



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2020 – 11:58 AM BST

PDB ID : 7BUM
Title : mcGAS bound with pGpA
Authors : Wang, B.; Su, X.D.
Deposited on : 2020-04-07
Resolution : 3.05 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

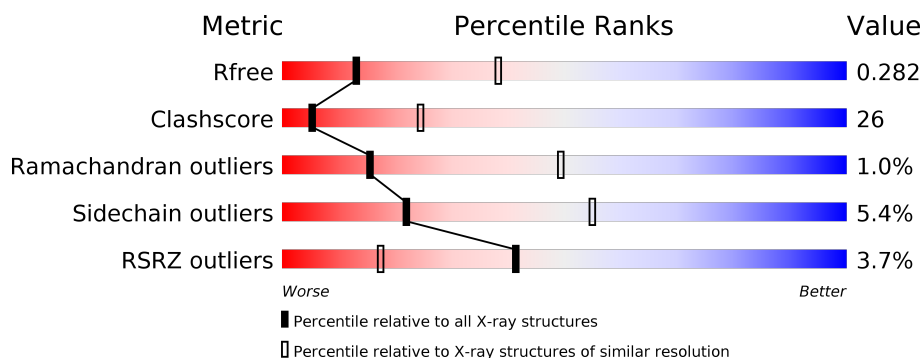
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.05 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	2752 (3.08-3.00)
Clashscore	141614	3096 (3.08-3.00)
Ramachandran outliers	138981	2986 (3.08-3.00)
Sidechain outliers	138945	2988 (3.08-3.00)
RSRZ outliers	127900	2636 (3.08-3.00)

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	507	<div> <div>2%</div> <div> <div></div> <div>38%</div> <div>31%</div> <div>•</div> <div>29%</div> </div> </div>
1	B	507	<div> <div>3%</div> <div> <div></div> <div>46%</div> <div>23%</div> <div>•</div> <div>29%</div> </div> </div>

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	A	A	602	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6099 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.

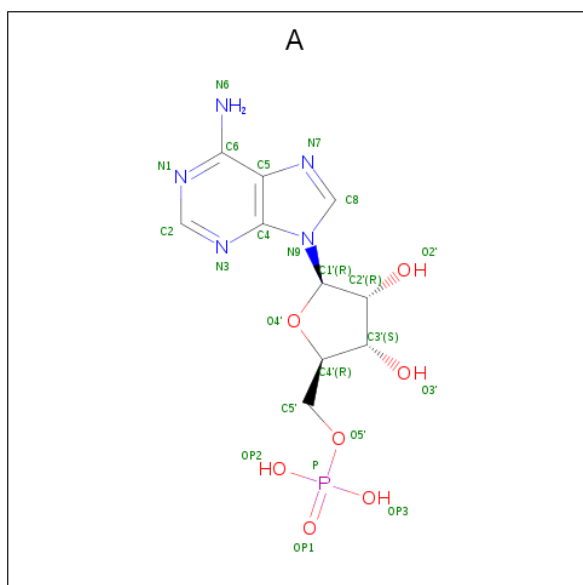
- Molecule 1 is a protein called Cyclic GMP-AMP synthase.

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	361	Total	C	N	O	S	0	0	0
			2984	1919	508	544	13			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by author).

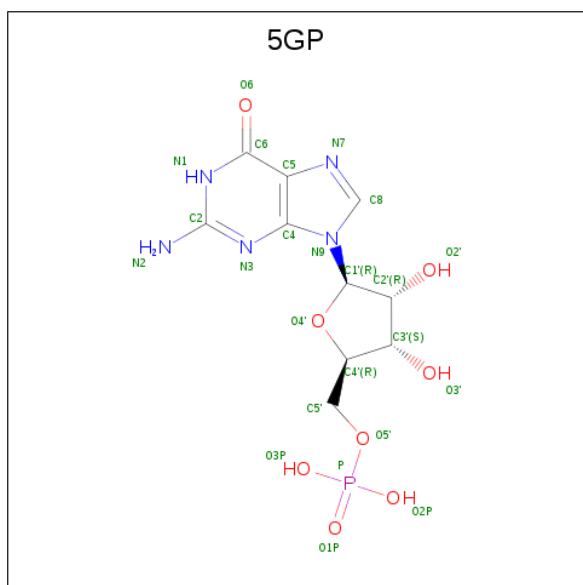
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is ADENOSINE-5'-MONOPHOSPHATE (three-letter code: A) (formula: C₁₀H₁₄N₅O₇P) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
3	B	1	Total	C	N	O	P	0	0
			22	10	5	6	1		

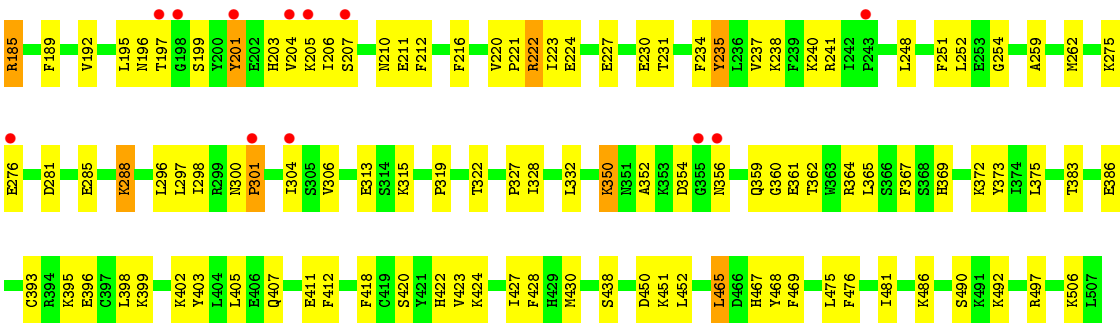
- Molecule 4 is GUANOSINE-5'-MONOPHOSPHATE (three-letter code: 5GP) (formula: $C_{10}H_{14}N_5O_8P$).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	21	Total	O	0	0
			21	21		
5	B	16	Total	O	0	0
			16	16		



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	48.34 Å 109.59 Å 75.87 Å 90.00° 93.93° 90.00°	Depositor
Resolution (Å)	30.82 – 3.05 30.82 – 3.05	Depositor EDS
% Data completeness (in resolution range)	95.3 (30.82-3.05) 95.3 (30.82-3.05)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 3.06 Å)	Xtriage
Refinement program	PHENIX 1.14	Depositor
R, R_{free}	0.217 , 0.280 0.218 , 0.282	Depositor DCC
R_{free} test set	1449 reflections (10.05%)	wwPDB-VP
Wilson B-factor (Å ²)	66.4	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 51.9	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	0.91	EDS
Total number of atoms	6099	wwPDB-VP
Average B, all atoms (Å ²)	68.0	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: ZN, 5GP

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	0.58	1/3049 (0.0%)	0.75	3/4093 (0.1%)
1	B	0.57	1/3049 (0.0%)	0.66	0/4093
All	All	0.57	2/6098 (0.0%)	0.71	3/8186 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	235	TYR	CE1-CZ	-5.68	1.31	1.38
1	A	397	CYS	CB-SG	-5.50	1.72	1.81

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	322	THR	CB-CA-C	-5.89	95.71	111.60
1	A	148	ASP	N-CA-C	-5.68	95.67	111.00
1	A	324	GLU	N-CA-CB	5.21	119.97	110.60

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	202	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2984	0	3034	129	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	22	0	12	1	0
3	B	22	0	12	0	0
4	A	24	0	11	0	0
4	B	24	0	11	2	0
5	A	21	0	0	14	0
5	B	16	0	0	10	0
All	All	6099	0	6114	321	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 26.

All (321) 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:204:VAL:CG2	1:A:398:LEU:HD23	1.59	1.30
1:B:259:ALA:HB3	1:B:360:GLY:O	1.50	1.12
1:A:169:GLU:HB2	1:B:201:TYR:OH	1.49	1.11
1:A:300:ASN:HB3	1:A:301:PRO:CD	1.77	1.11
1:A:293:ALA:HB2	1:A:309:ILE:HD13	1.31	1.10
1:A:300:ASN:CB	1:A:301:PRO:HD3	1.80	1.09
1:B:300:ASN:HB2	5:B:710:HOH:O	1.51	1.08
1:A:281:ASP:O	1:A:299:ARG:HB3	1.54	1.07
1:A:431:TRP:HB3	5:A:704:HOH:O	1.50	1.07
1:B:428:PHE:HB3	1:B:468:TYR:CE1	1.91	1.06
1:A:204:VAL:HG22	1:A:398:LEU:HD23	1.05	1.02
1:A:323:LYS:HA	1:A:341:ARG:NH2	1.75	1.01
1:A:477:SER:HB2	5:A:709:HOH:O	1.58	1.00
1:A:204:VAL:HG22	1:A:398:LEU:CD2	1.96	0.96
1:B:156:LYS:HA	1:B:159:LEU:HD23	1.44	0.96
1:B:156:LYS:HA	1:B:159:LEU:CD2	1.97	0.94
1:A:150:LEU:H	1:A:150:LEU:HD23	1.32	0.93
1:A:293:ALA:HB2	1:A:309:ILE:CD1	2.01	0.91
1:B:234:PHE:CD1	1:B:352:ALA:HB2	2.06	0.90
1:A:300:ASN:HB3	1:A:301:PRO:HD3	0.92	0.90
1:A:204:VAL:CG2	1:A:398:LEU:CD2	2.50	0.89
1:A:391:LYS:HG2	5:A:720:HOH:O	1.73	0.89
1:B:428:PHE:HB3	1:B:468:TYR:CD1	2.08	0.88
1:A:323:LYS:HA	1:A:341:ARG:HH21	1.37	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:GLU:HB2	1:B:201:TYR:HH	1.36	0.87
1:A:202:GLU:OE1	1:A:206:ILE:CG2	2.23	0.86
1:A:207:SER:OG	1:A:211:GLU:HB3	1.75	0.86
1:A:284:VAL:CG2	5:A:715:HOH:O	2.24	0.85
1:B:350:LYS:HD2	1:B:364:ARG:NH1	1.90	0.85
1:A:212:PHE:HB2	1:A:306:VAL:HG22	1.58	0.85
1:A:202:GLU:O	1:A:202:GLU:HG3	1.78	0.84
1:A:358:PHE:CE2	5:A:714:HOH:O	2.31	0.83
1:A:283:SER:HB2	1:A:299:ARG:NH1	1.94	0.83
1:A:327:PRO:HD2	1:A:468:TYR:CZ	2.13	0.83
1:A:271:LYS:O	1:A:275:LYS:HD3	1.79	0.82
1:B:465:LEU:HD22	1:B:476:PHE:CE2	2.15	0.81
1:A:182:MET:HE1	1:A:270:ILE:CG1	2.11	0.81
1:A:274:VAL:HA	1:A:277:ILE:CG2	2.10	0.81
1:A:210:ASN:OD1	1:A:304:ILE:HA	1.80	0.80
1:A:202:GLU:OE1	1:A:206:ILE:HG22	1.82	0.80
1:A:182:MET:HE1	1:A:270:ILE:HG13	1.63	0.80
1:B:465:LEU:HD23	1:B:476:PHE:CZ	2.17	0.80
1:B:350:LYS:CD	1:B:364:ARG:NH1	2.46	0.79
1:A:284:VAL:HG22	5:A:715:HOH:O	1.84	0.78
1:B:159:LEU:HD12	1:B:159:LEU:O	1.82	0.78
1:A:162:LYS:CE	1:B:201:TYR:O	2.32	0.78
1:B:465:LEU:CD2	1:B:476:PHE:CZ	2.67	0.78
1:A:212:PHE:O	1:A:306:VAL:HA	1.83	0.77
1:B:465:LEU:HD23	1:B:476:PHE:CE1	2.20	0.76
1:A:204:VAL:HG23	1:A:398:LEU:HB3	1.66	0.76
1:A:283:SER:HB2	1:A:299:ARG:HH12	1.51	0.75
1:B:328:ILE:HD12	1:B:332:LEU:HB2	1.70	0.74
1:B:160:LYS:HB3	1:B:164:ILE:HG12	1.70	0.74
1:B:156:LYS:CA	1:B:159:LEU:HD23	2.18	0.73
1:A:402:LYS:O	1:A:406:GLU:HG3	1.87	0.73
1:B:164:ILE:HD12	1:B:205:LYS:HD2	1.71	0.73
1:B:157:LEU:O	1:B:160:LYS:HD3	1.87	0.73
1:A:177:ARG:HD2	1:A:277:ILE:HD11	1.72	0.72
1:A:150:LEU:H	1:A:150:LEU:CD2	2.03	0.71
1:A:320:ILE:HD12	1:A:320:ILE:H	1.53	0.71
1:A:182:MET:CE	1:A:270:ILE:HG12	2.20	0.71
1:B:506:LYS:HB3	5:B:709:HOH:O	1.89	0.71
1:B:203:HIS:CE1	1:B:375:LEU:CD2	2.73	0.70
1:A:218:LEU:HD11	5:A:703:HOH:O	1.91	0.70
1:A:212:PHE:CB	1:A:306:VAL:HG22	2.20	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:HIS:HE1	1:B:375:LEU:CD2	2.04	0.70
1:A:161:ARG:O	1:A:164:ILE:HG13	1.92	0.70
1:B:203:HIS:CE1	1:B:375:LEU:HD23	2.27	0.69
1:A:162:LYS:HE2	1:B:201:TYR:O	1.92	0.69
1:A:274:VAL:HA	1:A:277:ILE:HG21	1.73	0.69
1:B:259:ALA:H	1:B:361:GLU:HA	1.57	0.69
1:A:244:ARG:HG3	1:A:244:ARG:O	1.93	0.68
1:A:295:THR:HA	1:A:307:ASP:HB3	1.75	0.68
1:A:188:GLU:HG3	1:A:247:PRO:HB2	1.75	0.67
1:A:298:ILE:HB	1:A:304:ILE:H	1.59	0.67
1:A:298:ILE:N	1:A:304:ILE:O	2.21	0.67
1:B:424:LYS:NZ	5:B:701:HOH:O	2.26	0.67
1:B:285:GLU:HB2	1:B:297:LEU:HG	1.76	0.66
1:A:149:LYS:O	1:A:153:VAL:HG13	1.95	0.66
1:B:450:ASP:OD2	1:B:497:ARG:NH2	2.21	0.66
1:A:274:VAL:C	1:A:277:ILE:HG22	2.16	0.66
1:A:358:PHE:CD2	5:A:714:HOH:O	2.49	0.66
1:A:274:VAL:HA	1:A:277:ILE:HG22	1.78	0.65
1:A:290:GLY:HA2	5:A:707:HOH:O	1.96	0.65
1:B:328:ILE:HD12	1:B:332:LEU:CB	2.25	0.65
1:B:402:LYS:HE3	1:B:423:VAL:HB	1.77	0.65
1:A:161:ARG:HB2	1:A:203:HIS:NE2	2.12	0.65
1:A:492:LYS:O	1:A:496:GLU:HB2	1.97	0.65
1:B:350:LYS:HD2	1:B:364:ARG:CZ	2.27	0.65
1:B:451:LYS:HD2	5:B:713:HOH:O	1.97	0.65
1:A:162:LYS:HE3	1:B:201:TYR:O	1.96	0.64
1:B:156:LYS:CA	1:B:159:LEU:CD2	2.74	0.64
1:A:280:ILE:HG22	1:A:281:ASP:H	1.62	0.64
1:B:158:ARG:O	1:B:158:ARG:HG2	1.97	0.64
1:B:252:LEU:HD13	1:B:254:GLY:O	1.97	0.64
1:A:265:LYS:HE2	1:A:269:ILE:HD11	1.80	0.64
1:A:293:ALA:CB	1:A:309:ILE:HD13	2.19	0.63
1:A:160:LYS:O	1:A:160:LYS:HD3	1.97	0.63
1:A:403:TYR:HD2	1:A:504:PHE:CE1	2.17	0.63
1:A:207:SER:OG	1:A:211:GLU:CB	2.45	0.63
1:A:287:GLU:O	1:A:287:GLU:HG3	1.97	0.63
1:A:206:ILE:HD11	3:A:602:A:O3'	1.99	0.63
1:B:465:LEU:CD2	1:B:476:PHE:CE2	2.80	0.63
1:A:403:TYR:CD2	1:A:504:PHE:CE1	2.87	0.62
1:A:199:SER:HB3	1:A:201:TYR:CD2	2.34	0.62
1:A:274:VAL:HG12	1:A:274:VAL:O	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:274:VAL:CA	1:A:277:ILE:HG22	2.30	0.61
1:A:381:GLU:O	1:A:384:CYS:HB2	1.99	0.61
1:A:190:LYS:HG2	5:A:711:HOH:O	2.00	0.61
1:B:350:LYS:HE3	4:B:603:5GP:H8	1.83	0.61
1:A:200:TYR:C	1:A:202:GLU:H	2.02	0.61
1:A:328:ILE:HG13	1:A:328:ILE:O	2.01	0.60
1:B:356:ASN:OD1	1:B:359:GLN:HG2	2.01	0.60
1:B:402:LYS:CE	1:B:423:VAL:HB	2.31	0.60
1:A:277:ILE:HG23	1:A:277:ILE:O	2.00	0.60
1:A:199:SER:HB3	1:A:201:TYR:CE2	2.37	0.60
1:A:208:ALA:N	1:A:209:PRO:CD	2.64	0.60
1:B:203:HIS:HE1	1:B:375:LEU:HD22	1.65	0.60
1:A:303:GLU:O	1:A:303:GLU:HG3	2.02	0.60
1:B:481:ILE:O	1:B:486:LYS:NZ	2.34	0.60
1:A:162:LYS:NZ	1:B:205:LYS:HE2	2.17	0.59
1:A:414:GLU:OE1	1:A:414:GLU:N	2.25	0.59
1:A:205:LYS:HG3	1:A:424:LYS:HZ1	1.67	0.59
1:B:156:LYS:HA	1:B:159:LEU:HD21	1.84	0.59
1:B:350:LYS:HE3	4:B:603:5GP:C8	2.34	0.58
1:B:197:THR:OG1	1:B:197:THR:O	2.20	0.58
1:B:424:LYS:HD3	5:B:705:HOH:O	2.04	0.58
1:A:169:GLU:CB	1:B:201:TYR:OH	2.40	0.58
1:B:155:ASP:O	1:B:159:LEU:HD23	2.04	0.58
1:A:182:MET:CE	1:A:270:ILE:CG1	2.77	0.57
1:A:327:PRO:HD2	1:A:468:TYR:CE1	2.38	0.57
1:A:424:LYS:HG2	1:A:428:PHE:CE2	2.39	0.57
1:A:415:LEU:HD22	1:A:489:LEU:HD13	1.85	0.57
1:B:149:LYS:O	1:B:153:VAL:HG23	2.03	0.57
1:B:424:LYS:CD	5:B:705:HOH:O	2.53	0.57
1:A:150:LEU:O	1:A:153:VAL:HG22	2.04	0.57
1:B:154:LEU:HD22	1:B:399:LYS:HE3	1.84	0.57
1:A:220:VAL:HG12	1:A:248:LEU:HD21	1.87	0.57
1:A:300:ASN:CB	1:A:301:PRO:CD	2.58	0.57
1:B:369:HIS:CD2	5:B:707:HOH:O	2.57	0.57
1:A:298:ILE:CG2	1:A:301:PRO:HD2	2.35	0.56
1:B:403:TYR:HE1	1:B:407:GLN:HE21	1.52	0.56
1:A:178:LEU:HD21	1:A:182:MET:CE	2.36	0.56
1:B:359:GLN:HB3	1:B:362:THR:HG23	1.86	0.56
1:A:205:LYS:HG3	1:A:424:LYS:NZ	2.21	0.56
1:A:407:GLN:HB3	1:A:503:ILE:CG1	2.35	0.56
1:B:204:VAL:O	1:B:204:VAL:HG13	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:223:ILE:HG13	1:B:238:LYS:O	2.04	0.56
1:A:241:ARG:O	1:A:241:ARG:HG2	2.05	0.56
1:A:292:PRO:O	1:A:309:ILE:HG23	2.06	0.56
1:A:204:VAL:HG23	1:A:398:LEU:CB	2.35	0.55
1:A:323:LYS:O	1:A:337:ARG:NH2	2.39	0.55
1:A:331:TRP:HB2	1:A:435:PRO:HB3	1.86	0.55
1:B:154:LEU:HD11	1:B:396:GLU:HG3	1.88	0.55
1:B:179:LEU:O	1:B:183:GLN:HG3	2.06	0.55
1:B:465:LEU:HB3	1:B:475:LEU:HB2	1.88	0.55
1:B:156:LYS:HD2	1:B:159:LEU:HD21	1.88	0.55
1:A:216:PHE:HD2	1:A:266:PHE:HZ	1.55	0.54
1:A:481:ILE:HB	1:A:486:LYS:HE3	1.90	0.54
1:B:223:ILE:CD1	1:B:238:LYS:O	2.56	0.54
1:A:277:ILE:O	1:A:277:ILE:HG13	2.08	0.54
1:B:169:GLU:O	1:B:173:LYS:HG2	2.09	0.53
1:B:223:ILE:HG12	1:B:237:VAL:HG13	1.91	0.53
1:B:383:THR:O	1:B:386:GLU:HB2	2.08	0.53
1:B:172:ASN:O	1:B:176:GLU:HB2	2.09	0.53
1:A:493:ILE:O	1:A:496:GLU:N	2.39	0.53
1:A:499:ASN:N	1:A:499:ASN:OD1	2.41	0.53
1:B:224:GLU:HG2	1:B:240:LYS:HE3	1.91	0.52
1:B:398:LEU:HB2	1:B:427:ILE:HD13	1.90	0.52
1:A:220:VAL:HG21	1:A:223:ILE:HD12	1.91	0.52
1:A:502:PRO:HA	1:A:505:ASP:OD2	2.08	0.52
1:A:387:SER:HB3	1:B:159:LEU:HD11	1.91	0.52
1:B:220:VAL:HG12	1:B:248:LEU:CD2	2.40	0.52
1:A:450:ASP:OD1	1:A:497:ARG:NE	2.42	0.52
1:A:251:PHE:CD2	1:A:262:MET:HG3	2.44	0.52
1:A:285:GLU:HG3	1:A:297:LEU:CD2	2.40	0.52
1:A:348:VAL:HG21	1:A:364:ARG:NH2	2.25	0.52
1:A:207:SER:CB	1:A:211:GLU:HB3	2.41	0.51
1:B:227:GLU:HA	1:B:235:TYR:CD1	2.45	0.51
1:A:178:LEU:HD21	1:A:182:MET:HE2	1.93	0.51
1:A:299:ARG:O	1:A:299:ARG:HG3	2.11	0.50
1:A:185:ARG:HG2	1:A:186:GLU:N	2.26	0.50
1:A:200:TYR:CD1	1:A:200:TYR:C	2.85	0.50
1:A:295:THR:HA	1:A:307:ASP:CB	2.40	0.50
1:A:181:ARG:HH11	1:A:181:ARG:HG2	1.77	0.50
1:B:224:GLU:CG	1:B:240:LYS:HE3	2.42	0.50
1:A:292:PRO:HB3	1:A:349:PRO:O	2.12	0.50
1:B:319:PRO:HG2	1:B:365:LEU:HD11	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:199:SER:O	1:B:203:HIS:N	2.40	0.49
1:B:418:PHE:HA	1:B:422:HIS:CE1	2.46	0.49
1:A:431:TRP:HD1	1:A:440:TRP:CZ2	2.30	0.49
1:A:212:PHE:HB2	1:A:305:SER:O	2.11	0.49
1:A:402:LYS:HD3	5:A:717:HOH:O	2.12	0.49
1:A:176:GLU:HG2	1:B:173:LYS:HD2	1.94	0.49
1:A:178:LEU:CD2	1:A:182:MET:CE	2.91	0.49
1:A:217:LYS:NZ	5:A:701:HOH:O	2.45	0.49
1:A:383:THR:O	1:A:389:GLY:HA3	2.12	0.49
1:A:212:PHE:N	1:A:305:SER:O	2.46	0.49
1:B:428:PHE:CB	1:B:468:TYR:CD1	2.87	0.49
1:A:201:TYR:CD1	1:A:201:TYR:C	2.87	0.48
1:A:402:LYS:HD3	1:A:420:SER:HB2	1.95	0.48
1:B:154:LEU:CD1	1:B:396:GLU:HG3	2.43	0.48
1:B:486:LYS:O	1:B:490:SER:OG	2.29	0.48
1:A:186:GLU:HG2	1:A:186:GLU:O	2.13	0.48
1:A:273:GLU:O	1:A:277:ILE:HB	2.13	0.48
1:B:328:ILE:HD12	1:B:332:LEU:HD12	1.96	0.48
1:A:401:MET:HB2	1:A:401:MET:HE2	1.62	0.48
1:B:275:LYS:HG2	1:B:275:LYS:O	2.13	0.48
1:B:234:PHE:CE1	1:B:352:ALA:HB2	2.48	0.47
1:B:206:ILE:CD1	1:B:395:LYS:HD3	2.44	0.47
1:A:295:THR:O	1:A:296:LEU:HD23	2.14	0.47
1:A:297:LEU:HB2	1:A:304:ILE:O	2.14	0.47
1:A:270:ILE:HD13	1:A:294:VAL:HG21	1.97	0.47
1:B:156:LYS:O	1:B:159:LEU:HG	2.15	0.47
1:A:407:GLN:HB3	1:A:503:ILE:HG12	1.96	0.47
1:A:204:VAL:HG23	1:A:398:LEU:CG	2.44	0.46
1:A:431:TRP:CB	5:A:704:HOH:O	2.28	0.46
1:B:174:VAL:HG21	5:B:715:HOH:O	2.14	0.46
1:B:411:GLU:HG3	1:B:412:PHE:CD2	2.50	0.46
1:A:178:LEU:CD2	1:A:182:MET:HE3	2.46	0.46
1:A:266:PHE:CE2	1:A:310:LEU:HB2	2.51	0.46
1:A:274:VAL:CA	1:A:277:ILE:CG2	2.86	0.46
1:A:280:ILE:HG22	1:A:281:ASP:N	2.31	0.46
1:A:298:ILE:HB	1:A:304:ILE:N	2.29	0.46
1:A:479:GLU:N	1:A:479:GLU:OE1	2.49	0.46
1:B:315:LYS:NZ	5:B:703:HOH:O	2.43	0.46
1:A:177:ARG:O	1:A:181:ARG:HG3	2.15	0.46
1:A:212:PHE:CE2	1:A:304:ILE:HG21	2.51	0.46
1:A:181:ARG:NH1	1:A:181:ARG:HG2	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:467:HIS:HD2	1:B:469:PHE:H	1.64	0.46
1:B:179:LEU:HD13	1:B:216:PHE:CZ	2.51	0.45
1:A:179:LEU:O	1:A:183:GLN:HG2	2.16	0.45
1:A:483:ARG:O	1:A:487:GLU:HG3	2.16	0.45
1:B:276:GLU:O	1:B:276:GLU:HG3	2.17	0.45
1:B:285:GLU:OE2	1:B:288:LYS:HE2	2.16	0.45
1:A:147:PRO:HB2	1:A:148:ASP:H	1.57	0.45
1:A:424:LYS:HB2	1:A:424:LYS:HE3	1.61	0.45
1:B:156:LYS:O	1:B:159:LEU:CD2	2.64	0.45
1:A:205:LYS:HE2	1:A:424:LYS:HZ3	1.82	0.45
1:A:298:ILE:O	1:A:303:GLU:HA	2.16	0.45
1:A:391:LYS:HE3	1:A:391:LYS:HB3	1.85	0.45
1:B:350:LYS:HD3	1:B:364:ARG:NH1	2.27	0.45
1:A:201:TYR:O	1:A:201:TYR:HD1	2.00	0.45
1:B:222:ARG:HG3	1:B:240:LYS:HG3	1.99	0.45
1:B:298:ILE:HB	1:B:304:ILE:HG13	1.99	0.45
1:A:178:LEU:HD21	1:A:182:MET:HE3	1.99	0.44
1:A:250:HIS:O	1:A:250:HIS:ND1	2.50	0.44
1:A:217:LYS:HA	1:A:311:ALA:O	2.18	0.44
1:B:298:ILE:HD13	5:B:715:HOH:O	2.17	0.44
1:A:447:SER:O	1:A:451:LYS:HB2	2.18	0.44
1:A:324:GLU:HA	1:A:324:GLU:OE2	2.16	0.44
1:A:407:GLN:HG3	1:A:503:ILE:HG13	1.99	0.44
1:B:189:PHE:O	1:B:192:VAL:HB	2.18	0.44
1:B:492:LYS:HD3	1:B:492:LYS:HA	1.82	0.44
1:A:412:PHE:CD2	1:A:488:PHE:HZ	2.35	0.44
1:A:162:LYS:HZ2	1:B:205:LYS:HE2	1.81	0.44
1:B:296:LEU:HB2	1:B:306:VAL:HB	1.99	0.44
1:B:322:THR:HG21	1:B:367:PHE:CZ	2.53	0.44
1:A:177:ARG:HD2	1:A:277:ILE:CD1	2.45	0.44
1:B:177:ARG:HH11	1:B:181:ARG:HH21	1.66	0.44
1:B:411:GLU:HG3	1:B:412:PHE:CE2	2.51	0.44
1:B:430:MET:HG3	1:B:452:LEU:HD13	2.00	0.44
1:A:326:LEU:HD11	1:A:374:ILE:HD11	2.00	0.44
1:A:295:THR:HG23	1:A:307:ASP:HB3	2.00	0.43
1:A:407:GLN:HB3	1:A:503:ILE:HG13	2.00	0.43
1:B:195:LEU:C	1:B:197:THR:H	2.22	0.43
1:B:231:THR:HG21	1:B:234:PHE:HB2	2.00	0.43
1:B:372:LYS:HG3	1:B:373:TYR:N	2.33	0.43
1:B:393:CYS:O	1:B:396:GLU:HB2	2.19	0.43
1:B:251:PHE:CD2	1:B:262:MET:HG3	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:LEU:C	1:B:197:THR:N	2.72	0.43
1:B:402:LYS:HE2	1:B:420:SER:HA	2.00	0.43
1:A:460:LEU:HD21	1:A:489:LEU:HD23	2.00	0.43
1:A:432:THR:OG1	1:A:468:TYR:OH	2.32	0.43
1:B:173:LYS:HA	1:B:173:LYS:HD3	1.83	0.43
1:A:294:VAL:O	1:A:294:VAL:HG23	2.19	0.42
1:A:460:LEU:HD22	1:A:486:LYS:HA	2.01	0.42
1:A:150:LEU:N	1:A:150:LEU:CD2	2.72	0.42
1:A:150:LEU:HD23	1:A:150:LEU:N	2.11	0.42
1:A:210:ASN:OD1	1:A:303:GLU:O	2.37	0.42
1:B:405:LEU:HG	1:B:405:LEU:O	2.17	0.42
1:A:403:TYR:CD2	1:A:504:PHE:HE1	2.35	0.42
1:A:320:ILE:CD1	1:A:320:ILE:H	2.19	0.42
1:B:210:ASN:HB2	1:B:211:GLU:OE1	2.19	0.42
1:B:230:GLU:O	1:B:230:GLU:HG2	2.19	0.42
1:B:223:ILE:CG1	1:B:238:LYS:O	2.68	0.42
1:A:225:LEU:HD11	1:A:345:PHE:CE2	2.54	0.41
1:A:167:ALA:HB1	1:A:298:ILE:HG13	2.02	0.41
1:A:185:ARG:HG2	1:A:186:GLU:H	1.85	0.41
1:B:223:ILE:HG12	1:B:237:VAL:CG1	2.49	0.41
1:A:283:SER:O	1:A:296:LEU:HB3	2.19	0.41
1:A:492:LYS:O	1:A:496:GLU:CB	2.66	0.41
1:B:164:ILE:CD1	1:B:205:LYS:HD2	2.45	0.41
1:A:203:HIS:HB3	1:A:399:LYS:HE2	2.02	0.41
1:A:284:VAL:HA	1:A:296:LEU:HD22	2.03	0.41
1:A:299:ARG:CG	1:A:299:ARG:O	2.69	0.41
1:A:328:ILE:HD12	1:A:332:LEU:HB2	2.01	0.41
1:A:458:GLU:O	1:A:462:THR:OG1	2.34	0.41
1:B:281:ASP:O	1:B:298:ILE:HG23	2.20	0.41
1:B:300:ASN:OD1	1:B:301:PRO:HD3	2.21	0.41
1:A:150:LEU:HD12	1:A:400:LEU:HD13	2.03	0.41
1:A:159:LEU:HD12	1:A:159:LEU:O	2.21	0.41
1:A:199:SER:OG	1:A:200:TYR:N	2.53	0.41
1:A:378:HIS:HD2	1:A:394:ARG:HD2	1.86	0.41
1:A:503:ILE:HD12	1:A:503:ILE:HA	1.89	0.41
1:B:185:ARG:O	1:B:185:ARG:HD2	2.21	0.41
1:B:204:VAL:HG21	1:B:212:PHE:HA	2.03	0.41
1:A:159:LEU:CG	1:A:159:LEU:O	2.69	0.41
1:A:431:TRP:CA	5:A:704:HOH:O	2.67	0.41
1:A:495:TYR:O	1:A:495:TYR:CG	2.74	0.41
1:B:424:LYS:O	1:B:427:ILE:HG22	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:492:LYS:O	1:A:496:GLU:HG2	2.21	0.40
1:A:188:GLU:N	1:A:188:GLU:OE1	2.47	0.40
1:A:233:ALA:O	1:A:364:ARG:HA	2.21	0.40
1:A:301:PRO:HB2	1:A:302:GLU:H	1.57	0.40
1:A:384:CYS:C	1:A:386:GLU:H	2.25	0.40
1:B:156:LYS:CA	1:B:159:LEU:HD21	2.47	0.40
1:B:220:VAL:HG22	1:B:313:GLU:O	2.21	0.40
1:A:167:ALA:O	1:A:170:THR:HB	2.22	0.40
1:A:210:ASN:HA	1:A:304:ILE:HG23	2.02	0.40
1:B:221:PRO:O	1:B:222:ARG:C	2.60	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	359/507 (71%)	308 (86%)	46 (13%)	5 (1%)	11	40
1	B	359/507 (71%)	325 (90%)	32 (9%)	2 (1%)	25	60
All	All	718/1014 (71%)	633 (88%)	78 (11%)	7 (1%)	15	49

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	301	PRO
1	A	207	SER
1	A	209	PRO
1	A	300	ASN
1	A	302	GLU
1	B	301	PRO
1	B	327	PRO

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	334/447 (75%)	311 (93%)	23 (7%)	15	45
1	B	334/447 (75%)	321 (96%)	13 (4%)	32	66
All	All	668/894 (75%)	632 (95%)	36 (5%)	22	55

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	148	ASP
1	A	149	LYS
1	A	150	LEU
1	A	151	LYS
1	A	158	ARG
1	A	161	ARG
1	A	162	LYS
1	A	163	ASP
1	A	165	SER
1	A	185	ARG
1	A	207	SER
1	A	246	ASN
1	A	267	ARG
1	A	287	GLU
1	A	354	ASP
1	A	359	GLN
1	A	361	GLU
1	A	372	LYS
1	A	409	LYS
1	A	441	ASP
1	A	473	PHE
1	A	483	ARG
1	A	499	ASN
1	B	156	LYS
1	B	177	ARG
1	B	185	ARG
1	B	196	ASN

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Mol	Chain	Res	Type
1	B	201	TYR
1	B	207	SER
1	B	222	ARG
1	B	241	ARG
1	B	288	LYS
1	B	350	LYS
1	B	354	ASP
1	B	438	SER
1	B	465	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	356	ASN
1	B	196	ASN
1	B	203	HIS
1	B	444	ASN
1	B	467	HIS

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 6 ligands modelled in this entry, 2 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	$\# Z > 2$
3	A	A	602	4	18,24,25	7.30	12 (66%)	18,35,38	1.90	4 (22%)
3	A	B	602	4	18,24,25	7.12	13 (72%)	18,35,38	1.75	2 (11%)
4	5GP	A	603	3	22,26,26	4.95	11 (50%)	27,40,40	1.92	6 (22%)
4	5GP	B	603	3	22,26,26	4.72	10 (45%)	27,40,40	1.88	5 (18%)

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	A	A	602	4	-	0/3/25/26	0/3/3/3
3	A	B	602	4	-	1/3/25/26	0/3/3/3
4	5GP	A	603	3	-	5/6/26/26	0/3/3/3
4	5GP	B	603	3	-	5/6/26/26	0/3/3/3

All (46) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	602	A	C2'-C1'	-19.16	1.24	1.53
3	A	602	A	C2'-C1'	-18.79	1.25	1.53
3	A	602	A	O4'-C1'	15.62	1.62	1.41
3	B	602	A	O4'-C1'	14.29	1.61	1.41
3	A	602	A	C4-N3	9.60	1.48	1.35
4	A	603	5GP	C4-N3	9.56	1.50	1.35
3	B	602	A	C4-N3	9.10	1.48	1.35
4	A	603	5GP	O4'-C1'	-8.99	1.28	1.41
4	A	603	5GP	C3'-C4'	-8.92	1.30	1.53
3	A	602	A	C3'-C4'	-8.89	1.30	1.53
4	B	603	5GP	C4-N3	8.68	1.49	1.35
4	B	603	5GP	C3'-C4'	-8.65	1.30	1.53
4	B	603	5GP	C6-C5	8.58	1.56	1.41
3	A	602	A	C2-N3	8.53	1.45	1.32
3	B	602	A	C2-N3	8.52	1.45	1.32
4	B	603	5GP	O4'-C1'	-8.47	1.29	1.41
3	B	602	A	C3'-C4'	-8.40	1.31	1.53
4	A	603	5GP	C6-C5	8.25	1.55	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	603	5GP	O4'-C4'	8.04	1.63	1.45
4	B	603	5GP	O4'-C4'	8.00	1.62	1.45
4	A	603	5GP	C6-N1	6.81	1.44	1.33
3	B	602	A	C2-N1	6.58	1.46	1.33
3	A	602	A	C2-N1	6.19	1.45	1.33
4	A	603	5GP	C2-N2	5.95	1.45	1.33
4	B	603	5GP	C6-N1	5.89	1.43	1.33
4	A	603	5GP	C2-N1	5.82	1.45	1.35
4	B	603	5GP	C2-N2	5.64	1.45	1.33
3	A	602	A	C2'-C3'	5.32	1.67	1.53
4	B	603	5GP	C2-N1	5.12	1.44	1.35
3	B	602	A	C2'-C3'	5.02	1.67	1.53
4	A	603	5GP	O3'-C3'	3.63	1.51	1.43
3	A	602	A	O4'-C4'	3.54	1.52	1.45
4	B	603	5GP	O3'-C3'	3.48	1.51	1.43
3	A	602	A	C6-N6	3.31	1.46	1.34
3	B	602	A	O4'-C4'	3.14	1.52	1.45
3	B	602	A	C6-N6	3.06	1.45	1.34
4	A	603	5GP	C2-N3	2.91	1.48	1.34
4	A	603	5GP	O2'-C2'	-2.85	1.36	1.43
3	A	602	A	C6-C5	2.79	1.53	1.43
3	B	602	A	C6-C5	2.47	1.52	1.43
4	B	603	5GP	C2-N3	2.39	1.45	1.34
3	A	602	A	C5'-C4'	2.34	1.58	1.51
3	B	602	A	O3'-C3'	2.25	1.48	1.43
3	A	602	A	O2'-C2'	2.24	1.48	1.43
3	B	602	A	C5'-C4'	2.20	1.58	1.51
3	B	602	A	O2'-C2'	2.14	1.48	1.43

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	602	A	N3-C2-N1	-5.99	119.32	128.68
3	A	602	A	N3-C2-N1	-5.53	120.03	128.68
4	A	603	5GP	N3-C2-N1	-5.25	120.22	127.22
4	B	603	5GP	N3-C2-N1	-5.20	120.28	127.22
4	A	603	5GP	C2-N3-C4	4.80	120.84	115.36
4	B	603	5GP	C2-N3-C4	4.07	120.01	115.36
4	B	603	5GP	C5-C6-N1	-3.66	118.43	123.43
4	A	603	5GP	C5-C6-N1	-3.35	118.85	123.43
3	A	602	A	O4'-C1'-C2'	-3.26	102.17	106.93
4	B	603	5GP	C6-N1-C2	3.06	120.80	115.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	A	C3'-C2'-C1'	2.83	105.25	100.98
3	B	602	A	O4'-C1'-C2'	-2.82	102.80	106.93
4	B	603	5GP	C4-C5-N7	-2.80	106.48	109.40
4	A	603	5GP	C4-C5-N7	-2.50	106.80	109.40
4	A	603	5GP	O4'-C1'-C2'	-2.39	103.43	106.93
4	A	603	5GP	C6-N1-C2	2.29	119.56	115.93
3	A	602	A	C4-C5-N7	-2.23	107.07	109.40

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	603	5GP	C5'-O5'-P-O3P
4	A	603	5GP	O4'-C4'-C5'-O5'
4	B	603	5GP	C5'-O5'-P-O2P
4	B	603	5GP	C5'-O5'-P-O3P
4	B	603	5GP	O4'-C4'-C5'-O5'
4	B	603	5GP	C3'-C4'-C5'-O5'
4	A	603	5GP	C3'-C4'-C5'-O5'
4	B	603	5GP	C5'-O5'-P-O1P
3	B	602	A	O4'-C4'-C5'-O5'
4	A	603	5GP	C5'-O5'-P-O1P
4	A	603	5GP	C5'-O5'-P-O2P

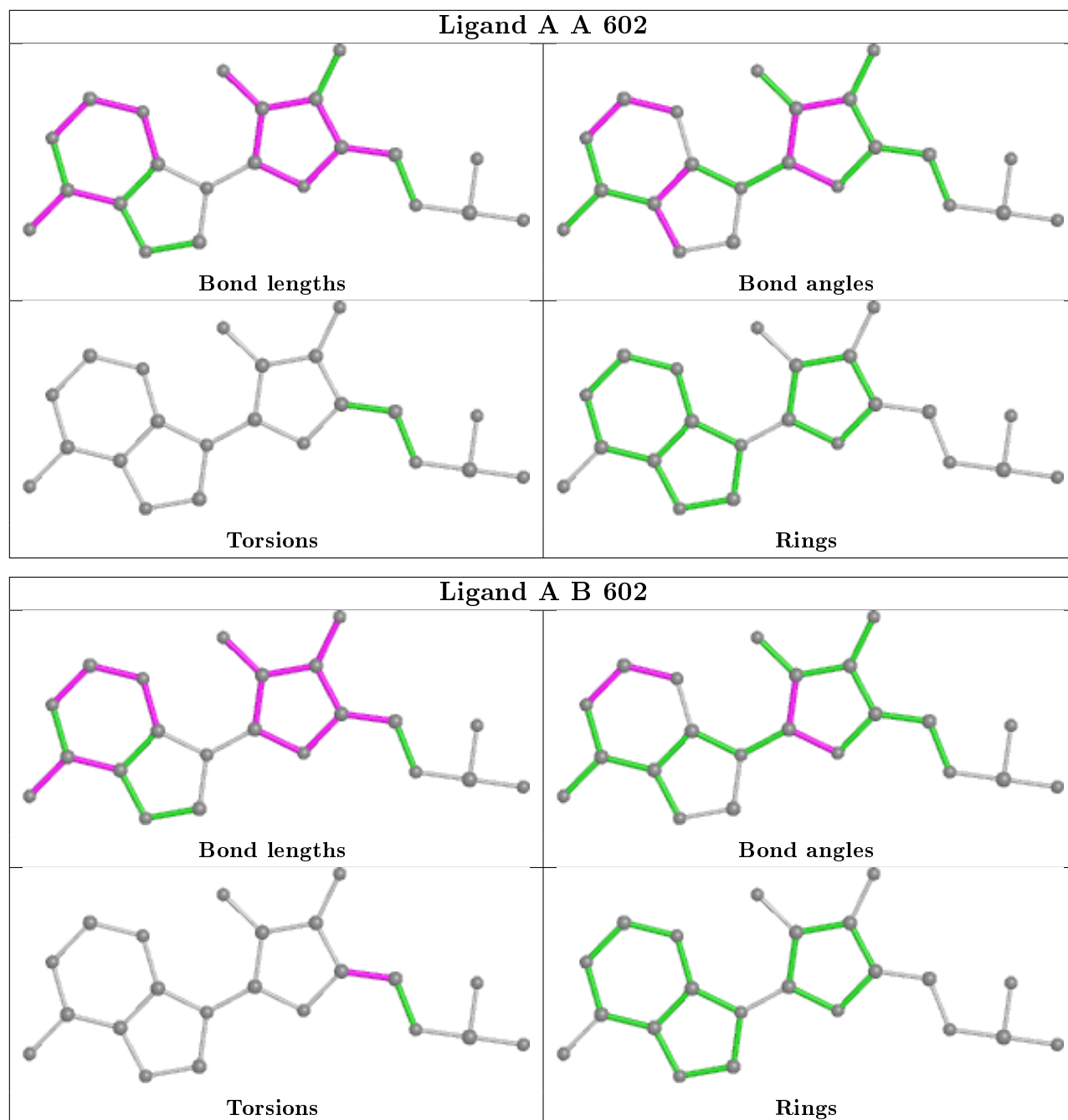
There are no ring outliers.

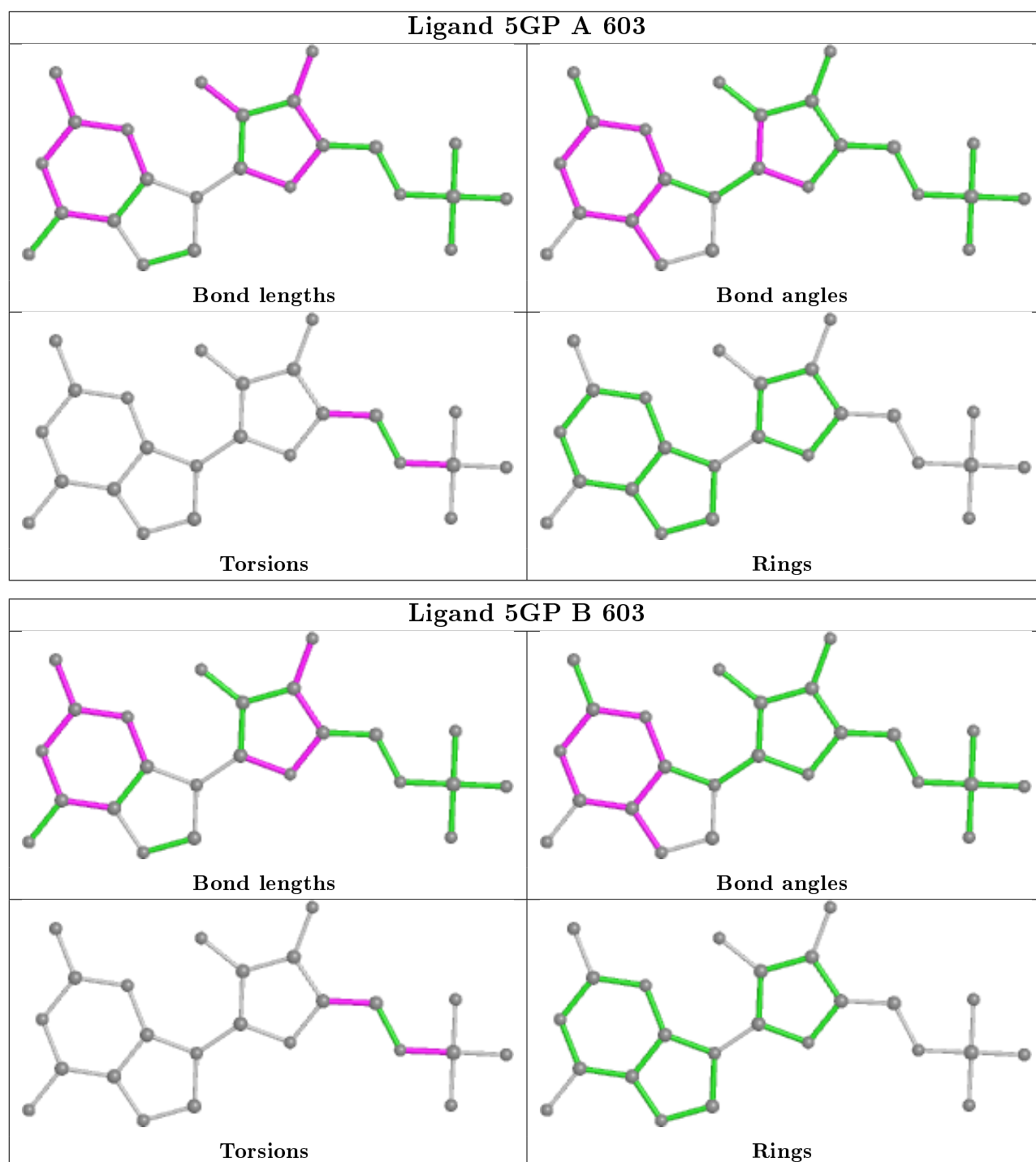
2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	A	1	0
4	B	603	5GP	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.





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	361/507 (71%)	-0.04	12 (3%)	46	20	27, 58, 130, 191	0
1	B	361/507 (71%)	-0.02	15 (4%)	36	14	26, 56, 140, 213	0
All	All	722/1014 (71%)	-0.03	27 (3%)	41	17	26, 57, 138, 213	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	162	LYS	8.2
1	B	355	GLY	4.3
1	B	161	ARG	4.3
1	A	302	GLU	4.3
1	A	303	GLU	3.8
1	B	207	SER	3.6
1	B	201	TYR	3.2
1	A	301	PRO	3.1
1	A	304	ILE	3.1
1	A	205	LYS	3.1
1	B	276	GLU	2.9
1	A	361	GLU	2.7
1	A	229	TYR	2.7
1	B	301	PRO	2.7
1	B	198	GLY	2.6
1	B	166	GLU	2.6
1	A	358	PHE	2.5
1	B	205	LYS	2.5
1	B	243	PRO	2.5
1	B	304	ILE	2.4
1	B	197	THR	2.4
1	A	356	ASN	2.2
1	A	436	GLN	2.2
1	B	204	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	164	ILE	2.1
1	A	163	ASP	2.1
1	B	356	ASN	2.1

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

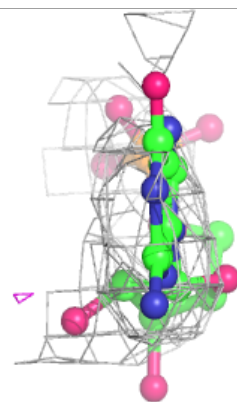
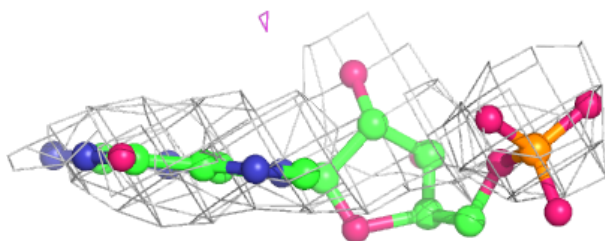
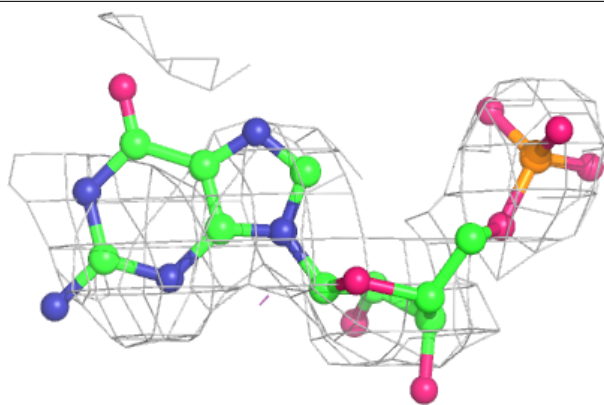
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	5GP	A	603	24/24	0.59	0.37	144,152,156,157	0
3	A	A	602	22/23	0.67	0.42	111,123,128,136	0
4	5GP	B	603	24/24	0.87	0.24	58,67,73,75	0
3	A	B	602	22/23	0.93	0.27	53,58,66,75	0
2	ZN	B	601	1/1	0.97	0.09	49,49,49,49	0
2	ZN	A	601	1/1	0.99	0.10	47,47,47,47	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.

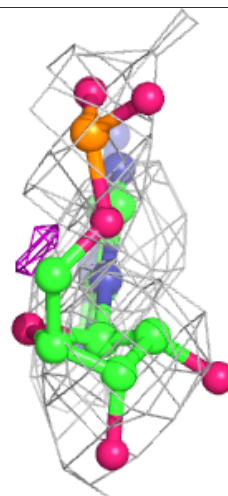
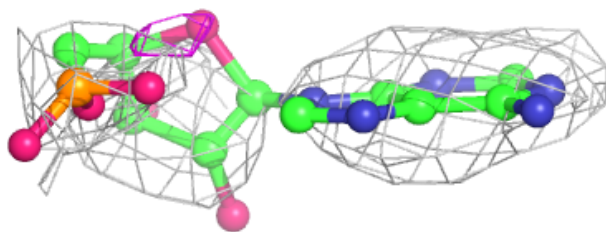
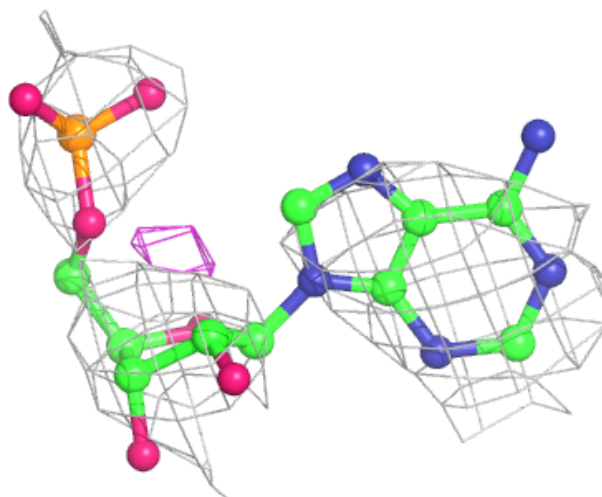
Electron density around 5GP A 603:

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



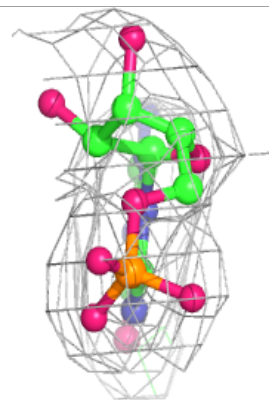
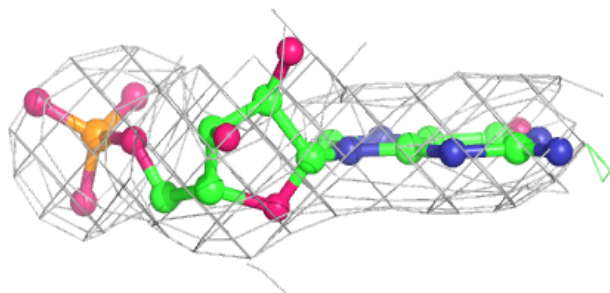
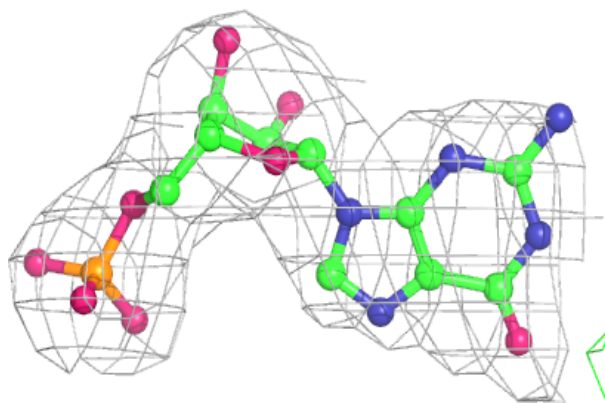
Electron density around A A 602:

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



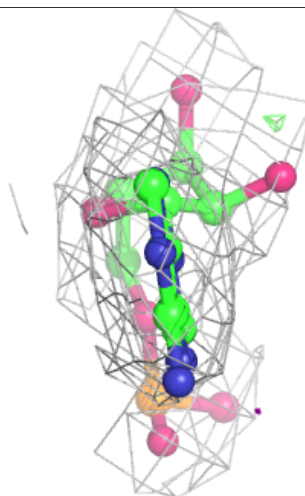
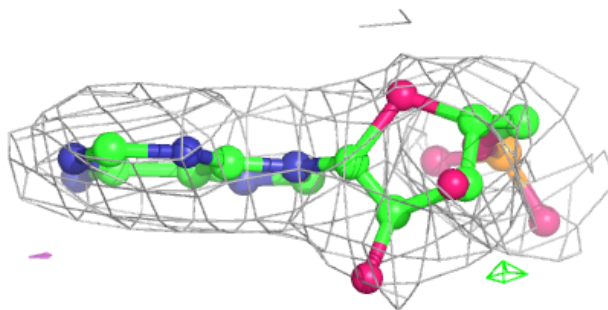
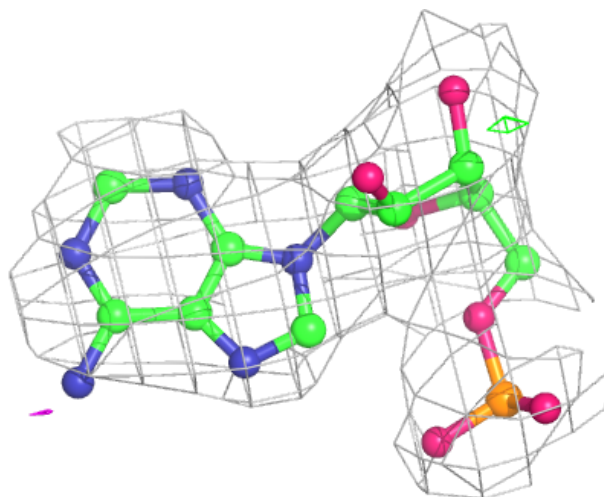
Electron density around 5GP B 603:

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



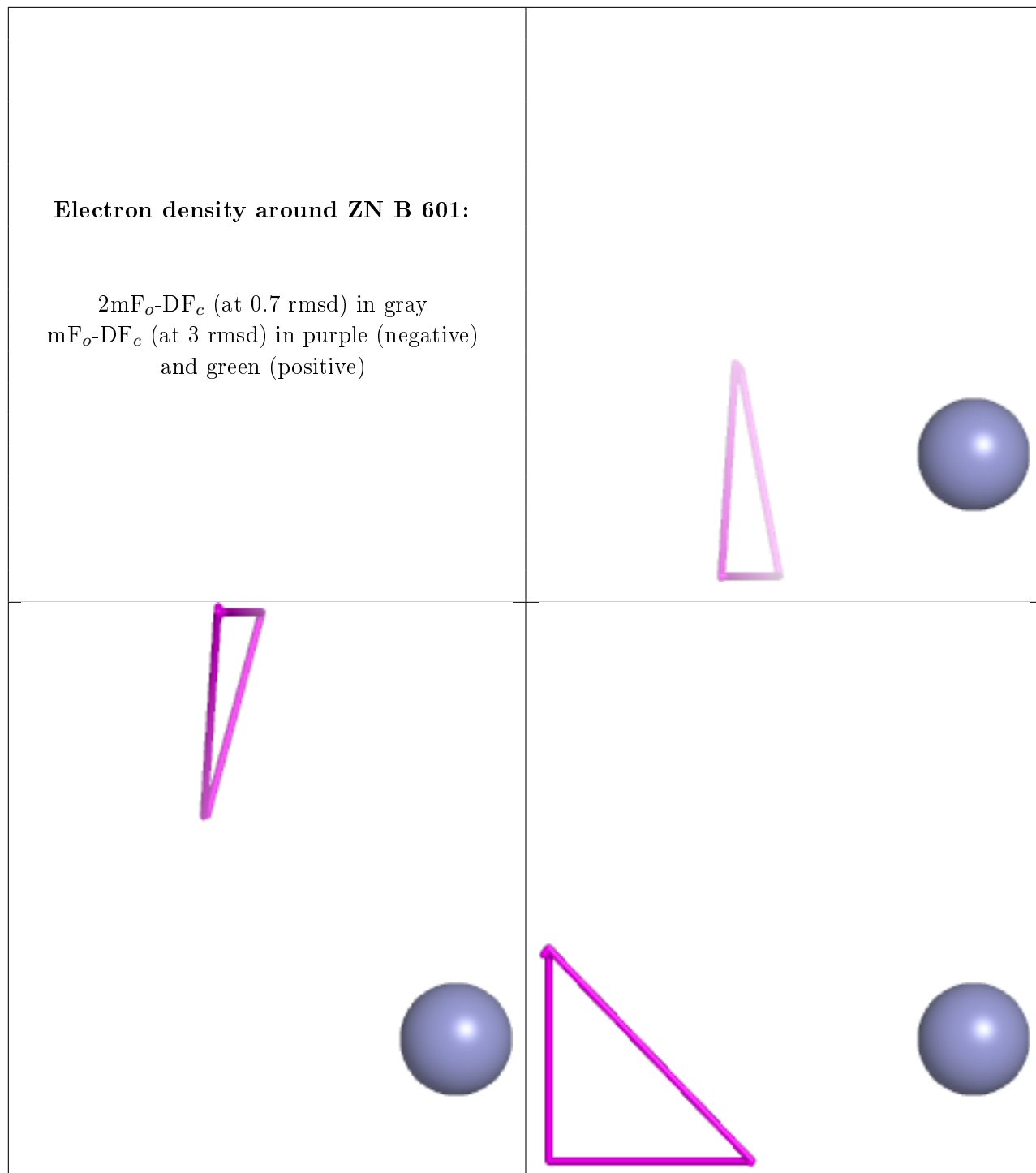
Electron density around A B 602:

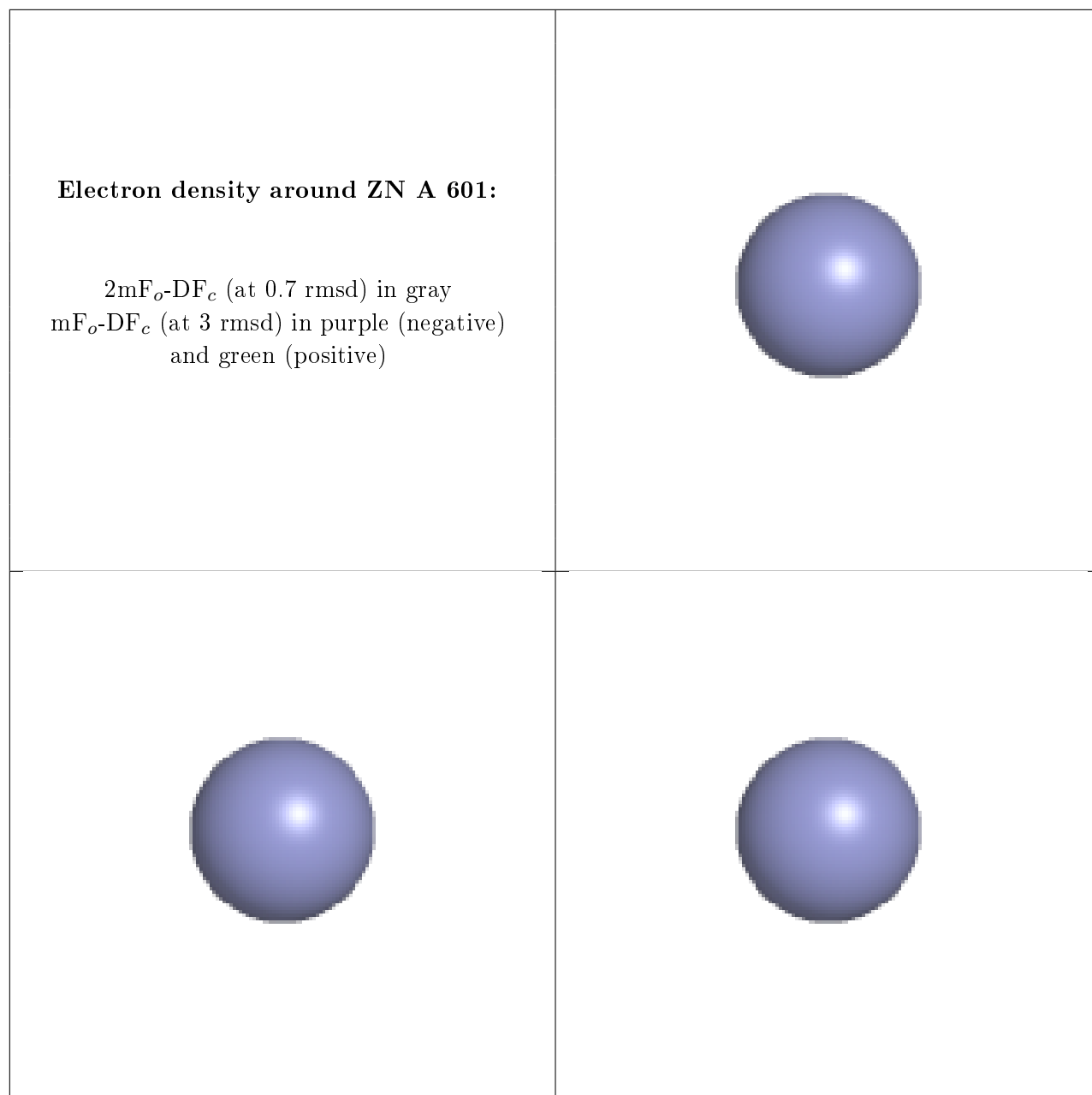
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ZN B 601:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.