



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 02:38 AM UTC

PDB ID : 4I50 / pdb_00004i50
Title : Crystal structure of tubulin-stathmin-TTL-Epothilone A complex
Authors : Prota, A.E.; Bargsten, K.; Zurwerra, D.; Field, J.J.; Diaz, J.F.; Altmann, K.H.; Steinmetz, M.O.
Deposited on : 2012-11-28
Resolution : 2.30 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

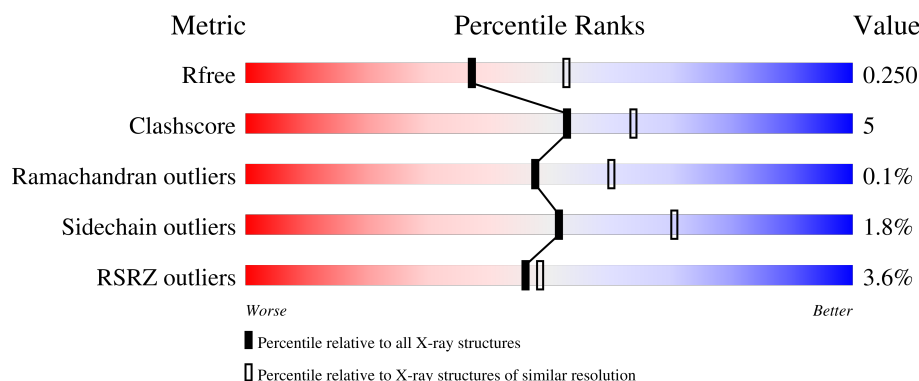
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	451	0% 88% 9% .
1	C	451	2% 87% 10% ..
2	B	445	3% 84% 10% 5%
2	D	445	2% 85% 11% ..
3	E	143	6% 74% 12% 14%

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Mol	Chain	Length	Quality of chain
4	F	384	7% 65% 12% • 21%

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 34573 atoms, of which 16917 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 Tubulin alpha-1B chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	436	Total	C	H	N	O	S	0	0	0
			6740	2155	3333	580	650	22			
1	C	440	Total	C	H	N	O	S	0	10	0
			6890	2201	3412	586	666	25			

- Molecule 2 is a protein called Tubulin beta-2B chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	422	Total	C	H	N	O	S	0	2	0
			6555	2092	3225	569	643	26			
2	D	430	Total	C	H	N	O	S	0	4	0
			6690	2132	3292	582	657	27			

- Molecule 3 is a protein called Stathmin-4.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	E	123	Total	C	H	N	O	S	0	0	0
			2047	625	1033	183	201	5			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	3	MET	ILE	cloning artifact	UNP P63043
E	4	ALA	SER	cloning artifact	UNP P63043

- Molecule 4 is a protein called Tubulin Tyrosine Ligase, TTL.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
4	F	303	Total	C	H	N	O	S	0	0	0
			4964	1597	2481	423	450	13			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	379	HIS	-	expression tag	UNP E1BQ43
F	380	HIS	-	expression tag	UNP E1BQ43
F	381	HIS	-	expression tag	UNP E1BQ43
F	382	HIS	-	expression tag	UNP E1BQ43
F	383	HIS	-	expression tag	UNP E1BQ43
F	384	HIS	-	expression tag	UNP E1BQ43

- Molecule 5 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total	C	H	N	O	P	0	0
			42	10	10	5	14	3		
5	C	1	Total	C	H	N	O	P	0	0
			42	10	10	5	14	3		

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

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

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	F	2	Total	Mg	0	0
			2	2		

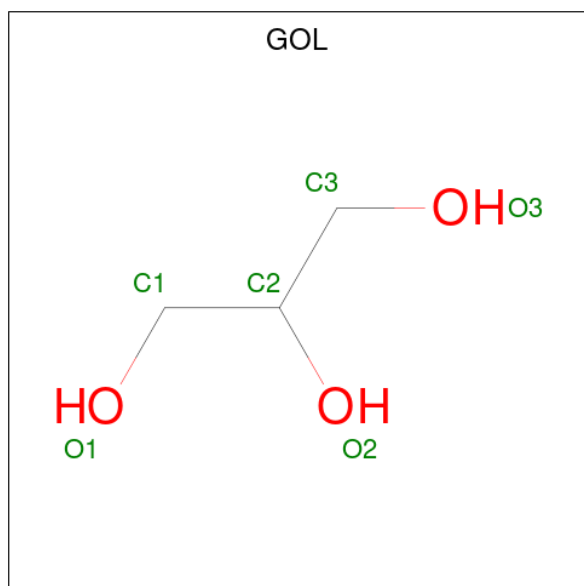
- Molecule 7 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	2	Total	Ca	0	0
			2	2		
7	B	1	Total	Ca	0	0
			1	1		
7	C	3	Total	Ca	0	0
			3	3		

- 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 GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



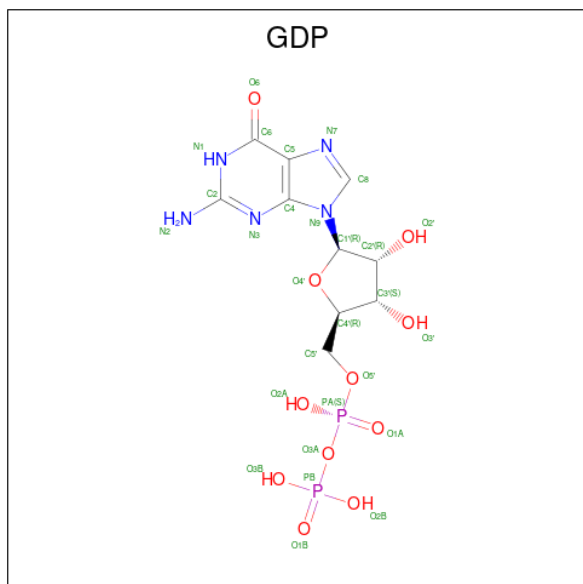
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	A	1	Total	C	H	O	0	0
			14	3	8	3		
9	A	1	Total	C	H	O	0	0
			14	3	8	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	D	1	Total	C	H	O	0	0
			14	3	8	3		

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



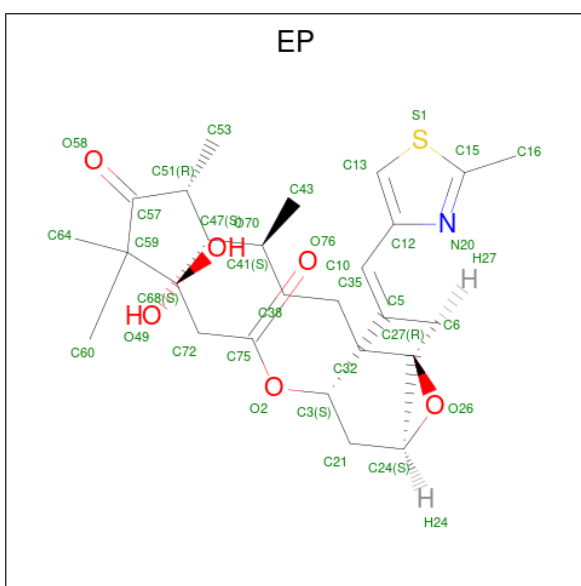
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
10	B	1	Total	C	H	N	O	P	0	0
			38	10	10	5	11	2		
10	D	1	Total	C	H	N	O	P	0	0
			38	10	10	5	11	2		

- Molecule 11 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
11	B	1	Total	C	H	N	O	S	0	0
			25	6	13	1	4	1		
11	D	1	Total	C	H	N	O	S	0	0
			25	6	13	1	4	1		

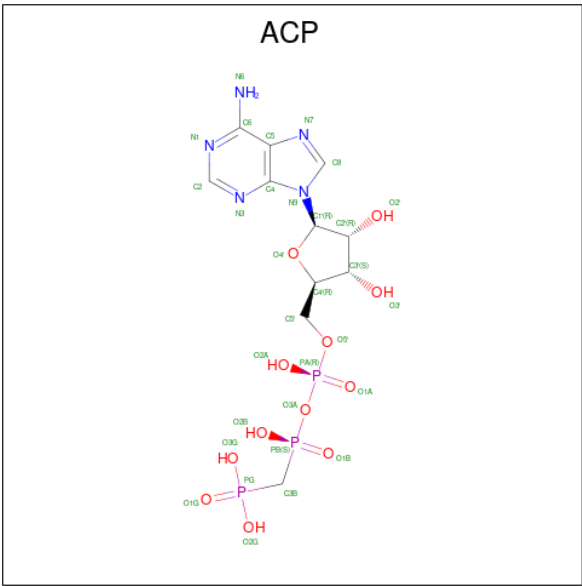
- Molecule 12 is EPOTHILONE A (CCD ID: EP) (formula: $C_{26}H_{39}NO_6S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
12	D	1	Total	C	H	N	O	S	0	0
			73	26	39	1	6	1		

- Molecule 13 is PHOSPHOMETHYLPHOSPHONIC ACID ADENYLATE ESTER (CCD

ID: ACP) (formula: C₁₁H₁₈N₅O₁₂P₃).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
13	F	1	Total	C	H	N	O	P	0	0
			43	11	12	5	12	3		

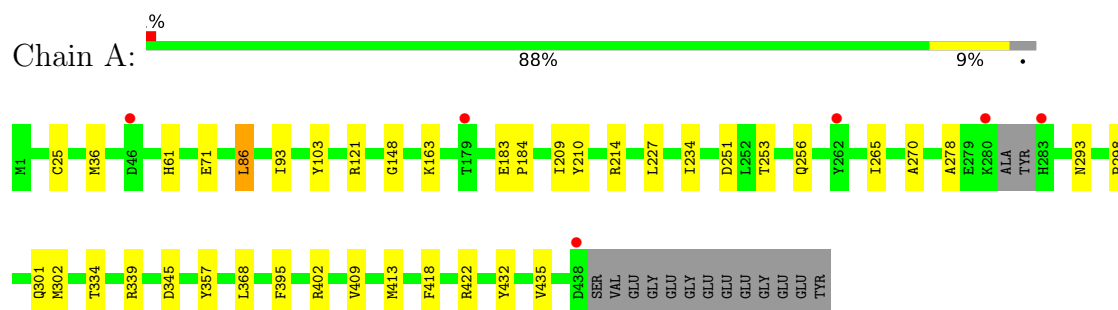
- Molecule 14 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
14	A	63	Total	O	0	0
			63	63		
14	B	50	Total	O	0	0
			50	50		
14	C	117	Total	O	0	0
			117	117		
14	D	40	Total	O	0	0
			40	40		
14	E	16	Total	O	0	0
			16	16		
14	F	20	Total	O	0	0
			20	20		

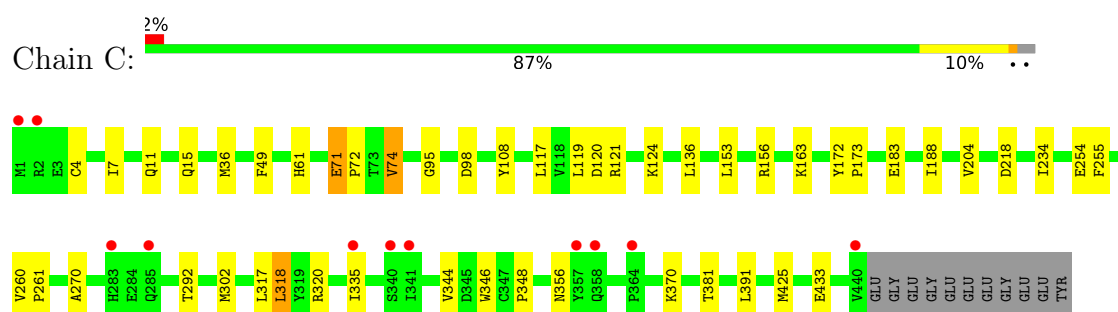
3 Residue-property plots [i](#)

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.

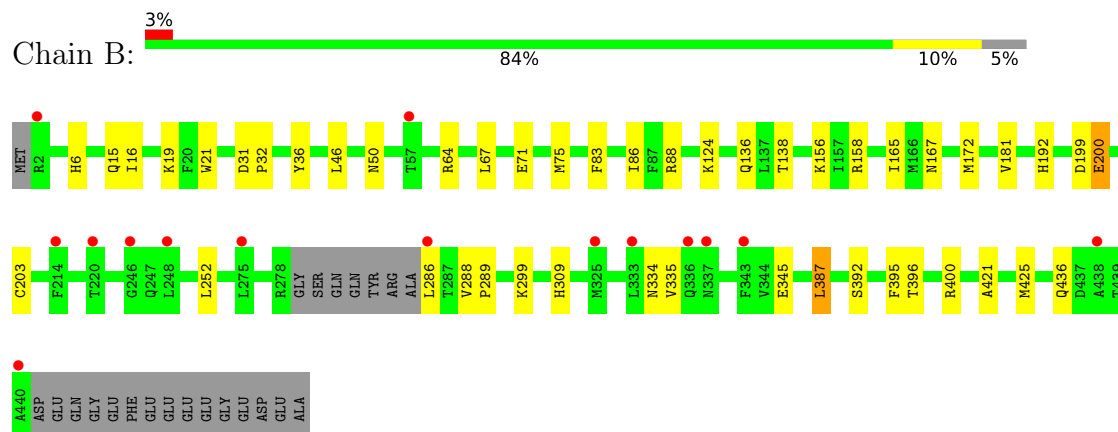
- Molecule 1: Tubulin alpha-1B chain



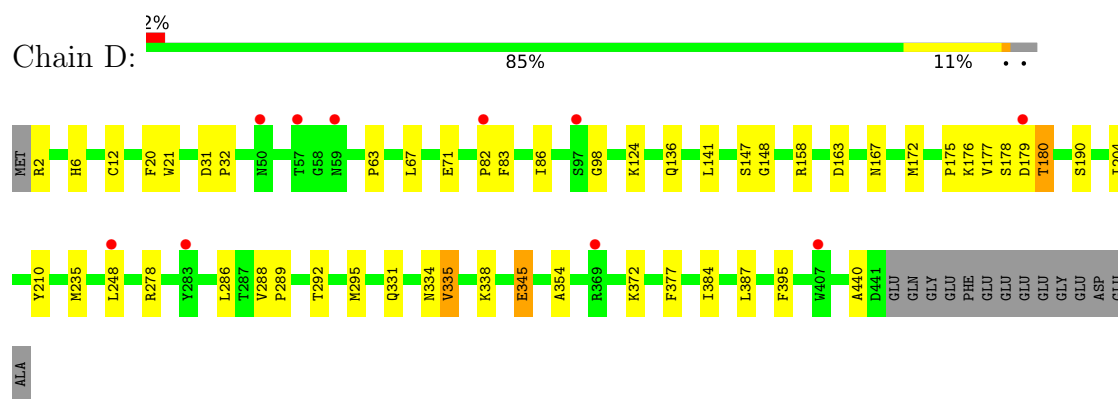
- Molecule 1: Tubulin alpha-1B chain



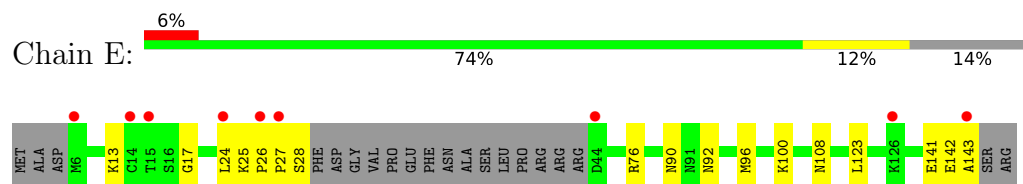
- Molecule 2: Tubulin beta-2B chain



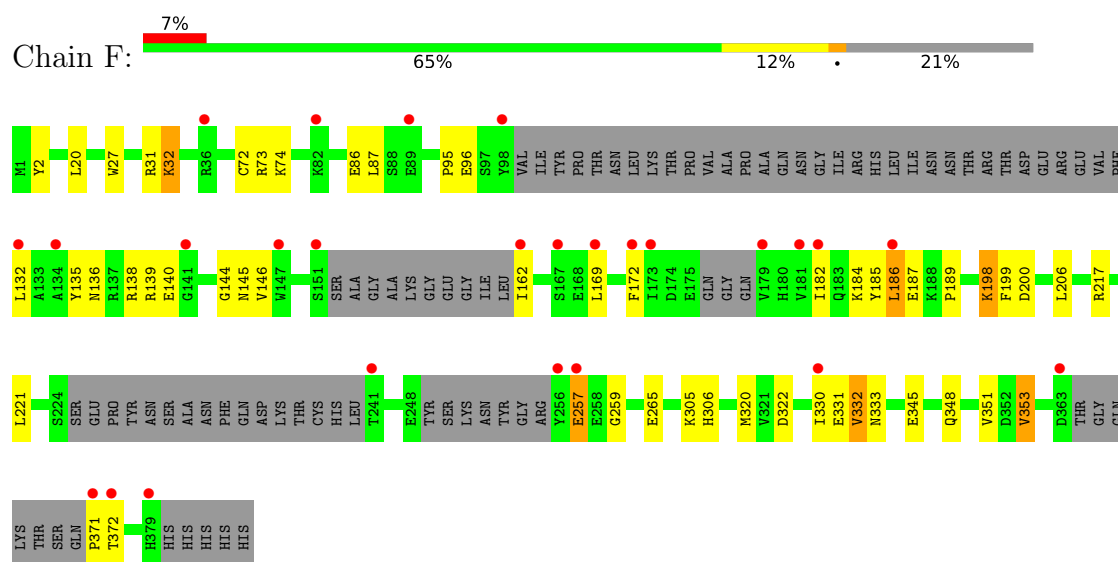
- Molecule 2: Tubulin beta-2B chain



• Molecule 3: Stathmin-4



• Molecule 4: Tubulin Tyrosine Ligase, TTL



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	103.62Å 155.13Å 180.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	77.56 – 2.30 77.56 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (77.56-2.30) 100.0 (77.56-2.30)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 2.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.191 , 0.246 0.197 , 0.250	Depositor DCC
R_{free} test set	6505 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	50.3	Xtriage
Anisotropy	0.261	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 61.6	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	34573	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.36% 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: ACP, CL, CA, MES, GDP, EP, GOL, GTP, 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	0.41	0/3483	0.72	0/4726
1	C	0.46	1/3588 (0.0%)	0.73	0/4871
2	B	0.41	0/3409	0.69	0/4617
2	D	0.39	0/3485	0.70	0/4720
3	E	0.38	0/1022	0.67	0/1356
4	F	0.33	0/2535	0.68	1/3418 (0.0%)
All	All	0.40	1/17522 (0.0%)	0.70	1/23708 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	74	VAL	C-N	-6.60	1.26	1.33

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	259	GLY	N-CA-C	-5.53	107.37	115.72

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	3407	3333	3320	21	0
1	C	3478	3412	3387	30	0
2	B	3330	3225	3215	35	0
2	D	3398	3292	3282	36	0
3	E	1014	1033	1029	12	0
4	F	2483	2481	2472	43	0
5	A	32	10	12	0	0
5	C	32	10	12	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
6	F	2	0	0	0	0
7	A	2	0	0	0	0
7	B	1	0	0	0	0
7	C	3	0	0	0	0
8	A	1	0	0	1	0
9	A	12	16	16	0	0
9	D	6	8	8	0	0
10	B	28	10	12	1	0
10	D	28	10	12	1	0
11	B	12	13	13	2	0
11	D	12	13	13	2	0
12	D	34	39	38	1	0
13	F	31	12	14	2	0
14	A	63	0	0	0	0
14	B	50	0	0	5	0
14	C	117	0	0	1	0
14	D	40	0	0	0	0
14	E	16	0	0	2	0
14	F	20	0	0	0	0
All	All	17656	16917	16855	172	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 5.

All (172) 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:C:234:ILE:HG21	1:C:302[B]:MET:SD	2.25	0.77
1:C:254:GLU:OE1	14:C:717:HOH:O	2.08	0.71
2:D:331:GLN:O	2:D:335:VAL:HG23	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:331:GLU:OE1	4:F:333:ASN:ND2	2.27	0.67
1:C:71:GLU:HB2	1:C:98:ASP:HB3	1.80	0.64
1:C:270:ALA:HB3	1:C:302[B]:MET:SD	2.37	0.63
1:C:4[A]:CYS:SG	1:C:136:LEU:HG	2.39	0.62
1:C:72:PRO:HD3	1:C:95:GLY:O	2.01	0.61
2:B:15:GLN:NE2	10:B:501:GDP:O6	2.32	0.61
2:B:158:ARG:CZ	11:B:504:MES:H21	2.30	0.61
12:D:504:EP:H13	12:D:504:EP:H62	1.82	0.60
2:D:136:GLN:HA	2:D:167:ASN:O	2.02	0.59
2:B:50:ASN:O	2:B:64:ARG:NH2	2.34	0.58
2:B:199:ASP:OD1	11:B:504:MES:H52	2.02	0.58
1:C:255:PHE:CZ	1:C:318:LEU:HD22	2.38	0.58
2:B:387:LEU:HD23	2:B:387:LEU:C	2.28	0.58
1:C:204:VAL:HG22	1:C:302[B]:MET:CE	2.34	0.57
1:C:292:THR:HG22	1:C:335:ILE:CD1	2.35	0.57
2:D:82:PRO:O	2:D:83:PHE:HB2	2.06	0.56
4:F:96:GLU:OE2	4:F:184:LYS:NZ	2.39	0.56
1:A:357:TYR:CZ	3:E:17:GLY:HA2	2.41	0.55
1:A:93:ILE:HD11	1:A:121:ARG:HG3	1.88	0.55
4:F:138:ARG:HD3	4:F:145:ASN:HA	1.87	0.55
2:B:181:VAL:HG23	1:C:348:PRO:CG	2.37	0.54
4:F:206:LEU:HD23	4:F:353:VAL:CG2	2.38	0.54
4:F:371:PRO:HA	4:F:372:THR:HB	1.90	0.54
1:A:293:ASN:OD1	1:A:339:ARG:NH1	2.42	0.53
2:D:163:ASP:HA	11:D:503:MES:C7	2.38	0.53
2:D:387:LEU:C	2:D:387:LEU:HD23	2.33	0.53
2:D:124:LYS:C	2:D:124:LYS:HD3	2.34	0.53
2:D:331:GLN:O	2:D:335:VAL:CG2	2.57	0.53
1:A:36:MET:HB3	1:A:61:HIS:CE1	2.44	0.53
4:F:371:PRO:CA	4:F:372:THR:HB	2.39	0.53
4:F:162:ILE:N	4:F:162:ILE:HD12	2.23	0.52
4:F:217:ARG:NH2	4:F:345:GLU:OE2	2.43	0.52
4:F:200:ASP:OD1	13:F:401:ACP:O3'	2.28	0.52
2:D:141:LEU:HD12	2:D:172:MET:SD	2.50	0.51
2:B:124:LYS:HD3	2:B:124:LYS:C	2.35	0.51
2:D:292:THR:CG2	2:D:335:VAL:HG21	2.41	0.51
2:B:172:MET:HG3	2:B:387:LEU:HD11	1.91	0.51
2:D:158:ARG:HG2	3:E:123:LEU:HD11	1.92	0.51
1:C:317:LEU:C	1:C:318:LEU:HD23	2.36	0.50
2:B:309:HIS:O	2:B:436:GLN:NE2	2.45	0.50
2:D:334:ASN:O	2:D:338:LYS:HG2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:163:ASP:HA	11:D:503:MES:H71	1.93	0.50
1:A:265:ILE:HG23	1:A:432:TYR:CE2	2.47	0.49
3:E:108:ASN:ND2	14:E:206:HOH:O	2.45	0.49
2:D:286:LEU:HD23	2:D:372:LYS:CG	2.41	0.49
2:D:83:PHE:O	2:D:86:ILE:HG22	2.13	0.49
3:E:141:GLU:O	3:E:143:ALA:N	2.43	0.49
4:F:198:LYS:CG	4:F:199:PHE:N	2.75	0.49
2:B:136:GLN:HA	2:B:167:ASN:O	2.12	0.49
4:F:257:GLU:O	4:F:257:GLU:CG	2.60	0.49
4:F:185:TYR:O	4:F:186:LEU:C	2.56	0.49
4:F:320:MET:HG2	4:F:330:ILE:HD11	1.95	0.49
1:A:270:ALA:HB3	1:A:302:MET:HG2	1.95	0.48
2:B:286:LEU:O	2:B:286:LEU:HG	2.13	0.48
2:D:248:LEU:HD23	2:D:354:ALA:HB2	1.95	0.48
2:D:67:LEU:N	2:D:67:LEU:HD12	2.28	0.48
2:D:286:LEU:HD23	2:D:372:LYS:HG2	1.95	0.48
4:F:146:VAL:CG1	4:F:162:ILE:CG2	2.92	0.48
1:A:298:PRO:HA	1:A:301:GLN:CD	2.38	0.48
1:A:345:ASP:OD1	1:A:345:ASP:N	2.46	0.48
2:D:31:ASP:HB2	2:D:32:PRO:CD	2.44	0.48
4:F:138:ARG:CZ	4:F:184:LYS:HE3	2.43	0.48
2:D:12:CYS:HB2	10:D:501:GDP:C8	2.49	0.47
2:D:177:VAL:HG12	2:D:177:VAL:O	2.14	0.47
4:F:144:GLY:HA3	4:F:145:ASN:HB2	1.96	0.47
2:D:345:GLU:HG3	2:D:440:ALA:HB2	1.95	0.47
1:A:210:TYR:CE1	1:A:214:ARG:HD2	2.50	0.47
2:B:6:HIS:CD2	2:B:21:TRP:HE1	2.33	0.47
2:B:192:HIS:CD2	2:B:421:ALA:HA	2.49	0.47
2:D:179:ASP:C	2:D:180:THR:CG2	2.88	0.47
2:B:67:LEU:HD12	2:B:67:LEU:N	2.30	0.47
2:B:395:PHE:CD2	2:B:395:PHE:C	2.92	0.47
4:F:132:LEU:O	4:F:135:TYR:N	2.47	0.47
2:D:295:MET:HE2	2:D:377:PHE:HB2	1.96	0.47
2:D:20:PHE:CD2	2:D:235:MET:HE2	2.50	0.46
4:F:182:ILE:O	4:F:182:ILE:HG23	2.15	0.46
1:C:173:PRO:HB3	1:C:183:GLU:OE2	2.16	0.46
2:B:396:THR:O	2:B:400:ARG:HG3	2.15	0.46
2:B:31:ASP:HB2	2:B:32:PRO:CD	2.46	0.46
1:A:25:CYS:SG	1:A:86:LEU:HD21	2.56	0.46
2:B:334:ASN:CB	14:B:623:HOH:O	2.64	0.46
2:B:165:ILE:HG21	2:B:252:LEU:HB3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:2:TYR:HB3	4:F:27:TRP:CZ3	2.51	0.46
4:F:73:ARG:O	4:F:74:LYS:C	2.58	0.46
1:A:183:GLU:N	1:A:184:PRO:CD	2.79	0.45
2:B:156:LYS:HD3	3:E:76:ARG:CZ	2.45	0.45
1:C:320:ARG:HA	1:C:356:ASN:O	2.16	0.45
2:D:21:TRP:CZ3	2:D:63:PRO:HB3	2.51	0.45
2:D:395:PHE:CD2	2:D:395:PHE:C	2.95	0.45
1:C:11:GLN:O	1:C:15:GLN:HG3	2.17	0.45
1:C:108:TYR:CD2	3:E:108:ASN:CG	2.95	0.45
2:D:147[A]:SER:OG	2:D:148:GLY:N	2.48	0.45
1:C:270:ALA:CB	1:C:302[B]:MET:SD	3.05	0.45
1:C:318:LEU:HD23	1:C:318:LEU:N	2.32	0.45
2:B:334:ASN:HB3	14:B:623:HOH:O	2.17	0.45
1:C:172:TYR:CE2	1:C:391:LEU:HD22	2.52	0.45
2:D:288:VAL:HB	2:D:289:PRO:HD3	1.98	0.45
1:C:163:LYS:HG2	3:E:90:ASN:OD1	2.17	0.45
1:A:209:ILE:HG22	1:A:227:LEU:HD22	1.99	0.44
2:B:71:GLU:HG2	14:B:612:HOH:O	2.16	0.44
2:D:292:THR:CG2	2:D:335:VAL:CG2	2.95	0.44
4:F:136:ASN:O	4:F:140:GLU:N	2.51	0.44
4:F:72:CYS:SG	4:F:332:VAL:HG23	2.57	0.44
1:A:357:TYR:CE1	3:E:17:GLY:HA2	2.53	0.44
4:F:187:GLU:C	4:F:189:PRO:HD3	2.43	0.44
2:D:175:PRO:HA	2:D:178:SER:HB3	1.98	0.44
4:F:138:ARG:HG2	4:F:144:GLY:HA3	2.00	0.43
4:F:257:GLU:O	4:F:257:GLU:HG3	2.18	0.43
2:B:288:VAL:N	2:B:289:PRO:CD	2.81	0.43
2:B:334:ASN:OD1	2:B:334:ASN:C	2.61	0.43
2:B:83:PHE:O	2:B:86:ILE:HG22	2.18	0.43
2:B:299:LYS:N	2:B:299:LYS:HD2	2.34	0.43
2:B:392:SER:HB2	2:B:425:MET:HE3	2.00	0.43
4:F:138:ARG:NH2	4:F:184:LYS:CE	2.82	0.43
1:C:36:MET:HB3	1:C:61:HIS:CE1	2.54	0.43
2:B:75:MET:N	14:B:642:HOH:O	2.47	0.43
1:C:188:ILE:HG13	1:C:425:MET:HG3	1.99	0.43
4:F:206:LEU:HD23	4:F:353:VAL:HG21	2.01	0.43
2:D:176:LYS:HD3	2:D:210:TYR:CD1	2.54	0.43
1:A:409:VAL:HA	1:A:413:MET:O	2.19	0.42
1:A:395:PHE:CD2	1:A:422:ARG:HD3	2.54	0.42
1:C:7:ILE:HG21	1:C:153:LEU:HD21	2.00	0.42
1:C:117:LEU:HD11	1:C:121:ARG:NH1	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:24:LEU:C	3:E:25:LYS:HG3	2.45	0.42
2:B:167:ASN:OD1	2:B:200:GLU:HG3	2.19	0.42
2:B:36:TYR:CD2	2:B:46:LEU:HD11	2.54	0.42
4:F:185:TYR:O	4:F:187:GLU:N	2.52	0.42
4:F:198:LYS:HG3	4:F:199:PHE:N	2.35	0.42
1:C:49:PHE:CD1	1:C:49:PHE:C	2.98	0.42
1:C:120[B]:ASP:OD2	1:C:124:LYS:NZ	2.50	0.42
4:F:138:ARG:NH2	4:F:184:LYS:HE2	2.34	0.42
1:A:278:ALA:HB2	1:A:368:LEU:C	2.45	0.42
4:F:20:LEU:CD2	4:F:348:GLN:OE1	2.68	0.42
4:F:139:ARG:HD2	4:F:139:ARG:HA	1.93	0.42
3:E:26:PRO:O	3:E:28:SER:N	2.52	0.41
1:A:163:LYS:HB2	8:A:504:CL:CL	2.56	0.41
1:C:11:GLN:HG3	1:C:74:VAL:HG21	2.02	0.41
1:A:251:ASP:OD1	1:A:253:THR:HG22	2.21	0.41
2:B:71:GLU:CG	14:B:612:HOH:O	2.68	0.41
1:C:344:VAL:HG21	1:C:346:TRP:CE2	2.56	0.41
3:E:13:LYS:NZ	14:E:214:HOH:O	2.53	0.41
4:F:221:LEU:HD12	4:F:221:LEU:N	2.34	0.41
4:F:172:PHE:CD2	4:F:172:PHE:C	2.98	0.41
4:F:322:ASP:OD1	4:F:322:ASP:C	2.64	0.41
4:F:169:LEU:HA	4:F:172:PHE:HB3	2.02	0.41
2:B:288:VAL:N	2:B:289:PRO:HD2	2.35	0.41
2:D:292:THR:HG21	2:D:335:VAL:CG2	2.50	0.41
3:E:92:ASN:O	3:E:96:MET:HG2	2.20	0.41
1:A:103:TYR:CE1	1:A:148:GLY:HA2	2.56	0.41
1:A:302:MET:N	1:A:302:MET:HE2	2.36	0.41
2:B:16:ILE:HD11	2:B:138:THR:HB	2.02	0.41
2:D:147[B]:SER:OG	2:D:190:SER:OG	2.30	0.41
4:F:138:ARG:CG	4:F:144:GLY:HA3	2.51	0.41
4:F:146:VAL:CG1	4:F:162:ILE:HG22	2.51	0.41
4:F:200:ASP:OD2	4:F:200:ASP:C	2.64	0.41
4:F:348:GLN:O	4:F:351:VAL:CG1	2.69	0.41
2:D:31:ASP:OD1	2:D:31:ASP:C	2.64	0.40
2:B:88:ARG:NH2	2:B:124:LYS:HD2	2.37	0.40
4:F:95:PRO:O	4:F:96:GLU:C	2.62	0.40
4:F:305:LYS:HG3	4:F:306:HIS:N	2.37	0.40
1:C:234:ILE:HD13	1:C:302[A]:MET:SD	2.60	0.40
1:C:260:VAL:HA	1:C:261:PRO:HD2	1.95	0.40
2:D:6:HIS:CD2	2:D:21:TRP:HE1	2.40	0.40
4:F:31:ARG:O	4:F:32:LYS:C	2.64	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:172:MET:CE	2:B:203:CYS:SG	3.09	0.40
2:D:71:GLU:HG3	2:D:98:GLY:CA	2.51	0.40
13:F:401:ACP:O2G	13:F:401:ACP:O2B	2.39	0.40
1:A:413:MET:CE	1:A:418:PHE:CE1	3.04	0.40
1:C:119:LEU:HD11	1:C:156:ARG:HB3	2.03	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	432/451 (96%)	423 (98%)	9 (2%)	0	100	100
1	C	448/451 (99%)	435 (97%)	13 (3%)	0	100	100
2	B	420/445 (94%)	403 (96%)	17 (4%)	0	100	100
2	D	432/445 (97%)	419 (97%)	13 (3%)	0	100	100
3	E	119/143 (83%)	115 (97%)	2 (2%)	2 (2%)	7	7
4	F	289/384 (75%)	271 (94%)	17 (6%)	1 (0%)	36	46
All	All	2140/2319 (92%)	2066 (96%)	71 (3%)	3 (0%)	48	60

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	E	142	GLU
4	F	32	LYS
3	E	27	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	368/379 (97%)	361 (98%)	7 (2%)	50	69
1	C	381/379 (100%)	375 (98%)	6 (2%)	55	73
2	B	367/383 (96%)	362 (99%)	5 (1%)	59	76
2	D	375/383 (98%)	368 (98%)	7 (2%)	50	69
3	E	110/127 (87%)	109 (99%)	1 (1%)	70	84
4	F	273/342 (80%)	265 (97%)	8 (3%)	37	55
All	All	1874/1993 (94%)	1840 (98%)	34 (2%)	51	70

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	GLU
1	A	86	LEU
1	A	234	ILE
1	A	256	GLN
1	A	334	THR
1	A	402	ARG
1	A	435	VAL
2	B	19	LYS
2	B	200	GLU
2	B	335	VAL
2	B	345	GLU
2	B	387	LEU
1	C	71	GLU
1	C	218	ASP
1	C	318	LEU
1	C	370	LYS
1	C	381	THR
1	C	433	GLU
2	D	2	ARG
2	D	180	THR
2	D	204	ILE
2	D	278	ARG

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Mol	Chain	Res	Type
2	D	335	VAL
2	D	345	GLU
2	D	384	ILE
3	E	100	LYS
4	F	86	GLU
4	F	87	LEU
4	F	186	LEU
4	F	198	LYS
4	F	257	GLU
4	F	265	GLU
4	F	332	VAL
4	F	353	VAL

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

Mol	Chain	Res	Type
1	A	101	ASN
1	A	393	HIS
2	B	349	ASN
2	B	385	GLN
1	C	15	GLN
1	C	133	GLN
1	C	216	ASN
1	C	393	HIS
2	D	8	GLN
2	D	50	ASN
2	D	229	HIS
2	D	436	GLN
3	E	12	ASN
3	E	108	ASN
4	F	269	GLN
4	F	310	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 24 ligands modelled in this entry, 13 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
9	GOL	D	505	-	5,5,5	0.36	0	5,5,5	0.18	0
10	GDP	B	501	6	29,30,30	1.18	2 (6%)	45,47,47	1.83	7 (15%)
5	GTP	A	501	6	33,34,34	1.03	3 (9%)	50,54,54	1.56	10 (20%)
12	EP	D	504	-	36,36,36	1.39	6 (16%)	48,53,53	1.64	9 (18%)
11	MES	B	504	-	12,12,12	2.37	1 (8%)	15,16,16	1.35	2 (13%)
11	MES	D	503	-	12,12,12	2.42	1 (8%)	15,16,16	1.42	3 (20%)
5	GTP	C	501	6	33,34,34	1.19	4 (12%)	50,54,54	1.53	9 (18%)
9	GOL	A	507	-	5,5,5	0.38	0	5,5,5	0.30	0
13	ACP	F	401	6	31,33,33	1.49	4 (12%)	47,52,52	2.33	12 (25%)
9	GOL	A	506	-	5,5,5	0.33	0	5,5,5	0.38	0
10	GDP	D	501	6	29,30,30	1.23	4 (13%)	45,47,47	1.63	5 (11%)

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
9	GOL	D	505	-	-	2/4/4/4	-
10	GDP	B	501	6	-	5/16/32/32	0/3/3/3
5	GTP	A	501	6	-	6/22/38/38	0/3/3/3
12	EP	D	504	-	-	4/50/55/55	0/3/3/3
11	MES	B	504	-	-	1/6/14/14	0/1/1/1
11	MES	D	503	-	-	3/6/14/14	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GTP	C	501	6	-	9/22/38/38	0/3/3/3
9	GOL	A	507	-	-	2/4/4/4	-
13	ACP	F	401	6	-	2/19/38/38	0/3/3/3
9	GOL	A	506	-	-	0/4/4/4	-
10	GDP	D	501	6	-	5/16/32/32	0/3/3/3

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	D	503	MES	C8-S	-8.16	1.66	1.77
11	B	504	MES	C8-S	-7.93	1.66	1.77
13	F	401	ACP	C6-N6	4.17	1.44	1.34
12	D	504	EP	C27-C24	3.84	1.53	1.46
10	D	501	GDP	C5-C4	3.06	1.47	1.38
5	C	501	GTP	PB-O3A	3.02	1.62	1.59
5	C	501	GTP	PA-O3A	3.02	1.62	1.59
12	D	504	EP	C12-N20	2.99	1.44	1.38
12	D	504	EP	C72-C68	2.96	1.57	1.53
13	F	401	ACP	C5-C4	2.94	1.44	1.39
13	F	401	ACP	PB-O1B	2.92	1.58	1.51
10	B	501	GDP	C5-C4	2.81	1.46	1.38
10	D	501	GDP	C6-N1	-2.76	1.33	1.38
10	B	501	GDP	C6-N1	-2.72	1.33	1.38
5	C	501	GTP	PB-O3B	2.71	1.62	1.59
5	A	501	GTP	PA-O3A	2.48	1.62	1.59
5	A	501	GTP	PB-O3B	2.39	1.62	1.59
13	F	401	ACP	C4-N9	2.19	1.42	1.37
12	D	504	EP	C21-C3	2.18	1.56	1.53
12	D	504	EP	C10-C12	2.12	1.49	1.41
10	D	501	GDP	PA-O3A	2.11	1.61	1.59
10	D	501	GDP	C4-N9	-2.11	1.32	1.38
5	C	501	GTP	C2-N3	2.08	1.38	1.33
12	D	504	EP	C51-C47	2.03	1.57	1.53
5	A	501	GTP	C2-N3	2.00	1.38	1.33

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	F	401	ACP	O2B-PB-O1B	7.66	134.89	109.95
12	D	504	EP	C27-O26-C24	5.99	64.99	60.67
13	F	401	ACP	O1B-PB-C3B	-5.87	93.60	109.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	B	501	GDP	C5-C4-N3	-5.80	119.15	128.39
13	F	401	ACP	C5-C4-N3	-5.40	119.29	126.72
10	D	501	GDP	C5-C4-N3	-5.22	120.08	128.39
10	B	501	GDP	C2-N3-C4	4.93	120.79	112.30
5	A	501	GTP	C2-N3-C4	4.79	120.55	112.30
5	C	501	GTP	C5-C4-N3	-4.70	120.92	128.39
5	A	501	GTP	C5-C4-N3	-4.66	120.98	128.39
10	D	501	GDP	C2-N3-C4	4.43	119.93	112.30
13	F	401	ACP	N3-C4-N9	4.31	134.50	127.17
5	C	501	GTP	C2-N3-C4	4.17	119.48	112.30
10	B	501	GDP	N9-C4-N3	4.04	134.04	125.95
10	D	501	GDP	C6-C5-N7	3.99	137.55	130.29
12	D	504	EP	C6-C5-C10	-3.98	119.20	124.96
11	B	504	MES	C5-N4-C3	3.97	117.39	108.84
13	F	401	ACP	N3-C2-N1	-3.95	122.60	128.58
13	F	401	ACP	C5-N7-C8	3.84	109.49	103.45
10	B	501	GDP	C6-C5-N7	3.81	137.22	130.29
10	D	501	GDP	N9-C4-N3	3.71	133.37	125.95
13	F	401	ACP	C2-N3-C4	3.46	120.27	111.83
5	C	501	GTP	N9-C8-N7	-3.13	107.60	113.40
5	C	501	GTP	N9-C4-N3	3.12	132.20	125.95
12	D	504	EP	O26-C24-C27	-3.09	57.47	59.66
13	F	401	ACP	C4-C5-N7	-3.03	107.12	110.58
11	D	503	MES	C5-N4-C3	3.00	115.31	108.84
12	D	504	EP	O26-C27-C24	-3.00	57.54	59.66
10	B	501	GDP	O2A-PA-O3A	2.95	115.25	107.27
10	B	501	GDP	C4-C5-N7	-2.93	106.03	110.67
5	C	501	GTP	C8-N7-C5	2.85	109.34	104.26
13	F	401	ACP	PB-O3A-PA	-2.85	123.08	132.37
10	D	501	GDP	C4-C5-N7	-2.84	106.17	110.67
13	F	401	ACP	N9-C8-N7	-2.82	109.93	113.94
5	A	501	GTP	N9-C8-N7	-2.81	108.19	113.40
11	D	503	MES	O1S-S-C8	2.76	110.90	106.73
12	D	504	EP	C53-C51-C47	-2.64	107.76	112.42
5	A	501	GTP	C2-N1-C6	-2.64	120.33	125.11
5	C	501	GTP	C2-N1-C6	-2.64	120.33	125.11
5	A	501	GTP	C8-N7-C5	2.62	108.92	104.26
12	D	504	EP	O2-C75-C72	2.61	116.11	111.43
5	A	501	GTP	C5-C6-N1	2.52	119.68	113.25
5	A	501	GTP	N9-C4-N3	2.49	130.94	125.95
12	D	504	EP	C16-C15-N20	2.48	128.97	123.74
5	A	501	GTP	O2B-PB-O3A	2.46	113.91	107.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	501	GTP	O2A-PA-O3A	2.42	113.80	107.27
13	F	401	ACP	O2A-PA-O1A	-2.40	101.28	112.44
5	C	501	GTP	C5-C6-N1	2.31	119.13	113.25
10	B	501	GDP	O3'-C3'-C4'	-2.30	104.48	111.08
11	D	503	MES	C6-C5-N4	-2.22	106.75	110.12
5	A	501	GTP	C4-C5-N7	-2.21	107.16	110.67
12	D	504	EP	C16-C15-S1	-2.18	116.14	121.91
12	D	504	EP	C3-O2-C75	2.16	119.49	116.25
5	A	501	GTP	C6-C5-N7	2.08	134.07	130.29
11	B	504	MES	O2S-S-C8	2.04	109.82	106.73
13	F	401	ACP	O2B-PB-C3B	-2.03	98.33	106.73
5	C	501	GTP	O6-C6-C5	-2.02	121.19	126.53

There are no chirality outliers.

All (39) 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	A	501	GTP	C5'-O5'-PA-O2A
5	C	501	GTP	C5'-O5'-PA-O3A
5	C	501	GTP	C5'-O5'-PA-O1A
5	C	501	GTP	C5'-O5'-PA-O2A
9	A	507	GOL	O1-C1-C2-C3
10	B	501	GDP	C5'-O5'-PA-O3A
10	B	501	GDP	C5'-O5'-PA-O1A
10	B	501	GDP	C5'-O5'-PA-O2A
10	D	501	GDP	C5'-O5'-PA-O3A
10	D	501	GDP	C5'-O5'-PA-O1A
11	B	504	MES	C8-C7-N4-C5
11	D	503	MES	C8-C7-N4-C5
11	D	503	MES	N4-C7-C8-S
12	D	504	EP	C12-C10-C5-C6
12	D	504	EP	C21-C3-C5-C10
13	F	401	ACP	O4'-C4'-C5'-O5'
13	F	401	ACP	C3'-C4'-C5'-O5'
9	D	505	GOL	O1-C1-C2-C3
9	A	507	GOL	O1-C1-C2-O2
11	D	503	MES	C8-C7-N4-C3
9	D	505	GOL	O1-C1-C2-O2
5	C	501	GTP	PB-O3B-PG-O1G
5	A	501	GTP	PB-O3B-PG-O2G

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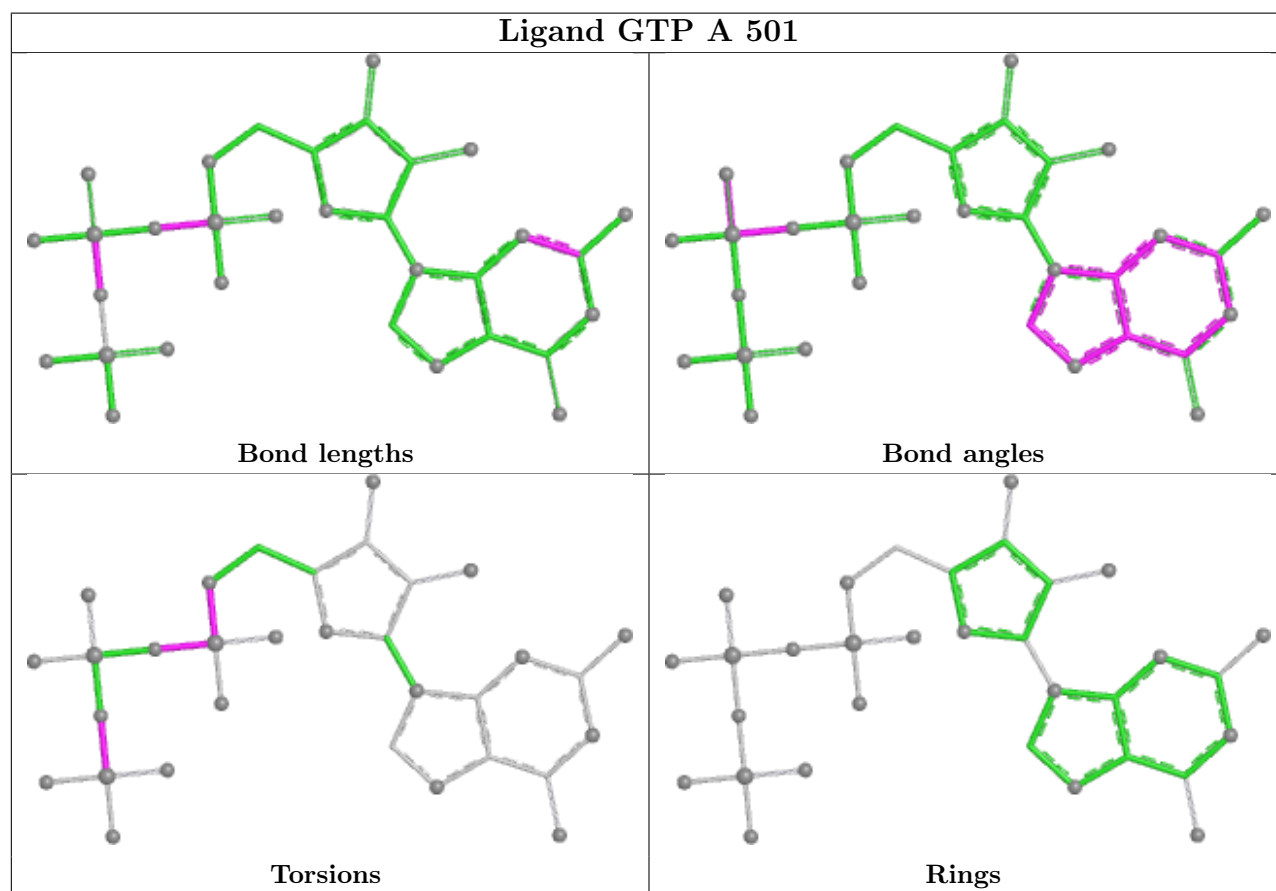
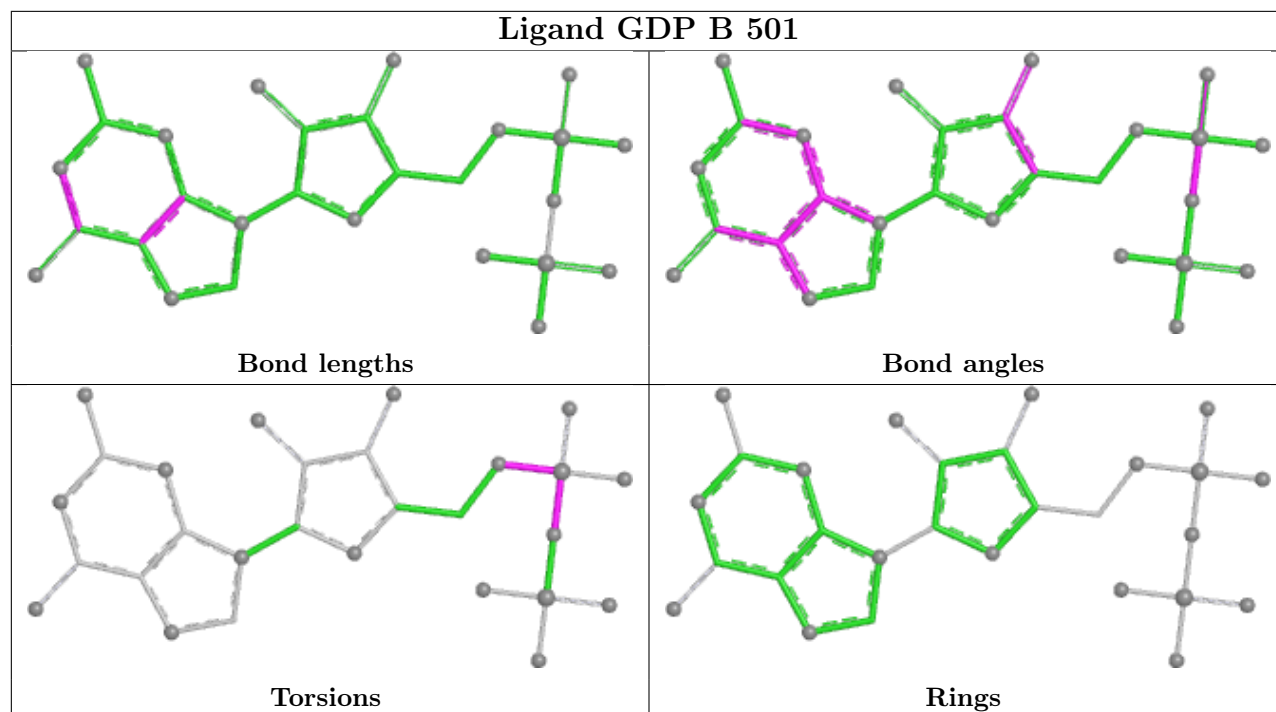
Mol	Chain	Res	Type	Atoms
10	D	501	GDP	C5'-O5'-PA-O2A
12	D	504	EP	O2-C3-C5-C10
10	D	501	GDP	PB-O3A-PA-O2A
5	C	501	GTP	PB-O3A-PA-O2A
10	B	501	GDP	PB-O3A-PA-O1A
10	B	501	GDP	PB-O3A-PA-O2A
5	C	501	GTP	C4'-C5'-O5'-PA
5	A	501	GTP	PB-O3B-PG-O3G
5	C	501	GTP	PB-O3B-PG-O2G
5	C	501	GTP	PB-O3B-PG-O3G
12	D	504	EP	C21-C3-C5-C6
5	C	501	GTP	PB-O3A-PA-O1A
5	A	501	GTP	PB-O3A-PA-O2A
10	D	501	GDP	PB-O3A-PA-O1A

There are no ring outliers.

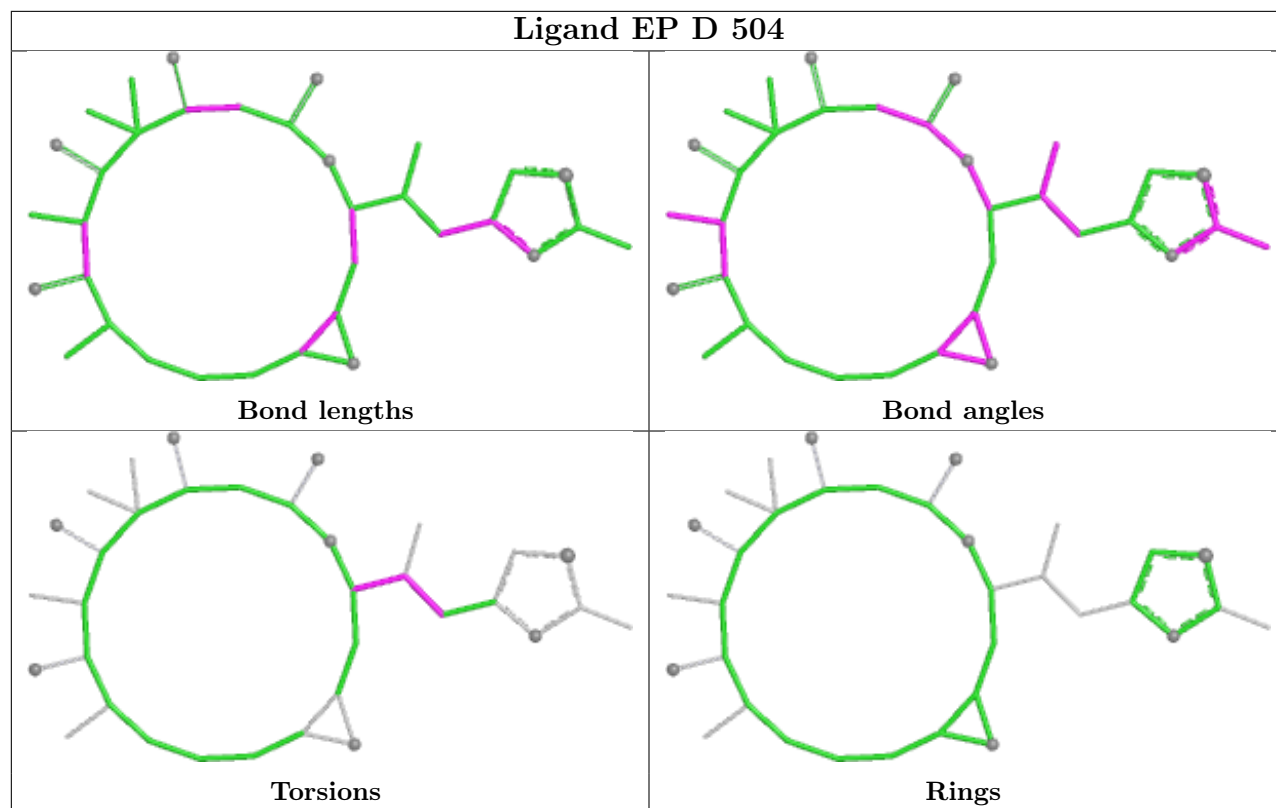
6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	B	501	GDP	1	0
12	D	504	EP	1	0
11	B	504	MES	2	0
11	D	503	MES	2	0
13	F	401	ACP	2	0
10	D	501	GDP	1	0

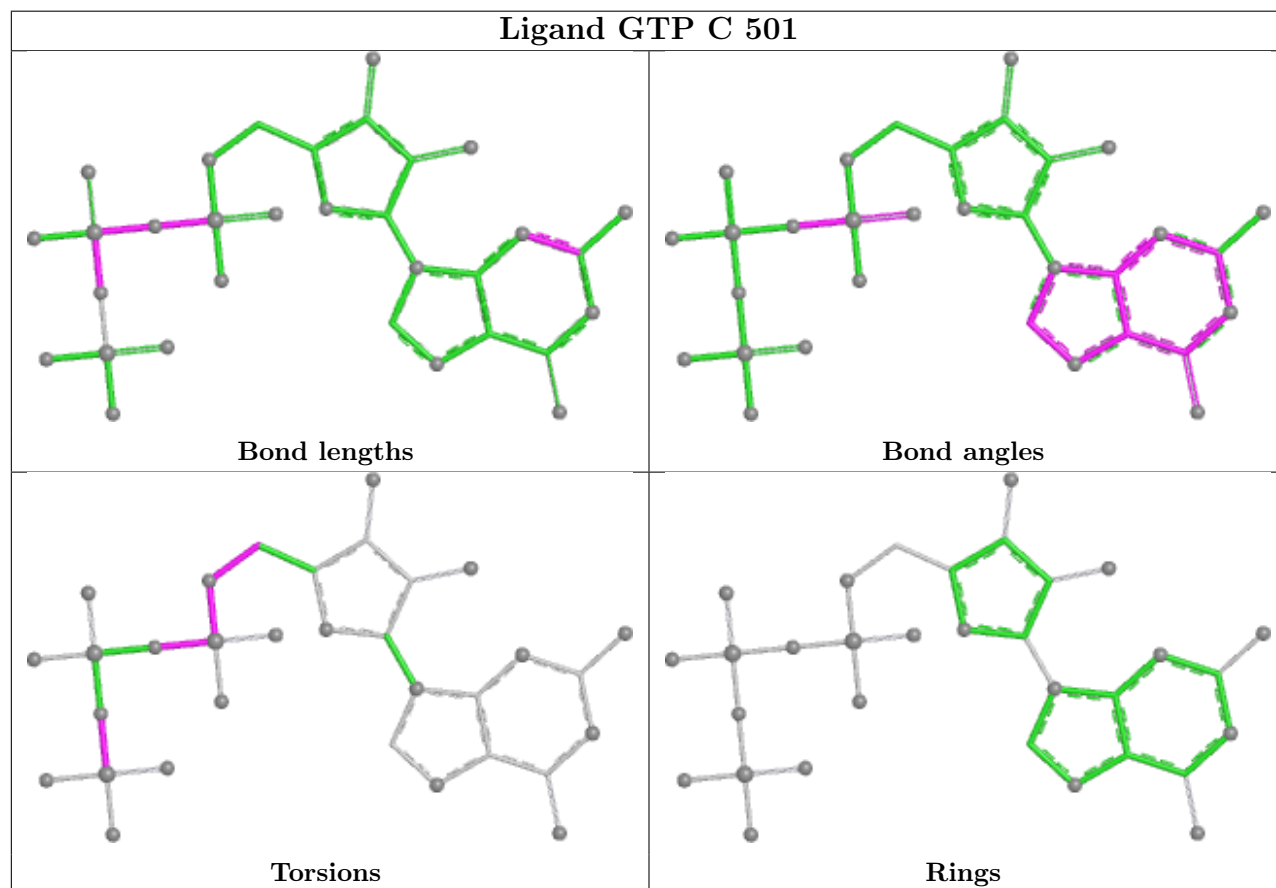
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

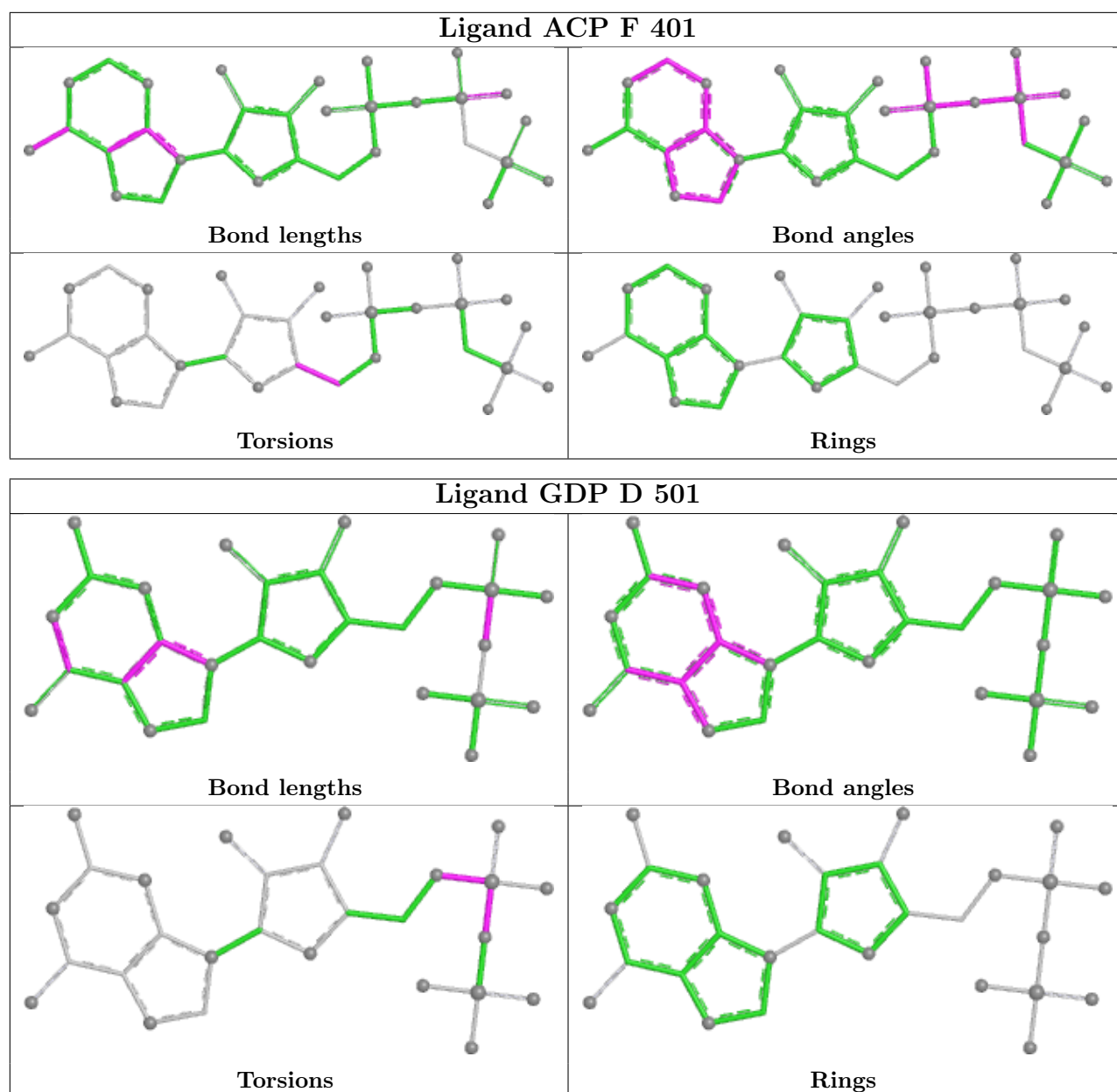


Ligand EP D 504



Ligand GTP C 501





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	436/451 (96%)	0.02	6 (1%) 73 75	33, 60, 104, 138	0
1	C	440/451 (97%)	-0.07	11 (2%) 58 60	24, 51, 91, 148	9 (2%)
2	B	422/445 (94%)	0.20	15 (3%) 46 48	25, 63, 108, 182	7 (1%)
2	D	430/445 (96%)	0.23	10 (2%) 61 63	34, 66, 110, 188	7 (1%)
3	E	123/143 (86%)	0.43	9 (7%) 21 23	42, 78, 122, 142	0
4	F	303/384 (78%)	0.65	26 (8%) 16 17	47, 87, 168, 244	0
All	All	2154/2319 (92%)	0.19	77 (3%) 46 48	24, 64, 121, 244	23 (1%)

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	440	VAL	5.7
2	B	286	LEU	5.0
3	E	143	ALA	4.7
4	F	132	LEU	4.5
4	F	172	PHE	4.1
4	F	182	ILE	3.8
2	B	57	THR	3.7
2	B	440	ALA	3.5
2	D	82	PRO	3.5
1	C	1	MET	3.4
1	C	283	HIS	3.4
4	F	371	PRO	3.3
4	F	134	ALA	3.3
4	F	141	GLY	3.3
3	E	14	CYS	3.3
2	D	369[A]	ARG	3.3
2	B	337	ASN	3.2
1	A	283	HIS	3.2
4	F	363	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
3	E	26	PRO	3.1
4	F	179	VAL	3.1
4	F	169	LEU	3.0
3	E	27	PRO	2.9
1	C	340	SER	2.9
4	F	162	ILE	2.9
4	F	256	TYR	2.8
2	D	179	ASP	2.8
1	C	357	TYR	2.8
2	D	59	ASN	2.8
3	E	24	LEU	2.8
4	F	98	TYR	2.8
2	D	57	THR	2.7
4	F	147	TRP	2.7
2	B	343	PHE	2.7
2	B	336	GLN	2.6
3	E	126	LYS	2.6
2	D	50	ASN	2.6
1	A	46	ASP	2.6
2	B	248	LEU	2.6
4	F	36	ARG	2.6
4	F	173	ILE	2.5
4	F	167	SER	2.5
4	F	241	THR	2.5
4	F	186	LEU	2.5
3	E	6	MET	2.5
4	F	257	GLU	2.4
2	B	333	LEU	2.4
1	A	179	THR	2.4
2	B	214	PHE	2.4
1	C	2	ARG	2.4
2	D	248	LEU	2.4
2	D	97	SER	2.3
4	F	151	SER	2.3
4	F	181	VAL	2.3
2	B	275	LEU	2.3
1	C	285	GLN	2.3
1	C	341	ILE	2.3
3	E	44	ASP	2.2
2	D	283	TYR	2.2
1	A	438	ASP	2.2
2	B	220	THR	2.2

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Mol	Chain	Res	Type	RSRZ
4	F	330	ILE	2.2
1	C	358	GLN	2.2
2	D	407	TRP	2.1
1	A	280	LYS	2.1
4	F	82	LYS	2.1
3	E	15	THR	2.1
2	B	2	ARG	2.1
1	A	262	TYR	2.1
2	B	246	GLY	2.1
2	B	438	ALA	2.1
4	F	379	HIS	2.1
1	C	364	PRO	2.1
2	B	325	MET	2.0
1	C	335	ILE	2.0
4	F	89	GLU	2.0
4	F	372	THR	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
7	CA	C	504	1/1	0.73	0.38	116,116,116,116	0
6	MG	F	403	1/1	0.75	0.23	97,97,97,97	0
6	MG	A	502	1/1	0.80	0.19	58,58,58,58	0
9	GOL	D	505	6/6	0.82	0.17	83,109,119,132	0
6	MG	F	402	1/1	0.84	0.20	126,126,126,126	0
7	CA	C	505	1/1	0.85	0.35	97,97,97,97	0
11	MES	D	503	12/12	0.85	0.17	87,108,149,155	0

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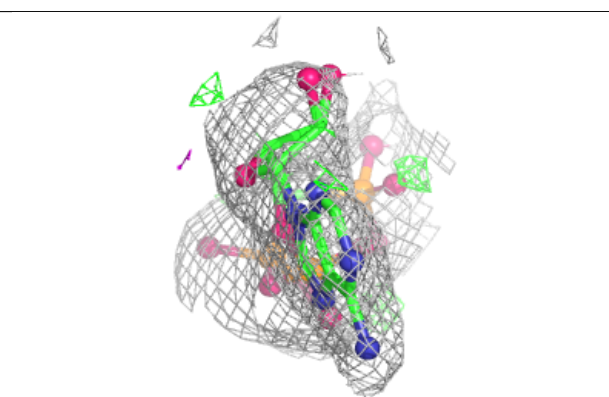
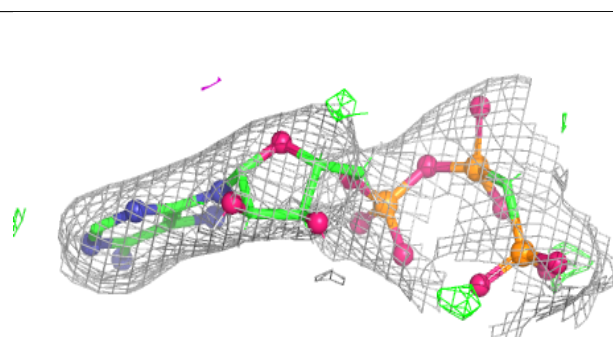
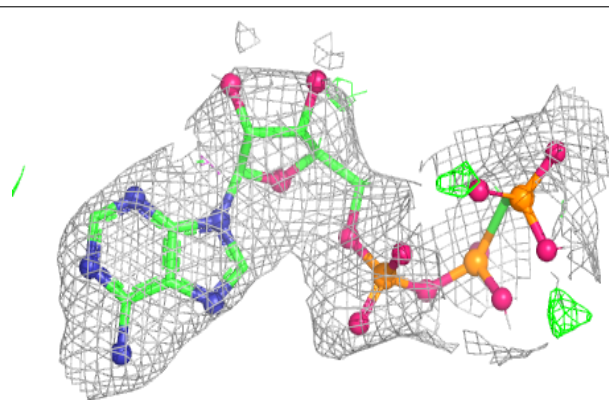
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
9	GOL	A	506	6/6	0.86	0.16	54,74,97,106	0
13	ACP	F	401	31/31	0.86	0.10	63,97,118,143	0
9	GOL	A	507	6/6	0.87	0.16	57,90,114,114	0
7	CA	B	503	1/1	0.88	0.27	91,91,91,91	0
6	MG	D	502	1/1	0.91	0.23	56,56,56,56	0
7	CA	A	505	1/1	0.91	0.23	97,97,97,97	0
7	CA	C	503	1/1	0.95	0.14	72,72,72,72	0
11	MES	B	504	12/12	0.95	0.09	52,67,89,94	0
10	GDP	D	501	28/28	0.96	0.07	41,53,69,74	0
12	EP	D	504	34/34	0.96	0.07	38,67,90,100	0
8	CL	A	504	1/1	0.96	0.09	73,73,73,73	0
5	GTP	A	501	32/32	0.98	0.05	28,42,56,60	0
6	MG	C	502	1/1	0.98	0.15	39,39,39,39	0
10	GDP	B	501	28/28	0.98	0.05	22,39,50,72	0
5	GTP	C	501	32/32	0.98	0.05	25,35,46,51	0
7	CA	A	503	1/1	0.99	0.10	72,72,72,72	0
6	MG	B	502	1/1	0.99	0.13	35,35,35,35	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.

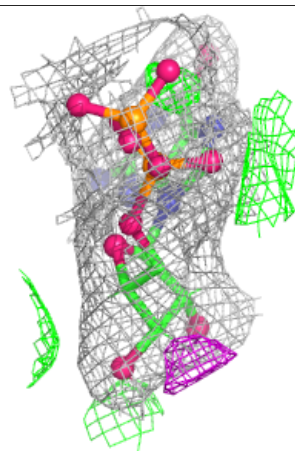
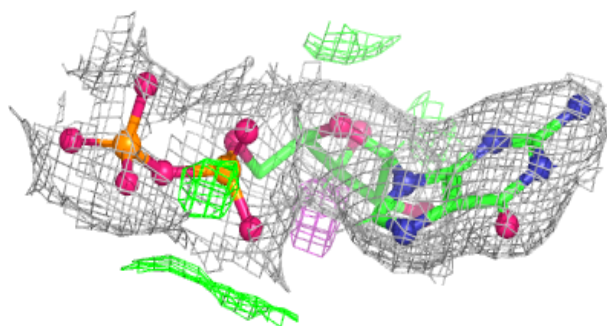
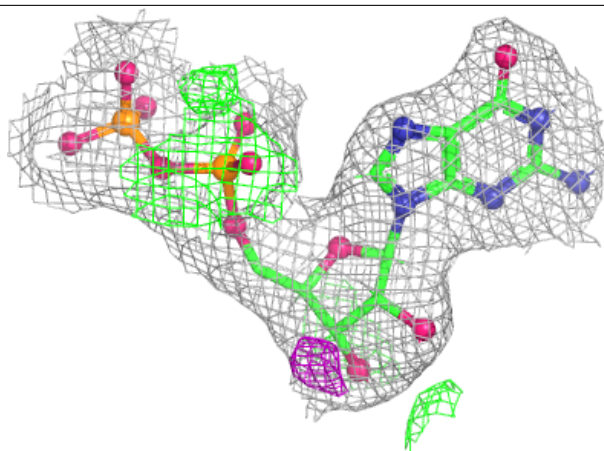
Electron density around ACP F 401:

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



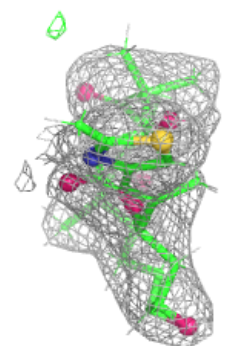
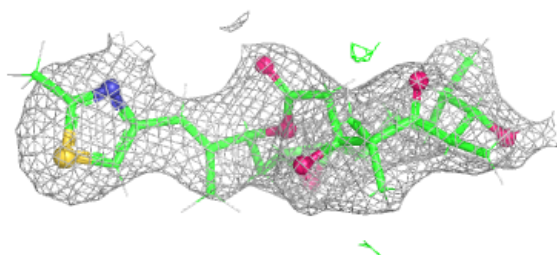
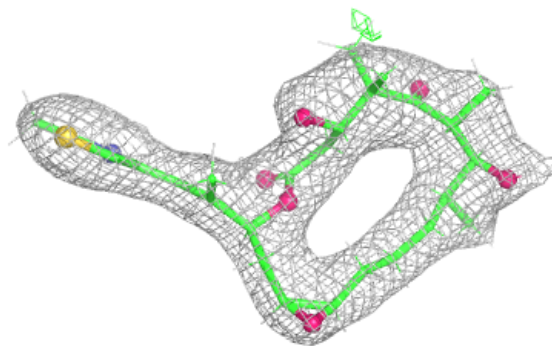
Electron density around GDP D 501:

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

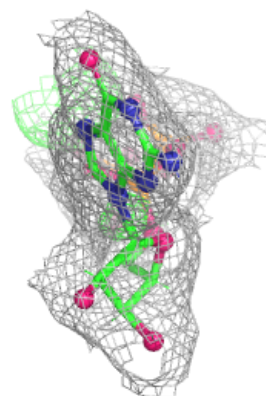
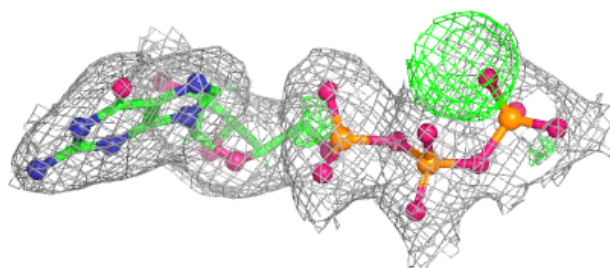
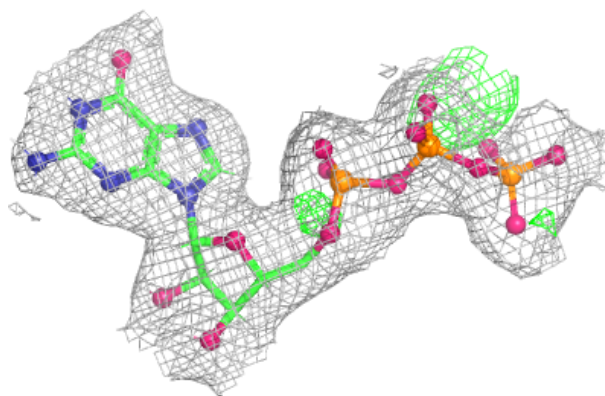


Electron density around EP D 504:

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

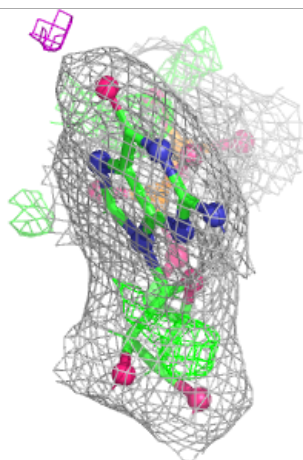
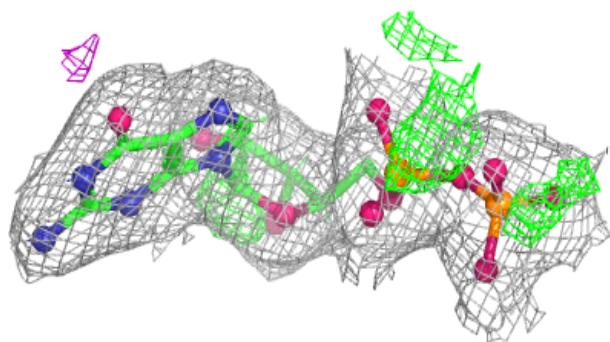
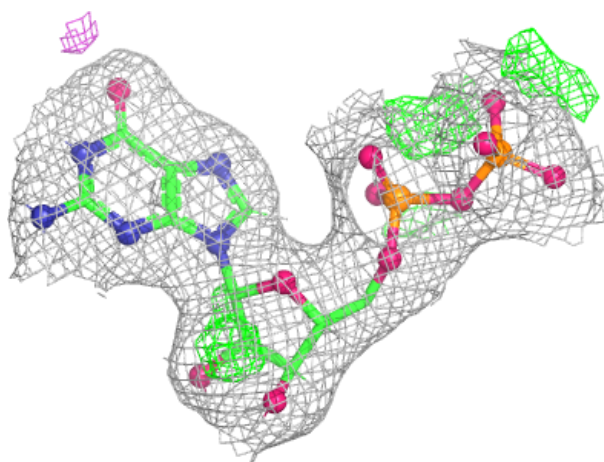
**Electron density around GTP A 501:**

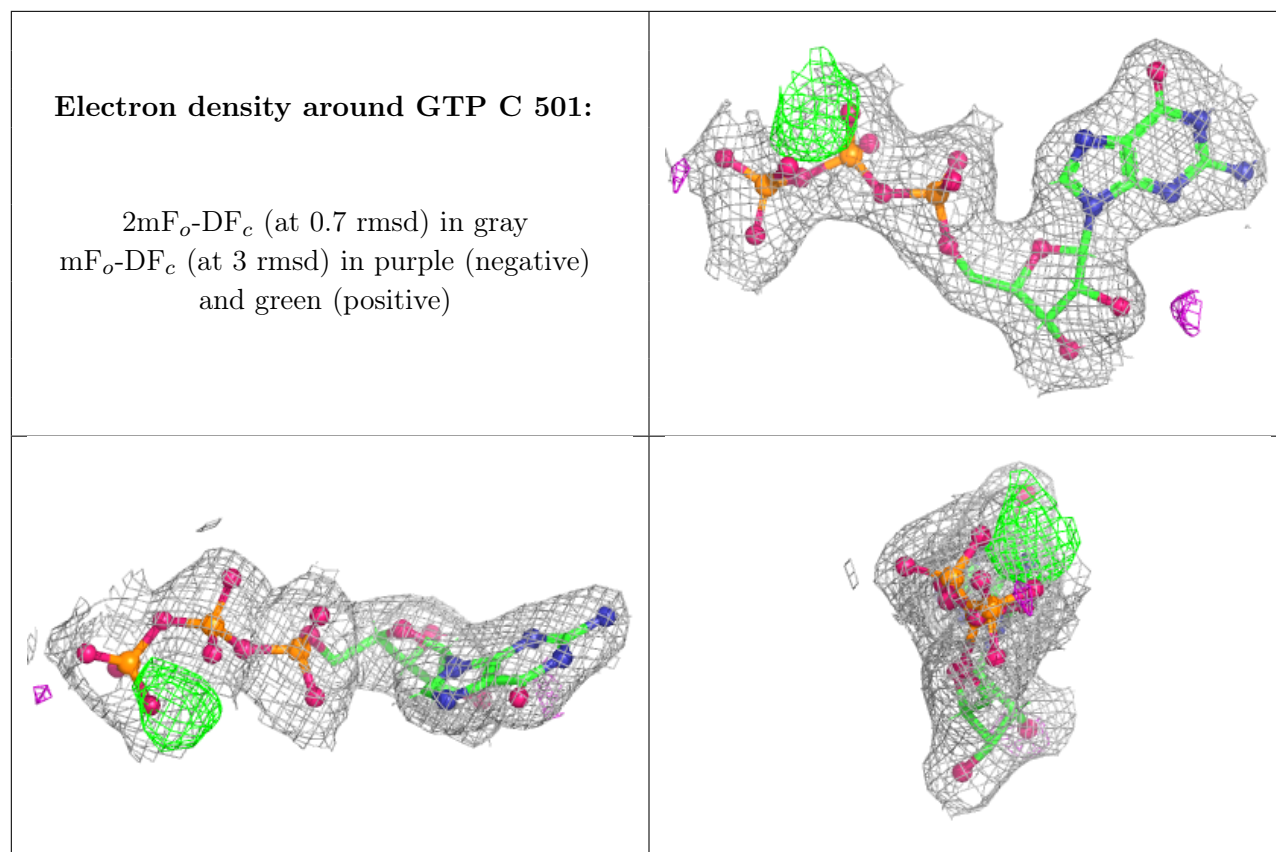
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around 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.