



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 08:43 PM UTC

PDB ID : 8YRK / pdb_00008yrk
Title : Tubulin-Compound KY216: stathmin-like domain complex
Authors : Xia, Y.Z.; He, X.Y.
Deposited on : 2024-03-21
Resolution : 2.74 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

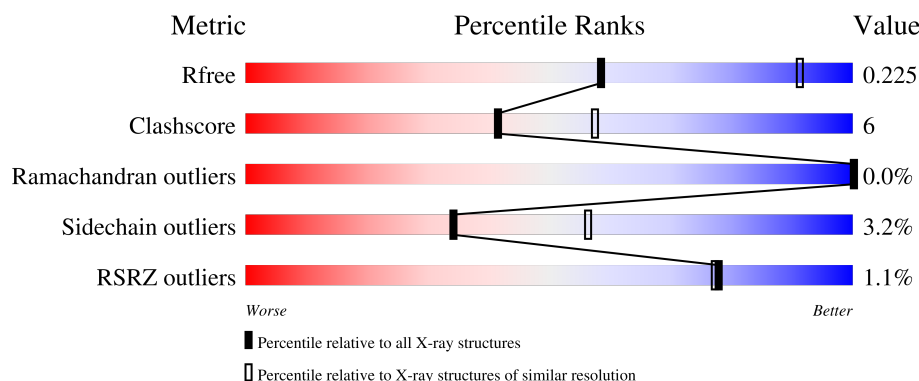
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.74 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	1819 (2.76-2.72)
Clashscore	190562	1866 (2.76-2.72)
Ramachandran outliers	187476	1830 (2.76-2.72)
Sidechain outliers	187428	1831 (2.76-2.72)
RSRZ outliers	180081	1819 (2.76-2.72)

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	450	% 83% 14% • •
1	C	450	% 83% 15% •
2	B	445	78% 17% 5%
2	D	445	% 75% 19% • 5%
3	E	143	% 68% 17% • 14%

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Mol	Chain	Length	Quality of chain
4	F	384	2% 70% 19% 10%

2 Entry composition

There are 15 unique types of molecules in this entry. The entry contains 17720 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 Detyrosinated tubulin alpha-1B chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	437	Total	C	N	O	S	0	4	0
			3431	2173	581	653	24			
1	C	440	Total	C	N	O	S	3	8	0
			3465	2193	585	663	24			

- Molecule 2 is a protein called Tubulin beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	424	Total	C	N	O	S	0	2	0
			3350	2107	572	645	26			
2	D	421	Total	C	N	O	S	0	0	0
			3306	2078	562	640	26			

- Molecule 3 is a protein called Stathmin-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	123	Total	C	N	O	S	0	2	0
			1026	633	186	202	5			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	3	MET	-	initiating methionine	UNP P63042
E	4	ALA	-	expression tag	UNP P63042

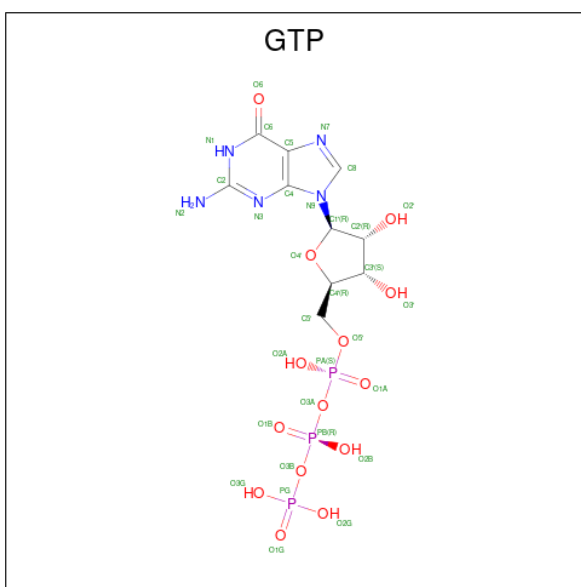
- Molecule 4 is a protein called Tubulin tyrosine ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	346	Total	C	N	O	S	0	4	0
			2859	1835	488	521	15			

There are 39 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	?	-	ALA	deletion	UNP A0A8V0Z8P0
F	?	-	GLU	deletion	UNP A0A8V0Z8P0
F	?	-	MET	deletion	UNP A0A8V0Z8P0
F	?	-	GLN	deletion	UNP A0A8V0Z8P0
F	?	-	GLN	deletion	UNP A0A8V0Z8P0
F	?	-	GLN	deletion	UNP A0A8V0Z8P0
F	?	-	LEU	deletion	UNP A0A8V0Z8P0
F	?	-	LEU	deletion	UNP A0A8V0Z8P0
F	?	-	GLU	deletion	UNP A0A8V0Z8P0
F	?	-	GLY	deletion	UNP A0A8V0Z8P0
F	?	-	ASP	deletion	UNP A0A8V0Z8P0
F	?	-	GLN	deletion	UNP A0A8V0Z8P0
F	?	-	THR	deletion	UNP A0A8V0Z8P0
F	?	-	LEU	deletion	UNP A0A8V0Z8P0
F	?	-	VAL	deletion	UNP A0A8V0Z8P0
F	?	-	LEU	deletion	UNP A0A8V0Z8P0
F	?	-	ALA	deletion	UNP A0A8V0Z8P0
F	?	-	SER	deletion	UNP A0A8V0Z8P0
F	?	-	SER	deletion	UNP A0A8V0Z8P0
F	?	-	THR	deletion	UNP A0A8V0Z8P0
F	?	-	HIS	deletion	UNP A0A8V0Z8P0
F	?	-	PRO	deletion	UNP A0A8V0Z8P0
F	?	-	GLU	deletion	UNP A0A8V0Z8P0
F	?	-	SER	deletion	UNP A0A8V0Z8P0
F	?	-	VAL	deletion	UNP A0A8V0Z8P0
F	?	-	ASP	deletion	UNP A0A8V0Z8P0
F	?	-	SER	deletion	UNP A0A8V0Z8P0
F	?	-	ASP	deletion	UNP A0A8V0Z8P0
F	?	-	LYS	deletion	UNP A0A8V0Z8P0
F	?	-	ASN	deletion	UNP A0A8V0Z8P0
F	?	-	HIS	deletion	UNP A0A8V0Z8P0
F	?	-	GLY	deletion	UNP A0A8V0Z8P0
F	?	-	PHE	deletion	UNP A0A8V0Z8P0
F	379	HIS	-	expression tag	UNP A0A8V0Z8P0
F	380	HIS	-	expression tag	UNP A0A8V0Z8P0
F	381	HIS	-	expression tag	UNP A0A8V0Z8P0
F	382	HIS	-	expression tag	UNP A0A8V0Z8P0
F	383	HIS	-	expression tag	UNP A0A8V0Z8P0
F	384	HIS	-	expression tag	UNP A0A8V0Z8P0

- Molecule 5 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).



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

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

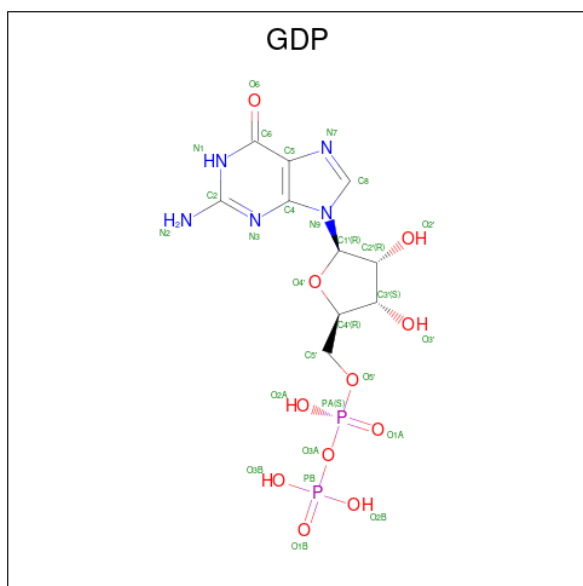
- Molecule 7 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total Ca 1 1	0	0
7	C	1	Total Ca 1 1	0	0

- Molecule 8 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	1	Total	Cl	0	0
			1	1		

- Molecule 9 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).



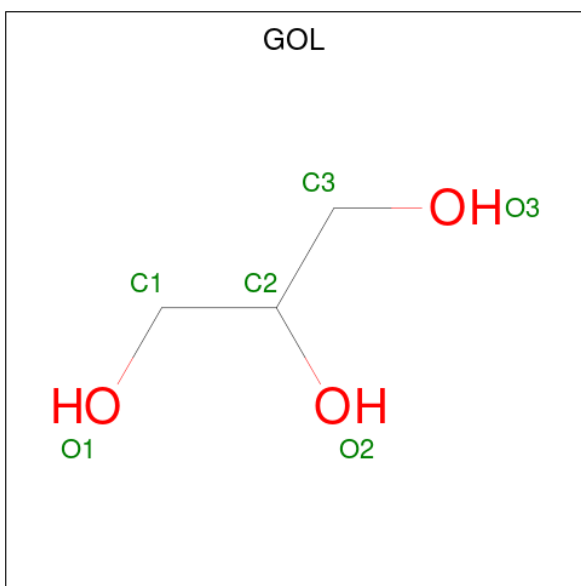
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	B	1	Total 28	C 10	N 5	O 11	P 2	0	0
9	D	1	Total 28	C 10	N 5	O 11	P 2	0	0

- Molecule 10 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
10	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
10	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

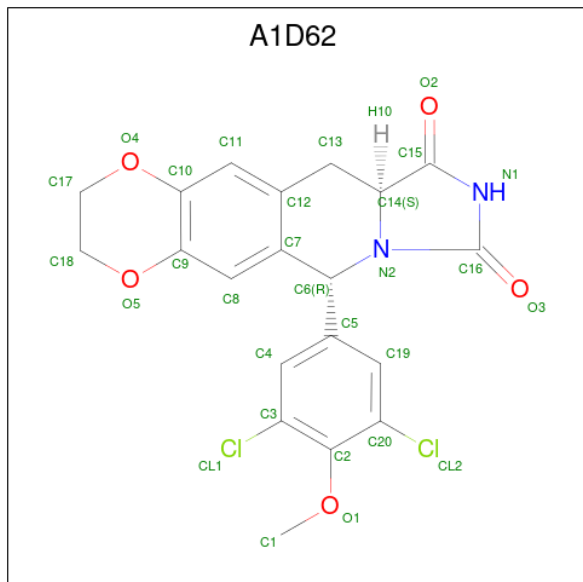
- Molecule 11 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

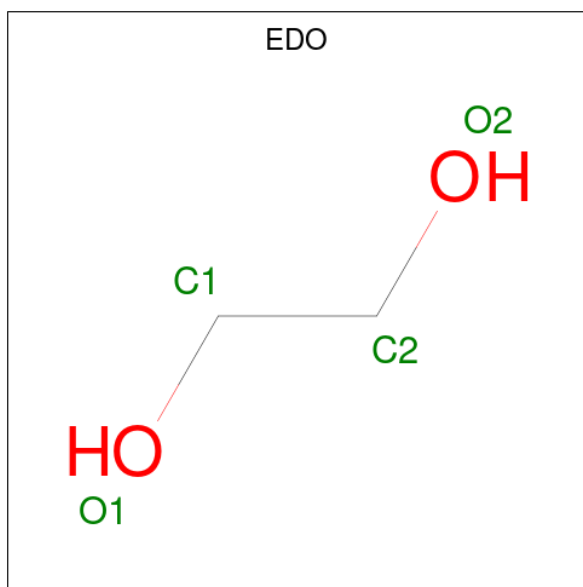
- Molecule 12 is (11R,16S)-11-(3,5-dichloro-4-methoxyphenyl)-4,7-dioxa-12,14-diazatetrac

cyclo[8.7.0.03,8.012,16]heptadeca-1,3(8),9-triene-13,15-dione (CCD ID: A1D62) (formula: $C_{20}H_{16}Cl_2N_2O_5$) (labeled as "Ligand of Interest" by depositor).



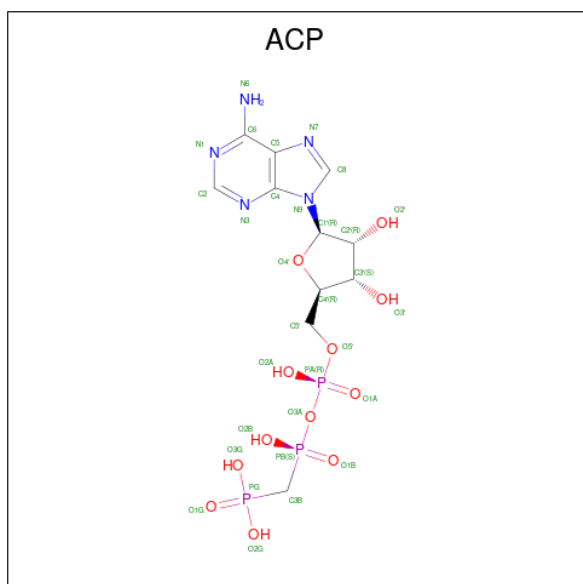
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
12	B	1	Total	C	Cl	N	O	0	0
			29	20	2	2	5		
12	D	1	Total	C	Cl	N	O	0	0
			29	20	2	2	5		

- Molecule 13 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
13	C	1	Total	C	O	0	0
			4	2	2		

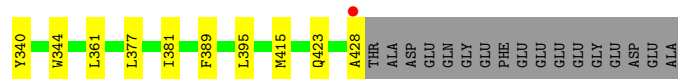
- Molecule 14 is PHOSPHOMETHYLPHOSPHONIC ACID ADENYLATE ESTER (CCD ID: ACP) (formula: $C_{11}H_{18}N_5O_{12}P_3$).



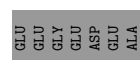
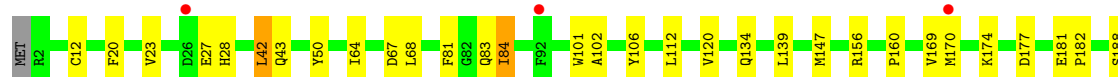
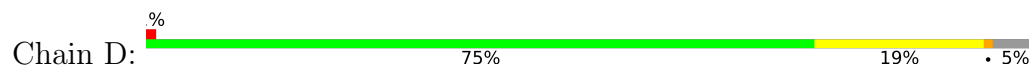
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
14	F	1	Total	C	N	O	P	0	0
			31	11	5	12	3		

- Molecule 15 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	A	5	Total	O	0	0
			5	5		
15	B	9	Total	O	0	0
			9	9		
15	C	12	Total	O	0	0
			12	12		
15	D	6	Total	O	0	0
			6	6		
15	E	1	Total	O	0	0
			1	1		



• Molecule 2: Tubulin beta chain



• Molecule 3: Stathmin-4



• Molecule 4: Tubulin tyrosine ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	104.10Å 158.98Å 179.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.54 – 2.74 43.54 – 2.74	Depositor EDS
% Data completeness (in resolution range)	96.0 (43.54-2.74) 95.9 (43.54-2.74)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.03 (at 2.73Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.173 , 0.226 0.173 , 0.225	Depositor DCC
R_{free} test set	3715 reflections (4.72%)	wwPDB-VP
Wilson B-factor (Å ²)	59.8	Xtriage
Anisotropy	0.133	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 52.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	17720	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.49% 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: EDO, A1D62, MES, GOL, ACP, CL, GTP, GDP, CA, 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	0.17	0/3521	0.34	0/4780
1	C	0.19	0/3564	0.38	0/4838
2	B	0.19	0/3430	0.37	0/4645
2	D	0.20	1/3379 (0.0%)	0.35	0/4578
3	E	0.18	0/1041	0.30	0/1382
4	F	0.14	0/2935	0.33	0/3964
All	All	0.18	1/17870 (0.0%)	0.35	0/24187

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	356	ILE	N-CA	5.54	1.50	1.46

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	3431	0	3352	40	0
1	C	3465	0	3383	36	0
2	B	3350	0	3237	48	0
2	D	3306	0	3182	54	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	1026	0	1042	14	0
4	F	2859	0	2841	40	0
5	A	32	0	12	2	0
5	C	32	0	12	1	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	D	2	0	0	0	0
7	A	1	0	0	0	0
7	C	1	0	0	0	0
8	A	1	0	0	0	0
9	B	28	0	12	0	0
9	D	28	0	12	2	0
10	B	24	0	24	6	0
11	B	6	0	8	1	0
12	B	29	0	0	1	0
12	D	29	0	0	3	0
13	C	4	0	6	0	0
14	F	31	0	14	1	0
15	A	5	0	0	0	0
15	B	9	0	0	0	0
15	C	12	0	0	0	0
15	D	6	0	0	1	0
15	E	1	0	0	0	0
All	All	17720	0	17137	224	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (224) 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)
9:D:502:GDP:O3B	15:D:601:HOH:O	1.85	0.93
1:A:336:LYS:HD2	3:E:24:LEU:HD13	1.62	0.81
1:A:161:TYR:HB3	1:A:164:LYS:HG3	1.64	0.78
2:B:114:ASP:HB3	11:B:505:GOL:H32	1.66	0.75
2:D:330:MET:HG3	2:D:351:THR:HG21	1.68	0.75
4:F:139:ARG:HG2	4:F:145:ASN:HD22	1.51	0.75
2:D:396:HIS:HA	2:D:399:THR:HG22	1.69	0.74
4:F:78:VAL:HG21	4:F:181:VAL:HG21	1.72	0.72
1:C:296:PHE:CZ	1:C:377:MET:HE1	2.26	0.70
2:B:274:THR:HG21	2:B:361:LEU:HD11	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:232:THR:HG21	2:D:300:MET:HG3	1.74	0.69
1:C:71:GLU:HG2	1:C:98:ASP:HB3	1.75	0.68
2:D:315:ALA:HB3	2:D:351:THR:HG22	1.77	0.67
2:D:246:LEU:HD11	12:D:504:A1D62:C15	2.24	0.67
12:B:506:A1D62:C1	12:B:506:A1D62:CL2	2.80	0.67
2:B:170:MET:HG3	2:B:377:LEU:HD21	1.77	0.65
12:D:504:A1D62:C1	12:D:504:A1D62:CL2	2.82	0.65
4:F:340:GLN:HA	4:F:343:TYR:HD2	1.62	0.65
2:D:42:LEU:HD13	2:D:243:PRO:HG2	1.78	0.65
2:D:238:THR:HG21	2:D:316:ILE:HG12	1.79	0.64
4:F:229:ASN:ND2	4:F:232:ASN:OD1	2.30	0.64
1:C:18:ASN:HD21	1:C:78:VAL:HG22	1.64	0.63
2:B:197:ASP:OD2	10:B:503:MES:H52	1.98	0.63
2:B:294:PHE:HB3	10:B:504:MES:H32	1.80	0.62
2:D:193:VAL:HG12	2:D:194:GLU:HG2	1.80	0.62
1:C:11:GLN:HG3	1:C:74:VAL:HG21	1.81	0.61
2:D:284:LEU:HD21	2:D:289:LEU:HD13	1.81	0.61
2:B:174:LYS:HD2	2:B:205:GLU:HG3	1.82	0.61
2:B:273:LEU:HD11	2:B:298:ASN:HA	1.84	0.60
4:F:148:ILE:HD13	4:F:162:ILE:HG12	1.85	0.59
2:D:102:ALA:HB2	2:D:403:MET:HE3	1.85	0.59
4:F:292:ARG:HH12	4:F:380:HIS:HB2	1.68	0.59
1:C:93:ILE:HG22	1:C:114:ILE:HD11	1.84	0.59
1:A:274:PRO:HB2	1:A:276:ILE:HD11	1.84	0.59
2:D:28:HIS:HA	2:D:43:GLN:HB3	1.85	0.59
2:D:332:ASN:O	2:D:336:LYS:HG2	2.03	0.58
2:D:378:PHE:HD1	2:D:415:MET:HE2	1.68	0.58
2:B:267:MET:HE1	2:B:303:CYS:HB2	1.87	0.57
1:C:18:ASN:ND2	1:C:78:VAL:HG22	2.20	0.56
1:C:172:TYR:HB3	1:C:205:ASP:HA	1.89	0.55
2:D:323:MET:HG2	2:D:353:VAL:HG11	1.89	0.55
1:A:275:VAL:HG13	1:A:368:LEU:HD21	1.89	0.55
4:F:14:TYR:HA	4:F:17:VAL:HG12	1.88	0.55
1:A:351:PHE:HE1	3:E:24:LEU:HD11	1.72	0.55
4:F:173:ILE:HA	4:F:176:GLN:HG2	1.89	0.55
1:C:7:ILE:HG21	1:C:153:LEU:HD21	1.88	0.54
4:F:240:LEU:HB2	4:F:245:ILE:HG21	1.89	0.54
2:B:2:ARG:O	2:B:131:GLN:NE2	2.35	0.54
2:B:156:ARG:CZ	10:B:503:MES:H21	2.38	0.54
2:B:31:ASP:HB2	2:B:32:PRO:HD2	1.89	0.54
4:F:136:ASN:O	4:F:140:GLU:HG2	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:247:LYS:HD2	4:F:253:TYR:CZ	2.42	0.54
1:C:4[A]:CYS:SG	1:C:136:LEU:HG	2.48	0.54
1:A:18:ASN:HD21	1:A:78:VAL:HG22	1.73	0.53
1:C:14:VAL:HG13	1:C:67:PHE:HD2	1.74	0.53
2:D:156:ARG:HG2	3:E:123:LEU:HD11	1.90	0.53
4:F:278:THR:HG23	4:F:281:ASN:H	1.74	0.53
4:F:163:SER:HB3	4:F:169:LEU:HG	1.92	0.52
4:F:248:GLU:HB3	4:F:249:TYR:HD1	1.75	0.52
4:F:102:PRO:HD3	4:F:179:VAL:HA	1.92	0.52
4:F:135:TYR:OH	4:F:165:GLU:HA	2.09	0.52
2:B:134:GLN:HA	2:B:165:ASN:O	2.10	0.52
1:C:341:ILE:HD13	1:C:351:PHE:HZ	1.75	0.52
1:A:36:MET:HB3	1:A:61:HIS:CE1	2.45	0.51
1:A:274:PRO:HB3	1:A:286:LEU:HD12	1.92	0.51
2:B:139:LEU:HD12	2:B:170:MET:SD	2.51	0.51
1:C:98:ASP:HB2	5:C:501:GTP:O1G	2.11	0.51
1:C:8:HIS:HB3	1:C:13:GLY:O	2.11	0.51
3:E:9:ILE:HG12	3:E:21:GLU:HB3	1.93	0.51
1:C:231:ILE:HA	1:C:234:ILE:HD12	1.92	0.51
2:D:238:THR:CG2	2:D:316:ILE:HG12	2.40	0.51
1:A:276:ILE:HG12	1:A:283:HIS:CE1	2.46	0.50
1:C:209:ILE:HG23	1:C:230:LEU:HD23	1.91	0.50
2:D:182:PRO:HG2	2:D:388:MET:HE2	1.94	0.50
2:D:170:MET:HE2	2:D:377:LEU:HD11	1.93	0.50
1:A:210:TYR:CE1	1:A:222:PRO:HD2	2.47	0.49
2:B:165:ASN:HD21	2:B:200:TYR:HE2	1.60	0.49
4:F:150:LYS:NZ	4:F:183:GLN:OE1	2.42	0.49
1:A:23:LEU:O	1:A:27:GLU:HG3	2.11	0.49
2:D:293:MET:HE3	2:D:294:PHE:HE1	1.76	0.49
3:E:127:ASP:O	3:E:131:GLU:HG2	2.12	0.49
2:B:306:ARG:HG2	2:B:340:TYR:HE2	1.78	0.49
4:F:148:ILE:HD11	4:F:160:ILE:HG13	1.93	0.49
3:E:131:GLU:O	3:E:135:LYS:HG2	2.12	0.49
2:D:101:TRP:HE3	2:D:403:MET:HE1	1.77	0.49
2:D:215:LEU:O	2:D:216:LYS:HG2	2.13	0.49
1:A:262:TYR:HE1	1:A:264:ARG:HB2	1.77	0.49
4:F:14:TYR:O	4:F:18:SER:OG	2.31	0.49
2:B:16[B]:ILE:HD13	2:B:229:VAL:HG11	1.94	0.49
3:E:58:GLU:O	3:E:62:LYS:HG3	2.12	0.49
2:D:101:TRP:CE3	2:D:403:MET:HE1	2.48	0.48
2:D:293:MET:HE3	2:D:294:PHE:CE1	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:68:LEU:HD12	2:B:97:ALA:HB2	1.96	0.48
4:F:195:GLY:HA3	4:F:197:ARG:HD3	1.95	0.48
1:A:152:LEU:HD11	3:E:54:LEU:HD11	1.95	0.48
2:D:160:PRO:HB2	3:E:116:LEU:HD11	1.95	0.47
1:A:112:LYS:HA	1:A:112:LYS:HD2	1.76	0.47
1:C:117:LEU:HD11	1:C:121:ARG:NH2	2.30	0.47
2:D:177:ASP:OD1	2:D:177:ASP:N	2.45	0.47
1:A:134:GLY:HA3	1:A:165:SER:O	2.15	0.47
1:C:254:GLU:HG2	1:C:352:LYS:HE2	1.95	0.47
4:F:226:GLU:HG3	4:F:237:THR:HB	1.95	0.47
1:C:210:TYR:CZ	1:C:222:PRO:HD2	2.50	0.47
1:A:56:THR:HG22	1:A:60:LYS:H	1.80	0.47
2:B:11:GLN:HA	2:B:72:THR:HG21	1.96	0.47
1:A:96:LYS:NZ	2:B:128:ASP:OD2	2.48	0.47
3:E:6:MET:SD	3:E:6:MET:N	2.88	0.46
2:D:106:TYR:CE2	2:D:403:MET:HE2	2.50	0.46
4:F:15:ALA:O	4:F:19:ARG:HG3	2.15	0.46
2:D:101:TRP:HE1	2:D:188:SER:HB3	1.80	0.46
2:D:285:THR:HG22	2:D:288:GLU:CD	2.41	0.46
1:A:15:GLN:NE2	5:A:501:GTP:O6	2.49	0.46
2:B:83:GLN:CD	2:B:83:GLN:H	2.24	0.46
4:F:226:GLU:OE2	4:F:250:SER:OG	2.33	0.46
4:F:252:ASN:OD1	4:F:252:ASN:N	2.48	0.46
2:B:95:SER:OG	2:B:108:GLU:HG2	2.16	0.46
2:B:389:PHE:HE1	2:B:395:LEU:HD11	1.80	0.45
2:D:211:CYS:HA	2:D:215:LEU:HB2	1.98	0.45
4:F:320:MET:HE3	4:F:320:MET:HB3	1.88	0.45
4:F:340:GLN:HA	4:F:343:TYR:CD2	2.48	0.45
1:A:56:THR:HG23	1:A:58:ALA:H	1.81	0.45
1:A:79:ARG:HH22	1:A:94:THR:HG23	1.81	0.45
1:A:98:ASP:HB2	5:A:501:GTP:O2G	2.16	0.45
1:A:188:ILE:HD12	1:A:425:MET:HG3	1.98	0.45
1:A:254:GLU:HG2	1:A:258:ASN:ND2	2.30	0.45
1:A:226:ASN:ND2	1:A:367:ASP:OD2	2.48	0.45
2:B:306:ARG:HE	10:B:504:MES:H81	1.81	0.45
4:F:243:HIS:HE1	4:F:247:LYS:HD3	1.82	0.45
2:D:267:MET:HG3	2:D:301:ALA:HB3	1.98	0.45
2:B:6:HIS:CD2	2:B:21:TRP:HE1	2.35	0.45
2:B:190:HIS:O	2:B:193:VAL:HG12	2.16	0.45
4:F:275[B]:LEU:HD13	4:F:325:LEU:HD11	1.99	0.45
2:D:50:TYR:OH	2:D:134:GLN:NE2	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:228:LEU:O	2:D:232:THR:HG23	2.16	0.45
4:F:7:ARG:HB2	4:F:42:GLY:HA2	1.99	0.45
2:D:12:CYS:HB2	9:D:502:GDP:C8	2.51	0.44
1:A:18:ASN:ND2	1:A:78:VAL:HG22	2.32	0.44
2:D:139:LEU:HD12	2:D:170:MET:SD	2.57	0.44
1:C:244:PHE:CD1	1:C:358:GLN:HG3	2.52	0.44
2:B:163:ILE:HG21	2:B:250:LEU:HB3	2.00	0.44
2:D:23:VAL:O	2:D:27:GLU:HG3	2.18	0.44
2:D:299:MET:HE3	2:D:305:PRO:HG3	1.98	0.44
1:A:166:LYS:HE2	1:A:197:HIS:O	2.17	0.44
3:E:106:GLU:HG3	3:E:107:SER:N	2.32	0.44
1:A:401:LYS:HG3	2:B:344:TRP:CE3	2.53	0.44
1:C:385:ALA:HA	1:C:388:TRP:CD1	2.52	0.44
2:D:102:ALA:CB	2:D:403:MET:HE3	2.47	0.44
2:B:335:ASN:OD1	4:F:36:ARG:HD3	2.17	0.44
4:F:284:LEU:O	4:F:288:LYS:HG3	2.17	0.44
2:D:81:PHE:O	2:D:84:ILE:HG22	2.17	0.43
2:D:310:TYR:CE1	2:D:367:PHE:HZ	2.36	0.43
4:F:314:LEU:HD22	4:F:350:ILE:HD11	2.00	0.43
2:B:304:ASP:HB3	2:B:307:HIS:ND1	2.32	0.43
2:B:67:ASP:O	2:B:92:PHE:HA	2.19	0.43
2:B:73:MET:HE3	2:B:90:PHE:HD2	1.83	0.43
2:B:152:ILE:HG23	2:B:164:MET:HG2	2.00	0.43
2:B:340:TYR:N	2:B:340:TYR:CD1	2.86	0.43
2:D:317:PHE:HB3	2:D:321:MET:SD	2.59	0.43
1:A:14:VAL:HG13	1:A:67:PHE:HD2	1.82	0.43
2:D:256:ASN:HB3	12:D:504:A1D62:C10	2.48	0.43
4:F:237:THR:HG21	4:F:250:SER:HA	2.00	0.43
1:A:88:HIS:ND1	1:A:90:GLU:HG2	2.33	0.43
1:A:188:ILE:HG23	1:A:425:MET:HG3	2.01	0.43
1:C:140:SER:HA	1:C:171:ILE:HB	2.01	0.43
1:C:141:PHE:O	1:C:147:SER:HB3	2.19	0.43
2:D:235:GLY:O	2:D:238:THR:HG22	2.19	0.43
4:F:217:ARG:NH2	4:F:374:ILE:HG22	2.33	0.43
2:B:322:SER:O	2:B:326:VAL:HG23	2.19	0.43
1:A:401:LYS:NZ	2:B:428:ALA:HB1	2.33	0.43
1:A:30:ILE:HG12	1:A:36:MET:HB2	2.01	0.43
1:C:317:LEU:HD21	1:C:377:MET:HE3	2.00	0.43
2:D:296:SER:HB2	2:D:305:PRO:HD2	2.01	0.43
1:C:368:LEU:HA	1:C:368:LEU:HD23	1.76	0.42
1:C:406:HIS:CG	2:D:261:PRO:HD3	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:81:ILE:HG12	4:F:87:LEU:HD13	2.01	0.42
2:B:104:GLY:O	2:B:109:GLY:HA3	2.19	0.42
1:C:155:GLU:HB3	3:E:101:LEU:HD22	2.00	0.42
2:D:20:PHE:CD1	2:D:233:MET:HE2	2.54	0.42
4:F:246:GLN:HA	4:F:250:SER:HB2	2.01	0.42
2:B:67:ASP:HA	2:B:143:THR:HG21	2.01	0.42
2:D:174:LYS:HE3	2:D:174:LYS:HB2	1.85	0.42
1:A:262:TYR:CE1	1:A:264:ARG:HB2	2.55	0.42
2:B:73:MET:HE2	2:B:92:PHE:HB3	2.02	0.42
2:B:294:PHE:CZ	2:B:330:MET:HE1	2.54	0.42
4:F:149:ALA:HA	4:F:182:ILE:HG22	2.01	0.42
1:A:221:ARG:CZ	2:B:323:MET:HG2	2.48	0.42
1:C:36:MET:HB3	1:C:61:HIS:NE2	2.35	0.42
1:C:107:HIS:HD2	1:C:151[A]:SER:OG	2.03	0.42
1:C:388:TRP:CE3	1:C:425:MET:HE1	2.55	0.42
4:F:37:PHE:CZ	4:F:40:MET:HE3	2.54	0.42
2:B:294:PHE:HZ	2:B:330:MET:HE1	1.85	0.42
2:D:169:VAL:HA	2:D:202:ILE:O	2.20	0.42
1:A:55:GLU:HG2	1:A:61:HIS:CD2	2.55	0.41
2:D:112:LEU:HD23	2:D:147:MET:HE1	2.02	0.41
4:F:79:LYS:HE3	4:F:79:LYS:HB2	1.86	0.41
2:B:19:LYS:HD2	2:B:19:LYS:HA	1.87	0.41
1:C:172:TYR:CE2	1:C:391:LEU:HD22	2.54	0.41
1:A:5:ILE:O	1:A:135:PHE:HA	2.20	0.41
2:D:23:VAL:HG21	2:D:230:SER:HB2	2.01	0.41
2:D:67:ASP:OD1	2:D:68:LEU:N	2.54	0.41
2:B:103:LYS:HA	2:B:107:THR:OG1	2.20	0.41
1:C:271:THR:HG21	1:C:295:CYS:O	2.21	0.41
1:A:370:LYS:H	1:A:370:LYS:HG2	1.57	0.41
2:D:64:ILE:HD13	2:D:120:VAL:HG22	2.03	0.41
1:A:265:ILE:HG23	1:A:432:TYR:CZ	2.56	0.41
1:C:174:ALA:O	1:C:178:SER:N	2.53	0.41
1:C:214:ARG:HG2	1:C:219:ILE:O	2.21	0.41
3:E:76:ARG:NH1	3:E:79:GLU:OE2	2.53	0.41
2:B:165:ASN:ND2	2:B:198:GLU:HG3	2.36	0.41
1:A:172:TYR:HB3	1:A:205:ASP:HA	2.03	0.40
2:B:215:LEU:HD22	2:B:275:SER:HB2	2.02	0.40
2:B:306:ARG:NE	10:B:504:MES:H81	2.36	0.40
2:B:210:ILE:HG23	2:B:273:LEU:HD13	2.02	0.40
10:B:503:MES:H82	10:B:503:MES:H51	1.73	0.40
1:C:233:GLN:HG3	1:C:368:LEU:HD12	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:F:401:ACP:H3B2	14:F:401:ACP:O2A	2.21	0.40
1:C:34:GLY:HA3	1:C:60:LYS:HG3	2.03	0.40
1:C:40:LYS:HA	1:C:40:LYS:HD3	1.92	0.40
2:D:181:GLU:N	2:D:182:PRO:HD2	2.36	0.40
4:F:243:HIS:CE1	4:F:247:LYS:HD3	2.57	0.40
1:A:348:PRO:HG3	3:E:27:PRO:HG3	2.03	0.40
2:D:330:MET:O	2:D:334:GLN:HG3	2.22	0.40
2:B:381:ILE:HG22	2:B:415:MET:HE1	2.04	0.40
2:D:392:LYS:N	2:D:392:LYS:HD2	2.36	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	439/450 (98%)	428 (98%)	11 (2%)	0	100	100
1	C	445/450 (99%)	435 (98%)	9 (2%)	1 (0%)	43	62
2	B	422/445 (95%)	408 (97%)	14 (3%)	0	100	100
2	D	417/445 (94%)	413 (99%)	4 (1%)	0	100	100
3	E	121/143 (85%)	120 (99%)	1 (1%)	0	100	100
4	F	342/384 (89%)	328 (96%)	14 (4%)	0	100	100
All	All	2186/2317 (94%)	2132 (98%)	53 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	131	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	372/378 (98%)	360 (97%)	12 (3%)	34	56
1	C	378/378 (100%)	373 (99%)	5 (1%)	61	75
2	B	369/383 (96%)	362 (98%)	7 (2%)	50	68
2	D	363/383 (95%)	351 (97%)	12 (3%)	33	55
3	E	112/127 (88%)	102 (91%)	10 (9%)	9	17
4	F	316/342 (92%)	300 (95%)	16 (5%)	21	39
All	All	1910/1991 (96%)	1848 (97%)	62 (3%)	34	56

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	THR
1	A	60	LYS
1	A	68	VAL
1	A	124	LYS
1	A	181	VAL
1	A	229	ARG
1	A	324	VAL
1	A	340	SER
1	A	349	THR
1	A	363	VAL
1	A	370	LYS
1	A	423	GLU
2	B	30	ILE
2	B	84	ILE
2	B	130	LEU
2	B	214	THR
2	B	218	THR
2	B	316	ILE
2	B	423	GLN
1	C	26	LEU
1	C	68	VAL
1	C	70	LEU

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Mol	Chain	Res	Type
1	C	335	ILE
1	C	361	THR
2	D	42	LEU
2	D	83	GLN
2	D	84	ILE
2	D	193	VAL
2	D	198	GLU
2	D	205	GLU
2	D	229	VAL
2	D	336	LYS
2	D	343	GLU
2	D	355	ASP
2	D	395	LEU
2	D	420	SER
3	E	9	ILE
3	E	22	VAL
3	E	47	LEU
3	E	92	ASN
3	E	95	LYS
3	E	100	LYS
3	E	106	GLU
3	E	119	MET
3	E	138	GLU
3	E	141	GLU
4	F	12	SER
4	F	18	SER
4	F	46	ARG
4	F	76	SER
4	F	99	VAL
4	F	161	LEU
4	F	162	ILE
4	F	173	ILE
4	F	193	GLU
4	F	200	ASP
4	F	223	THR
4	F	248	GLU
4	F	255[A]	ARG
4	F	255[B]	ARG
4	F	257	GLU
4	F	285	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	HIS
1	A	128	GLN
1	A	283	HIS
1	A	393	HIS
2	B	165	ASN
2	B	334	GLN
2	B	426	GLN
1	C	11	GLN
1	C	18	ASN
1	C	283	HIS
1	C	356	ASN
1	C	393	HIS
2	D	329	GLN
2	D	416	ASN
3	E	64	GLN
4	F	180	HIS
4	F	212	ASN
4	F	379	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 7 are monoatomic - leaving 11 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	GTP	C	501	6	33,34,34	0.99	3 (9%)	50,54,54	1.63	13 (26%)
12	A1D62	B	506	-	33,33,33	0.62	0	43,50,50	1.59	5 (11%)
9	GDP	B	501	6	29,30,30	1.13	2 (6%)	45,47,47	1.68	7 (15%)
10	MES	B	503	-	12,12,12	2.21	1 (8%)	15,16,16	1.85	4 (26%)
10	MES	B	504	-	12,12,12	2.35	1 (8%)	15,16,16	2.48	5 (33%)
12	A1D62	D	504	-	33,33,33	0.64	0	43,50,50	1.64	5 (11%)
13	EDO	C	503	-	3,3,3	0.48	0	2,2,2	0.24	0
5	GTP	A	501	6	33,34,34	0.94	1 (3%)	50,54,54	1.52	8 (16%)
9	GDP	D	502	6	29,30,30	1.15	3 (10%)	45,47,47	1.71	6 (13%)
11	GOL	B	505	-	5,5,5	0.95	0	5,5,5	0.99	0
14	ACP	F	401	-	31,33,33	0.84	2 (6%)	47,52,52	0.81	1 (2%)

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	GTP	C	501	6	-	5/22/38/38	0/3/3/3
12	A1D62	B	506	-	-	2/6/41/41	0/5/5/5
9	GDP	B	501	6	-	4/16/32/32	0/3/3/3
10	MES	B	503	-	-	3/6/14/14	0/1/1/1
10	MES	B	504	-	-	4/6/14/14	0/1/1/1
12	A1D62	D	504	-	-	2/6/41/41	0/5/5/5
13	EDO	C	503	-	-	0/1/1/1	-
5	GTP	A	501	6	-	8/22/38/38	0/3/3/3
9	GDP	D	502	6	-	3/16/32/32	0/3/3/3
11	GOL	B	505	-	-	2/4/4/4	-
14	ACP	F	401	-	-	7/19/38/38	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	B	504	MES	C8-S	-7.84	1.66	1.77
10	B	503	MES	C8-S	-7.40	1.67	1.77
14	F	401	ACP	PB-O3A	3.93	1.62	1.58
9	D	502	GDP	C5-C4	3.14	1.47	1.38
9	B	501	GDP	C5-C4	2.93	1.46	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	B	501	GDP	C6-N1	-2.62	1.33	1.38
9	D	502	GDP	C6-N1	-2.51	1.34	1.38
14	F	401	ACP	PA-O3A	2.27	1.61	1.59
5	C	501	GTP	PB-O3B	2.24	1.61	1.59
5	A	501	GTP	PA-O3A	2.22	1.61	1.59
5	C	501	GTP	PB-O3A	2.19	1.61	1.59
9	D	502	GDP	C5-N7	-2.06	1.34	1.39
5	C	501	GTP	C2-N3	2.03	1.38	1.33

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	D	504	A1D62	O1-C2-C3	-6.09	109.40	121.25
12	B	506	A1D62	O1-C2-C3	-6.02	109.52	121.25
9	D	502	GDP	C5-C4-N3	-5.96	118.91	128.39
9	B	501	GDP	C5-C4-N3	-5.53	119.58	128.39
10	B	504	MES	C5-N4-C3	5.34	120.34	108.84
12	D	504	A1D62	O1-C2-C20	5.27	131.51	121.25
12	B	506	A1D62	O1-C2-C20	5.20	131.37	121.25
14	F	401	ACP	PB-O3A-PA	-5.14	115.59	132.37
9	D	502	GDP	C2-N3-C4	4.88	120.70	112.30
9	B	501	GDP	C2-N3-C4	4.84	120.63	112.30
10	B	503	MES	C5-N4-C3	4.80	119.17	108.84
5	C	501	GTP	C2-N3-C4	4.72	120.43	112.30
5	C	501	GTP	C5-C4-N3	-4.66	120.97	128.39
5	A	501	GTP	C2-N3-C4	4.56	120.16	112.30
5	A	501	GTP	C5-C4-N3	-4.50	121.22	128.39
9	D	502	GDP	N9-C4-N3	4.41	134.77	125.95
9	B	501	GDP	N9-C4-N3	4.26	134.47	125.95
10	B	504	MES	C2-C3-N4	-3.95	104.12	110.12
9	B	501	GDP	C6-C5-N7	3.63	136.90	130.29
10	B	504	MES	C7-N4-C5	3.47	120.50	111.24
9	D	502	GDP	C6-C5-N7	3.29	136.28	130.29
10	B	504	MES	C6-C5-N4	-3.17	105.30	110.12
5	C	501	GTP	N9-C8-N7	-2.97	107.89	113.40
5	A	501	GTP	N9-C8-N7	-2.95	107.92	113.40
10	B	504	MES	C7-N4-C3	2.87	118.90	111.24
5	C	501	GTP	N9-C4-N3	2.83	131.60	125.95
5	C	501	GTP	C8-N7-C5	2.77	109.19	104.26
5	C	501	GTP	C2-N1-C6	-2.74	120.14	125.11
5	A	501	GTP	N9-C4-N3	2.74	131.43	125.95
12	D	504	A1D62	C18-O5-C9	2.67	118.37	113.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	501	GTP	C8-N7-C5	2.63	108.95	104.26
10	B	503	MES	C7-N4-C5	2.57	118.08	111.24
5	C	501	GTP	C5-C6-N1	2.53	119.69	113.25
5	A	501	GTP	C2-N1-C6	-2.53	120.53	125.11
9	D	502	GDP	C4-C5-N7	-2.51	106.70	110.67
12	B	506	A1D62	C18-O5-C9	2.48	118.04	113.67
12	D	504	A1D62	C17-O4-C10	2.47	118.02	113.67
9	B	501	GDP	C4-C5-N7	-2.47	106.76	110.67
5	C	501	GTP	O2B-PB-O3A	2.44	113.88	107.27
12	B	506	A1D62	C17-O4-C10	2.32	117.76	113.67
5	A	501	GTP	C5-C6-N1	2.31	119.14	113.25
5	A	501	GTP	O6-C6-C5	-2.28	120.52	126.53
12	B	506	A1D62	C12-C13-C14	-2.27	106.64	111.61
10	B	503	MES	O2S-S-C8	2.25	110.12	106.73
5	C	501	GTP	O2A-PA-O3A	2.24	113.33	107.27
5	C	501	GTP	O3G-PG-O3B	2.20	112.00	104.64
12	D	504	A1D62	C15-C14-N2	-2.19	100.22	101.79
10	B	503	MES	C6-C5-N4	-2.14	106.86	110.12
5	C	501	GTP	O6-C6-C5	-2.10	120.98	126.53
9	B	501	GDP	C5-C6-N1	2.07	118.53	113.25
9	B	501	GDP	O6-C6-C5	-2.05	121.11	126.53
5	C	501	GTP	C4-C5-N7	-2.05	107.43	110.67
9	D	502	GDP	O6-C6-C5	-2.01	121.22	126.53
5	C	501	GTP	N1-C2-N3	-2.01	119.65	123.32

There are no chirality outliers.

All (40) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	501	GTP	C5'-O5'-PA-O3A
5	A	501	GTP	C5'-O5'-PA-O1A
5	C	501	GTP	C5'-O5'-PA-O3A
5	C	501	GTP	C5'-O5'-PA-O1A
9	B	501	GDP	C5'-O5'-PA-O3A
9	B	501	GDP	C5'-O5'-PA-O1A
9	D	502	GDP	C5'-O5'-PA-O3A
9	D	502	GDP	C5'-O5'-PA-O1A
10	B	503	MES	C8-C7-N4-C5
10	B	504	MES	C7-C8-S-O1S
10	B	504	MES	C7-C8-S-O3S
12	B	506	A1D62	C20-C2-O1-C1
12	D	504	A1D62	C20-C2-O1-C1

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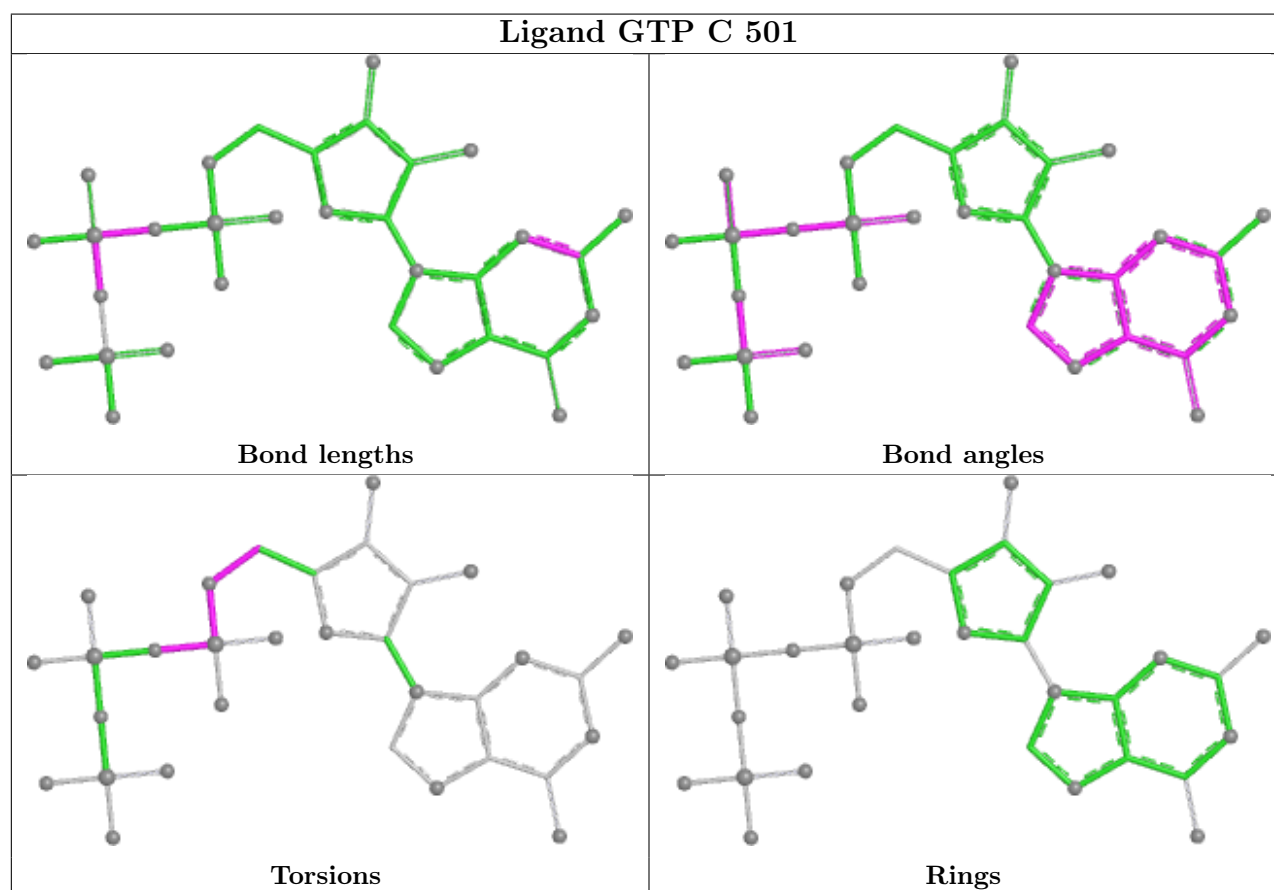
Mol	Chain	Res	Type	Atoms
12	D	504	A1D62	C3-C2-O1-C1
14	F	401	ACP	C5'-O5'-PA-O1A
14	F	401	ACP	C5'-O5'-PA-O2A
14	F	401	ACP	C5'-O5'-PA-O3A
14	F	401	ACP	O4'-C4'-C5'-O5'
14	F	401	ACP	C3'-C4'-C5'-O5'
11	B	505	GOL	O1-C1-C2-C3
12	B	506	A1D62	C3-C2-O1-C1
10	B	503	MES	C8-C7-N4-C3
10	B	504	MES	C8-C7-N4-C3
10	B	504	MES	C7-C8-S-O2S
5	A	501	GTP	PB-O3B-PG-O1G
5	A	501	GTP	C5'-O5'-PA-O2A
5	C	501	GTP	C5'-O5'-PA-O2A
9	B	501	GDP	C5'-O5'-PA-O2A
9	D	502	GDP	C5'-O5'-PA-O2A
14	F	401	ACP	PG-C3B-PB-O2B
5	C	501	GTP	PB-O3A-PA-O2A
11	B	505	GOL	O1-C1-C2-O2
10	B	503	MES	C7-C8-S-O1S
14	F	401	ACP	C4'-C5'-O5'-PA
5	A	501	GTP	PB-O3B-PG-O2G
5	A	501	GTP	PB-O3B-PG-O3G
5	A	501	GTP	PB-O3A-PA-O1A
5	A	501	GTP	PB-O3A-PA-O2A
5	C	501	GTP	C4'-C5'-O5'-PA
9	B	501	GDP	PB-O3A-PA-O1A

There are no ring outliers.

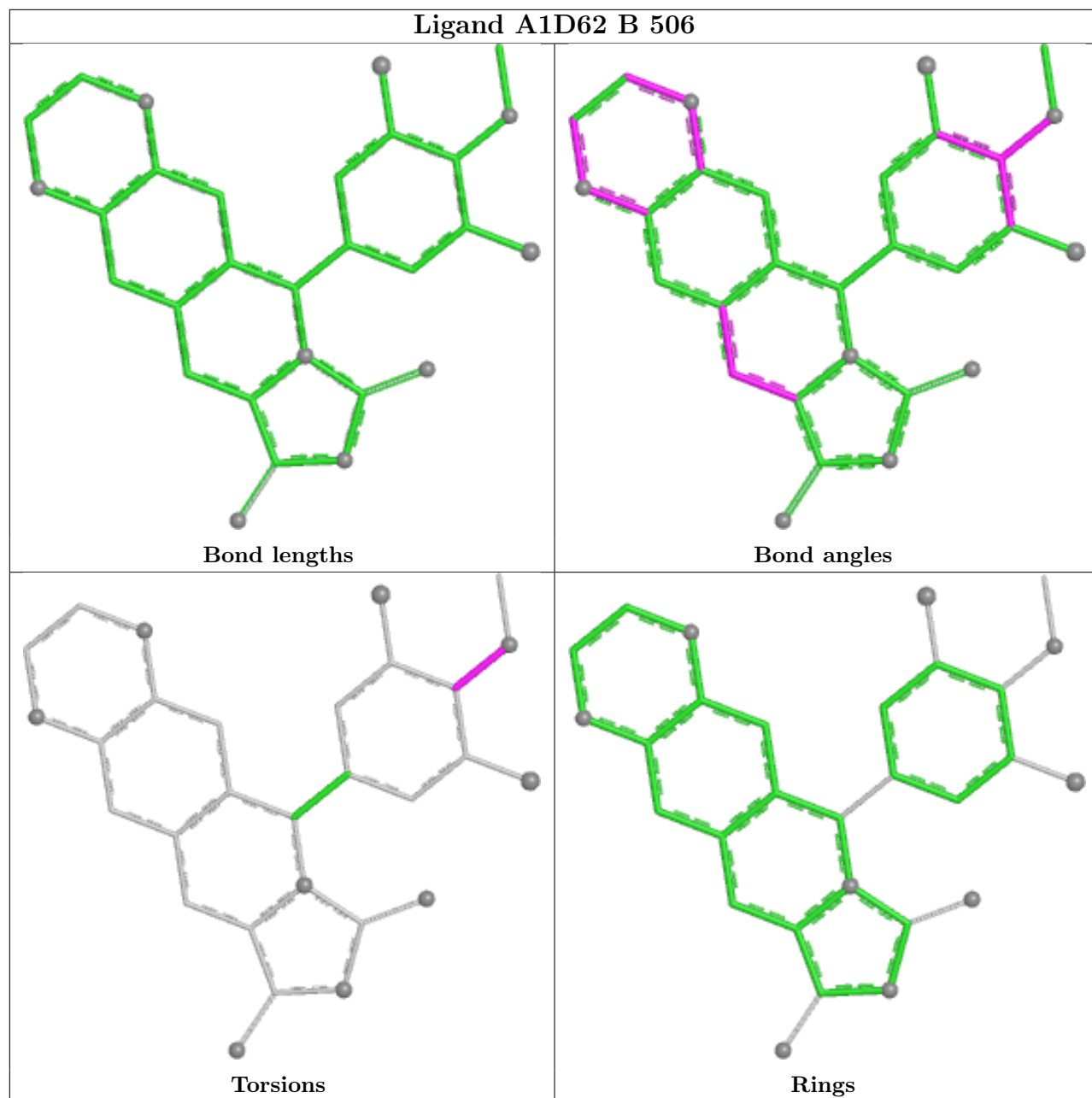
9 monomers are involved in 17 short contacts:

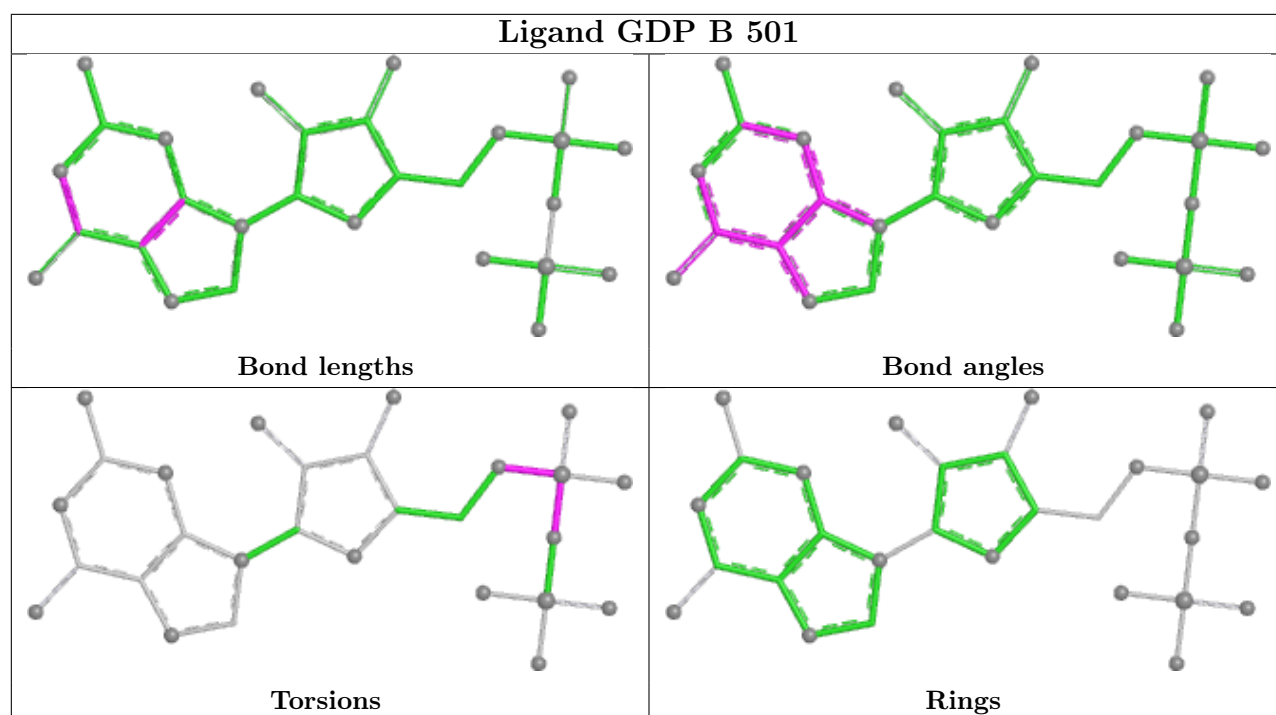
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	501	GTP	1	0
12	B	506	A1D62	1	0
10	B	503	MES	3	0
10	B	504	MES	3	0
12	D	504	A1D62	3	0
5	A	501	GTP	2	0
9	D	502	GDP	2	0
11	B	505	GOL	1	0
14	F	401	ACP	1	0

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.

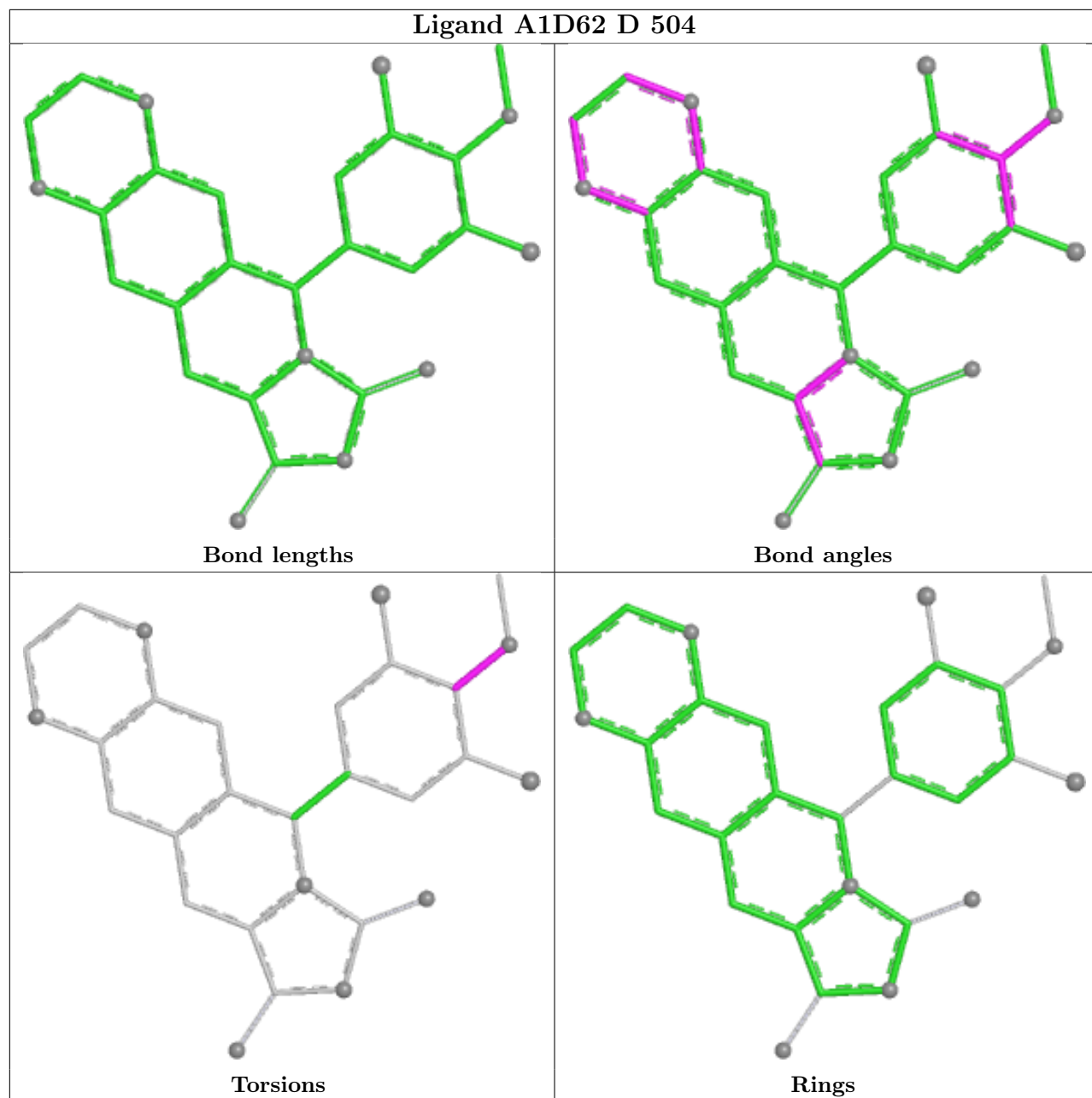


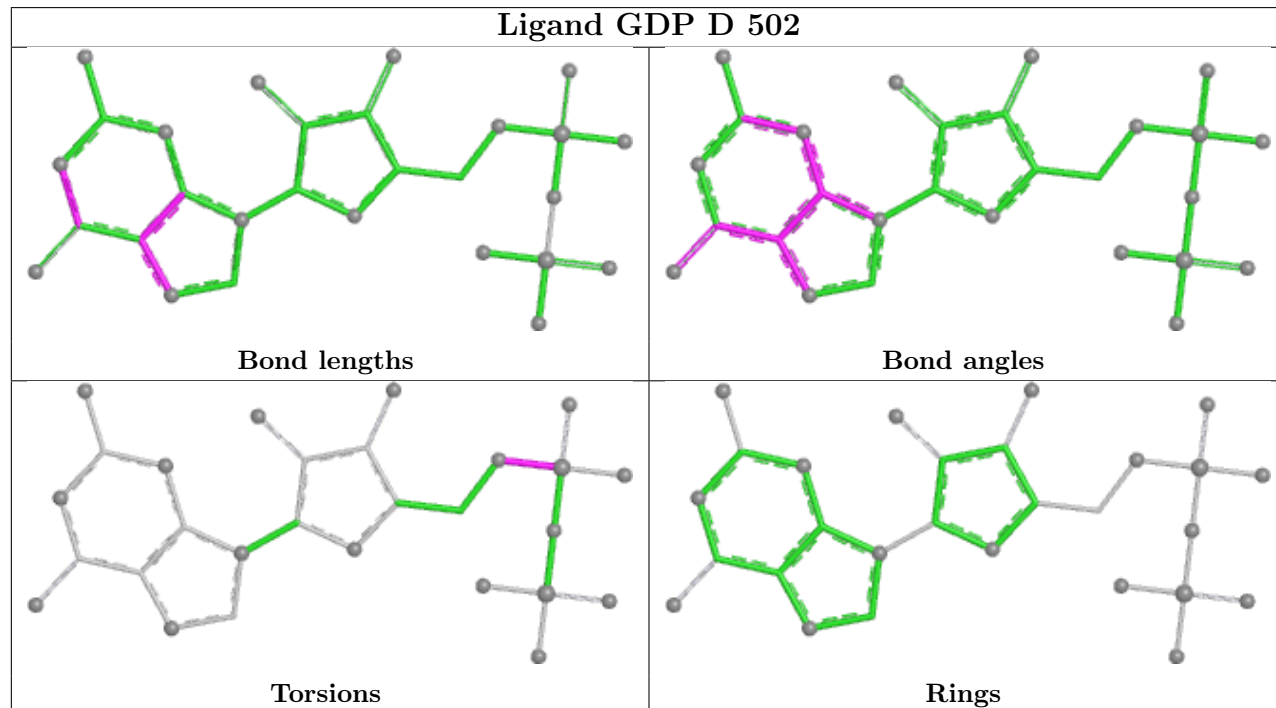
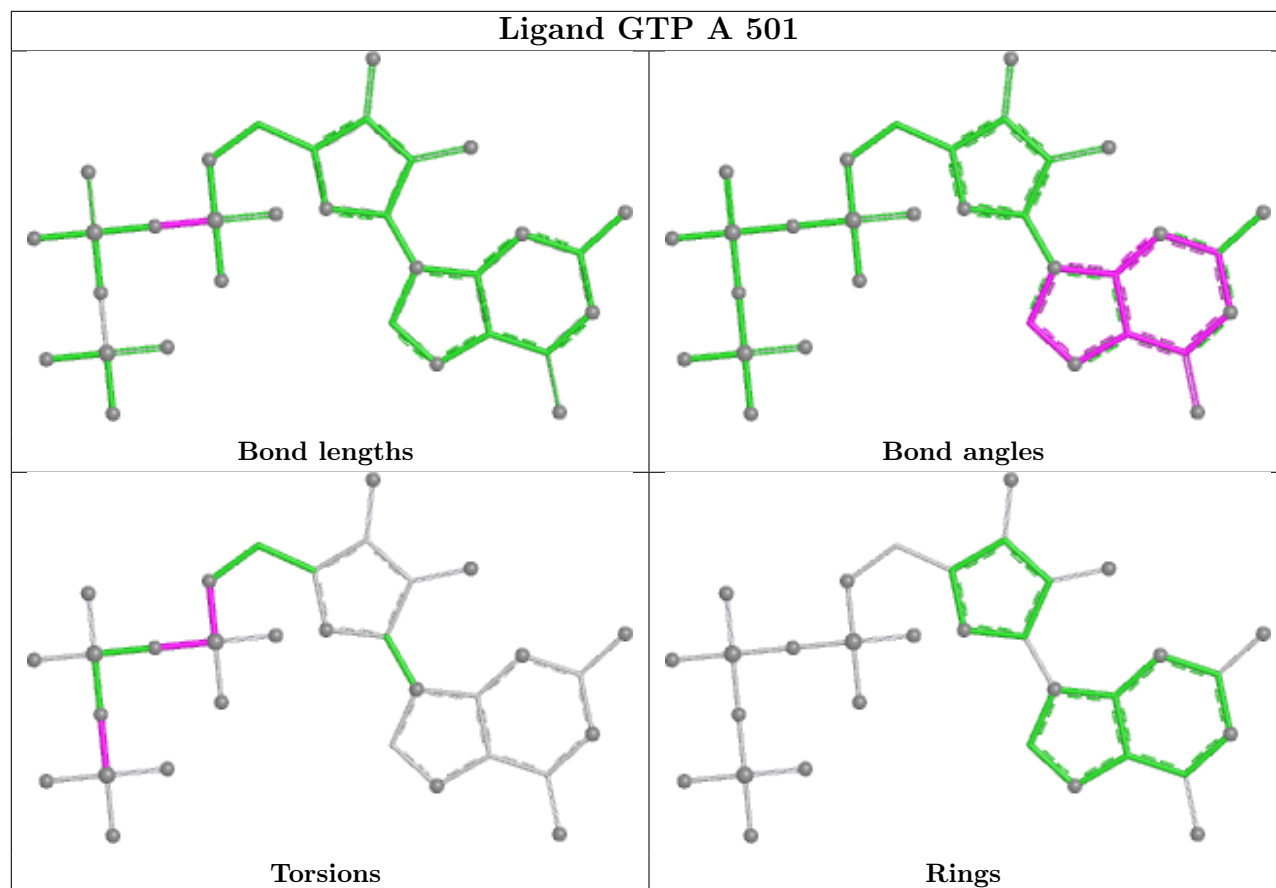
Ligand A1D62 B 506

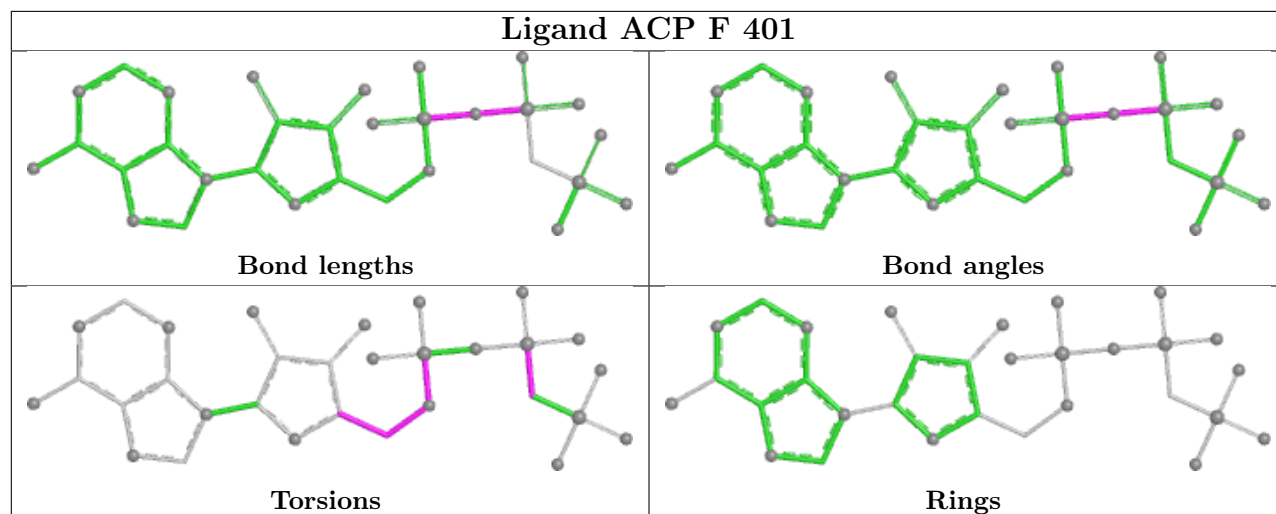




Ligand A1D62 D 504







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(Å²)	Q<0.9
1	A	437/450 (97%)	-0.25	3 (0%)	84	84	37, 61, 96, 134	4 (0%)
1	C	440/450 (97%)	-0.44	3 (0%)	84	84	28, 48, 75, 105	7 (1%)
2	B	424/445 (95%)	-0.25	2 (0%)	87	87	27, 57, 96, 136	2 (0%)
2	D	421/445 (94%)	-0.10	6 (1%)	73	72	42, 70, 110, 150	0
3	E	123/143 (86%)	0.06	1 (0%)	82	82	34, 75, 116, 149	2 (1%)
4	F	346/384 (90%)	0.26	9 (2%)	57	54	37, 91, 154, 175	4 (1%)
All	All	2191/2317 (94%)	-0.16	24 (1%)	78	77	27, 64, 118, 175	19 (0%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	428	ALA	3.6
1	A	437	VAL	3.3
2	D	170	MET	2.8
2	D	92	PHE	2.7
4	F	379	HIS	2.7
1	A	349	THR	2.7
2	D	26	ASP	2.6
2	D	283	ALA	2.6
3	E	24	LEU	2.5
4	F	158	GLU	2.4
1	C	340	SER	2.3
2	D	246	LEU	2.3
4	F	362	ALA	2.3
4	F	159	GLY	2.3
2	B	275	SER	2.2
4	F	186	LEU	2.2
4	F	249	TYR	2.2
1	A	112	LYS	2.1
4	F	233	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
4	F	169	LEU	2.1
1	C	440	VAL	2.1
4	F	167	SER	2.1
1	C	177	VAL	2.1
2	D	247	ASN	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](#)

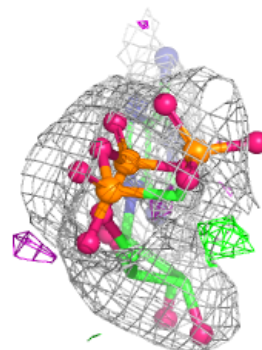
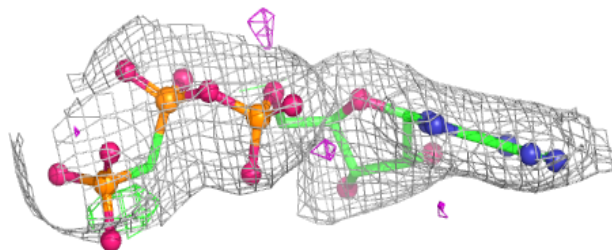
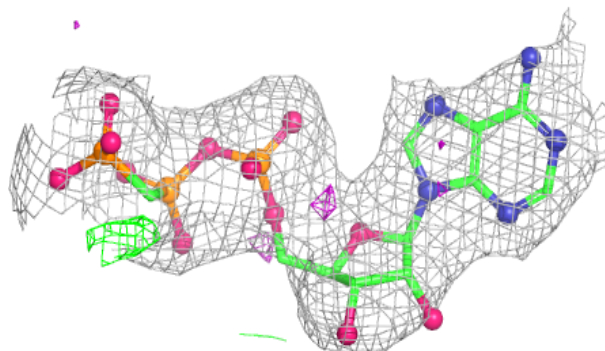
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
11	GOL	B	505	6/6	0.72	0.18	84,94,96,101	0
13	EDO	C	503	4/4	0.84	0.24	61,69,72,75	0
8	CL	A	504	1/1	0.87	0.10	86,86,86,86	0
10	MES	B	504	12/12	0.87	0.12	98,102,126,129	0
14	ACP	F	401	31/31	0.88	0.10	87,100,129,147	0
12	A1D62	D	504	29/29	0.91	0.16	55,81,119,161	0
6	MG	D	503	1/1	0.92	0.12	63,63,63,63	0
12	A1D62	B	506	29/29	0.95	0.12	49,64,96,159	0
9	GDP	D	502	28/28	0.95	0.09	51,64,77,78	0
10	MES	B	503	12/12	0.96	0.07	48,60,72,77	0
5	GTP	A	501	32/32	0.98	0.05	33,46,52,63	0
7	CA	A	503	1/1	0.98	0.03	94,94,94,94	0
5	GTP	C	501	32/32	0.98	0.06	31,38,49,55	0
9	GDP	B	501	28/28	0.98	0.06	31,44,51,55	0
6	MG	D	501	1/1	0.99	0.04	35,35,35,35	0
7	CA	C	502	1/1	0.99	0.03	58,58,58,58	0
6	MG	A	502	1/1	1.00	0.04	46,46,46,46	0
6	MG	B	502	1/1	1.00	0.05	50,50,50,50	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.

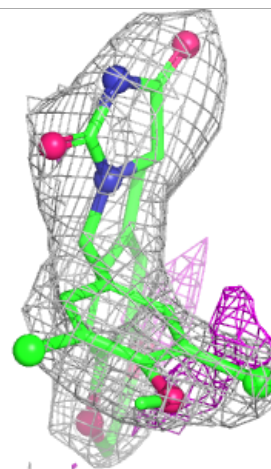
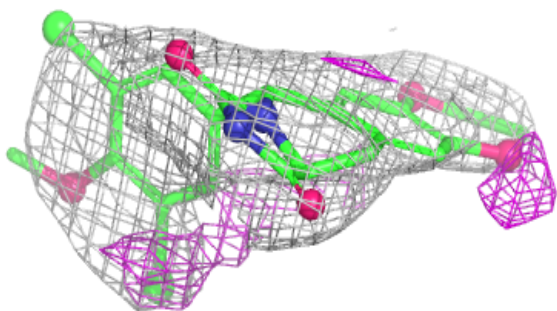
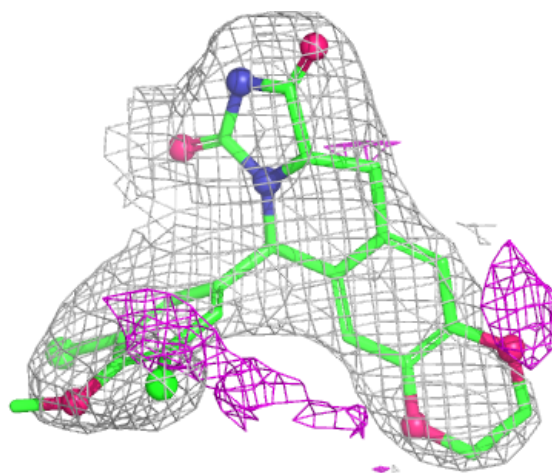
Electron density around ACP F 401:

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



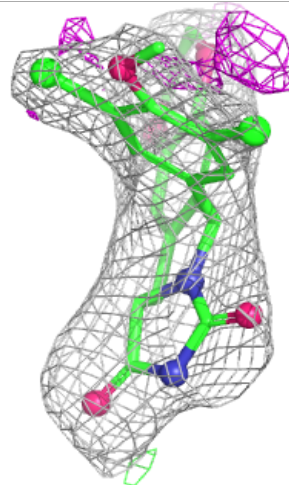
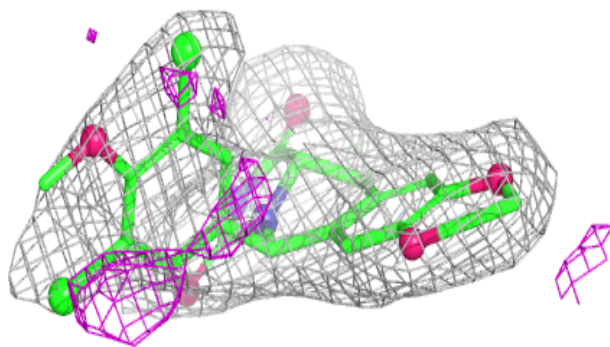
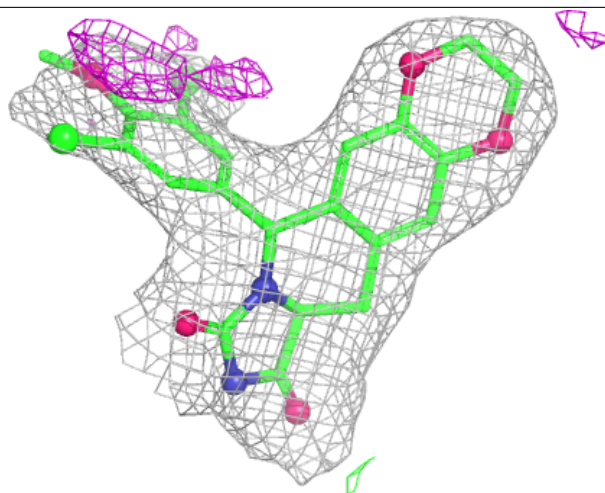
Electron density around A1D62 D 504:

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



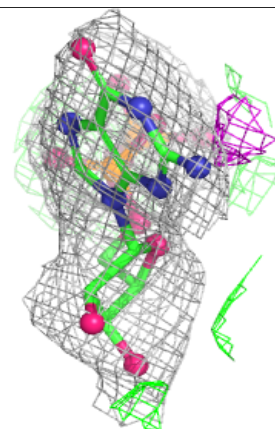
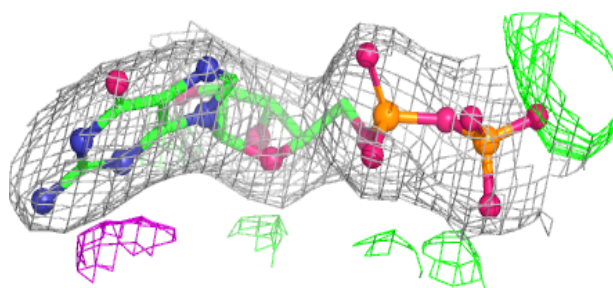
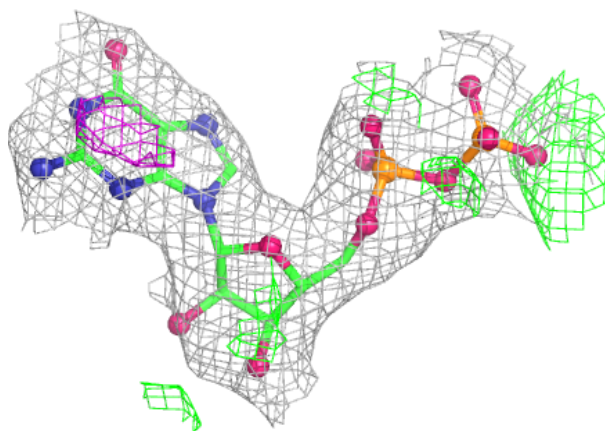
Electron density around A1D62 B 506:

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

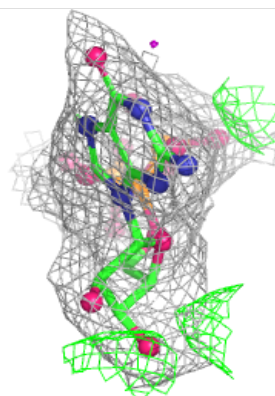
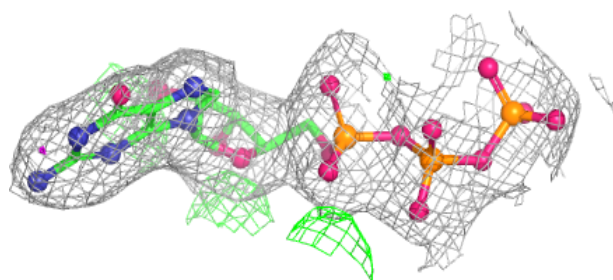
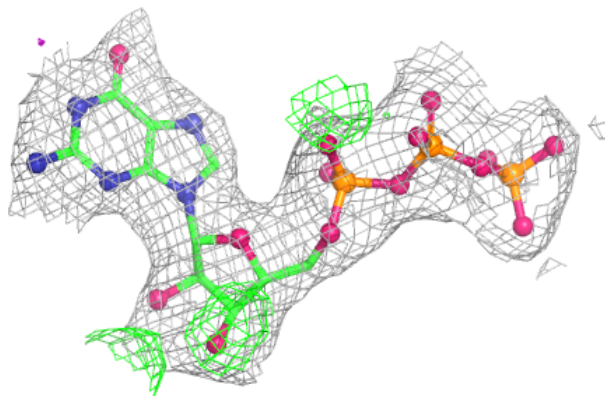


Electron density around GDP D 502:

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and green (positive)

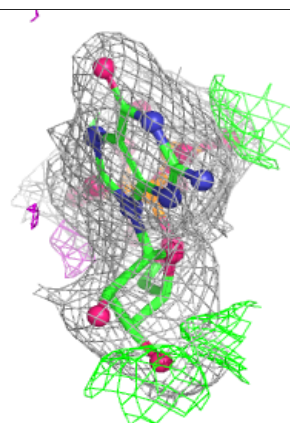
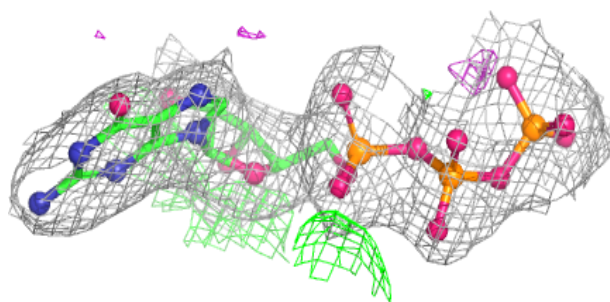
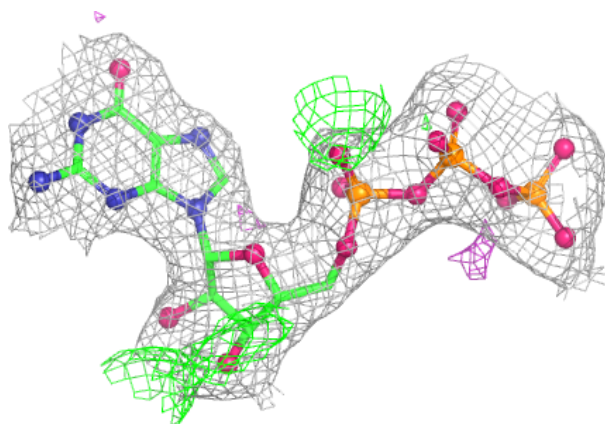
**Electron density around GTP A 501:**

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



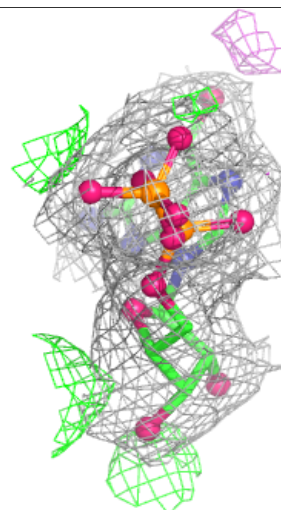
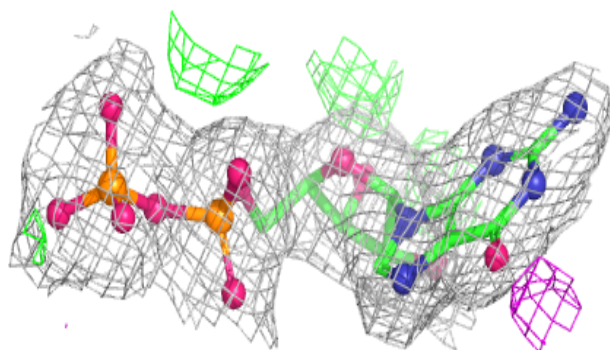
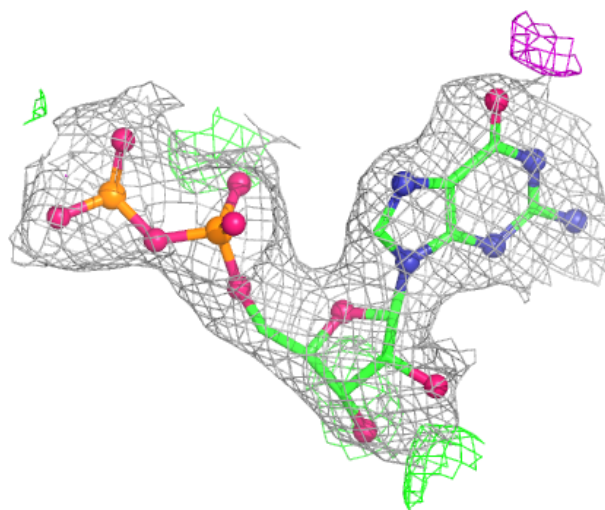
Electron density around GTP C 501:

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



Electron density around GDP B 501:

$2mF_o-DF_c$ (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.