



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:32 AM UTC

PDB ID : 9EVF / pdb\_00009evf  
Title : Citramalate lyase - ADP - pyruvate complex  
Authors : Ermler, U.; Berg, I.; Demmer, U.  
Deposited on : 2024-03-29  
Resolution : 1.40 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 : 4-5-2 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

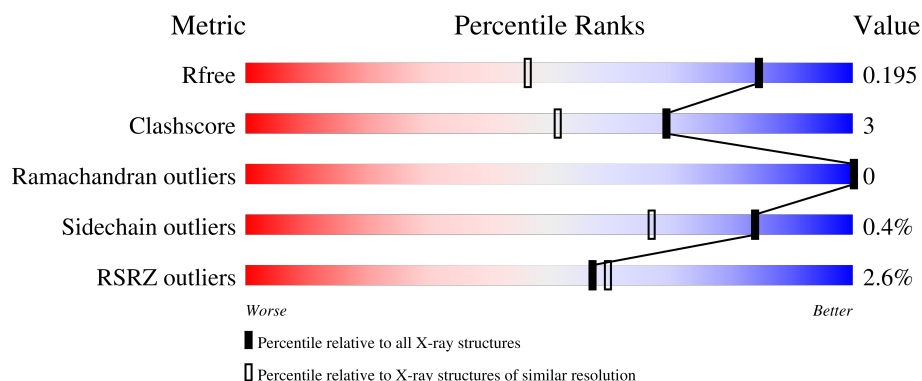
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.40 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))
$R_{free}$	180053	2563 (1.40-1.40)
Clashscore	190562	2660 (1.40-1.40)
Ramachandran outliers	187476	2611 (1.40-1.40)
Sidechain outliers	187428	2610 (1.40-1.40)
RSRZ outliers	180081	2561 (1.40-1.40)

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	456	<div> <div>2%</div> <div>91%</div> <div>9%</div> </div>
1	C	456	<div> <div>3%</div> <div>92%</div> <div>7%</div> </div>
2	B	105	<div> <div>4%</div> <div>93%</div> <div>6%</div> </div>
2	D	105	<div> <div>0%</div> <div>96%</div> <div>4%</div> </div>

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
5	PYR	A	504[A]	-	X	-	-
5	PYR	C	504[A]	-	X	-	-

## 2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 18501 atoms, of which 8776 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.

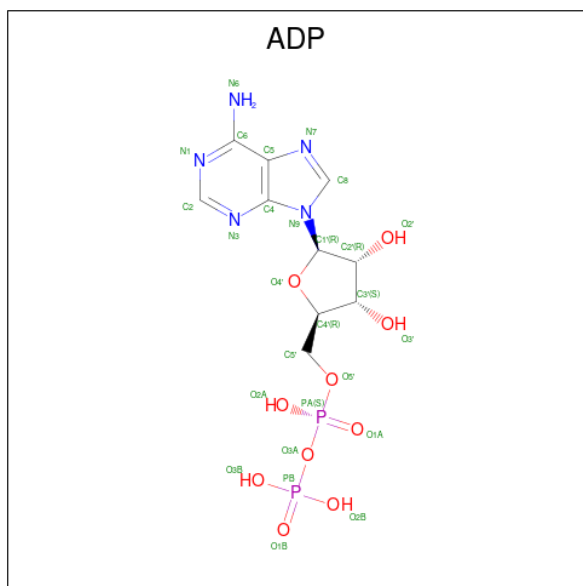
- Molecule 1 is a protein called Citramalate lyase, subunit A (cmlA).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	456	Total	C	H	N	O	S	0	21	0
			6983	2220	3472	597	673	21			
1	C	456	Total	C	H	N	O	S	0	17	0
			7002	2226	3478	601	675	22			

- Molecule 2 is a protein called Citramalate lyase, subunit B (cmlB).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	105	Total	C	H	N	O	S	0	6	0
			1756	558	884	151	160	3			
2	D	105	Total	C	H	N	O	S	0	0	0
			1702	543	856	146	154	3			

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ) (labeled as "Ligand of Interest" by depositor).

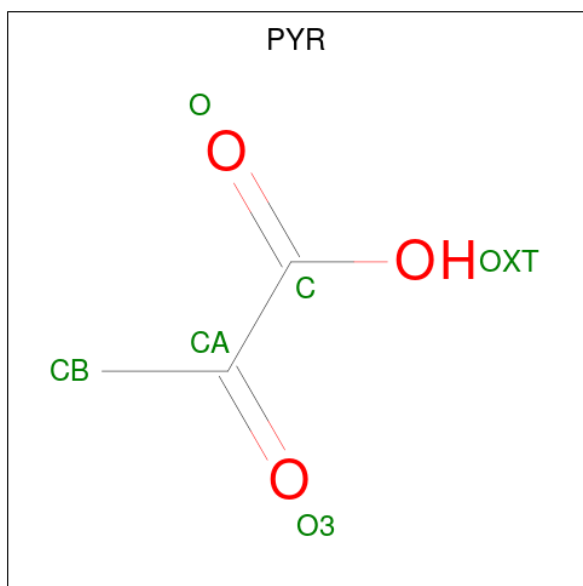


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	P	
			38	10	11	5	10	2	
3	C	1	Total	C	H	N	O	P	
			38	10	11	5	10	2	

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

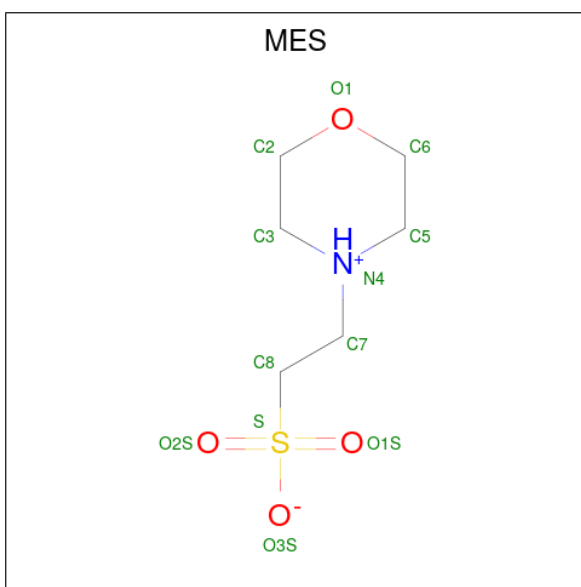
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mg		
			2	2	0	0
4	C	2	Total	Mg		
			2	2	0	0

- Molecule 5 is PYRUVIC ACID (CCD ID: PYR) (formula: C<sub>3</sub>H<sub>4</sub>O<sub>3</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O		
			18	6	6	6	0	1
5	C	1	Total	C	H	O		
			18	6	6	6	0	1

- Molecule 6 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: C<sub>6</sub>H<sub>13</sub>NO<sub>4</sub>S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
6	A	1	Total 25	C 6	H 13	N 1	O 4	S 1	0	1
6	A	1	Total 25	C 6	H 13	N 1	O 4	S 1	0	0
6	C	1	Total 25	C 6	H 13	N 1	O 4	S 1	0	0
6	D	1	Total 25	C 6	H 13	N 1	O 4	S 1	0	1

- Molecule 7 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	Cl	0	0
			1	1		

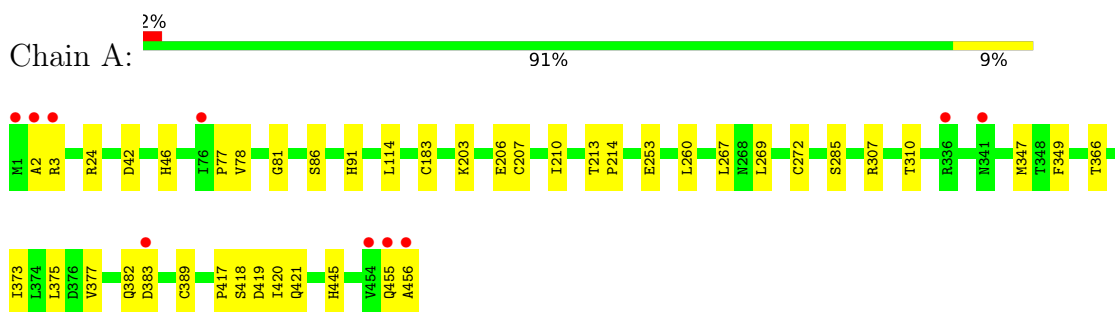
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	329	Total	O	0	3
			332	332		
8	B	67	Total	O	0	0
			67	67		
8	C	353	Total	O	0	3
			356	356		
8	D	86	Total	O	0	0
			86	86		

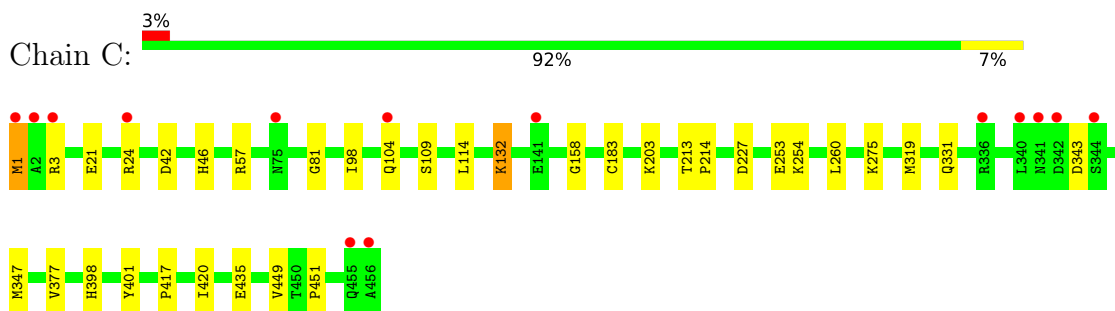
### 3 Residue-property plots [i](#)

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.

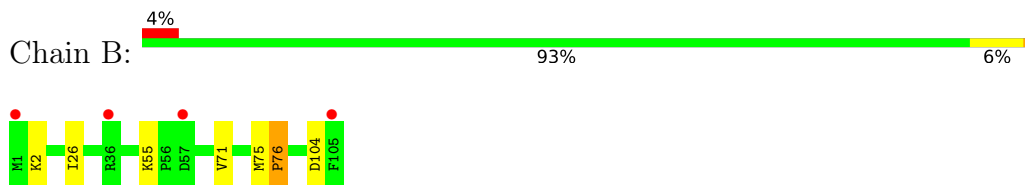
- Molecule 1: Citramalate lyase, subunit A (cmlA)



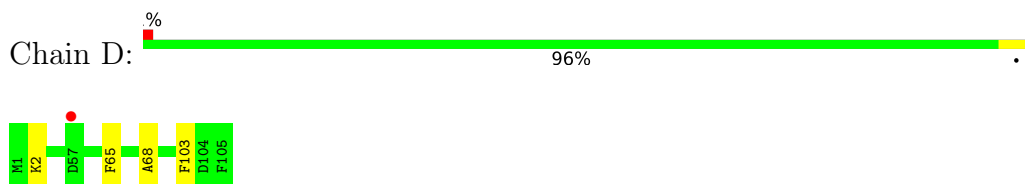
- Molecule 1: Citramalate lyase, subunit A (cmlA)



- Molecule 2: Citramalate lyase, subunit B (cmlB)



- Molecule 2: Citramalate lyase, subunit B (cmlB)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	107.21Å 108.34Å 98.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.10 – 1.40 38.10 – 1.40	Depositor EDS
% Data completeness (in resolution range)	98.7 (38.10-1.40) 98.9 (38.10-1.40)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.36 (at 1.40Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, $R_{free}$	0.178 , 0.195 0.178 , 0.195	Depositor DCC
$R_{free}$ test set	10923 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	19.0	Xtriage
Anisotropy	0.765	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 31.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.003 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	18501	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% 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 ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: MES, MG, ADP, PYR, CL

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	0.44	2/3668 (0.1%)	0.60	2/4981 (0.0%)
1	C	0.41	1/3651 (0.0%)	0.60	1/4952 (0.0%)
2	B	0.86	3/918 (0.3%)	1.07	10/1241 (0.8%)
2	D	0.37	0/864	0.59	0/1169
All	All	0.48	6/9101 (0.1%)	0.66	13/12343 (0.1%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	76	PRO	N-CA	17.25	1.68	1.47
2	B	55	LYS	C-N	11.52	1.51	1.33
2	B	75	MET	C-N	8.52	1.44	1.33
1	A	78[A]	VAL	N-CA	6.93	1.54	1.46
1	A	78[B]	VAL	N-CA	6.93	1.54	1.46
1	C	158	GLY	N-CA	5.03	1.52	1.45

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	75	MET	CA-C-N	15.33	135.55	119.76
2	B	75	MET	C-N-CA	15.33	135.55	119.76
2	B	75	MET	O-C-N	-9.85	113.87	121.55
2	B	55	LYS	CA-C-N	7.54	127.92	119.32
2	B	55	LYS	C-N-CA	7.54	127.92	119.32
2	B	76	PRO	CA-N-CD	-7.21	101.91	112.00
2	B	76	PRO	N-CA-CB	6.65	109.23	103.31
2	B	76	PRO	N-CD-CG	6.14	112.41	103.20
2	B	76	PRO	N-CA-C	-5.91	101.93	111.14
2	B	55	LYS	O-C-N	-5.69	115.06	121.43
1	A	77	PRO	CA-C-N	-5.52	115.91	123.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	77	PRO	C-N-CA	-5.52	115.91	123.14
1	C	343	ASP	CB-CA-C	5.03	119.96	109.95

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3511	3472	3371	29	0
1	C	3524	3478	3425	27	0
2	B	872	884	854	6	0
2	D	846	856	856	2	0
3	A	27	11	12	0	0
3	C	27	11	12	0	0
4	A	2	0	0	0	0
4	C	2	0	0	0	0
5	A	12	6	0	1	0
5	C	12	6	0	0	0
6	A	24	26	26	1	0
6	C	12	13	13	0	0
6	D	12	13	13	0	0
7	A	1	0	0	0	0
8	A	332	0	0	2	0
8	B	67	0	0	1	0
8	C	356	0	0	2	0
8	D	86	0	0	0	0
All	All	9725	8776	8582	61	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 (61) 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:B:76:PRO:N	2:B:76:PRO:CA	1.68	1.38
1:C:331:GLN:HG2	1:C:347[B]:MET:HE1	1.50	0.93
1:C:331:GLN:HG2	1:C:347[B]:MET:CE	2.03	0.88
1:A:310[B]:THR:HG22	1:A:419:ASP:HA	1.55	0.87
1:C:21:GLU:HG3	1:C:24[A]:ARG:HH22	1.42	0.84
1:A:455:GLN:OE1	1:A:456:ALA:N	2.17	0.77
2:B:76:PRO:N	2:B:76:PRO:C	2.49	0.70
1:A:310[B]:THR:HG23	1:A:389:CYS:SG	2.38	0.63
1:A:269:LEU:HD22	1:A:272[A]:CYS:SG	2.41	0.61
1:A:310[A]:THR:HG23	1:A:418:SER:O	2.00	0.61
1:C:21:GLU:HG3	1:C:24[A]:ARG:NH2	2.15	0.60
1:C:203:LYS:HE2	1:C:254:LYS:HE3	1.87	0.56
1:A:347:MET:HE3	8:A:604:HOH:O	2.03	0.56
2:B:2[B]:LYS:NZ	8:B:201:HOH:O	2.34	0.54
1:A:455:GLN:OE1	1:A:455:GLN:C	2.51	0.53
1:A:213:THR:HA	1:A:214:PRO:C	2.34	0.53
1:C:42:ASP:OD2	1:C:253:GLU:OE2	2.25	0.53
1:A:417:PRO:HD3	1:C:46:HIS:CG	2.43	0.53
2:B:26:ILE:HB	2:B:71:VAL:HG13	1.91	0.52
1:C:81:GLY:HA3	1:C:183[A]:CYS:SG	2.50	0.52
1:C:3:ARG:NE	1:C:3:ARG:HA	2.23	0.51
1:A:203:LYS:O	1:A:206:GLU:HG2	2.11	0.51
2:B:26:ILE:HB	2:B:71:VAL:CG1	2.40	0.50
1:C:347[A]:MET:HG2	1:C:377:VAL:HG22	1.92	0.50
1:C:203:LYS:HE2	1:C:254:LYS:CE	2.42	0.50
1:A:420:ILE:HG21	1:C:420:ILE:HG21	1.94	0.49
1:C:114:LEU:CD1	8:C:771:HOH:O	2.61	0.48
1:A:310[B]:THR:HG21	1:A:419:ASP:OD1	2.13	0.48
1:A:24:ARG:HH11	1:A:24:ARG:HG2	1.78	0.48
1:A:285:SER:OG	6:A:505[A]:MES:H22	2.13	0.47
2:D:65:PHE:HD2	2:D:68:ALA:HB3	1.79	0.47
1:C:3:ARG:CZ	1:C:3:ARG:HB2	2.45	0.47
1:C:254:LYS:HG3	1:C:260:LEU:HD22	1.97	0.47
1:A:46:HIS:CG	1:C:417:PRO:HD3	2.49	0.46
1:A:81:GLY:HA3	1:A:183[A]:CYS:SG	2.56	0.46
1:C:254:LYS:HA	1:C:254:LYS:HD3	1.80	0.45
1:A:86:SER:HB2	1:A:91:HIS:CE1	2.51	0.45
1:C:319:MET:HB2	1:C:401:TYR:CZ	2.51	0.45
5:A:504[B]:PYR:O	1:C:398:HIS:CE1	2.70	0.45
1:C:213:THR:HA	1:C:214:PRO:C	2.42	0.45
1:A:114[B]:LEU:HD12	1:A:445:HIS:NE2	2.32	0.44
1:C:109:SER:HA	1:C:449:VAL:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:3:ARG:HG2	1:C:3:ARG:HH11	1.82	0.44
1:A:382:GLN:NE2	1:A:421:GLN:HG3	2.33	0.43
1:C:114:LEU:HD12	8:C:771:HOH:O	2.18	0.43
1:A:2:ALA:HB3	8:A:638:HOH:O	2.18	0.43
1:A:260:LEU:HB2	1:A:267:LEU:HB2	2.00	0.43
1:A:349:PHE:CZ	1:A:375:LEU:HD13	2.53	0.43
1:A:42:ASP:OD2	1:A:253:GLU:OE2	2.37	0.42
2:B:2[A]:LYS:HD3	2:B:104:ASP:HA	2.01	0.42
1:C:132:LYS:HB3	1:C:435[A]:GLU:HB3	2.02	0.42
1:A:207:CYS:HB3	1:A:210:ILE:HD12	2.01	0.42
1:A:307:ARG:HA	1:A:377:VAL:O	2.20	0.41
1:A:3:ARG:N	1:A:3:ARG:HD2	2.35	0.41
1:C:57:ARG:HG3	1:C:98:ILE:HD11	2.01	0.41
1:C:1:MET:N	1:C:227:ASP:OD1	2.50	0.40
1:C:104[A]:GLN:HG2	1:C:451:PRO:O	2.22	0.40
2:D:2:LYS:HA	2:D:103:PHE:O	2.21	0.40

There are no symmetry-related clashes.

## 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	475/456 (104%)	465 (98%)	10 (2%)	0	100	100
1	C	471/456 (103%)	459 (98%)	12 (2%)	0	100	100
2	B	109/105 (104%)	107 (98%)	2 (2%)	0	100	100
2	D	103/105 (98%)	101 (98%)	2 (2%)	0	100	100
All	All	1158/1122 (103%)	1132 (98%)	26 (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	379/360 (105%)	377 (100%)	2 (0%)	81	61
1	C	377/360 (105%)	374 (99%)	3 (1%)	73	48
2	B	99/93 (106%)	99 (100%)	0	100	100
2	D	93/93 (100%)	93 (100%)	0	100	100
All	All	948/906 (105%)	943 (100%)	5 (0%)	84	61

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

Mol	Chain	Res	Type
1	A	383[A]	ASP
1	A	383[B]	ASP
1	C	1	MET
1	C	132	LYS
1	C	275	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	104	GLN
1	A	382	GLN
1	A	403	ASN
2	B	60	HIS
2	D	83	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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 5 are monoatomic - leaving 10 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	$\# Z  > 2$
6	MES	C	505	-	12,12,12	1.46	3 (25%)	15,16,16	2.09	4 (26%)
3	ADP	C	501	4	28,29,29	1.03	2 (7%)	43,45,45	0.62	1 (2%)
6	MES	A	505[A]	-	12,12,12	1.07	1 (8%)	15,16,16	2.06	5 (33%)
6	MES	A	506	-	12,12,12	1.26	1 (8%)	15,16,16	1.63	3 (20%)
6	MES	D	201[B]	-	12,12,12	1.08	1 (8%)	15,16,16	2.03	5 (33%)
5	PYR	C	504[A]	-	5,5,5	1.87	2 (40%)	3,6,6	1.76	1 (33%)
5	PYR	C	504[B]	-	5,5,5	1.33	0	3,6,6	1.85	1 (33%)
5	PYR	A	504[A]	-	5,5,5	1.79	1 (20%)	3,6,6	1.50	1 (33%)
5	PYR	A	504[B]	-	5,5,5	2.57	3 (60%)	3,6,6	0.67	0
3	ADP	A	501	4	28,29,29	0.96	2 (7%)	43,45,45	0.61	1 (2%)

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
6	MES	C	505	-	-	0/6/14/14	0/1/1/1
3	ADP	C	501	4	-	4/16/32/32	0/3/3/3
6	MES	A	505[A]	-	-	0/6/14/14	0/1/1/1
6	MES	A	506	-	-	0/6/14/14	0/1/1/1
6	MES	D	201[B]	-	-	0/6/14/14	0/1/1/1
5	PYR	C	504[A]	-	-	3/4/4/4	-
5	PYR	C	504[B]	-	-	0/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PYR	A	504[A]	-	-	4/4/4/4	-
5	PYR	A	504[B]	-	-	0/4/4/4	-
3	ADP	A	501	4	-	3/16/32/32	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	504[B]	PYR	O-C	3.86	1.32	1.22
3	C	501	ADP	PB-O3B	-3.84	1.40	1.54
5	A	504[A]	PYR	O3-CA	-3.41	1.15	1.23
5	A	504[B]	PYR	O3-CA	-3.07	1.16	1.23
6	A	505[A]	MES	C8-S	3.01	1.81	1.77
6	D	201[B]	MES	C8-S	2.99	1.81	1.77
3	A	501	ADP	PB-O3B	-2.86	1.44	1.54
5	C	504[A]	PYR	O3-CA	-2.85	1.16	1.23
5	A	504[B]	PYR	OXT-C	-2.72	1.23	1.30
6	C	505	MES	O1S-S	-2.72	1.37	1.45
3	A	501	ADP	PB-O2B	-2.65	1.45	1.54
6	A	506	MES	O3S-S	-2.33	1.38	1.47
5	C	504[A]	PYR	OXT-C	-2.29	1.24	1.30
6	C	505	MES	C8-S	2.17	1.80	1.77
3	C	501	ADP	PB-O2B	-2.11	1.47	1.54
6	C	505	MES	O3S-S	-2.08	1.39	1.47

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	201[B]	MES	O2S-S-C8	5.19	114.57	106.73
6	A	505[A]	MES	O1S-S-C8	5.16	114.52	106.73
6	C	505	MES	O1S-S-C8	4.72	113.86	106.73
6	C	505	MES	O3S-S-O1S	-4.17	100.97	111.40
6	A	505[A]	MES	O2S-S-C8	3.66	112.26	106.73
6	A	506	MES	O3S-S-C8	3.65	113.14	106.00
6	D	201[B]	MES	O1S-S-C8	3.58	112.13	106.73
6	C	505	MES	O3S-S-C8	3.04	111.95	106.00
6	A	506	MES	O1S-S-C8	2.98	111.23	106.73
5	C	504[B]	PYR	OXT-C-CA	2.91	121.65	113.59
3	C	501	ADP	O2A-PA-O3A	2.81	114.86	107.27
5	C	504[A]	PYR	OXT-C-CA	2.63	120.88	113.59
6	C	505	MES	O2S-S-C8	2.49	110.49	106.73
6	A	505[A]	MES	O2S-S-O1S	-2.40	106.02	113.82
6	D	201[B]	MES	O2S-S-O1S	-2.35	106.19	113.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	506	MES	O3S-S-O1S	-2.28	105.70	111.40
6	A	505[A]	MES	O3S-S-C8	2.24	110.38	106.00
5	A	504[A]	PYR	OXT-C-CA	2.22	119.76	113.59
3	A	501	ADP	O3B-PB-O2B	2.08	115.59	107.80
6	A	505[A]	MES	O3S-S-O1S	-2.07	106.23	111.40
6	D	201[B]	MES	O3S-S-C8	2.05	110.02	106.00
6	D	201[B]	MES	O3S-S-O2S	-2.00	106.39	111.40

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	504[A]	PYR	O-C-CA-O3
5	A	504[A]	PYR	OXT-C-CA-O3
5	A	504[A]	PYR	OXT-C-CA-CB
5	C	504[A]	PYR	OXT-C-CA-CB
5	A	504[A]	PYR	O-C-CA-CB
5	C	504[A]	PYR	O-C-CA-CB
3	C	501	ADP	PB-O3A-PA-O2A
3	A	501	ADP	PA-O3A-PB-O1B
3	A	501	ADP	PB-O3A-PA-O1A
3	C	501	ADP	PA-O3A-PB-O2B
3	C	501	ADP	PB-O3A-PA-O1A
5	C	504[A]	PYR	O-C-CA-O3
3	A	501	ADP	PB-O3A-PA-O2A
3	C	501	ADP	O4'-C4'-C5'-O5'

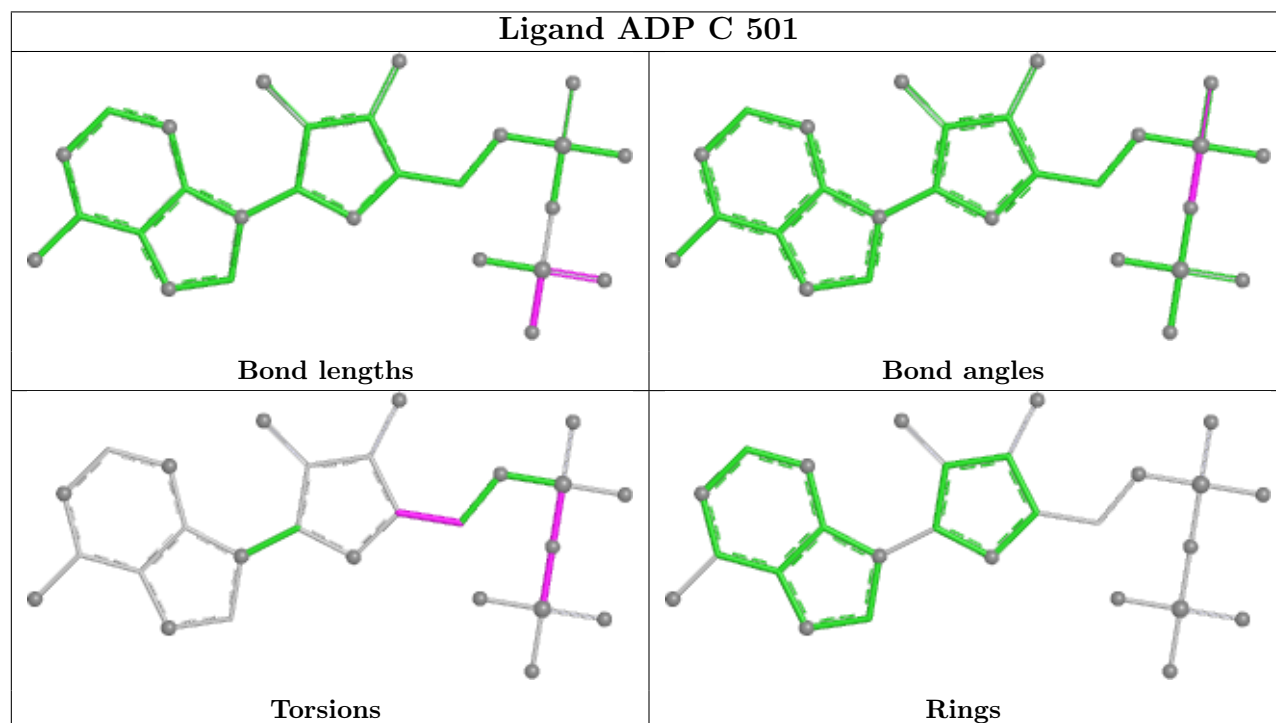
There are no ring outliers.

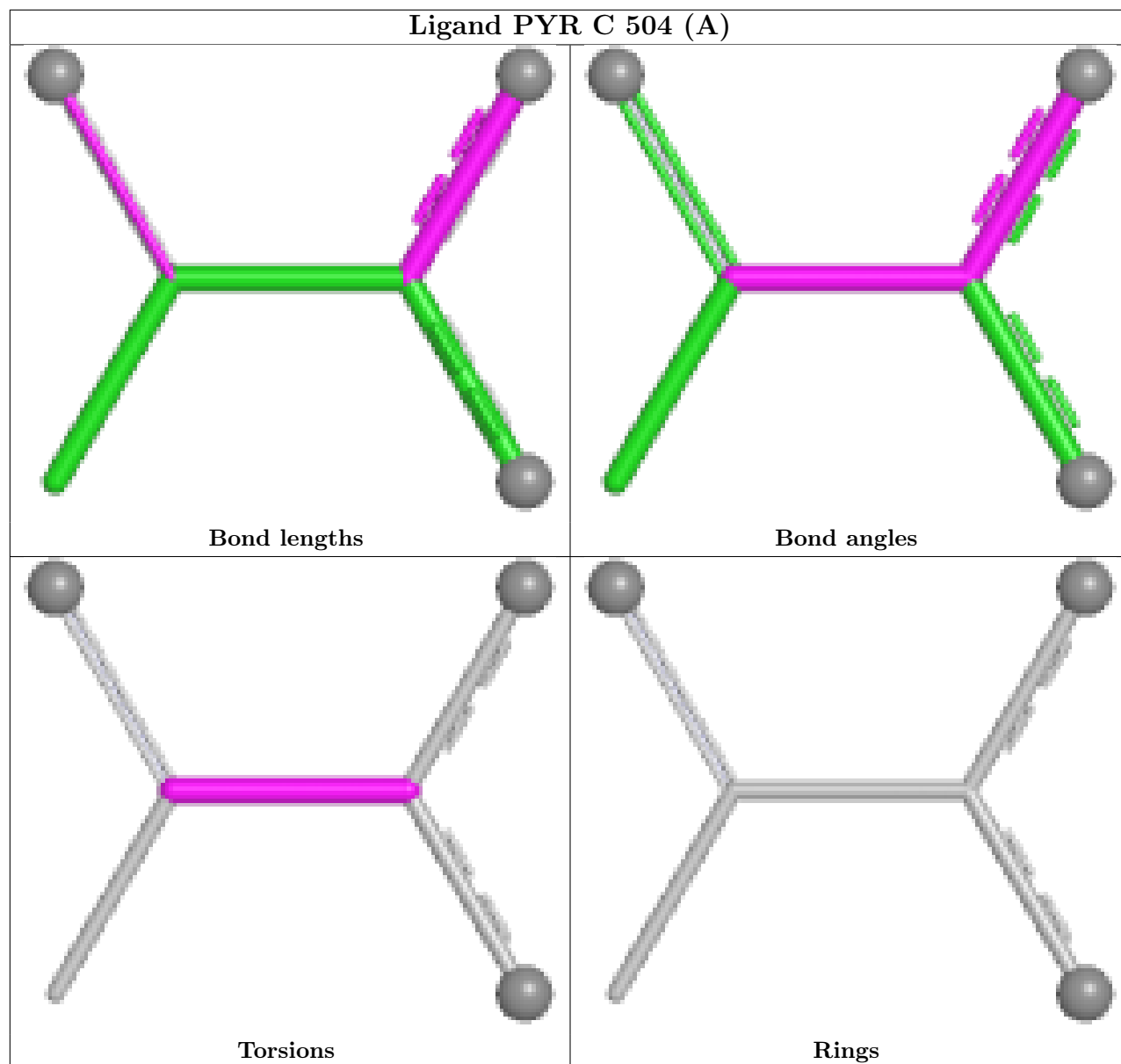
2 monomers are involved in 2 short contacts:

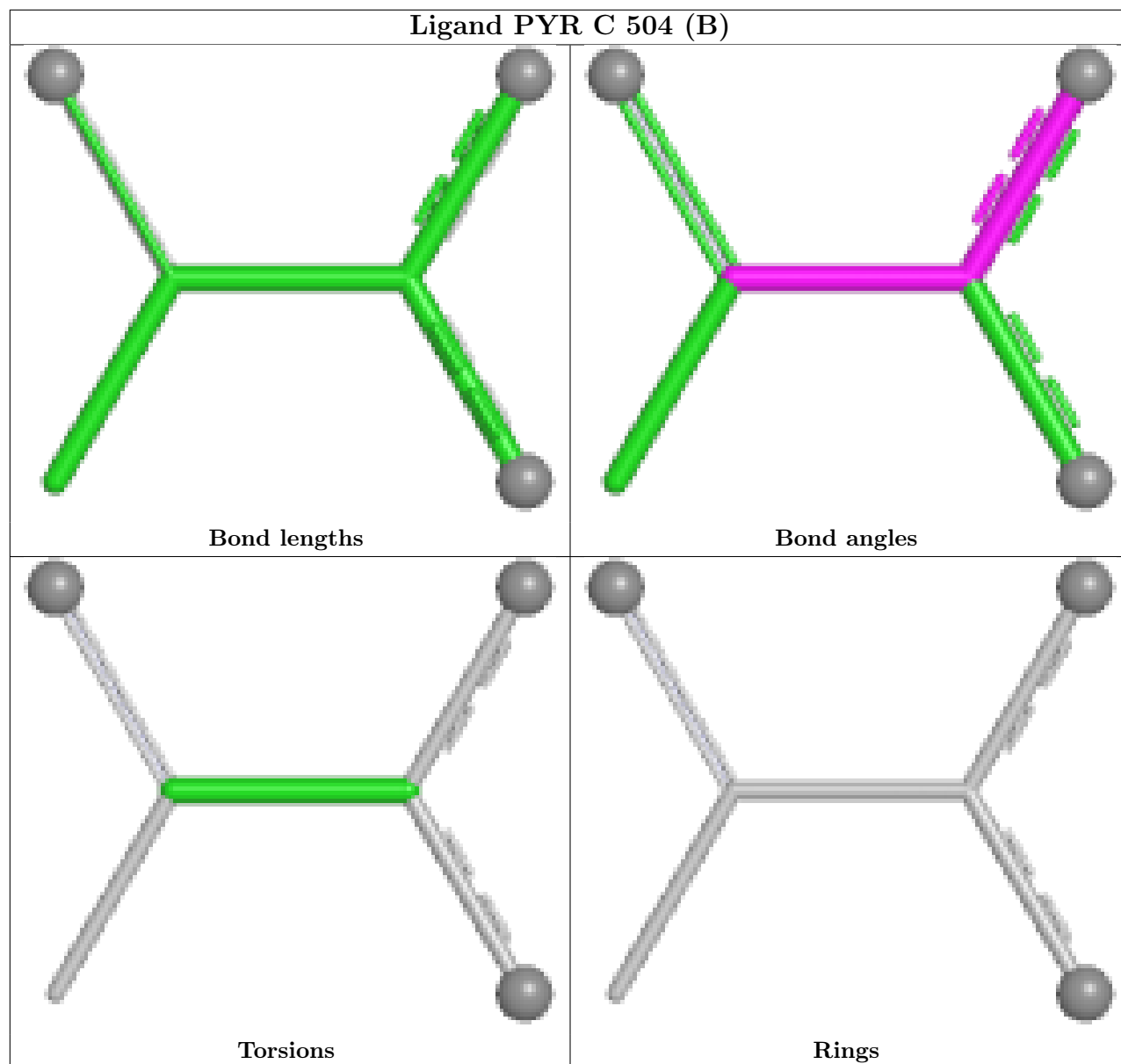
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	505[A]	MES	1	0
5	A	504[B]	PYR	1	0

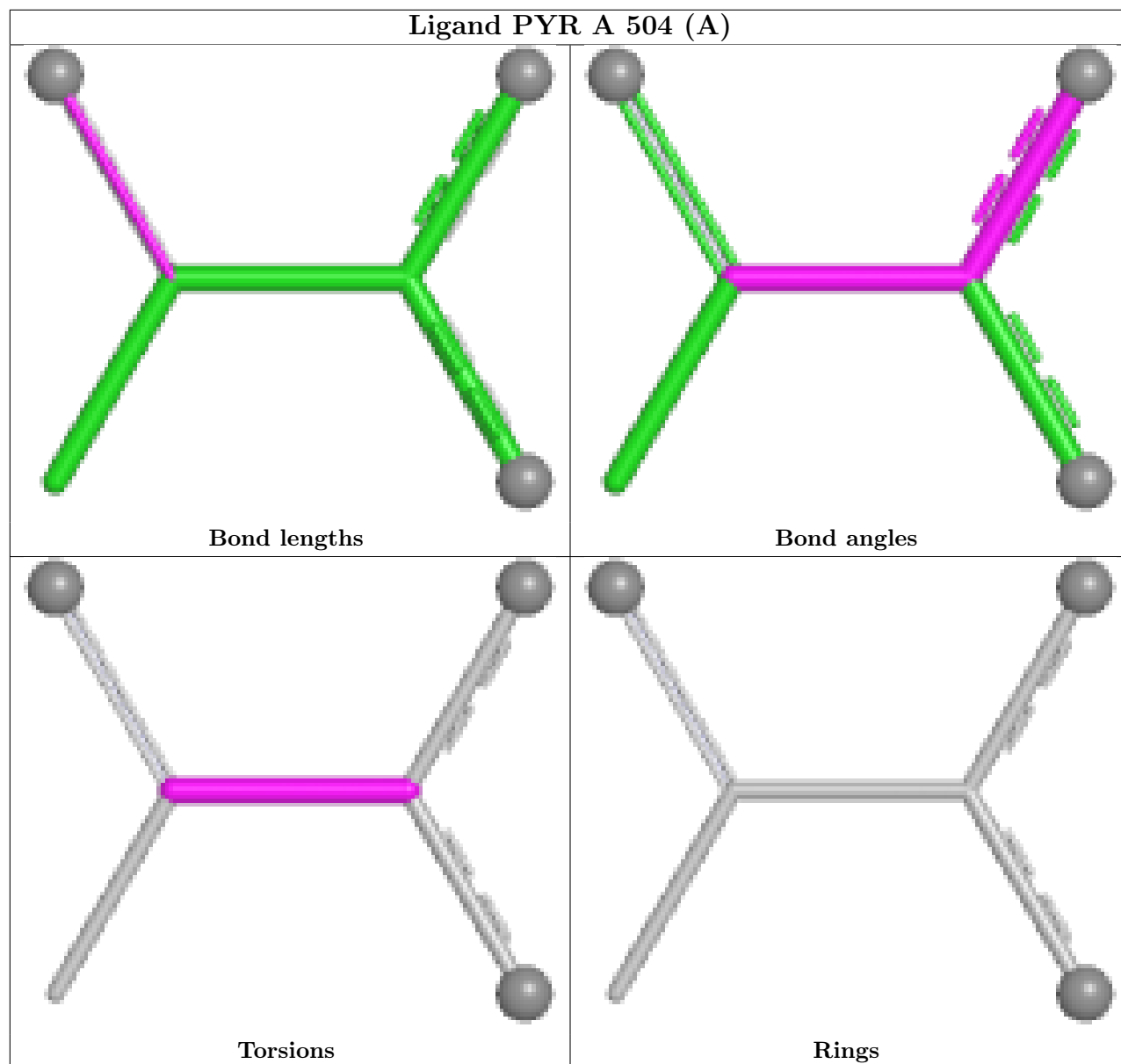
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

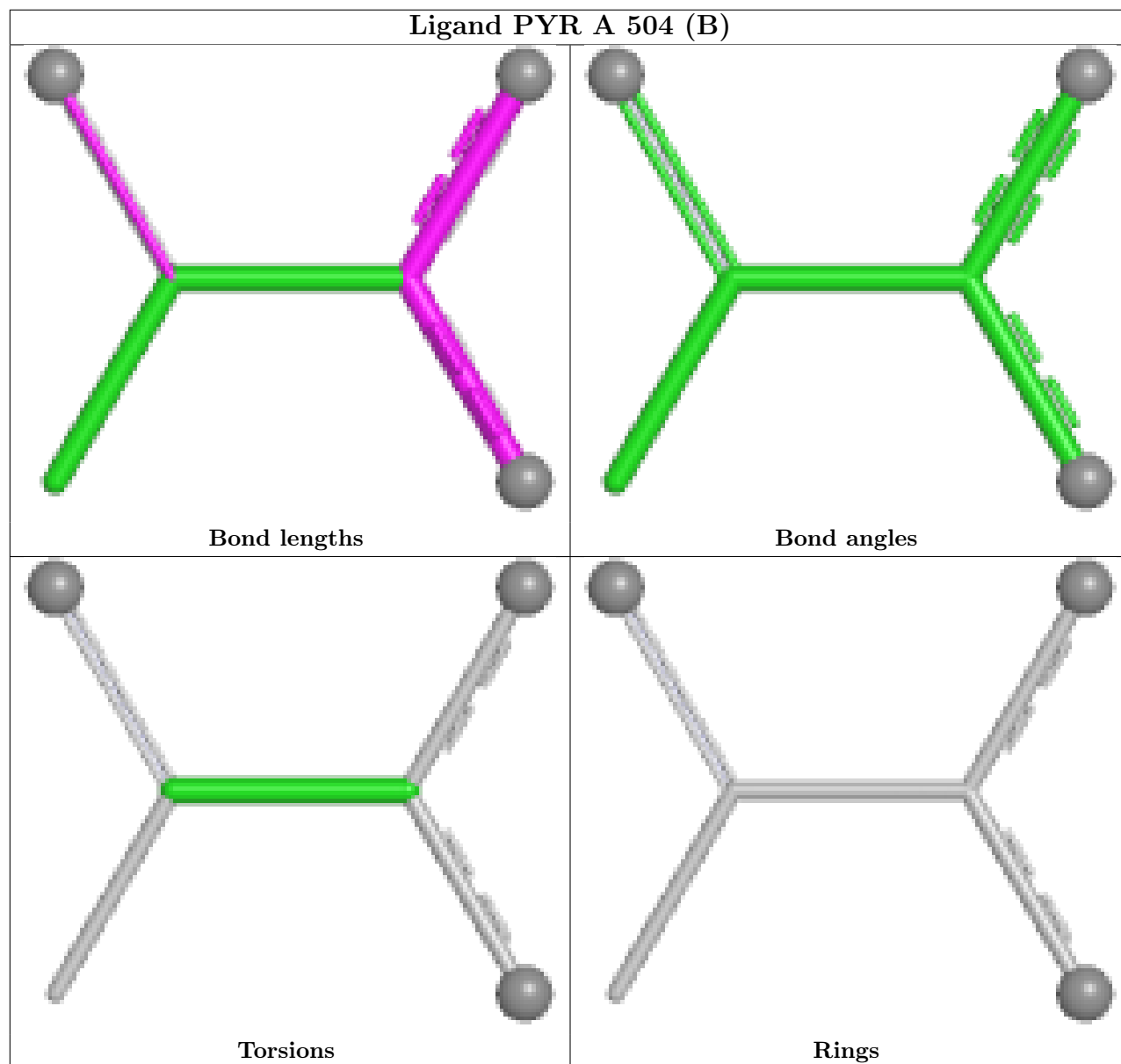
average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

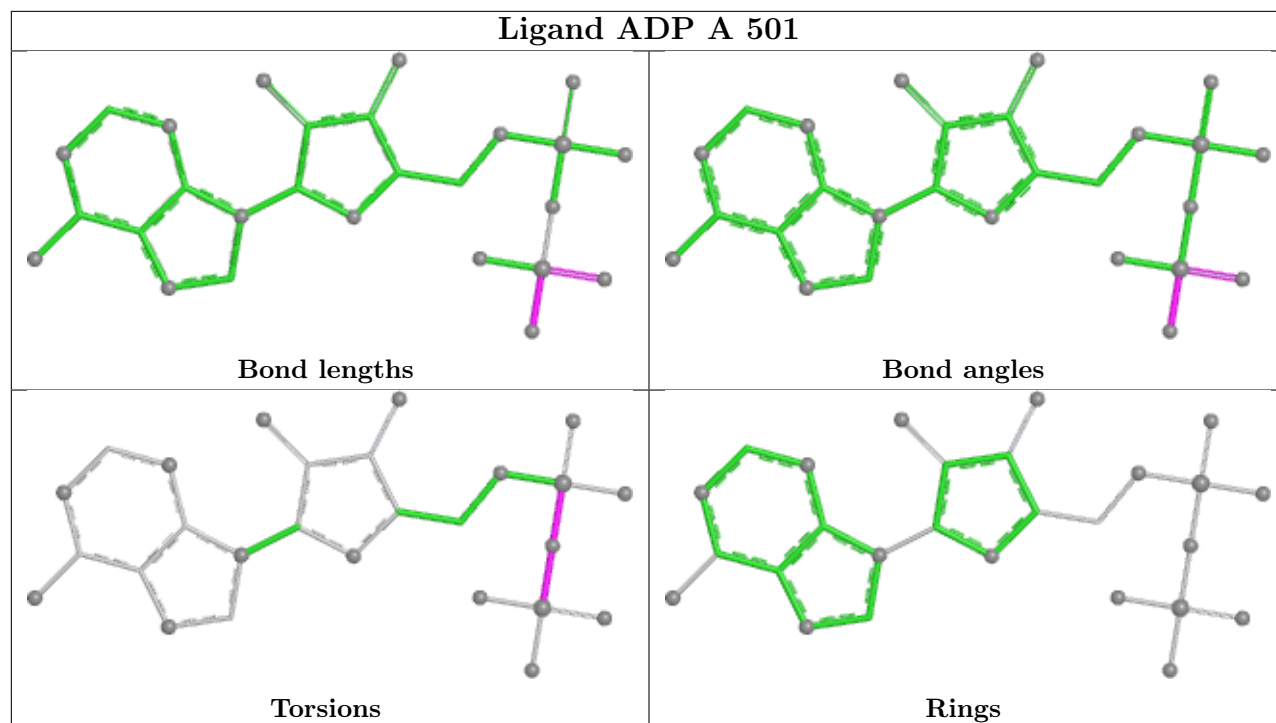












## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	456/456 (100%)	0.06	10 (2%) 62 65	11, 25, 39, 108	10 (2%)
1	C	456/456 (100%)	0.06	14 (3%) 51 54	11, 25, 37, 87	11 (2%)
2	B	105/105 (100%)	0.17	4 (3%) 44 46	16, 28, 38, 58	3 (2%)
2	D	105/105 (100%)	-0.05	1 (0%) 79 83	21, 25, 36, 45	0
All	All	1122/1122 (100%)	0.06	29 (2%) 57 60	11, 25, 38, 108	24 (2%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	2	ALA	8.0
1	A	456	ALA	6.9
1	C	1	MET	5.6
1	C	456	ALA	5.4
2	B	105	PHE	4.5
1	A	2	ALA	4.4
1	C	340	LEU	4.2
1	A	454	VAL	3.8
1	C	3	ARG	3.7
1	A	455	GLN	3.5
1	A	3	ARG	3.5
1	A	1	MET	3.4
1	C	342	ASP	3.3
2	B	1	MET	3.2
1	A	341	ASN	3.2
1	C	341	ASN	3.1
1	C	24[A]	ARG	2.9
1	C	455	GLN	2.8
1	A	336	ARG	2.7
2	B	57[A]	ASP	2.6
1	C	141[A]	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
2	D	57	ASP	2.5
1	C	336	ARG	2.5
1	A	383[A]	ASP	2.5
2	B	36	ARG	2.5
1	A	76	ILE	2.4
1	C	344	SER	2.3
1	C	75	ASN	2.2
1	C	104[A]	GLN	2.1

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

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

## 6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 6.4 Ligands [i](#)

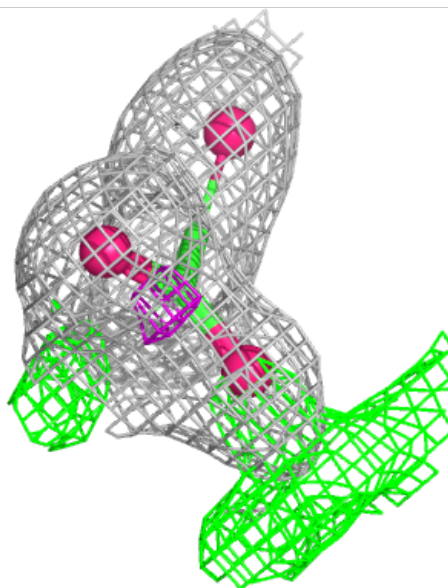
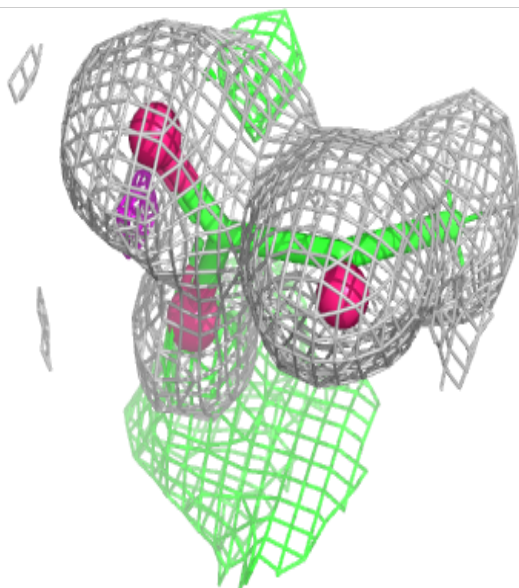
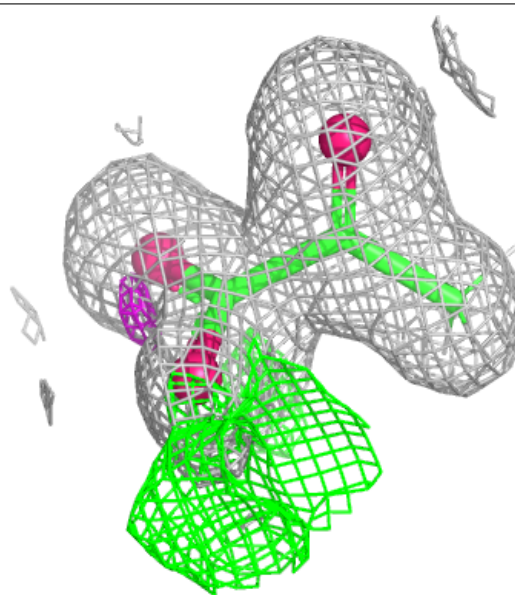
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	PYR	A	504[A]	6/6	0.81	0.20	19,31,35,39	9
5	PYR	A	504[B]	6/6	0.81	0.20	16,25,30,40	9
6	MES	A	505[A]	12/12	0.85	0.16	24,44,54,90	25
6	MES	D	201[B]	12/12	0.85	0.15	29,42,53,88	25
5	PYR	C	504[A]	6/6	0.88	0.13	13,16,26,40	9
5	PYR	C	504[B]	6/6	0.88	0.13	19,23,31,40	9
4	MG	A	502	1/1	0.90	0.09	19,19,19,19	0
4	MG	C	502	1/1	0.91	0.11	17,17,17,17	0
4	MG	A	503	1/1	0.93	0.09	18,18,18,18	0
4	MG	C	503	1/1	0.94	0.07	19,19,19,19	0
7	CL	A	507	1/1	0.95	0.11	28,28,28,28	0
6	MES	A	506	12/12	0.96	0.08	19,28,37,41	0
3	ADP	C	501	27/27	0.96	0.08	15,20,29,34	0
3	ADP	A	501	27/27	0.96	0.08	17,21,30,36	0
6	MES	C	505	12/12	0.97	0.08	20,30,39,44	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.

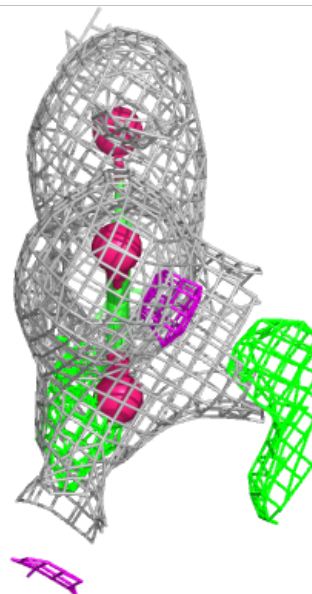
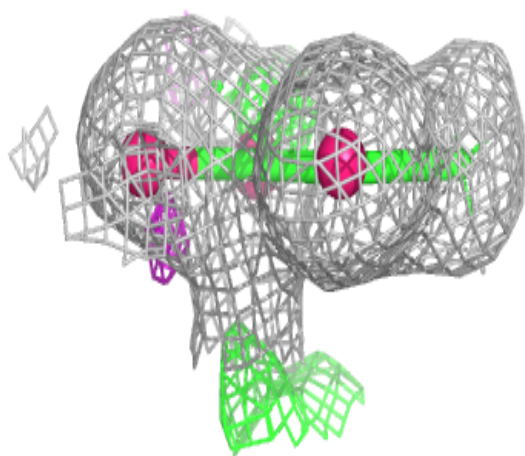
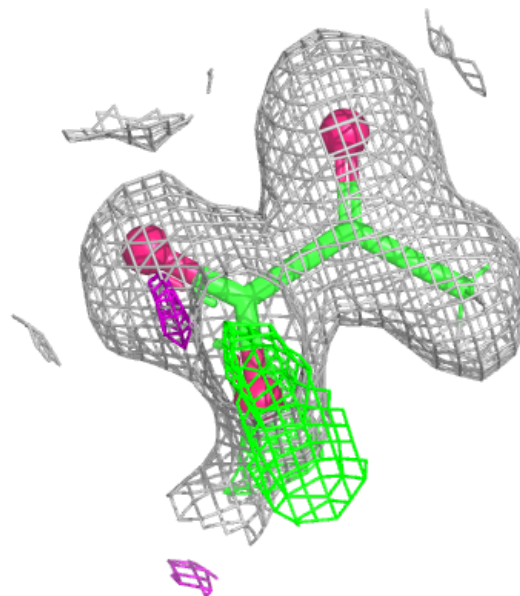
**Electron density around PYR A 504 (A):**

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



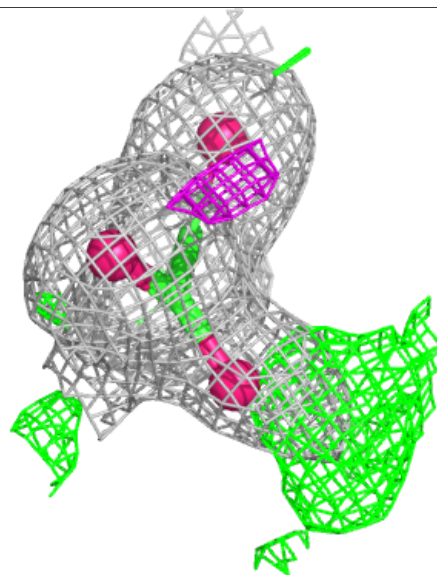
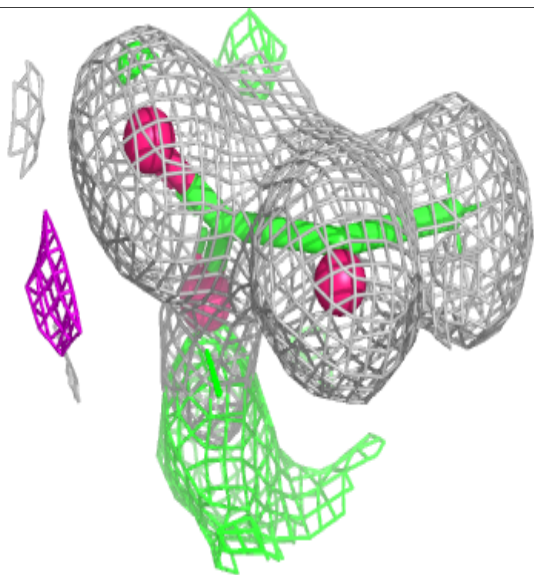
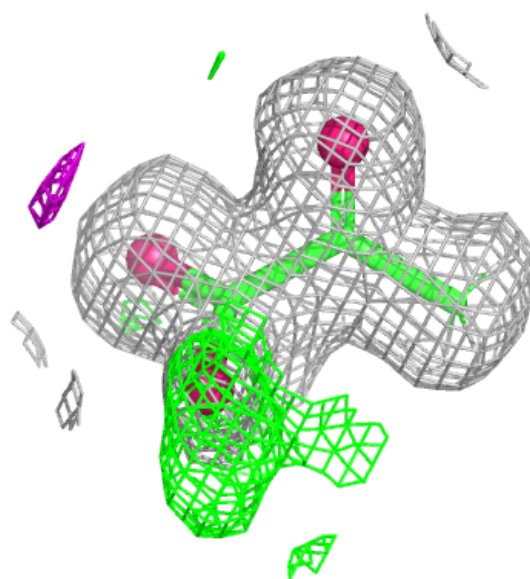
**Electron density around PYR A 504 (B):**

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



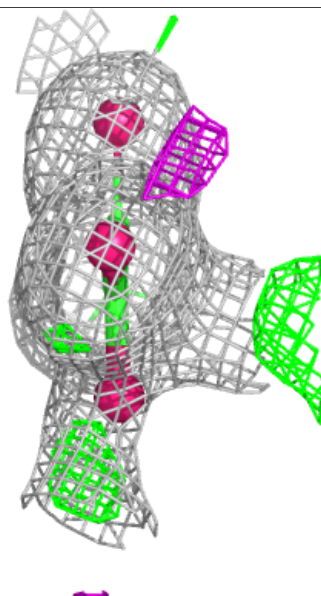
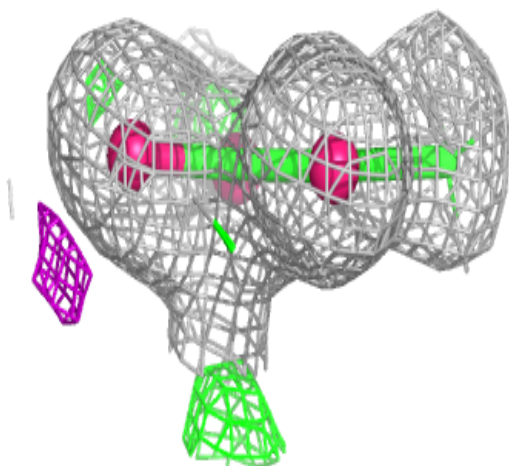
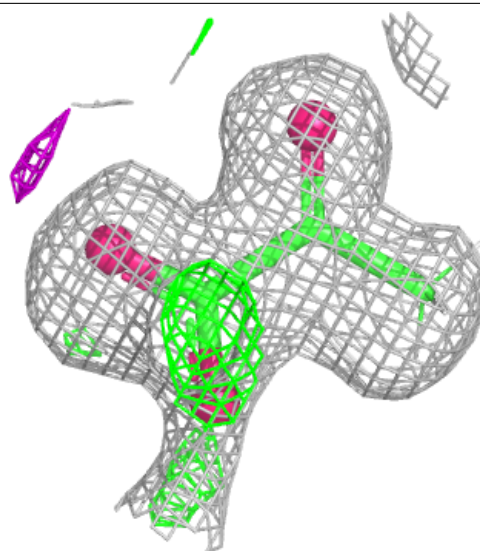
**Electron density around PYR C 504 (A):**

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



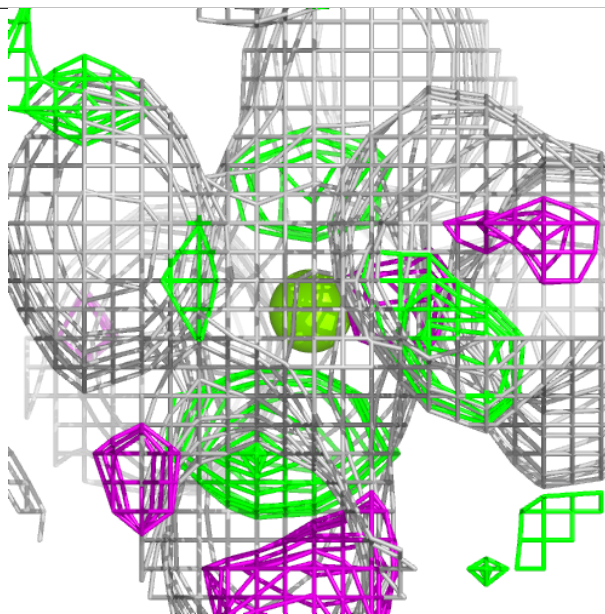
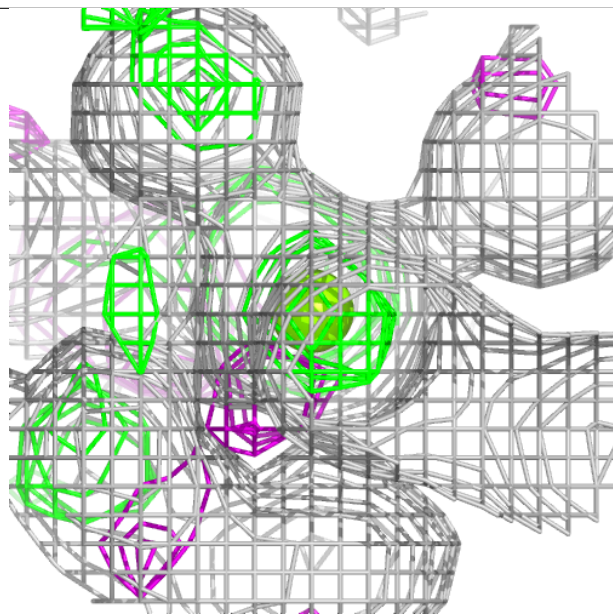
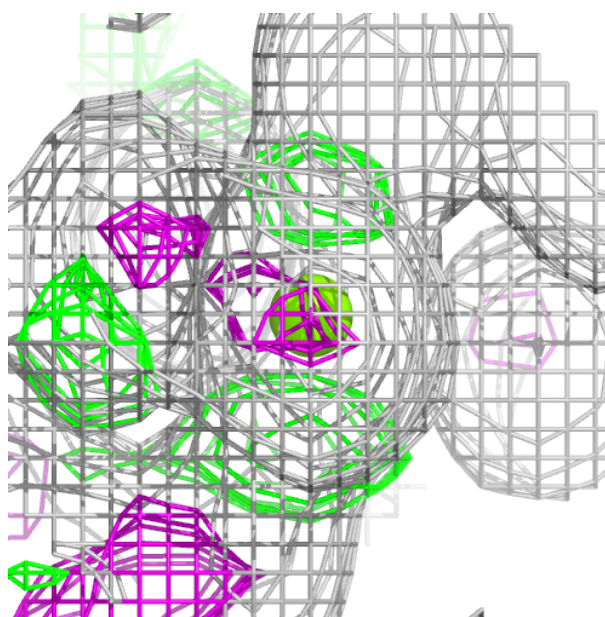
**Electron density around PYR C 504 (B):**

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and green (positive)



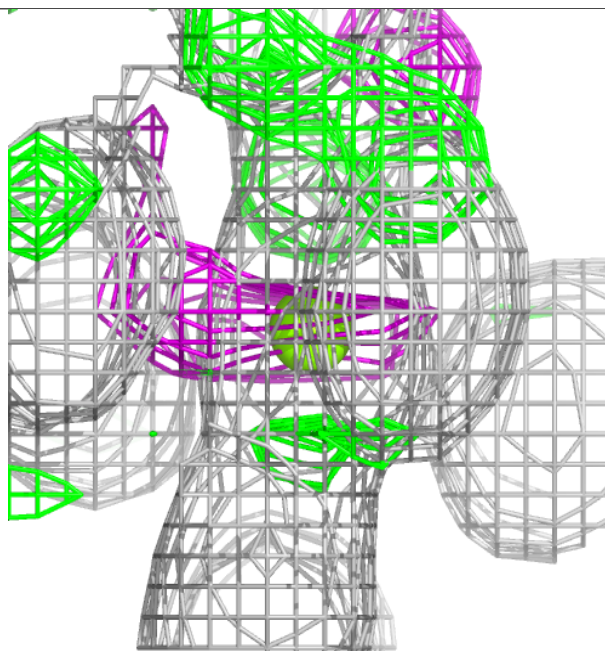
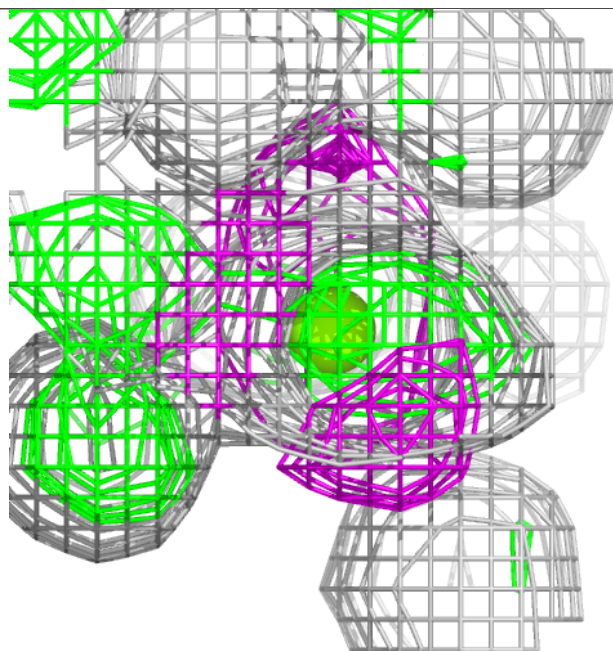
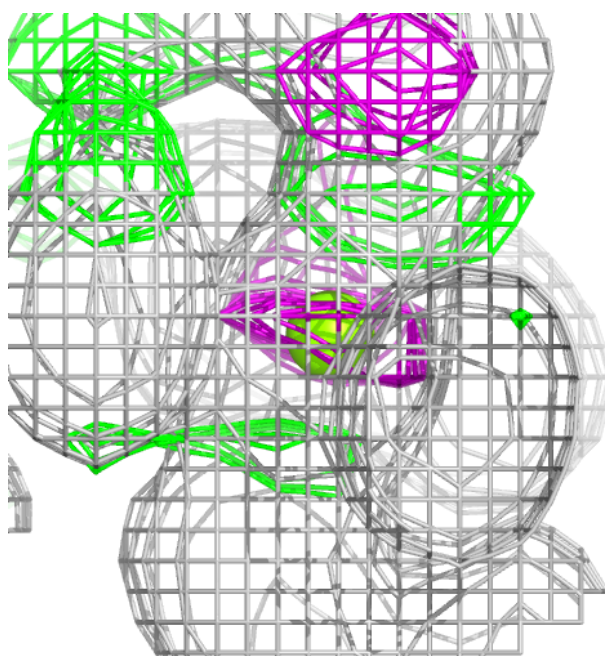
**Electron density around MG A 502:**

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and green (positive)



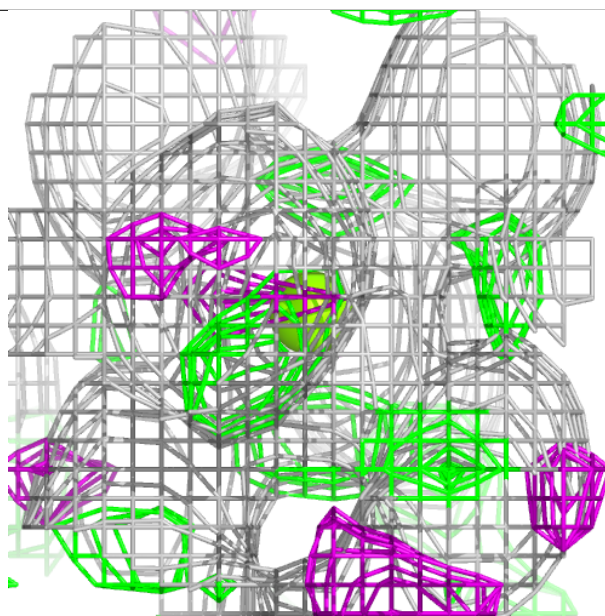
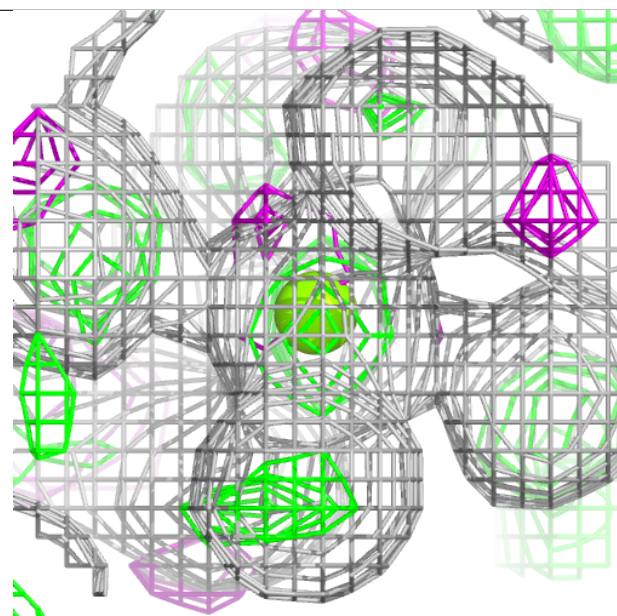
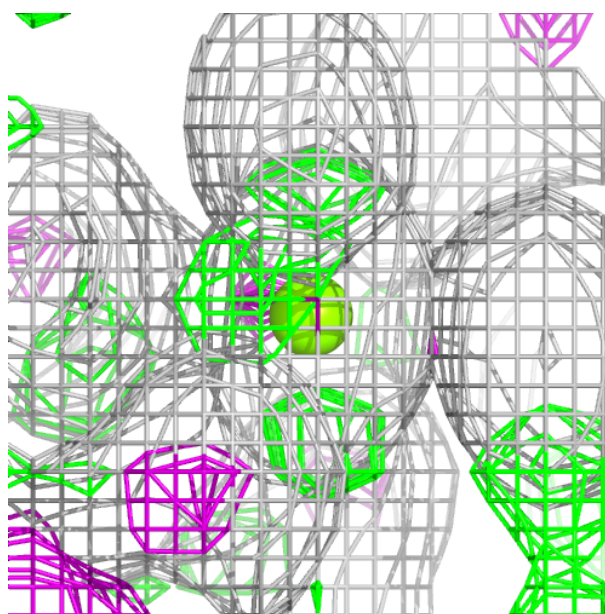
**Electron density around MG C 502:**

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and green (positive)



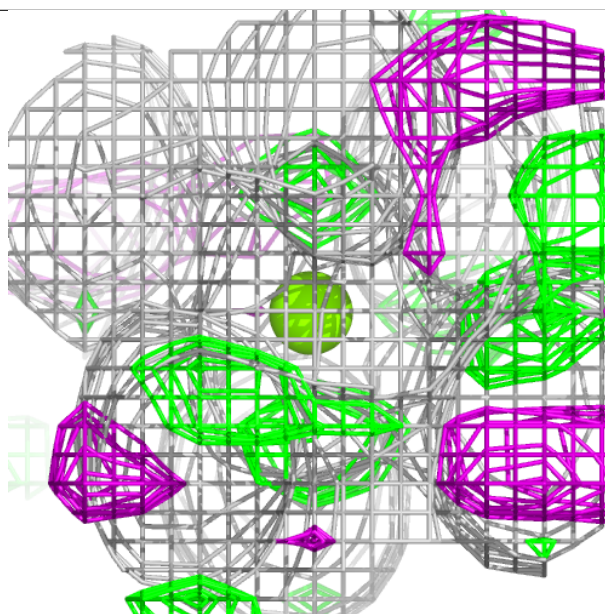
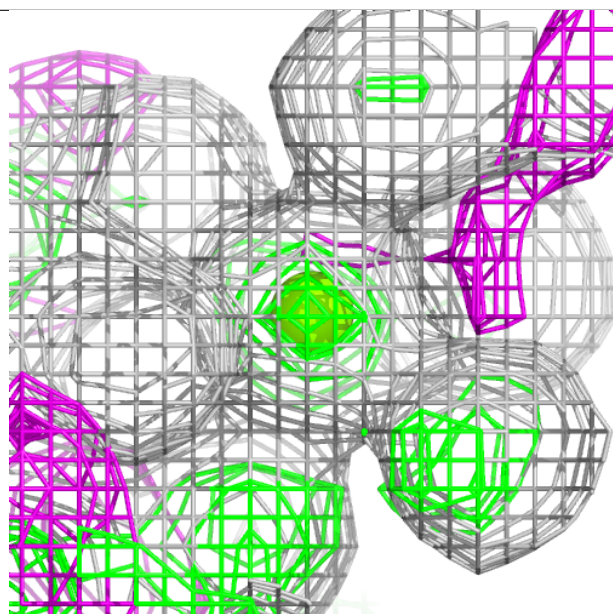
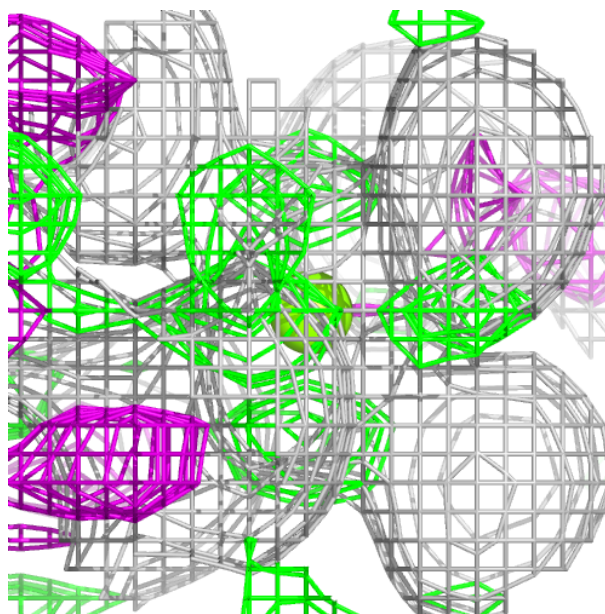
**Electron density around MG A 503:**

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and green (positive)



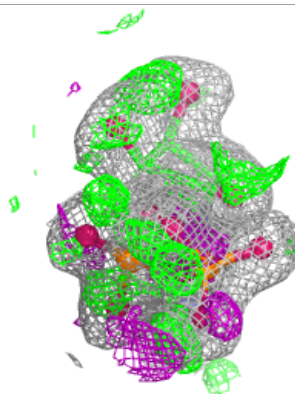
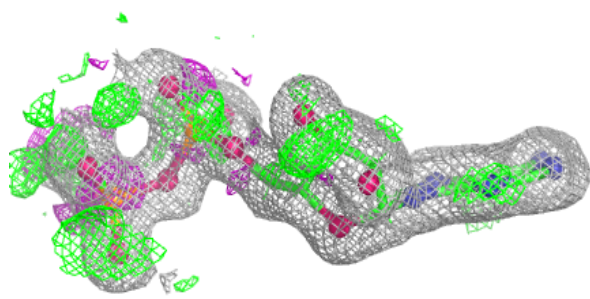
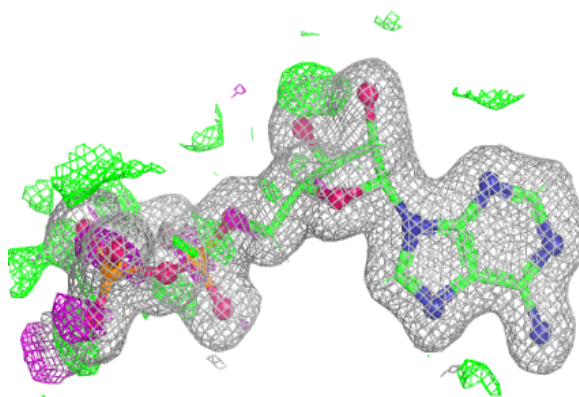
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and green (positive)

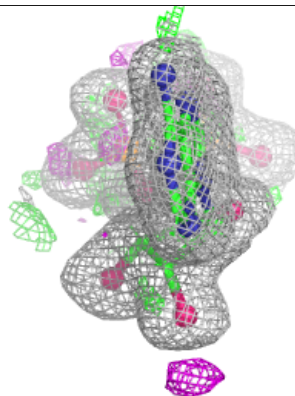
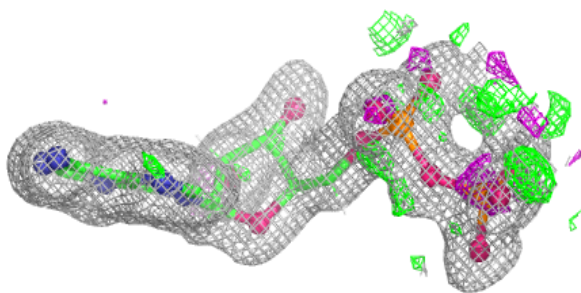
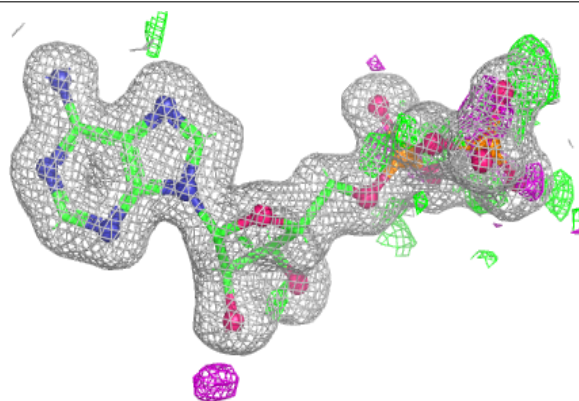


**Electron density around ADP C 501:**

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

**Electron density around ADP 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)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.