



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:58 PM UTC

PDB ID : 4Q7H / pdb_00004q7h
Title : Crystal structure of SAMHD1 catalytic core with GTP
Authors : Xiaohong, Q.; Yanhong, L.; Jia, K.; Xiaofang, Y.
Deposited on : 2014-04-25
Resolution : 2.59 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

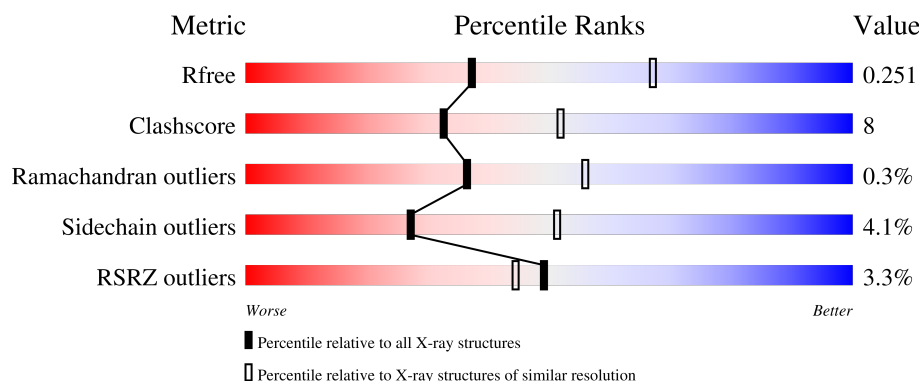
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	4770 (2.60-2.56)
Clashscore	190562	5124 (2.60-2.56)
Ramachandran outliers	187476	5046 (2.60-2.56)
Sidechain outliers	187428	5046 (2.60-2.56)
RSRZ outliers	180081	4770 (2.60-2.56)

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	526	2% 71% 14% • 13%
1	B	526	5% 70% 15% • 15%
1	C	526	2% 67% 16% • 15%
1	D	526	3% 65% 18% • 17%

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
3	GTP	B	902	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14424 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 Deoxynucleoside triphosphate triphosphohydrolase SAMHD1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	459	Total	C	N	O	S	0	0	0
			3646	2336	630	662	18			
1	B	449	Total	C	N	O	S	0	0	0
			3557	2277	616	646	18			
1	C	447	Total	C	N	O	S	0	0	0
			3572	2291	621	642	18			
1	D	439	Total	C	N	O	S	0	0	0
			3493	2237	605	634	17			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	101	HIS	-	expression tag	UNP Q9Y3Z3
A	102	HIS	-	expression tag	UNP Q9Y3Z3
A	103	HIS	-	expression tag	UNP Q9Y3Z3
A	104	HIS	-	expression tag	UNP Q9Y3Z3
A	105	HIS	-	expression tag	UNP Q9Y3Z3
A	106	HIS	-	expression tag	UNP Q9Y3Z3
A	107	GLY	-	expression tag	UNP Q9Y3Z3
A	108	SER	-	expression tag	UNP Q9Y3Z3
A	266	TYR	CYS	engineered mutation	UNP Q9Y3Z3
B	101	HIS	-	expression tag	UNP Q9Y3Z3
B	102	HIS	-	expression tag	UNP Q9Y3Z3
B	103	HIS	-	expression tag	UNP Q9Y3Z3
B	104	HIS	-	expression tag	UNP Q9Y3Z3
B	105	HIS	-	expression tag	UNP Q9Y3Z3
B	106	HIS	-	expression tag	UNP Q9Y3Z3
B	107	GLY	-	expression tag	UNP Q9Y3Z3
B	108	SER	-	expression tag	UNP Q9Y3Z3
B	266	TYR	CYS	engineered mutation	UNP Q9Y3Z3
C	101	HIS	-	expression tag	UNP Q9Y3Z3
C	102	HIS	-	expression tag	UNP Q9Y3Z3
C	103	HIS	-	expression tag	UNP Q9Y3Z3

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Chain	Residue	Modelled	Actual	Comment	Reference
C	104	HIS	-	expression tag	UNP Q9Y3Z3
C	105	HIS	-	expression tag	UNP Q9Y3Z3
C	106	HIS	-	expression tag	UNP Q9Y3Z3
C	107	GLY	-	expression tag	UNP Q9Y3Z3
C	108	SER	-	expression tag	UNP Q9Y3Z3
C	266	TYR	CYS	engineered mutation	UNP Q9Y3Z3
D	101	HIS	-	expression tag	UNP Q9Y3Z3
D	102	HIS	-	expression tag	UNP Q9Y3Z3
D	103	HIS	-	expression tag	UNP Q9Y3Z3
D	104	HIS	-	expression tag	UNP Q9Y3Z3
D	105	HIS	-	expression tag	UNP Q9Y3Z3
D	106	HIS	-	expression tag	UNP Q9Y3Z3
D	107	GLY	-	expression tag	UNP Q9Y3Z3
D	108	SER	-	expression tag	UNP Q9Y3Z3
D	266	TYR	CYS	engineered mutation	UNP Q9Y3Z3

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0

- Molecule 3 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃).



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	C	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

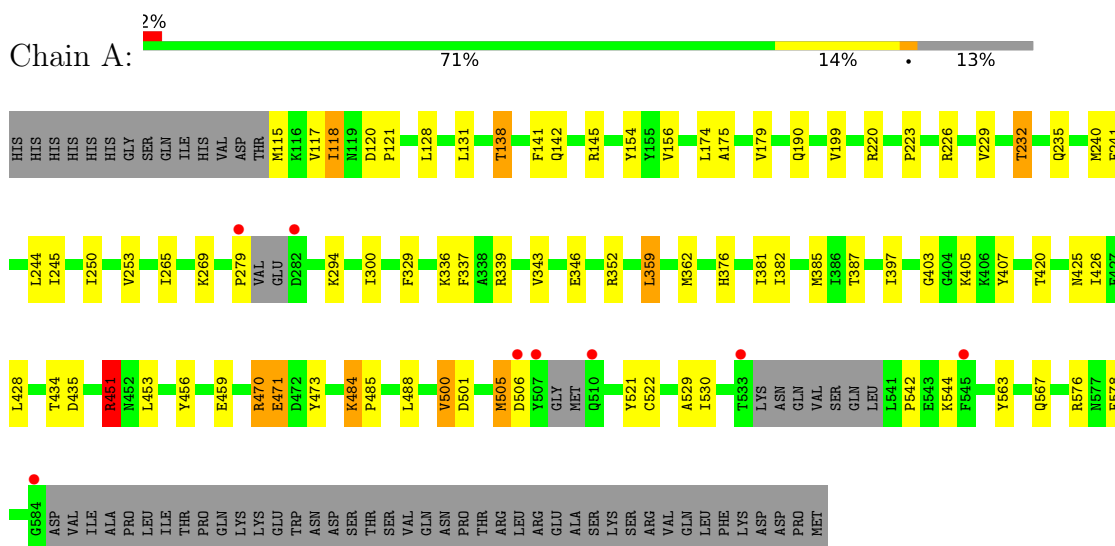
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	18	Total	O	0	0
			18	18		
4	B	18	Total	O	0	0
			18	18		
4	C	7	Total	O	0	0
			7	7		
4	D	13	Total	O	0	0
			13	13		

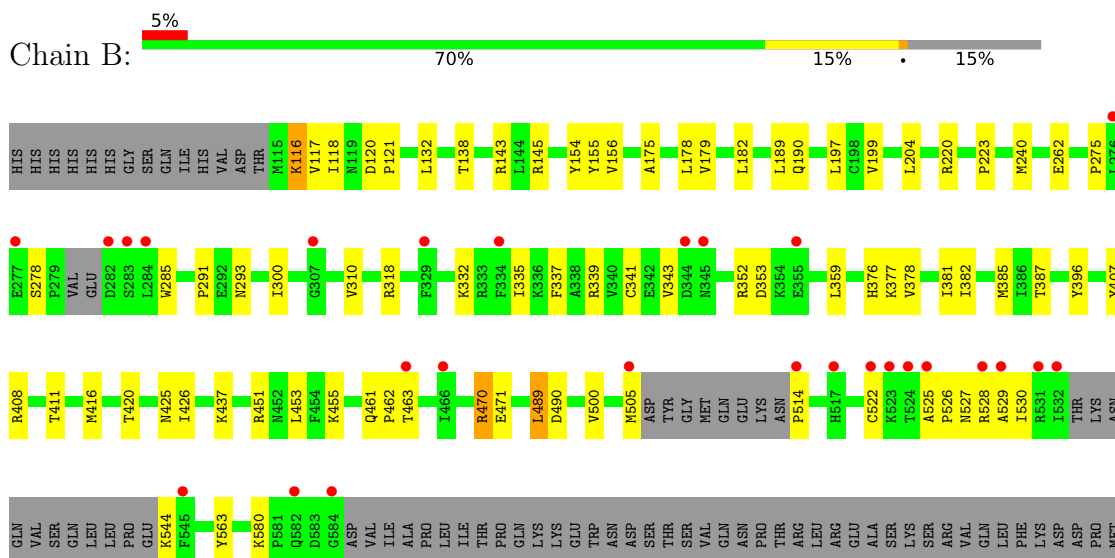
3 Residue-property plots [i](#)

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

• Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

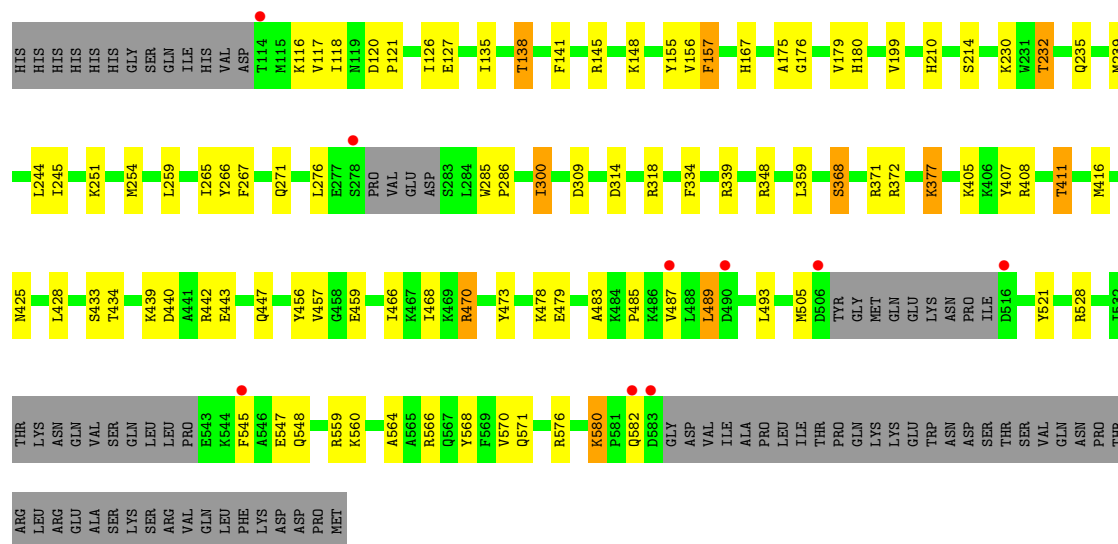


• Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

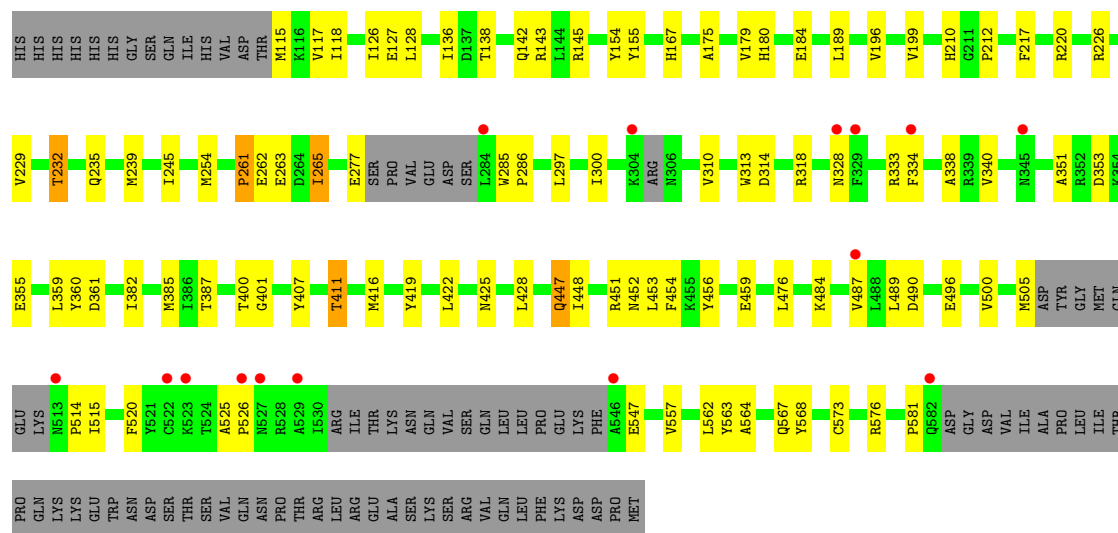


• Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1





● Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.94Å 142.76Å 199.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.57 – 2.59 34.57 – 2.59	Depositor EDS
% Data completeness (in resolution range)	94.5 (34.57-2.59) 88.5 (34.57-2.59)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 2.57Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.191 , 0.248 0.195 , 0.251	Depositor DCC
R_{free} test set	2000 reflections (2.74%)	wwPDB-VP
Wilson B-factor (Å ²)	43.9	Xtriage
Anisotropy	0.407	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 34.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14424	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.79% 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, ZN

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	0.58	0/3732	0.87	9/5052 (0.2%)
1	B	0.57	0/3640	0.90	8/4927 (0.2%)
1	C	0.54	0/3656	0.88	7/4946 (0.1%)
1	D	0.55	0/3574	0.88	6/4833 (0.1%)
All	All	0.56	0/14602	0.88	30/19758 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	526	PRO	N-CA-CB	10.12	110.11	102.73
1	D	526	PRO	N-CA-CB	7.61	110.53	103.46
1	A	279	PRO	N-CA-CB	6.66	110.33	103.00
1	B	514	PRO	N-CA-CB	6.62	110.29	103.00
1	D	261	PRO	N-CA-CB	6.50	110.08	103.25
1	A	435	ASP	CA-C-N	5.97	125.85	119.28
1	A	435	ASP	C-N-CA	5.97	125.85	119.28
1	D	514	PRO	N-CA-CB	5.94	109.21	103.51
1	B	489	LEU	N-CA-C	5.71	117.64	109.14
1	C	251	LYS	CA-C-N	-5.65	113.20	119.19
1	C	251	LYS	C-N-CA	-5.65	113.20	119.19
1	C	214	SER	CB-CA-C	-5.61	110.09	116.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	263	GLU	N-CA-C	-5.61	105.17	112.23
1	C	580	LYS	CA-C-N	5.56	126.79	119.84
1	C	580	LYS	C-N-CA	5.56	126.79	119.84
1	A	542	PRO	N-CA-CB	5.49	110.09	103.45
1	D	487	VAL	N-CA-C	5.39	111.98	106.21
1	B	500	VAL	CA-C-N	-5.36	115.15	122.93
1	B	500	VAL	C-N-CA	-5.36	115.15	122.93
1	A	451	ARG	N-CA-C	5.30	118.51	111.30
1	A	226	ARG	CA-C-N	5.23	124.90	119.56
1	A	226	ARG	C-N-CA	5.23	124.90	119.56
1	D	515	ILE	N-CA-C	5.21	115.98	110.72
1	B	285	TRP	CA-C-N	5.09	124.78	119.64
1	B	285	TRP	C-N-CA	5.09	124.78	119.64
1	A	484	LYS	CA-C-N	5.08	125.02	119.78
1	A	484	LYS	C-N-CA	5.08	125.02	119.78
1	B	116	LYS	N-CA-C	5.02	117.75	109.76
1	C	157	PHE	CA-C-N	5.01	125.28	119.47
1	C	157	PHE	C-N-CA	5.01	125.28	119.47

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	505	MET	Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3646	0	3491	55	0
1	B	3557	0	3406	52	0
1	C	3572	0	3450	68	0
1	D	3493	0	3330	57	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	32	0	12	2	0
3	B	32	0	11	9	0
3	C	32	0	12	8	0
4	A	18	0	0	0	0
4	B	18	0	0	0	0
4	C	7	0	0	0	0
4	D	13	0	0	0	0
All	All	14424	0	13712	223	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (223) 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:D:385:MET:HE1	1:D:453:LEU:HD22	1.46	0.97
1:C:232:THR:HG22	1:C:235:GLN:H	1.33	0.92
1:A:232:THR:HG22	1:A:235:GLN:H	1.37	0.89
1:D:447:GLN:HE21	1:D:452:ASN:HD22	1.19	0.88
1:B:381:ILE:HG22	1:B:385:MET:HE2	1.59	0.83
3:C:902:GTP:O1B	3:C:902:GTP:N2	2.12	0.83
1:D:489:LEU:HD21	1:D:567:GLN:HG3	1.61	0.83
1:C:116:LYS:NZ	3:C:902:GTP:H2'	1.95	0.81
1:C:505:MET:HE3	1:C:545:PHE:HB2	1.60	0.81
1:A:381:ILE:HG22	1:A:385:MET:HE2	1.65	0.77
1:D:117:VAL:HG22	1:D:127:GLU:HG2	1.66	0.76
1:B:343:VAL:HG22	1:B:529:ALA:HB2	1.68	0.75
1:B:408:ARG:H	1:B:411:THR:HG22	1.52	0.73
1:D:254:MET:HE1	1:D:265:ILE:HG13	1.71	0.72
1:A:145:ARG:NH2	1:B:154:TYR:O	2.24	0.71
1:C:425:ASN:OD1	1:D:425:ASN:ND2	2.24	0.70
1:C:479:GLU:OE1	1:C:576:ARG:NH1	2.25	0.70
1:D:232:THR:HG22	1:D:235:GLN:H	1.57	0.68
1:A:138:THR:HG22	1:A:141:PHE:H	1.58	0.68
1:B:382:ILE:HA	1:B:385:MET:HE3	1.75	0.68
1:C:334:PHE:HE2	1:C:359:LEU:HD11	1.59	0.68
1:D:447:GLN:NE2	1:D:452:ASN:HD22	1.90	0.68
1:D:407:TYR:HB3	1:D:411:THR:HG22	1.74	0.68
1:A:138:THR:HG21	1:A:244:LEU:HG	1.78	0.66
1:A:175:ALA:HB1	1:A:199:VAL:HG12	1.77	0.65
1:A:397:ILE:HG21	1:A:426:ILE:HD11	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:ARG:HG2	1:A:521:TYR:CZ	2.32	0.65
1:B:275:PRO:HB2	1:B:278:SER:HB3	1.79	0.65
1:A:385:MET:HE1	1:A:453:LEU:HD22	1.79	0.64
1:C:116:LYS:HZ1	3:C:902:GTP:H2'	1.59	0.64
1:A:451:ARG:NE	3:B:902:GTP:O1A	2.32	0.63
1:C:138:THR:HG22	1:C:141:PHE:H	1.61	0.62
1:A:382:ILE:HA	1:A:385:MET:HE3	1.82	0.61
1:B:522:CYS:HB2	1:B:530:ILE:HD13	1.81	0.61
1:C:145:ARG:NH2	1:D:154:TYR:O	2.34	0.61
1:D:179:VAL:HG12	1:D:196:VAL:HG22	1.82	0.61
1:D:476:LEU:HB2	1:D:500:VAL:HG11	1.83	0.60
1:B:385:MET:HE1	1:B:453:LEU:HD22	1.84	0.60
1:A:505:MET:HB2	1:A:506:ASP:HA	1.84	0.59
1:B:525:ALA:HB1	1:B:527:ASN:HD21	1.67	0.59
1:D:143:ARG:NH2	1:D:210:HIS:O	2.36	0.59
1:C:145:ARG:NH2	1:D:155:TYR:HA	2.18	0.58
1:C:407:TYR:HB3	1:C:411:THR:HG22	1.85	0.58
1:D:334:PHE:O	1:D:338:ALA:HB2	2.04	0.58
1:D:239:MET:HE3	1:D:416:MET:SD	2.44	0.58
1:C:334:PHE:CE2	1:C:359:LEU:HD11	2.39	0.57
1:C:505:MET:HG2	1:C:545:PHE:CD2	2.39	0.57
1:A:505:MET:N	1:A:506:ASP:OD1	2.33	0.57
1:B:120:ASP:OD1	1:B:318:ARG:NH2	2.38	0.56
1:A:425:ASN:OD1	1:B:425:ASN:ND2	2.38	0.56
1:A:154:TYR:O	1:B:145:ARG:NH2	2.38	0.56
1:D:328:ASN:HB2	1:D:361:ASP:HB3	1.86	0.56
1:C:138:THR:HG21	1:C:244:LEU:HG	1.86	0.56
1:A:339:ARG:HD3	1:A:521:TYR:CZ	2.41	0.56
1:C:116:LYS:HE2	3:C:902:GTP:H3'	1.86	0.56
1:C:116:LYS:HZ3	3:C:902:GTP:H2'	1.71	0.56
1:C:433:SER:O	1:C:442:ARG:NH1	2.39	0.55
1:C:117:VAL:HG22	1:C:127:GLU:HG2	1.88	0.55
1:C:468:ILE:H	1:C:548:GLN:HE22	1.52	0.55
1:C:266:TYR:CE2	1:C:276:LEU:HD12	2.42	0.55
1:B:143:ARG:HD2	1:B:420:THR:HA	1.89	0.54
1:C:155:TYR:HA	1:D:145:ARG:NH2	2.21	0.54
1:C:483:ALA:HB1	1:C:571:GLN:HG3	1.89	0.54
1:A:156:VAL:HG11	1:A:376:HIS:CE1	2.43	0.54
1:C:570:VAL:HG13	1:C:580:LYS:HE3	1.88	0.54
1:C:179:VAL:HG22	1:C:300:ILE:HD12	1.90	0.54
1:A:522:CYS:HB2	1:A:530:ILE:HD11	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:484:LYS:HE3	1:D:484:LYS:HA	1.90	0.53
1:B:132:LEU:HD22	1:B:204:LEU:HD22	1.90	0.53
1:B:544:LYS:N	1:D:360:TYR:HH	2.06	0.53
1:C:156:VAL:HG23	1:C:157:PHE:CD2	2.43	0.53
1:D:189:LEU:HD21	1:D:340:VAL:HB	1.90	0.53
1:A:336:LYS:HD2	1:A:337:PHE:CE1	2.44	0.52
1:D:142:GLN:OE1	1:D:145:ARG:HD2	2.09	0.52
1:B:117:VAL:O	3:B:902:GTP:O2'	2.28	0.52
1:D:220:ARG:HG2	1:D:387:THR:HG21	1.92	0.52
1:D:333:ARG:NH2	1:D:355:GLU:OE1	2.39	0.52
1:C:120:ASP:OD1	1:C:121:PRO:HD2	2.10	0.52
1:D:334:PHE:CE2	1:D:359:LEU:HD21	2.45	0.51
1:C:489:LEU:HD12	1:C:564:ALA:HB2	1.91	0.51
1:A:576:ARG:HG2	1:A:578:PHE:CE2	2.45	0.51
1:B:408:ARG:H	1:B:411:THR:CG2	2.23	0.51
1:C:368:SER:O	1:C:372:ARG:HB2	2.10	0.51
3:A:902:GTP:H5''	1:B:451:ARG:HE	1.75	0.51
1:B:116:LYS:NZ	3:B:902:GTP:H3'	2.26	0.50
1:C:439:LYS:NZ	1:C:440:ASP:OD1	2.38	0.50
1:B:116:LYS:HE3	3:B:902:GTP:H2'	1.92	0.50
1:D:573:CYS:SG	1:D:581:PRO:HG3	2.51	0.50
1:C:339:ARG:HD3	1:C:521:TYR:CZ	2.46	0.50
1:C:457:VAL:O	1:C:566:ARG:NH1	2.44	0.50
1:D:314:ASP:OD1	1:D:318:ARG:NH1	2.45	0.50
1:A:451:ARG:NH2	3:B:902:GTP:C6	2.80	0.50
1:B:156:VAL:HG11	1:B:376:HIS:CE1	2.47	0.50
1:C:138:THR:HG22	1:C:141:PHE:N	2.25	0.49
1:D:220:ARG:NH2	1:D:500:VAL:O	2.45	0.49
1:C:408:ARG:H	1:C:411:THR:HB	1.77	0.49
1:D:155:TYR:O	1:D:451:ARG:NH1	2.45	0.49
1:D:385:MET:HE3	1:D:448:ILE:HG12	1.94	0.49
1:A:329:PHE:CD1	1:A:362:MET:HB2	2.47	0.49
1:D:382:ILE:HA	1:D:385:MET:HE2	1.94	0.49
1:A:220:ARG:HG2	1:A:387:THR:HG21	1.94	0.49
1:C:405:LYS:HD3	1:C:407:TYR:OH	2.12	0.49
1:C:493:LEU:HB2	1:C:568:TYR:CE2	2.47	0.49
1:A:117:VAL:O	3:A:902:GTP:O2'	2.31	0.49
1:C:483:ALA:O	1:C:485:PRO:HD3	2.13	0.49
1:B:489:LEU:HD21	1:B:563:TYR:CE2	2.48	0.48
1:B:223:PRO:CG	1:B:470:ARG:HG2	2.43	0.48
1:A:501:ASP:OD1	1:A:501:ASP:N	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:351:ALA:O	1:D:520:PHE:HA	2.14	0.48
1:C:239:MET:HE2	1:C:416:MET:HE3	1.94	0.48
1:D:476:LEU:CB	1:D:500:VAL:HG11	2.44	0.47
1:D:212:PRO:HD2	1:D:217:PHE:CD1	2.49	0.47
1:B:175:ALA:HB2	1:B:310:VAL:HG21	1.95	0.47
1:B:240:MET:HE2	1:B:416:MET:SD	2.54	0.47
1:A:241:PHE:CE2	1:A:245:ILE:HD11	2.50	0.47
1:A:245:ILE:HG23	1:A:250:ILE:HB	1.96	0.47
1:B:116:LYS:HE2	3:B:902:GTP:O2A	2.14	0.47
1:B:175:ALA:O	1:B:179:VAL:HG22	2.15	0.46
1:C:116:LYS:HE2	3:C:902:GTP:C3'	2.45	0.46
1:D:385:MET:HG2	1:D:454:PHE:CE2	2.49	0.46
1:C:470:ARG:HG3	1:C:473:TYR:CZ	2.50	0.46
1:B:522:CYS:HB2	1:B:530:ILE:HG21	1.96	0.46
1:C:254:MET:HE3	1:C:259:LEU:HD23	1.97	0.46
1:A:470:ARG:HG3	1:A:473:TYR:CE2	2.49	0.46
1:A:505:MET:CB	1:A:506:ASP:HA	2.46	0.46
1:C:148:LYS:HE2	1:C:148:LYS:HB2	1.63	0.45
1:C:175:ALA:HB1	1:C:199:VAL:HG12	1.97	0.45
1:D:167:HIS:HA	1:D:318:ARG:NH1	2.30	0.45
1:D:261:PRO:O	1:D:265:ILE:HD12	2.16	0.45
1:A:405:LYS:HD3	1:A:407:TYR:OH	2.17	0.45
1:D:175:ALA:HB1	1:D:199:VAL:HG12	1.98	0.45
1:A:265:ILE:O	1:A:269:LYS:HG3	2.17	0.45
1:B:197:LEU:HD23	1:B:197:LEU:HA	1.84	0.45
1:B:120:ASP:CG	1:B:318:ARG:HH22	2.25	0.45
1:B:179:VAL:HG11	1:B:199:VAL:HG21	1.99	0.45
1:B:353:ASP:OD1	1:B:530:ILE:HD11	2.16	0.45
1:C:377:LYS:HD2	1:C:456:TYR:CE1	2.51	0.45
1:A:120:ASP:OD1	1:A:121:PRO:HD2	2.17	0.45
1:C:443:GLU:O	1:C:447:GLN:HG2	2.17	0.45
1:A:359:LEU:HD12	1:A:359:LEU:HA	1.80	0.45
1:B:116:LYS:CE	3:B:902:GTP:H3'	2.46	0.44
1:D:505:MET:HG2	1:D:547:GLU:O	2.17	0.44
1:A:220:ARG:HH21	1:A:501:ASP:HA	1.82	0.44
1:A:385:MET:HE1	1:A:453:LEU:HB3	1.99	0.44
1:C:245:ILE:HD12	1:C:265:ILE:HD13	1.98	0.44
1:D:310:VAL:HG12	1:D:313:TRP:CE3	2.51	0.44
1:D:562:LEU:HD23	1:D:562:LEU:HA	1.80	0.44
1:B:116:LYS:HE3	3:B:902:GTP:H3'	1.99	0.44
1:C:470:ARG:HG3	1:C:473:TYR:CE2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:456:TYR:OH	1:A:459:GLU:HB2	2.17	0.43
1:C:285:TRP:HA	1:C:286:PRO:HD3	1.73	0.43
1:B:341:CYS:HB2	1:B:528:ARG:HA	2.00	0.43
1:C:167:HIS:HA	1:C:318:ARG:HD2	2.00	0.43
1:D:353:ASP:OD1	1:D:353:ASP:N	2.52	0.43
1:A:179:VAL:HG21	1:A:199:VAL:HG11	2.00	0.43
1:B:376:HIS:HE1	1:B:378:VAL:HB	1.83	0.43
1:D:297:LEU:HD23	1:D:297:LEU:HA	1.79	0.43
1:C:428:LEU:HD13	1:D:425:ASN:HB2	2.01	0.43
1:C:466:ILE:O	1:C:466:ILE:HD12	2.18	0.43
1:A:118:ILE:HD12	1:A:128:LEU:HD11	2.00	0.43
1:C:138:THR:HG21	1:C:244:LEU:CG	2.49	0.43
1:B:178:LEU:HD13	1:B:335:ILE:HG12	2.01	0.43
1:D:447:GLN:HG3	1:D:452:ASN:HB2	2.01	0.43
1:C:425:ASN:HB2	1:D:428:LEU:HD13	2.01	0.42
1:A:232:THR:HB	1:A:235:GLN:CD	2.43	0.42
1:C:116:LYS:HE3	3:C:902:GTP:HN22	1.84	0.42
1:C:179:VAL:HG21	1:C:199:VAL:HG11	2.01	0.42
1:D:143:ARG:NH1	1:D:422:LEU:O	2.53	0.42
1:A:471:GLU:H	1:A:471:GLU:HG3	1.31	0.42
1:B:377:LYS:HE3	1:B:377:LYS:HB2	1.92	0.42
1:C:135:ILE:O	1:C:138:THR:HB	2.19	0.42
1:C:416:MET:HE2	1:C:416:MET:HA	2.02	0.42
1:B:489:LEU:HD12	1:B:489:LEU:HA	1.85	0.42
1:C:314:ASP:OD1	1:C:318:ARG:HD2	2.19	0.42
1:A:174:LEU:HD23	1:A:174:LEU:HA	1.81	0.42
1:A:232:THR:HB	1:A:235:GLN:NE2	2.34	0.42
1:A:131:LEU:HD13	1:A:253:VAL:HG12	2.02	0.42
1:B:291:PRO:HG2	1:B:293:ASN:OD1	2.20	0.42
1:B:396:TYR:CD1	1:B:437:LYS:HD3	2.55	0.42
1:C:176:GLY:O	1:C:180:HIS:HB2	2.20	0.42
1:B:461:GLN:HA	1:B:462:PRO:HD3	1.94	0.42
1:A:484:LYS:HA	1:A:485:PRO:HD3	1.77	0.41
1:B:359:LEU:HD23	1:B:359:LEU:HA	1.88	0.41
1:C:300:ILE:O	1:C:309:ASP:HB2	2.20	0.41
1:D:447:GLN:HE21	1:D:452:ASN:ND2	2.00	0.41
1:A:544:LYS:HD2	1:A:544:LYS:N	2.35	0.41
1:A:563:TYR:O	1:A:567:GLN:HG2	2.20	0.41
1:B:220:ARG:HG2	1:B:387:THR:HG21	2.01	0.41
1:C:116:LYS:HE3	3:C:902:GTP:O1B	2.20	0.41
1:D:401:GLY:HA3	1:D:407:TYR:CE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:493:LEU:HD23	1:C:493:LEU:HA	1.87	0.41
1:D:115:MET:HB2	1:D:128:LEU:O	2.21	0.41
1:C:339:ARG:HD3	1:C:521:TYR:CE2	2.55	0.41
1:C:407:TYR:HD2	1:C:411:THR:HG22	1.85	0.41
1:A:223:PRO:CG	1:A:470:ARG:HG2	2.51	0.41
1:C:267:PHE:O	1:C:271:GLN:HG3	2.20	0.41
1:D:416:MET:HE1	1:D:419:TYR:HD2	1.84	0.41
1:A:240:MET:HE1	1:A:420:THR:OG1	2.20	0.41
1:B:118:ILE:HD13	3:B:902:GTP:C2'	2.50	0.41
1:B:407:TYR:HD2	1:B:411:THR:HG23	1.86	0.41
1:C:470:ARG:HA	1:C:473:TYR:CE1	2.55	0.41
1:A:190:GLN:O	1:A:294:LYS:NZ	2.44	0.41
1:A:343:VAL:HG22	1:A:529:ALA:CB	2.51	0.41
1:A:428:LEU:HD23	1:A:428:LEU:HA	1.84	0.41
1:B:189:LEU:O	1:B:190:GLN:HB2	2.20	0.41
1:C:179:VAL:CG2	1:C:199:VAL:HG11	2.51	0.41
1:D:563:TYR:O	1:D:567:GLN:HG2	2.21	0.41
1:B:300:ILE:O	1:B:310:VAL:HG22	2.20	0.41
1:D:564:ALA:O	1:D:568:TYR:HD1	2.04	0.41
1:A:138:THR:HG21	1:A:244:LEU:CG	2.48	0.40
1:B:120:ASP:OD1	1:B:121:PRO:HD2	2.21	0.40
1:D:226:ARG:O	1:D:229:VAL:HG22	2.20	0.40
1:C:439:LYS:O	1:C:443:GLU:HG2	2.22	0.40
1:C:456:TYR:OH	1:C:459:GLU:HB2	2.21	0.40
1:D:180:HIS:O	1:D:184:GLU:HG2	2.20	0.40
1:B:337:PHE:CD2	1:B:352:ARG:HD2	2.56	0.40
1:D:456:TYR:OH	1:D:459:GLU:HB2	2.20	0.40
1:A:142:GLN:O	1:A:145:ARG:HG3	2.21	0.40
1:A:220:ARG:NH2	1:A:500:VAL:O	2.54	0.40
1:A:403:GLY:C	1:A:405:LYS:H	2.29	0.40
1:B:155:TYR:O	1:B:451:ARG:NH1	2.49	0.40
1:D:285:TRP:HA	1:D:286:PRO:HD3	1.82	0.40
1:B:182:LEU:HD22	1:B:339:ARG:HA	2.04	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	451/526 (86%)	436 (97%)	15 (3%)	0	100	100
1	B	441/526 (84%)	425 (96%)	15 (3%)	1 (0%)	43	63
1	C	439/526 (84%)	414 (94%)	23 (5%)	2 (0%)	24	44
1	D	429/526 (82%)	408 (95%)	19 (4%)	2 (0%)	24	44
All	All	1760/2104 (84%)	1683 (96%)	72 (4%)	5 (0%)	36	56

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	490	ASP
1	C	582	GLN
1	D	490	ASP
1	C	547	GLU
1	D	525	ALA

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	376/470 (80%)	362 (96%)	14 (4%)	30	55
1	B	366/470 (78%)	356 (97%)	10 (3%)	39	65
1	C	370/470 (79%)	350 (95%)	20 (5%)	20	40
1	D	357/470 (76%)	341 (96%)	16 (4%)	24	47
All	All	1469/1880 (78%)	1409 (96%)	60 (4%)	27	51

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	115	MET
1	A	118	ILE
1	A	138	THR
1	A	229	VAL
1	A	232	THR
1	A	300	ILE
1	A	346	GLU
1	A	359	LEU
1	A	434	THR
1	A	451	ARG
1	A	470	ARG
1	A	471	GLU
1	A	488	LEU
1	A	500	VAL
1	B	138	THR
1	B	262	GLU
1	B	332	LYS
1	B	426	ILE
1	B	455	LYS
1	B	463	THR
1	B	470	ARG
1	B	471	GLU
1	B	505	MET
1	B	580	LYS
1	C	118	ILE
1	C	126	ILE
1	C	138	THR
1	C	210	HIS
1	C	230	LYS
1	C	232	THR
1	C	300	ILE
1	C	348	ARG
1	C	368	SER
1	C	371	ARG
1	C	377	LYS
1	C	411	THR
1	C	434	THR
1	C	470	ARG
1	C	478	LYS
1	C	487	VAL
1	C	489	LEU
1	C	528	ARG

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Mol	Chain	Res	Type
1	C	559	ARG
1	C	560	LYS
1	D	118	ILE
1	D	126	ILE
1	D	136	ILE
1	D	138	THR
1	D	232	THR
1	D	245	ILE
1	D	262	GLU
1	D	265	ILE
1	D	277	GLU
1	D	300	ILE
1	D	400	THR
1	D	411	THR
1	D	447	GLN
1	D	496	GLU
1	D	557	VAL
1	D	576	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	125	HIS
1	A	200	GLN
1	A	235	GLN
1	A	321	HIS
1	A	370	HIS
1	B	200	GLN
1	B	243	HIS
1	B	328	ASN
1	C	345	ASN
1	C	425	ASN
1	C	504	ASN
1	C	548	GLN
1	D	235	GLN
1	D	327	ASN
1	D	364	HIS
1	D	425	ASN
1	D	447	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 ⓘ

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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	$\# Z > 2$
3	GTP	A	902	-	33,34,34	2.88	11 (33%)	50,54,54	1.76	11 (22%)
3	GTP	B	902	-	33,34,34	2.90	12 (36%)	50,54,54	2.03	13 (26%)
3	GTP	C	902	-	33,34,34	2.86	11 (33%)	50,54,54	1.83	13 (26%)

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	A	902	-	-	4/22/38/38	0/3/3/3
3	GTP	B	902	-	-	6/22/38/38	0/3/3/3
3	GTP	C	902	-	-	8/22/38/38	0/3/3/3

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	902	GTP	C3'-C4'	-8.71	1.30	1.53
3	B	902	GTP	C3'-C4'	-8.58	1.31	1.53
3	A	902	GTP	C3'-C4'	-8.34	1.31	1.53
3	A	902	GTP	O4'-C4'	7.48	1.61	1.45
3	B	902	GTP	O4'-C4'	7.09	1.60	1.45
3	C	902	GTP	O4'-C4'	7.07	1.60	1.45
3	A	902	GTP	C4-N3	5.49	1.46	1.34
3	B	902	GTP	C2-N3	5.47	1.46	1.33
3	C	902	GTP	C4-N3	5.20	1.46	1.34
3	B	902	GTP	C4-N3	5.17	1.46	1.34
3	A	902	GTP	C2-N3	5.14	1.45	1.33
3	C	902	GTP	C2-N3	4.78	1.44	1.33
3	B	902	GTP	C2-N2	4.29	1.44	1.34
3	C	902	GTP	O4'-C1'	-4.26	1.32	1.42
3	B	902	GTP	O4'-C1'	-4.22	1.32	1.42
3	A	902	GTP	C2-N2	4.03	1.43	1.34
3	A	902	GTP	O4'-C1'	-3.82	1.33	1.42
3	C	902	GTP	PB-O3B	3.79	1.63	1.59
3	C	902	GTP	O3'-C3'	3.74	1.52	1.43
3	A	902	GTP	PB-O3B	3.70	1.63	1.59
3	B	902	GTP	PB-O3B	3.59	1.63	1.59
3	C	902	GTP	C2-N2	3.56	1.42	1.34
3	B	902	GTP	O3'-C3'	3.50	1.51	1.43
3	C	902	GTP	PB-O3A	3.42	1.63	1.59
3	A	902	GTP	O3'-C3'	3.17	1.50	1.43
3	A	902	GTP	PB-O3A	3.04	1.62	1.59
3	B	902	GTP	PB-O3A	2.76	1.62	1.59
3	B	902	GTP	O2'-C2'	-2.61	1.36	1.43
3	B	902	GTP	C5-N7	-2.59	1.33	1.39
3	A	902	GTP	C2-N1	2.42	1.43	1.37
3	A	902	GTP	O2'-C2'	-2.40	1.37	1.43
3	C	902	GTP	O2'-C2'	-2.14	1.37	1.43
3	B	902	GTP	C5-C6	2.14	1.52	1.44
3	C	902	GTP	C5-C6	2.05	1.52	1.44

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	902	GTP	C5-C4-N3	-7.08	117.12	128.39
3	C	902	GTP	C5-C4-N3	-6.27	118.42	128.39
3	A	902	GTP	C5-C4-N3	-5.90	119.00	128.39
3	B	902	GTP	C2-N3-C4	5.26	121.37	112.30
3	C	902	GTP	C2-N3-C4	5.19	121.24	112.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	902	GTP	N9-C4-N3	4.93	135.81	125.95
3	A	902	GTP	C2-N3-C4	4.45	119.97	112.30
3	A	902	GTP	N9-C4-N3	3.87	133.70	125.95
3	C	902	GTP	N9-C4-N3	3.51	132.97	125.95
3	B	902	GTP	C2'-C1'-N9	3.13	121.95	113.25
3	B	902	GTP	C1'-N9-C8	-3.02	118.16	126.73
3	B	902	GTP	C3'-C2'-C1'	3.01	107.17	101.46
3	C	902	GTP	C2-N1-C6	-2.92	119.81	125.11
3	A	902	GTP	C2-N1-C6	-2.81	120.02	125.11
3	B	902	GTP	C6-C5-C4	2.77	123.00	118.83
3	B	902	GTP	N9-C8-N7	-2.72	108.35	113.40
3	C	902	GTP	C5-C6-N1	2.71	120.16	113.25
3	A	902	GTP	N9-C8-N7	-2.71	108.37	113.40
3	B	902	GTP	C1'-N9-C4	2.59	134.12	126.49
3	A	902	GTP	C2'-C3'-C4'	2.56	107.56	102.61
3	A	902	GTP	C3'-C2'-C1'	2.50	106.20	101.46
3	A	902	GTP	O6-C6-C5	-2.45	120.06	126.53
3	A	902	GTP	C8-N7-C5	2.40	108.54	104.26
3	C	902	GTP	C3'-C2'-C1'	2.38	105.96	101.46
3	B	902	GTP	C8-N7-C5	2.38	108.49	104.26
3	B	902	GTP	C2-N1-C6	-2.34	120.87	125.11
3	A	902	GTP	C5-C6-N1	2.33	119.18	113.25
3	C	902	GTP	C2'-C1'-N9	2.28	119.60	113.25
3	C	902	GTP	O3B-PB-O1B	-2.28	103.85	110.70
3	B	902	GTP	C2'-C3'-C4'	2.18	106.82	102.61
3	C	902	GTP	N9-C8-N7	-2.16	109.39	113.40
3	B	902	GTP	O6-C6-C5	-2.11	120.95	126.53
3	C	902	GTP	C8-N7-C5	2.08	107.96	104.26
3	A	902	GTP	O2G-PG-O3B	2.07	111.58	104.64
3	C	902	GTP	C4-C5-N7	-2.04	107.44	110.67
3	C	902	GTP	O2B-PB-O3B	2.03	112.76	107.27
3	C	902	GTP	O3G-PG-O3B	2.02	111.42	104.64

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	902	GTP	C5'-O5'-PA-O1A
3	C	902	GTP	PB-O3B-PG-O3G
3	C	902	GTP	C5'-O5'-PA-O3A
3	C	902	GTP	C5'-O5'-PA-O1A
3	A	902	GTP	O4'-C4'-C5'-O5'

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Mol	Chain	Res	Type	Atoms
3	A	902	GTP	C3'-C4'-C5'-O5'
3	B	902	GTP	O4'-C4'-C5'-O5'
3	B	902	GTP	C3'-C4'-C5'-O5'
3	A	902	GTP	C4'-C5'-O5'-PA
3	B	902	GTP	PB-O3A-PA-O2A
3	C	902	GTP	PB-O3A-PA-O2A
3	C	902	GTP	C5'-O5'-PA-O2A
3	B	902	GTP	PA-O3A-PB-O2B
3	C	902	GTP	PG-O3B-PB-O1B
3	B	902	GTP	PG-O3B-PB-O2B
3	C	902	GTP	PB-O3B-PG-O1G
3	B	902	GTP	C2'-C1'-N9-C4
3	C	902	GTP	C2'-C1'-N9-C4

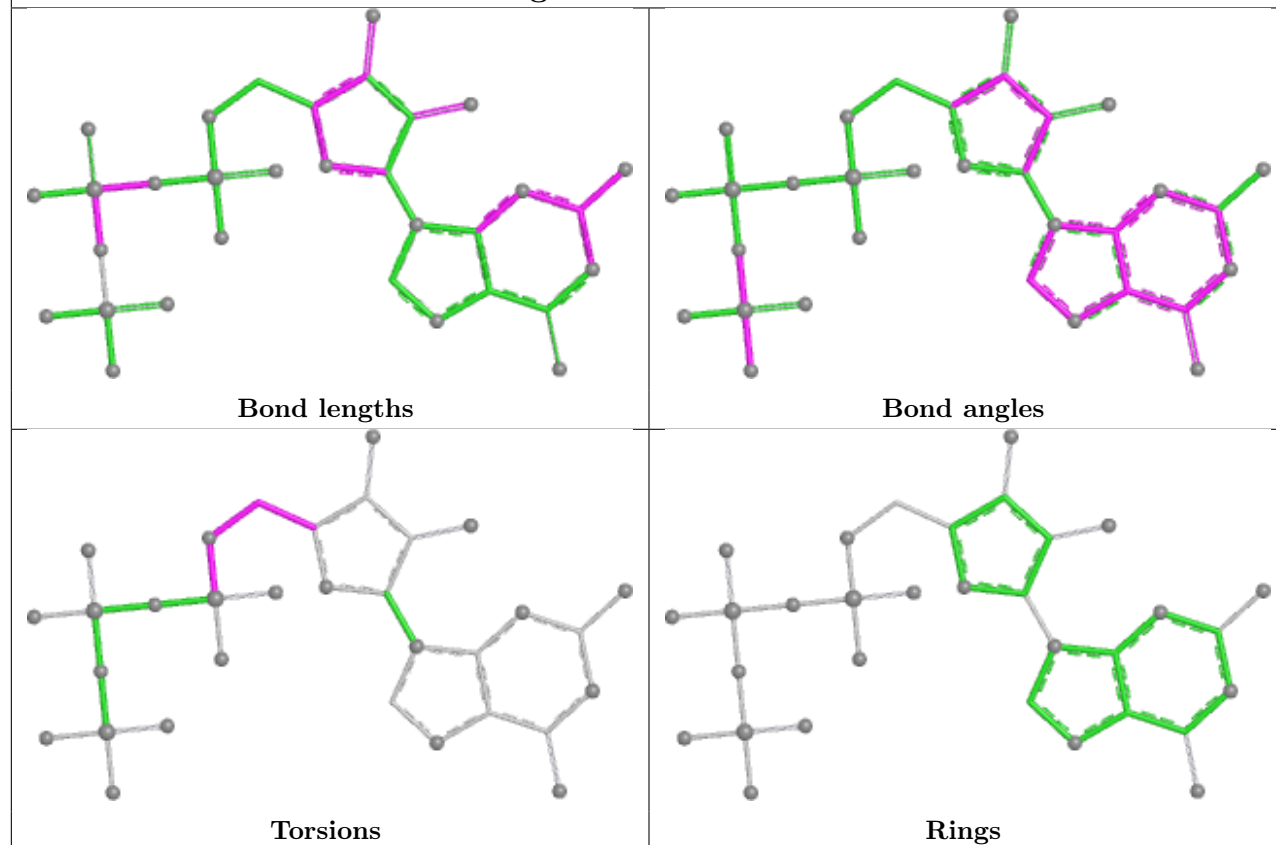
There are no ring outliers.

3 monomers are involved in 19 short contacts:

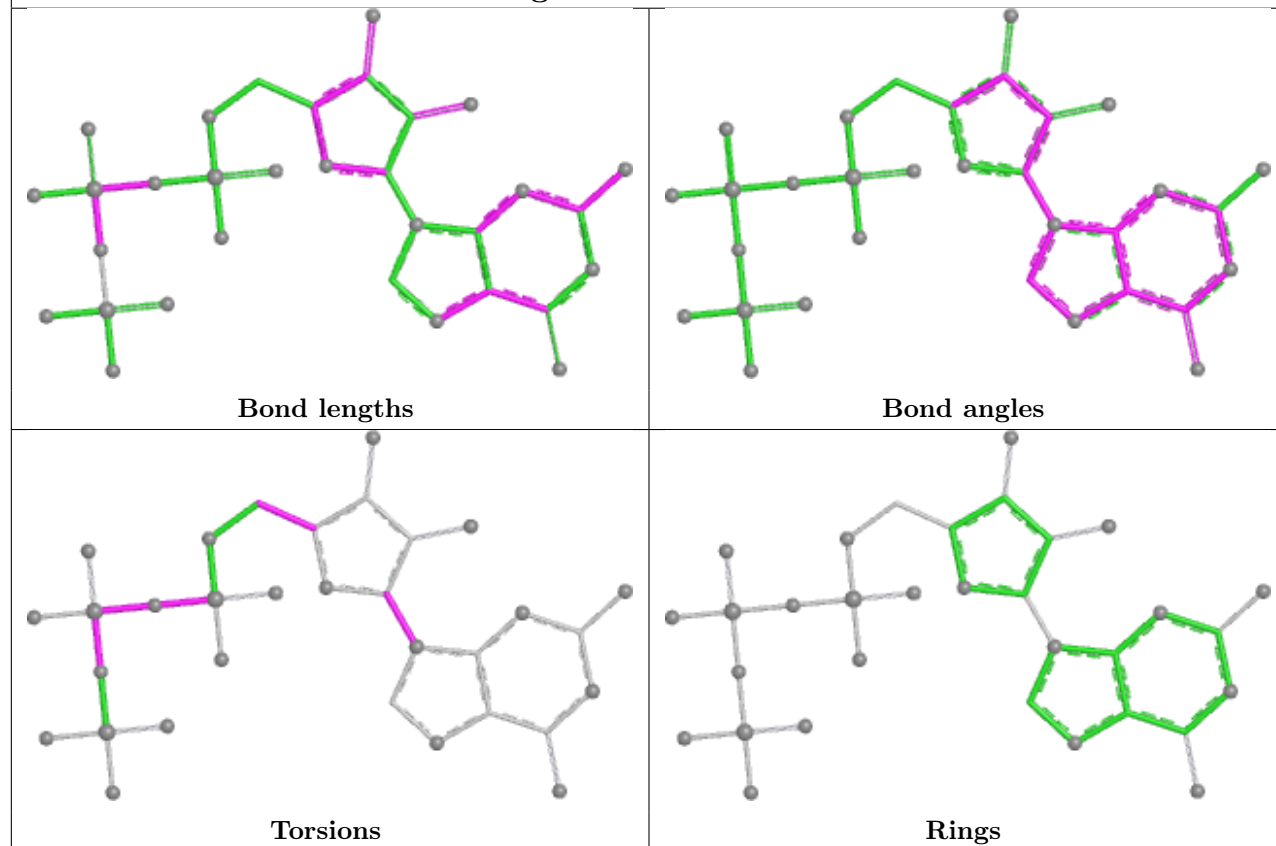
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	902	GTP	2	0
3	B	902	GTP	9	0
3	C	902	GTP	8	0

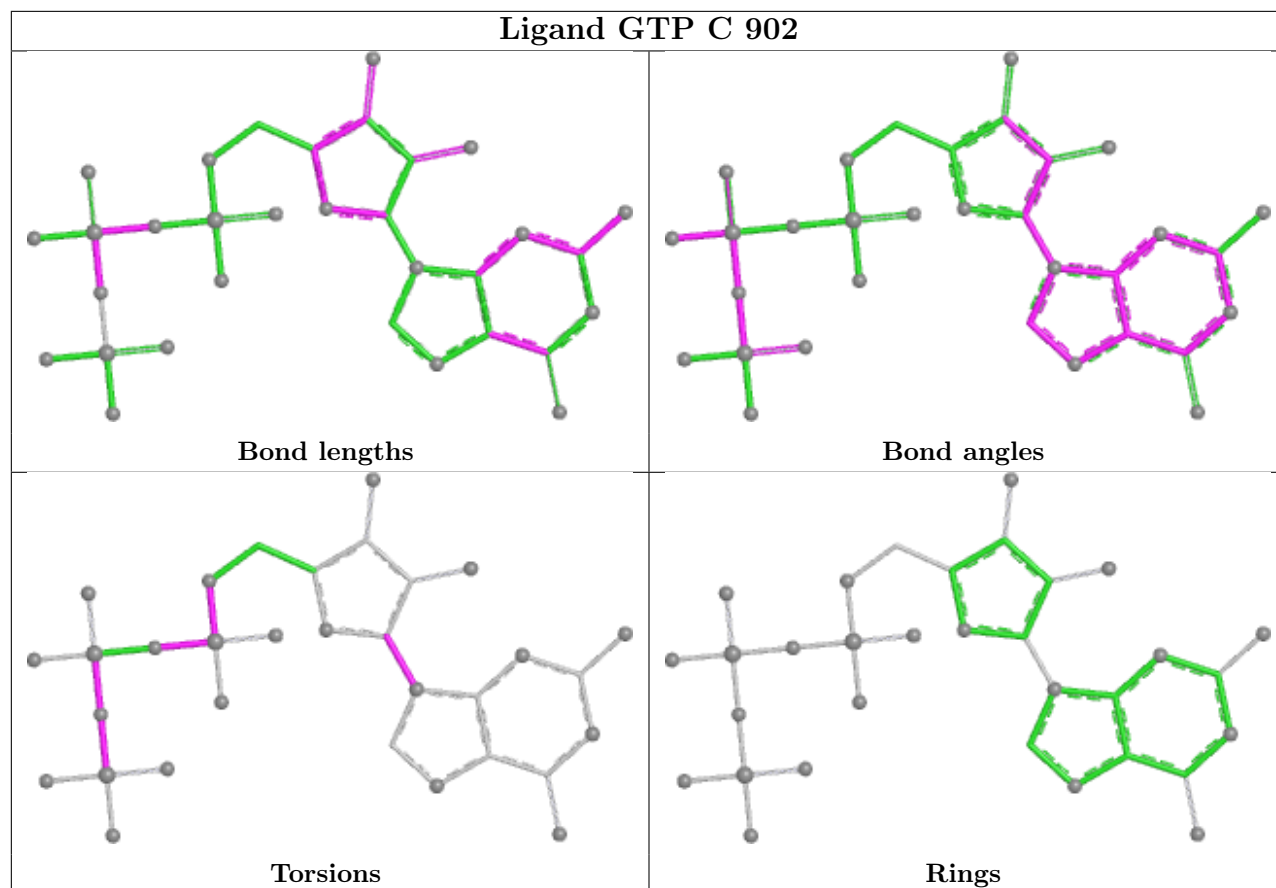
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 902



Ligand GTP B 902





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	459/526 (87%)	-0.16	8 (1%) 69 65	29, 46, 76, 106	0
1	B	449/526 (85%)	0.11	27 (6%) 27 23	31, 48, 91, 114	0
1	C	447/526 (84%)	0.04	9 (2%) 65 61	35, 51, 79, 99	0
1	D	439/526 (83%)	0.13	15 (3%) 48 43	34, 51, 84, 93	0
All	All	1794/2104 (85%)	0.03	59 (3%) 49 44	29, 49, 83, 114	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	516	ASP	4.8
1	A	507	TYR	4.3
1	C	506	ASP	4.1
1	B	505	MET	3.9
1	A	506	ASP	3.8
1	A	510	GLN	3.6
1	C	583	ASP	3.6
1	B	307	GLY	3.6
1	D	304	LYS	3.6
1	A	533	THR	3.4
1	B	531	ARG	3.4
1	C	114	THR	3.3
1	B	345	ASN	3.3
1	A	545	PHE	3.3
1	B	284	LEU	3.3
1	B	584	GLY	3.3
1	B	282	ASP	3.2
1	D	522	CYS	3.2
1	B	523	LYS	3.1
1	B	524	THR	3.1
1	B	532	ILE	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	546	ALA	3.1
1	B	329	PHE	3.0
1	B	277	GLU	2.9
1	B	545	PHE	2.9
1	B	355	GLU	2.9
1	A	584	GLY	2.8
1	B	522	CYS	2.7
1	C	582	GLN	2.7
1	B	344	ASP	2.7
1	C	545	PHE	2.6
1	A	282	ASP	2.6
1	D	329	PHE	2.6
1	B	528	ARG	2.6
1	D	284	LEU	2.6
1	D	328	ASN	2.6
1	C	490	ASP	2.6
1	D	345	ASN	2.5
1	B	466	ILE	2.5
1	D	582	GLN	2.5
1	B	514	PRO	2.5
1	B	276	LEU	2.4
1	B	582	GLN	2.4
1	A	279	PRO	2.4
1	B	525	ALA	2.4
1	C	278	SER	2.3
1	D	487	VAL	2.3
1	D	526	PRO	2.2
1	B	463	THR	2.2
1	D	334	PHE	2.2
1	D	513	ASN	2.2
1	D	523	LYS	2.2
1	B	283	SER	2.1
1	C	487	VAL	2.1
1	D	527	ASN	2.1
1	B	334	PHE	2.0
1	B	529	ALA	2.0
1	D	529	ALA	2.0
1	B	517	HIS	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 [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

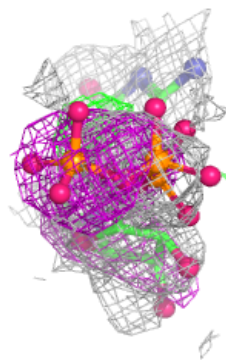
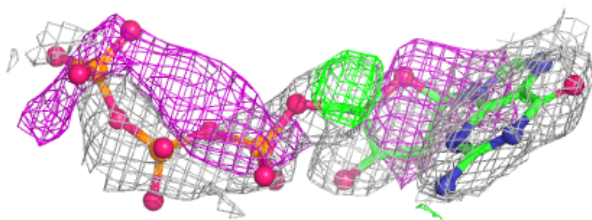
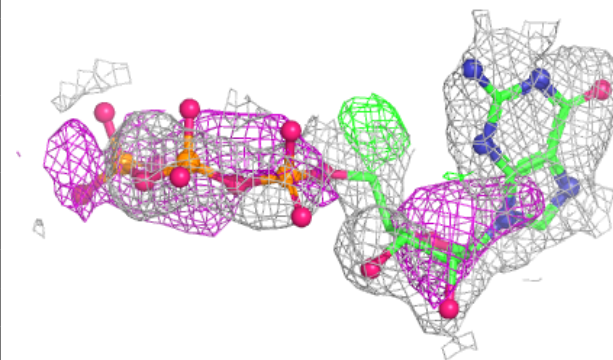
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GTP	A	902	32/32	0.57	0.16	53,80,148,151	0
3	GTP	B	902	32/32	0.63	0.15	48,71,103,119	0
3	GTP	C	902	32/32	0.66	0.15	54,71,109,117	0
2	ZN	C	901	1/1	0.98	0.06	42,42,42,42	0
2	ZN	D	901	1/1	0.99	0.05	39,39,39,39	0
2	ZN	B	901	1/1	0.99	0.04	39,39,39,39	0
2	ZN	A	901	1/1	1.00	0.05	34,34,34,34	0

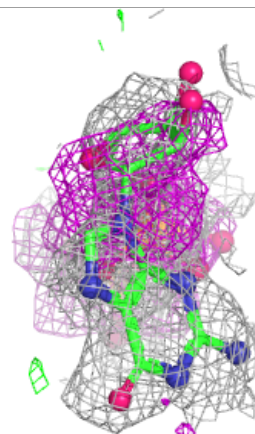
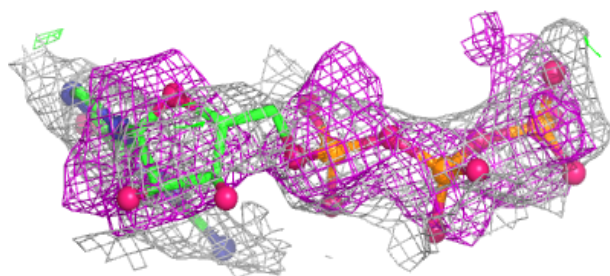
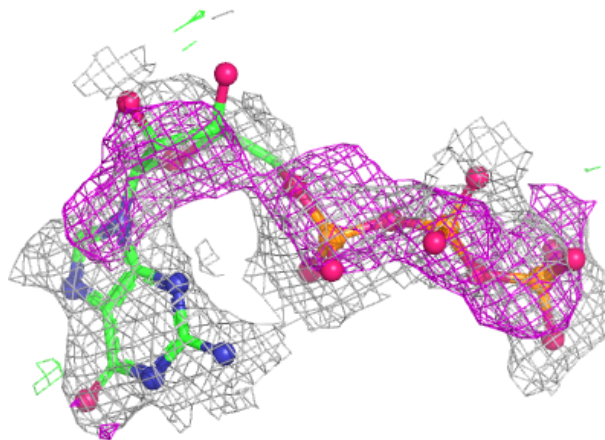
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around GTP A 902:

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

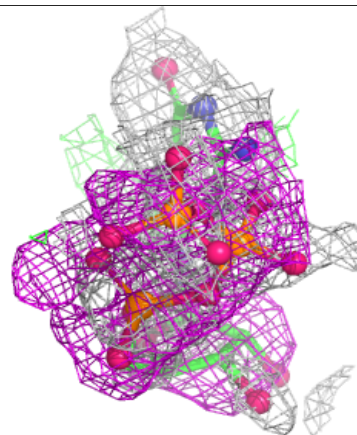
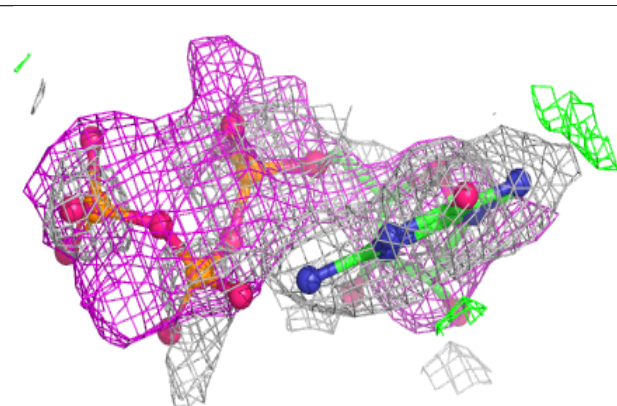
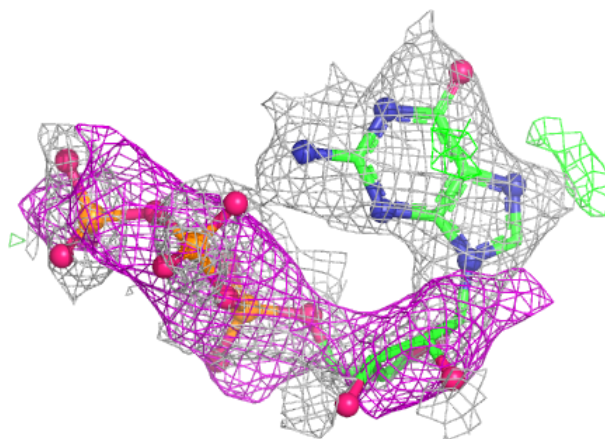
**Electron density around GTP B 902:**

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



Electron density around GTP C 902:

$2mF_o - DF_c$ (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.