



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:55 AM UTC

PDB ID : 3EA5 / pdb\_00003ea5  
Title : Kap95p Binding Induces the Switch Loops of RanGDP to adopt the GTP-bound Conformation: Implications for Nuclear Import Complex Assembly Dynamics  
Authors : Forwood, J.K.; Lonhienne, J.K.; Guncar, G.; Stewart, M.; Marfori, M.; Kobe, B.  
Deposited on : 2008-08-24  
Resolution : 2.50 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](mailto: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>.

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The following versions of software and data (see [references](#) ①) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtrriage (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

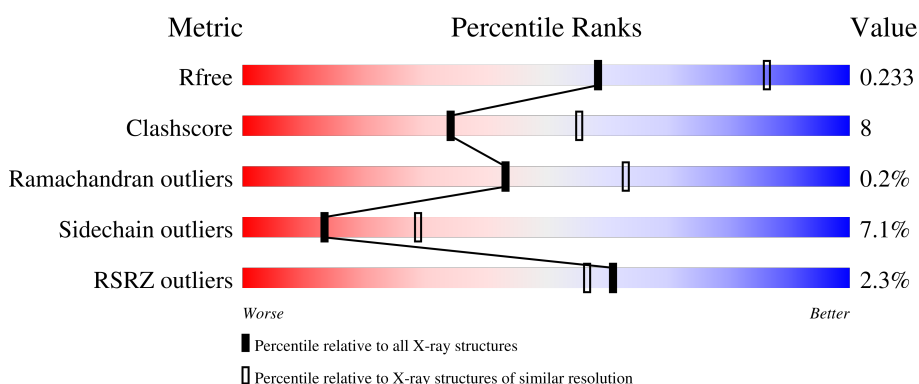
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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  $\geq 3$ , 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	216	 64% 13% 19% 2%
1	C	216	 67% 10% 22% 2%
2	B	861	 82% 15% 3%
2	D	861	 83% 13% 2%

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:

<b>Mol</b>	<b>Type</b>	<b>Chain</b>	<b>Res</b>	<b>Chirality</b>	<b>Geometry</b>	<b>Clashes</b>	<b>Electron density</b>
4	MG	A	222	-	-	X	-

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 16428 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 GTP-binding nuclear protein Ran.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	174	1415	918	248	244	5	0	0	0
1	C	169	1377	895	242	236	4	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	181	CYS	ALA	engineered mutation	UNP P62826
C	181	CYS	ALA	engineered mutation	UNP P62826

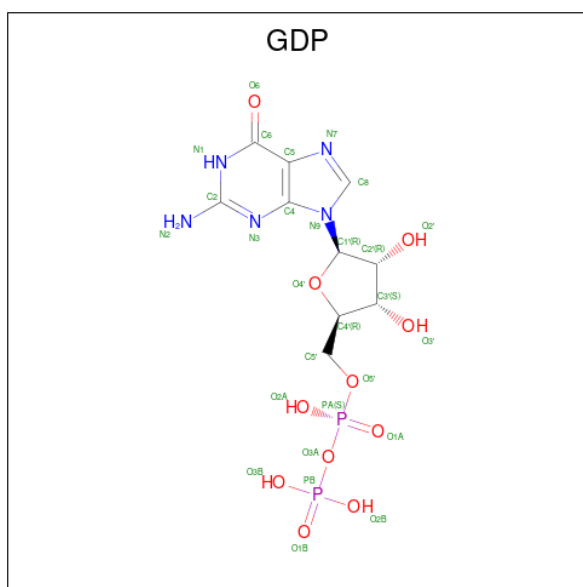
- Molecule 2 is a protein called Importin subunit beta-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	859	6642	4197	1099	1310	36	0	0	0
2	D	849	6570	4157	1086	1291	36	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	254	LYS	LEU	engineered mutation	UNP Q06142
D	254	LYS	LEU	engineered mutation	UNP Q06142

- Molecule 3 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula:  $C_{10}H_{15}N_5O_{11}P_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
3	A	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
3	C	1	Total	C	N	O	P	0	0
			28	10	5	11	2		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mg	0	0
			2	2		
4	C	2	Total	Mg	0	0
			2	2		

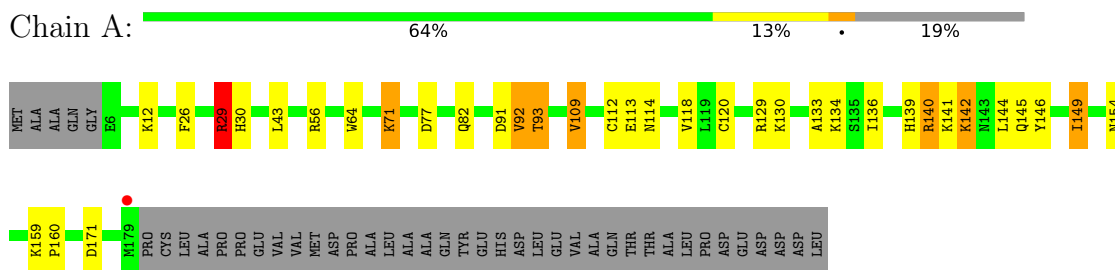
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	68	Total	O	0	0
			68	68		
5	B	178	Total	O	0	0
			178	178		
5	C	34	Total	O	0	0
			34	34		
5	D	84	Total	O	0	0
			84	84		

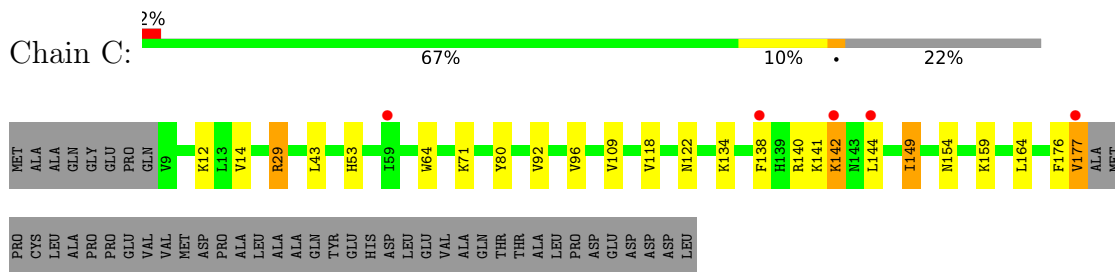
### 3 Residue-property plots

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

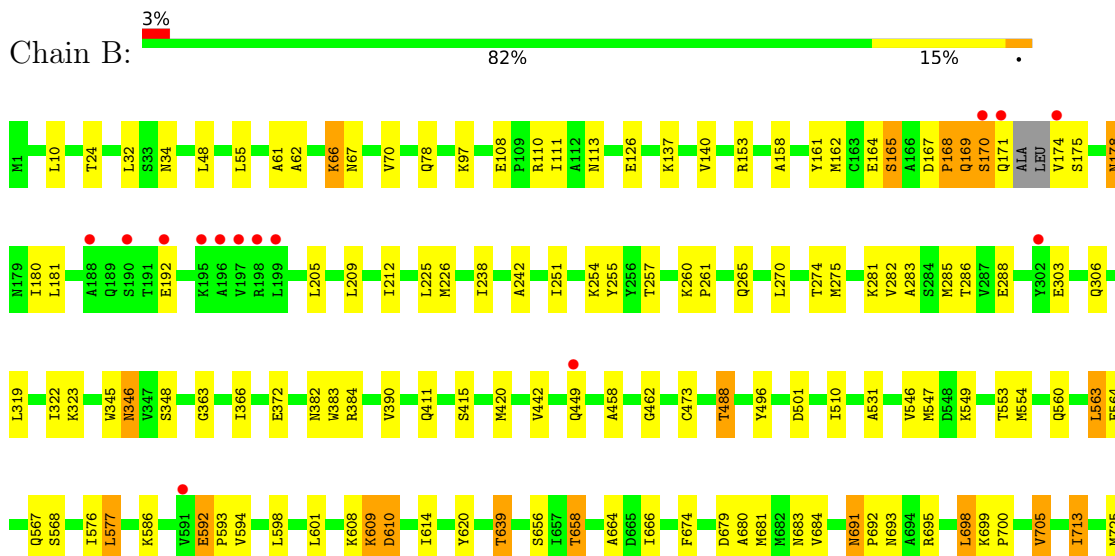
- Molecule 1: GTP-binding nuclear protein Ran



- Molecule 1: GTP-binding nuclear protein Ran

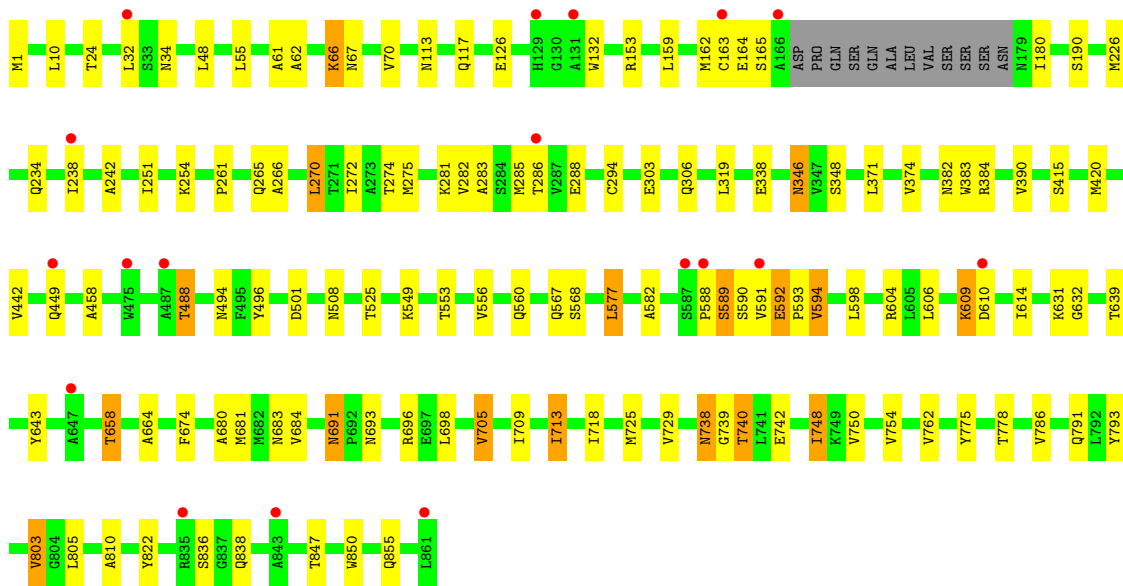
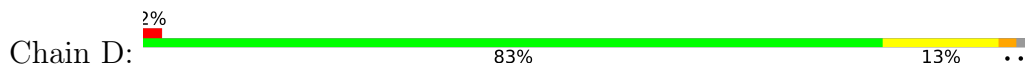


- Molecule 2: Importin subunit beta-1





• Molecule 2: Importin subunit beta-1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	110.85Å 127.81Å 171.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 20.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.50) 99.7 (20.00-2.50)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.78 (at 2.50Å)	Xtrriage
Refinement program	REFMAC 5.3.0037	Depositor
R, $R_{free}$	0.189 , 0.232 0.193 , 0.233	Depositor DCC
$R_{free}$ test set	4240 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	50.5	Xtrriage
Anisotropy	0.066	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 55.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	16428	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	68.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.60% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical 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: MG, GDP

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	0.92	1/1450 (0.1%)	1.00	3/1958 (0.2%)
1	C	0.74	0/1411	0.91	1/1905 (0.1%)
2	B	0.78	0/6755	0.92	5/9179 (0.1%)
2	D	0.70	2/6682 (0.0%)	0.89	2/9079 (0.0%)
All	All	0.76	3/16298 (0.0%)	0.92	11/22121 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	631	LYS	N-CA	7.20	1.55	1.46
2	D	718	ILE	CA-CB	5.46	1.57	1.53
1	A	109	VAL	CA-CB	5.21	1.60	1.54

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	838	GLN	N-CA-C	-6.50	105.27	113.72
1	C	142	LYS	N-CA-C	6.18	118.52	111.11
2	B	488	THR	CA-C-N	-5.90	112.46	119.84
2	B	488	THR	C-N-CA	-5.90	112.46	119.84
2	D	488	THR	CA-C-N	-5.71	112.70	119.84
2	D	488	THR	C-N-CA	-5.71	112.70	119.84
1	A	142	LYS	N-CA-C	5.62	117.85	111.11
2	B	372	GLU	N-CA-C	-5.17	105.65	111.28
2	B	175	SER	N-CA-C	5.15	118.40	110.42
1	A	29	ARG	NE-CZ-NH2	-5.03	114.67	119.20
1	A	114	ASN	N-CA-C	5.03	118.94	112.41

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	1415	0	1435	34	0
1	C	1377	0	1400	20	0
2	B	6642	0	6592	122	0
2	D	6570	0	6530	90	0
3	A	28	0	12	0	0
3	C	28	0	12	1	0
4	A	2	0	0	2	0
4	C	2	0	0	1	0
5	A	68	0	0	4	0
5	B	178	0	0	15	0
5	C	34	0	0	2	0
5	D	84	0	0	15	0
All	All	16428	0	15981	244	0

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

All (244) 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)
2:D:594:VAL:HG23	5:D:885:HOH:O	1.39	1.18
2:B:226:MET:HE1	2:B:251:ILE:HG21	1.30	1.13
2:D:226:MET:HE1	2:D:251:ILE:HG21	1.11	1.08
2:B:169:GLN:HE21	2:B:174:VAL:HG21	1.23	1.04
2:D:226:MET:CE	2:D:251:ILE:HG21	1.90	1.00
2:B:167:ASP:O	2:B:169:GLN:N	2.03	0.92
1:C:154:ASN:HD21	2:D:567:GLN:HE21	1.18	0.87
1:C:142:LYS:HE2	2:D:281:LYS:HZ2	1.40	0.86
2:D:371:LEU:O	5:D:926:HOH:O	1.93	0.85
1:C:29:ARG:HH11	2:D:567:GLN:HE22	1.21	0.85
2:B:257:THR:HG22	5:D:865:HOH:O	1.76	0.84
2:B:226:MET:CE	2:B:251:ILE:HG21	2.07	0.84
2:B:169:GLN:NE2	2:B:174:VAL:HG21	1.93	0.82
2:B:620:TYR:HE2	2:B:658:THR:HG21	1.46	0.81
2:B:639:THR:HG22	5:B:878:HOH:O	1.80	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:169:GLN:HE21	2:B:174:VAL:CG2	1.93	0.80
2:B:411:GLN:HG2	5:B:1000:HOH:O	1.82	0.79
2:D:742:GLU:HG2	5:D:921:HOH:O	1.83	0.79
2:B:140:VAL:HG21	2:B:178:ASN:HB3	1.64	0.79
2:D:226:MET:HE1	2:D:251:ILE:CG2	2.05	0.78
4:A:222:MG:MG	5:A:224:HOH:O	1.27	0.78
2:B:169:GLN:NE2	2:B:174:VAL:HG11	1.99	0.78
5:B:863:HOH:O	2:D:838:GLN:HG3	1.84	0.77
2:B:226:MET:HE1	2:B:251:ILE:HD13	1.65	0.76
1:C:141:LYS:HB2	1:C:144:LEU:HD12	1.66	0.76
2:B:169:GLN:HA	5:B:935:HOH:O	1.86	0.76
2:B:113:ASN:HD21	2:B:153:ARG:HH12	1.35	0.75
4:C:222:MG:MG	5:C:223:HOH:O	1.30	0.74
2:D:374:VAL:HB	5:D:926:HOH:O	1.90	0.72
2:B:609:LYS:HD3	5:B:933:HOH:O	1.89	0.71
2:B:261:PRO:O	2:B:265:GLN:HG2	1.92	0.70
1:A:149:ILE:HD11	1:A:160:PRO:HG3	1.74	0.69
2:D:371:LEU:C	5:D:926:HOH:O	2.35	0.69
2:D:588:PRO:C	2:D:590:SER:H	1.99	0.69
2:D:338:GLU:HG2	2:D:384:ARG:HH21	1.57	0.68
1:C:29:ARG:HH11	2:D:567:GLN:NE2	1.92	0.68
2:D:604:ARG:HD2	5:D:869:HOH:O	1.93	0.68
1:A:82:GLN:NE2	2:B:110:ARG:HH22	1.92	0.68
2:B:674:PHE:HB3	2:B:713:ILE:HD11	1.76	0.68
2:B:680:ALA:O	2:B:684:VAL:HG23	1.94	0.67
2:D:374:VAL:N	5:D:926:HOH:O	2.27	0.67
2:D:338:GLU:HG2	2:D:384:ARG:NH2	2.09	0.67
2:D:226:MET:HE1	2:D:251:ILE:HD13	1.77	0.66
2:B:169:GLN:HG2	2:B:174:VAL:HG21	1.77	0.66
2:B:608:LYS:O	5:B:961:HOH:O	2.14	0.66
2:D:508:ASN:OD1	2:D:549:LYS:NZ	2.28	0.65
2:D:691:ASN:HD22	2:D:693:ASN:H	1.44	0.65
4:A:222:MG:MG	5:A:223:HOH:O	1.40	0.64
2:B:169:GLN:HE21	2:B:174:VAL:HG11	1.62	0.64
2:D:113:ASN:HD21	2:D:153:ARG:HH12	1.45	0.64
2:B:577:LEU:HD13	2:B:614:ILE:HD11	1.78	0.64
1:A:154:ASN:HD21	2:B:567:GLN:HE21	1.44	0.64
2:B:729:VAL:HG22	2:B:778:THR:HG21	1.78	0.64
1:A:149:ILE:HD11	1:A:160:PRO:CG	2.28	0.64
1:C:154:ASN:HD21	2:D:567:GLN:NE2	1.94	0.64
3:C:220:GDP:O1B	5:C:223:HOH:O	2.15	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:LYS:NZ	5:A:281:HOH:O	2.31	0.62
2:B:108:GLU:HG3	5:B:905:HOH:O	1.99	0.61
2:D:658:THR:HG22	5:D:890:HOH:O	2.00	0.61
2:D:66:LYS:HE3	2:D:67:ASN:OD1	2.00	0.61
2:D:1:MET:N	5:D:872:HOH:O	2.34	0.60
2:B:281:LYS:O	2:B:285:MET:HG3	2.02	0.60
2:B:738:ASN:HD22	2:B:739:GLY:N	2.00	0.60
2:D:691:ASN:ND2	2:D:693:ASN:H	2.00	0.59
1:A:29:ARG:HH11	2:B:567:GLN:HE22	1.49	0.59
2:B:420:MET:HG3	2:B:458:ALA:HB1	1.83	0.59
2:B:729:VAL:HG22	2:B:778:THR:CG2	2.33	0.59
2:B:346:ASN:ND2	2:B:348:SER:OG	2.35	0.59
2:B:691:ASN:C	2:B:691:ASN:HD22	2.10	0.59
2:B:786:VAL:O	2:B:793:TYR:HB3	2.03	0.59
2:D:281:LYS:O	2:D:285:MET:HG3	2.03	0.58
2:B:66:LYS:HE3	2:B:67:ASN:OD1	2.02	0.58
2:D:261:PRO:O	2:D:265:GLN:HG2	2.04	0.58
2:D:738:ASN:HD22	2:D:739:GLY:N	2.00	0.57
2:D:346:ASN:ND2	2:D:348:SER:OG	2.37	0.57
1:A:136:ILE:HD13	1:A:146:TYR:CZ	2.39	0.57
2:B:609:LYS:O	2:B:610:ASP:HB2	2.03	0.57
2:B:691:ASN:ND2	2:B:693:ASN:H	2.02	0.57
2:B:226:MET:CE	2:B:251:ILE:HD13	2.33	0.57
2:D:226:MET:CE	2:D:251:ILE:HD13	2.35	0.57
2:D:681:MET:HE2	2:D:681:MET:HA	1.88	0.56
2:D:242:ALA:HB2	2:D:285:MET:HE2	1.86	0.56
1:A:91:ASP:OD1	1:A:93:THR:HB	2.06	0.56
2:B:275:MET:HA	2:B:286:THR:HG21	1.88	0.56
1:C:142:LYS:HE2	2:D:281:LYS:NZ	2.18	0.56
2:D:275:MET:HA	2:D:286:THR:HG21	1.88	0.56
2:B:283:ALA:O	2:B:286:THR:HG22	2.05	0.55
2:D:577:LEU:HD13	2:D:614:ILE:HD11	1.88	0.55
2:D:609:LYS:O	2:D:610:ASP:HB2	2.06	0.55
2:B:169:GLN:CG	2:B:174:VAL:HG21	2.36	0.55
2:D:786:VAL:O	2:D:793:TYR:HB3	2.06	0.55
2:B:169:GLN:HE21	2:B:174:VAL:CG1	2.20	0.55
1:C:53:HIS:O	1:C:176:PHE:O	2.25	0.54
2:D:382:ASN:HD22	2:D:383:TRP:N	2.05	0.54
2:D:691:ASN:HD22	2:D:691:ASN:C	2.15	0.54
2:D:163:CYS:SG	2:D:180:ILE:HG21	2.48	0.54
2:B:547:MET:HE1	2:B:594:VAL:HG21	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:803:VAL:HG12	2:D:847:THR:HG22	1.89	0.53
2:B:169:GLN:HE21	2:B:174:VAL:CB	2.21	0.53
1:A:77:ASP:OD2	2:B:66:LYS:NZ	2.41	0.53
2:B:382:ASN:HD22	2:B:383:TRP:N	2.06	0.53
2:D:234:GLN:NE2	2:D:270:LEU:HD13	2.23	0.53
1:A:142:LYS:CE	2:B:281:LYS:HD3	2.39	0.53
2:B:729:VAL:CG2	2:B:778:THR:HG21	2.39	0.53
1:A:56:ARG:NH1	1:A:171:ASP:OD2	2.42	0.53
2:B:554:MET:CE	2:B:601:LEU:HD22	2.38	0.53
2:B:664:ALA:HB2	2:B:705:VAL:HG23	1.91	0.53
1:C:142:LYS:CE	2:D:281:LYS:HZ2	2.17	0.53
2:B:725:MET:HE2	2:B:775:TYR:HB3	1.91	0.52
2:B:620:TYR:CE2	2:B:658:THR:HG21	2.35	0.52
2:D:632:GLY:N	5:D:877:HOH:O	2.42	0.52
2:B:738:ASN:ND2	2:B:740:THR:OG1	2.42	0.52
1:C:176:PHE:O	1:C:177:VAL:C	2.51	0.52
1:A:71:LYS:HB3	1:A:71:LYS:HZ3	1.75	0.52
2:B:837:GLY:N	5:B:996:HOH:O	2.39	0.52
2:B:78:GLN:NE2	5:B:1013:HOH:O	2.18	0.52
2:B:242:ALA:HB2	2:B:285:MET:HE2	1.92	0.52
1:A:142:LYS:HE2	2:B:281:LYS:HD3	1.92	0.52
1:A:71:LYS:CB	1:A:71:LYS:NZ	2.73	0.51
2:B:288:GLU:OE1	5:B:906:HOH:O	2.19	0.51
2:B:382:ASN:HD22	2:B:382:ASN:C	2.19	0.51
2:B:546:VAL:HB	2:B:576:ILE:HG23	1.92	0.51
2:D:680:ALA:O	2:D:684:VAL:HG23	2.10	0.51
2:D:420:MET:HG3	2:D:458:ALA:HB1	1.93	0.51
1:A:133:ALA:HA	1:A:136:ILE:HD12	1.93	0.51
1:A:112:CYS:O	1:A:113:GLU:C	2.51	0.50
1:A:12:LYS:HE3	1:A:64:TRP:CE2	2.47	0.50
1:A:82:GLN:NE2	2:B:110:ARG:NH2	2.58	0.50
2:B:531:ALA:O	2:B:586:LYS:HE2	2.11	0.50
2:B:691:ASN:HD22	2:B:693:ASN:H	1.58	0.50
2:B:382:ASN:ND2	2:B:384:ARG:H	2.09	0.50
2:D:664:ALA:HB2	2:D:705:VAL:HG23	1.94	0.50
2:B:158:ALA:O	2:B:162:MET:HE3	2.12	0.49
2:B:171:GLN:O	2:B:174:VAL:N	2.45	0.49
2:B:803:VAL:HG11	2:B:848:ALA:HA	1.94	0.49
2:D:48:LEU:HD13	2:D:62:ALA:HB2	1.94	0.49
1:A:142:LYS:HZ1	2:B:281:LYS:HD3	1.77	0.49
2:B:205:LEU:HD11	2:B:225:LEU:HD11	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:549:LYS:O	2:B:553:THR:HG23	2.12	0.49
2:B:681:MET:HE2	2:B:681:MET:HA	1.95	0.49
1:A:154:ASN:HD21	2:B:567:GLN:NE2	2.10	0.49
2:B:303:GLU:OE2	2:D:303:GLU:OE2	2.30	0.49
1:C:12:LYS:HE3	1:C:64:TRP:CE2	2.48	0.48
2:D:283:ALA:O	2:D:286:THR:HG22	2.13	0.48
2:B:161:TYR:O	2:B:165:SER:HB2	2.13	0.48
1:A:71:LYS:HB3	1:A:71:LYS:NZ	2.28	0.48
2:B:411:GLN:CG	5:B:1000:HOH:O	2.49	0.48
2:D:346:ASN:HD22	2:D:348:SER:H	1.59	0.48
2:D:382:ASN:HD22	2:D:382:ASN:C	2.21	0.48
1:A:26:PHE:CE1	1:A:30:HIS:HE1	2.31	0.48
2:D:738:ASN:ND2	2:D:740:THR:OG1	2.45	0.48
2:B:108:GLU:HB3	2:B:111:ILE:HD12	1.96	0.48
2:B:679:ASP:HB3	5:B:899:HOH:O	2.14	0.48
1:C:118:VAL:HG23	1:C:164:LEU:HD21	1.96	0.47
1:A:120:CYS:SG	1:A:149:ILE:HD13	2.54	0.47
1:A:141:LYS:HB2	1:A:144:LEU:HD12	1.96	0.47
2:B:181:LEU:HD23	2:B:181:LEU:C	2.39	0.47
2:B:592:GLU:N	2:B:593:PRO:HD2	2.28	0.47
1:A:120:CYS:SG	1:A:149:ILE:CD1	3.02	0.47
2:D:658:THR:CG2	5:D:890:HOH:O	2.62	0.47
2:B:346:ASN:HD22	2:B:348:SER:H	1.61	0.47
2:B:275:MET:CA	2:B:286:THR:HG21	2.45	0.47
2:B:656:SER:OG	2:B:698:LEU:HD21	2.15	0.47
2:D:592:GLU:N	2:D:593:PRO:HD2	2.29	0.47
2:D:159:LEU:HD22	2:D:180:ILE:HG23	1.98	0.46
2:D:729:VAL:HG22	2:D:778:THR:HG21	1.97	0.46
2:D:674:PHE:HB3	2:D:713:ILE:HD11	1.96	0.46
2:D:588:PRO:C	2:D:590:SER:N	2.69	0.46
1:C:141:LYS:HB2	1:C:144:LEU:CD1	2.40	0.46
2:D:274:THR:HG23	2:D:282:VAL:CG1	2.45	0.46
1:A:92:VAL:HG13	1:A:129:ARG:HB3	1.97	0.46
1:C:142:LYS:NZ	2:D:281:LYS:NZ	2.64	0.46
1:C:142:LYS:HZ1	2:D:281:LYS:HD3	1.81	0.45
1:C:14:VAL:HG11	1:C:80:TYR:HA	1.98	0.45
2:D:725:MET:HE2	2:D:775:TYR:HB3	1.99	0.45
2:B:167:ASP:O	2:B:168:PRO:C	2.59	0.45
2:B:322:ILE:O	2:B:323:LYS:C	2.60	0.45
2:B:837:GLY:CA	5:B:996:HOH:O	2.63	0.45
2:D:374:VAL:CA	5:D:926:HOH:O	2.64	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:126:GLU:OE1	2:D:132:TRP:HB2	2.17	0.45
1:A:82:GLN:HE21	2:B:110:ARG:HH12	1.65	0.44
2:D:606:LEU:HB3	2:D:643:TYR:CZ	2.52	0.44
2:B:180:ILE:HD12	5:B:980:HOH:O	2.15	0.44
2:B:363:GLY:O	2:B:366:ILE:HG22	2.17	0.44
2:B:674:PHE:HB3	2:B:713:ILE:CD1	2.46	0.44
1:A:29:ARG:HH11	2:B:567:GLN:NE2	2.14	0.44
1:A:136:ILE:HD13	1:A:146:TYR:CE2	2.52	0.44
2:B:803:VAL:CG1	2:B:848:ALA:HA	2.46	0.44
1:A:140:ARG:HG2	2:B:345:TRP:NE1	2.33	0.44
2:B:260:LYS:HB3	2:B:261:PRO:HD3	1.99	0.44
2:B:510:ILE:HG22	2:B:510:ILE:O	2.16	0.44
2:D:275:MET:CA	2:D:286:THR:HG21	2.48	0.44
2:D:762:VAL:HG21	2:D:805:LEU:HD22	2.00	0.44
2:B:274:THR:HG23	2:B:282:VAL:HG12	1.99	0.44
2:B:48:LEU:HD13	2:B:62:ALA:HB2	2.00	0.43
2:B:800:ARG:HD3	2:B:844:THR:HA	1.99	0.43
1:C:142:LYS:CE	2:D:281:LYS:NZ	2.78	0.43
2:D:117:GLN:NE2	5:D:912:HOH:O	2.49	0.43
2:B:462:GLY:HA3	2:B:473:CYS:SG	2.58	0.43
2:D:606:LEU:HB3	2:D:643:TYR:CE1	2.54	0.43
2:B:169:GLN:O	2:B:170:SER:CB	2.66	0.43
2:B:170:SER:HA	5:B:1023:HOH:O	2.19	0.43
2:D:729:VAL:HG22	2:D:778:THR:CG2	2.49	0.43
1:A:139:HIS:HB2	1:A:145:GLN:HA	2.01	0.42
2:B:167:ASP:O	2:B:169:GLN:CA	2.65	0.42
2:B:346:ASN:HD22	2:B:348:SER:N	2.17	0.42
2:D:346:ASN:HD22	2:D:348:SER:N	2.16	0.42
1:A:82:GLN:HE22	2:B:110:ARG:NH2	2.17	0.42
2:B:835:ARG:NH1	2:B:849:ARG:HG3	2.33	0.42
2:D:10:LEU:HD13	2:D:61:ALA:HB2	2.00	0.42
1:C:122:ASN:OD1	1:C:149:ILE:HG22	2.19	0.42
2:D:577:LEU:HD13	2:D:614:ILE:CD1	2.49	0.42
2:B:209:LEU:O	2:B:255:TYR:OH	2.30	0.42
2:B:738:ASN:ND2	2:B:743:ALA:CB	2.82	0.42
2:B:592:GLU:HG3	2:B:593:PRO:CD	2.50	0.42
2:B:225:LEU:C	2:B:225:LEU:HD23	2.45	0.42
2:D:162:MET:HE3	2:D:162:MET:HB2	1.89	0.42
2:D:374:VAL:CB	5:D:926:HOH:O	2.60	0.42
2:D:549:LYS:O	2:D:553:THR:HG23	2.20	0.42
2:B:10:LEU:HD13	2:B:61:ALA:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:274:THR:HG23	2:B:282:VAL:CG1	2.50	0.42
2:D:592:GLU:HG3	2:D:593:PRO:CD	2.50	0.42
2:D:748:ILE:HD13	2:D:791:GLN:HG2	2.02	0.41
1:A:159:LYS:O	1:A:160:PRO:C	2.63	0.41
2:B:97:LYS:CE	2:B:126:GLU:OE1	2.69	0.41
2:B:181:LEU:HD12	2:B:212:ILE:HD12	2.01	0.41
2:B:695:ARG:HG2	2:B:698:LEU:HD23	2.02	0.41
2:D:346:ASN:HD22	2:D:346:ASN:C	2.28	0.41
2:D:242:ALA:HA	2:D:285:MET:SD	2.60	0.41
2:D:750:VAL:O	2:D:754:VAL:HG23	2.21	0.41
1:C:138:PHE:CD2	1:C:138:PHE:C	2.97	0.41
2:D:810:ALA:HB2	2:D:822:TYR:CZ	2.55	0.41
2:B:592:GLU:HG3	2:B:593:PRO:HD3	2.02	0.41
2:B:691:ASN:HD22	2:B:692:PRO:N	2.19	0.41
2:D:709:ILE:O	2:D:713:ILE:HB	2.21	0.41
5:A:254:HOH:O	2:B:563:LEU:HD22	2.20	0.41
2:B:554:MET:CE	2:B:601:LEU:CD2	2.99	0.41
2:D:525:THR:HG23	2:D:582:ALA:HB2	2.01	0.41
2:B:699:LYS:HB3	2:B:700:PRO:CD	2.51	0.40
2:D:66:LYS:HE2	2:D:66:LYS:HB3	1.94	0.40
1:A:142:LYS:NZ	2:B:281:LYS:HD3	2.35	0.40
2:B:158:ALA:O	2:B:162:MET:HB2	2.21	0.40
2:B:698:LEU:HA	2:B:698:LEU:HD13	1.88	0.40
1:C:142:LYS:NZ	2:D:281:LYS:HZ1	2.20	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	172/216 (80%)	166 (96%)	6 (4%)	0	<a href="#">100</a> <a href="#">100</a>

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	167/216 (77%)	160 (96%)	7 (4%)	0	100	100
2	B	855/861 (99%)	827 (97%)	25 (3%)	3 (0%)	30	49
2	D	845/861 (98%)	821 (97%)	22 (3%)	2 (0%)	43	63
All	All	2039/2154 (95%)	1974 (97%)	60 (3%)	5 (0%)	43	63

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	168	PRO
2	B	169	GLN
2	B	170	SER
2	D	266	ALA
2	D	589	SER

### 5.3.2 Protein sidechains [i](#)

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	153/186 (82%)	142 (93%)	11 (7%)	13	28
1	C	149/186 (80%)	138 (93%)	11 (7%)	13	27
2	B	725/726 (100%)	675 (93%)	50 (7%)	14	30
2	D	715/726 (98%)	664 (93%)	51 (7%)	13	29
All	All	1742/1824 (96%)	1619 (93%)	123 (7%)	13	29

All (123) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	ARG
1	A	43	LEU
1	A	71	LYS
1	A	92	VAL
1	A	93	THR
1	A	109	VAL

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	118	VAL
1	A	130	LYS
1	A	134	LYS
1	A	140	ARG
1	A	149	ILE
2	B	24	THR
2	B	32	LEU
2	B	34	ASN
2	B	55	LEU
2	B	66	LYS
2	B	70	VAL
2	B	137	LYS
2	B	164	GLU
2	B	165	SER
2	B	178	ASN
2	B	192	GLU
2	B	238	ILE
2	B	254	LYS
2	B	270	LEU
2	B	306	GLN
2	B	319	LEU
2	B	346	ASN
2	B	390	VAL
2	B	415	SER
2	B	442	VAL
2	B	449	GLN
2	B	488	THR
2	B	496	TYR
2	B	501	ASP
2	B	560	GLN
2	B	563	LEU
2	B	564	GLU
2	B	568	SER
2	B	577	LEU
2	B	592	GLU
2	B	598	LEU
2	B	609	LYS
2	B	610	ASP
2	B	639	THR
2	B	658	THR
2	B	666	ILE
2	B	683	ASN

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
2	B	691	ASN
2	B	698	LEU
2	B	705	VAL
2	B	713	ILE
2	B	738	ASN
2	B	740	THR
2	B	741	LEU
2	B	748	ILE
2	B	755	LEU
2	B	803	VAL
2	B	835	ARG
2	B	836	SER
2	B	838	GLN
1	C	29	ARG
1	C	43	LEU
1	C	71	LYS
1	C	92	VAL
1	C	96	VAL
1	C	109	VAL
1	C	134	LYS
1	C	140	ARG
1	C	149	ILE
1	C	159	LYS
1	C	177	VAL
2	D	24	THR
2	D	32	LEU
2	D	34	ASN
2	D	55	LEU
2	D	66	LYS
2	D	70	VAL
2	D	164	GLU
2	D	165	SER
2	D	190	SER
2	D	238	ILE
2	D	254	LYS
2	D	270	LEU
2	D	272	ILE
2	D	288	GLU
2	D	294	CYS
2	D	306	GLN
2	D	319	LEU
2	D	346	ASN

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Mol	Chain	Res	Type
2	D	390	VAL
2	D	415	SER
2	D	442	VAL
2	D	449	GLN
2	D	488	THR
2	D	494	ASN
2	D	496	TYR
2	D	501	ASP
2	D	556	VAL
2	D	560	GLN
2	D	568	SER
2	D	577	LEU
2	D	589	SER
2	D	591	VAL
2	D	592	GLU
2	D	594	VAL
2	D	598	LEU
2	D	609	LYS
2	D	639	THR
2	D	658	THR
2	D	683	ASN
2	D	691	ASN
2	D	696	ARG
2	D	698	LEU
2	D	705	VAL
2	D	713	ILE
2	D	738	ASN
2	D	740	THR
2	D	748	ILE
2	D	803	VAL
2	D	836	SER
2	D	850	TRP
2	D	855	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (48) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	GLN
1	A	82	GLN
1	A	100	ASN
1	A	114	ASN
2	B	46	GLN

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
2	B	78	GLN
2	B	87	GLN
2	B	113	ASN
2	B	169	GLN
2	B	234	GLN
2	B	313	GLN
2	B	346	ASN
2	B	365	HIS
2	B	376	GLN
2	B	382	ASN
2	B	404	GLN
2	B	423	GLN
2	B	567	GLN
2	B	649	ASN
2	B	691	ASN
2	B	712	ASN
2	B	722	ASN
2	B	738	ASN
2	B	785	GLN
2	B	838	GLN
2	B	842	GLN
1	C	48	HIS
1	C	82	GLN
1	C	100	ASN
1	C	114	ASN
2	D	46	GLN
2	D	78	GLN
2	D	87	GLN
2	D	113	ASN
2	D	306	GLN
2	D	346	ASN
2	D	376	GLN
2	D	382	ASN
2	D	404	GLN
2	D	423	GLN
2	D	567	GLN
2	D	649	ASN
2	D	691	ASN
2	D	712	ASN
2	D	738	ASN
2	D	785	GLN
2	D	838	GLN

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Mol	Chain	Res	Type
2	D	842	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 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
3	GDP	C	220	4	29,30,30	1.24	4 (13%)	45,47,47	1.50	6 (13%)
3	GDP	A	220	4	29,30,30	1.27	4 (13%)	45,47,47	1.54	8 (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
3	GDP	C	220	4	-	6/16/32/32	0/3/3/3
3	GDP	A	220	4	-	4/16/32/32	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	220	GDP	C6-N1	-3.27	1.32	1.38
3	C	220	GDP	C5-C4	3.01	1.47	1.38
3	A	220	GDP	C5-N7	-2.90	1.33	1.39
3	A	220	GDP	C5-C4	2.82	1.46	1.38
3	C	220	GDP	C5-N7	-2.77	1.33	1.39
3	C	220	GDP	C6-N1	-2.67	1.33	1.38
3	C	220	GDP	C4-N9	-2.45	1.31	1.38
3	A	220	GDP	C4-N9	-2.28	1.32	1.38

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	220	GDP	C5-C4-N3	-4.94	120.53	128.39
3	C	220	GDP	C5-C4-N3	-4.62	121.04	128.39
3	A	220	GDP	N9-C4-N3	3.72	133.39	125.95
3	C	220	GDP	C2-N3-C4	3.48	118.29	112.30
3	C	220	GDP	N9-C4-N3	3.40	132.76	125.95
3	A	220	GDP	C2-N3-C4	3.29	117.97	112.30
3	C	220	GDP	C6-C5-N7	2.90	135.57	130.29
3	A	220	GDP	N2-C2-N3	-2.68	114.44	119.67
3	C	220	GDP	C4-C5-N7	-2.54	106.65	110.67
3	A	220	GDP	O2B-PB-O3A	2.48	112.96	104.64
3	A	220	GDP	C6-C5-N7	2.47	134.78	130.29
3	A	220	GDP	O6-C6-C5	-2.21	120.71	126.53
3	A	220	GDP	C4-C5-N7	-2.02	107.47	110.67
3	C	220	GDP	C8-N7-C5	2.01	107.84	104.26

There are no chirality outliers.

All (10) torsion outliers are listed below:

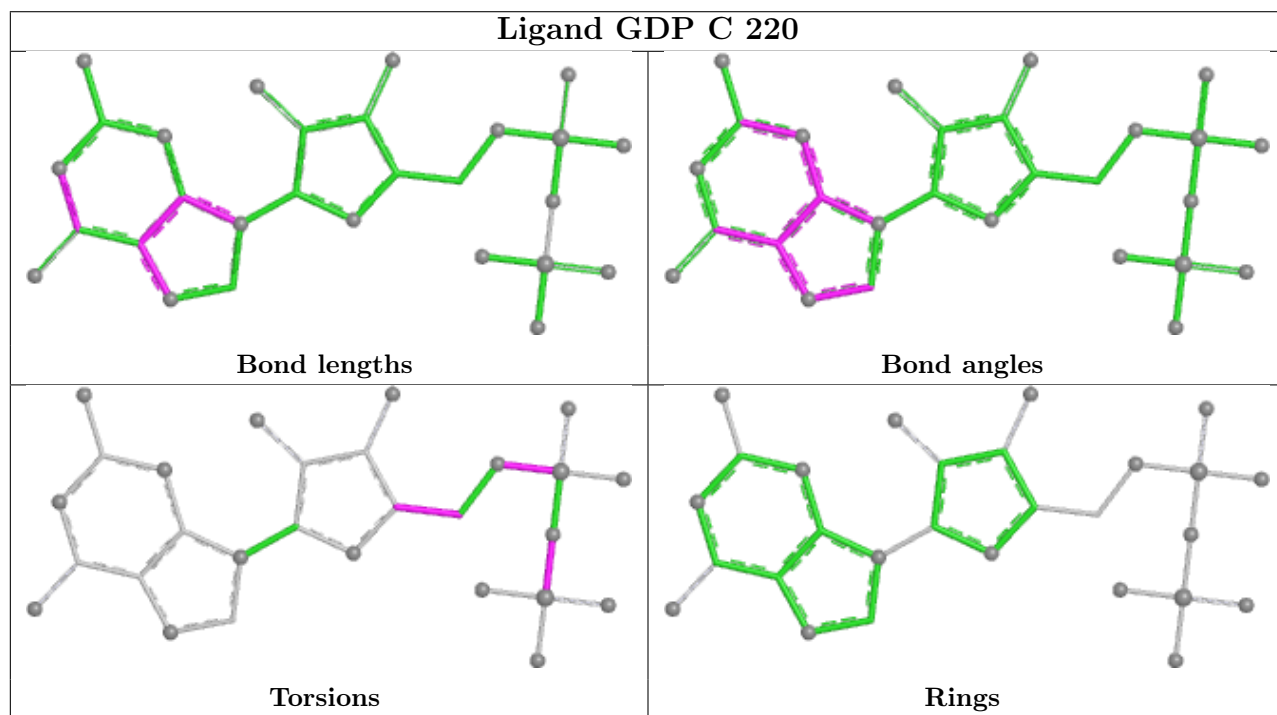
Mol	Chain	Res	Type	Atoms
3	A	220	GDP	PA-O3A-PB-O2B
3	C	220	GDP	PA-O3A-PB-O3B
3	C	220	GDP	C5'-O5'-PA-O1A
3	C	220	GDP	O4'-C4'-C5'-O5'
3	A	220	GDP	O4'-C4'-C5'-O5'
3	A	220	GDP	C5'-O5'-PA-O1A
3	C	220	GDP	C5'-O5'-PA-O3A
3	C	220	GDP	C3'-C4'-C5'-O5'
3	A	220	GDP	PA-O3A-PB-O1B
3	C	220	GDP	PA-O3A-PB-O2B

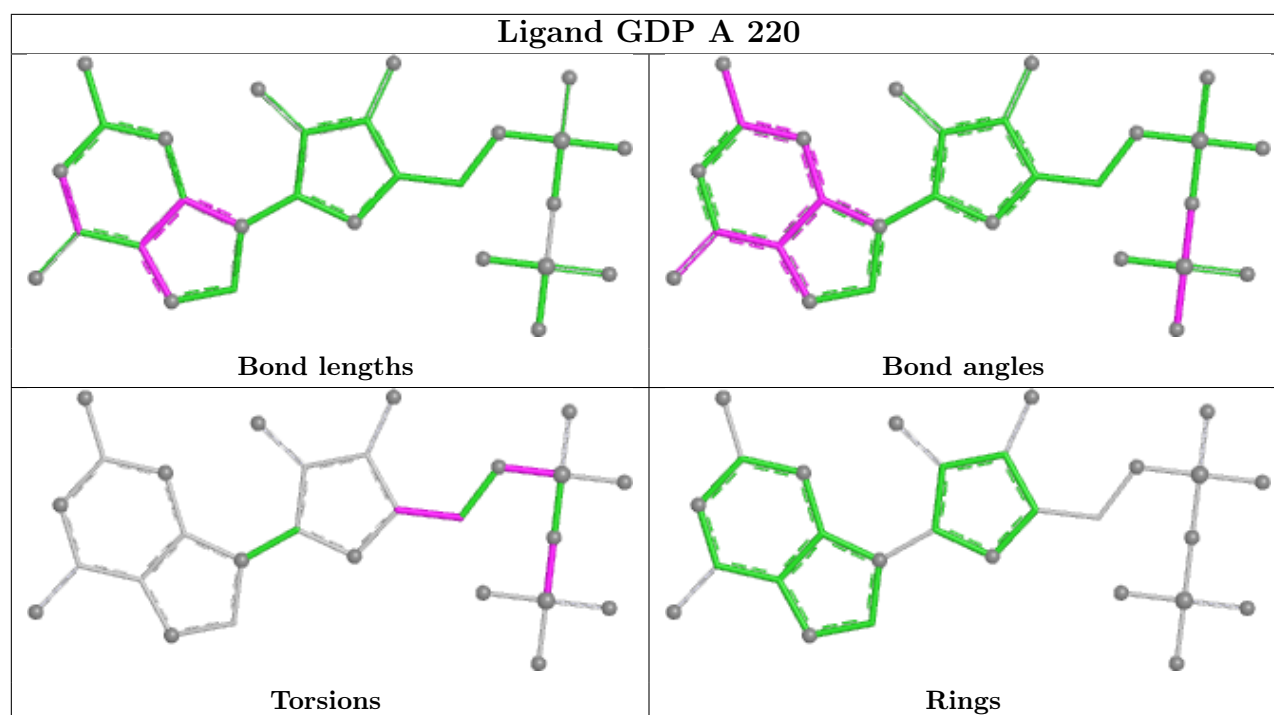
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	220	GDP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





## 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, 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	174/216 (80%)	-0.18	1 (0%) 85 83	52, 62, 88, 121	0
1	C	169/216 (78%)	0.10	5 (2%) 52 48	55, 65, 93, 111	0
2	B	859/861 (99%)	0.10	24 (2%) 55 50	51, 65, 83, 110	0
2	D	849/861 (98%)	0.28	18 (2%) 63 59	53, 67, 82, 99	0
All	All	2051/2154 (95%)	0.15	48 (2%) 61 57	51, 66, 84, 121	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	166	ALA	3.5
2	B	836	SER	3.4
2	B	841	SER	3.4
1	C	177	VAL	3.0
2	D	591	VAL	3.0
2	B	844	THR	3.0
2	D	32	LEU	2.8
2	B	843	ALA	2.8
2	B	838	GLN	2.7
2	B	839	LEU	2.7
2	B	196	ALA	2.6
1	C	144	LEU	2.5
2	D	587	SER	2.5
2	D	588	PRO	2.5
2	D	487	ALA	2.4
2	B	449	GLN	2.4
2	D	835	ARG	2.4
2	B	174	VAL	2.4
2	B	190	SER	2.4
2	D	861	LEU	2.4
1	C	59	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	199	LEU	2.4
2	B	197	VAL	2.4
2	B	171	GLN	2.3
2	D	610	ASP	2.3
1	C	142	LYS	2.3
2	B	195	LYS	2.3
2	D	286	THR	2.3
2	B	198	ARG	2.2
2	B	849	ARG	2.2
1	A	179	MET	2.2
2	D	238	ILE	2.2
2	B	834	THR	2.1
2	B	591	VAL	2.1
1	C	138	PHE	2.1
2	D	131	ALA	2.1
2	D	163	CYS	2.1
2	B	188	ALA	2.1
2	D	647	ALA	2.1
2	D	843	ALA	2.1
2	B	192	GLU	2.0
2	B	847	THR	2.0
2	B	302	TYR	2.0
2	B	170	SER	2.0
2	D	449	GLN	2.0
2	B	840	PHE	2.0
2	D	475	TRP	2.0
2	D	129	HIS	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 oligosaccharides in this entry.

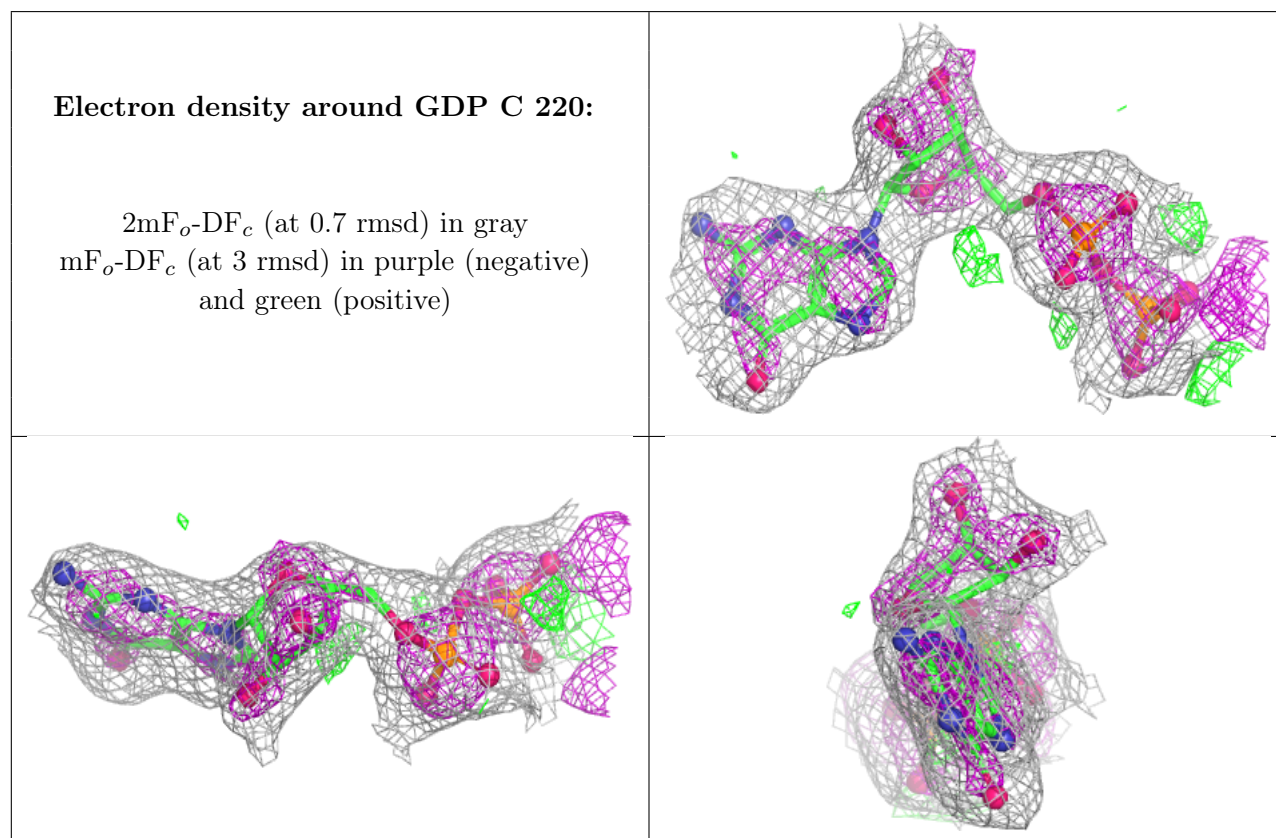
## 6.4 Ligands [i](#)

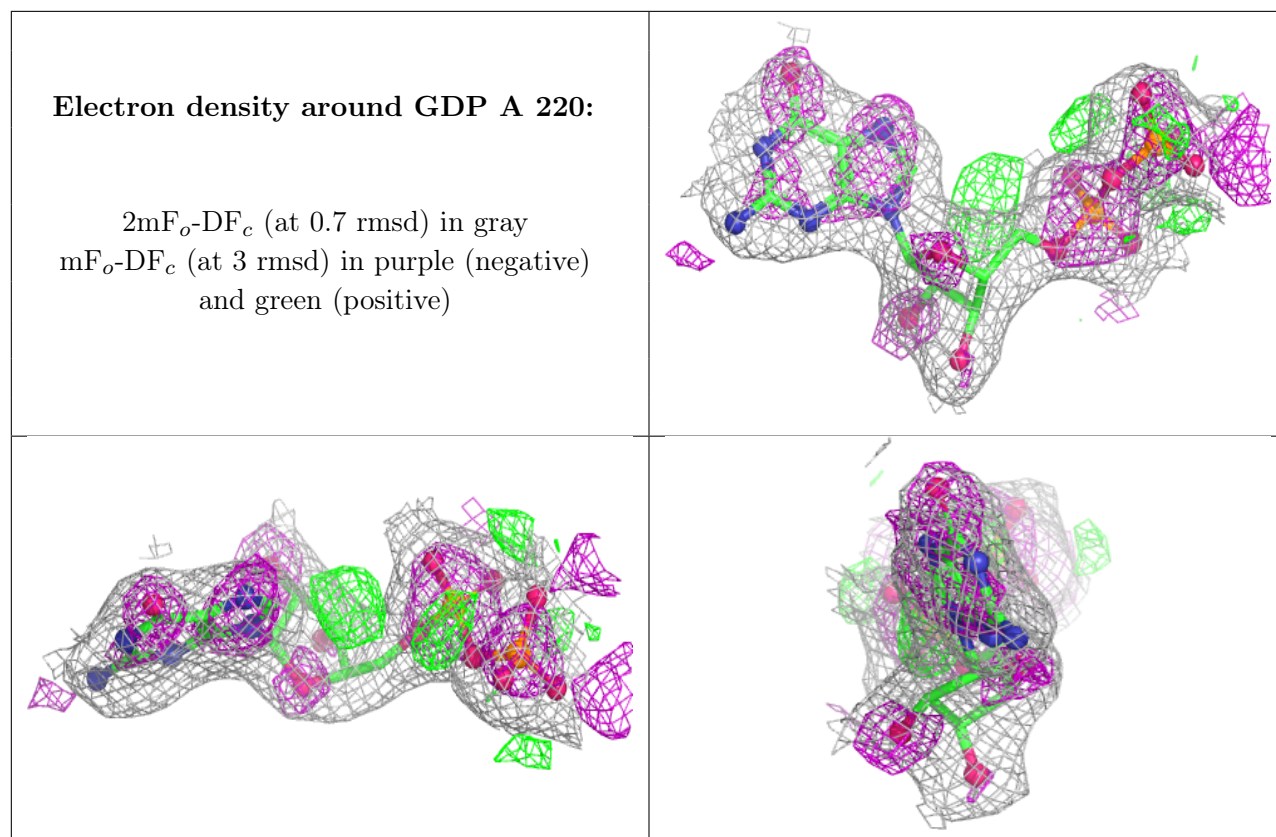
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 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	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	GDP	C	220	28/28	0.95	0.10	53,58,61,61	0
3	GDP	A	220	28/28	0.96	0.11	42,49,53,54	0
4	MG	C	221	1/1	0.96	0.10	55,55,55,55	0
4	MG	A	221	1/1	0.99	0.10	43,43,43,43	0
4	MG	C	222	1/1	0.99	0.12	58,58,58,58	0
4	MG	A	222	1/1	1.00	0.10	55,55,55,55	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.