



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 21, 2021 – 01:41 PM EDT

PDB ID : 5S5D
Title : Tubulin-Z32400357-complex
Authors : Muehlethaler, T.; Gioia, D.; Prota, A.E.; Sharpe, M.E.; Cavalli, A.; Steinmetz, M.O.
Deposited on : 2020-11-08
Resolution : 1.90 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.20
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.20

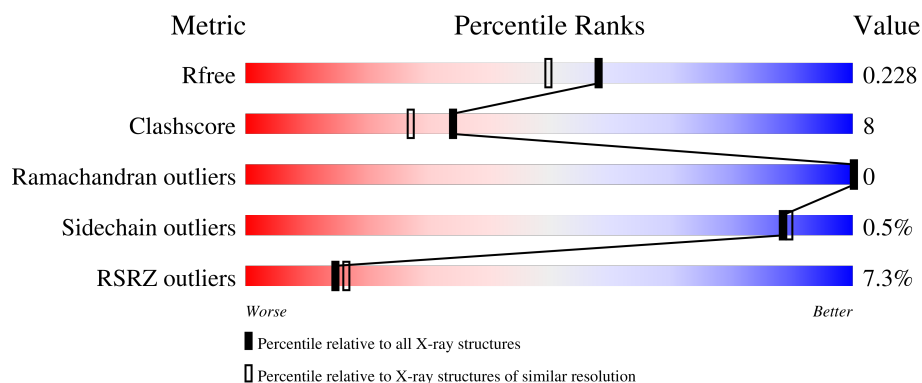
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 Å.

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




Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	<div> <div>4%</div> <div>79%</div> <div>18%</div> <div>.</div> </div>
1	C	451	<div> <div>2%</div> <div>83%</div> <div>14%</div> <div>.</div> </div>
2	B	445	<div> <div>7%</div> <div>77%</div> <div>18%</div> <div>5%</div> </div>
2	D	445	<div> <div>4%</div> <div>81%</div> <div>16%</div> <div>.</div> </div>
3	E	143	<div> <div>6%</div> <div>73%</div> <div>13%</div> <div>15%</div> </div>

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Mol	Chain	Length	Quality of chain
4	F	384	 A horizontal bar chart showing the quality of chain F. The bar is divided into four segments: red (20%), green (72%), yellow (18%), and grey (9%). The percentages are labeled above or below the corresponding segments.

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 18605 atoms, of which 36 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	437	Total	C	N	O	S	0	0	0
			3416	2163	581	650	22			
1	C	440	Total	C	N	O	S	0	1	0
			3443	2178	585	657	23			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	423	Total	C	N	O	S	1	1	0
			3349	2103	575	644	27			
2	D	431	Total	C	N	O	S	5	0	0
			3389	2126	580	656	27			

- Molecule 3 is a protein called Stathmin-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	122	Total	C	N	O	S	0	0	0
			1008	622	182	199	5			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	3	MET	-	initiating methionine	UNP P63043
E	4	ALA	-	expression tag	UNP P63043

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	348	Total	C	N	O	S	0	0	0
			2859	1833	491	521	14			

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 (three-letter code: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).



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

- Molecule 6 is MAGNESIUM ION (three-letter code: 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	1	Total	Mg	0	0
			1	1		

- Molecule 7 is CALCIUM ION (three-letter code: 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	2	Total	Ca	0	0
			2	2		

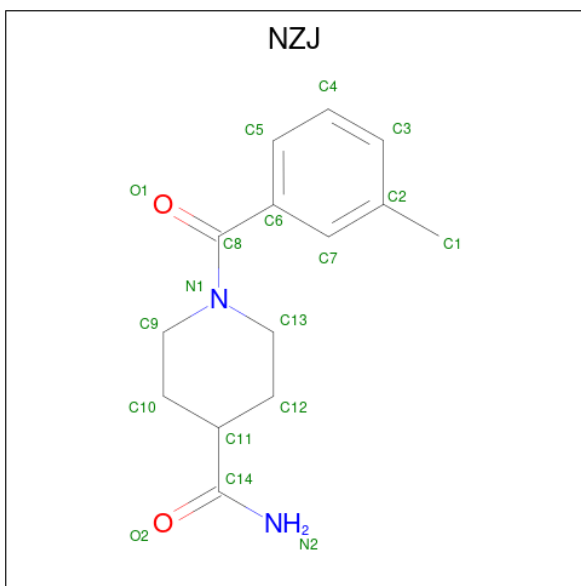
- Molecule 8 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).





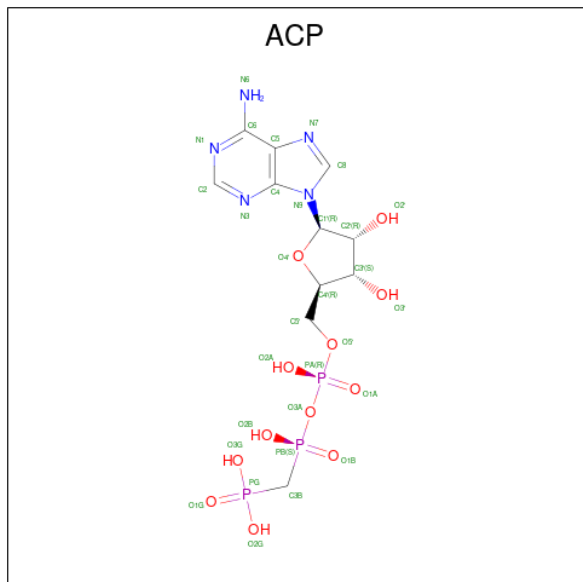
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 10 is 1-(3-methylbenzene-1-carbonyl)piperidine-4-carboxamide (three-letter code: NZJ) (formula: $C_{14}H_{18}N_2O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
10	B	1	Total	C	H	N	O	0	0
			36	14	18	2	2		
10	C	1	Total	C	H	N	O	0	0
			36	14	18	2	2		

- Molecule 11 is PHOSPHOMETHYLPHOSPHONIC ACID ADENYLATE ESTER (three-letter code: ACP) (formula: $C_{11}H_{18}N_5O_{12}P_3$).



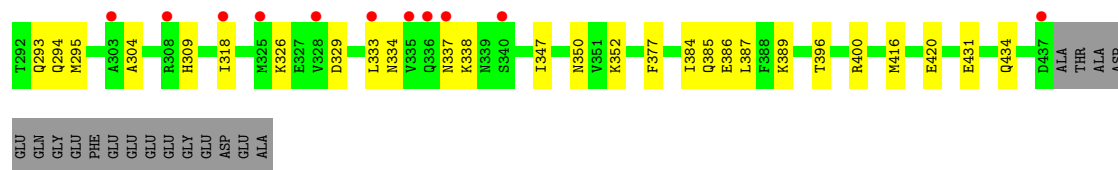
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
11	F	1	Total	C	N	O	P	0	0
			31	11	5	12	3		

- Molecule 12 is water.

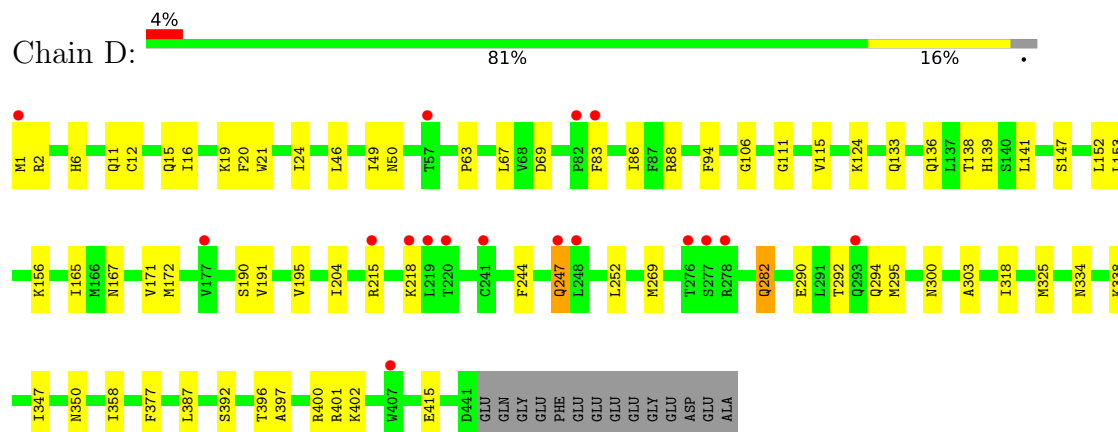
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	175	Total	O	0	0
			175	175		
12	B	165	Total	O	0	0
			165	165		
12	C	316	Total	O	0	0
			316	316		
12	D	128	Total	O	0	0
			128	128		
12	E	51	Total	O	0	0
			51	51		
12	F	61	Total	O	0	0
			61	61		

- Molecule 1: Tubulin alpha-1B chain

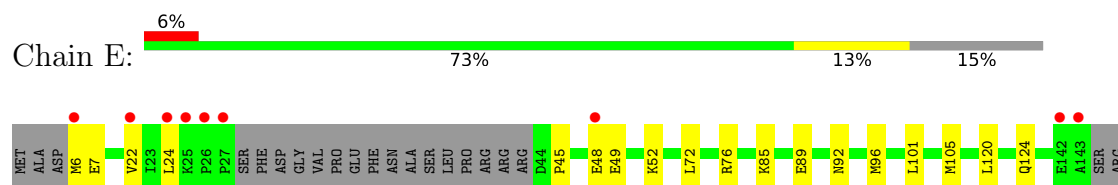




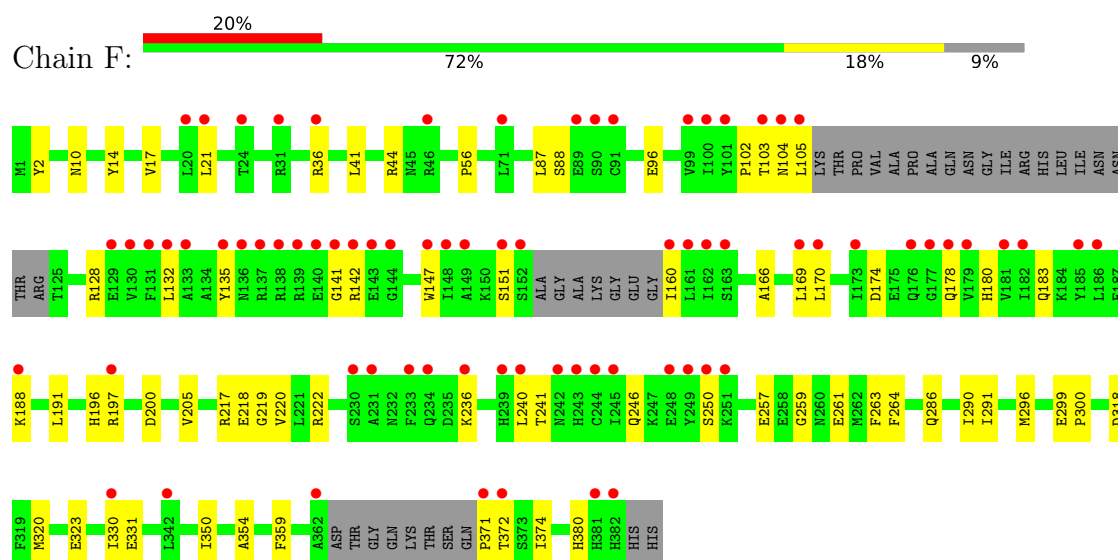
• Molecule 2: Tubulin beta-2B chain



• 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	105.07Å 159.08Å 179.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	63.42 – 1.90 118.97 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (63.42-1.90) 99.9 (118.97-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.14 (at 1.90Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.200 , 0.227 0.200 , 0.228	Depositor DCC
R_{free} test set	11889 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	43.1	Xtriage
Anisotropy	0.154	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 47.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	18605	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.41% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MES, GTP, CA, GDP, MG, NZJ, ACP

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	0.25	0/3494	0.42	0/4743
1	C	0.26	0/3521	0.44	1/4780 (0.0%)
2	B	0.26	0/3422	0.43	0/4630
2	D	0.25	0/3464	0.42	0/4692
3	E	0.24	0/1016	0.35	0/1348
4	F	0.24	0/2926	0.39	0/3954
All	All	0.25	0/17843	0.42	1/24147 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	171	ILE	C-N-CA	5.62	135.74	121.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	3416	0	3330	67	0
1	C	3443	0	3352	53	0
2	B	3349	0	3230	60	0
2	D	3389	0	3266	56	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	1008	0	1024	13	0
4	F	2859	0	2823	54	0
5	A	32	0	12	1	0
5	C	32	0	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	1	0	0	0	0
7	A	2	0	0	0	0
7	B	1	0	0	0	0
7	C	2	0	0	0	0
8	B	28	0	12	0	0
8	D	28	0	12	1	0
9	B	12	0	12	3	0
10	B	18	18	0	0	0
10	C	18	18	0	0	0
11	F	31	0	14	4	0
12	A	175	0	0	5	0
12	B	165	0	0	3	0
12	C	316	0	0	8	0
12	D	128	0	0	4	0
12	E	51	0	0	0	0
12	F	61	0	0	1	0
All	All	18569	36	17099	290	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 8.

All (290) 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:381:THR:HG22	1:A:383:ALA:H	1.20	1.06
4:F:102:PRO:HG2	4:F:105:LEU:HD13	1.47	0.95
1:C:221:ARG:HG3	2:D:325:MET:HG2	1.54	0.89
1:C:270:ALA:O	1:C:302:MET:HG2	1.74	0.87
4:F:241:THR:OG1	11:F:401:ACP:O3'	1.99	0.81
1:C:209:ILE:HD11	1:C:302:MET:HE3	1.62	0.81
1:C:209:ILE:HD11	1:C:302:MET:CE	2.11	0.80
1:C:234:ILE:HD13	1:C:302:MET:HE1	1.63	0.80
2:D:334:ASN:HD21	2:D:338:LYS:HE3	1.45	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:SER:OG	1:A:183:GLU:OE1	2.01	0.78
4:F:236:LYS:HB3	4:F:240:LEU:HD13	1.64	0.77
1:C:234:ILE:HD13	1:C:302:MET:CE	2.15	0.77
1:A:335:ILE:HG23	1:A:339:ARG:HG3	1.67	0.76
1:C:327:ASP:OD2	12:C:601:HOH:O	2.04	0.75
2:B:163:ASP:OD1	12:B:601:HOH:O	2.05	0.74
1:A:70:LEU:HD13	1:A:110:ILE:HG21	1.70	0.74
4:F:331:GLU:OE2	11:F:401:ACP:O3G	2.07	0.73
2:B:274:PRO:HB3	2:B:286:LEU:HD22	1.71	0.72
2:D:282:GLN:HA	2:D:282:GLN:HE21	1.56	0.71
4:F:102:PRO:HG2	4:F:105:LEU:CD1	2.21	0.70
4:F:318:ASP:OD2	11:F:401:ACP:O2G	2.10	0.70
2:B:83:PHE:O	2:B:86:ILE:HG22	1.92	0.69
1:A:71:GLU:OE2	1:A:73:THR:OG1	2.07	0.69
1:C:423:GLU:OE1	12:C:603:HOH:O	2.10	0.69
1:C:433:GLU:OE2	12:C:602:HOH:O	2.09	0.68
1:A:294:ALA:O	1:A:300:ASN:ND2	2.27	0.68
1:C:1:MET:O	12:C:604:HOH:O	2.11	0.68
4:F:128:ARG:NH2	4:F:174:ASP:OD1	2.27	0.68
4:F:10:ASN:HB2	4:F:44:ARG:HH22	1.58	0.67
1:C:423:GLU:OE2	12:C:605:HOH:O	2.12	0.67
4:F:197:ARG:NH1	4:F:257:GLU:OE2	2.27	0.67
3:E:48:GLU:HG2	3:E:52:LYS:HE3	1.77	0.66
1:C:367:ASP:OD1	12:C:606:HOH:O	2.12	0.66
4:F:371:PRO:HA	4:F:372:THR:O	1.95	0.66
1:A:336:LYS:HG3	3:E:24:LEU:HD13	1.77	0.66
4:F:371:PRO:HA	4:F:372:THR:HB	1.77	0.66
2:D:290:GLU:OE2	12:D:601:HOH:O	2.14	0.66
2:D:136:GLN:HA	2:D:167:ASN:O	1.96	0.65
2:D:21:TRP:CZ3	2:D:63:PRO:HB3	2.31	0.65
1:A:179:THR:HA	2:B:352:LYS:HD2	1.78	0.65
1:C:93:ILE:HD11	1:C:121:ARG:HG3	1.79	0.64
1:C:4[A]:CYS:SG	1:C:136:LEU:HG	2.37	0.64
2:D:269:MET:HG3	2:D:303:ALA:HB3	1.80	0.64
2:B:2:ARG:HB2	2:B:133:GLN:CG	2.28	0.64
2:B:141:LEU:HD12	2:B:172:MET:SD	2.39	0.63
2:B:337:ASN:OD1	4:F:36:ARG:HD3	1.98	0.63
2:D:83:PHE:O	2:D:86:ILE:HG22	1.99	0.62
2:D:334:ASN:ND2	2:D:338:LYS:HE3	2.13	0.62
2:D:397:ALA:HA	2:D:400:ARG:NH1	2.15	0.62
1:A:209:ILE:HD11	1:A:302:MET:SD	2.40	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:286:LEU:HD12	2:B:290:GLU:OE1	2.00	0.61
2:B:21:TRP:CZ3	2:B:63:PRO:HB3	2.36	0.61
1:C:76:ASP:O	1:C:80:THR:HG22	2.01	0.61
4:F:246:GLN:O	4:F:250:SER:HB3	2.01	0.61
2:D:106:GLY:O	2:D:111:GLY:HA3	2.01	0.61
4:F:151:SER:HB3	4:F:180:HIS:CD2	2.36	0.60
2:B:420:GLU:OE1	12:B:602:HOH:O	2.16	0.60
1:C:322:ASP:OD1	12:C:607:HOH:O	2.16	0.60
3:E:48:GLU:CG	3:E:52:LYS:HE3	2.32	0.60
1:A:66:VAL:HG23	1:A:125:LEU:HD12	1.84	0.59
2:D:1:MET:HG3	2:D:50:ASN:HB2	1.82	0.59
1:C:244:PHE:CD1	1:C:358:GLN:HG2	2.37	0.59
1:C:270:ALA:HB3	1:C:302:MET:HG3	1.84	0.58
2:B:396:THR:O	2:B:400:ARG:HG3	2.02	0.58
4:F:371:PRO:CA	4:F:372:THR:HB	2.32	0.58
2:B:88:ARG:NH1	2:B:90:ASP:HB2	2.19	0.58
1:A:335:ILE:HG23	1:A:339:ARG:CG	2.33	0.58
2:B:123:ARG:O	2:B:127:GLU:HG3	2.05	0.57
2:B:295:MET:CG	2:B:377:PHE:HB2	2.34	0.57
2:D:295:MET:HE2	2:D:377:PHE:HB2	1.85	0.57
1:C:419:SER:O	1:C:423:GLU:HG3	2.05	0.57
1:C:234:ILE:CD1	1:C:302:MET:HE1	2.35	0.57
2:B:69:ASP:O	2:B:94:PHE:HA	2.05	0.57
2:D:244:PHE:CE1	2:D:358:ILE:HD12	2.40	0.57
1:C:249:ASN:OD1	1:C:356:ASN:ND2	2.38	0.56
1:A:188:ILE:HD12	1:A:395:PHE:CD2	2.40	0.56
2:B:274:PRO:HB3	2:B:286:LEU:CD2	2.34	0.56
2:B:284:ARG:NH2	2:B:290:GLU:OE2	2.37	0.56
3:E:120:LEU:O	3:E:124:GLN:HG3	2.05	0.56
2:B:304:ALA:N	12:B:610:HOH:O	2.39	0.56
2:D:215:ARG:O	2:D:218:LYS:HE3	2.04	0.56
2:B:164:ARG:O	9:B:504:MES:H31	2.06	0.56
2:D:2:ARG:NH1	12:D:604:HOH:O	2.31	0.56
1:A:351:PHE:HE1	3:E:24:LEU:HD11	1.71	0.56
2:D:152:LEU:O	2:D:156:LYS:HG2	2.05	0.56
3:E:92:ASN:O	3:E:96:MET:HG2	2.06	0.56
4:F:10:ASN:CB	4:F:44:ARG:HH22	2.18	0.56
4:F:17:VAL:O	4:F:21:LEU:HG	2.06	0.56
3:E:85:LYS:O	3:E:89:GLU:HG3	2.06	0.56
2:D:69:ASP:O	2:D:94:PHE:HA	2.05	0.55
1:C:292:THR:HG22	1:C:335:ILE:CD1	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:GLN:CG	4:F:56:PRO:HB3	2.37	0.55
2:D:124:LYS:C	2:D:124:LYS:HD3	2.26	0.55
1:A:36:MET:HB3	1:A:61:HIS:CE1	2.41	0.55
3:E:101:LEU:O	3:E:105:MET:HG2	2.07	0.55
4:F:151:SER:HB3	4:F:180:HIS:CG	2.42	0.55
4:F:200:ASP:OD1	4:F:222:ARG:HB2	2.07	0.55
1:A:210:TYR:CE1	1:A:222:PRO:HD2	2.42	0.54
1:C:286:LEU:HA	1:C:290:GLU:OE1	2.06	0.54
2:D:397:ALA:O	2:D:401:ARG:NH1	2.40	0.54
2:B:36:TYR:CD1	2:B:46:LEU:HD21	2.43	0.54
1:A:430:LYS:O	1:A:434:GLU:HG3	2.07	0.54
2:B:136:GLN:HA	2:B:167:ASN:O	2.08	0.54
1:A:254:GLU:HG2	1:A:258:ASN:ND2	2.23	0.53
3:E:72:LEU:O	3:E:76:ARG:HG2	2.07	0.53
1:C:320:ARG:HA	1:C:356:ASN:O	2.08	0.53
4:F:132:LEU:HD21	4:F:170:LEU:HD11	1.91	0.53
1:A:88:HIS:HB2	1:A:89:PRO:HD2	1.90	0.53
4:F:371:PRO:HA	4:F:372:THR:C	2.28	0.53
1:A:187:SER:CB	1:A:391:LEU:HD21	2.38	0.53
1:C:1:MET:HE3	1:C:131:GLY:HA3	1.91	0.53
2:D:21:TRP:CE3	2:D:63:PRO:HB3	2.45	0.52
2:D:282:GLN:HA	2:D:282:GLN:NE2	2.23	0.52
1:C:167:LEU:HG	1:C:200:CYS:HB3	1.89	0.52
2:B:290:GLU:O	2:B:294:GLN:HG3	2.10	0.52
2:D:115:VAL:HG23	2:D:153:LEU:HD23	1.90	0.52
1:A:21:TRP:CZ3	1:A:63:PRO:HB3	2.45	0.52
2:B:124:LYS:HD3	2:B:124:LYS:C	2.30	0.52
1:A:214:ARG:HG2	1:A:219:ILE:O	2.10	0.52
2:D:167:ASN:ND2	12:D:612:HOH:O	2.43	0.52
4:F:2:TYR:CE1	4:F:359:PHE:HB3	2.45	0.52
4:F:296:MET:SD	4:F:380:HIS:HB2	2.50	0.51
4:F:320:MET:HG3	4:F:330:ILE:HD11	1.92	0.51
1:A:284:GLU:CD	1:A:284:GLU:H	2.14	0.51
1:A:220:GLU:OE1	2:B:326:LYS:HD2	2.10	0.51
2:D:1:MET:HG3	2:D:50:ASN:CB	2.40	0.51
1:C:66:VAL:HG23	1:C:125:LEU:CD1	2.41	0.51
4:F:147:TRP:HB2	4:F:169:LEU:CD1	2.41	0.51
2:B:187:ALA:O	2:B:191:VAL:HG23	2.11	0.50
1:C:88:HIS:HE1	1:C:90:GLU:HG3	1.75	0.50
2:D:6:HIS:CD2	2:D:21:TRP:HE1	2.29	0.50
2:B:106:GLY:O	2:B:111:GLY:HA3	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:172:MET:HG3	2:D:387:LEU:HD11	1.92	0.50
2:B:38:GLY:HA3	2:B:45:GLN:OE1	2.12	0.50
2:B:431:GLU:O	2:B:434:GLN:HG2	2.12	0.50
2:D:171:VAL:HA	2:D:204:ILE:O	2.12	0.50
2:B:334:ASN:OD1	2:B:338:LYS:HD3	2.11	0.50
2:B:347:ILE:HG22	2:B:350:ASN:HB3	1.92	0.50
1:A:265:ILE:HG23	1:A:432:TYR:CE1	2.48	0.49
2:B:2:ARG:HB2	2:B:133:GLN:NE2	2.27	0.49
1:C:88:HIS:CE1	1:C:90:GLU:HG3	2.46	0.49
4:F:219:GLY:HA3	4:F:264:PHE:CZ	2.47	0.49
2:B:199:ASP:OD1	9:B:504:MES:H62	2.12	0.49
2:D:141:LEU:HD12	2:D:172:MET:SD	2.52	0.49
1:C:172:TYR:HB3	1:C:205:ASP:HA	1.95	0.49
1:A:71:GLU:HG2	1:A:72:PRO:HD2	1.95	0.49
2:B:158:ARG:NH1	2:B:196:GLU:O	2.45	0.49
2:B:42:LEU:H	2:B:42:LEU:HD12	1.78	0.49
2:D:12:CYS:HB2	8:D:501:GDP:C8	2.48	0.49
1:A:74:VAL:HB	12:A:610:HOH:O	2.13	0.49
1:A:188:ILE:HD12	1:A:395:PHE:HB2	1.95	0.49
1:C:271:THR:HG21	1:C:295:CYS:O	2.12	0.49
1:C:11:GLN:HE22	2:D:247:GLN:NE2	2.11	0.49
4:F:205:VAL:HG21	4:F:291:ILE:HD13	1.95	0.48
2:B:318:ILE:N	2:B:318:ILE:HD12	2.28	0.48
1:C:270:ALA:C	1:C:302:MET:HG2	2.33	0.48
4:F:14:TYR:HB3	4:F:41:LEU:HD13	1.94	0.48
1:C:21:TRP:CZ3	1:C:63:PRO:HB3	2.48	0.48
2:B:295:MET:HG2	2:B:377:PHE:HB2	1.95	0.48
2:B:289:PRO:O	2:B:293:GLN:HG3	2.13	0.48
1:A:196:GLU:OE1	1:A:196:GLU:HA	2.14	0.48
2:D:165:ILE:HG21	2:D:252:LEU:HB3	1.95	0.48
4:F:141:GLY:O	4:F:142:ARG:HB2	2.13	0.48
2:B:385:GLN:OE1	2:B:389:LYS:HE3	2.14	0.48
2:B:88:ARG:HH11	2:B:90:ASP:HB2	1.78	0.47
2:B:172:MET:HG3	2:B:387:LEU:HD11	1.96	0.47
2:D:402:LYS:HE2	2:D:415:GLU:OE1	2.14	0.47
3:E:45:PRO:HA	3:E:49:GLU:OE1	2.14	0.47
4:F:103:THR:HG23	4:F:128:ARG:NH2	2.29	0.47
1:A:210:TYR:CZ	1:A:222:PRO:HD2	2.49	0.47
2:B:2:ARG:HB2	2:B:133:GLN:HG3	1.96	0.47
2:B:6:HIS:CD2	2:B:21:TRP:HE1	2.31	0.47
2:D:2:ARG:HB3	2:D:133:GLN:HG2	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:102:PRO:HB2	4:F:104:ASN:OD1	2.14	0.47
2:D:396:THR:O	2:D:400:ARG:HG2	2.15	0.47
2:D:1:MET:HE2	2:D:50:ASN:HB2	1.97	0.47
4:F:188:LYS:HD3	4:F:323:GLU:OE2	2.13	0.47
4:F:350:ILE:O	4:F:354:ALA:HB3	2.15	0.47
1:A:226:ASN:ND2	1:A:367:ASP:OD2	2.48	0.47
2:D:15:GLN:O	2:D:19:LYS:HG2	2.15	0.47
2:D:318:ILE:N	2:D:318:ILE:HD12	2.30	0.46
4:F:147:TRP:HB2	4:F:169:LEU:HD11	1.96	0.46
4:F:87:LEU:O	4:F:88:SER:OG	2.29	0.46
1:A:278:ALA:HA	1:A:369:ALA:HB2	1.98	0.46
9:B:504:MES:H51	9:B:504:MES:H81	1.55	0.46
1:A:100:ALA:HA	2:B:254:LYS:HG3	1.98	0.46
2:D:67:LEU:N	2:D:67:LEU:HD12	2.30	0.46
4:F:286:GLN:O	4:F:290:ILE:HG13	2.15	0.46
1:A:66:VAL:HG23	1:A:125:LEU:CD1	2.45	0.46
1:C:320:ARG:HG3	1:C:360:PRO:HG3	1.98	0.46
2:D:2:ARG:HB2	2:D:133:GLN:HE21	1.80	0.46
4:F:191:LEU:HD12	4:F:196:HIS:CE1	2.51	0.46
2:D:347:ILE:HG22	2:D:350:ASN:HB3	1.98	0.46
11:F:401:ACP:O3G	11:F:401:ACP:O1B	2.34	0.46
4:F:160:ILE:HD12	4:F:160:ILE:N	2.31	0.45
4:F:359:PHE:O	12:F:501:HOH:O	2.21	0.45
2:D:88:ARG:HD2	12:D:711:HOH:O	2.16	0.45
2:B:46:LEU:HA	2:B:49:ILE:HB	1.97	0.45
1:C:270:ALA:HB3	1:C:302:MET:CG	2.46	0.45
1:A:70:LEU:HB2	1:A:98:ASP:HA	1.97	0.45
1:A:176:GLN:HG3	4:F:56:PRO:HB3	1.98	0.45
1:A:21:TRP:CE3	1:A:63:PRO:HB3	2.52	0.45
2:D:2:ARG:HB3	2:D:133:GLN:CG	2.47	0.45
2:D:191:VAL:O	2:D:195:VAL:HG23	2.17	0.45
2:D:387:LEU:HD23	2:D:387:LEU:C	2.37	0.45
1:A:193:THR:HG23	12:A:653:HOH:O	2.16	0.45
1:C:341:ILE:HD13	1:C:351:PHE:HZ	1.81	0.45
1:A:207:GLU:OE2	1:A:304:LYS:HD2	2.17	0.44
2:B:269:MET:HG2	2:B:384:ILE:HD13	1.99	0.44
2:D:1:MET:CE	2:D:50:ASN:HB2	2.47	0.44
4:F:217:ARG:HG3	4:F:218:GLU:HG2	1.98	0.44
2:B:21:TRP:CE3	2:B:63:PRO:HB3	2.52	0.44
1:A:187:SER:HB3	1:A:391:LEU:HD21	1.98	0.44
1:A:304:LYS:HB2	12:A:619:HOH:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:320:MET:CG	4:F:330:ILE:HD11	2.47	0.44
1:A:300:ASN:HB3	12:A:601:HOH:O	2.17	0.44
2:B:223:THR:O	2:B:227:LEU:HD13	2.18	0.44
1:A:151:SER:HB2	1:A:193:THR:OG1	2.18	0.43
1:C:344:VAL:HG21	1:C:346:TRP:CE2	2.53	0.43
2:B:16:ILE:HD13	2:B:231:VAL:HG11	1.99	0.43
1:C:194:THR:O	1:C:194:THR:HG22	2.18	0.43
1:C:430:LYS:HE2	1:C:434:GLU:OE2	2.18	0.43
1:A:98:ASP:HB2	5:A:501:GTP:O2G	2.18	0.43
2:B:42:LEU:HD12	2:B:42:LEU:N	2.33	0.43
2:B:200:GLU:OE2	2:B:255:LEU:HG	2.19	0.43
1:A:188:ILE:HD11	1:A:392:ASP:HA	1.99	0.43
2:B:23:VAL:HG21	2:B:232:SER:HB3	2.00	0.43
1:A:176:GLN:HG3	4:F:56:PRO:HG3	2.00	0.43
1:A:292:THR:HG22	1:A:335:ILE:HD12	2.01	0.43
3:E:7:GLU:O	3:E:22:VAL:HA	2.18	0.43
1:A:100:ALA:HA	2:B:254:LYS:CG	2.49	0.43
1:C:192:HIS:CG	1:C:421:ALA:HA	2.54	0.43
2:D:20:PHE:CZ	2:D:24:ILE:HD13	2.53	0.43
2:D:147:SER:HB2	2:D:190:SER:OG	2.19	0.43
1:A:142:GLY:HA3	1:A:183:GLU:OE2	2.19	0.43
4:F:135:TYR:CZ	4:F:166:ALA:HB2	2.54	0.43
1:A:188:ILE:HD12	1:A:395:PHE:CB	2.49	0.42
1:A:351:PHE:HE1	3:E:24:LEU:CD1	2.32	0.42
2:B:416:MET:O	2:B:420:GLU:HG3	2.18	0.42
1:C:312:TYR:CD1	1:C:341:ILE:HG23	2.54	0.42
2:D:11:GLN:O	2:D:15:GLN:HG2	2.19	0.42
4:F:96:GLU:O	4:F:183:GLN:HA	2.19	0.42
1:A:209:ILE:HG23	1:A:230:LEU:HD23	2.00	0.42
4:F:205:VAL:CG2	4:F:291:ILE:HD13	2.50	0.42
4:F:259:GLY:O	4:F:261:GLU:HG3	2.20	0.42
1:A:3:GLU:OE1	1:A:3:GLU:N	2.51	0.42
2:D:115:VAL:HG23	2:D:153:LEU:CD2	2.50	0.42
2:D:392:SER:O	2:D:396:THR:HG22	2.20	0.42
4:F:299:GLU:HB3	4:F:300:PRO:HD3	2.02	0.42
1:A:233:GLN:HG3	1:A:368:LEU:CD1	2.49	0.42
2:B:67:LEU:HD12	2:B:67:LEU:N	2.35	0.42
1:C:66:VAL:HG12	1:C:68:VAL:HG23	2.01	0.42
2:D:16:ILE:HD11	2:D:138:THR:HB	2.02	0.42
1:A:132:LEU:O	1:A:164:LYS:NZ	2.53	0.41
1:A:241:SER:HB2	1:A:248:LEU:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:66:VAL:HG23	1:C:125:LEU:HD11	2.02	0.41
1:C:2:ARG:HD3	1:C:2:ARG:HA	1.73	0.41
1:C:93:ILE:CD1	1:C:121:ARG:HG3	2.49	0.41
2:D:294:GLN:HG2	2:D:300:ASN:ND2	2.34	0.41
4:F:178:GLN:N	4:F:178:GLN:OE1	2.52	0.41
2:B:333:LEU:O	2:B:337:ASN:ND2	2.53	0.41
1:C:36:MET:HB3	1:C:61:HIS:CE1	2.55	0.41
1:A:166:LYS:HE2	1:A:197:HIS:O	2.20	0.41
1:C:233:GLN:HG3	1:C:368:LEU:CD1	2.50	0.41
4:F:217:ARG:NH2	4:F:374:ILE:HA	2.35	0.41
2:B:165:ILE:HG21	2:B:252:LEU:HB3	2.01	0.41
2:B:247:GLN:O	2:B:250:ALA:HB2	2.21	0.41
1:A:320:ARG:HG3	12:A:681:HOH:O	2.20	0.41
1:C:133:GLN:NE2	1:C:253:THR:HG21	2.35	0.41
2:B:291:LEU:HD23	2:B:291:LEU:HA	1.92	0.41
1:A:71:GLU:HG2	1:A:72:PRO:CD	2.51	0.41
1:A:75:ILE:HB	1:A:94:THR:CG2	2.51	0.41
2:B:309:HIS:ND1	2:B:386:GLU:OE2	2.37	0.41
4:F:151:SER:HB3	4:F:180:HIS:CE1	2.56	0.41
1:C:196:GLU:HG2	12:C:702:HOH:O	2.20	0.41
1:C:66:VAL:HG23	1:C:125:LEU:HD12	2.03	0.40
2:D:46:LEU:HA	2:D:49:ILE:HB	2.03	0.40
2:B:329:ASP:O	2:B:333:LEU:HG	2.20	0.40
1:A:31:GLN:HB2	1:A:33:ASP:OD1	2.22	0.40
1:A:187:SER:HB2	1:A:391:LEU:HD21	2.03	0.40
1:A:217:LEU:HD21	1:A:368:LEU:HD23	2.02	0.40
1:C:11:GLN:HE22	2:D:247:GLN:CD	2.24	0.40
1:C:275:VAL:HG13	1:C:368:LEU:HD21	2.04	0.40
2:D:292:THR:O	2:D:295:MET:HG2	2.21	0.40
4:F:220:VAL:HG12	4:F:263:PHE:CE1	2.56	0.40
1:A:188:ILE:HD12	1:A:395:PHE:CG	2.57	0.40
1:A:227:LEU:O	1:A:231:ILE:HG13	2.21	0.40
1:A:329:ASN:HB3	3:E:6:MET:CE	2.51	0.40
1:A:419:SER:O	1:A:423:GLU:HG3	2.21	0.40
4:F:320:MET:HG2	4:F:330:ILE:HG13	2.02	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	435/451 (96%)	427 (98%)	8 (2%)	0	100	100
1	C	439/451 (97%)	430 (98%)	9 (2%)	0	100	100
2	B	418/445 (94%)	410 (98%)	8 (2%)	0	100	100
2	D	429/445 (96%)	421 (98%)	8 (2%)	0	100	100
3	E	118/143 (82%)	118 (100%)	0	0	100	100
4	F	340/384 (88%)	331 (97%)	9 (3%)	0	100	100
All	All	2179/2319 (94%)	2137 (98%)	42 (2%)	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	368/379 (97%)	365 (99%)	3 (1%)	81	82
1	C	372/379 (98%)	369 (99%)	3 (1%)	81	82
2	B	368/383 (96%)	367 (100%)	1 (0%)	92	93
2	D	372/383 (97%)	369 (99%)	3 (1%)	81	82
3	E	109/127 (86%)	109 (100%)	0	100	100
4	F	315/342 (92%)	315 (100%)	0	100	100
All	All	1904/1993 (96%)	1894 (100%)	10 (0%)	88	89

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	221	ARG
1	A	282	TYR
1	A	300	ASN
2	B	139	HIS
1	C	2	ARG
1	C	71	GLU
1	C	221	ARG
2	D	139	HIS
2	D	247	GLN
2	D	282	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	88	HIS
1	A	301	GLN
2	B	294	GLN
1	C	11	GLN
1	C	85	GLN
1	C	406	HIS
2	D	281	GLN
2	D	282	GLN
4	F	229	ASN
4	F	333	ASN
4	F	380	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 18 ligands modelled in this entry, 10 are monoatomic - leaving 8 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
8	GDP	D	501	6	24,30,30	1.17	2 (8%)	31,47,47	1.91	7 (22%)
9	MES	B	504	-	12,12,12	2.23	1 (8%)	14,16,16	2.00	5 (35%)
10	NZJ	B	505	-	19,19,19	1.01	2 (10%)	26,26,26	0.99	3 (11%)
8	GDP	B	501	6	24,30,30	1.13	2 (8%)	31,47,47	1.84	7 (22%)
5	GTP	A	501	6	26,34,34	1.00	1 (3%)	33,54,54	1.71	7 (21%)
5	GTP	C	501	6	26,34,34	0.97	1 (3%)	33,54,54	1.61	5 (15%)
10	NZJ	C	505	-	19,19,19	0.99	2 (10%)	26,26,26	1.23	4 (15%)
11	ACP	F	401	6	27,33,33	1.40	5 (18%)	32,52,52	1.47	4 (12%)

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
8	GDP	D	501	6	-	3/12/32/32	0/3/3/3
9	MES	B	504	-	-	4/6/14/14	0/1/1/1
10	NZJ	B	505	-	-	2/12/22/22	0/2/2/2
8	GDP	B	501	6	-	3/12/32/32	0/3/3/3
5	GTP	A	501	6	-	9/18/38/38	0/3/3/3
5	GTP	C	501	6	-	8/18/38/38	0/3/3/3
10	NZJ	C	505	-	-	2/12/22/22	0/2/2/2
11	ACP	F	401	6	-	9/15/38/38	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	B	504	MES	C8-S	-7.46	1.66	1.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	D	501	GDP	C6-C5	4.13	1.48	1.41
8	B	501	GDP	C6-C5	3.69	1.47	1.41
5	A	501	GTP	C6-N1	3.23	1.38	1.33
5	C	501	GTP	C6-N1	3.12	1.38	1.33
11	F	401	ACP	PG-O2G	2.94	1.61	1.54
11	F	401	ACP	PG-O3G	2.90	1.61	1.54
11	F	401	ACP	PB-O3A	2.73	1.61	1.58
11	F	401	ACP	C5-C4	2.54	1.47	1.40
10	C	505	NZJ	C8-N1	2.43	1.40	1.34
10	B	505	NZJ	C8-N1	2.41	1.40	1.34
8	D	501	GDP	C5-C4	2.39	1.47	1.40
8	B	501	GDP	C5-C4	2.29	1.47	1.40
11	F	401	ACP	PB-O2B	2.24	1.61	1.56
10	B	505	NZJ	C14-N2	2.17	1.38	1.32
10	C	505	NZJ	C14-N2	2.16	1.38	1.32

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	501	GTP	N3-C2-N1	-5.55	119.82	127.22
5	C	501	GTP	N3-C2-N1	-5.27	120.19	127.22
8	D	501	GDP	C2-N3-C4	4.83	120.87	115.36
8	B	501	GDP	C2-N3-C4	4.51	120.51	115.36
5	A	501	GTP	C2-N3-C4	4.33	120.30	115.36
8	B	501	GDP	C6-C5-C4	-4.29	116.71	120.80
8	D	501	GDP	C6-C5-C4	-4.20	116.79	120.80
8	B	501	GDP	C6-N1-C2	4.18	122.58	115.93
8	D	501	GDP	C6-N1-C2	4.14	122.50	115.93
11	F	401	ACP	PA-O3A-PB	-4.05	119.70	132.56
8	D	501	GDP	C5-C6-N1	-3.95	118.03	123.43
8	B	501	GDP	C5-C6-N1	-3.94	118.05	123.43
5	C	501	GTP	C2-N3-C4	3.91	119.83	115.36
9	B	504	MES	C5-N4-C3	3.63	117.00	108.83
8	B	501	GDP	N3-C2-N1	-3.51	122.54	127.22
11	F	401	ACP	C3'-C2'-C1'	3.46	106.19	100.98
8	D	501	GDP	N3-C2-N1	-3.42	122.66	127.22
11	F	401	ACP	N3-C2-N1	-3.12	123.80	128.68
5	C	501	GTP	C5-C6-N1	-3.02	119.30	123.43
9	B	504	MES	O1S-S-C8	2.99	110.52	106.92
10	B	505	NZJ	O2-C14-N2	-2.96	117.85	123.00
10	C	505	NZJ	O2-C14-N2	-2.89	117.97	123.00
9	B	504	MES	O2S-S-C8	2.85	110.34	106.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	501	GTP	C5-C6-N1	-2.79	119.61	123.43
9	B	504	MES	C6-C5-N4	-2.73	105.97	110.10
8	D	501	GDP	PA-O3A-PB	-2.72	123.48	132.83
5	C	501	GTP	C6-N1-C2	2.69	120.20	115.93
8	D	501	GDP	C4-C5-N7	-2.65	106.63	109.40
11	F	401	ACP	C4-C5-N7	-2.64	106.64	109.40
5	A	501	GTP	C6-N1-C2	2.64	120.13	115.93
10	B	505	NZJ	O1-C8-N1	-2.33	118.49	122.34
9	B	504	MES	C7-N4-C5	2.26	117.01	111.23
5	A	501	GTP	PB-O3B-PG	-2.25	125.12	132.83
5	C	501	GTP	PA-O3A-PB	-2.17	125.39	132.83
10	C	505	NZJ	O1-C8-N1	-2.16	118.76	122.34
8	B	501	GDP	C4-C5-N7	-2.16	107.15	109.40
5	A	501	GTP	N2-C2-N1	2.15	120.60	117.25
10	B	505	NZJ	C11-C14-N2	2.14	119.98	116.54
8	B	501	GDP	PA-O3A-PB	-2.11	125.60	132.83
5	A	501	GTP	PA-O3A-PB	-2.08	125.69	132.83
10	C	505	NZJ	C9-C10-C11	2.02	113.77	110.41
10	C	505	NZJ	C6-C8-N1	2.02	121.28	118.72

There are no chirality outliers.

All (40) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	501	GTP	C5'-O5'-PA-O1A
5	A	501	GTP	C5'-O5'-PA-O2A
5	C	501	GTP	C5'-O5'-PA-O1A
5	C	501	GTP	C5'-O5'-PA-O2A
8	B	501	GDP	C5'-O5'-PA-O1A
8	B	501	GDP	C5'-O5'-PA-O2A
8	D	501	GDP	C5'-O5'-PA-O1A
8	D	501	GDP	C5'-O5'-PA-O2A
9	B	504	MES	C8-C7-N4-C5
9	B	504	MES	C7-C8-S-O2S
9	B	504	MES	C7-C8-S-O3S
10	B	505	NZJ	C10-C11-C14-N2
10	B	505	NZJ	C10-C11-C14-O2
11	F	401	ACP	PB-C3B-PG-O1G
11	F	401	ACP	PB-C3B-PG-O2G
11	F	401	ACP	PB-C3B-PG-O3G
11	F	401	ACP	PG-C3B-PB-O1B
11	F	401	ACP	PG-C3B-PB-O2B

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Mol	Chain	Res	Type	Atoms
11	F	401	ACP	PG-C3B-PB-O3A
11	F	401	ACP	O4'-C4'-C5'-O5'
11	F	401	ACP	C3'-C4'-C5'-O5'
5	A	501	GTP	PB-O3B-PG-O1G
5	C	501	GTP	PB-O3B-PG-O1G
9	B	504	MES	C7-C8-S-O1S
10	C	505	NZJ	C10-C11-C14-N2
10	C	505	NZJ	C10-C11-C14-O2
5	A	501	GTP	C4'-C5'-O5'-PA
11	F	401	ACP	PB-O3A-PA-O2A
5	A	501	GTP	PB-O3B-PG-O2G
5	A	501	GTP	PB-O3B-PG-O3G
5	C	501	GTP	PB-O3B-PG-O2G
5	C	501	GTP	PB-O3B-PG-O3G
5	A	501	GTP	C5'-O5'-PA-O3A
5	C	501	GTP	C5'-O5'-PA-O3A
8	B	501	GDP	C5'-O5'-PA-O3A
8	D	501	GDP	C5'-O5'-PA-O3A
5	A	501	GTP	PB-O3A-PA-O1A
5	A	501	GTP	PB-O3A-PA-O2A
5	C	501	GTP	PB-O3A-PA-O1A
5	C	501	GTP	PB-O3A-PA-O2A

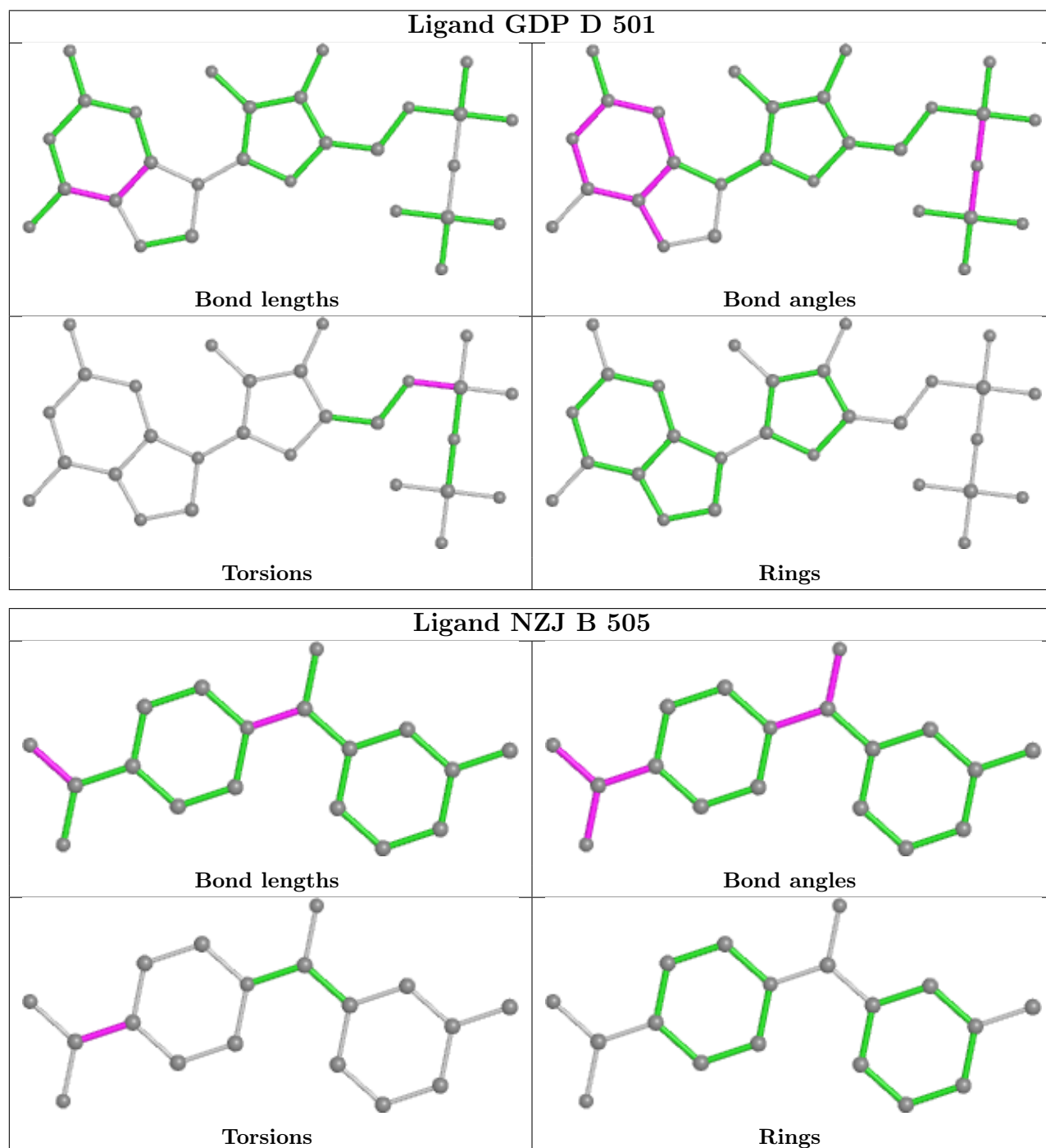
There are no ring outliers.

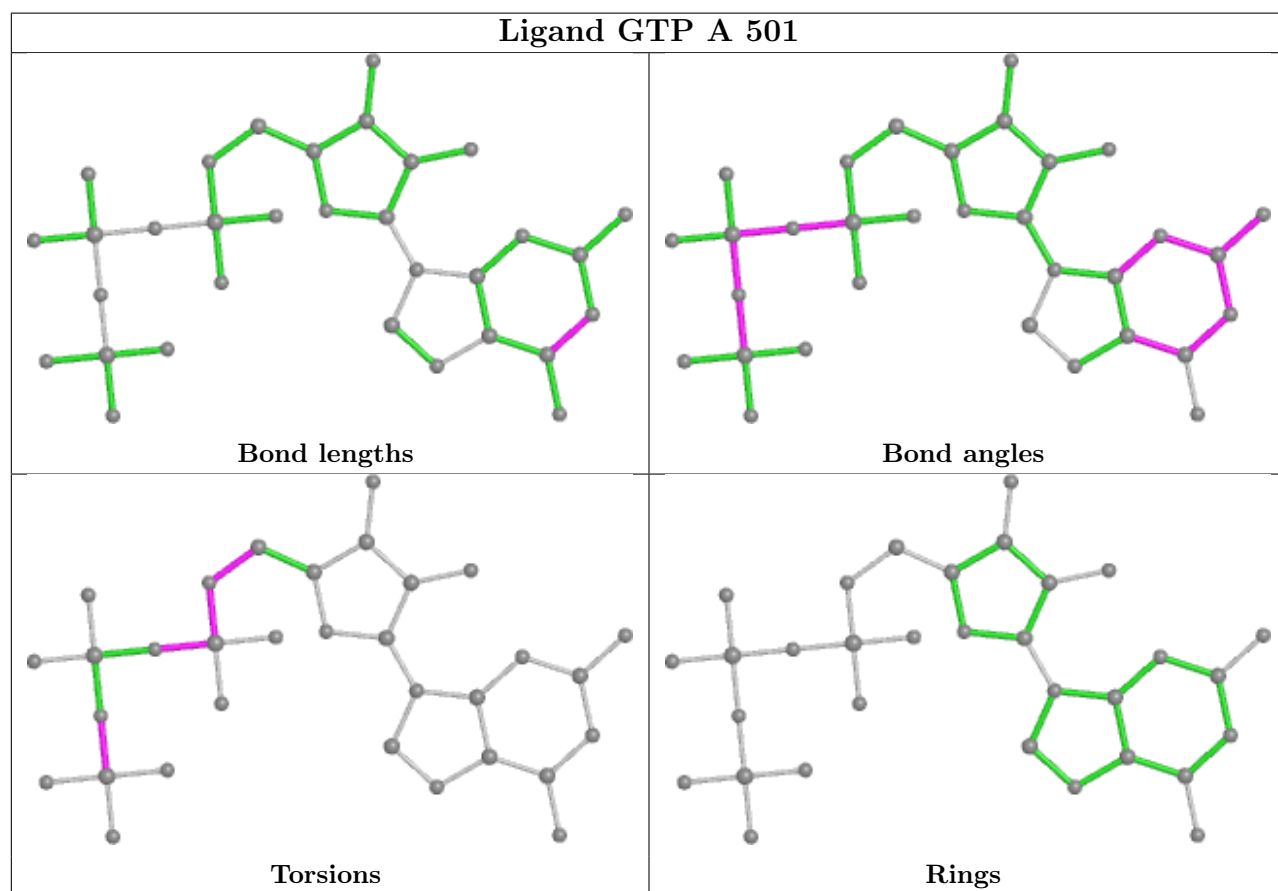
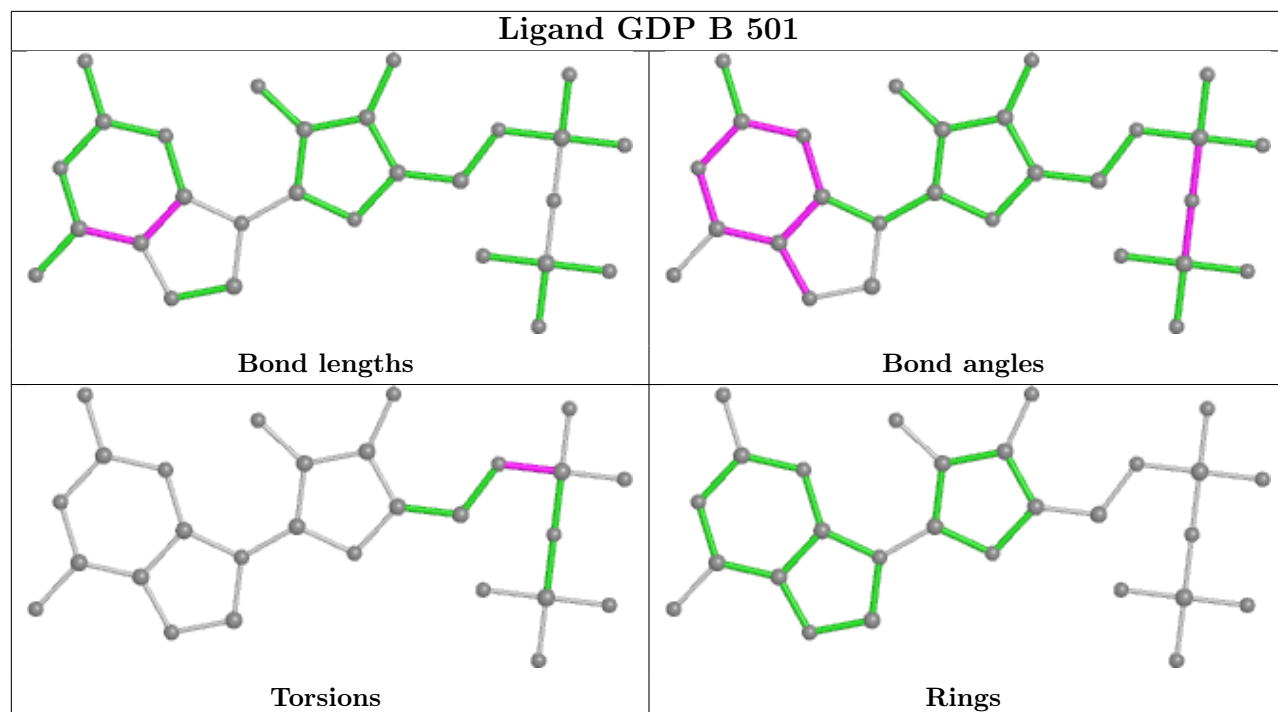
4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	D	501	GDP	1	0
9	B	504	MES	3	0
5	A	501	GTP	1	0
11	F	401	ACP	4	0

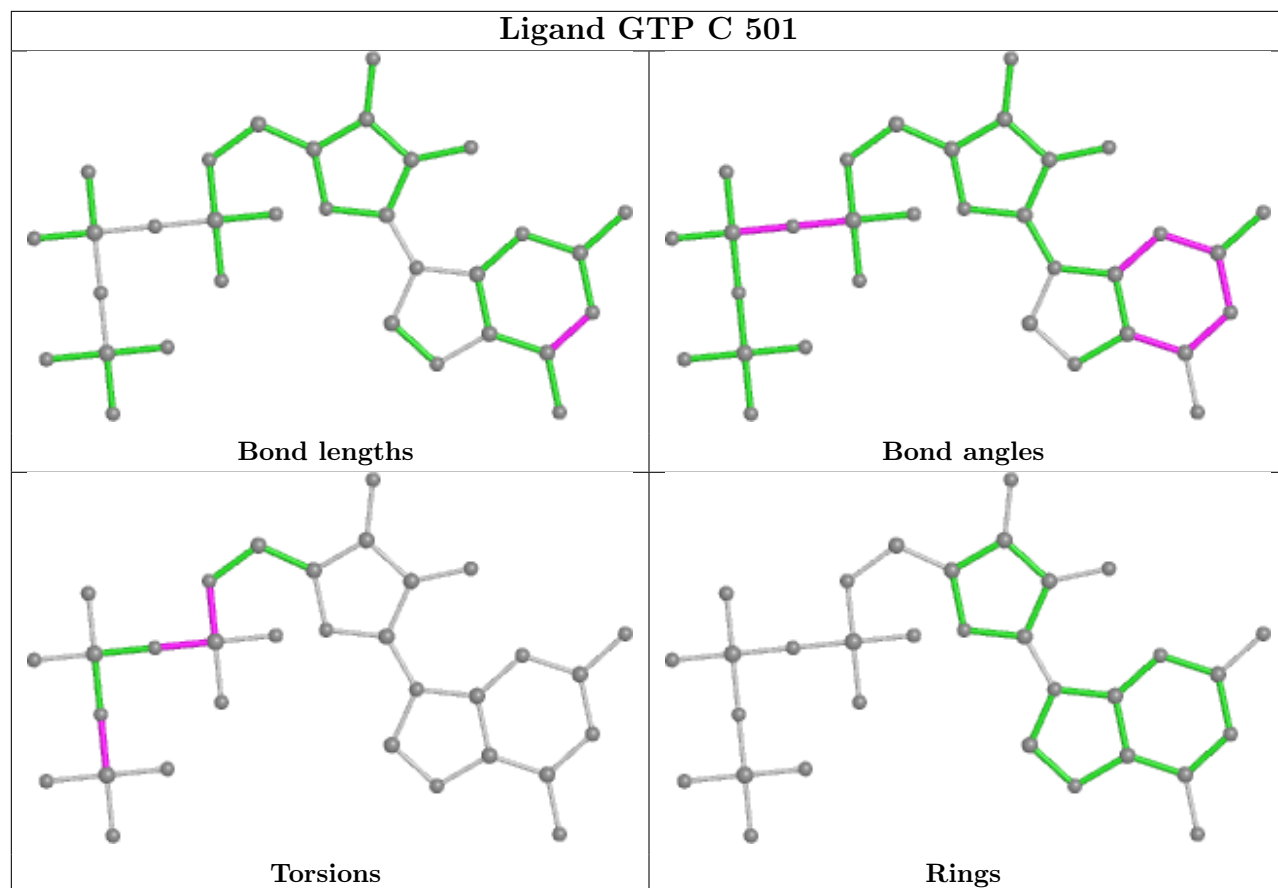
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

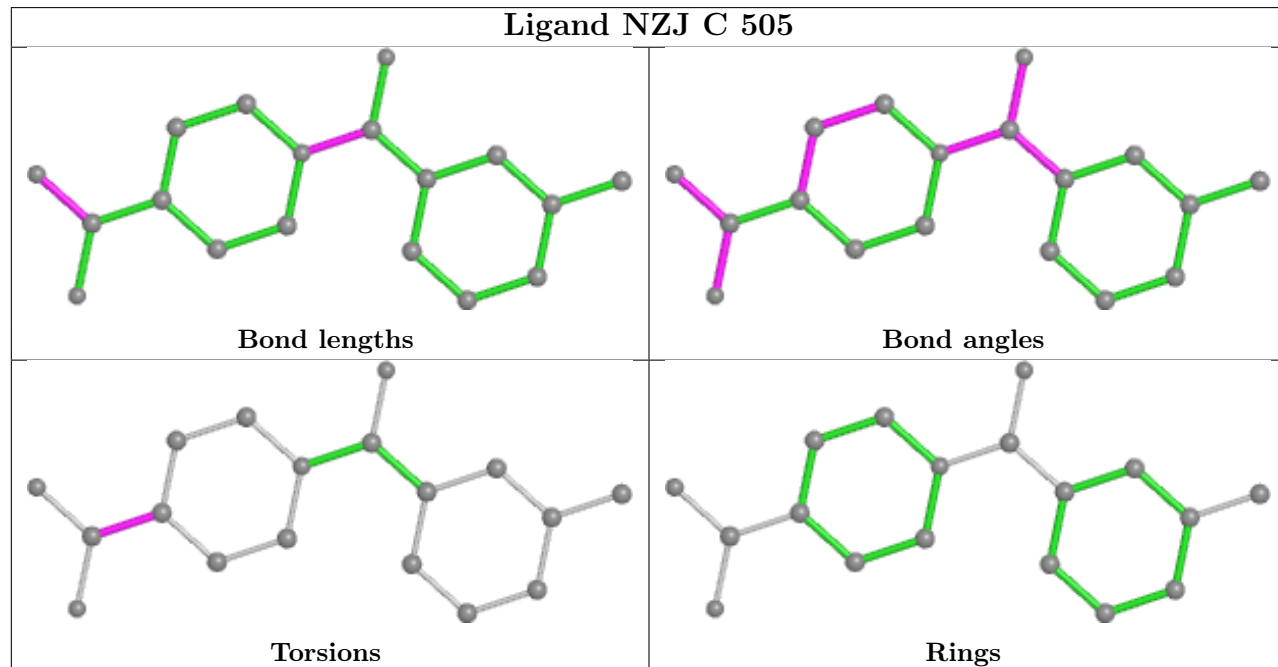


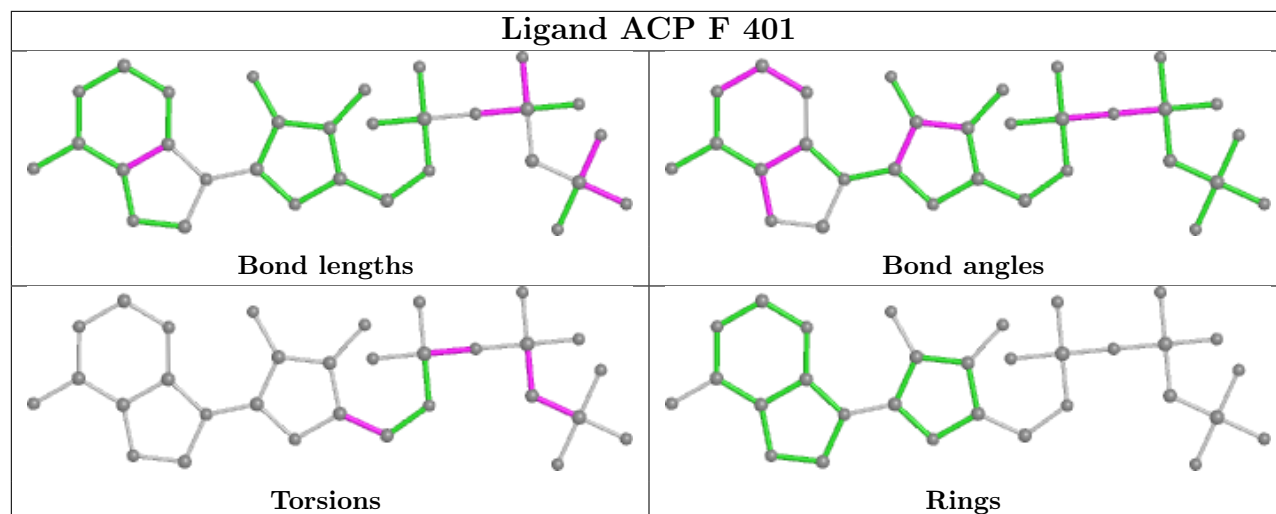


Ligand GTP C 501



Ligand NZJ C 505





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	437/451 (96%)	0.56	19 (4%)	35	38	34, 50, 83, 140	0
1	C	440/451 (97%)	0.56	11 (2%)	57	60	29, 40, 59, 103	0
2	B	423/445 (95%)	0.74	30 (7%)	16	17	29, 50, 90, 130	2 (0%)
2	D	431/445 (96%)	0.50	17 (3%)	39	42	34, 54, 83, 113	4 (0%)
3	E	122/143 (85%)	0.80	9 (7%)	14	16	38, 57, 102, 135	0
4	F	348/384 (90%)	1.30	75 (21%)	0	0	45, 76, 136, 166	0
All	All	2201/2319 (94%)	0.71	161 (7%)	15	16	29, 51, 102, 166	6 (0%)

All (161) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	F	173	ILE	8.8
3	E	26	PRO	7.6
4	F	372	THR	7.4
4	F	177	GLY	7.0
4	F	233	PHE	6.6
2	B	1	MET	6.5
2	B	333	LEU	6.3
4	F	105	LEU	6.2
1	A	179	THR	6.2
4	F	231	ALA	5.9
4	F	89	GLU	5.9
3	E	6	MET	5.6
1	A	346	TRP	5.6
4	F	142	ARG	5.6
2	B	276	THR	5.4
4	F	143	GLU	5.3
4	F	179	VAL	5.3
4	F	103	THR	5.3
4	F	130	VAL	5.2

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Mol	Chain	Res	Type	RSRZ
4	F	240	LEU	5.1
2	B	337	ASN	5.0
1	A	262	TYR	5.0
4	F	182	ILE	4.8
4	F	20	LEU	4.8
2	B	59	ASN	4.7
2	B	281	GLN	4.7
4	F	186	LEU	4.6
4	F	244	CYS	4.6
4	F	161	LEU	4.6
4	F	99	VAL	4.6
4	F	234	GLN	4.6
1	A	282	TYR	4.5
4	F	100	ILE	4.5
2	D	1	MET	4.3
1	C	340	SER	4.3
4	F	152	SER	4.3
4	F	169	LEU	4.2
4	F	90	SER	4.2
1	A	281	ALA	4.1
2	D	82	PRO	4.1
4	F	104	ASN	4.1
4	F	162	ILE	4.0
2	B	57	THR	3.9
4	F	371	PRO	3.9
2	B	325	MET	3.9
4	F	342	LEU	3.9
3	E	25	LYS	3.8
4	F	362	ALA	3.8
2	B	437	ASP	3.8
4	F	176	GLN	3.8
4	F	140	GLU	3.7
4	F	243	HIS	3.7
4	F	132	LEU	3.7
4	F	249	TYR	3.6
4	F	149	ALA	3.6
3	E	143	ALA	3.6
4	F	137	ARG	3.6
2	D	407	TRP	3.5
1	A	163	LYS	3.5
1	A	345	ASP	3.5
3	E	24	LEU	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	340	SER	3.5
2	D	276	THR	3.5
4	F	144	GLY	3.4
2	D	57	THR	3.4
4	F	330	ILE	3.3
2	B	283	TYR	3.3
2	D	215	ARG	3.3
4	F	250	SER	3.3
4	F	148	ILE	3.3
4	F	178	GLN	3.2
4	F	135	TYR	3.2
1	A	70	LEU	3.2
4	F	245	ILE	3.2
4	F	151	SER	3.2
1	C	302	MET	3.2
4	F	251	LYS	3.1
4	F	141	GLY	3.1
4	F	24	THR	3.0
3	E	48	GLU	3.0
4	F	160	ILE	3.0
2	B	308	ARG	3.0
4	F	138	ARG	2.9
2	B	285	ALA	2.9
1	A	335	ILE	2.9
2	D	83	PHE	2.9
2	D	278	ARG	2.9
1	C	218	ASP	2.9
4	F	239	HIS	2.9
4	F	101	TYR	2.8
2	B	245	PRO	2.8
2	B	335	VAL	2.8
4	F	181	VAL	2.8
2	B	58	GLY	2.8
4	F	242	ASN	2.8
2	B	246	GLY	2.8
1	C	308	ARG	2.8
4	F	91	CYS	2.8
1	C	286	LEU	2.8
2	B	318	ILE	2.7
4	F	163	SER	2.7
4	F	188	LYS	2.7
1	C	350	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
2	B	328	VAL	2.7
4	F	21	LEU	2.7
2	D	218	LYS	2.7
4	F	46	ARG	2.7
2	B	56	ALA	2.7
3	E	27	PRO	2.7
3	E	142	GLU	2.7
1	C	357	TYR	2.6
4	F	133	ALA	2.6
4	F	131	PHE	2.6
2	B	284	ARG	2.6
2	D	220	THR	2.6
2	D	277	SER	2.6
2	B	247	GLN	2.6
4	F	31	ARG	2.5
2	B	82	PRO	2.5
1	A	381	THR	2.5
2	B	340	SER	2.5
2	B	278	ARG	2.5
4	F	139	ARG	2.5
4	F	381	HIS	2.5
1	A	88	HIS	2.4
1	A	284	GLU	2.4
1	C	284	GLU	2.4
2	D	177	VAL	2.4
1	A	348	PRO	2.4
4	F	170	LEU	2.3
2	D	293	GLN	2.3
2	B	62	VAL	2.3
4	F	36	ARG	2.3
1	C	339	ARG	2.3
2	B	286	LEU	2.3
4	F	129	GLU	2.3
4	F	236	LYS	2.3
4	F	197	ARG	2.3
4	F	136	ASN	2.3
1	A	332	ILE	2.2
2	B	303	ALA	2.2
2	B	336	GLN	2.2
1	A	221	ARG	2.2
4	F	147	TRP	2.2
4	F	248	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
4	F	230	SER	2.2
2	D	219	LEU	2.2
1	C	283	HIS	2.2
2	B	241	CYS	2.2
2	B	36	TYR	2.1
1	C	295	CYS	2.1
3	E	22	VAL	2.1
2	D	247	GLN	2.1
2	D	241	CYS	2.1
4	F	382	HIS	2.1
1	A	362	VAL	2.1
1	A	285	GLN	2.1
4	F	71	LEU	2.1
4	F	185	TYR	2.0
2	D	248	LEU	2.0
1	A	349	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 monosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	CA	B	503	1/1	0.70	0.11	87,87,87,87	0
9	MES	B	504	12/12	0.76	0.20	62,69,88,93	0
6	MG	F	402	1/1	0.77	0.09	77,77,77,77	0
6	MG	D	502	1/1	0.79	0.11	51,51,51,51	0
7	CA	A	504	1/1	0.79	0.12	98,98,98,98	0
11	ACP	F	401	31/31	0.82	0.20	75,87,95,99	0
10	NZJ	B	505	18/18	0.91	0.15	39,48,62,62	36

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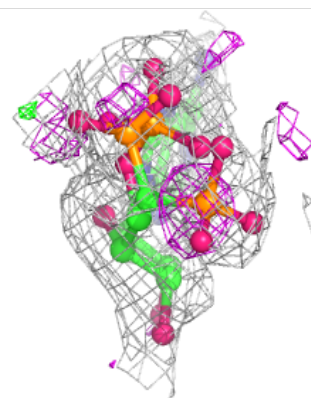
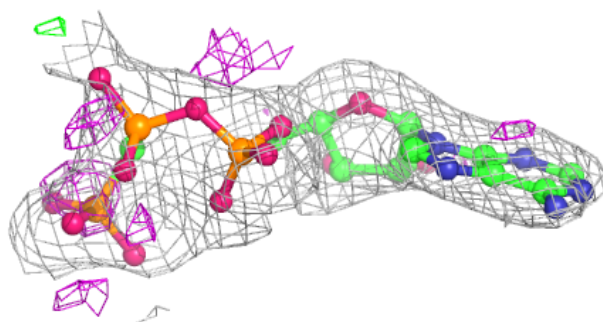
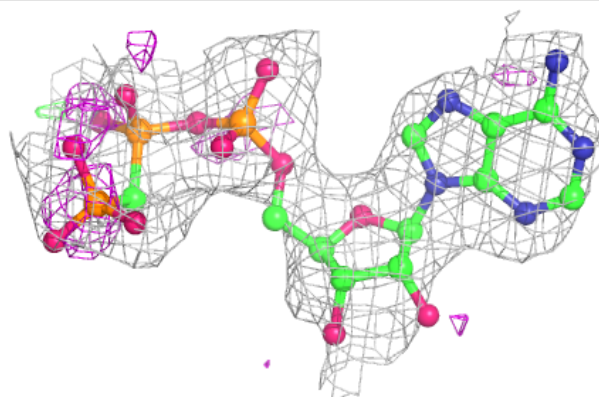
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
6	MG	C	502	1/1	0.92	0.09	32,32,32,32	0
8	GDP	D	501	28/28	0.95	0.15	45,50,58,60	0
10	NZJ	C	505	18/18	0.96	0.14	32,43,56,56	0
6	MG	A	502	1/1	0.96	0.08	37,37,37,37	0
7	CA	C	503	1/1	0.97	0.08	46,46,46,46	0
5	GTP	A	501	32/32	0.97	0.14	31,35,38,40	0
5	GTP	C	501	32/32	0.98	0.14	28,32,34,35	0
7	CA	C	504	1/1	0.98	0.10	45,45,45,45	0
8	GDP	B	501	28/28	0.98	0.14	30,33,38,41	0
7	CA	A	503	1/1	0.98	0.08	70,70,70,70	0
6	MG	B	502	1/1	1.00	0.27	28,28,28,28	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.

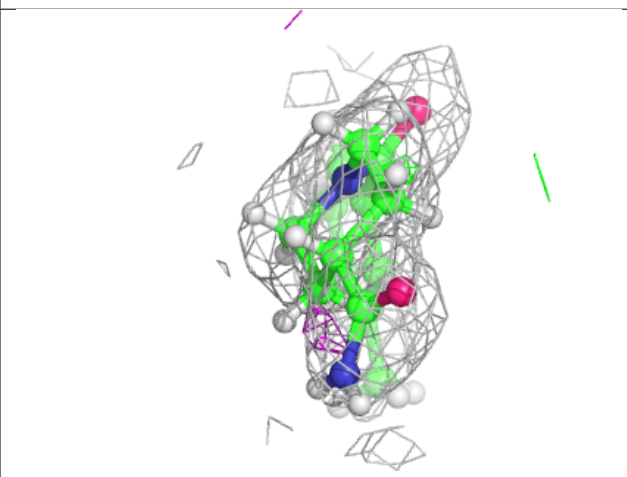
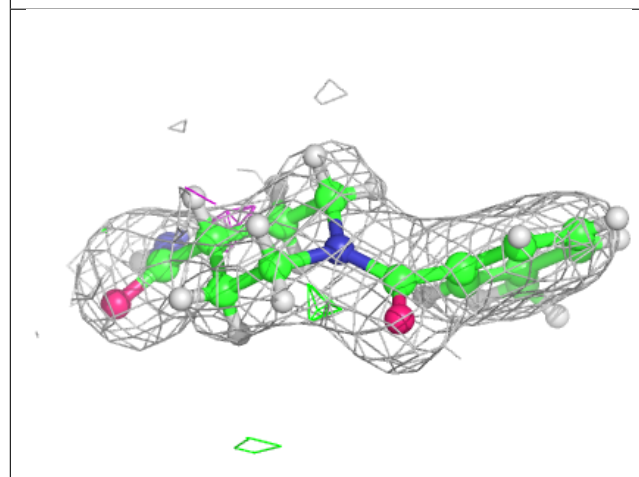
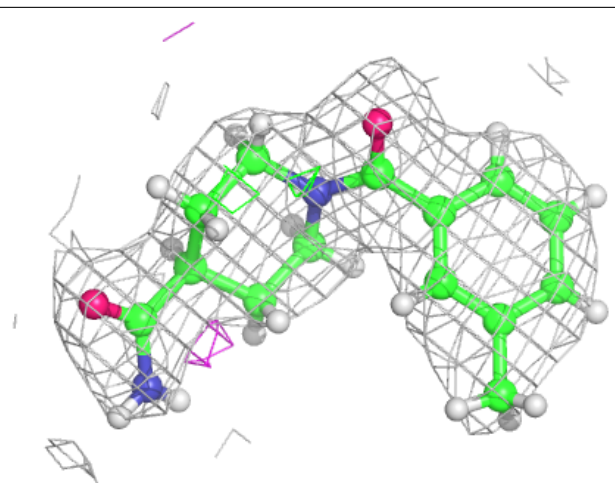
Electron density around ACP F 401:

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



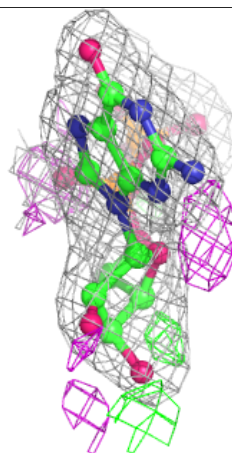
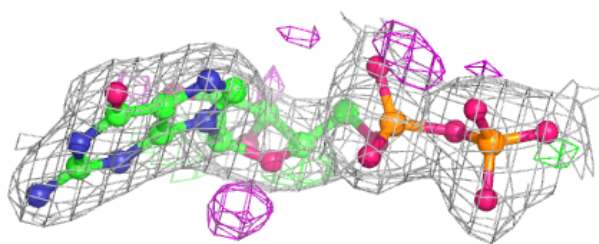
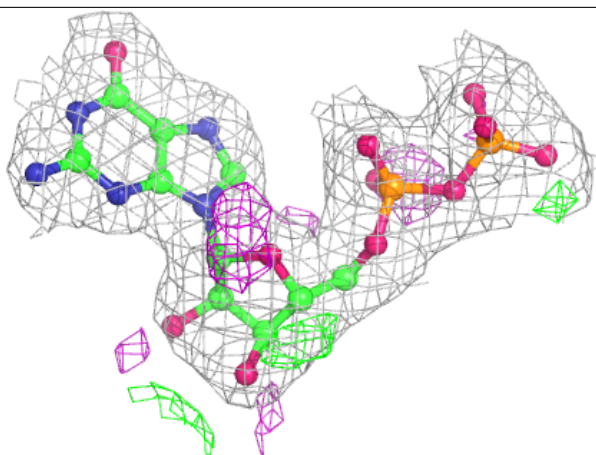
Electron density around NZJ B 505:

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

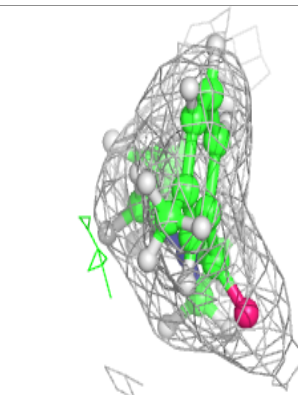
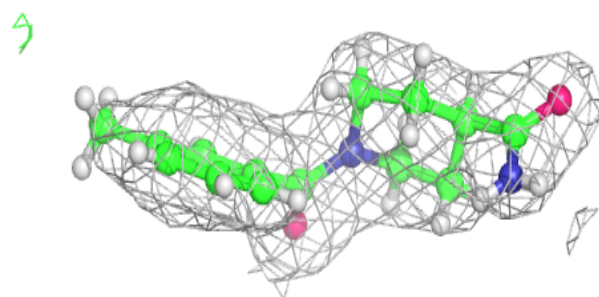
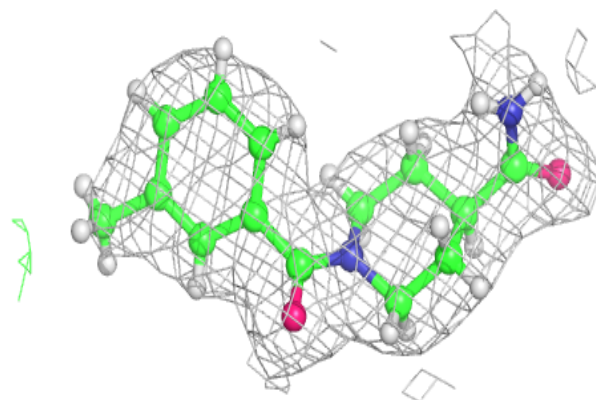


Electron density around GDP D 501:

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

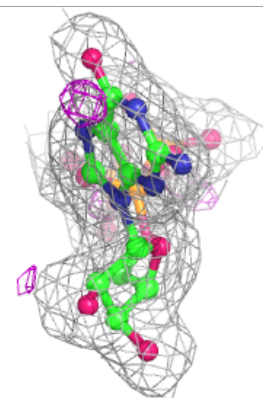
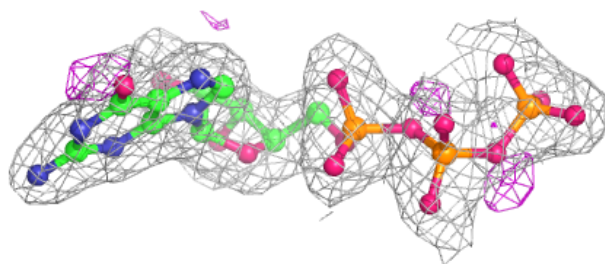
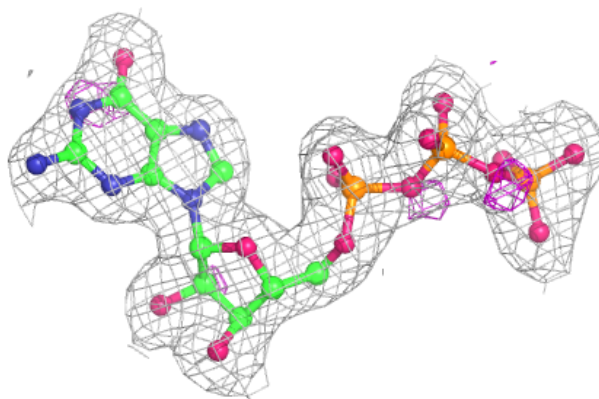
**Electron density around NZJ C 505:**

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

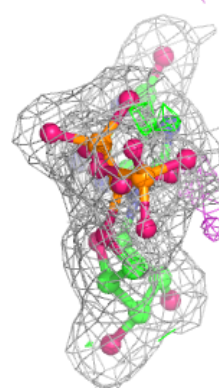
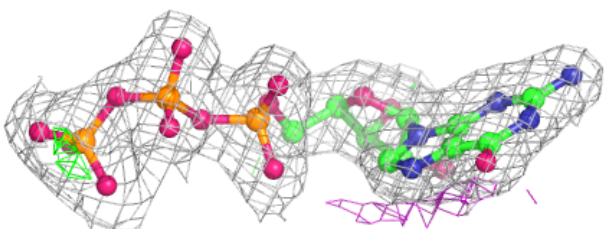
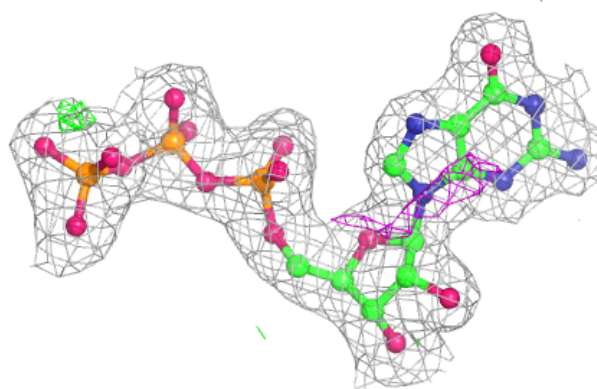


Electron density around GTP A 501:

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

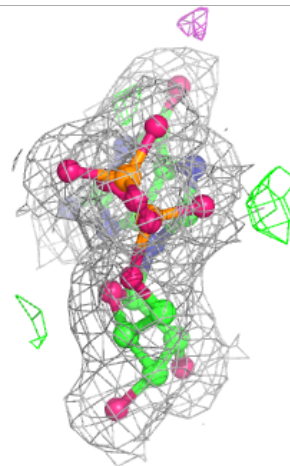
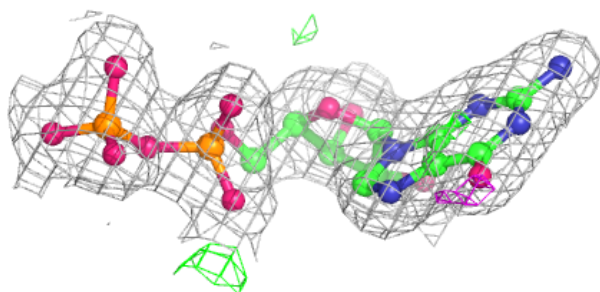
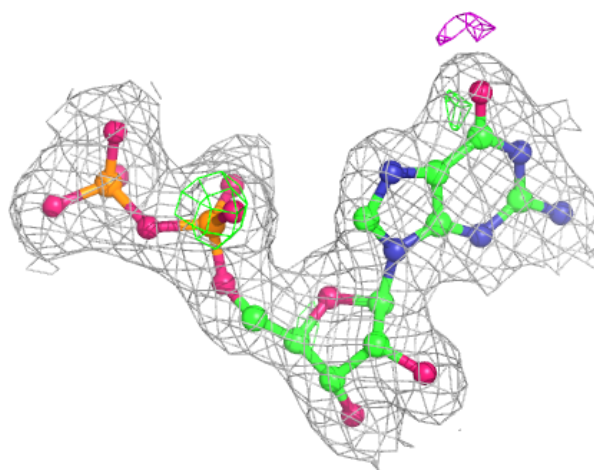
**Electron density around GTP C 501:**

$2mF_o-DF_c$ (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.