



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:49 PM UTC

PDB ID : 5S54 / pdb_00005s54
Title : Tubulin-Z2856434816-complex
Authors : Muehlethaler, T.; Gioia, D.; Protá, A.E.; Sharpe, M.E.; Cavalli, A.; Steinmetz, M.O.
Deposited on : 2020-11-08
Resolution : 2.40 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

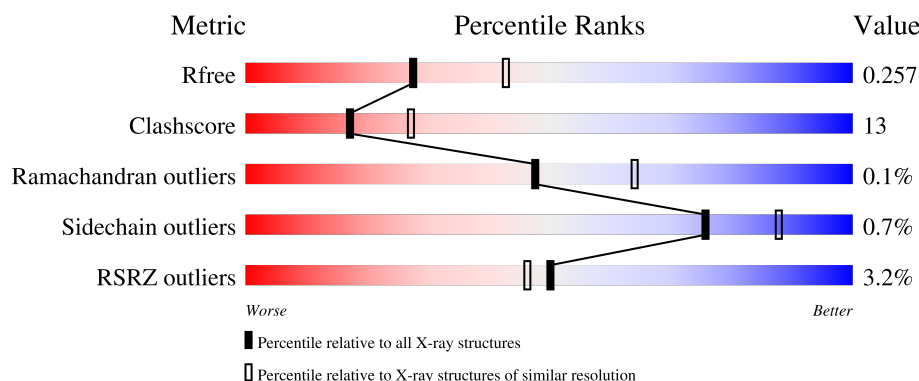
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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.



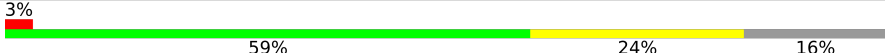
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	2% 69% 27% ..
1	C	451	2% 75% 23% .
2	B	445	3% 70% 25% 5%
2	D	445	3% 67% 29% .
3	E	143	8% 65% 20% 14%

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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: a small red segment at the beginning labeled '3%', followed by a large green segment labeled '59%', a yellow segment labeled '24%', and a small grey segment at the end labeled '16%'. The segments are separated by thin white lines.

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 17623 atoms, of which 12 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	438	Total	C	N	O	S	0	0	0
			3424	2167	582	653	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
			3339	2098	571	643	27			
2	D	426	Total	C	N	O	S	5	0	0
			3343	2098	570	648	27			

- Molecule 3 is a protein called Stathmin-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	123	Total	C	N	O	S	0	0	0
			1014	625	183	201	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

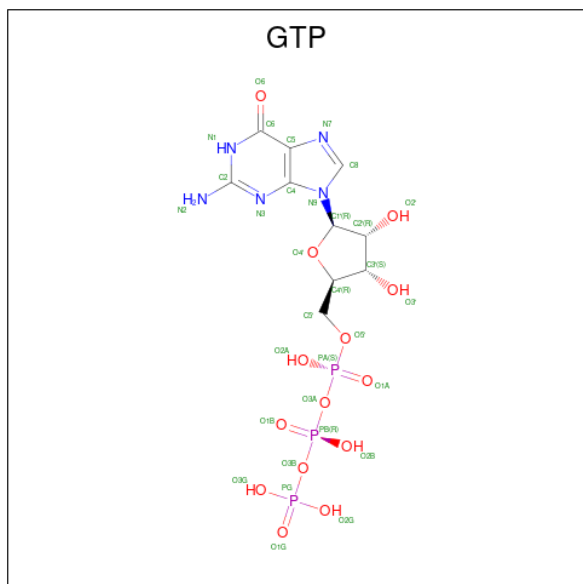
- Molecule 4 is a protein called Tubulin-Tyrosine Ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	322	Total	C	N	O	S	0	0	0
			2623	1688	447	476	12			

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: $C_{10}H_{16}N_5O_{14}P_3$).



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 (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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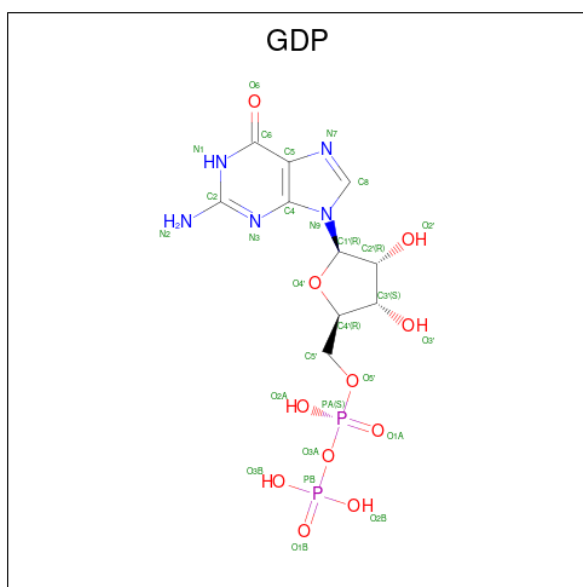
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	F	1	Total	Mg	0	0
			1	1		

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

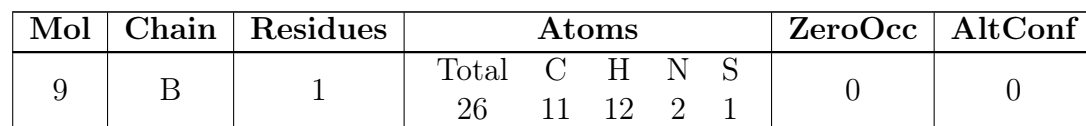
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	2	Total	Ca	0	0
			2	2		
7	B	1	Total	Ca	0	0
			1	1		
7	C	1	Total	Ca	0	0
			1	1		

- Molecule 8 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).



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

- Molecule 9 is 1-(pyridin-4-yl)-N-[(thiophen-2-yl)methyl]methanamine (CCD ID: WLS) (formula: C₁₁H₁₂N₂S) (labeled as "Ligand of Interest" by depositor).



- # ACP
-
- ORTEP diagram of the ACP molecule. The structure shows a pyrimidine ring system (N1, N3, N7, N8, C2, C4, C5, C6) linked to a ribose sugar (C1', C2', C3', C4', C5') via a glycosidic bond. The ribose sugar is further linked to a phosphate chain (P1, P2, P3) via a triphosphate group. The structure is labeled with atom names and displacement ellipsoids at the 50% probability level.

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

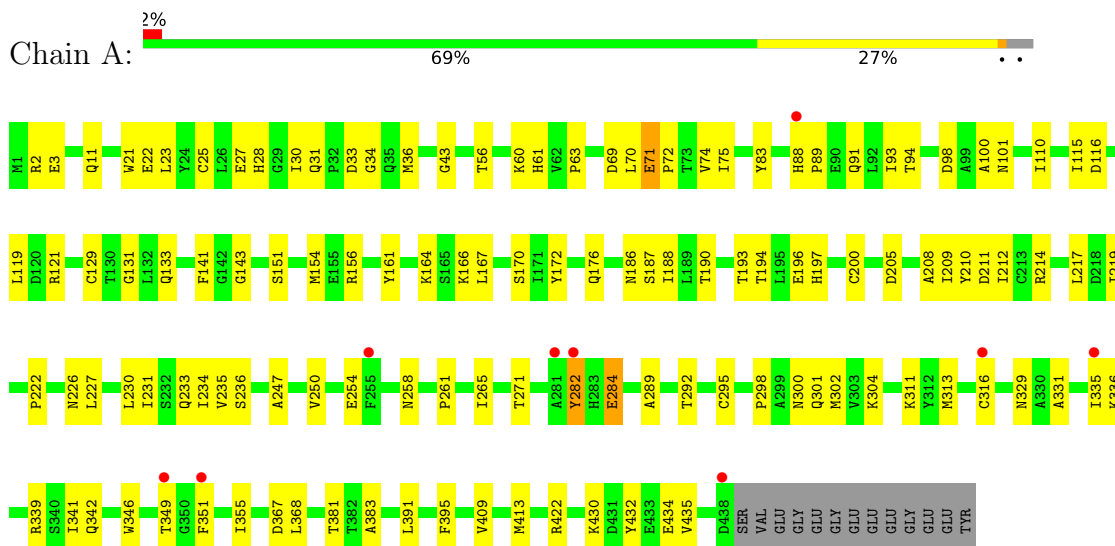
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	43	Total 43	O 43	0	0
11	B	48	Total 48	O 48	0	0
11	C	135	Total 135	O 135	0	0
11	D	17	Total 17	O 17	0	0
11	E	5	Total 5	O 5	0	0
11	F	3	Total 3	O 3	0	0

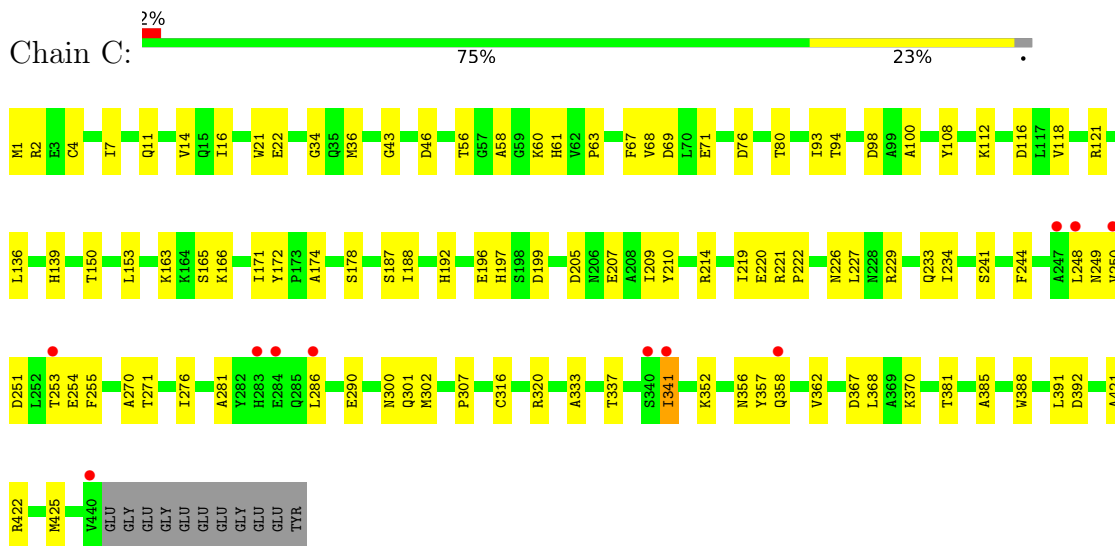
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Tubulin alpha-1B chain

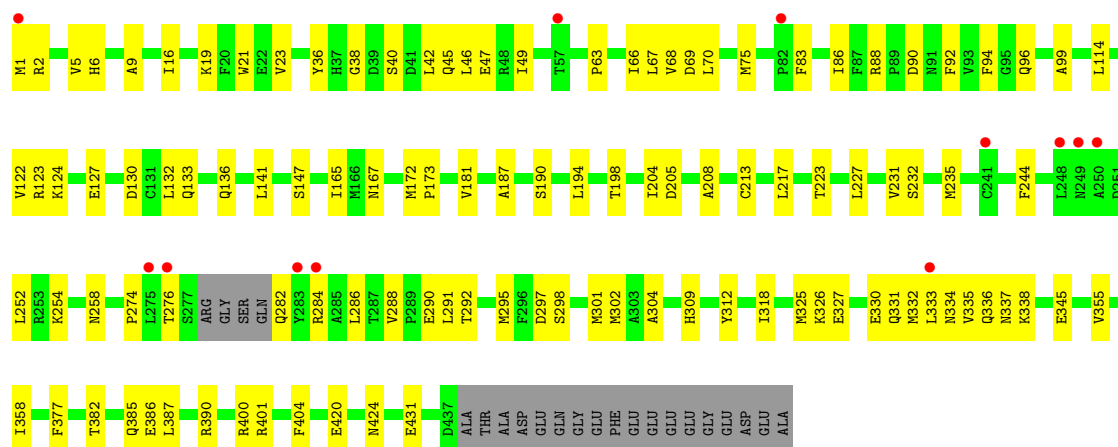


- Molecule 1: Tubulin alpha-1B chain

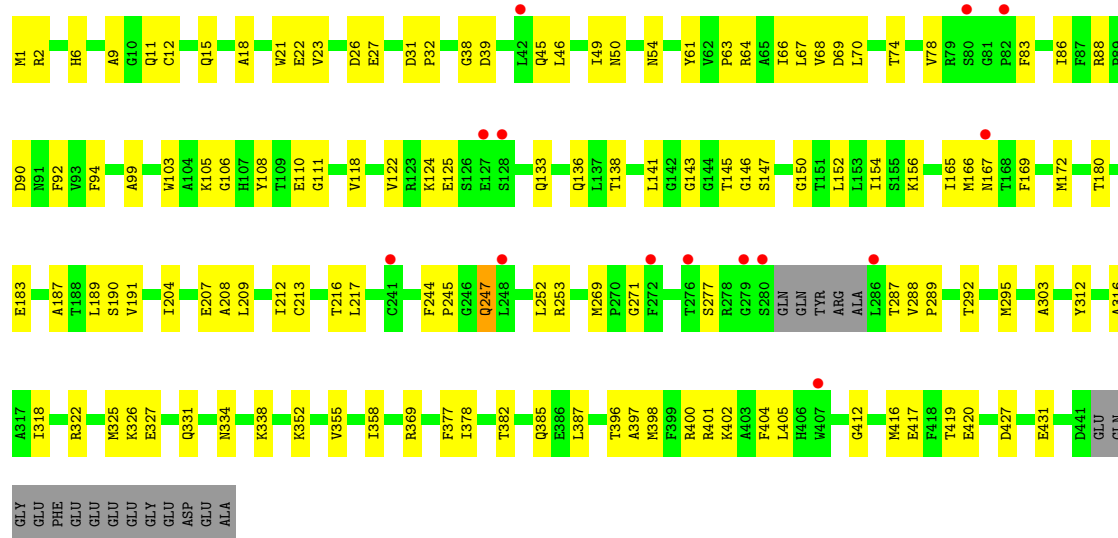


- Molecule 2: Tubulin beta-2B chain

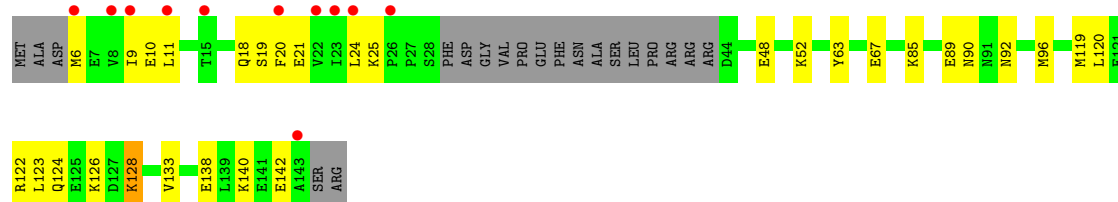




• Molecule 2: Tubulin beta-2B chain

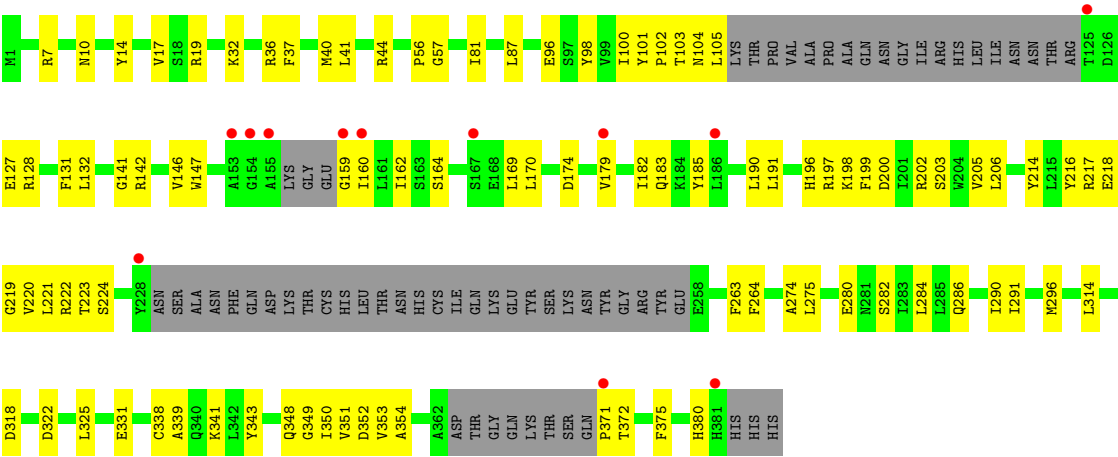


• Molecule 3: Stathmin-4



• Molecule 4: Tubulin-Tyrosine Ligase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	104.33Å 157.15Å 179.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	71.96 – 2.40 71.96 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.9 (71.96-2.40) 99.9 (71.96-2.40)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.15 (at 2.40Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.206 , 0.250 0.212 , 0.257	Depositor DCC
R_{free} test set	5756 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	68.5	Xtriage
Anisotropy	0.172	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 51.2	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	17623	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.83% 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: GTP, MG, GDP, ACP, CA, WLS

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	0.11	0/3502	0.30	0/4754
1	C	0.13	0/3521	0.32	0/4780
2	B	0.12	0/3413	0.29	0/4621
2	D	0.11	0/3416	0.28	0/4626
3	E	0.11	0/1022	0.25	0/1356
4	F	0.10	0/2682	0.30	0/3624
All	All	0.12	0/17556	0.29	0/23761

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3424	0	3334	98	0
1	C	3443	0	3352	75	0
2	B	3339	0	3214	90	0
2	D	3343	0	3222	91	0
3	E	1014	0	1029	27	0
4	F	2623	0	2609	70	0
5	A	32	0	12	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	32	0	12	1	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	1	0	0	0	0
8	B	28	0	12	0	0
8	D	28	0	12	3	0
9	B	14	12	0	1	0
10	F	31	0	14	3	0
11	A	43	0	0	0	0
11	B	48	0	0	4	0
11	C	135	0	0	6	0
11	D	17	0	0	1	0
11	E	5	0	0	0	0
11	F	3	0	0	0	0
All	All	17611	12	16822	429	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 13.

All (429) 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)
4:F:102:PRO:HG2	4:F:105:LEU:HD13	1.40	1.02
2:D:1:MET:HE2	2:D:50:ASN:HB2	1.49	0.94
1:A:336:LYS:HG3	3:E:24:LEU:HD13	1.51	0.90
2:D:334:ASN:HD21	2:D:338:LYS:HE3	1.35	0.89
1:C:93:ILE:HD11	1:C:121:ARG:HG3	1.56	0.88
2:B:274:PRO:HB3	2:B:286:LEU:HD22	1.57	0.85
2:D:217:LEU:HA	2:D:277:SER:HB3	1.58	0.84
2:D:83:PHE:O	2:D:86:ILE:HG22	1.80	0.82
2:B:424:ASN:HB3	11:B:636:HOH:O	1.81	0.81
4:F:331:GLU:OE2	10:F:401:ACP:O3G	1.98	0.81
2:D:397:ALA:O	2:D:401:ARG:NH1	2.13	0.81
1:A:209:ILE:HG22	1:A:227:LEU:HD22	1.64	0.78
2:B:96:GLN:HB3	1:C:1:MET:HE2	1.66	0.77
1:A:335:ILE:HG23	1:A:339:ARG:HG3	1.65	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:136:GLN:HA	2:D:167:ASN:O	1.86	0.75
1:C:76:ASP:O	1:C:80:THR:HG22	1.87	0.74
4:F:147:TRP:HB2	4:F:169:LEU:HD11	1.70	0.74
2:B:284:ARG:NH2	2:B:290:GLU:OE2	2.22	0.72
2:B:337:ASN:OD1	4:F:36:ARG:HD3	1.88	0.72
4:F:128:ARG:NH2	4:F:174:ASP:OD1	2.23	0.71
2:B:431:GLU:OE1	11:B:601:HOH:O	2.08	0.70
1:C:221:ARG:HG3	2:D:325:MET:HG2	1.73	0.70
8:D:501:GDP:O1A	11:D:601:HOH:O	2.09	0.70
2:D:295:MET:HE2	2:D:377:PHE:HB2	1.73	0.69
1:A:100:ALA:HA	2:B:254:LYS:HG3	1.74	0.69
1:A:271:THR:HG21	1:A:295:CYS:HA	1.74	0.69
2:D:397:ALA:HA	2:D:400:ARG:NH1	2.08	0.69
1:C:320:ARG:HA	1:C:356:ASN:O	1.93	0.68
1:A:381:THR:HG22	1:A:383:ALA:H	1.59	0.68
1:C:220:GLU:HB3	2:D:326:LYS:HD2	1.76	0.68
2:B:67:LEU:HD22	2:B:92:PHE:CE2	2.28	0.67
2:B:83:PHE:O	2:B:86:ILE:HG22	1.95	0.67
2:D:244:PHE:CE1	2:D:358:ILE:HD12	2.30	0.67
1:A:2:ARG:HB2	1:A:133:GLN:HE21	1.58	0.67
4:F:348:GLN:NE2	4:F:352:ASP:OD1	2.27	0.67
1:C:270:ALA:O	1:C:302:MET:HG2	1.94	0.67
2:D:334:ASN:ND2	2:D:338:LYS:HE3	2.10	0.67
1:A:166:LYS:HE2	1:A:197:HIS:O	1.94	0.66
1:C:209:ILE:HD11	1:C:302:MET:CE	2.24	0.66
1:C:209:ILE:HD11	1:C:302:MET:HE3	1.76	0.66
4:F:371:PRO:HA	4:F:372:THR:O	1.97	0.66
1:A:214:ARG:HG2	1:A:219:ILE:O	1.95	0.65
4:F:132:LEU:HD21	4:F:170:LEU:HD11	1.77	0.65
1:A:36:MET:HB3	1:A:61:HIS:CE1	2.32	0.65
2:B:172:MET:HG3	2:B:387:LEU:HD11	1.80	0.64
2:D:172:MET:HG3	2:D:387:LEU:HD11	1.79	0.64
3:E:48:GLU:HG2	3:E:52:LYS:HE3	1.79	0.64
4:F:102:PRO:HG2	4:F:105:LEU:CD1	2.23	0.64
1:C:4[A]:CYS:SG	1:C:136:LEU:HG	2.38	0.64
2:D:2:ARG:HB2	2:D:133:GLN:HE21	1.63	0.64
2:B:21:TRP:CZ3	2:B:63:PRO:HB3	2.33	0.64
1:C:234:ILE:HG12	1:C:302:MET:HE2	1.79	0.63
2:D:427:ASP:O	2:D:431:GLU:HG3	1.98	0.63
2:B:345:GLU:OE1	2:B:345:GLU:N	2.23	0.63
2:D:2:ARG:HB3	2:D:133:GLN:CG	2.28	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43:GLY:HA2	1:A:56:THR:O	1.99	0.63
4:F:371:PRO:HA	4:F:372:THR:C	2.24	0.63
1:A:88:HIS:CD2	1:A:91:GLN:HG3	2.34	0.62
1:C:188:ILE:HG13	1:C:425:MET:HG3	1.80	0.62
4:F:146:VAL:HG22	4:F:164:SER:HB3	1.80	0.62
2:D:46:LEU:HA	2:D:49:ILE:HB	1.80	0.62
1:A:34:GLY:HA3	1:A:60:LYS:HG3	1.82	0.62
2:D:23:VAL:O	2:D:27:GLU:HG3	2.00	0.62
1:C:214:ARG:HG2	1:C:219:ILE:O	2.00	0.62
4:F:314:LEU:HD22	4:F:350:ILE:HD11	1.81	0.61
4:F:14:TYR:HB3	4:F:41:LEU:HD13	1.80	0.61
1:C:165:SER:HA	1:C:199:ASP:OD2	2.01	0.61
4:F:205:VAL:HG21	4:F:291:ILE:HD13	1.81	0.61
4:F:10:ASN:HB2	4:F:44:ARG:HH22	1.65	0.61
2:B:147:SER:HG	2:B:190:SER:HG	1.47	0.61
4:F:198:LYS:HG2	4:F:199:PHE:H	1.66	0.60
1:A:188:ILE:HD12	1:A:395:PHE:CD2	2.37	0.60
4:F:349:GLY:O	4:F:353:VAL:HG22	2.01	0.60
1:C:1:MET:HE3	1:C:2:ARG:HH11	1.66	0.59
2:B:288:VAL:HG12	2:B:331:GLN:HG3	1.84	0.59
4:F:159:GLY:C	4:F:160:ILE:HD12	2.27	0.59
2:B:123:ARG:O	2:B:127:GLU:HG3	2.02	0.59
2:B:292:THR:HG22	2:B:335:VAL:HG21	1.84	0.59
2:B:431:GLU:HB2	11:B:601:HOH:O	2.02	0.59
1:A:329:ASN:HB3	3:E:6:MET:HE3	1.84	0.59
2:D:152:LEU:O	2:D:156:LYS:HG2	2.02	0.58
1:A:151:SER:HB2	1:A:193:THR:OG1	2.04	0.58
1:A:351:PHE:HE1	3:E:24:LEU:HD11	1.68	0.58
2:D:124:LYS:C	2:D:124:LYS:HD3	2.28	0.58
2:D:1:MET:CE	2:D:50:ASN:HB2	2.29	0.58
1:A:265:ILE:HG23	1:A:432:TYR:CZ	2.38	0.58
2:B:295:MET:CG	2:B:377:PHE:HB2	2.33	0.58
1:C:251:ASP:OD2	1:C:253:THR:OG1	2.21	0.58
2:B:2:ARG:HB2	2:B:133:GLN:HE21	1.69	0.58
4:F:371:PRO:CA	4:F:372:THR:HB	2.34	0.57
1:A:23:LEU:HD23	1:A:236:SER:HB2	1.87	0.57
1:A:292:THR:HG22	1:A:335:ILE:HD12	1.87	0.57
1:C:229:ARG:NH1	11:C:608:HOH:O	2.36	0.57
3:E:85:LYS:O	3:E:89:GLU:HG3	2.04	0.57
1:A:21:TRP:CZ3	1:A:63:PRO:HB3	2.39	0.57
3:E:10:GLU:O	3:E:20:PHE:HA	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:274:ALA:HB3	4:F:275:LEU:HD22	1.87	0.56
2:B:69:ASP:O	2:B:94:PHE:HA	2.04	0.56
2:B:96:GLN:CB	1:C:1:MET:HE2	2.35	0.56
2:B:276:THR:HG21	2:B:282:GLN:HA	1.85	0.56
2:D:26:ASP:OD2	2:D:369:ARG:HD2	2.04	0.56
1:A:250:VAL:HG12	1:A:254:GLU:OE1	2.06	0.56
2:D:12:CYS:HB2	8:D:501:GDP:C8	2.40	0.56
1:A:430:LYS:HE2	1:A:434:GLU:OE2	2.06	0.56
2:B:136:GLN:HA	2:B:167:ASN:O	2.06	0.56
2:B:194:LEU:HD22	2:B:198:THR:HG21	1.87	0.56
1:A:2:ARG:CB	1:A:133:GLN:HG3	2.35	0.56
1:C:196:GLU:HG2	11:C:628:HOH:O	2.05	0.55
3:E:119:MET:HA	3:E:122:ARG:NH2	2.21	0.55
2:B:141:LEU:HD12	2:B:172:MET:SD	2.47	0.55
2:D:269:MET:HG3	2:D:303:ALA:HB3	1.87	0.55
1:C:254:GLU:HG2	1:C:352:LYS:HE2	1.89	0.55
2:D:208:ALA:O	2:D:212:ILE:HG13	2.06	0.55
1:A:265:ILE:HG23	1:A:432:TYR:CE1	2.42	0.55
1:C:209:ILE:HG22	1:C:227:LEU:HD22	1.88	0.55
2:B:301:MET:HE1	2:B:312:TYR:OH	2.06	0.55
1:C:21:TRP:CZ3	1:C:63:PRO:HB3	2.42	0.55
1:C:174:ALA:HB1	1:C:207:GLU:HB2	1.89	0.54
1:A:71:GLU:HG2	1:A:72:PRO:HD2	1.90	0.54
2:D:22:GLU:HG2	2:D:83:PHE:CD1	2.42	0.54
2:D:154:ILE:HG23	2:D:166:MET:HG2	1.89	0.54
1:A:302:MET:N	1:A:302:MET:HE2	2.22	0.54
1:C:34:GLY:HA3	1:C:60:LYS:HG3	1.90	0.54
3:E:120:LEU:O	3:E:124:GLN:HG3	2.07	0.54
1:C:163:LYS:HG3	3:E:90:ASN:OD1	2.08	0.54
2:B:66:ILE:HD12	2:B:122:VAL:HG22	1.90	0.54
2:D:396:THR:O	2:D:400:ARG:HG2	2.08	0.54
1:A:2:ARG:HB3	1:A:133:GLN:HG3	1.89	0.53
1:C:286:LEU:HA	1:C:290:GLU:OE1	2.08	0.53
1:A:289:ALA:HA	1:A:331:ALA:CB	2.38	0.53
2:B:19:LYS:HB3	2:B:232:SER:OG	2.08	0.53
2:B:165:ILE:HG21	2:B:252:LEU:HB3	1.89	0.53
1:A:93:ILE:HD11	1:A:121:ARG:HG3	1.90	0.53
2:D:382:THR:O	2:D:385:GLN:HG2	2.07	0.53
1:A:172:TYR:HB3	1:A:205:ASP:HA	1.89	0.53
1:A:271:THR:CG2	1:A:295:CYS:HA	2.39	0.53
2:B:205:ASP:OD2	2:B:390:ARG:NH2	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:96:GLU:O	4:F:183:GLN:HA	2.08	0.53
10:F:401:ACP:O3G	10:F:401:ACP:O1B	2.26	0.53
4:F:216:TYR:CE2	4:F:218:GLU:HB2	2.43	0.53
1:A:187:SER:HB3	1:A:391:LEU:HD21	1.91	0.53
1:A:211:ASP:OD2	1:A:304:LYS:NZ	2.32	0.53
2:B:16:ILE:HD13	2:B:231:VAL:HG11	1.91	0.53
2:B:331:GLN:O	2:B:335:VAL:HG23	2.09	0.53
2:B:333:LEU:HD13	4:F:57:GLY:HA3	1.90	0.52
1:A:209:ILE:HG22	1:A:227:LEU:CD2	2.37	0.52
2:B:40:SER:OG	2:B:42:LEU:HD13	2.08	0.52
2:B:124:LYS:HD3	2:B:124:LYS:C	2.35	0.52
1:C:187:SER:HB3	1:C:391:LEU:HD21	1.91	0.52
2:D:9:ALA:HA	2:D:68:VAL:O	2.10	0.52
3:E:128:LYS:O	3:E:128:LYS:HD3	2.08	0.52
1:A:98:ASP:HB2	5:A:501:GTP:O2G	2.10	0.52
1:A:210:TYR:CE1	1:A:222:PRO:HD2	2.44	0.52
2:D:69:ASP:O	2:D:94:PHE:HA	2.10	0.52
2:D:118:VAL:O	2:D:122:VAL:HG23	2.10	0.52
1:C:249:ASN:OD1	1:C:356:ASN:ND2	2.40	0.52
2:D:105:LYS:HB3	2:D:110:GLU:HG2	1.92	0.52
1:A:335:ILE:CG2	1:A:339:ARG:HG3	2.39	0.52
2:B:2:ARG:HB2	2:B:133:GLN:NE2	2.25	0.52
2:D:145:THR:HB	8:D:501:GDP:O2B	2.09	0.52
2:D:67:LEU:HD22	2:D:92:PHE:CE2	2.45	0.52
1:A:101:ASN:HD22	2:B:258:ASN:HD21	1.58	0.51
2:D:26:ASP:CG	2:D:369:ARG:HD2	2.35	0.51
2:B:420:GLU:HB3	11:B:630:HOH:O	2.10	0.51
3:E:48:GLU:CG	3:E:52:LYS:HE3	2.40	0.51
4:F:286:GLN:O	4:F:290:ILE:HG13	2.10	0.51
1:A:176:GLN:CG	4:F:56:PRO:HB3	2.40	0.51
4:F:199:PHE:CB	4:F:223:THR:HG22	2.41	0.51
2:B:274:PRO:HB3	2:B:286:LEU:CD2	2.34	0.51
4:F:37:PHE:CZ	4:F:40:MET:HE3	2.44	0.51
4:F:205:VAL:CG2	4:F:291:ILE:HD13	2.40	0.51
1:A:247:ALA:HB3	3:E:19:SER:OG	2.11	0.51
1:A:298:PRO:HA	1:A:301:GLN:CD	2.35	0.51
1:A:69:ASP:O	1:A:94:THR:HA	2.11	0.51
1:A:196:GLU:OE1	1:A:196:GLU:HA	2.10	0.51
3:E:11:LEU:HD11	3:E:18:GLN:OE1	2.11	0.51
1:C:172:TYR:CE2	1:C:391:LEU:HD22	2.46	0.50
2:D:18:ALA:O	2:D:22:GLU:HG3	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:295:MET:HG3	2:B:377:PHE:HB2	1.94	0.50
4:F:341:LYS:HG2	4:F:341:LYS:O	2.12	0.50
2:B:88:ARG:NH1	2:B:90:ASP:HB2	2.26	0.50
2:B:204:ILE:HD13	2:B:231:VAL:HG13	1.92	0.50
2:D:2:ARG:HB3	2:D:133:GLN:HG3	1.92	0.50
1:C:166:LYS:HE2	1:C:197:HIS:O	2.11	0.50
1:C:226:ASN:ND2	1:C:367:ASP:OD2	2.44	0.50
4:F:17:VAL:HG13	4:F:351:VAL:HG22	1.94	0.50
1:C:233:GLN:HG3	1:C:368:LEU:CD1	2.41	0.50
2:B:6:HIS:CD2	2:B:21:TRP:HE1	2.29	0.50
2:D:6:HIS:CD2	2:D:21:TRP:HE1	2.29	0.50
2:D:67:LEU:N	2:D:67:LEU:HD12	2.27	0.50
2:D:141:LEU:HD12	2:D:172:MET:SD	2.51	0.50
2:D:316:ALA:HB3	2:D:378:ILE:HB	1.94	0.50
4:F:220:VAL:HG12	4:F:263:PHE:CE1	2.47	0.50
2:B:38:GLY:HA3	2:B:45:GLN:OE1	2.11	0.50
4:F:81:ILE:HA	4:F:87:LEU:HD12	1.93	0.50
2:B:325:MET:HE2	2:B:355:VAL:HB	1.92	0.50
2:D:70:LEU:HD12	2:D:99:ALA:HB2	1.94	0.49
2:D:288:VAL:HG12	2:D:331:GLN:HG3	1.93	0.49
4:F:100:ILE:HD12	4:F:128:ARG:HA	1.93	0.49
4:F:371:PRO:N	4:F:372:THR:HB	2.27	0.49
1:A:187:SER:CB	1:A:391:LEU:HD21	2.42	0.49
2:D:106:GLY:O	2:D:111:GLY:HA3	2.12	0.49
1:A:119:LEU:HD12	1:A:156:ARG:NH2	2.27	0.49
1:A:335:ILE:HG23	1:A:339:ARG:CG	2.39	0.49
1:A:355:ILE:HB	3:E:18:GLN:O	2.13	0.49
1:C:341:ILE:HD12	1:C:341:ILE:O	2.12	0.49
1:C:139:HIS:CD2	1:C:150:THR:HG21	2.47	0.49
2:D:180:THR:HB	2:D:183:GLU:HG3	1.93	0.49
3:E:138:GLU:O	3:E:142:GLU:HG3	2.13	0.49
4:F:198:LYS:HG2	4:F:199:PHE:N	2.27	0.49
2:B:327:GLU:O	2:B:331:GLN:HG2	2.13	0.49
4:F:199:PHE:HB2	4:F:223:THR:HG22	1.95	0.49
2:B:165:ILE:HD13	9:B:504:WLS:C11	2.43	0.49
2:D:398:MET:HE3	2:D:404:PHE:HD2	1.77	0.49
1:A:88:HIS:HB2	1:A:89:PRO:HD2	1.94	0.49
1:A:161:TYR:HB3	1:A:164:LYS:HD3	1.94	0.49
1:A:289:ALA:HA	1:A:331:ALA:HB1	1.94	0.48
1:C:22:GLU:OE1	11:C:601:HOH:O	2.20	0.48
1:A:11:GLN:HG3	1:A:74:VAL:HG21	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:GLN:HG3	4:F:56:PRO:HB3	1.95	0.48
2:B:297:ASP:OD1	2:B:298:SER:N	2.46	0.48
2:D:1:MET:HE2	2:D:50:ASN:CB	2.34	0.48
4:F:103:THR:HG23	4:F:128:ARG:NH2	2.27	0.48
1:C:11:GLN:HE22	2:D:247:GLN:NE2	2.10	0.48
1:C:248:LEU:HD12	1:C:357:TYR:OH	2.14	0.48
2:D:74:THR:O	2:D:78:VAL:HG23	2.14	0.48
2:D:103:TRP:CE3	2:D:189:LEU:HD13	2.47	0.48
1:C:36:MET:HB3	1:C:61:HIS:CE1	2.48	0.48
1:A:301:GLN:C	1:A:302:MET:HE2	2.38	0.48
1:C:43:GLY:HA2	1:C:56:THR:O	2.14	0.48
1:C:108:TYR:O	1:C:112:LYS:HG2	2.14	0.48
1:A:292:THR:HG22	1:A:335:ILE:CD1	2.43	0.48
4:F:10:ASN:CB	4:F:44:ARG:HH22	2.26	0.48
2:B:332:MET:O	2:B:336:GLN:HG3	2.12	0.48
2:D:387:LEU:HD23	2:D:387:LEU:C	2.39	0.48
1:A:351:PHE:HE1	3:E:24:LEU:CD1	2.27	0.48
2:B:325:MET:HE2	2:B:355:VAL:CB	2.42	0.48
3:E:123:LEU:HD23	3:E:126:LYS:HD3	1.96	0.47
4:F:199:PHE:CD1	4:F:221:LEU:HD23	2.49	0.47
2:B:286:LEU:HD12	2:B:290:GLU:OE1	2.14	0.47
4:F:191:LEU:HD12	4:F:196:HIS:CE1	2.48	0.47
1:C:16:ILE:CD1	1:C:171:ILE:HD11	2.44	0.47
2:D:180:THR:O	2:D:183:GLU:HG3	2.14	0.47
1:C:7:ILE:HG21	1:C:153:LEU:HD21	1.95	0.47
4:F:371:PRO:HA	4:F:372:THR:HB	1.95	0.47
1:A:188:ILE:HD12	1:A:395:PHE:HB2	1.96	0.47
1:A:209:ILE:HG23	1:A:230:LEU:HD23	1.94	0.47
2:B:286:LEU:HD23	2:B:291:LEU:CD2	2.45	0.47
2:B:420:GLU:O	2:B:424:ASN:ND2	2.48	0.47
1:A:188:ILE:HD12	1:A:395:PHE:CG	2.50	0.47
1:A:226:ASN:ND2	1:A:367:ASP:OD2	2.44	0.47
1:A:311:LYS:HA	1:A:342:GLN:O	2.15	0.47
1:C:21:TRP:CE3	1:C:63:PRO:HB3	2.49	0.47
1:C:93:ILE:HD11	1:C:121:ARG:CG	2.37	0.47
2:D:63:PRO:HD3	2:D:86:ILE:HG12	1.97	0.47
2:D:412:GLY:C	3:E:133:VAL:HG13	2.40	0.47
4:F:338:CYS:HB3	4:F:343:TYR:CE1	2.50	0.47
4:F:318:ASP:OD2	10:F:401:ACP:O2G	2.33	0.47
2:D:147:SER:HB2	2:D:190:SER:OG	2.14	0.47
1:A:231:ILE:O	1:A:235:VAL:HG23	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:46:ASP:OD1	1:C:46:ASP:N	2.46	0.47
2:B:2:ARG:HB2	2:B:133:GLN:CG	2.45	0.46
1:A:336:LYS:HG3	3:E:24:LEU:CD1	2.36	0.46
1:C:174:ALA:CB	1:C:207:GLU:HB2	2.45	0.46
4:F:104:ASN:C	4:F:105:LEU:HD12	2.40	0.46
1:A:31:GLN:HB2	1:A:33:ASP:OD1	2.15	0.46
1:A:71:GLU:HG2	1:A:72:PRO:CD	2.45	0.46
2:D:2:ARG:CB	2:D:133:GLN:HE21	2.28	0.46
1:A:3:GLU:OE1	1:A:129:CYS:HB3	2.15	0.46
2:B:147:SER:OG	2:B:190:SER:OG	2.21	0.46
2:B:66:ILE:CD1	2:B:122:VAL:HG22	2.45	0.46
2:B:325:MET:HE2	2:B:355:VAL:HG21	1.97	0.46
3:E:92:ASN:O	3:E:96:MET:HG2	2.14	0.46
4:F:162:ILE:HD13	4:F:185:TYR:CE1	2.50	0.46
4:F:202:ARG:HD3	4:F:203:SER:N	2.31	0.46
2:B:1:MET:HB2	2:B:130:ASP:OD2	2.15	0.46
2:B:114:LEU:O	2:B:114:LEU:HG	2.16	0.46
1:C:250:VAL:HG11	1:C:352:LYS:HE3	1.98	0.46
2:D:402:LYS:HB3	2:D:405:LEU:HD12	1.98	0.46
1:A:234:ILE:HD12	1:A:234:ILE:N	2.31	0.46
2:B:46:LEU:HA	2:B:49:ILE:HB	1.97	0.46
2:D:108:TYR:OH	2:D:417:GLU:OE2	2.22	0.46
1:A:209:ILE:CG2	1:A:227:LEU:HD22	2.42	0.45
1:A:395:PHE:CD1	1:A:422:ARG:HD3	2.50	0.45
2:D:416:MET:O	2:D:420:GLU:HG3	2.16	0.45
2:D:11:GLN:O	2:D:15:GLN:HG2	2.16	0.45
1:A:23:LEU:O	1:A:27:GLU:HG3	2.16	0.45
2:B:75:MET:HE3	2:B:92:PHE:CD2	2.52	0.45
1:C:100:ALA:CB	2:D:253:ARG:HD2	2.46	0.45
1:A:25:CYS:HB3	1:A:30:ILE:O	2.16	0.45
2:B:23:VAL:HG21	2:B:232:SER:HB3	1.99	0.45
1:C:172:TYR:HB3	1:C:205:ASP:HA	1.98	0.45
1:C:244:PHE:CD1	1:C:358:GLN:HG2	2.52	0.45
2:B:181:VAL:HG21	2:B:404:PHE:CZ	2.52	0.45
2:B:387:LEU:C	2:B:387:LEU:HD23	2.42	0.45
1:C:14:VAL:HG13	1:C:67:PHE:HD2	1.81	0.45
2:B:5:VAL:HG23	2:B:132:LEU:HD11	1.99	0.45
2:B:231:VAL:O	2:B:235:MET:HG3	2.17	0.45
1:C:116:ASP:HB2	11:C:671:HOH:O	2.15	0.45
1:A:188:ILE:HD12	1:A:395:PHE:CB	2.47	0.45
1:A:227:LEU:O	1:A:231:ILE:HG13	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:64:ARG:HG3	2:D:125:GLU:OE1	2.16	0.45
3:E:92:ASN:OD1	3:E:96:MET:HE2	2.16	0.45
4:F:217:ARG:HG3	4:F:218:GLU:HG2	1.98	0.45
4:F:219:GLY:HA3	4:F:264:PHE:CZ	2.52	0.45
1:A:154:MET:HG3	1:A:194:THR:HG23	1.99	0.44
2:D:88:ARG:NH1	2:D:90:ASP:HB2	2.32	0.44
2:D:146:GLY:O	2:D:150:GLY:HA3	2.17	0.44
4:F:214:TYR:HB3	4:F:375:PHE:HB3	1.99	0.44
2:B:67:LEU:N	2:B:67:LEU:HD12	2.31	0.44
2:D:292:THR:O	2:D:295:MET:HG2	2.17	0.44
4:F:197:ARG:HB2	4:F:224:SER:O	2.17	0.44
4:F:280:GLU:OE1	4:F:284:LEU:HD23	2.17	0.44
1:C:234:ILE:CG1	1:C:302:MET:HE2	2.45	0.44
2:D:31:ASP:HB2	2:D:32:PRO:CD	2.47	0.44
4:F:101:TYR:CD2	4:F:179:VAL:HG22	2.52	0.44
4:F:206:LEU:HD21	4:F:354:ALA:HB2	1.99	0.44
1:A:284:GLU:CD	1:A:284:GLU:H	2.24	0.44
1:C:93:ILE:CD1	1:C:121:ARG:HG3	2.37	0.44
2:D:31:ASP:HB2	2:D:32:PRO:HD2	1.99	0.44
1:A:265:ILE:HG21	1:A:313:MET:HE1	1.99	0.44
1:C:234:ILE:HD12	1:C:234:ILE:H	1.83	0.44
2:D:318:ILE:N	2:D:318:ILE:HD12	2.33	0.44
2:B:208:ALA:HB2	2:B:304:ALA:HB2	2.00	0.43
4:F:19:ARG:HD2	4:F:19:ARG:O	2.17	0.43
1:A:346:TRP:HZ2	1:A:435:VAL:HG13	1.81	0.43
2:B:42:LEU:N	2:B:42:LEU:HD12	2.33	0.43
1:A:409:VAL:HA	1:A:413:MET:O	2.18	0.43
2:B:46:LEU:O	2:B:49:ILE:HG22	2.18	0.43
1:C:69:ASP:O	1:C:94:THR:HA	2.18	0.43
1:C:241:SER:HA	1:C:249:ASN:OD1	2.18	0.43
2:B:309:HIS:HD1	2:B:386:GLU:CD	2.26	0.43
1:C:178:SER:OG	2:D:352:LYS:NZ	2.52	0.43
1:A:71:GLU:HG2	1:A:72:PRO:N	2.32	0.43
2:B:21:TRP:CE3	2:B:63:PRO:HB3	2.53	0.43
2:B:244:PHE:CD1	2:B:358:ILE:HD12	2.53	0.43
1:C:255:PHE:CE1	1:C:352:LYS:HG2	2.54	0.43
1:C:255:PHE:CD1	1:C:316:CYS:HB3	2.54	0.43
4:F:141:GLY:O	4:F:142:ARG:HB2	2.18	0.43
1:A:217:LEU:HD21	1:A:368:LEU:HD23	2.00	0.43
2:D:204:ILE:HG22	2:D:209:LEU:HD11	2.01	0.43
4:F:131:PHE:CE1	4:F:182:ILE:HG21	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:286:LEU:HD23	2:B:291:LEU:HD23	2.01	0.43
2:D:38:GLY:HA3	2:D:45:GLN:OE1	2.18	0.43
1:A:167:LEU:HG	1:A:200:CYS:HB3	2.00	0.43
1:A:208:ALA:O	1:A:212:ILE:HG13	2.19	0.43
1:A:233:GLN:HG3	1:A:368:LEU:HD12	2.00	0.43
2:D:66:ILE:HD12	2:D:122:VAL:HG22	2.00	0.43
1:C:229:ARG:NE	11:C:601:HOH:O	2.25	0.43
4:F:282:SER:HB2	4:F:325:LEU:HD13	1.99	0.43
1:A:141:PHE:CE1	1:A:170:SER:HB3	2.54	0.42
2:B:2:ARG:CB	2:B:133:GLN:HE21	2.31	0.42
2:D:165:ILE:HG21	2:D:252:LEU:HB3	2.00	0.42
1:A:254:GLU:HG2	1:A:258:ASN:ND2	2.34	0.42
1:A:329:ASN:HB3	3:E:6:MET:CE	2.48	0.42
2:B:382:THR:O	2:B:385:GLN:HG2	2.20	0.42
2:D:46:LEU:O	2:D:49:ILE:HG22	2.20	0.42
1:A:22:GLU:HG3	1:A:83:TYR:CE1	2.54	0.42
1:A:70:LEU:HD13	1:A:110:ILE:HG21	2.01	0.42
1:A:186:ASN:O	1:A:190:THR:HG22	2.20	0.42
2:B:325:MET:HE2	2:B:355:VAL:CG2	2.49	0.42
1:C:276:ILE:HD13	1:C:281:ALA:HB2	2.01	0.42
4:F:200:ASP:OD1	4:F:222:ARG:HB2	2.19	0.42
4:F:280:GLU:HA	4:F:284:LEU:HB3	2.01	0.42
1:A:2:ARG:HB2	1:A:133:GLN:HG3	2.01	0.42
1:A:75:ILE:HB	1:A:94:THR:CG2	2.49	0.42
2:B:318:ILE:N	2:B:318:ILE:HD12	2.35	0.42
1:C:71:GLU:HG2	1:C:98:ASP:HB3	2.00	0.42
1:C:362:VAL:HG22	1:C:370:LYS:HD3	2.00	0.42
2:D:287:THR:HB	2:D:289:PRO:HD2	2.00	0.42
4:F:263:PHE:CZ	4:F:341:LYS:HE2	2.54	0.42
1:A:336:LYS:HD2	1:A:341:ILE:HD12	2.01	0.42
1:C:385:ALA:HA	1:C:388:TRP:CD1	2.55	0.42
1:A:2:ARG:HB3	1:A:131:GLY:O	2.20	0.42
2:B:334:ASN:OD1	2:B:338:LYS:HD3	2.19	0.42
1:C:210:TYR:CZ	1:C:222:PRO:HD2	2.54	0.42
1:C:333:ALA:O	1:C:337:THR:HG23	2.20	0.42
2:D:295:MET:CE	2:D:377:PHE:HB2	2.47	0.42
4:F:220:VAL:HG11	4:F:339:ALA:HB2	2.01	0.42
1:A:28:HIS:O	1:A:36:MET:HE3	2.20	0.42
2:B:325:MET:HE3	2:B:325:MET:HB2	1.97	0.42
2:D:212:ILE:O	2:D:216:THR:HB	2.20	0.42
2:D:322:ARG:NH1	2:D:322:ARG:HB3	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:68:VAL:HG11	1:C:118:VAL:HG21	2.01	0.42
1:A:143:GLY:HA3	5:A:501:GTP:O3A	2.20	0.42
2:B:5:VAL:HG23	2:B:132:LEU:CD1	2.50	0.42
2:B:70:LEU:HD12	2:B:99:ALA:HB2	2.01	0.42
2:D:213:CYS:HA	2:D:217:LEU:HB2	2.02	0.42
2:B:400:ARG:HG3	2:B:401:ARG:HG2	2.01	0.41
1:C:271:THR:HG23	1:C:300:ASN:O	2.20	0.41
2:D:271:GLY:HA3	2:D:377:PHE:HB3	2.02	0.41
2:B:326:LYS:O	2:B:330:GLU:HG3	2.19	0.41
4:F:296:MET:SD	4:F:380:HIS:HB2	2.60	0.41
2:B:213:CYS:HA	2:B:217:LEU:HB2	2.03	0.41
1:C:234:ILE:HD12	1:C:234:ILE:N	2.35	0.41
2:D:327:GLU:O	2:D:331:GLN:HG2	2.19	0.41
2:B:75:MET:HE3	2:B:92:PHE:HD2	1.85	0.41
2:B:223:THR:O	2:B:227:LEU:HD13	2.20	0.41
1:C:192:HIS:CG	1:C:421:ALA:HA	2.55	0.41
1:A:115:ILE:HG23	1:A:116:ASP:N	2.35	0.41
1:C:58:ALA:HB3	11:C:613:HOH:O	2.20	0.41
1:C:98:ASP:HB2	5:C:501:GTP:O2G	2.20	0.41
2:B:36:TYR:CD1	2:B:46:LEU:HD21	2.56	0.41
2:B:173:PRO:HG3	2:B:187:ALA:HB2	2.03	0.41
2:D:312:TYR:CE1	2:D:377:PHE:HZ	2.39	0.41
3:E:63:TYR:O	3:E:67:GLU:HG2	2.20	0.41
4:F:190:LEU:HB2	4:F:322:ASP:O	2.20	0.41
4:F:220:VAL:HG12	4:F:263:PHE:CD1	2.56	0.41
2:D:143:GLY:O	2:D:147:SER:OG	2.37	0.41
2:D:187:ALA:O	2:D:191:VAL:HG23	2.20	0.41
2:D:398:MET:HE3	2:D:404:PHE:CD2	2.56	0.41
1:C:301:GLN:OE1	1:C:307:PRO:HG2	2.21	0.41
2:D:21:TRP:CE3	2:D:63:PRO:HB3	2.56	0.41
2:D:32:PRO:HA	2:D:83:PHE:CD2	2.56	0.41
4:F:280:GLU:HA	4:F:284:LEU:CB	2.51	0.41
1:C:392:ASP:HB3	1:C:422:ARG:CZ	2.51	0.41
2:D:54:ASN:O	2:D:61:TYR:HA	2.21	0.41
2:D:325:MET:HE2	2:D:355:VAL:HG21	2.02	0.41
1:A:349:THR:HB	3:E:25:LYS:HB3	2.04	0.40
2:D:138:THR:HG22	2:D:169:PHE:HB2	2.03	0.40
3:E:67:GLU:OE1	3:E:67:GLU:HA	2.20	0.40
2:B:9:ALA:HA	2:B:68:VAL:O	2.21	0.40
4:F:98:TYR:HA	4:F:127:GLU:OE1	2.20	0.40
4:F:372:THR:O	4:F:372:THR:HG22	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:198:LYS:O	4:F:199:PHE:HB3	2.21	0.40
3:E:9:ILE:HD11	3:E:21:GLU:OE1	2.22	0.40
1:A:22:GLU:HG3	1:A:83:TYR:HE1	1.87	0.40
3:E:140:LYS:HE2	3:E:140:LYS:HB3	1.88	0.40
4:F:7:ARG:NH1	4:F:32:LYS:O	2.55	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	436/451 (97%)	423 (97%)	11 (2%)	2 (0%)	24	37
1	C	439/451 (97%)	431 (98%)	8 (2%)	0	100	100
2	B	420/445 (94%)	407 (97%)	13 (3%)	0	100	100
2	D	422/445 (95%)	409 (97%)	12 (3%)	1 (0%)	43	58
3	E	119/143 (83%)	118 (99%)	1 (1%)	0	100	100
4	F	312/384 (81%)	297 (95%)	15 (5%)	0	100	100
All	All	2148/2319 (93%)	2085 (97%)	60 (3%)	3 (0%)	48	64

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	282	TYR
2	D	245	PRO
1	A	261	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	369/379 (97%)	364 (99%)	5 (1%)	59	79
1	C	372/379 (98%)	370 (100%)	2 (0%)	81	91
2	B	366/383 (96%)	364 (100%)	2 (0%)	81	91
2	D	368/383 (96%)	364 (99%)	4 (1%)	65	82
3	E	110/127 (87%)	109 (99%)	1 (1%)	70	85
4	F	287/342 (84%)	287 (100%)	0	100	100
All	All	1872/1993 (94%)	1858 (99%)	14 (1%)	76	88

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	GLU
1	A	282	TYR
1	A	284	GLU
1	A	300	ASN
1	A	316	CYS
2	B	47	GLU
2	B	302	MET
1	C	341	ILE
1	C	381	THR
2	D	39	ASP
2	D	207	GLU
2	D	247	GLN
2	D	419	THR
3	E	128	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	88	HIS
1	A	101	ASN

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Mol	Chain	Res	Type
1	A	283	HIS
1	A	329	ASN
2	B	15	GLN
2	B	193	GLN
2	B	282	GLN
2	B	294	GLN
2	B	424	ASN
1	C	11	GLN
1	C	393	HIS
2	D	294	GLN
2	D	300	ASN
2	D	331	GLN
3	E	103	GLN
3	E	136	ASN
4	F	180	HIS
4	F	269	GLN
4	F	306	HIS
4	F	348	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 9 are monoatomic - leaving 6 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	29,30,30	1.17	3 (10%)	45,47,47	1.70	6 (13%)
5	GTP	C	501	6	33,34,34	0.96	2 (6%)	50,54,54	1.50	10 (20%)
9	WLS	B	504	-	15,15,15	0.19	0	17,18,18	0.28	0
10	ACP	F	401	6	31,33,33	1.68	9 (29%)	47,52,52	1.87	10 (21%)
8	GDP	B	501	6	29,30,30	1.16	3 (10%)	45,47,47	1.72	6 (13%)
5	GTP	A	501	6	33,34,34	0.98	2 (6%)	50,54,54	1.52	9 (18%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	GDP	D	501	6	-	4/16/32/32	0/3/3/3
5	GTP	C	501	6	-	7/22/38/38	0/3/3/3
9	WLS	B	504	-	-	2/6/6/6	0/2/2/2
10	ACP	F	401	6	-	9/19/38/38	0/3/3/3
8	GDP	B	501	6	-	3/16/32/32	0/3/3/3
5	GTP	A	501	6	-	5/22/38/38	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	F	401	ACP	C5-C4	4.73	1.47	1.39
8	D	501	GDP	C5-C4	3.07	1.47	1.38
8	B	501	GDP	C5-C4	3.04	1.47	1.38
10	F	401	ACP	PG-O2G	2.99	1.61	1.55
10	F	401	ACP	PG-O3G	2.97	1.61	1.55
10	F	401	ACP	PB-O3A	2.76	1.61	1.58
10	F	401	ACP	C5-C6	2.67	1.48	1.41
8	D	501	GDP	C6-N1	-2.47	1.34	1.38
8	B	501	GDP	C6-N1	-2.37	1.34	1.38
5	C	501	GTP	PA-O3A	2.34	1.62	1.59
10	F	401	ACP	C8-N7	2.33	1.36	1.31
5	A	501	GTP	PA-O3A	2.30	1.62	1.59
5	A	501	GTP	PB-O3B	2.25	1.61	1.59
10	F	401	ACP	C5-N7	-2.25	1.35	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	501	GDP	PA-O3A	2.23	1.61	1.59
5	C	501	GTP	C2-N3	2.09	1.38	1.33
10	F	401	ACP	PA-O3A	2.09	1.61	1.59
10	F	401	ACP	PB-O2B	2.08	1.61	1.56
8	D	501	GDP	C5-N7	-2.03	1.35	1.39

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	D	501	GDP	C5-C4-N3	-5.89	119.02	128.39
8	B	501	GDP	C5-C4-N3	-5.75	119.23	128.39
10	F	401	ACP	C5-C4-N3	-5.73	118.83	126.72
8	B	501	GDP	C2-N3-C4	4.99	120.89	112.30
8	D	501	GDP	C2-N3-C4	4.87	120.69	112.30
5	A	501	GTP	C5-C4-N3	-4.65	120.98	128.39
10	F	401	ACP	N3-C4-N9	4.62	135.02	127.17
5	A	501	GTP	C2-N3-C4	4.56	120.16	112.30
5	C	501	GTP	C5-C4-N3	-4.55	121.14	128.39
5	C	501	GTP	C2-N3-C4	4.34	119.78	112.30
8	B	501	GDP	N9-C4-N3	4.33	134.61	125.95
8	D	501	GDP	N9-C4-N3	4.20	134.36	125.95
10	F	401	ACP	PB-O3A-PA	-3.97	119.41	132.37
10	F	401	ACP	C2-N3-C4	3.64	120.71	111.83
8	B	501	GDP	C6-C5-N7	3.45	136.57	130.29
8	D	501	GDP	C6-C5-N7	3.44	136.54	130.29
10	F	401	ACP	C4-C5-N7	-3.35	106.75	110.58
10	F	401	ACP	N3-C2-N1	-3.19	123.75	128.58
5	C	501	GTP	N9-C8-N7	-2.89	108.05	113.40
5	C	501	GTP	N9-C4-N3	2.86	131.66	125.95
5	A	501	GTP	N9-C4-N3	2.84	131.62	125.95
5	A	501	GTP	C2-N1-C6	-2.81	120.02	125.11
5	C	501	GTP	C2-N1-C6	-2.76	120.10	125.11
10	F	401	ACP	C4-N9-C8	2.76	108.63	105.74
8	D	501	GDP	C4-C5-N7	-2.75	106.31	110.67
5	A	501	GTP	N9-C8-N7	-2.69	108.40	113.40
10	F	401	ACP	C3'-C2'-C1'	2.61	106.40	101.46
5	C	501	GTP	C8-N7-C5	2.60	108.89	104.26
8	B	501	GDP	C4-C5-N7	-2.53	106.67	110.67
10	F	401	ACP	C5-N7-C8	2.46	107.31	103.45
5	A	501	GTP	C8-N7-C5	2.43	108.59	104.26
5	A	501	GTP	C5-C6-N1	2.43	119.43	113.25
5	A	501	GTP	O6-C6-C5	-2.38	120.25	126.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	501	GTP	C5-C6-N1	2.37	119.30	113.25
5	C	501	GTP	O6-C6-C5	-2.33	120.37	126.53
8	B	501	GDP	O6-C6-C5	-2.32	120.42	126.53
5	C	501	GTP	O2B-PB-O3A	2.10	112.94	107.27
5	C	501	GTP	O2A-PA-O3A	2.05	112.81	107.27
5	A	501	GTP	O2A-PA-O3A	2.03	112.77	107.27
10	F	401	ACP	C6-C5-N7	2.01	135.97	132.09
8	D	501	GDP	O6-C6-C5	-2.01	121.22	126.53

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	501	GTP	PB-O3B-PG-O2G
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
8	B	501	GDP	C5'-O5'-PA-O3A
8	B	501	GDP	C5'-O5'-PA-O1A
8	B	501	GDP	C5'-O5'-PA-O2A
8	D	501	GDP	C5'-O5'-PA-O3A
8	D	501	GDP	C5'-O5'-PA-O1A
8	D	501	GDP	C5'-O5'-PA-O2A
10	F	401	ACP	PB-C3B-PG-O1G
10	F	401	ACP	PB-C3B-PG-O2G
10	F	401	ACP	PB-C3B-PG-O3G
10	F	401	ACP	PG-C3B-PB-O1B
10	F	401	ACP	PG-C3B-PB-O2B
10	F	401	ACP	PG-C3B-PB-O3A
10	F	401	ACP	C5'-O5'-PA-O1A
10	F	401	ACP	C5'-O5'-PA-O2A
10	F	401	ACP	C5'-O5'-PA-O3A
9	B	504	WLS	N02-C01-C10-S14
9	B	504	WLS	N02-C01-C10-C11
5	C	501	GTP	PB-O3B-PG-O1G
5	A	501	GTP	C4'-C5'-O5'-PA
5	C	501	GTP	PB-O3B-PG-O2G
5	C	501	GTP	PB-O3B-PG-O3G
8	D	501	GDP	PA-O3A-PB-O2B

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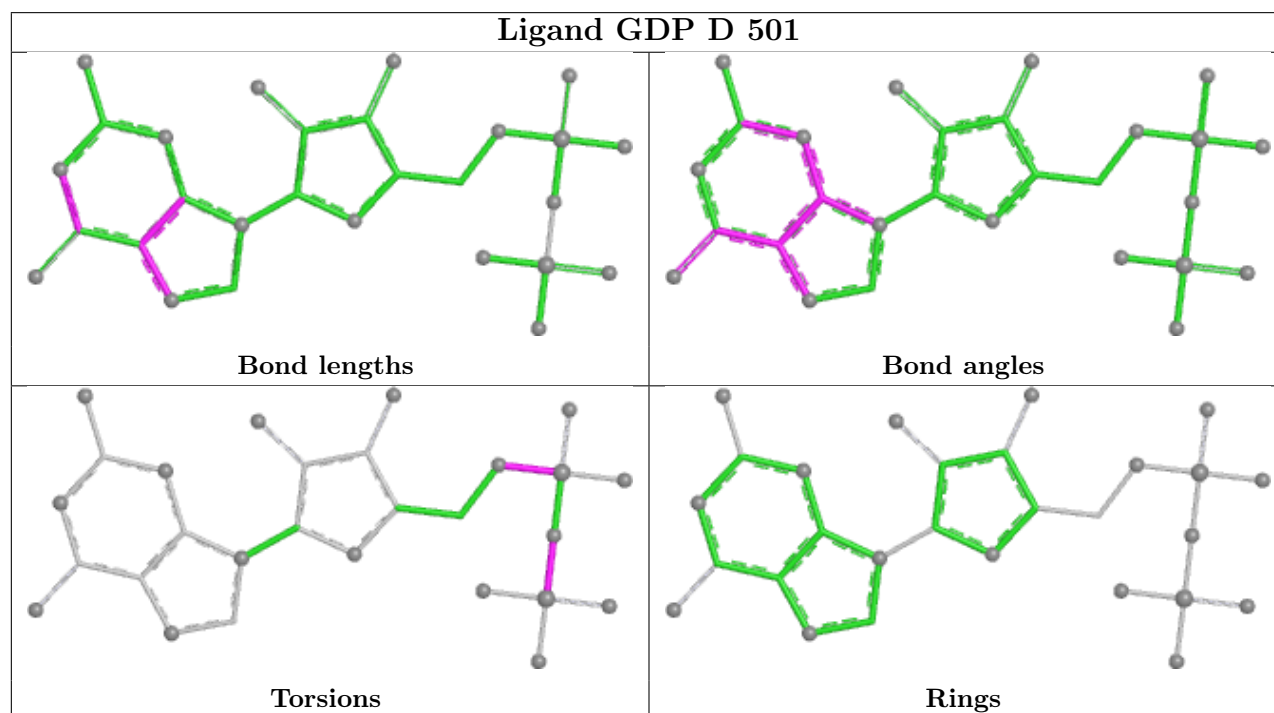
Mol	Chain	Res	Type	Atoms
5	C	501	GTP	PB-O3A-PA-O2A

There are no ring outliers.

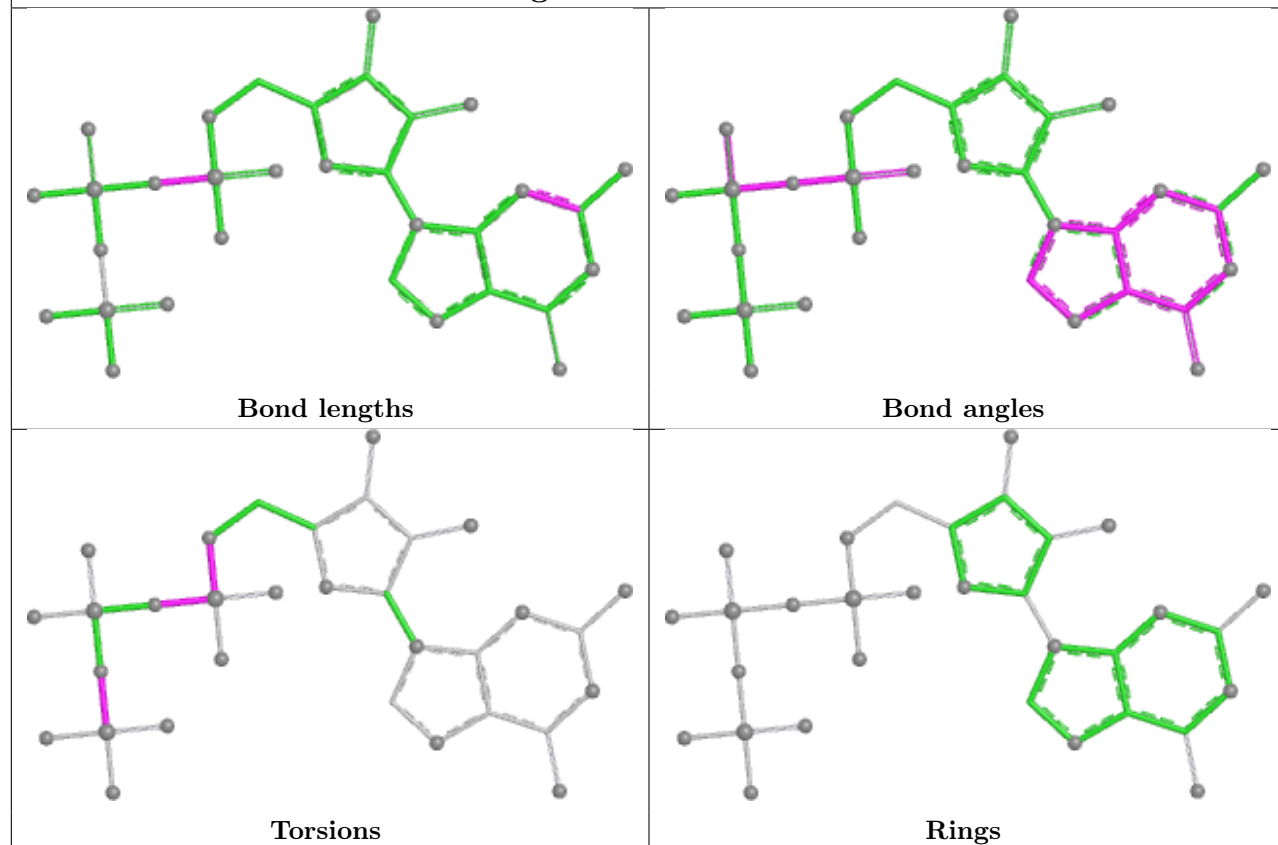
5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	D	501	GDP	3	0
5	C	501	GTP	1	0
9	B	504	WLS	1	0
10	F	401	ACP	3	0
5	A	501	GTP	2	0

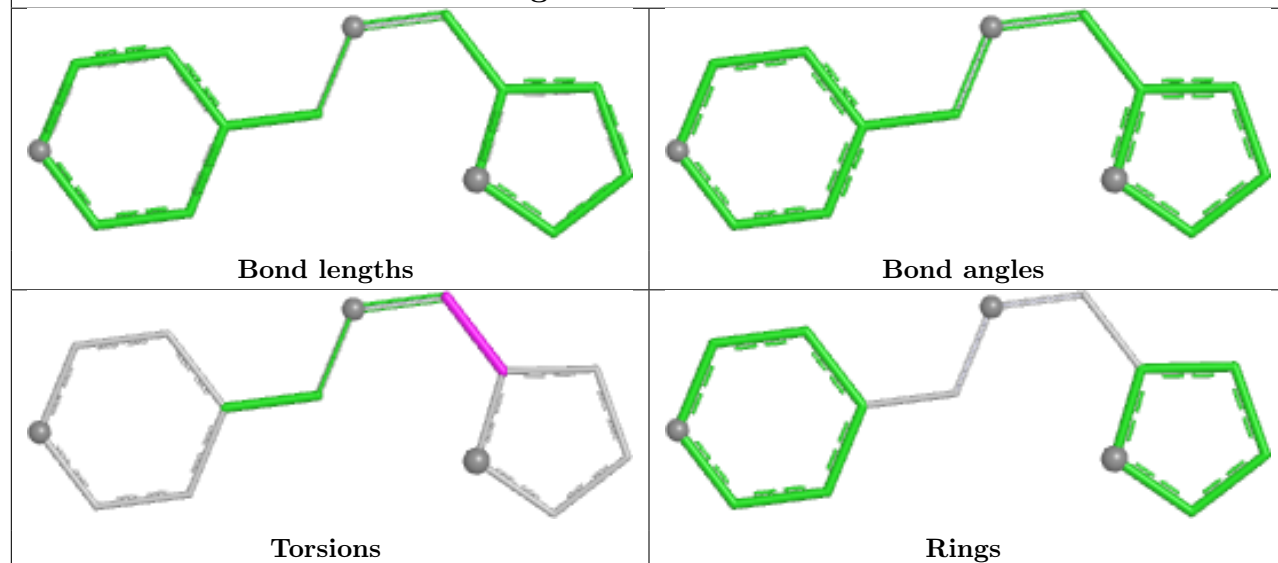
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



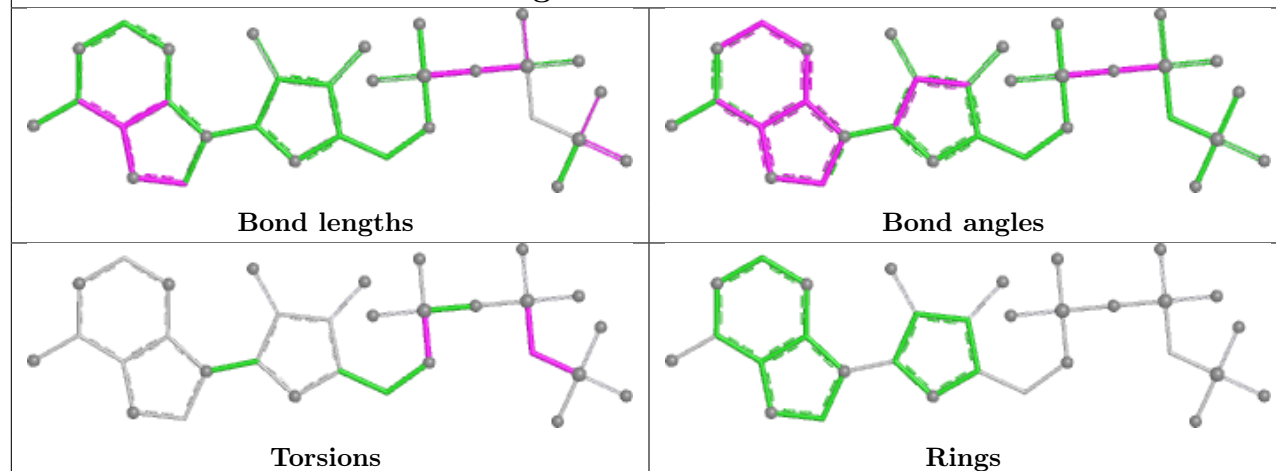
Ligand GTP C 501



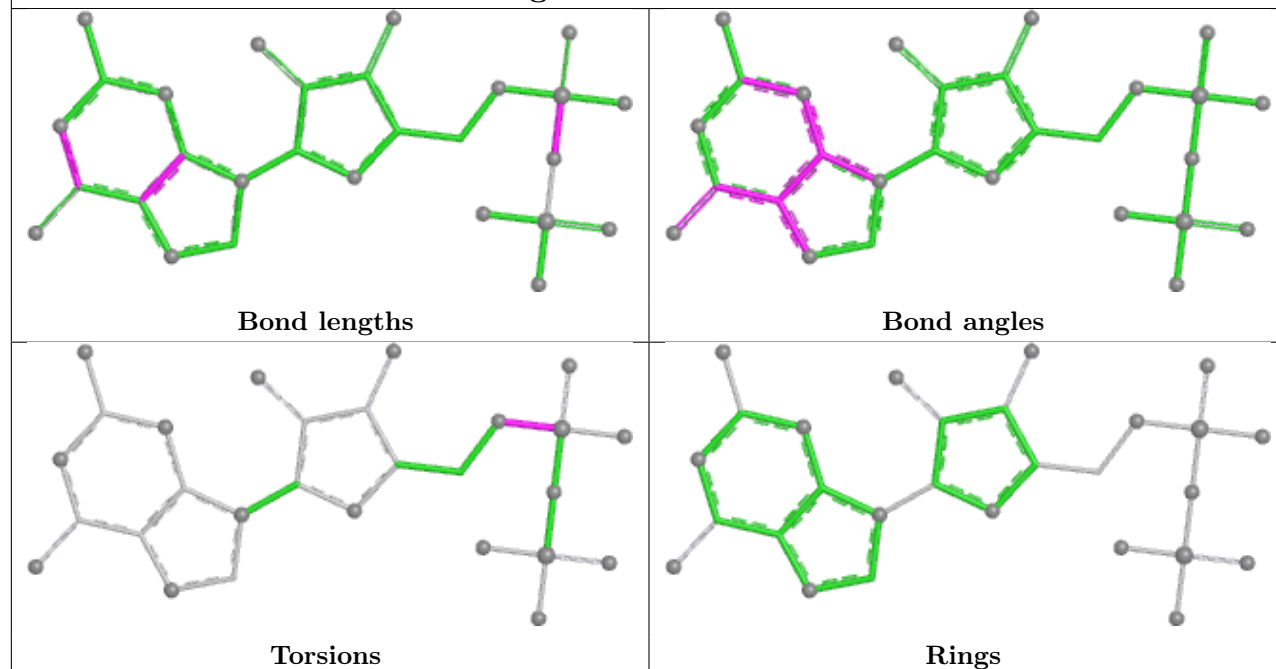
Ligand WLS B 504

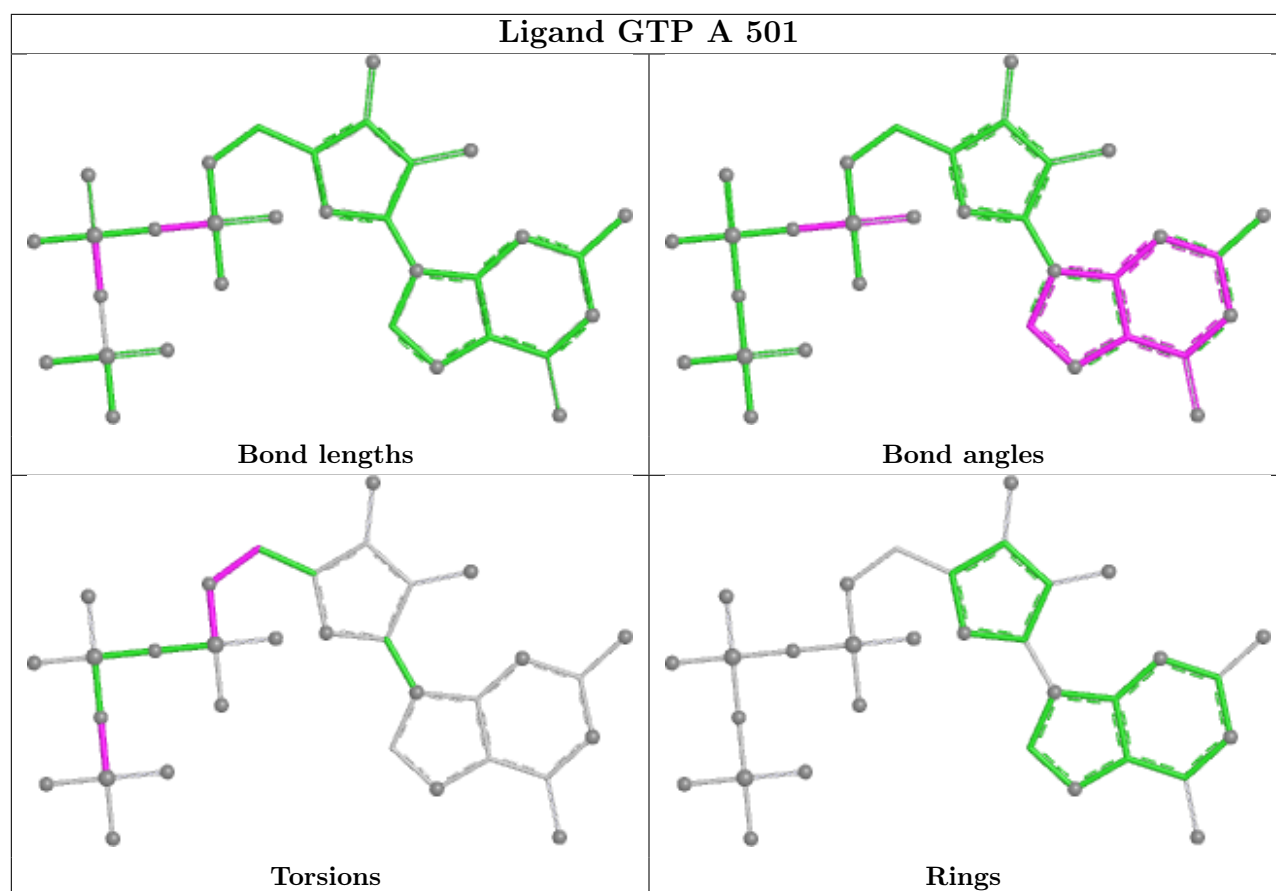


Ligand ACP F 401



Ligand GDP B 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	438/451 (97%)	0.30	9 (2%) 63 59	56, 81, 121, 202	0
1	C	440/451 (97%)	0.07	11 (2%) 58 54	33, 64, 97, 132	1 (0%)
2	B	423/445 (95%)	0.27	12 (2%) 55 51	36, 75, 126, 170	3 (0%)
2	D	426/445 (95%)	0.32	14 (3%) 49 45	59, 86, 124, 163	4 (0%)
3	E	123/143 (86%)	0.63	11 (8%) 15 12	61, 93, 145, 195	0
4	F	322/384 (83%)	0.46	12 (3%) 45 41	68, 100, 169, 225	0
All	All	2172/2319 (93%)	0.29	69 (3%) 50 46	33, 81, 134, 225	8 (0%)

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	F	155	ALA	4.8
2	B	249	ASN	4.7
2	B	248	LEU	4.7
1	C	440	VAL	4.1
1	A	349	THR	4.0
2	B	250	ALA	3.7
3	E	24	LEU	3.7
4	F	371	PRO	3.6
3	E	20	PHE	3.5
3	E	26	PRO	3.5
4	F	186	LEU	3.3
2	B	284	ARG	3.3
2	D	276	THR	3.3
3	E	22	VAL	3.2
4	F	159	GLY	3.2
4	F	167	SER	3.2
3	E	143	ALA	3.0
3	E	8	VAL	3.0
1	C	340	SER	3.0

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Mol	Chain	Res	Type	RSRZ
2	D	167	ASN	3.0
4	F	125	THR	3.0
2	B	57	THR	2.9
1	A	351	PHE	2.9
1	A	88	HIS	2.8
2	B	1	MET	2.8
1	C	248	LEU	2.8
3	E	23	ILE	2.7
2	B	241	CYS	2.7
2	B	333	LEU	2.7
2	D	272	PHE	2.6
4	F	160	ILE	2.6
1	C	341	ILE	2.6
2	B	275	LEU	2.6
2	D	280	SER	2.6
1	C	358	GLN	2.6
1	A	281	ALA	2.6
2	D	248	LEU	2.6
1	A	282	TYR	2.6
1	C	284	GLU	2.5
4	F	179	VAL	2.5
1	C	253	THR	2.5
3	E	9	ILE	2.4
1	C	247	ALA	2.4
2	D	241	CYS	2.4
2	D	82	PRO	2.3
3	E	11	LEU	2.3
2	B	283	TYR	2.3
4	F	154	GLY	2.3
1	A	438	ASP	2.3
1	A	316	CYS	2.3
4	F	153	ALA	2.3
1	C	283	HIS	2.3
4	F	381	HIS	2.2
2	D	127	GLU	2.2
1	C	250	VAL	2.2
4	F	228	TYR	2.2
1	C	286	LEU	2.2
2	D	42	LEU	2.2
2	D	286	LEU	2.2
2	B	82	PRO	2.2
2	D	407	TRP	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	276	THR	2.1
3	E	15	THR	2.1
3	E	6	MET	2.1
2	D	279	GLY	2.0
2	D	80	SER	2.0
2	D	128	SER	2.0
1	A	255	PHE	2.0
1	A	335	ILE	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

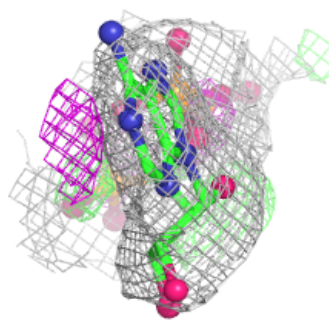
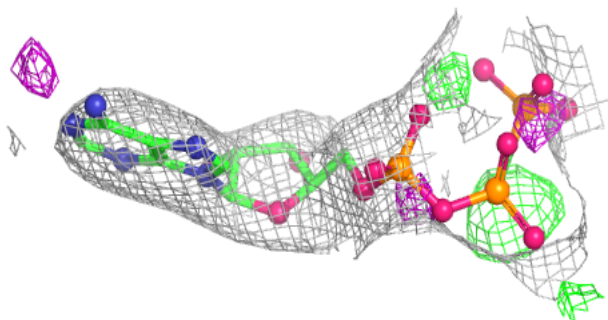
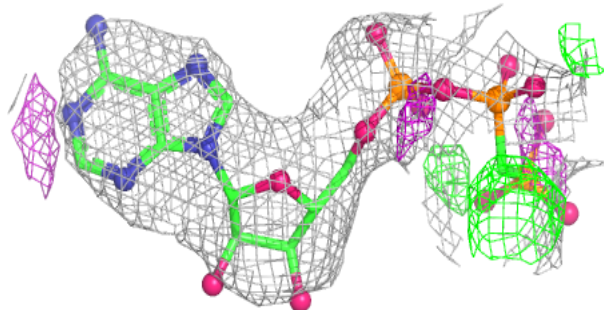
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	0.74	0.12	119,119,119,119	0
6	MG	F	402	1/1	0.77	0.16	122,122,122,122	0
10	ACP	F	401	31/31	0.79	0.13	136,157,166,166	0
9	WLS	B	504	14/14	0.91	0.10	72,87,101,115	0
8	GDP	D	501	28/28	0.93	0.09	70,77,89,95	0
6	MG	D	502	1/1	0.94	0.07	79,79,79,79	0
7	CA	A	504	1/1	0.95	0.08	106,106,106,106	0
5	GTP	A	501	32/32	0.97	0.06	52,61,70,70	0
8	GDP	B	501	28/28	0.98	0.06	49,55,62,69	0
7	CA	A	503	1/1	0.98	0.04	115,115,115,115	0
5	GTP	C	501	32/32	0.98	0.05	47,52,56,58	0
6	MG	B	502	1/1	0.98	0.07	59,59,59,59	0
6	MG	A	502	1/1	1.00	0.04	60,60,60,60	0
7	CA	C	503	1/1	1.00	0.02	94,94,94,94	0
6	MG	C	502	1/1	1.00	0.05	60,60,60,60	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.

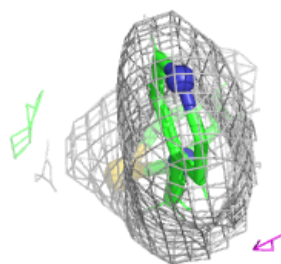
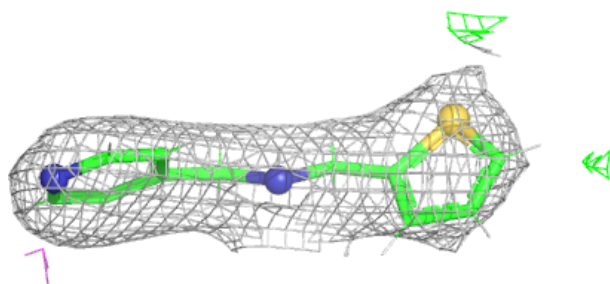
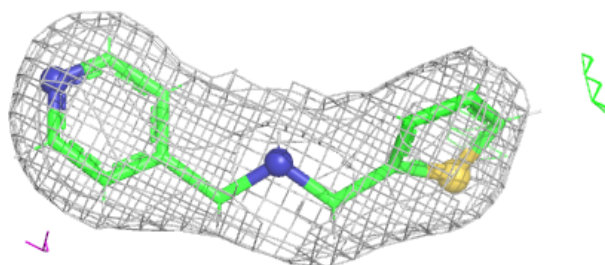
Electron density around ACP F 401:

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

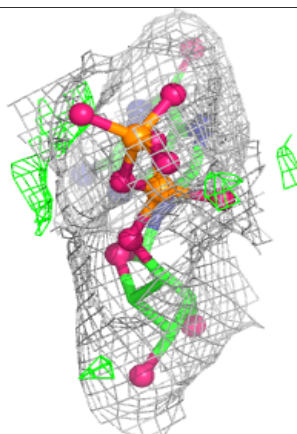
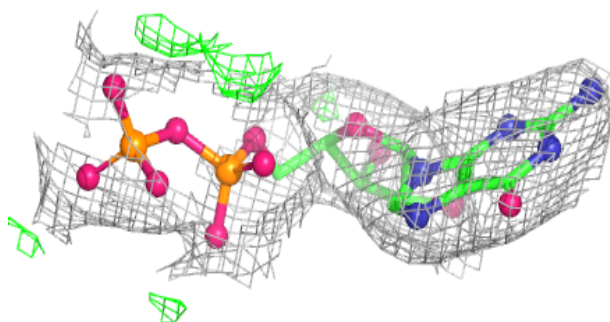
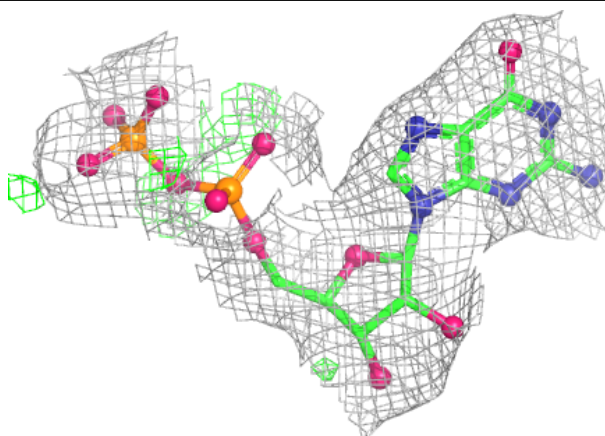


Electron density around WLS B 504:

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

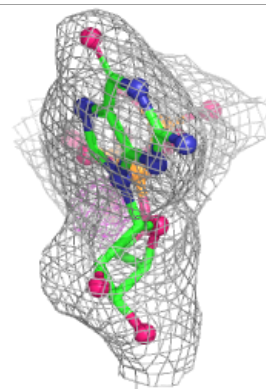
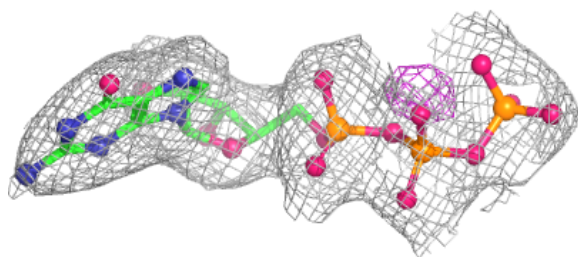
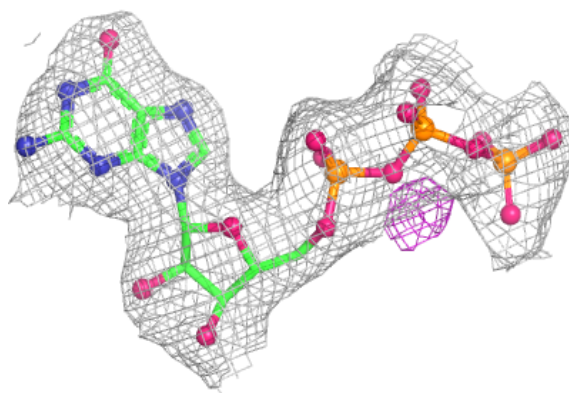
**Electron density around GDP D 501:**

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



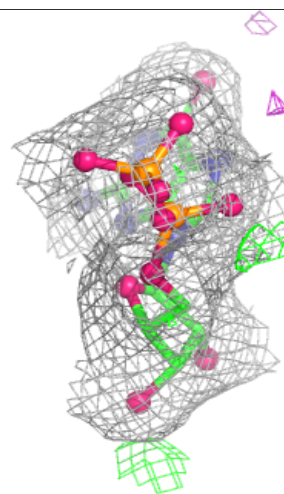
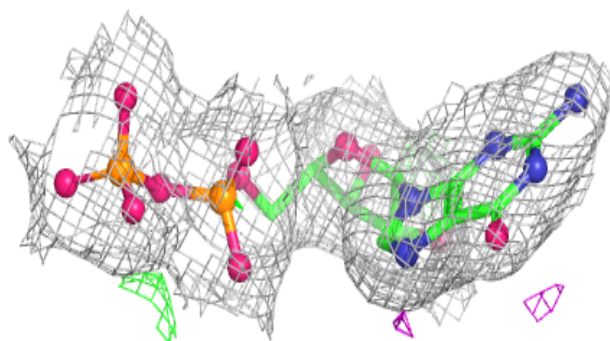
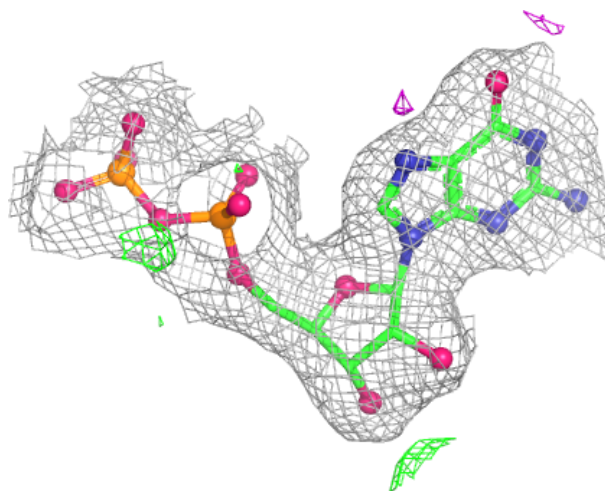
Electron density around GTP A 501:

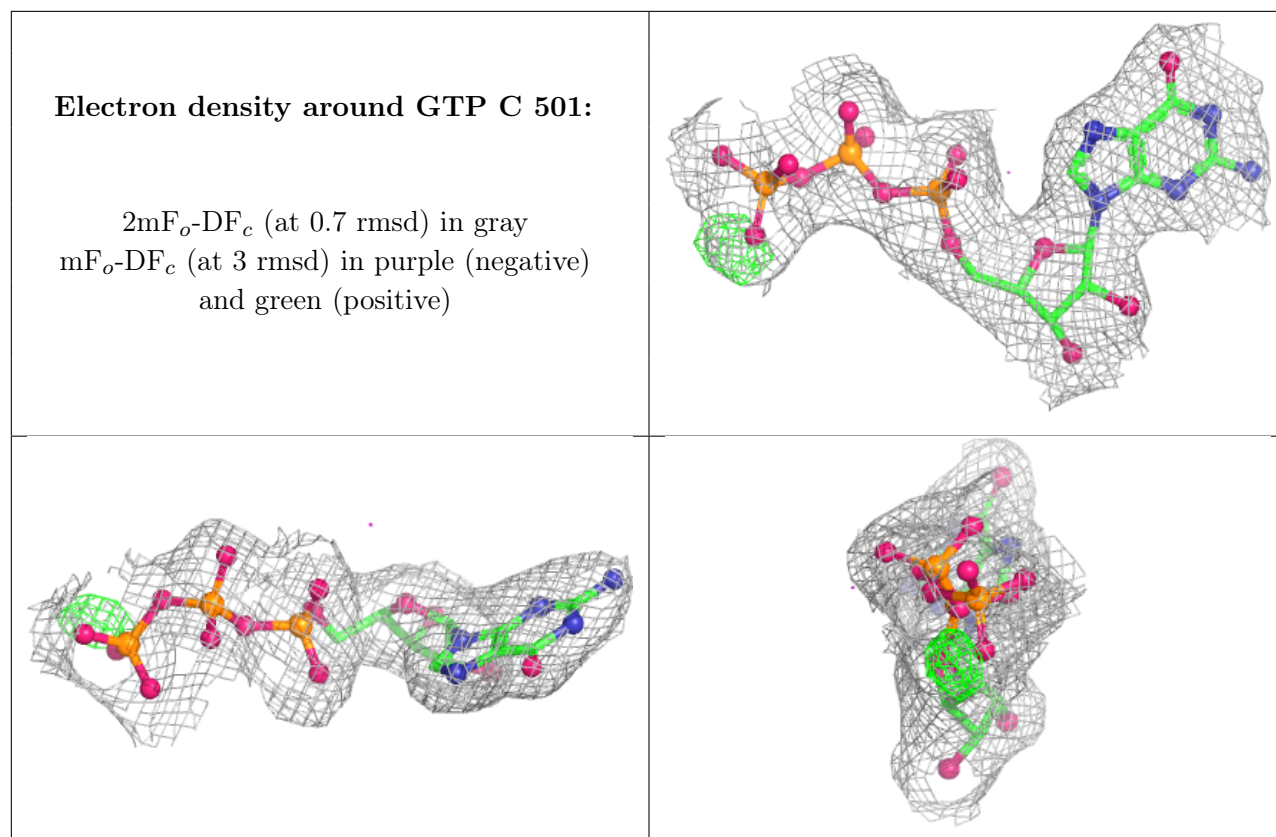
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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.