



# wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 10, 2023 – 02:31 AM EDT

PDB ID : 7MNZ  
Title : Crystal Structure of Nup358/RanBP2 Ran-binding domain 4 in complex with Ran-GPPNHP  
Authors : Bley, C.J.; Nie, S.; Mobbs, G.W.; Petrovic, S.; Gres, A.T.; Liu, X.; Mukherjee, S.; Harvey, S.; Huber, F.M.; Lin, D.H.; Brown, B.; Tang, A.W.; Rundlet, E.J.; Correia, A.R.; Chen, S.; Regmi, S.G.; Stevens, T.A.; Jette, C.A.; Dasso, M.; Patke, A.; Palazzo, A.F.; Kossiakoff, A.A.; Hoelz, A.  
Deposited on : 2021-05-01  
Resolution : 2.35 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35.1  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)

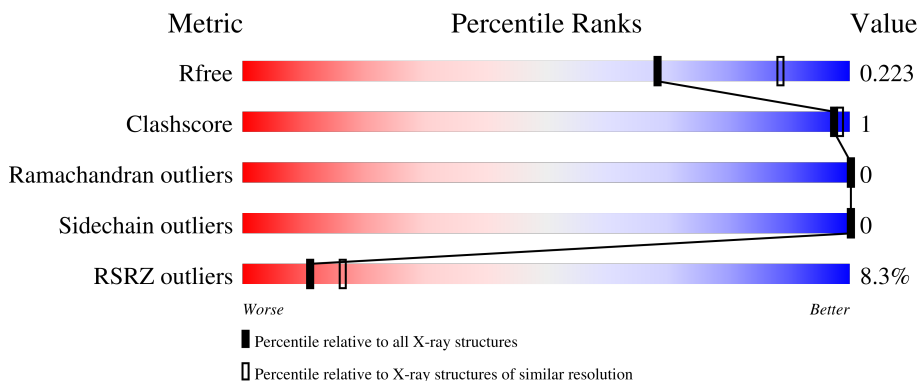
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.35 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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	216	 8% 92% 5%
1	C	216	 7% 91% 5%
1	E	216	 11% 92% 5%

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Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
 Validation Pipeline (wwPDB-VP) : 2.35.1

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Mol	Chain	Length	Quality of chain
1	G	216	<p>7% 94% 5%</p>
1	I	216	<p>8% 93% 6%</p>
1	K	216	<p>11% 93% 5%</p>
2	B	139	<p>8% 99%</p>
2	D	139	<p>6% 98% ..</p>
2	F	139	<p>5% 98% ..</p>
2	H	139	<p>6% 99% ..</p>
2	J	139	<p>6% 99% ..</p>
2	L	139	<p>10% 99% ..</p>

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 34162 atoms, of which 16802 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 GTP-binding nuclear protein Ran.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	206	3328	1073	1670	285	294	6	0	1	0
1	C	205	3311	1067	1661	284	293	6	0	2	0
1	E	206	3314	1068	1663	284	293	6	0	1	0
1	G	206	3312	1068	1661	284	293	6	0	0	0
1	I	204	3284	1058	1649	282	289	6	0	1	0
1	K	205	3314	1068	1665	284	291	6	0	1	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP P62826
C	0	SER	-	expression tag	UNP P62826
E	0	SER	-	expression tag	UNP P62826
G	0	SER	-	expression tag	UNP P62826
I	0	SER	-	expression tag	UNP P62826
K	0	SER	-	expression tag	UNP P62826

- Molecule 2 is a protein called E3 SUMO-protein ligase RanBP2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	B	138	2247	711	1120	197	214	5	0	0	0
2	D	138	2271	717	1133	201	215	5	0	1	0
2	F	138	2271	717	1133	201	215	5	0	1	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O	S			
2	H	138	2272	717	1134	201	215	5	0	2	0
2	J	138	2248	711	1121	197	214	5	0	0	0
2	L	138	2247	711	1120	197	214	5	0	0	0

There are 72 discrepancies between the modelled and reference sequences:

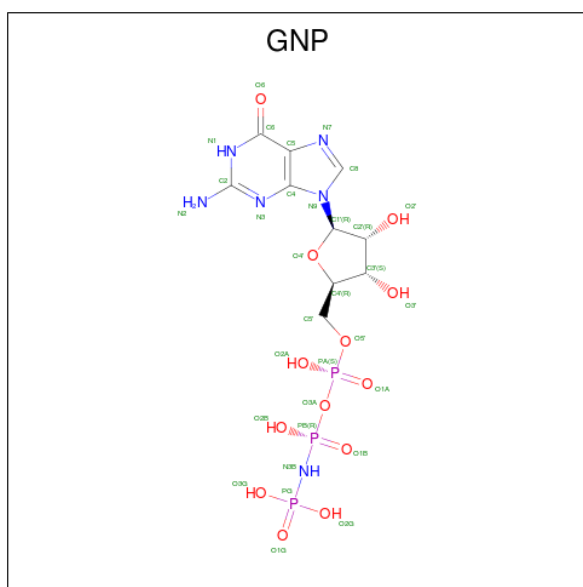
Chain	Residue	Modelled	Actual	Comment	Reference
B	2907	GLY	-	expression tag	UNP P49792
B	2908	PRO	-	expression tag	UNP P49792
B	2909	GLY	-	expression tag	UNP P49792
B	2910	SER	-	expression tag	UNP P49792
B	2962	GLN	TRP	engineered mutation	UNP P49792
B	2963	ASN	HIS	engineered mutation	UNP P49792
B	2964	TYR	THR	engineered mutation	UNP P49792
B	2965	ASP	MET	engineered mutation	UNP P49792
B	2966	ASN	LYS	engineered mutation	UNP P49792
B	2967	LYS	ASN	engineered mutation	UNP P49792
B	2968	GLN	TYR	engineered mutation	UNP P49792
B	2969	VAL	TYR	engineered mutation	UNP P49792
D	2907	GLY	-	expression tag	UNP P49792
D	2908	PRO	-	expression tag	UNP P49792
D	2909	GLY	-	expression tag	UNP P49792
D	2910	SER	-	expression tag	UNP P49792
D	2962	GLN	TRP	engineered mutation	UNP P49792
D	2963	ASN	HIS	engineered mutation	UNP P49792
D	2964	TYR	THR	engineered mutation	UNP P49792
D	2965	ASP	MET	engineered mutation	UNP P49792
D	2966	ASN	LYS	engineered mutation	UNP P49792
D	2967	LYS	ASN	engineered mutation	UNP P49792
D	2968	GLN	TYR	engineered mutation	UNP P49792
D	2969	VAL	TYR	engineered mutation	UNP P49792
F	2907	GLY	-	expression tag	UNP P49792
F	2908	PRO	-	expression tag	UNP P49792
F	2909	GLY	-	expression tag	UNP P49792
F	2910	SER	-	expression tag	UNP P49792
F	2962	GLN	TRP	engineered mutation	UNP P49792
F	2963	ASN	HIS	engineered mutation	UNP P49792
F	2964	TYR	THR	engineered mutation	UNP P49792
F	2965	ASP	MET	engineered mutation	UNP P49792

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Chain	Residue	Modelled	Actual	Comment	Reference
F	2966	ASN	LYS	engineered mutation	UNP P49792
F	2967	LYS	ASN	engineered mutation	UNP P49792
F	2968	GLN	TYR	engineered mutation	UNP P49792
F	2969	VAL	TYR	engineered mutation	UNP P49792
H	2907	GLY	-	expression tag	UNP P49792
H	2908	PRO	-	expression tag	UNP P49792
H	2909	GLY	-	expression tag	UNP P49792
H	2910	SER	-	expression tag	UNP P49792
H	2962	GLN	TRP	engineered mutation	UNP P49792
H	2963	ASN	HIS	engineered mutation	UNP P49792
H	2964	TYR	THR	engineered mutation	UNP P49792
H	2965	ASP	MET	engineered mutation	UNP P49792
H	2966	ASN	LYS	engineered mutation	UNP P49792
H	2967	LYS	ASN	engineered mutation	UNP P49792
H	2968	GLN	TYR	engineered mutation	UNP P49792
H	2969	VAL	TYR	engineered mutation	UNP P49792
J	2907	GLY	-	expression tag	UNP P49792
J	2908	PRO	-	expression tag	UNP P49792
J	2909	GLY	-	expression tag	UNP P49792
J	2910	SER	-	expression tag	UNP P49792
J	2962	GLN	TRP	engineered mutation	UNP P49792
J	2963	ASN	HIS	engineered mutation	UNP P49792
J	2964	TYR	THR	engineered mutation	UNP P49792
J	2965	ASP	MET	engineered mutation	UNP P49792
J	2966	ASN	LYS	engineered mutation	UNP P49792
J	2967	LYS	ASN	engineered mutation	UNP P49792
J	2968	GLN	TYR	engineered mutation	UNP P49792
J	2969	VAL	TYR	engineered mutation	UNP P49792
L	2907	GLY	-	expression tag	UNP P49792
L	2908	PRO	-	expression tag	UNP P49792
L	2909	GLY	-	expression tag	UNP P49792
L	2910	SER	-	expression tag	UNP P49792
L	2962	GLN	TRP	engineered mutation	UNP P49792
L	2963	ASN	HIS	engineered mutation	UNP P49792
L	2964	TYR	THR	engineered mutation	UNP P49792
L	2965	ASP	MET	engineered mutation	UNP P49792
L	2966	ASN	LYS	engineered mutation	UNP P49792
L	2967	LYS	ASN	engineered mutation	UNP P49792
L	2968	GLN	TYR	engineered mutation	UNP P49792
L	2969	VAL	TYR	engineered mutation	UNP P49792

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-GUANYLATE ESTER (three-letter code: GNP) (formula: C<sub>10</sub>H<sub>17</sub>N<sub>6</sub>O<sub>13</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	H	N	O			P
3	A	1	Total 44	C 10	H 12	N 6	O 13	P 3	0	0
3	C	1	Total 44	C 10	H 12	N 6	O 13	P 3	0	0
3	E	1	Total 44	C 10	H 12	N 6	O 13	P 3	0	0
3	G	1	Total 44	C 10	H 12	N 6	O 13	P 3	0	0
3	I	1	Total 44	C 10	H 12	N 6	O 13	P 3	0	0
3	K	1	Total 44	C 10	H 12	N 6	O 13	P 3	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
4	A	1	Total 1	Mg 1	0	0
4	C	1	Total 1	Mg 1	0	0
4	E	1	Total 1	Mg 1	0	0
4	G	1	Total 1	Mg 1	0	0
4	I	1	Total 1	Mg 1	0	0
4	K	1	Total 1	Mg 1	0	0

- Molecule 5 is water.

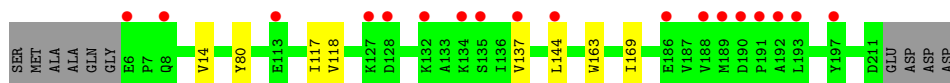
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	64	Total O 64 64	0	0
5	B	20	Total O 20 20	0	0
5	C	79	Total O 79 79	0	0
5	D	25	Total O 25 25	0	0
5	E	35	Total O 35 35	0	0
5	F	21	Total O 21 21	0	0
5	G	68	Total O 68 68	0	0
5	H	39	Total O 39 39	0	0
5	I	34	Total O 34 34	0	0
5	J	21	Total O 21 21	0	0
5	K	50	Total O 50 50	0	0
5	L	17	Total O 17 17	0	0



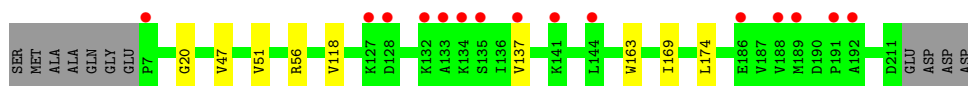
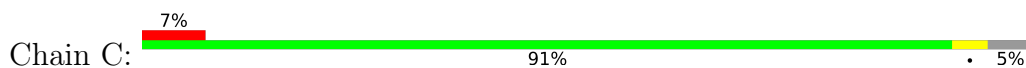
### 3 Residue-property plots [i](#)

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.

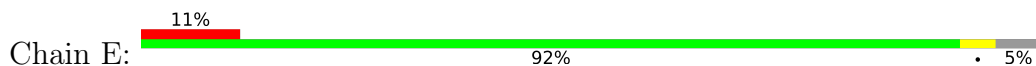
- Molecule 1: GTP-binding nuclear protein Ran



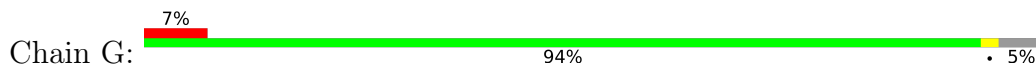
- Molecule 1: GTP-binding nuclear protein Ran



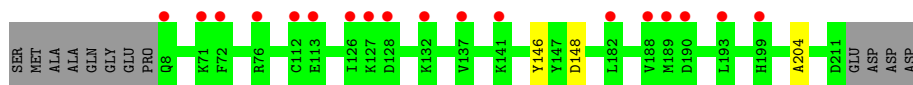
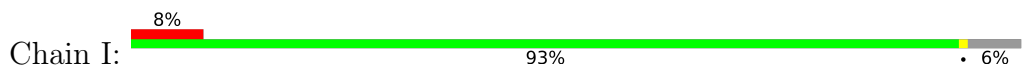
- Molecule 1: GTP-binding nuclear protein Ran



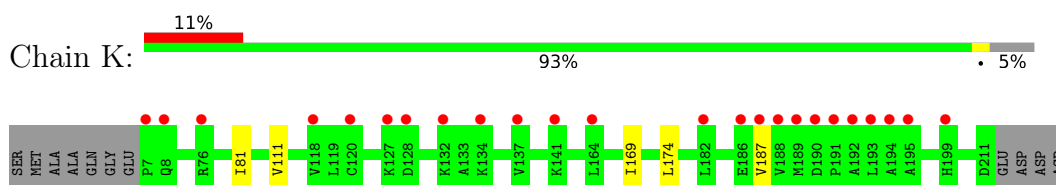
- Molecule 1: GTP-binding nuclear protein Ran



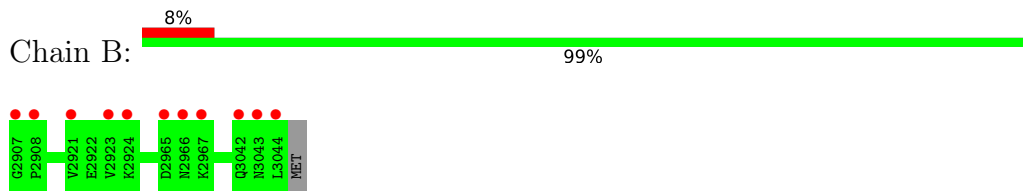
- Molecule 1: GTP-binding nuclear protein Ran



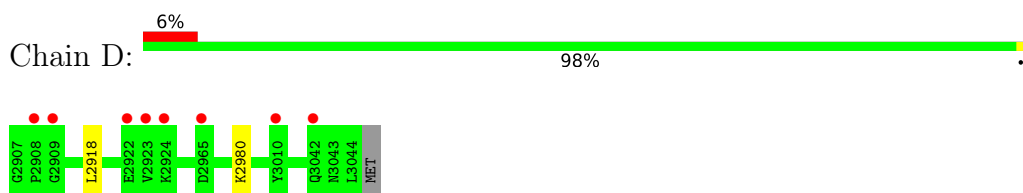
- Molecule 1: GTP-binding nuclear protein Ran



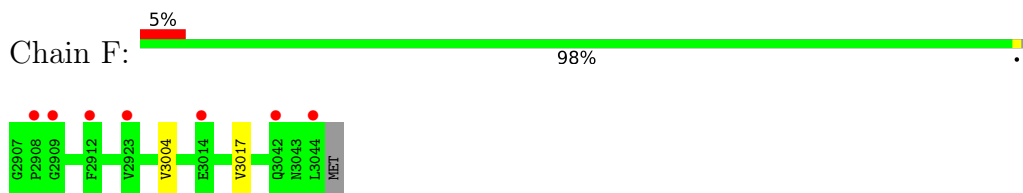
- Molecule 2: E3 SUMO-protein ligase RanBP2



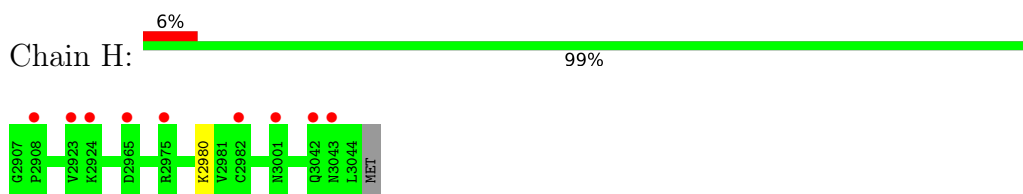
- Molecule 2: E3 SUMO-protein ligase RanBP2



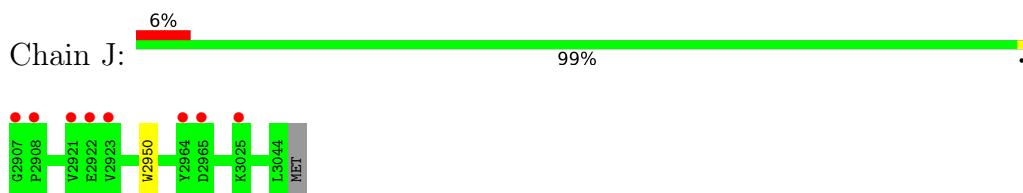
- Molecule 2: E3 SUMO-protein ligase RanBP2



- Molecule 2: E3 SUMO-protein ligase RanBP2

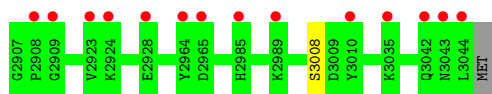


- Molecule 2: E3 SUMO-protein ligase RanBP2



- Molecule 2: E3 SUMO-protein ligase RanBP2





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	142.69Å 142.69Å 96.65Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.36 – 2.35 29.36 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.36-2.35) 99.9 (29.36-2.35)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.71 (at 2.36Å)	Xtriage
Refinement program	PHENIX 1.19.1	Depositor
R, $R_{free}$	0.183 , 0.225 0.182 , 0.223	Depositor DCC
$R_{free}$ test set	4428 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	47.5	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 39.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l 0.019 for h,-h-k,-l 0.000 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	34162	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 27.69 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.1024e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

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

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GNP, MG

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	0.25	0/1700	0.49	0/2307
1	C	0.25	0/1700	0.49	0/2306
1	E	0.25	0/1702	0.47	0/2308
1	G	0.25	0/1693	0.49	0/2297
1	I	0.25	0/1685	0.47	0/2284
1	K	0.25	0/1691	0.48	0/2294
2	B	0.24	0/1147	0.48	0/1544
2	D	0.25	0/1158	0.49	0/1558
2	F	0.25	0/1158	0.49	0/1558
2	H	0.24	0/1166	0.49	0/1569
2	J	0.24	0/1147	0.47	0/1544
2	L	0.24	0/1147	0.48	0/1544
All	All	0.25	0/17094	0.48	0/23113

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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	1658	1670	1669	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1650	1661	1649	8	0
1	E	1651	1663	1648	7	0
1	G	1651	1661	1661	3	0
1	I	1635	1649	1635	2	0
1	K	1649	1665	1664	4	0
2	B	1127	1120	1120	0	0
2	D	1138	1133	1132	2	0
2	F	1138	1133	1132	3	0
2	H	1138	1134	1129	1	0
2	J	1127	1121	1120	1	0
2	L	1127	1120	1120	1	0
3	A	32	12	13	0	0
3	C	32	12	13	1	0
3	E	32	12	13	0	0
3	G	32	12	13	1	0
3	I	32	12	13	0	0
3	K	32	12	13	0	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
4	E	1	0	0	0	0
4	G	1	0	0	0	0
4	I	1	0	0	0	0
4	K	1	0	0	0	0
5	A	64	0	0	0	0
5	B	20	0	0	0	0
5	C	79	0	0	0	0
5	D	25	0	0	0	0
5	E	35	0	0	0	0
5	F	21	0	0	0	0
5	G	68	0	0	0	0
5	H	39	0	0	0	0
5	I	34	0	0	0	0
5	J	21	0	0	0	0
5	K	50	0	0	0	0
5	L	17	0	0	0	0
All	All	17360	16802	16757	28	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 1.

The worst 5 of 28 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:E:31:LEU:HD11	1:E:48:HIS:HB3	1.78	0.65
1:E:201:LEU:HD11	2:F:3004:VAL:HG12	1.84	0.57
1:C:169:ILE:HD11	1:C:174:LEU:HD22	1.89	0.55
1:I:146:TYR:OH	1:I:148:ASP:OD1	2.19	0.54
1:K:169:ILE:HD11	1:K:174:LEU:HD13	1.90	0.54

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	205/216 (95%)	201 (98%)	4 (2%)	0	100	100
1	C	205/216 (95%)	199 (97%)	6 (3%)	0	100	100
1	E	205/216 (95%)	198 (97%)	7 (3%)	0	100	100
1	G	204/216 (94%)	198 (97%)	6 (3%)	0	100	100
1	I	203/216 (94%)	196 (97%)	7 (3%)	0	100	100
1	K	204/216 (94%)	197 (97%)	7 (3%)	0	100	100
2	B	136/139 (98%)	133 (98%)	3 (2%)	0	100	100
2	D	137/139 (99%)	134 (98%)	3 (2%)	0	100	100
2	F	137/139 (99%)	135 (98%)	2 (2%)	0	100	100
2	H	138/139 (99%)	136 (99%)	2 (1%)	0	100	100
2	J	136/139 (98%)	133 (98%)	3 (2%)	0	100	100
2	L	136/139 (98%)	134 (98%)	2 (2%)	0	100	100
All	All	2046/2130 (96%)	1994 (98%)	52 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	179/185 (97%)	179 (100%)	0	100	100
1	C	179/185 (97%)	179 (100%)	0	100	100
1	E	179/185 (97%)	179 (100%)	0	100	100
1	G	178/185 (96%)	178 (100%)	0	100	100
1	I	177/185 (96%)	177 (100%)	0	100	100
1	K	178/185 (96%)	178 (100%)	0	100	100
2	B	124/125 (99%)	124 (100%)	0	100	100
2	D	125/125 (100%)	125 (100%)	0	100	100
2	F	125/125 (100%)	125 (100%)	0	100	100
2	H	126/125 (101%)	126 (100%)	0	100	100
2	J	124/125 (99%)	124 (100%)	0	100	100
2	L	124/125 (99%)	124 (100%)	0	100	100
All	All	1818/1860 (98%)	1818 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.



## 5.6 Ligand geometry

Of 12 ligands modelled in this entry, 6 are monoatomic - leaving 6 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
3	GNP	G	301	4	29,34,34	1.42	4 (13%)	33,54,54	2.37	7 (21%)
3	GNP	C	301	4	29,34,34	1.42	4 (13%)	33,54,54	2.36	7 (21%)
3	GNP	A	301	4	29,34,34	1.40	4 (13%)	33,54,54	2.34	7 (21%)
3	GNP	I	301	4	29,34,34	1.43	4 (13%)	33,54,54	2.39	7 (21%)
3	GNP	E	301	4	29,34,34	1.44	4 (13%)	33,54,54	2.38	7 (21%)
3	GNP	K	301	4	29,34,34	1.42	4 (13%)	33,54,54	2.40	8 (24%)

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
3	GNP	G	301	4	-	3/14/38/38	0/3/3/3
3	GNP	C	301	4	-	3/14/38/38	0/3/3/3
3	GNP	A	301	4	-	6/14/38/38	0/3/3/3
3	GNP	I	301	4	-	7/14/38/38	0/3/3/3
3	GNP	E	301	4	-	4/14/38/38	0/3/3/3
3	GNP	K	301	4	-	3/14/38/38	0/3/3/3

The worst 5 of 24 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	301	GNP	C6-N1	4.08	1.40	1.33
3	K	301	GNP	C6-N1	4.05	1.40	1.33
3	C	301	GNP	C6-N1	4.03	1.40	1.33
3	G	301	GNP	C6-N1	4.02	1.40	1.33
3	A	301	GNP	C6-N1	4.00	1.40	1.33

The worst 5 of 43 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	301	GNP	C5-C6-N1	-9.03	111.08	123.43
3	K	301	GNP	C5-C6-N1	-9.03	111.08	123.43
3	C	301	GNP	C5-C6-N1	-8.97	111.16	123.43
3	A	301	GNP	C5-C6-N1	-8.96	111.18	123.43
3	E	301	GNP	C5-C6-N1	-8.94	111.21	123.43

There are no chirality outliers.

5 of 26 torsion outliers are listed below:

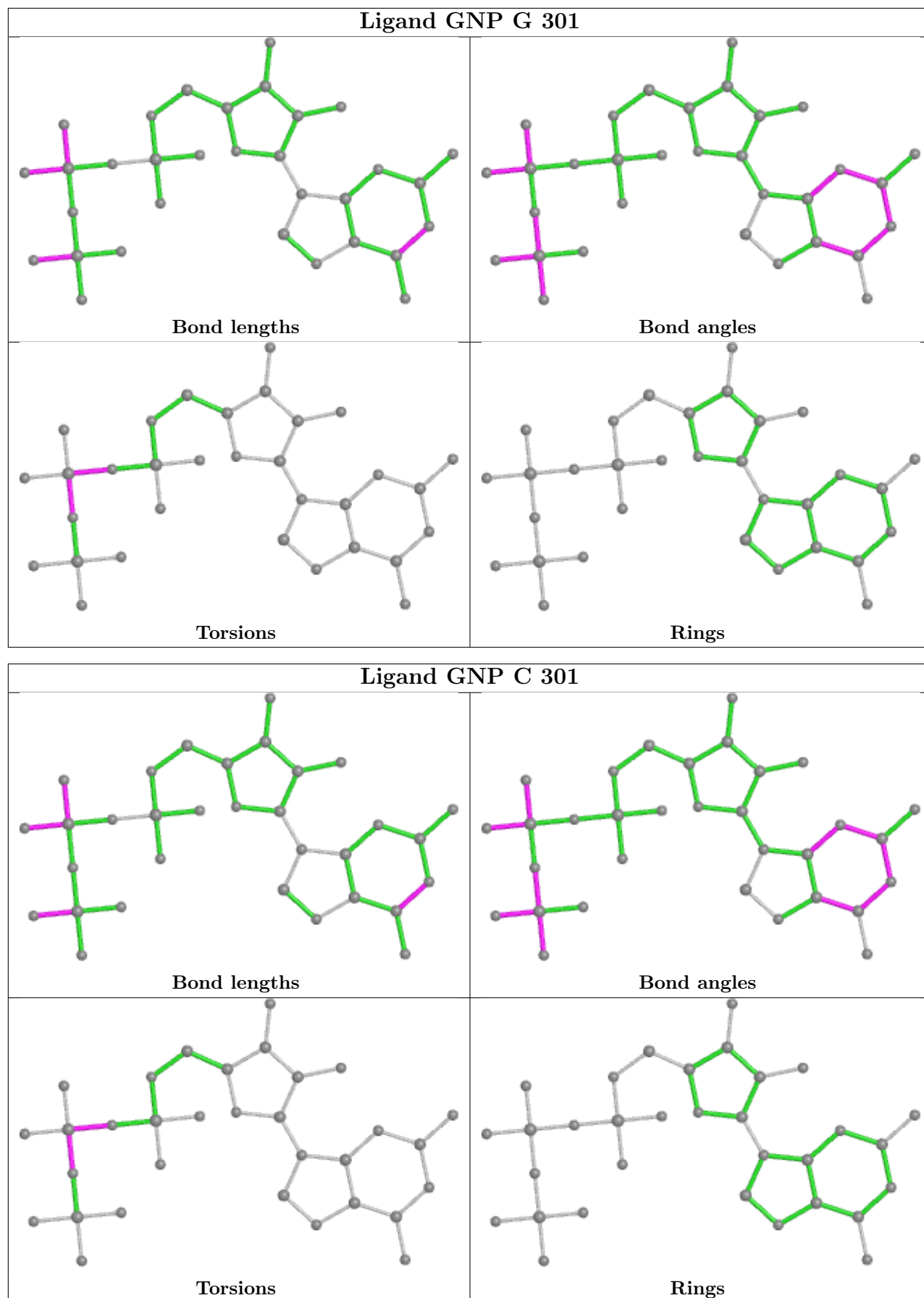
Mol	Chain	Res	Type	Atoms
3	A	301	GNP	PG-N3B-PB-O1B
3	A	301	GNP	PA-O3A-PB-O1B
3	A	301	GNP	PA-O3A-PB-O2B
3	A	301	GNP	C5'-O5'-PA-O3A
3	C	301	GNP	PG-N3B-PB-O1B

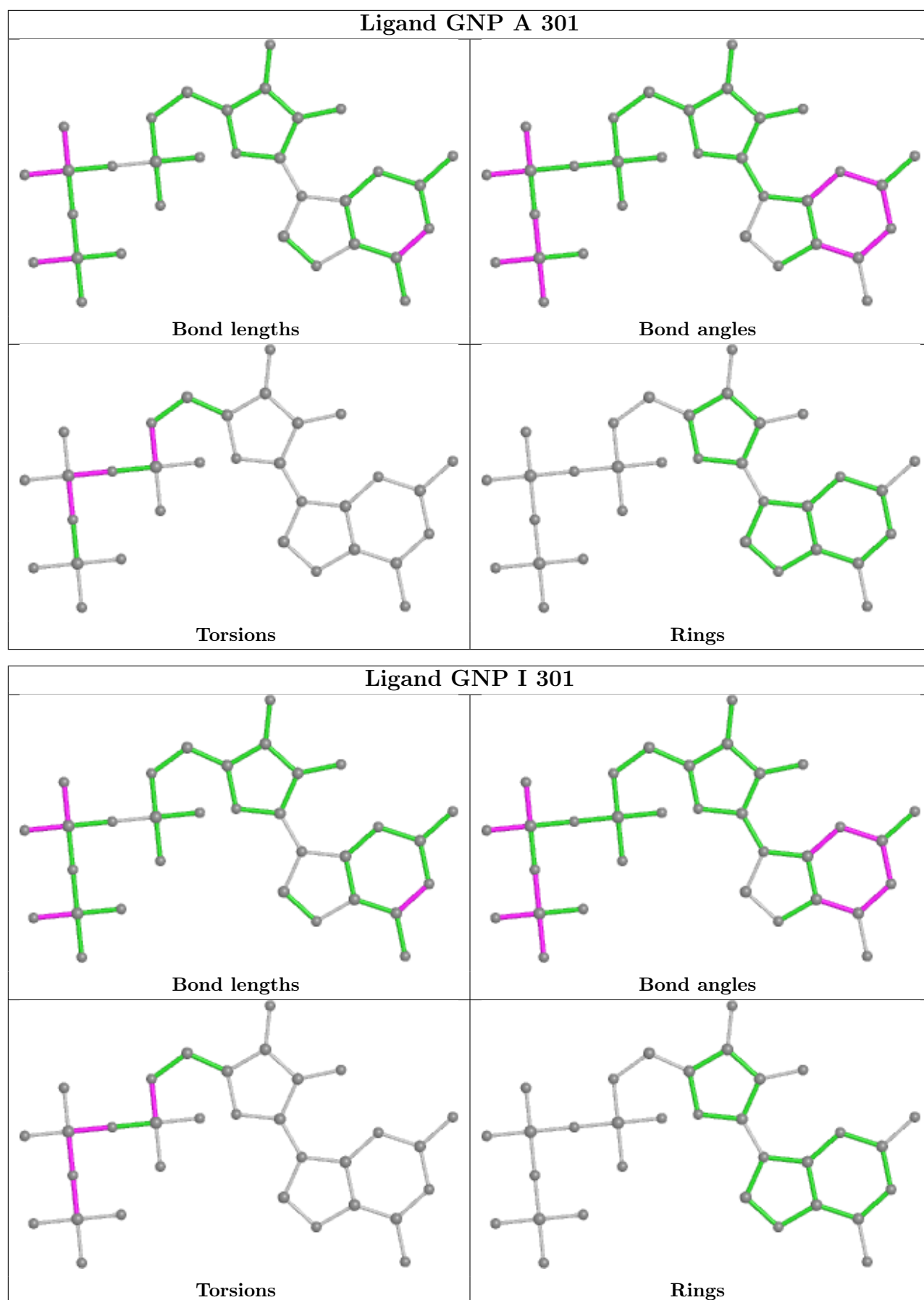
There are no ring outliers.

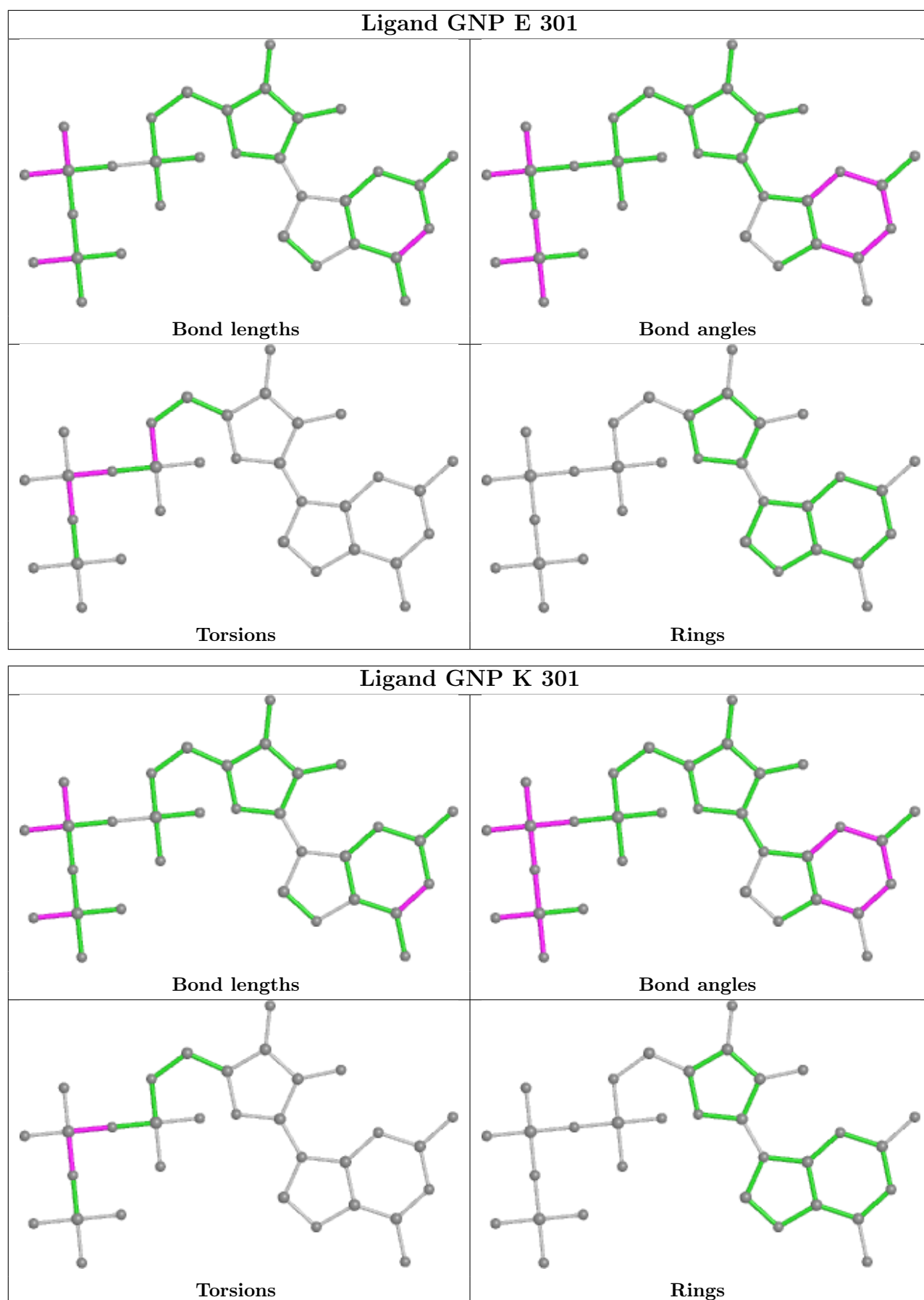
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	301	GNP	1	0
3	C	301	GNP	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	206/216 (95%)	0.41	18 (8%) 10 15	35, 50, 101, 126	0
1	C	205/216 (94%)	0.39	15 (7%) 15 22	31, 46, 98, 121	0
1	E	206/216 (95%)	0.66	23 (11%) 5 8	41, 58, 110, 134	0
1	G	206/216 (95%)	0.33	15 (7%) 15 22	32, 46, 95, 137	0
1	I	204/216 (94%)	0.52	18 (8%) 10 15	44, 58, 93, 124	0
1	K	205/216 (94%)	0.56	24 (11%) 4 7	40, 54, 102, 134	0
2	B	138/139 (99%)	0.48	11 (7%) 12 18	41, 63, 93, 119	0
2	D	138/139 (99%)	0.36	8 (5%) 23 33	36, 55, 85, 115	0
2	F	138/139 (99%)	0.27	7 (5%) 28 40	43, 59, 83, 93	0
2	H	138/139 (99%)	0.38	9 (6%) 18 27	35, 54, 81, 99	0
2	J	138/139 (99%)	0.35	8 (5%) 23 33	41, 63, 90, 122	0
2	L	138/139 (99%)	0.49	14 (10%) 7 11	44, 62, 89, 104	0
All	All	2060/2130 (96%)	0.44	170 (8%) 11 16	31, 56, 98, 137	0

The worst 5 of 170 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	189	MET	7.3
1	E	197	TYR	6.2
1	K	192	ALA	6.0
1	I	128	ASP	6.0
1	K	191	PRO	5.9

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

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

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

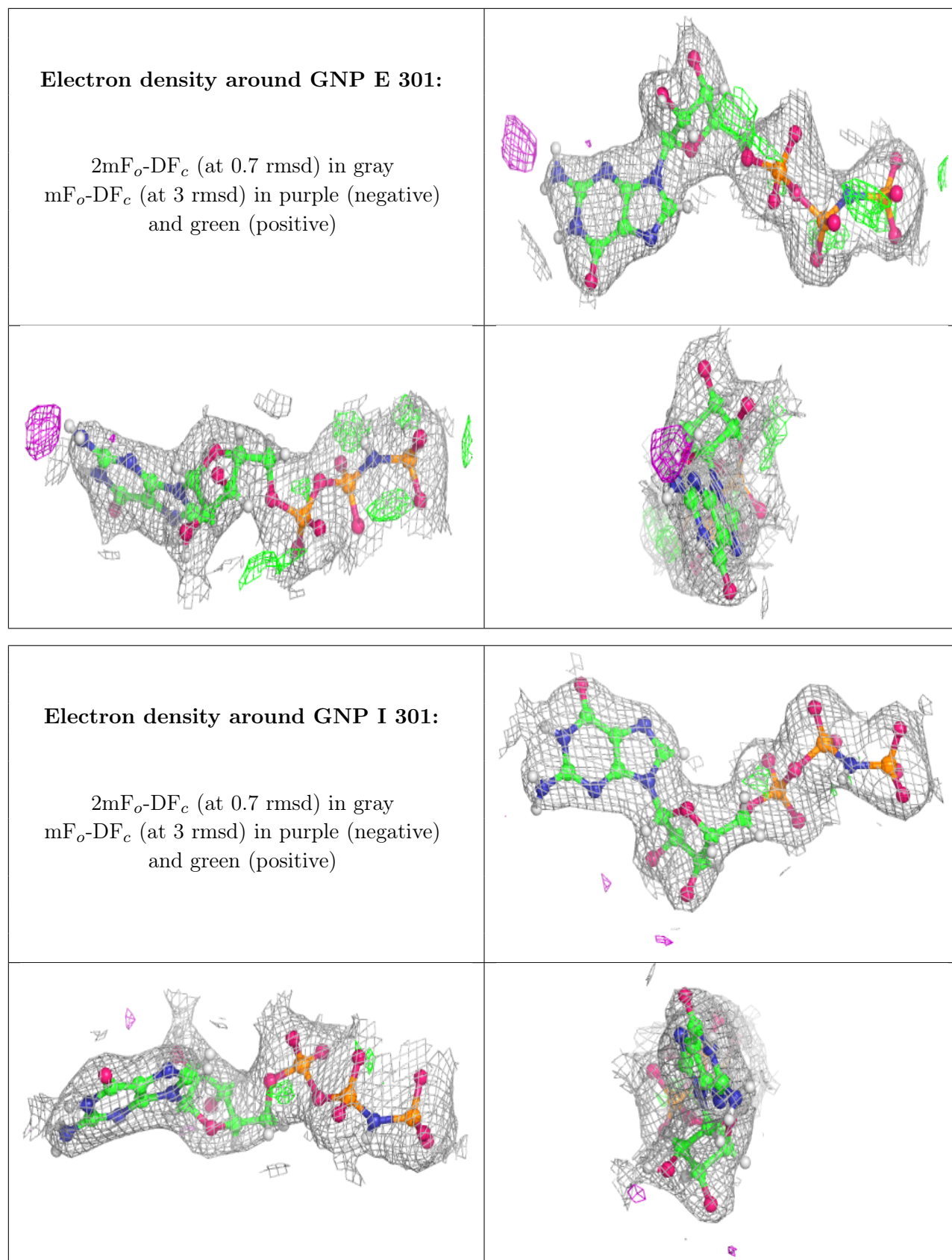
### 6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	MG	G	302	1/1	0.89	0.14	33,33,33,33	0
4	MG	E	302	1/1	0.91	0.13	49,49,49,49	0
4	MG	A	302	1/1	0.91	0.14	33,33,33,33	0
4	MG	C	302	1/1	0.96	0.20	36,36,36,36	0
4	MG	K	302	1/1	0.96	0.12	44,44,44,44	0
4	MG	I	302	1/1	0.97	0.09	41,41,41,41	0
3	GNP	E	301	32/32	0.97	0.15	39,49,62,65	0
3	GNP	I	301	32/32	0.98	0.12	37,48,58,60	0
3	GNP	K	301	32/32	0.98	0.14	34,47,60,67	0
3	GNP	C	301	32/32	0.98	0.15	30,39,50,52	0
3	GNP	A	301	32/32	0.98	0.17	31,42,49,60	0
3	GNP	G	301	32/32	0.99	0.13	28,36,48,48	0

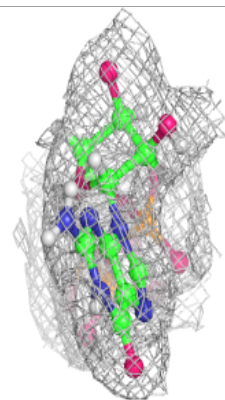
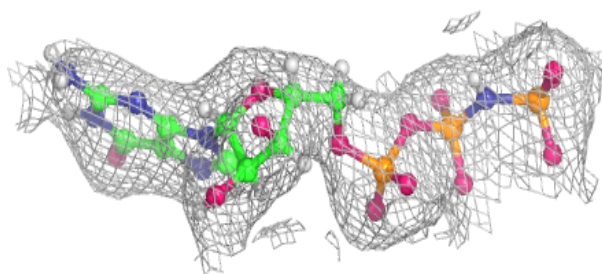
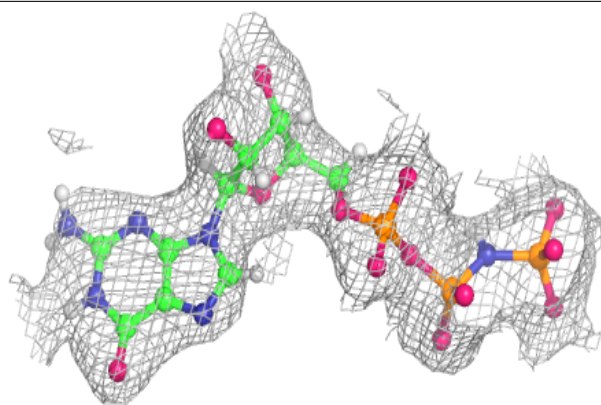
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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



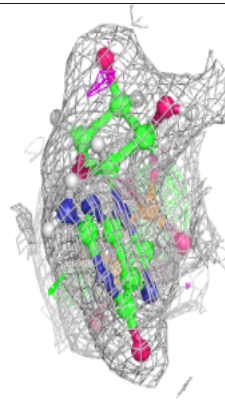
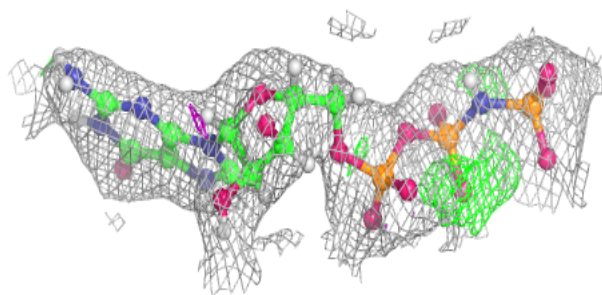
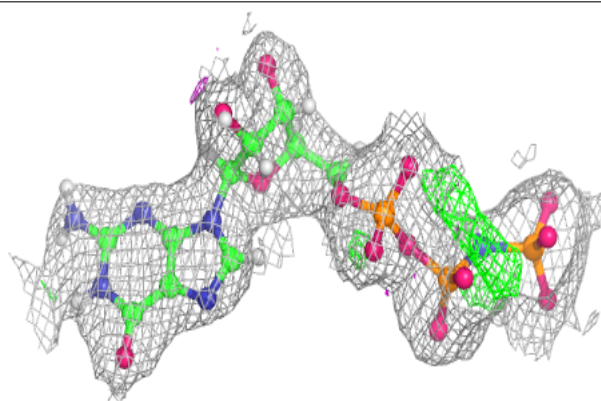


**Electron density around GNP K 301:**

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

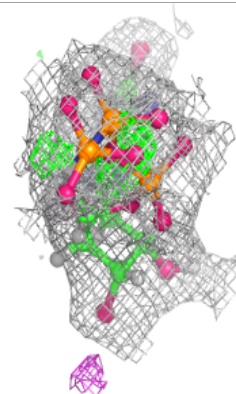
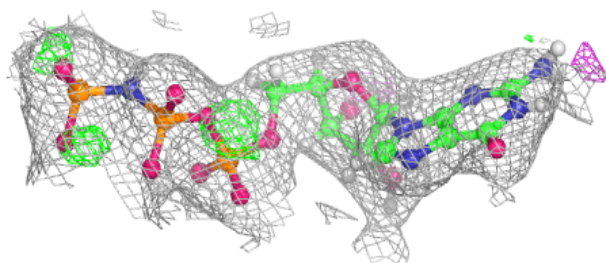
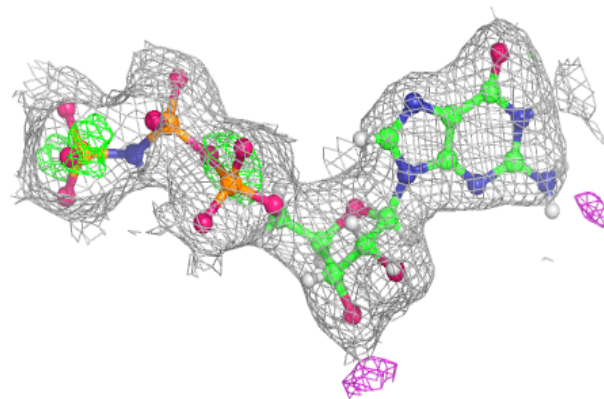
**Electron density around GNP C 301:**

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

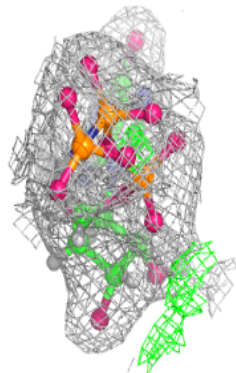
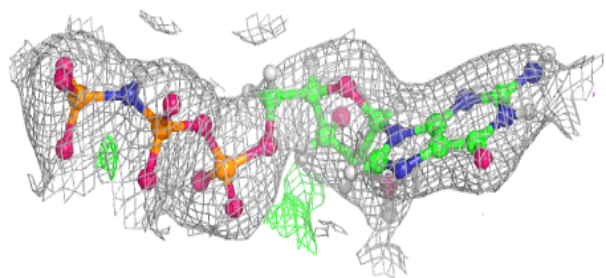
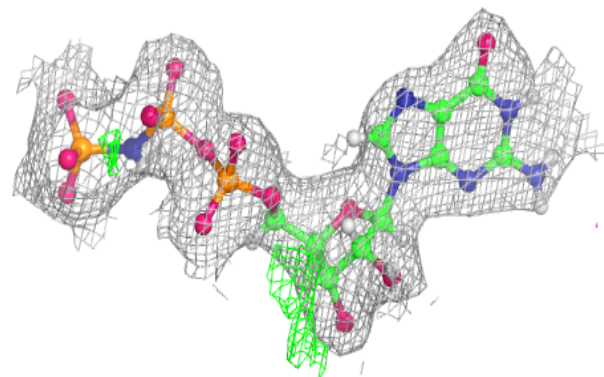


**Electron density around GNP A 301:**

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

**Electron density around GNP G 301:**

$2mF_o-DF_c$  (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.