



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:30 PM UTC

PDB ID : 5AO3 / pdb_00005ao3
Title : Crystal structure of human SAMHD1 (amino acid residues 115-626) bound to GTP
Authors : Schwefel, D.; Taylor, I.A.
Deposited on : 2015-09-09
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

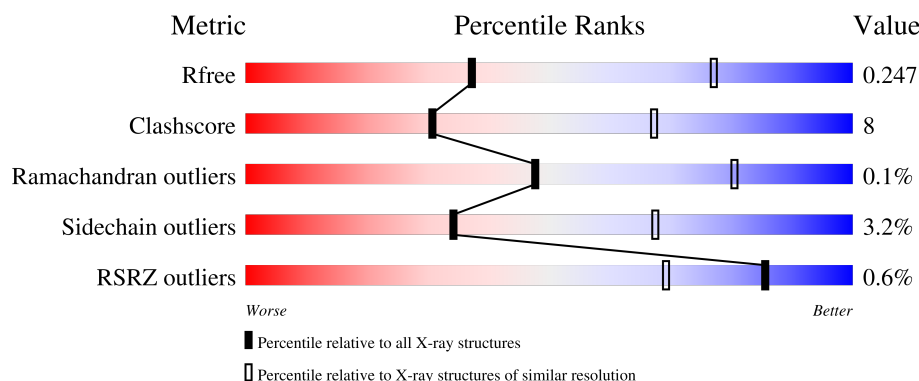
1 Overall quality at a glance

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	538	% 62% 18% • 18%
1	B	538	65% 15% • 19%
1	C	538	64% 16% • 18%
1	D	538	% 65% 16% • 18%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 14005 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 TRIPHOSPHO-HYDROLASE SAMHD1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	0	2	0
			3471	2228	598	626	19			
1	B	438	Total	C	N	O	S	0	1	0
			3446	2210	594	623	19			
1	C	439	Total	C	N	O	S	0	1	0
			3426	2200	590	617	19			
1	D	439	Total	C	N	O	S	0	1	0
			3470	2227	605	619	19			

There are 104 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	89	MET	-	expression tag	UNP Q9Y3Z3
A	90	ALA	-	expression tag	UNP Q9Y3Z3
A	91	SER	-	expression tag	UNP Q9Y3Z3
A	92	TRP	-	expression tag	UNP Q9Y3Z3
A	93	SER	-	expression tag	UNP Q9Y3Z3
A	94	HIS	-	expression tag	UNP Q9Y3Z3
A	95	PRO	-	expression tag	UNP Q9Y3Z3
A	96	GLN	-	expression tag	UNP Q9Y3Z3
A	97	PHE	-	expression tag	UNP Q9Y3Z3
A	98	GLU	-	expression tag	UNP Q9Y3Z3
A	99	LYS	-	expression tag	UNP Q9Y3Z3
A	100	GLY	-	expression tag	UNP Q9Y3Z3
A	101	ALA	-	expression tag	UNP Q9Y3Z3
A	102	LEU	-	expression tag	UNP Q9Y3Z3
A	103	GLU	-	expression tag	UNP Q9Y3Z3
A	104	VAL	-	expression tag	UNP Q9Y3Z3
A	105	LEU	-	expression tag	UNP Q9Y3Z3
A	106	PHE	-	expression tag	UNP Q9Y3Z3
A	107	GLN	-	expression tag	UNP Q9Y3Z3
A	108	GLY	-	expression tag	UNP Q9Y3Z3

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Chain	Residue	Modelled	Actual	Comment	Reference
A	109	PRO	-	expression tag	UNP Q9Y3Z3
A	110	GLY	-	expression tag	UNP Q9Y3Z3
A	111	TYR	-	expression tag	UNP Q9Y3Z3
A	112	GLN	-	expression tag	UNP Q9Y3Z3
A	113	ASP	-	expression tag	UNP Q9Y3Z3
A	114	PRO	-	expression tag	UNP Q9Y3Z3
B	89	MET	-	expression tag	UNP Q9Y3Z3
B	90	ALA	-	expression tag	UNP Q9Y3Z3
B	91	SER	-	expression tag	UNP Q9Y3Z3
B	92	TRP	-	expression tag	UNP Q9Y3Z3
B	93	SER	-	expression tag	UNP Q9Y3Z3
B	94	HIS	-	expression tag	UNP Q9Y3Z3
B	95	PRO	-	expression tag	UNP Q9Y3Z3
B	96	GLN	-	expression tag	UNP Q9Y3Z3
B	97	PHE	-	expression tag	UNP Q9Y3Z3
B	98	GLU	-	expression tag	UNP Q9Y3Z3
B	99	LYS	-	expression tag	UNP Q9Y3Z3
B	100	GLY	-	expression tag	UNP Q9Y3Z3
B	101	ALA	-	expression tag	UNP Q9Y3Z3
B	102	LEU	-	expression tag	UNP Q9Y3Z3
B	103	GLU	-	expression tag	UNP Q9Y3Z3
B	104	VAL	-	expression tag	UNP Q9Y3Z3
B	105	LEU	-	expression tag	UNP Q9Y3Z3
B	106	PHE	-	expression tag	UNP Q9Y3Z3
B	107	GLN	-	expression tag	UNP Q9Y3Z3
B	108	GLY	-	expression tag	UNP Q9Y3Z3
B	109	PRO	-	expression tag	UNP Q9Y3Z3
B	110	GLY	-	expression tag	UNP Q9Y3Z3
B	111	TYR	-	expression tag	UNP Q9Y3Z3
B	112	GLN	-	expression tag	UNP Q9Y3Z3
B	113	ASP	-	expression tag	UNP Q9Y3Z3
B	114	PRO	-	expression tag	UNP Q9Y3Z3
C	89	MET	-	expression tag	UNP Q9Y3Z3
C	90	ALA	-	expression tag	UNP Q9Y3Z3
C	91	SER	-	expression tag	UNP Q9Y3Z3
C	92	TRP	-	expression tag	UNP Q9Y3Z3
C	93	SER	-	expression tag	UNP Q9Y3Z3
C	94	HIS	-	expression tag	UNP Q9Y3Z3
C	95	PRO	-	expression tag	UNP Q9Y3Z3
C	96	GLN	-	expression tag	UNP Q9Y3Z3
C	97	PHE	-	expression tag	UNP Q9Y3Z3
C	98	GLU	-	expression tag	UNP Q9Y3Z3

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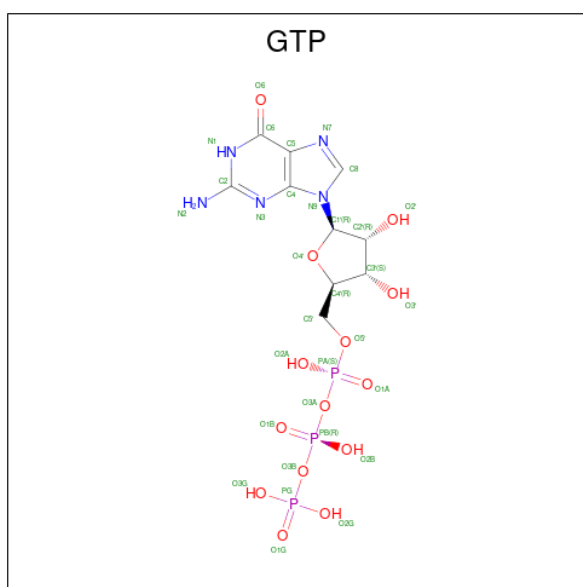
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Chain	Residue	Modelled	Actual	Comment	Reference
C	99	LYS	-	expression tag	UNP Q9Y3Z3
C	100	GLY	-	expression tag	UNP Q9Y3Z3
C	101	ALA	-	expression tag	UNP Q9Y3Z3
C	102	LEU	-	expression tag	UNP Q9Y3Z3
C	103	GLU	-	expression tag	UNP Q9Y3Z3
C	104	VAL	-	expression tag	UNP Q9Y3Z3
C	105	LEU	-	expression tag	UNP Q9Y3Z3
C	106	PHE	-	expression tag	UNP Q9Y3Z3
C	107	GLN	-	expression tag	UNP Q9Y3Z3
C	108	GLY	-	expression tag	UNP Q9Y3Z3
C	109	PRO	-	expression tag	UNP Q9Y3Z3
C	110	GLY	-	expression tag	UNP Q9Y3Z3
C	111	TYR	-	expression tag	UNP Q9Y3Z3
C	112	GLN	-	expression tag	UNP Q9Y3Z3
C	113	ASP	-	expression tag	UNP Q9Y3Z3
C	114	PRO	-	expression tag	UNP Q9Y3Z3
D	89	MET	-	expression tag	UNP Q9Y3Z3
D	90	ALA	-	expression tag	UNP Q9Y3Z3
D	91	SER	-	expression tag	UNP Q9Y3Z3
D	92	TRP	-	expression tag	UNP Q9Y3Z3
D	93	SER	-	expression tag	UNP Q9Y3Z3
D	94	HIS	-	expression tag	UNP Q9Y3Z3
D	95	PRO	-	expression tag	UNP Q9Y3Z3
D	96	GLN	-	expression tag	UNP Q9Y3Z3
D	97	PHE	-	expression tag	UNP Q9Y3Z3
D	98	GLU	-	expression tag	UNP Q9Y3Z3
D	99	LYS	-	expression tag	UNP Q9Y3Z3
D	100	GLY	-	expression tag	UNP Q9Y3Z3
D	101	ALA	-	expression tag	UNP Q9Y3Z3
D	102	LEU	-	expression tag	UNP Q9Y3Z3
D	103	GLU	-	expression tag	UNP Q9Y3Z3
D	104	VAL	-	expression tag	UNP Q9Y3Z3
D	105	LEU	-	expression tag	UNP Q9Y3Z3
D	106	PHE	-	expression tag	UNP Q9Y3Z3
D	107	GLN	-	expression tag	UNP Q9Y3Z3
D	108	GLY	-	expression tag	UNP Q9Y3Z3
D	109	PRO	-	expression tag	UNP Q9Y3Z3
D	110	GLY	-	expression tag	UNP Q9Y3Z3
D	111	TYR	-	expression tag	UNP Q9Y3Z3
D	112	GLN	-	expression tag	UNP Q9Y3Z3
D	113	ASP	-	expression tag	UNP Q9Y3Z3
D	114	PRO	-	expression tag	UNP Q9Y3Z3

- Molecule 2 is FE (III) ION (CCD ID: FE) (formula: Fe).

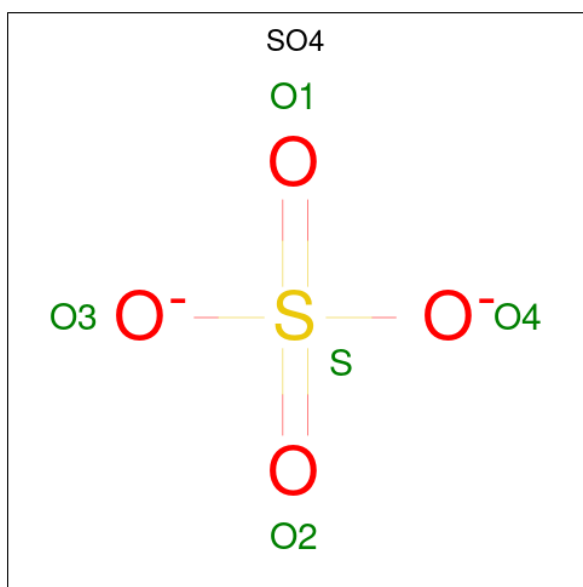
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Fe	0	0
			1	1		
2	B	1	Total	Fe	0	0
			1	1		
2	C	1	Total	Fe	0	0
			1	1		
2	D	1	Total	Fe	0	0
			1	1		

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

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 5 is water.

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

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	2	Total	O	0	0
			2	2		
5	C	2	Total	O	0	0
			2	2		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	81.92Å 95.73Å 97.59Å 91.37° 108.97° 115.33°	Depositor
Resolution (Å)	29.66 – 3.00 29.66 – 3.00	Depositor EDS
% Data completeness (in resolution range)	91.1 (29.66-3.00) 91.1 (29.66-3.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 3.00Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.8.4_1496)	Depositor
R, R_{free}	0.177 , 0.247 0.181 , 0.247	Depositor DCC
R_{free} test set	2275 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	69.7	Xtriage
Anisotropy	0.198	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 66.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14005	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.38% 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: SO4, FE, GTP

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.66	0/3562	1.03	9/4829 (0.2%)
1	B	0.65	1/3532 (0.0%)	1.01	9/4784 (0.2%)
1	C	0.58	0/3513	0.99	7/4763 (0.1%)
1	D	0.62	0/3558	0.99	5/4811 (0.1%)
All	All	0.63	1/14165 (0.0%)	1.01	30/19187 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	222	ILE	CA-CB	6.13	1.58	1.53

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	576	ARG	N-CA-C	-9.70	101.11	113.72
1	C	576	ARG	N-CA-C	-7.67	103.75	113.72
1	A	285	TRP	CA-C-N	6.92	126.63	119.64
1	A	285	TRP	C-N-CA	6.92	126.63	119.64
1	B	580	LYS	CA-C-N	6.71	126.70	119.78
1	B	580	LYS	C-N-CA	6.71	126.70	119.78
1	A	513	ASN	CA-C-N	6.46	127.91	119.84
1	A	513	ASN	C-N-CA	6.46	127.91	119.84
1	A	226	ARG	CA-C-N	6.24	126.14	119.28
1	A	226	ARG	C-N-CA	6.24	126.14	119.28
1	C	300	ILE	N-CA-C	6.18	116.36	110.74
1	A	458	GLY	N-CA-C	5.86	119.32	110.18
1	A	305	ARG	N-CA-C	5.82	117.49	111.03
1	D	525	ALA	CA-C-N	5.74	125.27	119.19
1	D	525	ALA	C-N-CA	5.74	125.27	119.19
1	C	305	ARG	N-CA-C	5.69	117.35	111.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	525	ALA	CA-C-N	5.51	125.03	119.19
1	C	525	ALA	C-N-CA	5.51	125.03	119.19
1	B	285	TRP	CA-C-N	5.49	125.18	119.64
1	B	285	TRP	C-N-CA	5.49	125.18	119.64
1	C	466	ILE	N-CA-C	5.42	116.04	108.89
1	B	476	LEU	CA-C-N	-5.35	113.39	119.28
1	B	476	LEU	C-N-CA	-5.35	113.39	119.28
1	C	464	GLY	N-CA-C	-5.29	103.83	112.83
1	A	222	ILE	CB-CA-C	-5.23	108.81	114.35
1	B	120	ASP	CA-C-N	5.12	124.73	119.56
1	B	120	ASP	C-N-CA	5.12	124.73	119.56
1	B	179	VAL	CB-CA-C	-5.11	105.43	111.97
1	D	222	ILE	CA-C-N	-5.08	113.70	119.28
1	D	222	ILE	C-N-CA	-5.08	113.70	119.28

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	3471	0	3299	65	0
1	B	3446	0	3277	56	0
1	C	3426	0	3241	55	0
1	D	3470	0	3348	54	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
3	A	32	0	12	4	0
3	C	32	0	12	1	0
3	D	64	0	24	2	0
4	A	15	0	0	0	0
4	B	20	0	0	2	0
4	C	10	0	0	0	0
4	D	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	1	0	0	1	0
5	B	2	0	0	2	0
5	C	2	0	0	0	0
All	All	14005	0	13213	226	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 (226) 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)
3:A:1585:GTP:H5''	1:C:451:ARG:HH21	1.35	0.89
1:B:371:ARG:NH1	4:B:1585:SO4:O1	2.12	0.83
1:A:505:MET:HE1	1:A:547:GLU:HB2	1.65	0.78
1:A:179:VAL:HG22	1:A:300:ILE:HD13	1.68	0.75
1:B:352:ARG:HB3	1:B:521:TYR:CZ	2.22	0.75
1:A:442[B]:ARG:HD2	1:A:446:LYS:HE3	1.68	0.74
1:B:402:ALA:HB2	1:B:417:GLU:CD	2.16	0.71
1:B:175:ALA:HB1	1:B:199:VAL:HG12	1.73	0.71
1:C:320:CYS:SG	1:C:327:ASN:HB2	2.30	0.70
1:A:175:ALA:HB1	1:A:199:VAL:HG12	1.71	0.69
1:D:497:ASP:OD2	1:D:556:LYS:NZ	2.26	0.69
1:B:116:LYS:NZ	3:D:1585:GTP:O1A	2.25	0.67
1:D:179:VAL:HG22	1:D:300:ILE:HD13	1.76	0.67
1:D:308:ILE:HG12	1:D:362:MET:HE1	1.75	0.67
1:C:175:ALA:HB1	1:C:199:VAL:HG12	1.76	0.67
1:A:428:LEU:HD13	1:C:425:ASN:HB2	1.76	0.67
1:B:179:VAL:HG22	1:B:300:ILE:HD13	1.77	0.66
1:A:513:ASN:HB3	1:A:515:ILE:HG22	1.78	0.65
1:D:442:ARG:HG2	1:D:446:LYS:HE3	1.78	0.64
1:B:337:PHE:HB3	1:B:352:ARG:HG2	1.79	0.64
1:D:220:ARG:HG2	1:D:387:THR:HG21	1.79	0.63
1:D:576:ARG:HG2	1:D:578:PHE:CE2	2.33	0.63
1:B:428:LEU:HD13	1:D:425:ASN:HB2	1.81	0.62
1:C:175:ALA:O	1:C:179:VAL:HG23	1.98	0.62
1:B:372:ARG:NH1	4:B:1586:SO4:O1	2.31	0.60
1:B:439:LYS:NZ	1:B:443:GLU:OE2	2.36	0.59
3:A:1585:GTP:O2B	1:C:378:VAL:HG21	2.02	0.58
1:A:431:LEU:HD13	1:A:445:LEU:HB3	1.86	0.58
1:B:566:ARG:NH2	1:B:582:GLN:O	2.37	0.58
1:D:275:PRO:HB3	1:D:286:PRO:HB2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:299:GLU:OE2	1:C:305:ARG:NH2	2.37	0.57
1:C:197:LEU:O	1:C:201:ILE:HG13	2.04	0.57
1:A:462:PRO:HA	1:A:578:PHE:CD1	2.39	0.57
1:C:397:ILE:HG21	1:C:426:ILE:HD11	1.86	0.56
1:D:175:ALA:O	1:D:179:VAL:HG23	2.05	0.56
1:D:360:TYR:HE1	1:D:515:ILE:HG13	1.70	0.56
1:D:355:GLU:O	1:D:355:GLU:HG3	2.06	0.56
1:A:381:ILE:HG23	1:A:454:PHE:HB2	1.88	0.55
1:D:431:LEU:HD23	1:D:432:TYR:CE1	2.41	0.55
1:B:285:TRP:CH2	1:B:287:TYR:HD2	2.23	0.55
1:B:576:ARG:HG2	1:B:578:PHE:CE2	2.41	0.55
1:C:144:LEU:HD23	1:C:147:ILE:HD12	1.88	0.55
1:B:408:ARG:HG2	1:B:408:ARG:HH11	1.72	0.55
1:A:303:ASN:ND2	1:A:306:ASN:OD1	2.40	0.54
1:C:116:LYS:NZ	3:C:1585:GTP:O2A	2.38	0.54
1:D:483:ALA:O	1:D:485:PRO:HD3	2.06	0.54
1:A:343:VAL:HG22	1:A:529:ALA:HB2	1.90	0.54
3:A:1585:GTP:H5"	1:C:451:ARG:NH2	2.15	0.54
1:D:175:ALA:HB1	1:D:199:VAL:HG12	1.89	0.54
1:A:500:VAL:HB	1:A:552:VAL:HG22	1.91	0.53
1:D:231:TRP:CD2	1:D:413:ILE:HD12	2.43	0.53
1:A:179:VAL:HG22	1:A:300:ILE:CD1	2.37	0.53
1:B:146:TYR:CD1	1:D:428:LEU:HD11	2.43	0.53
1:C:359:LEU:HD11	1:C:518:VAL:HG11	1.89	0.53
1:B:515:ILE:O	1:B:518:VAL:HG12	2.08	0.53
1:D:353:ASP:OD1	1:D:354:LYS:N	2.41	0.53
1:C:385:MET:HG2	1:C:454:PHE:CE2	2.43	0.53
1:D:232:THR:OG1	1:D:235:GLN:HG3	2.09	0.53
1:B:226:ARG:NE	1:B:229:VAL:HG21	2.25	0.52
1:A:456:TYR:OH	1:A:459:GLU:HB2	2.10	0.52
1:B:212:PRO:HD2	1:B:217:PHE:CD1	2.45	0.52
1:D:320:CYS:SG	1:D:327:ASN:HB2	2.50	0.52
1:D:559:ARG:HG2	1:D:560:LYS:N	2.24	0.52
1:B:392:LYS:HE2	1:B:444:ILE:HD11	1.90	0.52
1:C:275:PRO:HB3	1:C:286:PRO:HB2	1.92	0.52
1:C:566:ARG:NH2	1:C:582:GLN:O	2.44	0.51
1:A:175:ALA:HB1	1:A:199:VAL:CG1	2.38	0.51
1:C:522:CYS:HB2	1:C:530:ILE:HD11	1.91	0.51
1:A:150:LEU:HD12	1:A:160:ALA:HB1	1.92	0.51
1:C:180:HIS:O	1:C:184:GLU:HG2	2.11	0.51
1:A:123:HIS:CE1	1:A:331:TYR:HH	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:ARG:HD2	1:A:420:THR:HA	1.93	0.50
1:D:197:LEU:O	1:D:201:ILE:HG13	2.11	0.50
1:D:359:LEU:HD11	1:D:518:VAL:HG11	1.93	0.50
1:A:226:ARG:NE	1:A:229:VAL:HG21	2.26	0.50
1:C:179:VAL:HG22	1:C:300:ILE:HD13	1.93	0.50
1:C:473:TYR:CE1	1:C:502:VAL:HG21	2.47	0.50
1:D:369:LEU:HD13	1:D:374:TYR:OH	2.12	0.50
1:D:576:ARG:HG2	1:D:578:PHE:CZ	2.45	0.50
1:C:194:ARG:HD2	1:C:258:GLY:O	2.11	0.50
1:B:220:ARG:HG2	1:B:387:THR:HG21	1.94	0.50
1:D:360:TYR:CE1	1:D:515:ILE:HG13	2.47	0.49
1:C:144:LEU:HD22	1:C:164:ARG:CB	2.42	0.49
1:D:232:THR:HB	1:D:234:GLU:OE1	2.13	0.49
1:C:576:ARG:HG2	1:C:578:PHE:CE2	2.47	0.49
1:B:486:LYS:HA	1:B:487:VAL:C	2.38	0.49
1:D:156:VAL:HG11	1:D:376:HIS:CE1	2.48	0.49
1:A:171:VAL:HG21	1:A:206:HIS:CE1	2.48	0.49
1:A:212:PRO:HD2	1:A:217:PHE:CD1	2.47	0.49
1:C:500:VAL:HG22	1:C:552:VAL:HG22	1.94	0.49
1:C:392:LYS:HD2	1:C:444:ILE:HD11	1.95	0.48
1:A:232:THR:HB	1:A:234:GLU:OE1	2.13	0.48
1:B:428:LEU:HD22	1:B:432:TYR:CZ	2.47	0.48
1:A:178:LEU:HD23	1:A:300:ILE:HG12	1.94	0.48
1:A:505:MET:HE1	1:A:547:GLU:CB	2.41	0.48
1:B:156:VAL:HG11	1:B:376:HIS:CE1	2.48	0.48
1:B:126:ILE:HG23	1:B:173:TYR:CD1	2.48	0.48
1:A:175:ALA:O	1:A:179:VAL:HG23	2.13	0.48
1:B:444:ILE:O	1:B:447[B]:GLN:HB2	2.14	0.48
1:A:330:ASP:OD2	1:A:333:ARG:HB2	2.14	0.48
1:B:226:ARG:HE	1:B:229:VAL:HG21	1.78	0.48
1:B:563:TYR:O	1:B:567:GLN:HG2	2.13	0.48
1:C:120:ASP:OD1	1:C:318:ARG:NH2	2.47	0.48
1:C:442:ARG:O	1:C:446:LYS:HG3	2.13	0.48
1:C:169:LEU:HD23	1:C:204:LEU:HD11	1.96	0.48
1:D:425:ASN:ND2	1:D:429:GLU:OE2	2.47	0.48
1:A:562:LEU:HA	1:A:562:LEU:HD23	1.60	0.47
1:A:566:ARG:NH2	1:A:582:GLN:O	2.46	0.47
1:C:385:MET:HG2	1:C:454:PHE:CD2	2.49	0.47
1:D:223:PRO:HG3	1:D:470:ARG:HG2	1.97	0.47
1:D:223:PRO:CG	1:D:470:ARG:HG2	2.45	0.47
1:A:228:GLU:H	1:A:228:GLU:HG3	1.21	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:359:LEU:HD13	1:B:518:VAL:HG11	1.95	0.47
1:C:462:PRO:HA	1:C:578:PHE:CD1	2.49	0.47
1:D:443:GLU:O	1:D:447[A]:GLN:HG2	2.15	0.47
1:D:150:LEU:HD12	1:D:160:ALA:HB1	1.95	0.47
1:D:143:ARG:HD2	1:D:420:THR:HA	1.96	0.47
1:C:470:ARG:HG3	1:C:473:TYR:CE2	2.50	0.47
1:A:259:LEU:O	1:A:261:PRO:HD3	2.15	0.46
1:C:489:LEU:HD13	1:C:564:ALA:HA	1.96	0.46
1:A:131:LEU:HD23	1:A:197:LEU:HD13	1.97	0.46
1:C:489:LEU:HD22	1:C:563:TYR:HE2	1.80	0.46
1:B:183:GLY:HA2	1:B:191:ILE:HD12	1.97	0.46
1:C:144:LEU:HD22	1:C:164:ARG:HB2	1.98	0.46
1:D:385:MET:HG2	1:D:454:PHE:CD2	2.50	0.46
1:B:501:ASP:HB2	1:B:553:TYR:CE2	2.50	0.46
1:D:470:ARG:HG3	1:D:473:TYR:CE2	2.50	0.46
1:D:231:TRP:CE3	1:D:413:ILE:HD12	2.49	0.46
1:A:167:HIS:O	1:A:171:VAL:HG23	2.15	0.46
1:B:393:ALA:HB2	1:B:441:ALA:CB	2.46	0.46
1:B:562:LEU:HD23	1:B:562:LEU:HA	1.77	0.46
1:A:356:VAL:HG23	1:A:515:ILE:HD11	1.98	0.46
1:C:143:ARG:HD2	1:C:420:THR:HA	1.97	0.46
1:D:462:PRO:HB3	1:D:578:PHE:CE1	2.50	0.46
1:A:333:ARG:NH2	1:A:358:ASN:HB2	2.31	0.46
1:C:308:ILE:HG12	1:C:362:MET:HE1	1.98	0.46
1:B:431:LEU:HD13	1:B:445:LEU:HB3	1.98	0.45
1:D:462:PRO:HA	1:D:578:PHE:CD1	2.51	0.45
1:D:515:ILE:HG23	1:D:520:PHE:HE1	1.81	0.45
1:A:523:LYS:O	1:A:526:PRO:HD3	2.16	0.45
1:A:207:ASP:HB2	5:A:2001:HOH:O	2.17	0.45
1:A:315:TYR:HA	1:A:318:ARG:HB3	1.98	0.45
1:A:563:TYR:O	1:A:567:GLN:HG2	2.16	0.45
1:C:426:ILE:O	1:C:430:ILE:HG13	2.17	0.45
1:D:566:ARG:O	1:D:570:VAL:HG23	2.16	0.45
1:A:359:LEU:HA	1:A:359:LEU:HD23	1.61	0.45
1:C:150:LEU:HD12	1:C:160:ALA:HB1	1.99	0.44
1:C:462:PRO:HG3	1:C:550:ILE:HD11	1.98	0.44
1:A:388:ASP:HB3	1:A:444:ILE:HD13	2.00	0.44
1:B:141:PHE:CE2	1:B:204:LEU:HG	2.53	0.44
1:B:179:VAL:HG22	1:B:300:ILE:CD1	2.46	0.44
1:C:115:MET:HE3	1:C:129:HIS:N	2.33	0.44
1:C:353:ASP:OD1	1:C:354:LYS:N	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:VAL:O	3:A:1585:GTP:O2'	2.36	0.44
1:A:115:MET:HE3	1:A:127:GLU:HG2	2.00	0.44
1:C:203:GLY:O	1:C:206:HIS:ND1	2.42	0.44
1:B:178:LEU:HD23	1:B:300:ILE:HG23	1.99	0.44
1:A:179:VAL:CG2	1:A:199:VAL:HG11	2.48	0.44
1:A:576:ARG:HG2	1:A:578:PHE:CE2	2.52	0.44
1:A:462:PRO:HG3	1:A:548:GLN:OE1	2.18	0.43
1:D:480:VAL:HG22	1:D:572:TRP:CD2	2.52	0.43
1:C:259:LEU:O	1:C:261:PRO:HD3	2.19	0.43
1:C:223:PRO:CG	1:C:470:ARG:HG2	2.48	0.43
1:D:120:ASP:OD1	1:D:318:ARG:NH2	2.51	0.43
1:D:459:GLU:OE2	1:D:549:LEU:HD13	2.18	0.43
1:A:197:LEU:O	1:A:201:ILE:HG13	2.19	0.43
1:B:330:ASP:OD2	1:B:333:ARG:HB2	2.18	0.43
1:B:451:ARG:HD2	1:B:451:ARG:HA	1.88	0.43
1:A:397:ILE:HD11	1:A:430:ILE:HG12	2.00	0.43
1:C:132:LEU:HD23	1:C:132:LEU:HA	1.70	0.43
1:A:360:TYR:CD1	1:A:360:TYR:C	2.96	0.43
1:C:174:LEU:HB3	1:C:310:VAL:HB	2.00	0.43
1:A:468:ILE:HG21	1:A:476:LEU:HD11	2.01	0.43
1:B:369:LEU:HD23	1:B:369:LEU:HA	1.93	0.43
1:C:566:ARG:O	1:C:570:VAL:HG23	2.19	0.43
1:A:352:ARG:HB2	1:A:521:TYR:CZ	2.54	0.42
1:A:480:VAL:HG22	1:A:572:TRP:CG	2.54	0.42
1:B:146:TYR:C	1:B:147:ILE:HG13	2.44	0.42
1:D:197:LEU:HD23	1:D:197:LEU:HA	1.73	0.42
1:D:492:LYS:HA	1:D:568:TYR:OH	2.19	0.42
1:C:427:PHE:CE2	1:C:445:LEU:HD22	2.54	0.42
1:A:139:PRO:HD3	1:C:450:TYR:CE1	2.54	0.42
1:D:385:MET:HE3	1:D:454:PHE:CE1	2.54	0.42
1:B:207:ASP:HB2	5:B:2001:HOH:O	2.18	0.42
1:A:369:LEU:O	1:A:373:ALA:HB3	2.20	0.42
1:B:259:LEU:O	1:B:261:PRO:HD3	2.19	0.42
1:C:197:LEU:HD23	1:C:197:LEU:HA	1.74	0.42
1:A:132:LEU:HD22	1:A:204:LEU:HD22	2.02	0.42
1:C:483:ALA:O	1:C:485:PRO:HD3	2.18	0.42
1:B:417:GLU:H	1:B:417:GLU:HG2	1.45	0.42
1:A:224:LEU:HD23	1:A:224:LEU:HA	1.82	0.42
1:B:388:ASP:O	1:B:392:LYS:HG3	2.19	0.42
1:D:312:LYS:HA	1:D:315:TYR:CE2	2.55	0.42
1:A:443:GLU:O	1:A:447[B]:GLN:HG3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:459:GLU:OE2	1:B:549:LEU:HD13	2.20	0.42
1:A:442[B]:ARG:CD	1:A:446:LYS:HE3	2.43	0.41
1:B:120:ASP:OD1	1:B:318:ARG:NH2	2.53	0.41
1:B:242:GLU:OE1	1:B:269:LYS:NZ	2.35	0.41
1:B:303:ASN:ND2	1:B:306:ASN:OD1	2.53	0.41
1:C:312:LYS:HD3	1:C:315:TYR:OH	2.20	0.41
1:B:116:LYS:NZ	3:D:1585:GTP:HN22	2.17	0.41
1:B:553:TYR:OH	5:B:2002:HOH:O	2.20	0.41
1:A:285:TRP:HA	1:A:286:PRO:HD3	1.77	0.41
1:D:275:PRO:CB	1:D:286:PRO:HB2	2.50	0.41
1:A:120:ASP:OD1	1:A:318:ARG:NH2	2.53	0.41
1:B:226:ARG:NH1	1:B:411:THR:HA	2.35	0.41
1:B:359:LEU:HA	1:B:359:LEU:HD23	1.79	0.41
1:B:143:ARG:HD2	1:B:420:THR:HA	2.02	0.41
1:D:456:TYR:OH	1:D:459:GLU:HB2	2.21	0.41
1:D:505:MET:HG3	1:D:547:GLU:HB2	2.03	0.41
1:A:226:ARG:HE	1:A:229:VAL:HG21	1.86	0.41
1:B:175:ALA:O	1:B:179:VAL:HG23	2.20	0.41
1:B:275:PRO:HD3	1:B:287:TYR:CE1	2.56	0.41
1:D:522:CYS:SG	1:D:523:LYS:N	2.94	0.41
1:A:388:ASP:O	1:A:392:LYS:HG3	2.21	0.41
1:A:126:ILE:HG23	1:A:173:TYR:HB2	2.02	0.40
1:A:453:LEU:O	1:A:455:LYS:NZ	2.52	0.40
1:B:333:ARG:NH2	1:B:358:ASN:HB2	2.36	0.40
1:C:359:LEU:HD23	1:C:359:LEU:HA	1.87	0.40
1:A:438:LEU:O	1:A:442[B]:ARG:HB2	2.21	0.40
1:D:515:ILE:HG23	1:D:520:PHE:CE1	2.56	0.40
1:C:222:ILE:HB	1:C:223:PRO:HD3	2.03	0.40
1:D:486:LYS:HA	1:D:487:VAL:C	2.46	0.40
1:A:480:VAL:HG22	1:A:572:TRP:CD2	2.56	0.40
1:C:571:GLN:O	1:C:574:ALA:HB3	2.22	0.40
1:D:204:LEU:HD12	1:D:204:LEU:HA	1.91	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	434/538 (81%)	422 (97%)	12 (3%)	0	100	100
1	B	431/538 (80%)	424 (98%)	6 (1%)	1 (0%)	43	76
1	C	432/538 (80%)	423 (98%)	8 (2%)	1 (0%)	43	76
1	D	432/538 (80%)	423 (98%)	9 (2%)	0	100	100
All	All	1729/2152 (80%)	1692 (98%)	35 (2%)	2 (0%)	48	80

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	485	PRO
1	B	485	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	355/478 (74%)	344 (97%)	11 (3%)	35	68
1	B	352/478 (74%)	344 (98%)	8 (2%)	44	74
1	C	347/478 (73%)	336 (97%)	11 (3%)	34	67
1	D	359/478 (75%)	344 (96%)	15 (4%)	26	61
All	All	1413/1912 (74%)	1368 (97%)	45 (3%)	34	67

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	117	VAL
1	A	228	GLU
1	A	326	GLN
1	A	349	ILE

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Mol	Chain	Res	Type
1	A	387	THR
1	A	470	ARG
1	A	491	VAL
1	A	500	VAL
1	A	505	MET
1	A	507	TYR
1	A	547	GLU
1	B	117	VAL
1	B	288	LYS
1	B	348	ARG
1	B	399	ILE
1	B	417	GLU
1	B	426	ILE
1	B	445	LEU
1	B	470	ARG
1	C	115	MET
1	C	117	VAL
1	C	180	HIS
1	C	260	ILE
1	C	276	LEU
1	C	326	GLN
1	C	387	THR
1	C	399	ILE
1	C	443	GLU
1	C	470	ARG
1	C	474	GLU
1	D	117	VAL
1	D	210	HIS
1	D	284	LEU
1	D	348	ARG
1	D	387	THR
1	D	397	ILE
1	D	439	LYS
1	D	445	LEU
1	D	461	GLN
1	D	463	THR
1	D	470	ARG
1	D	471	GLU
1	D	474	GLU
1	D	505	MET
1	D	559	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	200	GLN
1	A	243	HIS
1	A	256	GLN
1	A	425	ASN
1	A	504	ASN
1	B	163	ASN
1	B	326	GLN
1	B	327	ASN
1	C	149	GLN
1	C	210	HIS
1	C	248	ASN
1	C	327	ASN
1	C	328	ASN
1	C	364	HIS
1	C	425	ASN
1	D	200	GLN
1	D	303	ASN
1	D	327	ASN
1	D	380	ASN
1	D	425	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 4 are monoatomic - leaving 15 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
4	SO4	B	1587	-	4,4,4	0.22	0	6,6,6	0.11	0
4	SO4	A	1588	2	4,4,4	0.29	0	6,6,6	0.24	0
4	SO4	D	1587	-	4,4,4	0.24	0	6,6,6	0.11	0
3	GTP	C	1585	-	33,34,34	0.97	3 (9%)	50,54,54	1.86	9 (18%)
4	SO4	A	1586	-	4,4,4	0.25	0	6,6,6	0.21	0
3	GTP	A	1585	-	33,34,34	1.01	2 (6%)	50,54,54	1.90	7 (14%)
4	SO4	D	1588	2	4,4,4	0.31	0	6,6,6	0.23	0
3	GTP	D	1585	-	33,34,34	1.00	2 (6%)	50,54,54	1.77	9 (18%)
4	SO4	C	1587	2	4,4,4	0.36	0	6,6,6	0.28	0
3	GTP	D	1586	-	33,34,34	0.87	0	50,54,54	1.71	9 (18%)
4	SO4	B	1586	-	4,4,4	0.23	0	6,6,6	0.14	0
4	SO4	C	1586	-	4,4,4	0.22	0	6,6,6	0.10	0
4	SO4	B	1588	2	4,4,4	0.33	0	6,6,6	0.14	0
4	SO4	B	1585	-	4,4,4	0.25	0	6,6,6	0.14	0
4	SO4	A	1587	-	4,4,4	0.23	0	6,6,6	0.15	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GTP	D	1586	-	-	5/22/38/38	0/3/3/3
3	GTP	D	1585	-	-	8/22/38/38	0/3/3/3
3	GTP	C	1585	-	-	9/22/38/38	0/3/3/3
3	GTP	A	1585	-	-	6/22/38/38	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1585	GTP	PB-O3B	2.65	1.62	1.59
3	D	1585	GTP	PB-O3B	2.63	1.62	1.59
3	D	1585	GTP	C2-N3	2.47	1.39	1.33
3	A	1585	GTP	C2-N3	2.21	1.38	1.33
3	C	1585	GTP	PB-O3A	2.12	1.61	1.59
3	C	1585	GTP	PA-O3A	2.05	1.61	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1585	GTP	PA-O3A	2.05	1.61	1.59

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1585	GTP	C5-C4-N3	-6.42	118.17	128.39
3	C	1585	GTP	C5-C4-N3	-6.09	118.69	128.39
3	C	1585	GTP	C2-N3-C4	5.93	122.51	112.30
3	D	1585	GTP	C5-C4-N3	-5.86	119.06	128.39
3	D	1586	GTP	C5-C4-N3	-5.75	119.24	128.39
3	A	1585	GTP	C2-N3-C4	5.63	122.00	112.30
3	D	1586	GTP	C2-N3-C4	5.29	121.41	112.30
3	D	1585	GTP	C2-N3-C4	5.22	121.29	112.30
3	A	1585	GTP	N9-C4-N3	4.55	135.06	125.95
3	D	1585	GTP	N9-C4-N3	3.74	133.44	125.95
3	D	1586	GTP	N9-C4-N3	3.69	133.34	125.95
3	C	1585	GTP	C2-N1-C6	-3.53	118.70	125.11
3	A	1585	GTP	C2-N1-C6	-3.35	119.03	125.11
3	C	1585	GTP	N9-C4-N3	3.31	132.58	125.95
3	A	1585	GTP	C5-C6-N1	3.18	121.36	113.25
3	D	1585	GTP	C2-N1-C6	-3.03	119.62	125.11
3	C	1585	GTP	C8-N7-C5	3.00	109.61	104.26
3	D	1585	GTP	C8-N7-C5	2.94	109.49	104.26
3	C	1585	GTP	N2-C2-N1	2.92	122.92	116.76
3	D	1586	GTP	C2-N1-C6	-2.90	119.85	125.11
3	C	1585	GTP	C5-C6-N1	2.76	120.27	113.25
3	D	1585	GTP	C3'-C2'-C1'	2.75	106.67	101.46
3	C	1585	GTP	C4-C5-N7	-2.69	106.41	110.67
3	A	1585	GTP	C8-N7-C5	2.64	108.96	104.26
3	D	1585	GTP	C5-C6-N1	2.64	119.96	113.25
3	C	1585	GTP	N9-C8-N7	-2.46	108.83	113.40
3	D	1586	GTP	C5-C6-N1	2.42	119.42	113.25
3	A	1585	GTP	N9-C8-N7	-2.39	108.97	113.40
3	D	1586	GTP	O6-C6-C5	-2.36	120.29	126.53
3	D	1585	GTP	N9-C8-N7	-2.36	109.03	113.40
3	D	1586	GTP	C8-N7-C5	2.30	108.36	104.26
3	D	1585	GTP	C4-C5-N7	-2.21	107.17	110.67
3	D	1586	GTP	N2-C2-N1	2.15	121.30	116.76
3	D	1586	GTP	N9-C8-N7	-2.01	109.67	113.40

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1585	GTP	PB-O3B-PG-O3G
3	A	1585	GTP	C5'-O5'-PA-O3A
3	A	1585	GTP	C5'-O5'-PA-O1A
3	A	1585	GTP	C5'-O5'-PA-O2A
3	C	1585	GTP	C5'-O5'-PA-O3A
3	C	1585	GTP	C5'-O5'-PA-O1A
3	D	1585	GTP	C5'-O5'-PA-O3A
3	D	1585	GTP	C5'-O5'-PA-O1A
3	D	1585	GTP	C5'-O5'-PA-O2A
3	D	1586	GTP	C5'-O5'-PA-O3A
3	D	1586	GTP	C5'-O5'-PA-O1A
3	D	1586	GTP	C5'-O5'-PA-O2A
3	D	1586	GTP	O4'-C4'-C5'-O5'
3	C	1585	GTP	O4'-C4'-C5'-O5'
3	C	1585	GTP	C3'-C4'-C5'-O5'
3	C	1585	GTP	C4'-C5'-O5'-PA
3	C	1585	GTP	C5'-O5'-PA-O2A
3	C	1585	GTP	PG-O3B-PB-O2B
3	D	1585	GTP	PB-O3A-PA-O1A
3	C	1585	GTP	C2'-C1'-N9-C8
3	D	1585	GTP	C2'-C1'-N9-C8
3	A	1585	GTP	PB-O3B-PG-O2G
3	A	1585	GTP	C2'-C1'-N9-C8
3	D	1585	GTP	C2'-C1'-N9-C4
3	C	1585	GTP	PG-O3B-PB-O1B
3	D	1585	GTP	PG-O3B-PB-O2B
3	D	1586	GTP	C2'-C1'-N9-C4
3	D	1585	GTP	PB-O3A-PA-O2A

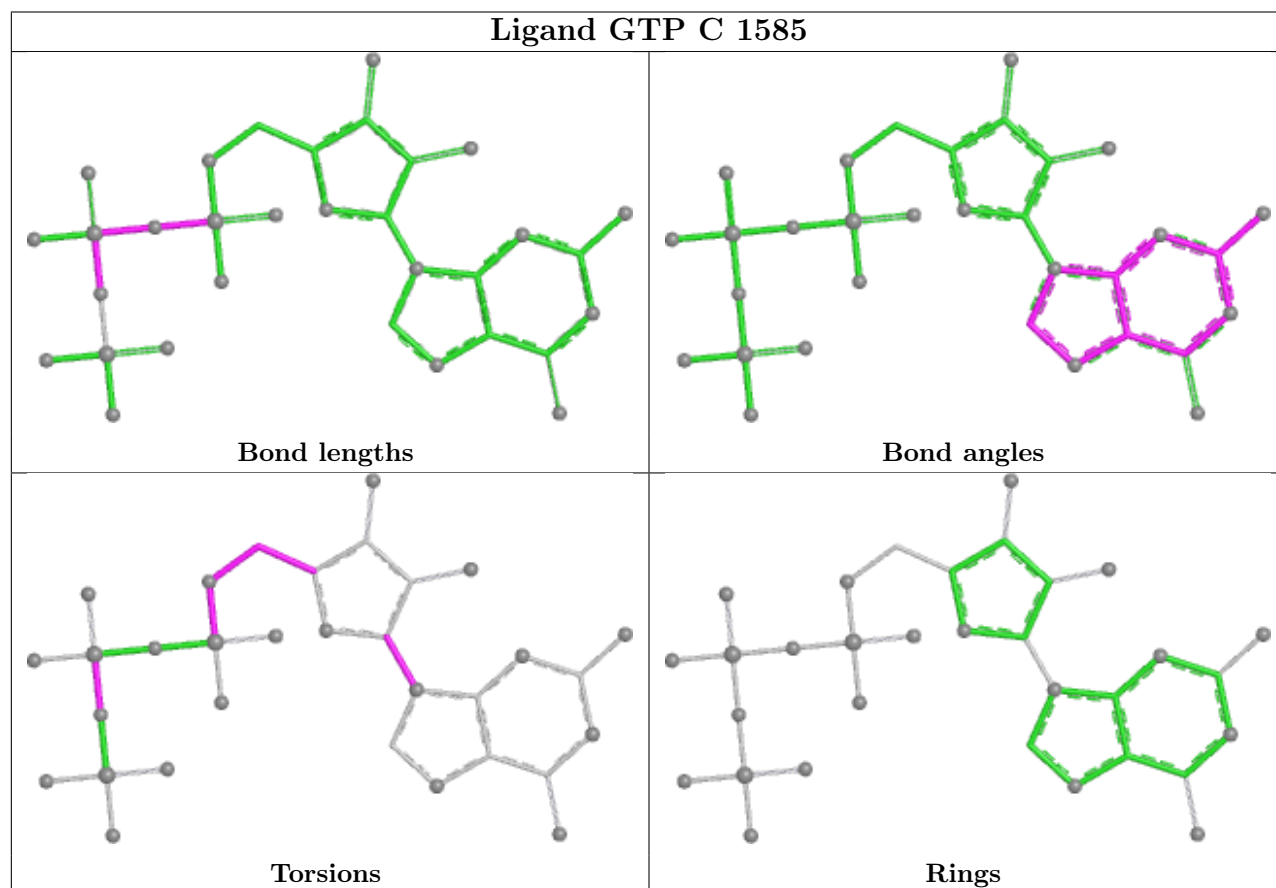
There are no ring outliers.

5 monomers are involved in 9 short contacts:

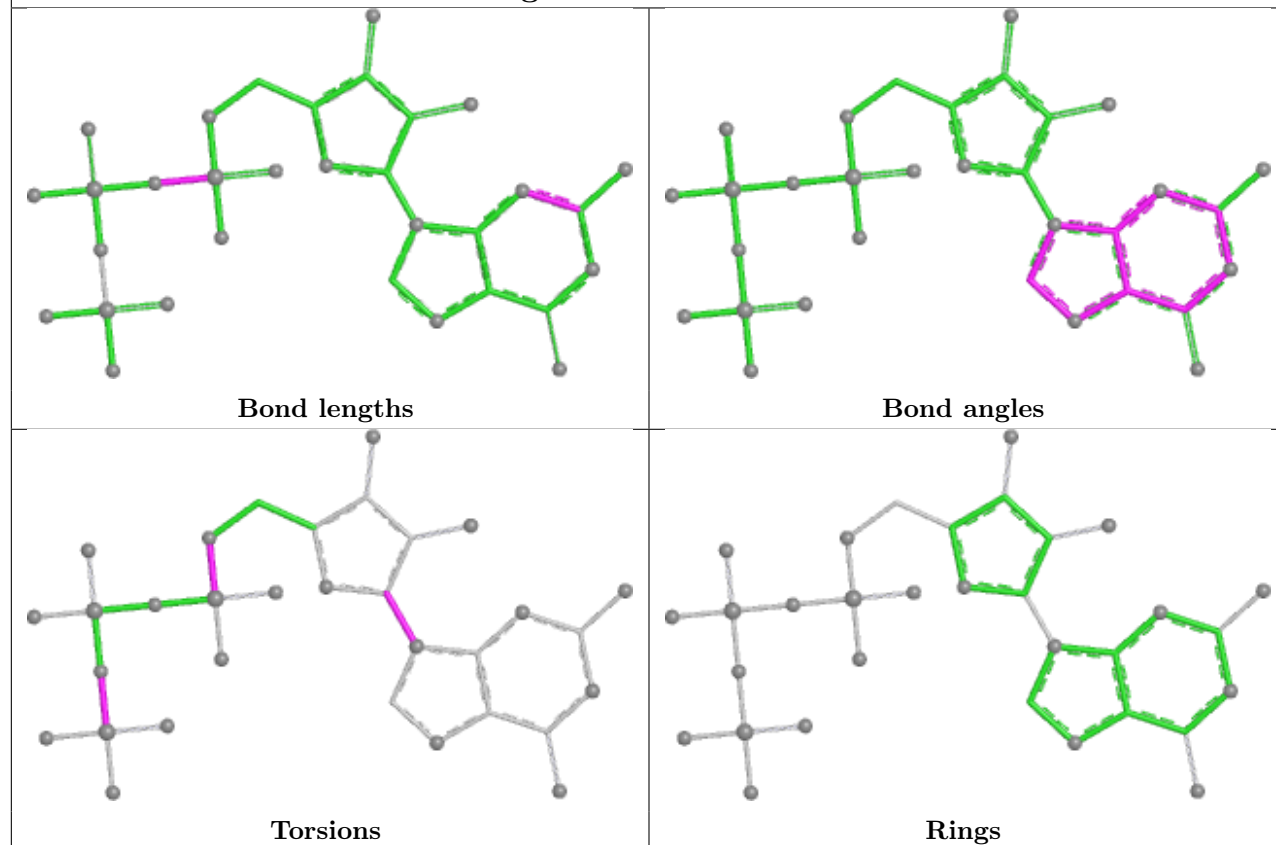
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1585	GTP	1	0
3	A	1585	GTP	4	0
3	D	1585	GTP	2	0
4	B	1586	SO4	1	0
4	B	1585	SO4	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is

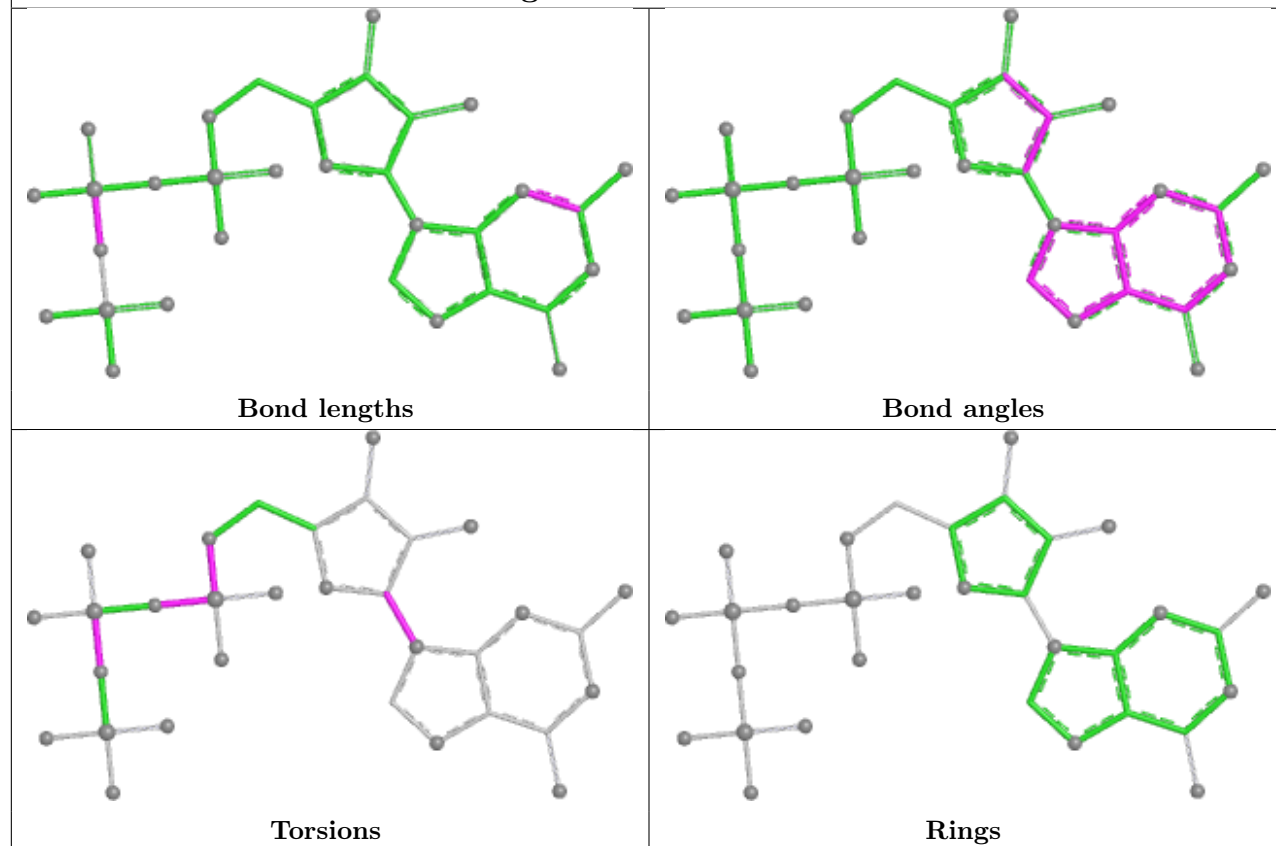
within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

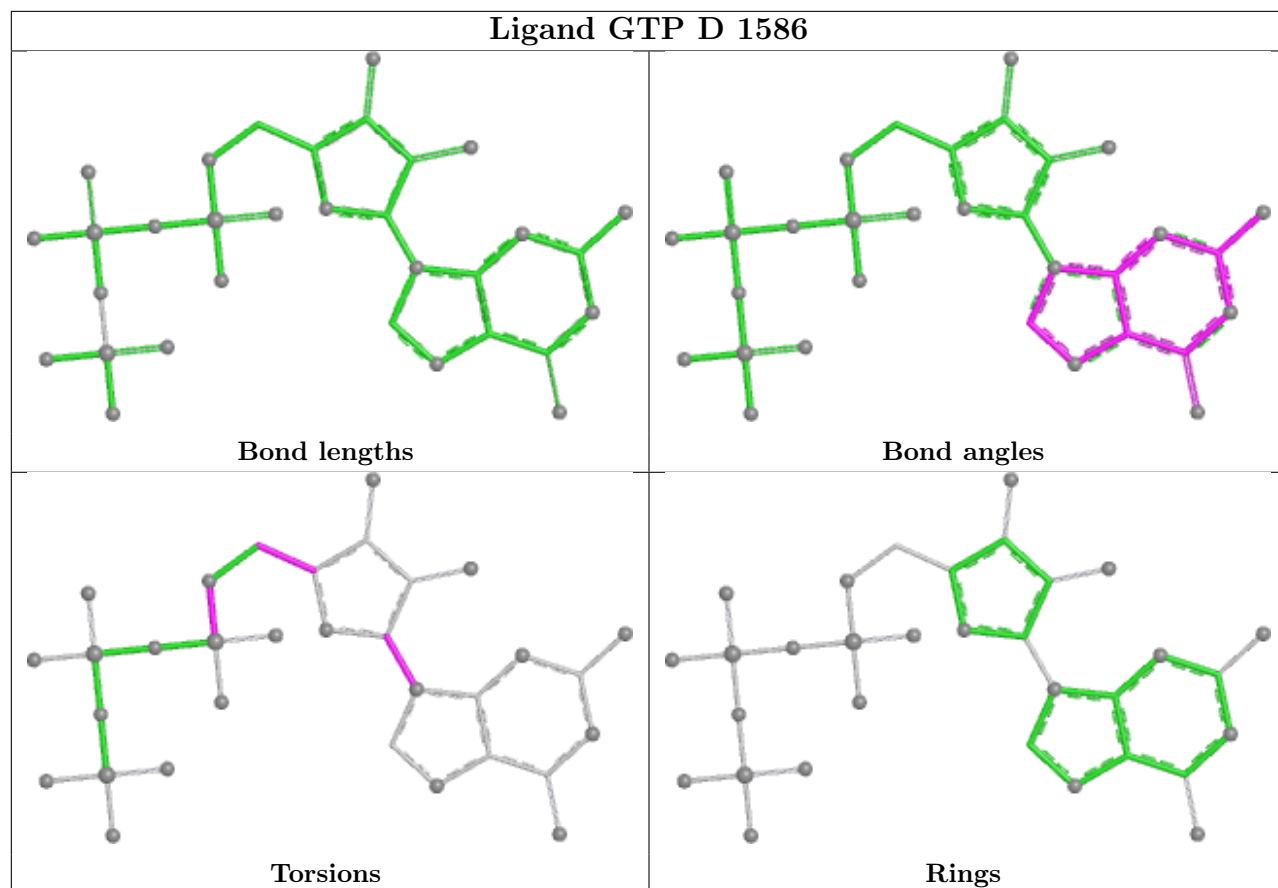


Ligand GTP A 1585



Ligand GTP D 1585





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 [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	440/538 (81%)	-0.37	5 (1%) 78 57	17, 58, 114, 202	2 (0%)
1	B	438/538 (81%)	-0.42	1 (0%) 91 83	23, 60, 109, 177	1 (0%)
1	C	439/538 (81%)	-0.31	2 (0%) 87 72	31, 67, 117, 170	1 (0%)
1	D	439/538 (81%)	-0.31	3 (0%) 84 66	29, 63, 112, 175	1 (0%)
All	All	1756/2152 (81%)	-0.35	11 (0%) 85 69	17, 62, 114, 202	5 (0%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	345	ASN	3.5
1	A	506	ASP	3.4
1	D	114	PRO	2.9
1	D	583	ASP	2.6
1	C	276	LEU	2.6
1	A	507	TYR	2.6
1	A	299	GLU	2.5
1	D	515	ILE	2.2
1	A	449	GLU	2.2
1	B	127	GLU	2.2
1	C	516	ASP	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands

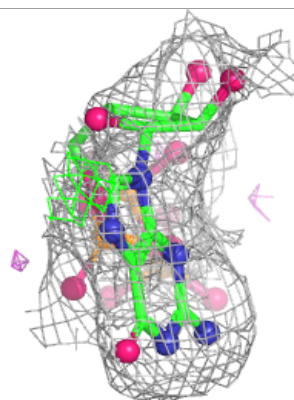
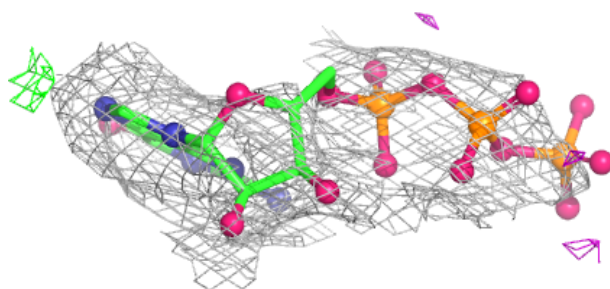
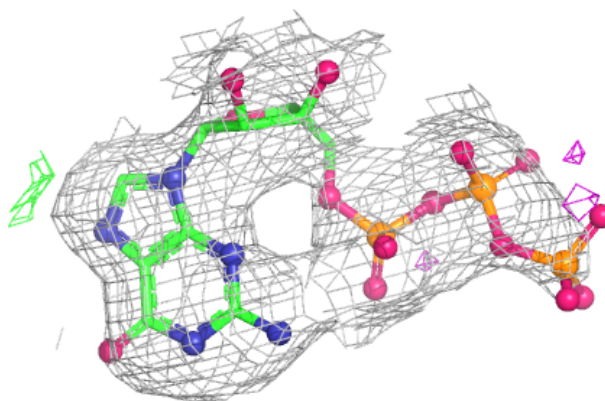
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists 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
4	SO4	A	1587	5/5	0.77	0.12	114,120,121,125	0
4	SO4	B	1586	5/5	0.77	0.21	135,135,139,140	0
3	GTP	D	1585	32/32	0.80	0.12	58,116,197,199	0
4	SO4	D	1587	5/5	0.80	0.15	129,135,137,142	0
4	SO4	B	1587	5/5	0.82	0.24	133,136,139,143	0
4	SO4	C	1586	5/5	0.84	0.19	109,115,116,124	0
3	GTP	A	1585	32/32	0.85	0.11	64,109,163,181	0
4	SO4	A	1586	5/5	0.88	0.15	128,130,135,136	0
4	SO4	C	1587	5/5	0.90	0.11	106,106,106,106	0
4	SO4	B	1588	5/5	0.90	0.12	97,97,97,97	0
4	SO4	D	1588	5/5	0.90	0.13	112,112,112,112	0
3	GTP	C	1585	32/32	0.91	0.10	42,86,155,195	0
3	GTP	D	1586	32/32	0.92	0.09	50,89,183,200	0
4	SO4	B	1585	5/5	0.93	0.09	115,121,124,126	0
4	SO4	A	1588	5/5	0.95	0.11	93,93,93,93	0
2	FE	D	1584	1/1	1.00	0.02	30,30,30,30	0
2	FE	A	1584	1/1	1.00	0.01	30,30,30,30	0
2	FE	B	1584	1/1	1.00	0.02	32,32,32,32	0
2	FE	C	1584	1/1	1.00	0.02	36,36,36,36	0

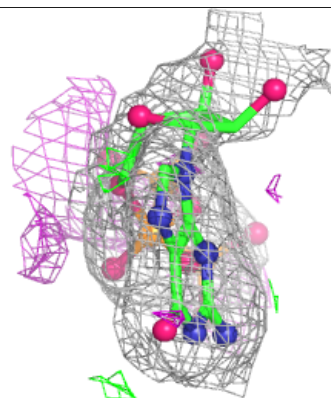
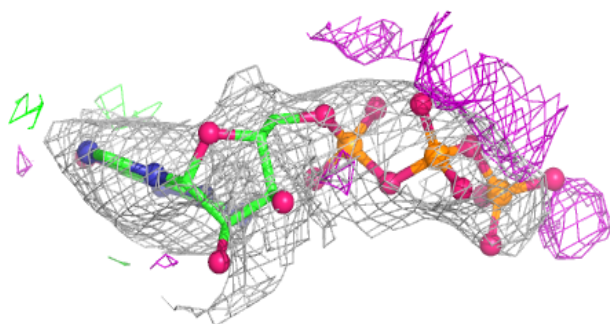
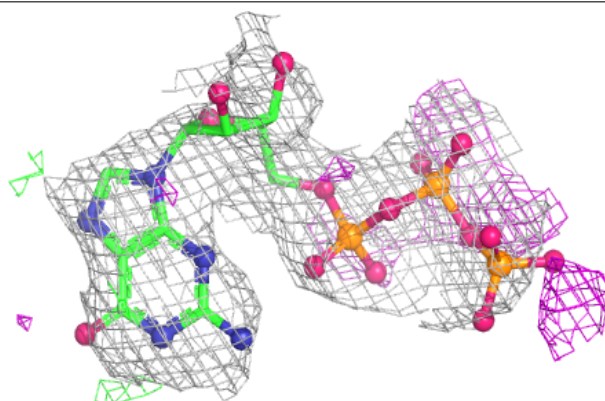
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around GTP D 1585:

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

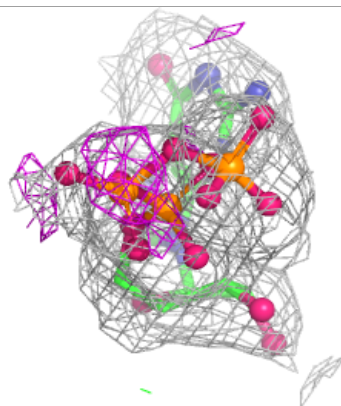
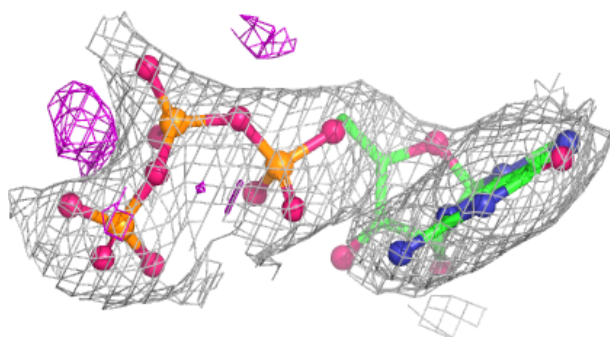
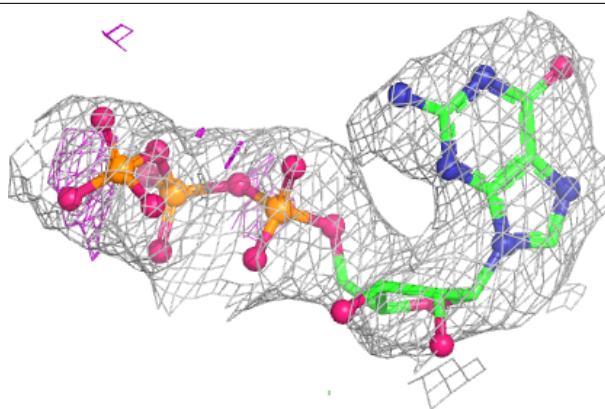
**Electron density around GTP A 1585:**

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

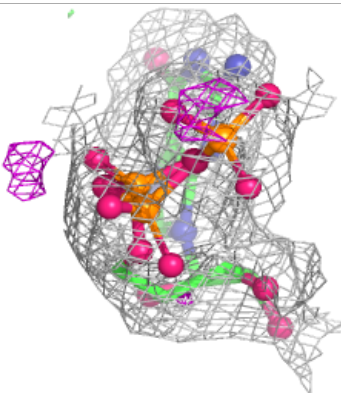
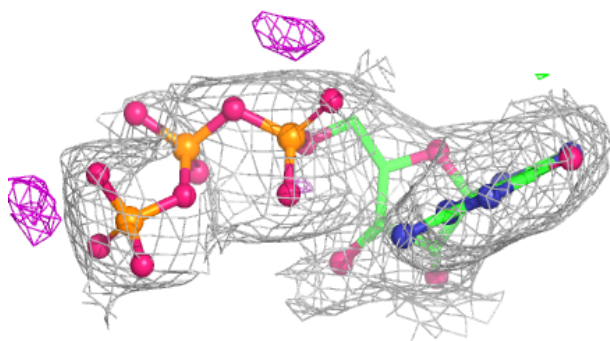
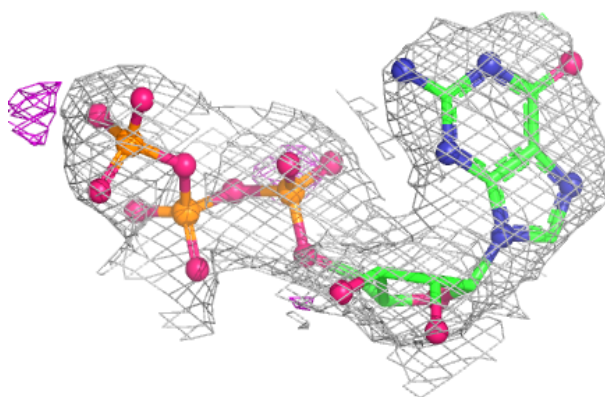


Electron density around GTP C 1585:

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

**Electron density around GTP D 1586:**

$2mF_o-DF_c$ (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.