



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 10, 2023 – 06:48 PM EDT

PDB ID : 4K81
Title : Crystal structure of the Grb14 RA and PH domains in complex with GTP-loaded H-Ras
Authors : Qamra, R.; Hubbard, S.R.
Deposited on : 2013-04-17
Resolution : 2.40 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 : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.35.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35.1

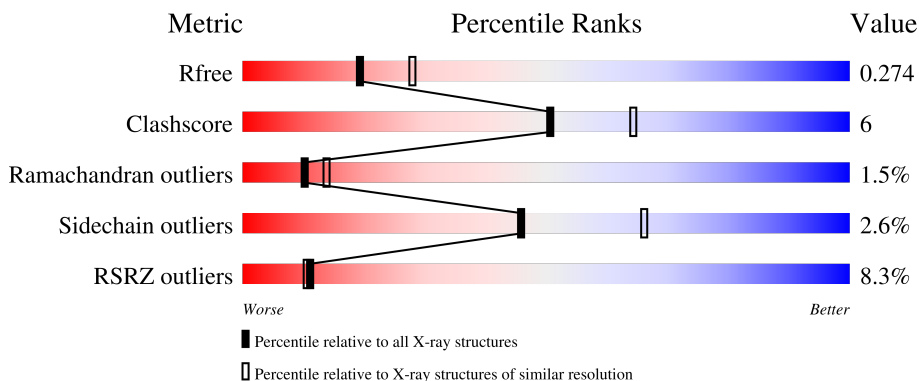
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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.



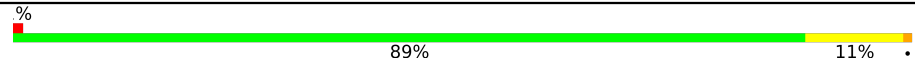

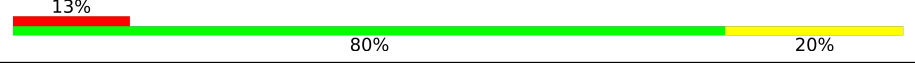
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	258	
1	C	258	
1	E	258	
1	G	258	
2	B	171	

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Mol	Chain	Length	Quality of chain
2	D	171	 <p>% 89% 11% .</p>
2	F	171	 <p>2% 81% 19% .</p>
2	H	171	 <p>13% 80% 20%</p>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 13867 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 Growth factor receptor-bound protein 14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	244	2023	1312	340	361	10	0	0	0
1	C	244	2023	1312	340	361	10	0	0	0
1	E	244	2023	1312	340	361	10	0	0	0
1	G	244	2023	1312	340	361	10	0	0	0

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	99	GLY	-	expression tag	UNP Q14449
A	100	HIS	-	expression tag	UNP Q14449
A	101	MET	-	expression tag	UNP Q14449
A	102	ALA	-	expression tag	UNP Q14449
A	103	SER	-	expression tag	UNP Q14449
A	104	GLY	-	expression tag	UNP Q14449
A	105	SER	-	expression tag	UNP Q14449
A	272	ALA	LYS	engineered mutation	UNP Q14449
A	273	ALA	GLU	engineered mutation	UNP Q14449
C	99	GLY	-	expression tag	UNP Q14449
C	100	HIS	-	expression tag	UNP Q14449
C	101	MET	-	expression tag	UNP Q14449
C	102	ALA	-	expression tag	UNP Q14449
C	103	SER	-	expression tag	UNP Q14449
C	104	GLY	-	expression tag	UNP Q14449
C	105	SER	-	expression tag	UNP Q14449
C	272	ALA	LYS	engineered mutation	UNP Q14449
C	273	ALA	GLU	engineered mutation	UNP Q14449
E	99	GLY	-	expression tag	UNP Q14449
E	100	HIS	-	expression tag	UNP Q14449
E	101	MET	-	expression tag	UNP Q14449

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Chain	Residue	Modelled	Actual	Comment	Reference
E	102	ALA	-	expression tag	UNP Q14449
E	103	SER	-	expression tag	UNP Q14449
E	104	GLY	-	expression tag	UNP Q14449
E	105	SER	-	expression tag	UNP Q14449
E	272	ALA	LYS	engineered mutation	UNP Q14449
E	273	ALA	GLU	engineered mutation	UNP Q14449
G	99	GLY	-	expression tag	UNP Q14449
G	100	HIS	-	expression tag	UNP Q14449
G	101	MET	-	expression tag	UNP Q14449
G	102	ALA	-	expression tag	UNP Q14449
G	103	SER	-	expression tag	UNP Q14449
G	104	GLY	-	expression tag	UNP Q14449
G	105	SER	-	expression tag	UNP Q14449
G	272	ALA	LYS	engineered mutation	UNP Q14449
G	273	ALA	GLU	engineered mutation	UNP Q14449

- Molecule 2 is a protein called GTPase HRas.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	171	1352	842	233	269	8	0	0	0
2	D	171	1352	842	233	269	8	0	0	0
2	F	171	1352	842	233	269	8	0	0	0
2	H	171	1352	842	233	269	8	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

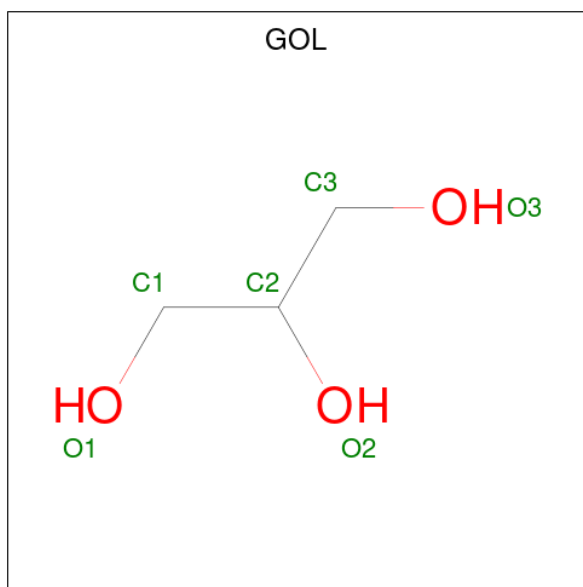
Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	GLY	-	expression tag	UNP P01112
B	-3	ALA	-	expression tag	UNP P01112
B	-2	MET	-	expression tag	UNP P01112
B	-1	GLY	-	expression tag	UNP P01112
B	0	SER	-	expression tag	UNP P01112
B	12	VAL	GLY	engineered mutation	UNP P01112
D	-4	GLY	-	expression tag	UNP P01112
D	-3	ALA	-	expression tag	UNP P01112
D	-2	MET	-	expression tag	UNP P01112
D	-1	GLY	-	expression tag	UNP P01112
D	0	SER	-	expression tag	UNP P01112

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Chain	Residue	Modelled	Actual	Comment	Reference
D	12	VAL	GLY	engineered mutation	UNP P01112
F	-4	GLY	-	expression tag	UNP P01112
F	-3	ALA	-	expression tag	UNP P01112
F	-2	MET	-	expression tag	UNP P01112
F	-1	GLY	-	expression tag	UNP P01112
F	0	SER	-	expression tag	UNP P01112
F	12	VAL	GLY	engineered mutation	UNP P01112
H	-4	GLY	-	expression tag	UNP P01112
H	-3	ALA	-	expression tag	UNP P01112
H	-2	MET	-	expression tag	UNP P01112
H	-1	GLY	-	expression tag	UNP P01112
H	0	SER	-	expression tag	UNP P01112
H	12	VAL	GLY	engineered mutation	UNP P01112

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



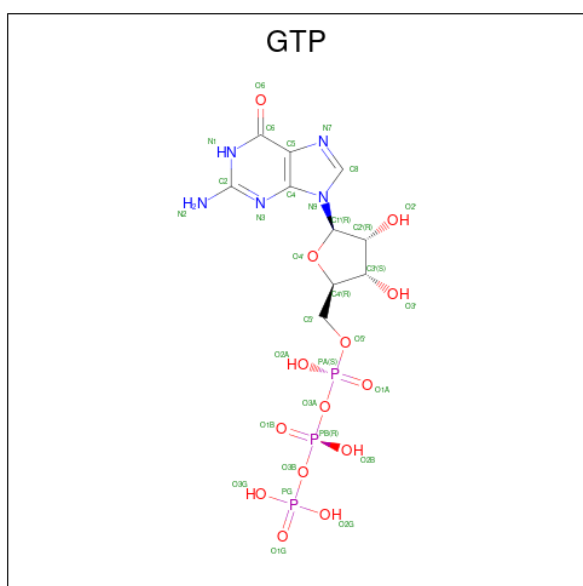
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	6	3	3	0	0
3	A	1	6	3	3	0	0
3	A	1	6	3	3	0	0
3	A	1	6	3	3	0	0
3	B	1	6	3	3	0	0

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

- Molecule 4 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
4	D	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
4	F	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
4	H	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Mg	0	0
			1	1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	D	1	Total Mg 1 1	0	0
5	F	1	Total Mg 1 1	0	0
5	H	1	Total Mg 1 1	0	0

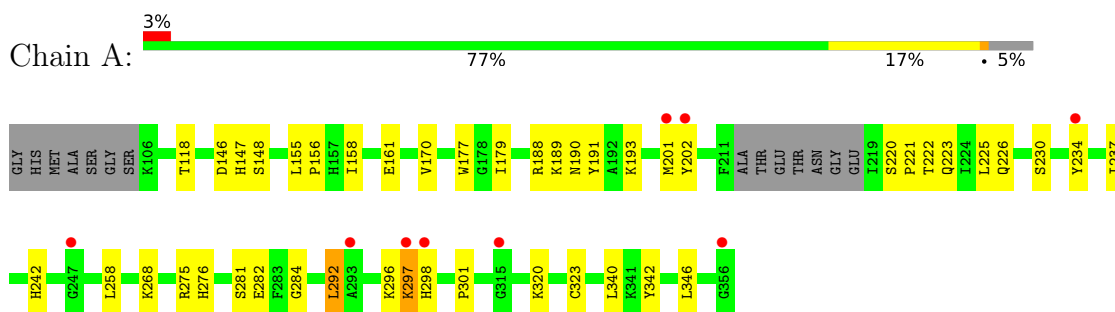
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	46	Total O 46 46	0	0
6	B	24	Total O 24 24	0	0
6	C	25	Total O 25 25	0	0
6	D	21	Total O 21 21	0	0
6	E	33	Total O 33 33	0	0
6	F	16	Total O 16 16	0	0
6	G	10	Total O 10 10	0	0
6	H	6	Total O 6 6	0	0

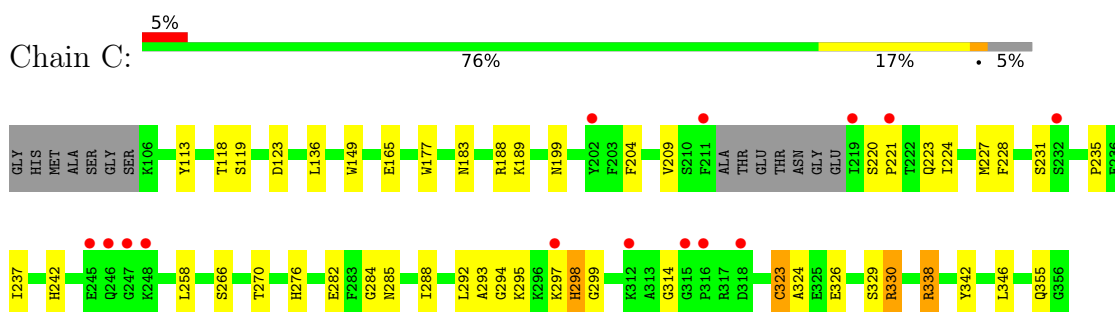
3 Residue-property plots [i](#)

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.

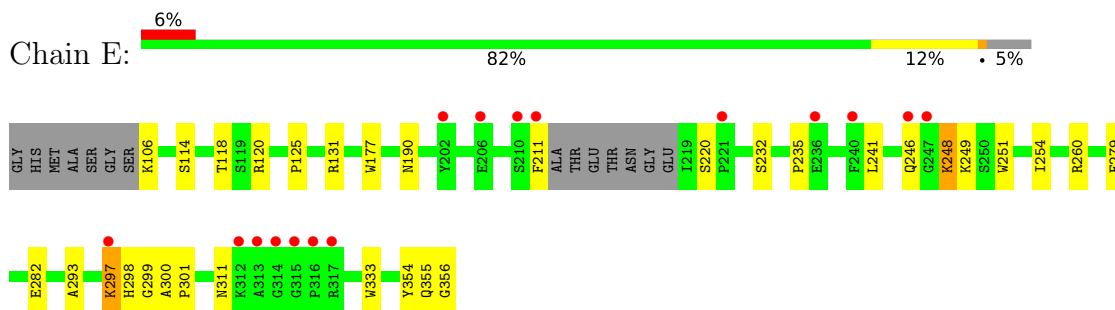
- Molecule 1: Growth factor receptor-bound protein 14



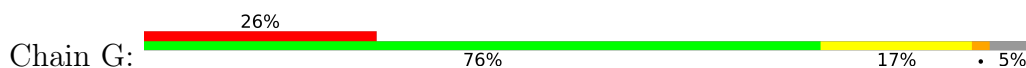
- Molecule 1: Growth factor receptor-bound protein 14

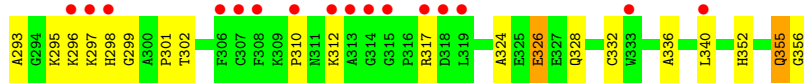
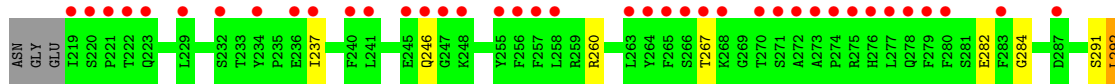


- Molecule 1: Growth factor receptor-bound protein 14

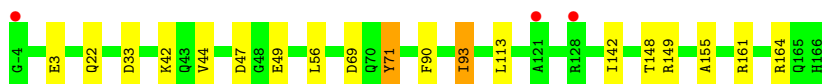
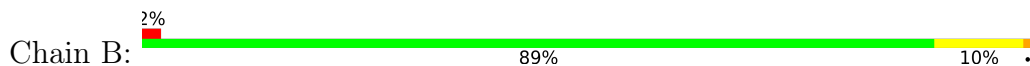


- Molecule 1: Growth factor receptor-bound protein 14

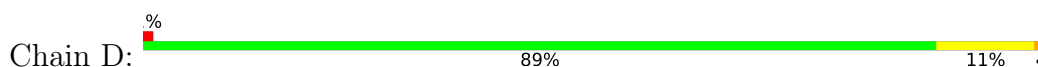




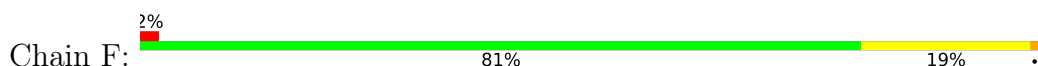
• Molecule 2: GTPase HRas



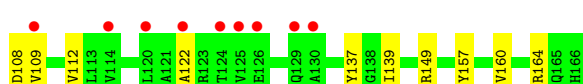
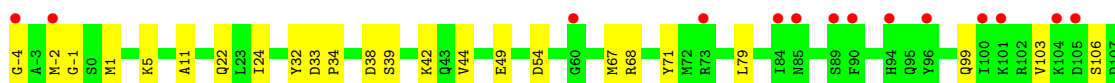
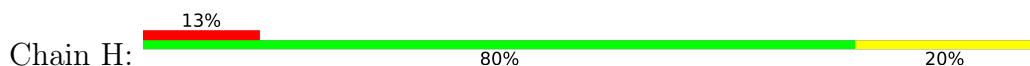
• Molecule 2: GTPase HRas



• Molecule 2: GTPase HRas



• Molecule 2: GTPase HRas



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.73Å 115.59Å 103.11Å 90.00° 96.75° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 46.81 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.7 (50.00-2.40) 99.6 (46.81-2.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 2.39Å)	Xtrriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.219 , 0.277 0.217 , 0.274	Depositor DCC
R_{free} test set	3628 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	51.7	Xtrriage
Anisotropy	0.368	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 49.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13867	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.27% 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: MG, GTP, 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	0.62	0/2083	0.68	0/2808
1	C	0.56	1/2083 (0.0%)	0.64	1/2808 (0.0%)
1	E	0.58	2/2083 (0.1%)	0.64	0/2808
1	G	0.51	1/2083 (0.0%)	0.53	0/2808
2	B	0.48	0/1371	0.63	0/1850
2	D	0.51	0/1371	0.66	0/1850
2	F	0.43	0/1371	0.59	0/1850
2	H	0.36	0/1371	0.53	0/1850
All	All	0.52	4/13816 (0.0%)	0.62	1/18632 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	177	TRP	CD2-CE2	5.25	1.47	1.41
1	G	177	TRP	CD2-CE2	5.23	1.47	1.41
1	C	149	TRP	CD2-CE2	5.22	1.47	1.41
1	E	251	TRP	CD2-CE2	5.01	1.47	1.41

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	338	ARG	NE-CZ-NH2	-5.27	117.67	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2023	0	1974	31	0
1	C	2023	0	1974	28	0
1	E	2023	0	1974	15	0
1	G	2023	0	1974	24	0
2	B	1352	0	1326	14	0
2	D	1352	0	1326	16	0
2	F	1352	0	1325	19	0
2	H	1352	0	1326	19	0
3	A	24	0	32	4	0
3	B	6	0	8	1	0
3	C	12	0	16	0	0
3	E	12	0	16	0	0
4	B	32	0	12	0	0
4	D	32	0	12	0	0
4	F	32	0	12	1	0
4	H	32	0	12	0	0
5	B	1	0	0	0	0
5	D	1	0	0	0	0
5	F	1	0	0	0	0
5	H	1	0	0	0	0
6	A	46	0	0	6	0
6	B	24	0	0	2	0
6	C	25	0	0	0	0
6	D	21	0	0	0	0
6	E	33	0	0	1	0
6	F	16	0	0	2	0
6	G	10	0	0	0	0
6	H	6	0	0	1	0
All	All	13867	0	13319	154	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 6.

All (154) 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:342:TYR:HB2	6:A:540:HOH:O	1.56	1.03
1:C:330:ARG:HH11	1:C:330:ARG:HG2	1.28	0.96
1:E:356:GLY:HA2	6:E:507:HOH:O	1.72	0.86
2:D:84:ILE:HD12	2:D:123:ARG:HG2	1.63	0.80
1:A:346:LEU:HG	6:A:540:HOH:O	1.82	0.78
2:B:49:GLU:HB2	2:B:164:ARG:HH22	1.48	0.77
2:D:22:GLN:O	2:D:149:ARG:NH1	2.19	0.75
2:B:3:GLU:O	2:D:-4:GLY:HA3	1.86	0.75
1:E:297:LYS:C	1:E:299:GLY:H	1.91	0.74
2:H:106:SER:HB3	2:H:109:VAL:HG23	1.74	0.69
2:H:137:TYR:HB2	2:H:139:ILE:HG12	1.74	0.69
1:E:120:ARG:HG3	2:F:38:ASP:OD1	1.92	0.69
1:A:275:ARG:HB2	6:A:525:HOH:O	1.92	0.68
2:H:68:ARG:HA	2:H:71:TYR:CE2	2.29	0.67
2:B:93:ILE:HD13	2:B:113:LEU:HD11	1.76	0.67
1:G:237:ILE:HD13	1:G:336:ALA:HB1	1.77	0.65
1:C:118:THR:HB	2:D:39:SER:O	1.97	0.64
6:B:318:HOH:O	1:C:294:GLY:HA2	1.98	0.62
1:C:330:ARG:HH11	1:C:330:ARG:CG	2.09	0.61
2:F:0:SER:OG	2:H:-2:MET:HG3	2.02	0.60
1:G:260:ARG:HA	1:G:340:LEU:HD11	1.84	0.60
2:B:142:ILE:HD12	2:B:155:ALA:HA	1.82	0.60
1:A:220:SER:HB3	1:A:223:GLN:HB2	1.84	0.60
1:G:297:LYS:C	1:G:299:GLY:H	2.05	0.60
2:B:47:ASP:HB3	1:C:295:LYS:HE3	1.86	0.58
1:C:227:MET:O	1:C:235:PRO:HD3	2.03	0.58
2:F:3:GLU:O	2:H:-4:GLY:HA3	2.03	0.58
1:G:352:HIS:O	1:G:355:GLN:NE2	2.37	0.57
1:A:242:HIS:CG	1:A:301:PRO:HG2	2.39	0.57
1:C:165:GLU:OE2	1:C:338:ARG:HD2	2.05	0.56
1:A:268:LYS:HG3	1:A:276:HIS:CD2	2.41	0.56
1:C:330:ARG:HG2	1:C:330:ARG:NH1	2.08	0.56
1:G:174:LEU:O	1:G:177:TRP:HB2	2.06	0.56
2:F:84:ILE:HD13	2:F:123:ARG:HG2	1.89	0.55
2:F:117:LYS:HB3	2:F:120:LEU:HD12	1.89	0.55
2:H:32:TYR:CE1	2:H:34:PRO:HG3	2.41	0.55
1:E:297:LYS:C	1:E:299:GLY:N	2.59	0.54
1:C:242:HIS:HB2	1:C:323:CYS:HB3	1.90	0.53
1:A:190:ASN:HD21	3:A:402:GOL:H11	1.72	0.53
1:G:310:PRO:HG2	1:G:317:ARG:HA	1.90	0.53
2:H:22:GLN:CG	2:H:149:ARG:HG3	2.39	0.52
1:C:266:SER:HA	1:C:276:HIS:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:241:LEU:HD21	1:E:333:TRP:CD1	2.44	0.52
1:A:296:LYS:O	1:A:298:HIS:N	2.37	0.51
1:A:242:HIS:HB2	1:A:323:CYS:HB3	1.93	0.51
1:G:205:PRO:HD2	1:G:208:MET:HG3	1.92	0.51
2:F:131:GLN:O	2:F:135:ARG:HB2	2.10	0.51
2:B:90:PHE:O	2:B:93:ILE:HG13	2.10	0.51
1:E:232:SER:O	1:E:260:ARG:HD3	2.10	0.50
1:E:106:LYS:O	1:E:125:PRO:HA	2.11	0.50
1:C:355:GLN:HA	1:C:355:GLN:OE1	2.12	0.50
4:F:200:GTP:O2G	6:F:310:HOH:O	2.18	0.50
2:D:68:ARG:HA	2:D:71:TYR:CE2	2.47	0.49
1:A:223:GLN:O	1:A:226:GLN:HG2	2.13	0.49
2:H:11:ALA:HB1	6:H:304:HOH:O	2.12	0.49
2:D:84:ILE:HD11	2:D:118:CYS:HA	1.95	0.48
2:B:49:GLU:HB2	2:B:164:ARG:NH2	2.23	0.48
2:B:148:THR:O	2:B:149:ARG:HB2	2.13	0.48
1:C:188:ARG:HG2	1:C:189:LYS:N	2.29	0.48
1:G:153:GLU:O	1:G:161:GLU:HA	2.13	0.48
1:A:146:ASP:HB3	1:A:148:SER:H	1.77	0.48
1:C:123:ASP:O	1:G:356:GLY:HA2	2.14	0.48
1:E:279:PHE:HZ	1:E:282:GLU:HB2	1.79	0.47
1:C:220:SER:HB2	1:C:223:GLN:HB2	1.97	0.47
1:E:131:ARG:NH1	1:E:355:GLN:HG3	2.30	0.47
2:F:20:THR:O	2:F:24:ILE:HG12	2.15	0.47
2:F:149:ARG:NH2	2:F:153:GLU:OE2	2.41	0.47
1:G:155:LEU:HD12	1:G:160:VAL:HG23	1.95	0.47
1:A:147:HIS:HA	3:A:401:GOL:H32	1.96	0.47
1:C:342:TYR:HB2	1:C:346:LEU:HG	1.97	0.47
1:A:281:SER:HB2	1:A:320:LYS:HD3	1.96	0.47
2:D:41:ARG:NH1	2:D:52:LEU:HD21	2.29	0.47
1:G:119:SER:H	2:H:39:SER:HB3	1.80	0.47
1:A:296:LYS:C	1:A:298:HIS:H	2.19	0.46
2:D:148:THR:O	2:D:149:ARG:HB2	2.15	0.46
2:D:157:TYR:N	2:D:157:TYR:CD1	2.83	0.46
3:A:401:GOL:H31	6:A:532:HOH:O	2.15	0.46
1:G:297:LYS:C	1:G:299:GLY:N	2.68	0.46
2:H:106:SER:CB	2:H:109:VAL:HG23	2.45	0.46
2:B:22:GLN:CG	2:B:149:ARG:HG2	2.46	0.46
1:G:326:GLU:HG3	1:G:328:GLN:H	1.81	0.46
1:C:165:GLU:OE1	1:C:342:TYR:CE2	2.68	0.46
1:A:220:SER:O	1:A:223:GLN:HB3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:161:GLU:OE2	1:A:188:ARG:NH2	2.49	0.45
1:G:160:VAL:HG12	1:G:291:SER:HA	1.97	0.45
1:A:201:MET:HB3	1:A:221:PRO:HA	1.98	0.45
1:A:201:MET:HG3	1:A:202:TYR:CD2	2.52	0.45
2:F:76:GLU:OE2	2:H:4:GLY:HA2	2.16	0.45
1:A:342:TYR:HB2	1:A:346:LEU:HG	1.98	0.45
2:F:22:GLN:O	2:F:149:ARG:NH1	2.47	0.44
1:E:190:ASN:OD1	1:E:190:ASN:C	2.55	0.44
2:F:-2:MET:HA	2:H:1:MET:O	2.17	0.44
1:A:177:TRP:HZ3	3:A:404:GOL:H12	1.82	0.44
2:B:161:ARG:HD2	6:B:321:HOH:O	2.18	0.44
2:F:39:SER:HA	2:F:55:ILE:O	2.16	0.44
2:D:116:ASN:CG	2:D:117:LYS:H	2.21	0.44
1:A:222:THR:HA	1:A:225:LEU:HB3	2.00	0.44
1:A:193:LYS:CD	6:A:536:HOH:O	2.66	0.43
1:A:296:LYS:HG3	1:A:297:LYS:N	2.32	0.43
2:D:22:GLN:NE2	2:D:149:ARG:HG3	2.33	0.43
1:C:284:GLY:O	1:C:285:ASN:HB2	2.17	0.43
1:E:211:PHE:CE1	1:E:235:PRO:HB3	2.52	0.43
2:D:84:ILE:CD1	2:D:118:CYS:HA	2.47	0.43
2:F:84:ILE:C	2:F:84:ILE:HD12	2.39	0.43
2:H:5:LYS:HA	2:H:54:ASP:HB3	2.00	0.43
2:F:74:THR:HB	6:F:301:HOH:O	2.17	0.43
1:G:149:TRP:CZ2	1:G:189:LYS:HE2	2.54	0.43
1:E:248:LYS:HE2	1:E:249:LYS:HE2	2.01	0.43
1:G:282:GLU:CD	1:G:284:GLY:H	2.21	0.43
2:H:44:VAL:HG11	2:H:157:TYR:OH	2.19	0.43
1:A:156:PRO:HD2	6:A:522:HOH:O	2.18	0.43
1:C:227:MET:HG2	1:C:235:PRO:HG3	1.99	0.43
1:G:302:THR:OG1	1:G:324:ALA:O	2.36	0.43
1:A:155:LEU:HD12	1:A:158:ILE:HD11	2.00	0.42
1:A:237:ILE:HB	1:A:258:LEU:HB3	2.01	0.42
1:C:113:TYR:CD1	1:C:119:SER:HB3	2.54	0.42
1:C:297:LYS:C	1:C:299:GLY:H	2.22	0.42
1:G:120:ARG:HG3	2:H:38:ASP:OD1	2.19	0.42
1:G:126:SER:HA	1:G:171:ILE:HG13	2.01	0.42
1:A:282:GLU:HG3	1:A:284:GLY:H	1.85	0.42
1:C:237:ILE:HB	1:C:258:LEU:HB3	2.00	0.42
2:D:41:ARG:HH11	2:D:52:LEU:HD21	1.84	0.42
1:E:300:ALA:HA	1:E:301:PRO:HD3	1.80	0.42
1:A:189:LYS:HE2	1:A:191:TYR:OH	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:177:TRP:CD2	1:C:183:ASN:HB2	2.54	0.42
2:F:161:ARG:NH2	1:G:296:LYS:HD2	2.35	0.42
2:D:22:GLN:HG3	2:D:149:ARG:HG2	2.00	0.42
2:B:33:ASP:OD1	1:E:354:TYR:HE1	2.03	0.42
1:G:174:LEU:HD11	1:G:185:LEU:HD11	2.02	0.42
1:C:165:GLU:HG2	1:C:346:LEU:CD2	2.51	0.41
2:F:69:ASP:O	2:F:73:ARG:HG3	2.20	0.41
2:H:160:VAL:O	2:H:164:ARG:HG3	2.20	0.41
2:B:56:LEU:HD23	2:B:71:TYR:HB2	2.02	0.41
1:G:292:LEU:HB2	1:G:293:ALA:H	1.66	0.41
1:A:282:GLU:HA	1:A:282:GLU:OE1	2.20	0.41
2:B:42:LYS:HE3	2:B:44:VAL:HG12	2.02	0.41
2:B:49:GLU:CB	2:B:164:ARG:HH22	2.27	0.41
2:F:49:GLU:HB3	2:F:164:ARG:NH2	2.36	0.41
2:H:99:GLN:O	2:H:103:VAL:HG23	2.20	0.41
2:D:157:TYR:N	2:D:157:TYR:HD1	2.17	0.41
1:A:234:TYR:CZ	1:A:340:LEU:HB2	2.56	0.41
1:C:295:LYS:HB2	1:C:298:HIS:CD2	2.56	0.41
1:C:324:ALA:HB1	1:C:329:SER:HB3	2.03	0.41
2:D:116:ASN:CG	2:D:117:LYS:N	2.75	0.41
1:G:205:PRO:HD3	1:G:332:CYS:SG	2.61	0.40
1:G:209:VAL:HG12	1:G:210:SER:N	2.36	0.40
1:C:204:PHE:HB3	1:C:209:VAL:HG23	2.02	0.40
1:C:220:SER:HA	1:C:221:PRO:HD3	1.97	0.40
1:C:224:ILE:O	1:C:228:PHE:HD2	2.04	0.40
1:E:118:THR:HB	2:F:39:SER:O	2.22	0.40
2:F:126:GLU:N	2:F:129:GLN:OE1	2.33	0.40
2:H:24:ILE:CD1	2:H:42:LYS:HB2	2.51	0.40
2:H:79:LEU:HD12	2:H:112:VAL:HB	2.03	0.40
1:A:118:THR:HG21	3:B:201:GOL:C1	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	240/258 (93%)	223 (93%)	15 (6%)	2 (1%)	19	29
1	C	240/258 (93%)	222 (92%)	14 (6%)	4 (2%)	9	11
1	E	240/258 (93%)	222 (92%)	13 (5%)	5 (2%)	7	8
1	G	240/258 (93%)	213 (89%)	19 (8%)	8 (3%)	4	3
2	B	169/171 (99%)	163 (96%)	6 (4%)	0	100	100
2	D	169/171 (99%)	166 (98%)	2 (1%)	1 (1%)	25	36
2	F	169/171 (99%)	161 (95%)	6 (4%)	2 (1%)	13	19
2	H	169/171 (99%)	160 (95%)	7 (4%)	2 (1%)	13	19
All	All	1636/1716 (95%)	1530 (94%)	82 (5%)	24 (2%)	10	14

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	293	ALA
1	E	248	LYS
1	C	298	HIS
2	D	-3	ALA
1	E	293	ALA
1	E	298	HIS
1	G	246	GLN
2	H	-1	GLY
2	H	122	ALA
1	C	314	GLY
1	G	295	LYS
1	G	301	PRO
1	G	312	LYS
1	A	292	LEU
1	A	297	LYS
1	E	246	GLN
1	G	200	PRO
1	G	292	LEU
1	G	298	HIS
1	E	297	LYS
2	F	-1	GLY
2	F	37	GLU
1	G	355	GLN
1	C	292	LEU

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	219/228 (96%)	215 (98%)	4 (2%)	59 76
1	C	219/228 (96%)	210 (96%)	9 (4%)	30 48
1	E	219/228 (96%)	215 (98%)	4 (2%)	59 76
1	G	219/228 (96%)	215 (98%)	4 (2%)	59 76
2	B	147/147 (100%)	144 (98%)	3 (2%)	55 74
2	D	147/147 (100%)	143 (97%)	4 (3%)	44 65
2	F	147/147 (100%)	141 (96%)	6 (4%)	30 48
2	H	147/147 (100%)	143 (97%)	4 (3%)	44 65
All	All	1464/1500 (98%)	1426 (97%)	38 (3%)	46 66

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	170	VAL
1	A	179	ILE
1	A	230	SER
1	A	292	LEU
2	B	69	ASP
2	B	71	TYR
2	B	93	ILE
1	C	136	LEU
1	C	199	ASN
1	C	231	SER
1	C	270	THR
1	C	282	GLU
1	C	288	ILE
1	C	323	CYS
1	C	326	GLU
1	C	330	ARG
2	D	2	THR
2	D	51	CYS
2	D	71	TYR

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Mol	Chain	Res	Type
2	D	93	ILE
1	E	114	SER
1	E	220	SER
1	E	254	ILE
1	E	311	ASN
2	F	31	GLU
2	F	71	TYR
2	F	87	THR
2	F	107	ASP
2	F	129	GLN
2	F	166	HIS
1	G	142	HIS
1	G	165	GLU
1	G	267	THR
1	G	326	GLU
2	H	33	ASP
2	H	49	GLU
2	H	67	MET
2	H	108	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	345	GLN
1	A	349	ASN
1	C	298	HIS
1	C	345	GLN
1	C	349	ASN
1	E	207	HIS
1	E	246	GLN
1	E	311	ASN
1	E	345	GLN
1	E	349	ASN
1	E	352	HIS
2	F	61	GLN
2	F	150	GLN
1	G	328	GLN
1	G	349	ASN
1	G	355	GLN

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 17 ligands modelled in this entry, 4 are monoatomic - leaving 13 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	GOL	C	400	-	5,5,5	0.38	0	5,5,5	0.60	0
3	GOL	E	400	-	5,5,5	0.31	0	5,5,5	0.51	0
3	GOL	A	403	-	5,5,5	0.59	0	5,5,5	0.64	0
3	GOL	C	401	-	5,5,5	0.21	0	5,5,5	0.49	0
3	GOL	A	402	-	5,5,5	0.36	0	5,5,5	0.46	0
4	GTP	B	202	5	26,34,34	1.01	1 (3%)	32,54,54	1.06	3 (9%)
3	GOL	B	201	-	5,5,5	0.43	0	5,5,5	0.58	0
3	GOL	E	401	-	5,5,5	0.29	0	5,5,5	0.90	0
4	GTP	D	200	5	26,34,34	0.94	1 (3%)	32,54,54	1.17	4 (12%)
3	GOL	A	401	-	5,5,5	0.40	0	5,5,5	0.68	0
4	GTP	H	200	5	26,34,34	0.89	1 (3%)	32,54,54	1.24	4 (12%)
4	GTP	F	200	5	26,34,34	0.99	1 (3%)	32,54,54	1.14	4 (12%)
3	GOL	A	404	-	5,5,5	0.61	0	5,5,5	0.89	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	GOL	C	400	-	-	2/4/4/4	-
3	GOL	E	400	-	-	2/4/4/4	-
3	GOL	A	403	-	-	0/4/4/4	-
3	GOL	C	401	-	-	2/4/4/4	-
3	GOL	A	402	-	-	4/4/4/4	-
4	GTP	B	202	5	-	2/18/38/38	0/3/3/3
3	GOL	B	201	-	-	2/4/4/4	-
3	GOL	E	401	-	-	2/4/4/4	-
4	GTP	D	200	5	-	2/18/38/38	0/3/3/3
3	GOL	A	401	-	-	2/4/4/4	-
4	GTP	H	200	5	-	2/18/38/38	0/3/3/3
4	GTP	F	200	5	-	2/18/38/38	0/3/3/3
3	GOL	A	404	-	-	0/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	202	GTP	C6-N1	-3.05	1.33	1.37
4	F	200	GTP	C6-N1	-2.33	1.34	1.37
4	D	200	GTP	C6-N1	-2.30	1.34	1.37
4	H	200	GTP	C6-N1	-2.18	1.34	1.37

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	200	GTP	PA-O3A-PB	-3.19	121.89	132.83
4	D	200	GTP	PA-O3A-PB	-2.72	123.48	132.83
4	F	200	GTP	C8-N7-C5	2.58	107.91	102.99
4	F	200	GTP	PA-O3A-PB	-2.53	124.14	132.83
4	F	200	GTP	PB-O3B-PG	-2.48	124.33	132.83
4	D	200	GTP	PB-O3B-PG	-2.40	124.58	132.83
4	H	200	GTP	PB-O3B-PG	-2.37	124.69	132.83
4	H	200	GTP	C5-C6-N1	2.36	118.12	113.95
4	B	202	GTP	O3G-PG-O2G	2.30	116.41	107.64
4	D	200	GTP	C8-N7-C5	2.29	107.36	102.99
4	B	202	GTP	PA-O3A-PB	-2.29	124.96	132.83
4	B	202	GTP	PB-O3B-PG	-2.22	125.20	132.83
4	H	200	GTP	C8-N7-C5	2.21	107.20	102.99
4	D	200	GTP	C5-C6-N1	2.15	117.75	113.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	200	GTP	O3G-PG-O2G	2.11	115.69	107.64

There are no chirality outliers.

All (24) torsion outliers are listed below:

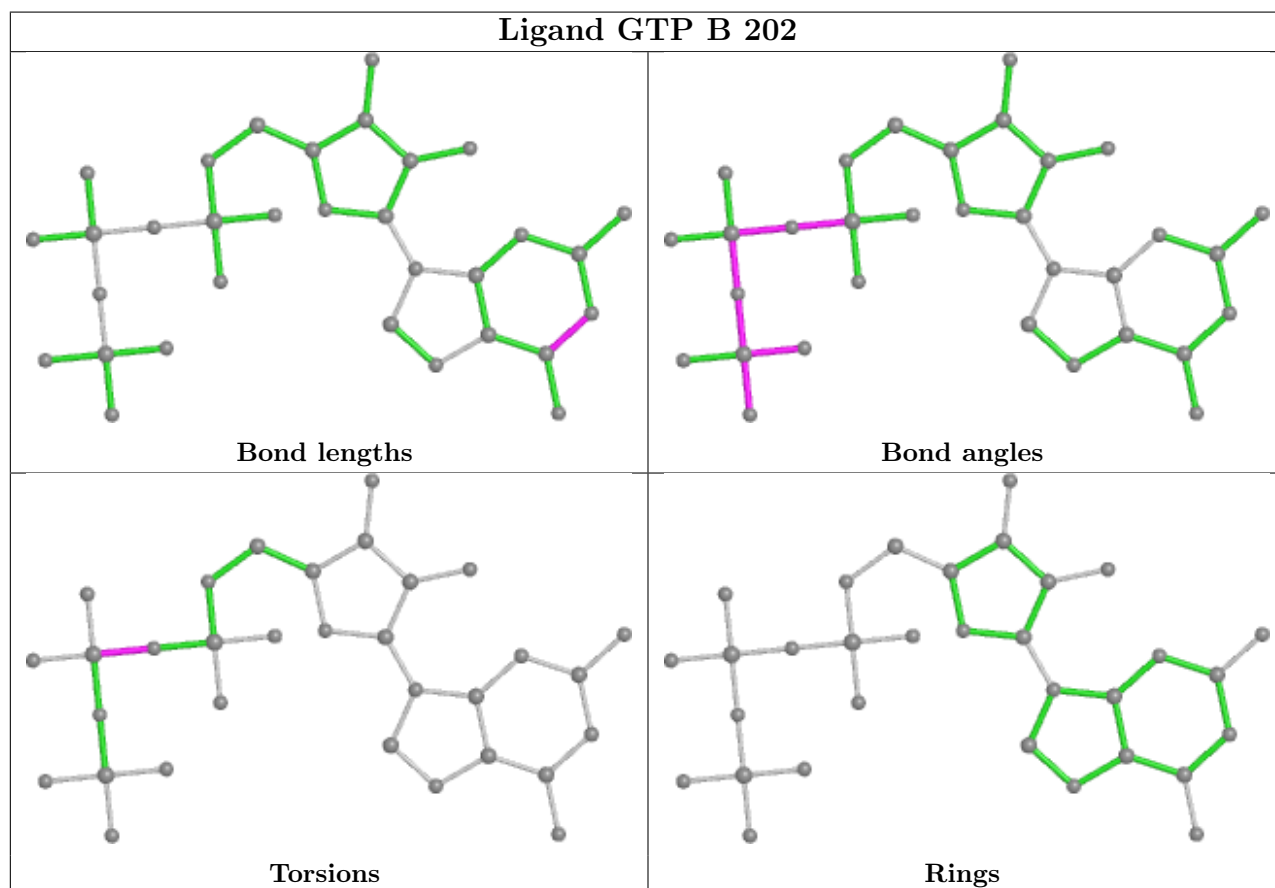
Mol	Chain	Res	Type	Atoms
3	A	401	GOL	C1-C2-C3-O3
3	A	402	GOL	O1-C1-C2-C3
3	A	402	GOL	C1-C2-C3-O3
3	B	201	GOL	O1-C1-C2-C3
3	C	400	GOL	C1-C2-C3-O3
3	C	401	GOL	O1-C1-C2-C3
3	A	401	GOL	O2-C2-C3-O3
3	A	402	GOL	O1-C1-C2-O2
3	E	401	GOL	O1-C1-C2-C3
3	A	402	GOL	O2-C2-C3-O3
3	B	201	GOL	O1-C1-C2-O2
3	C	400	GOL	O2-C2-C3-O3
3	E	401	GOL	O1-C1-C2-O2
3	C	401	GOL	O1-C1-C2-O2
4	B	202	GTP	PA-O3A-PB-O2B
4	D	200	GTP	PA-O3A-PB-O2B
4	F	200	GTP	PA-O3A-PB-O2B
4	H	200	GTP	PA-O3A-PB-O1B
4	H	200	GTP	PA-O3A-PB-O2B
3	E	400	GOL	O2-C2-C3-O3
4	B	202	GTP	PA-O3A-PB-O1B
4	D	200	GTP	PA-O3A-PB-O1B
4	F	200	GTP	PA-O3A-PB-O1B
3	E	400	GOL	C1-C2-C3-O3

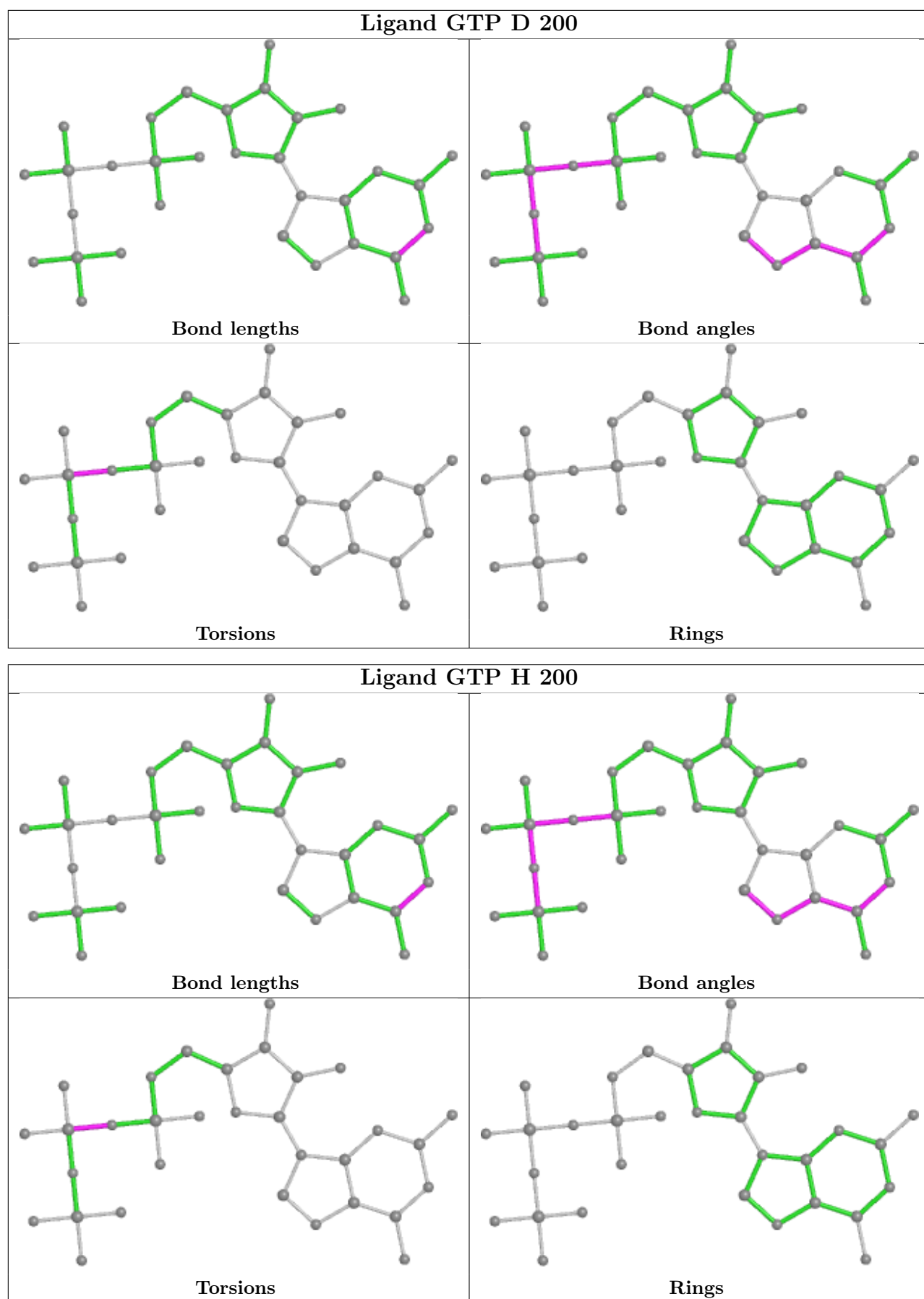
There are no ring outliers.

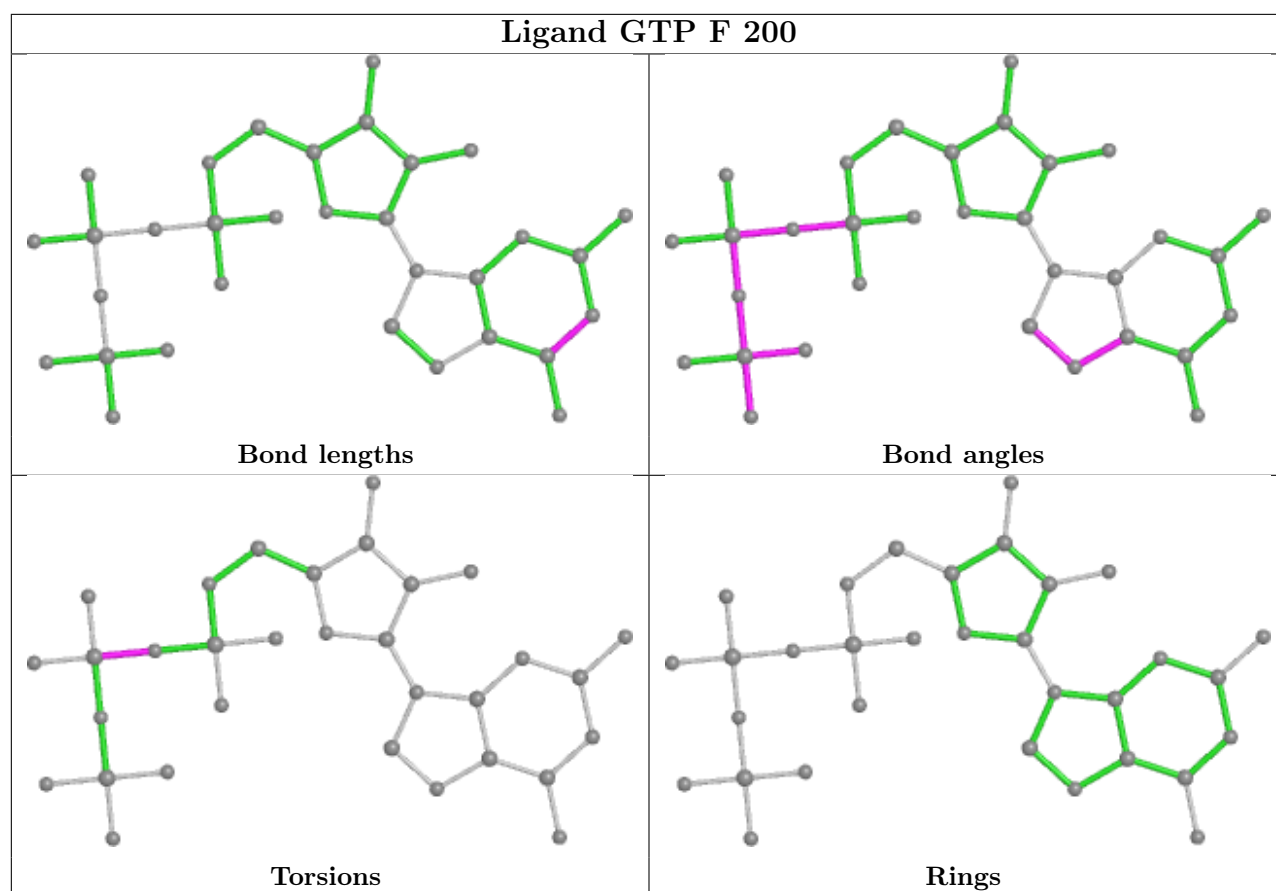
5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	402	GOL	1	0
3	B	201	GOL	1	0
3	A	401	GOL	2	0
4	F	200	GTP	1	0
3	A	404	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	244/258 (94%)	0.19	9 (3%) 41 41	26, 46, 87, 107	0
1	C	244/258 (94%)	0.33	14 (5%) 23 22	34, 57, 110, 127	0
1	E	244/258 (94%)	0.37	16 (6%) 18 17	29, 55, 95, 136	0
1	G	244/258 (94%)	1.45	68 (27%) 0 0	43, 93, 159, 187	0
2	B	171/171 (100%)	0.12	3 (1%) 68 66	29, 52, 85, 112	0
2	D	171/171 (100%)	0.02	1 (0%) 89 88	29, 47, 73, 103	0
2	F	171/171 (100%)	0.30	4 (2%) 60 58	38, 63, 95, 121	0
2	H	171/171 (100%)	1.06	23 (13%) 3 2	47, 90, 156, 196	0
All	All	1660/1716 (96%)	0.50	138 (8%) 11 10	26, 60, 129, 196	0

All (138) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	257	PHE	9.2
1	G	271	SER	8.5
1	G	297	LYS	8.2
1	G	246	GLN	7.0
1	G	265	PHE	6.9
1	G	315	GLY	6.9
1	G	264	TYR	6.7
1	E	313	ALA	6.4
1	G	207	HIS	6.3
1	G	247	GLY	6.2
2	F	-1	GLY	6.1
1	G	266	SER	6.0
1	C	211	PHE	6.0
1	E	246	GLN	6.0
1	G	270	THR	5.9
1	E	211	PHE	5.9

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Mol	Chain	Res	Type	RSRZ
1	E	316	PRO	5.9
1	E	315	GLY	5.8
1	G	219	ILE	5.8
1	G	245	GLU	5.6
1	G	277	LEU	5.4
1	G	248	LYS	5.4
1	G	240	PHE	5.4
1	C	246	GLN	5.3
1	G	221	PRO	5.2
2	H	90	PHE	5.1
2	H	130	ALA	5.1
2	H	109	VAL	5.0
1	G	268	LYS	5.0
1	E	312	LYS	4.9
1	G	313	ALA	4.8
1	G	272	ALA	4.8
1	A	297	LYS	4.8
2	H	96	TYR	4.8
2	H	-2	MET	4.6
2	H	125	VAL	4.5
1	G	333	TRP	4.4
1	G	276	HIS	4.4
1	G	280	PHE	4.4
1	G	273	ALA	4.3
1	E	297	LYS	4.3
1	G	296	LYS	4.3
2	H	126	GLU	4.2
1	G	308	PHE	4.1
1	E	317	ARG	4.0
1	G	209	VAL	3.9
2	H	104	LYS	3.8
1	G	298	HIS	3.8
2	H	100	ILE	3.7
1	C	245	GLU	3.7
1	C	297	LYS	3.6
1	E	314	GLY	3.5
2	B	-4	GLY	3.5
1	C	202	TYR	3.4
2	H	120	LEU	3.4
1	C	312	LYS	3.4
2	F	90	PHE	3.3
1	G	275	ARG	3.3

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Mol	Chain	Res	Type	RSRZ
1	G	220	SER	3.3
2	D	-4	GLY	3.3
1	G	310	PRO	3.2
2	H	122	ALA	3.2
1	G	287	ASP	3.2
2	H	60	GLY	3.2
1	C	221	PRO	3.2
1	G	208	MET	3.2
2	H	105	ASP	3.2
2	H	129	GLN	3.1
1	G	256	PHE	3.1
2	H	84	ILE	3.1
1	A	298	HIS	3.1
1	G	267	THR	3.1
1	G	236	GLU	3.1
1	G	210	SER	3.1
1	G	258	LEU	3.1
1	A	234	TYR	3.0
1	G	106	LYS	3.0
1	G	202	TYR	3.0
1	A	202	TYR	3.0
1	G	158	ILE	2.9
1	G	234	TYR	2.9
1	G	229	LEU	2.9
2	B	121	ALA	2.9
1	G	222	THR	2.9
2	H	89	SER	2.9
1	G	307	CYS	2.8
1	C	232	SER	2.8
1	G	312	LYS	2.8
1	E	240	PHE	2.8
1	C	316	PRO	2.8
1	G	318	ASP	2.8
1	G	279	PHE	2.7
1	A	356	GLY	2.7
1	G	314	GLY	2.7
1	G	241	LEU	2.7
1	C	219	ILE	2.7
1	G	317	ARG	2.6
2	H	73	ARG	2.6
1	E	210	SER	2.6
2	B	128	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
2	H	101	LYS	2.5
1	G	237	ILE	2.5
1	E	247	GLY	2.5
1	G	283	PHE	2.5
2	F	0	SER	2.5
1	G	223	GLN	2.5
1	G	255	TYR	2.5
1	G	206	GLU	2.5
1	A	247	GLY	2.5
1	G	306	PHE	2.4
1	C	315	GLY	2.4
1	G	278	GLN	2.4
2	H	-4	GLY	2.4
1	A	293	ALA	2.4
1	A	315	GLY	2.4
1	E	206	GLU	2.4
1	G	232	SER	2.4
2	H	94	HIS	2.3
1	E	202	TYR	2.3
2	H	85	ASN	2.3
1	C	247	GLY	2.3
1	G	180	GLU	2.3
1	G	108	GLN	2.2
2	F	84	ILE	2.2
2	H	114	VAL	2.2
1	C	248	LYS	2.2
1	G	263	LEU	2.2
1	E	236	GLU	2.2
1	A	201	MET	2.1
1	G	115	GLU	2.1
1	G	274	PRO	2.1
1	G	319	LEU	2.1
1	E	221	PRO	2.1
1	G	200	PRO	2.1
1	G	211	PHE	2.1
1	G	340	LEU	2.1
1	C	318	ASP	2.0
2	H	124	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 [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

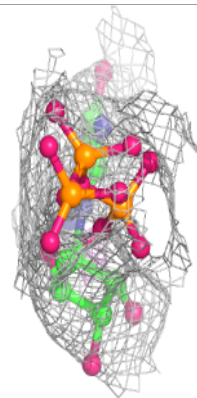
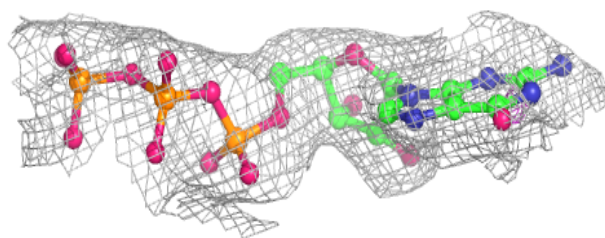
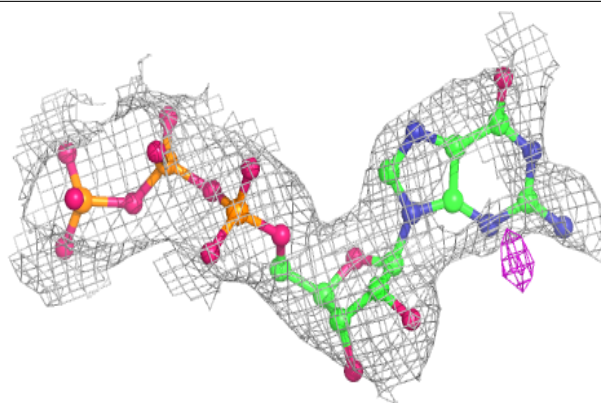
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GOL	C	400	6/6	0.75	0.38	65,67,68,70	0
3	GOL	A	403	6/6	0.78	0.34	49,51,55,56	0
3	GOL	A	404	6/6	0.84	0.15	44,50,54,55	0
5	MG	H	201	1/1	0.84	0.21	63,63,63,63	0
3	GOL	E	401	6/6	0.87	0.20	41,43,44,45	0
5	MG	D	201	1/1	0.90	0.18	45,45,45,45	0
3	GOL	C	401	6/6	0.91	0.18	40,43,44,44	0
3	GOL	B	201	6/6	0.91	0.19	46,54,55,61	0
3	GOL	A	402	6/6	0.92	0.18	32,34,34,37	0
3	GOL	A	401	6/6	0.93	0.25	46,47,48,51	0
4	GTP	H	200	32/32	0.94	0.18	69,80,91,93	0
3	GOL	E	400	6/6	0.95	0.15	41,43,44,47	0
5	MG	B	203	1/1	0.97	0.16	40,40,40,40	0
4	GTP	F	200	32/32	0.97	0.15	39,56,66,66	0
5	MG	F	201	1/1	0.97	0.22	54,54,54,54	0
4	GTP	B	202	32/32	0.97	0.15	34,41,52,53	0
4	GTP	D	200	32/32	0.98	0.13	38,46,50,51	0

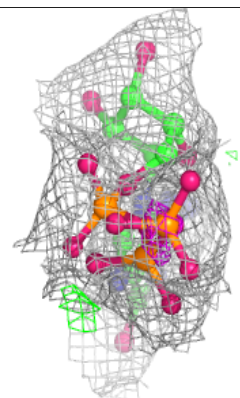
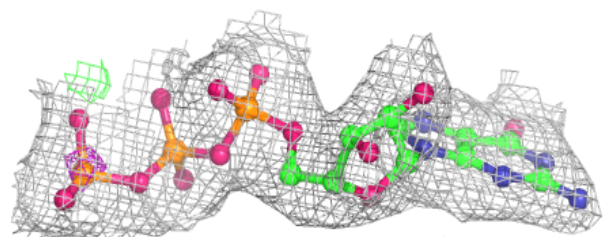
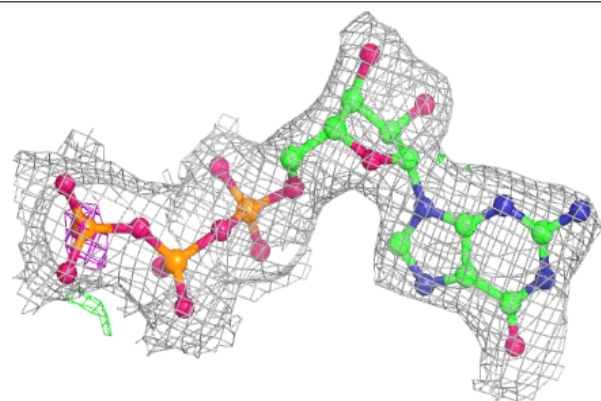
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around GTP H 200:

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

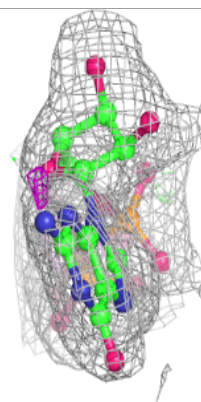
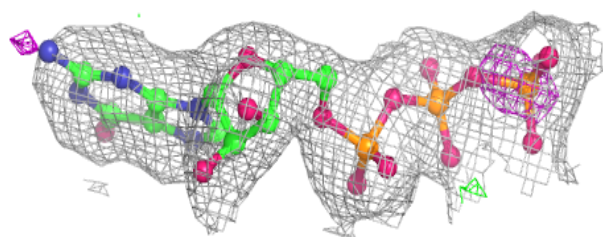
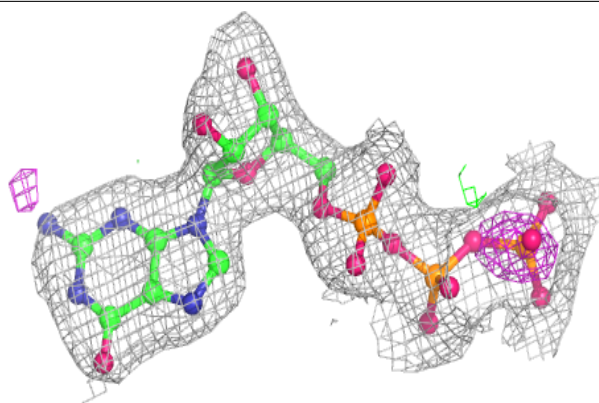
**Electron density around GTP F 200:**

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

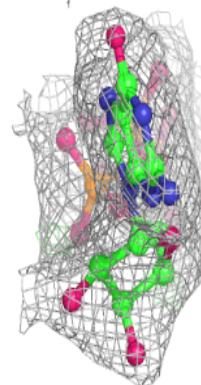
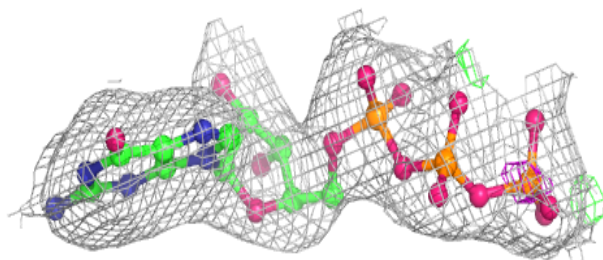
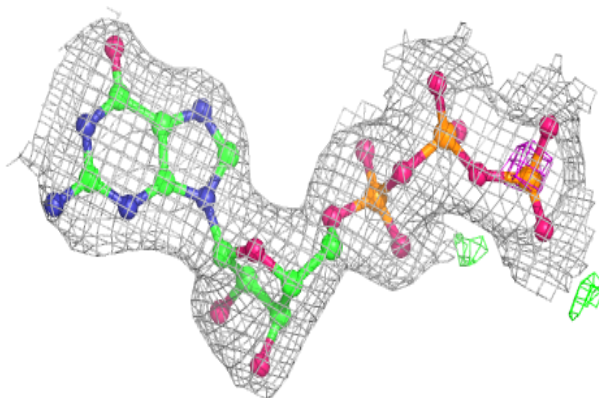


Electron density around GTP B 202:

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

**Electron density around GTP D 200:**

$2mF_o-DF_c$ (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.