



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 06:11 pm BST

PDB ID : 4LHZ
Title : Crystal structure of GTP-bound Rab8:Rabin8
Authors : Guo, Z.; Hou, X.M.; Goody, R.S.; Itzen, A.
Deposited on : 2013-07-01
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

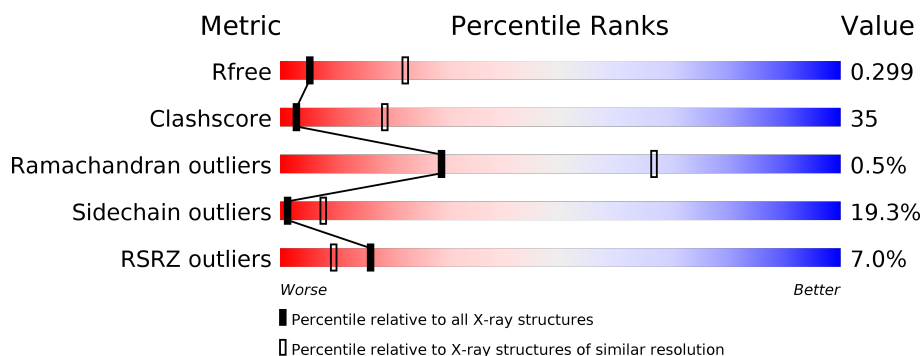
1 Overall quality at a glance

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}	130704	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (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 on 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	186	<div> <div>39%</div> <div>44%</div> <div>11%</div> <div>6%</div> </div>
1	B	186	<div> <div>39%</div> <div>38%</div> <div>16%</div> <div>6%</div> </div>
2	C	78	<div>8%</div> <div>65%</div> <div>26%</div> <div>8%</div>
2	D	78	<div>13%</div> <div>74%</div> <div>14%</div> <div>9%</div>
2	E	78	<div>13%</div> <div>68%</div> <div>24%</div> <div>5%</div>
2	F	78	<div>22%</div> <div>78%</div> <div>15%</div> <div></div>

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GTP	A	201	-	-	X	-
3	GTP	B	201	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4799 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 Ras-related protein Rab-8A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	175	Total	C	N	O	S	0	0	0
			1390	888	235	260	7			
1	B	175	Total	C	N	O	S	0	0	0
			1389	885	233	264	7			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP P61006
A	0	HIS	-	EXPRESSION TAG	UNP P61006
B	-1	GLY	-	EXPRESSION TAG	UNP P61006
B	0	HIS	-	EXPRESSION TAG	UNP P61006

- Molecule 2 is a protein called Rab-3A-interacting protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	72	Total	C	N	O	S	0	0	0
			481	293	89	97	2			
2	D	71	Total	C	N	O	S	0	0	0
			469	284	83	100	2			
2	E	76	Total	C	N	O	S	0	0	0
			498	303	92	101	2			
2	F	76	Total	C	N	O	S	0	0	0
			508	307	96	103	2			

There are 8 discrepancies between the modelled and reference sequences:

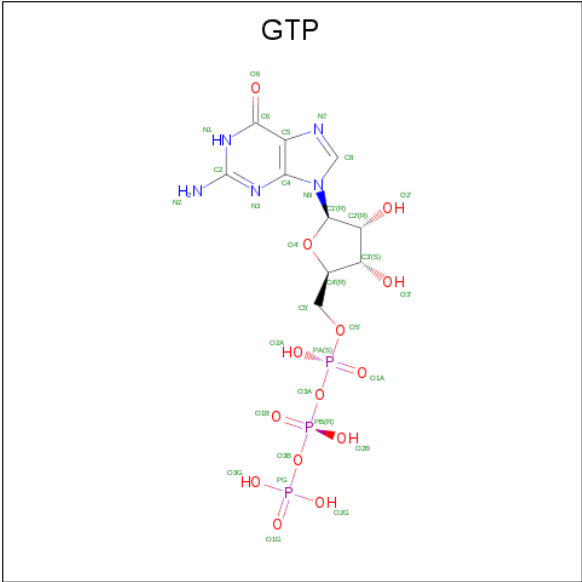
Chain	Residue	Modelled	Actual	Comment	Reference
C	155	GLY	-	EXPRESSION TAG	UNP Q96QF0
C	156	PRO	-	EXPRESSION TAG	UNP Q96QF0
D	155	GLY	-	EXPRESSION TAG	UNP Q96QF0
D	156	PRO	-	EXPRESSION TAG	UNP Q96QF0

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Chain	Residue	Modelled	Actual	Comment	Reference
E	155	GLY	-	EXPRESSION TAG	UNP Q96QF0
E	156	PRO	-	EXPRESSION TAG	UNP Q96QF0
F	155	GLY	-	EXPRESSION TAG	UNP Q96QF0
F	156	PRO	-	EXPRESSION TAG	UNP Q96QF0

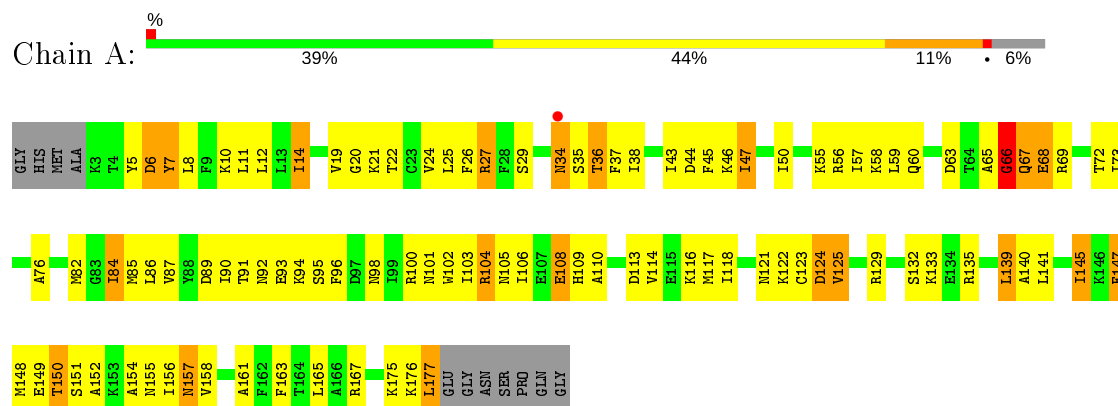
- Molecule 3 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃).



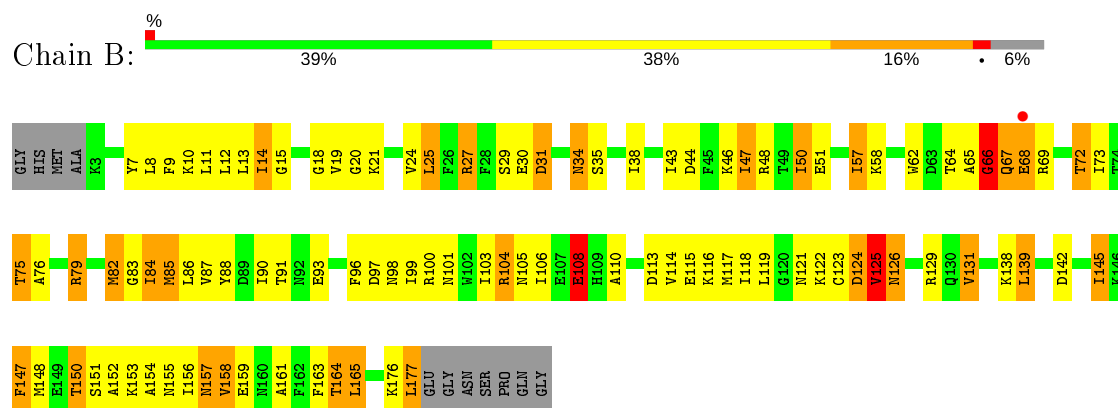
3 Residue-property plots

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

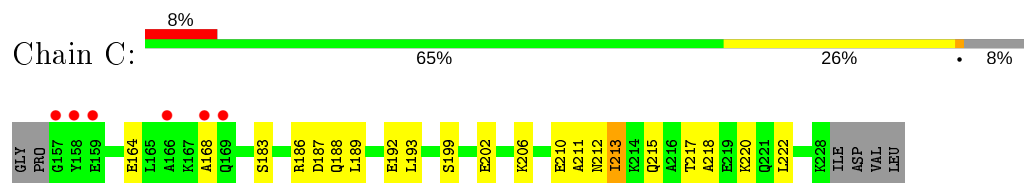
• Molecule 1: Ras-related protein Rab-8A



• Molecule 1: Ras-related protein Rab-8A



• Molecule 2: Rab-3A-interacting protein



• Molecule 2: Rab-3A-interacting protein

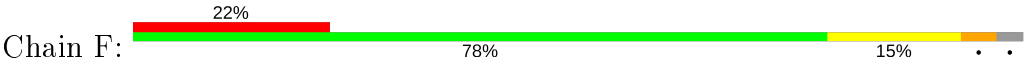




• Molecule 2: Rab-3A-interacting protein



• Molecule 2: Rab-3A-interacting protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	81.34Å 165.56Å 167.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.99 – 3.20 19.96 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.5 (19.99-3.20) 99.5 (19.96-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.58 (at 3.22Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.278 , 0.309 0.269 , 0.299	Depositor DCC
R_{free} test set	954 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	79.3	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 82.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4799	wwPDB-VP
Average B, all atoms (Å ²)	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.78% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	1.00	0/1411	1.03	2/1895 (0.1%)
1	B	0.91	0/1410	0.99	2/1896 (0.1%)
2	C	0.65	0/482	0.66	0/651
2	D	0.60	0/470	0.62	0/638
2	E	0.48	0/499	0.56	0/674
2	F	0.60	0/509	0.60	0/687
All	All	0.82	0/4781	0.87	4/6441 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	7	TYR	CB-CA-C	-5.50	99.39	110.40
1	B	66	GLY	N-CA-C	5.32	126.40	113.10
1	A	66	GLY	N-CA-C	5.31	126.36	113.10
1	B	125	VAL	N-CA-C	-5.11	97.19	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	6	ASP	Mainchain
1	B	108	GLU	Mainchain

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	1390	0	1388	125	0
1	B	1389	0	1372	154	0
2	C	481	0	409	27	0
2	D	469	0	380	22	0
2	E	498	0	411	16	0
2	F	508	0	423	20	0
3	A	32	0	12	10	0
3	B	32	0	12	11	0
All	All	4799	0	4407	321	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 35.

All (321) 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:A:14:ILE:CG2	1:A:84:ILE:HD11	1.44	1.46
1:A:14:ILE:CG2	1:A:84:ILE:CD1	2.22	1.17
1:B:14:ILE:CG2	1:B:84:ILE:HD13	1.73	1.16
1:B:108:GLU:OE2	1:B:108:GLU:O	1.68	1.12
1:B:19:VAL:HG11	1:B:87:VAL:HG12	1.31	1.09
1:B:14:ILE:HG21	1:B:84:ILE:HD13	1.30	1.07
1:B:176:LYS:C	1:B:177:LEU:HD22	1.74	1.06
1:B:86:LEU:HD21	1:B:106:ILE:HD12	1.36	1.05
1:A:72:THR:HG22	2:F:189:LEU:HD12	1.39	1.02
1:A:72:THR:CG2	2:F:189:LEU:CD1	2.39	1.01
1:A:56:ARG:O	1:A:57:ILE:HD13	1.59	1.00
1:A:72:THR:CG2	2:F:189:LEU:HD12	1.93	0.98
1:A:90:ILE:CD1	1:A:123:CYS:HA	1.93	0.98
1:B:124:ASP:HB3	1:B:151:SER:OG	1.66	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:186:ARG:HD3	2:D:185:VAL:HG21	1.45	0.96
1:A:14:ILE:HG23	1:A:84:ILE:HD11	0.98	0.94
1:B:84:ILE:HD12	1:B:84:ILE:C	1.88	0.93
1:B:19:VAL:CG1	1:B:87:VAL:HG12	1.99	0.91
1:B:161:ALA:O	1:B:164:THR:HG22	1.73	0.88
1:B:72:THR:HG23	2:D:189:LEU:CD1	2.03	0.88
1:A:14:ILE:HG22	1:A:84:ILE:CD1	2.00	0.88
1:A:56:ARG:C	1:A:57:ILE:HD13	1.94	0.87
1:A:14:ILE:HG22	1:A:84:ILE:HD11	1.52	0.86
1:B:121:ASN:HA	1:B:150:THR:HG22	1.57	0.84
1:A:150:THR:CG2	1:A:151:SER:N	2.42	0.83
1:A:150:THR:HG23	1:A:151:SER:N	1.94	0.83
1:B:84:ILE:HD12	1:B:85:MET:N	1.95	0.82
1:B:117:MET:CE	1:B:119:LEU:HD11	2.11	0.81
1:A:89:ASP:OD1	1:A:91:THR:HB	1.80	0.81
2:C:186:ARG:CD	2:D:185:VAL:HG21	2.10	0.80
1:A:14:ILE:HG23	1:A:84:ILE:CD1	1.94	0.78
1:B:25:LEU:HD11	1:B:46:LYS:HB2	1.65	0.77
1:B:177:LEU:N	1:B:177:LEU:HD22	1.97	0.77
2:E:203:GLU:OE1	2:E:203:GLU:HA	1.84	0.77
1:B:117:MET:CE	1:B:161:ALA:HB1	2.14	0.77
1:B:14:ILE:CD1	1:B:106:ILE:HD11	2.14	0.76
1:B:14:ILE:HG21	1:B:84:ILE:CD1	2.13	0.76
1:A:72:THR:HG22	2:F:189:LEU:CD1	2.07	0.76
1:B:117:MET:HE2	1:B:119:LEU:HD11	1.67	0.76
1:A:121:ASN:OD1	1:A:150:THR:HG22	1.86	0.76
1:B:164:THR:CG2	1:B:165:LEU:N	2.48	0.76
1:B:153:LYS:HG3	3:B:201:GTP:C6	2.21	0.75
1:B:38:ILE:HD11	1:B:65:ALA:HB2	1.68	0.75
1:A:8:LEU:HD11	1:A:60:GLN:NE2	2.02	0.74
1:B:72:THR:HG22	1:B:73:ILE:N	2.00	0.74
1:A:93:GLU:OE2	1:A:135:ARG:NH2	2.21	0.74
1:B:25:LEU:HD22	1:B:25:LEU:O	1.87	0.73
2:C:193:LEU:CD2	2:D:193:LEU:HG	2.18	0.73
1:B:148:MET:HE1	1:B:164:THR:HG21	1.72	0.72
1:B:50:ILE:HG12	1:B:163:PHE:CZ	2.24	0.72
1:A:90:ILE:HD11	1:A:123:CYS:HA	1.72	0.72
1:B:24:VAL:HG13	1:B:158:VAL:HG12	1.71	0.71
1:B:79:ARG:CZ	2:C:186:ARG:HH21	2.04	0.71
1:A:21:LYS:HB2	3:A:201:GTP:O1B	1.91	0.71
1:B:72:THR:CG2	1:B:73:ILE:N	2.54	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:79:ARG:NH2	2:C:186:ARG:HE	1.88	0.70
1:A:72:THR:CG2	2:F:189:LEU:HD13	2.20	0.70
1:B:108:GLU:OE2	1:B:108:GLU:C	2.29	0.70
1:A:12:LEU:O	1:A:84:ILE:HD12	1.92	0.70
1:B:27:ARG:CZ	1:B:155:ASN:HB2	2.22	0.69
1:A:29:SER:CB	1:A:46:LYS:HG2	2.22	0.69
1:A:19:VAL:CG1	1:A:87:VAL:HG12	2.22	0.69
1:A:63:ASP:C	1:A:63:ASP:OD1	2.29	0.69
2:C:215:GLN:HG3	2:D:214:LYS:HB2	1.73	0.69
1:B:124:ASP:C	1:B:124:ASP:OD2	2.29	0.68
1:B:117:MET:HE1	1:B:148:MET:HE1	1.76	0.68
1:B:72:THR:CG2	2:D:189:LEU:HD12	2.23	0.68
1:A:7:TYR:CD2	1:A:57:ILE:HD12	2.29	0.68
1:B:66:GLY:O	1:B:67:GLN:C	2.31	0.68
1:A:27:ARG:CZ	1:A:155:ASN:HB2	2.24	0.68
1:B:47:ILE:HD12	2:C:202:GLU:OE2	1.93	0.67
1:B:14:ILE:CG2	1:B:84:ILE:CD1	2.62	0.67
1:A:124:ASP:HB3	1:A:151:SER:OG	1.94	0.67
1:B:86:LEU:HD21	1:B:106:ILE:CD1	2.19	0.67
1:B:106:ILE:O	1:B:110:ALA:HB3	1.95	0.67
1:A:106:ILE:O	1:A:110:ALA:HB3	1.95	0.67
1:A:86:LEU:HD21	1:A:106:ILE:HD12	1.76	0.66
1:B:121:ASN:OD1	1:B:150:THR:CG2	2.43	0.66
1:A:118:ILE:HG13	1:A:145:ILE:HD11	1.77	0.66
1:B:72:THR:CG2	2:D:189:LEU:CD1	2.73	0.66
2:C:215:GLN:HG3	2:D:214:LYS:CB	2.26	0.66
1:A:132:SER:OG	1:A:135:ARG:HB2	1.95	0.66
1:A:93:GLU:O	1:A:96:PHE:N	2.28	0.66
1:B:27:ARG:NE	1:B:155:ASN:HB2	2.11	0.66
1:B:13:LEU:O	1:B:64:THR:HG22	1.96	0.66
1:A:14:ILE:HG22	1:A:84:ILE:HD13	1.77	0.65
1:B:164:THR:HG22	1:B:165:LEU:N	2.12	0.65
1:B:88:TYR:HB3	1:B:99:ILE:HD11	1.77	0.65
1:B:164:THR:HG22	1:B:165:LEU:H	1.61	0.64
1:B:31:ASP:OD2	1:B:46:LYS:CE	2.46	0.64
1:B:159:GLU:O	1:B:163:PHE:HD1	1.81	0.64
1:A:22:THR:N	3:A:201:GTP:O1B	2.28	0.64
2:C:222:LEU:HD23	2:D:222:LEU:HG	1.78	0.63
1:A:121:ASN:OD1	1:A:150:THR:CG2	2.45	0.63
1:A:66:GLY:O	1:A:67:GLN:C	2.31	0.63
1:A:72:THR:HB	2:F:192:GLU:OE2	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:148:MET:CE	1:B:164:THR:HG21	2.29	0.63
1:A:90:ILE:HD13	1:A:123:CYS:HA	1.76	0.63
1:A:27:ARG:NE	1:A:155:ASN:HB2	2.14	0.62
1:B:117:MET:HE1	1:B:161:ALA:HB1	1.80	0.62
1:A:72:THR:HB	2:F:192:GLU:CD	2.20	0.62
1:A:101:ASN:O	1:A:104:ARG:HB2	1.98	0.61
1:B:19:VAL:HG11	1:B:87:VAL:CG1	2.21	0.61
1:A:38:ILE:HD11	1:A:65:ALA:HB2	1.81	0.61
2:C:164:GLU:O	2:C:168:ALA:N	2.24	0.61
1:A:72:THR:HG23	2:F:189:LEU:CD1	2.30	0.61
1:A:6:ASP:OD2	1:A:55:LYS:HB3	2.02	0.60
1:B:72:THR:HG23	2:D:189:LEU:HD12	1.79	0.60
1:A:19:VAL:HG11	1:A:87:VAL:HG12	1.83	0.60
1:A:20:GLY:HA2	3:A:201:GTP:PA	2.42	0.59
2:C:193:LEU:HD21	2:D:193:LEU:HG	1.84	0.59
2:E:175:LYS:NZ	2:F:176:ASP:CB	2.66	0.59
1:B:117:MET:HE3	1:B:119:LEU:HD11	1.84	0.59
1:B:121:ASN:OD1	1:B:150:THR:HG22	2.03	0.59
1:B:176:LYS:O	1:B:177:LEU:C	2.41	0.59
2:E:170:ARG:O	2:E:174:LEU:N	2.30	0.58
1:A:163:PHE:O	1:A:167:ARG:HG3	2.04	0.58
1:B:122:LYS:NZ	3:B:201:GTP:O4'	2.34	0.58
1:A:36:THR:OG1	1:A:37:PHE:N	2.34	0.58
1:B:117:MET:CE	1:B:148:MET:HE1	2.34	0.58
2:E:181:ARG:O	2:E:185:VAL:HG13	2.03	0.58
1:B:96:PHE:CG	1:B:131:VAL:HG11	2.39	0.58
1:B:118:ILE:HG13	1:B:145:ILE:HD11	1.85	0.57
1:B:7:TYR:HB3	1:B:57:ILE:HD13	1.86	0.57
1:A:122:LYS:HG2	3:A:201:GTP:C6	2.39	0.57
1:A:50:ILE:HD13	1:A:59:LEU:HD11	1.86	0.57
1:B:117:MET:HE1	1:B:161:ALA:CB	2.35	0.57
1:B:117:MET:HE1	1:B:161:ALA:CA	2.34	0.57
1:B:122:LYS:HB3	1:B:125:VAL:CG1	2.35	0.57
2:E:175:LYS:HZ2	2:F:176:ASP:CB	2.17	0.57
1:B:124:ASP:HB3	1:B:151:SER:CB	2.36	0.56
1:A:45:PHE:CE1	1:A:47:ILE:HD12	2.40	0.56
1:B:72:THR:HG23	2:D:189:LEU:HD13	1.84	0.56
1:A:68:GLU:OE1	1:A:69:ARG:N	2.36	0.55
1:B:68:GLU:OE1	1:B:69:ARG:N	2.36	0.55
2:C:164:GLU:O	2:C:168:ALA:CB	2.54	0.55
1:B:31:ASP:OD2	1:B:46:LYS:HE3	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:PHE:CD2	1:B:57:ILE:HD11	2.42	0.55
1:A:72:THR:HB	2:F:192:GLU:OE1	2.06	0.55
1:A:122:LYS:HD3	3:A:201:GTP:C4	2.42	0.55
1:B:126:ASN:ND2	1:B:129:ARG:HH11	2.05	0.55
1:B:29:SER:O	1:B:30:GLU:HB2	2.07	0.55
1:B:124:ASP:HB3	1:B:151:SER:HG	1.70	0.54
2:E:211:ALA:HB2	2:F:208:VAL:HG13	1.88	0.54
1:A:117:MET:HE1	1:A:148:MET:SD	2.46	0.54
1:A:5:TYR:CD1	1:A:58:LYS:HE3	2.43	0.54
1:B:155:ASN:CG	1:B:155:ASN:O	2.46	0.54
1:B:126:ASN:HD21	1:B:129:ARG:HH11	1.54	0.54
1:B:19:VAL:CG1	1:B:19:VAL:O	2.51	0.54
1:B:159:GLU:O	1:B:163:PHE:CD1	2.60	0.54
1:B:177:LEU:H	1:B:177:LEU:HD13	1.73	0.54
1:B:72:THR:HG22	1:B:73:ILE:H	1.71	0.54
1:B:8:LEU:HD12	1:B:58:LYS:CB	2.38	0.53
1:A:34:ASN:H	1:A:34:ASN:ND2	2.07	0.53
1:B:34:ASN:ND2	1:B:34:ASN:H	2.07	0.53
1:A:12:LEU:HG	1:A:84:ILE:HD13	1.89	0.53
1:A:155:ASN:CG	1:A:155:ASN:O	2.46	0.53
1:A:5:TYR:HD1	1:A:58:LYS:HE3	1.74	0.53
1:B:117:MET:HE2	1:B:161:ALA:HB1	1.88	0.53
1:B:14:ILE:HG23	1:B:84:ILE:HD13	1.83	0.53
1:A:150:THR:HG23	1:A:151:SER:H	1.73	0.53
1:B:82:MET:O	1:B:114:VAL:HG23	2.09	0.53
1:B:124:ASP:OD2	1:B:125:VAL:N	2.42	0.53
1:A:148:MET:HG2	1:A:157:ASN:OD1	2.08	0.53
1:B:114:VAL:O	1:B:114:VAL:HG13	2.09	0.53
1:B:90:ILE:CD1	1:B:123:CYS:HA	2.38	0.53
1:A:44:ASP:HA	2:E:201:PHE:CE2	2.44	0.52
1:A:26:PHE:CZ	1:A:36:THR:HB	2.45	0.52
1:B:148:MET:SD	1:B:157:ASN:OD1	2.67	0.52
1:B:164:THR:HG23	1:B:165:LEU:N	2.23	0.52
1:B:38:ILE:CD1	1:B:65:ALA:HB2	2.39	0.52
2:C:168:ALA:HB1	2:D:168:ALA:HB1	1.92	0.52
1:A:72:THR:HG23	2:F:189:LEU:HD13	1.92	0.52
1:A:140:ALA:HB2	1:A:147:PHE:HB2	1.92	0.52
1:A:177:LEU:HD22	1:A:177:LEU:N	2.25	0.52
1:B:177:LEU:N	1:B:177:LEU:HD13	2.25	0.52
2:E:213:ILE:O	2:E:216:ALA:HB3	2.10	0.52
1:A:114:VAL:HG13	1:A:114:VAL:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:86:LEU:HD11	1:B:106:ILE:CD1	2.39	0.51
2:C:186:ARG:O	2:C:187:ASP:C	2.48	0.51
1:B:50:ILE:HG12	1:B:163:PHE:HZ	1.76	0.51
1:A:122:LYS:HD3	3:A:201:GTP:C2	2.45	0.51
1:B:21:LYS:HB2	3:B:201:GTP:O1B	2.11	0.51
1:A:122:LYS:C	1:A:124:ASP:H	2.13	0.50
1:A:14:ILE:CG2	1:A:84:ILE:HD13	2.30	0.50
1:A:19:VAL:O	1:A:19:VAL:HG13	2.10	0.50
1:B:159:GLU:HG3	1:B:163:PHE:CE1	2.46	0.50
1:A:154:ALA:HB3	1:A:156:ILE:HG13	1.94	0.50
1:A:34:ASN:HD22	1:A:34:ASN:H	1.59	0.50
1:B:116:LYS:HZ3	1:B:116:LYS:HB2	1.77	0.50
1:A:133:LYS:HD3	1:A:149:GLU:CD	2.32	0.50
1:A:36:THR:OG1	1:A:38:ILE:HG22	2.12	0.50
1:B:14:ILE:HD13	1:B:86:LEU:HG	1.94	0.50
1:B:152:ALA:N	3:B:201:GTP:O6	2.45	0.50
1:B:86:LEU:CD2	1:B:106:ILE:HD12	2.26	0.50
1:B:86:LEU:HD11	1:B:106:ILE:HD13	1.94	0.50
1:B:154:ALA:HB3	1:B:156:ILE:HG13	1.94	0.50
1:A:118:ILE:CG1	1:A:145:ILE:HD11	2.41	0.50
1:A:19:VAL:HG22	1:A:89:ASP:N	2.27	0.50
1:B:98:ASN:O	1:B:99:ILE:C	2.49	0.50
1:B:72:THR:HG22	2:D:192:GLU:OE2	2.11	0.50
1:B:12:LEU:HD12	1:B:62:TRP:O	2.12	0.49
2:F:185:VAL:O	2:F:186:ARG:C	2.51	0.49
1:B:117:MET:HE1	1:B:161:ALA:HA	1.93	0.49
2:C:215:GLN:HA	2:D:215:GLN:HG2	1.94	0.49
2:C:186:ARG:HG2	2:D:185:VAL:HG11	1.95	0.49
1:A:108:GLU:OE2	1:A:109:HIS:ND1	2.46	0.49
1:B:34:ASN:HD22	1:B:34:ASN:H	1.59	0.49
2:E:222:LEU:O	2:E:226:GLN:CB	2.61	0.49
2:E:175:LYS:O	2:E:179:CYS:N	2.36	0.49
1:B:117:MET:CE	1:B:148:MET:CE	2.90	0.48
1:B:19:VAL:HG12	1:B:19:VAL:O	2.11	0.48
1:B:14:ILE:HD13	1:B:84:ILE:HD11	1.96	0.48
1:B:79:ARG:NH2	2:C:186:ARG:HH21	2.12	0.48
1:B:122:LYS:HB3	1:B:125:VAL:HG13	1.96	0.48
1:A:122:LYS:HD3	3:A:201:GTP:N3	2.29	0.48
1:A:7:TYR:CD2	1:A:57:ILE:CD1	2.97	0.48
1:B:21:LYS:NZ	3:B:201:GTP:O2B	2.36	0.48
1:B:21:LYS:HZ1	3:B:201:GTP:PB	2.35	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:ILE:O	1:A:76:ALA:HB3	2.14	0.47
1:B:83:GLY:HA2	1:B:115:GLU:O	2.14	0.47
1:A:123:CYS:SG	1:A:150:THR:O	2.71	0.47
1:A:26:PHE:HZ	1:A:36:THR:HB	1.79	0.47
1:A:14:ILE:HD11	1:A:102:TRP:HB3	1.97	0.47
1:A:92:ASN:O	1:A:95:SER:OG	2.33	0.47
1:A:129:ARG:NH2	1:A:149:GLU:OE2	2.48	0.47
2:C:217:THR:O	2:C:220:LYS:N	2.47	0.47
1:A:34:ASN:O	1:A:36:THR:HG22	2.14	0.47
1:A:96:PHE:CZ	1:A:139:LEU:HD12	2.50	0.47
1:A:7:TYR:HD2	1:A:57:ILE:CD1	2.28	0.46
1:A:89:ASP:CG	1:A:91:THR:HB	2.35	0.46
1:B:14:ILE:HD13	1:B:106:ILE:HD11	1.94	0.46
1:B:50:ILE:HD12	1:B:51:GLU:H	1.81	0.46
2:C:222:LEU:HD11	2:D:221:GLN:HB2	1.97	0.46
1:A:29:SER:OG	1:A:46:LYS:HG2	2.15	0.46
2:C:217:THR:O	2:C:218:ALA:C	2.53	0.46
1:A:19:VAL:O	1:A:19:VAL:CG1	2.61	0.46
1:A:100:ARG:O	1:A:103:ILE:HB	2.16	0.46
1:A:10:LYS:H	1:A:82:MET:CE	2.29	0.46
1:A:125:VAL:O	1:A:125:VAL:HG22	2.15	0.46
1:B:104:ARG:O	1:B:105:ASN:C	2.55	0.46
1:A:122:LYS:O	1:A:124:ASP:N	2.49	0.45
1:B:147:PHE:C	1:B:147:PHE:CD2	2.90	0.45
2:E:193:LEU:HD13	2:F:189:LEU:HG	1.98	0.45
1:A:96:PHE:CE1	1:A:139:LEU:HD12	2.51	0.45
1:B:43:ILE:HG22	1:B:44:ASP:N	2.31	0.45
1:A:84:ILE:HD12	1:A:85:MET:N	2.32	0.45
1:B:79:ARG:HH21	2:C:186:ARG:HE	1.62	0.45
1:B:84:ILE:HD11	1:B:106:ILE:HD11	1.99	0.45
1:B:153:LYS:HG3	3:B:201:GTP:N1	2.30	0.45
1:B:124:ASP:CB	1:B:151:SER:OG	2.51	0.45
1:B:15:GLY:O	1:B:21:LYS:HE2	2.17	0.45
2:E:181:ARG:HB2	2:E:181:ARG:CZ	2.46	0.44
1:A:104:ARG:O	1:A:105:ASN:C	2.55	0.44
1:B:12:LEU:HB3	1:B:84:ILE:HB	1.99	0.44
1:B:84:ILE:HD11	1:B:86:LEU:HG	1.98	0.44
1:B:87:VAL:HG22	1:B:119:LEU:HB2	2.00	0.44
1:B:93:GLU:OE1	1:B:97:ASP:OD1	2.35	0.44
2:C:211:ALA:HB2	2:D:208:VAL:HG13	1.98	0.44
2:F:196:LEU:HD11	2:F:200:LEU:CD2	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:ILE:CG2	1:B:48:ARG:N	2.79	0.44
1:B:9:PHE:HA	1:B:82:MET:SD	2.58	0.44
1:A:114:VAL:HG13	1:A:116:LYS:HG3	2.00	0.44
1:A:45:PHE:HZ	1:A:60:GLN:HG2	1.83	0.44
2:F:200:LEU:HA	2:F:200:LEU:HD13	1.76	0.44
2:C:206:LYS:O	2:C:210:GLU:HG3	2.18	0.44
1:A:47:ILE:HD13	2:E:202:GLU:OE2	2.18	0.43
1:B:139:LEU:O	1:B:142:ASP:HB2	2.18	0.43
1:B:21:LYS:NZ	3:B:201:GTP:PB	2.91	0.43
2:D:189:LEU:HD12	2:D:189:LEU:HA	1.75	0.43
1:B:121:ASN:CA	1:B:150:THR:HG22	2.40	0.43
1:B:121:ASN:OD1	1:B:150:THR:HG23	2.18	0.43
1:B:27:ARG:HD3	1:B:152:ALA:O	2.19	0.43
2:E:180:GLU:O	2:E:184:LYS:HB2	2.19	0.43
1:B:18:GLY:N	3:B:201:GTP:O2G	2.48	0.43
1:B:35:SER:HB3	2:D:207:MET:SD	2.59	0.43
1:A:92:ASN:HD21	1:A:94:LYS:NZ	2.17	0.43
1:A:140:ALA:HB1	1:A:145:ILE:O	2.19	0.42
1:B:98:ASN:O	1:B:101:ASN:N	2.52	0.42
1:B:86:LEU:HD22	1:B:103:ILE:HG13	2.02	0.42
3:A:201:GTP:O3A	3:A:201:GTP:O2G	2.37	0.42
1:A:19:VAL:HG13	1:A:87:VAL:HG12	1.99	0.42
1:A:91:THR:CG2	1:A:91:THR:O	2.68	0.42
1:B:122:LYS:O	1:B:125:VAL:HG13	2.20	0.42
1:B:20:GLY:HA2	3:B:201:GTP:PA	2.59	0.42
1:A:150:THR:HG22	1:A:151:SER:N	2.26	0.42
1:A:117:MET:HE2	1:A:161:ALA:HB1	2.00	0.42
1:B:10:LYS:O	1:B:11:LEU:HD12	2.20	0.42
1:B:91:THR:O	1:B:91:THR:CG2	2.68	0.42
1:A:117:MET:HE3	1:A:148:MET:HB2	2.02	0.41
1:A:145:ILE:HD13	1:A:145:ILE:HG21	1.84	0.41
1:B:114:VAL:CG1	1:B:114:VAL:O	2.68	0.41
1:B:75:THR:HG22	1:B:76:ALA:N	2.33	0.41
2:D:196:LEU:O	2:D:196:LEU:HD12	2.20	0.41
1:A:176:LYS:O	1:A:176:LYS:CG	2.68	0.41
1:A:176:LYS:O	1:A:176:LYS:HG3	2.20	0.41
1:B:27:ARG:CD	1:B:152:ALA:O	2.69	0.41
1:A:22:THR:HB	3:A:201:GTP:O2A	2.21	0.41
1:A:98:ASN:O	1:A:101:ASN:HB2	2.21	0.41
1:B:148:MET:HE1	1:B:161:ALA:HA	2.01	0.41
1:A:20:GLY:HA3	3:A:201:GTP:C8	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:212:ASN:O	2:C:213:ILE:C	2.58	0.41
2:F:188:GLN:O	2:F:189:LEU:C	2.58	0.41
2:E:165:LEU:O	2:E:169:GLN:CB	2.69	0.41
1:B:114:VAL:HG13	1:B:116:LYS:HG3	2.01	0.41
1:B:153:LYS:HG3	3:B:201:GTP:C5	2.53	0.41
1:B:176:LYS:O	1:B:177:LEU:HD22	2.14	0.41
2:C:164:GLU:O	2:C:168:ALA:HB3	2.20	0.41
1:A:43:ILE:HG22	2:E:201:PHE:CZ	2.55	0.41
1:A:5:TYR:HD2	1:A:7:TYR:N	2.18	0.41
1:B:73:ILE:HG12	2:C:193:LEU:HD11	2.02	0.41
1:A:73:ILE:CD1	2:F:192:GLU:HB3	2.51	0.41
2:F:227:GLY:O	2:F:231:VAL:N	2.54	0.41
1:A:114:VAL:O	1:A:114:VAL:CG1	2.68	0.40
2:C:186:ARG:HD2	2:D:185:VAL:HG11	2.02	0.40
1:A:122:LYS:C	1:A:124:ASP:N	2.74	0.40
1:B:50:ILE:O	1:B:57:ILE:HG23	2.20	0.40
1:A:27:ARG:HD3	1:A:152:ALA:O	2.21	0.40
1:B:122:LYS:O	1:B:123:CYS:C	2.60	0.40
1:B:31:ASP:OD2	1:B:46:LYS:HE2	2.21	0.40
1:A:92:ASN:C	1:A:92:ASN:OD1	2.59	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	173/186 (93%)	164 (95%)	8 (5%)	1 (1%)	25	64
1	B	173/186 (93%)	161 (93%)	11 (6%)	1 (1%)	25	64
2	C	70/78 (90%)	62 (89%)	7 (10%)	1 (1%)	11	46
2	D	69/78 (88%)	68 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	E	74/78 (95%)	72 (97%)	2 (3%)	0	100	100
2	F	74/78 (95%)	69 (93%)	5 (7%)	0	100	100
All	All	633/684 (92%)	596 (94%)	34 (5%)	3 (0%)	29	67

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	66	GLY
1	B	66	GLY
2	C	213	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	146/159 (92%)	119 (82%)	27 (18%)	1	8
1	B	146/159 (92%)	112 (77%)	34 (23%)	1	3
2	C	36/66 (54%)	31 (86%)	5 (14%)	3	16
2	D	35/66 (53%)	32 (91%)	3 (9%)	10	38
2	E	35/66 (53%)	27 (77%)	8 (23%)	1	4
2	F	37/66 (56%)	30 (81%)	7 (19%)	1	8
All	All	435/582 (75%)	351 (81%)	84 (19%)	1	8

All (84) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LEU
1	A	14	ILE
1	A	24	VAL
1	A	25	LEU
1	A	27	ARG
1	A	34	ASN
1	A	35	SER

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Mol	Chain	Res	Type
1	A	36	THR
1	A	47	ILE
1	A	67	GLN
1	A	68	GLU
1	A	84	ILE
1	A	104	ARG
1	A	108	GLU
1	A	113	ASP
1	A	124	ASP
1	A	125	VAL
1	A	139	LEU
1	A	141	LEU
1	A	145	ILE
1	A	147	PHE
1	A	150	THR
1	A	157	ASN
1	A	158	VAL
1	A	165	LEU
1	A	175	LYS
1	A	177	LEU
1	B	14	ILE
1	B	25	LEU
1	B	27	ARG
1	B	31	ASP
1	B	34	ASN
1	B	47	ILE
1	B	50	ILE
1	B	57	ILE
1	B	67	GLN
1	B	68	GLU
1	B	72	THR
1	B	75	THR
1	B	79	ARG
1	B	82	MET
1	B	84	ILE
1	B	85	MET
1	B	100	ARG
1	B	104	ARG
1	B	108	GLU
1	B	113	ASP
1	B	124	ASP
1	B	125	VAL

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Mol	Chain	Res	Type
1	B	126	ASN
1	B	131	VAL
1	B	138	LYS
1	B	139	LEU
1	B	145	ILE
1	B	147	PHE
1	B	150	THR
1	B	157	ASN
1	B	158	VAL
1	B	164	THR
1	B	165	LEU
1	B	177	LEU
2	C	183	SER
2	C	188	GLN
2	C	189	LEU
2	C	192	GLU
2	C	199	SER
2	D	189	LEU
2	D	193	LEU
2	D	199	SER
2	E	178	GLU
2	E	179	CYS
2	E	180	GLU
2	E	181	ARG
2	E	194	GLU
2	E	195	GLU
2	E	196	LEU
2	E	203	GLU
2	F	179	CYS
2	F	186	ARG
2	F	189	LEU
2	F	191	GLN
2	F	193	LEU
2	F	200	LEU
2	F	206	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	ASN
1	A	60	GLN
1	A	67	GLN

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Mol	Chain	Res	Type
1	B	34	ASN
1	B	67	GLN
1	B	101	ASN
1	B	126	ASN
2	F	188	GLN
2	F	215	GLN
2	F	221	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

2 ligands 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	GTP	B	201	-	26,34,34	1.12	1 (3%)	33,54,54	1.95	8 (24%)
3	GTP	A	201	-	26,34,34	1.01	2 (7%)	33,54,54	1.72	7 (21%)

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	B	201	-	-	1/18/38/38	0/3/3/3
3	GTP	A	201	-	-	2/18/38/38	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	201	GTP	C6-N1	4.01	1.40	1.33
3	A	201	GTP	C6-C5	-2.56	1.36	1.41
3	A	201	GTP	C6-N1	2.01	1.36	1.33

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	201	GTP	PB-O3B-PG	-5.19	115.02	132.83
3	B	201	GTP	N3-C2-N1	-5.03	120.51	127.22
3	A	201	GTP	N3-C2-N1	-4.24	121.57	127.22
3	B	201	GTP	C2-N3-C4	3.86	119.77	115.36
3	B	201	GTP	PA-O3A-PB	-3.83	119.70	132.83
3	B	201	GTP	O2G-PG-O3B	3.68	116.96	104.64
3	B	201	GTP	C5-C6-N1	-3.52	118.62	123.43
3	B	201	GTP	PB-O3B-PG	-3.50	120.83	132.83
3	A	201	GTP	PA-O3A-PB	-3.34	121.36	132.83
3	B	201	GTP	C6-N1-C2	2.59	120.05	115.93
3	A	201	GTP	C2-N3-C4	2.50	118.21	115.36
3	A	201	GTP	C5-C6-N1	-2.46	120.07	123.43
3	B	201	GTP	N2-C2-N3	2.17	121.33	117.79
3	A	201	GTP	C6-N1-C2	2.15	119.34	115.93
3	A	201	GTP	O3G-PG-O2G	2.05	115.47	107.64

There are no chirality outliers.

All (3) torsion outliers are listed below:

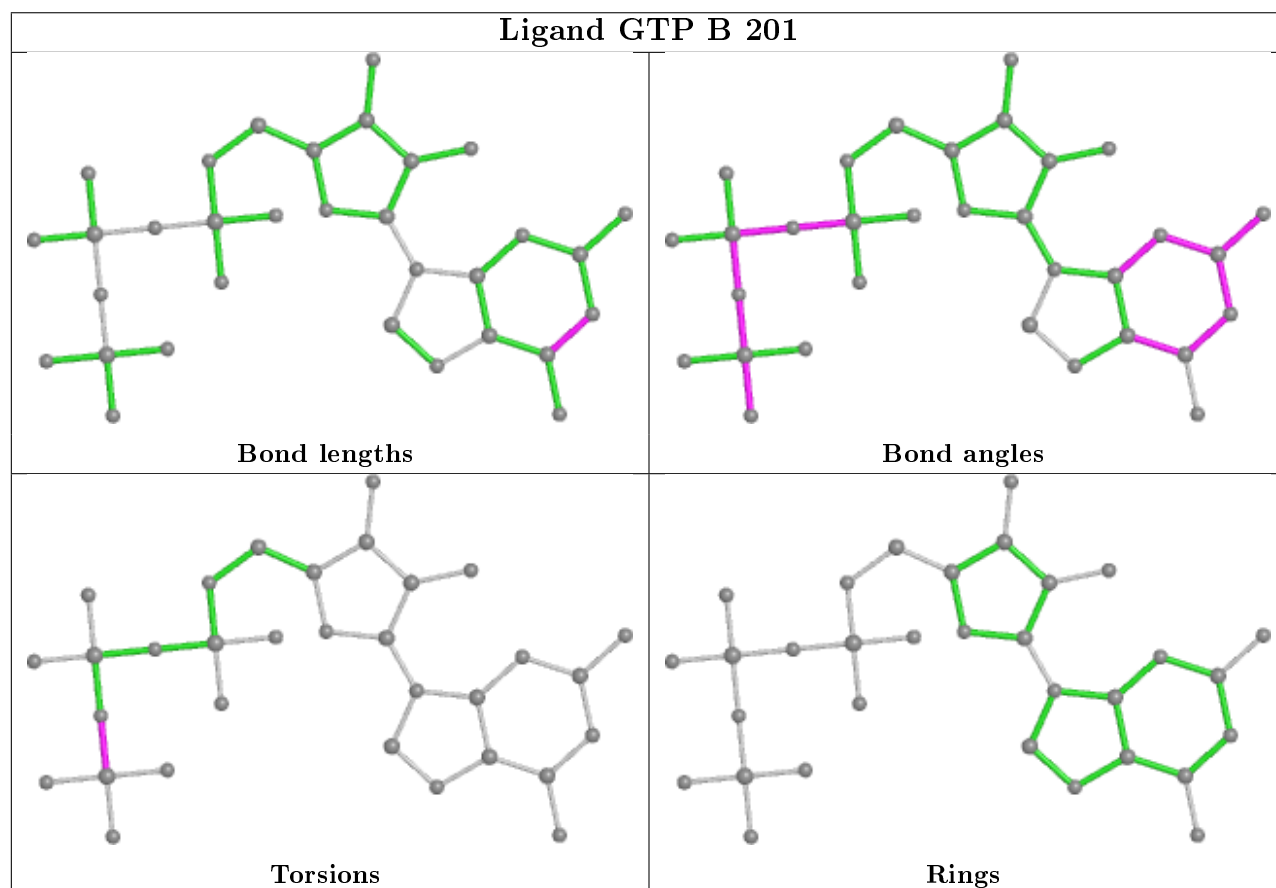
Mol	Chain	Res	Type	Atoms
3	A	201	GTP	PB-O3B-PG-O2G
3	B	201	GTP	PB-O3B-PG-O1G
3	A	201	GTP	PB-O3B-PG-O1G

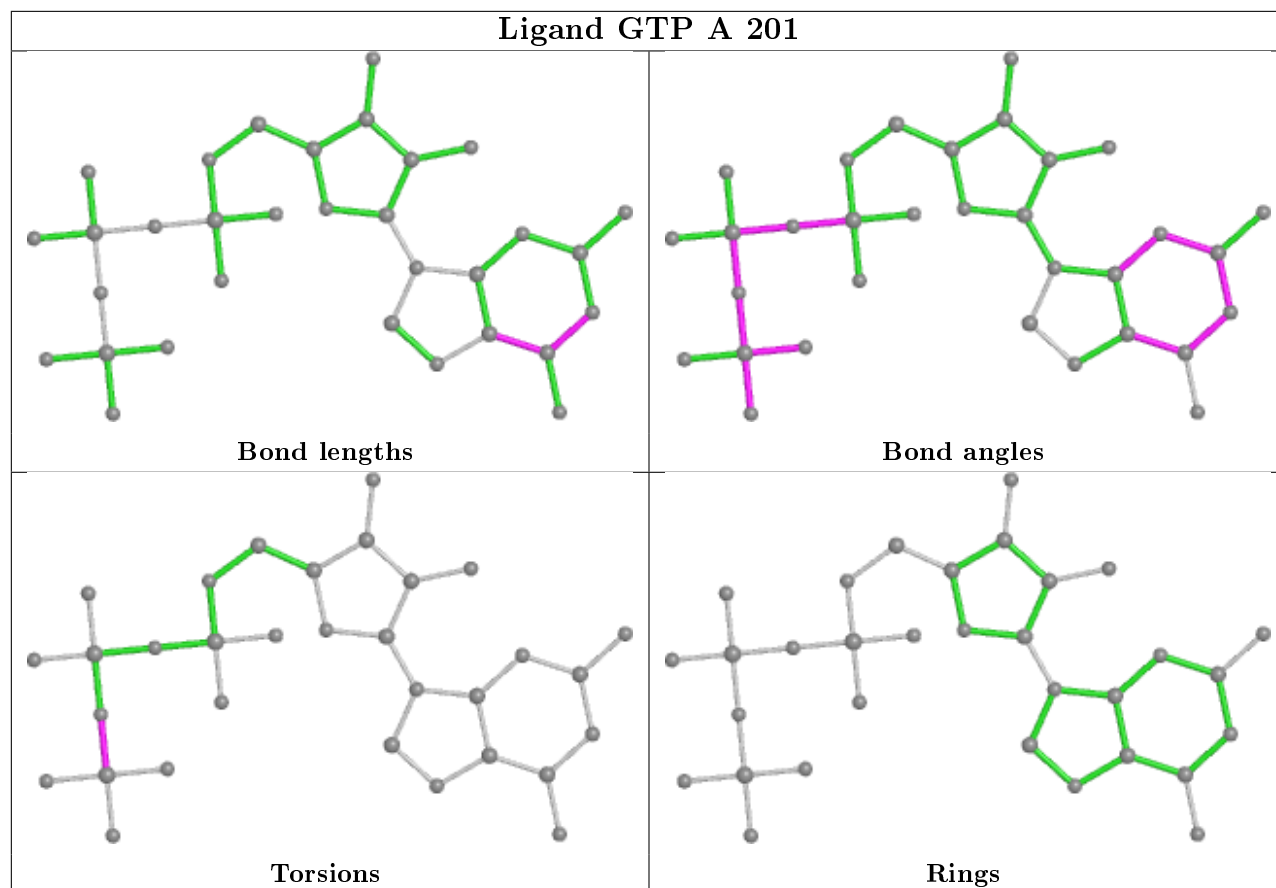
There are no ring outliers.

2 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	201	GTP	11	0
3	A	201	GTP	10	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.





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	175/186 (94%)	-0.05	1 (0%) 89 83	37, 67, 121, 154	0
1	B	175/186 (94%)	-0.03	1 (0%) 89 83	41, 74, 128, 174	0
2	C	72/78 (92%)	0.34	6 (8%) 11 6	44, 110, 181, 209	0
2	D	71/78 (91%)	0.52	10 (14%) 2 2	38, 117, 188, 202	0
2	E	76/78 (97%)	0.43	10 (13%) 3 2	40, 115, 236, 276	0
2	F	76/78 (97%)	0.76	17 (22%) 0 0	40, 126, 245, 265	0
All	All	645/684 (94%)	0.21	45 (6%) 16 9	37, 81, 201, 276	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	157	GLY	7.3
2	F	160	ARG	6.8
2	F	161	LEU	5.3
2	F	159	GLU	5.3
2	E	232	LEU	5.0
2	F	228	LYS	4.6
2	D	159	GLU	4.6
2	F	162	LYS	4.5
2	D	161	LEU	4.0
2	C	158	TYR	4.0
2	D	227	GLY	4.0
2	F	225	ALA	3.8
2	E	229	ILE	3.8
2	D	164	GLU	3.7
2	F	163	GLU	3.6
2	D	160	ARG	3.4
2	C	166	ALA	3.3
2	E	227	GLY	3.1
2	F	224	GLU	3.1

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Mol	Chain	Res	Type	RSRZ
2	E	163	GLU	3.0
2	D	157	GLY	2.9
2	C	159	GLU	2.9
2	F	230	ASP	2.9
2	E	160	ARG	2.9
2	D	158	TYR	2.9
2	E	166	ALA	2.9
2	D	162	LYS	2.8
2	E	159	GLU	2.8
2	D	224	GLU	2.8
2	F	166	ALA	2.7
2	F	169	GLN	2.7
1	A	34	ASN	2.7
1	B	68	GLU	2.5
2	D	166	ALA	2.5
2	E	228	LYS	2.5
2	F	157	GLY	2.4
2	C	168	ALA	2.4
2	F	231	VAL	2.4
2	E	157	GLY	2.3
2	F	226	GLN	2.3
2	E	230	ASP	2.3
2	F	229	ILE	2.2
2	F	168	ALA	2.1
2	F	165	LEU	2.0
2	C	169	GLN	2.0

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](#)

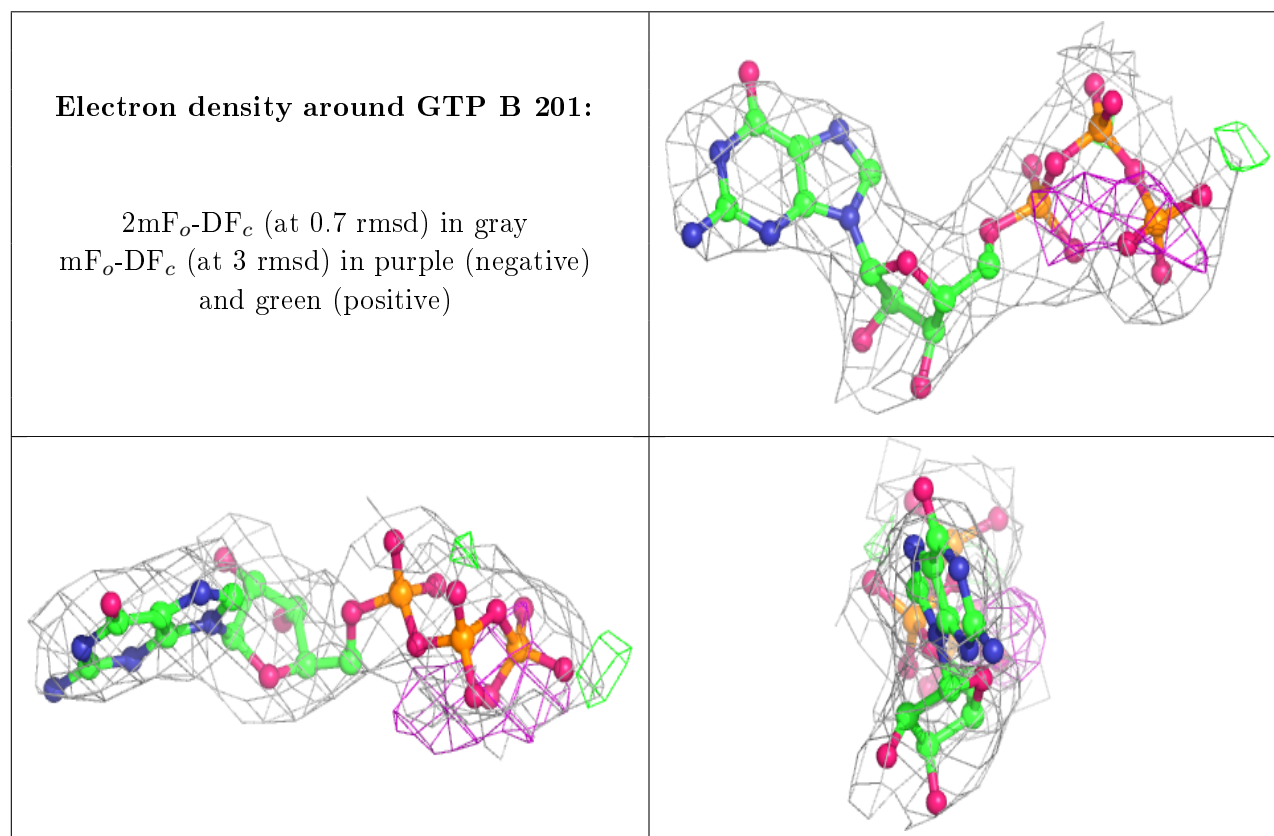
There are no carbohydrates in this entry.

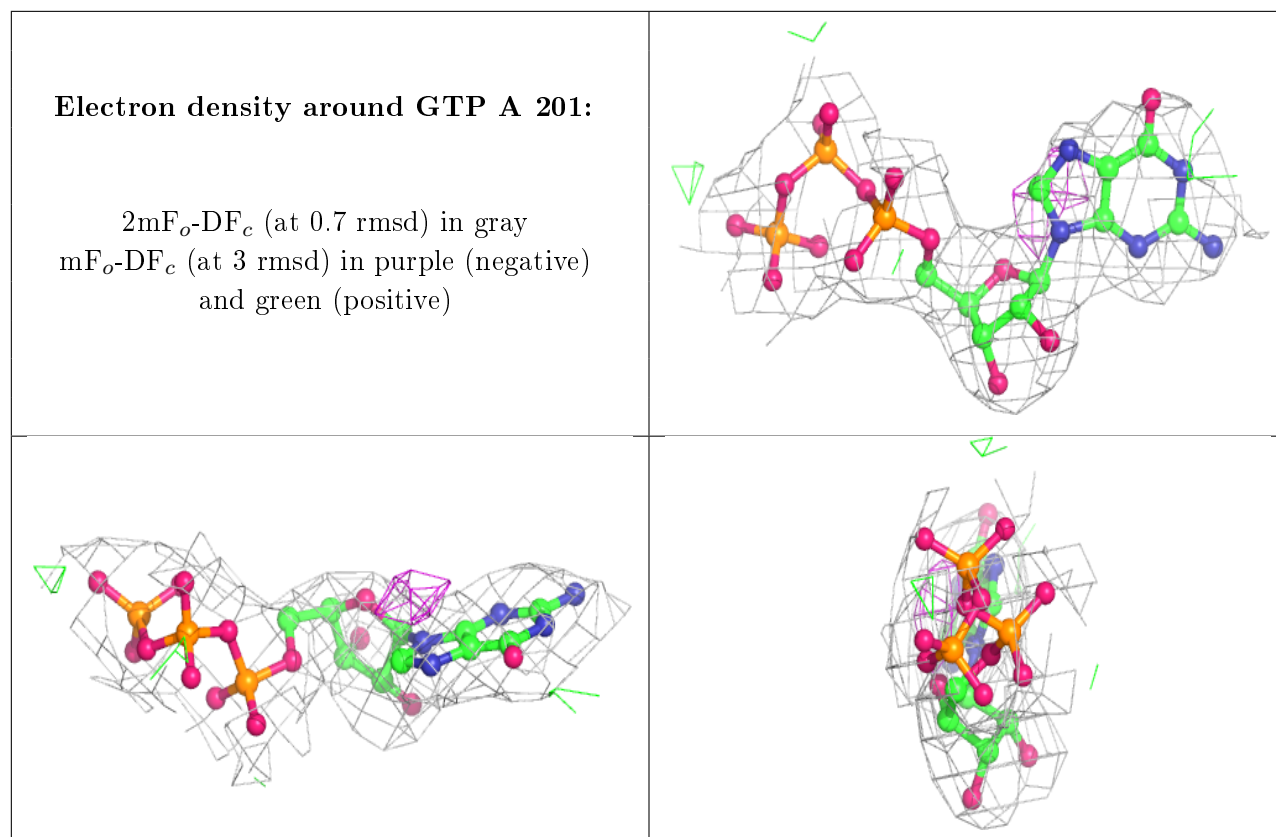
6.4 Ligands [i](#)

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(\AA^2)	Q<0.9
3	GTP	B	201	32/32	0.94	0.23	55,75,77,78	0
3	GTP	A	201	32/32	0.96	0.15	47,51,59,61	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.