



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 03:47 AM UTC

PDB ID : 5IQD / pdb_00005iqd
Title : Aminoglycoside Phosphotransferase (2'')-Ia (CTD of AAC(6')-Ie/APH(2'')-Ia)
in complex with GMPPNP, Magnesium, and Ribostamycin
Authors : Caldwell, S.J.; Berghuis, A.M.
Deposited on : 2016-03-10
Resolution : 2.20 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

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

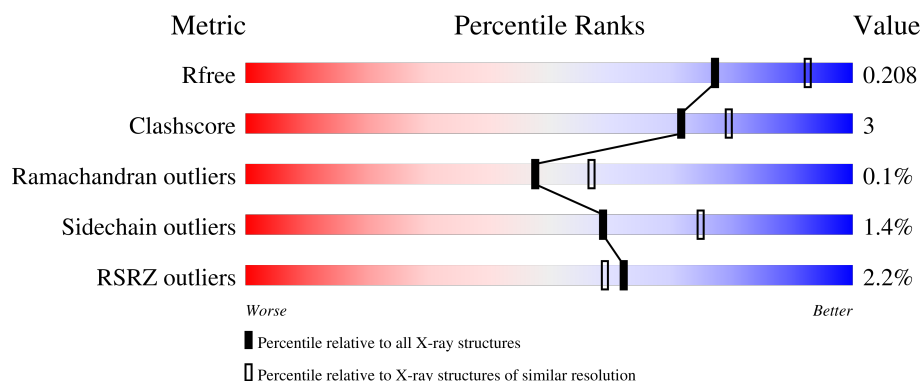
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	305	3% 89% 7% ..
1	B	305	2% 88% 9% ..
1	C	305	% 88% 7% ..
1	D	305	3% 89% 8% ..

2 Entry composition [i](#)

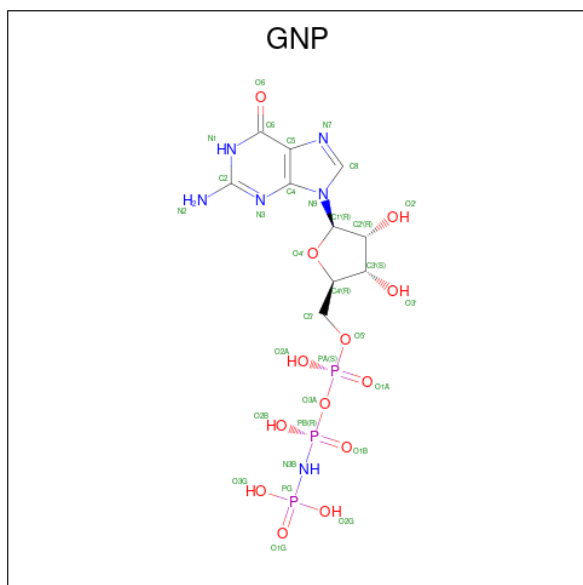
There are 6 unique types of molecules in this entry. The entry contains 11182 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Bifunctional AAC/APH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	S	0	4	0
			2447	1559	381	496	11			
1	B	299	Total	C	N	O	S	0	3	0
			2478	1581	382	504	11			
1	C	292	Total	C	N	O	S	0	3	0
			2425	1548	377	489	11			
1	D	297	Total	C	N	O	S	0	3	0
			2453	1566	383	493	11			

- Molecule 2 is PHOSPHOAMINOPHOSPHONIC ACID-GUANYLATE ESTER (CCD ID: GNP) (formula: $C_{10}H_{17}N_6O_{13}P_3$).



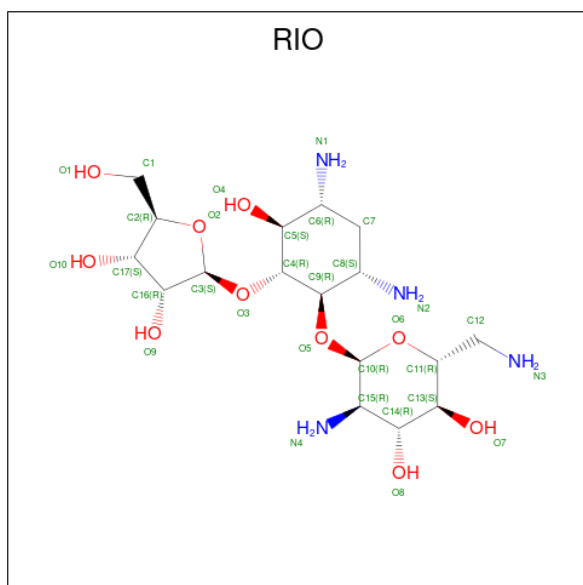
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
2	B	1	Total	C	N	O	P	0	1
			41	10	7	19	5		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
2	D	1	Total	C	N	O	P	0	1
			41	10	7	19	5		

- Molecule 3 is RIBOSTAMYCIN (CCD ID: RIO) (formula: $C_{17}H_{34}N_4O_{10}$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			31	17	4	10		
3	B	1	Total	C	N	O	0	0
			31	17	4	10		
3	C	1	Total	C	N	O	0	0
			31	17	4	10		
3	D	1	Total	C	N	O	0	0
			31	17	4	10		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	2	Total 2	Mg 2	0	0

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total 1	Cl 1	0	0
5	B	1	Total 1	Cl 1	0	0
5	C	1	Total 1	Cl 1	0	0

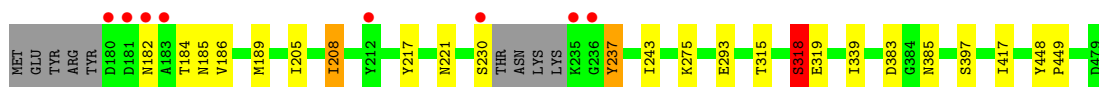
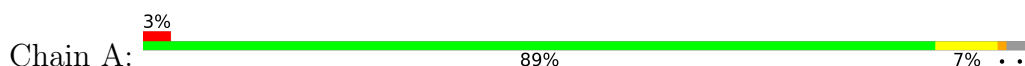
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	279	Total 279	O 279	0	0
6	B	325	Total 325	O 325	0	0
6	C	275	Total 275	O 275	0	0
6	D	219	Total 219	O 219	0	0

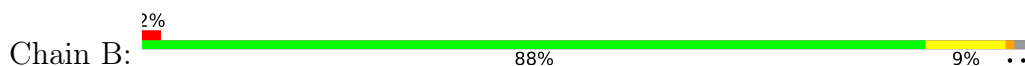
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

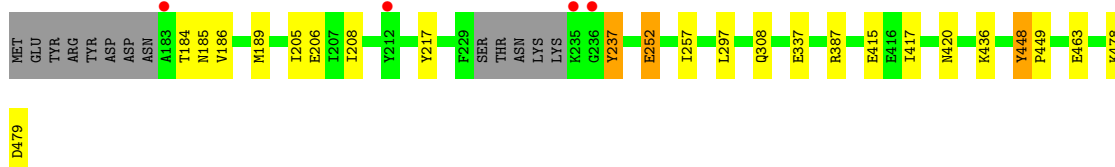
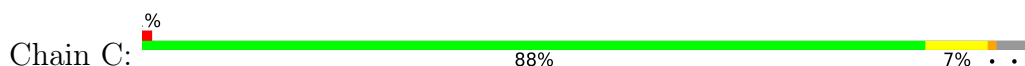
- Molecule 1: Bifunctional AAC/APH



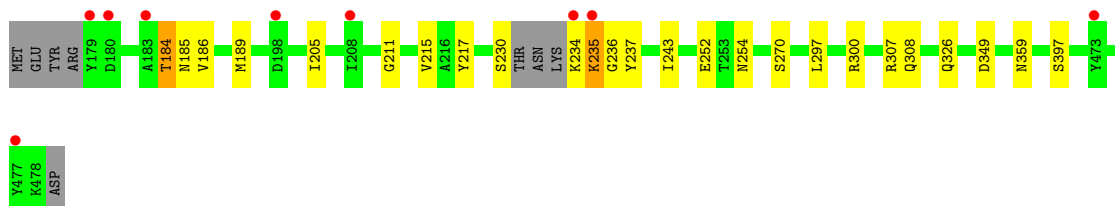
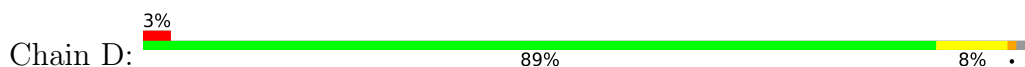
- Molecule 1: Bifunctional AAC/APH



- Molecule 1: Bifunctional AAC/APH



- Molecule 1: Bifunctional AAC/APH



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	90.22Å 99.67Å 93.40Å 90.00° 104.90° 90.00°	Depositor
Resolution (Å)	90.26 – 2.20 90.26 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (90.26-2.20) 99.9 (90.26-2.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.162 , 0.202 0.170 , 0.208	Depositor DCC
R_{free} test set	4078 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	38.5	Xtriage
Anisotropy	0.120	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.017 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	11182	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% 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: CL, MG, GNP, RIO

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	1.12	1/2498 (0.0%)	1.00	8/3369 (0.2%)
1	B	1.06	0/2528	0.95	3/3410 (0.1%)
1	C	1.07	2/2474 (0.1%)	0.99	3/3335 (0.1%)
1	D	0.96	2/2502 (0.1%)	0.91	1/3373 (0.0%)
All	All	1.05	5/10002 (0.0%)	0.96	15/13487 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	236	GLY	CA-C	5.78	1.58	1.51
1	C	448	TYR	C-O	-5.42	1.19	1.24
1	D	236	GLY	N-CA	5.40	1.51	1.44
1	A	237	TYR	N-CA	5.31	1.53	1.46
1	C	257	ILE	C-O	-5.14	1.19	1.24

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	208	ILE	CB-CA-C	-8.57	101.53	110.88
1	B	208	ILE	CB-CA-C	-7.35	102.87	110.88
1	A	417	ILE	N-CA-C	6.47	116.61	110.53
1	A	319	GLU	CB-CG-CD	6.04	122.87	112.60
1	D	237	TYR	N-CA-C	5.95	120.16	113.02
1	A	339	ILE	N-CA-C	5.91	118.68	112.83
1	B	417	ILE	N-CA-C	5.76	115.94	110.53
1	A	318	SER	CA-C-N	5.60	128.34	120.28
1	A	318	SER	C-N-CA	5.60	128.34	120.28
1	B	339	ILE	N-CA-C	5.54	118.31	112.83
1	A	237	TYR	N-CA-C	5.50	122.51	110.80
1	C	417	ILE	N-CA-C	5.45	115.66	110.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	237	TYR	N-CA-C	5.34	121.66	113.28
1	C	463	GLU	CB-CG-CD	5.14	121.35	112.60
1	A	385	ASN	CB-CA-C	-5.05	102.03	110.56

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	2447	0	2346	14	1
1	B	2478	0	2373	17	1
1	C	2425	0	2334	22	0
1	D	2453	0	2361	21	0
2	A	32	0	13	0	0
2	B	41	0	2	0	0
2	C	32	0	13	0	0
2	D	41	0	2	0	0
3	A	31	0	34	0	0
3	B	31	0	34	0	0
3	C	31	0	34	1	0
3	D	31	0	34	0	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
4	C	2	0	0	0	0
4	D	2	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	1	0
6	A	279	0	0	4	0
6	B	325	0	0	5	1
6	C	275	0	0	8	3
6	D	219	0	0	7	0
All	All	11182	0	9580	65	3

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 (65) 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)
5:C:802:CL:CL	6:C:1107:HOH:O	2.29	0.86
1:D:359:ASN:OD1	6:D:1050:HOH:O	1.99	0.80
1:C:184:THR:HG21	1:D:215:VAL:HG21	1.61	0.80
1:C:237:TYR:HA	6:C:1384:HOH:O	1.80	0.79
1:B:349[A]:ASP:OD1	6:B:1135:HOH:O	2.07	0.71
1:D:211:GLY:HA3	6:D:1091:HOH:O	1.89	0.70
1:C:185:ASN:OD1	1:D:184:THR:OG1	2.09	0.70
1:A:185:ASN:ND2	1:B:179:TYR:OH	2.26	0.68
1:D:185:ASN:ND2	1:D:217:TYR:OH	2.28	0.65
1:B:478:LYS:NZ	6:B:1272:HOH:O	2.29	0.65
1:C:184:THR:CG2	1:D:215:VAL:HG21	2.27	0.65
1:D:270:SER:OG	6:D:1212:HOH:O	2.15	0.65
1:B:463:GLU:OE1	6:B:1001:HOH:O	2.14	0.65
1:B:298:LEU:HA	6:B:1268:HOH:O	2.00	0.62
1:D:307[A]:ARG:NH2	6:D:1005:HOH:O	2.32	0.62
1:A:243:ILE:HD13	1:A:397:SER:HB2	1.83	0.59
1:C:237:TYR:CA	6:C:1384:HOH:O	2.45	0.58
1:D:254:ASN:HB2	6:D:1500:HOH:O	2.04	0.57
1:B:467:ASN:ND2	6:B:1434:HOH:O	2.38	0.56
1:B:243:ILE:HD13	1:B:397:SER:HB2	1.89	0.55
1:C:436:LYS:NZ	6:C:1113:HOH:O	2.39	0.55
1:C:186:VAL:HG13	1:C:205:ILE:HG23	1.87	0.55
1:D:186:VAL:HG13	1:D:205:ILE:HG23	1.89	0.55
1:D:326:GLN:NE2	6:D:1013:HOH:O	2.33	0.54
1:B:186:VAL:HG13	1:B:205:ILE:HG23	1.89	0.54
1:A:186:VAL:HG13	1:A:205:ILE:HG23	1.90	0.53
1:B:252:GLU:O	1:B:308:GLN:NE2	2.34	0.52
1:C:252:GLU:O	1:C:308:GLN:NE2	2.31	0.52
1:D:349:ASP:OD1	6:D:1171:HOH:O	2.19	0.52
1:A:182:ASN:HA	1:B:181:ASP:OD1	2.10	0.52
1:B:189:MET:HG2	1:B:217:TYR:CE1	2.44	0.52
1:D:243:ILE:HD13	1:D:397[A]:SER:HB2	1.92	0.51
1:A:184:THR:HG21	1:B:215:VAL:HG21	1.91	0.51
1:C:184:THR:CG2	1:D:185:ASN:OD1	2.59	0.50
1:D:252:GLU:O	1:D:308:GLN:NE2	2.36	0.50
1:C:189:MET:HG2	1:C:217:TYR:CE1	2.47	0.49
1:D:189:MET:HG2	1:D:217:TYR:CE1	2.47	0.49
1:C:186:VAL:CG1	1:C:205:ILE:HG23	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:230:SER:O	1:B:231:THR:HB	2.14	0.47
1:A:221[B]:ASN:ND2	6:A:1199:HOH:O	2.39	0.46
1:A:448:TYR:HB3	1:A:449:PRO:HD3	1.98	0.46
1:A:189:MET:HG2	1:A:217:TYR:CE1	2.50	0.46
6:A:1217:HOH:O	1:B:479:ASP:N	2.44	0.46
1:C:448:TYR:HB3	1:C:449:PRO:HD3	1.97	0.46
1:B:186:VAL:CG1	1:B:205:ILE:HG23	2.47	0.45
1:D:297:LEU:HD23	1:D:300:ARG:HH11	1.81	0.45
1:A:186:VAL:CG1	1:A:205:ILE:HG23	2.47	0.44
1:C:337:GLU:HG2	6:C:1341:HOH:O	2.17	0.43
1:D:186:VAL:CG1	1:D:205:ILE:HG23	2.47	0.43
1:C:420:ASN:ND2	6:C:1031:HOH:O	2.51	0.43
1:A:315:THR:O	1:A:318:SER:OG	2.33	0.43
1:C:308:GLN:HB2	6:C:1191:HOH:O	2.19	0.43
1:C:478:LYS:O	1:C:479:ASP:CB	2.67	0.43
1:A:184:THR:CG2	1:B:185:ASN:ND2	2.82	0.42
1:C:415:GLU:OE1	3:C:600:RIO:N2	2.52	0.42
1:A:275:LYS:NZ	6:A:1190:HOH:O	2.51	0.42
1:C:297:LEU:HD23	1:C:387:ARG:HD3	2.02	0.42
1:B:234:LYS:O	1:B:235:LYS:HB2	2.20	0.41
1:D:234:LYS:O	1:D:235:LYS:HB2	2.20	0.41
1:C:184:THR:HG22	1:D:185:ASN:OD1	2.21	0.41
1:C:206:GLU:HB3	6:C:1328:HOH:O	2.21	0.41
1:C:478:LYS:O	1:C:479:ASP:HB2	2.21	0.41
1:A:237:TYR:HA	6:A:1384:HOH:O	2.22	0.40
1:A:383:ASP:OD1	1:A:383:ASP:C	2.64	0.40
1:C:184:THR:HB	1:D:185:ASN:OD1	2.21	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:293:GLU:OE1	6:C:1080:HOH:O[2_747]	2.11	0.09
6:B:1391:HOH:O	6:C:1240:HOH:O[2_746]	2.13	0.07
1:B:288:SER:O	6:C:1219:HOH:O[2_646]	2.18	0.02

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	296/305 (97%)	285 (96%)	11 (4%)	0	100	100
1	B	298/305 (98%)	286 (96%)	12 (4%)	0	100	100
1	C	291/305 (95%)	280 (96%)	11 (4%)	0	100	100
1	D	296/305 (97%)	285 (96%)	10 (3%)	1 (0%)	36	42
All	All	1181/1220 (97%)	1136 (96%)	44 (4%)	1 (0%)	48	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	235	LYS

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	271/281 (96%)	268 (99%)	3 (1%)	65	79
1	B	275/281 (98%)	267 (97%)	8 (3%)	37	51
1	C	269/281 (96%)	267 (99%)	2 (1%)	76	87
1	D	271/281 (96%)	269 (99%)	2 (1%)	76	87
All	All	1086/1124 (97%)	1071 (99%)	15 (1%)	59	75

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	208	ILE
1	A	230	SER
1	A	318	SER
1	B	182	ASN
1	B	184	THR
1	B	208	ILE
1	B	230	SER
1	B	231	THR
1	B	285	GLU
1	B	463	GLU
1	B	470	LYS
1	C	208	ILE
1	C	252	GLU
1	D	184	THR
1	D	230	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	182	ASN
1	A	308	GLN
1	A	341	ASN
1	A	420	ASN
1	A	431	ASN
1	A	467	ASN
1	B	467	ASN
1	C	327	ASN
1	C	386	ASN
1	D	296	ASN
1	D	326	GLN
1	D	420	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](#)

Of 21 ligands modelled in this entry, 11 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
2	GNP	D	500[A]	4	34,34,34	0.85	1 (2%)	47,54,54	0.97	4 (8%)
3	RIO	A	600	-	32,33,33	0.47	0	46,49,49	1.10	3 (6%)
3	RIO	D	600	-	32,33,33	0.37	0	46,49,49	1.11	5 (10%)
2	GNP	A	500	4	34,34,34	1.45	3 (8%)	47,54,54	1.16	5 (10%)
2	GNP	B	500[B]	4	34,34,34	0.73	0	47,54,54	1.02	3 (6%)
3	RIO	C	600	-	32,33,33	0.44	0	46,49,49	1.04	2 (4%)
3	RIO	B	600	-	32,33,33	0.43	0	46,49,49	1.20	3 (6%)
2	GNP	C	500	4	34,34,34	1.07	2 (5%)	47,54,54	0.86	2 (4%)
2	GNP	D	500[B]	4	34,34,34	0.86	2 (5%)	47,54,54	0.69	1 (2%)
2	GNP	B	500[A]	4	34,34,34	0.77	1 (2%)	47,54,54	0.81	2 (4%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GNP	D	500[A]	4	-	2/18/38/38	0/3/3/3
3	RIO	A	600	-	-	3/12/68/68	0/3/3/3
3	RIO	D	600	-	-	3/12/68/68	0/3/3/3
2	GNP	A	500	4	-	3/18/38/38	0/3/3/3
2	GNP	B	500[B]	4	-	4/18/38/38	0/3/3/3
3	RIO	C	600	-	-	2/12/68/68	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	RIO	B	600	-	-	2/12/68/68	0/3/3/3
2	GNP	C	500	4	-	5/18/38/38	0/3/3/3
2	GNP	D	500[B]	4	-	3/18/38/38	0/3/3/3
2	GNP	B	500[A]	4	-	2/18/38/38	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	GNP	PB-O2B	5.32	1.71	1.56
2	C	500	GNP	PB-O3A	-4.93	1.53	1.59
2	A	500	GNP	PG-O2G	4.60	1.69	1.56
2	A	500	GNP	PB-O3A	-3.79	1.54	1.59
2	D	500[B]	GNP	PB-O3A	2.72	1.62	1.59
2	D	500[A]	GNP	PB-O3A	2.60	1.62	1.59
2	D	500[B]	GNP	PB-O2B	2.29	1.62	1.56
2	B	500[A]	GNP	PB-O2B	2.20	1.62	1.56
2	C	500	GNP	PG-O2G	2.16	1.62	1.56

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500[B]	GNP	O1G-PG-N3B	-5.12	104.23	111.77
3	B	600	RIO	O3-C3-O2	-4.42	106.85	111.37
2	D	500[A]	GNP	O1G-PG-N3B	3.82	117.40	111.77
2	A	500	GNP	O1G-PG-N3B	-3.51	106.60	111.77
2	A	500	GNP	O2G-PG-O1G	-3.37	105.00	113.45
3	C	600	RIO	O3-C4-C9	3.30	115.79	107.42
2	A	500	GNP	O2G-PG-O3G	3.09	115.88	107.59
2	C	500	GNP	O1G-PG-N3B	-3.05	107.27	111.77
3	D	600	RIO	C4-C9-C8	-3.02	105.70	111.31
2	A	500	GNP	O2'-C2'-C1'	-2.96	99.90	110.10
3	B	600	RIO	O3-C4-C9	2.77	114.46	107.42
2	D	500[A]	GNP	O3A-PB-N3B	2.63	113.89	106.59
2	D	500[B]	GNP	O2B-PB-O3A	2.43	112.75	104.64
3	A	600	RIO	C3-O3-C4	-2.42	112.24	117.98
3	A	600	RIO	O3-C3-O2	2.41	113.83	111.37
3	B	600	RIO	C4-C9-C8	-2.33	106.99	111.31
3	A	600	RIO	C4-C9-C8	-2.30	107.05	111.31
2	B	500[B]	GNP	O2A-PA-O3A	2.29	113.47	107.27
2	B	500[A]	GNP	O2B-PB-O1B	-2.28	104.97	109.87
2	C	500	GNP	O2B-PB-O3A	2.27	112.21	104.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	500[A]	GNP	O3A-PA-O1A	2.19	117.30	110.70
3	C	600	RIO	O5-C10-C15	2.18	111.64	108.08
3	D	600	RIO	C10-O5-C9	2.17	123.12	117.98
2	D	500[A]	GNP	O2G-PG-O3G	-2.16	101.78	107.59
3	D	600	RIO	O5-C9-C4	2.13	112.82	107.42
2	B	500[A]	GNP	O2'-C2'-C1'	-2.04	103.07	110.10
2	B	500[B]	GNP	O2'-C2'-C1'	-2.04	103.07	110.10
3	D	600	RIO	O5-C9-C8	-2.03	104.33	109.18
2	A	500	GNP	O3'-C3'-C2'	-2.01	105.36	111.82
3	D	600	RIO	C3-O3-C4	-2.01	113.22	117.98

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	500	GNP	PG-N3B-PB-O1B
2	B	500[A]	GNP	PG-N3B-PB-O1B
2	B	500[B]	GNP	PG-N3B-PB-O1B
2	D	500[A]	GNP	PB-N3B-PG-O1G
2	D	500[A]	GNP	PG-N3B-PB-O1B
2	D	500[B]	GNP	PB-N3B-PG-O1G
2	D	500[B]	GNP	PG-N3B-PB-O1B
3	B	600	RIO	O6-C10-O5-C9
3	C	600	RIO	O6-C10-O5-C9
3	A	600	RIO	C4-C9-O5-C10
3	D	600	RIO	C4-C9-O5-C10
3	A	600	RIO	O6-C10-O5-C9
3	B	600	RIO	C4-C9-O5-C10
3	A	600	RIO	C8-C9-O5-C10
3	D	600	RIO	C8-C9-O5-C10
3	C	600	RIO	C4-C9-O5-C10
2	C	500	GNP	PG-N3B-PB-O3A
2	A	500	GNP	PB-O3A-PA-O1A
2	B	500[B]	GNP	PB-O3A-PA-O2A
3	D	600	RIO	O6-C10-O5-C9
2	B	500[B]	GNP	PB-O3A-PA-O1A
2	C	500	GNP	PB-O3A-PA-O1A
2	D	500[B]	GNP	PB-O3A-PA-O1A
2	C	500	GNP	PB-O3A-PA-O2A
2	C	500	GNP	PG-N3B-PB-O1B
2	B	500[A]	GNP	PA-O3A-PB-O1B
2	A	500	GNP	PG-N3B-PB-O3A

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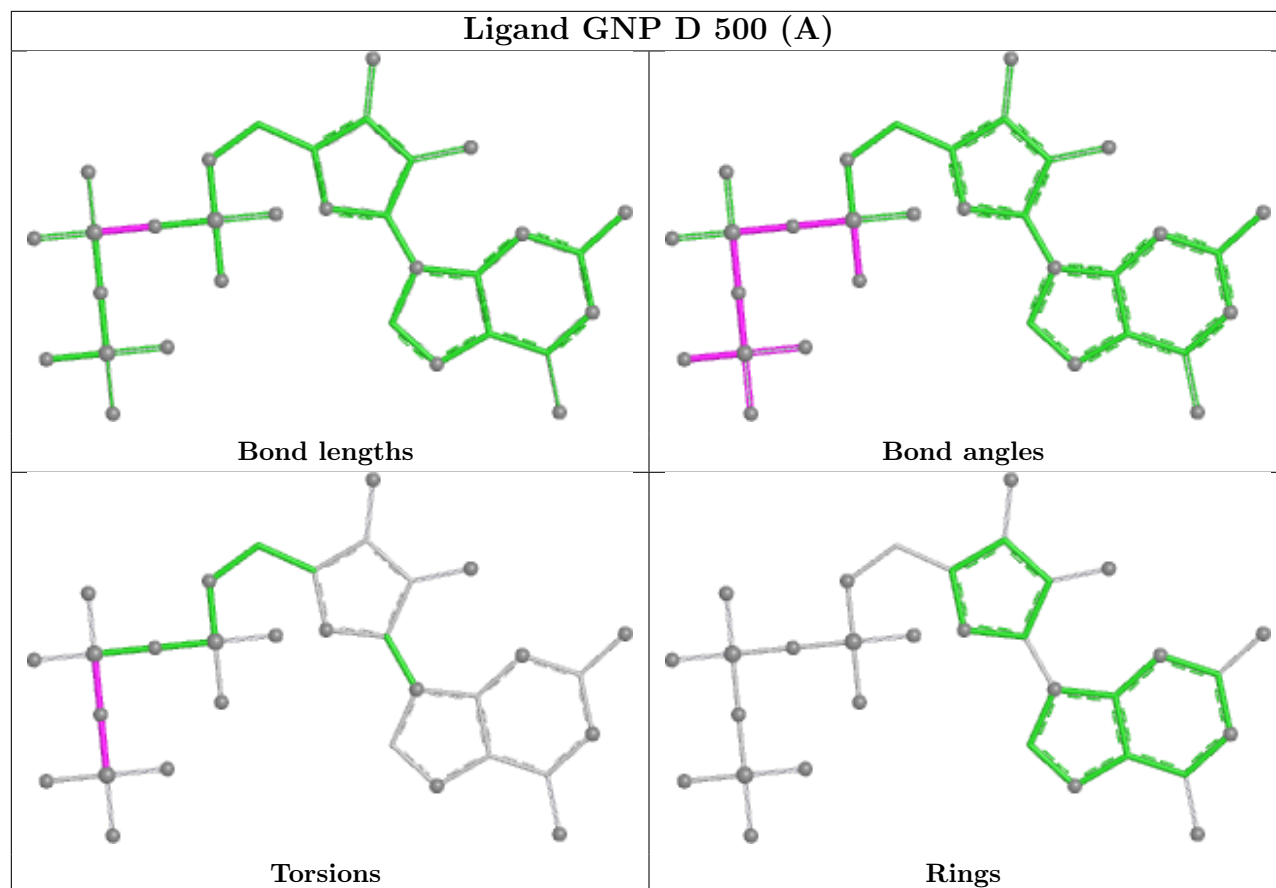
Mol	Chain	Res	Type	Atoms
2	B	500[B]	GNP	PG-N3B-PB-O3A
2	C	500	GNP	PB-N3B-PG-O1G

There are no ring outliers.

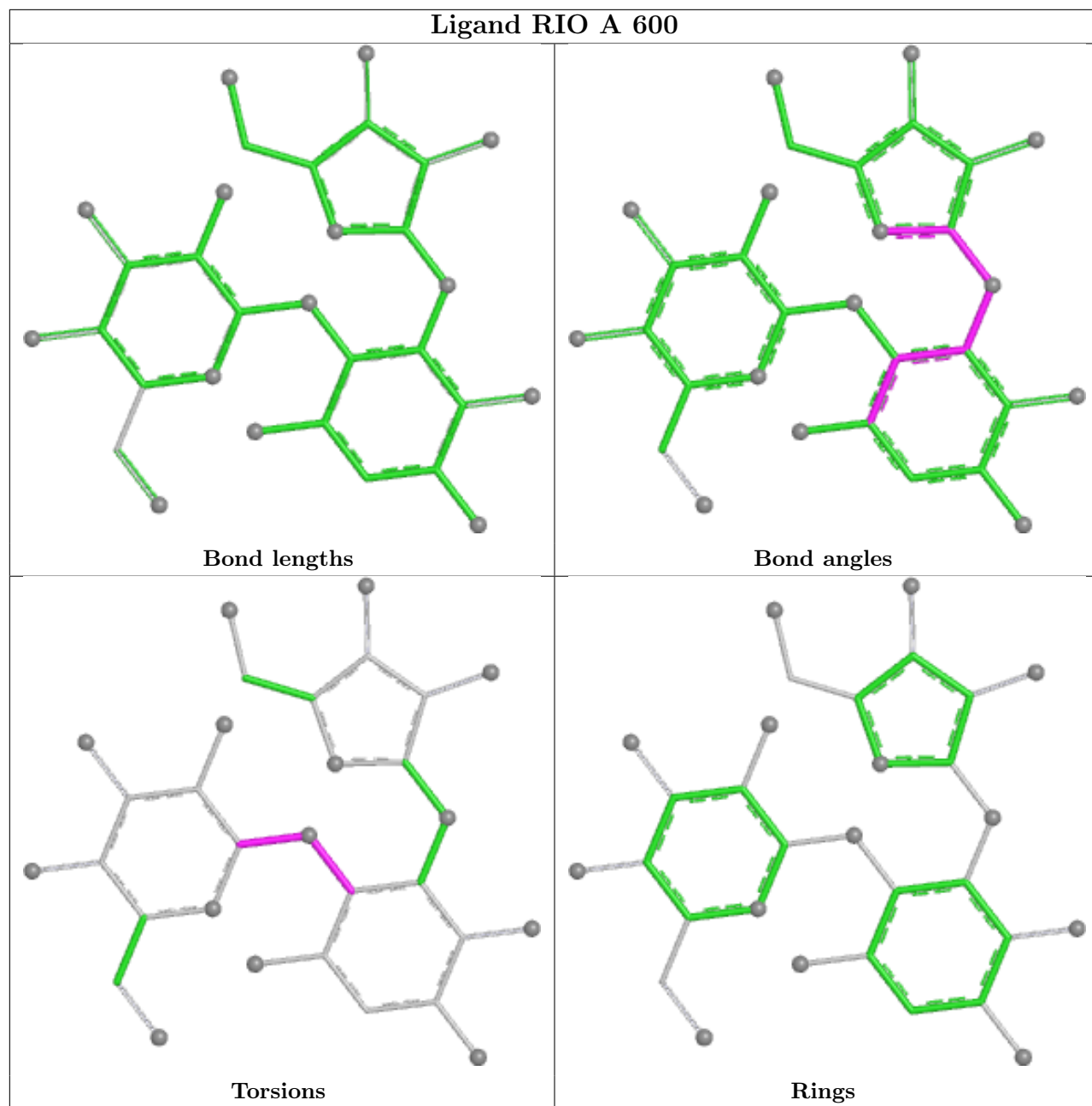
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	600	RIO	1	0

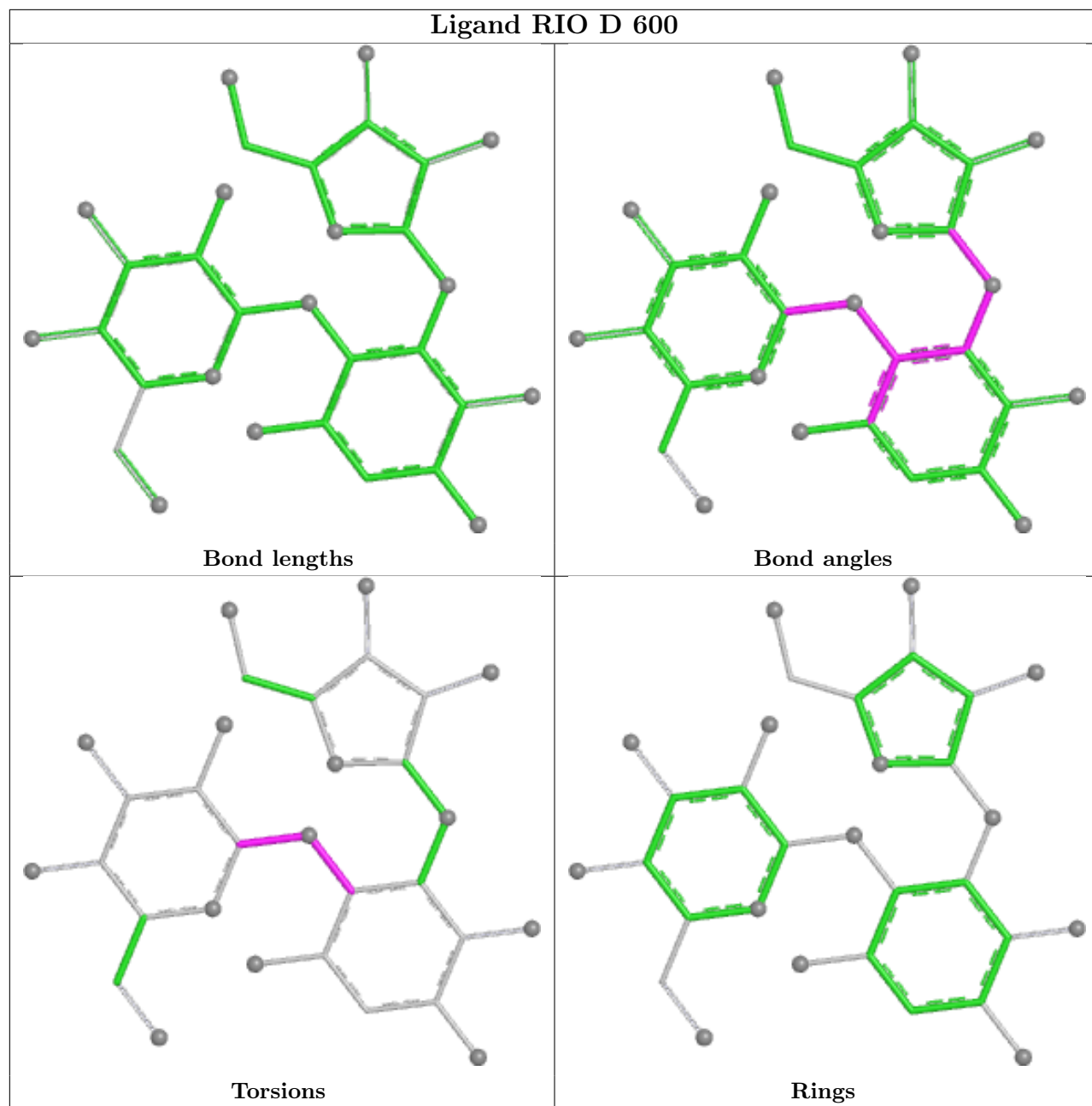
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

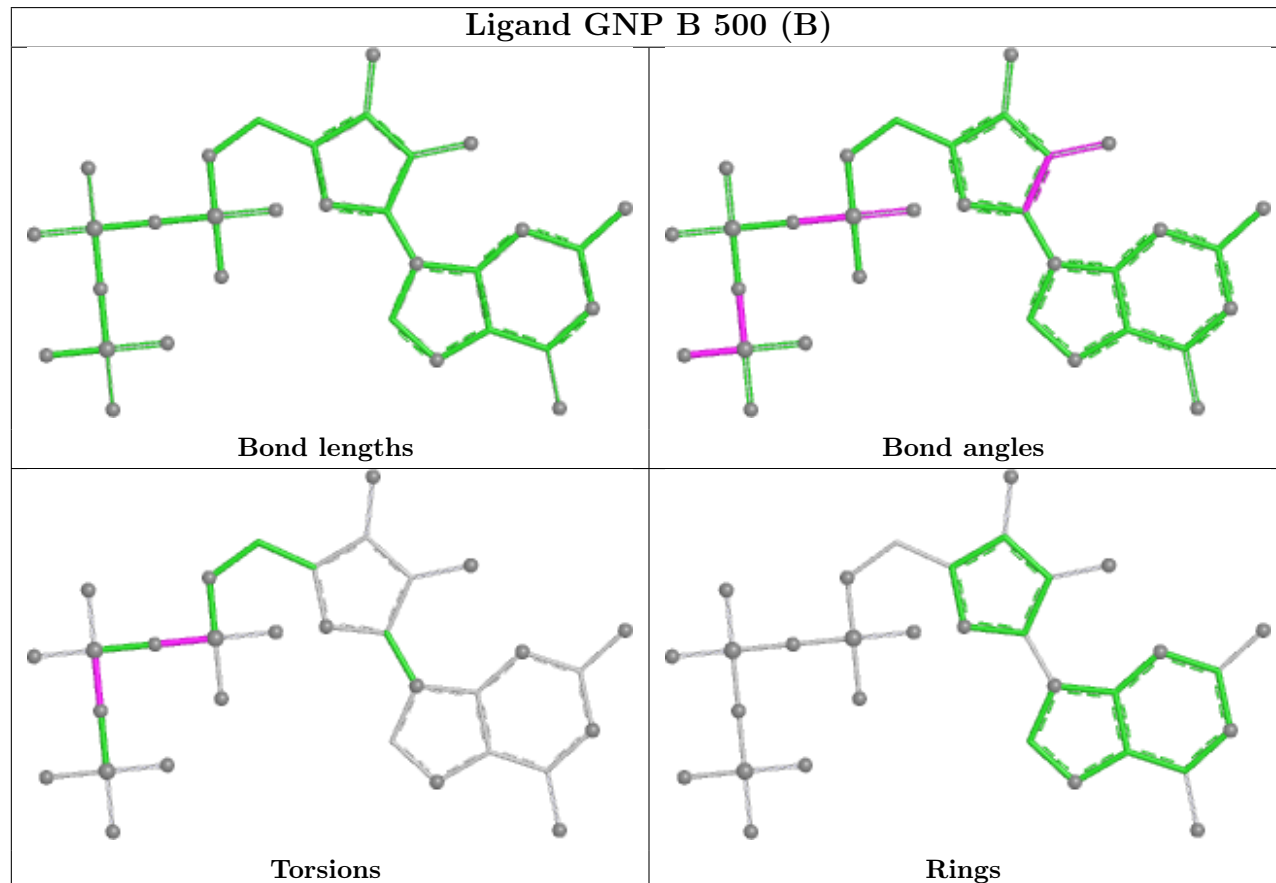
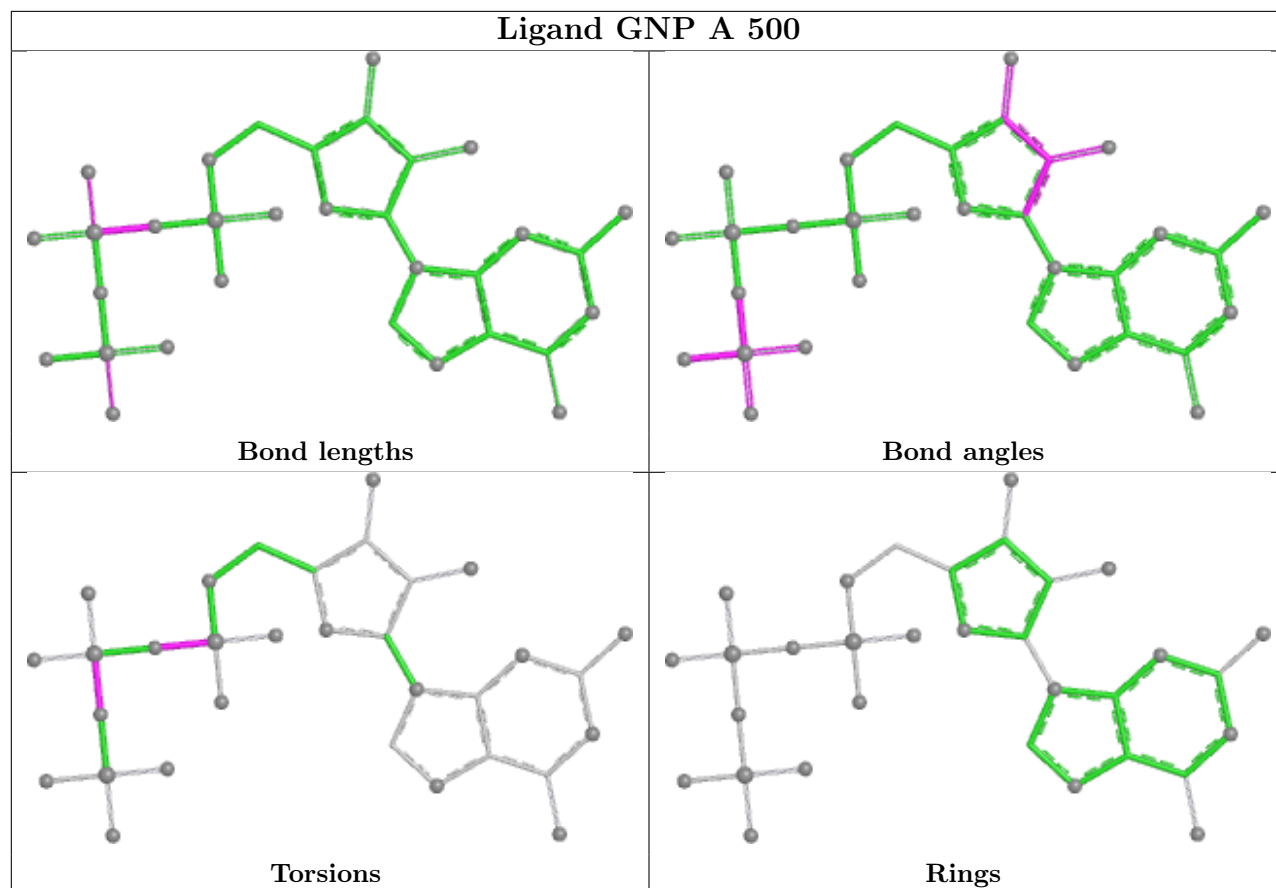


Ligand RIO A 600

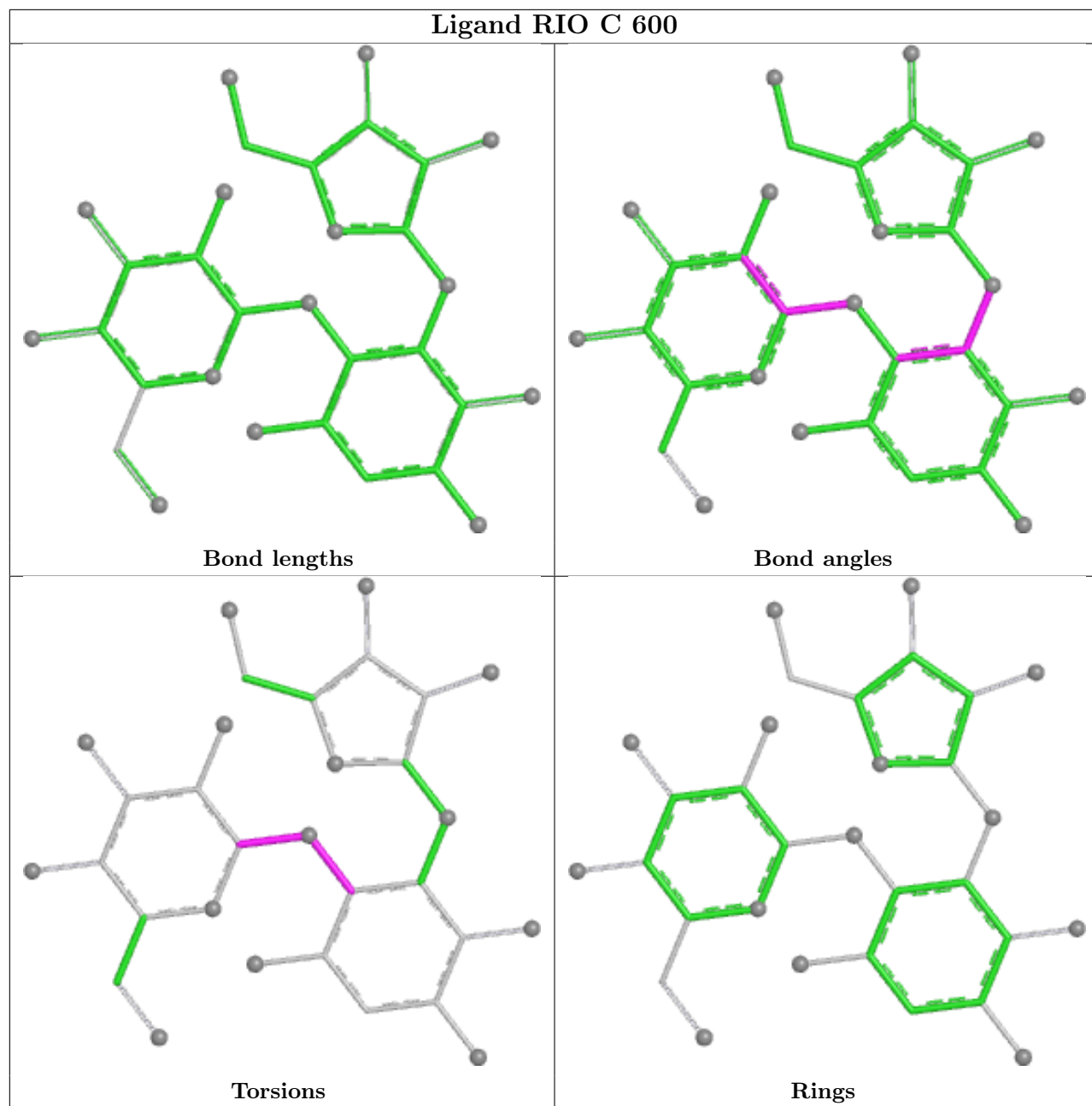


Ligand RIO D 600

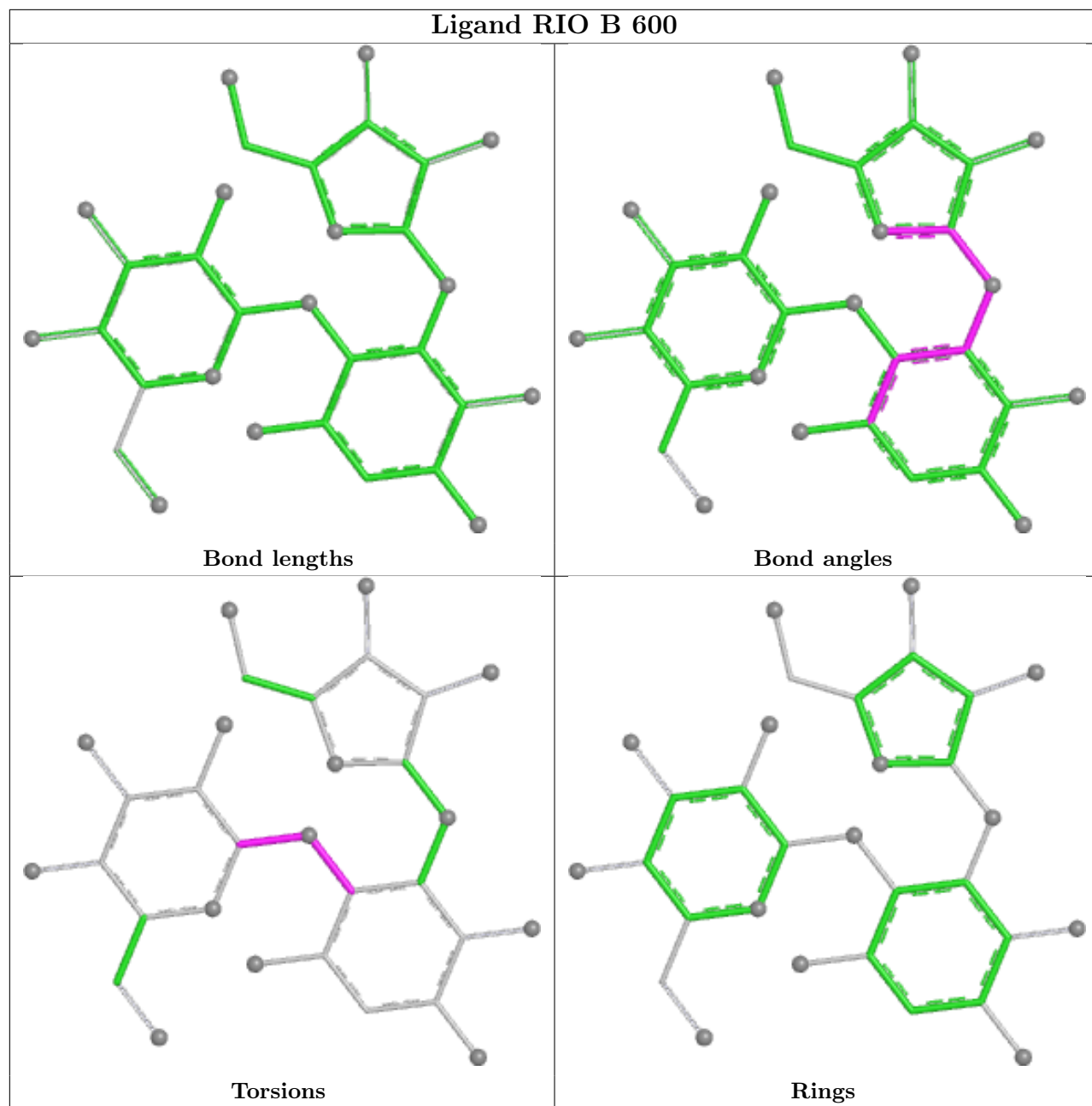


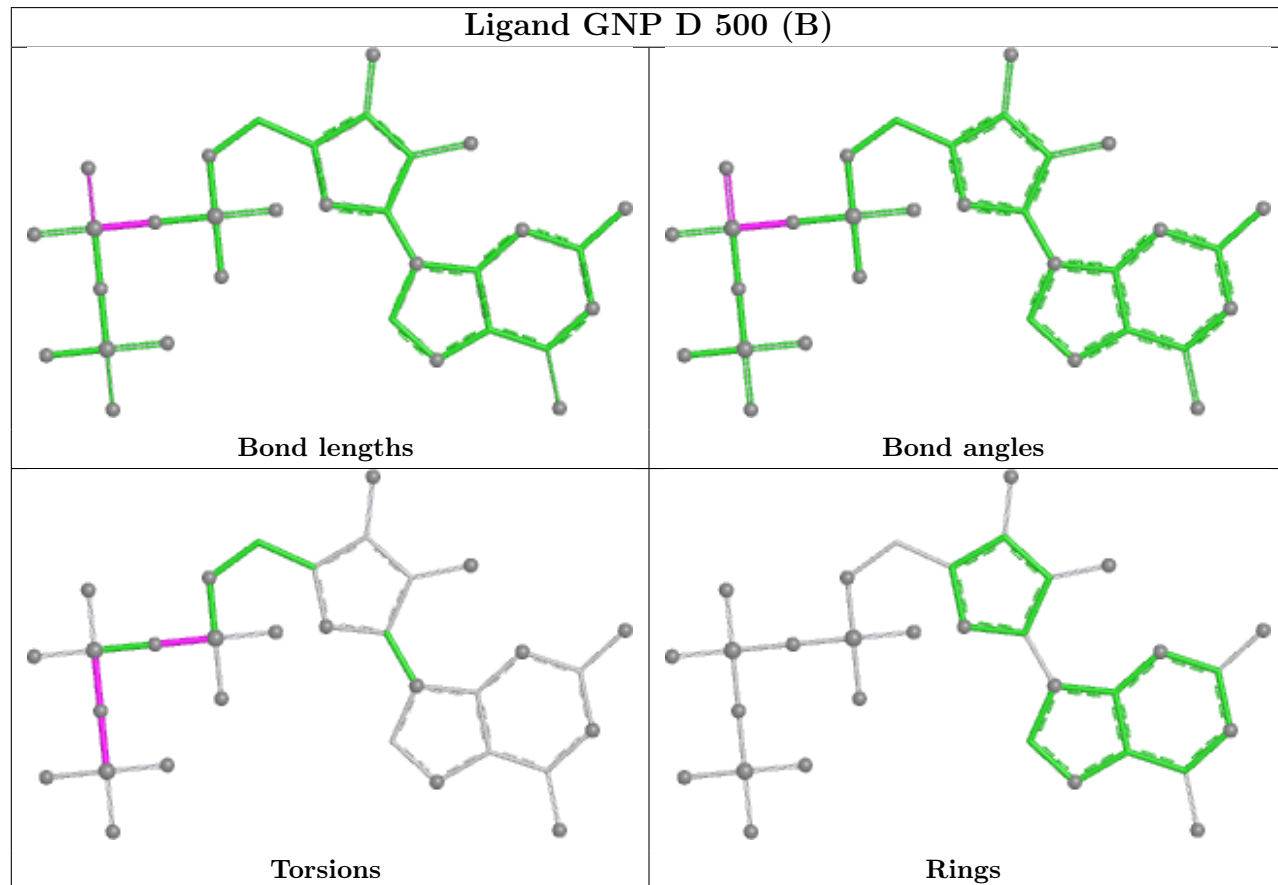
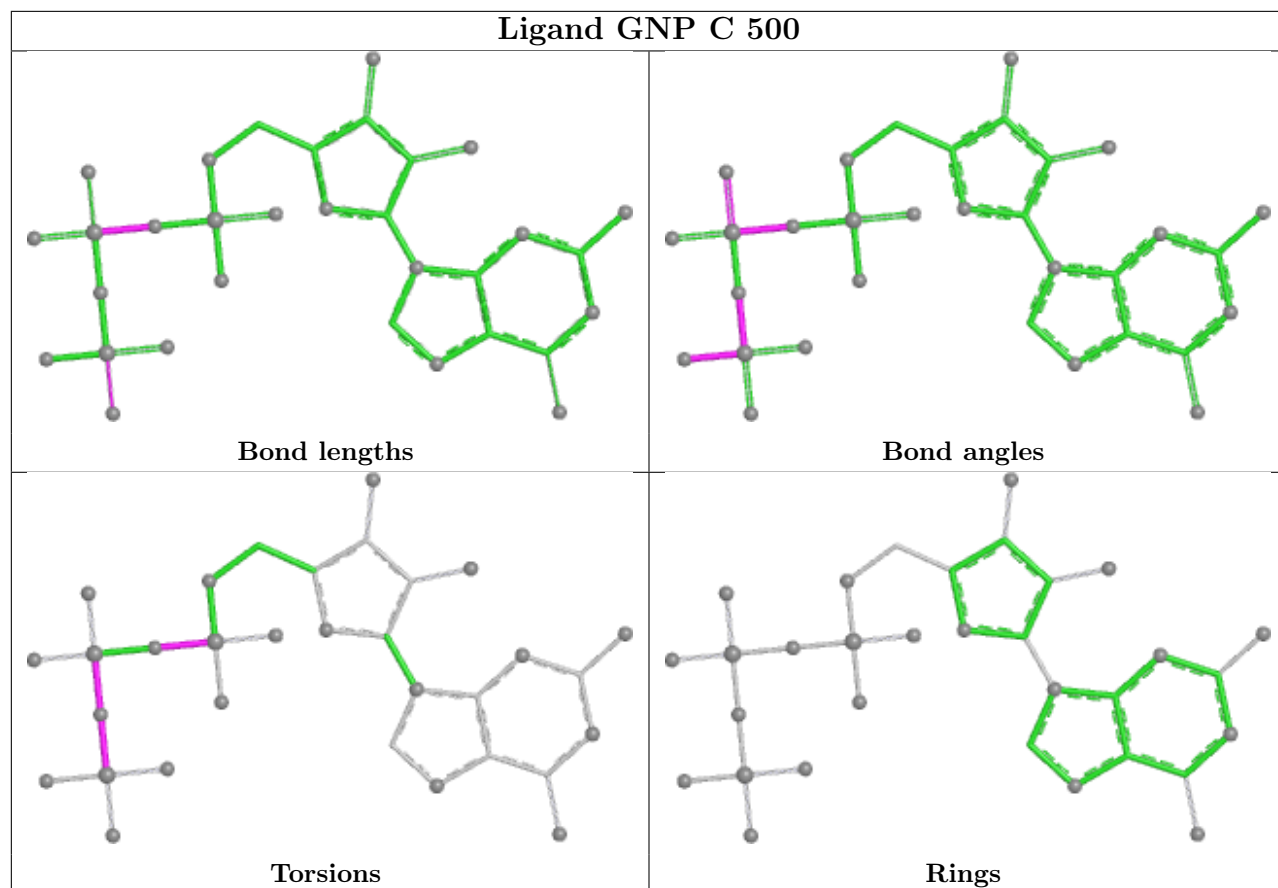


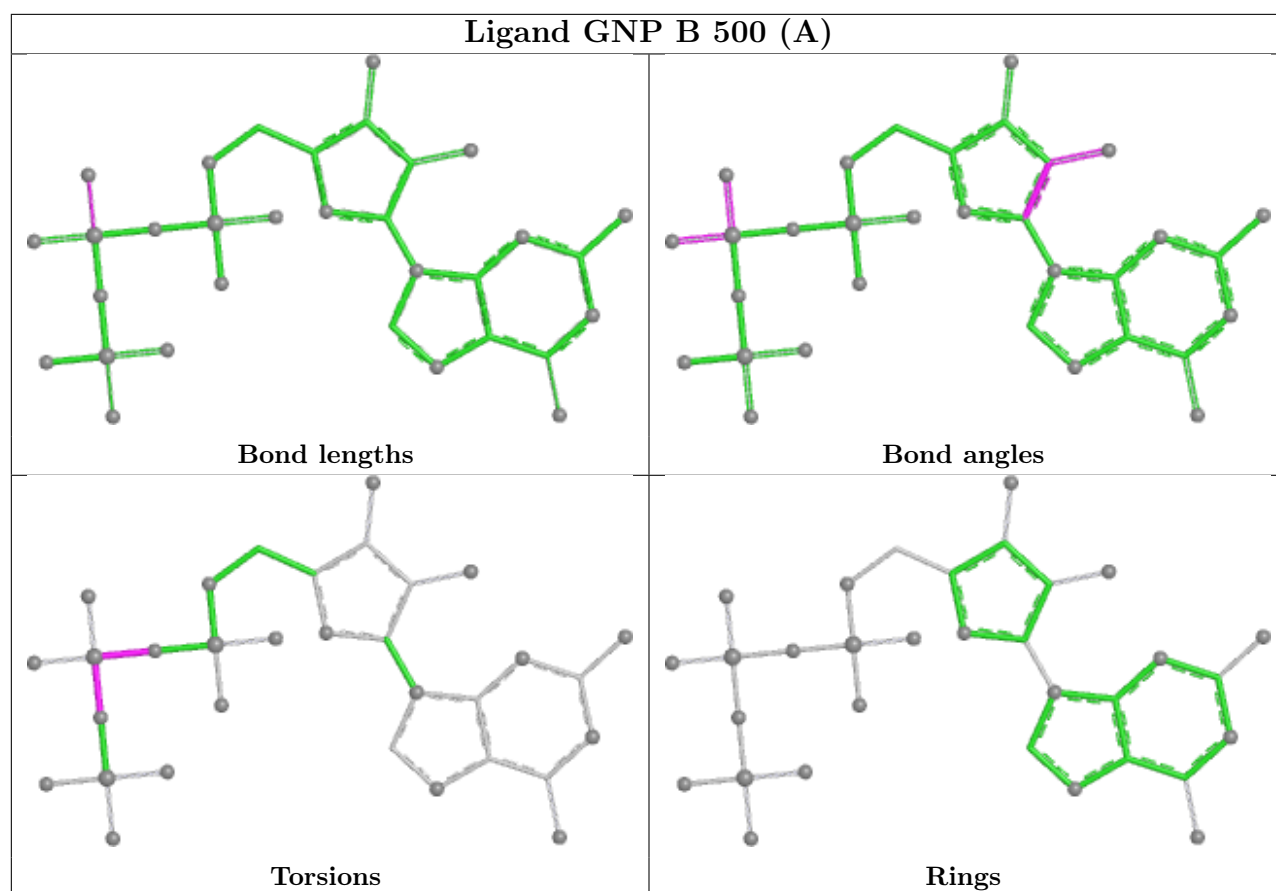
Ligand RIO C 600



Ligand RIO B 600







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	296/305 (97%)	-0.34	8 (2%) 56 53	19, 41, 64, 121	4 (1%)
1	B	299/305 (98%)	-0.36	5 (1%) 69 66	19, 40, 72, 103	3 (1%)
1	C	292/305 (95%)	-0.32	4 (1%) 73 71	21, 42, 73, 96	3 (1%)
1	D	297/305 (97%)	-0.08	9 (3%) 52 49	20, 50, 91, 126	3 (1%)
All	All	1184/1220 (97%)	-0.28	26 (2%) 62 59	19, 43, 77, 126	13 (1%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	236	GLY	4.8
1	C	235	LYS	4.1
1	B	479	ASP	4.1
1	C	183	ALA	3.6
1	B	234	LYS	3.4
1	D	235	LYS	3.4
1	D	234	LYS	3.4
1	D	477	TYR	3.4
1	A	235	LYS	3.4
1	A	181	ASP	3.0
1	B	235	LYS	3.0
1	D	473	TYR	3.0
1	C	212	TYR	3.0
1	A	180	ASP	2.9
1	D	179	TYR	2.8
1	D	180	ASP	2.8
1	B	231	THR	2.8
1	D	198	ASP	2.7
1	D	183	ALA	2.7
1	A	182	ASN	2.7
1	C	236	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	183	ALA	2.5
1	D	208	ILE	2.5
1	A	212	TYR	2.2
1	A	230	SER	2.1
1	B	179	TYR	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, 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
3	RIO	C	600	31/31	0.84	0.13	64,70,80,88	0
3	RIO	A	600	31/31	0.93	0.08	41,49,56,59	0
5	CL	C	802	1/1	0.94	0.11	63,63,63,63	0
5	CL	B	802	1/1	0.95	0.10	72,72,72,72	0
3	RIO	D	600	31/31	0.95	0.07	46,50,60,65	0
3	RIO	B	600	31/31	0.96	0.06	35,38,46,47	0
5	CL	A	802	1/1	0.96	0.09	62,62,62,62	0
2	GNP	B	500[A]	32/32	0.98	0.05	29,34,37,40	9
2	GNP	B	500[B]	32/32	0.98	0.05	29,34,37,40	9
2	GNP	C	500	32/32	0.98	0.05	35,39,43,44	0
4	MG	A	700	1/1	0.98	0.04	35,35,35,35	0
2	GNP	D	500[A]	32/32	0.98	0.05	34,39,45,48	9
2	GNP	D	500[B]	32/32	0.98	0.05	34,39,41,42	9
2	GNP	A	500	32/32	0.98	0.05	29,34,38,40	0
4	MG	A	702	1/1	0.99	0.04	37,37,37,37	0
4	MG	C	700	1/1	0.99	0.02	34,34,34,34	0
4	MG	D	702	1/1	0.99	0.02	43,43,43,43	0
4	MG	B	702	1/1	1.00	0.02	36,36,36,36	0

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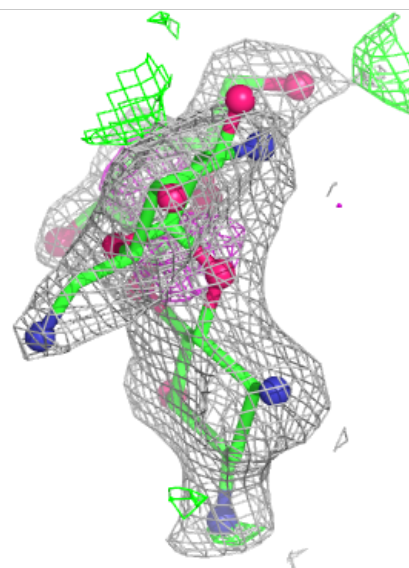
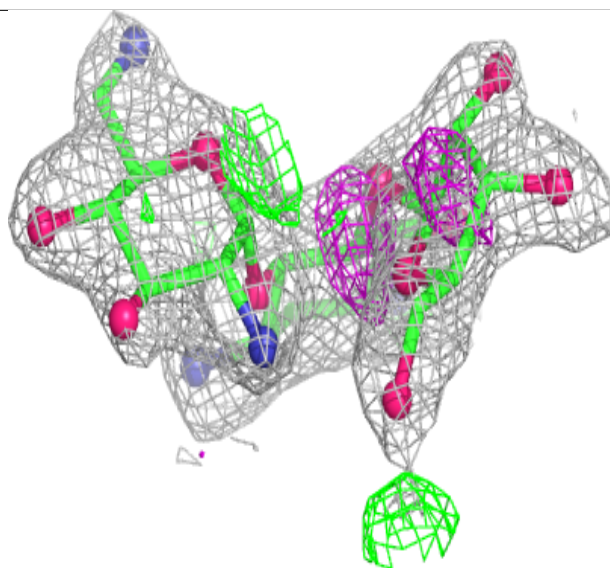
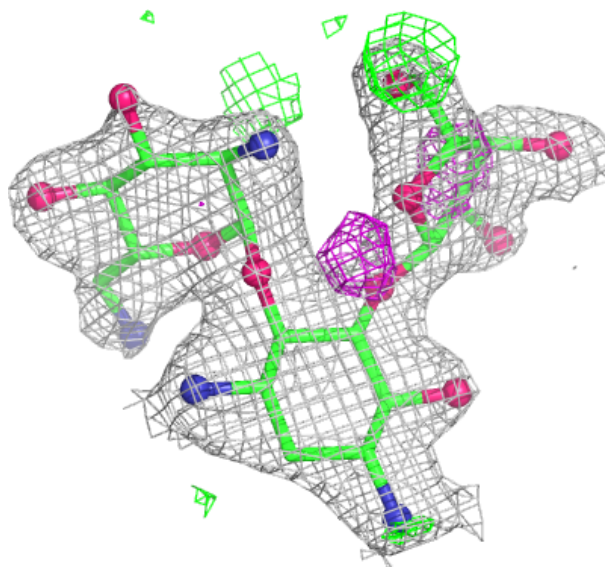
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MG	B	700	1/1	1.00	0.02	30,30,30,30	0
4	MG	C	702	1/1	1.00	0.04	38,38,38,38	0
4	MG	D	700	1/1	1.00	0.02	31,31,31,31	0

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.

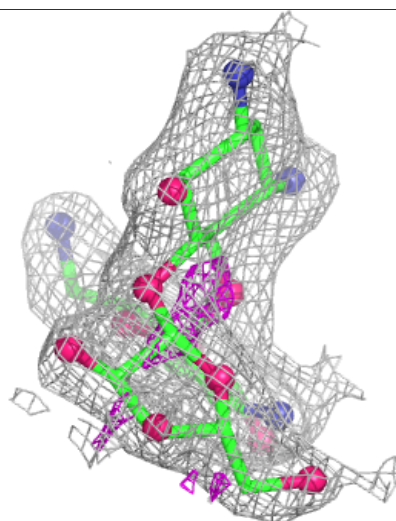
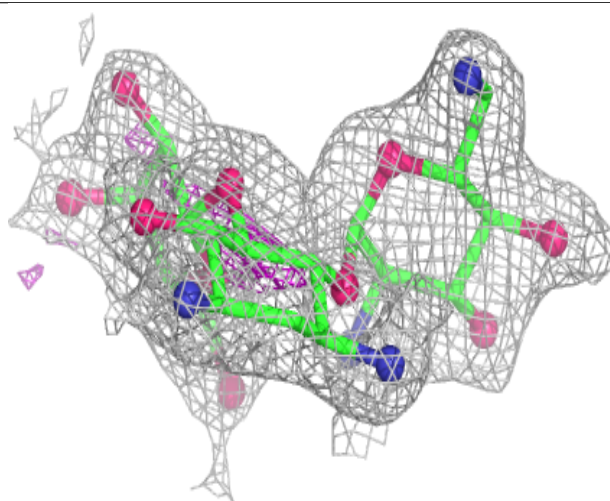
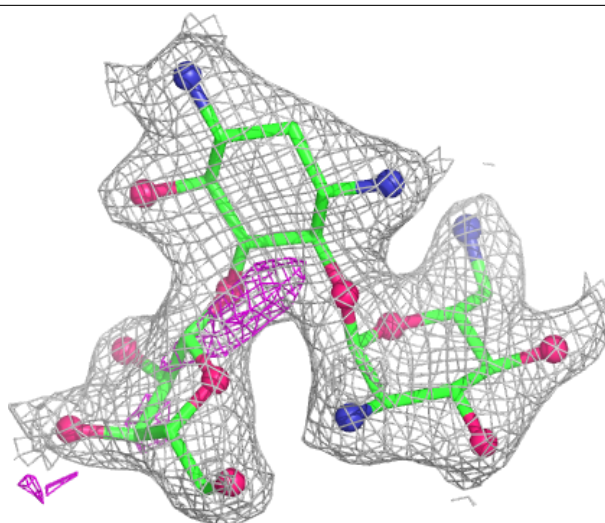
Electron density around RIO C 600:

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



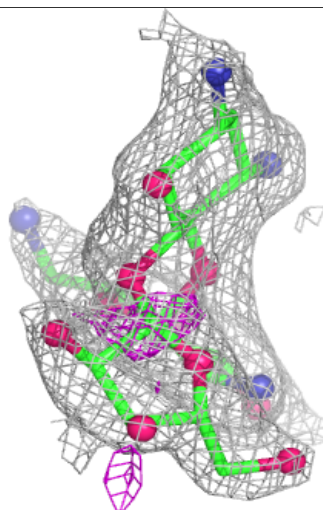
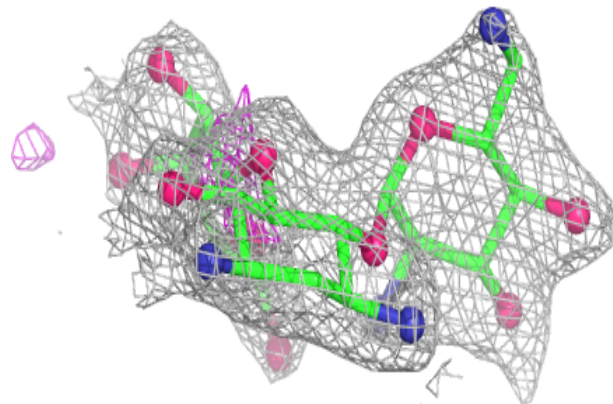
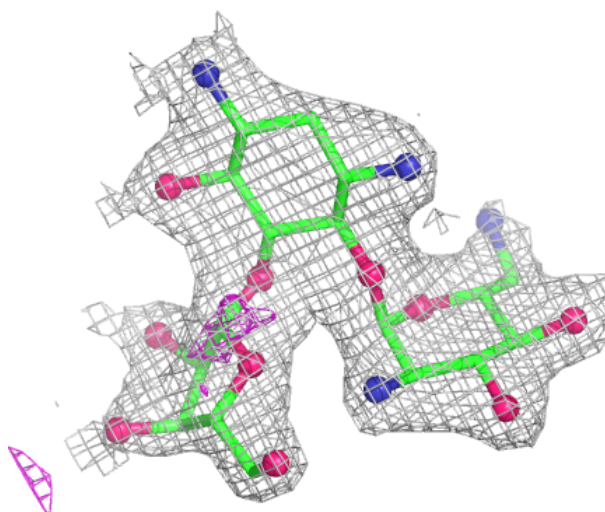
Electron density around RIO A 600:

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



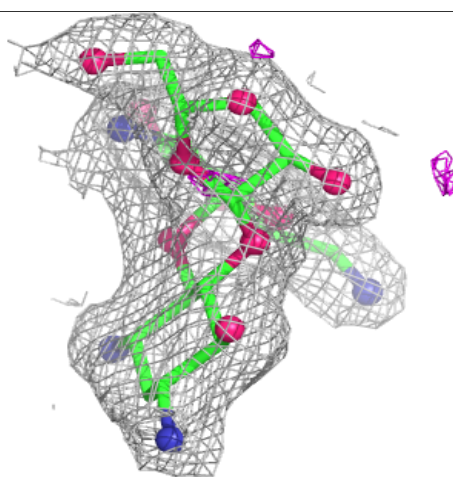
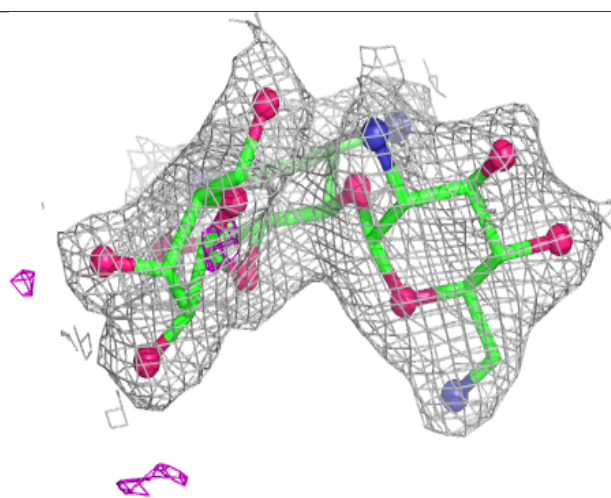
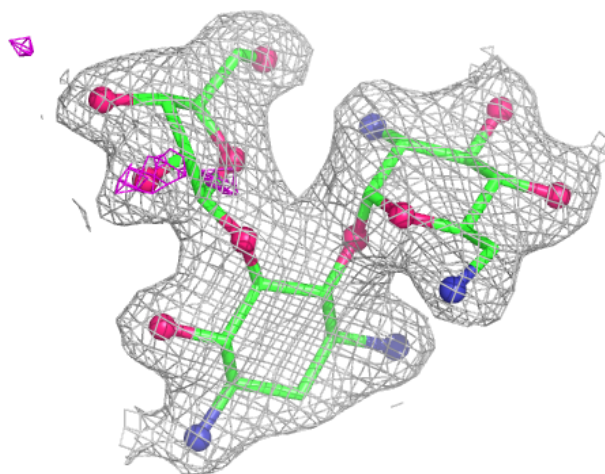
Electron density around RIO D 600:

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



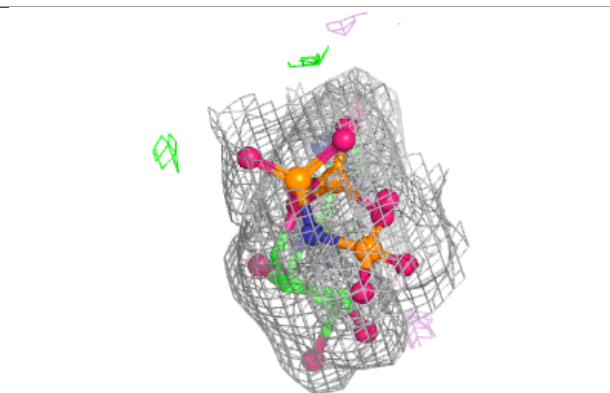
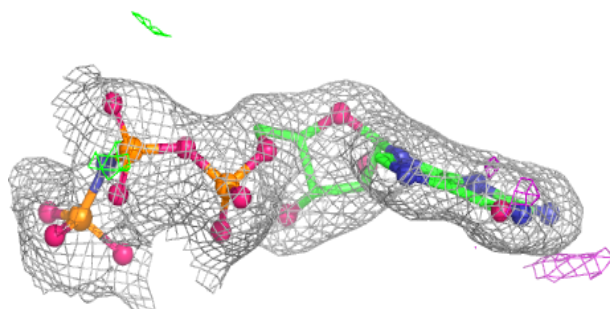
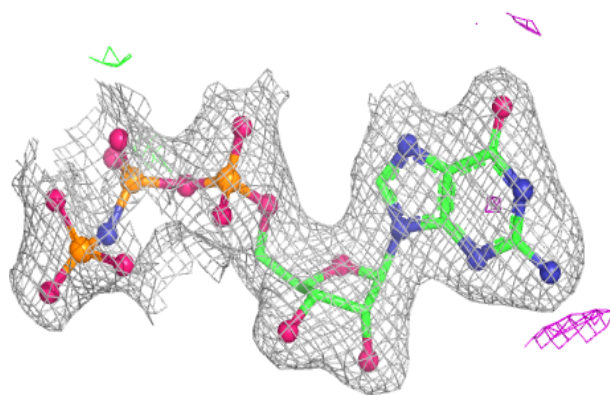
Electron density around RIO B 600:

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

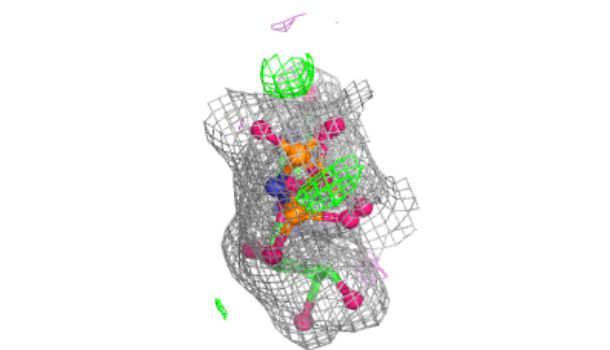
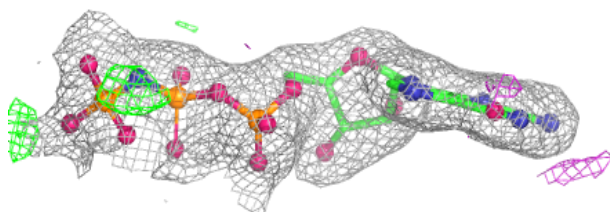
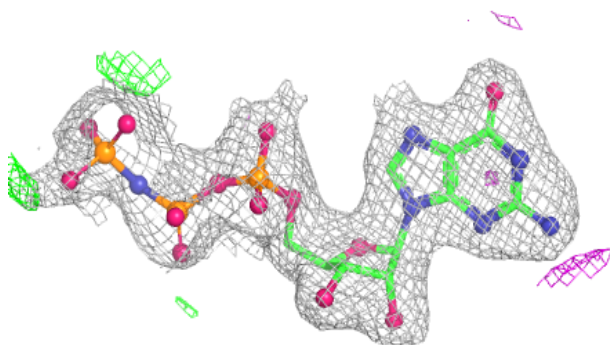


Electron density around GNP B 500 (A):

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

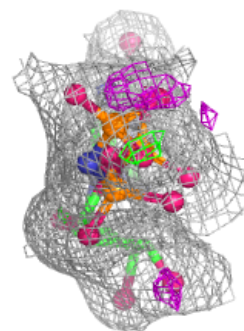
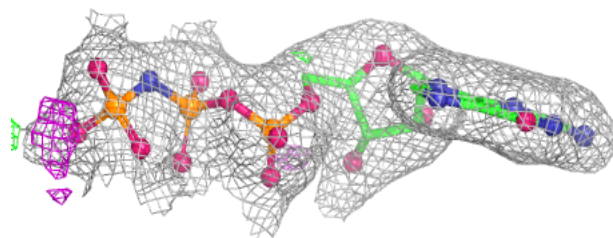
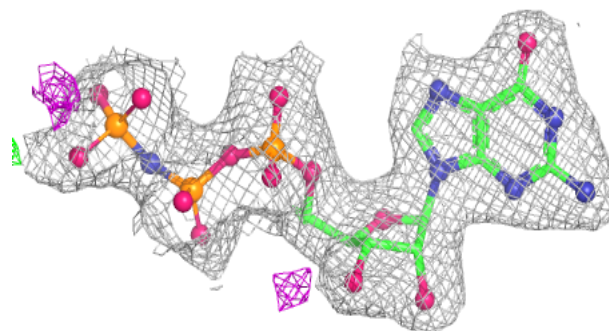
**Electron density around GNP B 500 (B):**

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

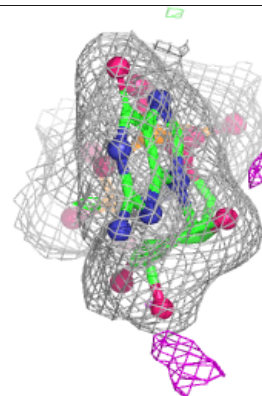
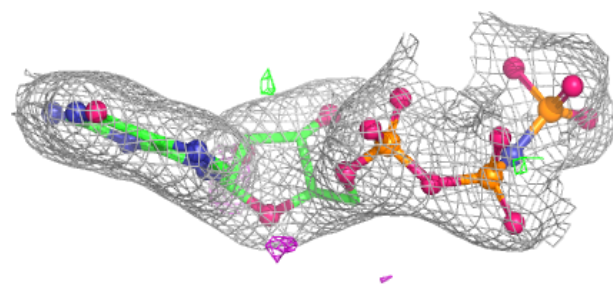
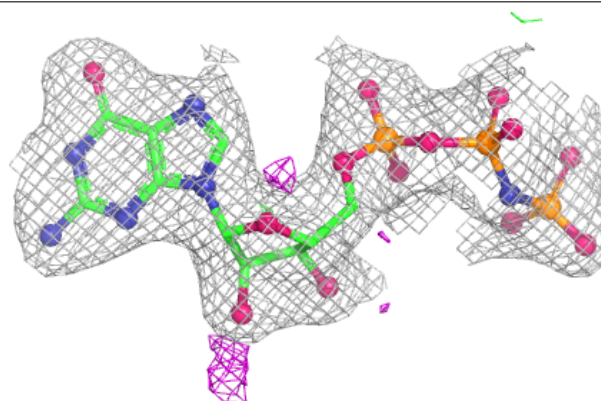


Electron density around GNP C 500:

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

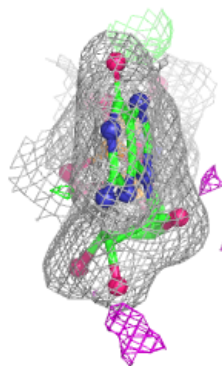
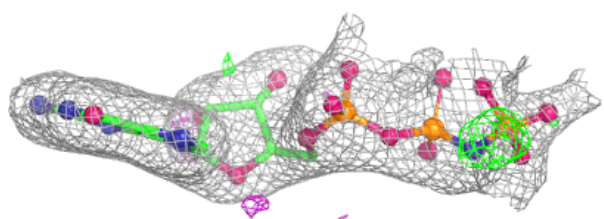
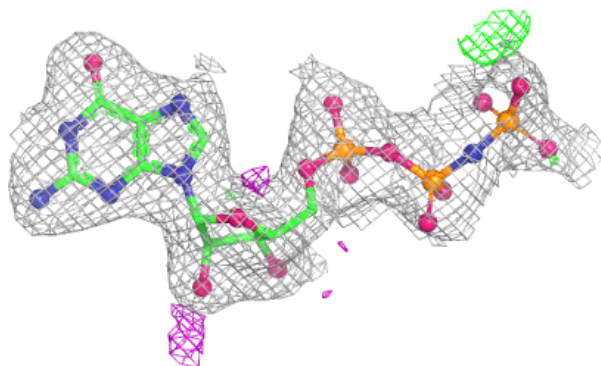
**Electron density around GNP D 500 (A):**

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

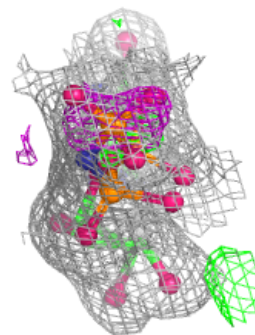
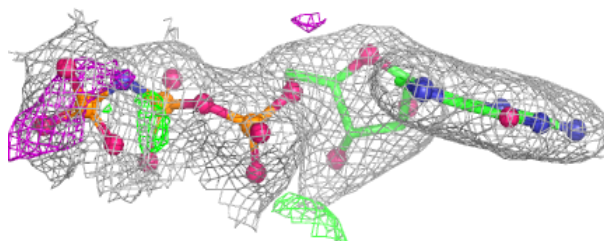
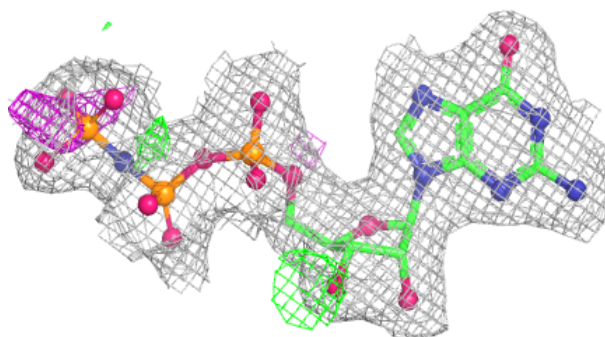


Electron density around GNP D 500 (B):

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

**Electron density around GNP A 500:**

$2mF_o-DF_c$ (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.