



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2014 – 04:14 PM GMT

PDB ID : 4C0Q  
Title : Transportin 3 in complex with Ran(Q69L)GTP  
Authors : Maertens, G.; Hare, S.; Cherepanov, P.  
Deposited on : 2013-08-06  
Resolution : 3.42 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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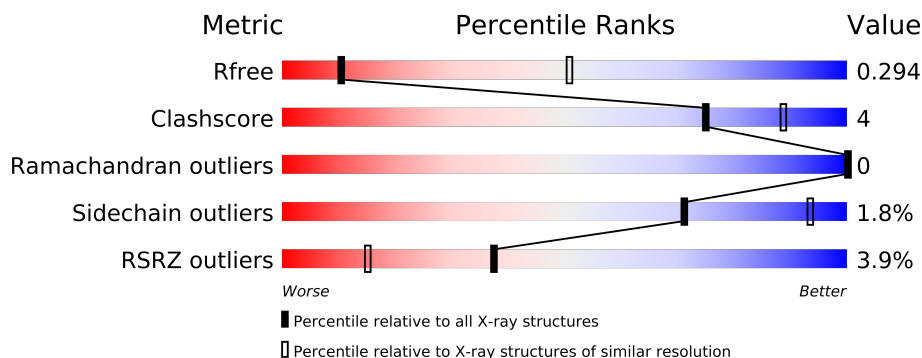
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.16 November 2013  
Xtriage (Phenix) : dev-1323  
EDS : trunk22714  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk22714

# 1 Overall quality at a glance

The reported resolution of this entry is 3.42 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))
$R_{free}$	66092	1013 (3.54-3.30)
Clashscore	79885	1270 (3.54-3.30)
Ramachandran outliers	78287	1232 (3.54-3.30)
Sidechain outliers	78261	1232 (3.54-3.30)
RSRZ outliers	66119	1013 (3.54-3.30)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	923	
1	B	923	
2	C	215	
2	D	215	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	MG	C	1178	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15766 atoms, of which 0 are hydrogen and 0 are deuterium.

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 TRANSPORTIN-3.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	871	Total	C	N	O	S	Se	0	0	0
			6927	4420	1177	1277	31	22			
1	B	749	Total	C	N	O	S	Se	0	0	0
			5979	3838	1007	1088	26	20			

- Molecule 2 is a protein called GTP-BINDING NUCLEAR PROTEIN RAN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	169	Total	C	N	O	S	0	0	0
			1376	896	241	235	4			
2	D	175	Total	C	N	O	S	0	0	0
			1418	923	247	243	5			

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 MAGNESIUM ION (three-letter code: MG) (formula: Mg).

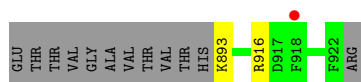
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

- Molecule 4 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula: C<sub>10</sub>H<sub>16</sub>N<sub>5</sub>O<sub>14</sub>P<sub>3</sub>).



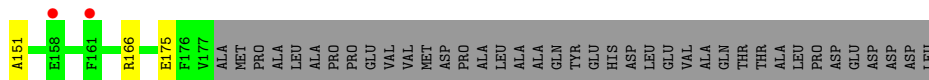
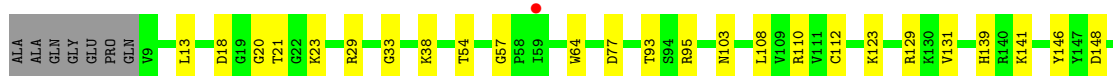
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	C	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
4	D	1	Total	C	N	O	P	0	0
			32	10	5	14	3		





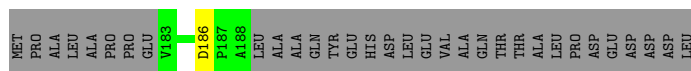
- Molecule 2: GTP-BINDING NUCLEAR PROTEIN RAN

Chain C:



- Molecule 2: GTP-BINDING NUCLEAR PROTEIN RAN

Chain D:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.95Å 93.51Å 104.71Å 78.43° 68.29° 68.28°	Depositor
Resolution (Å)	39.95 – 3.42 39.94 – 3.42	Depositor EDS
% Data completeness (in resolution range)	99.1 (39.95-3.42) 99.2 (39.94-3.42)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.14 (at 3.40Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.268 , 0.290 0.270 , 0.294	Depositor DCC
$R_{free}$ test set	1766 reflections (5.05%)	DCC
Wilson B-factor (Å <sup>2</sup> )	135.4	Xtriage
Anisotropy	0.400	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 55.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 34990 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	15766	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	137.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.75% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GTP, 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	0.22	0/7040	0.41	0/9520
1	B	0.22	0/6074	0.41	0/8200
2	C	0.23	0/1410	0.41	0/1904
2	D	0.24	0/1452	0.42	0/1961
All	All	0.22	0/15976	0.41	0/21585

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6927	0	0	25	0
1	B	5979	0	0	16	0
2	C	1376	0	0	14	0
2	D	1418	0	0	13	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	C	32	0	12	2	0
4	D	32	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	15766	0	24	63	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 4.

All (63) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:671:ARG:NH1	1:A:710:GLU:OE2	2.09	0.85
1:A:917:ASP:OD1	2:D:56:ARG:NH2	2.11	0.84
1:A:68:LYS:NZ	2:C:77:ASP:OD2	2.21	0.74
1:B:68:LYS:NZ	2:D:77:ASP:OD2	2.25	0.69
1:B:340:GLU:OE1	1:B:402:ARG:NH2	2.26	0.68
1:A:157:ARG:NH1	2:C:103:ASN:OD1	2.27	0.67
2:C:29:ARG:NH2	2:C:33:GLY:O	2.27	0.66
2:C:146:TYR:OH	2:C:148:ASP:OD1	2.17	0.62
2:D:70:GLU:OE2	2:D:76:ARG:NH2	2.34	0.61
2:D:56:ARG:NH1	2:D:171:ASP:OD2	2.34	0.60
2:D:29:ARG:NH1	2:D:154:ASN:OD1	2.35	0.59
1:B:247:ASP:OD1	1:B:297:ARG:NH2	2.35	0.59
2:D:166:ARG:NH2	2:D:175:GLU:OE2	2.36	0.58
2:D:18:ASP:OD1	2:D:104:TRP:NE1	2.37	0.57
1:A:247:ASP:OD1	1:A:297:ARG:NH2	2.38	0.57
1:A:456:ASN:ND2	1:A:495:ASN:OD1	2.38	0.57
2:C:13:LEU:O	2:C:64:TRP:N	2.38	0.57
2:C:23:LYS:N	4:C:1179:GTP:O2B	2.38	0.57
1:A:692:TYR:OH	1:A:732:THR:OG1	2.23	0.57
2:D:29:ARG:NH2	2:D:33:GLY:O	2.38	0.56
1:B:68:LYS:NZ	1:B:107:GLN:OE1	2.40	0.54
1:A:521:LYS:O	1:A:525:ASN:ND2	2.40	0.54
1:B:556:GLU:OE1	1:B:556:GLU:N	2.40	0.54
1:A:176:TYR:O	1:A:179:THR:OG1	2.26	0.54
2:D:53:HIS:O	2:D:177:VAL:N	2.39	0.53
1:A:556:GLU:N	1:A:556:GLU:OE1	2.42	0.53
1:B:582:SER:O	1:B:586:SER:OG	2.27	0.53
1:B:235:ASP:OD2	1:B:286:ARG:NH2	2.42	0.52
1:A:831:GLN:OE1	1:A:874:ASN:ND2	2.41	0.52
1:A:395:ASP:N	1:A:395:ASP:OD1	2.42	0.52
2:C:108:LEU:O	2:C:112:CYS:N	2.43	0.52
1:B:176:TYR:O	1:B:179:THR:OG1	2.29	0.51
2:D:146:TYR:OH	2:D:148:ASP:OD1	2.29	0.51
1:A:613:ASP:OD1	1:A:661:ARG:NE	2.43	0.51
1:A:152:GLU:OE1	2:C:110:ARG:NH2	2.44	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:233:GLN:NE2	1:A:275:THR:OG1	2.45	0.49
1:B:79:PRO:O	1:B:82:SER:OG	2.30	0.49
1:A:805:THR:O	1:A:818:ARG:NH1	2.45	0.49
1:B:831:GLN:OE1	1:B:874:ASN:ND2	2.45	0.49
1:B:395:ASP:OD1	1:B:395:ASP:N	2.46	0.48
1:B:232:LEU:O	1:B:295:TYR:OH	2.31	0.48
2:D:18:ASP:OD2	2:D:100:ASN:ND2	2.47	0.48
1:A:313:THR:O	1:A:313:THR:OG1	2.32	0.48
1:B:313:THR:OG1	1:B:313:THR:O	2.32	0.47
1:A:257:GLU:OE1	1:A:257:GLU:N	2.47	0.47
1:A:232:LEU:O	1:A:295:TYR:OH	2.33	0.47
2:C:20:GLY:O	2:C:123:LYS:NZ	2.48	0.46
2:D:108:LEU:O	2:D:112:CYS:N	2.48	0.46
2:C:166:ARG:NH2	2:C:175:GLU:OE2	2.48	0.46
1:B:257:GLU:OE1	1:B:257:GLU:N	2.48	0.46
1:A:79:PRO:O	1:A:82:SER:OG	2.33	0.46
2:C:151:ALA:N	4:C:1179:GTP:O6	2.51	0.43
2:C:129:ARG:NE	2:C:131:VAL:O	2.52	0.43
2:C:18:ASP:O	2:C:21:THR:OG1	2.37	0.42
2:D:13:LEU:O	2:D:64:TRP:N	2.52	0.42
1:B:325:GLU:OE2	1:B:371:ARG:NH1	2.53	0.41
1:A:779:TRP:O	1:A:783:SER:N	2.53	0.41
2:C:54:THR:OG1	2:C:57:GLY:O	2.38	0.41
1:A:618:ILE:O	1:A:622:THR:OG1	2.39	0.41
1:A:277:GLU:O	1:A:281:HIS:ND1	2.55	0.40
1:A:68:LYS:NZ	1:A:107:GLN:OE1	2.54	0.40
1:A:478:ARG:NH1	1:A:512:GLU:OE1	2.54	0.40
1:B:394:ASP:OD1	1:B:397:GLY:N	2.55	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	857/923 (93%)	834 (97%)	23 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	717/923 (78%)	701 (98%)	16 (2%)	0	100	100
2	C	167/215 (78%)	163 (98%)	4 (2%)	0	100	100
2	D	171/215 (80%)	166 (97%)	5 (3%)	0	100	100
All	All	1912/2276 (84%)	1864 (98%)	48 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	779/798 (98%)	768 (99%)	11 (1%)	78	95
1	B	671/798 (84%)	661 (98%)	10 (2%)	76	95
2	C	149/184 (81%)	144 (97%)	5 (3%)	49	87
2	D	154/184 (84%)	148 (96%)	6 (4%)	43	84
All	All	1753/1964 (89%)	1721 (98%)	32 (2%)	71	93

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	GLN
1	A	198	LEU
1	A	209	PHE
1	A	282	MSE
1	A	395	ASP
1	A	459	LEU
1	A	546	ARG
1	A	584	LEU
1	A	667	ARG
1	A	759	PHE
1	A	916	ARG
1	B	65	GLN
1	B	198	LEU
1	B	209	PHE

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Mol	Chain	Res	Type
1	B	395	ASP
1	B	584	LEU
1	B	667	ARG
1	B	759	PHE
1	B	837	LEU
1	B	893	LYS
1	B	916	ARG
2	C	38	LYS
2	C	93	THR
2	C	95	ARG
2	C	139	HIS
2	C	141	LYS
2	D	38	LYS
2	D	93	THR
2	D	95	ARG
2	D	139	HIS
2	D	141	LYS
2	D	186	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates 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	C	1179	3	34,34,34	1.06	2 (5%)	52,54,54	2.76	9 (17%)
4	GTP	D	1190	3	34,34,34	1.06	2 (5%)	52,54,54	2.69	9 (17%)

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	C	1179	3	-	0/22/38/38	0/3/3/3
4	GTP	D	1190	3	-	0/22/38/38	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	1179	GTP	C2-N3	4.26	1.38	1.33
4	D	1190	GTP	C2-N3	4.06	1.38	1.33
4	D	1190	GTP	C5-N7	-2.29	1.35	1.38
4	C	1179	GTP	C5-N7	-2.21	1.35	1.38

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	1179	GTP	C6-C5-N7	-15.29	132.08	134.14
4	D	1190	GTP	C6-C5-N7	-14.25	132.22	134.14
4	D	1190	GTP	PA-O3A-PB	-5.66	116.27	131.93
4	C	1179	GTP	C5-C4-N3	-5.22	120.06	126.07
4	D	1190	GTP	C5-C4-N3	-5.00	120.31	126.07
4	C	1179	GTP	PA-O3A-PB	-4.70	118.90	131.93
4	D	1190	GTP	PB-O3B-PG	-4.66	119.02	131.93
4	C	1179	GTP	PB-O3B-PG	-4.54	119.37	131.93
4	C	1179	GTP	C2-N3-C4	4.22	120.36	115.30
4	D	1190	GTP	C2-N3-C4	4.13	120.25	115.30
4	C	1179	GTP	N3-C4-N9	4.05	132.84	126.91
4	D	1190	GTP	N3-C4-N9	3.74	132.39	126.91
4	C	1179	GTP	C3'-C2'-C1'	3.14	105.85	100.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1190	GTP	C3'-C2'-C1'	3.11	105.80	100.92
4	D	1190	GTP	O3A-PB-O3B	3.01	107.78	101.66
4	D	1190	GTP	C4'-O4'-C1'	2.74	112.73	109.72
4	C	1179	GTP	C4'-O4'-C1'	2.72	112.70	109.72
4	C	1179	GTP	O3A-PB-O3B	2.27	106.29	101.66

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	871/923 (94%)	0.29	16 (1%) 65 30	59, 130, 186, 206	0
1	B	749/923 (81%)	0.59	57 (7%) 14 6	92, 172, 210, 232	0
2	C	169/215 (78%)	0.33	3 (1%) 65 30	62, 104, 148, 158	0
2	D	175/215 (81%)	0.24	0 100 100	65, 89, 119, 134	0
All	All	1964/2276 (86%)	0.40	76 (3%) 37 15	59, 140, 201, 232	0

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	688	MSE	5.8
1	A	488	MSE	5.8
1	B	662	CYS	5.0
1	B	488	MSE	4.6
1	B	724	MSE	4.6
1	A	417	MSE	4.4
1	B	918	PHE	3.7
1	B	459	LEU	3.5
1	B	219	MSE	3.5
1	B	590	MSE	3.4
1	B	615	LEU	3.2
1	A	5	LYS	3.2
1	B	310	ILE	3.1
1	B	765	VAL	3.1
1	B	863	VAL	3.1
1	B	842	CYS	3.1
1	B	431	ASN	3.0
1	B	702	TYR	3.0
1	B	365	PHE	3.0
1	A	769	ARG	3.0
1	B	318	LEU	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	414	ILE	2.9
1	B	410	LEU	2.9
2	C	59	ILE	2.9
1	B	322	ARG	2.9
1	B	828	GLN	2.9
1	B	444	MSE	2.9
1	B	306	PHE	2.9
2	C	161	PHE	2.8
1	B	658	ILE	2.8
1	B	5	LYS	2.8
1	B	78	LEU	2.8
1	B	11	VAL	2.8
1	B	670	VAL	2.7
1	B	197	MSE	2.7
1	A	572	LEU	2.6
1	B	802	LEU	2.5
1	B	67	MSE	2.5
1	B	804	HIS	2.5
1	B	826	MSE	2.5
1	B	308	GLU	2.4
1	A	410	LEU	2.4
1	B	269	LEU	2.4
1	B	324	LEU	2.4
1	B	777	LEU	2.4
1	B	799	LEU	2.4
1	B	806	GLY	2.4
1	B	657	ARG	2.4
1	A	53	ILE	2.3
1	A	444	MSE	2.3
1	B	184	LEU	2.3
1	A	459	LEU	2.3
1	B	838	LEU	2.3
2	C	158	GLU	2.3
1	B	592	LEU	2.3
1	B	141	PRO	2.3
1	B	498	PHE	2.3
1	A	60	CYS	2.2
1	B	458	THR	2.2
1	B	75	PHE	2.2
1	B	829	LEU	2.2
1	A	8	LEU	2.2
1	A	11	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	97	LEU	2.2
1	A	310	ILE	2.2
1	B	168	GLU	2.2
1	B	752	LEU	2.1
1	B	413	LEU	2.1
1	B	44	TRP	2.1
1	A	419	CYS	2.1
1	B	608	PRO	2.1
1	B	751	ASP	2.1
1	A	46	ILE	2.0
1	B	666	LEU	2.0
1	B	843	PHE	2.0
1	A	826	MSE	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q < 0.9
3	MG	C	1178	1/1	0.36	2.60	76,76,76,76	0
4	GTP	C	1179	32/32	0.27	0.69	66,80,93,95	0
4	GTP	D	1190	32/32	0.24	-0.33	63,73,86,87	0
3	MG	D	1189	1/1	0.19	-0.69	68,68,68,68	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.