



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 10:13 PM EDT

PDB ID : 6CZF / pdb_00006czf
Title : The structure of E. coli PurF in complex with ppGpp-Mg
Authors : Wang, B.; Grant, R.A.; Laub, M.T.
Deposited on : 2018-04-09
Resolution : 1.95 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

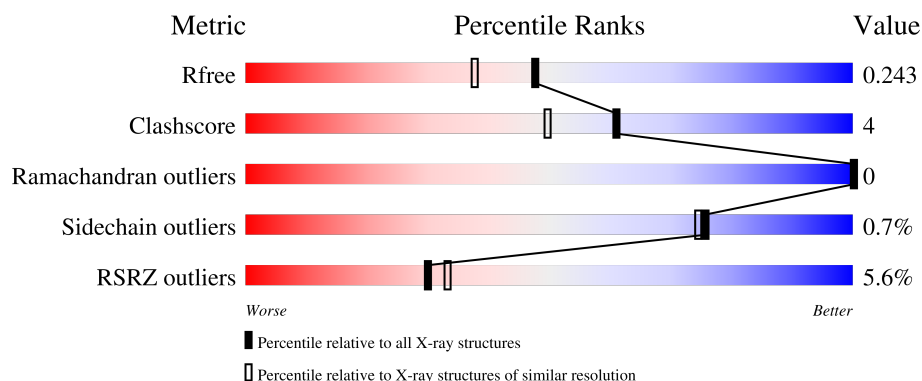
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	498	7% 86% 13%
1	B	498	3% 93% 6%
1	C	498	8% 87% 13%
1	D	498	4% 91% 8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 31549 atoms, of which 15444 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.

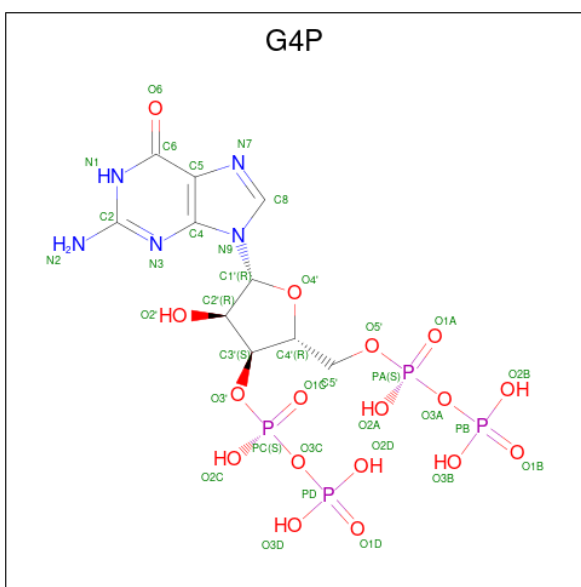
- Molecule 1 is a protein called Amidophosphoribosyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	498	Total	C	H	N	O	S	0	0	0
			7783	2453	3872	700	741	17			
1	B	495	Total	C	H	N	O	S	0	0	0
			7711	2434	3831	692	737	17			
1	C	498	Total	C	H	N	O	S	0	0	0
			7791	2455	3876	700	743	17			
1	D	494	Total	C	H	N	O	S	0	0	0
			7692	2428	3821	692	734	17			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	1
			2	2		
2	D	1	Total	Mg	0	1
			2	2		

- Molecule 3 is GUANOSINE-5',3'-TETRAPHOSPHATE (CCD ID: G4P) (formula: C₁₀H₁₇N₅O₁₇P₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total 94	C 20	H 22	N 10	O 34	P 8	0	1
3	D	1	Total 94	C 20	H 22	N 10	O 34	P 8	0	1

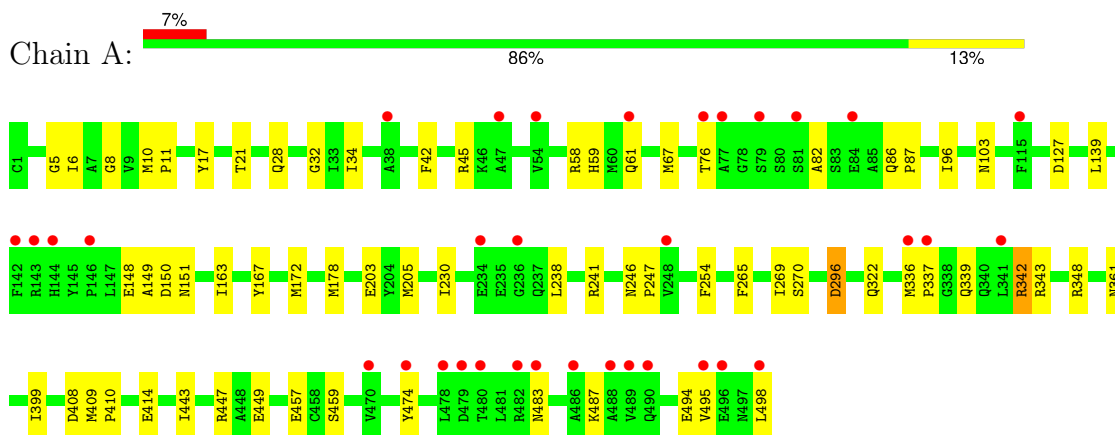
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	78	Total O 78 78	0	0
4	B	113	Total O 113 113	0	0
4	C	74	Total O 74 74	0	0
4	D	115	Total O 115 115	0	0

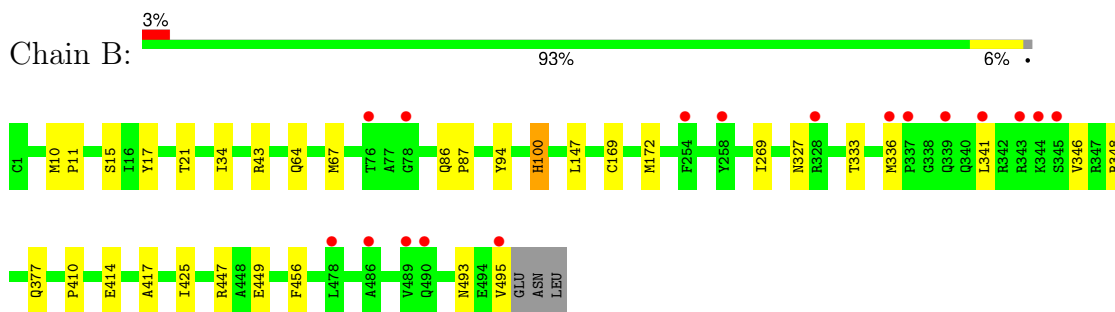
3 Residue-property plots [i](#)

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.

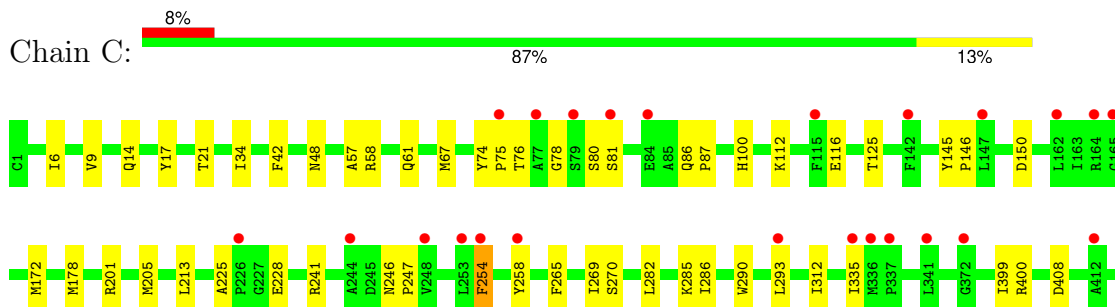
- Molecule 1: Amidophosphoribosyltransferase

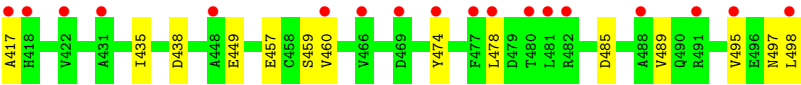


- Molecule 1: Amidophosphoribosyltransferase

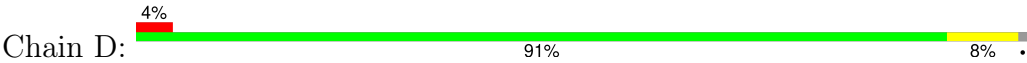


- Molecule 1: Amidophosphoribosyltransferase





● Molecule 1: Amidophosphoribosyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	107.29Å 115.34Å 156.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.64 – 1.95 48.64 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.5 (48.64-1.95) 90.4 (48.64-1.95)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.63 (at 1.95Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.216 , 0.241 0.218 , 0.243	Depositor DCC
R_{free} test set	1987 reflections (1.41%)	wwPDB-VP
Wilson B-factor (Å ²)	37.0	Xtriage
Anisotropy	0.252	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 24.8	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	0.96	EDS
Total number of atoms	31549	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 86.10 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.0387e-07. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: G4P, MG

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	0.28	0/3979	0.54	0/5386
1	B	0.27	0/3948	0.54	0/5346
1	C	0.27	0/3983	0.53	0/5391
1	D	0.27	0/3939	0.55	0/5333
All	All	0.27	0/15849	0.54	0/21456

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	3911	3872	3872	44	0
1	B	3880	3831	3831	24	0
1	C	3915	3876	3876	47	0
1	D	3871	3821	3821	26	0
2	A	2	0	0	0	0
2	D	2	0	0	0	0
3	A	72	22	22	1	0
3	D	72	22	22	0	0
4	A	78	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	113	0	0	1	0
4	C	74	0	0	1	0
4	D	115	0	0	0	0
All	All	16105	15444	15444	131	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 4.

All (131) 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:495:VAL:HG21	1:C:61:GLN:HB3	1.40	1.03
1:A:495:VAL:HG21	1:C:61:GLN:CB	2.11	0.81
1:A:61:GLN:HB3	1:C:495:VAL:HG21	1.63	0.79
1:B:341:LEU:H	1:B:341:LEU:HD23	1.51	0.75
1:B:336:MET:HA	1:B:336:MET:HE2	1.69	0.74
1:A:342:ARG:HD3	1:B:15:SER:OG	1.86	0.74
1:C:150:ASP:OD1	1:C:241:ARG:NH2	2.28	0.67
1:A:150:ASP:OD1	1:A:241:ARG:NH2	2.31	0.63
1:A:296:ASP:HB2	1:A:361:ASN:O	2.00	0.62
1:C:254:PHE:CE1	1:C:460:VAL:HG21	2.34	0.61
1:C:58:ARG:O	1:C:61:GLN:HG3	1.99	0.61
1:A:342:ARG:CD	1:B:15:SER:OG	2.49	0.61
1:A:45:ARG:HG2	1:A:59:HIS:HB3	1.83	0.61
1:D:10:MET:HE3	1:D:195:ARG:NH2	2.18	0.59
1:C:254:PHE:CD1	1:C:460:VAL:HG21	2.38	0.58
1:A:103:ASN:HA	1:A:127:ASP:OD2	2.04	0.58
1:A:269:ILE:HG23	1:A:449:GLU:HB2	1.87	0.57
1:A:76:THR:HG22	1:A:408:ASP:HB2	1.86	0.56
1:C:285:LYS:NZ	1:C:438:ASP:OD1	2.37	0.56
1:C:282:LEU:O	1:C:286:ILE:HG13	2.06	0.55
1:C:254:PHE:CD2	1:C:258:TYR:HE2	2.25	0.55
1:C:282:LEU:HD23	1:C:312:ILE:HD11	1.89	0.54
1:B:493:ASN:O	1:C:498:LEU:HB2	2.07	0.54
1:C:497:ASN:O	1:C:498:LEU:C	2.50	0.54
1:A:17:TYR:O	1:A:21:THR:HG23	2.08	0.53
1:D:341:LEU:HD23	1:D:341:LEU:H	1.73	0.53
1:C:78:GLY:N	1:C:485:ASP:OD2	2.41	0.53
1:A:399:ILE:HG21	1:A:409:MET:HE1	1.91	0.53
1:C:225:ALA:HB3	1:C:228:GLU:HB2	1.90	0.53
1:C:6:ILE:C	1:C:178:MET:HE1	2.34	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:ILE:C	1:A:178:MET:HE1	2.35	0.52
1:C:265:PHE:CE2	1:C:270:SER:HB2	2.44	0.52
1:D:269:ILE:HD13	1:D:446:VAL:HG12	1.91	0.52
1:A:148:GLU:O	1:A:151:ASN:N	2.43	0.52
1:A:336:MET:HA	1:A:336:MET:HE2	1.91	0.51
1:B:495:VAL:O	1:C:497:ASN:HA	2.10	0.51
1:C:76:THR:HG22	1:C:408:ASP:HB2	1.92	0.51
1:C:400:ARG:NH1	1:C:417:ALA:O	2.43	0.51
1:A:58:ARG:O	1:A:61:GLN:HG3	2.09	0.51
1:D:205:MET:CG	1:D:213:LEU:HD22	2.41	0.51
1:A:399:ILE:CG2	1:A:409:MET:HE1	2.41	0.51
1:C:269:ILE:HG23	1:C:449:GLU:HB2	1.93	0.51
1:A:8:GLY:HA2	1:A:203:GLU:HG2	1.91	0.50
1:A:103:ASN:ND2	1:A:474:TYR:OH	2.40	0.50
1:A:34:ILE:HD11	1:A:42:PHE:HB3	1.93	0.49
1:A:96:ILE:HD13	1:A:139:LEU:HD12	1.94	0.49
1:B:410:PRO:HB2	1:B:414:GLU:HB2	1.95	0.48
1:C:290:TRP:CG	1:C:293:LEU:HD12	2.49	0.48
1:A:483:ASN:O	1:A:487:LYS:HG3	2.13	0.48
1:C:17:TYR:O	1:C:21:THR:HG23	2.13	0.48
1:A:343:ARG:HD3	1:A:348:ARG:HG2	1.96	0.48
1:C:457:GLU:OE2	1:C:459:SER:OG	2.28	0.48
1:B:34:ILE:HA	1:B:43:ARG:O	2.14	0.47
1:D:269:ILE:HG23	1:D:449:GLU:HB2	1.96	0.47
1:C:474:TYR:CE2	1:C:478:LEU:HD11	2.49	0.47
1:C:254:PHE:CD2	1:C:258:TYR:CE2	3.02	0.47
4:A:671:HOH:O	1:B:333:THR:HG22	2.14	0.47
1:B:86:GLN:HB3	1:B:87:PRO:HA	1.95	0.47
1:C:205:MET:HE2	1:C:213:LEU:HD23	1.96	0.47
1:D:86:GLN:HB3	1:D:87:PRO:HA	1.96	0.47
1:D:94:TYR:CE2	1:D:147:LEU:HD11	2.50	0.47
1:B:67:MET:HG2	1:B:172:MET:SD	2.55	0.46
1:D:251:PRO:HG2	1:D:456:PHE:CD1	2.50	0.46
1:B:327:ASN:HB3	4:B:682:HOH:O	2.14	0.46
1:D:171:ALA:HB3	1:D:179:VAL:HG22	1.96	0.46
1:C:78:GLY:HA2	1:C:125:THR:OG1	2.16	0.46
1:D:336:MET:HE2	1:D:336:MET:HA	1.98	0.46
1:B:269:ILE:HG23	1:B:449:GLU:HB2	1.98	0.45
1:B:348:ARG:CZ	1:B:348:ARG:HA	2.46	0.45
1:C:86:GLN:HB3	1:C:87:PRO:HA	1.98	0.45
1:C:67:MET:HG2	1:C:172:MET:SD	2.57	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:285:LYS:HE2	1:C:435:ILE:HD11	1.99	0.45
1:D:10:MET:HB2	1:D:11:PRO:HD2	1.99	0.45
1:C:485:ASP:O	1:C:489:VAL:HG23	2.17	0.45
1:A:336:MET:HB3	1:A:339:GLN:HB2	1.98	0.44
1:B:346:VAL:HG11	1:B:377:GLN:HB3	1.99	0.44
1:C:285:LYS:HG2	1:C:435:ILE:HD11	1.99	0.44
1:B:447:ARG:HG3	1:B:456:PHE:HE1	1.82	0.44
4:C:566:HOH:O	1:D:327:ASN:HB3	2.16	0.44
1:D:93:PRO:HB2	1:D:94:TYR:CD2	2.52	0.44
1:C:74:TYR:CD1	1:C:75:PRO:HD2	2.52	0.44
1:D:67:MET:HE1	1:D:177:GLY:HA2	2.00	0.44
1:A:457:GLU:OE2	1:A:459:SER:OG	2.35	0.44
1:B:10:MET:HB2	1:B:11:PRO:HD2	2.00	0.44
1:B:100:HIS:HB2	1:B:169:CYS:SG	2.58	0.44
1:A:494:GLU:O	1:A:498:LEU:HG	2.18	0.43
1:D:67:MET:HG2	1:D:172:MET:SD	2.58	0.43
1:D:342:ARG:O	1:D:342:ARG:HG3	2.18	0.43
1:D:10:MET:HB2	1:D:11:PRO:CD	2.48	0.43
1:C:34:ILE:HD11	1:C:42:PHE:HB3	2.00	0.43
1:D:379:ILE:HD13	1:D:431:ALA:HB2	2.01	0.43
1:A:322:GLN:HG3	4:A:676:HOH:O	2.17	0.43
1:C:112:LYS:HZ3	1:C:116:GLU:CD	2.24	0.43
1:A:265:PHE:CE2	1:A:270:SER:HB2	2.53	0.43
1:B:346:VAL:HG22	1:B:346:VAL:O	2.18	0.43
1:A:32:GLY:HA3	1:A:87:PRO:HD3	2.00	0.43
1:A:246:ASN:N	1:A:247:PRO:CD	2.82	0.43
1:C:14:GLN:NE2	1:D:348:ARG:HH22	2.16	0.43
1:C:254:PHE:HD2	1:C:258:TYR:CE2	2.36	0.43
1:A:495:VAL:HG11	1:C:57:ALA:O	2.19	0.43
1:D:17:TYR:O	1:D:21:THR:HG23	2.18	0.43
1:A:230:ILE:HG23	1:A:238:LEU:CD1	2.49	0.42
1:D:336:MET:O	1:D:339:GLN:CB	2.67	0.42
1:B:341:LEU:HD23	1:B:341:LEU:N	2.28	0.42
1:C:399:ILE:O	1:C:417:ALA:HB3	2.18	0.42
1:D:285:LYS:HG2	1:D:435:ILE:HD11	2.02	0.42
1:A:410:PRO:HB2	1:A:414:GLU:HB2	2.01	0.42
1:B:64:GLN:HG2	1:D:491:ARG:NH2	2.35	0.42
1:C:246:ASN:N	1:C:247:PRO:CD	2.82	0.42
1:C:335:ILE:HD12	1:C:335:ILE:H	1.83	0.42
1:A:6:ILE:O	1:A:178:MET:HE1	2.20	0.42
1:C:80:SER:HB3	1:C:485:ASP:CG	2.44	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:94:TYR:CE2	1:B:147:LEU:HD11	2.55	0.42
1:D:346:VAL:O	1:D:346:VAL:HG22	2.20	0.42
1:A:163:ILE:HD11	1:A:167:TYR:CZ	2.54	0.41
1:C:145:TYR:CD1	1:C:146:PRO:HA	2.55	0.41
1:C:48:ASN:ND2	1:C:81:SER:O	2.43	0.41
1:C:457:GLU:CD	1:C:459:SER:HG	2.28	0.41
1:A:67:MET:HG2	1:A:172:MET:SD	2.60	0.41
1:A:82:ALA:C	3:A:502[A]:G4P:HN21	2.28	0.41
1:B:17:TYR:O	1:B:21:THR:HG23	2.20	0.41
1:B:417:ALA:HA	1:B:425:ILE:HD11	2.01	0.41
1:A:86:GLN:HB3	1:A:87:PRO:HA	2.02	0.41
1:A:148:GLU:O	1:A:149:ALA:C	2.63	0.41
1:C:9:VAL:HG11	1:C:201:ARG:HB3	2.02	0.41
1:D:193:GLY:HA3	1:D:213:LEU:HD13	2.03	0.41
1:D:253:LEU:HD12	1:D:253:LEU:HA	1.94	0.41
1:A:443:ILE:O	1:A:447:ARG:HB2	2.20	0.40
1:A:5:GLY:O	1:A:205:MET:HA	2.21	0.40
1:A:10:MET:HB2	1:A:11:PRO:CD	2.51	0.40
1:A:336:MET:HA	1:A:337:PRO:HD2	1.98	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	496/498 (100%)	477 (96%)	19 (4%)	0	100	100
1	B	493/498 (99%)	478 (97%)	15 (3%)	0	100	100
1	C	496/498 (100%)	469 (95%)	27 (5%)	0	100	100
1	D	492/498 (99%)	481 (98%)	11 (2%)	0	100	100
All	All	1977/1992 (99%)	1905 (96%)	72 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	416/417 (100%)	412 (99%)	4 (1%)	68	64
1	B	412/417 (99%)	411 (100%)	1 (0%)	87	87
1	C	417/417 (100%)	415 (100%)	2 (0%)	81	82
1	D	410/417 (98%)	405 (99%)	5 (1%)	63	57
All	All	1655/1668 (99%)	1643 (99%)	12 (1%)	76	74

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	GLN
1	A	254	PHE
1	A	296	ASP
1	A	342	ARG
1	B	100	HIS
1	C	100	HIS
1	C	254	PHE
1	D	100	HIS
1	D	131	LEU
1	D	235	GLU
1	D	263	ASP
1	D	336	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	103	ASN
1	A	185	ASN
1	A	200	ASN
1	A	246	ASN
1	A	437	GLN

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Mol	Chain	Res	Type
1	B	322	GLN
1	B	418	HIS
1	C	66	ASN
1	D	64	GLN
1	D	91	ASN
1	D	322	GLN
1	D	418	HIS
1	D	454	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 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	G4P	D	502[B]	2	36,38,38	1.15	4 (11%)	56,61,61	1.80	13 (23%)
3	G4P	A	502[A]	2	36,38,38	1.45	5 (13%)	56,61,61	1.73	14 (25%)
3	G4P	A	502[B]	2	36,38,38	1.21	4 (11%)	56,61,61	1.68	12 (21%)
3	G4P	D	502[A]	2	36,38,38	1.20	3 (8%)	56,61,61	1.84	12 (21%)

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	G4P	D	502[B]	2	-	5/27/43/43	0/3/3/3
3	G4P	A	502[A]	2	-	5/27/43/43	0/3/3/3
3	G4P	A	502[B]	2	-	4/27/43/43	0/3/3/3
3	G4P	D	502[A]	2	-	5/27/43/43	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502[A]	G4P	PC-O3C	5.51	1.65	1.59
3	A	502[B]	G4P	PC-O3C	3.15	1.62	1.59
3	D	502[A]	G4P	C5-C4	2.94	1.46	1.38
3	D	502[A]	G4P	PC-O3C	2.92	1.62	1.59
3	D	502[B]	G4P	C5-C4	2.81	1.46	1.38
3	A	502[A]	G4P	C5-C4	2.81	1.46	1.38
3	A	502[B]	G4P	C5-C4	2.80	1.46	1.38
3	A	502[B]	G4P	C6-N1	-2.75	1.33	1.38
3	D	502[B]	G4P	PC-O3C	2.67	1.62	1.59
3	D	502[A]	G4P	C6-N1	-2.53	1.34	1.38
3	D	502[B]	G4P	C6-N1	-2.36	1.34	1.38
3	A	502[A]	G4P	PA-O3A	2.23	1.61	1.59
3	A	502[A]	G4P	C6-N1	-2.20	1.34	1.38
3	A	502[A]	G4P	C4-N9	-2.16	1.32	1.38
3	D	502[B]	G4P	C4-N9	-2.10	1.32	1.38
3	A	502[B]	G4P	C4-N9	-2.10	1.32	1.38

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	502[A]	G4P	C5-C4-N3	-5.90	118.99	128.39
3	D	502[B]	G4P	C5-C4-N3	-5.41	119.78	128.39
3	A	502[A]	G4P	C5-C4-N3	-5.40	119.80	128.39
3	D	502[A]	G4P	C2-N3-C4	5.35	121.51	112.30
3	D	502[B]	G4P	C2-N3-C4	5.05	121.00	112.30
3	A	502[B]	G4P	C5-C4-N3	-4.84	120.69	128.39
3	A	502[A]	G4P	C2-N3-C4	4.80	120.57	112.30
3	D	502[A]	G4P	C6-C5-N7	4.41	138.31	130.29
3	D	502[B]	G4P	C6-C5-N7	4.32	138.15	130.29
3	A	502[B]	G4P	C2-N3-C4	4.20	119.54	112.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502[A]	G4P	N9-C4-N3	4.12	134.19	125.95
3	A	502[B]	G4P	C6-C5-N7	4.09	137.74	130.29
3	D	502[A]	G4P	N9-C4-N3	3.91	133.78	125.95
3	D	502[B]	G4P	N9-C4-N3	3.72	133.39	125.95
3	A	502[B]	G4P	N9-C4-N3	3.66	133.28	125.95
3	A	502[A]	G4P	C6-C5-N7	3.65	136.93	130.29
3	D	502[A]	G4P	C4-C5-N7	-3.21	105.58	110.67
3	D	502[A]	G4P	O4'-C4'-C5'	-3.19	99.12	109.33
3	D	502[B]	G4P	C4-C5-N7	-2.95	105.99	110.67
3	A	502[B]	G4P	C4-C5-N7	-2.88	106.11	110.67
3	D	502[B]	G4P	O4'-C4'-C5'	-2.88	100.12	109.33
3	A	502[B]	G4P	N9-C8-N7	-2.62	108.54	113.40
3	A	502[A]	G4P	O4'-C1'-N9	2.62	114.29	108.36
3	A	502[A]	G4P	C4-C5-N7	-2.60	106.55	110.67
3	D	502[A]	G4P	O4'-C1'-N9	2.58	114.20	108.36
3	A	502[B]	G4P	O4'-C1'-N9	2.54	114.11	108.36
3	A	502[A]	G4P	O4'-C4'-C5'	-2.50	101.34	109.33
3	A	502[B]	G4P	C8-N9-C4	2.49	110.69	106.03
3	A	502[B]	G4P	O4'-C4'-C5'	-2.44	101.52	109.33
3	A	502[B]	G4P	C8-N7-C5	2.44	108.60	104.26
3	D	502[A]	G4P	O2C-PC-O3C	2.42	113.80	107.27
3	A	502[B]	G4P	C1'-N9-C4	-2.36	119.50	126.49
3	D	502[B]	G4P	O2C-PC-O3C	2.35	113.63	107.27
3	A	502[A]	G4P	O6-C6-C5	-2.28	120.53	126.53
3	D	502[A]	G4P	C8-N7-C5	2.25	108.27	104.26
3	D	502[B]	G4P	O2A-PA-O3A	2.24	113.33	107.27
3	D	502[B]	G4P	O3B-PB-O2B	2.23	116.16	107.80
3	A	502[A]	G4P	C8-N9-C4	2.22	110.18	106.03
3	A	502[A]	G4P	C2-N1-C6	-2.20	121.13	125.11
3	D	502[A]	G4P	C5-C6-N1	2.18	118.79	113.25
3	D	502[B]	G4P	C5-C6-N1	2.16	118.76	113.25
3	A	502[A]	G4P	N9-C8-N7	-2.16	109.40	113.40
3	A	502[B]	G4P	O3B-PB-O2B	2.15	115.85	107.80
3	D	502[B]	G4P	C1'-N9-C4	-2.14	120.16	126.49
3	D	502[A]	G4P	C1'-N9-C4	-2.11	120.25	126.49
3	A	502[A]	G4P	C5-C6-N1	2.11	118.61	113.25
3	D	502[B]	G4P	O4'-C1'-N9	2.05	113.01	108.36
3	D	502[A]	G4P	O2A-PA-O3A	2.04	112.79	107.27
3	A	502[A]	G4P	C1'-N9-C4	-2.04	120.47	126.49
3	D	502[B]	G4P	C8-N7-C5	2.04	107.89	104.26
3	A	502[A]	G4P	C8-N7-C5	2.01	107.84	104.26

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502[A]	G4P	PC-O3C-PD-O2D
3	D	502[A]	G4P	PA-O3A-PB-O2B
3	A	502[A]	G4P	C3'-C4'-C5'-O5'
3	D	502[A]	G4P	C3'-C4'-C5'-O5'
3	D	502[B]	G4P	C3'-C4'-C5'-O5'
3	A	502[A]	G4P	O4'-C4'-C5'-O5'
3	A	502[B]	G4P	O4'-C4'-C5'-O5'
3	A	502[B]	G4P	C3'-C4'-C5'-O5'
3	D	502[A]	G4P	O4'-C4'-C5'-O5'
3	D	502[B]	G4P	O4'-C4'-C5'-O5'
3	A	502[B]	G4P	PC-O3C-PD-O2D
3	D	502[B]	G4P	C5'-O5'-PA-O1A
3	A	502[B]	G4P	C4'-C5'-O5'-PA
3	A	502[A]	G4P	C4'-C5'-O5'-PA
3	D	502[A]	G4P	C4'-C5'-O5'-PA
3	D	502[B]	G4P	C4'-C5'-O5'-PA
3	A	502[A]	G4P	PA-O3A-PB-O2B
3	D	502[B]	G4P	PA-O3A-PB-O2B
3	D	502[A]	G4P	C3'-O3'-PC-O1C

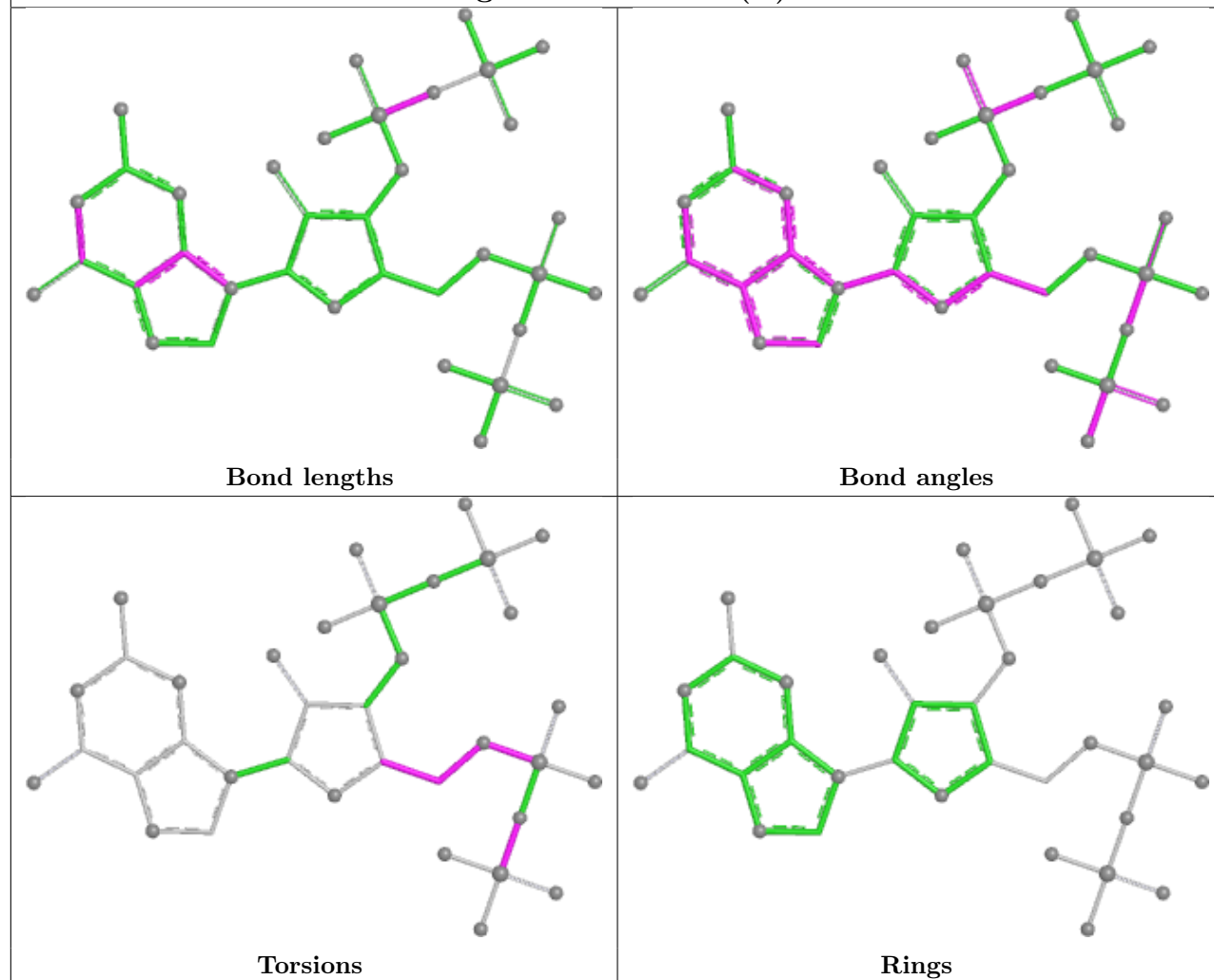
There are no ring outliers.

1 monomer is involved in 1 short contact:

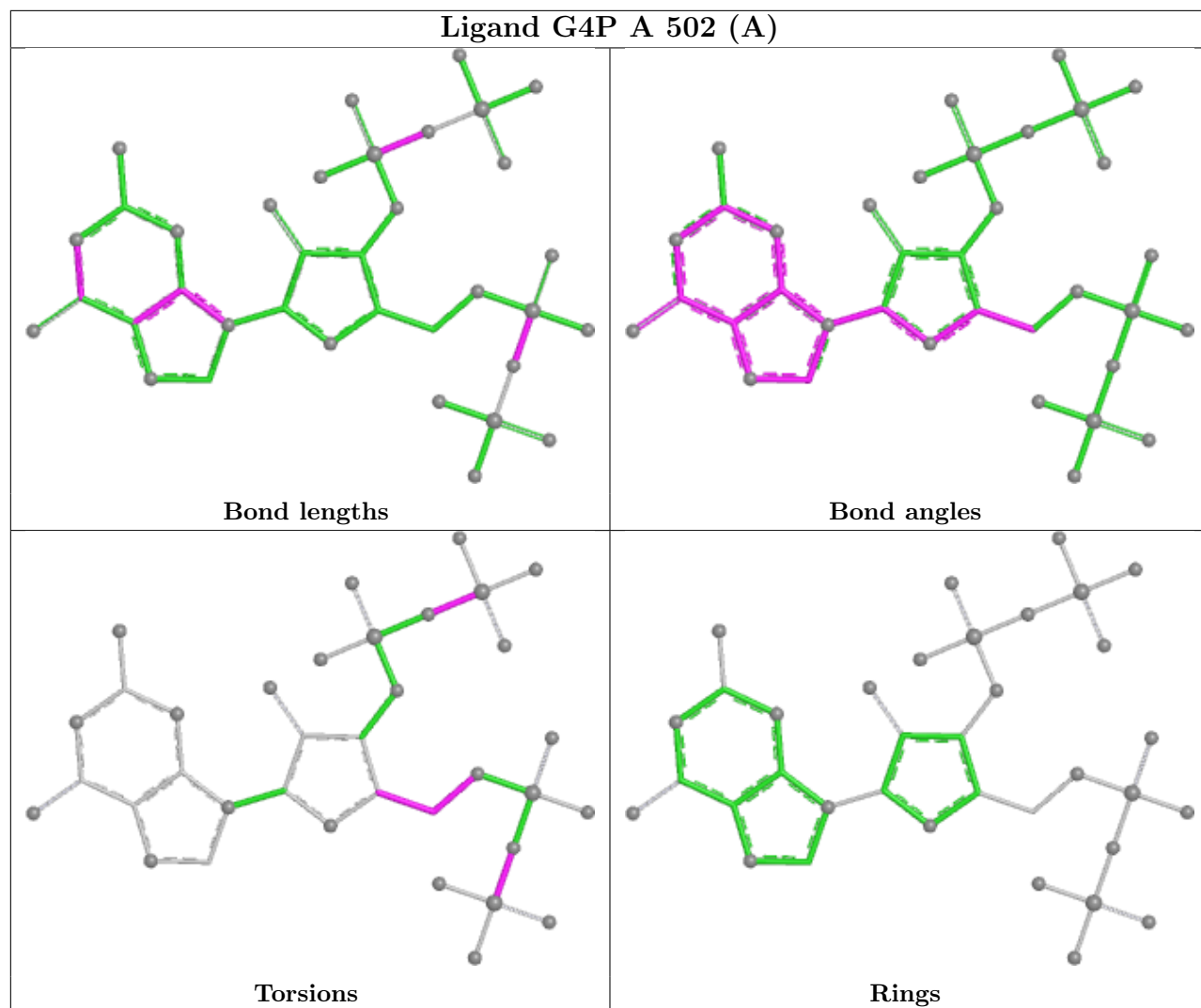
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502[A]	G4P	1	0

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

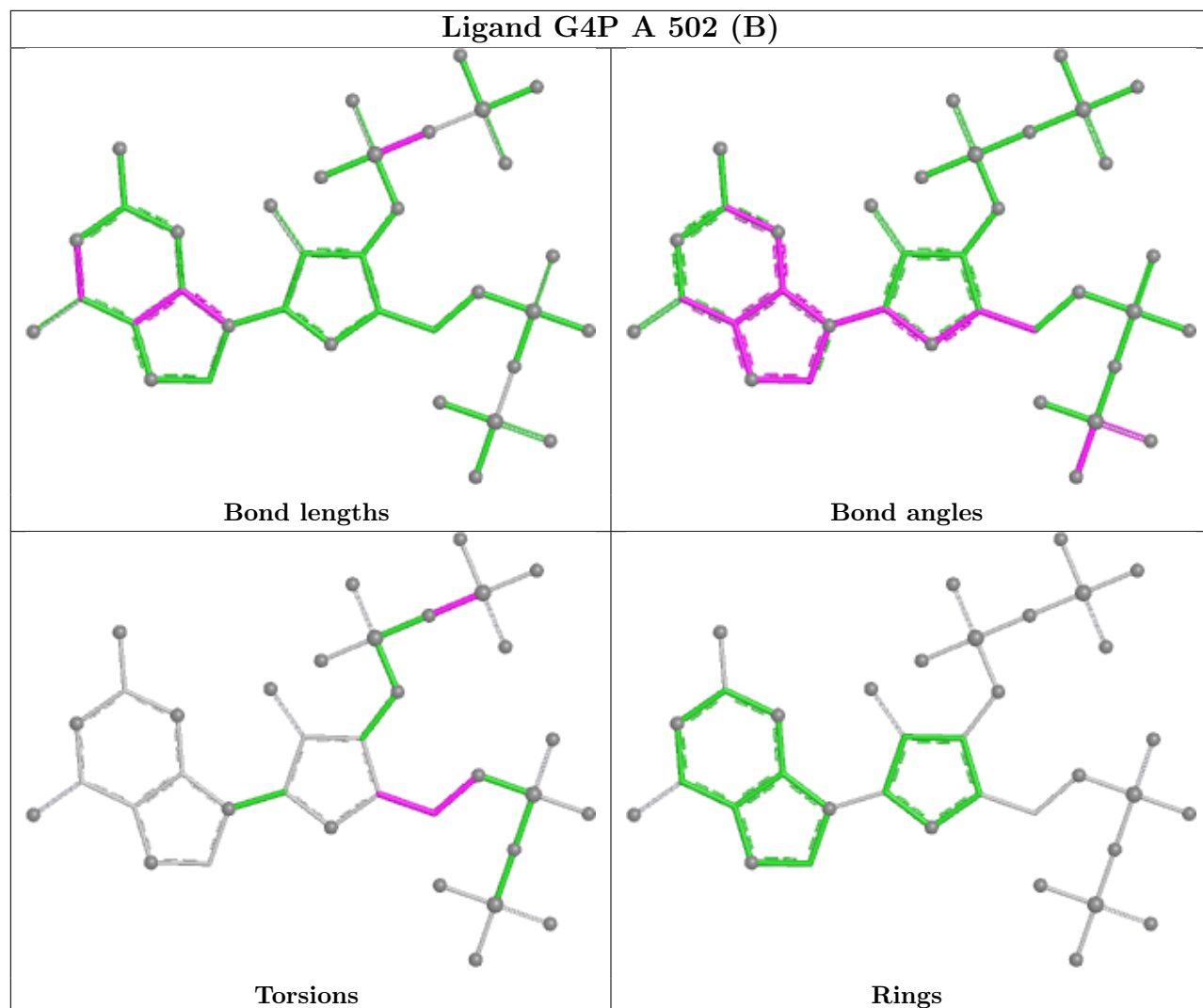
Ligand G4P D 502 (B)

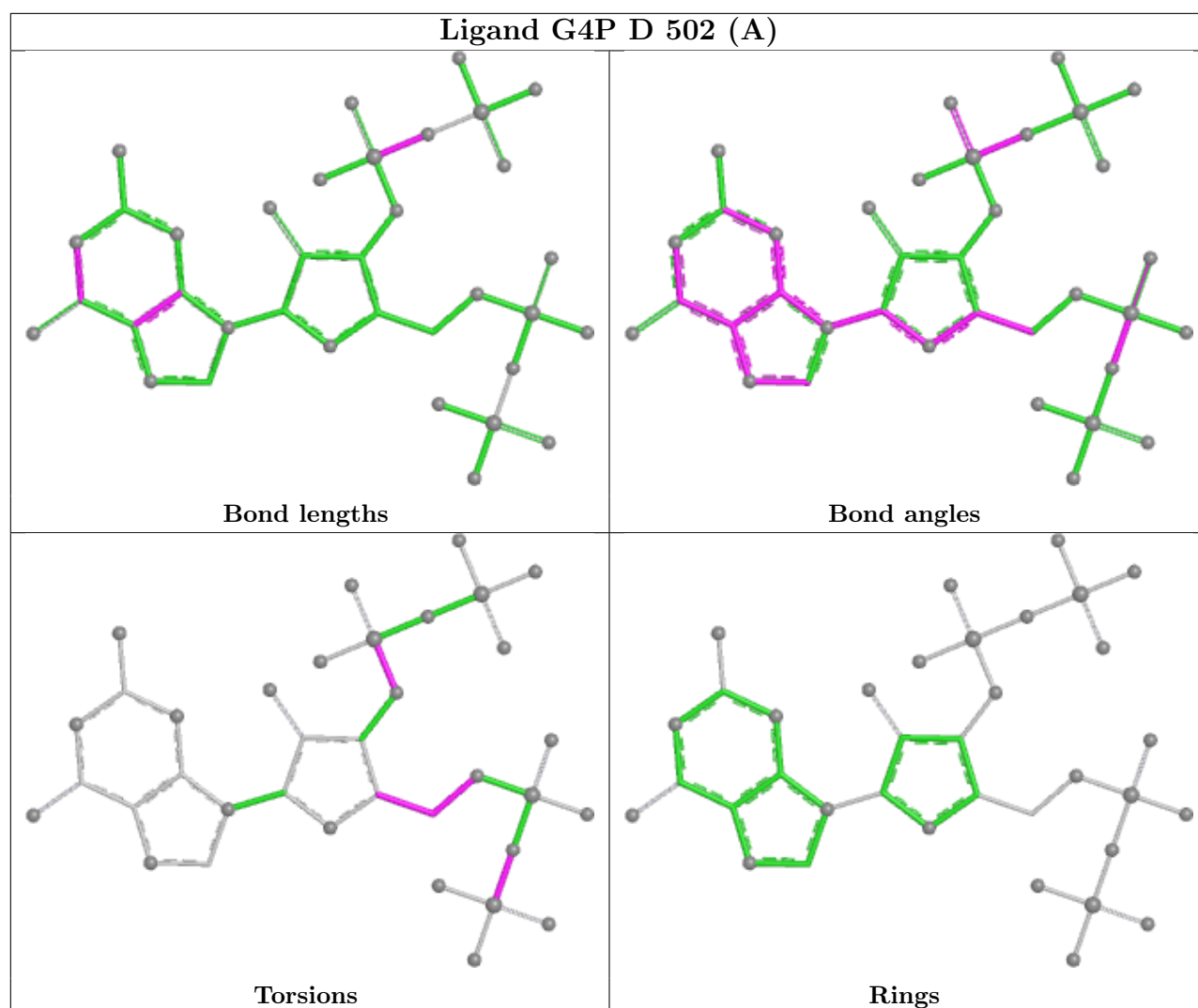


Ligand G4P A 502 (A)



Ligand G4P A 502 (B)





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	498/498 (100%)	0.59	34 (6%)	23	26	34, 52, 77, 94	0
1	B	495/498 (99%)	0.09	17 (3%)	48	54	28, 43, 77, 99	0
1	C	498/498 (100%)	0.84	42 (8%)	17	19	36, 56, 84, 93	0
1	D	494/498 (99%)	0.17	19 (3%)	44	49	30, 45, 80, 103	0
All	All	1985/1992 (99%)	0.42	112 (5%)	30	33	28, 49, 80, 103	0

All (112) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	498	LEU	5.5
1	A	495	VAL	4.4
1	C	495	VAL	4.3
1	C	480	THR	3.8
1	A	498	LEU	3.7
1	B	344	LYS	3.7
1	B	258	TYR	3.7
1	C	248	VAL	3.6
1	C	477	PHE	3.6
1	D	337	PRO	3.5
1	C	466	VAL	3.5
1	B	495	VAL	3.5
1	B	343	ARG	3.4
1	D	341	LEU	3.4
1	C	488	ALA	3.3
1	A	144	HIS	3.2
1	D	489	VAL	3.2
1	D	258	TYR	3.2
1	B	341	LEU	3.2
1	B	254	PHE	3.2
1	B	336	MET	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	337	PRO	3.1
1	C	478	LEU	3.1
1	D	344	LYS	3.1
1	D	343	ARG	2.9
1	A	77	ALA	2.9
1	A	496	GLU	2.9
1	A	474	TYR	2.9
1	A	486	ALA	2.8
1	D	1	CYS	2.8
1	A	234	GLU	2.8
1	C	84	GLU	2.8
1	C	79	SER	2.7
1	C	77	ALA	2.7
1	C	417	ALA	2.7
1	D	477	PHE	2.7
1	B	339	GLN	2.7
1	A	248	VAL	2.7
1	A	38	ALA	2.7
1	A	142	PHE	2.7
1	A	76	THR	2.7
1	C	481	LEU	2.7
1	C	341	LEU	2.6
1	A	490	GLN	2.6
1	C	469	ASP	2.6
1	A	81	SER	2.6
1	B	76	THR	2.6
1	B	328	ARG	2.5
1	A	488	ALA	2.5
1	C	142	PHE	2.5
1	A	79	SER	2.5
1	A	482	ARG	2.5
1	A	483	ASN	2.4
1	D	490	GLN	2.4
1	A	480	THR	2.4
1	A	236	GLY	2.4
1	B	78	GLY	2.4
1	D	478	LEU	2.4
1	B	337	PRO	2.4
1	A	84	GLU	2.4
1	C	460	VAL	2.4
1	C	164	ARG	2.3
1	D	342	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	491	ARG	2.3
1	D	348	ARG	2.3
1	D	76	THR	2.3
1	C	258	TYR	2.3
1	B	489	VAL	2.3
1	C	482	ARG	2.2
1	D	254	PHE	2.2
1	B	490	GLN	2.2
1	C	147	LEU	2.2
1	C	81	SER	2.2
1	B	486	ALA	2.2
1	C	418	HIS	2.2
1	A	470	VAL	2.2
1	C	244	ALA	2.2
1	C	254	PHE	2.2
1	C	474	TYR	2.2
1	C	162	LEU	2.2
1	A	146	PRO	2.2
1	A	61	GLN	2.1
1	A	478	LEU	2.1
1	B	478	LEU	2.1
1	C	293	LEU	2.1
1	D	336	MET	2.1
1	C	448	ALA	2.1
1	A	115	PHE	2.1
1	A	54	VAL	2.1
1	D	491	ARG	2.1
1	A	47	ALA	2.1
1	C	431	ALA	2.1
1	D	339	GLN	2.1
1	C	115	PHE	2.1
1	C	165	GLY	2.1
1	D	78	GLY	2.1
1	C	336	MET	2.1
1	C	422	VAL	2.1
1	A	143	ARG	2.1
1	B	345	SER	2.0
1	C	335	ILE	2.0
1	D	487	LYS	2.0
1	C	226	PRO	2.0
1	A	341	LEU	2.0
1	C	253	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	412	ALA	2.0
1	A	479	ASP	2.0
1	A	336	MET	2.0
1	C	372	GLY	2.0
1	A	337	PRO	2.0
1	C	75	PRO	2.0
1	A	489	VAL	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

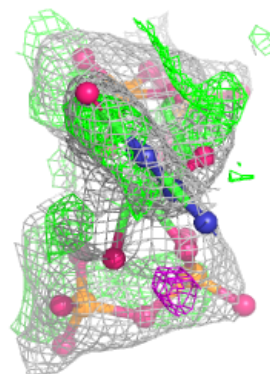
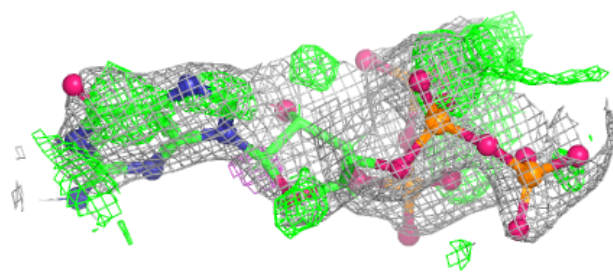
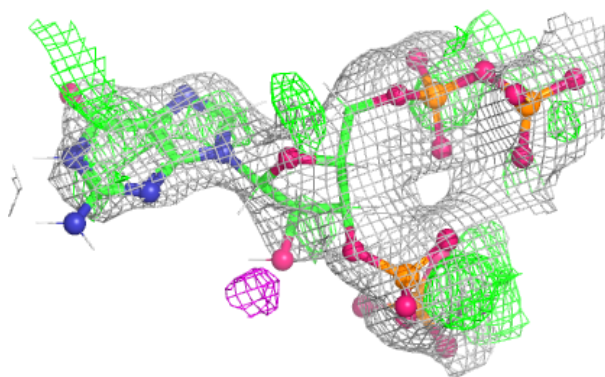
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	G4P	A	502[A]	36/36	0.87	0.18	28,37,51,62	47
3	G4P	A	502[B]	36/36	0.87	0.18	28,36,49,57	47
3	G4P	D	502[A]	36/36	0.92	0.14	26,36,51,62	47
3	G4P	D	502[B]	36/36	0.92	0.14	25,36,49,55	47
2	MG	A	501[A]	1/1	0.95	0.07	36,36,36,36	1
2	MG	A	501[B]	1/1	0.95	0.07	38,38,38,38	1
2	MG	D	501[A]	1/1	0.96	0.09	37,37,37,37	1
2	MG	D	501[B]	1/1	0.96	0.09	37,37,37,37	1

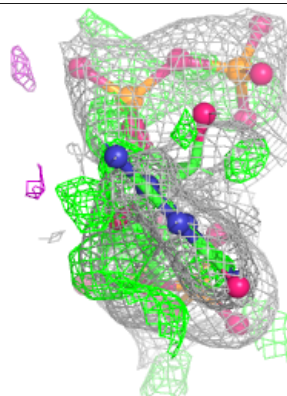
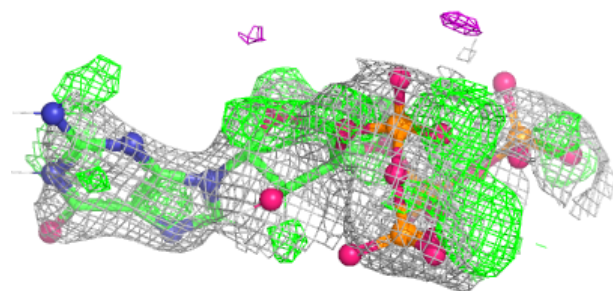
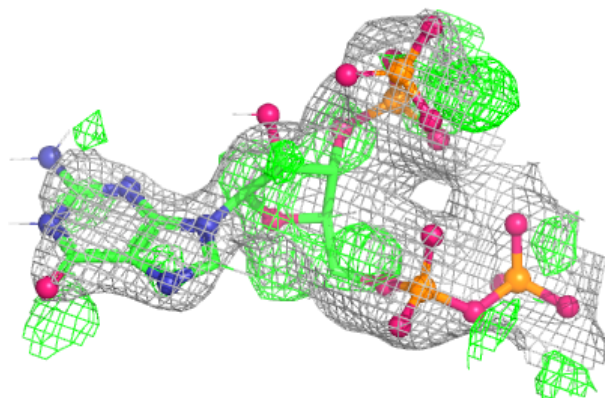
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around G4P A 502 (A):

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

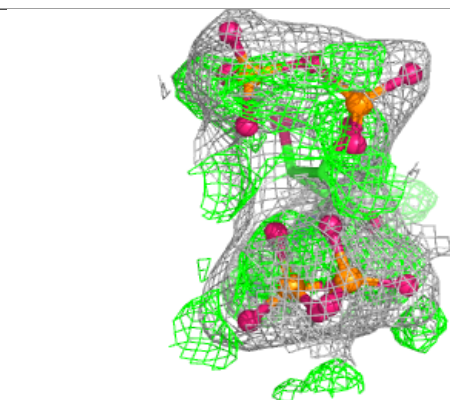
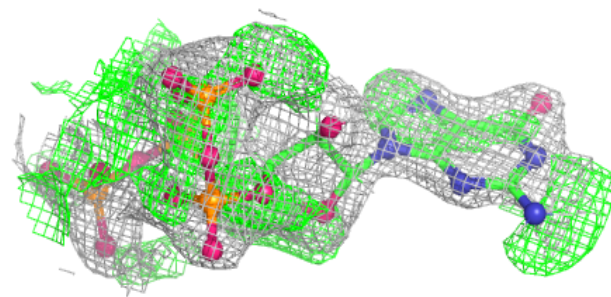
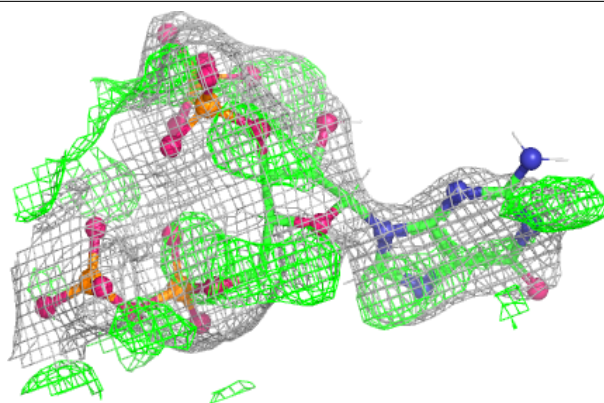
**Electron density around G4P A 502 (B):**

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

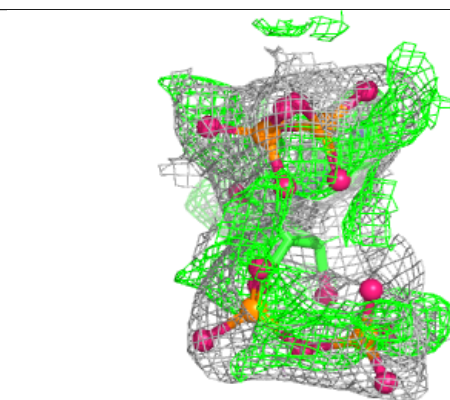
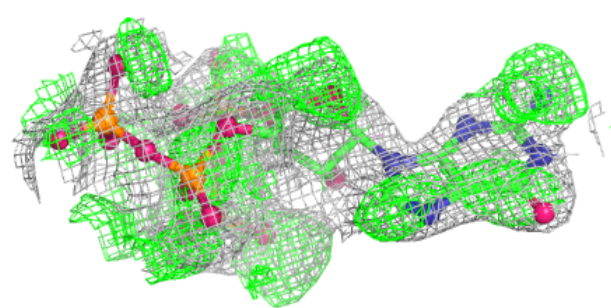
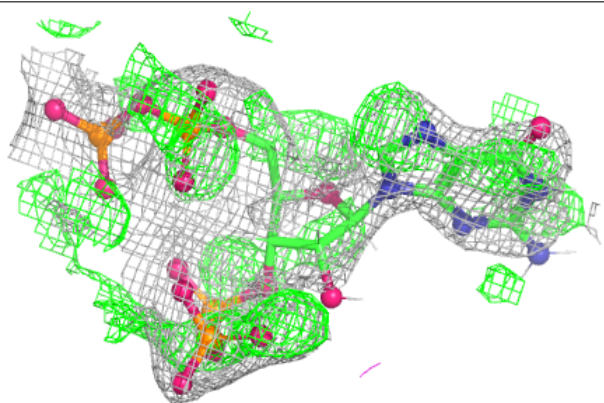


Electron density around G4P D 502 (A):

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

**Electron density around G4P D 502 (B):**

$2mF_o-DF_c$ (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.