



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 05:51 AM UTC

PDB ID : 6Y32 / pdb_00006y32
Title : Structure of the GTPase heterodimer of human SRP54 and SRalpha
Authors : Juair, K.D.; Becker, M.M.M.; Wild, K.; Sinning, I.
Deposited on : 2020-02-17
Resolution : 2.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

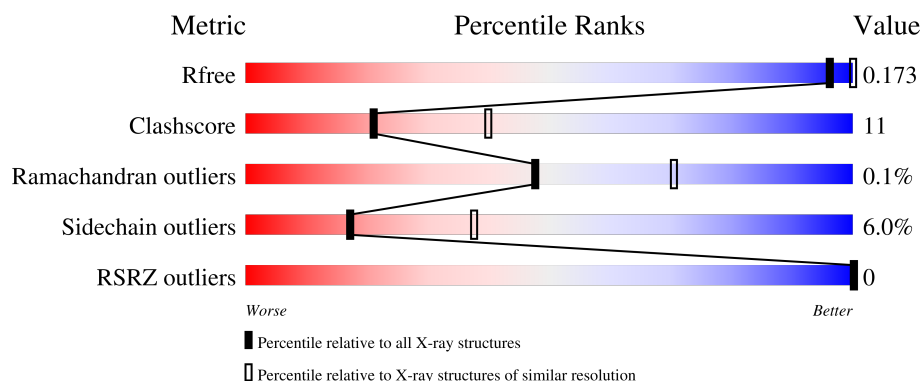
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	304	
1	C	304	
1	E	304	
1	G	304	
2	B	315	

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Mol	Chain	Length	Quality of chain
2	D	315	68%26%
2	F	315	70%24%
2	H	315	73%21% • 5%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 18356 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	0	0
			2120	1353	354	401	12			
1	C	268	Total	C	N	O	S	0	0	0
			2075	1326	348	390	11			
1	E	268	Total	C	N	O	S	0	0	0
			2075	1326	348	390	11			
1	G	267	Total	C	N	O	S	0	0	0
			2067	1322	346	388	11			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	initiating methionine	UNP P61011
A	-6	GLY	-	expression tag	UNP P61011
A	-5	HIS	-	expression tag	UNP P61011
A	-4	HIS	-	expression tag	UNP P61011
A	-3	HIS	-	expression tag	UNP P61011
A	-2	HIS	-	expression tag	UNP P61011
A	-1	HIS	-	expression tag	UNP P61011
A	0	HIS	-	expression tag	UNP P61011
C	-7	MET	-	initiating methionine	UNP P61011
C	-6	GLY	-	expression tag	UNP P61011
C	-5	HIS	-	expression tag	UNP P61011
C	-4	HIS	-	expression tag	UNP P61011
C	-3	HIS	-	expression tag	UNP P61011
C	-2	HIS	-	expression tag	UNP P61011
C	-1	HIS	-	expression tag	UNP P61011
C	0	HIS	-	expression tag	UNP P61011
E	-7	MET	-	initiating methionine	UNP P61011
E	-6	GLY	-	expression tag	UNP P61011
E	-5	HIS	-	expression tag	UNP P61011
E	-4	HIS	-	expression tag	UNP P61011
E	-3	HIS	-	expression tag	UNP P61011

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	HIS	-	expression tag	UNP P61011
E	-1	HIS	-	expression tag	UNP P61011
E	0	HIS	-	expression tag	UNP P61011
G	-7	MET	-	initiating methionine	UNP P61011
G	-6	GLY	-	expression tag	UNP P61011
G	-5	HIS	-	expression tag	UNP P61011
G	-4	HIS	-	expression tag	UNP P61011
G	-3	HIS	-	expression tag	UNP P61011
G	-2	HIS	-	expression tag	UNP P61011
G	-1	HIS	-	expression tag	UNP P61011
G	0	HIS	-	expression tag	UNP P61011

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	302	Total	C	N	O	S	0	0	0
			2297	1444	407	430	16			
2	D	305	Total	C	N	O	S	0	0	0
			2322	1460	412	434	16			
2	F	304	Total	C	N	O	S	0	0	0
			2316	1456	411	433	16			
2	H	300	Total	C	N	O	S	0	0	0
			2284	1436	404	428	16			

There are 32 discrepancies between the modelled and reference sequences:

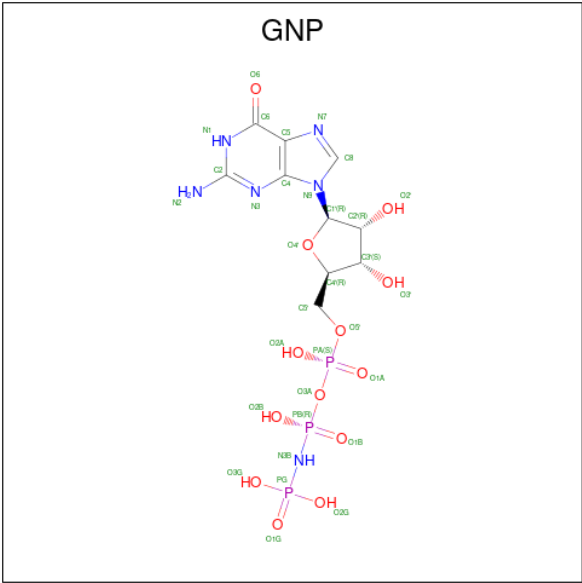
Chain	Residue	Modelled	Actual	Comment	Reference
B	330	MET	-	initiating methionine	UNP P08240
B	331	GLY	-	expression tag	UNP P08240
B	639	HIS	-	expression tag	UNP P08240
B	640	HIS	-	expression tag	UNP P08240
B	641	HIS	-	expression tag	UNP P08240
B	642	HIS	-	expression tag	UNP P08240
B	643	HIS	-	expression tag	UNP P08240
B	644	HIS	-	expression tag	UNP P08240
D	330	MET	-	initiating methionine	UNP P08240
D	331	GLY	-	expression tag	UNP P08240
D	639	HIS	-	expression tag	UNP P08240
D	640	HIS	-	expression tag	UNP P08240
D	641	HIS	-	expression tag	UNP P08240
D	642	HIS	-	expression tag	UNP P08240
D	643	HIS	-	expression tag	UNP P08240

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Chain	Residue	Modelled	Actual	Comment	Reference
D	644	HIS	-	expression tag	UNP P08240
F	330	MET	-	initiating methionine	UNP P08240
F	331	GLY	-	expression tag	UNP P08240
F	639	HIS	-	expression tag	UNP P08240
F	640	HIS	-	expression tag	UNP P08240
F	641	HIS	-	expression tag	UNP P08240
F	642	HIS	-	expression tag	UNP P08240
F	643	HIS	-	expression tag	UNP P08240
F	644	HIS	-	expression tag	UNP P08240
H	330	MET	-	initiating methionine	UNP P08240
H	331	GLY	-	expression tag	UNP P08240
H	639	HIS	-	expression tag	UNP P08240
H	640	HIS	-	expression tag	UNP P08240
H	641	HIS	-	expression tag	UNP P08240
H	642	HIS	-	expression tag	UNP P08240
H	643	HIS	-	expression tag	UNP P08240
H	644	HIS	-	expression tag	UNP P08240

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-GUANYLATE ESTER (CCD ID: GNP) (formula: C₁₀H₁₇N₆O₁₃P₃) (labeled as "Ligand of Interest" by depositor).



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		

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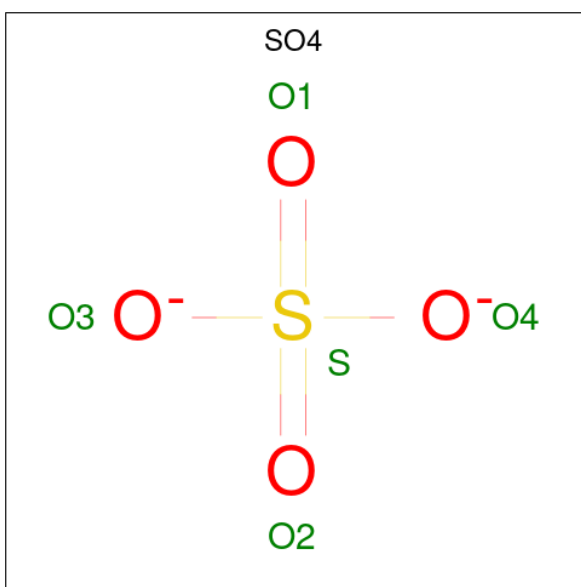
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	C	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
3	D	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
3	E	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
3	F	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
3	G	1	Total	C	N	O	P	0	0
			32	10	6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

- Molecule 4 is MAGNESIUM ION (CCD ID: 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 SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		
5	D	1	Total	O	S	0	0
			5	4	1		
5	F	1	Total	O	S	0	0
			5	4	1		
5	F	1	Total	O	S	0	0
			5	4	1		
5	F	1	Total	O	S	0	0
			5	4	1		
5	G	1	Total	O	S	0	0
			5	4	1		
5	H	1	Total	O	S	0	0
			5	4	1		
5	H	1	Total	O	S	0	0
			5	4	1		
5	H	1	Total	O	S	0	0
			5	4	1		
5	H	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is water.

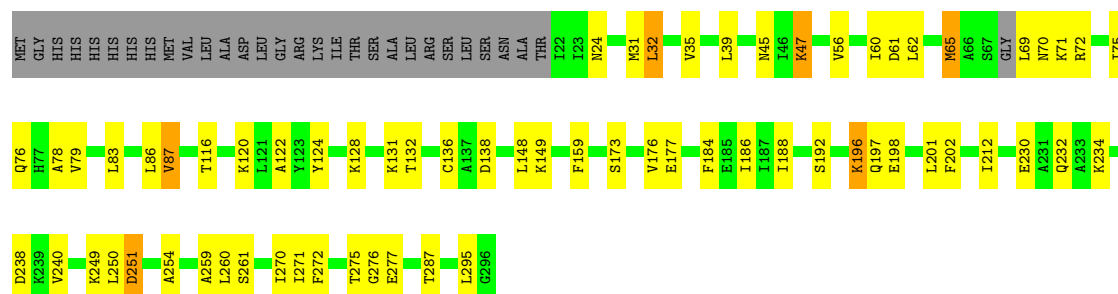
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	56	Total 56	O 56	0	0
6	B	80	Total 80	O 80	0	0
6	C	47	Total 47	O 47	0	0
6	D	50	Total 50	O 50	0	0
6	E	49	Total 49	O 49	0	0
6	F	55	Total 55	O 55	0	0
6	G	45	Total 45	O 45	0	0
6	H	84	Total 84	O 84	0	0

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.

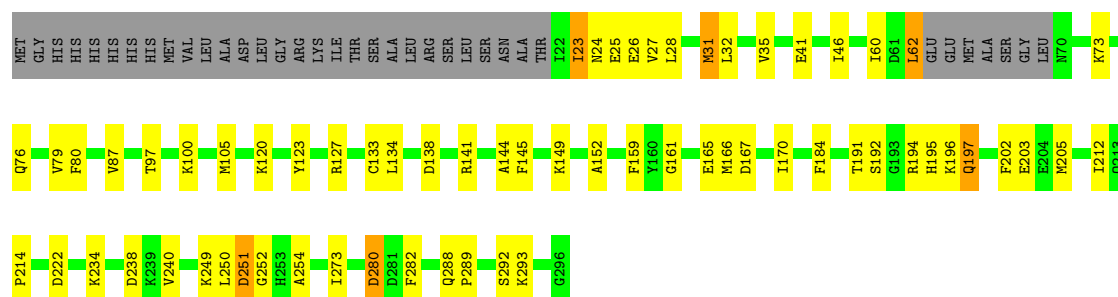
- Molecule 1: Signal recognition particle 54 kDa protein

Chain A: 



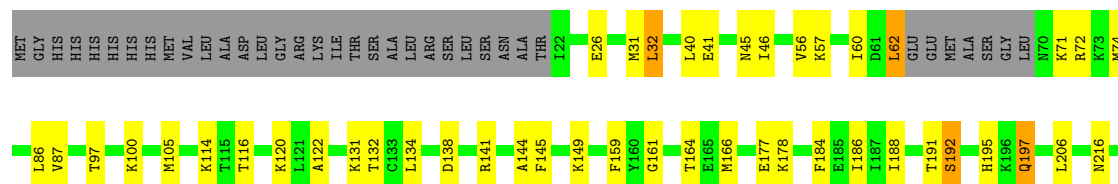
- Molecule 1: Signal recognition particle 54 kDa protein

Chain C: 



- Molecule 1: Signal recognition particle 54 kDa protein

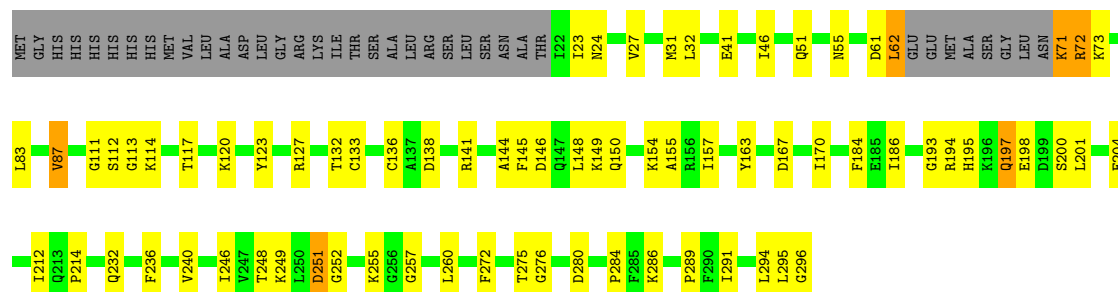
Chain E: 





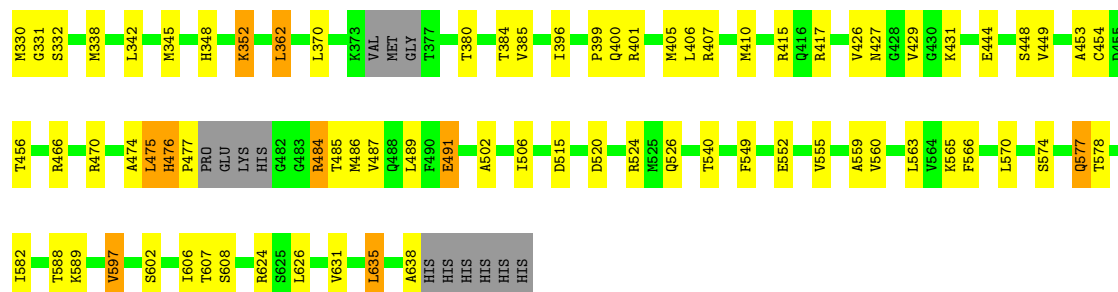
- Molecule 1: Signal recognition particle 54 kDa protein

Chain G: 63% 23% 12%



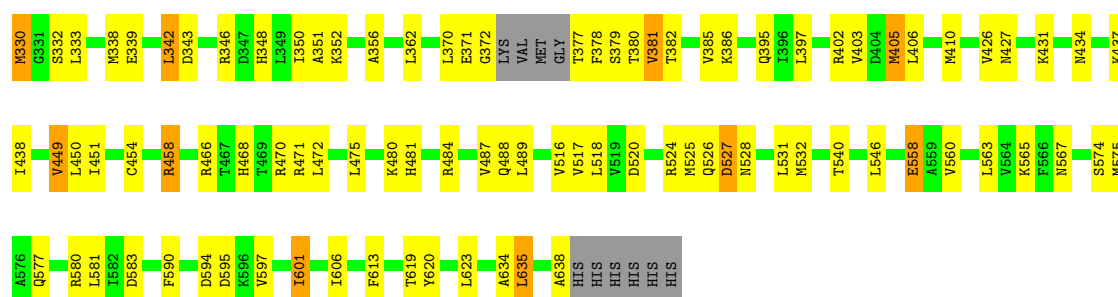
- Molecule 2: Signal recognition particle receptor subunit alpha

Chain B: 71% 22%



- Molecule 2: Signal recognition particle receptor subunit alpha

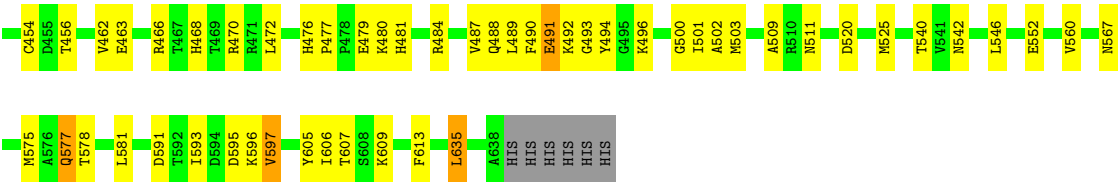
Chain D: 68% 26%



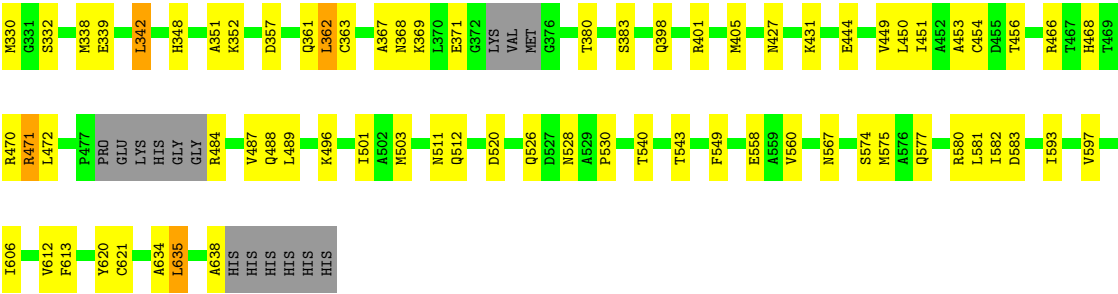
- Molecule 2: Signal recognition particle receptor subunit alpha

Chain F: 70% 24%





• Molecule 2: Signal recognition particle receptor subunit alpha



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	155.82Å 70.63Å 172.57Å 90.00° 116.81° 90.00°	Depositor
Resolution (Å)	49.55 – 2.60 49.55 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.5 (49.55-2.60) 99.4 (49.55-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.149 , 0.184 0.151 , 0.173	Depositor DCC
R_{free} test set	5161 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	44.9	Xtriage
Anisotropy	0.822	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 38.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	0.254 for h,-k,-h-l	Xtriage
Reported twinning fraction	0.260 for h,-k,-h-l	Depositor
Outliers	0 of 103282 reflections	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	18356	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.04% 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

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	0.34	0/2153	0.52	0/2899
1	C	0.32	0/2108	0.51	0/2839
1	E	0.33	0/2108	0.48	0/2839
1	G	0.30	0/2100	0.49	0/2828
2	B	0.40	0/2323	0.64	0/3136
2	D	0.38	0/2351	0.59	0/3177
2	F	0.36	0/2345	0.54	0/3168
2	H	0.38	0/2310	0.56	0/3120
All	All	0.36	0/17798	0.55	0/24006

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	2120	0	2165	42	0
1	C	2075	0	2123	41	0
1	E	2075	0	2123	44	0
1	G	2067	0	2117	51	0
2	B	2297	0	2361	62	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	2322	0	2379	64	0
2	F	2316	0	2376	57	0
2	H	2284	0	2346	46	0
3	A	32	0	13	3	0
3	B	32	0	13	1	0
3	C	32	0	13	1	0
3	D	32	0	13	2	0
3	E	32	0	13	3	0
3	F	32	0	13	1	0
3	G	32	0	13	6	0
3	H	32	0	13	2	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	B	15	0	0	1	0
5	C	5	0	0	0	0
5	D	5	0	0	1	0
5	F	15	0	0	0	0
5	G	5	0	0	1	0
5	H	25	0	0	1	0
6	A	56	0	0	0	0
6	B	80	0	0	6	0
6	C	47	0	0	2	0
6	D	50	0	0	4	0
6	E	49	0	0	4	0
6	F	55	0	0	4	0
6	G	45	0	0	2	0
6	H	84	0	0	6	0
All	All	18356	0	18094	395	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 11.

All (395) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:560:VAL:HG13	2:F:606:ILE:CD1	1.47	1.44
2:F:560:VAL:HG13	2:F:606:ILE:HD11	1.14	1.13
2:F:560:VAL:CG1	2:F:606:ILE:HD11	1.87	1.03
2:F:560:VAL:HG13	2:F:606:ILE:HD13	1.37	1.00
2:F:560:VAL:CG1	2:F:606:ILE:CD1	2.41	0.97
2:D:330:MET:N	2:D:330:MET:SD	2.45	0.89
2:H:580:ARG:NH2	2:H:583:ASP:OD2	2.13	0.81
2:F:597:VAL:HG12	2:F:635:LEU:HD13	1.61	0.81
2:B:540:THR:HG21	2:B:574:SER:HA	1.62	0.80
2:B:399:PRO:O	6:B:801:HOH:O	2.01	0.79
2:D:330:MET:N	2:D:332:SER:H	1.82	0.78
2:B:563:LEU:HD13	2:B:606:ILE:HD11	1.65	0.77
2:B:415:ARG:NH2	6:B:803:HOH:O	2.18	0.77
1:E:141:ARG:HB3	1:E:144:ALA:HB2	1.67	0.76
2:F:477:PRO:HG2	2:F:480:LYS:HB2	1.69	0.74
2:B:352:LYS:HD3	2:B:635:LEU:HD12	1.69	0.73
2:D:352:LYS:HE2	2:D:634:ALA:O	1.89	0.73
2:H:348:HIS:CE1	2:H:638:ALA:HB2	2.25	0.72
2:B:484:ARG:HE	2:B:485:THR:H	1.37	0.72
2:B:427:ASN:OD1	6:B:802:HOH:O	2.08	0.71
1:G:163:TYR:N	5:G:303:SO4:O4	2.23	0.71
1:E:141:ARG:NH1	3:E:301:GNP:O3G	2.25	0.69
1:G:255:LYS:NZ	1:G:296:GLY:OXT	2.26	0.69
1:G:24:ASN:OD1	1:G:62:LEU:HG	1.92	0.69
1:A:277:GLU:OE1	3:A:301:GNP:O2'	2.10	0.69
2:B:444:GLU:HG2	2:D:619:THR:HB	1.75	0.68
1:G:31:MET:HE1	1:G:72:ARG:HB3	1.76	0.68
2:F:342:LEU:HG	2:F:363:CYS:HB3	1.76	0.68
2:F:378:PHE:HB3	2:F:381:VAL:HB	1.75	0.68
2:H:342:LEU:HG	2:H:363:CYS:HB3	1.76	0.68
2:F:407:ARG:HH12	2:F:411:ASP:HB2	1.59	0.68
2:H:348:HIS:HE1	2:H:638:ALA:HB2	1.56	0.67
2:B:407:ARG:NE	1:C:203:GLU:HG3	2.10	0.67
1:C:141:ARG:HB3	1:C:144:ALA:HB2	1.76	0.66
2:F:417:ARG:NH1	6:F:803:HOH:O	2.29	0.66
2:F:330:MET:HB3	2:F:335:ARG:NH1	2.11	0.66
2:D:348:HIS:NE2	2:D:638:ALA:HB3	2.11	0.65
1:G:198:GLU:OE2	1:G:200:SER:OG	2.14	0.65
2:D:580:ARG:NH2	2:D:583:ASP:OD2	2.30	0.65
2:D:527:ASP:OD2	6:D:801:HOH:O	2.15	0.65
2:H:580:ARG:NH1	6:H:805:HOH:O	2.30	0.64
2:B:475:LEU:HD13	2:D:471:ARG:HD3	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:87:VAL:HG21	1:G:291:ILE:HD11	1.79	0.64
2:B:475:LEU:HB3	2:B:477:PRO:HG3	1.79	0.64
2:F:330:MET:HB3	2:F:335:ARG:HH11	1.63	0.64
1:G:154:LYS:NZ	6:G:402:HOH:O	2.24	0.64
2:F:362:LEU:HD11	2:F:396:ILE:HD11	1.81	0.63
2:H:401:ARG:NH1	6:H:807:HOH:O	2.32	0.63
1:G:71:LYS:NZ	1:G:72:ARG:HG3	2.14	0.63
2:B:624:ARG:NH1	5:B:704:SO4:O1	2.31	0.63
1:G:141:ARG:NH1	3:G:301:GNP:O2G	2.31	0.63
1:C:165:GLU:OE2	6:C:401:HOH:O	2.16	0.62
2:D:377:THR:HB	2:D:380:THR:OG1	2.00	0.62
1:A:24:ASN:HB3	1:A:62:LEU:HD13	1.82	0.61
1:G:248:THR:OG1	1:G:249:LYS:HG3	2.02	0.60
2:H:560:VAL:HG13	2:H:606:ILE:HG21	1.83	0.60
1:A:47:LYS:NZ	1:A:230:GLU:OE2	2.33	0.60
2:D:397:LEU:HD21	2:D:601:ILE:HD12	1.83	0.60
2:F:332:SER:O	2:F:336:GLU:HB2	2.02	0.60
1:C:196:LYS:NZ	2:D:594:ASP:OD2	2.25	0.60
1:A:249:LYS:NZ	2:B:427:ASN:O	2.34	0.60
1:A:251:ASP:HB3	1:A:275:THR:O	2.02	0.59
1:E:206:LEU:HB2	1:E:239:LYS:HE3	1.85	0.59
2:D:560:VAL:HG13	2:D:606:ILE:HG21	1.83	0.59
2:F:454:CYS:HB2	2:F:520:ASP:O	2.03	0.59
2:D:574:SER:OG	2:D:575:MET:N	2.34	0.58
2:B:453:ALA:HB2	2:B:489:LEU:HD11	1.84	0.58
2:B:606:ILE:HD12	2:B:607:THR:N	2.18	0.58
2:F:450:LEU:HD11	2:F:490:PHE:HB2	1.84	0.58
2:H:496:LYS:HB2	2:H:501:ILE:HD11	1.85	0.58
2:F:496:LYS:HB2	2:F:501:ILE:HD11	1.86	0.58
1:C:23:ILE:HD12	1:C:23:ILE:H	1.68	0.58
1:C:120:LYS:HD3	1:C:282:PHE:HB2	1.85	0.58
2:H:597:VAL:HG12	2:H:635:LEU:HD13	1.87	0.57
1:G:83:LEU:HD11	1:G:295:LEU:HD13	1.86	0.57
2:H:454:CYS:HB2	2:H:520:ASP:O	2.05	0.57
2:H:466:ARG:HD3	2:H:470:ARG:NH1	2.19	0.57
2:H:369:LYS:NZ	6:H:802:HOH:O	2.24	0.56
1:E:32:LEU:HD11	1:E:56:VAL:HG11	1.87	0.56
2:F:491:GLU:HG2	2:F:493:GLY:H	1.70	0.56
2:D:466:ARG:HD3	2:D:470:ARG:CZ	2.36	0.56
2:D:352:LYS:HZ2	2:D:638:ALA:HB2	1.70	0.56
1:G:167:ASP:HB3	1:G:170:ILE:HD12	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:400:GLN:OE1	2:B:608:SER:O	2.24	0.56
2:B:560:VAL:HG13	2:B:606:ILE:HG21	1.88	0.56
1:C:145:PHE:CE1	1:C:161:GLY:HA3	2.41	0.56
1:G:251:ASP:HB3	1:G:275:THR:O	2.06	0.55
3:A:301:GNP:O4'	2:B:427:ASN:HB3	2.06	0.55
2:H:549:PHE:HB2	2:H:582:ILE:HD13	1.87	0.55
2:B:577:GLN:CD	2:B:577:GLN:H	2.12	0.55
1:A:86:LEU:HB3	1:A:260:LEU:HD22	1.88	0.55
1:C:250:LEU:HA	1:C:254:ALA:HB3	1.89	0.55
2:F:450:LEU:HD12	2:F:488:GLN:O	2.07	0.55
1:E:195:HIS:CD2	1:E:197:GLN:HG2	2.41	0.55
1:A:250:LEU:HA	1:A:254:ALA:HB3	1.89	0.55
1:E:72:ARG:NH1	6:E:404:HOH:O	2.30	0.55
2:D:352:LYS:HD3	2:D:635:LEU:HD12	1.89	0.54
2:F:462:VAL:HG11	2:F:491:GLU:HG3	1.90	0.54
1:E:31:MET:SD	1:E:72:ARG:HD3	2.48	0.54
2:B:429:VAL:HG12	2:B:552:GLU:HG2	1.90	0.54
2:B:484:ARG:NE	2:B:485:THR:H	2.04	0.54
2:F:466:ARG:HD3	2:F:470:ARG:CZ	2.37	0.54
3:G:301:GNP:O4'	2:H:427:ASN:HB3	2.08	0.54
1:E:288:GLN:O	1:E:292:SER:OG	2.17	0.54
2:F:405:MET:HE1	2:F:613:PHE:HE2	1.73	0.54
2:B:417:ARG:NH1	6:B:812:HOH:O	2.39	0.53
1:C:288:GLN:O	1:C:292:SER:OG	2.21	0.53
1:G:136:CYS:HB2	1:G:148:LEU:HD22	1.89	0.53
1:C:35:VAL:HG11	1:C:79:VAL:HG11	1.90	0.53
1:C:105:MET:HE3	1:C:191:THR:HG21	1.90	0.53
1:E:251:ASP:N	1:E:251:ASP:OD1	2.41	0.53
1:A:87:VAL:HG13	1:A:260:LEU:HD21	1.90	0.53
2:D:524:ARG:HH11	2:D:531:LEU:HB3	1.73	0.53
2:H:352:LYS:NZ	2:H:634:ALA:O	2.41	0.53
5:H:706:SO4:O1	6:H:801:HOH:O	2.15	0.53
1:E:62:LEU:HD21	1:E:71:LYS:HE3	1.91	0.53
1:G:51:GLN:O	1:G:55:ASN:ND2	2.41	0.53
2:H:621:CYS:HB2	6:H:817:HOH:O	2.09	0.52
1:C:250:LEU:HD12	1:C:273:ILE:HB	1.91	0.52
1:E:45:ASN:ND2	1:E:230:GLU:OE2	2.41	0.52
1:E:138:ASP:CG	1:E:192:SER:HB3	2.35	0.52
1:G:71:LYS:HZ2	1:G:72:ARG:HG3	1.73	0.52
1:G:123:TYR:CZ	1:G:127:ARG:HD2	2.45	0.52
1:A:198:GLU:HG2	1:A:201:LEU:HD12	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:560:VAL:HA	2:D:606:ILE:HD13	1.92	0.51
2:H:453:ALA:HB2	2:H:489:LEU:HD11	1.92	0.51
2:H:540:THR:HG21	2:H:575:MET:HG2	1.92	0.51
2:D:613:PHE:HB3	2:D:623:LEU:HD11	1.92	0.51
1:A:35:VAL:HG21	1:A:79:VAL:HG21	1.92	0.51
1:E:248:THR:OG1	1:E:249:LYS:HG3	2.10	0.51
1:G:212:ILE:HG13	1:G:214:PRO:HD3	1.92	0.51
1:C:60:ILE:HG22	1:C:62:LEU:HB2	1.92	0.51
1:C:138:ASP:OD1	1:C:192:SER:HB3	2.11	0.51
2:B:348:HIS:NE2	2:B:638:ALA:HB3	2.26	0.51
1:G:249:LYS:NZ	2:H:427:ASN:O	2.41	0.51
1:E:197:GLN:OE1	1:E:232:GLN:NE2	2.44	0.51
2:H:540:THR:CG2	2:H:575:MET:HG2	2.41	0.51
2:F:456:THR:HG21	2:F:501:ILE:HD12	1.93	0.51
2:F:540:THR:HG21	2:F:575:MET:H	1.76	0.51
1:G:198:GLU:HG2	1:G:201:LEU:HG	1.92	0.51
2:D:352:LYS:HE3	2:D:595:ASP:O	2.11	0.50
1:E:122:ALA:HB2	1:E:188:ILE:HD12	1.93	0.50
1:E:138:ASP:OD1	1:E:192:SER:HB3	2.11	0.50
2:D:526:GLN:HB2	2:D:565:LYS:HB2	1.93	0.50
3:C:301:GNP:O4'	2:D:427:ASN:HB3	2.11	0.50
1:A:136:CYS:HB2	1:A:148:LEU:HD13	1.93	0.50
1:E:230:GLU:HB2	1:E:261:SER:HB3	1.93	0.50
1:G:155:ALA:O	1:G:157:ILE:HG13	2.11	0.50
1:A:32:LEU:HD11	1:A:56:VAL:HG11	1.92	0.50
2:D:450:LEU:HD12	2:D:488:GLN:O	2.11	0.50
2:H:357:ASP:OD1	2:H:361:GLN:NE2	2.43	0.50
2:H:450:LEU:HD12	2:H:488:GLN:O	2.12	0.50
2:D:352:LYS:NZ	2:D:595:ASP:OD2	2.45	0.49
3:E:301:GNP:H5'1	6:E:406:HOH:O	2.12	0.49
2:H:398:GLN:O	2:H:398:GLN:HG2	2.13	0.49
2:B:526:GLN:HB2	2:B:565:LYS:HB2	1.94	0.49
1:E:105:MET:HE3	1:E:191:THR:HG21	1.92	0.49
1:G:112:SER:O	1:G:249:LYS:HE3	2.12	0.49
2:B:577:GLN:NE2	6:B:806:HOH:O	2.34	0.49
1:C:23:ILE:O	1:C:27:VAL:HG23	2.12	0.49
2:F:552:GLU:OE1	6:F:801:HOH:O	2.20	0.49
1:G:145:PHE:O	1:G:149:LYS:HB2	2.13	0.49
2:H:352:LYS:HZ2	2:H:638:ALA:H	1.60	0.49
1:A:138:ASP:OD1	1:A:192:SER:HB3	2.13	0.49
2:B:577:GLN:HA	1:E:292:SER:HB3	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:87:VAL:HG21	1:E:287:THR:HG23	1.94	0.49
1:G:286:LYS:O	1:G:289:PRO:HD2	2.12	0.49
2:F:479:GLU:HG2	2:F:480:LYS:N	2.27	0.49
2:B:602:SER:O	2:B:606:ILE:HG13	2.12	0.49
2:H:574:SER:OG	2:H:575:MET:N	2.44	0.49
1:A:87:VAL:HG21	1:A:287:THR:HG23	1.94	0.49
2:F:357:ASP:O	2:F:361:GLN:HG3	2.12	0.49
1:A:31:MET:HE1	1:A:76:GLN:HG3	1.94	0.49
1:E:149:LYS:HG3	1:E:159:PHE:CE2	2.48	0.49
1:E:216:ASN:OD1	1:E:243:ALA:HB3	2.12	0.49
2:B:484:ARG:NE	2:B:485:THR:O	2.46	0.48
2:B:549:PHE:HB2	2:B:582:ILE:HD13	1.95	0.48
1:G:113:GLY:HA2	3:G:301:GNP:H5'2	1.95	0.48
2:B:338:MET:O	2:B:342:LEU:HG	2.12	0.48
1:C:97:THR:HB	1:C:100:LYS:HG2	1.95	0.48
2:F:362:LEU:HD12	2:F:362:LEU:HA	1.66	0.48
2:B:477:PRO:C	2:B:486:MET:HE2	2.39	0.48
1:G:87:VAL:CG2	1:G:291:ILE:HD11	2.43	0.48
1:E:31:MET:HE2	1:E:72:ARG:HB3	1.96	0.48
2:F:387:GLN:O	2:F:391:GLU:HG3	2.12	0.48
2:D:454:CYS:HB2	2:D:520:ASP:O	2.13	0.48
2:F:407:ARG:NH1	2:F:411:ASP:HB2	2.27	0.48
1:G:195:HIS:CD2	1:G:197:GLN:HG2	2.49	0.48
2:B:555:VAL:HG13	2:B:559:ALA:HB2	1.95	0.48
1:C:249:LYS:NZ	2:D:427:ASN:O	2.46	0.48
2:F:406:LEU:O	2:F:410:MET:HG3	2.13	0.48
1:G:46:ILE:HG21	2:H:351:ALA:HB2	1.96	0.48
1:A:196:LYS:HB3	1:A:232:GLN:OE1	2.14	0.48
2:B:474:ALA:HB3	2:D:475:LEU:HD23	1.96	0.48
1:A:32:LEU:HD12	1:A:32:LEU:HA	1.62	0.47
2:B:477:PRO:HD3	6:B:804:HOH:O	2.14	0.47
2:D:382:THR:O	2:D:386:LYS:HB2	2.14	0.47
1:G:204:GLU:OE1	6:G:401:HOH:O	2.20	0.47
1:A:131:LYS:HD3	1:A:184:PHE:HA	1.97	0.47
1:C:46:ILE:HG21	2:D:351:ALA:HB2	1.96	0.47
1:C:280:ASP:OD1	1:C:280:ASP:N	2.41	0.47
1:A:271:ILE:HG13	1:A:272:PHE:CD2	2.50	0.47
3:A:301:GNP:H8	3:A:301:GNP:H5'2	1.95	0.47
2:D:370:LEU:HD23	2:D:370:LEU:HA	1.67	0.47
1:E:288:GLN:HG3	1:E:289:PRO:HD3	1.96	0.47
1:C:167:ASP:HB3	1:C:170:ILE:HD12	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:138:ASP:HB3	1:G:144:ALA:HB1	1.96	0.47
2:F:468:HIS:O	2:F:472:LEU:HG	2.15	0.47
1:C:166:MET:HB3	1:C:166:MET:HE2	1.56	0.47
2:D:338:MET:HG2	2:D:342:LEU:HD22	1.96	0.47
2:F:567:ASN:OD1	2:F:581:LEU:HB3	2.15	0.47
2:F:577:GLN:O	2:F:578:THR:OG1	2.27	0.47
1:G:194:ARG:NH1	1:G:201:LEU:O	2.48	0.47
1:C:60:ILE:HG22	1:C:62:LEU:H	1.78	0.47
2:F:450:LEU:HD23	2:F:509:ALA:HB2	1.97	0.47
2:B:338:MET:HE3	2:B:342:LEU:HD11	1.95	0.46
2:D:378:PHE:O	2:D:381:VAL:HG23	2.14	0.46
2:D:517:VAL:O	2:D:518:LEU:HD23	2.14	0.46
2:B:466:ARG:HD3	2:B:470:ARG:CZ	2.45	0.46
1:E:221:MET:HG3	1:E:247:VAL:HG22	1.98	0.46
2:D:397:LEU:HD23	2:D:397:LEU:HA	1.76	0.46
2:D:405:MET:HE3	2:D:405:MET:HB2	1.74	0.46
2:H:338:MET:SD	2:H:367:ALA:HA	2.56	0.46
1:A:176:VAL:HG21	1:A:212:ILE:HB	1.97	0.46
2:B:400:GLN:HB3	2:B:401:ARG:H	1.52	0.46
1:C:196:LYS:HA	1:C:202:PHE:CE2	2.50	0.46
2:F:596:LYS:NZ	6:F:814:HOH:O	2.48	0.46
2:B:502:ALA:O	2:B:506:ILE:HG13	2.16	0.46
2:D:468:HIS:O	2:D:472:LEU:HG	2.16	0.46
1:E:131:LYS:HG2	1:E:184:PHE:CD1	2.51	0.46
1:E:134:LEU:HD23	1:E:188:ILE:HB	1.97	0.46
1:E:177:GLU:HG3	1:E:178:LYS:N	2.31	0.46
2:B:407:ARG:HE	1:C:203:GLU:HG3	1.80	0.46
1:E:132:THR:HG22	1:E:186:ILE:HB	1.97	0.46
1:E:197:GLN:HG3	2:F:591:ASP:O	2.16	0.46
2:F:595:ASP:O	2:F:597:VAL:N	2.48	0.46
2:F:500:GLY:HA2	2:F:503:MET:HE3	1.97	0.45
1:A:47:LYS:HB3	1:A:47:LYS:HE3	1.76	0.45
1:A:234:LYS:HE2	1:A:238:ASP:OD2	2.17	0.45
2:F:454:CYS:SG	2:F:502:ALA:HB2	2.56	0.45
1:A:198:GLU:HG2	1:A:201:LEU:HG	1.99	0.45
1:A:198:GLU:HG2	1:A:201:LEU:CD1	2.46	0.45
2:B:407:ARG:CZ	1:C:203:GLU:HG3	2.46	0.45
1:C:251:ASP:OD1	1:C:251:ASP:N	2.49	0.45
2:D:370:LEU:CD1	2:D:385:VAL:HG12	2.46	0.45
2:F:451:ILE:O	2:F:489:LEU:HD12	2.16	0.45
1:C:123:TYR:CZ	1:C:127:ARG:HD2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:405:MET:HE1	2:H:613:PHE:HE2	1.81	0.45
2:B:345:MET:HA	2:B:345:MET:HE2	1.99	0.45
1:C:149:LYS:HE3	1:C:159:PHE:CZ	2.52	0.45
2:D:480:LYS:HD2	2:D:481:HIS:CE1	2.51	0.45
2:B:597:VAL:CG1	2:B:631:VAL:HG13	2.47	0.45
1:E:120:LYS:HD3	1:E:282:PHE:HB2	1.99	0.45
1:A:196:LYS:HA	1:A:202:PHE:CE2	2.52	0.45
1:G:133:CYS:HB3	1:G:184:PHE:CE1	2.51	0.45
2:B:555:VAL:CG1	2:B:559:ALA:HB2	2.47	0.45
1:E:149:LYS:HG3	1:E:159:PHE:CZ	2.52	0.45
1:G:146:ASP:O	1:G:150:GLN:HG3	2.17	0.44
2:H:348:HIS:HE1	2:H:638:ALA:CB	2.26	0.44
2:D:377:THR:O	2:D:381:VAL:HG22	2.17	0.44
2:B:454:CYS:HB2	2:B:520:ASP:O	2.17	0.44
2:D:352:LYS:HZ1	2:D:595:ASP:CG	2.25	0.44
2:H:451:ILE:O	2:H:489:LEU:HD12	2.17	0.44
1:E:57:LYS:O	1:E:60:ILE:HG22	2.16	0.44
1:E:166:MET:HE3	1:E:166:MET:HB3	1.80	0.44
1:A:122:ALA:HB2	1:A:188:ILE:HD12	2.00	0.44
1:A:132:THR:HG22	1:A:186:ILE:HB	1.99	0.44
2:B:444:GLU:CG	2:D:619:THR:HB	2.47	0.44
2:F:421:VAL:HG22	2:F:546:LEU:HB3	2.00	0.44
2:H:468:HIS:O	2:H:472:LEU:HG	2.18	0.44
1:C:222:ASP:C	1:C:222:ASP:OD1	2.61	0.44
1:E:265:ALA:O	6:E:401:HOH:O	2.21	0.44
2:H:431:LYS:NZ	3:H:701:GNP:O3G	2.48	0.44
1:G:132:THR:HG22	1:G:186:ILE:HB	2.00	0.44
1:A:31:MET:SD	1:A:72:ARG:HA	2.58	0.43
2:B:426:VAL:HA	2:B:524:ARG:O	2.18	0.43
1:A:83:LEU:HD11	1:A:295:LEU:HD13	2.01	0.43
2:F:357:ASP:OD1	2:F:361:GLN:NE2	2.50	0.43
1:E:32:LEU:HD12	1:E:32:LEU:HA	1.71	0.43
1:E:86:LEU:HD23	1:E:86:LEU:HA	1.81	0.43
1:A:173:SER:O	1:A:177:GLU:HG2	2.18	0.43
1:A:259:ALA:HB1	1:A:270:ILE:HD11	1.99	0.43
1:C:28:LEU:O	1:C:31:MET:N	2.51	0.43
1:C:149:LYS:HE3	1:C:159:PHE:HZ	1.82	0.43
2:D:528:ASN:O	2:D:532:MET:HG2	2.19	0.43
2:F:476:HIS:HD2	2:F:481:HIS:CE1	2.37	0.43
1:G:71:LYS:HB3	1:G:71:LYS:HE3	1.76	0.43
2:F:463:GLU:OE2	2:F:494:TYR:OH	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:368:ASN:O	2:H:371:GLU:HG2	2.19	0.43
2:D:426:VAL:HG22	6:D:802:HOH:O	2.19	0.43
2:D:434:ASN:O	2:D:438:ILE:HG13	2.19	0.43
2:B:606:ILE:HD12	2:B:607:THR:HG23	2.00	0.43
1:E:62:LEU:HD11	1:E:71:LYS:HE3	2.01	0.43
1:E:116:THR:HG22	1:E:120:LYS:HE3	2.01	0.43
1:G:280:ASP:OD1	1:G:280:ASP:N	2.47	0.43
2:H:362:LEU:HD12	2:H:362:LEU:HA	1.79	0.43
2:D:431:LYS:NZ	3:D:701:GNP:O2G	2.47	0.43
2:D:449:VAL:HA	2:D:516:VAL:O	2.18	0.43
2:H:577:GLN:HB2	6:H:806:HOH:O	2.19	0.43
1:C:149:LYS:HG3	1:C:159:PHE:CZ	2.53	0.43
2:D:525:MET:HB3	6:D:801:HOH:O	2.18	0.43
2:D:590:PHE:CZ	2:D:597:VAL:HG23	2.54	0.43
2:F:405:MET:HE3	2:F:405:MET:HB2	1.77	0.43
1:G:257:GLY:HA2	1:G:294:LEU:HD11	1.99	0.43
2:B:491:GLU:C	2:B:491:GLU:CD	2.86	0.43
1:C:28:LEU:HA	1:C:31:MET:HE2	2.00	0.42
2:D:563:LEU:HD23	2:D:563:LEU:HA	1.90	0.42
1:G:61:ASP:OD2	1:G:62:LEU:N	2.52	0.42
1:G:141:ARG:HB3	1:G:144:ALA:HB2	2.01	0.42
1:G:32:LEU:HD12	1:G:32:LEU:HA	1.62	0.42
1:A:149:LYS:HG3	1:A:159:PHE:CE2	2.55	0.42
2:B:405:MET:HE3	2:B:405:MET:HB2	1.85	0.42
2:D:350:ILE:HG12	2:D:356:ALA:HA	2.01	0.42
2:F:415:ARG:NH2	6:F:808:HOH:O	2.36	0.42
1:G:286:LYS:HD3	1:G:286:LYS:HA	1.84	0.42
2:B:362:LEU:HD12	2:B:362:LEU:HA	1.70	0.42
1:A:65:MET:HB2	1:A:71:LYS:NZ	2.35	0.42
1:C:134:LEU:HD12	1:C:152:ALA:HB2	2.01	0.42
1:A:39:LEU:HA	1:A:39:LEU:HD23	1.75	0.42
2:B:362:LEU:HD11	2:B:396:ILE:CD1	2.50	0.42
1:G:111:GLY:HA2	3:G:301:GNP:HNB3	1.84	0.42
3:G:301:GNP:O3'	3:H:701:GNP:O1G	2.34	0.42
2:H:330:MET:HA	2:H:332:SER:N	2.34	0.42
1:C:194:ARG:HD2	1:C:205:MET:SD	2.59	0.42
1:C:195:HIS:CD2	1:C:197:GLN:HG2	2.55	0.42
2:D:470:ARG:NH1	5:D:703:SO4:O3	2.53	0.42
2:D:567:ASN:OD1	2:D:581:LEU:HB3	2.20	0.42
1:E:145:PHE:CE1	1:E:161:GLY:HA3	2.55	0.42
1:A:230:GLU:HB2	1:A:261:SER:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:288:GLN:HB3	1:C:289:PRO:HD3	2.02	0.42
2:D:451:ILE:O	2:D:489:LEU:HD12	2.19	0.42
6:E:430:HOH:O	2:F:525:MET:HE1	2.20	0.42
1:A:120:LYS:NZ	1:A:276:GLY:O	2.37	0.41
2:B:330:MET:HA	2:B:331:GLY:C	2.45	0.41
2:B:566:PHE:O	2:B:570:LEU:HG	2.20	0.41
2:B:380:THR:O	2:B:384:THR:N	2.50	0.41
2:B:448:SER:O	2:B:515:ASP:HB2	2.20	0.41
2:D:406:LEU:O	2:D:410:MET:HG3	2.20	0.41
1:E:40:LEU:HD23	1:E:40:LEU:HA	1.91	0.41
1:G:232:GLN:O	1:G:236:PHE:HD1	2.03	0.41
2:H:528:ASN:OD1	2:H:530:PRO:HD2	2.20	0.41
2:H:612:VAL:HG12	2:H:613:PHE:CD2	2.55	0.41
1:C:80:PHE:HZ	1:C:288:GLN:HG2	1.85	0.41
2:F:397:LEU:O	2:F:399:PRO:HD3	2.20	0.41
1:E:114:LYS:NZ	3:E:301:GNP:O1B	2.45	0.41
2:B:476:HIS:N	2:B:477:PRO:HD3	2.35	0.41
1:C:133:CYS:HB3	1:C:184:PHE:CE1	2.56	0.41
1:E:40:LEU:HD21	1:E:46:ILE:HG13	2.03	0.41
1:A:69:LEU:HB2	2:H:503:MET:SD	2.61	0.41
2:B:606:ILE:HD12	2:B:607:THR:H	1.83	0.41
1:G:117:THR:HG21	1:G:246:ILE:HG21	2.02	0.41
2:H:456:THR:HG21	2:H:501:ILE:HD12	2.01	0.41
1:A:45:ASN:OD1	1:A:47:LYS:HB3	2.20	0.41
1:A:116:THR:HG22	1:A:120:LYS:HE3	2.03	0.41
2:D:403:VAL:N	6:D:814:HOH:O	2.52	0.41
2:F:397:LEU:HD23	2:F:397:LEU:HA	1.85	0.41
2:F:607:THR:HB	2:F:609:LYS:HG2	2.03	0.41
2:B:431:LYS:NZ	3:B:701:GNP:O2G	2.41	0.41
2:B:577:GLN:H	2:B:577:GLN:NE2	2.19	0.41
6:C:411:HOH:O	2:D:558:GLU:HB2	2.21	0.41
2:D:458:ARG:CZ	2:D:458:ARG:HB3	2.50	0.41
2:F:409:ILE:HG12	2:F:419:TYR:CD1	2.56	0.41
2:F:431:LYS:NZ	3:F:701:GNP:O1G	2.51	0.41
1:A:56:VAL:HG23	1:A:78:ALA:HB1	2.03	0.41
2:B:588:THR:OG1	2:B:589:LYS:HG3	2.20	0.41
2:D:371:GLU:HA	2:D:372:GLY:HA2	1.54	0.41
2:F:503:MET:HA	2:F:542:ASN:OD1	2.21	0.41
1:G:23:ILE:HG13	1:G:27:VAL:HG23	2.03	0.41
1:A:60:ILE:HD13	1:A:75:ILE:HG12	2.02	0.40
2:D:381:VAL:O	2:D:385:VAL:HG13	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:491:GLU:HG2	2:F:492:LYS:N	2.36	0.40
2:H:526:GLN:NE2	2:H:558:GLU:OE2	2.54	0.40
2:H:560:VAL:HA	2:H:606:ILE:HD13	2.03	0.40
2:B:406:LEU:O	2:B:410:MET:HG3	2.21	0.40
2:D:402:ARG:HH11	2:D:402:ARG:HB2	1.87	0.40
2:D:437:LYS:HG2	2:D:620:TYR:HA	2.03	0.40
2:H:471:ARG:HD2	2:H:620:TYR:HE2	1.85	0.40
2:H:567:ASN:OD1	2:H:581:LEU:HB3	2.22	0.40
1:C:234:LYS:HE2	1:C:238:ASP:OD2	2.21	0.40
2:F:362:LEU:HD13	2:F:605:TYR:HE1	1.87	0.40
1:G:114:LYS:H	3:G:301:GNP:PB	2.45	0.40
2:B:370:LEU:HD12	2:B:385:VAL:HG12	2.02	0.40
1:C:212:ILE:HG13	1:C:214:PRO:HD3	2.03	0.40
2:D:343:ASP:CG	2:D:346:ARG:HH12	2.30	0.40
1:E:97:THR:HB	1:E:100:LYS:HE2	2.03	0.40
1:G:198:GLU:HG3	1:G:201:LEU:H	1.87	0.40
1:G:272:PHE:CD1	1:G:284:PRO:HA	2.57	0.40
2:H:511:ASN:HD22	2:H:512:GLN:NE2	2.20	0.40
2:H:597:VAL:CG1	2:H:635:LEU:HD13	2.50	0.40
1:A:124:TYR:O	1:A:128:LYS:HG2	2.22	0.40
2:D:352:LYS:HZ2	2:D:638:ALA:CB	2.35	0.40
2:D:431:LYS:H	3:D:701:GNP:PB	2.45	0.40
1:G:87:VAL:HG22	1:G:260:LEU:HD21	2.02	0.40
1:G:114:LYS:HE3	1:G:193:GLY:HA2	2.02	0.40
1:G:120:LYS:NZ	1:G:276:GLY:O	2.40	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	270/304 (89%)	267 (99%)	3 (1%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	264/304 (87%)	255 (97%)	8 (3%)	1 (0%)	30	51
1	E	264/304 (87%)	257 (97%)	7 (3%)	0	100	100
1	G	263/304 (86%)	258 (98%)	4 (2%)	1 (0%)	30	51
2	B	296/315 (94%)	284 (96%)	12 (4%)	0	100	100
2	D	301/315 (96%)	286 (95%)	15 (5%)	0	100	100
2	F	300/315 (95%)	289 (96%)	11 (4%)	0	100	100
2	H	294/315 (93%)	285 (97%)	9 (3%)	0	100	100
All	All	2252/2476 (91%)	2181 (97%)	69 (3%)	2 (0%)	48	70

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	252	GLY
1	G	252	GLY

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	229/253 (90%)	219 (96%)	10 (4%)	25	50
1	C	224/253 (88%)	208 (93%)	16 (7%)	13	31
1	E	224/253 (88%)	212 (95%)	12 (5%)	20	42
1	G	223/253 (88%)	214 (96%)	9 (4%)	28	55
2	B	247/259 (95%)	232 (94%)	15 (6%)	17	37
2	D	249/259 (96%)	229 (92%)	20 (8%)	11	25
2	F	249/259 (96%)	231 (93%)	18 (7%)	13	30
2	H	246/259 (95%)	233 (95%)	13 (5%)	20	43
All	All	1891/2048 (92%)	1778 (94%)	113 (6%)	17	37

All (113) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	LEU
1	A	47	LYS
1	A	61	ASP
1	A	65	MET
1	A	70	ASN
1	A	87	VAL
1	A	196	LYS
1	A	197	GLN
1	A	240	VAL
1	A	251	ASP
2	B	332	SER
2	B	352	LYS
2	B	362	LEU
2	B	449	VAL
2	B	456	THR
2	B	475	LEU
2	B	476	HIS
2	B	484	ARG
2	B	487	VAL
2	B	491	GLU
2	B	577	GLN
2	B	578	THR
2	B	597	VAL
2	B	626	LEU
2	B	635	LEU
1	C	23	ILE
1	C	24	ASN
1	C	25	GLU
1	C	26	GLU
1	C	31	MET
1	C	32	LEU
1	C	41	GLU
1	C	62	LEU
1	C	73	LYS
1	C	76	GLN
1	C	87	VAL
1	C	197	GLN
1	C	240	VAL
1	C	251	ASP
1	C	280	ASP
1	C	293	LYS
2	D	330	MET
2	D	333	LEU

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Mol	Chain	Res	Type
2	D	339	GLU
2	D	342	LEU
2	D	362	LEU
2	D	379	SER
2	D	381	VAL
2	D	395	GLN
2	D	405	MET
2	D	449	VAL
2	D	458	ARG
2	D	484	ARG
2	D	487	VAL
2	D	527	ASP
2	D	540	THR
2	D	546	LEU
2	D	558	GLU
2	D	577	GLN
2	D	601	ILE
2	D	635	LEU
1	E	26	GLU
1	E	32	LEU
1	E	41	GLU
1	E	62	LEU
1	E	74	MET
1	E	164	THR
1	E	192	SER
1	E	197	GLN
1	E	240	VAL
1	E	251	ASP
1	E	288	GLN
1	E	293	LYS
2	F	336	GLU
2	F	342	LEU
2	F	362	LEU
2	F	378	PHE
2	F	381	VAL
2	F	384	THR
2	F	386	LYS
2	F	387	GLN
2	F	402	ARG
2	F	449	VAL
2	F	484	ARG
2	F	487	VAL

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Mol	Chain	Res	Type
2	F	491	GLU
2	F	511	ASN
2	F	577	GLN
2	F	593	ILE
2	F	597	VAL
2	F	635	LEU
1	G	41	GLU
1	G	62	LEU
1	G	71	LYS
1	G	72	ARG
1	G	73	LYS
1	G	87	VAL
1	G	197	GLN
1	G	240	VAL
1	G	251	ASP
2	H	339	GLU
2	H	342	LEU
2	H	362	LEU
2	H	380	THR
2	H	383	SER
2	H	444	GLU
2	H	449	VAL
2	H	471	ARG
2	H	484	ARG
2	H	487	VAL
2	H	543	THR
2	H	593	ILE
2	H	635	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	210	ASN
1	C	24	ASN
1	C	55	ASN
1	C	288	GLN
1	E	55	ASN
2	F	476	HIS
2	F	481	HIS
2	H	348	HIS
2	H	512	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 30 ligands modelled in this entry, 8 are monoatomic - leaving 22 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	$\# Z > 2$
5	SO4	B	703	-	4,4,4	0.25	0	6,6,6	0.27	0
3	GNP	D	701	4	34,34,34	1.15	4 (11%)	47,54,54	0.99	4 (8%)
5	SO4	F	703	-	4,4,4	0.29	0	6,6,6	0.43	0
3	GNP	G	301	4	34,34,34	1.28	6 (17%)	47,54,54	1.09	4 (8%)
5	SO4	H	707	-	4,4,4	0.24	0	6,6,6	0.40	0
5	SO4	H	704	-	4,4,4	0.23	0	6,6,6	0.07	0
5	SO4	H	705	-	4,4,4	0.26	0	6,6,6	0.09	0
5	SO4	G	303	-	4,4,4	0.32	0	6,6,6	0.61	0
5	SO4	H	703	-	4,4,4	0.27	0	6,6,6	0.36	0
3	GNP	H	701	4	34,34,34	1.12	4 (11%)	47,54,54	1.07	1 (2%)
5	SO4	B	705	-	4,4,4	0.27	0	6,6,6	0.16	0
5	SO4	C	303	-	4,4,4	0.20	0	6,6,6	0.12	0
5	SO4	F	705	-	4,4,4	0.27	0	6,6,6	0.42	0
3	GNP	B	701	4	34,34,34	1.27	5 (14%)	47,54,54	1.07	4 (8%)
3	GNP	E	301	4	34,34,34	1.28	5 (14%)	47,54,54	0.88	3 (6%)
5	SO4	F	704	-	4,4,4	0.27	0	6,6,6	0.17	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GNP	A	301	4	34,34,34	1.39	8 (23%)	47,54,54	1.07	6 (12%)
5	SO4	H	706	-	4,4,4	0.24	0	6,6,6	0.30	0
3	GNP	F	701	4	34,34,34	1.14	5 (14%)	47,54,54	0.85	3 (6%)
3	GNP	C	301	4	34,34,34	1.22	5 (14%)	47,54,54	0.93	2 (4%)
5	SO4	D	703	-	4,4,4	0.26	0	6,6,6	0.36	0
5	SO4	B	704	-	4,4,4	0.23	0	6,6,6	0.22	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	A	301	4	-	4/18/38/38	0/3/3/3
3	GNP	B	701	4	-	3/18/38/38	0/3/3/3
3	GNP	D	701	4	-	3/18/38/38	0/3/3/3
3	GNP	E	301	4	-	3/18/38/38	0/3/3/3
3	GNP	F	701	4	-	8/18/38/38	0/3/3/3
3	GNP	G	301	4	-	8/18/38/38	0/3/3/3
3	GNP	C	301	4	-	4/18/38/38	0/3/3/3
3	GNP	H	701	4	-	2/18/38/38	0/3/3/3

All (42) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	301	GNP	PG-O1G	3.73	1.51	1.46
3	G	301	GNP	PG-N3B	3.63	1.72	1.63
3	E	301	GNP	PG-N3B	3.62	1.72	1.63
3	E	301	GNP	PB-O3A	3.53	1.63	1.59
3	E	301	GNP	PB-O1B	3.45	1.51	1.46
3	A	301	GNP	PG-N3B	3.38	1.72	1.63
3	B	701	GNP	PG-O1G	3.32	1.51	1.46
3	G	301	GNP	PB-O3A	3.15	1.63	1.59
3	C	301	GNP	PG-N3B	3.10	1.71	1.63
3	B	701	GNP	PB-O3A	3.04	1.62	1.59
3	A	301	GNP	PG-O1G	3.04	1.50	1.46
3	B	701	GNP	PG-N3B	2.99	1.71	1.63
3	F	701	GNP	PB-O2B	-2.94	1.49	1.56
3	D	701	GNP	PB-O3A	2.89	1.62	1.59
3	A	301	GNP	PB-O1B	2.87	1.50	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	301	GNP	PA-O2A	-2.86	1.42	1.55
3	F	701	GNP	PG-O1G	2.77	1.50	1.46
3	H	701	GNP	PB-O2B	-2.77	1.49	1.56
3	D	701	GNP	PG-O1G	2.75	1.50	1.46
3	E	301	GNP	PG-O1G	2.70	1.50	1.46
3	D	701	GNP	PB-O2B	-2.69	1.49	1.56
3	F	701	GNP	PB-O3A	2.69	1.62	1.59
3	C	301	GNP	PB-O3A	2.65	1.62	1.59
3	H	701	GNP	PG-N3B	2.62	1.70	1.63
3	A	301	GNP	PA-O3A	-2.56	1.56	1.59
3	C	301	GNP	PB-O1B	2.53	1.50	1.46
3	D	701	GNP	PB-O1B	2.53	1.50	1.46
3	A	301	GNP	PB-O2B	-2.53	1.50	1.56
3	G	301	GNP	PB-O1B	2.51	1.50	1.46
3	F	701	GNP	PG-N3B	2.49	1.69	1.63
3	B	701	GNP	PB-O2B	-2.43	1.50	1.56
3	H	701	GNP	PB-O3A	2.40	1.62	1.59
3	G	301	GNP	PB-O2B	-2.39	1.50	1.56
3	C	301	GNP	PB-O2B	-2.39	1.50	1.56
3	G	301	GNP	PG-O1G	2.32	1.49	1.46
3	F	701	GNP	PB-O1B	2.28	1.49	1.46
3	A	301	GNP	PB-O3A	2.24	1.61	1.59
3	A	301	GNP	PA-O5'	-2.22	1.50	1.59
3	B	701	GNP	PA-O2A	-2.22	1.45	1.55
3	E	301	GNP	PB-O2B	-2.21	1.50	1.56
3	H	701	GNP	PA-O3A	-2.19	1.57	1.59
3	G	301	GNP	PA-O3A	-2.14	1.57	1.59

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	701	GNP	O1G-PG-N3B	-5.45	103.75	111.77
3	G	301	GNP	O3G-PG-O1G	-3.08	105.72	113.45
3	B	701	GNP	O2G-PG-O1G	-3.08	105.73	113.45
3	B	701	GNP	O3A-PB-N3B	-3.07	98.08	106.59
3	G	301	GNP	O2A-PA-O5'	-3.00	93.96	107.57
3	E	301	GNP	O2G-PG-O1G	-2.93	106.11	113.45
3	G	301	GNP	O2A-PA-O1A	2.88	125.86	112.44
3	C	301	GNP	O1G-PG-N3B	-2.75	107.73	111.77
3	C	301	GNP	O2G-PG-O1G	-2.67	106.75	113.45
3	D	701	GNP	O1G-PG-N3B	-2.67	107.84	111.77
3	A	301	GNP	O2A-PA-O5'	-2.62	95.69	107.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	701	GNP	O2B-PB-O1B	-2.45	104.61	109.87
3	G	301	GNP	O5'-C5'-C4'	-2.40	100.81	108.99
3	D	701	GNP	O3A-PB-N3B	-2.33	100.13	106.59
3	A	301	GNP	O2A-PA-O1A	2.32	123.25	112.44
3	F	701	GNP	O3G-PG-O1G	-2.30	107.69	113.45
3	D	701	GNP	O2A-PA-O1A	2.27	123.01	112.44
3	A	301	GNP	O1G-PG-N3B	-2.26	108.45	111.77
3	A	301	GNP	O3G-PG-O1G	-2.25	107.81	113.45
3	E	301	GNP	O2A-PA-O1A	2.25	122.90	112.44
3	D	701	GNP	O2B-PB-O1B	-2.21	105.13	109.87
3	F	701	GNP	O2A-PA-O1A	2.17	122.53	112.44
3	A	301	GNP	O1B-PB-N3B	2.16	114.95	111.77
3	A	301	GNP	O2B-PB-O1B	-2.08	105.41	109.87
3	B	701	GNP	O2B-PB-O3A	2.06	111.53	104.64
3	E	301	GNP	O2B-PB-O1B	-2.01	105.56	109.87
3	F	701	GNP	O1G-PG-N3B	-2.01	108.82	111.77

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	301	GNP	PB-N3B-PG-O1G
3	A	301	GNP	PG-N3B-PB-O1B
3	A	301	GNP	PG-N3B-PB-O3A
3	B	701	GNP	PB-N3B-PG-O1G
3	B	701	GNP	PG-N3B-PB-O1B
3	B	701	GNP	PG-N3B-PB-O3A
3	C	301	GNP	PG-N3B-PB-O1B
3	C	301	GNP	PG-N3B-PB-O3A
3	D	701	GNP	PB-N3B-PG-O1G
3	D	701	GNP	PG-N3B-PB-O1B
3	D	701	GNP	PG-N3B-PB-O3A
3	E	301	GNP	PB-N3B-PG-O1G
3	E	301	GNP	PG-N3B-PB-O1B
3	E	301	GNP	PG-N3B-PB-O3A
3	F	701	GNP	PB-N3B-PG-O1G
3	F	701	GNP	PG-N3B-PB-O1B
3	F	701	GNP	PG-N3B-PB-O3A
3	F	701	GNP	C5'-O5'-PA-O1A
3	F	701	GNP	C5'-O5'-PA-O2A
3	G	301	GNP	PB-N3B-PG-O1G
3	G	301	GNP	PG-N3B-PB-O1B

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Mol	Chain	Res	Type	Atoms
3	H	701	GNP	PG-N3B-PB-O1B
3	H	701	GNP	PG-N3B-PB-O3A
3	F	701	GNP	C3'-C4'-C5'-O5'
3	F	701	GNP	O4'-C4'-C5'-O5'
3	G	301	GNP	O4'-C4'-C5'-O5'
3	G	301	GNP	C3'-C4'-C5'-O5'
3	F	701	GNP	C5'-O5'-PA-O3A
3	G	301	GNP	C5'-O5'-PA-O3A
3	G	301	GNP	C5'-O5'-PA-O1A
3	G	301	GNP	C5'-O5'-PA-O2A
3	C	301	GNP	PA-O3A-PB-O2B
3	C	301	GNP	PA-O3A-PB-O1B
3	A	301	GNP	O4'-C4'-C5'-O5'
3	G	301	GNP	PG-N3B-PB-O3A

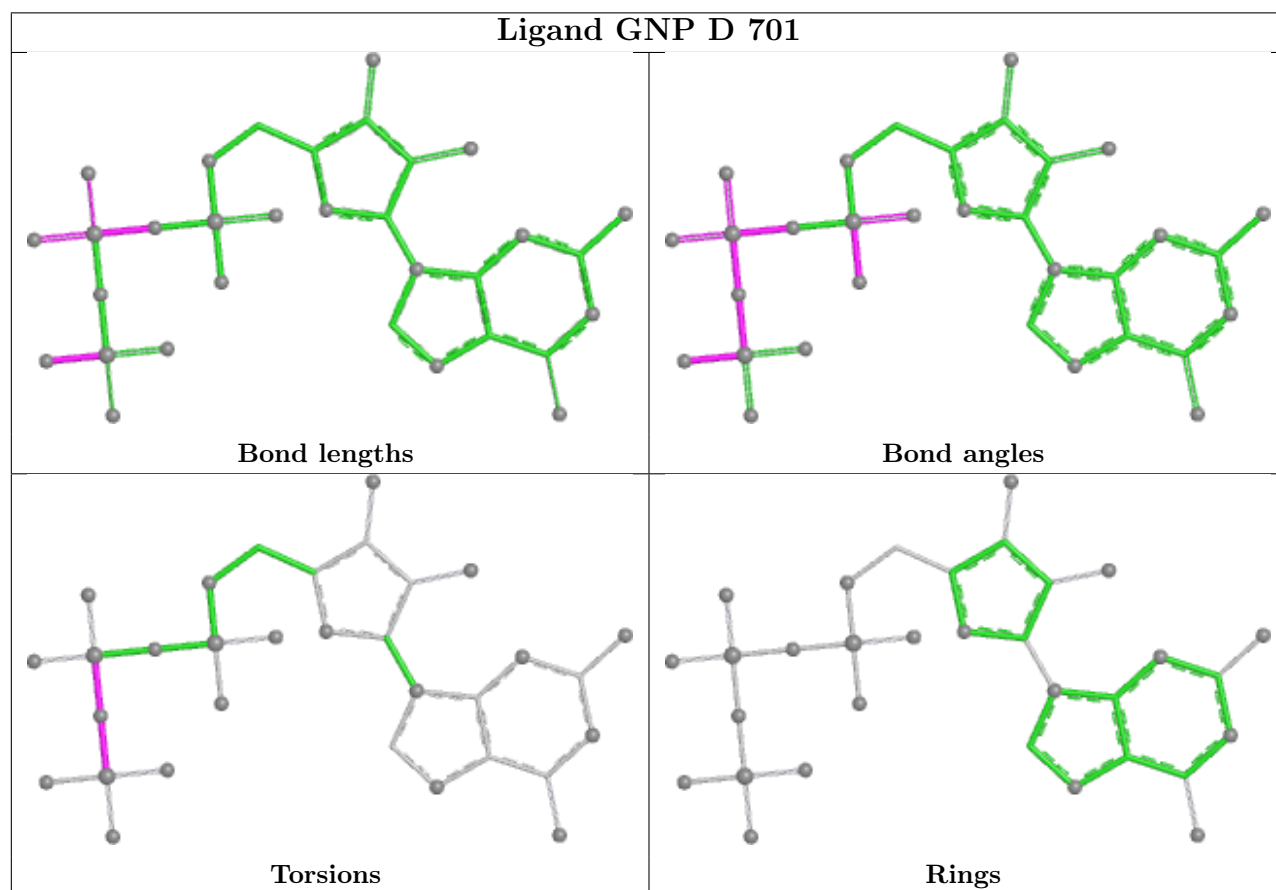
There are no ring outliers.

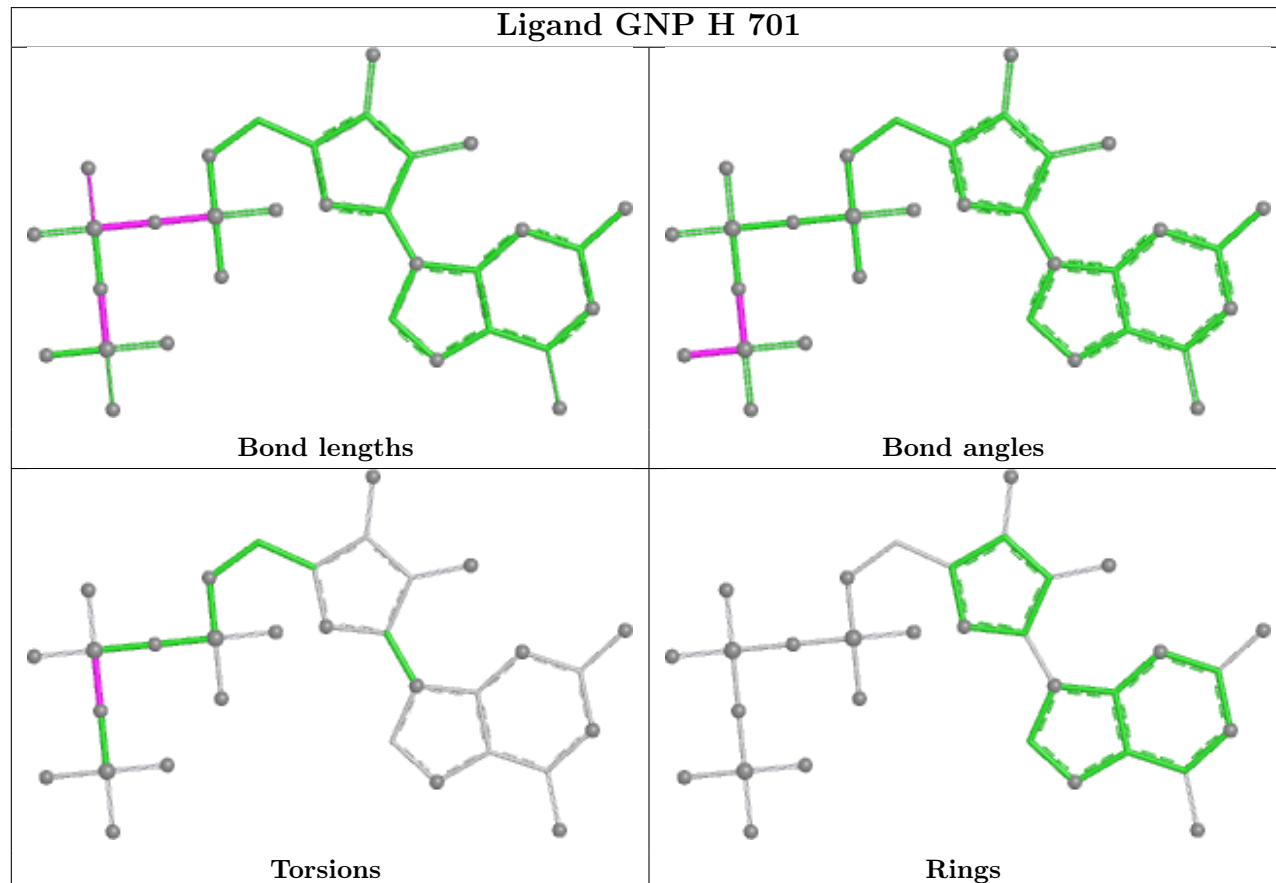
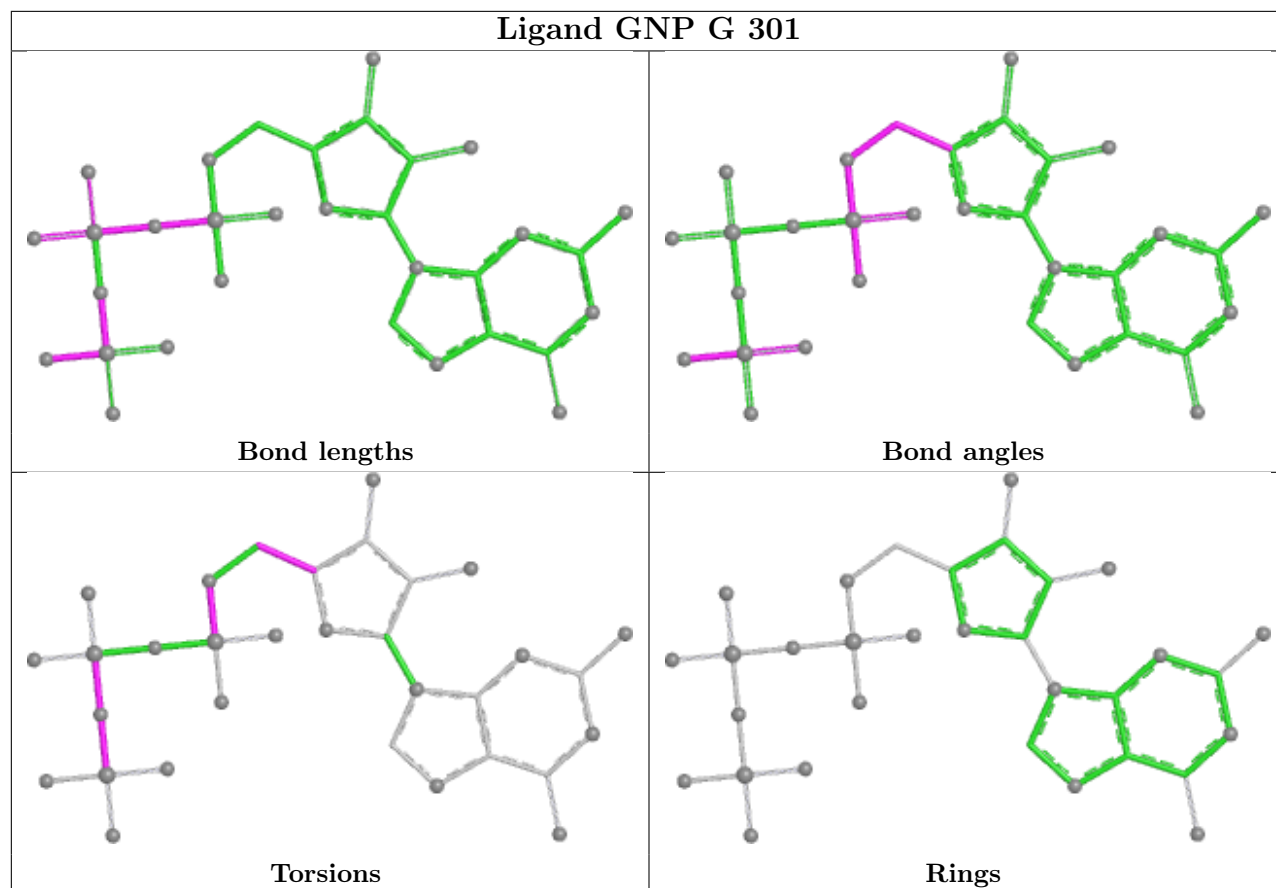
12 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	701	GNP	2	0
3	G	301	GNP	6	0
5	G	303	SO4	1	0
3	H	701	GNP	2	0
3	B	701	GNP	1	0
3	E	301	GNP	3	0
3	A	301	GNP	3	0
5	H	706	SO4	1	0
3	F	701	GNP	1	0
3	C	301	GNP	1	0
5	D	703	SO4	1	0
5	B	704	SO4	1	0

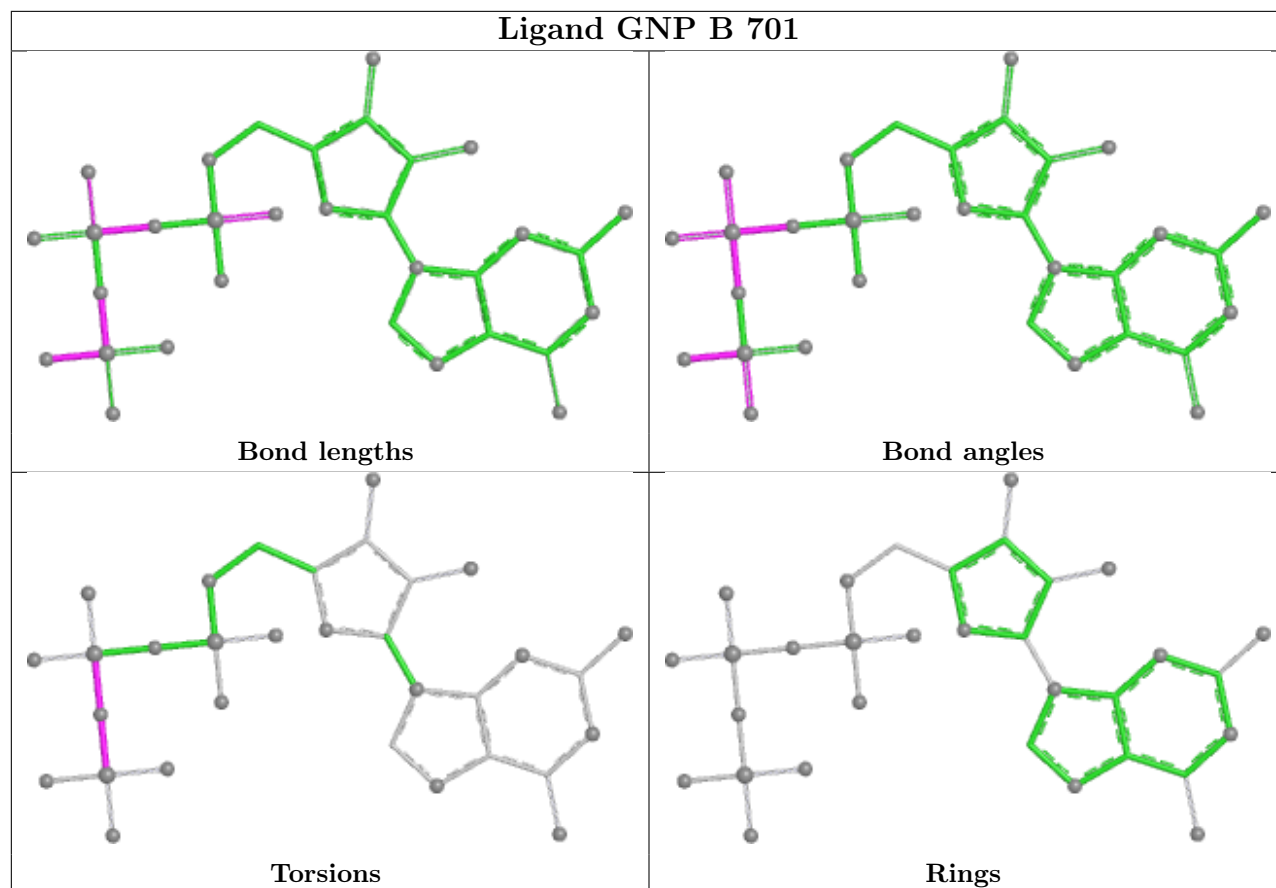
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier.

The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

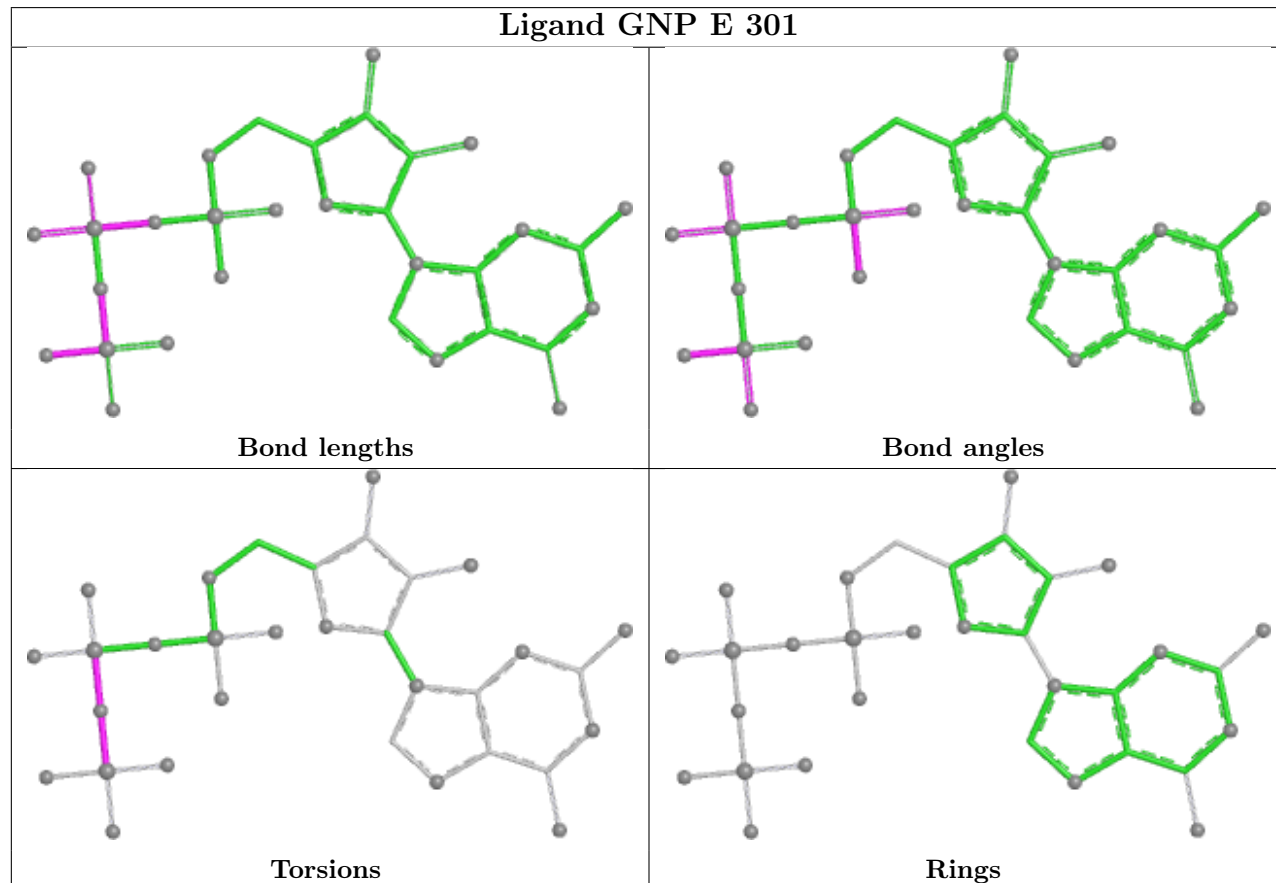


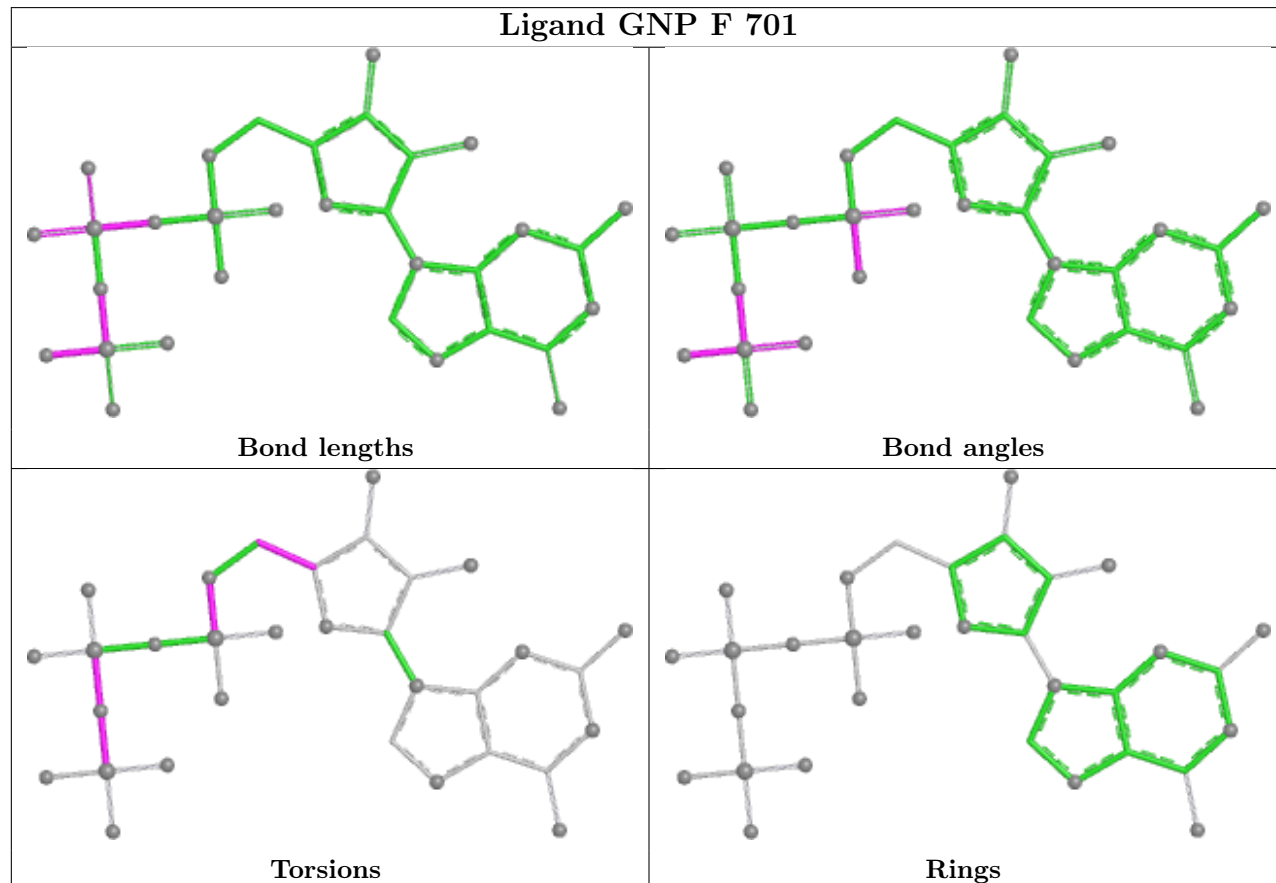
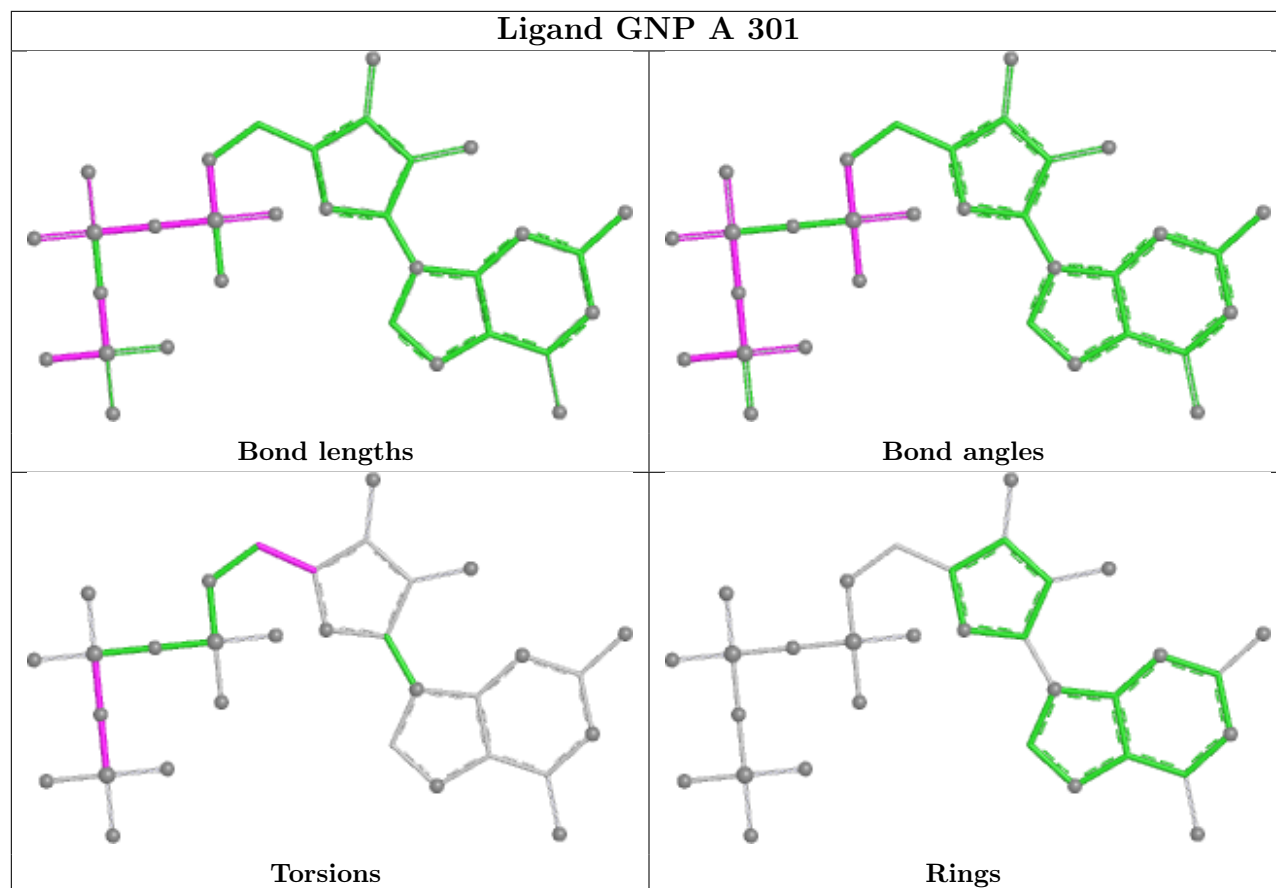


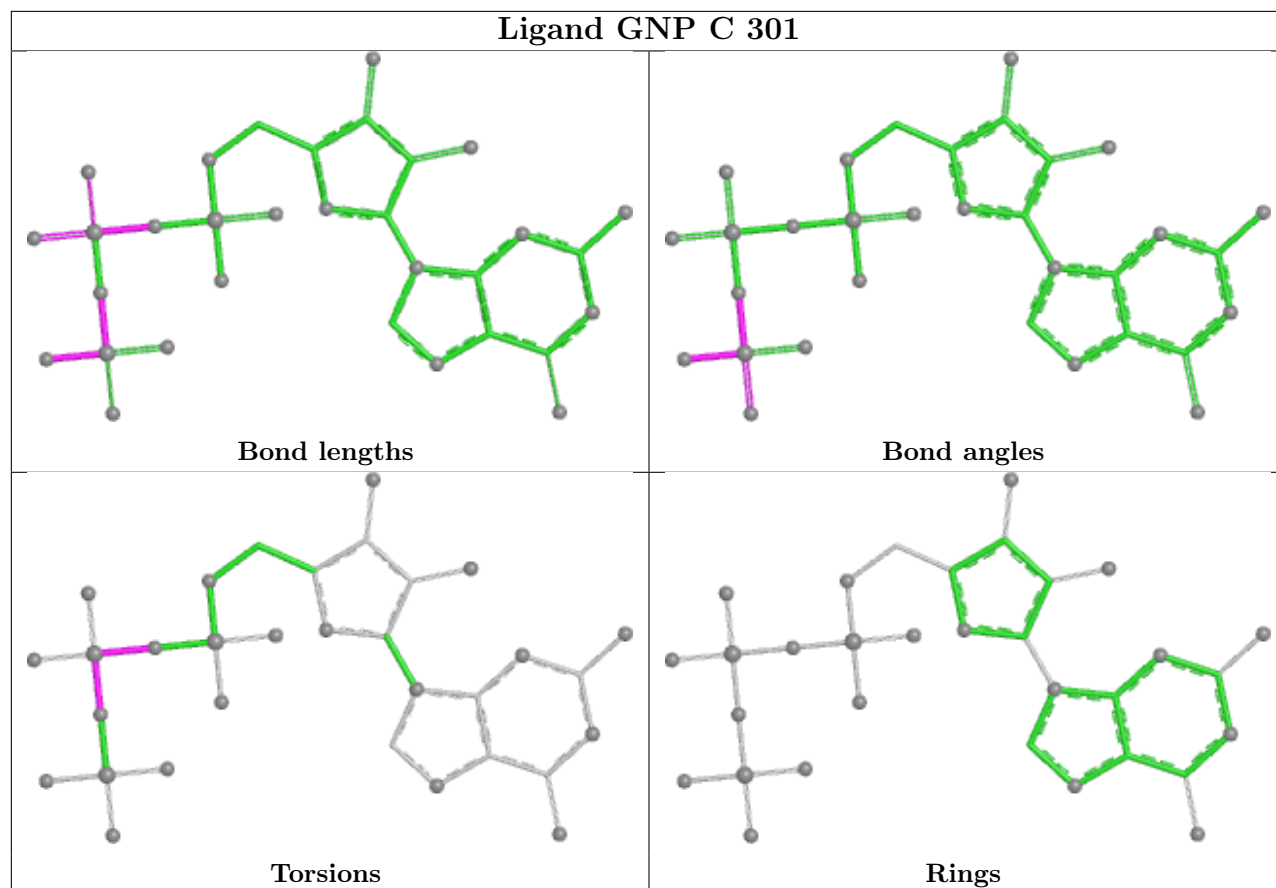
Ligand GNP B 701



Ligand GNP E 301







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, 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	274/304 (90%)	-1.63	0 100 100	38, 59, 95, 118	0
1	C	268/304 (88%)	-1.61	0 100 100	37, 57, 98, 145	0
1	E	268/304 (88%)	-1.64	0 100 100	32, 57, 97, 124	0
1	G	267/304 (87%)	-1.60	0 100 100	35, 61, 99, 139	0
2	B	302/315 (95%)	-1.66	0 100 100	34, 46, 91, 130	0
2	D	305/315 (96%)	-1.61	0 100 100	31, 52, 98, 149	0
2	F	304/315 (96%)	-1.64	0 100 100	29, 53, 98, 139	0
2	H	300/315 (95%)	-1.67	0 100 100	34, 48, 92, 130	0
All	All	2288/2476 (92%)	-1.63	0 100 100	29, 54, 97, 149	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

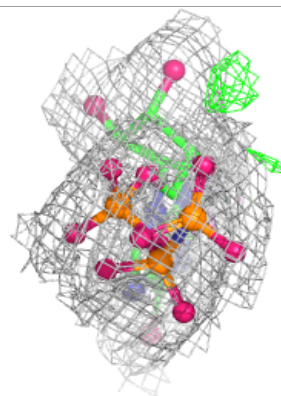
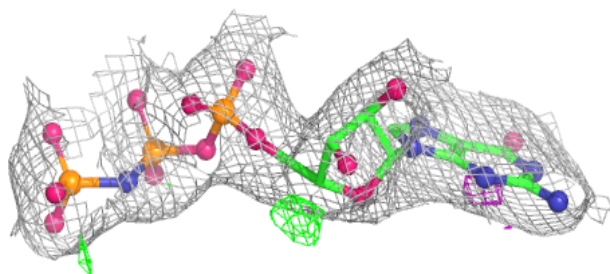
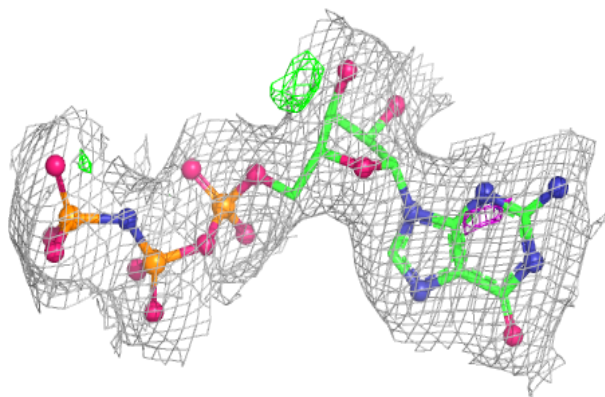
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	SO4	H	705	5/5	0.94	0.04	122,125,126,126	0
5	SO4	F	705	5/5	0.98	0.03	140,142,143,144	0
5	SO4	C	303	5/5	0.98	0.03	126,127,127,127	0
5	SO4	H	706	5/5	0.98	0.07	122,125,126,129	0
5	SO4	H	707	5/5	0.98	0.03	113,114,115,117	0
5	SO4	F	704	5/5	0.99	0.04	84,87,88,91	0
5	SO4	B	705	5/5	0.99	0.05	105,108,108,113	0
5	SO4	G	303	5/5	0.99	0.04	122,124,126,131	0
5	SO4	H	703	5/5	0.99	0.04	90,91,92,93	0
5	SO4	H	704	5/5	0.99	0.05	176,176,177,178	0
5	SO4	B	704	5/5	0.99	0.02	111,111,112,112	0
5	SO4	D	703	5/5	0.99	0.03	119,121,123,125	0
5	SO4	F	703	5/5	0.99	0.02	83,84,87,87	0
4	MG	F	702	1/1	1.00	0.01	31,31,31,31	0
4	MG	G	302	1/1	1.00	0.01	30,30,30,30	0
4	MG	H	702	1/1	1.00	0.02	29,29,29,29	0
5	SO4	B	703	5/5	1.00	0.03	74,77,78,78	0
3	GNP	A	301	32/32	1.00	0.02	13,36,44,50	0
3	GNP	B	701	32/32	1.00	0.02	21,32,42,45	0
3	GNP	C	301	32/32	1.00	0.02	27,39,51,52	0
3	GNP	D	701	32/32	1.00	0.02	23,34,45,62	0
3	GNP	E	301	32/32	1.00	0.02	21,40,50,51	0
3	GNP	F	701	32/32	1.00	0.03	19,36,55,98	0
3	GNP	G	301	32/32	1.00	0.02	20,35,49,53	0
3	GNP	H	701	32/32	1.00	0.02	17,36,47,76	0
4	MG	A	302	1/1	1.00	0.01	35,35,35,35	0
4	MG	B	702	1/1	1.00	0.01	30,30,30,30	0
4	MG	C	302	1/1	1.00	0.01	32,32,32,32	0
4	MG	D	702	1/1	1.00	0.02	34,34,34,34	0
4	MG	E	302	1/1	1.00	0.01	38,38,38,38	0

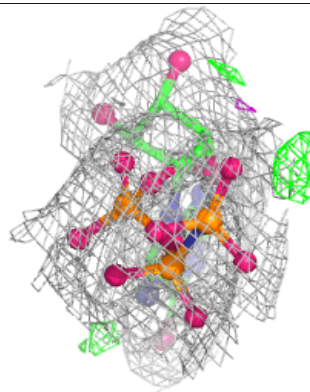
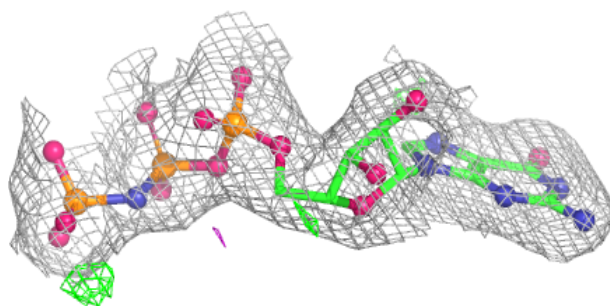
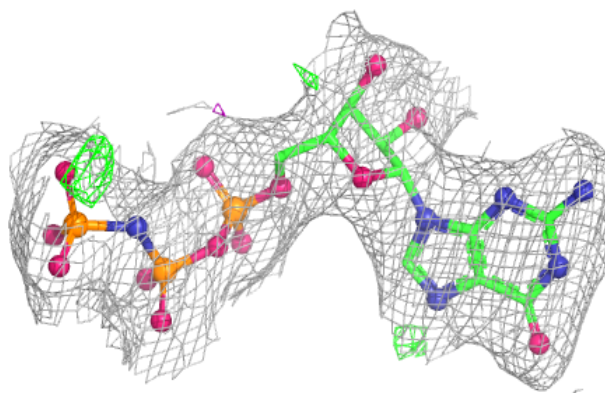
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around GNP 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)

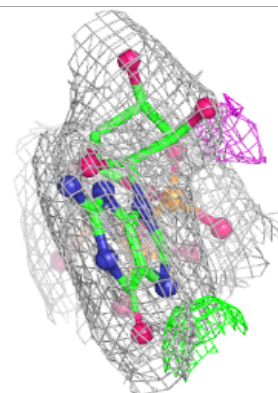
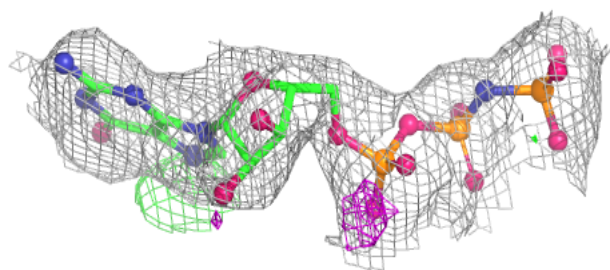
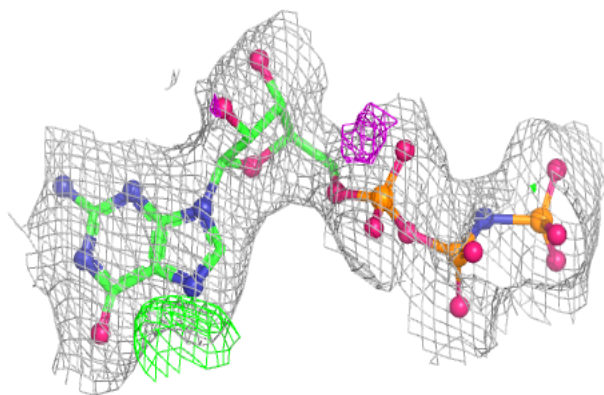
**Electron density around GNP B 701:**

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

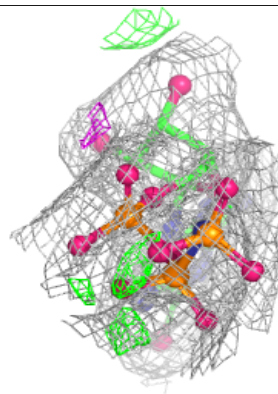
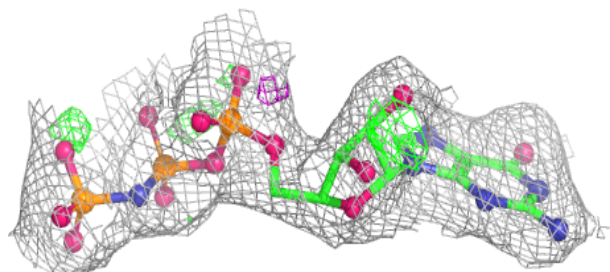
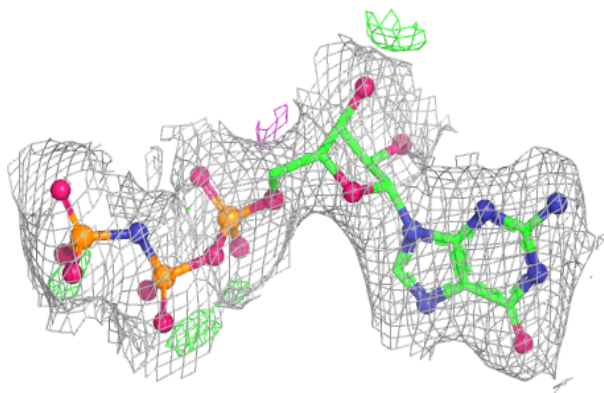


Electron density around GNP C 301:

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

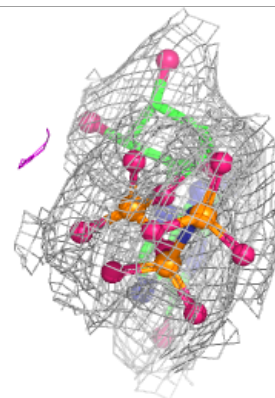
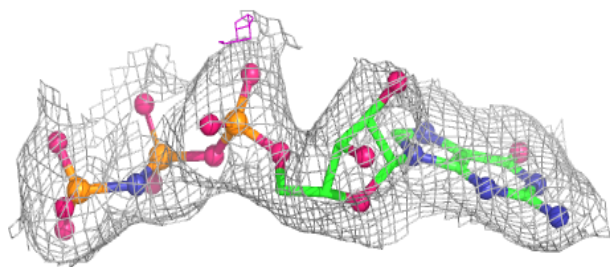
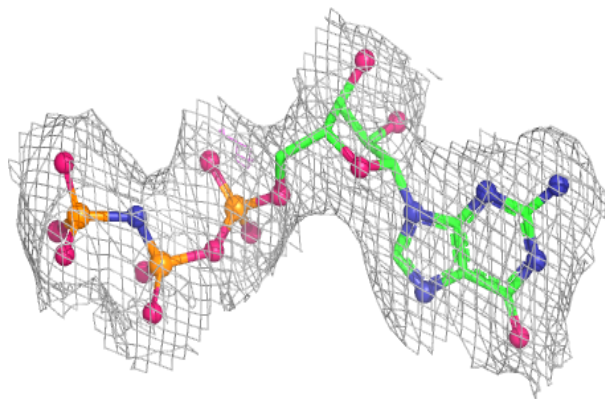
**Electron density around GNP D 701:**

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

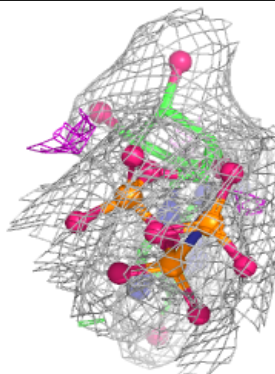
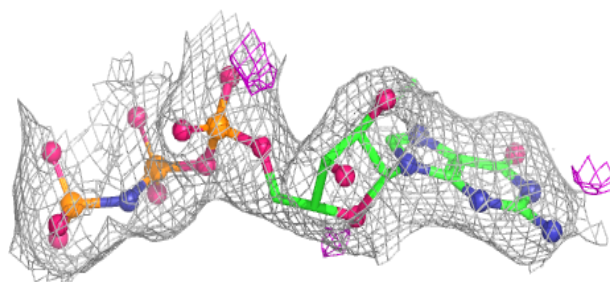
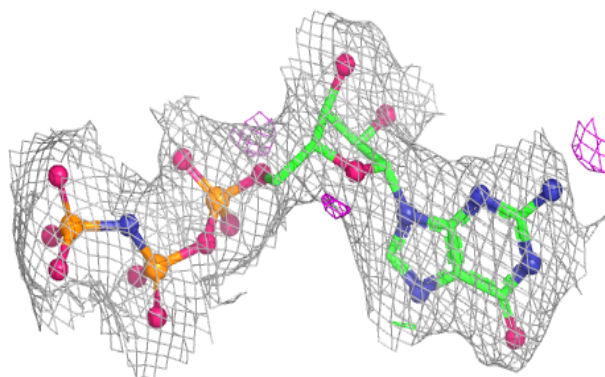


Electron density around GNP E 301:

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

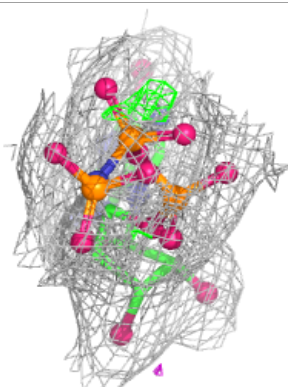
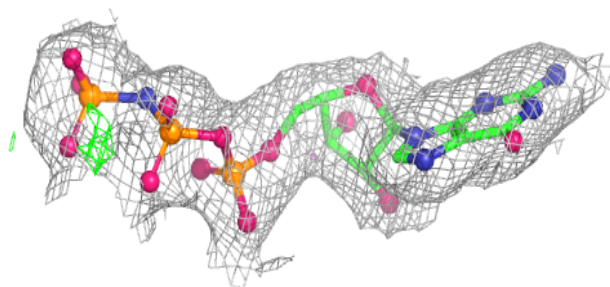
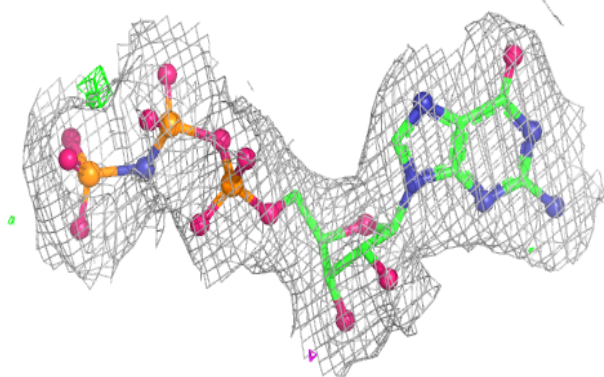
**Electron density around GNP F 701:**

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

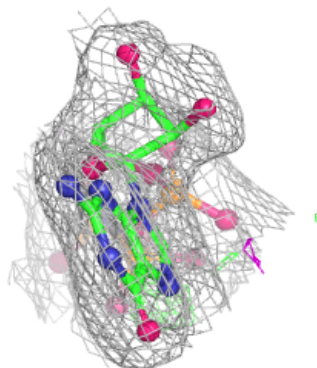
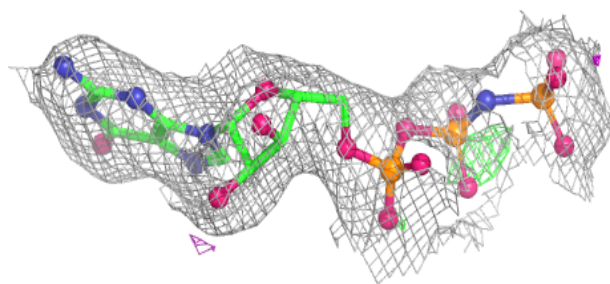
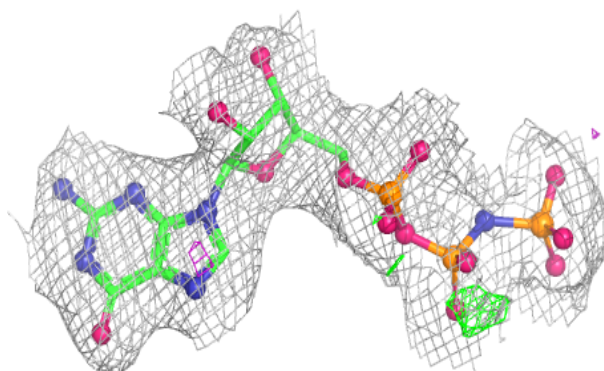


Electron density around GNP G 301:

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

**Electron density around GNP H 701:**

$2mF_o-DF_c$ (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.