



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 27, 2024 – 06:34 pm BST

PDB ID : 5L3S
Title : Structure of the GTPase heterodimer of crenarchaeal SRP54 and FtsY
Authors : Bange, G.; Wild, K.; Sinning, I.
Deposited on : 2016-05-24
Resolution : 1.90 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:

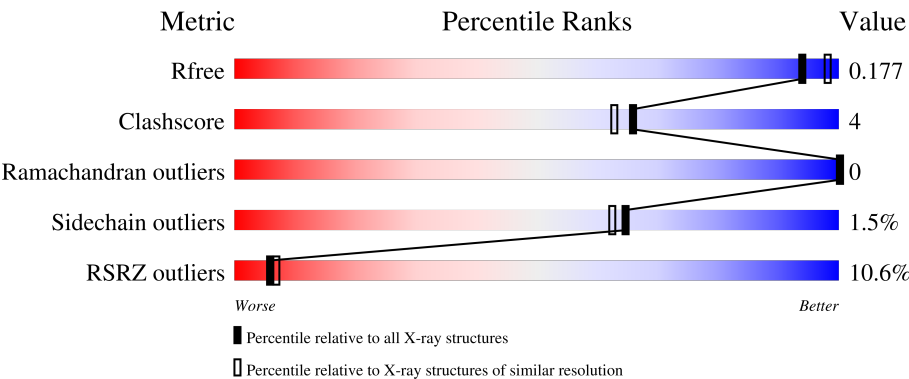
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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.



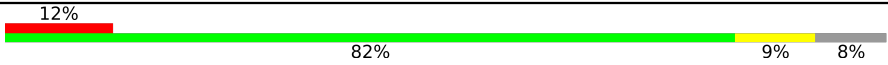


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	298	<div><div>7%</div><div>83%9%8%</div></div>
1	C	298	<div><div>19%</div><div>78%12%9%</div></div>
1	E	298	<div><div>11%</div><div>83%7%9%</div></div>
1	G	298	<div><div>5%</div><div>83%10%7%</div></div>
2	B	296	<div><div>6%</div><div>87%9%. </div></div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	D	296	
2	F	296	
2	H	296	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	G	A	303	-	-	-	X
6	GOL	H	404	-	-	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 19033 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 Signal recognition particle 54 kDa protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	0	2	0
			2142	1380	353	404	5			
1	C	270	Total	C	N	O	S	0	2	0
			2110	1360	346	399	5			
1	E	271	Total	C	N	O	S	0	1	0
			2115	1362	349	399	5			
1	G	277	Total	C	N	O	S	0	0	0
			2150	1383	355	407	5			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	expression tag	UNP Q97ZE7
A	-4	HIS	-	expression tag	UNP Q97ZE7
A	-3	HIS	-	expression tag	UNP Q97ZE7
A	-2	HIS	-	expression tag	UNP Q97ZE7
A	-1	HIS	-	expression tag	UNP Q97ZE7
A	0	HIS	-	expression tag	UNP Q97ZE7
C	-5	HIS	-	expression tag	UNP Q97ZE7
C	-4	HIS	-	expression tag	UNP Q97ZE7
C	-3	HIS	-	expression tag	UNP Q97ZE7
C	-2	HIS	-	expression tag	UNP Q97ZE7
C	-1	HIS	-	expression tag	UNP Q97ZE7
C	0	HIS	-	expression tag	UNP Q97ZE7
E	-5	HIS	-	expression tag	UNP Q97ZE7
E	-4	HIS	-	expression tag	UNP Q97ZE7
E	-3	HIS	-	expression tag	UNP Q97ZE7
E	-2	HIS	-	expression tag	UNP Q97ZE7
E	-1	HIS	-	expression tag	UNP Q97ZE7
E	0	HIS	-	expression tag	UNP Q97ZE7
G	-5	HIS	-	expression tag	UNP Q97ZE7
G	-4	HIS	-	expression tag	UNP Q97ZE7
G	-3	HIS	-	expression tag	UNP Q97ZE7

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
G	-2	HIS	-	expression tag	UNP Q97ZE7
G	-1	HIS	-	expression tag	UNP Q97ZE7
G	0	HIS	-	expression tag	UNP Q97ZE7

- Molecule 2 is a protein called Signal recognition particle receptor FtsY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	283	Total	C	N	O	S	0	1	0
			2235	1429	374	429	3			
2	D	271	Total	C	N	O	S	0	1	0
			2128	1361	352	412	3			
2	F	283	Total	C	N	O	S	0	0	0
			2228	1426	371	428	3			
2	H	278	Total	C	N	O	S	0	2	0
			2198	1406	367	422	3			

There are 28 discrepancies between the modelled and reference sequences:

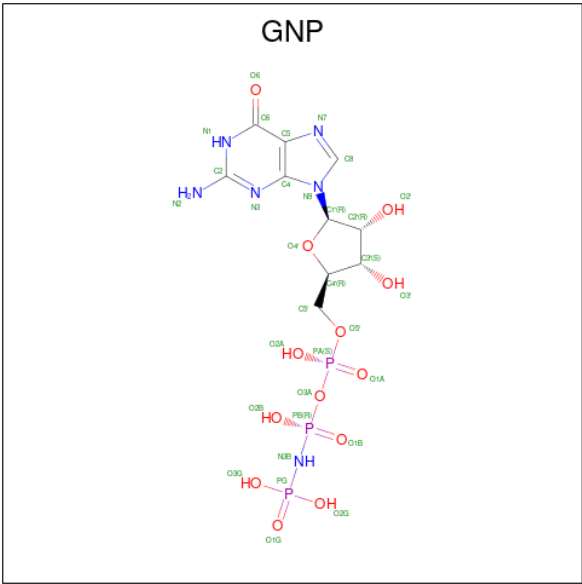
Chain	Residue	Modelled	Actual	Comment	Reference
B	73	HIS	-	expression tag	UNP P27414
B	74	HIS	-	expression tag	UNP P27414
B	75	HIS	-	expression tag	UNP P27414
B	76	HIS	-	expression tag	UNP P27414
B	77	HIS	-	expression tag	UNP P27414
B	78	HIS	-	expression tag	UNP P27414
B	279	LYS	ARG	conflict	UNP P27414
D	73	HIS	-	expression tag	UNP P27414
D	74	HIS	-	expression tag	UNP P27414
D	75	HIS	-	expression tag	UNP P27414
D	76	HIS	-	expression tag	UNP P27414
D	77	HIS	-	expression tag	UNP P27414
D	78	HIS	-	expression tag	UNP P27414
D	279	LYS	ARG	conflict	UNP P27414
F	73	HIS	-	expression tag	UNP P27414
F	74	HIS	-	expression tag	UNP P27414
F	75	HIS	-	expression tag	UNP P27414
F	76	HIS	-	expression tag	UNP P27414
F	77	HIS	-	expression tag	UNP P27414
F	78	HIS	-	expression tag	UNP P27414
F	279	LYS	ARG	conflict	UNP P27414
H	73	HIS	-	expression tag	UNP P27414
H	74	HIS	-	expression tag	UNP P27414

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
H	75	HIS	-	expression tag	UNP P27414
H	76	HIS	-	expression tag	UNP P27414
H	77	HIS	-	expression tag	UNP P27414
H	78	HIS	-	expression tag	UNP P27414
H	279	LYS	ARG	conflict	UNP P27414

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-GUANYLATE ESTER (three-letter code: GNP) (formula: C₁₀H₁₇N₆O₁₃P₃).

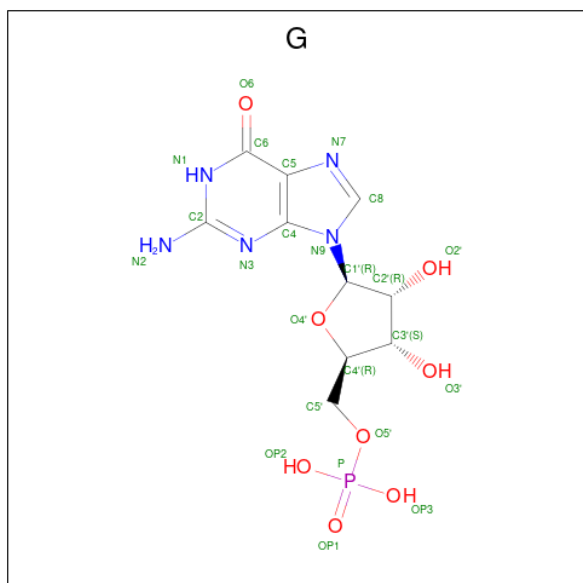


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
3	B	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
3	C	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
3	D	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
3	E	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
3	F	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
3	G	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
3	H	1	Total	C	N	O	P	0	0
			32	10	6	13	3		

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		
4	B	1	Total	Mg	0	0
			1	1		
4	C	1	Total	Mg	0	0
			1	1		
4	D	1	Total	Mg	0	0
			1	1		
4	E	1	Total	Mg	0	0
			1	1		
4	F	1	Total	Mg	0	0
			1	1		
4	G	1	Total	Mg	0	0
			1	1		
4	H	1	Total	Mg	0	0
			1	1		

- Molecule 5 is GUANOSINE-5'-MONOPHOSPHATE (three-letter code: G) (formula: C₁₀H₁₄N₅O₈P).



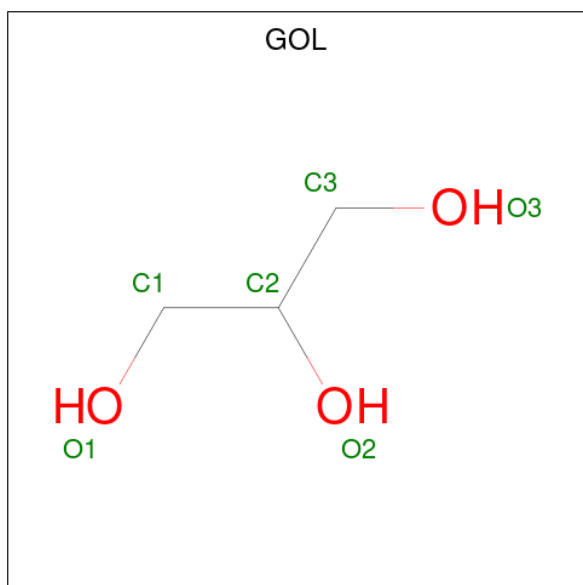
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			24	10	5	8	1		
5	C	1	Total	C	N	O	P	0	0
			24	10	5	8	1		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	E	1	Total	C	N	O	P	0	0
			24	10	5	8	1		
5	G	1	Total	C	N	O	P	0	0
			24	10	5	8	1		

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



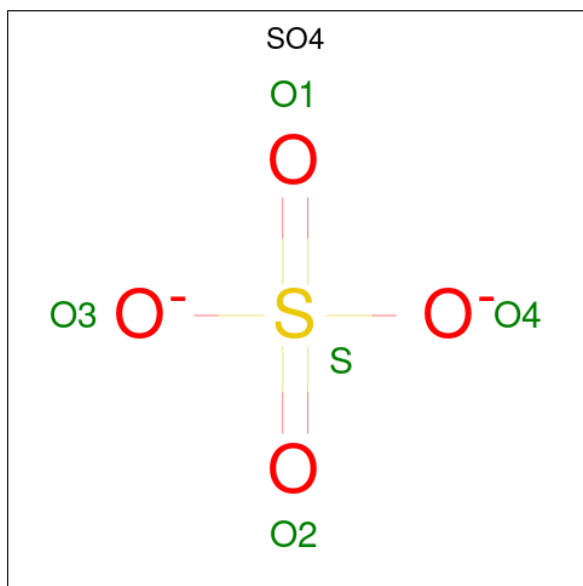
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	B	1	Total	C	O	0	0
			6	3	3		
6	B	1	Total	C	O	0	0
			6	3	3		
6	B	1	Total	C	O	0	0
			6	3	3		
6	C	1	Total	C	O	0	0
			6	3	3		
6	C	1	Total	C	O	0	0
			6	3	3		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	D	1	Total	C	O	0	0
			6	3	3		
6	E	1	Total	C	O	0	0
			6	3	3		
6	E	1	Total	C	O	0	0
			6	3	3		
6	F	1	Total	C	O	0	0
			6	3	3		
6	G	1	Total	C	O	0	0
			6	3	3		
6	G	1	Total	C	O	0	0
			6	3	3		
6	G	1	Total	C	O	0	0
			6	3	3		
6	H	1	Total	C	O	0	0
			6	3	3		
6	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	O	S	0	0
			5	4	1		
7	A	1	Total	O	S	0	0
			5	4	1		

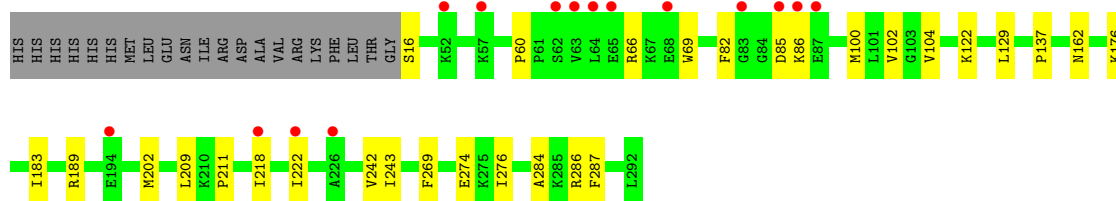
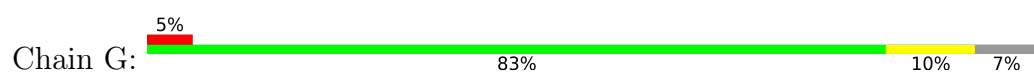
Continued on next page...

Continued from previous page...

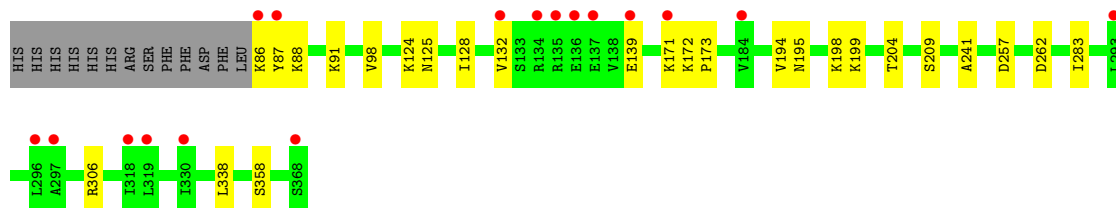
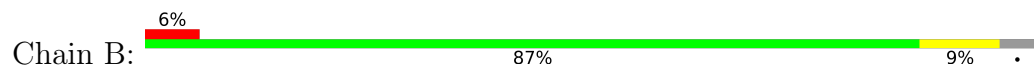
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	O	S	0	0
			5	4	1		
7	B	1	Total	O	S	0	0
			5	4	1		
7	B	1	Total	O	S	0	0
			5	4	1		
7	C	1	Total	O	S	0	0
			5	4	1		
7	E	1	Total	O	S	0	0
			5	4	1		
7	G	1	Total	O	S	0	0
			5	4	1		
7	H	1	Total	O	S	0	0
			5	4	1		

- Molecule 8 is water.

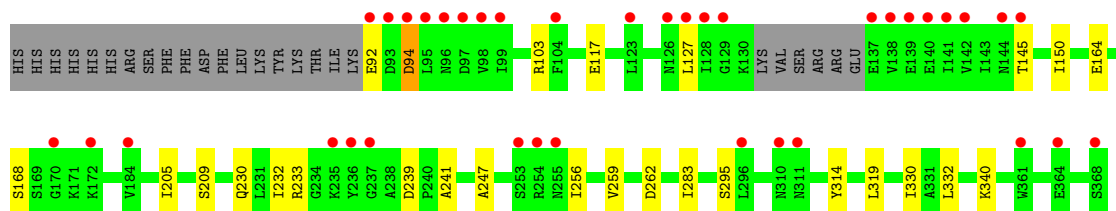
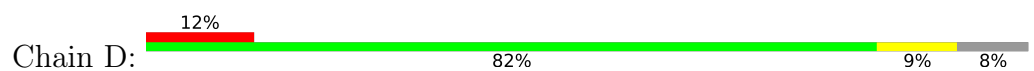
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	193	Total	O	0	0
			193	193		
8	B	190	Total	O	0	0
			190	190		
8	C	128	Total	O	0	0
			128	128		
8	D	135	Total	O	0	0
			135	135		
8	E	141	Total	O	0	0
			141	141		
8	F	106	Total	O	0	0
			106	106		
8	G	168	Total	O	0	0
			168	168		
8	H	153	Total	O	0	0
			153	153		



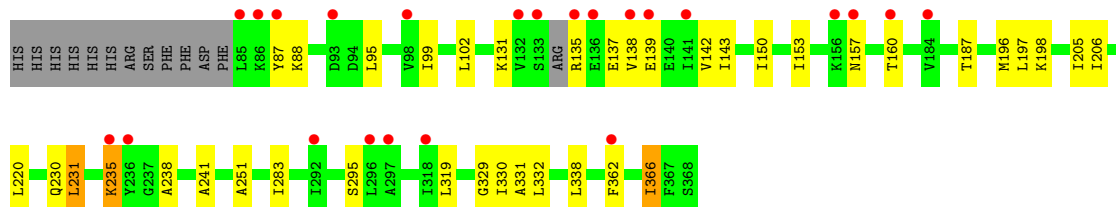
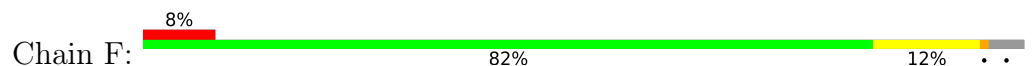
- Molecule 2: Signal recognition particle receptor FtsY



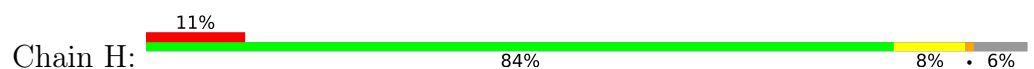
- Molecule 2: Signal recognition particle receptor FtsY



- Molecule 2: Signal recognition particle receptor FtsY



- Molecule 2: Signal recognition particle receptor FtsY





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.20Å 199.67Å 124.87Å 90.00° 92.95° 90.00°	Depositor
Resolution (Å)	37.87 – 1.90 37.87 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.3 (37.87-1.90) 89.8 (37.87-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.24 (at 1.89Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, R_{free}	0.173 , 0.216 0.175 , 0.177	Depositor DCC
R_{free} test set	3836 reflections (2.00%)	wwPDB-VP
Wilson B-factor (Å ²)	22.0	Xtriage
Anisotropy	0.691	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 62.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.068 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	19033	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.08% 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: GNP, MG, SO4, GOL

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	0.37	0/2181	0.52	0/2934
1	C	0.33	0/2150	0.47	0/2894
1	E	0.34	0/2152	0.48	0/2896
1	G	0.35	0/2186	0.49	0/2946
2	B	0.37	0/2268	0.51	0/3059
2	D	0.33	0/2159	0.48	0/2915
2	F	0.32	0/2257	0.49	0/3043
2	H	0.37	0/2233	0.50	0/3010
All	All	0.35	0/17586	0.49	0/23697

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	2142	0	2217	17	0
1	C	2110	0	2180	22	0
1	E	2115	0	2185	14	0
1	G	2150	0	2222	20	0
2	B	2235	0	2310	17	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	2128	0	2184	15	0
2	F	2228	0	2300	21	0
2	H	2198	0	2268	18	0
3	A	32	0	13	0	0
3	B	32	0	13	0	0
3	C	32	0	13	0	0
3	D	32	0	13	0	0
3	E	32	0	13	0	0
3	F	32	0	13	0	0
3	G	32	0	13	0	0
3	H	32	0	13	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	0
4	H	1	0	0	0	0
5	A	24	0	12	0	0
5	C	24	0	12	1	0
5	E	24	0	12	0	0
5	G	24	0	12	1	0
6	A	18	0	24	3	0
6	B	24	0	32	1	0
6	C	12	0	16	3	0
6	D	6	0	8	0	0
6	E	12	0	16	0	0
6	F	6	0	8	0	0
6	G	18	0	24	2	0
6	H	12	0	16	1	0
7	A	15	0	0	0	0
7	B	10	0	0	0	0
7	C	5	0	0	1	0
7	E	5	0	0	0	0
7	G	5	0	0	0	0
7	H	5	0	0	0	0
8	A	193	0	0	4	0
8	B	190	0	0	5	3
8	C	128	0	0	3	0
8	D	135	0	0	2	0
8	E	141	0	0	2	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	F	106	0	0	2	3
8	G	168	0	0	6	0
8	H	153	0	0	6	0
All	All	19033	0	18162	145	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 4.

All (145) 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:202:MET:SD	8:A:567:HOH:O	2.33	0.86
1:C:153[B]:GLN:NE2	1:E:153:GLN:OE1	2.18	0.76
2:B:91:LYS:NZ	8:B:502:HOH:O	2.18	0.75
2:F:88:LYS:HG3	2:F:135:ARG:HB2	1.70	0.73
1:A:286:ARG:NH1	6:A:305:GOL:O2	2.23	0.72
2:H:114:GLU:OE2	8:H:501:HOH:O	2.09	0.69
1:G:162:ASN:OD1	8:G:401:HOH:O	2.11	0.68
1:G:102:VAL:HB	1:G:202:MET:HE2	1.74	0.68
1:A:159[A]:ASN:OD1	8:A:401:HOH:O	2.14	0.65
2:B:125:ASN:OD1	8:B:501:HOH:O	2.13	0.65
2:F:196:MET:SD	8:F:603:HOH:O	2.54	0.65
2:D:103:ARG:NH1	2:D:117:GLU:OE1	2.29	0.64
2:D:205:ILE:HG22	2:D:230:GLN:HB3	1.80	0.64
2:H:123:LEU:HD22	2:H:146:LEU:HG	1.80	0.63
1:G:176:LYS:NZ	8:G:406:HOH:O	2.33	0.61
6:C:304:GOL:O1	8:C:401:HOH:O	2.16	0.61
2:F:205:ILE:HG22	2:F:230:GLN:HB3	1.82	0.61
2:F:329:GLY:HA2	2:F:366:ILE:HD11	1.82	0.61
2:H:102:LEU:HG	2:H:146:LEU:HD12	1.83	0.61
2:F:153:ILE:O	2:F:157:ASN:ND2	2.32	0.61
2:F:187:THR:HG23	2:F:220:LEU:HD13	1.84	0.60
1:A:24:ASP:OD1	1:A:52:LYS:NZ	2.34	0.60
1:E:100:MET:HE3	1:E:211:PRO:HG3	1.84	0.60
2:H:187:THR:HG23	2:H:220:LEU:HD13	1.84	0.59
1:C:232:ARG:NH2	7:C:306:SO4:O4	2.35	0.59
1:G:286:ARG:NH1	6:G:306:GOL:O2	2.36	0.58
1:C:167:ALA:HB1	1:C:209:LEU:HD11	1.86	0.58
2:H:158:GLN:NE2	8:H:504:HOH:O	2.35	0.57
2:F:198:LYS:NZ	8:F:502:HOH:O	2.36	0.57
2:F:241:ALA:HB1	2:F:283:ILE:HD13	1.85	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:218:ILE:HD11	1:C:222:ILE:HG21	1.87	0.56
2:B:306:ARG:NE	8:B:505:HOH:O	2.26	0.56
1:C:24:ASP:OD1	1:C:52:LYS:NZ	2.31	0.56
2:H:295:SER:HB3	2:H:319:LEU:HD12	1.87	0.56
2:D:319:LEU:HD21	2:D:330:ILE:HG21	1.87	0.55
1:C:51:ILE:HD13	1:C:74:VAL:HG22	1.89	0.55
2:B:124:LYS:O	2:B:128:ILE:HG12	2.07	0.55
2:D:314:TYR:CE2	2:D:340:LYS:HD3	2.42	0.54
1:C:16:SER:HB2	1:C:21:LYS:HE3	1.90	0.53
6:A:306:GOL:O1	8:A:402:HOH:O	2.19	0.53
2:H:139:GLU:O	2:H:143:ILE:HG12	2.09	0.53
2:H:124:LYS:O	2:H:128:ILE:HG12	2.10	0.52
2:H:267:MET:HE1	8:H:518:HOH:O	2.09	0.52
2:H:232:ILE:HG13	2:H:247:ALA:HA	1.91	0.52
1:C:291:ILE:O	1:C:292:LEU:HB2	2.10	0.52
2:D:241:ALA:HB1	2:D:283:ILE:HD12	1.90	0.52
1:A:54:ARG:NH1	1:A:77:GLU:OE2	2.41	0.52
1:C:286:ARG:NH2	8:C:412:HOH:O	2.43	0.52
1:A:260:VAL:HG21	1:A:267:ILE:HD11	1.91	0.51
1:E:54:ARG:HH11	1:E:73:ILE:HG23	1.75	0.51
1:A:243:ILE:HG12	1:A:269:PHE:HB2	1.92	0.51
2:D:164:GLU:O	2:D:168:SER:OG	2.22	0.51
1:C:270:ILE:HD13	6:C:304:GOL:H2	1.92	0.51
1:A:86:LYS:NZ	8:A:414:HOH:O	2.45	0.50
2:D:205:ILE:HD11	2:D:259:VAL:HG22	1.94	0.50
2:D:233:ARG:NH2	8:D:506:HOH:O	2.40	0.50
1:E:218:ILE:HD11	1:E:222:ILE:HG21	1.94	0.50
2:D:295:SER:HB3	2:D:319:LEU:HD22	1.94	0.50
2:H:285:LYS:NZ	8:H:507:HOH:O	2.39	0.50
1:G:243:ILE:HG12	1:G:269:PHE:HB2	1.94	0.50
1:A:22:ALA:HB1	1:A:65:GLU:HB3	1.92	0.50
2:F:235:LYS:HD2	2:F:238:ALA:HB2	1.93	0.50
2:D:150:ILE:HG23	2:D:332:LEU:HD13	1.93	0.50
2:F:205:ILE:HD11	2:F:251:ALA:HB2	1.93	0.49
1:A:215:ILE:HG12	1:A:241:SER:HB2	1.95	0.49
2:D:205:ILE:HG23	2:D:256:ILE:HG21	1.95	0.49
1:G:287:PHE:HB2	6:G:306:GOL:H11	1.94	0.49
1:A:92:PRO:HG3	1:A:181:ILE:HD11	1.95	0.49
1:C:234:HIS:HB2	1:C:263:THR:HG22	1.95	0.48
2:F:331:ALA:HB3	2:F:362:PHE:CE1	2.48	0.48
1:A:42:LYS:HB3	6:A:304:GOL:H32	1.95	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:88:LYS:N	2:B:132:VAL:O	2.42	0.48
2:D:92:GLU:HG2	2:D:94:ASP:HB2	1.96	0.48
1:G:82:PHE:HB2	8:G:452:HOH:O	2.13	0.48
2:F:138:VAL:O	2:F:142:VAL:HG23	2.14	0.48
1:A:234:HIS:HB2	1:A:263:THR:HG22	1.95	0.48
2:F:139:GLU:O	2:F:143:ILE:HG12	2.13	0.48
1:E:19:TYR:OH	1:E:55:LEU:O	2.23	0.47
2:F:295:SER:HB3	2:F:319:LEU:HD22	1.96	0.47
1:G:85:ASP:HA	1:G:284:ALA:HB3	1.95	0.47
1:C:100:MET:HE3	1:C:211:PRO:HG3	1.96	0.47
2:F:319:LEU:HD21	2:F:330:ILE:HG21	1.96	0.47
1:C:215:ILE:HG12	1:C:241:SER:HB2	1.95	0.47
2:H:199:LYS:NZ	8:H:505:HOH:O	2.36	0.46
1:C:243:ILE:HG12	1:C:269:PHE:HB2	1.98	0.46
1:G:122:LYS:NZ	8:G:415:HOH:O	2.49	0.46
2:B:257:ASP:OD1	6:B:405:GOL:O2	2.33	0.46
1:C:153[A]:GLN:NE2	8:C:407:HOH:O	2.38	0.46
1:G:100:MET:CE	1:G:209:LEU:HD12	2.46	0.46
1:G:100:MET:HE3	1:G:211:PRO:HG3	1.97	0.45
2:D:232:ILE:HG13	2:D:247:ALA:HA	1.98	0.45
1:G:16:SER:N	8:G:417:HOH:O	2.49	0.45
1:G:218:ILE:HD11	1:G:222:ILE:HG21	1.98	0.45
2:F:198:LYS:HE2	2:F:198:LYS:HB3	1.84	0.45
1:A:66:ARG:HG3	1:A:69:TRP:H	1.81	0.45
1:C:202:MET:HE3	1:C:202:MET:HB2	1.86	0.45
2:F:206:ILE:HB	2:F:231:LEU:HD23	1.99	0.45
2:B:241:ALA:HB1	2:B:283:ILE:HD13	1.99	0.44
1:E:169:LYS:NZ	8:E:412:HOH:O	2.48	0.44
2:B:204[A]:THR:HG22	8:B:520:HOH:O	2.17	0.44
1:C:244:ILE:HD11	1:C:267:ILE:HD13	2.00	0.44
1:G:137:PRO:HA	5:G:303:G:O3'	2.17	0.44
2:B:172:LYS:HA	2:B:173:PRO:HA	1.77	0.44
1:A:22:ALA:CB	1:A:65:GLU:HB3	2.48	0.43
2:B:194:VAL:HG13	2:B:204[A]:THR:HG21	1.99	0.43
2:F:87:TYR:CE1	2:F:131:LYS:HD2	2.54	0.43
2:H:102:LEU:HD21	2:H:143:ILE:HD13	2.00	0.43
2:B:195:ASN:O	2:B:199:LYS:HG2	2.19	0.43
1:C:137:PRO:HA	5:C:303:G:O3'	2.19	0.43
1:G:104:VAL:HA	1:G:189:ARG:O	2.19	0.43
1:G:274:GLU:OE1	8:G:402:HOH:O	2.22	0.43
2:D:127:LEU:HD21	2:D:145:THR:HG21	2.01	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:86:LYS:HE2	2:B:87:TYR:CZ	2.54	0.43
1:E:100:MET:HE3	1:E:211:PRO:CG	2.49	0.43
2:H:351:TYR:CD1	6:H:403:GOL:H32	2.54	0.43
2:F:150:ILE:HG23	2:F:332:LEU:HD13	2.01	0.42
2:B:306:ARG:HG2	2:B:338:LEU:HD21	2.02	0.42
1:A:101:LEU:HD23	1:A:215:ILE:HB	2.01	0.42
6:C:305:GOL:H31	8:D:607:HOH:O	2.19	0.42
2:D:209:SER:HB2	2:D:262:ASP:O	2.20	0.41
1:E:174:PHE:HB3	1:E:179:MET:HG2	2.01	0.41
2:B:98:VAL:HB	2:B:139:GLU:OE1	2.21	0.41
1:C:27:ILE:HG23	1:C:48:THR:HB	2.03	0.41
1:C:40:ASN:O	1:C:44:VAL:HG23	2.19	0.41
1:E:94:LYS:O	1:E:95:LEU:HD23	2.20	0.41
1:E:100:MET:HE2	1:E:209:LEU:HD12	2.02	0.41
2:B:171:LYS:HA	2:B:171:LYS:HD3	1.94	0.41
1:C:19:TYR:O	1:C:23:VAL:HG23	2.21	0.41
1:E:134:VAL:HG22	8:E:443:HOH:O	2.20	0.41
2:H:124:LYS:HB2	2:H:124:LYS:HE3	1.84	0.41
1:E:234:HIS:HB2	1:E:263:THR:HG22	2.03	0.41
2:F:95:LEU:O	2:F:99:ILE:HG12	2.21	0.41
2:B:209:SER:HB2	2:B:262:ASP:O	2.21	0.41
1:C:191:GLY:HA2	1:C:195:GLU:O	2.21	0.41
1:G:82:PHE:HE1	1:G:284:ALA:HB1	1.85	0.41
2:H:215:ALA:HA	2:H:218:GLU:OE1	2.21	0.41
2:B:198:LYS:HG2	8:B:520:HOH:O	2.21	0.41
2:H:125:ASN:O	8:H:503:HOH:O	2.22	0.40
1:A:81:LEU:HG	1:A:261:VAL:HG22	2.02	0.40
2:F:160:THR:OG1	1:G:86:LYS:NZ	2.55	0.40
1:G:60:PRO:HB3	1:G:69:TRP:CE3	2.56	0.40
1:G:129:LEU:HD23	1:G:183:ILE:HB	2.04	0.40
2:H:138:VAL:O	2:H:142:VAL:HG23	2.21	0.40
1:E:69:TRP:O	1:E:72:SER:OG	2.33	0.40
1:E:167:ALA:HB1	1:E:209:LEU:HD11	2.03	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)
8:B:501:HOH:O	8:F:534:HOH:O[2_546]	2.13	0.07
8:B:502:HOH:O	8:F:599:HOH:O[2_546]	2.13	0.07
8:B:649:HOH:O	8:F:593:HOH:O[2_546]	2.19	0.01

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	270/298 (91%)	267 (99%)	3 (1%)	0	100	100
1	C	268/298 (90%)	267 (100%)	1 (0%)	0	100	100
1	E	268/298 (90%)	268 (100%)	0	0	100	100
1	G	275/298 (92%)	273 (99%)	2 (1%)	0	100	100
2	B	282/296 (95%)	276 (98%)	6 (2%)	0	100	100
2	D	268/296 (90%)	262 (98%)	6 (2%)	0	100	100
2	F	279/296 (94%)	273 (98%)	6 (2%)	0	100	100
2	H	276/296 (93%)	270 (98%)	6 (2%)	0	100	100
All	All	2186/2376 (92%)	2156 (99%)	30 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	232/252 (92%)	230 (99%)	2 (1%)	78	79
1	C	228/252 (90%)	227 (100%)	1 (0%)	91	91
1	E	228/252 (90%)	225 (99%)	3 (1%)	69	68
1	G	233/252 (92%)	230 (99%)	3 (1%)	69	68
2	B	247/259 (95%)	246 (100%)	1 (0%)	91	91
2	D	235/259 (91%)	233 (99%)	2 (1%)	78	79

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	F	246/259 (95%)	239 (97%)	7 (3%)	43	36
2	H	243/259 (94%)	234 (96%)	9 (4%)	34	25
All	All	1892/2044 (93%)	1864 (98%)	28 (2%)	65	62

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	GLU
1	A	94	LYS
2	B	358	SER
1	C	141	ASP
2	D	94	ASP
2	D	239	ASP
1	E	94	LYS
1	E	141	ASP
1	E	275	LYS
2	F	102	LEU
2	F	137	GLU
2	F	197	LEU
2	F	231	LEU
2	F	235	LYS
2	F	338	LEU
2	F	366	ILE
1	G	66	ARG
1	G	242	VAL
1	G	276	ILE
2	H	102	LEU
2	H	134	ARG
2	H	146	LEU
2	H	197	LEU
2	H	220	LEU
2	H	231	LEU
2	H	307	HIS
2	H	319	LEU
2	H	366	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	D	157	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	E	177	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 47 ligands modelled in this entry, 8 are monoatomic - leaving 39 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	F	401	4	29,34,34	1.77	6 (20%)	33,54,54	2.26	8 (24%)
6	GOL	G	304	-	5,5,5	0.35	0	5,5,5	0.97	0
6	GOL	B	406	-	5,5,5	0.34	0	5,5,5	0.39	0
3	GNP	E	301	4	29,34,34	1.68	7 (24%)	33,54,54	2.42	10 (30%)
6	GOL	G	305	-	5,5,5	0.40	0	5,5,5	0.51	0
6	GOL	E	304	-	5,5,5	0.34	0	5,5,5	0.30	0
3	GNP	G	301	4	29,34,34	1.65	6 (20%)	33,54,54	2.36	9 (27%)
7	SO4	C	306	-	4,4,4	0.15	0	6,6,6	0.05	0
6	GOL	D	403	-	5,5,5	0.43	0	5,5,5	0.48	0
7	SO4	B	407	-	4,4,4	0.16	0	6,6,6	0.13	0
5	G	E	303	-	22,26,26	1.02	2 (9%)	26,40,40	0.83	1 (3%)
3	GNP	A	301	4	29,34,34	1.67	5 (17%)	33,54,54	2.51	9 (27%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	GOL	B	403	-	5,5,5	0.40	0	5,5,5	0.47	0
3	GNP	H	401	4	29,34,34	1.60	5 (17%)	33,54,54	2.10	5 (15%)
6	GOL	B	404	-	5,5,5	0.30	0	5,5,5	0.37	0
3	GNP	B	401	4	29,34,34	1.78	5 (17%)	33,54,54	2.04	5 (15%)
6	GOL	A	306	-	5,5,5	0.35	0	5,5,5	0.32	0
5	G	A	303	-	22,26,26	1.03	2 (9%)	26,40,40	0.81	1 (3%)
7	SO4	A	308	-	4,4,4	0.13	0	6,6,6	0.13	0
3	GNP	C	301	4	29,34,34	1.80	6 (20%)	33,54,54	2.36	9 (27%)
7	SO4	A	307	-	4,4,4	0.14	0	6,6,6	0.05	0
7	SO4	G	307	-	4,4,4	0.14	0	6,6,6	0.07	0
7	SO4	E	306	-	4,4,4	0.14	0	6,6,6	0.05	0
6	GOL	H	403	-	5,5,5	0.51	0	5,5,5	0.54	0
6	GOL	H	404	-	5,5,5	0.37	0	5,5,5	0.16	0
7	SO4	B	408	-	4,4,4	0.14	0	6,6,6	0.05	0
6	GOL	F	403	-	5,5,5	0.40	0	5,5,5	0.34	0
6	GOL	G	306	-	5,5,5	0.35	0	5,5,5	0.32	0
5	G	G	303	-	22,26,26	1.05	3 (13%)	26,40,40	0.95	2 (7%)
7	SO4	A	309	-	4,4,4	0.15	0	6,6,6	0.13	0
6	GOL	E	305	-	5,5,5	0.35	0	5,5,5	0.55	0
3	GNP	D	401	4	29,34,34	1.67	4 (13%)	33,54,54	2.18	7 (21%)
5	G	C	303	-	22,26,26	1.06	2 (9%)	26,40,40	0.97	1 (3%)
6	GOL	A	304	-	5,5,5	0.34	0	5,5,5	0.28	0
6	GOL	C	304	-	5,5,5	0.31	0	5,5,5	0.30	0
7	SO4	H	405	-	4,4,4	0.16	0	6,6,6	0.14	0
6	GOL	A	305	-	5,5,5	0.36	0	5,5,5	0.50	0
6	GOL	C	305	-	5,5,5	0.34	0	5,5,5	0.39	0
6	GOL	B	405	-	5,5,5	0.43	0	5,5,5	0.29	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GNP	F	401	4	-	3/14/38/38	0/3/3/3
6	GOL	G	304	-	-	2/4/4/4	-
6	GOL	B	406	-	-	2/4/4/4	-
3	GNP	E	301	4	-	2/14/38/38	0/3/3/3
6	GOL	G	305	-	-	0/4/4/4	-
6	GOL	E	304	-	-	4/4/4/4	-

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GNP	G	301	4	-	2/14/38/38	0/3/3/3
6	GOL	D	403	-	-	2/4/4/4	-
5	G	E	303	-	-	2/6/26/26	0/3/3/3
3	GNP	A	301	4	-	2/14/38/38	0/3/3/3
6	GOL	B	403	-	-	2/4/4/4	-
3	GNP	H	401	4	-	3/14/38/38	0/3/3/3
6	GOL	B	404	-	-	0/4/4/4	-
3	GNP	B	401	4	-	3/14/38/38	0/3/3/3
6	GOL	A	306	-	-	2/4/4/4	-
5	G	A	303	-	-	1/6/26/26	0/3/3/3
3	GNP	C	301	4	-	2/14/38/38	0/3/3/3
6	GOL	H	403	-	-	3/4/4/4	-
6	GOL	H	404	-	-	2/4/4/4	-
6	GOL	F	403	-	-	2/4/4/4	-
6	GOL	G	306	-	-	4/4/4/4	-
5	G	G	303	-	-	2/6/26/26	0/3/3/3
6	GOL	E	305	-	-	2/4/4/4	-
3	GNP	D	401	4	-	3/14/38/38	0/3/3/3
5	G	C	303	-	-	2/6/26/26	0/3/3/3
6	GOL	A	304	-	-	4/4/4/4	-
6	GOL	C	304	-	-	2/4/4/4	-
6	GOL	A	305	-	-	0/4/4/4	-
6	GOL	C	305	-	-	0/4/4/4	-
6	GOL	B	405	-	-	0/4/4/4	-

All (53) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	401	GNP	PB-O3A	-5.75	1.51	1.59
3	F	401	GNP	PB-O3A	-5.69	1.51	1.59
3	D	401	GNP	PB-O3A	-5.06	1.52	1.59
3	A	301	GNP	PG-O1G	4.99	1.54	1.46
3	C	301	GNP	PG-O1G	4.53	1.53	1.46
3	E	301	GNP	PG-O1G	4.34	1.53	1.46
3	H	401	GNP	PB-O3A	-4.20	1.53	1.59
3	B	401	GNP	C6-N1	4.14	1.40	1.33
3	C	301	GNP	C6-N1	4.14	1.40	1.33
3	A	301	GNP	C6-N1	4.12	1.40	1.33

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	301	GNP	C6-N1	4.11	1.40	1.33
3	F	401	GNP	C6-N1	4.09	1.40	1.33
3	G	301	GNP	PB-O3A	-4.09	1.53	1.59
3	G	301	GNP	C6-N1	4.00	1.40	1.33
3	D	401	GNP	C6-N1	3.90	1.39	1.33
3	C	301	GNP	PB-O2B	-3.89	1.46	1.56
3	H	401	GNP	C6-N1	3.82	1.39	1.33
3	E	301	GNP	PB-O3A	-3.67	1.54	1.59
3	D	401	GNP	PB-O2B	-3.60	1.47	1.56
3	C	301	GNP	PB-O3A	-3.57	1.54	1.59
3	G	301	GNP	PB-O2B	-3.48	1.47	1.56
3	B	401	GNP	PB-O2B	-3.43	1.47	1.56
3	B	401	GNP	PG-O1G	3.40	1.51	1.46
3	H	401	GNP	PB-O2B	-3.36	1.47	1.56
3	F	401	GNP	PG-O1G	3.23	1.51	1.46
3	G	301	GNP	PG-O1G	3.23	1.51	1.46
3	A	301	GNP	PB-O3A	-3.20	1.55	1.59
3	H	401	GNP	PG-O1G	3.16	1.51	1.46
3	A	301	GNP	PB-O2B	-3.08	1.48	1.56
3	D	401	GNP	PG-O1G	3.03	1.51	1.46
3	F	401	GNP	PB-O2B	-2.94	1.48	1.56
3	E	301	GNP	PB-O2B	-2.60	1.49	1.56
3	C	301	GNP	PG-O2G	-2.56	1.49	1.56
3	E	301	GNP	PB-O1B	2.48	1.50	1.46
5	A	303	G	C5-C6	-2.43	1.42	1.47
5	E	303	G	C5-C6	-2.41	1.42	1.47
5	C	303	G	C5-C6	-2.40	1.42	1.47
3	G	301	GNP	PG-O2G	-2.34	1.50	1.56
5	E	303	G	C8-N7	-2.33	1.31	1.35
5	G	303	G	C5-C6	-2.32	1.42	1.47
3	G	301	GNP	PB-O1B	2.31	1.49	1.46
5	C	303	G	C8-N7	-2.31	1.31	1.35
3	E	301	GNP	PG-O2G	-2.30	1.50	1.56
3	F	401	GNP	PG-O3G	-2.29	1.50	1.56
3	A	301	GNP	PG-O2G	-2.23	1.50	1.56
5	A	303	G	C8-N7	-2.23	1.31	1.35
3	H	401	GNP	PB-O1B	2.22	1.49	1.46
3	C	301	GNP	PB-O1B	2.15	1.49	1.46
5	G	303	G	C8-N7	-2.13	1.31	1.35
3	E	301	GNP	PG-O3G	-2.11	1.51	1.56
3	B	401	GNP	C2-N1	2.10	1.39	1.35
3	F	401	GNP	C2-N1	2.08	1.39	1.35

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	G	303	G	P-OP3	-2.04	1.47	1.54

All (67) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	401	GNP	C5-C6-N1	-8.35	112.00	123.43
3	E	301	GNP	C5-C6-N1	-8.34	112.02	123.43
3	A	301	GNP	C5-C6-N1	-8.27	112.11	123.43
3	D	401	GNP	C5-C6-N1	-8.19	112.22	123.43
3	B	401	GNP	C5-C6-N1	-8.14	112.30	123.43
3	F	401	GNP	C5-C6-N1	-8.01	112.47	123.43
3	G	301	GNP	C5-C6-N1	-7.92	112.60	123.43
3	C	301	GNP	C5-C6-N1	-7.70	112.90	123.43
3	E	301	GNP	C2-N1-C6	5.83	125.19	115.93
3	A	301	GNP	C2-N1-C6	5.75	125.07	115.93
3	D	401	GNP	C2-N1-C6	5.63	124.88	115.93
3	H	401	GNP	C2-N1-C6	5.63	124.87	115.93
3	G	301	GNP	C2-N1-C6	5.45	124.59	115.93
3	C	301	GNP	C2-N1-C6	5.40	124.51	115.93
3	F	401	GNP	C2-N1-C6	5.37	124.47	115.93
3	B	401	GNP	C2-N1-C6	5.08	124.00	115.93
3	E	301	GNP	O2B-PB-O1B	5.03	120.46	109.92
3	A	301	GNP	O1B-PB-N3B	-4.64	104.94	111.77
3	G	301	GNP	O1G-PG-N3B	-4.55	105.06	111.77
3	A	301	GNP	O2B-PB-O1B	4.30	118.93	109.92
3	F	401	GNP	O2B-PB-O1B	4.15	118.63	109.92
3	F	401	GNP	O1B-PB-N3B	-4.02	105.86	111.77
3	D	401	GNP	O2B-PB-O1B	3.84	117.97	109.92
3	C	301	GNP	O2B-PB-O1B	3.81	117.91	109.92
3	C	301	GNP	O3G-PG-O1G	-3.71	104.13	113.45
3	G	301	GNP	O2B-PB-O1B	3.65	117.57	109.92
3	A	301	GNP	N3-C2-N1	-3.41	122.68	127.22
3	A	301	GNP	O1G-PG-N3B	-3.41	106.75	111.77
3	A	301	GNP	O3G-PG-O1G	-3.24	105.30	113.45
3	E	301	GNP	O3G-PG-O1G	-3.19	105.44	113.45
3	G	301	GNP	N3-C2-N1	-3.17	122.99	127.22
3	E	301	GNP	N3-C2-N1	-3.16	123.01	127.22
3	C	301	GNP	N3-C2-N1	-3.15	123.02	127.22
3	A	301	GNP	C4-C5-C6	-3.12	117.82	120.80
3	C	301	GNP	O2G-PG-O3G	3.11	115.92	107.64
3	C	301	GNP	C4-C5-C6	-3.10	117.83	120.80
3	C	301	GNP	O1G-PG-N3B	-3.08	107.23	111.77

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	401	GNP	N3-C2-N1	-3.06	123.14	127.22
3	E	301	GNP	C4-C5-C6	-2.85	118.07	120.80
3	D	401	GNP	N3-C2-N1	-2.83	123.45	127.22
3	G	301	GNP	O2G-PG-O3G	2.80	115.09	107.64
3	B	401	GNP	C2-N3-C4	-2.79	112.17	115.36
3	G	301	GNP	C4-C5-C6	-2.70	118.22	120.80
3	A	301	GNP	O2G-PG-O3G	2.69	114.81	107.64
3	H	401	GNP	C2-N3-C4	-2.66	112.32	115.36
3	E	301	GNP	O1B-PB-N3B	-2.65	107.88	111.77
3	G	301	GNP	O3G-PG-O1G	-2.61	106.89	113.45
3	B	401	GNP	N3-C2-N1	-2.56	123.80	127.22
3	D	401	GNP	C4-C5-C6	-2.54	118.37	120.80
3	E	301	GNP	O1G-PG-N3B	-2.51	108.08	111.77
3	F	401	GNP	C4-C5-C6	-2.50	118.41	120.80
3	D	401	GNP	C2-N3-C4	-2.48	112.52	115.36
3	H	401	GNP	N3-C2-N1	-2.47	123.93	127.22
3	G	301	GNP	O1B-PB-N3B	-2.45	108.16	111.77
3	E	301	GNP	O2G-PG-O3G	2.41	114.07	107.64
3	E	301	GNP	C2-N3-C4	-2.37	112.65	115.36
3	D	401	GNP	O1G-PG-N3B	-2.36	108.30	111.77
5	G	303	G	OP2-P-OP1	2.26	119.52	110.68
5	G	303	G	O6-C6-C5	2.24	128.74	124.37
3	B	401	GNP	O1G-PG-N3B	-2.21	108.52	111.77
3	F	401	GNP	C2-N3-C4	-2.19	112.86	115.36
5	A	303	G	O6-C6-C5	2.17	128.62	124.37
5	C	303	G	O6-C6-C5	2.15	128.57	124.37
3	F	401	GNP	O1G-PG-N3B	-2.14	108.62	111.77
3	H	401	GNP	O2B-PB-O1B	2.13	114.39	109.92
5	E	303	G	O6-C6-C5	2.10	128.48	124.37
3	C	301	GNP	O1B-PB-N3B	-2.05	108.75	111.77

There are no chirality outliers.

All (60) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	301	GNP	PG-N3B-PB-O1B
3	A	301	GNP	PG-N3B-PB-O3A
3	B	401	GNP	PB-N3B-PG-O1G
3	B	401	GNP	PG-N3B-PB-O1B
3	B	401	GNP	PG-N3B-PB-O3A
3	C	301	GNP	PG-N3B-PB-O1B
3	C	301	GNP	PG-N3B-PB-O3A

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
3	D	401	GNP	PB-N3B-PG-O1G
3	D	401	GNP	PG-N3B-PB-O1B
3	E	301	GNP	PG-N3B-PB-O1B
3	E	301	GNP	PG-N3B-PB-O3A
3	F	401	GNP	PB-N3B-PG-O1G
3	F	401	GNP	PG-N3B-PB-O1B
3	F	401	GNP	PG-N3B-PB-O3A
3	G	301	GNP	PG-N3B-PB-O1B
3	G	301	GNP	PG-N3B-PB-O3A
3	H	401	GNP	PB-N3B-PG-O1G
3	H	401	GNP	PG-N3B-PB-O1B
3	H	401	GNP	PG-N3B-PB-O3A
5	C	303	G	C4'-C5'-O5'-P
5	G	303	G	C4'-C5'-O5'-P
6	A	304	GOL	O1-C1-C2-C3
6	A	304	GOL	C1-C2-C3-O3
6	A	306	GOL	O1-C1-C2-C3
6	B	403	GOL	O1-C1-C2-C3
6	B	406	GOL	O1-C1-C2-C3
6	D	403	GOL	O1-C1-C2-C3
6	E	305	GOL	O1-C1-C2-C3
6	F	403	GOL	O1-C1-C2-C3
6	G	304	GOL	C1-C2-C3-O3
6	G	306	GOL	O1-C1-C2-C3
6	G	306	GOL	C1-C2-C3-O3
6	H	403	GOL	O1-C1-C2-C3
6	H	404	GOL	O1-C1-C2-C3
6	A	304	GOL	O1-C1-C2-O2
6	D	403	GOL	O1-C1-C2-O2
5	E	303	G	C4'-C5'-O5'-P
6	E	304	GOL	C1-C2-C3-O3
6	B	403	GOL	O1-C1-C2-O2
6	B	406	GOL	O1-C1-C2-O2
6	E	304	GOL	O2-C2-C3-O3
6	F	403	GOL	O1-C1-C2-O2
5	A	303	G	C4'-C5'-O5'-P
6	A	306	GOL	O1-C1-C2-O2
6	E	305	GOL	O1-C1-C2-O2
6	G	306	GOL	O1-C1-C2-O2
6	G	306	GOL	O2-C2-C3-O3
6	H	403	GOL	O1-C1-C2-O2
6	H	404	GOL	O1-C1-C2-O2

Continued on next page...

Continued from previous page...

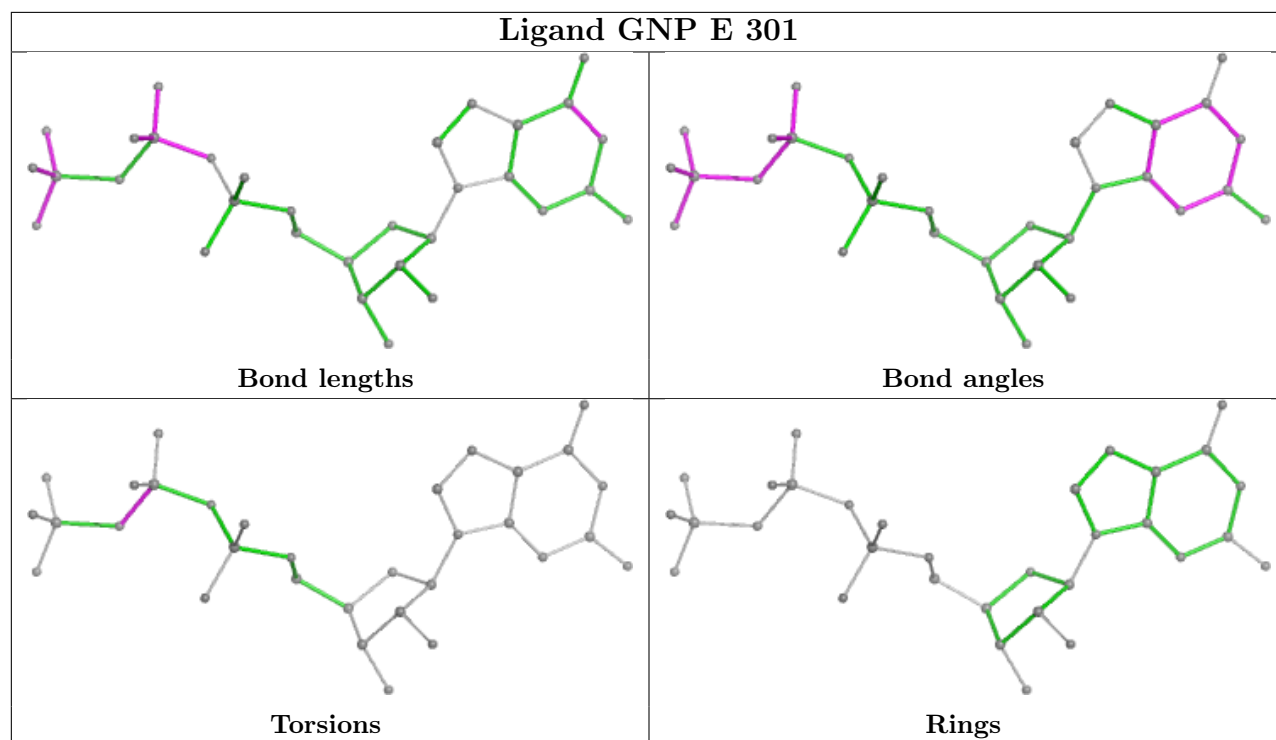
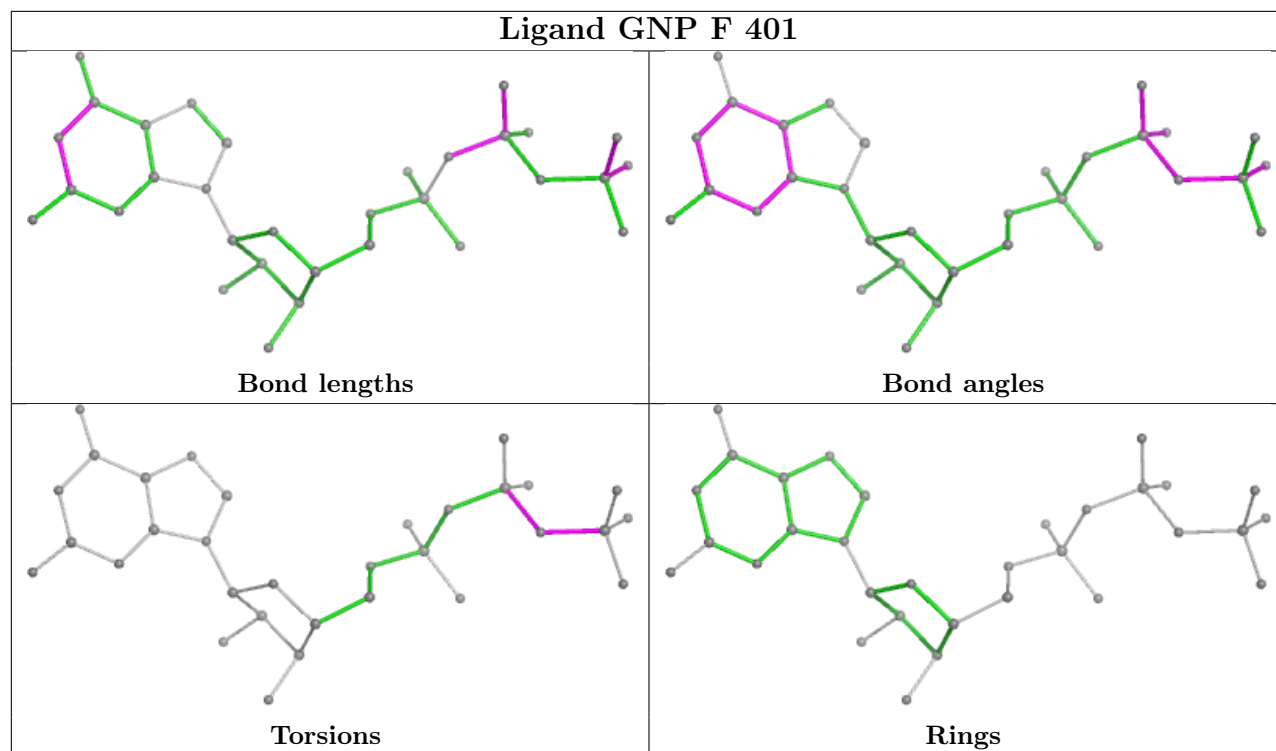
Mol	Chain	Res	Type	Atoms
6	A	304	GOL	O2-C2-C3-O3
6	E	304	GOL	O1-C1-C2-O2
6	C	304	GOL	C1-C2-C3-O3
3	D	401	GNP	PG-N3B-PB-O3A
6	G	304	GOL	O2-C2-C3-O3
6	C	304	GOL	O2-C2-C3-O3
5	E	303	G	O4'-C4'-C5'-O5'
5	G	303	G	O4'-C4'-C5'-O5'
6	E	304	GOL	O1-C1-C2-C3
5	C	303	G	O4'-C4'-C5'-O5'
6	H	403	GOL	O2-C2-C3-O3

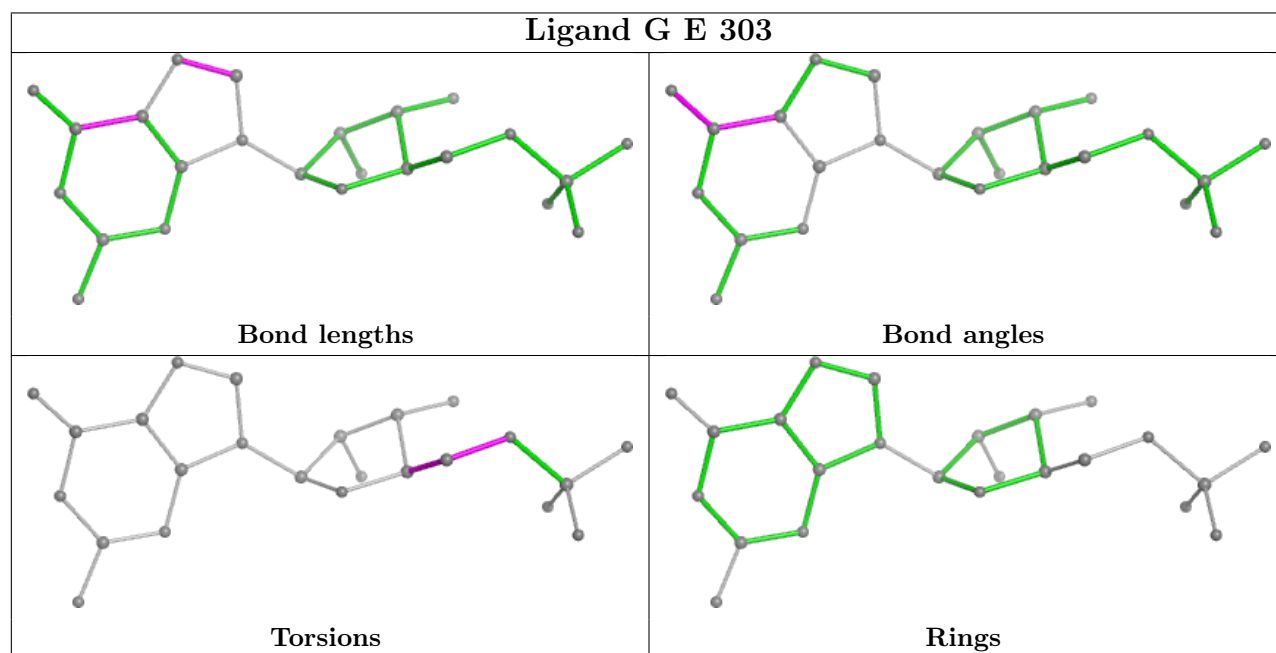
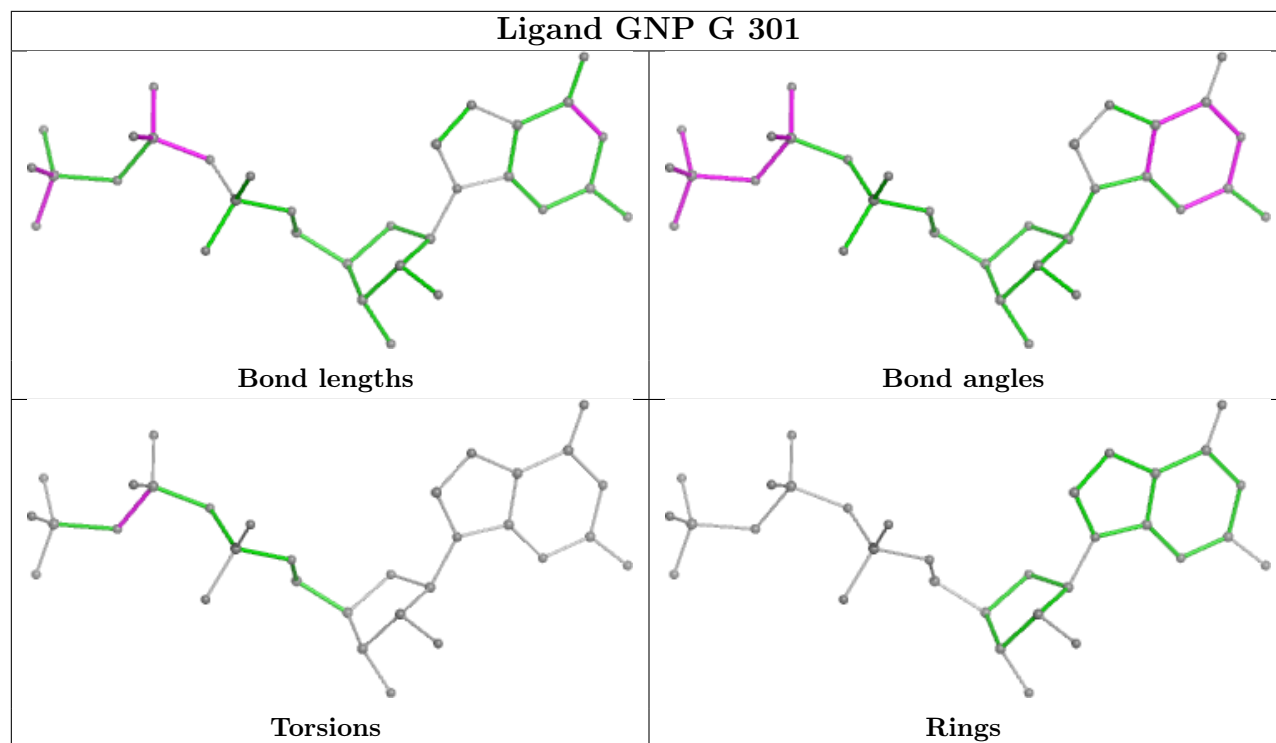
There are no ring outliers.

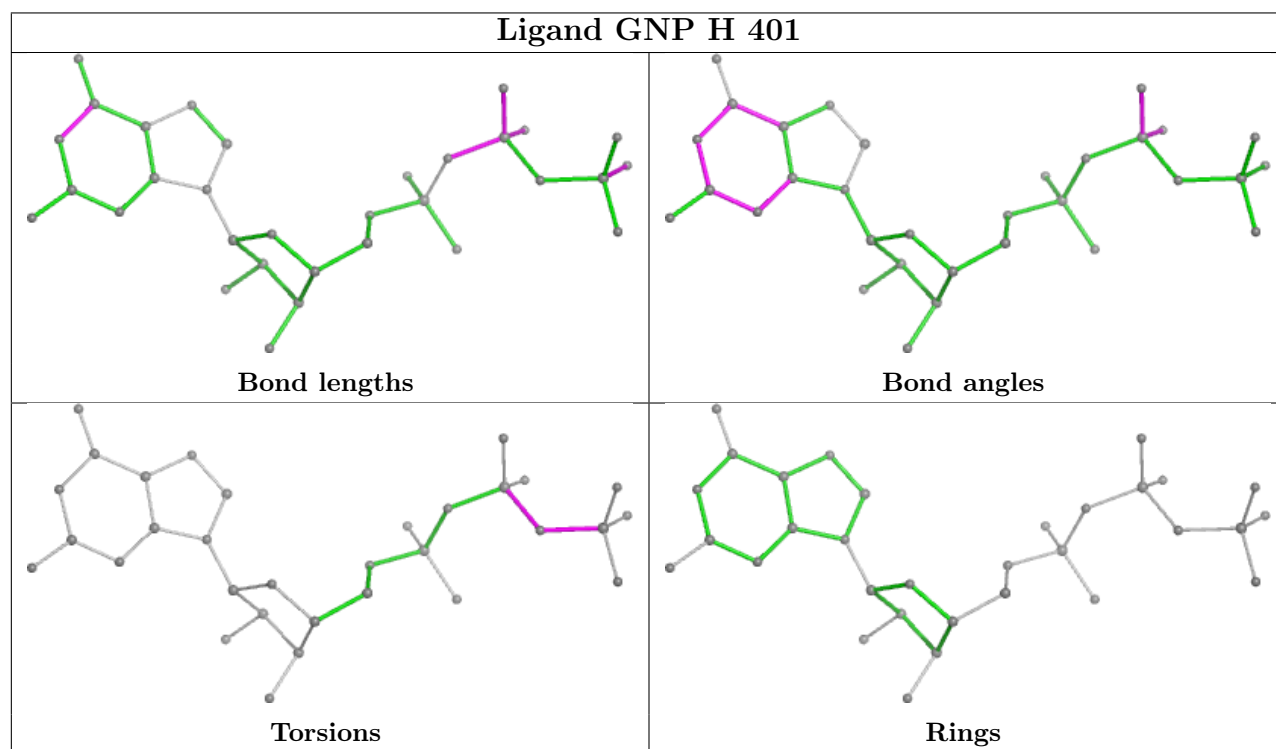
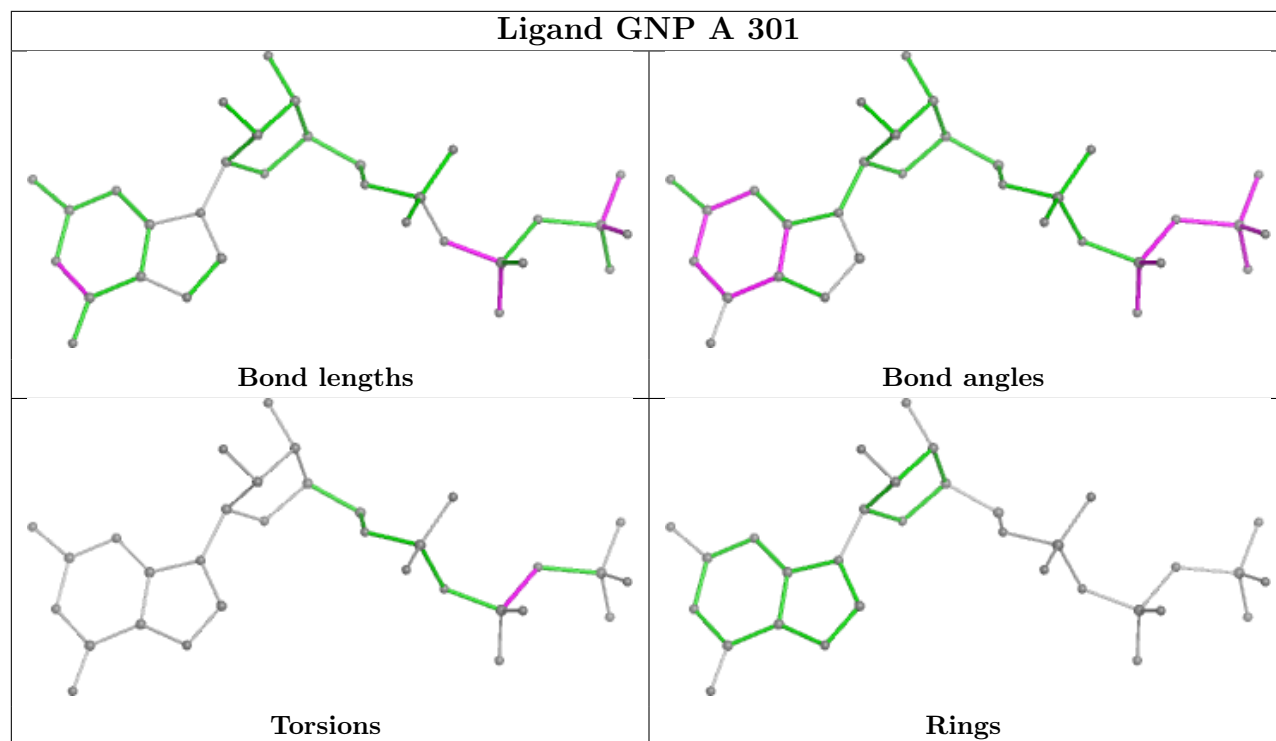
11 monomers are involved in 13 short contacts:

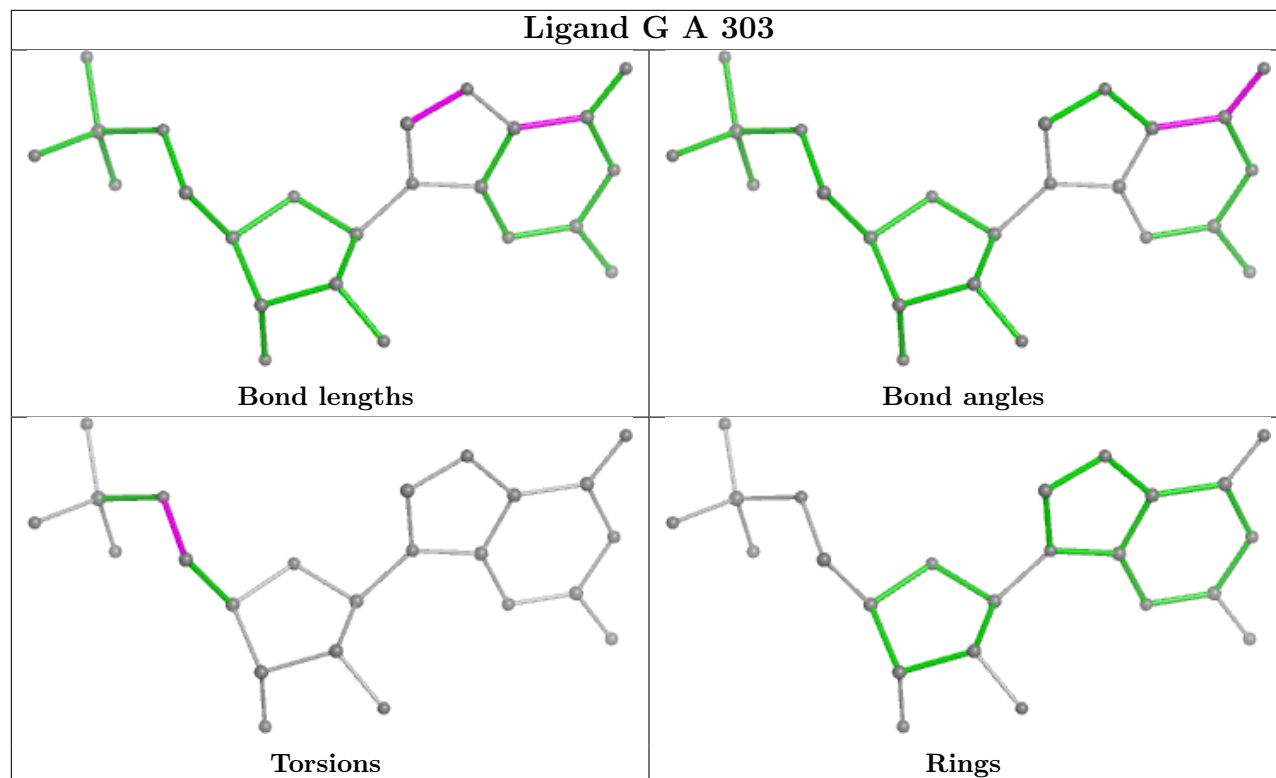
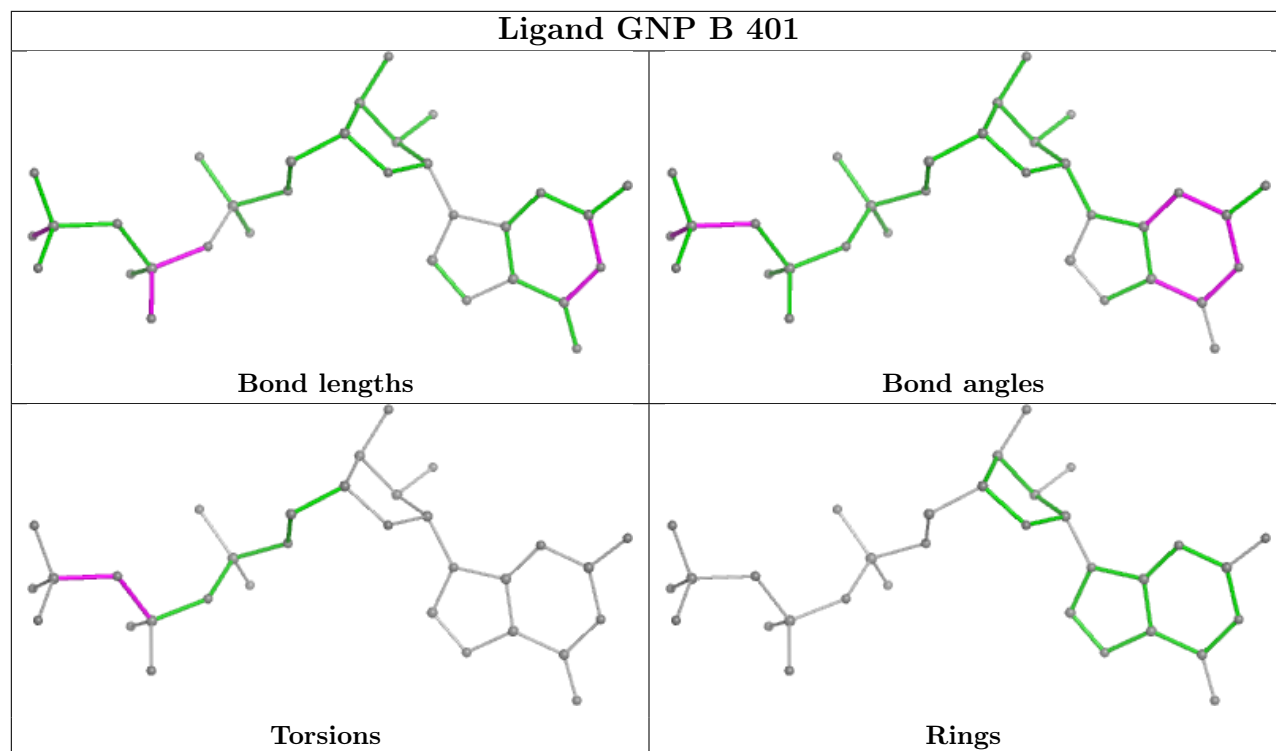
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	306	SO4	1	0
6	A	306	GOL	1	0
6	H	403	GOL	1	0
6	G	306	GOL	2	0
5	G	303	G	1	0
5	C	303	G	1	0
6	A	304	GOL	1	0
6	C	304	GOL	2	0
6	A	305	GOL	1	0
6	C	305	GOL	1	0
6	B	405	GOL	1	0

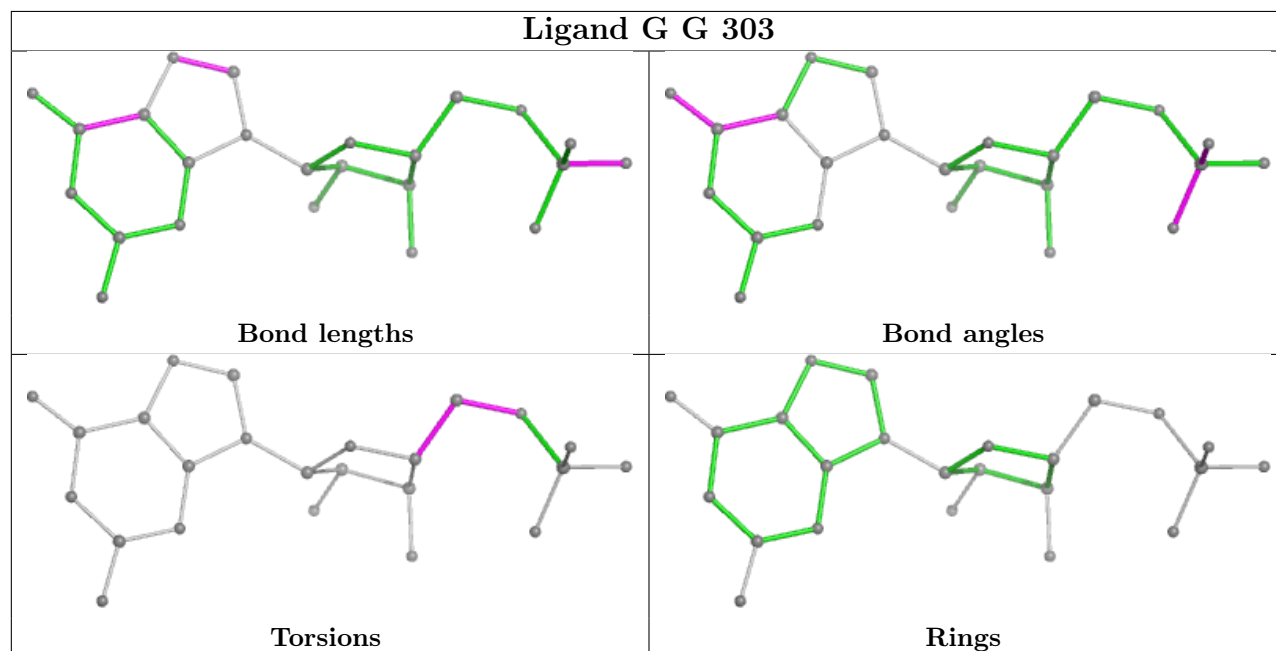
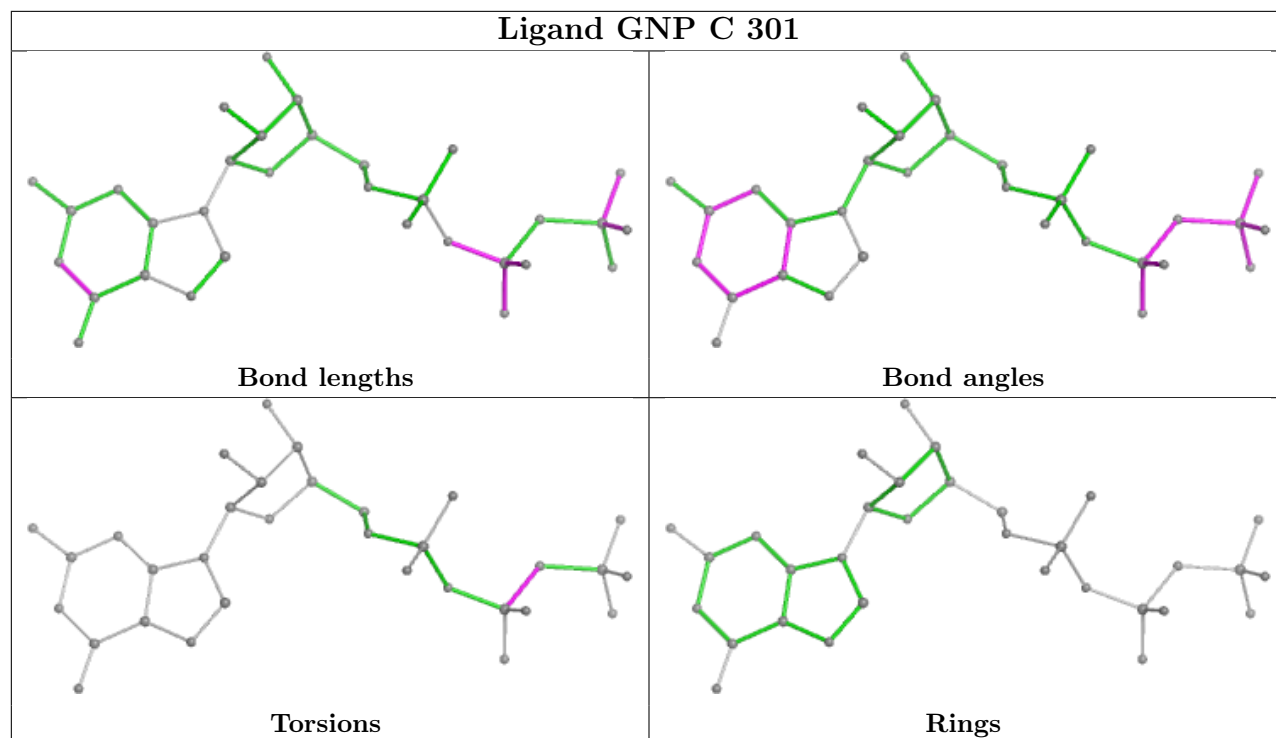
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

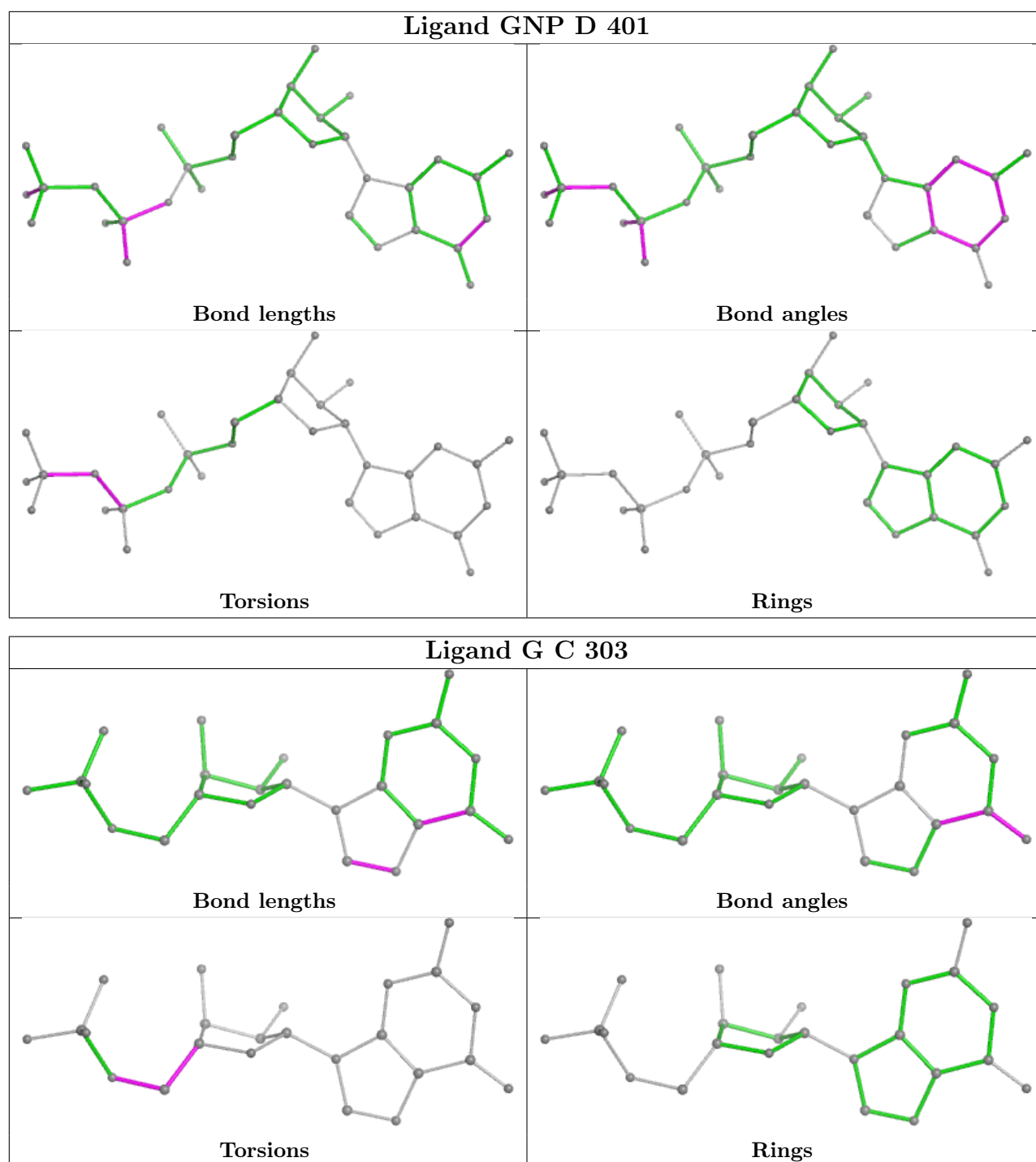












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	274/298 (91%)	0.41	21 (7%) 13 15	18, 34, 73, 113	0
1	C	270/298 (90%)	1.17	57 (21%) 1 0	22, 44, 106, 135	0
1	E	271/298 (90%)	0.48	32 (11%) 4 5	22, 39, 97, 126	0
1	G	277/298 (92%)	0.27	15 (5%) 25 29	18, 38, 70, 112	0
2	B	283/296 (95%)	0.29	17 (6%) 21 24	19, 33, 64, 107	0
2	D	271/296 (91%)	0.73	37 (13%) 3 3	23, 44, 100, 120	0
2	F	283/296 (95%)	0.60	23 (8%) 12 13	23, 46, 79, 123	0
2	H	278/296 (93%)	0.50	32 (11%) 4 5	19, 36, 90, 117	0
All	All	2207/2376 (92%)	0.55	234 (10%) 6 7	18, 39, 89, 135	0

All (234) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	236	TYR	9.8
1	G	63	VAL	9.4
1	C	19	TYR	8.9
2	H	368	SER	8.7
2	H	132	VAL	8.2
2	D	236	TYR	8.1
1	C	67	LYS	8.0
1	C	73	ILE	8.0
2	D	95	LEU	7.8
1	C	17	THR	7.8
1	C	23	VAL	7.4
2	D	138	VAL	7.2
2	H	90	ILE	7.0
1	C	58	GLU	7.0
2	H	95	LEU	6.9
1	E	70	PHE	6.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	56	ASN	6.5
1	A	63	VAL	6.3
1	C	49	ALA	6.2
2	D	94	ASP	6.0
2	D	128	ILE	6.0
1	C	55	LEU	6.0
2	F	87	TYR	5.9
1	C	57	LYS	5.9
2	B	87	TYR	5.9
1	A	65	GLU	5.8
1	C	71	ILE	5.8
1	C	70	PHE	5.6
1	C	74	VAL	5.5
1	C	285	LYS	5.3
2	D	139	GLU	5.2
2	D	97	ASP	5.2
1	C	69	TRP	5.2
2	H	138	VAL	5.2
2	H	127	LEU	5.2
2	H	128	ILE	5.1
2	F	132	VAL	5.1
1	C	59	LYS	5.1
1	C	76	ASP	5.0
1	E	66	ARG	5.0
1	C	53	GLU	5.0
1	C	20	GLU	5.0
2	D	141	ILE	5.0
2	H	91	LYS	4.9
1	G	64	LEU	4.9
2	H	93	ASP	4.7
2	H	97	ASP	4.7
1	C	50	LYS	4.7
1	E	16	SER	4.7
2	F	86	LYS	4.6
1	E	23	VAL	4.6
1	C	222	ILE	4.5
2	B	134	ARG	4.5
2	H	137	GLU	4.5
1	C	84	GLY	4.5
2	D	368	SER	4.5
2	H	94	ASP	4.4
2	H	236[A]	TYR	4.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	72	SER	4.2
1	C	22	ALA	4.2
2	F	138	VAL	4.1
1	C	18	PRO	4.1
1	C	51	ILE	4.1
1	C	28	LYS	4.1
1	C	54	ARG	4.1
1	C	75	TYR	3.9
2	H	92	GLU	3.9
1	A	222	ILE	3.9
1	C	218	ILE	3.9
2	D	361	TRP	3.8
2	D	235	LYS	3.8
1	E	69	TRP	3.8
2	H	141	ILE	3.8
1	E	68	GLU	3.8
2	B	132	VAL	3.8
2	D	142	VAL	3.8
1	G	62	SER	3.7
2	H	129	GLY	3.7
2	B	296	LEU	3.7
1	C	47	LEU	3.7
2	D	184	VAL	3.7
2	H	131	LYS	3.6
1	C	24	ASP	3.6
2	D	364	GLU	3.6
1	C	21	LYS	3.6
2	B	135	ARG	3.6
2	F	296	LEU	3.5
2	B	184	VAL	3.5
2	D	98	VAL	3.4
1	G	86	LYS	3.4
2	H	98	VAL	3.4
2	D	104	PHE	3.4
1	C	217	VAL	3.4
1	E	67	LYS	3.4
2	D	170	GLY	3.3
1	A	64	LEU	3.3
1	C	104	VAL	3.3
2	D	99	ILE	3.2
2	D	127	LEU	3.2
1	E	45	PHE	3.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	B	297	ALA	3.2
1	G	85	ASP	3.2
2	D	140	GLU	3.2
2	H	135	ARG	3.1
2	F	235	LYS	3.1
2	H	139	GLU	3.1
1	E	24	ASP	3.1
1	E	28	LYS	3.1
1	C	79	SER	3.0
2	D	92	GLU	3.0
2	B	319	LEU	3.0
2	H	121	GLU	3.0
2	D	311	ASN	3.0
1	C	80	LYS	3.0
1	C	194	GLU	3.0
1	E	85	ASP	3.0
1	G	83	GLY	2.9
2	F	85	LEU	2.9
2	H	133	SER	2.9
2	B	293	LEU	2.9
1	E	18	PRO	2.9
1	E	53	GLU	2.9
2	F	135	ARG	2.8
2	H	126	ASN	2.8
1	A	218	ILE	2.8
1	A	83	GLY	2.8
1	C	68	GLU	2.8
2	F	362	PHE	2.8
1	E	17	THR	2.8
1	C	52	LYS	2.7
2	D	237	GLY	2.7
1	E	54	ARG	2.7
1	G	68	GLU	2.7
1	C	286	ARG	2.7
2	D	255	ASN	2.7
1	A	87	GLU	2.7
1	E	135[A]	TYR	2.7
2	F	141	ILE	2.7
2	B	137	GLU	2.7
2	F	184	VAL	2.6
2	H	184	VAL	2.6
2	F	93	ASP	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	F	160	THR	2.6
1	C	86	LYS	2.6
1	E	56	ASN	2.6
1	E	49	ALA	2.6
1	G	218	ILE	2.6
2	H	96	ASN	2.6
2	F	136	GLU	2.6
1	A	81	LEU	2.5
1	C	196	THR	2.5
1	C	25	GLU	2.5
2	B	136	GLU	2.5
1	A	16	SER	2.5
1	C	107	SER	2.5
1	E	22	ALA	2.5
2	B	86	LYS	2.5
1	A	67	LYS	2.5
2	F	139	GLU	2.5
2	H	124	LYS	2.5
1	G	65	GLU	2.5
2	H	99	ILE	2.5
1	E	52	LYS	2.5
2	F	133	SER	2.5
2	D	96	ASN	2.4
1	C	77	GLU	2.4
2	B	330	ILE	2.4
1	G	87	GLU	2.4
2	D	253	SER	2.4
1	E	286	ARG	2.4
2	B	318	ILE	2.4
2	D	172	LYS	2.4
1	C	220	ALA	2.3
1	E	57	LYS	2.3
1	A	69	TRP	2.3
1	C	16	SER	2.3
1	C	102	VAL	2.3
1	E	47	LEU	2.3
2	D	296	LEU	2.3
1	C	85	ASP	2.3
1	C	87	GLU	2.3
1	E	26	PHE	2.3
1	A	86	LYS	2.3
2	H	142	VAL	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	71	ILE	2.3
1	G	222	ILE	2.3
1	C	78	LEU	2.3
1	C	26	PHE	2.3
1	C	82	PHE	2.3
1	G	226	ALA	2.3
2	B	368	SER	2.3
2	H	123	LEU	2.3
2	D	137	GLU	2.2
1	G	57	LYS	2.2
1	A	216	LEU	2.2
1	A	135[A]	TYR	2.2
1	A	220	ALA	2.2
1	E	256	ALA	2.2
2	F	297	ALA	2.2
1	A	217	VAL	2.2
1	E	73	ILE	2.2
1	E	20	GLU	2.2
1	C	135[A]	TYR	2.2
1	A	260	VAL	2.2
1	E	59	LYS	2.2
1	A	292	LEU	2.2
2	H	134	ARG	2.2
2	B	171	LYS	2.2
2	F	157	ASN	2.2
2	B	139	GLU	2.2
1	E	222	ILE	2.2
1	G	52	LYS	2.2
2	D	129	GLY	2.2
1	E	218	ILE	2.1
2	H	172	LYS	2.1
2	D	93	ASP	2.1
2	D	310	ASN	2.1
2	F	318	ILE	2.1
2	D	254	ARG	2.1
1	C	113	ALA	2.1
2	H	143	ILE	2.1
1	E	58	GLU	2.1
2	D	126	ASN	2.1
1	A	257	LEU	2.1
2	D	123	LEU	2.1
1	A	66	ARG	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	D	145	THR	2.0
2	F	98	VAL	2.0
2	F	292	ILE	2.0
1	G	194	GLU	2.0
2	F	156	LYS	2.0
2	D	144	ASN	2.0
1	A	197	LYS	2.0
1	C	245	THR	2.0

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 monosaccharides 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
6	GOL	H	404	6/6	0.49	0.45	74,84,86,87	0
5	G	A	303	24/24	0.55	0.43	84,97,107,109	24
6	GOL	A	304	6/6	0.60	0.24	54,63,71,73	0
6	GOL	C	305	6/6	0.62	0.29	67,76,78,79	0
5	G	E	303	24/24	0.70	0.36	58,83,95,99	24
6	GOL	B	403	6/6	0.78	0.21	45,47,52,52	0
6	GOL	F	403	6/6	0.80	0.20	50,55,58,58	0
6	GOL	B	406	6/6	0.80	0.27	65,72,74,77	0
7	SO4	C	306	5/5	0.81	0.29	108,109,110,111	0
7	SO4	G	307	5/5	0.81	0.35	103,104,104,107	0
6	GOL	C	304	6/6	0.82	0.16	46,51,56,59	0
5	G	C	303	24/24	0.82	0.30	53,83,110,112	0
7	SO4	E	306	5/5	0.83	0.27	134,135,136,137	0
6	GOL	G	306	6/6	0.83	0.14	39,52,53,56	0
7	SO4	B	408	5/5	0.84	0.25	119,121,122,122	0

Continued on next page...

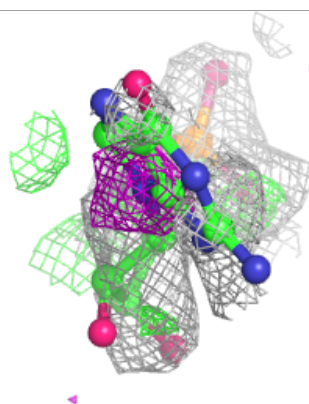
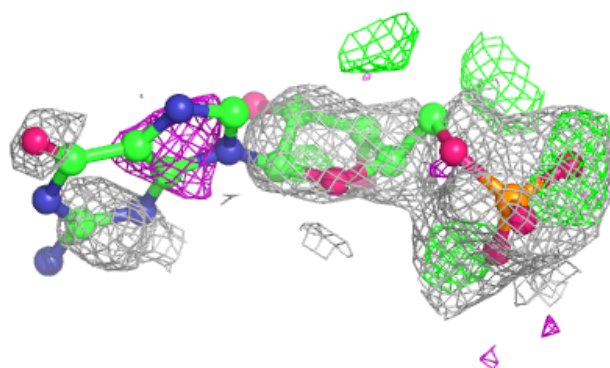
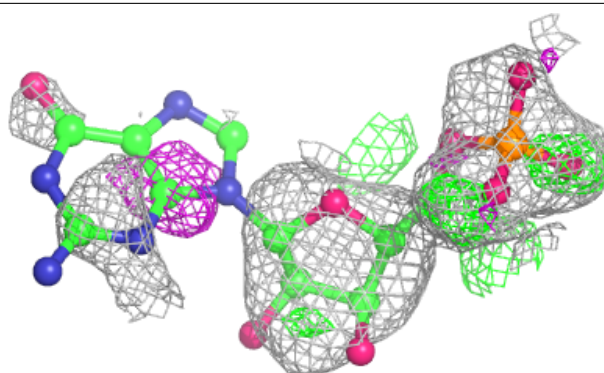
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
6	GOL	B	404	6/6	0.84	0.17	46,58,60,61	0
5	G	G	303	24/24	0.85	0.29	35,80,100,105	24
6	GOL	B	405	6/6	0.86	0.17	40,45,52,54	0
6	GOL	A	306	6/6	0.87	0.20	56,59,63,64	0
6	GOL	E	304	6/6	0.88	0.13	45,46,55,63	0
6	GOL	A	305	6/6	0.88	0.15	29,35,43,45	0
6	GOL	G	305	6/6	0.88	0.18	45,54,60,60	0
6	GOL	G	304	6/6	0.90	0.14	30,37,51,61	0
6	GOL	D	403	6/6	0.90	0.16	40,43,47,48	0
6	GOL	H	403	6/6	0.91	0.13	35,40,43,45	0
6	GOL	E	305	6/6	0.91	0.25	41,55,59,63	0
7	SO4	A	309	5/5	0.94	0.23	52,66,81,84	0
7	SO4	A	307	5/5	0.95	0.27	59,67,71,72	0
7	SO4	B	407	5/5	0.95	0.15	38,44,60,65	0
7	SO4	A	308	5/5	0.95	0.16	56,58,62,71	0
4	MG	B	402	1/1	0.97	0.12	19,19,19,19	0
4	MG	H	402	1/1	0.97	0.06	21,21,21,21	0
4	MG	G	302	1/1	0.98	0.09	22,22,22,22	0
3	GNP	B	401	32/32	0.98	0.14	12,18,23,24	0
3	GNP	D	401	32/32	0.98	0.12	17,23,29,31	0
3	GNP	E	301	32/32	0.98	0.10	16,23,30,31	0
3	GNP	F	401	32/32	0.98	0.12	18,22,27,30	0
3	GNP	A	301	32/32	0.98	0.14	13,20,23,24	0
4	MG	C	302	1/1	0.98	0.09	25,25,25,25	0
4	MG	D	402	1/1	0.98	0.08	22,22,22,22	0
7	SO4	H	405	5/5	0.98	0.16	58,60,68,70	0
3	GNP	H	401	32/32	0.99	0.12	16,22,25,26	0
4	MG	E	302	1/1	0.99	0.04	25,25,25,25	0
4	MG	F	402	1/1	0.99	0.08	24,24,24,24	0
4	MG	A	302	1/1	0.99	0.12	20,20,20,20	0
3	GNP	C	301	32/32	0.99	0.15	17,24,28,28	0
3	GNP	G	301	32/32	0.99	0.10	14,20,24,27	0

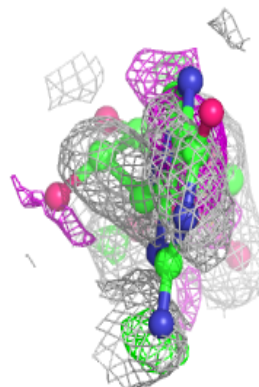
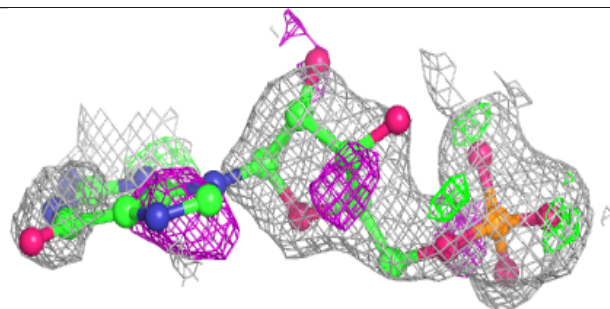
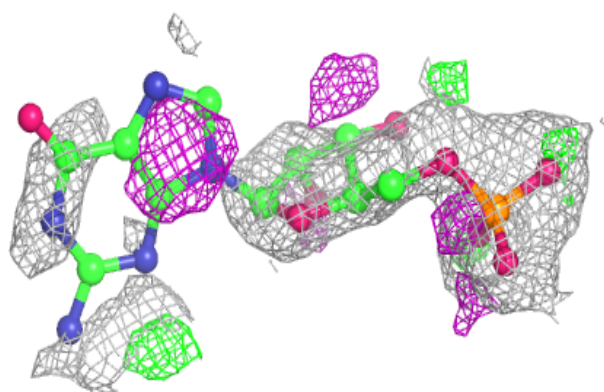
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around G A 303:

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

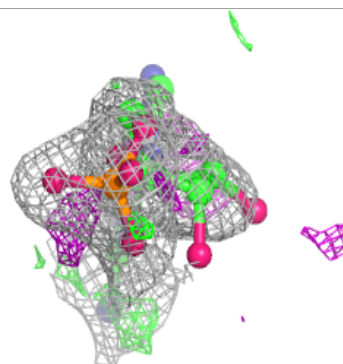
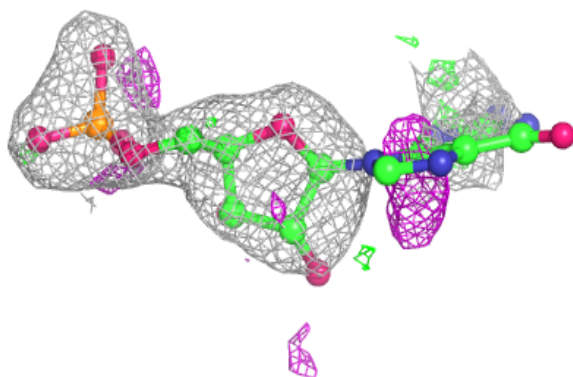
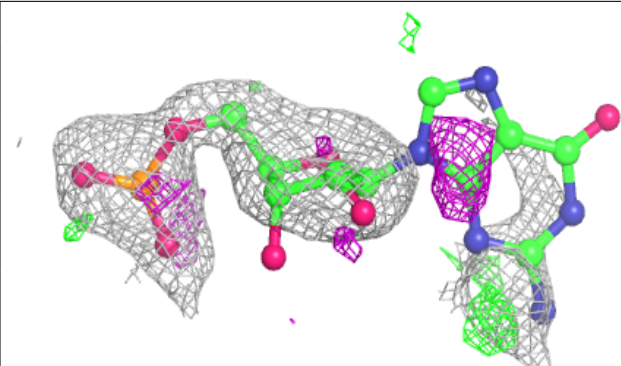
**Electron density around G E 303:**

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

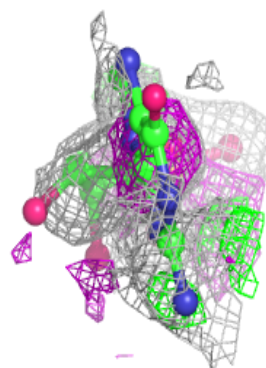
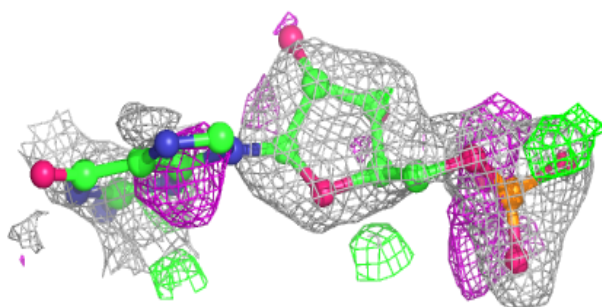
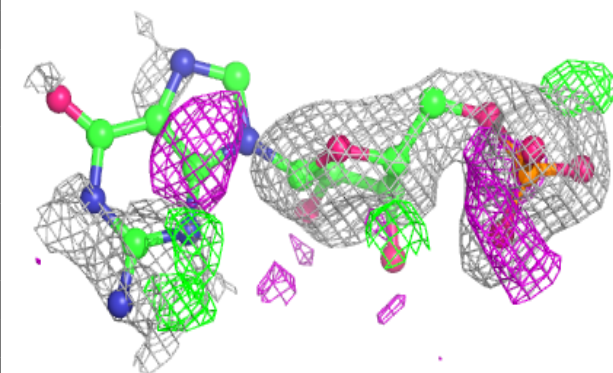


Electron density around G C 303:

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

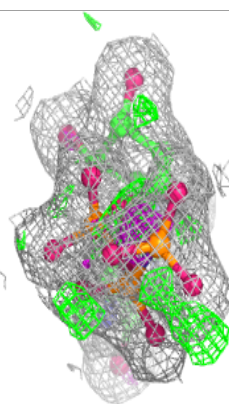
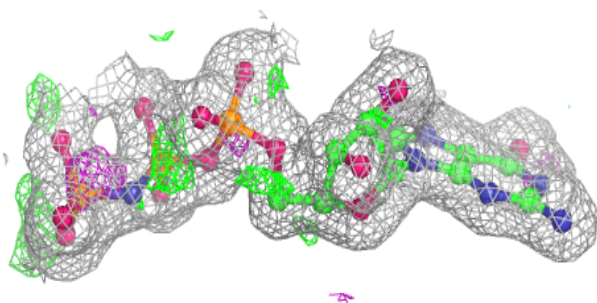
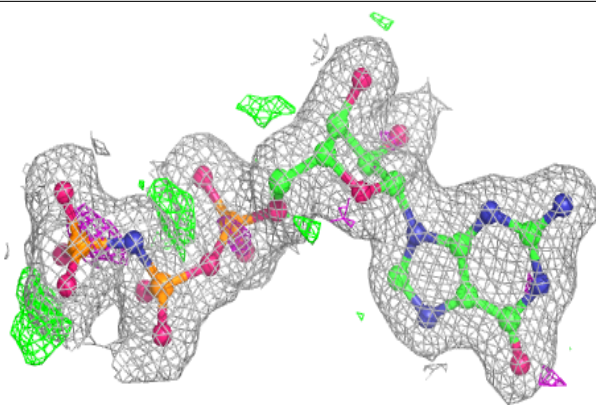
**Electron density around G G 303:**

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

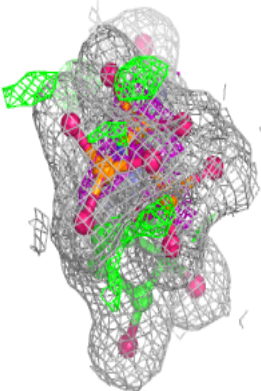
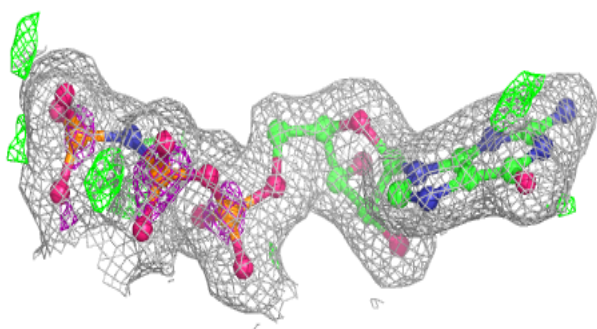
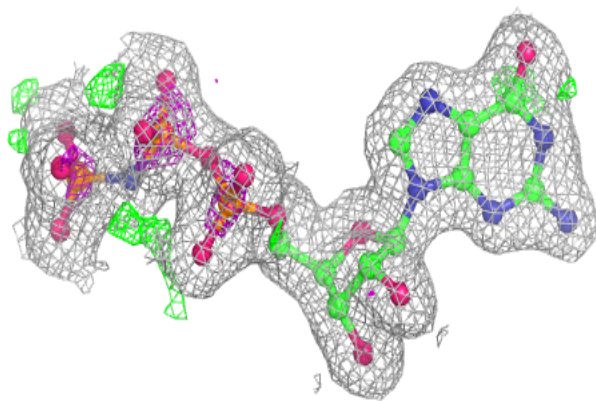


Electron density around GNP B 401:

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

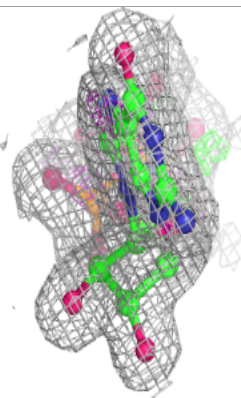
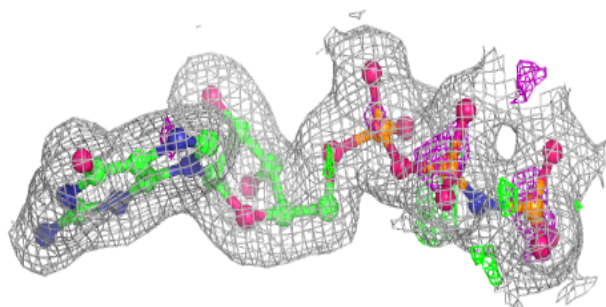
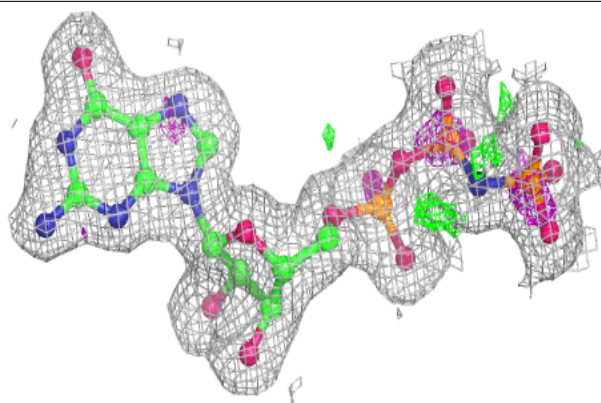
**Electron density around GNP D 401:**

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

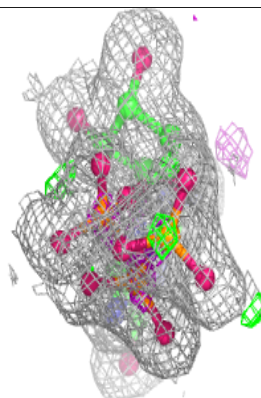
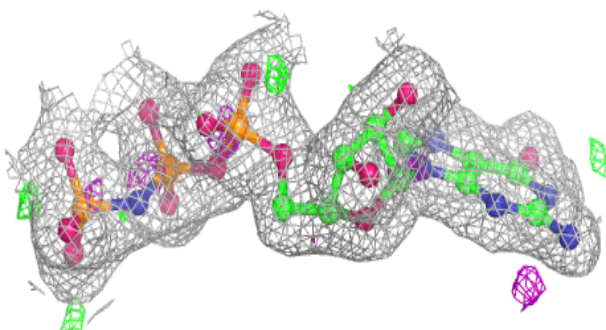
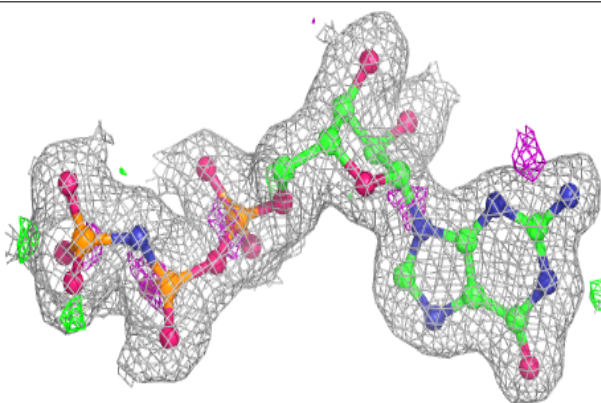


Electron density around GNP E 301:

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

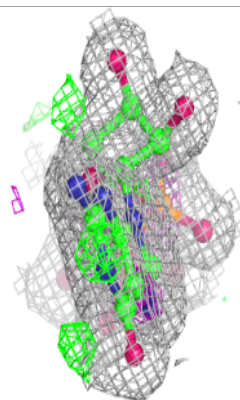
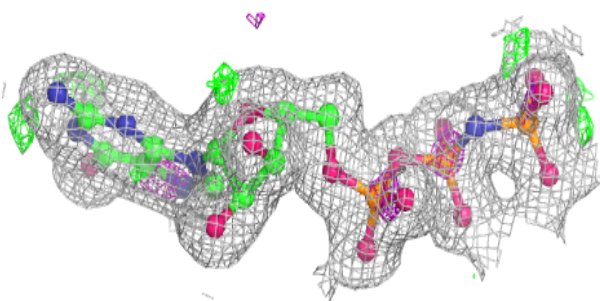
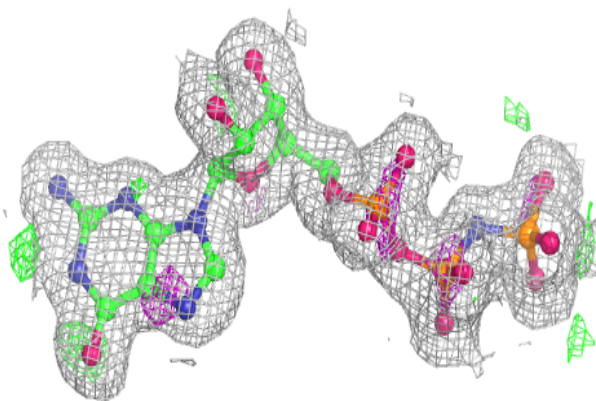
**Electron density around GNP F 401:**

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

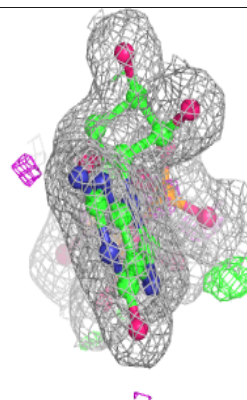
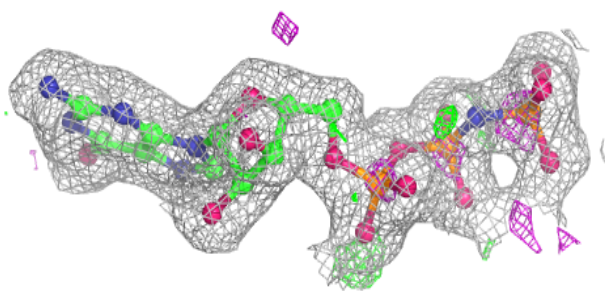
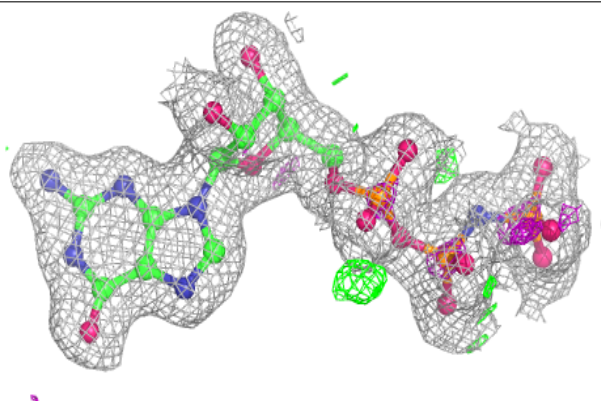


Electron density around GNP A 301:

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

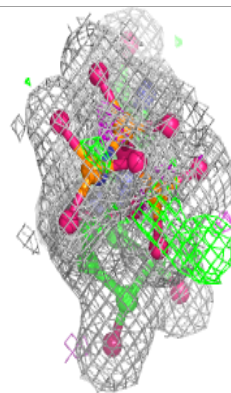
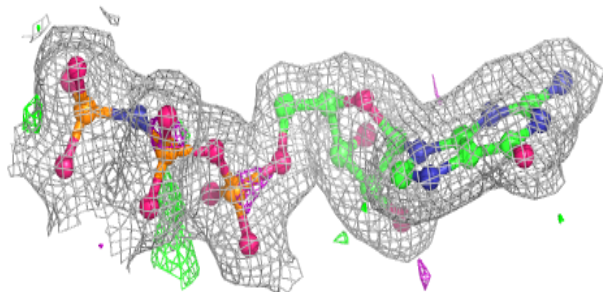
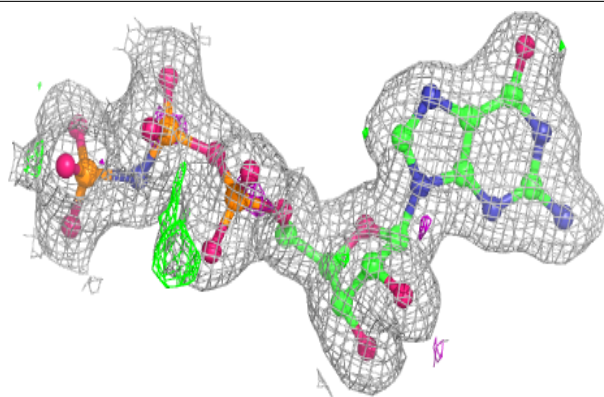
**Electron density around GNP H 401:**

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

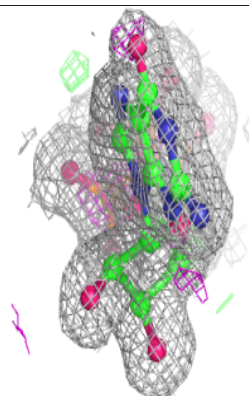
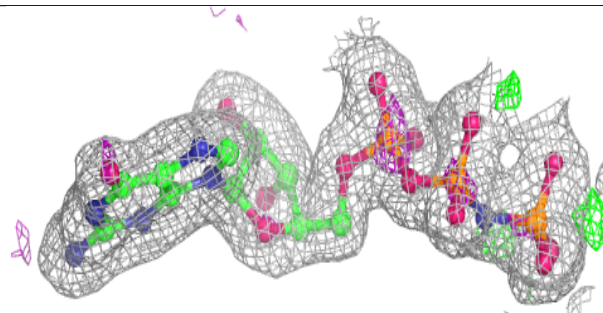
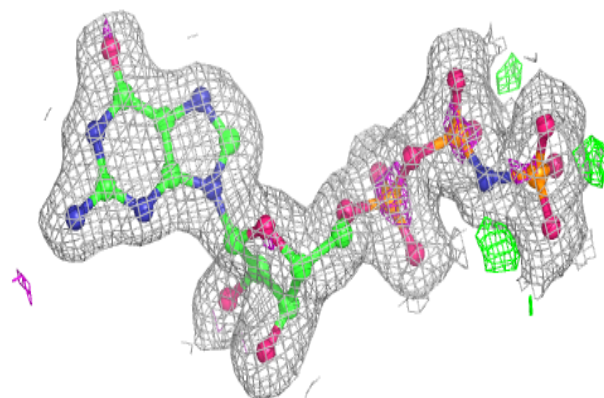


Electron density around GNP C 301:

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

**Electron density around GNP 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.