



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 09:34 AM EDT

PDB ID : 5XOX / pdb_00005xox
Title : Crystal structure of tRNA(His) guanylyltransferase from *Saccharomyces cerevisiae*
Authors : Lee, K.; Lee, E.H.; Son, J.; Hwang, K.Y.
Deposited on : 2017-05-31
Resolution : 3.00 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:

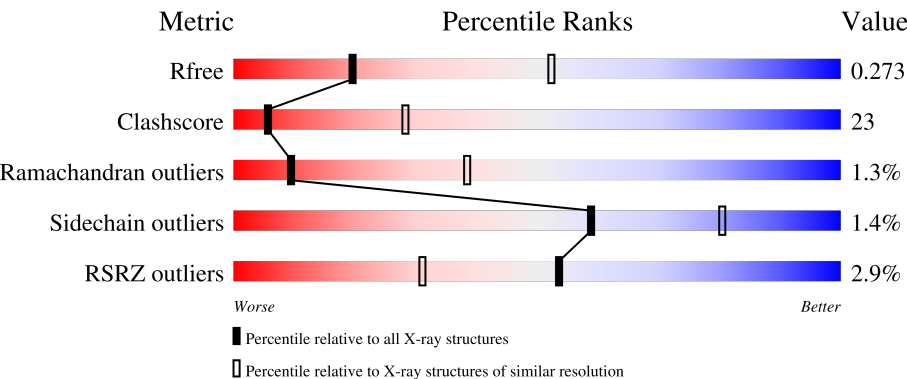
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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.



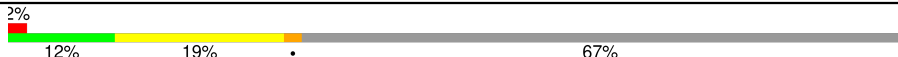
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	61% 30% 6%
1	B	243	54% 38% 6%
1	C	243	3% 56% 37% 6%
1	D	243	4% 58% 33% 6%
1	E	243	3% 55% 37% 6%

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Mol	Chain	Length	Quality of chain
1	F	243	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	PO4	A	304	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10372 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 tRNA(His) guanylyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	229	Total	C	N	O	S	0	0	0
			1898	1238	310	342	8			
1	B	229	Total	C	N	O	S	0	0	0
			1898	1238	310	342	8			
1	C	229	Total	C	N	O	S	0	0	0
			1898	1238	310	342	8			
1	D	229	Total	C	N	O	S	0	0	0
			1898	1238	310	342	8			
1	E	229	Total	C	N	O	S	0	0	0
			1898	1238	310	342	8			
1	F	80	Total	C	N	O	S	0	0	0
			687	455	110	120	2			

There are 36 discrepancies between the modelled and reference sequences:

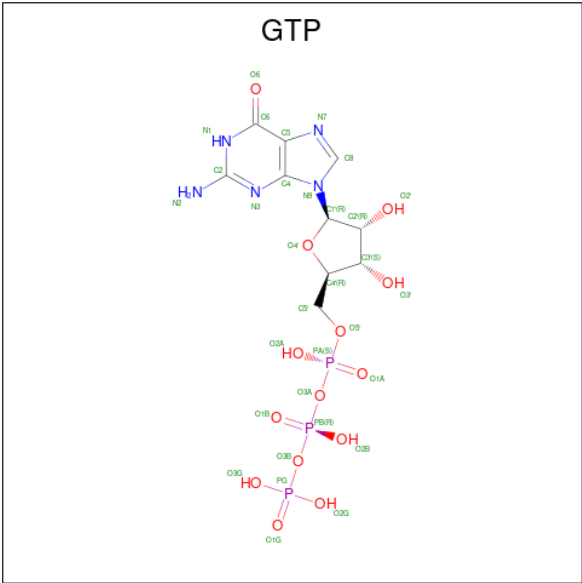
Chain	Residue	Modelled	Actual	Comment	Reference
A	238	HIS	-	expression tag	UNP P53215
A	239	HIS	-	expression tag	UNP P53215
A	240	HIS	-	expression tag	UNP P53215
A	241	HIS	-	expression tag	UNP P53215
A	242	HIS	-	expression tag	UNP P53215
A	243	HIS	-	expression tag	UNP P53215
B	238	HIS	-	expression tag	UNP P53215
B	239	HIS	-	expression tag	UNP P53215
B	240	HIS	-	expression tag	UNP P53215
B	241	HIS	-	expression tag	UNP P53215
B	242	HIS	-	expression tag	UNP P53215
B	243	HIS	-	expression tag	UNP P53215
C	238	HIS	-	expression tag	UNP P53215
C	239	HIS	-	expression tag	UNP P53215
C	240	HIS	-	expression tag	UNP P53215
C	241	HIS	-	expression tag	UNP P53215
C	242	HIS	-	expression tag	UNP P53215

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Chain	Residue	Modelled	Actual	Comment	Reference
C	243	HIS	-	expression tag	UNP P53215
D	238	HIS	-	expression tag	UNP P53215
D	239	HIS	-	expression tag	UNP P53215
D	240	HIS	-	expression tag	UNP P53215
D	241	HIS	-	expression tag	UNP P53215
D	242	HIS	-	expression tag	UNP P53215
D	243	HIS	-	expression tag	UNP P53215
E	238	HIS	-	expression tag	UNP P53215
E	239	HIS	-	expression tag	UNP P53215
E	240	HIS	-	expression tag	UNP P53215
E	241	HIS	-	expression tag	UNP P53215
E	242	HIS	-	expression tag	UNP P53215
E	243	HIS	-	expression tag	UNP P53215
F	238	HIS	-	expression tag	UNP P53215
F	239	HIS	-	expression tag	UNP P53215
F	240	HIS	-	expression tag	UNP P53215
F	241	HIS	-	expression tag	UNP P53215
F	242	HIS	-	expression tag	UNP P53215
F	243	HIS	-	expression tag	UNP P53215

- Molecule 2 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃).



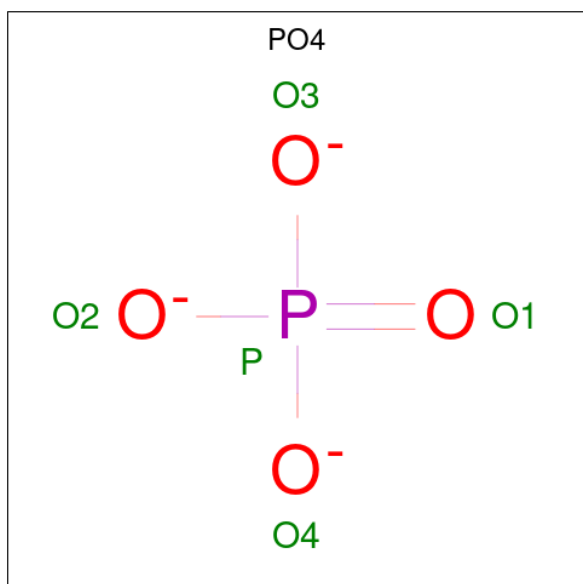
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	D	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	E	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Mg	0	0
			2	2		
3	B	2	Total	Mg	0	0
			2	2		
3	C	2	Total	Mg	0	0
			2	2		
3	D	2	Total	Mg	0	0
			2	2		
3	E	2	Total	Mg	0	0
			2	2		

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		

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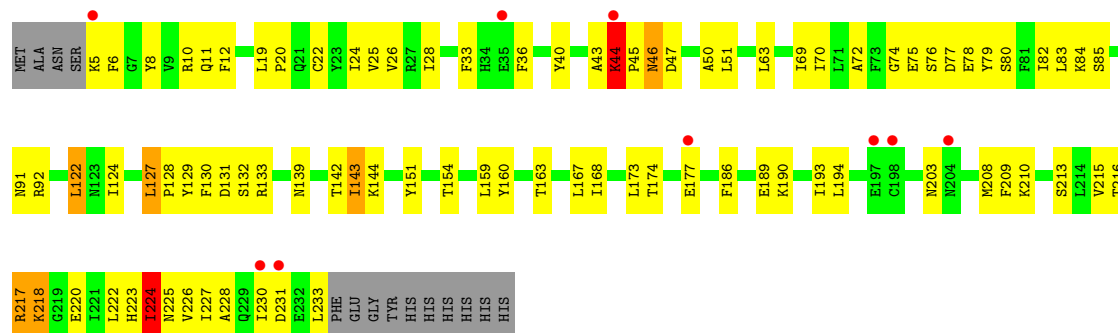
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	P	0	0
			5	4	1		
4	C	1	Total	O	P	0	0
			5	4	1		
4	D	1	Total	O	P	0	0
			5	4	1		
4	E	1	Total	O	P	0	0
			5	4	1		

- Molecule 1: tRNA(His) guanylyltransferase

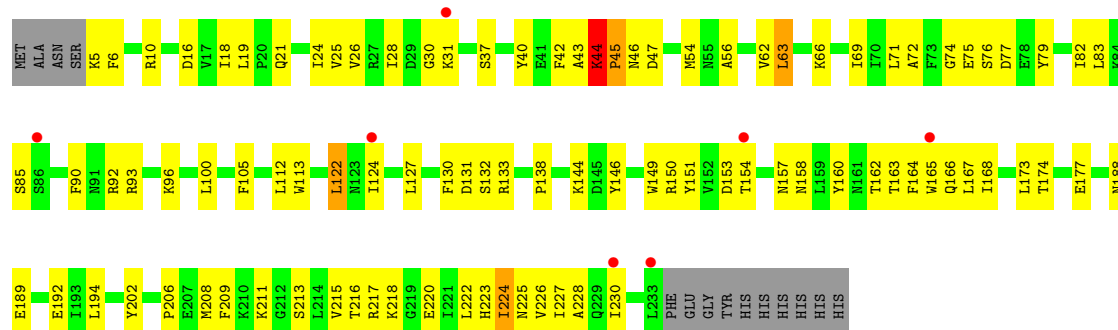




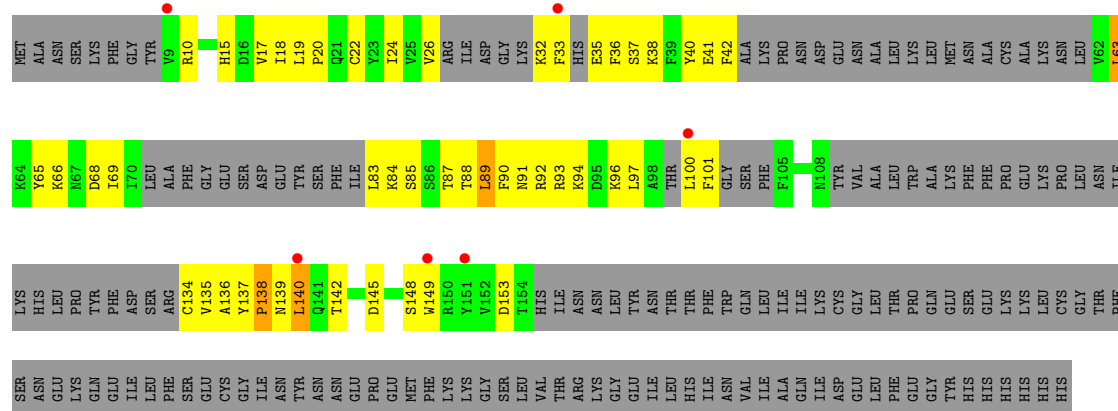
• Molecule 1: tRNA(His) guanylyltransferase



• Molecule 1: tRNA(His) guanylyltransferase



• Molecule 1: tRNA(His) guanylyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	I 41	Depositor
Cell constants a, b, c, α , β , γ	164.70Å 164.70Å 275.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.84 – 3.00 29.84 – 3.00	Depositor EDS
% Data completeness (in resolution range)	93.4 (29.84-3.00) 93.3 (29.84-3.00)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.72 (at 3.00Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.239 , 0.274 0.239 , 0.273	Depositor DCC
R_{free} test set	2026 reflections (2.12%)	wwPDB-VP
Wilson B-factor (Å ²)	71.8	Xtriage
Anisotropy	0.018	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 83.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.066 for h,-k,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10372	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.72% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.85	1/1947 (0.1%)	1.23	13/2628 (0.5%)
1	B	0.80	2/1947 (0.1%)	1.24	12/2628 (0.5%)
1	C	0.78	1/1947 (0.1%)	1.12	10/2628 (0.4%)
1	D	0.84	2/1947 (0.1%)	1.19	17/2628 (0.6%)
1	E	0.72	0/1947	1.14	15/2628 (0.6%)
1	F	0.73	0/698	1.21	5/932 (0.5%)
All	All	0.79	6/10433 (0.1%)	1.19	72/14072 (0.5%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	224	ILE	CA-C	6.35	1.60	1.52
1	A	224	ILE	CA-C	6.18	1.60	1.52
1	B	224	ILE	CA-C	5.73	1.60	1.52
1	C	224	ILE	CA-C	5.37	1.59	1.52
1	B	227	ILE	CA-CB	-5.05	1.47	1.53
1	D	224	ILE	CA-CB	5.05	1.61	1.54

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	44	LYS	CA-C-N	-15.96	99.89	119.84
1	B	44	LYS	C-N-CA	-15.96	99.89	119.84
1	A	44	LYS	CA-C-N	-14.44	101.78	119.84
1	A	44	LYS	C-N-CA	-14.44	101.78	119.84
1	E	44	LYS	CA-C-N	-10.27	107.99	119.28
1	E	44	LYS	C-N-CA	-10.27	107.99	119.28
1	B	137	TYR	CA-C-N	9.13	129.37	119.87
1	B	137	TYR	C-N-CA	9.13	129.37	119.87
1	F	137	TYR	CA-C-N	8.89	130.96	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	137	TYR	C-N-CA	8.89	130.96	119.84
1	D	44	LYS	CA-C-N	-7.75	110.16	119.84
1	D	44	LYS	C-N-CA	-7.75	110.16	119.84
1	D	19	LEU	CA-C-N	7.52	127.57	119.90
1	D	19	LEU	C-N-CA	7.52	127.57	119.90
1	E	45	PRO	CA-C-N	-7.44	112.24	122.95
1	E	45	PRO	C-N-CA	-7.44	112.24	122.95
1	E	122	LEU	N-CA-C	7.17	121.29	112.54
1	B	66	LYS	N-CA-C	-7.13	104.38	113.23
1	E	226	VAL	N-CA-C	-7.08	98.28	108.48
1	A	228	ALA	N-CA-C	-6.96	104.39	113.17
1	A	77	ASP	N-CA-C	-6.93	104.40	112.92
1	E	228	ALA	N-CA-C	-6.85	104.54	113.17
1	C	223	HIS	N-CA-C	6.75	119.35	107.61
1	C	226	VAL	N-CA-C	-6.70	98.47	108.12
1	C	228	ALA	N-CA-C	-6.59	105.27	113.38
1	B	122	LEU	N-CA-C	6.45	120.98	113.18
1	D	226	VAL	CA-C-N	-6.43	114.82	122.22
1	D	226	VAL	C-N-CA	-6.43	114.82	122.22
1	B	16	ASP	N-CA-C	6.43	119.30	110.24
1	A	216	THR	N-CA-C	-6.38	100.47	110.36
1	C	19	LEU	CA-C-N	6.36	126.31	119.76
1	C	19	LEU	C-N-CA	6.36	126.31	119.76
1	C	214	LEU	N-CA-C	6.27	118.93	108.20
1	B	226	VAL	CA-C-N	-6.21	114.71	122.48
1	B	226	VAL	C-N-CA	-6.21	114.71	122.48
1	E	223	HIS	N-CA-C	6.19	118.77	108.13
1	A	224	ILE	N-CA-C	6.16	122.16	109.34
1	A	226	VAL	CA-C-N	-6.06	114.91	122.48
1	A	226	VAL	C-N-CA	-6.06	114.91	122.48
1	D	224	ILE	N-CA-C	6.05	121.92	109.34
1	D	122	LEU	N-CA-C	5.99	120.42	113.18
1	D	77	ASP	N-CA-C	-5.96	105.66	113.17
1	E	16	ASP	N-CA-C	5.93	118.60	110.24
1	D	226	VAL	N-CA-C	-5.89	99.16	107.75
1	B	228	ALA	N-CA-C	-5.87	105.77	113.17
1	F	135	VAL	N-CA-C	5.85	116.03	107.37
1	D	218	LYS	CB-CA-C	-5.83	109.84	116.54
1	B	223	HIS	N-CA-C	5.81	117.52	107.93
1	E	77	ASP	N-CA-C	-5.80	105.86	113.17
1	A	43	ALA	N-CA-C	5.78	117.71	108.41
1	E	44	LYS	N-CA-C	5.75	122.52	109.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	153	ASP	N-CA-C	5.72	115.83	108.34
1	A	122	LEU	N-CA-C	5.69	120.06	113.18
1	D	143	ILE	N-CA-C	-5.67	104.99	110.72
1	D	223	HIS	N-CA-C	5.64	117.31	107.49
1	E	226	VAL	CA-C-N	-5.62	115.45	122.48
1	E	226	VAL	C-N-CA	-5.62	115.45	122.48
1	A	44	LYS	C-N-CD	5.53	147.68	125.00
1	D	217	ARG	N-CA-C	-5.51	102.86	110.35
1	B	224	ILE	N-CA-C	5.43	120.64	109.34
1	D	228	ALA	N-CA-C	-5.37	106.41	113.17
1	C	74	GLY	N-CA-C	5.34	120.81	111.25
1	E	66	LYS	N-CA-C	-5.33	106.77	113.28
1	E	63	LEU	N-CA-C	-5.29	105.21	110.97
1	A	205	GLU	CA-C-N	5.25	125.50	119.83
1	A	205	GLU	C-N-CA	5.25	125.50	119.83
1	F	36	PHE	N-CA-C	-5.21	106.42	112.89
1	D	218	LYS	N-CA-C	5.08	116.03	108.31
1	C	87	THR	N-CA-C	5.07	117.78	110.28
1	C	226	VAL	CA-C-N	-5.04	116.18	122.48
1	C	226	VAL	C-N-CA	-5.04	116.18	122.48
1	D	46	ASN	N-CA-CB	-5.04	102.56	110.77

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	1898	0	1887	85	0
1	B	1898	0	1887	84	1
1	C	1898	0	1887	91	0
1	D	1898	0	1887	77	0
1	E	1898	0	1887	91	1
1	F	687	0	684	67	0
2	A	32	0	12	4	0
2	B	32	0	12	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	32	0	12	5	0
2	D	32	0	12	5	0
2	E	32	0	12	2	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
3	E	2	0	0	0	0
4	A	5	0	0	3	0
4	B	5	0	0	0	0
4	C	5	0	0	1	0
4	D	5	0	0	1	0
4	E	5	0	0	1	0
All	All	10372	0	10179	480	1

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 23.

All (480) 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:F:69:ILE:CG2	1:F:89:LEU:HD11	1.22	1.58
1:F:87:THR:CG2	1:F:89:LEU:HG	1.61	1.31
1:F:69:ILE:CG2	1:F:89:LEU:CD1	2.14	1.23
1:F:87:THR:HG23	1:F:89:LEU:HG	1.20	1.18
1:C:44:LYS:CB	1:C:45:PRO:HD2	1.70	1.15
1:F:69:ILE:HG21	1:F:89:LEU:CD1	1.74	1.15
1:C:44:LYS:HB2	1:C:45:PRO:HD2	1.23	1.14
1:C:44:LYS:HB3	1:C:45:PRO:CD	1.78	1.11
1:C:44:LYS:CB	1:C:45:PRO:CD	2.25	1.11
1:F:69:ILE:HG21	1:F:89:LEU:HD11	1.17	1.10
1:F:69:ILE:HG22	1:F:89:LEU:HD11	1.20	1.06
1:E:44:LYS:HB3	1:E:45:PRO:CD	1.90	1.01
1:F:87:THR:CG2	1:F:89:LEU:CG	2.43	0.96
1:E:75:GLU:HG3	1:E:154:THR:HG22	1.46	0.94
1:F:69:ILE:HG23	1:F:89:LEU:HD11	1.47	0.94
1:A:44:LYS:HA	1:A:44:LYS:CE	1.94	0.94
1:C:44:LYS:HB3	1:C:45:PRO:HD3	1.51	0.92
1:A:44:LYS:HA	1:A:44:LYS:NZ	1.85	0.92
1:F:94:LYS:HZ1	1:F:134:CYS:N	1.66	0.91
1:F:87:THR:HG21	1:F:89:LEU:CD1	2.02	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:8:TYR:O	1:C:11:GLN:HB2	1.72	0.90
1:F:38:LYS:NZ	1:F:40:TYR:O	2.06	0.88
1:D:163:THR:HG22	1:D:194:LEU:HD22	1.57	0.86
1:C:215:VAL:HA	1:C:220:GLU:HB3	1.58	0.86
1:C:43:ALA:HB3	1:C:47:ASP:HA	1.58	0.83
1:B:215:VAL:HG13	1:B:220:GLU:HB3	1.58	0.83
4:A:304:PO4:O1	1:B:133:ARG:NH2	2.12	0.83
1:C:10:ARG:HH12	1:E:192:GLU:CD	1.86	0.82
1:A:44:LYS:HD2	1:A:165:TRP:CD2	2.16	0.81
1:F:83:LEU:HD21	1:F:89:LEU:HD12	1.61	0.81
1:E:76:SER:HB3	2:E:301:GTP:H8	1.46	0.81
1:E:124:ILE:O	1:E:127:LEU:HD13	1.81	0.80
1:E:76:SER:HB3	2:E:301:GTP:C8	2.16	0.79
1:F:38:LYS:HZ2	1:F:40:TYR:C	1.91	0.79
1:A:79:TYR:OH	1:A:215:VAL:HG23	1.85	0.77
1:E:44:LYS:HB3	1:E:45:PRO:HD3	1.66	0.77
1:A:72:ALA:HB1	1:A:79:TYR:HE1	1.48	0.76
1:C:131:ASP:OD1	1:C:132:SER:N	2.19	0.76
1:F:65:TYR:HB2	1:F:69:ILE:HD11	1.67	0.76
1:E:144:LYS:NZ	1:E:230:ILE:HD13	2.00	0.75
1:D:33:PHE:N	2:D:302:GTP:O2B	2.14	0.75
1:F:69:ILE:HG21	1:F:89:LEU:HD13	1.69	0.75
1:C:144:LYS:NZ	1:C:230:ILE:HD13	2.00	0.75
1:A:144:LYS:NZ	1:A:230:ILE:HD13	2.01	0.74
1:E:224:ILE:HG12	1:E:225:ASN:H	1.52	0.74
1:A:33:PHE:N	2:A:301:GTP:O2B	2.21	0.74
1:E:85:SER:O	1:E:92:ARG:NH2	2.20	0.74
1:D:224:ILE:HG12	1:D:225:ASN:H	1.54	0.73
1:E:215:VAL:HG13	1:E:220:GLU:HB3	1.70	0.73
1:F:84:LYS:HZ2	1:F:85:SER:HB3	1.52	0.73
1:D:8:TYR:O	1:D:11:GLN:HB2	1.89	0.72
1:F:87:THR:HG23	1:F:89:LEU:CG	2.11	0.72
1:D:174:THR:N	1:D:177:GLU:OE2	2.18	0.72
1:E:44:LYS:CE	1:E:44:LYS:HA	2.17	0.72
1:D:215:VAL:HG13	1:D:220:GLU:HB3	1.71	0.72
1:C:33:PHE:N	2:C:301:GTP:O2B	2.16	0.71
1:E:43:ALA:HB3	1:E:47:ASP:HA	1.72	0.71
1:A:95:ASP:O	1:A:99:THR:OG1	2.10	0.70
1:C:44:LYS:O	2:C:301:GTP:N2	2.25	0.70
1:F:38:LYS:NZ	1:F:41:GLU:HA	2.07	0.70
1:C:189:GLU:O	1:C:193:ILE:HG12	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:194:LEU:HD23	1:E:202:TYR:HB2	1.74	0.69
1:A:72:ALA:HB1	1:A:79:TYR:CE1	2.27	0.69
1:D:28:ILE:O	1:D:78:GLU:HA	1.92	0.69
1:C:40:TYR:OH	1:C:122:LEU:O	2.11	0.68
1:B:48:GLU:HG2	1:B:52:LYS:HE2	1.75	0.68
1:E:43:ALA:N	1:E:47:ASP:OD2	2.23	0.68
1:C:43:ALA:HB3	1:C:47:ASP:CA	2.24	0.68
1:C:151:TYR:CZ	1:C:224:ILE:HA	2.28	0.67
1:B:33:PHE:N	2:B:302:GTP:O1B	2.26	0.67
1:D:24:ILE:HB	1:D:83:LEU:HB2	1.76	0.67
1:B:224:ILE:HG12	1:B:225:ASN:H	1.59	0.67
1:C:163:THR:HG22	1:C:194:LEU:HD22	1.77	0.67
1:C:76:SER:HB3	2:C:301:GTP:C8	2.29	0.67
1:F:138:PRO:HG2	1:F:142:THR:HG21	1.75	0.67
1:D:129:TYR:C	1:D:130:PHE:HD1	2.02	0.67
1:E:138:PRO:O	1:F:139:ASN:ND2	2.28	0.66
1:A:62:VAL:HA	1:A:69:ILE:HD12	1.76	0.66
1:A:144:LYS:HZ2	1:A:230:ILE:HD13	1.60	0.66
1:B:151:TYR:CZ	1:B:224:ILE:HA	2.30	0.66
1:F:87:THR:HG21	1:F:89:LEU:HD12	1.76	0.66
1:A:44:LYS:HB3	1:A:45:PRO:CD	2.25	0.66
1:B:163:THR:HG22	1:B:194:LEU:HD22	1.76	0.66
1:F:69:ILE:HG22	1:F:89:LEU:CD1	2.04	0.66
1:B:91:ASN:O	1:B:92:ARG:HB2	1.96	0.66
1:C:48:GLU:HG2	1:C:52:LYS:HE2	1.77	0.66
1:D:85:SER:O	1:D:92:ARG:NH2	2.29	0.66
1:A:96:LYS:O	1:A:100:LEU:HB2	1.95	0.65
1:D:163:THR:HG21	1:D:190:LYS:HB3	1.77	0.65
1:E:62:VAL:HG12	1:E:69:ILE:HG21	1.78	0.65
1:F:94:LYS:NZ	1:F:134:CYS:N	2.45	0.65
1:C:85:SER:O	1:C:92:ARG:NH2	2.30	0.65
1:B:144:LYS:HZ1	1:B:230:ILE:HD13	1.62	0.64
1:A:151:TYR:CZ	1:A:224:ILE:HA	2.31	0.64
1:D:70:ILE:HD11	1:D:84:LYS:HA	1.80	0.64
1:B:96:LYS:O	1:B:100:LEU:HB2	1.98	0.64
1:F:87:THR:O	1:F:92:ARG:NH1	2.31	0.63
1:B:40:TYR:OH	1:B:122:LEU:O	2.17	0.63
1:D:76:SER:HB3	2:D:302:GTP:H8	1.62	0.63
1:E:113:TRP:CD1	1:E:122:LEU:HD22	2.33	0.63
1:E:5:LYS:N	1:E:5:LYS:HD3	2.14	0.63
1:F:32:LYS:NZ	1:F:33:PHE:O	2.32	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:174:THR:N	1:B:177:GLU:OE2	2.28	0.62
1:F:38:LYS:HZ3	1:F:41:GLU:HA	1.62	0.62
1:A:79:TYR:OH	1:A:215:VAL:CG2	2.47	0.62
1:E:40:TYR:OH	1:E:122:LEU:O	2.12	0.62
1:B:76:SER:HB3	2:B:302:GTP:H8	1.65	0.62
1:C:163:THR:HG21	1:C:190:LYS:HB3	1.80	0.62
1:D:44:LYS:O	1:D:46:ASN:N	2.33	0.62
1:A:44:LYS:HA	1:A:44:LYS:HZ3	1.61	0.61
1:D:167:LEU:HB3	1:D:173:LEU:HD22	1.81	0.61
1:E:131:ASP:OD1	1:E:132:SER:N	2.33	0.61
1:A:129:TYR:C	1:A:130:PHE:HD1	2.07	0.61
1:D:51:LEU:HB3	1:D:208:MET:HE1	1.82	0.61
1:D:144:LYS:NZ	1:D:230:ILE:HD13	2.15	0.61
1:E:96:LYS:O	1:E:100:LEU:HB2	2.00	0.61
1:E:45:PRO:HB2	1:E:46:ASN:HD22	1.66	0.61
1:D:174:THR:OG1	1:D:177:GLU:OE1	2.18	0.61
1:E:44:LYS:HD2	1:E:165:TRP:CZ2	2.36	0.61
1:A:224:ILE:HG12	1:A:225:ASN:H	1.66	0.61
1:D:217:ARG:C	1:D:218:LYS:HD3	2.26	0.61
1:A:90:PHE:HB3	1:A:96:LYS:HD3	1.81	0.60
1:E:30:GLY:HA2	1:E:130:PHE:HA	1.84	0.60
1:A:30:GLY:HA3	1:A:33:PHE:CE1	2.35	0.60
1:C:24:ILE:HB	1:C:83:LEU:HB2	1.83	0.60
1:E:18:ILE:HG22	1:E:92:ARG:HH11	1.66	0.60
1:C:113:TRP:CG	1:C:122:LEU:HD21	2.36	0.60
1:D:43:ALA:HB3	1:D:47:ASP:HA	1.84	0.59
1:A:77:ASP:OD2	2:A:301:GTP:O1A	2.20	0.59
1:A:40:TYR:OH	1:A:122:LEU:O	2.21	0.59
1:C:215:VAL:HA	1:C:220:GLU:CB	2.29	0.59
1:C:19:LEU:O	1:C:92:ARG:NH1	2.36	0.59
1:C:72:ALA:HA	1:C:80:SER:O	2.02	0.59
1:F:17:VAL:HG13	1:F:92:ARG:HB2	1.84	0.59
1:C:144:LYS:HZ1	1:C:230:ILE:HD13	1.67	0.58
1:C:144:LYS:HZ2	1:C:230:ILE:HD13	1.68	0.58
1:E:188:ASN:OD1	1:E:189:GLU:N	2.36	0.58
1:A:113:TRP:CD2	1:A:122:LEU:HD21	2.38	0.58
1:D:44:LYS:O	1:D:45:PRO:C	2.45	0.58
1:A:30:GLY:HA3	1:A:33:PHE:CD1	2.39	0.58
1:C:10:ARG:NH1	1:E:192:GLU:OE1	2.36	0.58
1:D:131:ASP:OD1	1:D:132:SER:N	2.36	0.58
1:C:44:LYS:O	2:C:301:GTP:C2	2.56	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:84:LYS:NZ	1:F:85:SER:HB3	2.17	0.58
1:B:211:LYS:HB3	1:B:224:ILE:H	1.67	0.58
1:F:97:LEU:O	1:F:101:PHE:HB3	2.03	0.58
1:A:215:VAL:HA	1:A:220:GLU:HB3	1.86	0.58
1:A:12:PHE:CE1	1:C:20:PRO:HB2	2.38	0.58
1:A:163:THR:HG22	1:A:194:LEU:HD22	1.84	0.58
1:B:124:ILE:O	1:B:127:LEU:HD13	2.04	0.58
1:B:217:ARG:HB2	1:B:218:LYS:HD3	1.84	0.58
1:D:213:SER:HA	1:D:222:LEU:HD23	1.86	0.58
1:F:68:ASP:O	1:F:87:THR:OG1	2.22	0.58
1:B:217:ARG:C	1:B:218:LYS:HD3	2.29	0.57
1:E:151:TYR:CZ	1:E:224:ILE:HA	2.39	0.57
1:B:76:SER:HB3	2:B:302:GTP:C8	2.39	0.57
1:E:217:ARG:C	1:E:218:LYS:HD3	2.30	0.57
1:B:151:TYR:HE1	1:B:211:LYS:HA	1.68	0.57
1:B:167:LEU:HB3	1:B:173:LEU:HD22	1.87	0.57
1:C:69:ILE:HG12	1:C:83:LEU:CD2	2.35	0.57
1:E:44:LYS:HA	1:E:44:LYS:HE2	1.84	0.57
1:E:24:ILE:HB	1:E:83:LEU:HB2	1.87	0.57
1:E:144:LYS:HZ1	1:E:230:ILE:HD13	1.70	0.57
1:C:129:TYR:C	1:C:130:PHE:HD1	2.12	0.56
1:A:44:LYS:HB3	1:A:45:PRO:HD3	1.87	0.56
1:A:72:ALA:CB	1:A:79:TYR:HE1	2.18	0.56
1:E:43:ALA:CB	1:E:47:ASP:OD2	2.53	0.56
1:D:45:PRO:C	1:D:46:ASN:HD22	2.14	0.56
1:F:83:LEU:CD2	1:F:89:LEU:HD12	2.35	0.56
1:F:145:ASP:HA	1:F:148:SER:HB2	1.88	0.56
1:A:174:THR:OG1	1:A:177:GLU:OE1	2.22	0.56
1:F:18:ILE:HD11	1:F:94:LYS:HG2	1.87	0.55
1:A:93:ARG:NH2	4:A:304:PO4:O3	2.40	0.55
1:C:70:ILE:HG22	1:C:71:LEU:HD23	1.88	0.55
1:E:144:LYS:HZ2	1:E:230:ILE:HD13	1.71	0.55
1:C:79:TYR:HB3	1:C:81:PHE:CE1	2.42	0.55
1:F:18:ILE:HG21	1:F:24:ILE:HG12	1.87	0.55
1:F:89:LEU:O	1:F:90:PHE:C	2.49	0.55
1:A:215:VAL:HG12	1:A:216:THR:O	2.07	0.55
1:A:113:TRP:CG	1:A:122:LEU:HD21	2.41	0.55
1:B:105:PHE:CE1	1:B:130:PHE:HE2	2.24	0.55
1:A:44:LYS:HD2	1:A:165:TRP:CE3	2.42	0.55
1:F:140:LEU:HD13	1:F:140:LEU:H	1.72	0.55
1:D:76:SER:HB3	2:D:302:GTP:C8	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:144:LYS:HZ2	1:D:230:ILE:HD13	1.72	0.55
1:F:19:LEU:O	1:F:92:ARG:NH2	2.40	0.55
1:E:26:VAL:HA	1:E:133:ARG:O	2.06	0.55
1:F:68:ASP:HB2	1:F:89:LEU:HD21	1.88	0.55
1:F:15:HIS:HB3	1:F:93:ARG:HD2	1.89	0.54
1:A:30:GLY:HA2	1:A:130:PHE:HA	1.90	0.54
1:C:225:ASN:O	1:C:225:ASN:ND2	2.40	0.54
1:B:129:TYR:C	1:B:130:PHE:HD1	2.15	0.54
1:E:206:PRO:HD2	1:E:209:PHE:CD2	2.43	0.54
1:A:208:MET:CB	1:A:222:LEU:HD13	2.38	0.54
1:C:192:GLU:HA	1:C:192:GLU:OE2	2.07	0.54
1:F:87:THR:HG22	1:F:89:LEU:HG	1.78	0.54
1:F:63:LEU:HD11	1:F:66:LYS:HB2	1.89	0.54
1:C:18:ILE:CD1	1:C:94:LYS:HG2	2.38	0.54
1:E:44:LYS:HB3	1:E:45:PRO:HD2	1.88	0.54
1:E:150:ARG:O	1:E:154:THR:HG23	2.07	0.53
1:A:18:ILE:HG22	1:A:92:ARG:HH11	1.74	0.53
1:F:63:LEU:HD11	1:F:66:LYS:HD3	1.90	0.53
1:A:189:GLU:O	1:A:193:ILE:HG12	2.08	0.53
1:C:213:SER:HA	1:C:222:LEU:CD2	2.38	0.53
1:A:144:LYS:HZ1	1:A:230:ILE:HD13	1.73	0.53
1:B:150:ARG:O	1:B:154:THR:HG23	2.09	0.53
1:D:82:ILE:HD12	1:D:143:ILE:HD11	1.91	0.53
1:B:70:ILE:HD11	1:B:84:LYS:HA	1.91	0.53
1:D:91:ASN:O	1:D:92:ARG:HB2	2.09	0.53
1:D:10:ARG:NH2	4:D:301:PO4:O1	2.42	0.53
1:D:189:GLU:O	1:D:193:ILE:HG12	2.08	0.53
1:A:208:MET:HB2	1:A:222:LEU:HD13	1.90	0.52
1:B:113:TRP:CG	1:B:122:LEU:HD21	2.44	0.52
1:D:75:GLU:HG3	1:D:154:THR:HG22	1.90	0.52
1:C:18:ILE:HD12	1:C:94:LYS:HG2	1.92	0.52
1:C:224:ILE:HG12	1:C:225:ASN:H	1.73	0.52
1:D:203:ASN:HA	1:D:210:LYS:NZ	2.25	0.52
1:B:189:GLU:O	1:B:193:ILE:HG12	2.09	0.52
1:C:216:THR:OG1	1:C:217:ARG:N	2.43	0.52
1:D:47:ASP:HB3	1:D:50:ALA:HB3	1.91	0.52
1:B:123:ASN:CG	1:B:124:ILE:H	2.18	0.52
1:A:213:SER:HA	1:A:222:LEU:HD23	1.90	0.52
1:D:168:ILE:HG23	1:D:173:LEU:O	2.10	0.52
1:E:43:ALA:HB3	1:E:47:ASP:CA	2.38	0.52
1:A:140:LEU:O	1:A:143:ILE:HG22	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:26:VAL:HA	1:D:133:ARG:O	2.10	0.52
1:B:45:PRO:HG3	1:B:165:TRP:HD1	1.74	0.52
1:C:66:LYS:O	1:C:217:ARG:NH2	2.43	0.52
1:D:231:ASP:O	1:D:233:LEU:HD13	2.10	0.52
1:E:206:PRO:HD2	1:E:209:PHE:HD2	1.74	0.52
1:D:44:LYS:HA	1:D:44:LYS:HE2	1.92	0.51
1:E:133:ARG:HH22	4:E:304:PO4:P	2.33	0.51
1:A:194:LEU:HD23	1:A:202:TYR:HB2	1.92	0.51
1:B:113:TRP:CD2	1:B:122:LEU:HD21	2.46	0.51
1:C:224:ILE:HG23	1:C:225:ASN:N	2.26	0.51
1:D:213:SER:HA	1:D:222:LEU:CD2	2.40	0.51
1:E:71:LEU:O	1:E:82:ILE:HB	2.11	0.51
1:E:44:LYS:CE	1:E:44:LYS:CA	2.86	0.51
1:E:173:LEU:HG	1:E:177:GLU:HG3	1.91	0.51
1:A:227:ILE:O	1:A:230:ILE:HG22	2.11	0.50
1:C:74:GLY:HA3	1:C:79:TYR:CE2	2.46	0.50
1:A:150:ARG:O	1:A:154:THR:HG23	2.11	0.50
1:E:54:MET:HE3	1:E:79:TYR:HD2	1.76	0.50
1:A:85:SER:O	1:A:92:ARG:NH2	2.45	0.50
1:F:18:ILE:HG23	1:F:136:ALA:HB2	1.94	0.50
1:F:26:VAL:HG22	1:F:97:LEU:HD21	1.93	0.50
1:E:25:VAL:HG22	1:E:82:ILE:HD13	1.93	0.50
1:B:35:GLU:H	1:B:35:GLU:CD	2.19	0.50
1:D:159:LEU:O	1:D:163:THR:HG23	2.12	0.50
1:D:215:VAL:HG12	1:D:216:THR:O	2.12	0.49
1:E:211:LYS:HD2	1:E:224:ILE:H	1.77	0.49
1:A:213:SER:HA	1:A:222:LEU:CD2	2.41	0.49
1:D:40:TYR:OH	1:D:122:LEU:O	2.27	0.49
1:B:33:PHE:HA	1:B:36:PHE:HB3	1.94	0.49
1:C:62:VAL:HG12	1:C:69:ILE:HG21	1.94	0.49
1:A:99:THR:HG21	1:B:132:SER:HB3	1.95	0.49
1:C:103:SER:CB	1:D:130:PHE:H	2.25	0.49
1:B:216:THR:HG23	1:B:217:ARG:O	2.13	0.49
1:B:174:THR:OG1	1:B:177:GLU:OE1	2.19	0.49
1:E:215:VAL:HG12	1:E:216:THR:O	2.13	0.49
1:C:215:VAL:HG12	1:C:216:THR:O	2.13	0.49
1:C:46:ASN:HB3	1:C:209:PHE:CE1	2.49	0.48
1:C:151:TYR:CE1	1:C:224:ILE:HA	2.48	0.48
1:D:72:ALA:HA	1:D:80:SER:O	2.13	0.48
1:F:88:THR:O	1:F:88:THR:OG1	2.28	0.48
1:A:192:GLU:OE2	1:A:192:GLU:HA	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:45:PRO:C	1:E:46:ASN:HD22	2.22	0.48
1:E:163:THR:HG22	1:E:194:LEU:HD22	1.95	0.48
1:D:203:ASN:HA	1:D:210:LYS:HZ2	1.79	0.48
1:D:227:ILE:O	1:D:230:ILE:HG22	2.14	0.48
1:E:217:ARG:HB2	1:E:218:LYS:HD3	1.96	0.48
1:C:8:TYR:HA	1:C:11:GLN:OE1	2.14	0.48
1:C:113:TRP:CD1	1:C:122:LEU:HD21	2.49	0.48
1:D:151:TYR:CZ	1:D:224:ILE:HA	2.48	0.48
1:A:39:PHE:HD1	1:A:40:TYR:CD1	2.32	0.48
1:A:215:VAL:HA	1:A:220:GLU:CB	2.44	0.48
1:B:55:ASN:HD21	1:B:213:SER:HB3	1.78	0.48
1:E:43:ALA:HB3	1:E:47:ASP:CB	2.43	0.48
1:E:215:VAL:HG22	1:E:220:GLU:HB3	1.96	0.48
1:A:20:PRO:HB2	1:C:12:PHE:CE1	2.49	0.47
1:F:66:LYS:O	1:F:69:ILE:HG12	2.14	0.47
1:A:28:ILE:O	1:A:78:GLU:HA	2.15	0.47
1:C:194:LEU:HD12	1:C:194:LEU:HA	1.66	0.47
1:E:213:SER:HA	1:E:222:LEU:HD22	1.96	0.47
1:B:73:PHE:CZ	1:B:154:THR:HG21	2.49	0.47
1:B:100:LEU:HA	1:B:100:LEU:HD23	1.78	0.47
1:B:225:ASN:O	1:B:225:ASN:ND2	2.46	0.47
1:D:186:PHE:H	1:D:189:GLU:HB2	1.79	0.47
1:F:35:GLU:N	1:F:37:SER:HG	2.12	0.47
1:B:45:PRO:HG3	1:B:165:TRP:CD1	2.49	0.47
1:C:77:ASP:OD2	2:C:301:GTP:O2A	2.32	0.47
1:F:18:ILE:HG22	1:F:19:LEU:N	2.30	0.47
1:A:129:TYR:C	1:A:130:PHE:CD1	2.92	0.47
1:C:131:ASP:OD1	1:C:131:ASP:C	2.56	0.47
1:A:139:ASN:O	1:A:142:THR:OG1	2.29	0.47
1:B:24:ILE:HB	1:B:83:LEU:HB2	1.97	0.47
1:C:213:SER:HA	1:C:222:LEU:HD23	1.97	0.47
1:D:46:ASN:OD1	1:D:209:PHE:CG	2.67	0.47
1:A:106:THR:O	1:A:110:VAL:HG23	2.15	0.47
1:C:5:LYS:CG	1:C:6:PHE:H	2.27	0.47
1:D:5:LYS:CG	1:D:6:PHE:H	2.28	0.47
1:D:33:PHE:O	2:D:302:GTP:H1'	2.15	0.47
1:B:12:PHE:CE1	1:D:20:PRO:HB2	2.50	0.46
1:E:194:LEU:HD12	1:E:194:LEU:HA	1.78	0.46
1:B:44:LYS:HD3	1:B:44:LYS:HA	1.50	0.46
1:B:62:VAL:HA	1:B:69:ILE:HD12	1.98	0.46
1:E:19:LEU:O	1:E:92:ARG:NH1	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:100:LEU:O	1:F:101:PHE:HB2	2.15	0.46
1:C:28:ILE:O	1:C:78:GLU:HA	2.15	0.46
1:C:70:ILE:HD11	1:C:84:LYS:HA	1.96	0.46
1:E:158:ASN:O	1:E:162:THR:OG1	2.19	0.46
1:A:131:ASP:OD1	1:A:132:SER:N	2.48	0.46
1:A:188:ASN:OD1	1:A:189:GLU:N	2.49	0.46
1:B:151:TYR:CE1	1:B:224:ILE:HA	2.51	0.46
1:B:224:ILE:HG23	1:B:225:ASN:N	2.30	0.46
1:D:217:ARG:HB2	1:D:218:LYS:HD3	1.97	0.46
1:A:43:ALA:HB3	1:A:47:ASP:HA	1.98	0.46
1:B:131:ASP:OD1	1:B:132:SER:N	2.49	0.46
1:B:69:ILE:HG12	1:B:83:LEU:CD2	2.46	0.46
1:C:35:GLU:H	1:C:35:GLU:CD	2.23	0.46
1:C:113:TRP:CD1	1:C:122:LEU:CD2	2.99	0.46
1:D:163:THR:HA	1:D:194:LEU:HD13	1.97	0.46
1:A:76:SER:HB3	2:A:301:GTP:C8	2.51	0.46
1:C:10:ARG:NH1	1:E:192:GLU:OE2	2.34	0.46
1:C:10:ARG:NH2	4:C:304:PO4:O3	2.44	0.46
1:C:44:LYS:HA	1:C:44:LYS:HD3	1.82	0.46
1:D:124:ILE:O	1:D:127:LEU:HD13	2.16	0.46
1:F:19:LEU:HB3	1:F:22:CYS:SG	2.55	0.46
1:D:216:THR:HG23	1:D:217:ARG:O	2.16	0.45
1:E:46:ASN:OD1	1:E:209:PHE:CG	2.69	0.45
1:E:144:LYS:HE2	1:E:227:ILE:HG21	1.98	0.45
1:E:211:LYS:HB3	1:E:224:ILE:H	1.81	0.45
1:A:215:VAL:HG13	1:A:220:GLU:HB3	1.98	0.45
1:B:71:LEU:HD13	1:B:214:LEU:HD11	1.97	0.45
1:E:45:PRO:CB	1:E:46:ASN:HD22	2.29	0.45
1:E:227:ILE:O	1:E:230:ILE:HG22	2.16	0.45
1:A:130:PHE:H	1:B:103:SER:HB3	1.82	0.45
1:B:186:PHE:O	1:B:190:LYS:HG3	2.15	0.45
1:C:113:TRP:CD2	1:C:122:LEU:HD21	2.51	0.45
1:E:208:MET:HE2	1:E:208:MET:HB3	1.88	0.45
1:F:84:LYS:HG3	1:F:85:SER:N	2.31	0.45
1:B:208:MET:CA	1:B:222:LEU:HD13	2.47	0.45
1:C:33:PHE:HE1	1:C:130:PHE:HE1	1.64	0.45
1:E:71:LEU:HD12	1:E:72:ALA:N	2.32	0.45
1:F:87:THR:HG22	1:F:89:LEU:H	1.82	0.45
1:B:55:ASN:HD21	1:B:213:SER:CB	2.30	0.45
1:B:217:ARG:HG3	1:B:232:GLU:OE2	2.17	0.45
2:D:302:GTP:O3A	2:D:302:GTP:H4'	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:152:VAL:O	1:C:156:ILE:HG12	2.16	0.45
1:A:45:PRO:HG3	1:A:165:TRP:CD1	2.52	0.45
1:C:61:LEU:HA	1:C:104:PHE:HE2	1.81	0.45
1:D:44:LYS:C	1:D:46:ASN:N	2.75	0.45
1:A:33:PHE:O	1:A:37:SER:OG	2.28	0.45
1:E:37:SER:HB2	1:E:42:PHE:HB2	1.98	0.45
1:E:174:THR:O	1:E:177:GLU:HG2	2.16	0.45
1:E:160:TYR:CD1	1:E:160:TYR:C	2.95	0.44
1:E:167:LEU:HD23	1:E:167:LEU:HA	1.86	0.44
1:A:131:ASP:OD1	1:A:131:ASP:C	2.60	0.44
1:B:70:ILE:HD12	1:B:82:ILE:HG22	1.98	0.44
1:E:44:LYS:HD2	1:E:165:TRP:CH2	2.51	0.44
1:B:174:THR:HB	1:B:175:PRO:HD2	2.00	0.44
1:D:46:ASN:OD1	1:D:209:PHE:CD1	2.71	0.44
1:B:20:PRO:HB2	1:D:12:PHE:CE2	2.52	0.44
1:D:139:ASN:O	1:D:142:THR:OG1	2.34	0.44
1:D:224:ILE:CG1	1:D:225:ASN:H	2.26	0.44
1:A:44:LYS:HD2	1:A:165:TRP:CG	2.52	0.44
1:A:100:LEU:HD23	1:A:100:LEU:HA	1.73	0.44
1:B:200:ILE:HD12	1:B:200:ILE:HA	1.86	0.44
1:D:44:LYS:HB3	1:D:45:PRO:CD	2.47	0.44
1:A:17:VAL:HG12	1:A:93:ARG:HA	2.00	0.44
1:E:43:ALA:HB3	1:E:47:ASP:CG	2.42	0.44
1:E:75:GLU:CG	1:E:154:THR:HG22	2.33	0.44
1:A:54:MET:HE3	1:A:54:MET:HB3	1.92	0.44
1:C:82:ILE:HG23	1:C:143:ILE:HD11	2.00	0.44
1:D:70:ILE:HD12	1:D:82:ILE:HG22	2.00	0.44
1:F:19:LEU:O	1:F:22:CYS:HB2	2.17	0.44
1:B:29:ASP:OD1	1:B:77:ASP:OD2	2.35	0.43
1:B:177:GLU:O	1:B:181:LYS:HB2	2.19	0.43
1:C:82:ILE:HD12	1:C:143:ILE:HD11	2.00	0.43
1:C:227:ILE:O	1:C:230:ILE:HG22	2.18	0.43
1:D:160:TYR:HD2	1:D:190:LYS:HE2	1.83	0.43
1:A:205:GLU:HA	1:A:206:PRO:HD3	1.83	0.43
1:B:68:ASP:OD2	1:B:87:THR:OG1	2.25	0.43
1:C:140:LEU:HA	1:C:143:ILE:HG22	2.01	0.43
1:E:216:THR:OG1	1:E:217:ARG:N	2.51	0.43
1:F:66:LYS:HA	1:F:69:ILE:O	2.18	0.43
1:B:19:LEU:HA	1:B:20:PRO:HD3	1.84	0.43
1:B:213:SER:HA	1:B:222:LEU:CD2	2.48	0.43
1:C:97:LEU:HA	1:C:97:LEU:HD23	1.77	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:150:ARG:O	1:C:154:THR:HG23	2.18	0.43
1:E:21:GLN:HA	1:E:85:SER:O	2.17	0.43
1:C:99:THR:HG21	1:D:132:SER:O	2.18	0.43
1:E:56:ALA:HB1	1:E:112:LEU:HD13	2.00	0.43
1:E:90:PHE:O	1:E:93:ARG:HB2	2.18	0.43
1:F:24:ILE:HG23	1:F:94:LYS:HZ2	1.83	0.43
1:A:211:LYS:HB3	1:A:224:ILE:H	1.84	0.43
1:A:69:ILE:HG12	1:A:83:LEU:CD2	2.49	0.43
1:B:215:VAL:HG12	1:B:216:THR:O	2.19	0.43
1:D:51:LEU:CB	1:D:208:MET:HE1	2.48	0.43
1:E:105:PHE:CD1	1:E:130:PHE:HE2	2.37	0.43
1:F:19:LEU:HA	1:F:20:PRO:HD2	1.83	0.43
1:F:93:ARG:HG2	1:F:96:LYS:HG2	2.00	0.43
4:A:304:PO4:P	1:B:133:ARG:HH21	2.42	0.43
1:C:122:LEU:HD22	1:C:122:LEU:HA	1.65	0.43
1:D:69:ILE:HG12	1:D:83:LEU:HD23	2.01	0.43
1:D:74:GLY:N	1:D:79:TYR:CE1	2.87	0.43
1:F:38:LYS:HD3	1:F:42:PHE:N	2.33	0.43
1:F:96:LYS:HA	1:F:96:LYS:HD3	1.50	0.43
1:A:76:SER:HB3	2:A:301:GTP:H8	1.84	0.43
1:B:194:LEU:HD12	1:B:194:LEU:HA	1.81	0.43
1:B:195:PHE:HE2	1:F:10:ARG:HH12	1.67	0.43
1:C:54:MET:HE3	1:C:54:MET:HB3	1.64	0.43
1:E:146:TYR:O	1:E:149:TRP:HB3	2.19	0.43
1:A:200:ILE:HD12	1:A:200:ILE:HA	1.78	0.42
1:B:45:PRO:HD3	1:B:165:TRP:CD1	2.54	0.42
1:B:139:ASN:O	1:B:142:THR:OG1	2.36	0.42
1:B:164:PHE:O	1:B:168:ILE:HG12	2.19	0.42
1:D:44:LYS:HA	1:D:44:LYS:CE	2.48	0.42
1:D:22:CYS:O	1:D:85:SER:HA	2.19	0.42
1:A:113:TRP:CD1	1:A:113:TRP:C	2.96	0.42
1:B:27:ARG:O	1:B:132:SER:HA	2.19	0.42
1:B:68:ASP:O	1:B:83:LEU:HA	2.20	0.42
1:C:57:CYS:HB2	1:C:105:PHE:CE1	2.54	0.42
1:D:194:LEU:HD12	1:D:194:LEU:HA	1.68	0.42
1:E:5:LYS:HG2	1:E:6:PHE:H	1.84	0.42
1:B:113:TRP:CD1	1:B:113:TRP:C	2.97	0.42
1:C:40:TYR:OH	1:C:126:HIS:O	2.37	0.42
1:E:62:VAL:HA	1:E:69:ILE:HD12	2.00	0.42
1:E:164:PHE:O	1:E:168:ILE:HG12	2.19	0.42
1:A:58:ALA:O	1:A:62:VAL:HG13	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:TRP:HB2	1:B:6:PHE:HB3	2.01	0.42
1:B:144:LYS:NZ	1:B:230:ILE:HD13	2.31	0.42
1:C:217:ARG:O	1:C:218:LYS:HG2	2.19	0.42
1:D:33:PHE:HA	1:D:36:PHE:HB3	2.01	0.42
1:E:153:ASP:O	1:E:157:ASN:HB2	2.20	0.42
1:D:25:VAL:HG22	1:D:82:ILE:CD1	2.50	0.42
1:E:28:ILE:HG22	1:E:130:PHE:HB3	2.02	0.42
1:F:26:VAL:CG2	1:F:97:LEU:HD21	2.49	0.42
1:F:91:ASN:C	1:F:93:ARG:H	2.27	0.42
1:A:151:TYR:HE1	1:A:211:LYS:HA	1.85	0.42
1:C:123:ASN:CG	1:C:124:ILE:H	2.25	0.42
1:A:169:ILE:HG22	1:A:170:LYS:HG2	2.01	0.42
1:B:54:MET:HB3	1:B:54:MET:HE3	1.73	0.42
1:A:44:LYS:HA	1:A:44:LYS:HE2	1.91	0.41
1:E:18:ILE:HG22	1:E:92:ARG:NH1	2.33	0.41
1:E:167:LEU:O	1:E:173:LEU:HB2	2.19	0.41
1:A:122:LEU:HA	1:A:122:LEU:HD22	1.80	0.41
1:C:100:LEU:HD23	1:C:100:LEU:HA	1.90	0.41
1:F:91:ASN:O	1:F:91:ASN:ND2	2.38	0.41
1:B:159:LEU:HD12	1:B:159:LEU:HA	1.80	0.41
1:B:23:TYR:CB	1:B:143:ILE:HG13	2.51	0.41
1:B:115:LYS:HE3	1:B:115:LYS:HB2	1.93	0.41
1:E:162:THR:O	1:E:166:GLN:HG2	2.20	0.41
1:A:43:ALA:H	1:A:47:ASP:HB2	1.85	0.41
1:C:19:LEU:HD12	1:C:20:PRO:HD2	2.03	0.41
1:C:62:VAL:HA	1:C:69:ILE:HD12	2.02	0.41
1:D:69:ILE:HG12	1:D:83:LEU:CD2	2.51	0.41
1:D:127:LEU:HG	1:D:128:PRO:HD2	2.02	0.41
1:D:208:MET:HA	1:D:222:LEU:HD13	2.01	0.41
1:E:74:GLY:HA3	1:E:79:TYR:CE2	2.56	0.41
1:F:138:PRO:HB2	1:F:139:ASN:H	1.78	0.41
1:C:151:TYR:CE2	1:C:224:ILE:HG13	2.55	0.41
1:C:208:MET:HE2	1:C:208:MET:HB3	1.89	0.41
1:D:82:ILE:HG23	1:D:143:ILE:HD11	2.02	0.41
1:E:5:LYS:HG2	1:E:6:PHE:N	2.35	0.41
1:E:44:LYS:O	1:E:45:PRO:C	2.58	0.41
1:B:124:ILE:HD12	1:B:124:ILE:HG23	1.87	0.40
1:A:19:LEU:HD12	1:A:20:PRO:HD2	2.03	0.40
1:A:103:SER:HB3	1:B:130:PHE:H	1.87	0.40
1:B:35:GLU:CD	1:B:35:GLU:N	2.79	0.40
1:D:230:ILE:HG23	1:D:231:ASP:OD1	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:145:ASP:O	1:F:149:TRP:CB	2.70	0.40
1:C:19:LEU:HA	1:C:20:PRO:HD3	1.81	0.40
1:C:177:GLU:O	1:C:181:LYS:CB	2.69	0.40
1:B:130:PHE:N	1:B:130:PHE:CD1	2.88	0.40
1:C:43:ALA:CB	1:C:47:ASP:HA	2.42	0.40
1:A:224:ILE:CG1	1:A:225:ASN:H	2.32	0.40
1:B:195:PHE:HE2	1:F:10:ARG:NH1	2.19	0.40
1:B:208:MET:CB	1:B:222:LEU:HD13	2.51	0.40
1:E:224:ILE:HG23	1:E:225:ASN:N	2.37	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:184:GLY:O	1:E:10:ARG:NH2[6_445]	2.09	0.11

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	227/243 (93%)	211 (93%)	13 (6%)	3 (1%)	9	38
1	B	227/243 (93%)	209 (92%)	15 (7%)	3 (1%)	9	38
1	C	227/243 (93%)	212 (93%)	12 (5%)	3 (1%)	9	38
1	D	227/243 (93%)	213 (94%)	12 (5%)	2 (1%)	14	48
1	E	227/243 (93%)	206 (91%)	19 (8%)	2 (1%)	14	48
1	F	64/243 (26%)	51 (80%)	11 (17%)	2 (3%)	3	19
All	All	1199/1458 (82%)	1102 (92%)	82 (7%)	15 (1%)	9	38

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	44	LYS
1	C	44	LYS
1	D	44	LYS
1	E	44	LYS
1	C	224	ILE
1	A	224	ILE
1	B	44	LYS
1	B	224	ILE
1	D	224	ILE
1	E	224	ILE
1	F	89	LEU
1	F	138	PRO
1	C	43	ALA
1	A	45	PRO
1	B	152	VAL

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	210/222 (95%)	206 (98%)	4 (2%)	50	76
1	B	210/222 (95%)	207 (99%)	3 (1%)	59	80
1	C	210/222 (95%)	208 (99%)	2 (1%)	68	84
1	D	210/222 (95%)	208 (99%)	2 (1%)	68	84
1	E	210/222 (95%)	207 (99%)	3 (1%)	59	80
1	F	78/222 (35%)	76 (97%)	2 (3%)	40	72
All	All	1128/1332 (85%)	1112 (99%)	16 (1%)	59	80

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	PHE
1	A	44	LYS
1	A	63	LEU
1	A	127	LEU

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Mol	Chain	Res	Type
1	B	63	LEU
1	B	79	TYR
1	B	127	LEU
1	C	63	LEU
1	C	127	LEU
1	D	63	LEU
1	D	127	LEU
1	E	31	LYS
1	E	44	LYS
1	E	63	LEU
1	F	63	LEU
1	F	140	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	55	ASN
1	A	141	GLN
1	A	225	ASN
1	B	55	ASN
1	C	225	ASN
1	D	46	ASN
1	D	139	ASN
1	D	158	ASN
1	E	46	ASN
1	E	60	ASN
1	E	126	HIS
1	F	139	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 10 are monoatomic - leaving 10 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
2	GTP	B	302	3	33,34,34	1.05	3 (9%)	50,54,54	1.69	11 (22%)
4	PO4	A	304	-	4,4,4	0.89	0	6,6,6	0.73	0
2	GTP	D	302	3	33,34,34	0.90	1 (3%)	50,54,54	1.65	13 (26%)
2	GTP	A	301	3	33,34,34	1.37	4 (12%)	50,54,54	1.64	11 (22%)
4	PO4	B	301	-	4,4,4	0.66	0	6,6,6	0.62	0
4	PO4	C	304	-	4,4,4	0.99	0	6,6,6	0.86	0
2	GTP	E	301	3	33,34,34	1.35	4 (12%)	50,54,54	1.57	11 (22%)
4	PO4	E	304	-	4,4,4	0.88	0	6,6,6	0.64	0
2	GTP	C	301	3	33,34,34	1.19	4 (12%)	50,54,54	1.71	12 (24%)
4	PO4	D	301	-	4,4,4	0.94	0	6,6,6	1.05	1 (16%)

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
2	GTP	B	302	3	-	6/22/38/38	0/3/3/3
2	GTP	D	302	3	-	7/22/38/38	0/3/3/3
2	GTP	A	301	3	-	6/22/38/38	0/3/3/3
2	GTP	E	301	3	-	4/22/38/38	0/3/3/3
2	GTP	C	301	3	-	7/22/38/38	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	GTP	PB-O3B	4.43	1.64	1.59
2	E	301	GTP	PA-O3A	3.82	1.63	1.59
2	C	301	GTP	PB-O3B	3.79	1.63	1.59
2	E	301	GTP	PB-O3A	3.61	1.63	1.59
2	A	301	GTP	PB-O3A	3.59	1.63	1.59
2	E	301	GTP	PB-O3B	3.18	1.62	1.59
2	C	301	GTP	PB-O3A	2.77	1.62	1.59
2	A	301	GTP	PA-O3A	2.71	1.62	1.59
2	B	302	GTP	PA-O3A	2.55	1.62	1.59
2	E	301	GTP	C2-N3	2.42	1.39	1.33
2	A	301	GTP	C2-N3	2.42	1.39	1.33
2	C	301	GTP	PA-O3A	2.37	1.62	1.59
2	B	302	GTP	PB-O3B	2.18	1.61	1.59
2	C	301	GTP	C2-N3	2.12	1.38	1.33
2	D	302	GTP	PB-O3B	2.05	1.61	1.59
2	B	302	GTP	PB-O3A	2.02	1.61	1.59

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	301	GTP	C2-N3-C4	5.09	121.07	112.30
2	A	301	GTP	C2-N3-C4	5.00	120.91	112.30
2	E	301	GTP	C5-C4-N3	-4.89	120.60	128.39
2	B	302	GTP	C2-N3-C4	4.88	120.70	112.30
2	D	302	GTP	C2-N3-C4	4.84	120.64	112.30
2	A	301	GTP	C5-C4-N3	-4.74	120.84	128.39
2	D	302	GTP	C5-C4-N3	-4.65	120.98	128.39
2	C	301	GTP	C5-C4-N3	-4.48	121.26	128.39
2	E	301	GTP	C2-N3-C4	4.33	119.76	112.30
2	B	302	GTP	C5-C4-N3	-3.90	122.18	128.39
2	B	302	GTP	N2-C2-N1	3.66	124.50	116.76
2	E	301	GTP	N9-C4-N3	3.34	132.64	125.95
2	D	302	GTP	C2'-C1'-N9	-3.20	104.35	113.25
2	C	301	GTP	O3G-PG-O3B	3.10	115.03	104.64
2	B	302	GTP	O3G-PG-O3B	3.01	114.73	104.64
2	E	301	GTP	O6-C6-C5	-2.84	119.04	126.53
2	C	301	GTP	C5-C6-N1	2.83	120.45	113.25
2	D	302	GTP	C3'-C2'-C1'	2.79	106.74	101.46
2	C	301	GTP	O6-C6-C5	-2.76	119.24	126.53
2	D	302	GTP	N2-C2-N1	2.75	122.56	116.76
2	E	301	GTP	C2-N1-C6	-2.73	120.16	125.11
2	C	301	GTP	C2-N1-C6	-2.71	120.20	125.11
2	B	302	GTP	O6-C6-C5	-2.70	119.39	126.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	GTP	C3'-C2'-C1'	2.66	106.49	101.46
2	A	301	GTP	N9-C4-N3	2.65	131.26	125.95
2	B	302	GTP	N1-C2-N3	-2.59	118.58	123.32
2	C	301	GTP	C3'-C2'-C1'	2.57	106.33	101.46
2	C	301	GTP	N9-C4-N3	2.57	131.09	125.95
2	C	301	GTP	O4'-C1'-N9	2.56	114.16	108.36
2	B	302	GTP	C5-C6-N1	2.53	119.69	113.25
2	D	302	GTP	C2-N1-C6	-2.52	120.53	125.11
2	A	301	GTP	N2-C2-N1	2.50	122.04	116.76
2	B	302	GTP	C3'-C2'-C1'	2.49	106.17	101.46
2	A	301	GTP	N9-C8-N7	-2.48	108.80	113.40
2	A	301	GTP	C8-N7-C5	2.48	108.68	104.26
2	C	301	GTP	N2-C2-N1	2.43	121.89	116.76
2	D	302	GTP	O6-C6-C5	-2.42	120.15	126.53
2	C	301	GTP	N1-C2-N3	-2.41	118.90	123.32
2	A	301	GTP	C2-N1-C6	-2.40	120.75	125.11
2	A	301	GTP	C5-C6-N1	2.39	119.33	113.25
2	B	302	GTP	C2-N1-C6	-2.37	120.82	125.11
2	A	301	GTP	N1-C2-N3	-2.34	119.04	123.32
2	D	302	GTP	C5-C6-N1	2.34	119.21	113.25
2	D	302	GTP	N9-C4-N3	2.34	130.63	125.95
2	E	301	GTP	O3G-PG-O3B	2.29	112.32	104.64
2	C	301	GTP	N9-C8-N7	-2.28	109.18	113.40
2	B	302	GTP	N9-C8-N7	-2.26	109.20	113.40
2	E	301	GTP	C5-C6-N1	2.20	118.84	113.25
2	D	302	GTP	O4'-C1'-N9	2.18	113.31	108.36
2	A	301	GTP	C4-C5-N7	-2.16	107.24	110.67
2	E	301	GTP	N2-C2-N1	2.16	121.32	116.76
4	D	301	PO4	O4-P-O3	2.16	114.62	107.91
2	E	301	GTP	C3'-C2'-C1'	2.13	105.49	101.46
2	D	302	GTP	N1-C2-N3	-2.13	119.43	123.32
2	E	301	GTP	N9-C8-N7	-2.12	109.48	113.40
2	B	302	GTP	O2A-PA-O3A	2.10	112.96	107.27
2	E	301	GTP	C8-N7-C5	2.10	108.00	104.26
2	D	302	GTP	N9-C8-N7	-2.10	109.51	113.40
2	D	302	GTP	C8-N7-C5	2.02	107.86	104.26

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	GTP	C4'-C5'-O5'-PA

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Mol	Chain	Res	Type	Atoms
2	B	302	GTP	C4'-C5'-O5'-PA
2	D	302	GTP	C4'-C5'-O5'-PA
2	B	302	GTP	O4'-C4'-C5'-O5'
2	E	301	GTP	O4'-C4'-C5'-O5'
2	D	302	GTP	C3'-C4'-C5'-O5'
2	E	301	GTP	C3'-C4'-C5'-O5'
2	A	301	GTP	O4'-C4'-C5'-O5'
2	B	302	GTP	C3'-C4'-C5'-O5'
2	C	301	GTP	O4'-C4'-C5'-O5'
2	C	301	GTP	C3'-C4'-C5'-O5'
2	D	302	GTP	O4'-C4'-C5'-O5'
2	A	301	GTP	C3'-C4'-C5'-O5'
2	B	302	GTP	PA-O3A-PB-O1B
2	C	301	GTP	C4'-C5'-O5'-PA
2	E	301	GTP	C4'-C5'-O5'-PA
2	A	301	GTP	PA-O3A-PB-O3B
2	D	302	GTP	PA-O3A-PB-O3B
2	C	301	GTP	PA-O3A-PB-O3B
2	B	302	GTP	PG-O3B-PB-O2B
2	D	302	GTP	PG-O3B-PB-O1B
2	A	301	GTP	PA-O3A-PB-O1B
2	C	301	GTP	PG-O3B-PB-O1B
2	C	301	GTP	PA-O3A-PB-O1B
2	C	301	GTP	PA-O3A-PB-O2B
2	D	302	GTP	PA-O3A-PB-O1B
2	A	301	GTP	PG-O3B-PB-O2B
2	B	302	GTP	PG-O3B-PB-O1B
2	D	302	GTP	PG-O3B-PB-O2B
2	E	301	GTP	PB-O3A-PA-O2A

There are no ring outliers.

9 monomers are involved in 25 short contacts:

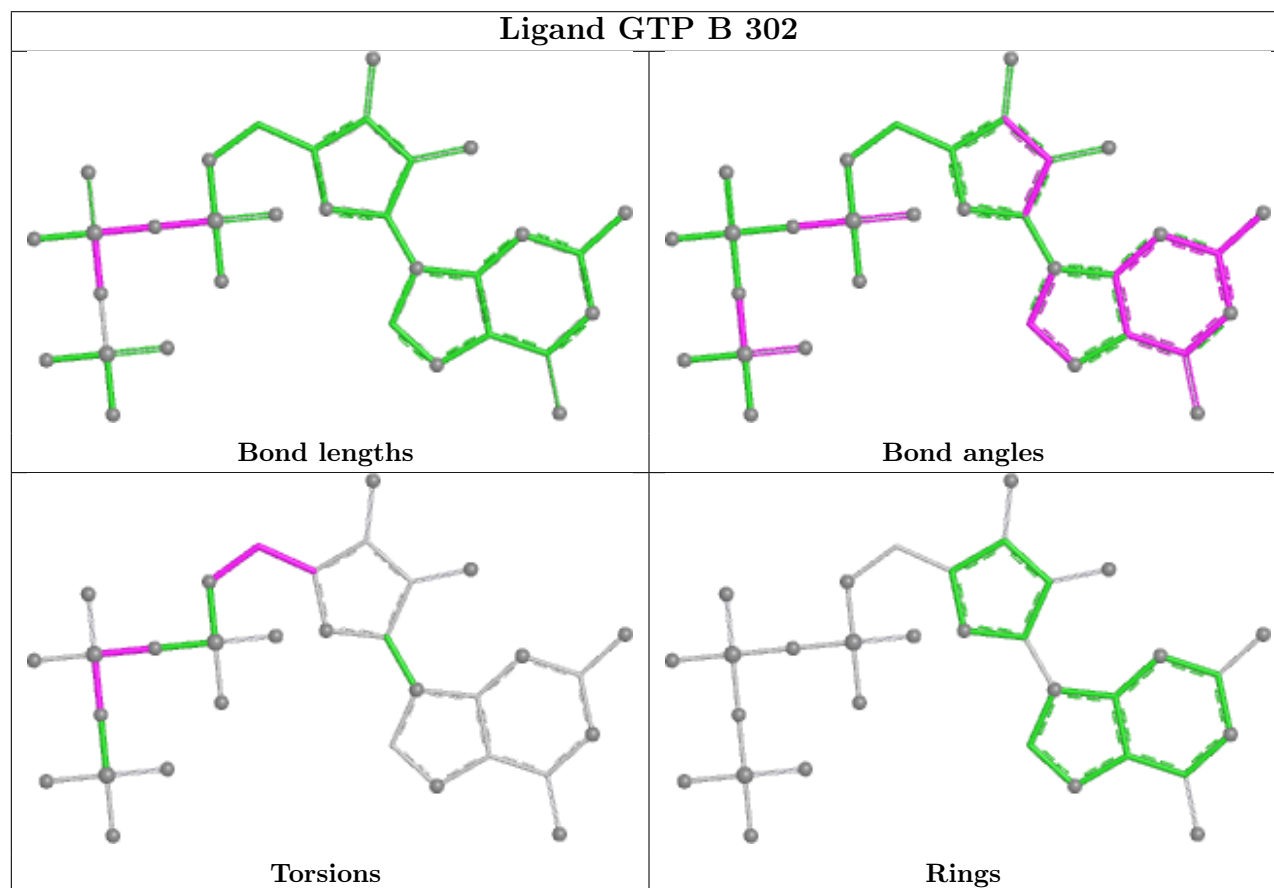
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	302	GTP	3	0
4	A	304	PO4	3	0
2	D	302	GTP	5	0
2	A	301	GTP	4	0
4	C	304	PO4	1	0
2	E	301	GTP	2	0
4	E	304	PO4	1	0
2	C	301	GTP	5	0

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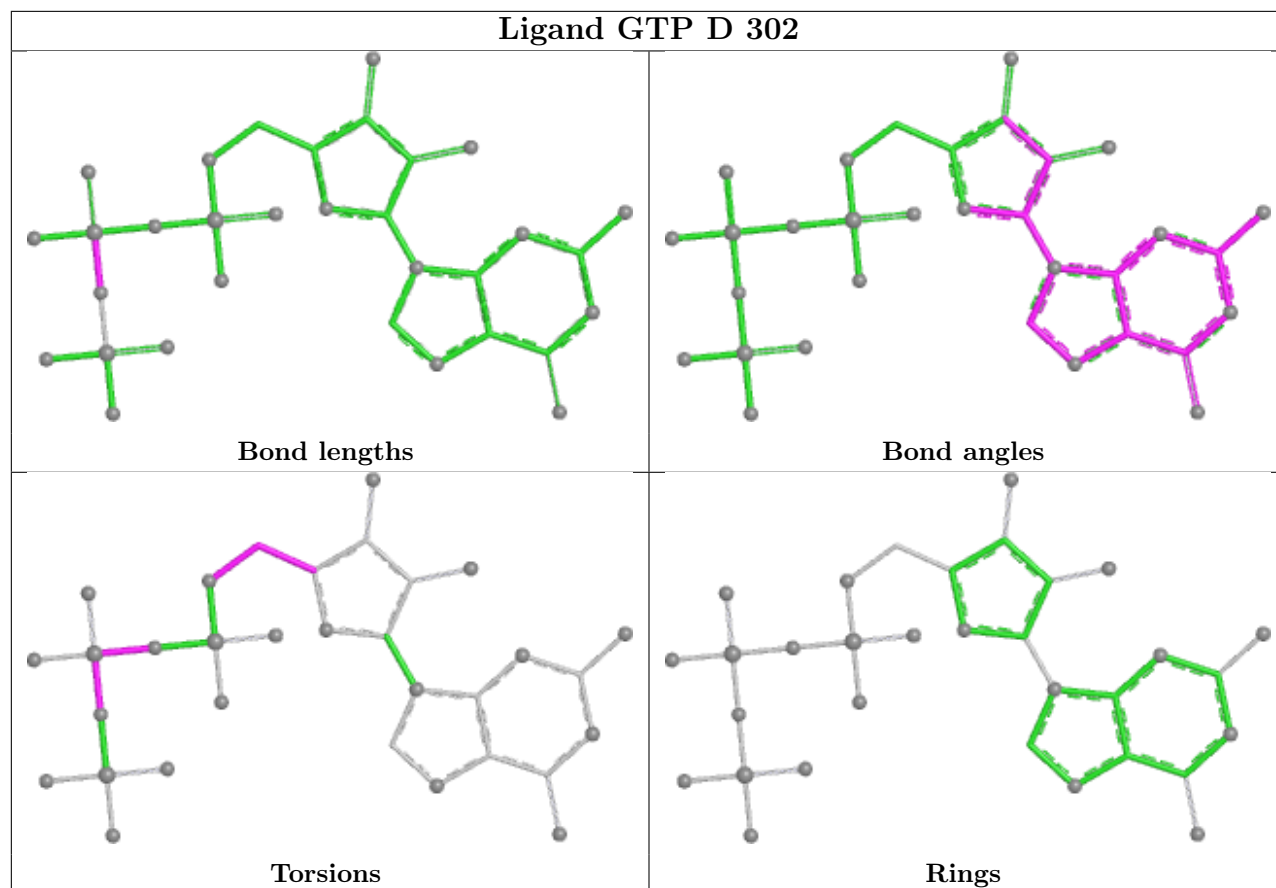
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	301	PO4	1	0

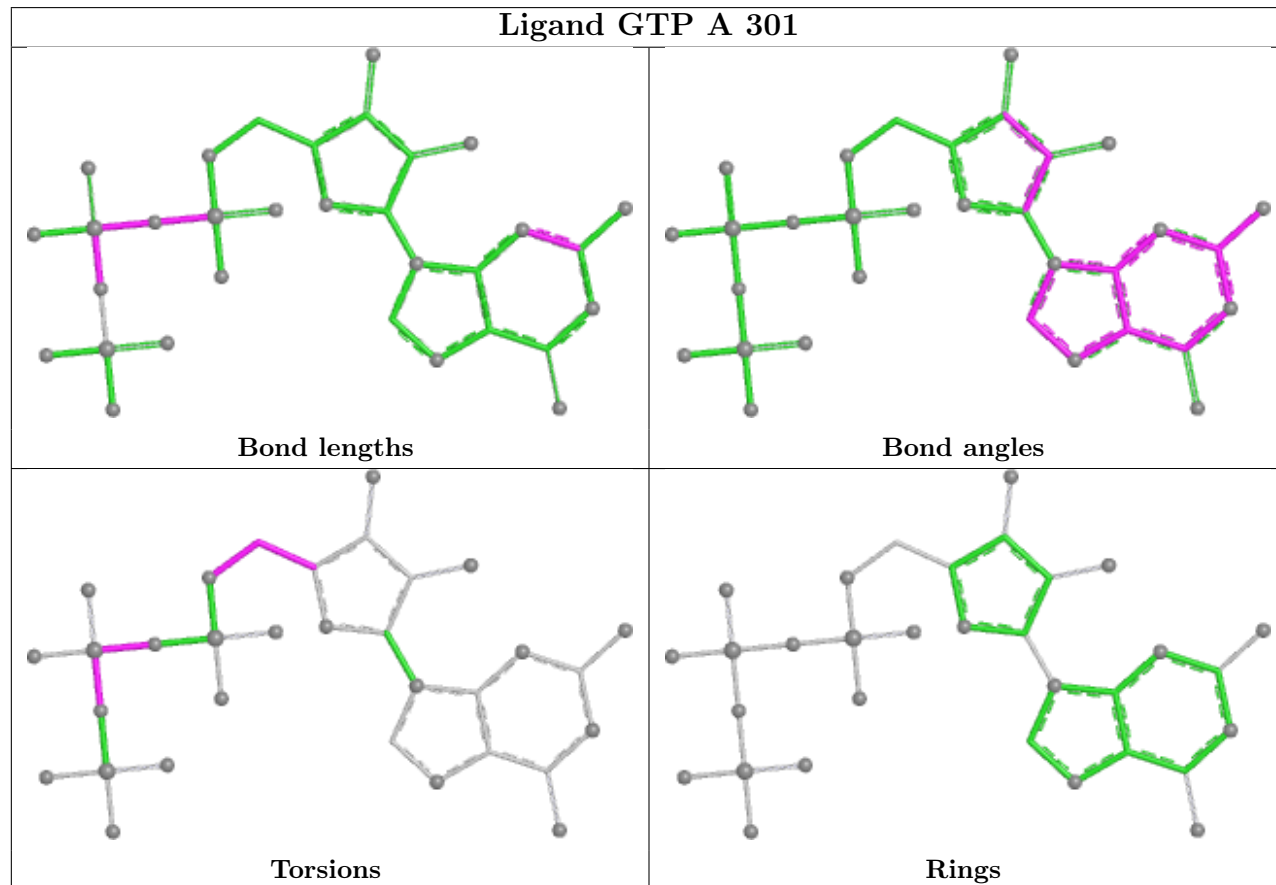
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



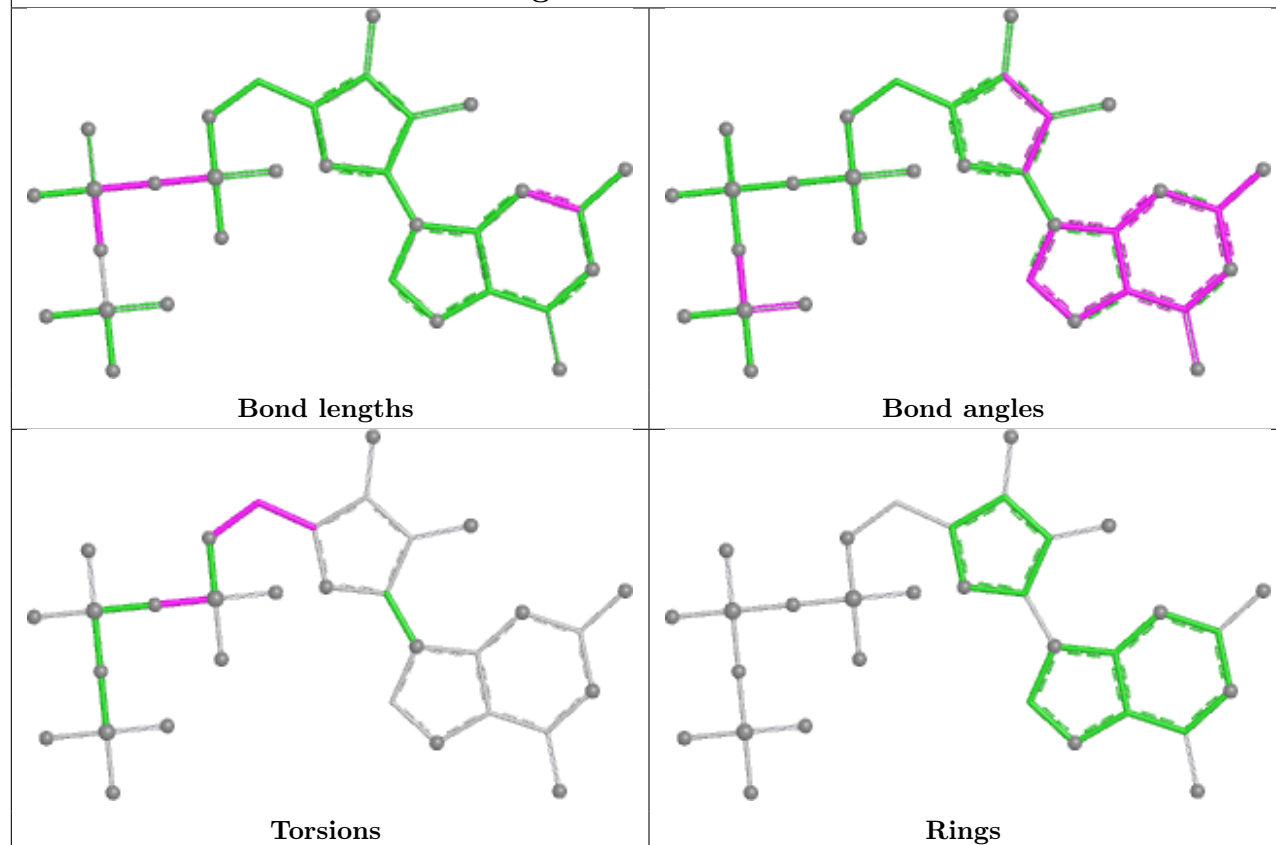
Ligand GTP D 302



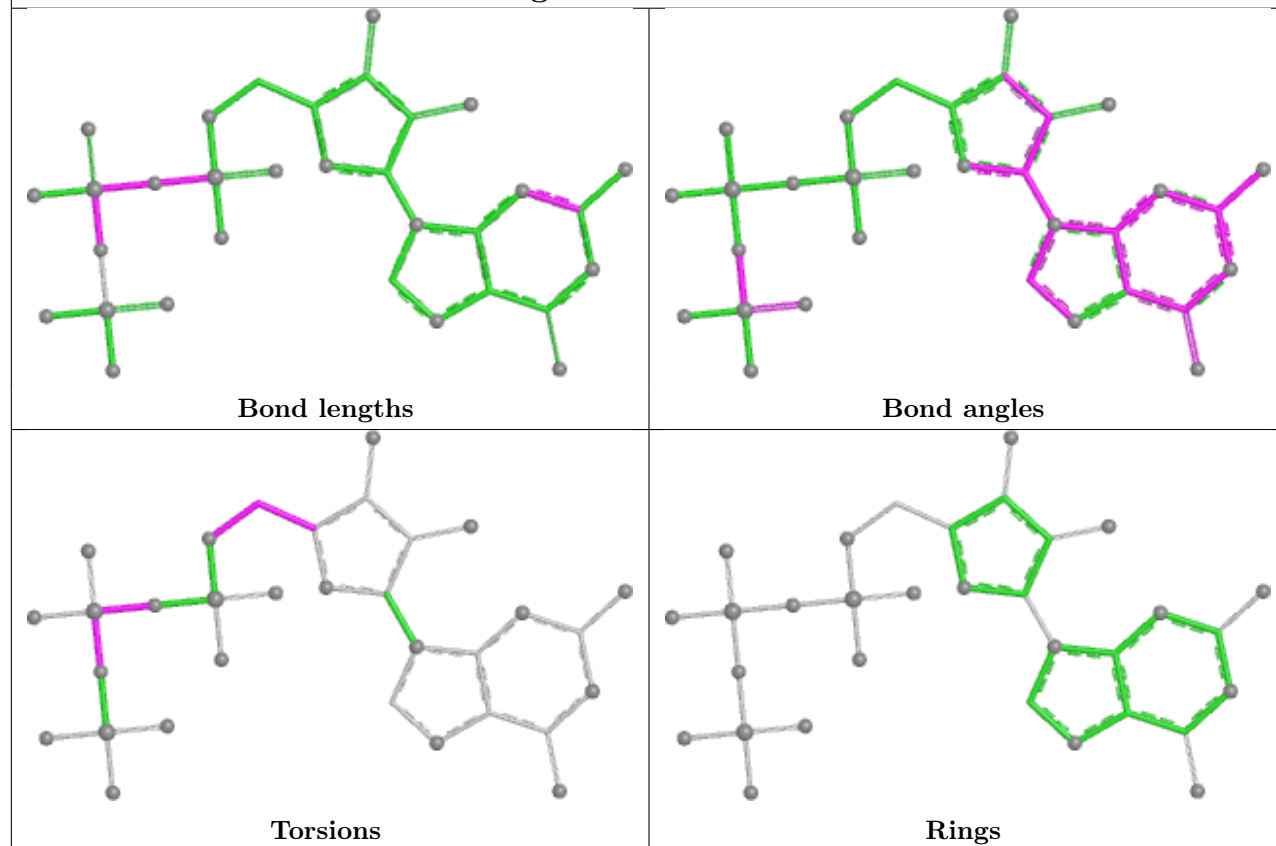
Ligand GTP A 301



Ligand GTP E 301



Ligand GTP C 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

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	229/243 (94%)	-0.32	3 (1%) 75 53	42, 77, 114, 166	0
1	B	229/243 (94%)	-0.25	3 (1%) 75 53	48, 96, 137, 172	0
1	C	229/243 (94%)	-0.19	8 (3%) 47 27	51, 87, 151, 184	0
1	D	229/243 (94%)	-0.12	9 (3%) 43 24	45, 88, 151, 191	0
1	E	229/243 (94%)	0.13	7 (3%) 51 30	87, 124, 162, 218	0
1	F	80/243 (32%)	0.42	6 (7%) 20 10	101, 137, 173, 177	0
All	All	1225/1458 (84%)	-0.11	36 (2%) 53 31	42, 100, 157, 218	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	5	LYS	4.5
1	C	233	LEU	4.4
1	D	35	GLU	4.3
1	E	233	LEU	4.0
1	A	233	LEU	4.0
1	E	230	ILE	3.8
1	C	223	HIS	3.7
1	E	31	LYS	3.6
1	A	230	ILE	3.5
1	F	9	VAL	3.2
1	D	204	ASN	3.2
1	B	5	LYS	3.2
1	D	231	ASP	3.0
1	D	197	GLU	2.8
1	E	165	TRP	2.7
1	D	198	CYS	2.7
1	C	227	ILE	2.7
1	C	155	HIS	2.7
1	D	230	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	177	GLU	2.5
1	C	211	LYS	2.5
1	B	197	GLU	2.5
1	E	124	ILE	2.5
1	D	44	LYS	2.4
1	F	100	LEU	2.3
1	F	140	LEU	2.3
1	B	233	LEU	2.3
1	C	45	PRO	2.2
1	F	33	PHE	2.2
1	C	175	PRO	2.2
1	F	151	TYR	2.2
1	C	230	ILE	2.2
1	E	154	THR	2.1
1	E	86	SER	2.1
1	F	149	TRP	2.1
1	A	41	GLU	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
2	GTP	E	301	32/32	0.86	0.10	113,137,159,162	0
4	PO4	E	304	5/5	0.87	0.07	99,111,118,122	0
2	GTP	C	301	32/32	0.92	0.09	104,114,128,142	0
2	GTP	B	302	32/32	0.93	0.08	97,124,143,147	0
3	MG	B	303	1/1	0.93	0.19	100,100,100,100	0
4	PO4	B	301	5/5	0.93	0.12	74,75,85,87	0

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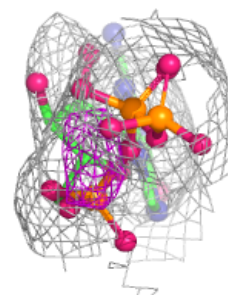
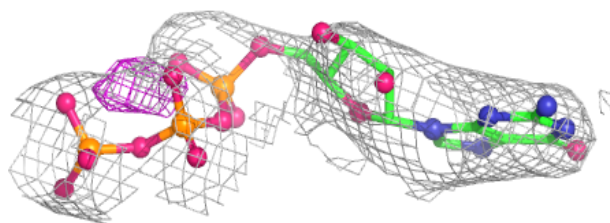
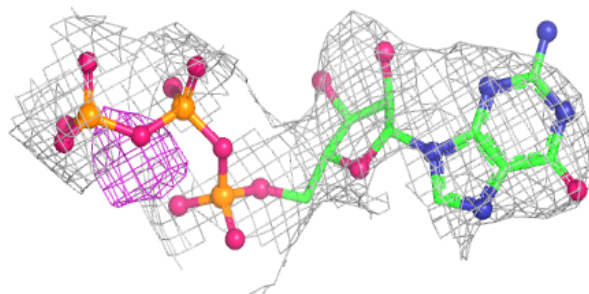
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GTP	A	301	32/32	0.93	0.07	65,94,107,112	0
4	PO4	C	304	5/5	0.95	0.07	69,73,81,86	0
4	PO4	A	304	5/5	0.95	0.06	57,93,102,104	0
3	MG	A	302	1/1	0.96	0.17	72,72,72,72	0
2	GTP	D	302	32/32	0.96	0.06	77,102,119,134	0
3	MG	C	302	1/1	0.96	0.08	95,95,95,95	0
4	PO4	D	301	5/5	0.96	0.06	66,67,68,93	0
3	MG	D	303	1/1	0.96	0.06	91,91,91,91	0
3	MG	E	302	1/1	0.97	0.06	112,112,112,112	0
3	MG	C	303	1/1	0.97	0.09	108,108,108,108	0
3	MG	E	303	1/1	0.98	0.04	106,106,106,106	0
3	MG	D	304	1/1	0.99	0.02	99,99,99,99	0
3	MG	B	304	1/1	0.99	0.05	104,104,104,104	0
3	MG	A	303	1/1	1.00	0.03	73,73,73,73	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.

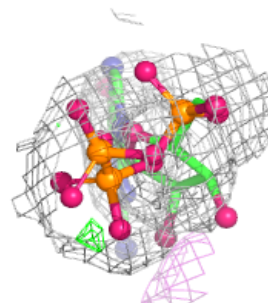
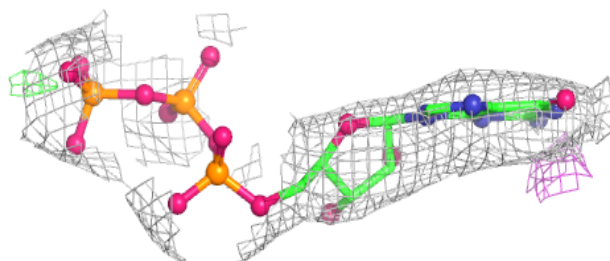
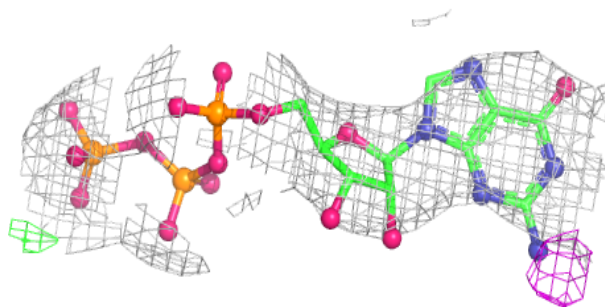
Electron density around GTP E 301:

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

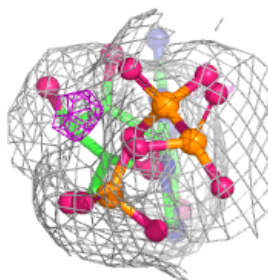
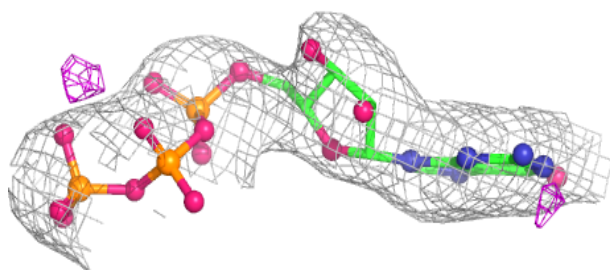
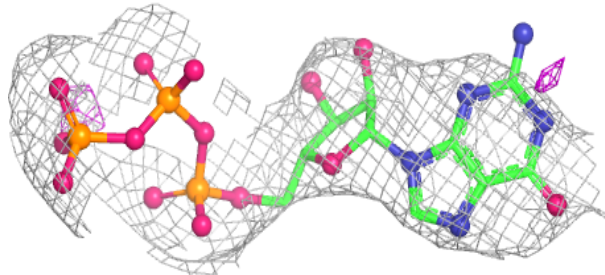


Electron density around GTP C 301:

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and green (positive)

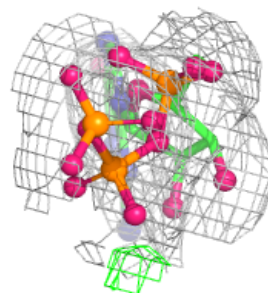
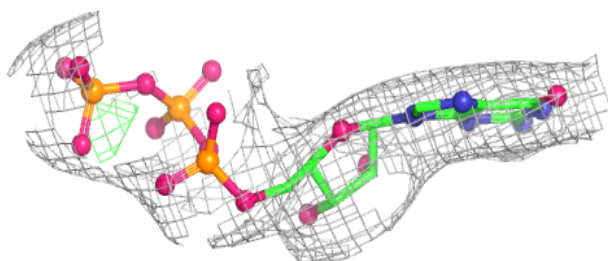
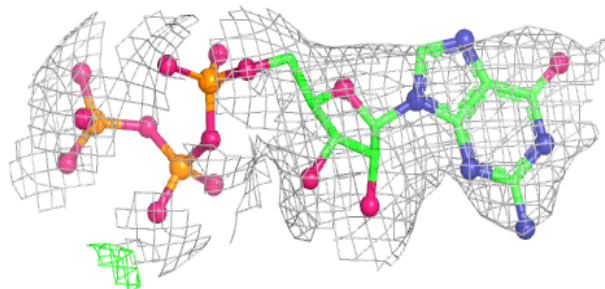
**Electron density around GTP B 302:**

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

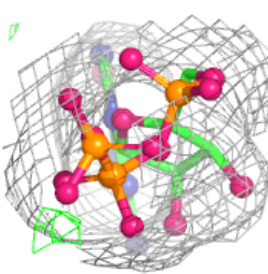
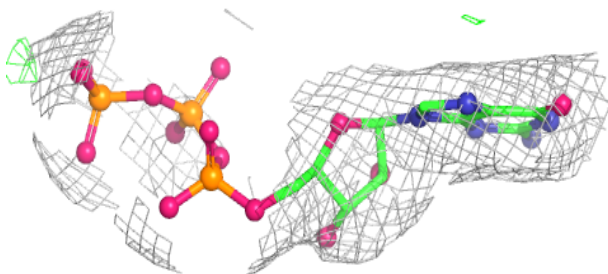
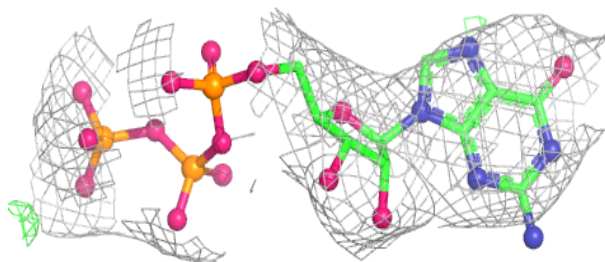


Electron density around GTP A 301:

$2mF_o-DF_c$ (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.