



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2020 – 12:33 PM BST

PDB ID : 7BUJ
Title : mcGAS bound with pppGpG
Authors : Wang, B.; Su, X.D.
Deposited on : 2020-04-07
Resolution : 2.13 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13
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)	:	2.13

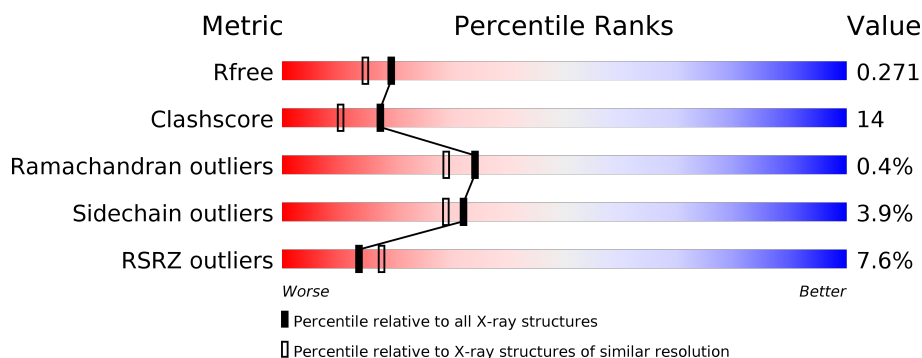
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.13 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	2523 (2.16-2.12)
Clashscore	141614	2653 (2.16-2.12)
Ramachandran outliers	138981	2618 (2.16-2.12)
Sidechain outliers	138945	2617 (2.16-2.12)
RSRZ outliers	127900	2485 (2.16-2.12)

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 on 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	447	
1	B	447	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6269 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.

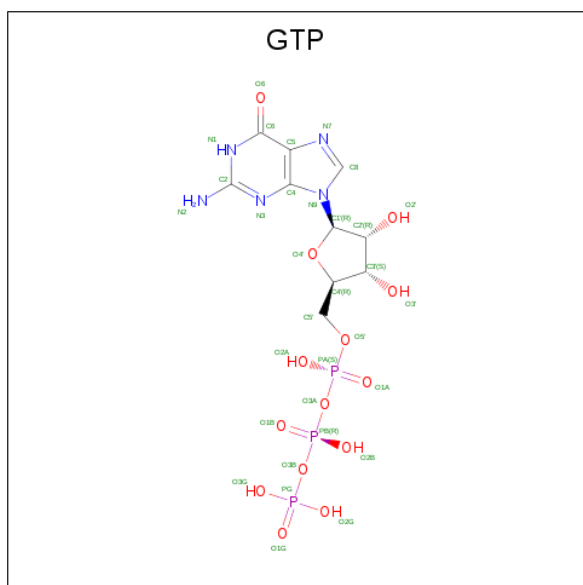
- Molecule 1 is a protein called Cyclic GMP-AMP synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	361	Total	C	N	O	S	0	0	0
			2984	1919	508	544	13			
1	B	362	Total	C	N	O	S	0	0	0
			2993	1924	509	547	13			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).

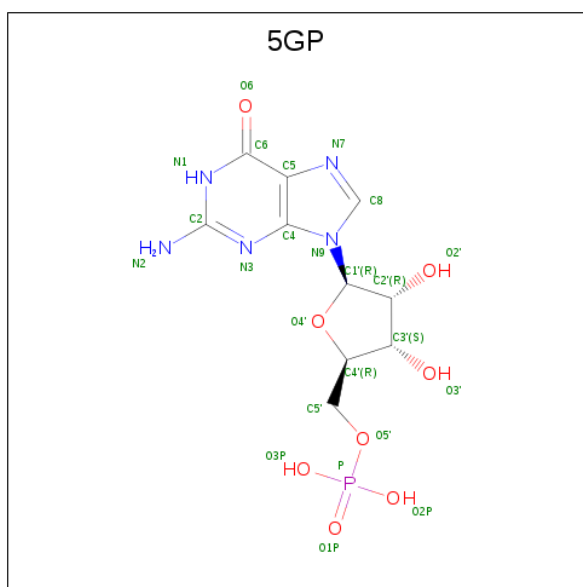


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
3	B	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

- Molecule 5 is GUANOSINE-5'-MONOPHOSPHATE (three-letter code: 5GP) (formula: C₁₀H₁₄N₅O₈P) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
5	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	91	Total	O	0	0
			91	91		

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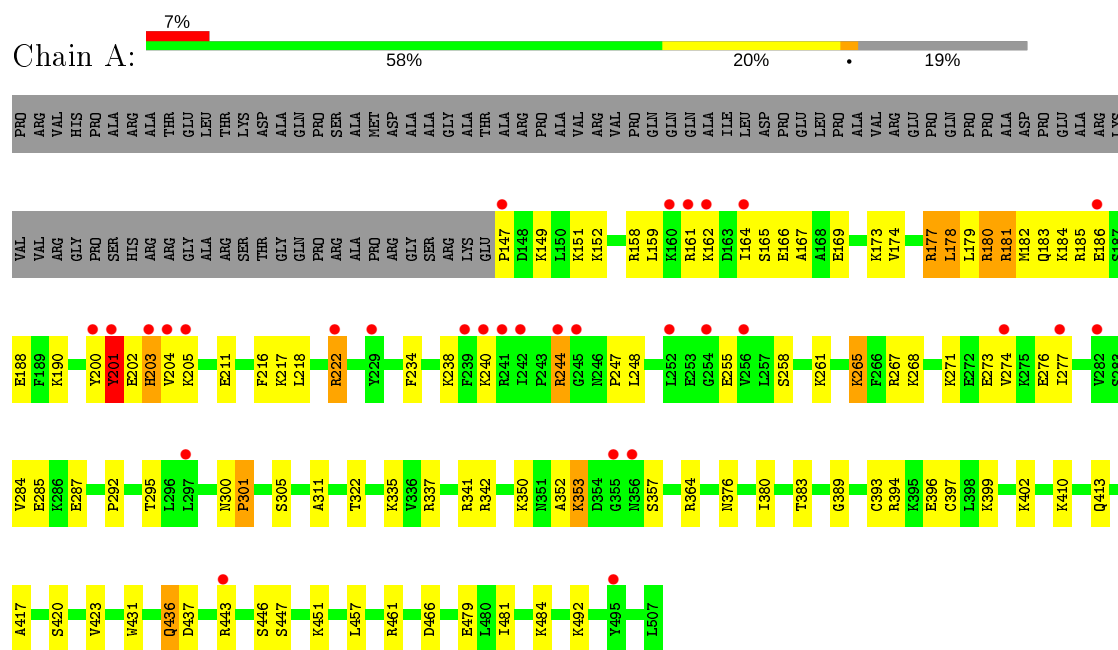
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	85	Total	O	0	0
			85	85		

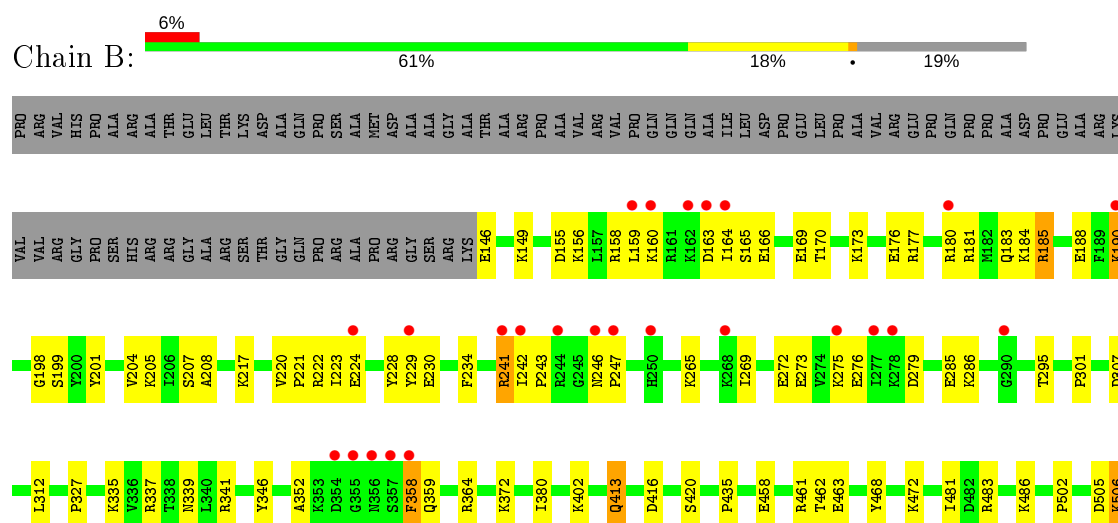
3 Residue-property plots [i](#)

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: Cyclic GMP-AMP synthase



• Molecule 1: Cyclic GMP-AMP synthase



L507

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	47.80 Å 109.63 Å 75.50 Å 90.00° 93.52° 90.00°	Depositor
Resolution (Å)	47.71 – 2.13 47.71 – 2.13	Depositor EDS
% Data completeness (in resolution range)	96.8 (47.71-2.13) 96.8 (47.71-2.13)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.14 Å)	Xtriage
Refinement program	PHENIX 1.14	Depositor
R, R_{free}	0.216 , 0.278 0.217 , 0.271	Depositor DCC
R_{free} test set	1983 reflections (4.72%)	wwPDB-VP
Wilson B-factor (Å ²)	49.7	Xtriage
Anisotropy	0.434	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 54.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6269	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.80% 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: GTP, ZN, 5GP, MN

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	0.60	0/3049	0.66	0/4093
1	B	0.50	0/3058	0.59	0/4106
All	All	0.55	0/6107	0.63	0/8199

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	2984	0	3034	100	0
1	B	2993	0	3039	73	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	32	0	11	2	0
3	B	32	0	11	0	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
5	A	23	0	12	1	0
5	B	23	0	11	0	0
6	A	91	0	0	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	85	0	0	7	0
All	All	6269	0	6118	169	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 14.

All (169) 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:300:ASN:CB	1:A:301:PRO:HD3	1.54	1.27
1:B:229:TYR:HB3	1:B:230:GLU:OE1	1.35	1.26
1:A:300:ASN:HB3	1:A:301:PRO:CD	1.66	1.21
1:A:159:LEU:HB2	1:A:203:HIS:CE1	1.93	1.02
1:B:352:ALA:O	1:B:359:GLN:HG3	1.68	0.93
1:A:177:ARG:HH12	1:B:180:ARG:CD	1.83	0.92
1:A:300:ASN:CB	1:A:301:PRO:CD	2.33	0.90
1:A:222:ARG:HD2	1:A:240:LYS:HD2	1.53	0.87
1:A:177:ARG:NE	1:A:180:ARG:HD2	1.93	0.83
1:A:161:ARG:HH11	1:A:164:ILE:HG21	1.45	0.81
1:B:158:ARG:HH22	1:B:205:LYS:HZ1	1.25	0.81
1:A:181:ARG:HH21	1:A:273:GLU:HB2	1.53	0.74
1:A:300:ASN:HB3	1:A:301:PRO:HD3	0.77	0.74
1:A:177:ARG:HH12	1:B:180:ARG:HD2	1.54	0.72
1:B:180:ARG:HE	1:B:184:LYS:HE3	1.56	0.71
1:B:229:TYR:CB	1:B:230:GLU:OE1	2.29	0.70
1:B:241:ARG:NH1	6:B:701:HOH:O	2.22	0.70
1:B:335:LYS:O	1:B:339:ASN:ND2	2.24	0.70
1:B:181:ARG:HH12	1:B:276:GLU:HB2	1.57	0.69
1:A:201:TYR:HA	1:A:204:VAL:HG23	1.75	0.69
1:A:397:CYS:SG	6:A:778:HOH:O	2.49	0.69
1:A:353:LYS:HD3	1:A:353:LYS:N	2.07	0.69
1:B:146:GLU:HB2	1:B:149:LYS:HB3	1.76	0.68
1:A:177:ARG:HH12	1:B:180:ARG:HD3	1.58	0.67
1:A:186:GLU:HG2	1:A:186:GLU:O	1.95	0.67
1:A:161:ARG:HA	1:A:164:ILE:HD12	1.75	0.67
1:B:461:ARG:NH2	1:B:462:THR:HG23	2.09	0.67
1:B:164:ILE:HD12	1:B:204:VAL:HG21	1.78	0.65
1:B:177:ARG:HH22	1:B:279:ASP:HB2	1.62	0.65
1:B:463:GLU:HG2	1:B:486:LYS:HD3	1.79	0.65
1:A:177:ARG:NH2	6:A:704:HOH:O	2.24	0.64
1:B:228:TYR:HE1	1:B:358:PHE:CD1	2.15	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:TYR:HA	1:A:204:VAL:CG2	2.29	0.62
1:B:458:GLU:CD	1:B:461:ARG:HH21	2.02	0.62
1:B:160:LYS:HB2	1:B:163:ASP:HB2	1.82	0.62
1:A:169:GLU:O	1:A:173:LYS:HG2	2.00	0.61
1:B:185:ARG:NH2	1:B:272:GLU:OE1	2.33	0.61
1:A:159:LEU:HA	6:A:702:HOH:O	1.99	0.61
1:B:173:LYS:O	1:B:177:ARG:HG3	2.00	0.61
1:A:267:ARG:CZ	1:A:287:GLU:HB3	2.31	0.60
1:A:188:GLU:N	1:A:188:GLU:OE1	2.27	0.60
1:B:181:ARG:NH2	1:B:273:GLU:HB3	2.17	0.60
1:B:461:ARG:HH22	1:B:462:THR:HG23	1.65	0.60
1:A:265:LYS:HD2	1:A:265:LYS:N	2.17	0.60
1:B:502:PRO:O	1:B:505:ASP:HB2	2.02	0.59
1:B:159:LEU:HB3	1:B:164:ILE:HD11	1.84	0.59
1:A:402:LYS:NZ	1:A:420:SER:O	2.26	0.59
1:B:156:LYS:NZ	6:B:704:HOH:O	2.35	0.59
1:B:435:PRO:HB2	6:B:726:HOH:O	2.02	0.58
1:B:507:LEU:N	1:B:507:LEU:HD22	2.18	0.58
1:A:180:ARG:NH1	6:A:707:HOH:O	2.30	0.58
1:A:410:LYS:NZ	6:A:709:HOH:O	2.37	0.58
1:B:217:LYS:NZ	6:B:705:HOH:O	2.37	0.58
1:A:204:VAL:O	1:A:204:VAL:HG12	2.04	0.57
1:B:243:PRO:HG2	1:B:246:ASN:HB2	1.86	0.57
1:B:341:ARG:HH11	1:B:341:ARG:HB3	1.68	0.57
1:A:200:TYR:C	1:A:202:GLU:H	2.06	0.57
1:B:265:LYS:O	1:B:269:ILE:HD12	2.04	0.57
1:A:376:ASN:ND2	6:A:710:HOH:O	2.38	0.57
1:A:201:TYR:C	1:A:201:TYR:HD1	2.09	0.55
1:A:174:VAL:O	1:A:178:LEU:HB2	2.05	0.55
1:A:178:LEU:O	1:A:182:MET:HG3	2.07	0.55
1:A:181:ARG:NH2	1:A:273:GLU:HB2	2.21	0.54
1:A:177:ARG:NH1	1:B:180:ARG:CD	2.64	0.54
1:A:167:ALA:HB3	1:A:200:TYR:CE2	2.43	0.53
1:B:472:LYS:HB2	6:B:722:HOH:O	2.08	0.53
1:B:285:GLU:HB3	1:B:295:THR:OG1	2.09	0.53
1:A:201:TYR:CD1	1:A:201:TYR:C	2.82	0.53
1:A:177:ARG:NH1	1:B:180:ARG:HD3	2.22	0.53
1:A:322:THR:OG1	1:A:341:ARG:NH2	2.42	0.53
1:A:159:LEU:HD22	6:A:702:HOH:O	2.07	0.53
1:A:201:TYR:O	1:A:201:TYR:HD1	1.90	0.53
1:A:147:PRO:N	1:A:446:SER:HG	2.06	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:341:ARG:NH1	1:B:341:ARG:HB3	2.25	0.51
3:A:602:GTP:O1A	5:A:605:5GP:H3'	2.10	0.51
1:A:265:LYS:HA	1:A:268:LYS:HB3	1.93	0.51
1:A:353:LYS:CD	1:A:353:LYS:N	2.72	0.51
1:B:188:GLU:HG3	1:B:247:PRO:HB2	1.93	0.51
1:A:161:ARG:NH1	1:A:164:ILE:HG21	2.23	0.50
1:A:177:ARG:CZ	1:A:180:ARG:HD2	2.40	0.50
1:A:222:ARG:HG3	1:A:240:LYS:HG3	1.91	0.50
1:A:205:LYS:O	1:A:399:LYS:HB2	2.12	0.50
1:B:242:ILE:HD12	1:B:242:ILE:H	1.76	0.50
1:B:158:ARG:NH2	1:B:205:LYS:HZ1	2.02	0.50
1:B:402:LYS:HD3	1:B:420:SER:HA	1.94	0.49
1:B:169:GLU:O	1:B:173:LYS:HG2	2.12	0.49
1:B:220:VAL:HG11	1:B:312:LEU:HD22	1.95	0.49
1:B:198:GLY:HA2	1:B:372:LYS:HZ2	1.78	0.49
1:A:217:LYS:HA	1:A:311:ALA:O	2.13	0.49
1:A:201:TYR:HB2	1:A:204:VAL:HB	1.95	0.48
1:A:265:LYS:HG3	1:A:268:LYS:HD3	1.94	0.48
1:A:276:GLU:O	1:A:277:ILE:HD12	2.13	0.48
1:A:394:ARG:HB3	1:A:431:TRP:CZ2	2.49	0.48
1:A:337:ARG:O	1:A:341:ARG:HG3	2.13	0.48
1:A:274:VAL:HG21	1:A:284:VAL:HG23	1.94	0.47
1:A:457:LEU:O	1:A:461:ARG:HG3	2.14	0.47
1:B:413:GLN:HG2	1:B:416:ASP:OD2	2.14	0.47
1:B:160:LYS:O	1:B:164:ILE:HG12	2.14	0.47
1:B:228:TYR:HE1	1:B:358:PHE:CE1	2.33	0.47
1:B:234:PHE:CD1	1:B:352:ALA:HB2	2.50	0.47
1:A:292:PRO:HD3	1:A:350:LYS:HD3	1.96	0.47
1:B:380:ILE:HB	6:B:726:HOH:O	2.13	0.47
1:A:402:LYS:HE3	1:A:420:SER:OG	2.15	0.47
1:A:380:ILE:HG21	1:A:436:GLN:HE21	1.80	0.47
1:B:181:ARG:HH12	1:B:276:GLU:CB	2.26	0.47
1:A:162:LYS:HE3	1:A:166:GLU:HG3	1.96	0.46
1:A:181:ARG:HD2	1:A:184:LYS:CD	2.46	0.46
1:B:286:LYS:HD2	1:B:286:LYS:HA	1.72	0.46
1:A:285:GLU:HG2	1:A:295:THR:HB	1.97	0.46
1:A:201:TYR:CA	1:A:204:VAL:HB	2.46	0.46
1:A:181:ARG:HD2	1:A:184:LYS:HD3	1.97	0.45
1:A:218:LEU:HD21	1:A:248:LEU:HD21	1.98	0.45
1:B:221:PRO:O	1:B:223:ILE:HG23	2.15	0.45
1:A:162:LYS:HA	1:A:165:SER:HB2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:GLU:OE2	1:B:265:LYS:HD3	2.17	0.45
1:B:483:ARG:HA	1:B:486:LYS:HD2	1.98	0.45
1:A:200:TYR:C	1:A:202:GLU:N	2.69	0.45
1:A:201:TYR:CD1	1:A:201:TYR:O	2.69	0.45
1:A:188:GLU:HA	1:A:247:PRO:HB2	1.98	0.45
1:B:228:TYR:CE1	1:B:358:PHE:CD1	3.02	0.45
1:A:300:ASN:HB2	1:A:301:PRO:HD3	1.78	0.45
1:A:258:SER:OG	1:A:261:LYS:HG3	2.17	0.44
1:A:271:LYS:NZ	6:A:716:HOH:O	2.50	0.44
1:A:364:ARG:NH2	3:A:602:GTP:O6	2.46	0.44
1:A:222:ARG:HD2	1:A:240:LYS:CD	2.37	0.44
1:A:238:LYS:HA	1:A:255:GLU:O	2.18	0.44
1:B:160:LYS:HB2	1:B:163:ASP:CB	2.47	0.44
1:B:337:ARG:O	1:B:341:ARG:HG3	2.17	0.44
1:B:234:PHE:CZ	1:B:364:ARG:HG3	2.52	0.44
1:A:185:ARG:HD2	1:A:190:LYS:NZ	2.32	0.44
1:A:159:LEU:HB2	1:A:203:HIS:HE1	1.67	0.44
1:A:211:GLU:HG2	1:A:305:SER:OG	2.18	0.43
1:A:393:CYS:HA	1:A:396:GLU:OE2	2.17	0.43
1:A:201:TYR:C	1:A:204:VAL:H	2.21	0.43
1:A:479:GLU:CD	1:A:479:GLU:H	2.22	0.43
1:B:199:SER:OG	1:B:201:TYR:HB2	2.19	0.43
1:B:234:PHE:CE1	1:B:364:ARG:HG3	2.54	0.43
1:A:149:LYS:HA	1:A:152:LYS:HD3	2.01	0.43
1:B:402:LYS:HE3	1:B:402:LYS:HB3	1.79	0.43
1:A:492:LYS:HD3	1:A:492:LYS:HA	1.93	0.42
1:B:295:THR:HG22	1:B:307:ASP:OD1	2.19	0.42
1:A:394:ARG:HD3	1:A:431:TRP:CD1	2.54	0.42
1:A:151:LYS:HB2	1:A:151:LYS:HE2	1.76	0.42
1:A:417:ALA:HB1	1:A:481:ILE:HG12	2.01	0.42
1:B:165:SER:O	1:B:169:GLU:HG3	2.19	0.42
1:A:179:LEU:O	1:A:183:GLN:HG3	2.20	0.42
1:A:466:ASP:OD1	6:A:701:HOH:O	2.21	0.42
1:B:312:LEU:O	1:B:346:TYR:HA	2.20	0.42
1:A:202:GLU:C	1:A:204:VAL:N	2.73	0.42
1:A:402:LYS:NZ	1:A:423:VAL:HB	2.35	0.42
1:B:506:LYS:O	1:B:506:LYS:CG	2.67	0.42
1:A:179:LEU:HD13	1:A:216:PHE:CZ	2.54	0.42
1:B:183:GLN:O	1:B:190:LYS:HE3	2.21	0.41
1:A:447:SER:O	1:A:451:LYS:HG3	2.20	0.41
1:B:481:ILE:HB	1:B:486:LYS:HE3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:TYR:O	1:A:202:GLU:N	2.54	0.41
1:A:234:PHE:CG	1:A:352:ALA:HB2	2.55	0.41
1:B:166:GLU:O	1:B:170:THR:HG22	2.20	0.41
1:A:436:GLN:NE2	1:A:437:ASP:H	2.19	0.41
1:B:230:GLU:OE1	1:B:230:GLU:N	2.49	0.41
1:B:380:ILE:HD12	6:B:726:HOH:O	2.20	0.41
1:A:244:ARG:HG3	1:A:244:ARG:H	1.72	0.41
1:A:413:GLN:HB2	1:A:413:GLN:HE21	1.59	0.40
1:B:506:LYS:O	1:B:506:LYS:HG3	2.21	0.40
1:A:383:THR:HB	1:A:389:GLY:HA3	2.03	0.40
1:B:207:SER:OG	1:B:208:ALA:N	2.55	0.40
1:A:335:LYS:HE2	1:A:335:LYS:HB2	1.92	0.40
1:B:327:PRO:HD2	1:B:468:TYR:CZ	2.57	0.40
1:A:158:ARG:O	6:A:702:HOH:O	2.22	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	359/447 (80%)	341 (95%)	16 (4%)	2 (1%)	25	17
1	B	360/447 (80%)	338 (94%)	21 (6%)	1 (0%)	41	36
All	All	719/894 (80%)	679 (94%)	37 (5%)	3 (0%)	34	29

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	301	PRO
1	B	301	PRO
1	A	201	TYR

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	334/400 (84%)	319 (96%)	15 (4%)	27	23
1	B	335/400 (84%)	324 (97%)	11 (3%)	38	35
All	All	669/800 (84%)	643 (96%)	26 (4%)	32	29

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	177	ARG
1	A	178	LEU
1	A	180	ARG
1	A	181	ARG
1	A	201	TYR
1	A	203	HIS
1	A	222	ARG
1	A	244	ARG
1	A	265	LYS
1	A	342	ARG
1	A	353	LYS
1	A	357	SER
1	A	436	GLN
1	A	443	ARG
1	A	484	LYS
1	B	155	ASP
1	B	176	GLU
1	B	185	ARG
1	B	190	LYS
1	B	222	ARG
1	B	224	GLU
1	B	241	ARG
1	B	275	LYS
1	B	358	PHE
1	B	413	GLN
1	B	506	LYS

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

Mol	Chain	Res	Type
1	A	413	GLN
1	A	436	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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 10 ligands modelled in this entry, 6 are monoatomic - leaving 4 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	5GP	A	605	3	18,25,26	5.02	12 (66%)	20,37,40	1.85	4 (20%)
3	GTP	B	602	5,4	26,34,34	4.26	18 (69%)	33,54,54	1.58	5 (15%)
5	5GP	B	605	3	18,25,26	5.12	13 (72%)	20,37,40	2.03	5 (25%)
3	GTP	A	602	5,4	26,34,34	4.83	10 (38%)	33,54,54	1.93	9 (27%)

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
5	5GP	A	605	3	-	2/3/25/26	0/3/3/3
3	GTP	B	602	5,4	-	0/18/38/38	0/3/3/3
5	5GP	B	605	3	-	1/3/25/26	0/3/3/3
3	GTP	A	602	5,4	-	1/18/38/38	0/3/3/3

All (53) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	GTP	O4'-C1'	14.91	1.61	1.41
3	B	602	GTP	O4'-C1'	11.90	1.57	1.41
3	B	602	GTP	O4'-C4'	-9.04	1.24	1.45
5	A	605	5GP	O4'-C1'	-8.91	1.28	1.41
5	A	605	5GP	C3'-C4'	-8.81	1.30	1.53
5	B	605	5GP	O4'-C1'	-8.79	1.28	1.41
3	A	602	GTP	C4-N3	8.74	1.49	1.35
5	B	605	5GP	C4-N3	8.43	1.48	1.35
5	B	605	5GP	C3'-C4'	-8.42	1.31	1.53
3	A	602	GTP	C6-C5	8.21	1.55	1.41
5	A	605	5GP	C4-N3	8.17	1.48	1.35
5	B	605	5GP	O4'-C4'	7.63	1.62	1.45
5	A	605	5GP	C6-C5	7.34	1.54	1.41
5	A	605	5GP	O4'-C4'	7.32	1.61	1.45
3	A	602	GTP	O4'-C4'	-6.99	1.29	1.45
5	B	605	5GP	C6-C5	6.98	1.53	1.41
3	B	602	GTP	C4-N3	6.77	1.46	1.35
5	B	605	5GP	C6-N1	6.73	1.44	1.33
3	A	602	GTP	C2-N2	6.71	1.47	1.33
3	B	602	GTP	C2'-C3'	-6.35	1.36	1.53
3	A	602	GTP	C6-N1	6.19	1.43	1.33
3	A	602	GTP	C2'-C3'	-5.89	1.37	1.53
5	B	605	5GP	C2-N1	5.69	1.45	1.35
5	B	605	5GP	C2-N2	5.59	1.45	1.33
5	A	605	5GP	C2-N1	5.45	1.45	1.35
3	A	602	GTP	C3'-C4'	5.44	1.66	1.53
5	A	605	5GP	C6-N1	5.35	1.42	1.33
5	A	605	5GP	C2-N2	5.17	1.44	1.33
3	B	602	GTP	C6-C5	4.92	1.49	1.41
3	B	602	GTP	C2-N2	4.88	1.43	1.33
3	A	602	GTP	C2-N1	4.87	1.44	1.35
3	B	602	GTP	C3'-C4'	4.48	1.64	1.53
5	A	605	5GP	O3'-C3'	3.80	1.51	1.43
3	B	602	GTP	C5-C4	-3.72	1.31	1.40
3	B	602	GTP	O6-C6	-3.57	1.15	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	602	GTP	C6-N1	3.32	1.38	1.33
5	B	605	5GP	O3'-C3'	3.12	1.50	1.43
3	B	602	GTP	C2-N1	2.67	1.40	1.35
5	B	605	5GP	O2'-C2'	-2.67	1.36	1.43
5	B	605	5GP	C2-N3	2.55	1.46	1.34
3	B	602	GTP	PG-O3G	-2.54	1.45	1.54
5	A	605	5GP	C2-N3	2.53	1.46	1.34
3	B	602	GTP	PB-O2B	-2.52	1.43	1.55
3	B	602	GTP	PG-O2G	-2.45	1.45	1.54
3	A	602	GTP	C2-N3	2.42	1.46	1.34
5	A	605	5GP	O5'-C5'	-2.36	1.39	1.44
3	B	602	GTP	PA-O1A	-2.27	1.42	1.50
5	B	605	5GP	O6-C6	-2.25	1.18	1.24
3	B	602	GTP	O2'-C2'	2.20	1.48	1.43
5	B	605	5GP	O5'-C5'	-2.20	1.39	1.44
3	B	602	GTP	C2-N3	2.18	1.44	1.34
5	A	605	5GP	O2'-C2'	-2.13	1.38	1.43
3	B	602	GTP	PA-O2A	-2.09	1.45	1.55

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	605	5GP	N3-C2-N1	-5.70	119.62	127.22
3	A	602	GTP	N3-C2-N1	-5.47	119.92	127.22
3	B	602	GTP	N3-C2-N1	-4.65	121.02	127.22
5	B	605	5GP	C2-N3-C4	4.61	120.62	115.36
5	A	605	5GP	N3-C2-N1	-4.58	121.11	127.22
3	A	602	GTP	C2-N3-C4	4.24	120.19	115.36
3	B	602	GTP	C2-N3-C4	3.95	119.86	115.36
3	A	602	GTP	C3'-C2'-C1'	3.46	106.18	100.98
5	A	605	5GP	C5-C6-N1	-3.44	118.72	123.43
3	A	602	GTP	C5-C6-N1	-3.16	119.11	123.43
3	B	602	GTP	O2G-PG-O3B	2.96	114.55	104.64
5	A	605	5GP	C2-N3-C4	2.84	118.61	115.36
3	B	602	GTP	PB-O3B-PG	-2.77	123.31	132.83
3	A	602	GTP	C6-N1-C2	2.74	120.29	115.93
5	A	605	5GP	C6-N1-C2	2.69	120.21	115.93
5	B	605	5GP	C5-C6-N1	-2.49	120.02	123.43
3	A	602	GTP	O4'-C1'-C2'	-2.25	103.64	106.93
3	A	602	GTP	N2-C2-N1	2.23	120.72	117.25
3	A	602	GTP	O2G-PG-O3B	2.23	112.10	104.64
5	B	605	5GP	C6-N1-C2	2.16	119.36	115.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	602	GTP	C3'-C2'-C1'	-2.15	97.73	100.98
5	B	605	5GP	N2-C2-N1	2.10	120.51	117.25
3	A	602	GTP	C4-C5-N7	-2.01	107.30	109.40

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	605	5GP	O4'-C4'-C5'-O5'
5	A	605	5GP	C3'-C4'-C5'-O5'
3	A	602	GTP	PA-O3A-PB-O1B
5	B	605	5GP	O4'-C4'-C5'-O5'

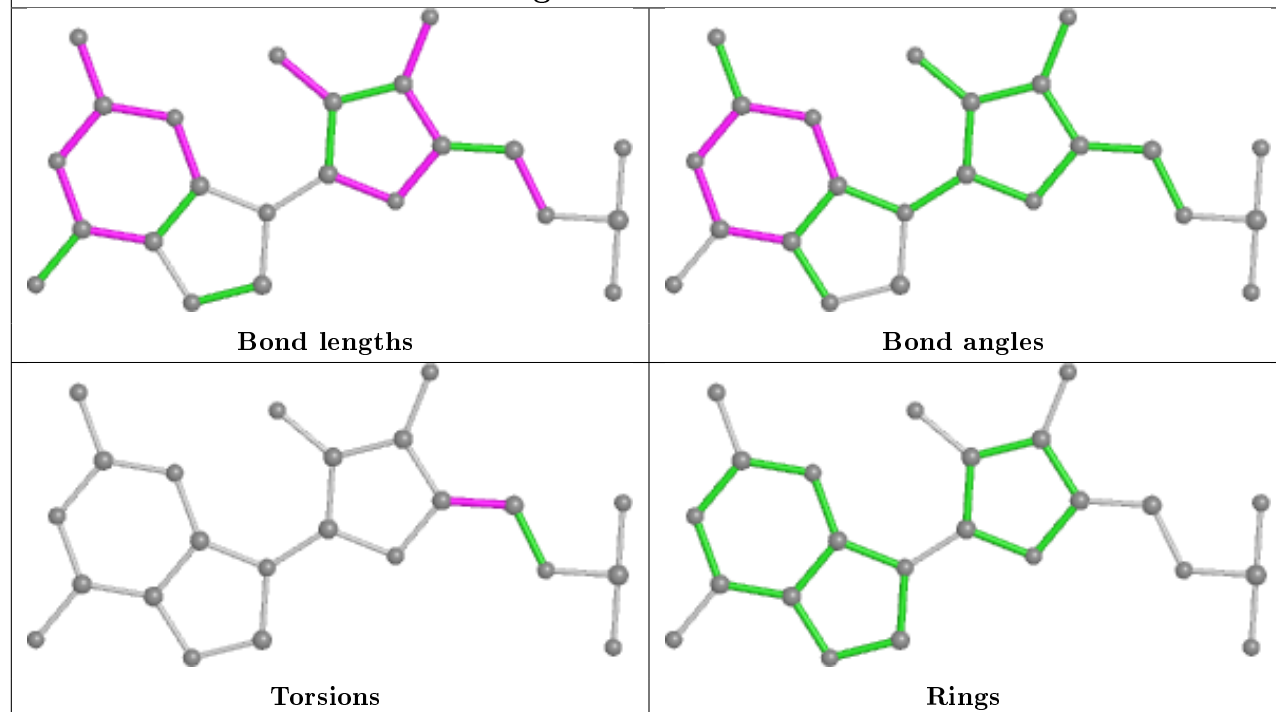
There are no ring outliers.

2 monomers are involved in 2 short contacts:

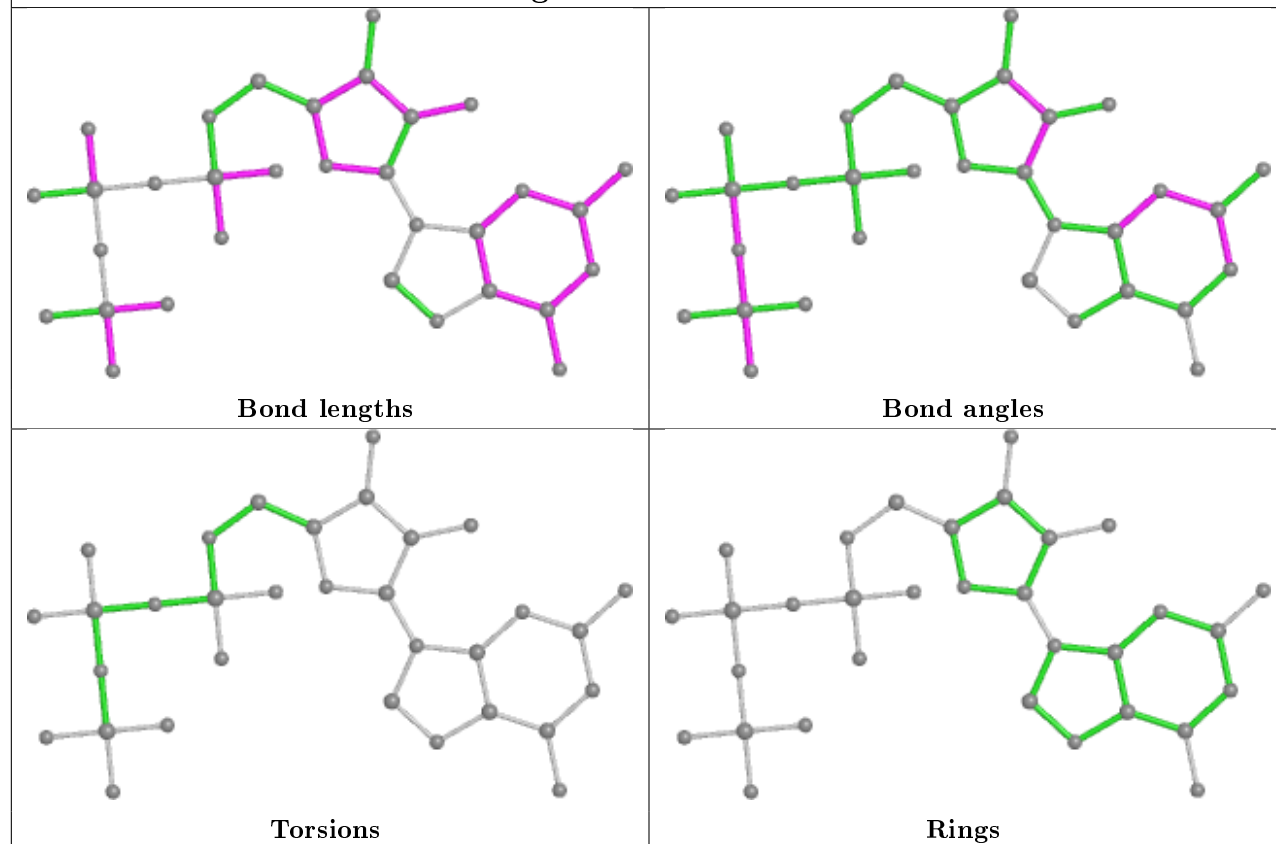
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	605	5GP	1	0
3	A	602	GTP	2	0

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.

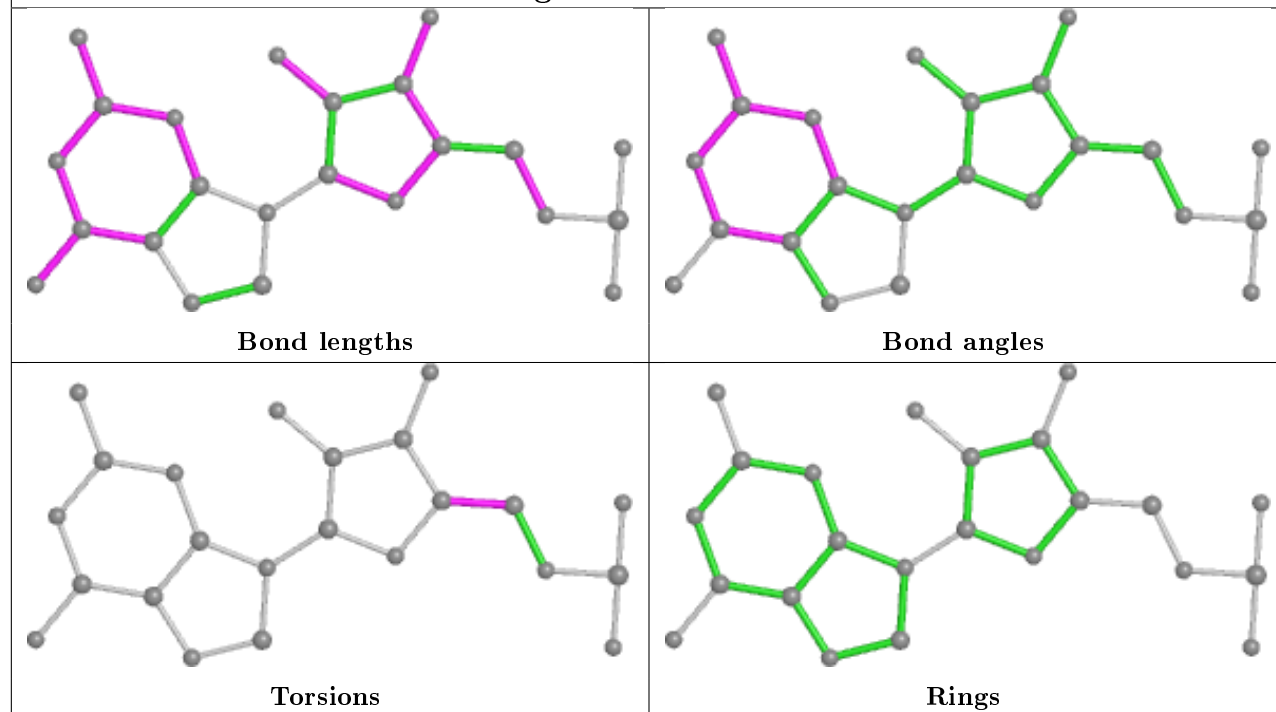
Ligand 5GP A 605



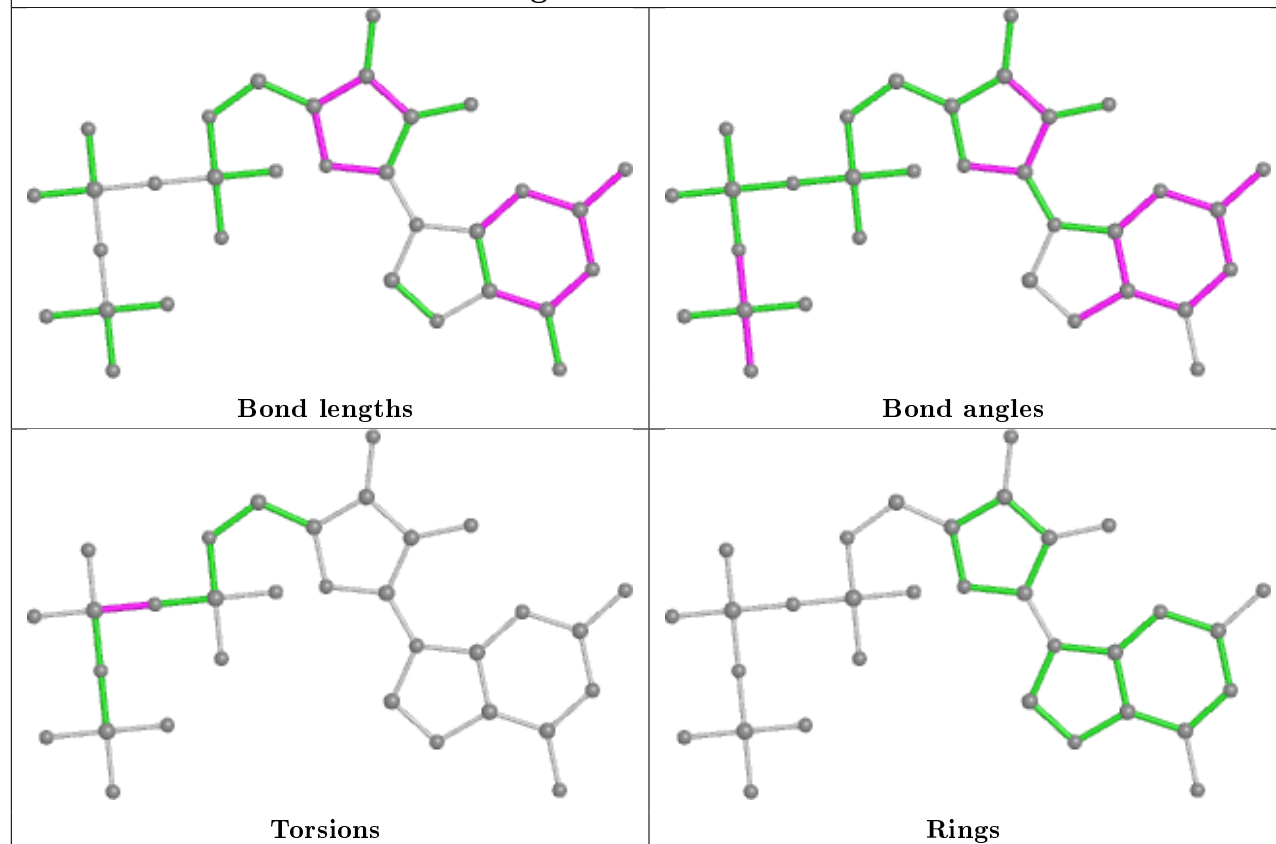
Ligand GTP B 602



Ligand 5GP B 605



Ligand GTP A 602



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	361/447 (80%)	0.71	30 (8%) 11 14	35, 61, 105, 147	0
1	B	362/447 (80%)	0.59	25 (6%) 16 21	38, 59, 105, 145	0
All	All	723/894 (80%)	0.65	55 (7%) 13 17	35, 59, 105, 147	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	205	LYS	7.0
1	A	240	LYS	7.0
1	B	357	SER	6.9
1	B	354	ASP	6.9
1	A	245	GLY	6.6
1	B	356	ASN	6.0
1	A	239	PHE	5.8
1	A	203	HIS	5.5
1	B	358	PHE	5.3
1	A	252	LEU	5.2
1	A	204	VAL	4.9
1	A	162	LYS	4.6
1	A	229	TYR	4.5
1	B	275	LYS	4.3
1	A	164	ILE	4.3
1	A	355	GLY	4.2
1	A	200	TYR	4.0
1	B	355	GLY	4.0
1	A	356	ASN	4.0
1	A	201	TYR	3.9
1	A	222	ARG	3.7
1	A	147	PRO	3.6
1	B	277	ILE	3.5
1	B	290	GLY	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	190	LYS	3.4
1	B	163	ASP	3.4
1	A	241	ARG	3.4
1	B	159	LEU	3.3
1	A	277	ILE	3.3
1	B	229	TYR	3.2
1	B	162	LYS	3.2
1	B	278	LYS	3.1
1	A	244	ARG	3.0
1	A	160	LYS	2.9
1	A	242	ILE	2.9
1	A	443	ARG	2.9
1	B	160	LYS	2.8
1	B	164	ILE	2.6
1	A	297	LEU	2.6
1	A	274	VAL	2.5
1	B	247	PRO	2.5
1	A	161	ARG	2.4
1	A	495	TYR	2.4
1	B	241	ARG	2.3
1	B	180	ARG	2.3
1	B	224	GLU	2.3
1	B	246	ASN	2.3
1	B	244	ARG	2.2
1	A	282	VAL	2.2
1	A	186	GLU	2.2
1	A	256	VAL	2.1
1	B	250	HIS	2.1
1	A	254	GLY	2.1
1	B	242	ILE	2.0
1	B	268	LYS	2.0

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

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

6.3 Carbohydrates ⓘ

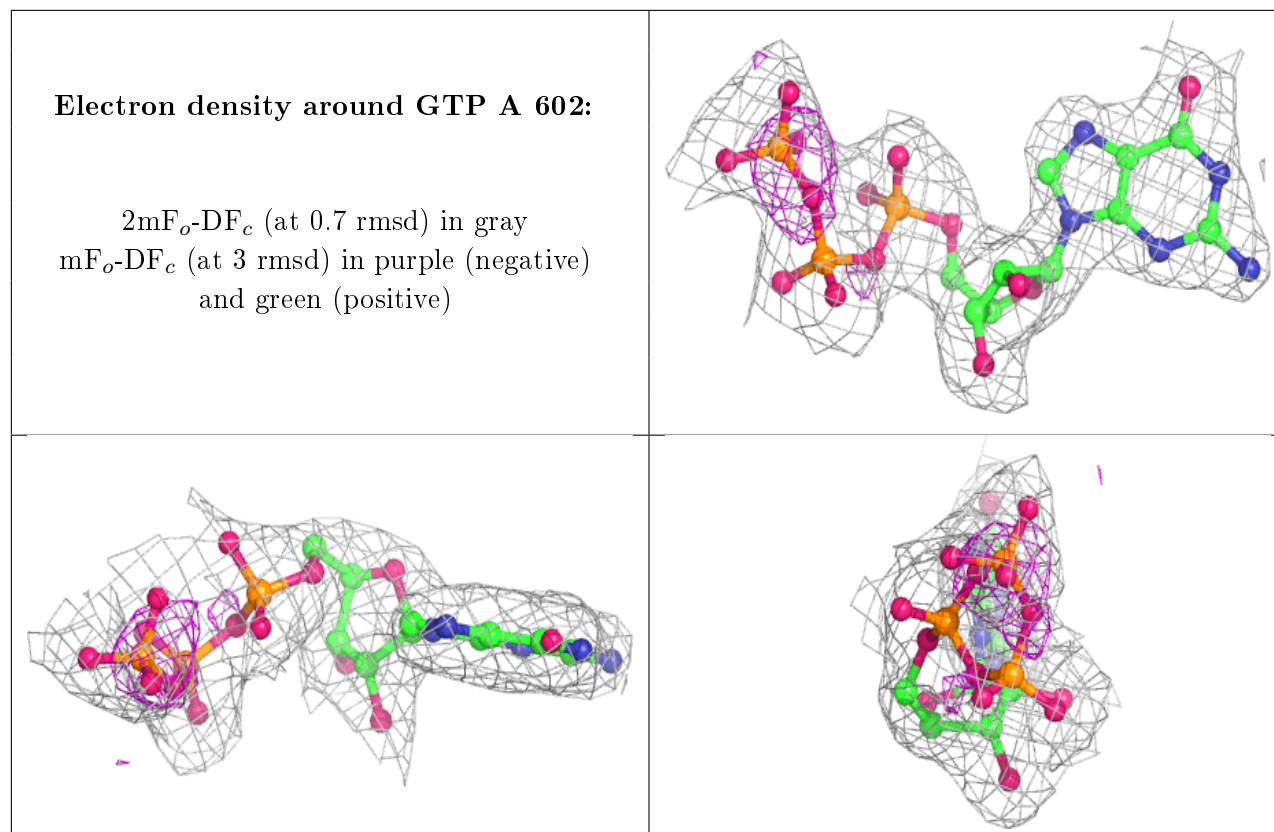
There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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

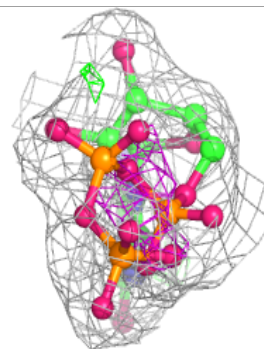
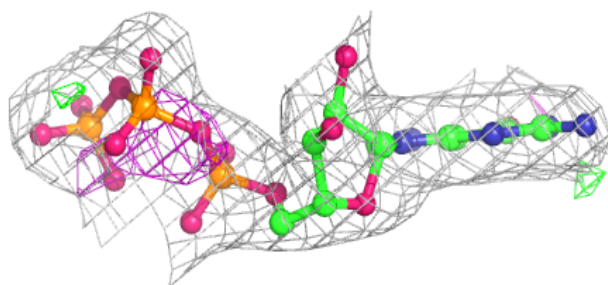
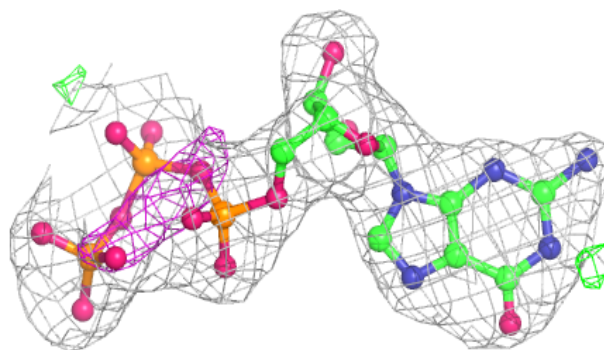
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MN	B	603	1/1	0.81	0.09	122,122,122,122	0
3	GTP	A	602	32/32	0.89	0.14	49,58,69,71	0
3	GTP	B	602	32/32	0.90	0.10	61,68,80,81	0
4	MN	B	604	1/1	0.90	0.05	106,106,106,106	0
4	MN	A	603	1/1	0.90	0.05	94,94,94,94	0
4	MN	A	604	1/1	0.91	0.07	105,105,105,105	0
5	5GP	A	605	23/24	0.93	0.14	48,52,55,57	0
5	5GP	B	605	23/24	0.95	0.12	55,58,62,63	0
2	ZN	A	601	1/1	0.99	0.15	46,46,46,46	0
2	ZN	B	601	1/1	1.00	0.11	45,45,45,45	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.

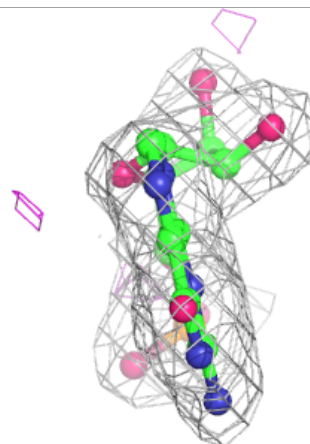
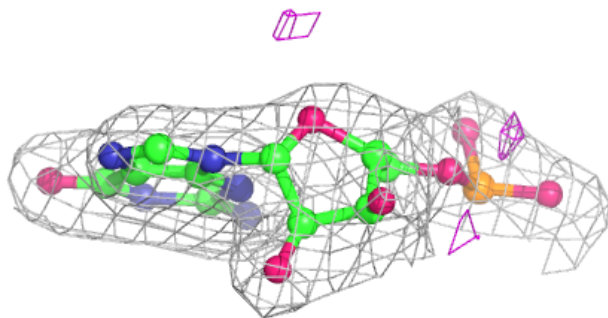
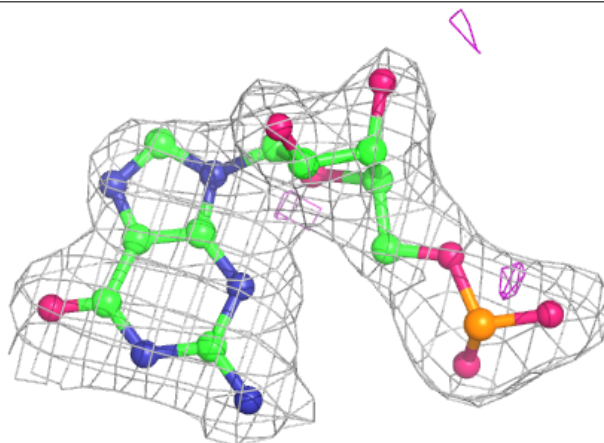


Electron density around GTP B 602:

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

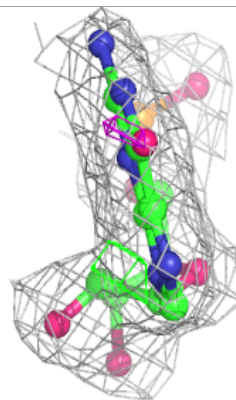
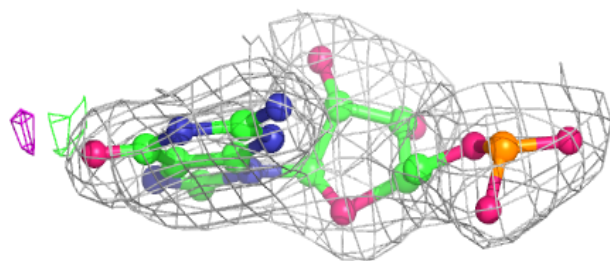
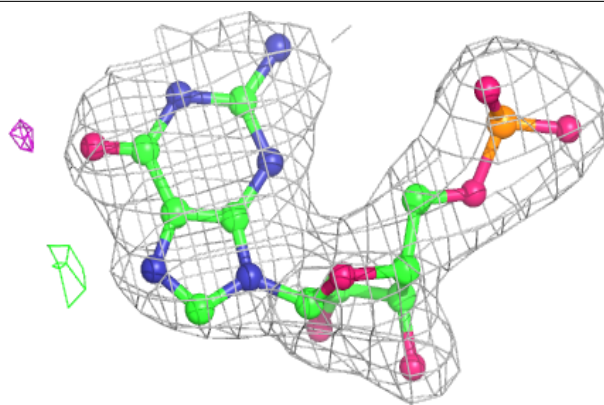
**Electron density around 5GP A 605:**

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



Electron density around 5GP B 605:

$2mF_o-DF_c$ (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.