



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 07:44 AM UTC

PDB ID : 7LZ4 / pdb_00007lz4
Title : Crystal structure of A211D mutant of Protein Kinase A RIa subunit, a Carney Complex mutation
Authors : Del Rio, J.; Wu, J.; Taylor, S.S.
Deposited on : 2021-03-08
Resolution : 4.16 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

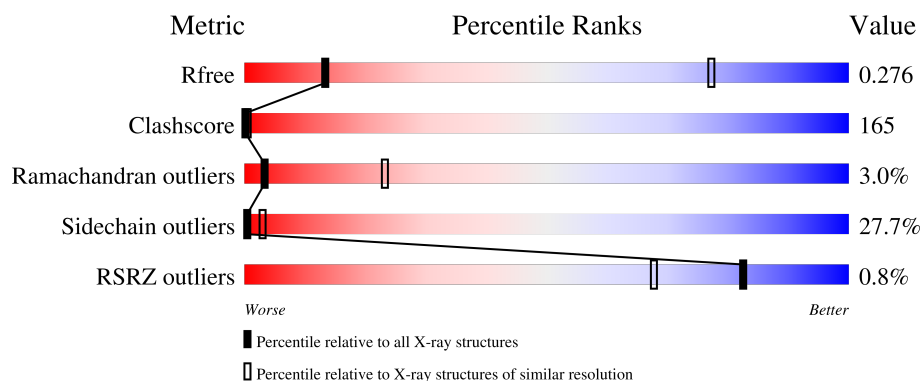
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.16 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.



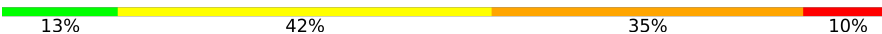
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1077 (4.50-3.82)
Clashscore	190562	1019 (4.48-3.84)
Ramachandran outliers	187476	1035 (4.50-3.82)
Sidechain outliers	187428	1022 (4.50-3.82)
RSRZ outliers	180081	1074 (4.50-3.82)

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	269	2% 12% 47% 31% 9%
1	B	269	% 11% 48% 33% 7%
1	C	269	% 12% 44% 32% 12%
1	D	269	% 11% 51% 29% 8%
1	E	269	% 11% 55% 29%

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Mol	Chain	Length	Quality of chain
1	F	269	
1	G	269	
1	H	269	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	CMP	A	401	-	-	X	-
2	CMP	A	402	-	-	X	-
2	CMP	B	401	-	-	X	-
2	CMP	B	402	-	-	X	-
2	CMP	C	401	-	-	X	-
2	CMP	C	402	-	-	X	-
2	CMP	D	401	-	-	X	-
2	CMP	D	402	-	-	X	-
2	CMP	E	401	-	-	X	-
2	CMP	E	402	-	-	X	-
2	CMP	F	401	-	-	X	-
2	CMP	F	402	-	-	X	-
2	CMP	G	401	-	-	X	-
2	CMP	G	402	-	-	X	-
2	CMP	H	401	-	-	X	-
2	CMP	H	402	-	-	X	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 17138 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.

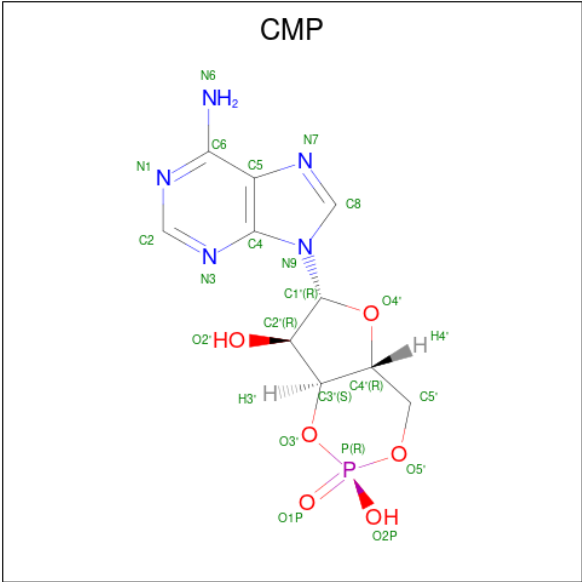
- Molecule 1 is a protein called cAMP-dependent protein kinase type I-alpha regulatory subunit, N-terminally processed.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	0	0	0
			2100	1333	358	401	8			
1	B	268	Total	C	N	O	S	0	0	0
			2100	1333	358	401	8			
1	C	268	Total	C	N	O	S	0	0	0
			2100	1333	358	401	8			
1	D	266	Total	C	N	O	S	0	0	0
			2086	1325	356	397	8			
1	E	268	Total	C	N	O	S	0	0	0
			2100	1333	358	401	8			
1	F	268	Total	C	N	O	S	0	0	0
			2100	1333	358	401	8			
1	G	268	Total	C	N	O	S	0	0	0
			2100	1333	358	401	8			
1	H	268	Total	C	N	O	S	0	0	0
			2100	1333	358	401	8			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	211	ASP	ALA	engineered mutation	UNP P00514
B	211	ASP	ALA	engineered mutation	UNP P00514
C	211	ASP	ALA	engineered mutation	UNP P00514
D	211	ASP	ALA	engineered mutation	UNP P00514
E	211	ASP	ALA	engineered mutation	UNP P00514
F	211	ASP	ALA	engineered mutation	UNP P00514
G	211	ASP	ALA	engineered mutation	UNP P00514
H	211	ASP	ALA	engineered mutation	UNP P00514

- Molecule 2 is ADENOSINE-3',5'-CYCLIC-MONOPHOSPHATE (CCD ID: CMP) (formula: $C_{10}H_{12}N_5O_6P$) (labeled as "Ligand of Interest" by depositor).

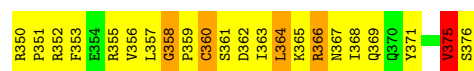


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	A	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	B	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	B	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	C	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	C	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	D	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	D	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	E	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	E	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	F	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	F	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	G	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	G	1	Total	C	N	O	P	0	0
			22	10	5	6	1		

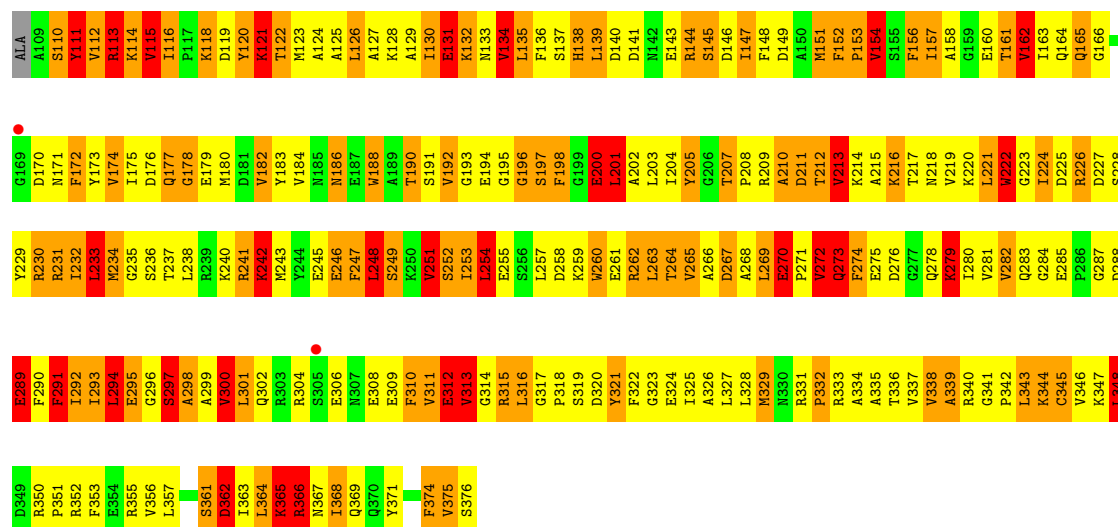
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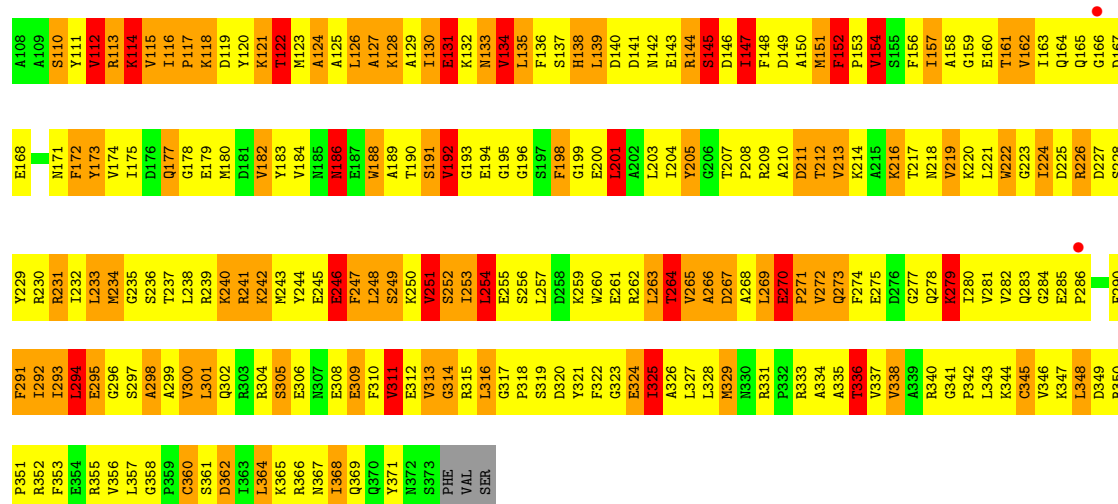
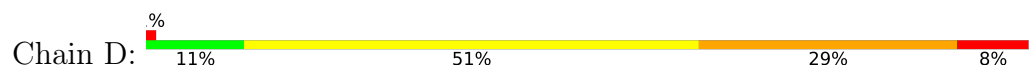
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	H	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	H	1	Total	C	N	O	P	0	0
			22	10	5	6	1		



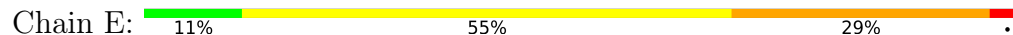
- Molecule 1: cAMP-dependent protein kinase type I-alpha regulatory subunit, N-terminally processed

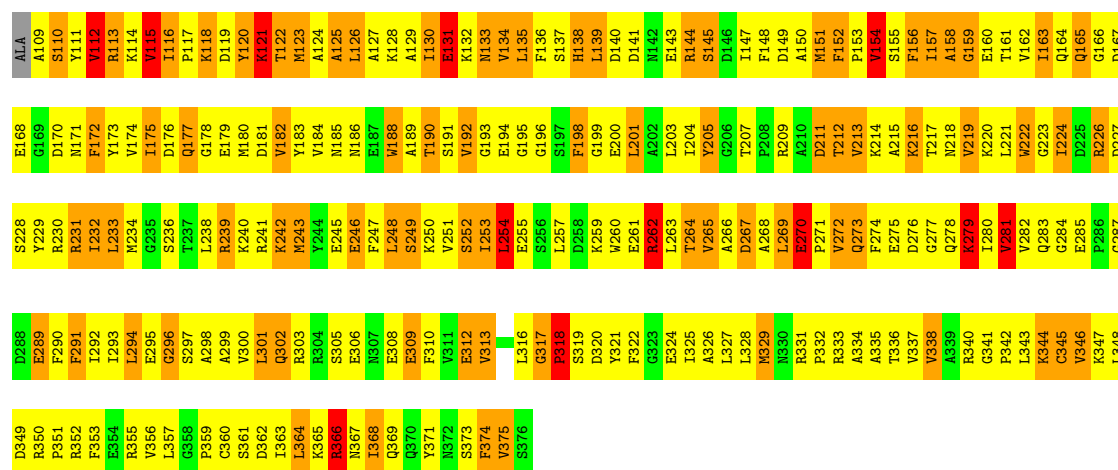


- Molecule 1: cAMP-dependent protein kinase type I-alpha regulatory subunit, N-terminally processed



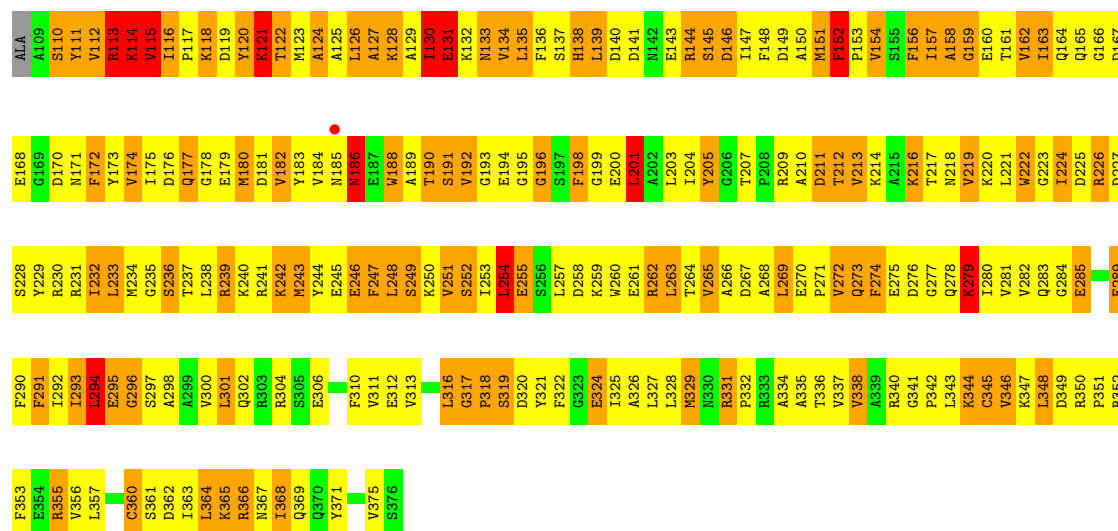
- Molecule 1: cAMP-dependent protein kinase type I-alpha regulatory subunit, N-terminally processed





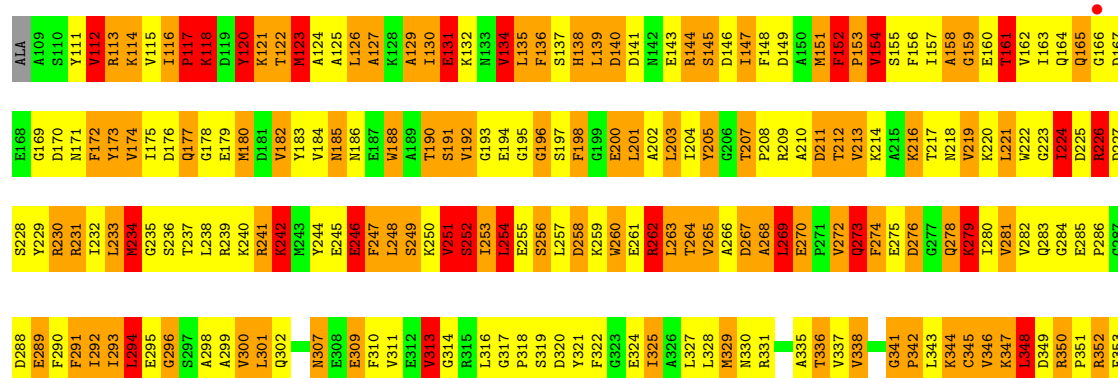
- Molecule 1: cAMP-dependent protein kinase type I-alpha regulatory subunit, N-terminally processed

Chain F: 13% 48% 34% .



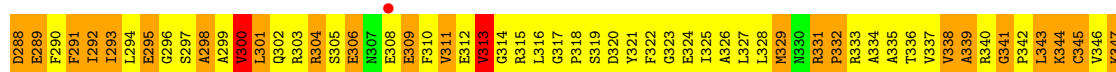
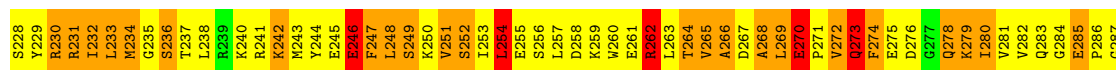
- Molecule 1: cAMP-dependent protein kinase type I-alpha regulatory subunit, N-terminally processed

Chain G: 13% 42% 35% 10%





- Molecule 1: cAMP-dependent protein kinase type I-alpha regulatory subunit, N-terminally processed



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	176.79Å 176.79Å 345.50Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.25 – 4.16 48.25 – 4.16	Depositor EDS
% Data completeness (in resolution range)	99.3 (48.25-4.16) 99.4 (48.25-4.16)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.69 (at 4.14Å)	Xtriage
Refinement program	PHENIX 1.7 _650	Depositor
R, R_{free}	0.221 , 0.270 0.228 , 0.276	Depositor DCC
R_{free} test set	1539 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	125.7	Xtriage
Anisotropy	0.440	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 158.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.418 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	17138	wwPDB-VP
Average B, all atoms (Å ²)	135.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% 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: CMP

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	1.72	33/2138 (1.5%)	1.97	93/2887 (3.2%)
1	B	1.68	30/2138 (1.4%)	1.93	75/2887 (2.6%)
1	C	1.57	21/2138 (1.0%)	2.01	90/2887 (3.1%)
1	D	1.57	20/2124 (0.9%)	1.91	71/2869 (2.5%)
1	E	1.40	19/2138 (0.9%)	1.77	54/2887 (1.9%)
1	F	1.51	23/2138 (1.1%)	1.88	77/2887 (2.7%)
1	G	1.90	58/2138 (2.7%)	2.31	126/2887 (4.4%)
1	H	1.56	22/2138 (1.0%)	1.91	74/2887 (2.6%)
All	All	1.62	226/17090 (1.3%)	1.97	660/23078 (2.9%)

The worst 5 of 226 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	112	VAL	CA-CB	19.77	1.81	1.54
1	A	112	VAL	CA-C	17.31	1.74	1.52
1	G	272	VAL	CA-CB	-13.99	1.45	1.55
1	F	192	VAL	CA-CB	-13.03	1.37	1.54
1	B	113	ARG	CA-C	12.57	1.68	1.52

The worst 5 of 660 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	121	LYS	CA-CB-CG	19.01	152.12	114.10
1	C	115	VAL	CB-CA-C	-16.25	89.87	110.99
1	G	115	VAL	N-CA-C	14.97	129.07	108.11
1	G	121	LYS	CB-CA-C	-14.94	86.00	110.79
1	C	279	LYS	CD-CE-NZ	-14.40	65.81	111.90

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	2100	0	2069	752	0
1	B	2100	0	2071	729	0
1	C	2100	0	2069	808	0
1	D	2086	0	2060	726	0
1	E	2100	0	2071	703	0
1	F	2100	0	2071	708	0
1	G	2100	0	2071	619	0
1	H	2100	0	2071	760	0
2	A	44	0	22	40	0
2	B	44	0	22	55	0
2	C	44	0	22	36	0
2	D	44	0	22	37	0
2	E	44	0	22	35	0
2	F	44	0	22	29	0
2	G	44	0	22	30	0
2	H	44	0	22	38	0
All	All	17138	0	16729	5572	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 165.

The worst 5 of 5572 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:112:VAL:CA	1:A:112:VAL:C	1.74	1.54
1:A:112:VAL:CA	1:A:112:VAL:CB	1.81	1.53
2:C:402:CMP:H2	2:C:402:CMP:C2	0.97	1.50
2:A:401:CMP:H2	2:A:401:CMP:C2	0.97	1.49
2:A:402:CMP:H2	2:A:402:CMP:C2	0.97	1.49

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	266/269 (99%)	225 (85%)	32 (12%)	9 (3%)	3	21
1	B	266/269 (99%)	218 (82%)	40 (15%)	8 (3%)	3	23
1	C	266/269 (99%)	221 (83%)	34 (13%)	11 (4%)	2	18
1	D	264/269 (98%)	226 (86%)	31 (12%)	7 (3%)	4	26
1	E	266/269 (99%)	219 (82%)	41 (15%)	6 (2%)	5	29
1	F	266/269 (99%)	224 (84%)	37 (14%)	5 (2%)	6	32
1	G	266/269 (99%)	224 (84%)	31 (12%)	11 (4%)	2	18
1	H	266/269 (99%)	225 (85%)	35 (13%)	6 (2%)	5	29
All	All	2126/2152 (99%)	1782 (84%)	281 (13%)	63 (3%)	3	23

5 of 63 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	112	VAL
1	A	252	SER
1	A	253	ILE
1	B	112	VAL
1	B	252	SER

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	222/229 (97%)	164 (74%)	58 (26%)	0	4

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	222/229 (97%)	159 (72%)	63 (28%)	0	2
1	C	222/229 (97%)	152 (68%)	70 (32%)	0	2
1	D	220/229 (96%)	154 (70%)	66 (30%)	0	2
1	E	222/229 (97%)	165 (74%)	57 (26%)	0	4
1	F	222/229 (97%)	161 (72%)	61 (28%)	0	3
1	G	222/229 (97%)	169 (76%)	53 (24%)	1	5
1	H	222/229 (97%)	159 (72%)	63 (28%)	0	2
All	All	1774/1832 (97%)	1283 (72%)	491 (28%)	0	3

5 of 491 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	249	SER
1	H	151	MET
1	E	233	LEU
1	H	135	LEU
1	H	273	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 24 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	372	ASN
1	F	372	ASN
1	F	302	GLN
1	G	273	GLN
1	B	372	ASN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

16 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
2	CMP	D	402	-	25,25,25	2.04	8 (32%)	37,39,39	2.34	12 (32%)
2	CMP	C	401	-	25,25,25	1.95	5 (20%)	37,39,39	2.87	13 (35%)
2	CMP	D	401	-	25,25,25	2.10	5 (20%)	37,39,39	2.55	12 (32%)
2	CMP	E	401	-	25,25,25	2.13	9 (36%)	37,39,39	2.53	13 (35%)
2	CMP	F	402	-	25,25,25	2.18	5 (20%)	37,39,39	2.31	10 (27%)
2	CMP	H	402	-	25,25,25	2.03	8 (32%)	37,39,39	2.21	11 (29%)
2	CMP	A	402	-	25,25,25	2.21	8 (32%)	37,39,39	3.11	13 (35%)
2	CMP	C	402	-	25,25,25	2.12	8 (32%)	37,39,39	2.66	10 (27%)
2	CMP	G	401	-	25,25,25	2.20	6 (24%)	37,39,39	2.30	11 (29%)
2	CMP	B	402	-	25,25,25	2.33	6 (24%)	37,39,39	2.55	12 (32%)
2	CMP	E	402	-	25,25,25	2.18	5 (20%)	37,39,39	2.15	10 (27%)
2	CMP	G	402	-	25,25,25	1.94	8 (32%)	37,39,39	2.80	12 (32%)
2	CMP	A	401	-	25,25,25	1.88	5 (20%)	37,39,39	2.65	13 (35%)
2	CMP	H	401	-	25,25,25	2.38	6 (24%)	37,39,39	3.27	14 (37%)
2	CMP	B	401	-	25,25,25	1.96	6 (24%)	37,39,39	2.89	14 (37%)
2	CMP	F	401	-	25,25,25	2.11	9 (36%)	37,39,39	2.27	13 (35%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	CMP	D	402	-	-	0/4/31/31	0/4/4/4
2	CMP	C	401	-	-	0/4/31/31	0/4/4/4
2	CMP	D	401	-	-	0/4/31/31	0/4/4/4
2	CMP	E	401	-	-	0/4/31/31	0/4/4/4
2	CMP	F	402	-	-	3/4/31/31	0/4/4/4
2	CMP	H	402	-	-	3/4/31/31	0/4/4/4
2	CMP	A	402	-	-	1/4/31/31	0/4/4/4
2	CMP	C	402	-	-	0/4/31/31	0/4/4/4
2	CMP	G	401	-	-	1/4/31/31	0/4/4/4
2	CMP	B	402	-	-	0/4/31/31	0/4/4/4
2	CMP	E	402	-	-	2/4/31/31	0/4/4/4
2	CMP	G	402	-	-	1/4/31/31	0/4/4/4
2	CMP	A	401	-	-	0/4/31/31	0/4/4/4
2	CMP	H	401	-	-	1/4/31/31	0/4/4/4
2	CMP	B	401	-	-	0/4/31/31	0/4/4/4
2	CMP	F	401	-	-	0/4/31/31	0/4/4/4

The worst 5 of 107 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	402	CMP	P-O3'	6.88	1.69	1.57
2	H	401	CMP	P-O3'	6.86	1.69	1.57
2	G	401	CMP	C5-C4	6.12	1.50	1.39
2	D	401	CMP	P-O3'	5.93	1.67	1.57
2	E	402	CMP	C5-C4	5.67	1.49	1.39

The worst 5 of 193 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	401	CMP	C5-C4-N3	-11.41	111.01	126.72
2	A	402	CMP	C5-C4-N3	-9.75	113.30	126.72
2	C	401	CMP	C5-C4-N3	-9.67	113.41	126.72
2	D	401	CMP	C5-C4-N3	-8.50	115.02	126.72
2	C	402	CMP	C5-C4-N3	-8.28	115.32	126.72

There are no chirality outliers.

5 of 12 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	402	CMP	C2'-C1'-N9-C4
2	F	402	CMP	C2'-C1'-N9-C4
2	E	402	CMP	C2'-C1'-N9-C8

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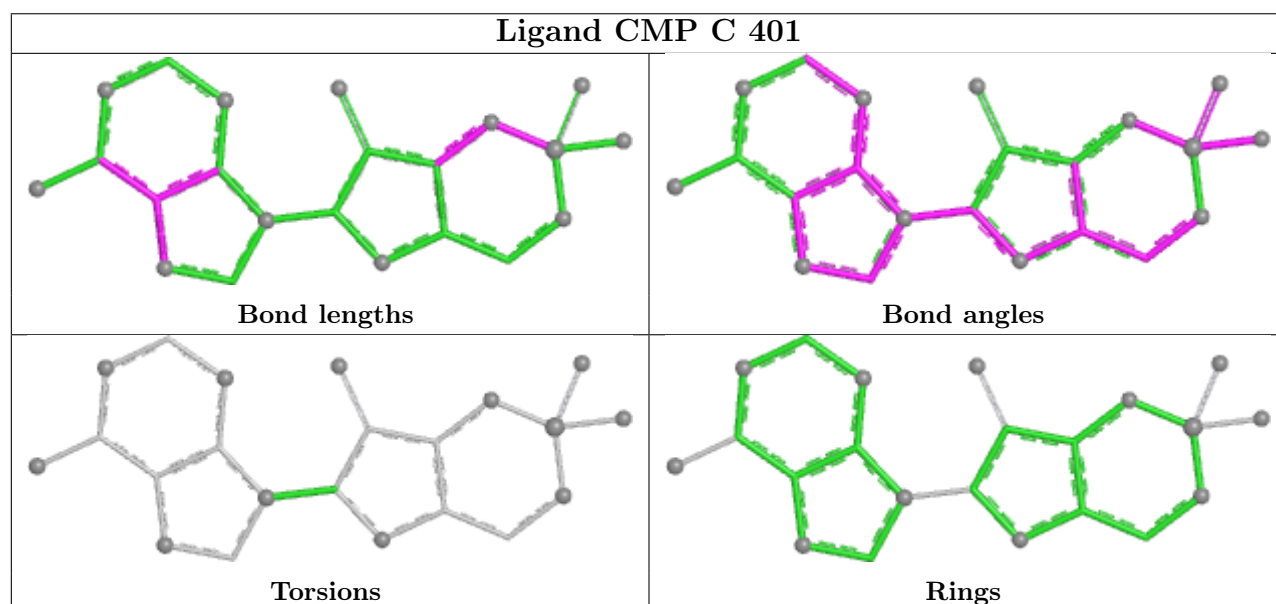
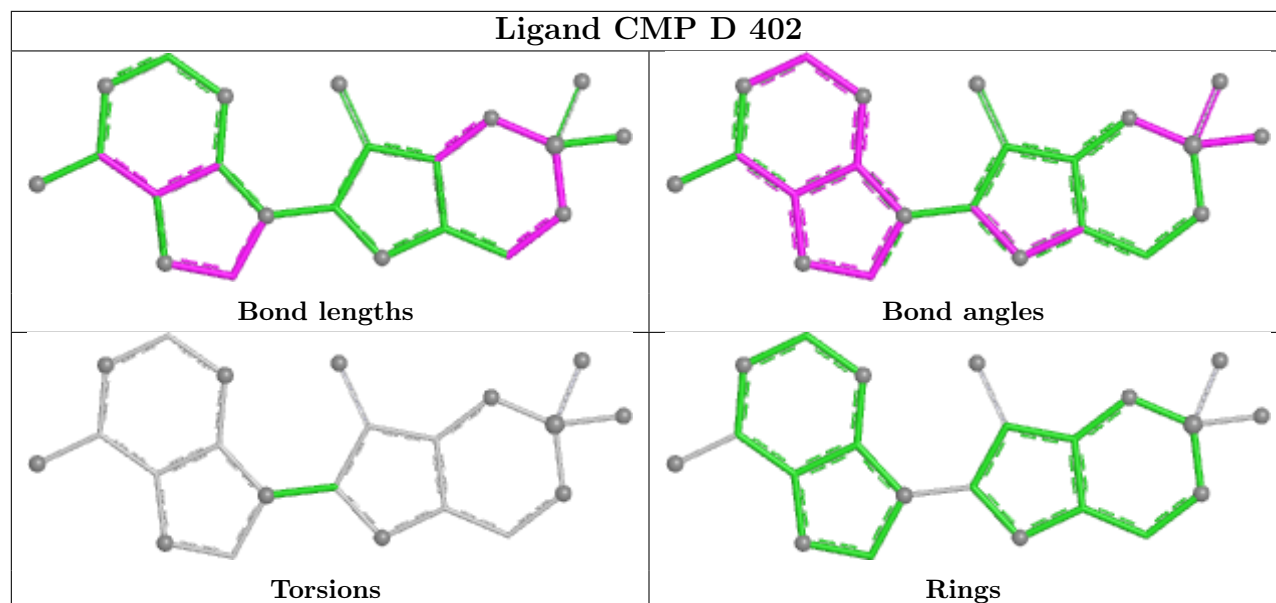
Mol	Chain	Res	Type	Atoms
2	F	402	CMP	C2'-C1'-N9-C8
2	H	402	CMP	C2'-C1'-N9-C8

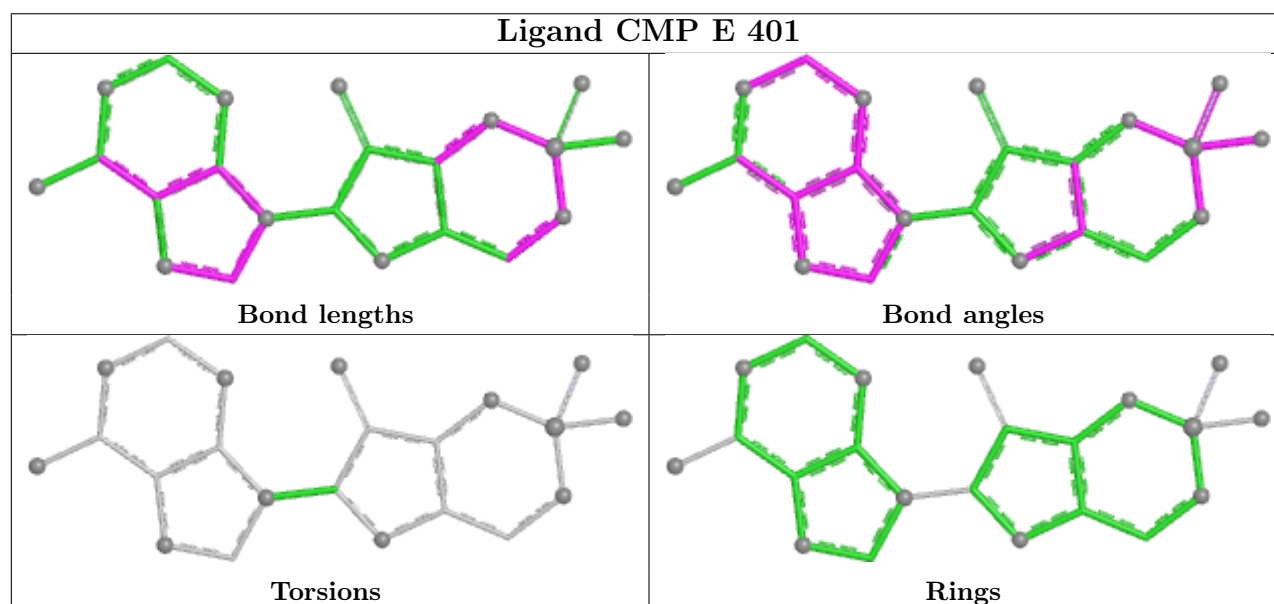
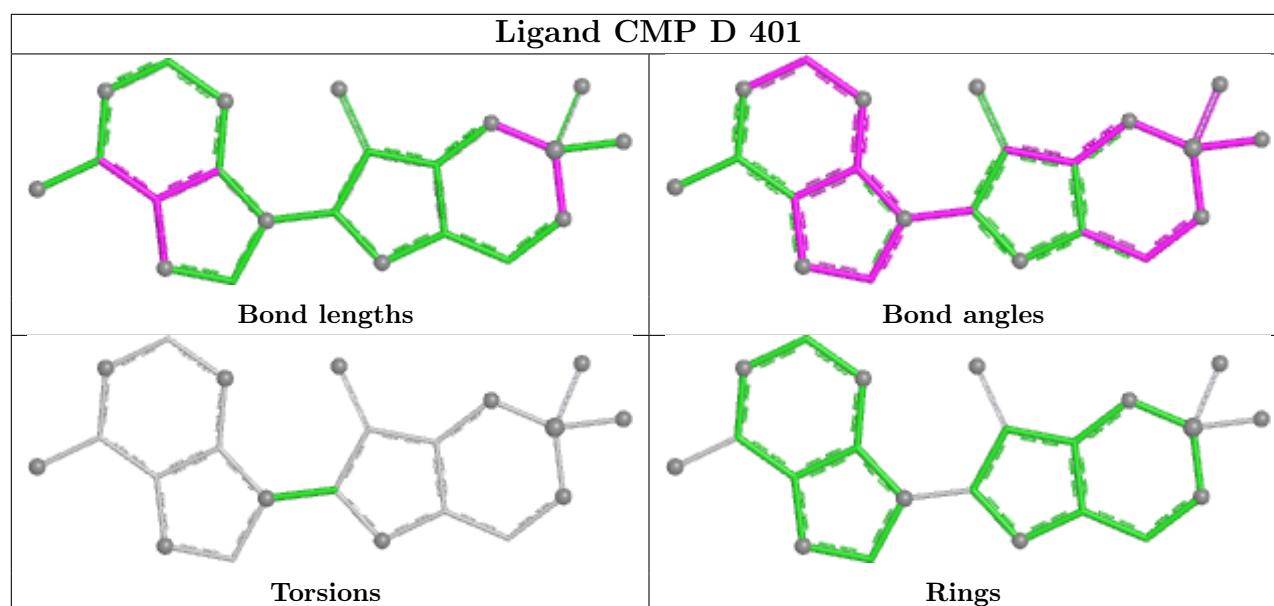
There are no ring outliers.

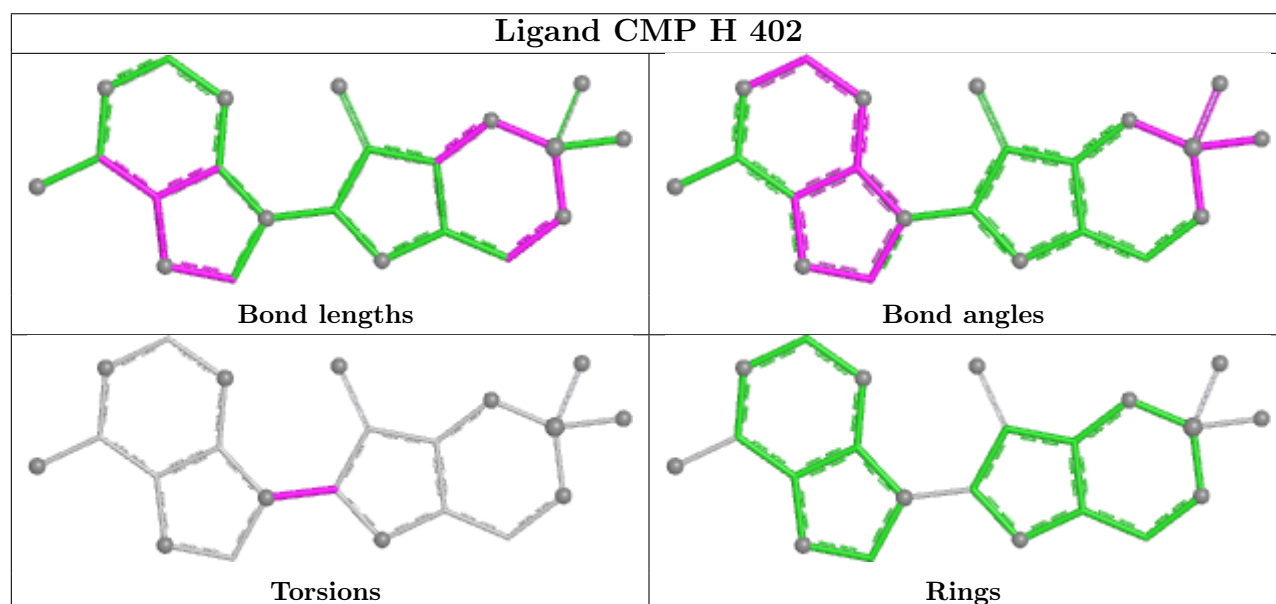
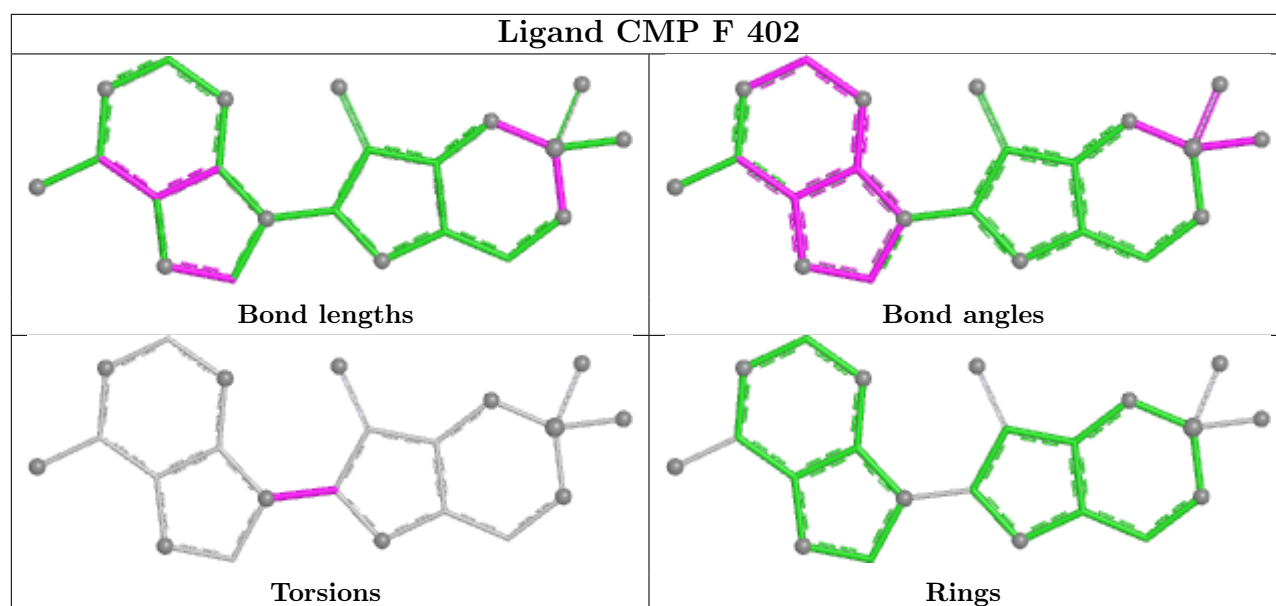
16 monomers are involved in 300 short contacts:

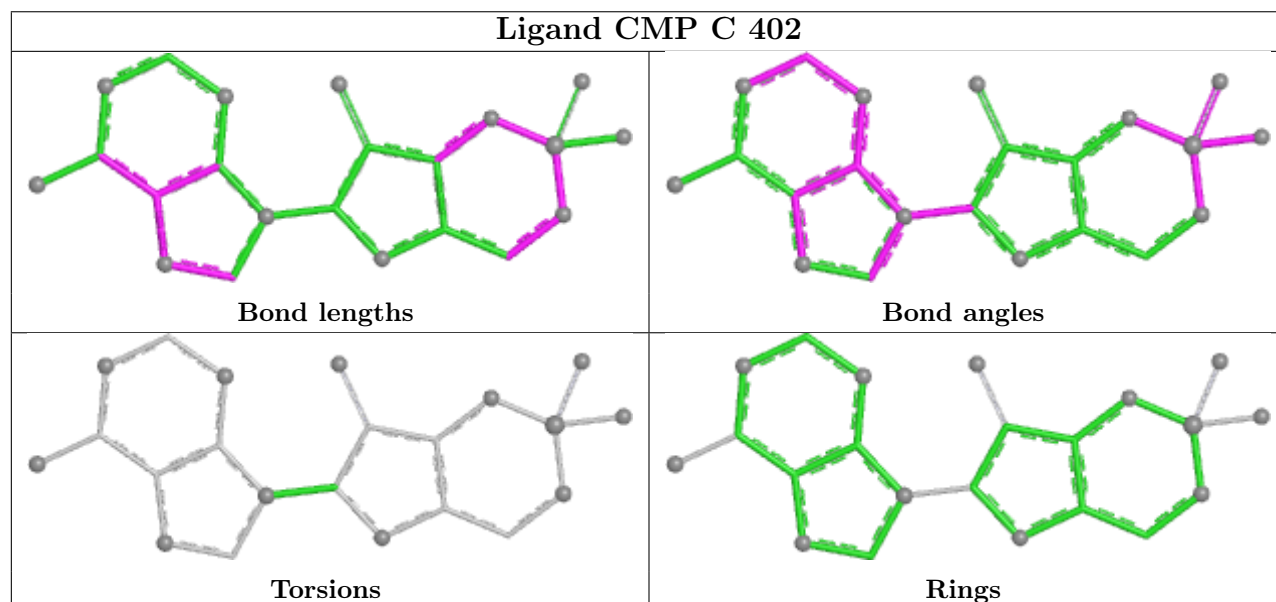
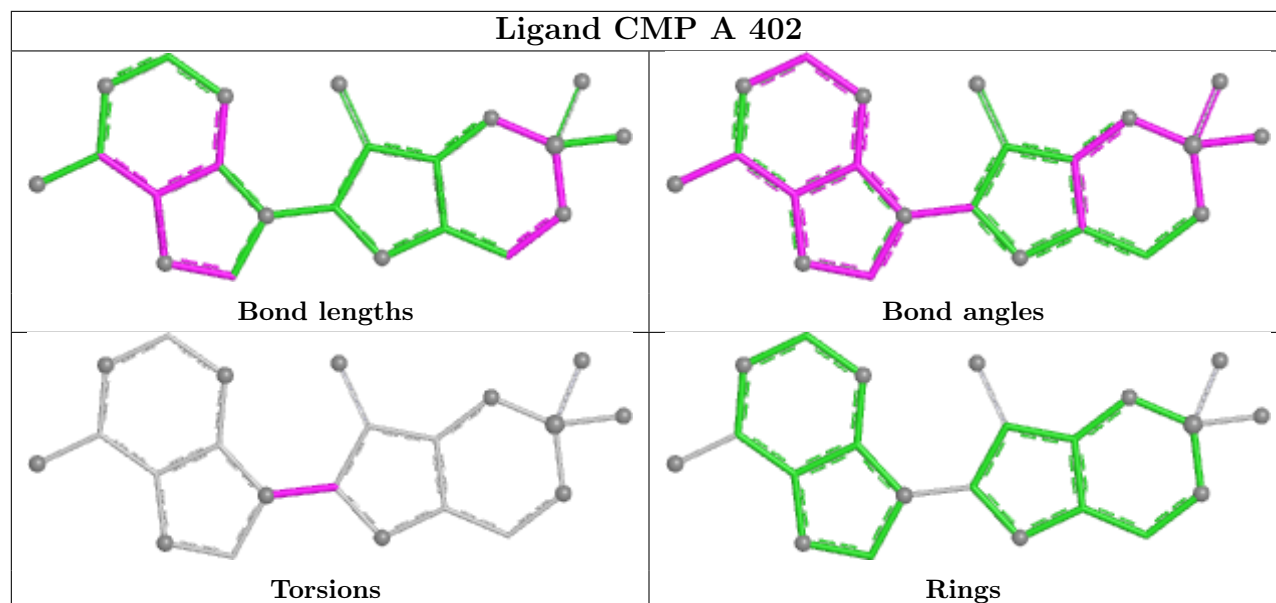
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	402	CMP	21	0
2	C	401	CMP	17	0
2	D	401	CMP	16	0
2	E	401	CMP	24	0
2	F	402	CMP	7	0
2	H	402	CMP	16	0
2	A	402	CMP	15	0
2	C	402	CMP	19	0
2	G	401	CMP	19	0
2	B	402	CMP	21	0
2	E	402	CMP	11	0
2	G	402	CMP	11	0
2	A	401	CMP	25	0
2	H	401	CMP	22	0
2	B	401	CMP	34	0
2	F	401	CMP	22	0

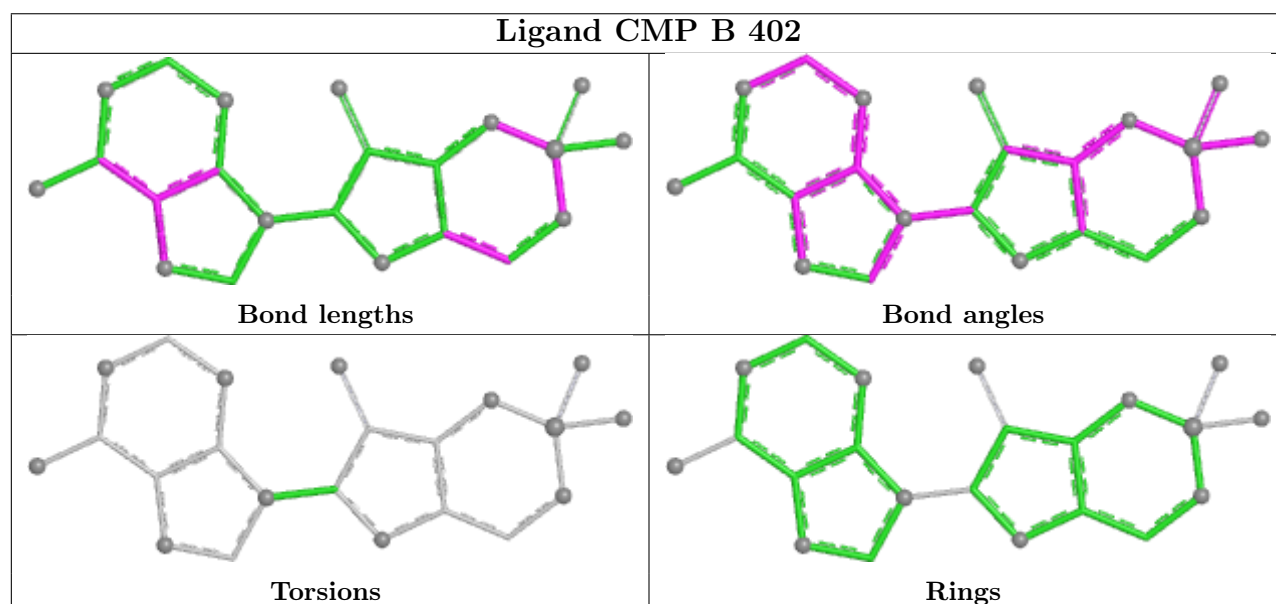
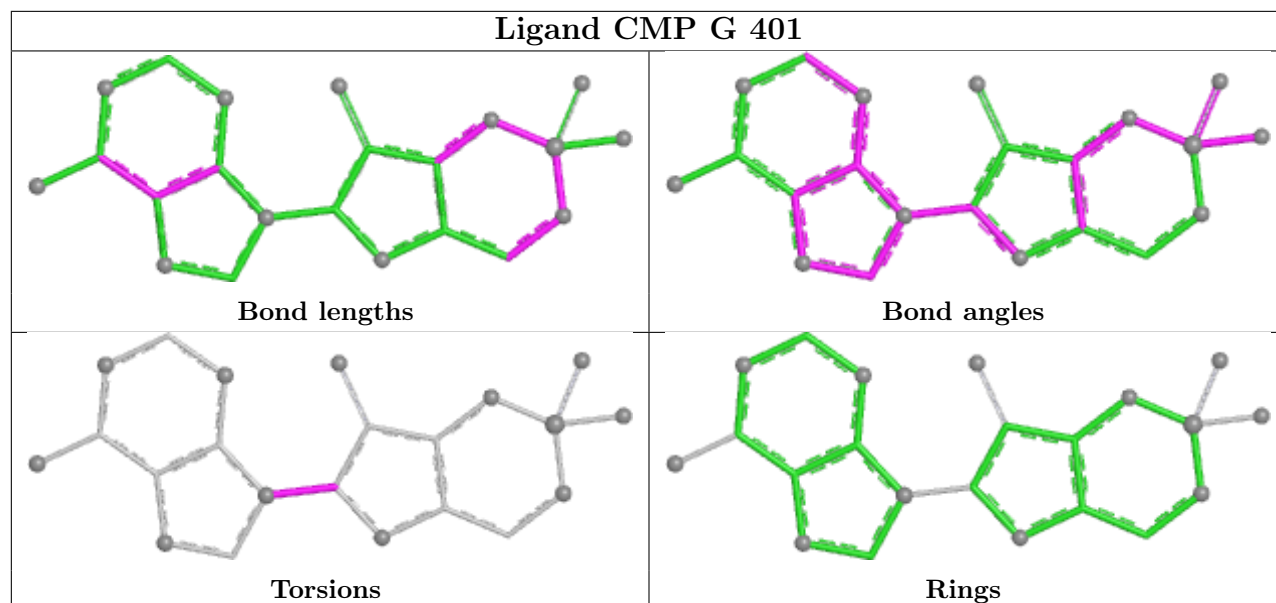
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

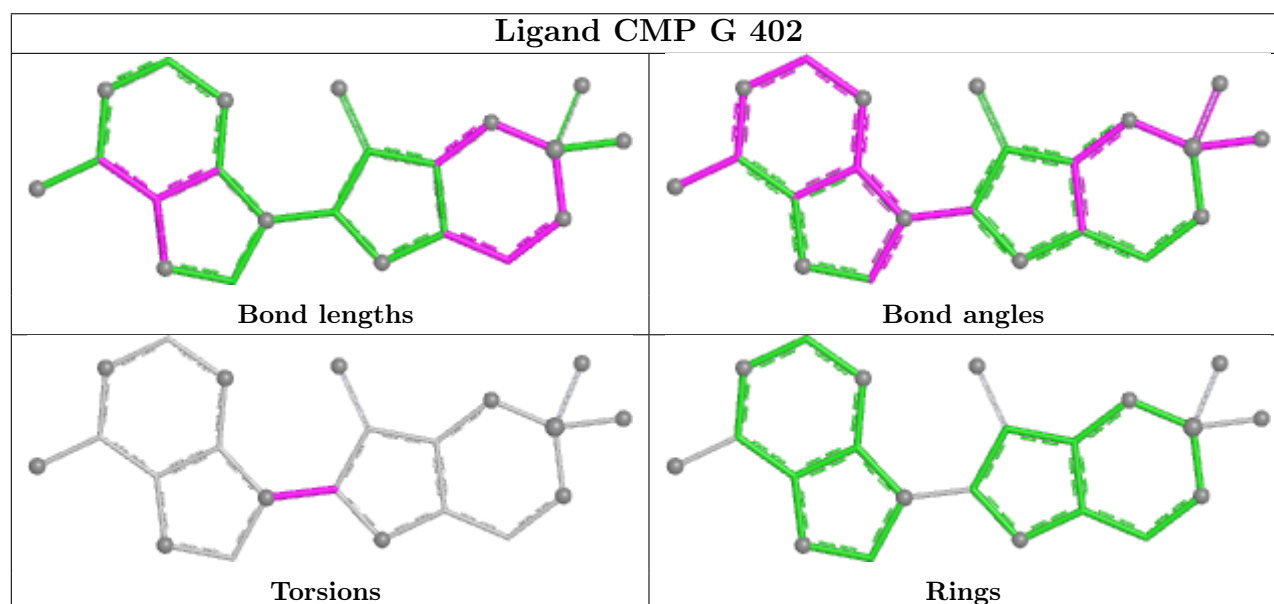
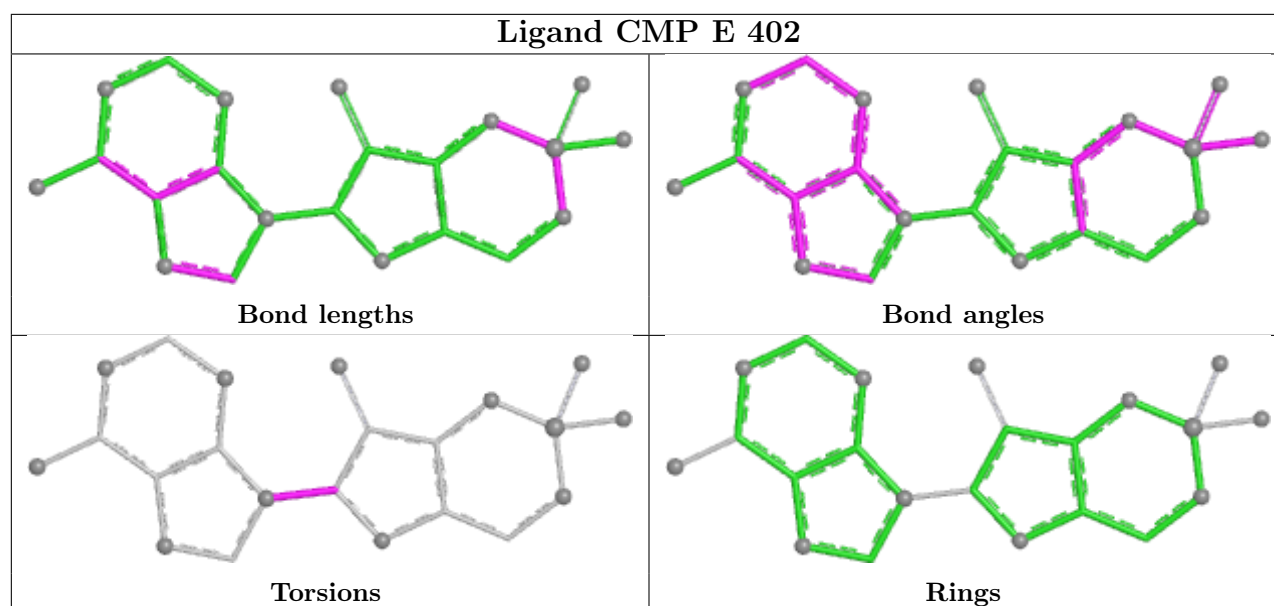


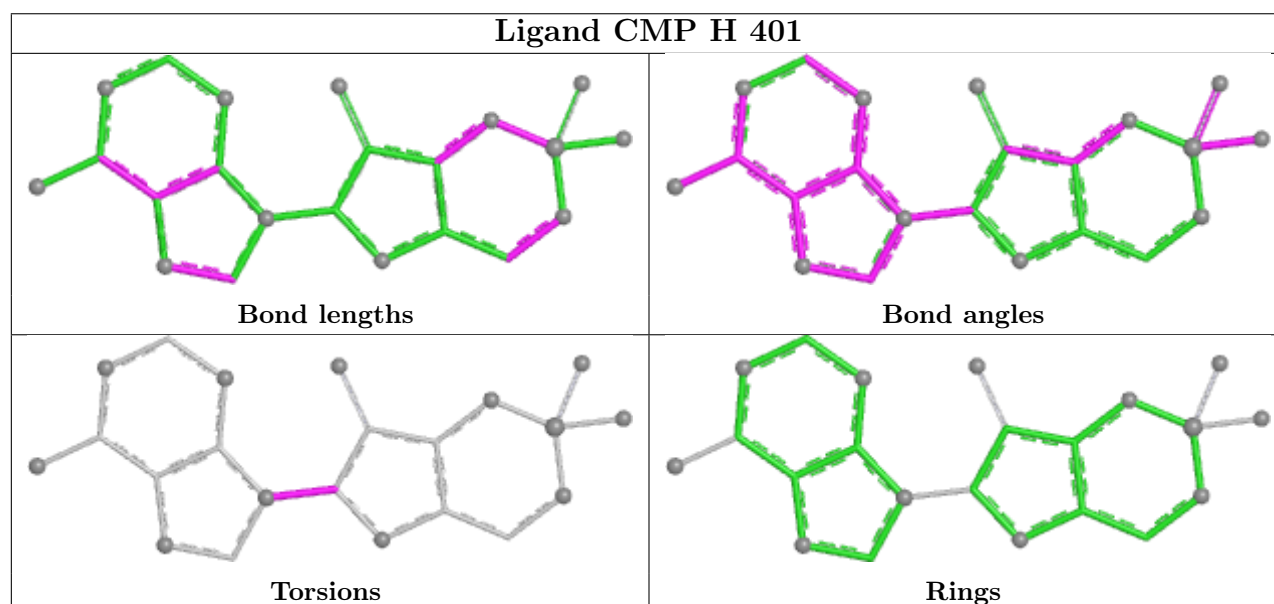
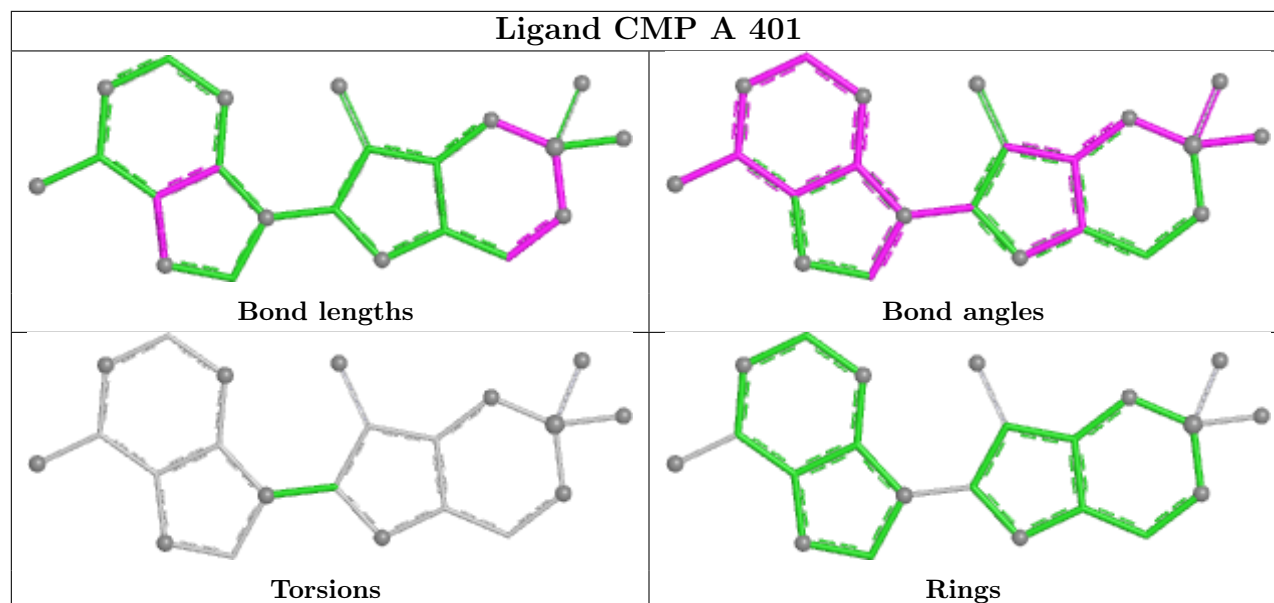


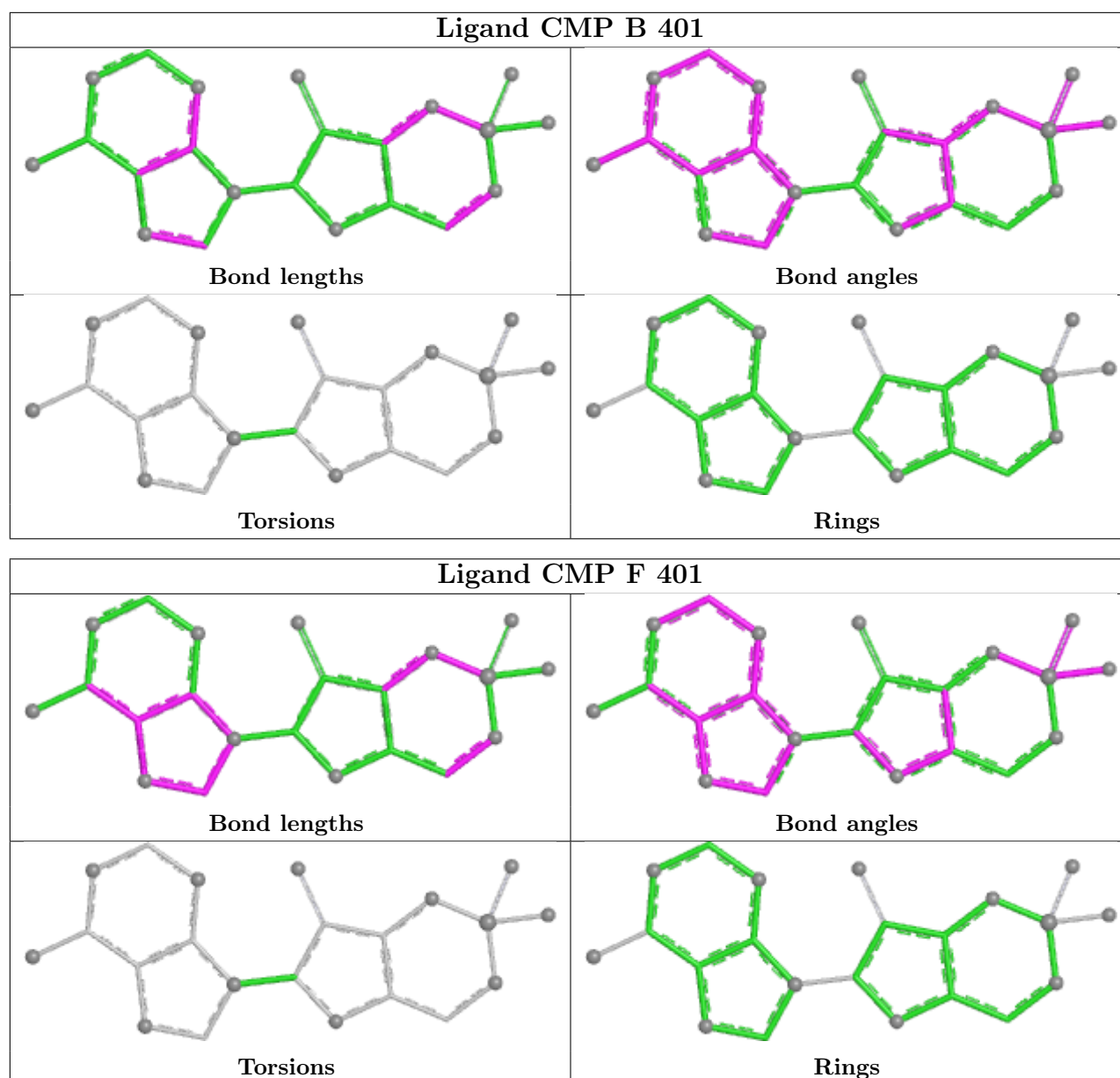












5.7 Other polymers [i](#)

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, 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	268/269 (99%)	-0.69	5 (1%) 66 51	70, 133, 202, 260	0
1	B	268/269 (99%)	-0.74	4 (1%) 72 56	70, 133, 202, 266	0
1	C	268/269 (99%)	-0.72	2 (0%) 84 70	71, 131, 201, 286	0
1	D	266/269 (98%)	-0.73	2 (0%) 82 68	75, 133, 194, 246	0
1	E	268/269 (99%)	-0.75	0 100 100	75, 139, 215, 259	0
1	F	268/269 (99%)	-0.69	1 (0%) 88 76	75, 138, 212, 261	0
1	G	268/269 (99%)	-0.70	1 (0%) 88 76	77, 135, 201, 268	0
1	H	268/269 (99%)	-0.70	3 (1%) 78 62	71, 132, 199, 274	0
All	All	2142/2152 (99%)	-0.72	18 (0%) 82 68	70, 134, 205, 286	0

The worst 5 of 18 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	166	GLY	4.2
1	F	185	ASN	3.6
1	H	169	GLY	3.2
1	D	166	GLY	2.7
1	A	168	GLU	2.6

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

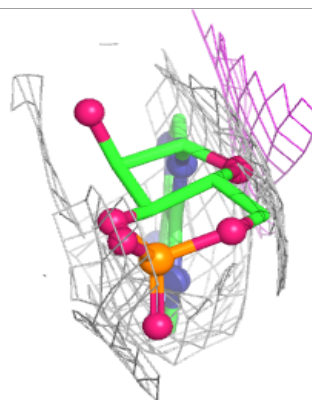
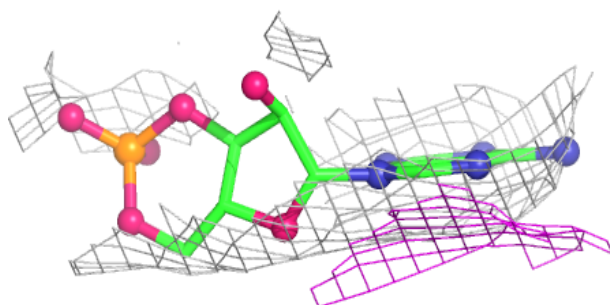
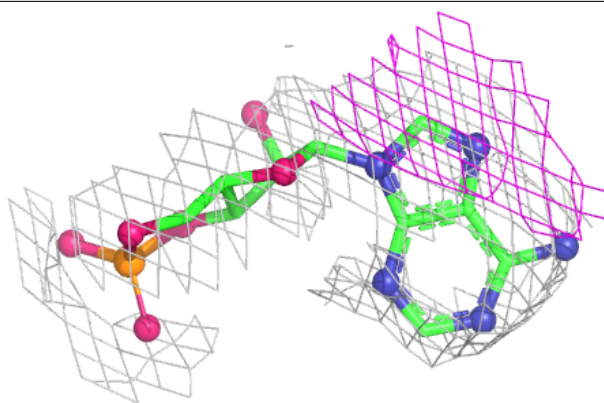
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	CMP	A	401	22/22	0.99	0.05	77,102,123,128	0
2	CMP	A	402	22/22	0.99	0.06	79,151,186,198	0
2	CMP	B	401	22/22	0.99	0.04	70,102,127,132	0
2	CMP	C	401	22/22	0.99	0.05	77,107,140,155	0
2	CMP	C	402	22/22	0.99	0.03	79,130,177,213	0
2	CMP	D	401	22/22	0.99	0.05	84,111,128,131	0
2	CMP	E	401	22/22	0.99	0.04	73,108,130,140	0
2	CMP	E	402	22/22	0.99	0.04	118,201,228,242	0
2	CMP	F	401	22/22	0.99	0.04	69,96,132,143	0
2	CMP	F	402	22/22	0.99	0.04	137,185,218,227	0
2	CMP	G	401	22/22	0.99	0.04	79,112,144,145	0
2	CMP	H	401	22/22	0.99	0.05	74,105,125,133	0
2	CMP	H	402	22/22	0.99	0.04	81,139,165,200	0
2	CMP	G	402	22/22	1.00	0.04	101,148,200,208	0
2	CMP	D	402	22/22	1.00	0.04	112,143,193,210	0
2	CMP	B	402	22/22	1.00	0.04	86,150,175,194	0

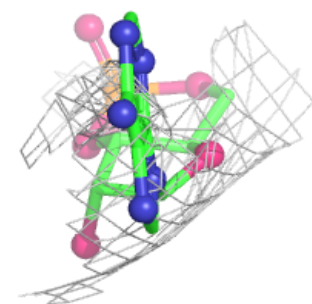
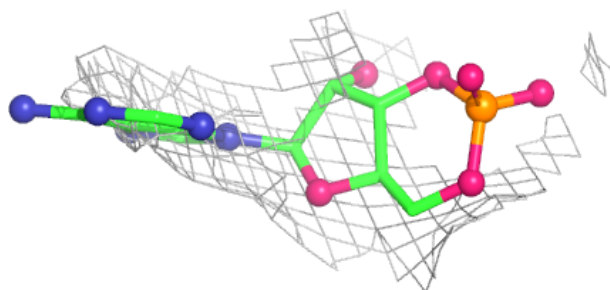
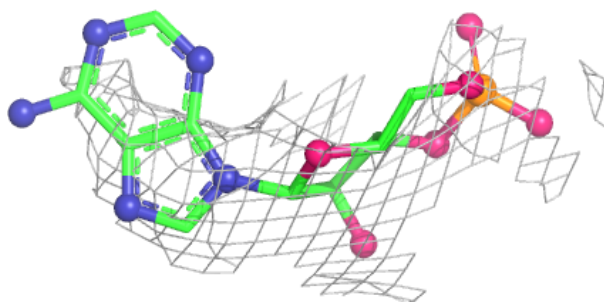
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around CMP A 401:

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

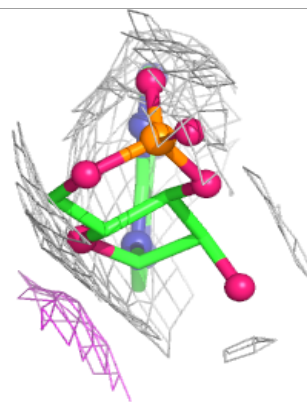
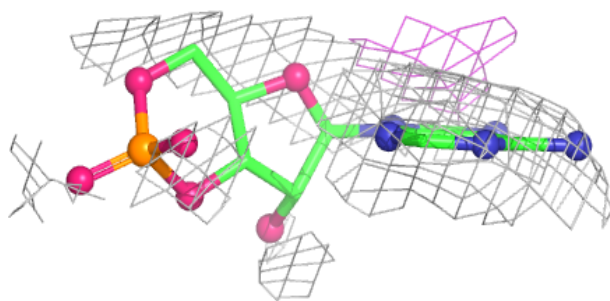
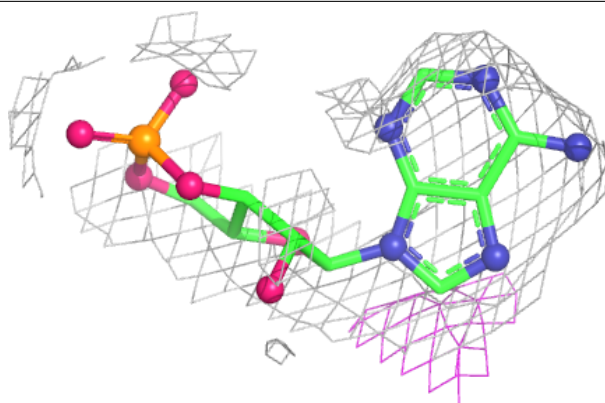
**Electron density around CMP A 402:**

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and green (positive)

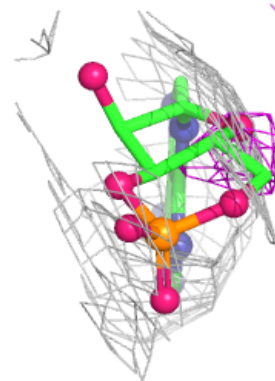
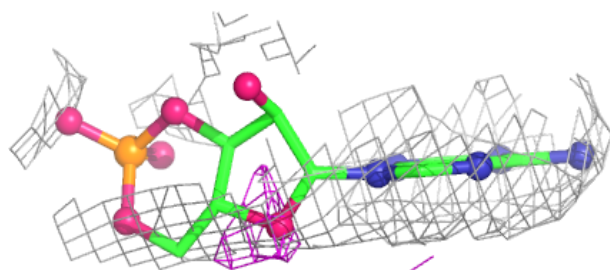
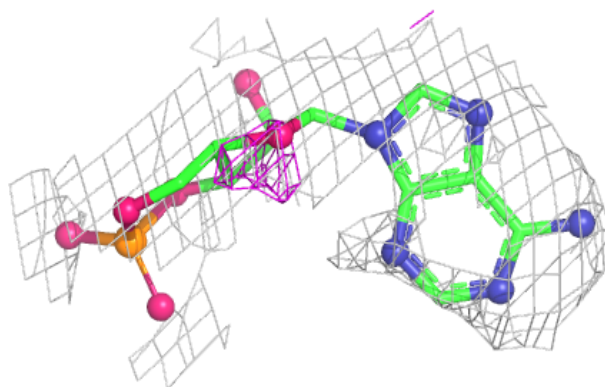


Electron density around CMP B 401:

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

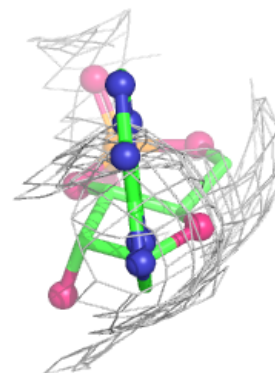
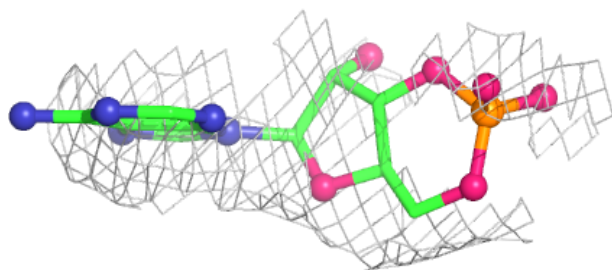
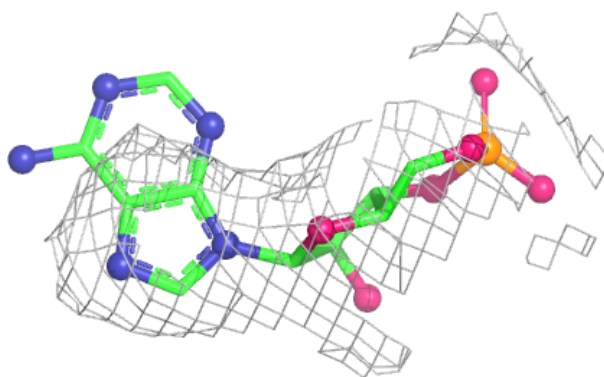
**Electron density around CMP C 401:**

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and green (positive)

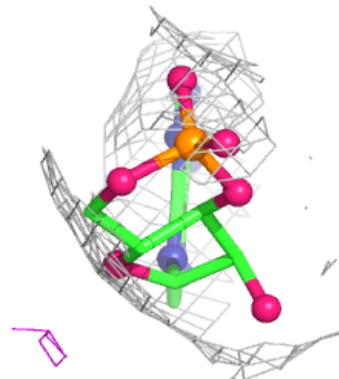
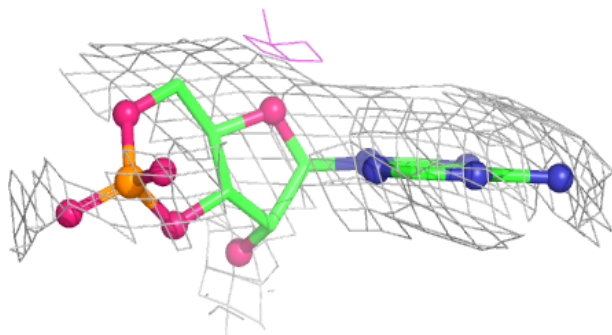
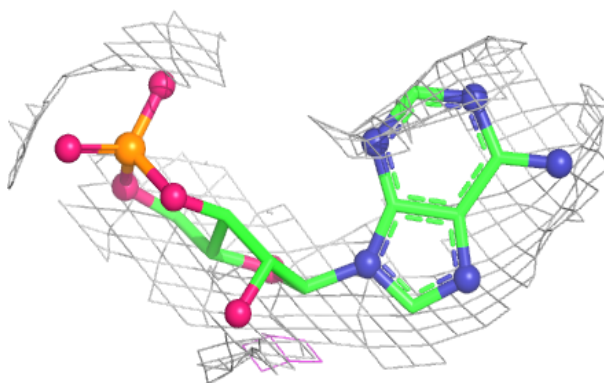


Electron density around CMP C 402:

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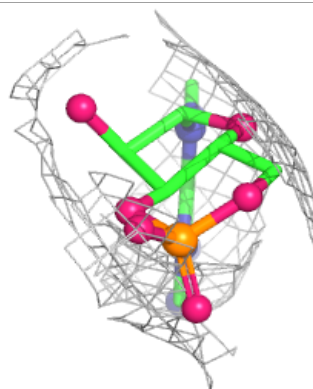
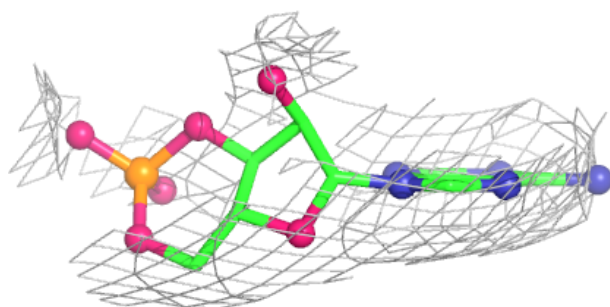
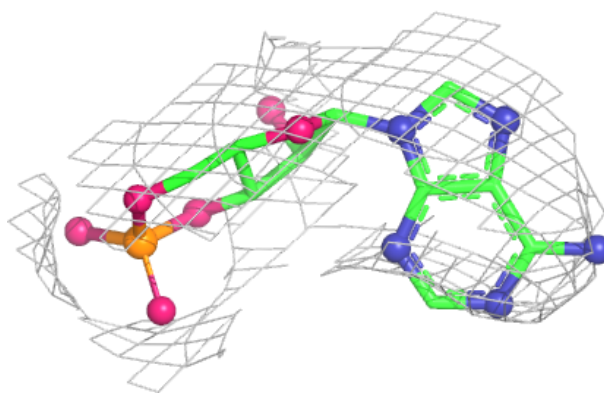
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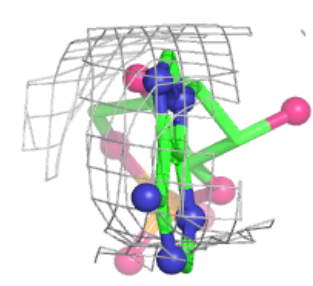
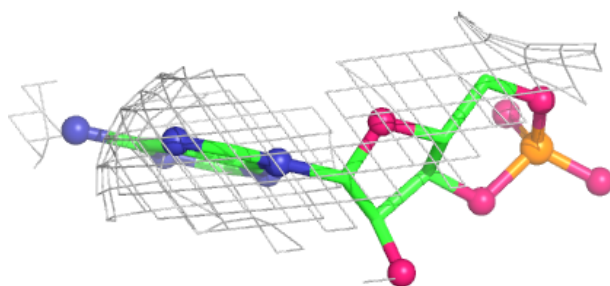
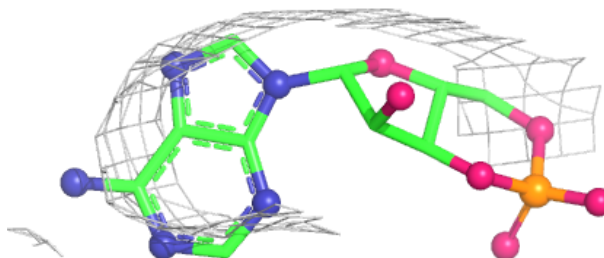


Electron density around CMP E 401:

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and green (positive)

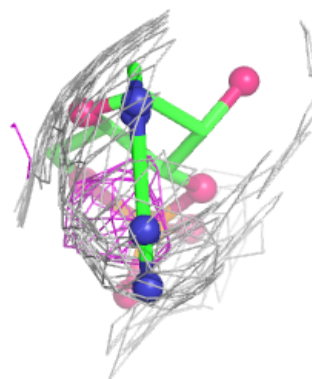
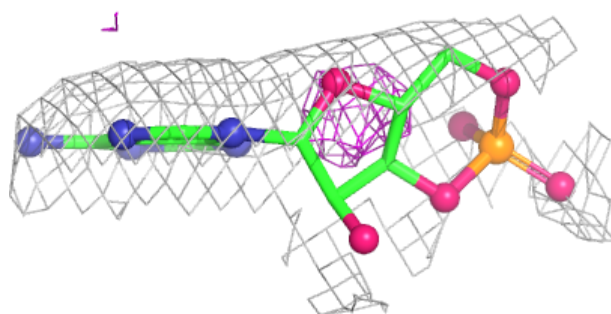
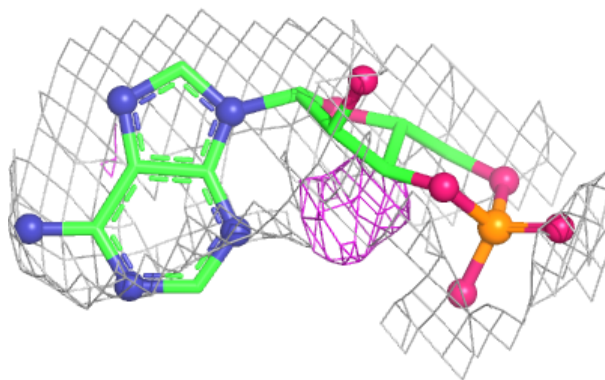
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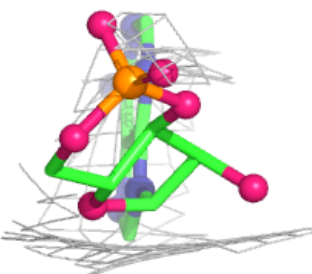
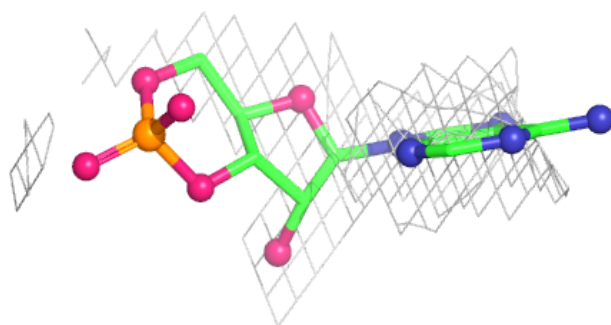
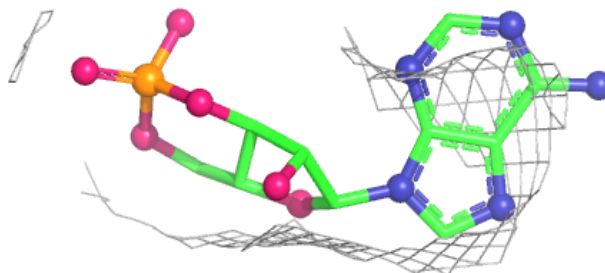


Electron density around CMP F 401:

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and green (positive)

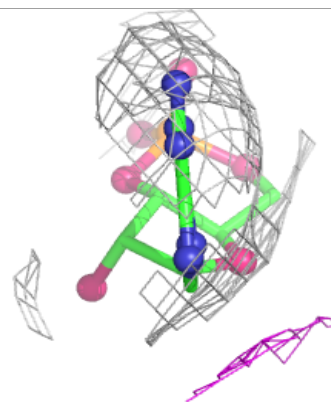
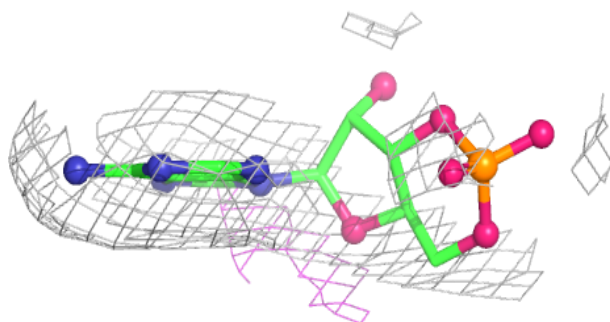
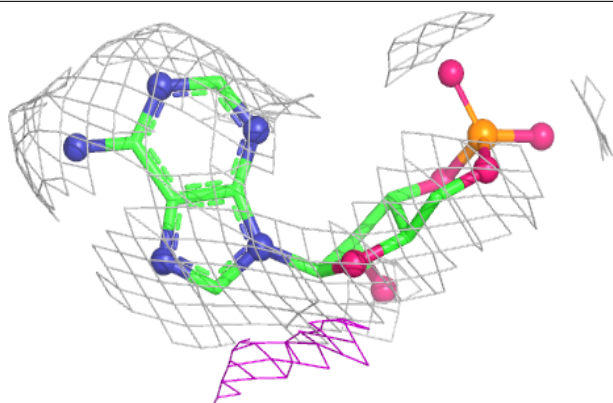
**Electron density around CMP F 402:**

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and green (positive)

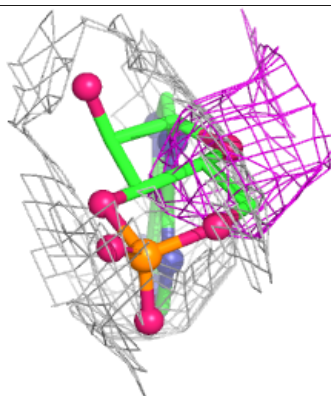
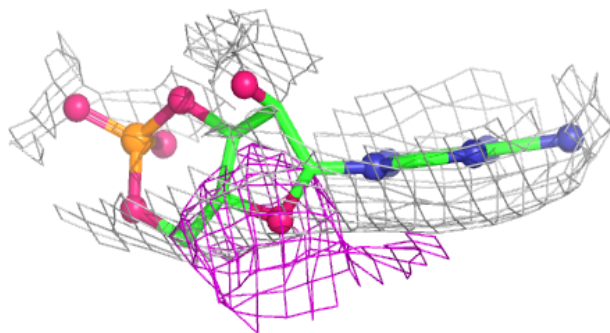
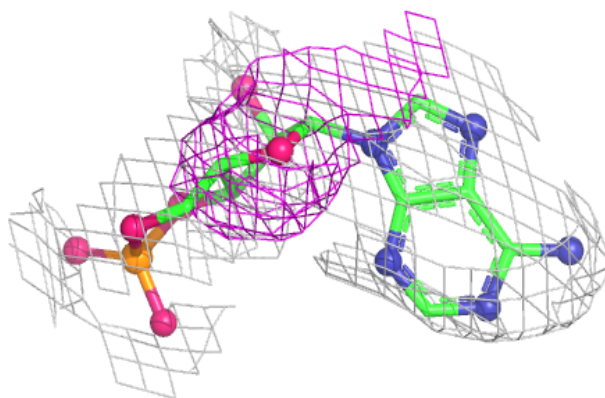


Electron density around CMP G 401:

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and green (positive)

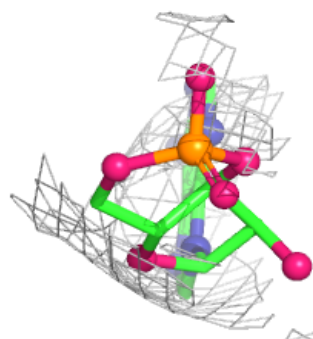
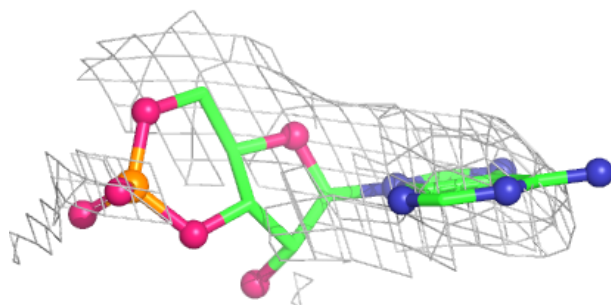
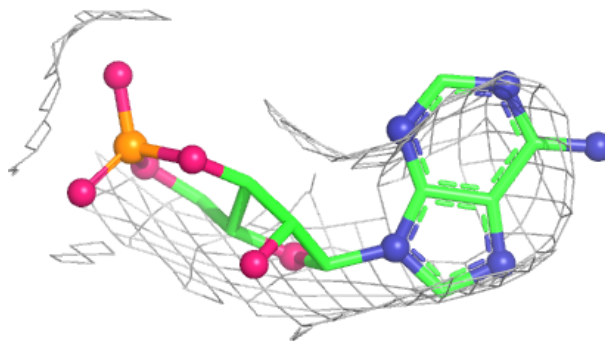
**Electron density around CMP H 401:**

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

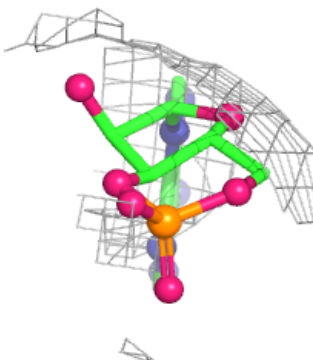
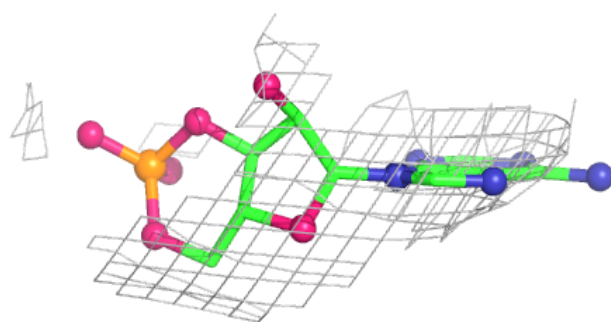
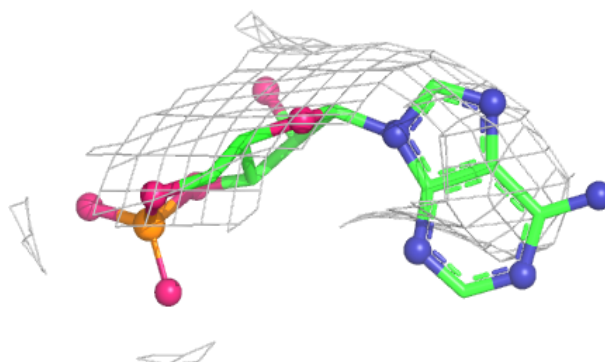


Electron density around CMP H 402:

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

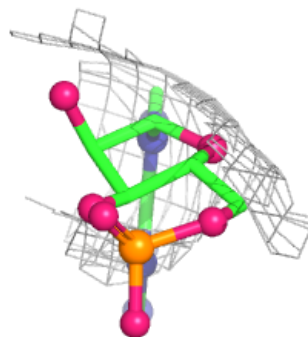
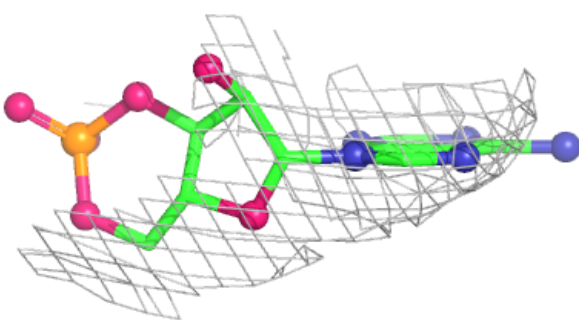
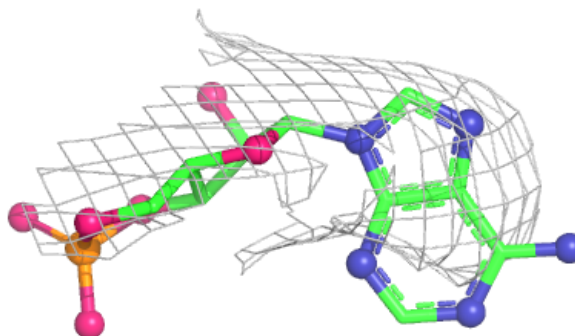
**Electron density around CMP G 402:**

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

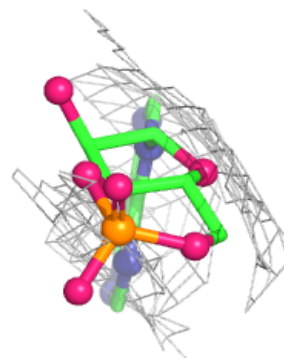
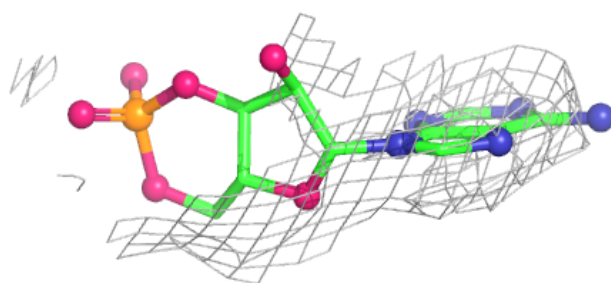
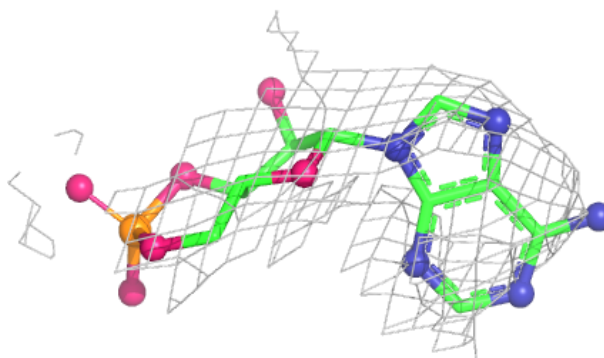


Electron density around CMP D 402:

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

**Electron density around CMP B 402:**

$2mF_o-DF_c$ (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.