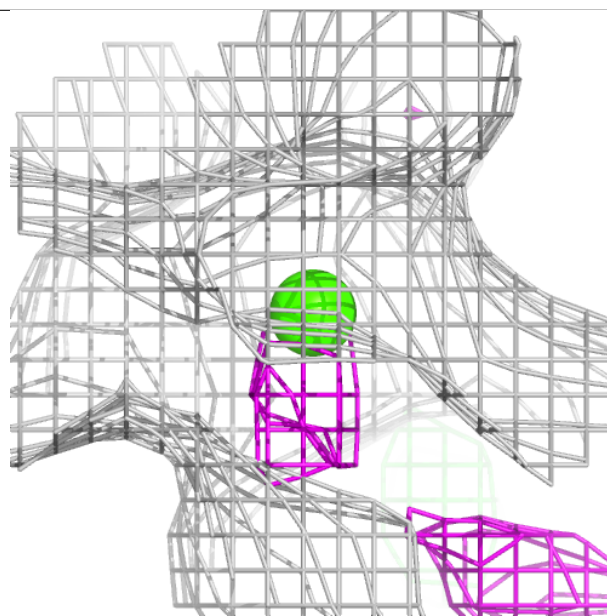
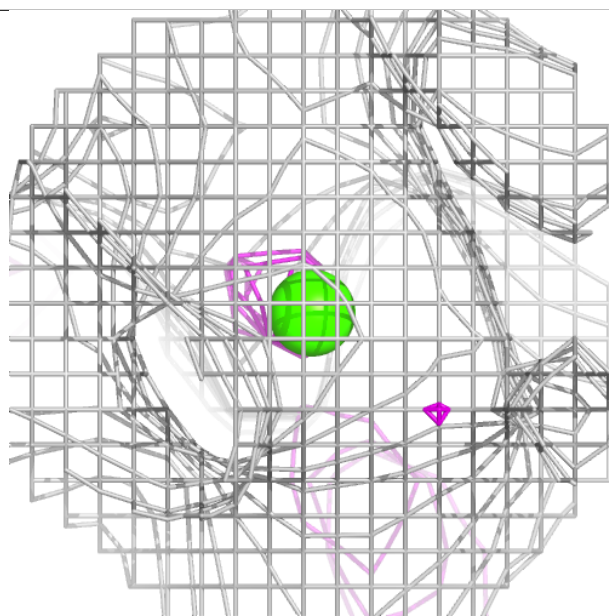
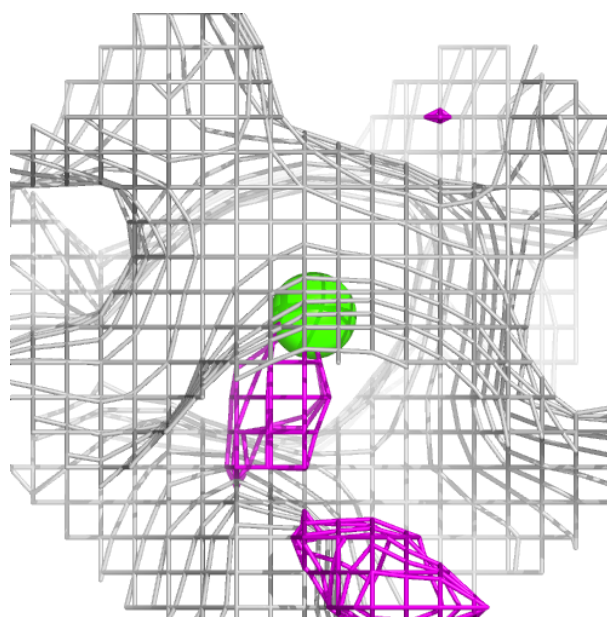
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





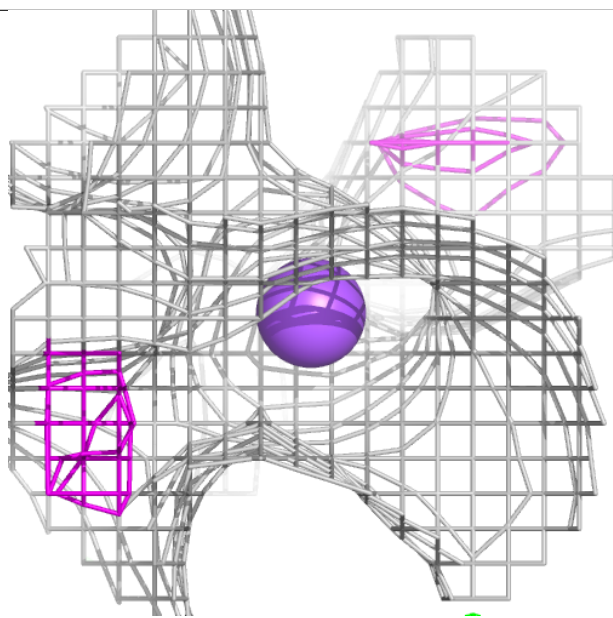
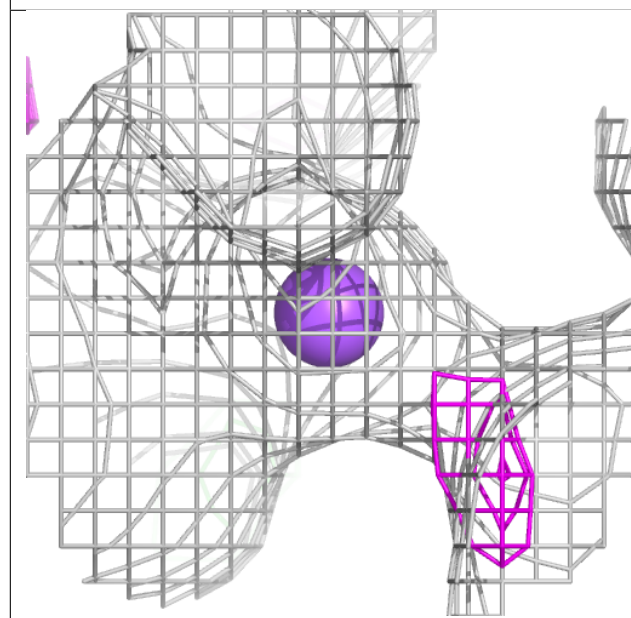
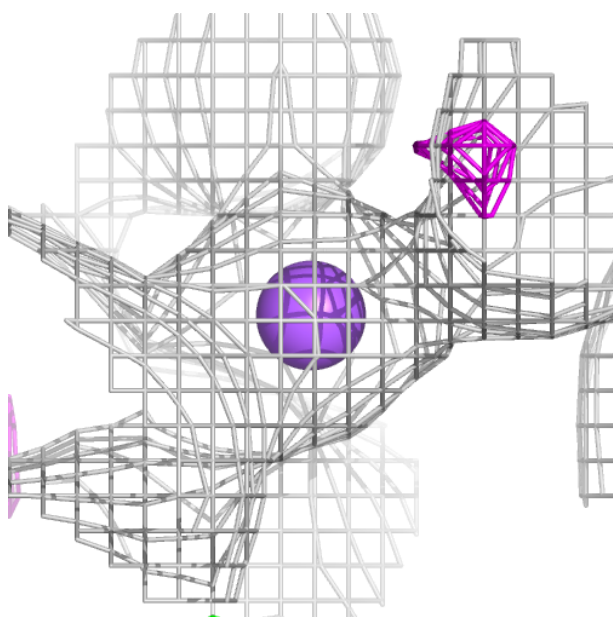
**Electron density around CA A 601:**

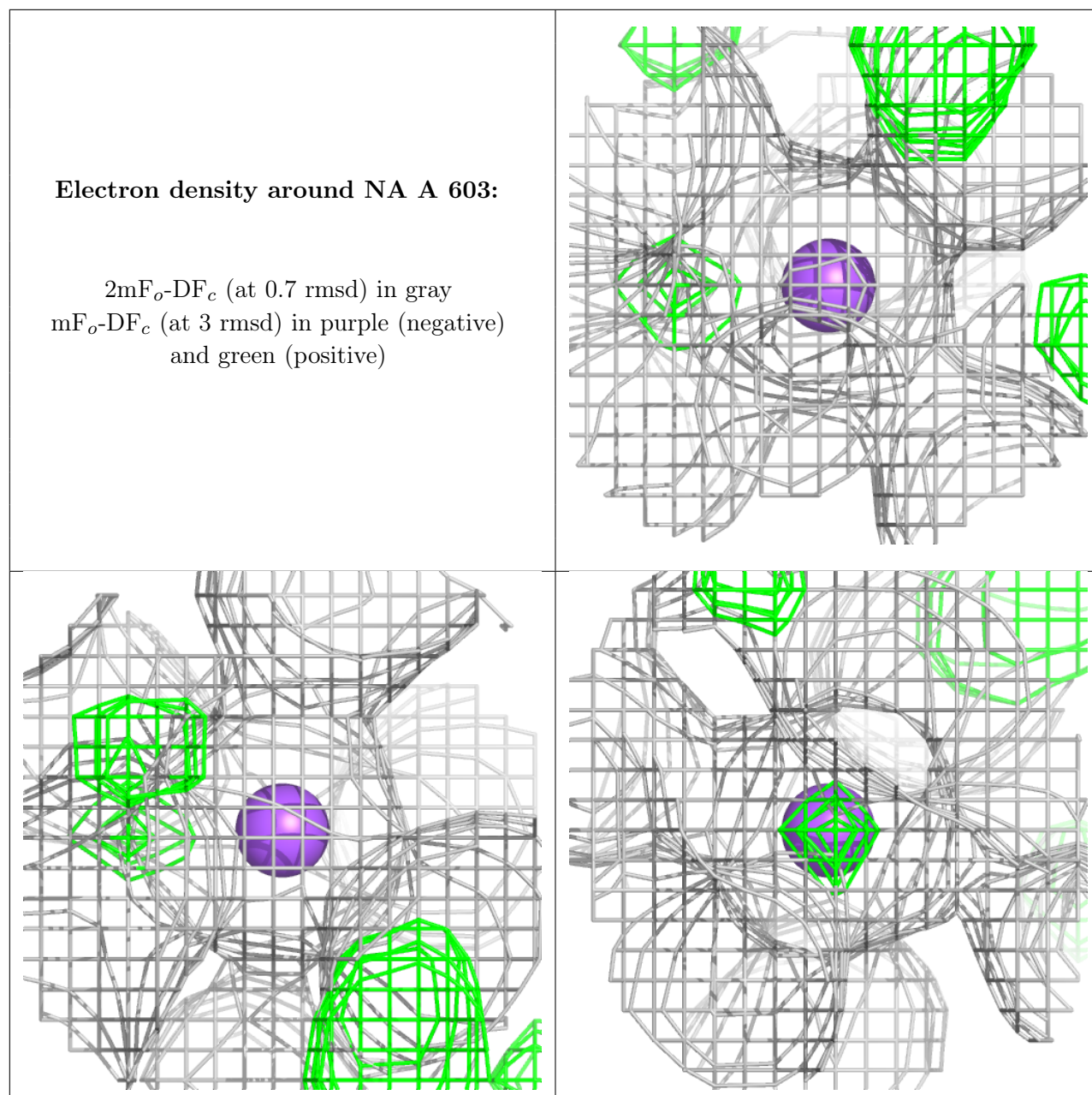
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around NA A 602:**

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





## 6.5 Other polymers ⓘ

There are no such residues in this entry.