



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 07:28 PM UTC

PDB ID : 5DLQ / pdb_00005dlq
Title : Crystal structure of RanGTP-Exportin 4-eIF5A complex
Authors : Aksu, M.; Trakhanov, S.; Gorlich, D.
Deposited on : 2015-09-07
Resolution : 3.20 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

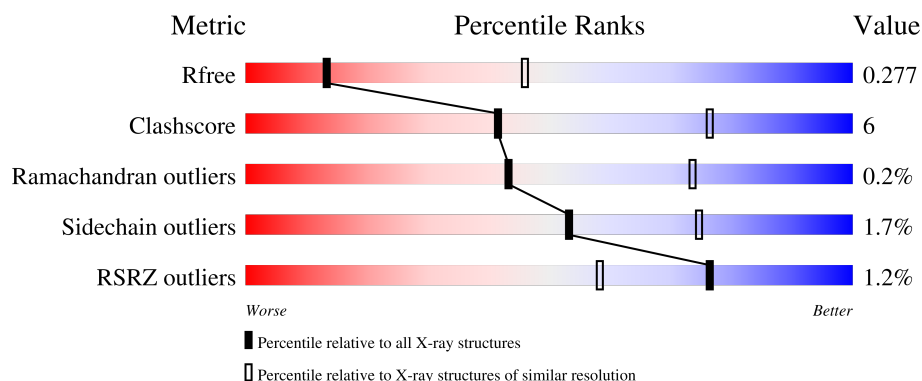
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	1113	
1	B	1113	
2	C	176	
2	D	176	
3	E	141	

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Mol	Chain	Length	Quality of chain
3	F	141	% 80% 16% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 20688 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 Exportin-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	1052	Total	C	N	O	S	0	0	0
			8199	5231	1351	1569	48			
1	A	983	Total	C	N	O	S	0	0	0
			7584	4849	1247	1443	45			

There are 78 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	PRO	deletion	UNP Q9ESJ0
B	?	-	LYS	deletion	UNP Q9ESJ0
B	?	-	LEU	deletion	UNP Q9ESJ0
B	?	-	GLY	deletion	UNP Q9ESJ0
B	?	-	ARG	deletion	UNP Q9ESJ0
B	?	-	HIS	deletion	UNP Q9ESJ0
B	?	-	TYR	deletion	UNP Q9ESJ0
B	?	-	ILE	deletion	UNP Q9ESJ0
B	?	-	ALA	deletion	UNP Q9ESJ0
B	?	-	MET	deletion	UNP Q9ESJ0
B	?	-	PHE	deletion	UNP Q9ESJ0
B	?	-	GLU	deletion	UNP Q9ESJ0
B	?	-	SER	deletion	UNP Q9ESJ0
B	?	-	SER	deletion	UNP Q9ESJ0
B	?	-	GLN	deletion	UNP Q9ESJ0
B	?	-	ASN	deletion	UNP Q9ESJ0
B	?	-	VAL	deletion	UNP Q9ESJ0
B	?	-	LEU	deletion	UNP Q9ESJ0
B	?	-	LEU	deletion	UNP Q9ESJ0
B	?	-	LYS	deletion	UNP Q9ESJ0
B	?	-	SER	deletion	UNP Q9ESJ0
B	?	-	ASP	deletion	UNP Q9ESJ0
B	?	-	THR	deletion	UNP Q9ESJ0
B	?	-	ASP	deletion	UNP Q9ESJ0
B	?	-	GLU	deletion	UNP Q9ESJ0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	VAL	deletion	UNP Q9ESJ0
B	?	-	PHE	deletion	UNP Q9ESJ0
B	?	-	ARG	deletion	UNP Q9ESJ0
B	?	-	GLY	deletion	UNP Q9ESJ0
B	?	-	HIS	deletion	UNP Q9ESJ0
B	?	-	GLU	deletion	UNP Q9ESJ0
B	?	-	PRO	deletion	UNP Q9ESJ0
B	?	-	GLY	deletion	UNP Q9ESJ0
B	?	-	GLN	deletion	UNP Q9ESJ0
B	?	-	ALA	deletion	UNP Q9ESJ0
B	?	-	ALA	deletion	UNP Q9ESJ0
B	?	-	GLY	deletion	UNP Q9ESJ0
B	?	-	ARG	deletion	UNP Q9ESJ0
B	1085	TYR	HIS	engineered mutation	UNP Q9ESJ0
A	?	-	PRO	deletion	UNP Q9ESJ0
A	?	-	LYS	deletion	UNP Q9ESJ0
A	?	-	LEU	deletion	UNP Q9ESJ0
A	?	-	GLY	deletion	UNP Q9ESJ0
A	?	-	ARG	deletion	UNP Q9ESJ0
A	?	-	HIS	deletion	UNP Q9ESJ0
A	?	-	TYR	deletion	UNP Q9ESJ0
A	?	-	ILE	deletion	UNP Q9ESJ0
A	?	-	ALA	deletion	UNP Q9ESJ0
A	?	-	MET	deletion	UNP Q9ESJ0
A	?	-	PHE	deletion	UNP Q9ESJ0
A	?	-	GLU	deletion	UNP Q9ESJ0
A	?	-	SER	deletion	UNP Q9ESJ0
A	?	-	SER	deletion	UNP Q9ESJ0
A	?	-	GLN	deletion	UNP Q9ESJ0
A	?	-	ASN	deletion	UNP Q9ESJ0
A	?	-	VAL	deletion	UNP Q9ESJ0
A	?	-	LEU	deletion	UNP Q9ESJ0
A	?	-	LEU	deletion	UNP Q9ESJ0
A	?	-	LYS	deletion	UNP Q9ESJ0
A	?	-	SER	deletion	UNP Q9ESJ0
A	?	-	ASP	deletion	UNP Q9ESJ0
A	?	-	THR	deletion	UNP Q9ESJ0
A	?	-	ASP	deletion	UNP Q9ESJ0
A	?	-	GLU	deletion	UNP Q9ESJ0
A	?	-	VAL	deletion	UNP Q9ESJ0
A	?	-	PHE	deletion	UNP Q9ESJ0
A	?	-	ARG	deletion	UNP Q9ESJ0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	GLY	deletion	UNP Q9ESJ0
A	?	-	HIS	deletion	UNP Q9ESJ0
A	?	-	GLU	deletion	UNP Q9ESJ0
A	?	-	PRO	deletion	UNP Q9ESJ0
A	?	-	GLY	deletion	UNP Q9ESJ0
A	?	-	GLN	deletion	UNP Q9ESJ0
A	?	-	ALA	deletion	UNP Q9ESJ0
A	?	-	ALA	deletion	UNP Q9ESJ0
A	?	-	GLY	deletion	UNP Q9ESJ0
A	?	-	ARG	deletion	UNP Q9ESJ0
A	1085	TYR	HIS	engineered mutation	UNP Q9ESJ0

- Molecule 2 is a protein called GTP-binding nuclear protein Ran.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	173	Total	C	N	O	S	0	0	0
			1368	890	238	236	4			
2	D	172	Total	C	N	O	S	0	0	0
			1373	894	236	239	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	69	LEU	GLN	engineered mutation	UNP P62826
D	69	LEU	GLN	engineered mutation	UNP P62826

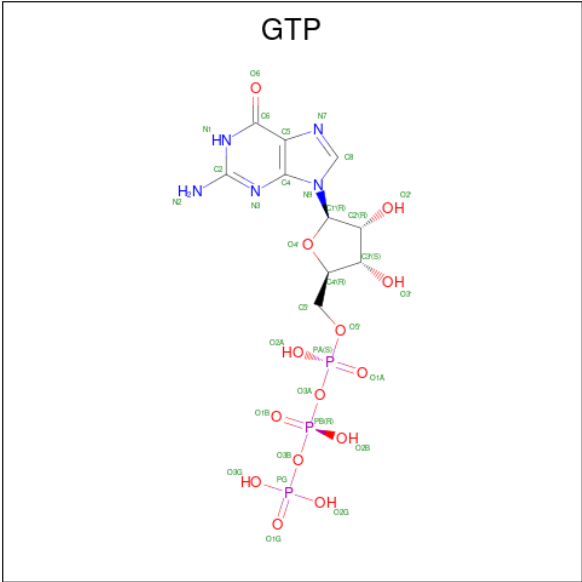
- Molecule 3 is a protein called Eukaryotic translation initiation factor 5A-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	138	Total	C	N	O	S	0	0	0
			1058	663	180	206	9			
3	E	137	Total	C	N	O	S	0	0	0
			1040	651	177	203	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	14	GLY	-	expression tag	UNP P63241
E	14	GLY	-	expression tag	UNP P63241

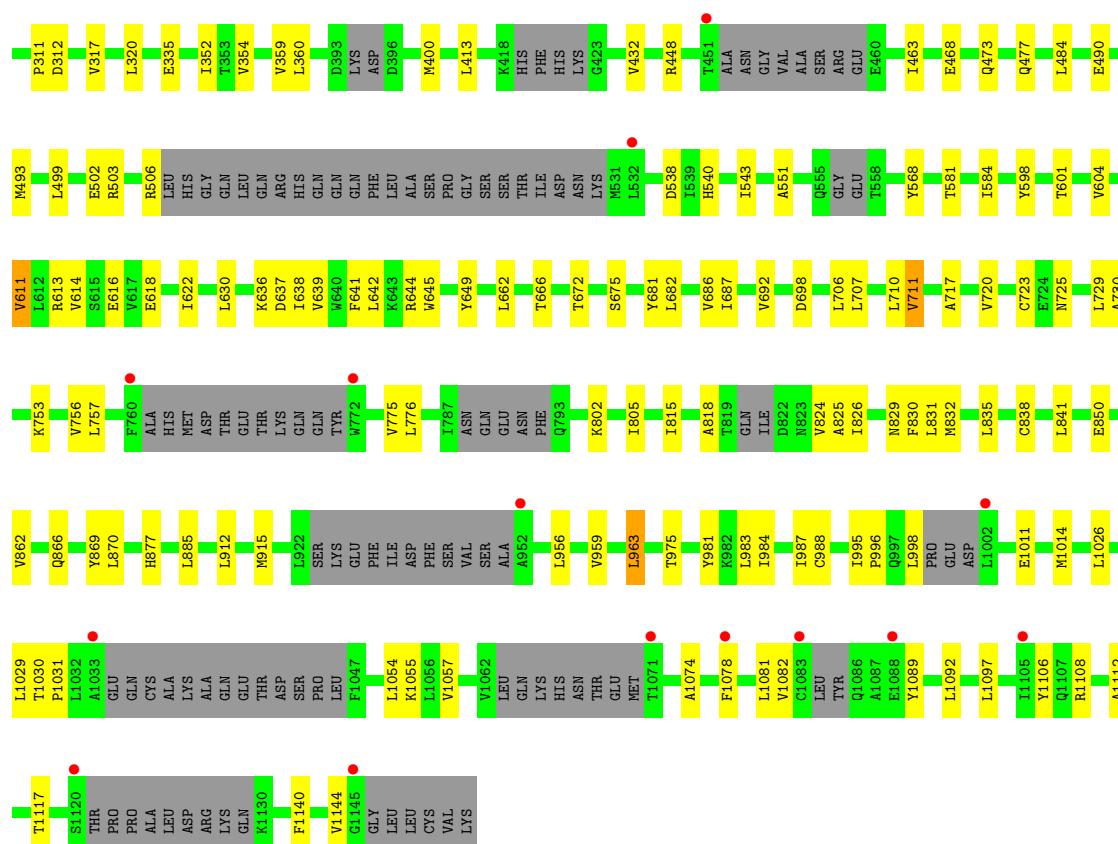
- Molecule 4 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃).



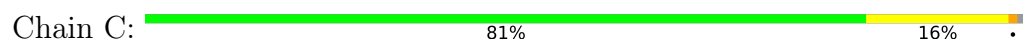
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	C	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
4	D	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

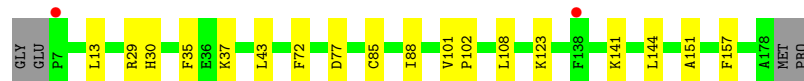
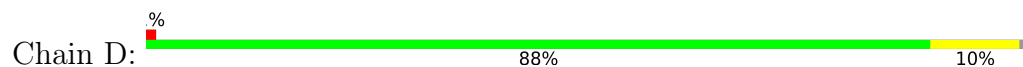
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	1	Total	Mg	0	0
			1	1		
5	D	1	Total	Mg	0	0
			1	1		



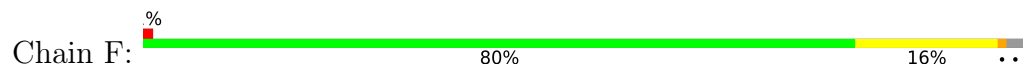
- Molecule 2: GTP-binding nuclear protein Ran



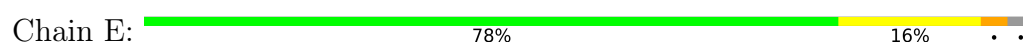
- Molecule 2: GTP-binding nuclear protein Ran



- Molecule 3: Eukaryotic translation initiation factor 5A-1



- Molecule 3: Eukaryotic translation initiation factor 5A-1



GLY	SER	A16	M20	E42	K50	K55	V59	I63	P74	S75	T76	H77	N83	N87	Q85	D96	G97	Y98	L101	R109	L112	R113	L114	P115	L134	I135	T136	W137	L138	E144	A148	I149	K150	A151	M152	ALA	LYS
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4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	98.62Å 98.62Å 726.86Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.41 – 3.20 49.41 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.41-3.20) 100.0 (49.41-3.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 3.19Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.221 , 0.270 0.232 , 0.277	Depositor DCC
R_{free} test set	3495 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	106.3	Xtriage
Anisotropy	0.409	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 79.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.038 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	20688	wwPDB-VP
Average B, all atoms (Å ²)	118.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.24% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GTP, 5CT, 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	0.11	0/7708	0.28	0/10473
1	B	0.11	0/8341	0.31	0/11336
2	C	0.11	0/1403	0.32	0/1904
2	D	0.10	0/1408	0.28	0/1908
3	E	0.09	0/1038	0.31	0/1398
3	F	0.11	0/1057	0.36	0/1423
All	All	0.11	0/20955	0.30	0/28442

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	7584	0	7444	110	0
1	B	8199	0	8095	100	0
2	C	1368	0	1353	18	0
2	D	1373	0	1369	13	0
3	E	1040	0	1036	14	0
3	F	1058	0	1057	14	0
4	C	32	0	12	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	32	0	12	1	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
All	All	20688	0	20378	262	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 6.

All (262) 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:B:1079:TYR:HH	1:B:1126:ASP:N	1.77	0.83
3:E:136:THR:HB	3:E:148:ALA:HB3	1.63	0.79
3:F:136:THR:HB	3:F:148:ALA:HB3	1.65	0.76
3:F:20:MET:HE2	3:F:81:VAL:HG21	1.71	0.73
1:B:91:GLU:OE2	1:B:94:ARG:NH1	2.22	0.72
1:B:210:SER:O	1:B:284:LYS:NZ	2.22	0.72
1:B:135:LYS:O	1:B:139:HIS:ND1	2.22	0.70
1:B:622:ILE:HD12	1:B:692:VAL:HG21	1.74	0.69
1:A:622:ILE:HD12	1:A:692:VAL:HG21	1.75	0.67
1:A:400:MET:HE3	1:A:477:GLN:HE22	1.58	0.67
3:F:37:PRO:HD3	3:F:141:MET:HE2	1.78	0.64
1:A:672:THR:HG23	1:A:675:SER:H	1.62	0.64
1:B:645:TRP:CD1	1:B:649:TYR:HB2	2.33	0.63
1:B:959:VAL:O	1:B:963:LEU:HB2	1.98	0.63
1:B:820:GLN:H	1:B:823:ASN:HB2	1.64	0.63
1:A:210:SER:O	1:A:284:LYS:NZ	2.32	0.62
1:A:285:ILE:HD12	1:A:291:MET:HE2	1.82	0.62
3:E:114:LEU:HD23	3:E:115:PRO:HD2	1.83	0.61
1:A:191:GLN:NE2	1:A:238:PHE:O	2.33	0.61
1:B:911:LEU:HD13	1:B:966:MET:HE2	1.83	0.60
1:B:553:ASP:O	1:B:555:GLN:N	2.33	0.60
1:B:721:ILE:HD13	1:B:761:ALA:HB3	1.83	0.60
1:B:988:CYS:HA	1:B:995:ILE:HD11	1.82	0.59
1:B:391:VAL:HG11	1:B:449:ASN:HA	1.85	0.59
1:B:352:ILE:HD13	1:B:413:LEU:HD12	1.85	0.58
1:A:645:TRP:CD1	1:A:649:TYR:HB2	2.37	0.58
1:B:636:LYS:HD3	1:B:698:ASP:HB3	1.85	0.58
1:B:825:ALA:O	1:B:829:ASN:ND2	2.37	0.58
1:B:824:VAL:HG11	1:B:869:TYR:HB2	1.86	0.58
1:A:956:LEU:HD13	1:A:998:LEU:HD21	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:LYS:HD2	1:A:138:PHE:CE2	2.38	0.57
1:A:1112:ALA:HB1	1:A:1140:PHE:HD1	1.68	0.57
1:A:581:THR:HA	1:A:601:THR:HG23	1.86	0.56
1:B:306:HIS:HB3	1:B:354:VAL:HG13	1.87	0.56
1:A:83:VAL:O	1:A:1108:ARG:NH1	2.38	0.56
1:A:1026:LEU:HD13	1:A:1074:ALA:HB2	1.88	0.56
1:A:835:LEU:HD21	1:A:862:VAL:HG21	1.86	0.56
1:A:956:LEU:HD21	1:A:987:ILE:HG21	1.88	0.56
1:A:618:GLU:HG3	1:A:630:LEU:HD13	1.87	0.56
1:B:320:LEU:HD21	1:B:360:LEU:HD23	1.88	0.55
3:E:55:LYS:HA	3:E:74:PRO:HA	1.87	0.55
1:B:83:VAL:O	1:B:1108:ARG:NH1	2.39	0.55
1:A:584:ILE:HG23	1:A:613:ARG:HH22	1.71	0.55
1:A:45:ARG:O	1:A:80:ARG:NH1	2.40	0.55
1:B:639:VAL:HG13	1:B:706:LEU:HD12	1.89	0.54
3:F:20:MET:HE3	3:F:25:LEU:HD23	1.88	0.54
1:A:832:MET:HE2	1:A:877:HIS:HB3	1.90	0.54
1:A:753:LYS:HA	1:A:818:ALA:HB2	1.89	0.54
1:B:116:LEU:HD22	1:B:159:ILE:HG12	1.90	0.54
1:B:707:LEU:O	1:B:711:VAL:HG23	2.07	0.54
1:B:672:THR:HG23	1:B:675:SER:H	1.74	0.53
2:D:151:ALA:HA	2:D:157:PHE:CE2	2.44	0.53
1:A:320:LEU:HD21	1:A:360:LEU:HD23	1.91	0.53
1:A:468:GLU:HG3	1:A:473:GLN:HB2	1.91	0.53
3:F:42:GLU:HB3	3:F:59:VAL:HG13	1.92	0.52
1:A:448:ARG:NH1	1:A:538:ASP:OD1	2.41	0.52
1:B:831:LEU:HD13	1:B:862:VAL:HG22	1.91	0.52
2:D:43:LEU:HA	2:D:72:PHE:HB3	1.92	0.52
3:F:75:SER:C	3:F:77:HIS:H	2.17	0.52
2:D:123:LYS:HG2	4:D:217:GTP:C6	2.44	0.52
1:B:285:ILE:HD12	1:B:291:MET:HE2	1.92	0.52
1:B:1016:SER:OG	1:B:1017:MET:N	2.42	0.52
1:B:899:ILE:HG12	3:F:119:LEU:HD13	1.92	0.52
3:E:75:SER:C	3:E:77:HIS:H	2.18	0.52
1:A:191:GLN:HG2	1:A:240:PRO:HD3	1.92	0.51
1:B:981:TYR:O	1:B:985:THR:OG1	2.23	0.51
2:C:39:TYR:CD1	4:C:217:GTP:H5"	2.45	0.51
1:B:611:VAL:HG21	1:B:641:PHE:CZ	2.45	0.51
1:B:819:THR:HG21	1:B:866:GLN:HG2	1.93	0.51
3:E:98:TYR:CE1	3:E:113:ARG:HG2	2.46	0.51
1:A:98:LEU:HD21	1:A:119:VAL:HG11	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:29:ARG:HG2	2:D:157:PHE:CE1	2.46	0.51
1:A:137:ILE:H	1:A:137:ILE:HD12	1.74	0.51
1:A:838:CYS:HA	1:A:841:LEU:HB2	1.92	0.51
1:A:12:ILE:HG12	1:A:57:ILE:HG12	1.92	0.51
1:B:346:SER:O	1:B:350:ASN:ND2	2.44	0.50
1:A:825:ALA:O	1:A:829:ASN:ND2	2.44	0.50
1:B:553:ASP:OD1	1:B:644:ARG:NH2	2.42	0.50
1:B:966:MET:HE1	1:B:976:LEU:HD11	1.93	0.50
1:A:187:LYS:O	1:A:191:GLN:HB2	2.11	0.50
2:C:15:LEU:HD13	2:C:23:LYS:HB3	1.92	0.50
3:F:55:LYS:HA	3:F:74:PRO:HA	1.93	0.50
1:A:707:LEU:O	1:A:711:VAL:HG23	2.11	0.50
1:A:725:ASN:O	1:A:729:LEU:HG	2.12	0.50
1:B:847:ASN:O	3:F:86:ARG:HD2	2.11	0.50
3:F:17:THR:HG22	3:F:82:PRO:HA	1.93	0.50
1:A:199:PHE:HB2	1:A:233:VAL:HG11	1.94	0.50
1:B:821:ILE:HG13	1:B:869:TYR:CE1	2.47	0.49
1:A:502:GLU:OE2	1:A:506:ARG:NE	2.45	0.49
1:A:756:VAL:HG21	1:A:815:ILE:HG12	1.94	0.49
1:A:831:LEU:HD13	1:A:862:VAL:HG12	1.94	0.49
1:A:776:LEU:HB3	1:A:830:PHE:CZ	2.47	0.49
1:A:988:CYS:HA	1:A:995:ILE:HD11	1.93	0.49
2:C:35:PHE:CE2	2:C:37:LYS:HG2	2.48	0.49
1:A:611:VAL:HG21	1:A:641:PHE:CE2	2.48	0.49
1:B:611:VAL:HG21	1:B:641:PHE:CE2	2.47	0.49
1:B:753:LYS:O	1:B:757:LEU:HB2	2.13	0.49
3:E:101:LEU:HD12	3:E:112:LEU:HD23	1.94	0.49
1:B:286:ARG:HE	1:B:334:ILE:HG23	1.78	0.49
1:A:717:ALA:HA	1:A:720:VAL:HG12	1.95	0.49
1:A:639:VAL:HG13	1:A:706:LEU:HD12	1.95	0.49
1:A:912:LEU:HD11	1:A:975:THR:HG22	1.95	0.48
1:B:84:LEU:HA	1:B:1108:ARG:HH12	1.77	0.48
1:B:94:ARG:NH2	1:B:140:GLU:OE1	2.45	0.48
1:A:642:LEU:HB2	1:A:706:LEU:HD11	1.95	0.48
3:F:89:PHE:CE1	3:F:103:GLN:HG2	2.49	0.48
1:B:263:GLU:HA	1:B:266:ARG:HB2	1.96	0.48
1:A:824:VAL:C	1:A:826:ILE:H	2.21	0.48
2:D:35:PHE:HE2	2:D:37:LYS:HG2	1.78	0.48
2:C:123:LYS:HG2	4:C:217:GTP:C6	2.49	0.48
1:A:824:VAL:HG11	1:A:869:TYR:HB2	1.95	0.48
2:C:95:ARG:HG2	2:C:99:LYS:HE3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:271:ASP:O	1:A:274:VAL:HG22	2.13	0.48
1:B:824:VAL:C	1:B:826:ILE:H	2.21	0.48
1:B:1112:ALA:HB1	1:B:1140:PHE:CD1	2.49	0.48
1:A:1011:GLU:HA	1:A:1014:MET:HE2	1.96	0.48
1:A:68:PHE:CD1	1:A:114:GLN:HG3	2.49	0.47
2:C:77:ASP:OD1	2:C:77:ASP:N	2.47	0.47
1:A:753:LYS:O	1:A:757:LEU:HB2	2.13	0.47
2:C:29:ARG:NH1	3:F:42:GLU:OE1	2.45	0.47
1:B:773:THR:HA	1:B:777:GLN:HB2	1.97	0.47
1:A:802:LYS:HA	1:A:805:ILE:HG22	1.97	0.47
1:B:584:ILE:HG23	1:B:613:ARG:HH22	1.80	0.47
1:A:866:GLN:O	1:A:870:LEU:HG	2.15	0.47
1:B:682:LEU:HD12	1:B:710:LEU:HD21	1.97	0.47
2:C:151:ALA:HA	2:C:157:PHE:CE2	2.50	0.47
1:A:263:GLU:HA	1:A:266:ARG:HB2	1.97	0.47
1:A:286:ARG:NH2	1:A:335:GLU:O	2.44	0.47
1:B:288:ASP:C	1:B:290:ASP:H	2.23	0.47
1:A:224:GLN:HG2	1:A:291:MET:HG2	1.97	0.47
1:B:486:ARG:NH2	1:B:552:ASP:OD1	2.42	0.47
1:A:78:VAL:HG11	1:A:122:ILE:HD11	1.97	0.47
1:B:922:LEU:HG	1:B:955:VAL:HG21	1.97	0.46
1:A:730:ALA:HB1	1:A:775:VAL:HG22	1.97	0.46
2:D:77:ASP:OD1	2:D:77:ASP:N	2.48	0.46
1:B:1003:PHE:O	1:B:1007:MET:HG2	2.15	0.46
1:A:686:VAL:HG21	1:A:707:LEU:HD13	1.97	0.46
1:A:288:ASP:C	1:A:290:ASP:H	2.23	0.46
1:A:15:LEU:HD22	1:A:41:PHE:CE1	2.50	0.46
1:A:84:LEU:HD23	1:A:1108:ARG:HH11	1.81	0.46
1:A:306:HIS:HB3	1:A:354:VAL:HG13	1.97	0.46
1:A:614:VAL:HG12	1:A:638:ILE:HD11	1.98	0.46
1:B:952:ALA:HB3	1:B:994:LYS:HZ2	1.79	0.46
3:E:87:ASN:O	3:E:136:THR:HA	2.15	0.46
1:B:1071:THR:HG22	1:B:1073:ALA:H	1.81	0.46
1:A:682:LEU:HD12	1:A:710:LEU:HD21	1.97	0.46
1:B:756:VAL:HG21	1:B:815:ILE:HG12	1.97	0.46
1:A:116:LEU:HD22	1:A:159:ILE:HG12	1.97	0.46
1:B:1112:ALA:HB1	1:B:1140:PHE:HD1	1.81	0.45
2:C:113:GLU:HB3	2:C:114:ASN:H	1.67	0.45
1:A:493:MET:HE1	1:A:604:VAL:HA	1.97	0.45
1:B:717:ALA:HA	1:B:720:VAL:HG12	1.97	0.45
1:A:68:PHE:CE1	1:A:114:GLN:HG3	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:ILE:HG23	1:A:413:LEU:HD13	1.98	0.45
1:A:1030:THR:OG1	1:A:1031:PRO:HD3	2.17	0.45
1:B:84:LEU:HD23	1:B:1108:ARG:HH11	1.82	0.45
1:A:490:GLU:HG2	1:A:568:TYR:CE1	2.52	0.45
1:A:723:CYS:O	1:A:725:ASN:N	2.42	0.45
1:B:828:PHE:CE2	1:B:874:LYS:HB3	2.51	0.45
1:A:584:ILE:HD12	1:A:601:THR:HG21	1.97	0.45
1:A:1097:LEU:HB3	1:A:1106:TYR:HE1	1.82	0.45
2:C:35:PHE:HE2	2:C:37:LYS:HG2	1.82	0.45
1:A:805:ILE:HD11	1:A:841:LEU:HG	1.98	0.45
1:A:959:VAL:O	1:A:963:LEU:HB2	2.16	0.45
1:B:98:LEU:HD21	1:B:119:VAL:HG11	1.97	0.45
1:B:595:ILE:HG22	1:B:598:TYR:HB3	1.97	0.45
1:B:618:GLU:HG3	1:B:630:LEU:HD13	1.99	0.45
1:A:1140:PHE:CE1	1:A:1144:VAL:HG21	2.51	0.45
1:A:1055:LYS:HE2	1:A:1092:LEU:HD21	1.98	0.45
1:A:618:GLU:O	1:A:622:ILE:HG12	2.17	0.44
2:D:141:LYS:HD2	2:D:144:LEU:HD12	1.99	0.44
3:E:95:GLN:C	3:E:97:GLY:H	2.25	0.44
1:B:154:THR:HG23	1:B:225:ARG:HD3	1.98	0.44
1:A:1089:TYR:HE2	1:A:1117:THR:HG22	1.82	0.44
2:D:13:LEU:HA	2:D:85:CYS:O	2.17	0.44
2:C:13:LEU:HA	2:C:85:CYS:O	2.18	0.44
2:C:89:MET:HE3	2:C:89:MET:HB2	1.90	0.44
1:A:304:SER:HA	1:A:354:VAL:HG21	1.99	0.44
1:A:611:VAL:HG21	1:A:641:PHE:CZ	2.53	0.44
1:B:286:ARG:HG3	1:B:340:GLU:OE2	2.18	0.44
1:B:187:LYS:O	1:B:191:GLN:HG3	2.17	0.44
1:A:981:TYR:HA	1:A:984:ILE:HG12	1.99	0.44
1:B:342:VAL:HG23	1:B:398:VAL:HG13	1.99	0.44
2:D:35:PHE:CE2	2:D:37:LYS:HG2	2.53	0.44
1:B:608:LEU:HB3	1:B:645:TRP:CH2	2.52	0.44
1:B:510:GLN:HG2	1:B:532:LEU:HD13	1.99	0.43
1:A:317:VAL:HG22	1:A:359:VAL:HG13	2.00	0.43
1:B:1054:LEU:HD13	1:B:1081:LEU:HB3	2.00	0.43
1:B:1030:THR:HG23	1:B:1080:THR:HG21	2.01	0.43
2:C:68:GLY:N	4:C:217:GTP:O1G	2.46	0.43
1:A:154:THR:HG23	1:A:225:ARG:HD3	2.01	0.43
1:B:199:PHE:HB2	1:B:233:VAL:HG11	1.99	0.43
1:A:25:PRO:HA	1:A:26:PRO:HD3	1.92	0.43
1:A:126:GLY:HA3	1:A:132:ILE:HG13	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:42:GLU:HB2	3:E:59:VAL:HG13	1.99	0.43
1:B:164:LEU:HD11	1:B:195:LEU:HD13	2.01	0.43
1:B:905:GLU:OE1	2:C:37:LYS:HD3	2.19	0.43
1:B:914:ILE:HG21	1:B:962:ILE:HG21	1.99	0.43
1:B:1034:GLU:OE1	1:B:1034:GLU:N	2.50	0.43
1:A:432:VAL:HG21	1:A:484:LEU:HD13	2.00	0.43
1:B:995:ILE:HB	1:B:996:PRO:HD3	2.01	0.43
1:A:616:GLU:HG2	1:A:681:TYR:OH	2.19	0.43
1:B:584:ILE:HG13	1:B:601:THR:HG21	2.00	0.43
1:B:687:ILE:HD11	1:B:729:LEU:HD22	2.01	0.42
1:A:1029:LEU:HD23	1:A:1029:LEU:HA	1.83	0.42
1:B:650:LEU:HD23	1:B:650:LEU:HA	1.86	0.42
1:A:540:HIS:ND1	1:A:637:ASP:OD2	2.29	0.42
3:E:63:ILE:HG13	3:E:138:LEU:HD21	2.02	0.42
1:B:141:VAL:HG13	1:B:156:ALA:HB1	2.01	0.42
1:B:879:TYR:CE1	1:B:921:LEU:HG	2.54	0.42
3:E:109:ARG:NH2	3:E:144:GLU:OE1	2.53	0.42
1:B:679:ILE:HD11	1:B:720:VAL:HB	2.01	0.42
1:A:1078:PHE:O	1:A:1082:VAL:HG23	2.19	0.42
1:B:616:GLU:HG2	1:B:681:TYR:OH	2.19	0.42
1:B:770:GLN:NE2	1:B:774:GLU:OE1	2.53	0.42
1:B:1084:LEU:HD23	1:B:1084:LEU:HA	1.93	0.42
2:D:151:ALA:HA	2:D:157:PHE:HE2	1.84	0.42
3:E:134:LEU:HB3	3:E:150:LYS:HB2	2.01	0.42
1:B:263:GLU:HG2	1:B:266:ARG:HE	1.85	0.42
1:B:1030:THR:HB	1:B:1031:PRO:HD3	2.01	0.42
1:B:1140:PHE:O	1:B:1144:VAL:N	2.48	0.42
1:B:65:TYR:CD1	2:C:81:ILE:HD12	2.54	0.42
1:A:29:VAL:O	1:A:30:SER:OG	2.29	0.42
1:B:683:LEU:HD12	1:B:683:LEU:HA	1.88	0.42
1:A:995:ILE:HB	1:A:996:PRO:HD3	2.02	0.42
1:B:499:LEU:O	1:B:503:ARG:HG2	2.20	0.42
1:B:678:ILE:O	1:B:682:LEU:HG	2.20	0.42
1:A:499:LEU:O	1:A:503:ARG:HG2	2.20	0.42
1:B:866:GLN:O	1:B:870:LEU:HG	2.20	0.41
2:C:101:VAL:N	2:C:102:PRO:HD2	2.35	0.41
1:A:662:LEU:O	1:A:666:THR:HG22	2.20	0.41
2:C:11:PHE:CG	2:C:168:LEU:HD13	2.55	0.41
1:A:543:ILE:HG12	1:A:611:VAL:HG12	2.02	0.41
1:A:850:GLU:H	1:A:850:GLU:CD	2.28	0.41
1:B:80:ARG:HG3	1:B:1146:GLY:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1146:GLY:O	1:B:1148:LEU:N	2.54	0.41
1:B:1013:GLY:HA2	1:B:1017:MET:HE3	2.02	0.41
1:A:636:LYS:HD3	1:A:698:ASP:HB3	2.01	0.41
3:F:98:TYR:CE1	3:F:113:ARG:HG2	2.56	0.41
1:A:220:SER:O	1:A:224:GLN:HG3	2.21	0.41
1:B:336:ILE:O	1:B:340:GLU:HB2	2.20	0.41
1:B:785:ARG:O	1:B:789:GLN:HG3	2.21	0.41
2:C:12:LYS:HE3	2:C:64:TRP:CE2	2.56	0.41
1:A:195:LEU:HD11	1:A:233:VAL:HG13	2.02	0.41
1:A:584:ILE:HD11	1:A:598:TYR:HB2	2.03	0.41
1:A:1054:LEU:HD13	1:A:1081:LEU:HB3	2.03	0.41
1:B:1091:GLU:O	1:B:1095:THR:OG1	2.30	0.41
1:A:622:ILE:HG23	1:A:692:VAL:HG11	2.02	0.41
1:A:687:ILE:HD11	1:A:729:LEU:CD1	2.51	0.41
1:A:1057:VAL:HG12	1:A:1078:PHE:CZ	2.55	0.41
1:B:1052:HIS:O	1:B:1052:HIS:ND1	2.50	0.40
2:D:29:ARG:HG2	2:D:157:PHE:CZ	2.56	0.40
3:F:152:MET:SD	3:F:152:MET:N	2.94	0.40
1:A:551:ALA:HB3	1:A:644:ARG:HD3	2.02	0.40
1:A:915:MET:HG3	1:A:983:LEU:HD12	2.03	0.40
2:D:88:ILE:HD11	2:D:108:LEU:HD12	2.03	0.40
1:B:188:ARG:CZ	1:A:311:PRO:HB3	2.51	0.40
1:A:776:LEU:HD23	1:A:830:PHE:CE2	2.57	0.40
2:D:101:VAL:N	2:D:102:PRO:HD2	2.36	0.40
3:E:20:MET:HE3	3:E:20:MET:HB2	1.91	0.40
1:A:140:GLU:O	1:A:144:LEU:HD13	2.22	0.40
1:A:463:ILE:HD12	3:E:50:5CT:H22	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	951/1113 (85%)	915 (96%)	35 (4%)	1 (0%)	48	79
1	B	1036/1113 (93%)	985 (95%)	47 (4%)	4 (0%)	30	62
2	C	171/176 (97%)	167 (98%)	4 (2%)	0	100	100
2	D	170/176 (97%)	166 (98%)	4 (2%)	0	100	100
3	E	134/141 (95%)	127 (95%)	6 (4%)	1 (1%)	18	52
3	F	135/141 (96%)	128 (95%)	7 (5%)	0	100	100
All	All	2597/2860 (91%)	2488 (96%)	103 (4%)	6 (0%)	43	73

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	554	THR
1	A	312	ASP
1	B	214	ASN
3	E	96	ASP
1	B	1121	THR
1	B	899	ILE

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	825/994 (83%)	819 (99%)	6 (1%)	76	83
1	B	901/994 (91%)	885 (98%)	16 (2%)	51	74
2	C	143/154 (93%)	137 (96%)	6 (4%)	26	60
2	D	146/154 (95%)	145 (99%)	1 (1%)	76	83
3	E	111/118 (94%)	107 (96%)	4 (4%)	31	63
3	F	114/118 (97%)	109 (96%)	5 (4%)	25	59
All	All	2240/2532 (88%)	2202 (98%)	38 (2%)	53	75

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	263	GLU
1	B	558	THR
1	B	576	VAL
1	B	595	ILE
1	B	611	VAL
1	B	697	GLN
1	B	711	VAL
1	B	715	GLU
1	B	767	THR
1	B	791	ASN
1	B	858	VAL
1	B	885	LEU
1	B	963	LEU
1	B	1064	GLN
1	B	1121	THR
1	B	1131	MET
2	C	14	VAL
2	C	15	LEU
2	C	30	HIS
2	C	63	VAL
2	C	69	LEU
2	C	109	VAL
3	F	59	VAL
3	F	76	THR
3	F	83	ASN
3	F	114	LEU
3	F	152	MET
1	A	63	VAL
1	A	263	GLU
1	A	611	VAL
1	A	711	VAL
1	A	885	LEU
1	A	963	LEU
2	D	30	HIS
3	E	59	VAL
3	E	83	ASN
3	E	112	LEU
3	E	114	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	515	GLN

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Mol	Chain	Res	Type
1	B	777	GLN
1	B	781	GLN
1	B	836	ASN
1	B	920	ASN
2	C	143	ASN
1	A	191	GLN
1	A	434	ASN
1	A	477	GLN
1	A	777	GLN
1	A	837	ASN
1	A	864	HIS
2	D	10	GLN
2	D	62	ASN
2	D	143	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	5CT	E	50	3	13,14,15	0.92	0	8,15,17	0.79	1 (12%)
3	5CT	F	50	3	13,14,15	0.92	0	8,15,17	0.72	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	5CT	E	50	3	-	6/13/14/16	-
3	5CT	F	50	3	-	8/13/14/16	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	E	50	5CT	C1-NZ-CE	-2.01	108.99	113.38

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	F	50	5CT	NZ-C1-C2-C3
3	F	50	5CT	NZ-C1-C2-O1
3	F	50	5CT	C1-C2-C3-C4
3	E	50	5CT	NZ-C1-C2-C3
3	E	50	5CT	NZ-C1-C2-O1
3	E	50	5CT	C1-C2-C3-C4
3	E	50	5CT	O1-C2-C3-C4
3	E	50	5CT	CD-CE-NZ-C1
3	F	50	5CT	CD-CE-NZ-C1
3	F	50	5CT	CE-CD-CG-CB
3	F	50	5CT	CA-CB-CG-CD
3	E	50	5CT	CA-CB-CG-CD
3	F	50	5CT	C2-C3-C4-N1
3	F	50	5CT	O1-C2-C3-C4

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	50	5CT	1	0

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	GTP	D	217	5	33,34,34	1.25	2 (6%)	50,54,54	1.58	9 (18%)
4	GTP	C	217	5	33,34,34	0.92	0	50,54,54	1.56	9 (18%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GTP	D	217	5	-	4/22/38/38	0/3/3/3
4	GTP	C	217	5	-	4/22/38/38	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	217	GTP	PA-O1A	3.11	1.61	1.50
4	D	217	GTP	PB-O1B	3.10	1.61	1.50

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	217	GTP	C5-C4-N3	-5.32	119.92	128.39
4	C	217	GTP	C5-C4-N3	-5.20	120.11	128.39
4	D	217	GTP	C2-N3-C4	5.02	120.94	112.30
4	C	217	GTP	C2-N3-C4	4.89	120.73	112.30
4	D	217	GTP	N9-C4-N3	4.03	134.02	125.95
4	C	217	GTP	N9-C4-N3	3.94	133.84	125.95
4	C	217	GTP	C8-N7-C5	2.96	109.53	104.26
4	D	217	GTP	C8-N7-C5	2.89	109.41	104.26
4	C	217	GTP	C2-N1-C6	-2.61	120.38	125.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	217	GTP	C2-N1-C6	-2.59	120.41	125.11
4	C	217	GTP	C4-C5-N7	-2.22	107.15	110.67
4	C	217	GTP	C6-C5-N7	2.21	134.31	130.29
4	C	217	GTP	O6-C6-C5	-2.19	120.74	126.53
4	D	217	GTP	C6-C5-N7	2.18	134.26	130.29
4	D	217	GTP	C4-C5-N7	-2.18	107.22	110.67
4	C	217	GTP	N9-C8-N7	-2.12	109.47	113.40
4	D	217	GTP	O6-C6-C5	-2.12	120.94	126.53
4	D	217	GTP	N9-C8-N7	-2.04	109.62	113.40

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	217	GTP	O4'-C4'-C5'-O5'
4	D	217	GTP	C3'-C4'-C5'-O5'
4	C	217	GTP	PB-O3B-PG-O3G
4	C	217	GTP	PB-O3B-PG-O2G
4	D	217	GTP	PB-O3B-PG-O3G
4	C	217	GTP	PG-O3B-PB-O2B
4	D	217	GTP	PG-O3B-PB-O2B
4	C	217	GTP	PG-O3B-PB-O1B

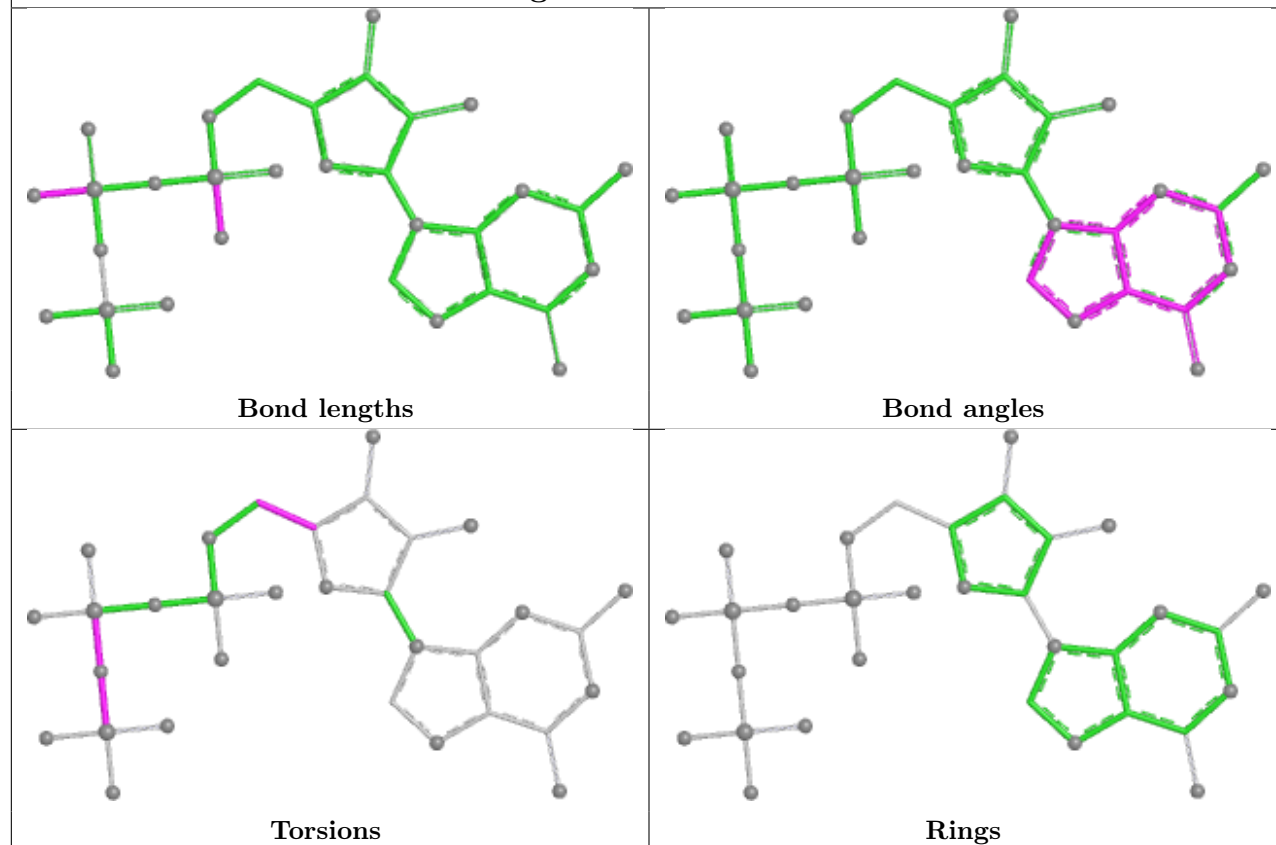
There are no ring outliers.

2 monomers are involved in 4 short contacts:

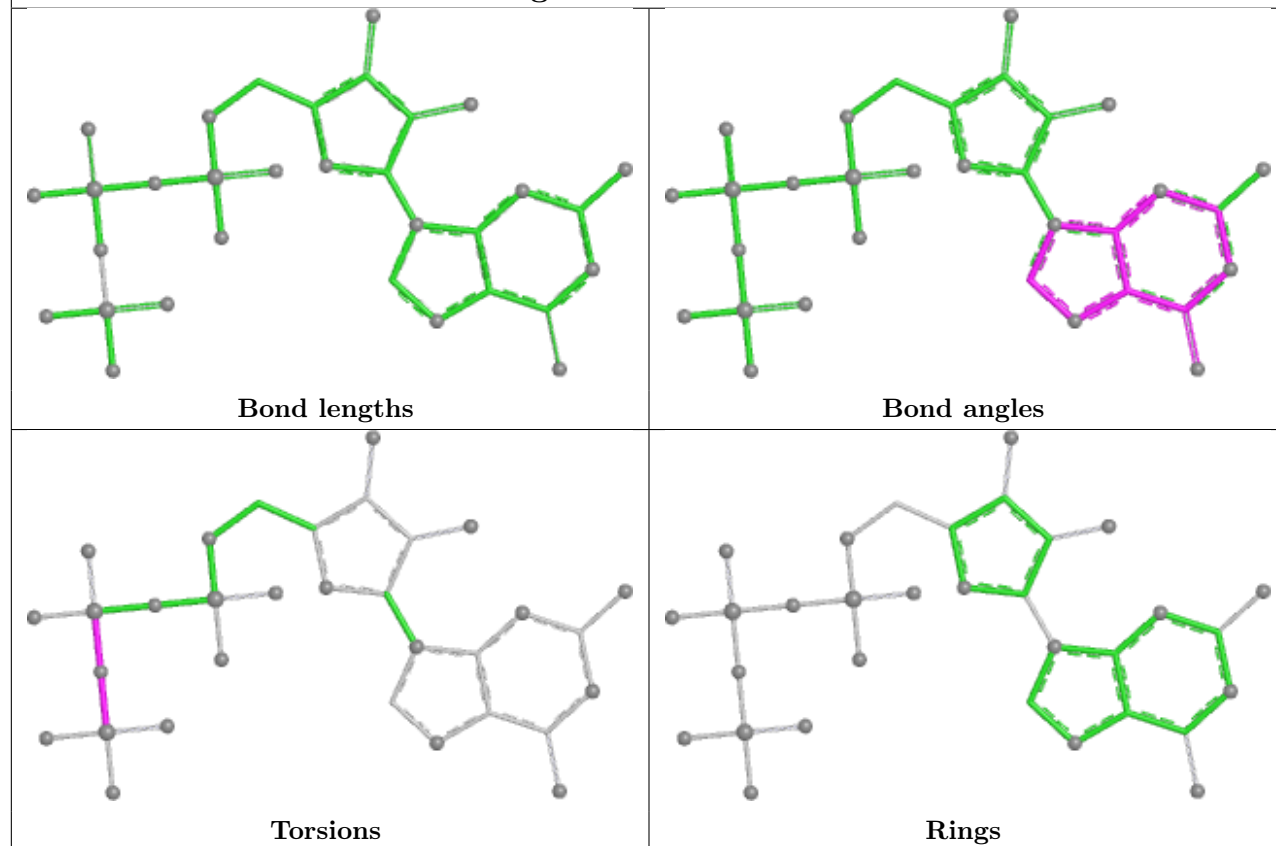
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	217	GTP	1	0
4	C	217	GTP	3	0

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 217



Ligand GTP C 217



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	983/1113 (88%)	-0.06	16 (1%) 70 51	95, 133, 173, 194	0
1	B	1052/1113 (94%)	-0.22	12 (1%) 78 61	68, 108, 158, 180	0
2	C	173/176 (98%)	-0.40	0 100 100	63, 83, 116, 130	0
2	D	172/176 (97%)	-0.32	2 (1%) 76 58	84, 102, 136, 149	0
3	E	136/141 (96%)	-0.16	0 100 100	108, 126, 146, 152	0
3	F	137/141 (97%)	-0.27	1 (0%) 84 69	71, 94, 133, 149	0
All	All	2653/2860 (92%)	-0.18	31 (1%) 76 58	63, 118, 165, 194	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1002	LEU	3.6
1	A	532	LEU	3.4
1	A	1033	ALA	3.4
3	F	76	THR	3.3
1	B	1062	VAL	3.2
1	B	1126	ASP	3.2
1	B	1067	ASN	3.1
1	A	1120	SER	3.1
1	A	952	ALA	3.0
1	A	1088	GLU	3.0
1	A	451	THR	2.9
1	B	1148	LEU	2.8
1	A	182	PHE	2.7
1	B	1046	LEU	2.6
1	A	1105	ILE	2.6
1	A	9	PRO	2.5
1	B	516	GLN	2.5
1	A	1083	CYS	2.4
1	A	1078	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
2	D	138	PHE	2.4
1	B	392	LEU	2.3
1	B	1127	ARG	2.3
1	B	457	SER	2.3
1	A	1145	GLY	2.2
2	D	7	PRO	2.2
1	B	129	ASP	2.2
1	A	772	TRP	2.1
1	A	1071	THR	2.1
1	A	760	PHE	2.1
1	B	924	LYS	2.0
1	B	1069	GLU	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	5CT	E	50	15/16	0.90	0.11	105,111,117,118	0
3	5CT	F	50	15/16	0.94	0.10	89,90,92,92	0

6.3 Carbohydrates [i](#)

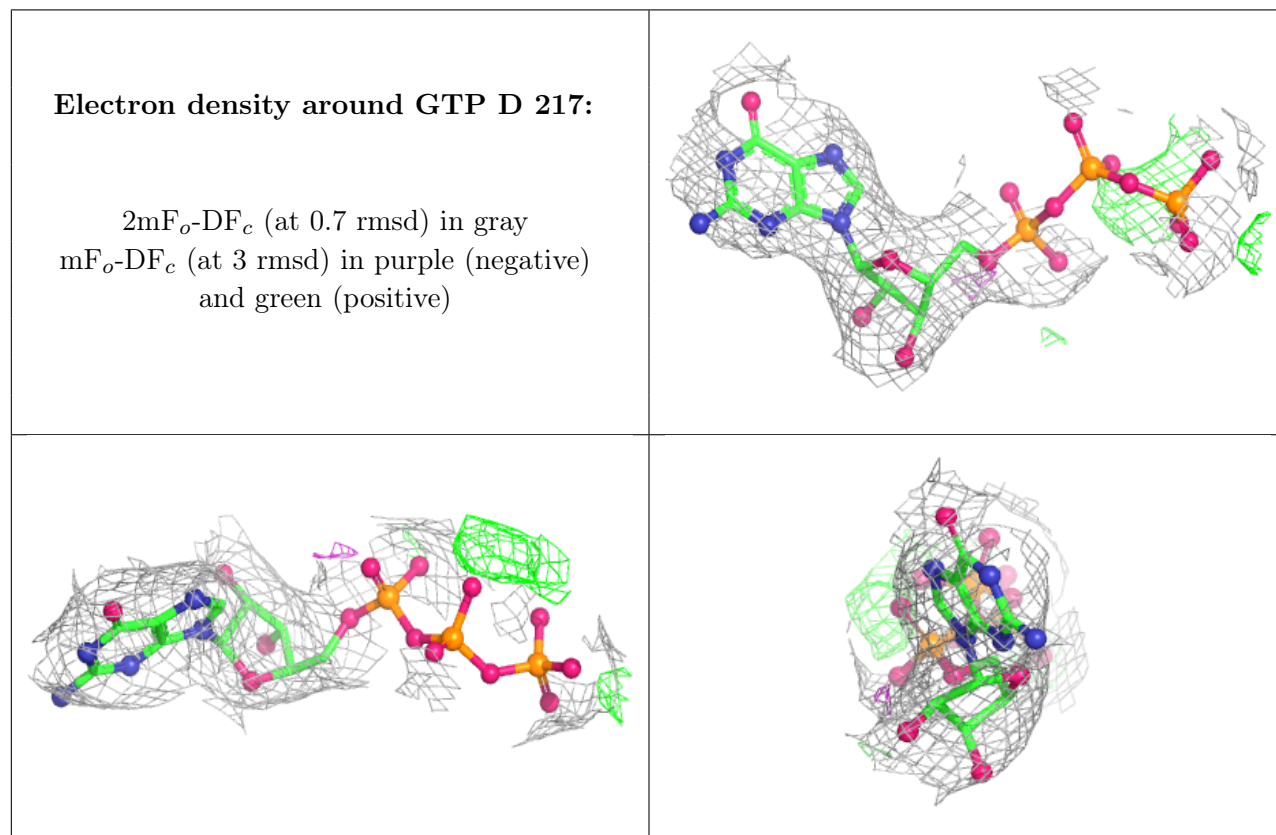
There are no oligosaccharides in this entry.

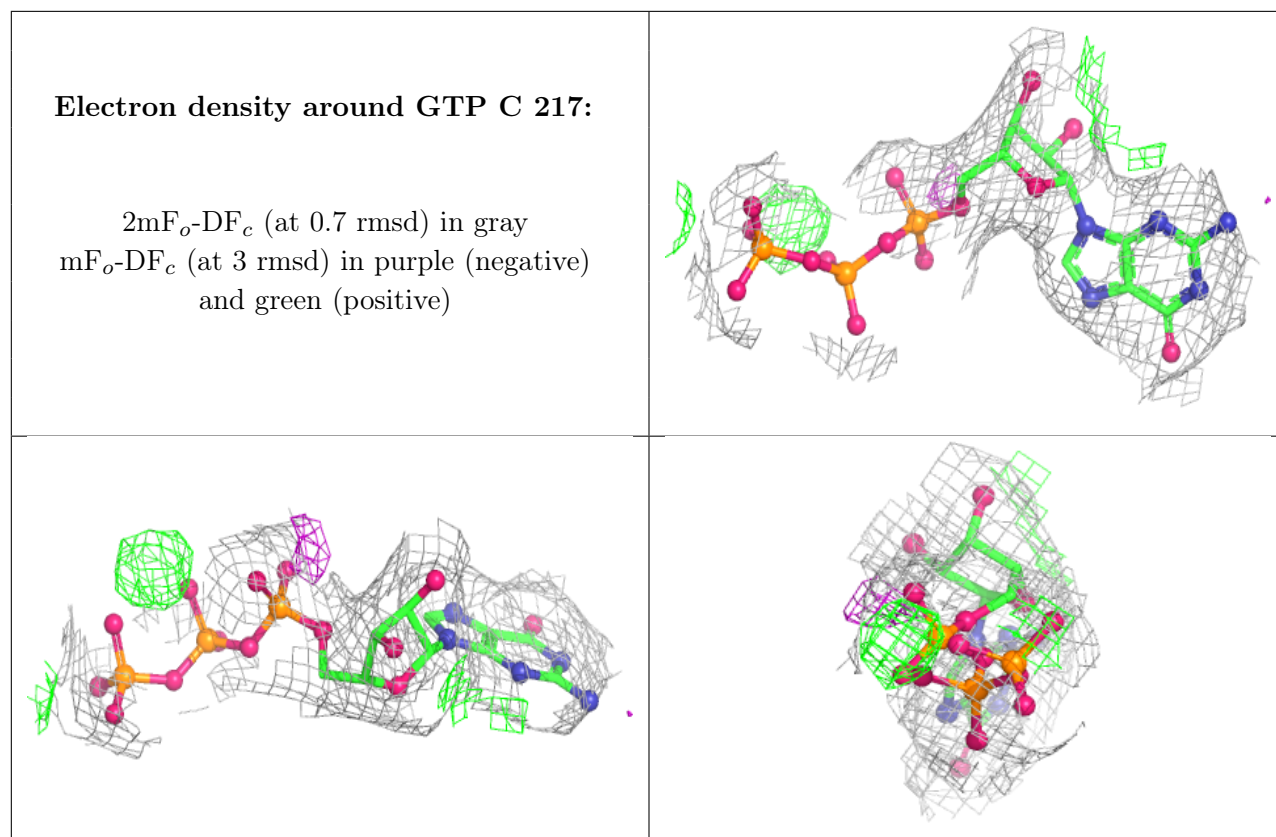
6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	MG	D	218	1/1	0.94	0.13	82,82,82,82	0
4	GTP	D	217	32/32	0.96	0.06	81,87,101,104	0
4	GTP	C	217	32/32	0.97	0.06	58,63,74,85	0
5	MG	C	218	1/1	0.99	0.07	58,58,58,58	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.