



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 03:15 PM UTC

PDB ID : 1JLR / pdb_00001jlr
Title : STRUCTURE OF THE URACIL PHOSPHORIBOSYLTRANSFERASE
GTP COMPLEX 2 MUTANT C128V
Authors : Schumacher, M.A.; Bashor, C.J.; Otsu, K.; Zu, S.; Parry, R.; Ulmman, B.;
Brennan, R.G.
Deposited on : 2001-07-16
Resolution : 2.45 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

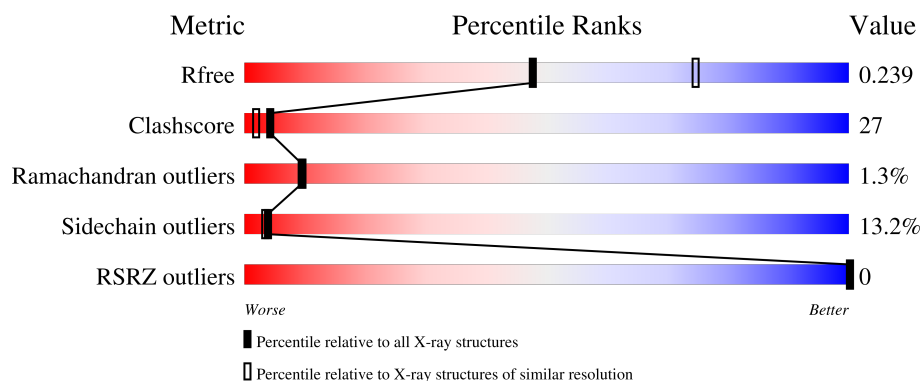
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

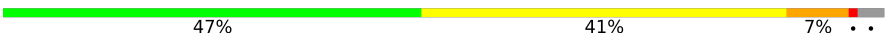
The reported resolution of this entry is 2.45 Å.

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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	243	
1	B	243	
1	C	243	
1	D	243	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7823 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Uracil Phosphoribosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	235	Total	C	N	O	S	0	0	0
			1874	1199	320	342	13			
1	B	235	Total	C	N	O	S	0	0	0
			1874	1199	320	342	13			
1	D	235	Total	C	N	O	S	0	0	0
			1874	1199	320	342	13			
1	C	235	Total	C	N	O	S	0	0	0
			1874	1199	320	342	13			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	84	GLN	GLU	conflict	UNP Q26998
A	128	VAL	CYS	engineered mutation	UNP Q26998
A	157	GLU	ASP	conflict	UNP Q26998
B	84	GLN	GLU	conflict	UNP Q26998
B	128	VAL	CYS	engineered mutation	UNP Q26998
B	157	GLU	ASP	conflict	UNP Q26998
D	84	GLN	GLU	conflict	UNP Q26998
D	128	VAL	CYS	engineered mutation	UNP Q26998
D	157	GLU	ASP	conflict	UNP Q26998
C	84	GLN	GLU	conflict	UNP Q26998
C	128	VAL	CYS	engineered mutation	UNP Q26998
C	157	GLU	ASP	conflict	UNP Q26998

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).



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

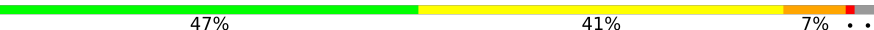
- Molecule 4 is water.

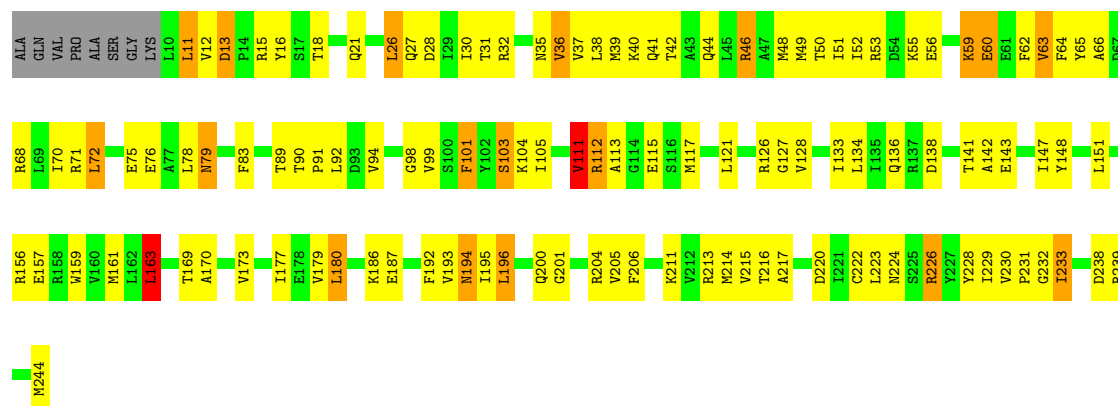
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	38	Total	O	0	0
			38	38		
4	B	33	Total	O	0	0
			33	33		
4	D	45	Total	O	0	0
			45	45		
4	C	43	Total	O	0	0
			43	43		

3 Residue-property plots

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.

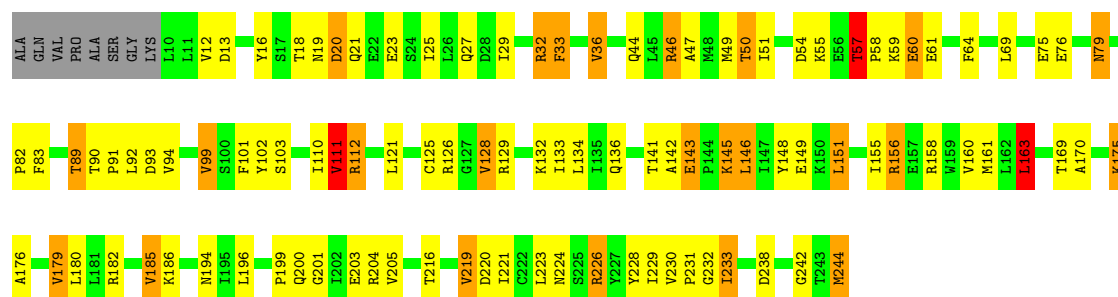
• Molecule 1: Uracil Phosphoribosyltransferase

Chain A: 



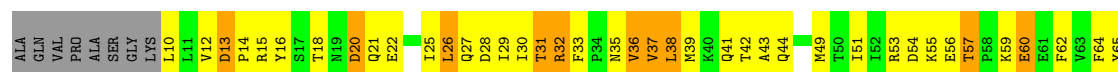
• Molecule 1: Uracil Phosphoribosyltransferase

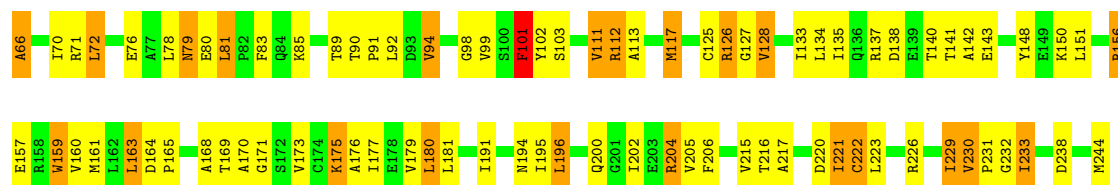
Chain B: 



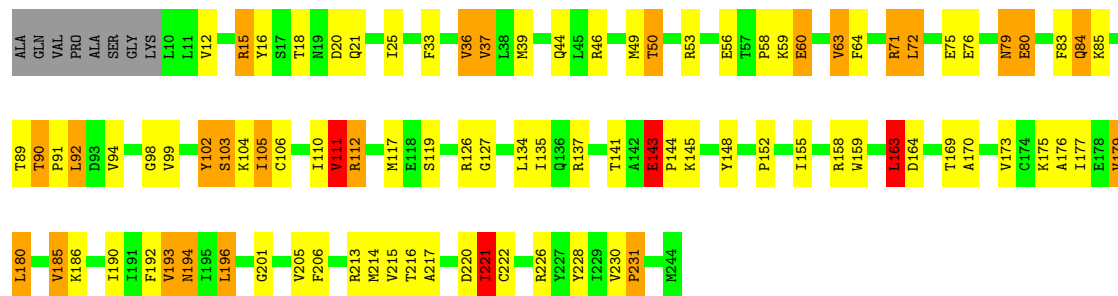
• Molecule 1: Uracil Phosphoribosyltransferase

Chain D: 





• Molecule 1: Uracil Phosphoribosyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.60Å 141.80Å 71.40Å 90.00° 115.00° 90.00°	Depositor
Resolution (Å)	16.14 – 2.45 16.14 – 2.45	Depositor EDS
% Data completeness (in resolution range)	86.6 (16.14-2.45) 86.4 (16.14-2.45)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 2.43Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.184 , 0.242 0.183 , 0.239	Depositor DCC
R_{free} test set	3437 reflections (9.83%)	wwPDB-VP
Wilson B-factor (Å ²)	20.1	Xtriage
Anisotropy	0.099	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 16.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.468 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7823	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.36% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, GTP

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	0.48	0/1905	0.99	9/2573 (0.3%)
1	B	0.46	0/1905	1.02	9/2573 (0.3%)
1	C	0.47	0/1905	1.02	11/2573 (0.4%)
1	D	0.47	1/1905 (0.1%)	1.01	13/2573 (0.5%)
All	All	0.47	1/7620 (0.0%)	1.01	42/10292 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	117	MET	SD-CE	-5.40	1.66	1.79

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	103	SER	N-CA-C	8.18	123.86	113.55
1	D	222	CYS	N-CA-C	7.49	118.13	108.24
1	C	221	ILE	N-CA-C	7.45	118.04	110.82
1	C	102	TYR	N-CA-C	-7.25	101.50	110.41
1	B	103	SER	N-CA-C	7.04	121.55	112.26
1	C	33	PHE	CA-C-N	6.95	126.93	119.28
1	C	33	PHE	C-N-CA	6.95	126.93	119.28
1	B	221	ILE	N-CA-C	6.76	116.86	110.30
1	A	101	PHE	N-CA-C	-6.62	100.68	110.48
1	A	159	TRP	N-CA-C	-6.57	98.92	109.96
1	B	33	PHE	CA-C-N	6.47	126.23	119.05
1	B	33	PHE	C-N-CA	6.47	126.23	119.05
1	B	111	VAL	N-CA-C	6.42	122.70	109.34
1	B	57	THR	CA-C-N	6.39	126.30	119.85
1	B	57	THR	C-N-CA	6.39	126.30	119.85
1	C	143	GLU	CA-C-N	6.37	127.37	119.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	143	GLU	C-N-CA	6.37	127.37	119.98
1	D	159	TRP	N-CA-C	-6.28	99.46	109.76
1	D	57	THR	CA-C-N	6.14	126.49	119.92
1	D	57	THR	C-N-CA	6.14	126.49	119.92
1	A	13	ASP	CA-C-N	5.88	125.36	119.24
1	A	13	ASP	C-N-CA	5.88	125.36	119.24
1	D	221	ILE	N-CA-C	5.81	117.74	111.58
1	B	110	ILE	N-CA-C	-5.67	98.88	107.73
1	A	222	CYS	N-CA-C	5.62	115.53	108.45
1	D	103	SER	N-CA-C	5.60	120.18	113.23
1	B	163	LEU	N-CA-C	5.54	117.56	108.52
1	C	163	LEU	N-CA-C	5.43	118.09	109.24
1	D	35	ASN	N-CA-C	5.38	119.83	113.16
1	C	104	LYS	N-CA-C	5.34	117.82	108.52
1	A	59	LYS	N-CA-C	5.34	117.79	111.33
1	D	66	ALA	N-CA-C	-5.33	105.37	111.07
1	A	104	LYS	N-CA-C	-5.29	99.16	108.20
1	A	163	LEU	N-CA-C	5.28	117.34	108.73
1	D	13	ASP	CA-C-N	5.24	124.86	119.05
1	D	13	ASP	C-N-CA	5.24	124.86	119.05
1	D	101	PHE	N-CA-C	5.15	118.06	110.59
1	C	111	VAL	N-CA-C	5.03	119.81	109.34
1	A	111	VAL	N-CA-C	5.03	119.80	109.34
1	D	33	PHE	CA-C-N	5.02	124.81	119.28
1	D	33	PHE	C-N-CA	5.02	124.81	119.28
1	C	15	ARG	N-CA-C	5.02	118.87	112.34

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	1874	0	1934	132	0
1	B	1874	0	1934	108	0
1	C	1874	0	1934	117	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1874	0	1934	132	0
2	A	10	0	0	0	0
2	B	10	0	0	1	0
2	C	10	0	0	1	0
2	D	10	0	0	0	0
3	A	32	0	12	2	0
3	B	32	0	12	2	0
3	C	32	0	12	0	0
3	D	32	0	12	2	0
4	A	38	0	0	7	0
4	B	33	0	0	2	0
4	C	43	0	0	3	0
4	D	45	0	0	4	0
All	All	7823	0	7784	418	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 27.

All (418) 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:90:THR:HG22	1:A:92:LEU:H	1.11	1.08
1:C:90:THR:HG23	1:C:92:LEU:H	1.26	0.99
1:A:49:MET:HE3	1:A:53:ARG:HE	1.28	0.97
1:B:156:ARG:HH11	1:B:156:ARG:H	1.10	0.96
1:B:231:PRO:O	1:D:90:THR:HG21	1.66	0.95
1:A:15:ARG:HD2	1:C:83:PHE:HB2	1.53	0.90
1:B:145:LYS:HE2	1:B:145:LYS:HA	1.57	0.86
1:C:141:THR:HG22	1:C:143:GLU:HG3	1.58	0.86
1:D:90:THR:HG22	1:D:92:LEU:H	1.42	0.84
1:A:90:THR:HG21	1:C:231:PRO:O	1.77	0.84
1:C:46:ARG:O	1:C:50:THR:HG23	1.76	0.83
1:D:54:ASP:O	1:D:57:THR:HG22	1.78	0.83
1:B:99:VAL:HG13	1:D:16:TYR:HB3	1.59	0.83
1:A:50:THR:CG2	1:C:98:GLY:HA3	2.10	0.82
1:B:59:LYS:HE2	1:C:127:GLY:O	1.80	0.80
1:A:44:GLN:HE21	1:C:83:PHE:HZ	1.28	0.79
1:A:99:VAL:H	1:C:50:THR:HG21	1.47	0.79
1:A:90:THR:HG22	1:A:92:LEU:N	1.96	0.78
1:C:49:MET:HE2	1:C:53:ARG:HH11	1.47	0.78
1:A:64:PHE:CZ	1:C:126:ARG:HD2	2.18	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:27:GLN:O	1:D:31:THR:HG22	1.84	0.77
1:C:49:MET:HE3	1:C:53:ARG:HD2	1.65	0.77
1:A:66:ALA:O	1:A:70:ILE:HG12	1.85	0.76
1:B:82:PRO:HB2	1:B:102:TYR:HD2	1.50	0.76
1:B:233:ILE:HD13	1:B:233:ILE:O	1.85	0.76
1:A:16:TYR:HB3	1:C:99:VAL:HG13	1.67	0.76
1:A:49:MET:HE3	1:A:53:ARG:NE	2.00	0.75
1:C:163:LEU:HD12	1:C:193:VAL:HG22	1.69	0.75
1:B:51:ILE:O	1:B:57:THR:HG21	1.86	0.74
1:A:46:ARG:O	1:A:50:THR:HG23	1.88	0.74
1:A:111:VAL:HG21	1:A:134:LEU:HB2	1.69	0.74
1:D:70:ILE:HD11	1:D:195:ILE:HD11	1.70	0.74
1:A:70:ILE:CD1	1:A:195:ILE:HD11	2.18	0.74
1:D:200:GLN:NE2	1:D:223:LEU:HD12	2.03	0.73
1:B:21:GLN:O	1:B:25:ILE:HG12	1.87	0.73
1:D:70:ILE:HD12	1:D:117:MET:HG2	1.68	0.73
1:D:156:ARG:HD3	4:D:816:HOH:O	1.87	0.73
1:B:125:CYS:O	1:B:128:VAL:HG22	1.88	0.72
1:A:50:THR:HG21	1:C:98:GLY:HA3	1.69	0.72
1:C:20:ASP:O	1:C:21:GLN:HG2	1.88	0.72
1:A:92:LEU:HD13	1:C:230:VAL:HG12	1.70	0.72
1:B:99:VAL:CG1	1:D:16:TYR:HB3	2.20	0.72
1:A:231:PRO:O	1:C:90:THR:HG21	1.91	0.71
1:A:111:VAL:HG11	1:A:134:LEU:HD13	1.70	0.71
1:D:233:ILE:HD13	1:D:233:ILE:O	1.90	0.71
1:B:90:THR:HG23	1:B:92:LEU:H	1.54	0.71
1:D:117:MET:CE	1:D:165:PRO:HD3	2.20	0.71
1:B:133:ILE:HD11	1:B:151:LEU:HD11	1.71	0.71
1:B:79:ASN:HD21	1:D:42:THR:HB	1.55	0.70
1:C:141:THR:CG2	1:C:143:GLU:HG3	2.22	0.70
1:A:126:ARG:HD2	1:C:64:PHE:CZ	2.27	0.70
1:A:70:ILE:HD12	1:A:117:MET:HG2	1.74	0.70
1:A:173:VAL:O	1:A:177:ILE:HG13	1.92	0.70
1:C:117:MET:HE1	1:C:193:VAL:HG23	1.73	0.70
1:D:26:LEU:O	1:D:30:ILE:HG12	1.93	0.69
1:C:49:MET:O	1:C:53:ARG:HG3	1.92	0.69
1:C:230:VAL:HA	1:C:231:PRO:C	2.17	0.69
1:A:26:LEU:O	1:A:30:ILE:HG12	1.93	0.69
1:B:244:MET:HE3	1:B:244:MET:H	1.58	0.69
1:D:12:VAL:HG11	1:D:18:THR:CG2	2.23	0.68
1:D:194:ASN:O	1:D:216:THR:HA	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:VAL:HG11	1:A:18:THR:CG2	2.23	0.68
1:D:62:PHE:HE1	1:D:233:ILE:HD11	1.57	0.68
1:A:59:LYS:O	1:A:63:VAL:HG13	1.94	0.67
1:C:163:LEU:HD12	1:C:193:VAL:CG2	2.24	0.67
1:B:145:LYS:CE	1:B:146:LEU:H	2.07	0.67
1:B:230:VAL:HA	1:B:231:PRO:C	2.19	0.67
1:D:36:VAL:HG12	1:D:206:PHE:CE2	2.29	0.67
1:A:42:THR:HB	1:C:79:ASN:HD21	1.60	0.67
1:A:70:ILE:HD11	1:A:195:ILE:HD11	1.76	0.67
1:A:186:LYS:HE2	1:A:186:LYS:HA	1.77	0.67
1:B:231:PRO:HA	1:D:94:VAL:HG11	1.77	0.67
1:A:244:MET:HB2	4:A:823:HOH:O	1.94	0.67
1:A:15:ARG:HG2	1:A:15:ARG:HH11	1.59	0.67
1:A:90:THR:HG23	1:A:91:PRO:HD2	1.78	0.66
1:A:244:MET:HE3	1:A:244:MET:HA	1.78	0.66
1:D:173:VAL:O	1:D:177:ILE:HG12	1.94	0.66
1:D:238:ASP:HB3	1:D:244:MET:HB3	1.77	0.66
1:A:133:ILE:HD11	1:A:151:LEU:HD21	1.76	0.66
1:B:112:ARG:HB3	2:B:800:PO4:O1	1.94	0.66
1:B:64:PHE:CZ	1:D:126:ARG:HD2	2.30	0.66
1:C:152:PRO:HG2	1:C:155:ILE:HB	1.77	0.65
1:D:133:ILE:HD11	1:D:151:LEU:HD21	1.78	0.65
1:A:127:GLY:O	1:D:59:LYS:HE2	1.97	0.65
1:A:50:THR:HG22	1:C:98:GLY:HA3	1.78	0.65
1:A:98:GLY:HA3	1:C:50:THR:CG2	2.26	0.65
1:C:60:GLU:CD	1:C:60:GLU:H	2.04	0.65
1:A:89:THR:HG22	4:A:833:HOH:O	1.96	0.64
1:D:90:THR:HG23	1:D:91:PRO:HD2	1.80	0.64
1:A:44:GLN:NE2	1:C:83:PHE:HZ	1.96	0.63
1:A:187:GLU:HG2	1:A:211:LYS:HB2	1.79	0.63
1:D:117:MET:HE1	1:D:165:PRO:N	2.13	0.63
1:B:201:GLY:O	1:B:205:VAL:HG12	1.98	0.63
1:A:231:PRO:HA	1:C:94:VAL:HG21	1.80	0.63
1:D:70:ILE:CD1	1:D:195:ILE:HD11	2.29	0.63
1:B:180:LEU:O	1:B:185:VAL:HG13	1.99	0.62
1:C:201:GLY:O	1:C:205:VAL:HG23	1.99	0.62
1:B:175:LYS:O	1:B:179:VAL:HG12	1.99	0.62
1:D:12:VAL:HG11	1:D:18:THR:HG23	1.82	0.62
1:A:12:VAL:HG11	1:A:18:THR:HG21	1.82	0.61
1:D:62:PHE:CE1	1:D:233:ILE:HD11	2.34	0.61
1:B:29:ILE:CD1	1:B:219:VAL:HG21	2.31	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:179:VAL:O	1:B:182:ARG:HG3	2.01	0.61
1:D:26:LEU:HG	1:D:38:LEU:HD11	1.82	0.61
1:C:49:MET:HE2	1:C:53:ARG:NH1	2.15	0.61
1:B:50:THR:HG21	1:D:99:VAL:H	1.65	0.61
1:B:148:TYR:OH	1:C:112:ARG:HG3	2.01	0.61
1:B:75:GLU:HG3	1:D:72:LEU:HD11	1.82	0.60
1:D:37:VAL:HG22	1:D:215:VAL:HG13	1.83	0.60
1:D:168:ALA:HB1	1:D:223:LEU:HD21	1.83	0.60
1:B:156:ARG:HH11	1:B:156:ARG:N	1.91	0.60
1:D:135:ILE:HD12	1:D:175:LYS:HG2	1.84	0.60
1:B:141:THR:HG22	1:B:143:GLU:HB2	1.82	0.60
1:B:60:GLU:H	1:B:60:GLU:CD	2.07	0.60
1:A:98:GLY:HA3	1:C:50:THR:HG21	1.84	0.59
1:C:59:LYS:O	1:C:63:VAL:HG13	2.03	0.59
1:A:138:ASP:HB3	1:A:143:GLU:H	1.68	0.58
1:B:47:ALA:O	1:B:51:ILE:HG12	2.02	0.58
3:B:301:GTP:O1G	1:C:59:LYS:HE3	2.02	0.58
1:C:49:MET:CE	1:C:53:ARG:HD2	2.32	0.58
1:B:141:THR:CG2	1:B:143:GLU:HB2	2.34	0.58
1:A:55:LYS:HG2	1:C:91:PRO:HA	1.85	0.58
1:C:80:GLU:HG3	1:C:213:ARG:HD3	1.86	0.58
1:B:226:ARG:HE	1:B:228:TYR:HE1	1.50	0.58
1:A:105:ILE:N	4:A:814:HOH:O	2.30	0.57
1:B:176:ALA:O	1:B:179:VAL:HG13	2.04	0.57
1:B:58:PRO:HB2	1:B:60:GLU:OE2	2.03	0.57
1:D:229:ILE:O	1:D:232:GLY:HA3	2.04	0.57
1:B:145:LYS:HE2	1:B:146:LEU:H	1.67	0.57
1:A:15:ARG:HG2	1:A:15:ARG:NH1	2.19	0.57
1:B:155:ILE:HG22	1:B:156:ARG:HH12	1.70	0.57
1:C:37:VAL:CG1	1:C:215:VAL:HG22	2.35	0.57
1:A:16:TYR:CD2	1:A:46:ARG:HG3	2.39	0.57
1:D:49:MET:HE3	1:D:53:ARG:HE	1.69	0.57
1:D:66:ALA:O	1:D:70:ILE:HG12	2.05	0.57
1:D:60:GLU:CD	1:D:60:GLU:H	2.11	0.56
1:B:194:ASN:O	1:B:216:THR:HA	2.05	0.56
1:A:32:ARG:HG2	1:A:32:ARG:HH11	1.69	0.56
1:B:49:MET:HE2	1:B:69:LEU:HD11	1.87	0.56
1:A:229:ILE:O	1:A:232:GLY:HA3	2.05	0.56
1:D:202:ILE:O	1:D:205:VAL:HG12	2.05	0.56
1:A:230:VAL:HG22	1:C:92:LEU:HD22	1.88	0.56
1:B:149:GLU:OE2	1:B:151:LEU:HD22	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:230:VAL:HA	1:D:231:PRO:C	2.29	0.56
1:A:94:VAL:HG11	1:C:231:PRO:HA	1.86	0.56
1:B:46:ARG:O	1:B:50:THR:CG2	2.54	0.56
1:C:53:ARG:NH2	1:C:220:ASP:OD2	2.39	0.56
1:A:60:GLU:CD	1:A:60:GLU:H	2.14	0.55
1:B:112:ARG:HD3	1:C:148:TYR:OH	2.06	0.55
1:D:161:MET:O	1:D:163:LEU:HD13	2.07	0.55
1:D:13:ASP:HB2	1:D:41:GLN:O	2.06	0.55
1:D:13:ASP:OD1	1:D:15:ARG:HG2	2.06	0.55
1:B:89:THR:O	1:D:55:LYS:HG2	2.06	0.55
1:D:125:CYS:HB2	1:D:128:VAL:CG2	2.36	0.55
1:A:70:ILE:HD13	1:A:195:ILE:HD11	1.89	0.55
1:A:138:ASP:HB3	1:A:142:ALA:N	2.21	0.55
1:D:196:LEU:CD2	1:D:229:ILE:HG21	2.37	0.55
1:A:16:TYR:CG	1:A:46:ARG:HG3	2.42	0.55
1:A:39:MET:HG2	1:A:76:GLU:HG2	1.89	0.55
1:B:125:CYS:HB3	1:B:128:VAL:HG13	1.88	0.55
1:A:163:LEU:HD12	1:A:193:VAL:HG22	1.89	0.55
1:D:137:ARG:HD3	1:D:142:ALA:O	2.07	0.55
1:B:112:ARG:NH2	1:B:238:ASP:OD2	2.40	0.54
1:C:53:ARG:HH22	1:C:220:ASP:CG	2.15	0.54
1:B:226:ARG:NE	1:B:228:TYR:HE1	2.05	0.54
1:C:192:PHE:CE1	1:C:214:MET:HE3	2.43	0.54
1:A:64:PHE:CE1	1:C:126:ARG:HD2	2.43	0.54
1:D:141:THR:O	1:D:142:ALA:HB3	2.07	0.54
1:C:39:MET:HG2	1:C:217:ALA:HA	1.89	0.54
1:C:143:GLU:OE2	1:C:145:LYS:HE2	2.07	0.54
1:C:111:VAL:O	1:C:112:ARG:HB2	2.08	0.54
1:A:138:ASP:HB3	1:A:142:ALA:H	1.72	0.54
1:B:90:THR:HG21	1:D:231:PRO:O	2.07	0.54
1:A:111:VAL:CG1	1:A:134:LEU:HD13	2.36	0.54
1:B:238:ASP:HB3	1:B:244:MET:HB3	1.89	0.54
1:D:125:CYS:CB	1:D:128:VAL:HG22	2.38	0.54
1:C:12:VAL:HG11	1:C:18:THR:CG2	2.38	0.54
1:A:35:ASN:HB2	4:A:813:HOH:O	2.08	0.54
1:B:186:LYS:HA	1:B:186:LYS:HE2	1.90	0.54
1:C:36:VAL:HG12	1:C:206:PHE:CE2	2.42	0.54
1:A:90:THR:HG23	1:A:91:PRO:CD	2.38	0.53
1:B:161:MET:O	1:B:163:LEU:HD13	2.08	0.53
1:A:117:MET:HE1	1:A:193:VAL:HG23	1.90	0.53
4:A:821:HOH:O	1:C:89:THR:HG22	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83:PHE:HZ	1:D:44:GLN:HE21	1.55	0.53
1:D:117:MET:HE2	1:D:165:PRO:HD3	1.90	0.53
1:B:13:ASP:HB3	1:B:16:TYR:CE2	2.44	0.53
1:A:192:PHE:CE1	1:A:214:MET:HE3	2.44	0.53
1:A:136:GLN:HB2	1:A:147:ILE:HG13	1.91	0.53
1:B:141:THR:O	1:B:142:ALA:HB3	2.09	0.53
1:D:51:ILE:HG21	1:D:65:TYR:CD2	2.44	0.53
1:A:138:ASP:CB	1:A:143:GLU:H	2.22	0.52
1:B:233:ILE:HD13	1:B:233:ILE:C	2.34	0.52
1:D:83:PHE:CE1	1:D:101:PHE:HB3	2.44	0.52
1:A:138:ASP:O	1:A:142:ALA:HA	2.08	0.52
1:B:79:ASN:ND2	1:D:42:THR:HB	2.23	0.52
1:B:54:ASP:HB3	1:B:57:THR:CG2	2.39	0.52
1:D:169:THR:O	1:D:170:ALA:HB3	2.10	0.52
1:D:80:GLU:HB2	1:D:191:ILE:HD13	1.92	0.52
1:D:117:MET:CE	1:D:165:PRO:CD	2.88	0.52
1:C:12:VAL:HG13	1:C:16:TYR:OH	2.08	0.52
1:C:135:ILE:HD13	1:C:175:LYS:HB3	1.92	0.52
1:A:79:ASN:OD1	1:C:44:GLN:HB2	2.10	0.52
1:B:242:GLY:HA2	1:B:244:MET:HE2	1.92	0.52
1:D:99:VAL:O	1:D:99:VAL:HG23	2.10	0.52
1:C:144:PRO:HB2	1:C:175:LYS:HD3	1.90	0.52
1:D:125:CYS:O	1:D:128:VAL:CG2	2.58	0.52
1:C:21:GLN:O	1:C:25:ILE:HG12	2.09	0.52
1:B:16:TYR:HB3	1:D:99:VAL:HG11	1.92	0.52
1:C:58:PRO:HB2	1:C:60:GLU:OE1	2.10	0.52
1:A:46:ARG:HG2	1:A:46:ARG:HH11	1.74	0.52
1:A:83:PHE:HB2	4:C:825:HOH:O	2.08	0.51
1:A:78:LEU:HD13	3:A:303:GTP:O6	2.10	0.51
1:A:99:VAL:N	1:C:50:THR:HG21	2.20	0.51
1:A:27:GLN:O	1:A:31:THR:HG22	2.11	0.51
1:D:168:ALA:HB1	1:D:223:LEU:CD2	2.41	0.51
1:C:90:THR:CG2	1:C:94:VAL:HG22	2.41	0.51
1:B:64:PHE:CE2	1:D:126:ARG:HD2	2.45	0.51
1:D:111:VAL:CG1	1:D:134:LEU:HD13	2.41	0.51
1:B:200:GLN:NE2	1:B:223:LEU:HG	2.26	0.51
1:C:196:LEU:HD21	1:C:220:ASP:OD1	2.10	0.51
1:C:194:ASN:O	1:C:216:THR:HA	2.10	0.51
1:A:196:LEU:HD21	1:A:220:ASP:CG	2.36	0.50
1:B:141:THR:HG22	1:B:143:GLU:H	1.76	0.50
1:D:99:VAL:C	4:D:803:HOH:O	2.53	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:VAL:HG12	1:A:206:PHE:CE2	2.45	0.50
1:D:196:LEU:HD21	1:D:229:ILE:HG21	1.92	0.50
1:A:148:TYR:OH	1:D:112:ARG:HD3	2.12	0.50
4:B:810:HOH:O	1:D:89:THR:HG22	2.12	0.50
1:C:196:LEU:HD21	1:C:220:ASP:CG	2.37	0.50
1:B:229:ILE:O	1:B:232:GLY:HA3	2.11	0.50
1:D:156:ARG:HG3	1:D:157:GLU:HG3	1.94	0.50
1:C:221:ILE:HG13	1:C:222:CYS:N	2.25	0.50
1:C:37:VAL:HG13	1:C:215:VAL:HA	1.94	0.50
1:A:53:ARG:O	1:C:91:PRO:HD3	2.11	0.49
1:C:112:ARG:HB3	2:C:802:PO4:O1	2.12	0.49
1:D:78:LEU:HD13	3:D:302:GTP:O6	2.12	0.49
1:D:85:LYS:HA	1:D:99:VAL:HG12	1.93	0.49
1:A:83:PHE:HZ	1:C:44:GLN:HE22	1.61	0.49
1:B:46:ARG:O	1:B:50:THR:HG23	2.12	0.49
1:B:101:PHE:HE2	1:D:44:GLN:NE2	2.11	0.49
1:B:99:VAL:HG13	1:D:16:TYR:CB	2.37	0.49
1:A:112:ARG:NH1	1:D:150:LYS:HG2	2.27	0.49
1:A:156:ARG:NH1	1:A:157:GLU:HG2	2.28	0.49
1:D:125:CYS:HB2	1:D:128:VAL:HG22	1.94	0.49
1:B:179:VAL:HA	1:B:182:ARG:HG2	1.94	0.49
1:C:15:ARG:HB3	4:C:825:HOH:O	2.11	0.48
1:A:156:ARG:NH1	1:A:156:ARG:HG3	2.27	0.48
1:B:55:LYS:HG2	1:D:89:THR:O	2.11	0.48
1:D:81:LEU:HD23	3:D:302:GTP:HN21	1.78	0.48
1:D:141:THR:HG22	1:D:143:GLU:HB2	1.95	0.48
1:B:90:THR:HG23	1:B:92:LEU:N	2.25	0.48
1:B:91:PRO:HD3	1:D:53:ARG:O	2.13	0.48
1:A:55:LYS:HG3	1:C:90:THR:C	2.38	0.48
1:A:226:ARG:HG3	1:A:226:ARG:HH11	1.79	0.48
1:D:90:THR:HG23	1:D:91:PRO:CD	2.43	0.48
1:C:110:ILE:HD12	1:C:164:ASP:HB2	1.95	0.48
1:B:32:ARG:HG2	1:B:32:ARG:HH11	1.78	0.48
1:C:36:VAL:HG12	1:C:206:PHE:HE2	1.79	0.48
1:D:112:ARG:O	1:D:113:ALA:C	2.55	0.48
1:C:90:THR:HG22	1:C:94:VAL:HG22	1.96	0.48
1:C:186:LYS:N	1:C:186:LYS:HD2	2.28	0.48
1:D:59:LYS:NZ	4:D:806:HOH:O	2.47	0.48
1:A:143:GLU:HG3	1:A:143:GLU:O	2.14	0.48
1:B:133:ILE:CD1	1:B:151:LEU:HD11	2.43	0.48
1:B:82:PRO:HB2	1:B:102:TYR:CD2	2.40	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:49:MET:CE	1:C:53:ARG:HH11	2.20	0.48
1:D:15:ARG:HG3	1:D:43:ALA:HB2	1.95	0.48
1:A:59:LYS:HE2	1:D:127:GLY:O	2.14	0.47
1:A:112:ARG:O	1:A:113:ALA:C	2.56	0.47
1:C:144:PRO:HB2	1:C:175:LYS:CD	2.44	0.47
1:A:136:GLN:HB2	1:A:147:ILE:CG1	2.44	0.47
1:B:44:GLN:HB2	1:D:79:ASN:OD1	2.13	0.47
1:C:176:ALA:O	1:C:179:VAL:HG22	2.14	0.47
1:A:42:THR:HB	1:C:79:ASN:ND2	2.27	0.47
1:B:111:VAL:HB	1:B:134:LEU:HD13	1.95	0.47
1:B:121:LEU:HG	1:B:128:VAL:HG21	1.95	0.47
1:A:55:LYS:HD2	1:C:90:THR:O	2.14	0.47
1:A:161:MET:O	1:A:163:LEU:HD13	2.15	0.47
1:A:201:GLY:O	1:A:205:VAL:HG12	2.14	0.47
1:B:175:LYS:O	1:B:175:LYS:HD3	2.15	0.47
1:C:20:ASP:O	1:C:21:GLN:CG	2.60	0.47
1:C:226:ARG:HD2	1:C:228:TYR:CE2	2.49	0.47
1:B:169:THR:O	1:B:170:ALA:HB3	2.15	0.47
1:D:138:ASP:O	1:D:142:ALA:HA	2.15	0.47
1:D:90:THR:HG22	1:D:91:PRO:N	2.29	0.46
1:B:145:LYS:HE2	1:B:145:LYS:CA	2.36	0.46
1:B:61:GLU:OE1	1:D:126:ARG:NH2	2.44	0.46
1:A:50:THR:HG21	1:C:99:VAL:H	1.80	0.46
1:B:12:VAL:HG11	1:B:18:THR:HG21	1.97	0.46
1:A:39:MET:HE2	1:A:217:ALA:HB2	1.98	0.46
1:B:46:ARG:O	1:B:50:THR:HG22	2.16	0.46
1:A:103:SER:N	4:A:838:HOH:O	2.35	0.46
1:D:26:LEU:HG	1:D:38:LEU:CD1	2.44	0.46
1:A:62:PHE:CD1	1:A:62:PHE:C	2.92	0.46
1:A:75:GLU:HG3	1:C:72:LEU:HD11	1.97	0.46
1:D:25:ILE:O	1:D:29:ILE:HG12	2.16	0.46
1:D:39:MET:HE2	1:D:217:ALA:HB2	1.97	0.46
1:B:69:LEU:C	1:B:69:LEU:HD23	2.41	0.46
1:B:92:LEU:O	1:B:93:ASP:HB2	2.16	0.46
1:D:226:ARG:HG2	1:D:226:ARG:HH11	1.81	0.46
1:A:126:ARG:HD2	1:C:64:PHE:CE1	2.50	0.46
1:D:112:ARG:NH2	1:D:238:ASP:OD2	2.50	0.45
1:B:129:ARG:NH1	3:B:301:GTP:O1B	2.49	0.45
1:D:125:CYS:O	1:D:128:VAL:HG22	2.17	0.45
1:C:84:GLN:CA	1:C:84:GLN:HE21	2.29	0.45
1:A:200:GLN:CD	1:A:223:LEU:HD22	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:VAL:HA	1:A:231:PRO:C	2.42	0.45
1:C:185:VAL:HG21	1:C:190:ILE:HD11	1.98	0.45
1:C:71:ARG:CG	1:C:71:ARG:HH11	2.30	0.45
1:C:84:GLN:HE21	1:C:84:GLN:HA	1.81	0.45
1:C:180:LEU:O	1:C:185:VAL:HG13	2.16	0.45
1:A:90:THR:CG2	1:A:91:PRO:N	2.80	0.45
3:A:303:GTP:PA	4:A:814:HOH:O	2.75	0.45
1:A:13:ASP:HB2	1:A:41:GLN:O	2.17	0.45
1:D:22:GLU:O	1:D:26:LEU:HB2	2.17	0.45
1:D:90:THR:CG2	1:D:91:PRO:HD2	2.47	0.45
1:A:169:THR:O	1:A:170:ALA:HB3	2.17	0.45
1:B:126:ARG:HD3	1:D:64:PHE:CE2	2.52	0.45
1:D:111:VAL:HG11	1:D:134:LEU:HD13	1.99	0.45
1:B:61:GLU:CD	1:D:126:ARG:HH22	2.24	0.44
1:D:202:ILE:O	1:D:205:VAL:CG1	2.65	0.44
1:C:46:ARG:O	1:C:50:THR:CG2	2.56	0.44
1:A:141:THR:O	1:A:142:ALA:HB3	2.17	0.44
1:A:226:ARG:HG3	1:A:226:ARG:NH1	2.32	0.44
1:B:242:GLY:C	1:B:244:MET:HE2	2.42	0.44
1:A:52:ILE:O	1:A:239:ARG:NH2	2.50	0.44
1:D:156:ARG:HB2	1:D:156:ARG:HH11	1.83	0.44
1:C:119:SER:HB2	4:C:826:HOH:O	2.17	0.44
1:D:13:ASP:HA	1:D:14:PRO:HD3	1.89	0.44
1:D:39:MET:HG2	1:D:76:GLU:HG2	1.99	0.44
1:D:148:TYR:OH	1:D:150:LYS:HE2	2.18	0.44
1:D:226:ARG:HG2	1:D:226:ARG:NH1	2.33	0.44
1:C:169:THR:O	1:C:170:ALA:HB3	2.17	0.44
1:D:28:ASP:O	1:D:32:ARG:HB2	2.17	0.44
1:D:36:VAL:HG12	1:D:206:PHE:CZ	2.53	0.44
1:A:16:TYR:HB3	1:C:99:VAL:CG1	2.41	0.43
1:A:121:LEU:HG	1:A:128:VAL:HG11	2.01	0.43
1:A:28:ASP:OD1	1:A:32:ARG:HD3	2.18	0.43
1:A:48:MET:HE3	1:A:68:ARG:HB3	2.01	0.43
1:B:54:ASP:O	1:B:57:THR:HG23	2.18	0.43
1:D:51:ILE:O	1:D:57:THR:HG21	2.18	0.43
1:C:16:TYR:CE2	1:C:46:ARG:HG3	2.53	0.43
1:A:224:ASN:C	1:A:226:ARG:H	2.26	0.43
1:D:90:THR:CG2	1:D:91:PRO:CD	2.97	0.43
1:A:163:LEU:HD12	1:A:193:VAL:CG2	2.48	0.43
1:D:102:TYR:CE2	1:D:159:TRP:CD1	3.06	0.43
1:D:179:VAL:HG23	1:D:180:LEU:N	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:20:ASP:C	1:C:21:GLN:HG2	2.44	0.43
1:C:37:VAL:HG11	1:C:215:VAL:HG22	1.98	0.43
1:A:83:PHE:HZ	1:C:44:GLN:NE2	2.16	0.43
1:C:137:ARG:NH1	1:C:144:PRO:HD3	2.34	0.43
1:A:156:ARG:HG3	1:A:156:ARG:HH11	1.84	0.43
1:B:76:GLU:O	1:B:79:ASN:HB2	2.18	0.43
1:A:26:LEU:HG	1:A:38:LEU:CD2	2.49	0.43
1:A:46:ARG:HB3	1:C:99:VAL:HG22	2.01	0.43
1:B:233:ILE:O	1:D:91:PRO:HG2	2.18	0.43
1:B:16:TYR:HB3	1:D:99:VAL:CG1	2.49	0.43
1:B:50:THR:CG2	1:D:98:GLY:HA3	2.49	0.43
1:B:156:ARG:H	1:B:156:ARG:HD2	1.83	0.42
1:B:199:PRO:O	1:B:203:GLU:HG3	2.19	0.42
1:C:221:ILE:CG1	1:C:222:CYS:N	2.81	0.42
1:C:226:ARG:O	1:C:226:ARG:HD3	2.19	0.42
1:A:112:ARG:CZ	1:D:150:LYS:HG2	2.49	0.42
1:A:186:LYS:HE2	1:A:186:LYS:CA	2.47	0.42
1:C:84:GLN:HA	1:C:84:GLN:NE2	2.34	0.42
1:B:224:ASN:HB3	1:B:228:TYR:H	1.85	0.42
1:A:12:VAL:HG11	1:A:18:THR:HG23	2.02	0.42
1:B:129:ARG:HD2	1:B:158:ARG:CZ	2.49	0.42
1:B:132:LYS:O	1:B:133:ILE:HD13	2.19	0.42
1:B:220:ASP:HB3	1:B:230:VAL:O	2.20	0.42
1:C:105:ILE:HG12	1:C:106:CYS:N	2.34	0.42
1:C:173:VAL:O	1:C:177:ILE:HG13	2.20	0.42
1:D:171:GLY:HA2	1:D:204:ARG:NH2	2.34	0.42
1:A:112:ARG:O	1:A:115:GLU:N	2.53	0.42
1:B:231:PRO:HA	1:D:94:VAL:CG1	2.48	0.42
1:B:46:ARG:HG2	4:B:836:HOH:O	2.19	0.42
1:D:117:MET:HE1	1:D:164:ASP:C	2.45	0.42
1:A:112:ARG:NH2	1:A:238:ASP:OD2	2.50	0.41
1:B:230:VAL:HA	1:B:232:GLY:N	2.35	0.41
1:A:72:LEU:HD11	1:C:75:GLU:HG3	2.02	0.41
1:B:29:ILE:HD12	1:B:219:VAL:HG21	2.00	0.41
1:D:20:ASP:O	1:D:21:GLN:HB2	2.21	0.41
1:D:222:CYS:SG	1:D:230:VAL:HG13	2.60	0.41
1:B:90:THR:OG1	1:B:91:PRO:HD2	2.20	0.41
1:D:220:ASP:HB2	1:D:229:ILE:CG2	2.50	0.41
1:D:196:LEU:HD22	1:D:229:ILE:HG21	2.02	0.41
1:A:11:LEU:HA	1:A:11:LEU:HD23	1.59	0.41
1:A:16:TYR:CB	1:C:99:VAL:HG13	2.45	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:90:THR:CG2	1:D:91:PRO:N	2.82	0.41
1:A:90:THR:CG2	1:A:92:LEU:HB2	2.50	0.41
1:C:102:TYR:CE1	1:C:159:TRP:CD1	3.09	0.41
1:C:193:VAL:CG2	1:C:193:VAL:O	2.68	0.41
1:D:179:VAL:CG2	1:D:180:LEU:N	2.83	0.41
1:D:205:VAL:HG13	1:D:206:PHE:N	2.36	0.41
1:D:244:MET:HB2	4:D:823:HOH:O	2.21	0.41
1:A:55:LYS:HD3	1:C:89:THR:HG23	2.02	0.41
1:A:134:LEU:HG	1:A:147:ILE:HB	2.03	0.41
1:A:194:ASN:O	1:A:216:THR:HA	2.21	0.41
1:B:90:THR:CG2	1:B:94:VAL:HG22	2.51	0.41
1:D:176:ALA:O	1:D:180:LEU:HD22	2.21	0.41
1:A:32:ARG:HG2	1:A:32:ARG:NH1	2.36	0.41
1:A:180:LEU:HD12	1:A:180:LEU:HA	1.82	0.41
1:D:12:VAL:HG13	1:D:16:TYR:OH	2.21	0.41
1:B:155:ILE:HG22	1:B:156:ARG:NH1	2.34	0.40
1:B:58:PRO:HG2	1:B:61:GLU:HG2	2.03	0.40
1:D:10:LEU:CB	1:D:30:ILE:HG13	2.51	0.40
1:D:125:CYS:HB3	1:D:128:VAL:HG22	2.02	0.40
1:D:220:ASP:HB3	1:D:230:VAL:O	2.21	0.40
1:C:111:VAL:CG1	1:C:134:LEU:HD13	2.51	0.40
1:C:117:MET:CE	1:C:193:VAL:HG23	2.47	0.40
1:A:51:ILE:HG21	1:A:65:TYR:CD2	2.56	0.40
1:A:228:TYR:HB3	1:C:92:LEU:HD21	2.03	0.40
1:A:233:ILE:O	1:A:233:ILE:HG12	2.20	0.40
1:C:76:GLU:O	1:C:79:ASN:HB2	2.22	0.40
1:B:33:PHE:O	1:B:36:VAL:HG12	2.22	0.40
1:C:84:GLN:O	1:C:85:LYS:C	2.65	0.40
1:C:144:PRO:O	1:C:175:LYS:HD3	2.20	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	233/243 (96%)	214 (92%)	15 (6%)	4 (2%)	7	6
1	B	233/243 (96%)	218 (94%)	12 (5%)	3 (1%)	9	9
1	C	233/243 (96%)	220 (94%)	10 (4%)	3 (1%)	9	9
1	D	233/243 (96%)	218 (94%)	13 (6%)	2 (1%)	14	17
All	All	932/972 (96%)	870 (93%)	50 (5%)	12 (1%)	9	9

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	112	ARG
1	A	111	VAL
1	A	112	ARG
1	B	112	ARG
1	D	112	ARG
1	A	103	SER
1	A	233	ILE
1	B	20	ASP
1	C	111	VAL
1	B	111	VAL
1	D	111	VAL
1	C	231	PRO

5.3.2 Protein sidechains [i](#)

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	206/212 (97%)	183 (89%)	23 (11%)	6	5
1	B	206/212 (97%)	175 (85%)	31 (15%)	3	2
1	C	206/212 (97%)	181 (88%)	25 (12%)	5	4
1	D	206/212 (97%)	176 (85%)	30 (15%)	3	2
All	All	824/848 (97%)	715 (87%)	109 (13%)	4	3

All (109) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LEU
1	A	21	GLN
1	A	26	LEU
1	A	36	VAL
1	A	37	VAL
1	A	40	LYS
1	A	46	ARG
1	A	56	GLU
1	A	60	GLU
1	A	63	VAL
1	A	71	ARG
1	A	72	LEU
1	A	79	ASN
1	A	101	PHE
1	A	163	LEU
1	A	179	VAL
1	A	180	LEU
1	A	194	ASN
1	A	196	LEU
1	A	204	ARG
1	A	213	ARG
1	A	215	VAL
1	A	226	ARG
1	B	19	ASN
1	B	20	ASP
1	B	23	GLU
1	B	27	GLN
1	B	32	ARG
1	B	36	VAL
1	B	46	ARG
1	B	50	THR
1	B	57	THR
1	B	60	GLU
1	B	79	ASN
1	B	89	THR
1	B	99	VAL
1	B	128	VAL
1	B	136	GLN
1	B	143	GLU
1	B	145	LYS
1	B	146	LEU
1	B	151	LEU
1	B	156	ARG

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Mol	Chain	Res	Type
1	B	160	VAL
1	B	163	LEU
1	B	175	LYS
1	B	179	VAL
1	B	185	VAL
1	B	196	LEU
1	B	204	ARG
1	B	219	VAL
1	B	226	ARG
1	B	233	ILE
1	B	244	MET
1	D	20	ASP
1	D	26	LEU
1	D	31	THR
1	D	32	ARG
1	D	36	VAL
1	D	37	VAL
1	D	38	LEU
1	D	56	GLU
1	D	60	GLU
1	D	71	ARG
1	D	72	LEU
1	D	79	ASN
1	D	81	LEU
1	D	94	VAL
1	D	101	PHE
1	D	126	ARG
1	D	128	VAL
1	D	140	THR
1	D	156	ARG
1	D	160	VAL
1	D	163	LEU
1	D	175	LYS
1	D	180	LEU
1	D	181	LEU
1	D	196	LEU
1	D	204	ARG
1	D	221	ILE
1	D	229	ILE
1	D	230	VAL
1	D	233	ILE
1	C	36	VAL

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Mol	Chain	Res	Type
1	C	37	VAL
1	C	50	THR
1	C	56	GLU
1	C	60	GLU
1	C	63	VAL
1	C	71	ARG
1	C	72	LEU
1	C	79	ASN
1	C	80	GLU
1	C	84	GLN
1	C	90	THR
1	C	92	LEU
1	C	103	SER
1	C	105	ILE
1	C	143	GLU
1	C	158	ARG
1	C	163	LEU
1	C	179	VAL
1	C	180	LEU
1	C	185	VAL
1	C	193	VAL
1	C	194	ASN
1	C	196	LEU
1	C	221	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	27	GLN
1	A	44	GLN
1	A	200	GLN
1	B	19	ASN
1	B	21	GLN
1	B	44	GLN
1	B	136	GLN
1	D	27	GLN
1	D	35	ASN
1	D	41	GLN
1	D	44	GLN
1	D	136	GLN
1	D	200	GLN

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Mol	Chain	Res	Type
1	C	21	GLN
1	C	35	ASN
1	C	44	GLN
1	C	84	GLN
1	C	136	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

12 ligands are modelled in this entry.

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	$\# Z > 2$
2	PO4	A	805	-	4,4,4	1.78	2 (50%)	6,6,6	0.45	0
2	PO4	C	803	-	4,4,4	1.71	0	6,6,6	0.49	0
3	GTP	A	303	-	33,34,34	1.95	6 (18%)	50,54,54	1.38	5 (10%)
2	PO4	A	806	-	4,4,4	1.79	2 (50%)	6,6,6	0.46	0
3	GTP	B	301	-	33,34,34	1.84	6 (18%)	50,54,54	1.38	5 (10%)
2	PO4	D	799	-	4,4,4	1.76	0	6,6,6	0.44	0
3	GTP	D	302	-	33,34,34	1.88	6 (18%)	50,54,54	1.36	5 (10%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GTP	C	300	-	33,34,34	1.80	6 (18%)	50,54,54	1.34	5 (10%)
2	PO4	C	802	-	4,4,4	1.79	1 (25%)	6,6,6	0.46	0
2	PO4	D	801	-	4,4,4	1.63	0	6,6,6	0.47	0
2	PO4	B	800	-	4,4,4	1.79	1 (25%)	6,6,6	0.47	0
2	PO4	B	804	-	4,4,4	1.73	1 (25%)	6,6,6	0.49	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GTP	B	301	-	-	3/22/38/38	0/3/3/3
3	GTP	A	303	-	-	3/22/38/38	0/3/3/3
3	GTP	C	300	-	-	6/22/38/38	0/3/3/3
3	GTP	D	302	-	-	7/22/38/38	0/3/3/3

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	302	GTP	PA-O3A	6.13	1.66	1.59
3	A	303	GTP	PA-O3A	5.70	1.65	1.59
3	A	303	GTP	PB-O3B	-5.49	1.53	1.59
3	B	301	GTP	PA-O3A	4.94	1.64	1.59
3	C	300	GTP	PA-O3A	4.85	1.64	1.59
3	C	300	GTP	PB-O3B	-4.82	1.54	1.59
3	B	301	GTP	PB-O3B	-4.69	1.54	1.59
3	B	301	GTP	C6-N1	4.40	1.47	1.38
3	A	303	GTP	C6-N1	4.27	1.46	1.38
3	C	300	GTP	C6-N1	4.12	1.46	1.38
3	D	302	GTP	PB-O3B	-4.10	1.55	1.59
3	D	302	GTP	C6-N1	3.94	1.46	1.38
3	B	301	GTP	C4-N3	3.72	1.42	1.34
3	A	303	GTP	C4-N3	3.61	1.42	1.34
3	D	302	GTP	C4-N3	3.61	1.42	1.34
3	C	300	GTP	C4-N3	3.57	1.42	1.34
3	D	302	GTP	PA-O5'	-3.19	1.46	1.59
3	C	300	GTP	PA-O5'	-2.96	1.47	1.59
3	B	301	GTP	PA-O5'	-2.96	1.47	1.59
3	A	303	GTP	PA-O5'	-2.58	1.49	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	301	GTP	C5-N7	-2.21	1.34	1.39
3	A	303	GTP	C5-N7	-2.11	1.34	1.39
2	B	804	PO4	P-O3	-2.10	1.48	1.54
2	C	802	PO4	P-O3	-2.10	1.48	1.54
3	C	300	GTP	C5-N7	-2.09	1.34	1.39
3	D	302	GTP	C5-N7	-2.09	1.34	1.39
2	B	800	PO4	P-O4	-2.06	1.48	1.54
2	A	805	PO4	P-O2	-2.06	1.48	1.54
2	A	806	PO4	P-O2	-2.05	1.48	1.54
2	A	806	PO4	P-O3	-2.04	1.48	1.54
2	A	805	PO4	P-O4	-2.02	1.48	1.54

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	302	GTP	O2G-PG-O3B	5.65	123.60	104.64
3	B	301	GTP	O2G-PG-O3B	5.59	123.37	104.64
3	A	303	GTP	O2G-PG-O3B	5.56	123.28	104.64
3	C	300	GTP	O2G-PG-O3B	5.52	123.15	104.64
3	A	303	GTP	O2B-PB-O3B	-3.51	97.78	107.27
3	B	301	GTP	O2B-PB-O3B	-3.41	98.07	107.27
3	C	300	GTP	O2B-PB-O3B	-3.21	98.60	107.27
3	D	302	GTP	O2B-PB-O3B	-2.80	99.69	107.27
3	D	302	GTP	O3G-PG-O1G	-2.77	100.03	110.83
3	C	300	GTP	O3G-PG-O1G	-2.64	100.54	110.83
3	A	303	GTP	C3'-C2'-C1'	2.60	106.38	101.46
3	B	301	GTP	O2G-PG-O1G	-2.59	100.74	110.83
3	B	301	GTP	O3G-PG-O1G	-2.59	100.75	110.83
3	A	303	GTP	O3G-PG-O1G	-2.58	100.77	110.83
3	C	300	GTP	C3'-C2'-C1'	2.55	106.29	101.46
3	D	302	GTP	O2G-PG-O1G	-2.54	100.92	110.83
3	A	303	GTP	O2G-PG-O1G	-2.53	101.00	110.83
3	B	301	GTP	C3'-C2'-C1'	2.51	106.21	101.46
3	D	302	GTP	C3'-C2'-C1'	2.46	106.12	101.46
3	C	300	GTP	O2G-PG-O1G	-2.46	101.25	110.83

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	302	GTP	C5'-O5'-PA-O3A
3	D	302	GTP	C3'-C4'-C5'-O5'

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Mol	Chain	Res	Type	Atoms
3	C	300	GTP	C5'-O5'-PA-O3A
3	C	300	GTP	C5'-O5'-PA-O2A
3	D	302	GTP	O4'-C4'-C5'-O5'
3	A	303	GTP	O4'-C4'-C5'-O5'
3	C	300	GTP	O4'-C4'-C5'-O5'
3	A	303	GTP	C4'-C5'-O5'-PA
3	A	303	GTP	C3'-C4'-C5'-O5'
3	B	301	GTP	O4'-C4'-C5'-O5'
3	B	301	GTP	C4'-C5'-O5'-PA
3	C	300	GTP	C3'-C4'-C5'-O5'
3	D	302	GTP	C5'-O5'-PA-O1A
3	C	300	GTP	PB-O3B-PG-O1G
3	C	300	GTP	C4'-C5'-O5'-PA
3	B	301	GTP	C3'-C4'-C5'-O5'
3	D	302	GTP	PA-O3A-PB-O1B
3	D	302	GTP	PA-O3A-PB-O2B
3	D	302	GTP	C2'-C1'-N9-C4

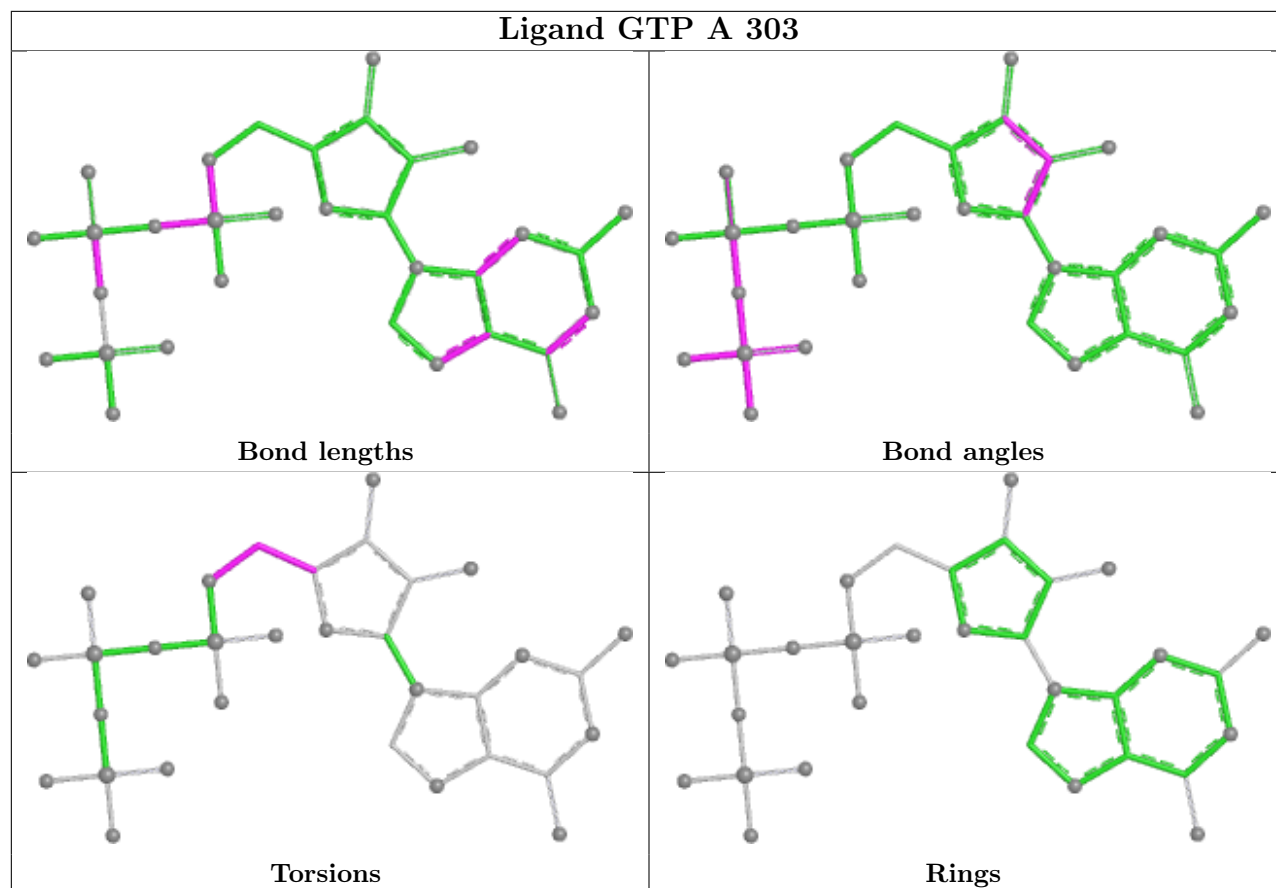
There are no ring outliers.

5 monomers are involved in 8 short contacts:

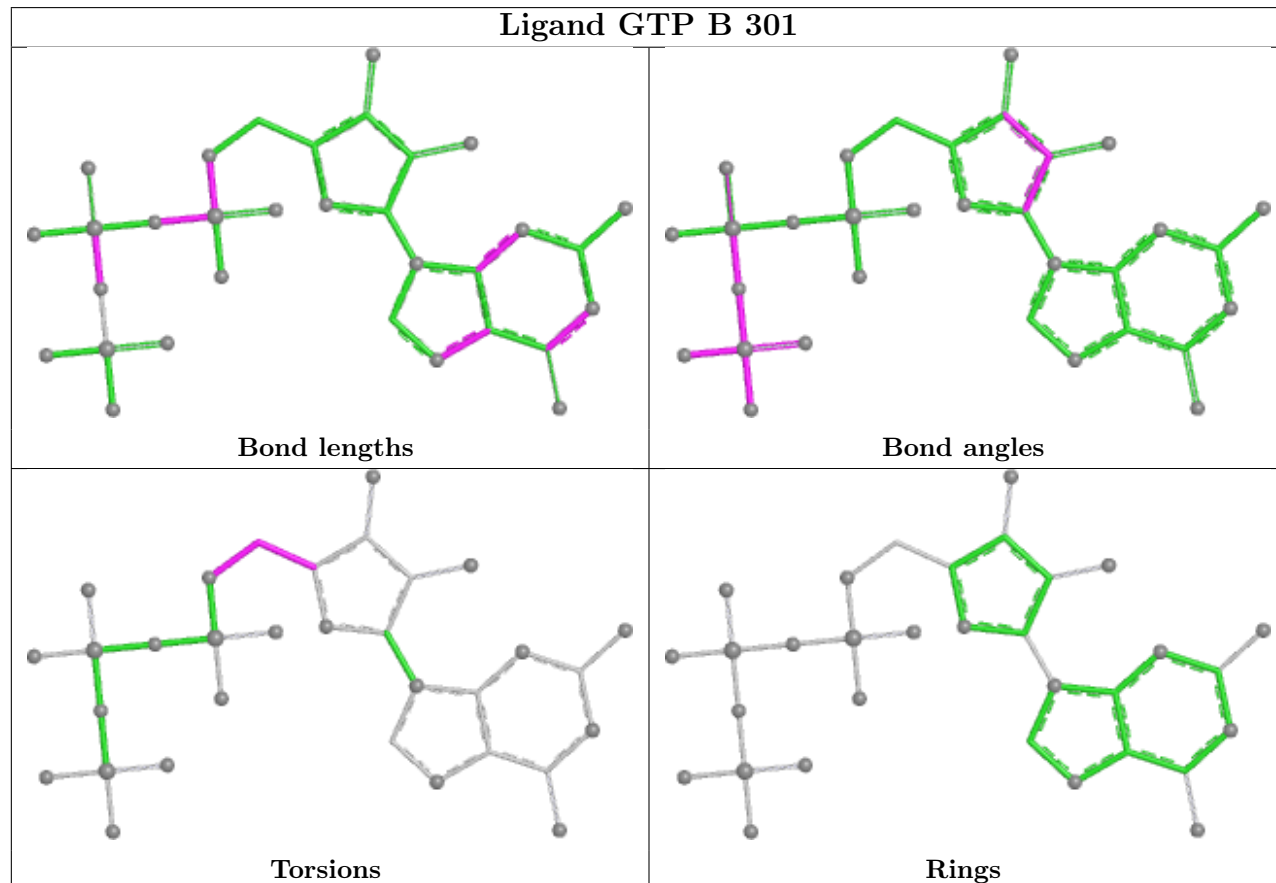
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	303	GTP	2	0
3	B	301	GTP	2	0
3	D	302	GTP	2	0
2	C	802	PO4	1	0
2	B	800	PO4	1	0

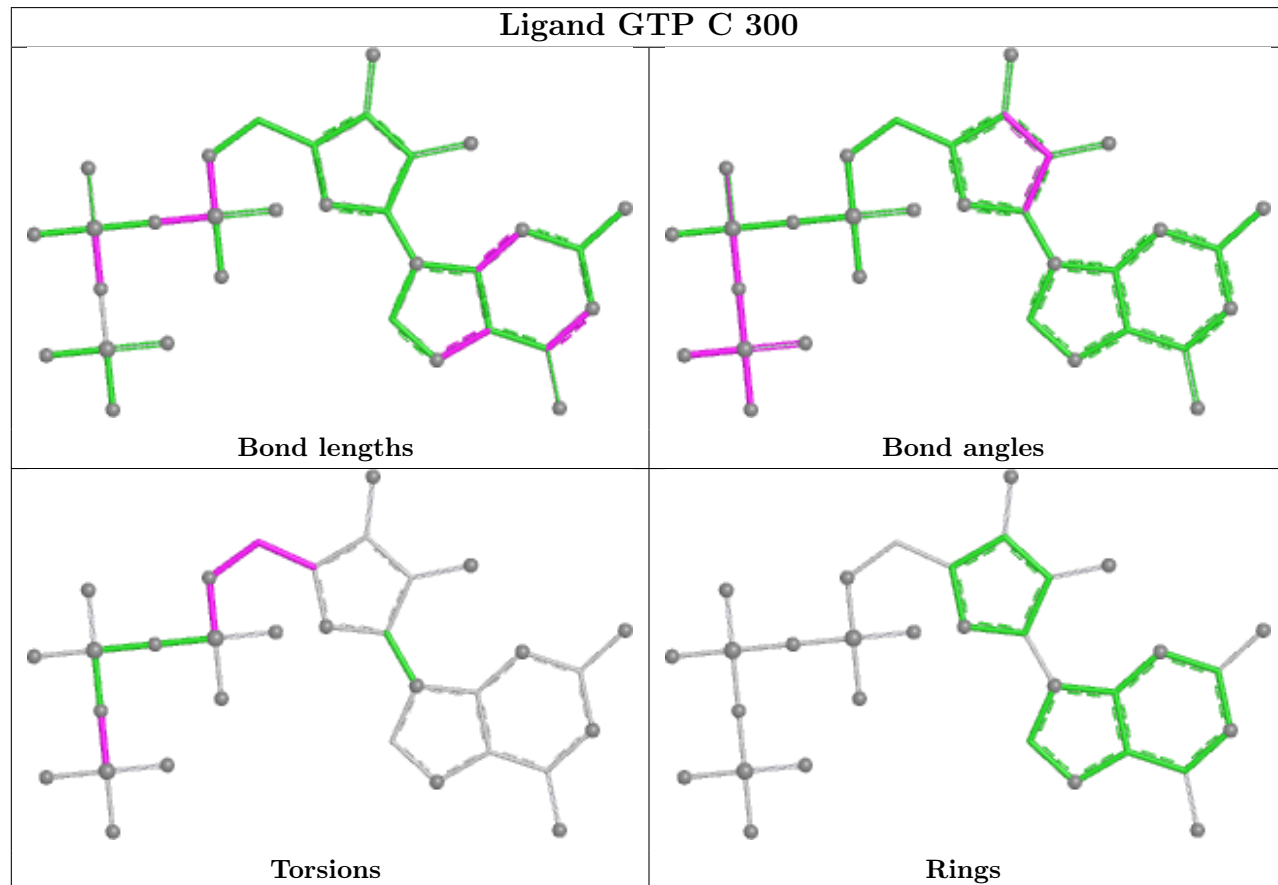
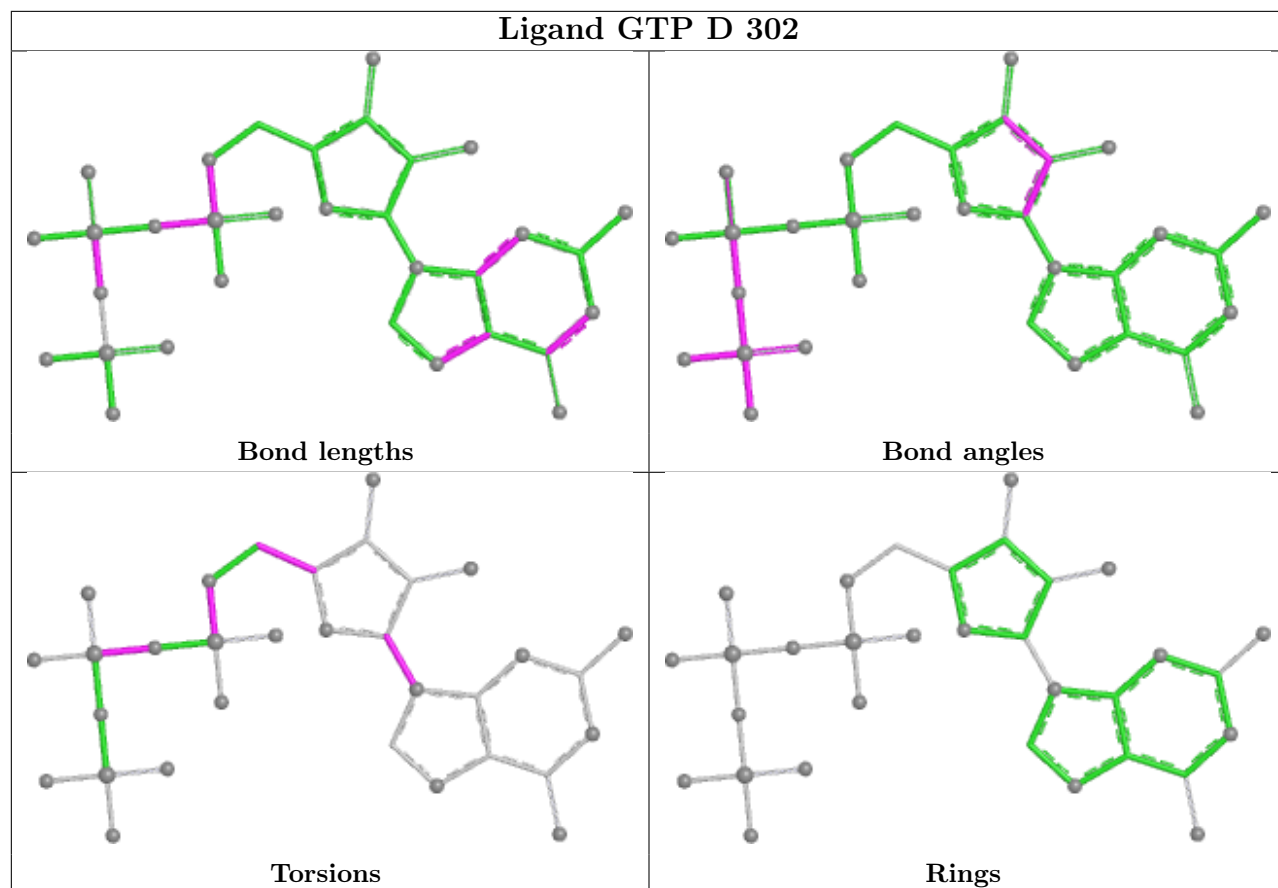
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

Ligand GTP A 303



Ligand GTP B 301





5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	235/243 (96%)	-1.80	0 100 100	4, 18, 64, 84	1 (0%)
1	B	235/243 (96%)	-1.80	0 100 100	4, 19, 52, 75	1 (0%)
1	C	235/243 (96%)	-1.81	0 100 100	4, 20, 50, 78	1 (0%)
1	D	235/243 (96%)	-1.76	0 100 100	5, 18, 67, 82	1 (0%)
All	All	940/972 (96%)	-1.79	0 100 100	4, 19, 61, 84	4 (0%)

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

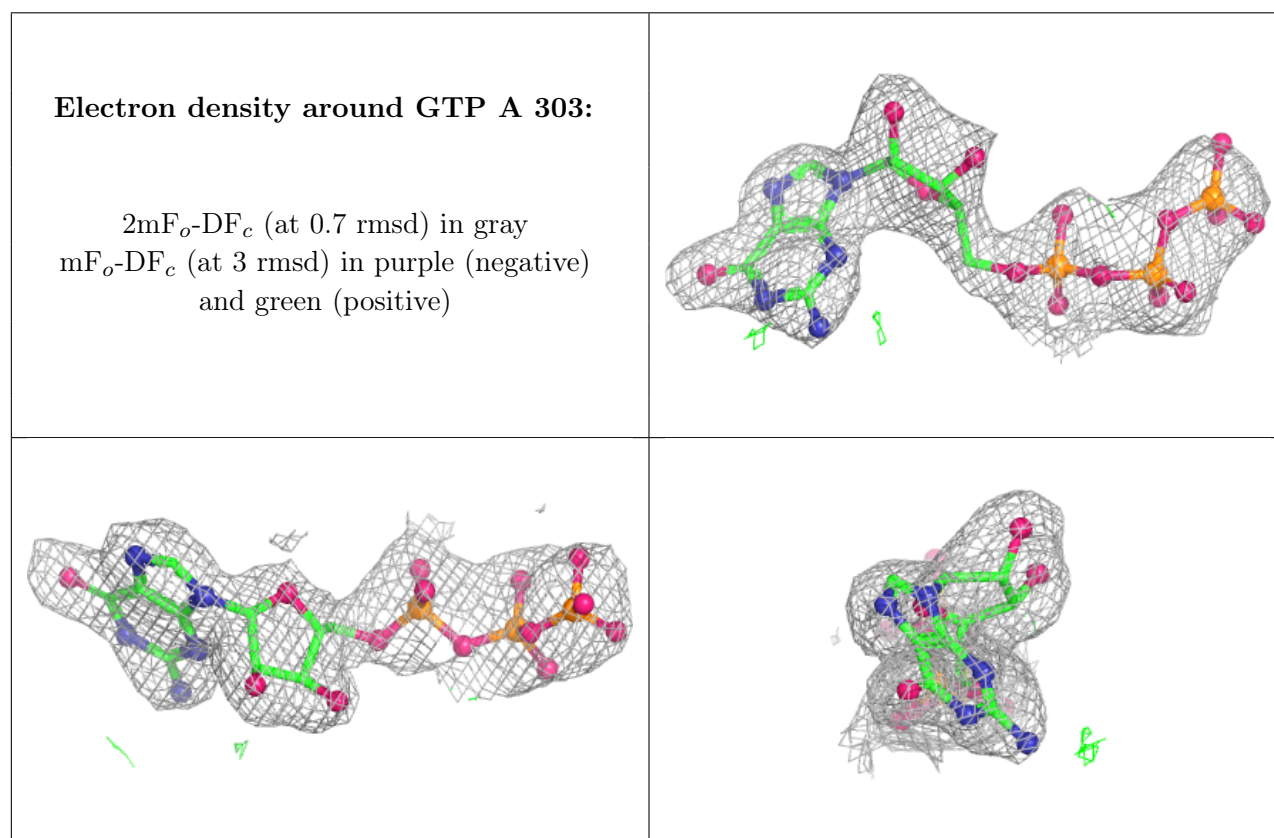
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	PO4	A	805	5/5	0.98	0.05	85,87,88,89	0
2	PO4	D	799	5/5	0.98	0.04	72,73,77,77	0
2	PO4	C	802	5/5	0.98	0.06	98,98,100,100	0
2	PO4	B	800	5/5	0.99	0.06	88,89,91,92	0

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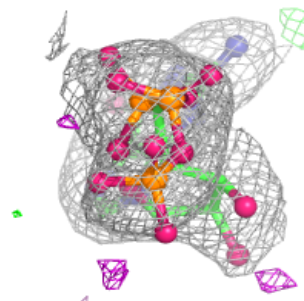
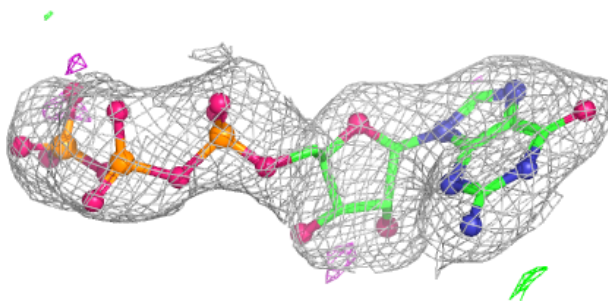
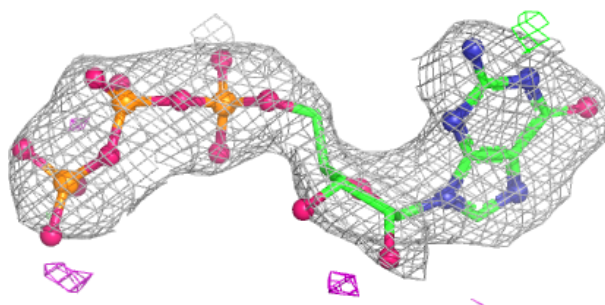
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GTP	A	303	32/32	0.99	0.03	43,53,75,77	0
3	GTP	B	301	32/32	0.99	0.03	34,57,86,86	0
3	GTP	D	302	32/32	0.99	0.03	30,53,79,80	0
3	GTP	C	300	32/32	0.99	0.03	53,58,85,85	0
2	PO4	A	806	5/5	1.00	0.02	31,36,38,38	0
2	PO4	D	801	5/5	1.00	0.02	29,29,34,34	0
2	PO4	B	804	5/5	1.00	0.02	39,40,42,45	0
2	PO4	C	803	5/5	1.00	0.02	36,38,43,44	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.

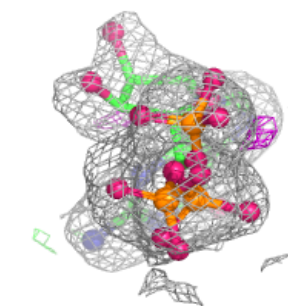
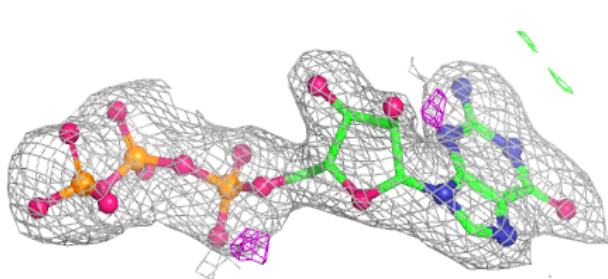
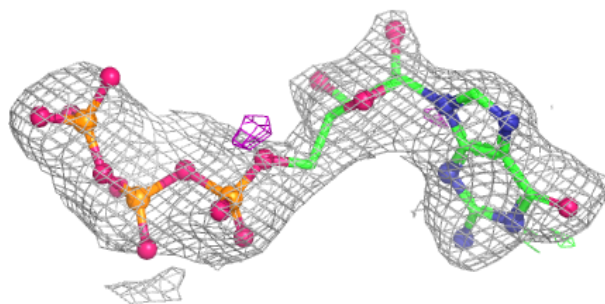


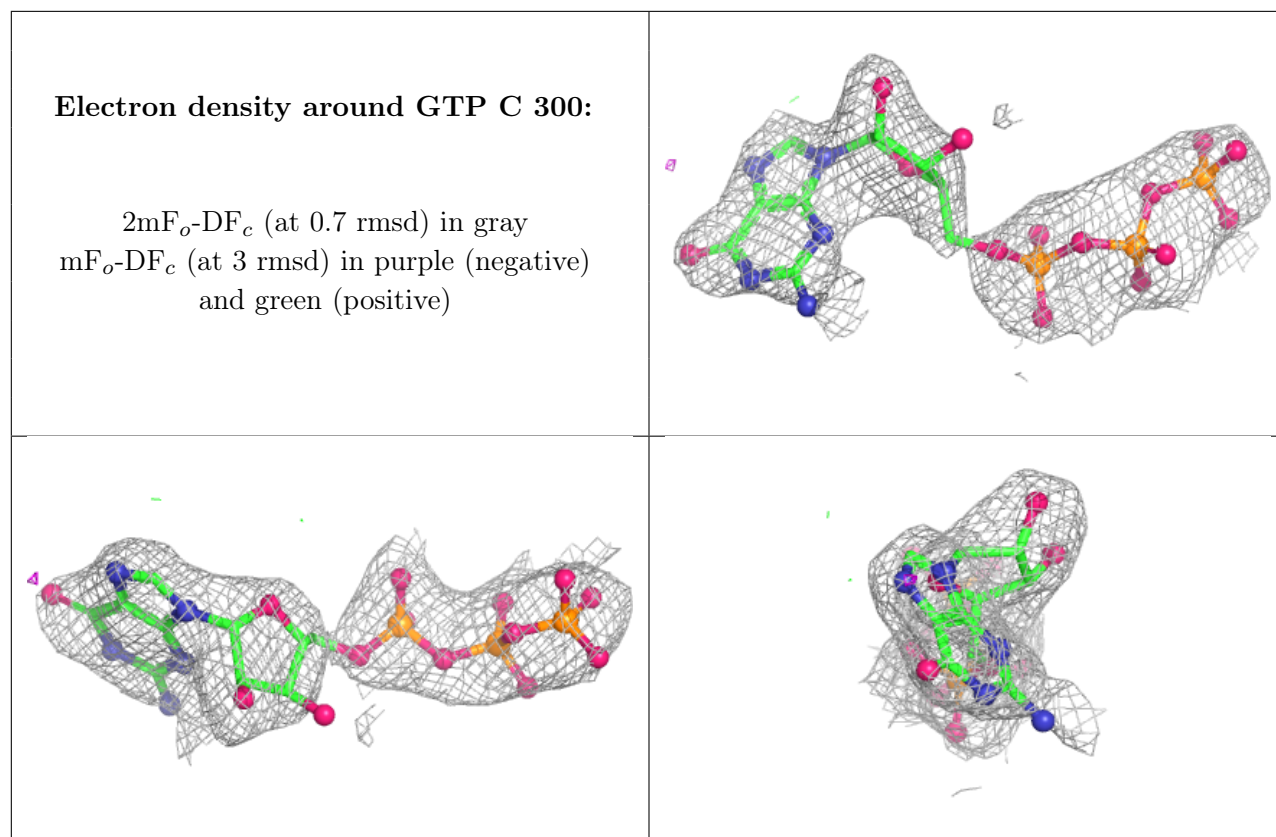
Electron density around GTP B 301:

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

**Electron density around GTP D 302:**

$2mF_o-DF_c$ (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.