



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 27, 2026 – 05:12 PM EDT

PDB ID : 3KUD / pdb_00003kud
Title : Complex of Ras-GDP with RafRBD(A85K)
Authors : Filchtinski, D.; Sharabi, O.; Rueppel, A.; Vetter, I.R.; Herrmann, C.; Shifman, J.M.
Deposited on : 2009-11-27
Resolution : 2.15 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

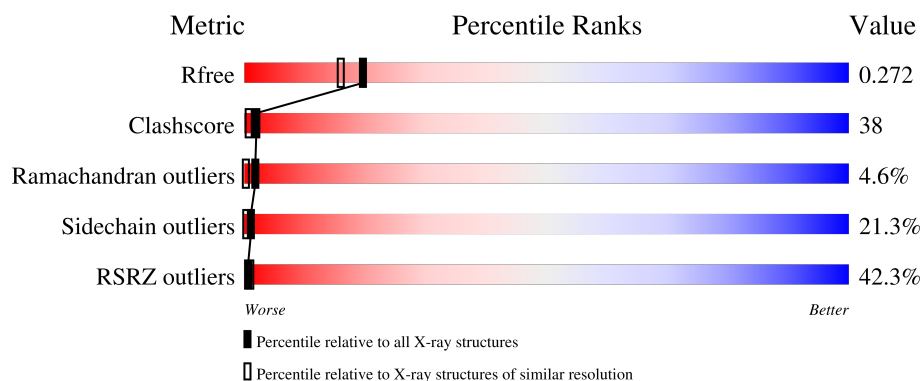
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	166	
2	B	81	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2005 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 GTPase HRas.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	165	Total	C	N	O	S	0	0	0
			1312	818	225	262	7			

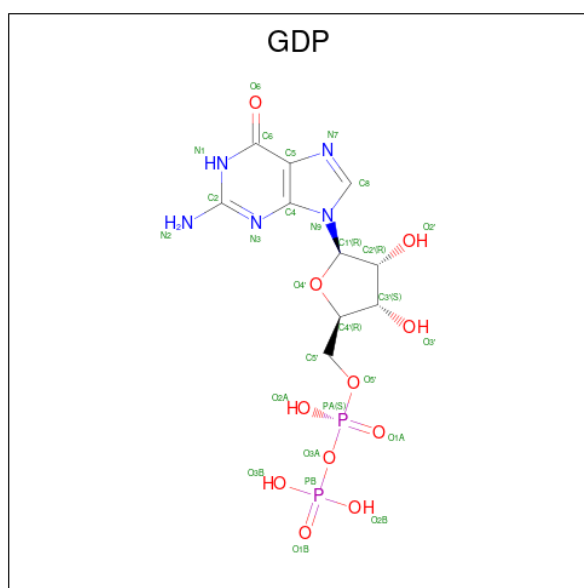
- Molecule 2 is a protein called RAF proto-oncogene serine/threonine-protein kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	76	Total	C	N	O	S	0	2	0
			625	395	119	106	5			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	85	LYS	ALA	engineered mutation	UNP P04049

- 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		

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		

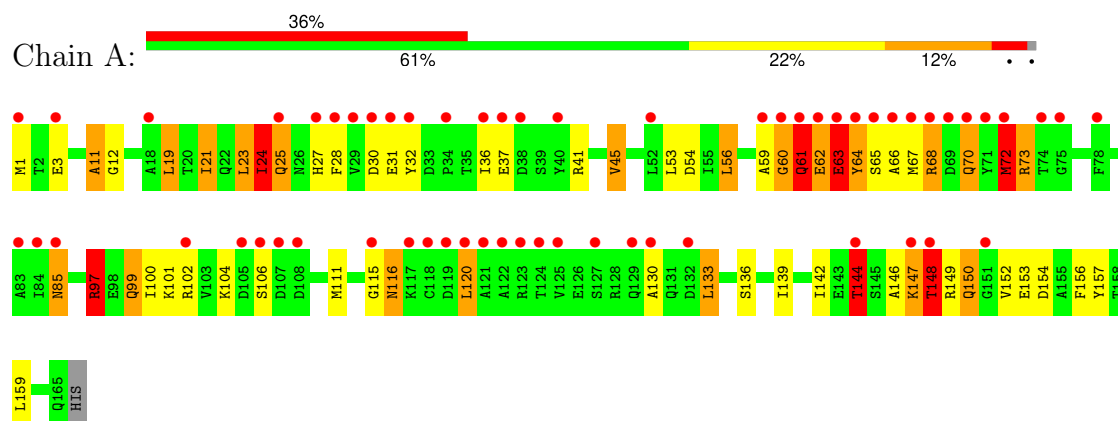
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	26	Total	O	0	0
			26	26		
5	B	13	Total	O	0	0
			13	13		

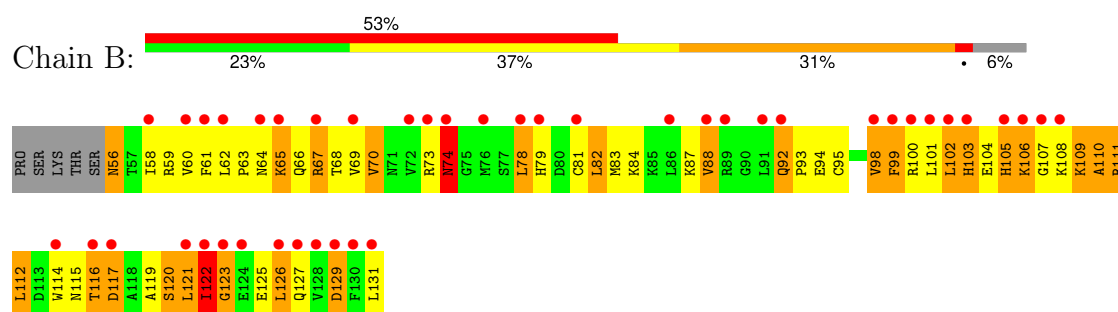
3 Residue-property plots [i](#)

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: GTPase HRas



• Molecule 2: RAF proto-oncogene serine/threonine-protein kinase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	91.13Å 91.13Å 277.98Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.68 – 2.15 19.68 – 2.15	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.68-2.15) 99.6 (19.68-2.15)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.15Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.237 , 0.264 0.249 , 0.272	Depositor DCC
R_{free} test set	1230 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	47.3	Xtriage
Anisotropy	0.351	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 67.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	2005	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.17% 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: GDP, 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	1.63	14/1330 (1.1%)	1.66	23/1795 (1.3%)
2	B	1.48	5/640 (0.8%)	1.53	9/857 (1.1%)
All	All	1.58	19/1970 (1.0%)	1.62	32/2652 (1.2%)

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	3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	25	GLN	CG-CD	9.03	1.74	1.52
1	A	27	HIS	CE1-NE2	-7.40	1.25	1.32
2	B	122	ILE	CA-CB	7.10	1.64	1.54
1	A	152	VAL	CA-C	6.76	1.61	1.52
1	A	116	ASN	C-O	6.32	1.31	1.23
1	A	23	LEU	CA-C	6.30	1.60	1.52
1	A	24	ILE	CA-CB	5.94	1.62	1.54
1	A	130	ALA	CA-CB	5.81	1.62	1.53
1	A	153	GLU	N-CA	5.80	1.53	1.46
2	B	111	ARG	C-O	5.71	1.30	1.23
2	B	122	ILE	CA-C	5.61	1.59	1.52
1	A	147	LYS	C-O	-5.43	1.17	1.24
1	A	25	GLN	CD-NE2	5.32	1.44	1.33
2	B	126	LEU	CA-C	5.23	1.59	1.52
1	A	45	VAL	CA-CB	5.20	1.60	1.54
2	B	116	THR	CA-CB	5.13	1.62	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	142	ILE	CG1-CD1	5.12	1.71	1.51
1	A	150	GLN	CA-C	5.07	1.59	1.52
1	A	120	LEU	N-CA	5.04	1.52	1.45

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	27	HIS	CE1-NE2-CD2	-12.62	96.39	109.00
1	A	27	HIS	ND1-CE1-NE2	9.51	117.91	108.40
2	B	74	ASN	N-CA-C	8.81	123.43	109.50
1	A	157	TYR	N-CA-C	-8.68	101.29	112.23
1	A	148	THR	N-CA-CB	-7.87	99.74	110.67
2	B	70	VAL	N-CA-C	7.61	118.86	107.75
1	A	61	GLN	N-CA-C	-7.34	103.75	114.39
1	A	45	VAL	CB-CA-C	7.17	120.96	110.77
1	A	148	THR	N-CA-C	-6.89	105.14	113.97
2	B	111	ARG	N-CA-C	6.76	120.29	110.28
1	A	64	TYR	N-CA-C	-6.58	92.56	111.00
1	A	139	ILE	CA-C-N	-6.57	113.87	120.31
1	A	139	ILE	C-N-CA	-6.57	113.87	120.31
1	A	27	HIS	CG-CD2-NE2	6.51	113.71	107.20
2	B	108	LYS	N-CA-C	6.11	119.36	109.40
1	A	25	GLN	N-CA-C	-6.01	104.40	113.89
1	A	144	THR	N-CA-CB	-5.94	100.48	111.53
1	A	70	GLN	N-CA-C	-5.90	98.24	110.80
2	B	87	LYS	N-CA-C	5.86	118.45	111.71
2	B	109	LYS	N-CA-C	5.66	117.95	108.73
1	A	133	LEU	N-CA-C	-5.62	105.06	111.07
1	A	154	ASP	CB-CA-C	-5.56	102.16	110.88
1	A	97	ARG	CG-CD-NE	5.55	124.21	112.00
1	A	62	GLU	N-CA-C	-5.52	99.03	110.80
1	A	61	GLN	CB-CA-C	5.52	118.32	109.27
1	A	99	GLN	N-CA-C	-5.51	104.96	110.97
1	A	27	HIS	CB-CG-CD2	5.28	138.07	131.20
1	A	115	GLY	N-CA-C	-5.25	100.73	113.18
1	A	27	HIS	ND1-CG-CD2	-5.15	100.95	106.10
2	B	110	ALA	N-CA-C	5.09	116.61	108.41
2	B	92	GLN	CA-C-N	5.02	125.04	119.32
2	B	92	GLN	C-N-CA	5.02	125.04	119.32

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	61	GLN	Peptide
1	A	62	GLU	Peptide
1	A	63	GLU	Peptide

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	1312	0	1288	53	0
2	B	625	0	660	98	0
3	A	28	0	12	1	0
4	A	1	0	0	0	0
5	A	26	0	0	1	1
5	B	13	0	0	4	0
All	All	2005	0	1960	151	1

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 38.

All (151) 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:25:GLN:CD	1:A:25:GLN:CG	1.74	1.59
1:A:61:GLN:O	1:A:63:GLU:HA	1.25	1.29
2:B:67[A]:ARG:NH1	2:B:67[A]:ARG:HG2	1.50	1.09
2:B:100:ARG:NH2	2:B:116:THR:HG21	1.70	1.06
2:B:67[A]:ARG:HH11	2:B:67[A]:ARG:CG	1.68	1.05
2:B:67[B]:ARG:HH11	2:B:67[B]:ARG:HG3	0.95	1.05
2:B:100:ARG:HH21	2:B:116:THR:HG21	1.24	1.00
2:B:99:PHE:HZ	2:B:129:ASP:HB2	1.26	1.00
2:B:67[B]:ARG:HE	2:B:69:VAL:HG22	1.26	0.97
2:B:101:LEU:HD21	2:B:109:LYS:HG2	1.46	0.97
1:A:148:THR:HG22	1:A:150:GLN:H	1.32	0.94
1:A:61:GLN:C	1:A:63:GLU:HA	1.95	0.91
1:A:61:GLN:O	1:A:63:GLU:CA	2.18	0.91
3:A:170:GDP:O3B	5:A:501:HOH:O	1.86	0.91
2:B:100:ARG:NH2	2:B:116:THR:CG2	2.33	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:67[B]:ARG:HH11	2:B:67[B]:ARG:CG	1.84	0.90
1:A:63:GLU:HB3	1:A:64:TYR:CD2	2.08	0.88
2:B:67[B]:ARG:HG3	2:B:67[B]:ARG:NH1	1.77	0.88
2:B:100:ARG:HG3	2:B:112:LEU:HD21	1.56	0.88
2:B:99:PHE:CZ	2:B:129:ASP:HB2	2.11	0.85
2:B:101:LEU:CD2	2:B:109:LYS:HG2	2.05	0.85
1:A:19:LEU:HD11	1:A:144:THR:HG23	1.58	0.84
2:B:67[A]:ARG:HG2	2:B:67[A]:ARG:HH11	0.74	0.84
2:B:67[B]:ARG:HE	2:B:69:VAL:CG2	1.91	0.83
2:B:100:ARG:HH21	2:B:116:THR:CG2	1.93	0.80
1:A:99:GLN:NE2	1:A:102:ARG:HH22	1.80	0.80
2:B:98:VAL:O	2:B:112:LEU:HB2	1.84	0.77
2:B:64:ASN:O	2:B:65:LYS:HB2	1.84	0.76
1:A:148:THR:CG2	1:A:150:GLN:H	1.97	0.76
1:A:67:MET:SD	1:A:68:ARG:HG3	2.26	0.76
2:B:59:ARG:HH21	2:B:125:GLU:HG3	1.51	0.75
1:A:41:ARG:HD3	1:A:54:ASP:OD1	1.87	0.75
2:B:103:HIS:CE1	5:B:502:HOH:O	2.39	0.74
1:A:99:GLN:HE22	1:A:102:ARG:HH22	1.35	0.74
1:A:19:LEU:HD11	1:A:144:THR:CG2	2.17	0.73
2:B:62:LEU:O	5:B:505:HOH:O	2.06	0.73
1:A:59:ALA:O	1:A:60:GLY:O	2.08	0.72
2:B:101:LEU:CD2	2:B:109:LYS:CG	2.72	0.68
2:B:78[B]:LEU:CD1	2:B:112:LEU:HD12	2.23	0.68
2:B:103:HIS:HA	2:B:107:GLY:H	1.59	0.68
2:B:99:PHE:HB2	5:B:529:HOH:O	1.94	0.68
2:B:78[A]:LEU:HD21	2:B:121:LEU:HD11	1.77	0.67
2:B:81:CYS:SG	2:B:82:LEU:HD12	2.34	0.67
1:A:116:ASN:OD1	1:A:144:THR:CG2	2.42	0.67
2:B:67[B]:ARG:NE	2:B:69:VAL:CG2	2.58	0.67
1:A:61:GLN:C	1:A:63:GLU:CA	2.66	0.66
2:B:56:ASN:HA	2:B:122:ILE:HG23	1.78	0.66
2:B:60:VAL:HG21	2:B:82:LEU:HD11	1.78	0.65
2:B:78[A]:LEU:HD21	2:B:121:LEU:CD1	2.26	0.65
1:A:19:LEU:HD13	1:A:146:ALA:HB2	1.78	0.65
1:A:37:GLU:OE2	2:B:67[B]:ARG:NH1	2.30	0.65
2:B:101:LEU:HD12	2:B:125:GLU:OE2	1.98	0.64
2:B:67[B]:ARG:HH21	2:B:69:VAL:HG21	1.62	0.63
1:A:73:ARG:HH11	1:A:104:LYS:HZ1	1.45	0.63
1:A:85:ASN:H	1:A:85:ASN:HD22	1.45	0.63
2:B:100:ARG:CG	2:B:112:LEU:HD21	2.28	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:ASN:HA	1:A:144:THR:HG22	1.82	0.62
2:B:111:ARG:C	2:B:112:LEU:HD23	2.25	0.61
2:B:78[B]:LEU:HD13	2:B:112:LEU:HD12	1.81	0.61
2:B:100:ARG:HB2	2:B:110:ALA:HB3	1.82	0.61
2:B:100:ARG:O	2:B:110:ALA:HB3	2.01	0.60
2:B:112:LEU:HD23	2:B:112:LEU:N	2.16	0.60
2:B:103:HIS:HB3	2:B:106:LYS:HZ2	1.65	0.60
2:B:59:ARG:HH11	2:B:59:ARG:HG3	1.67	0.60
2:B:67[B]:ARG:NE	2:B:69:VAL:HG22	2.06	0.60
2:B:59:ARG:HG3	2:B:59:ARG:NH1	2.16	0.59
2:B:122:ILE:O	2:B:122:ILE:CG2	2.49	0.59
2:B:58:ILE:HB	2:B:70:VAL:HG23	1.84	0.58
2:B:79:HIS:O	2:B:83:MET:HB2	2.02	0.58
2:B:100:ARG:NE	2:B:112:LEU:HD21	2.18	0.58
2:B:67[B]:ARG:HH21	2:B:69:VAL:CG2	2.16	0.57
2:B:100:ARG:HH22	2:B:116:THR:CG2	2.16	0.56
2:B:122:ILE:O	2:B:122:ILE:HG22	2.06	0.56
1:A:116:ASN:OD1	1:A:144:THR:HG23	2.05	0.56
2:B:102:LEU:HD11	2:B:110:ALA:HB2	1.88	0.56
1:A:12:GLY:HA2	1:A:60:GLY:HA3	1.88	0.56
2:B:100:ARG:HG3	2:B:112:LEU:CD2	2.34	0.56
1:A:73:ARG:NH1	1:A:104:LYS:HE3	2.21	0.55
1:A:25:GLN:CD	1:A:25:GLN:CB	2.72	0.55
1:A:11:ALA:HA	1:A:61:GLN:HG3	1.87	0.54
2:B:78[B]:LEU:HD11	2:B:112:LEU:HD12	1.89	0.53
2:B:59:ARG:HD2	2:B:59:ARG:C	2.34	0.53
2:B:59:ARG:NH2	2:B:125:GLU:HG3	2.21	0.53
2:B:61:PHE:O	2:B:127:GLN:HA	2.09	0.53
1:A:116:ASN:OD1	1:A:144:THR:HG22	2.09	0.52
1:A:100:ILE:HD13	1:A:111:MET:HE2	1.92	0.52
2:B:100:ARG:HE	2:B:112:LEU:HD21	1.74	0.52
2:B:100:ARG:HG2	2:B:126:LEU:CD2	2.40	0.51
1:A:65:SER:O	1:A:67:MET:N	2.40	0.51
1:A:73:ARG:HH11	1:A:104:LYS:NZ	2.09	0.51
1:A:149:ARG:O	1:A:150:GLN:C	2.53	0.51
2:B:101:LEU:CD2	5:B:529:HOH:O	2.57	0.51
1:A:99:GLN:NE2	1:A:102:ARG:NH2	2.56	0.50
2:B:78[A]:LEU:HD12	2:B:112:LEU:HD12	1.93	0.50
1:A:61:GLN:HA	1:A:63:GLU:O	2.12	0.50
2:B:100:ARG:CD	2:B:112:LEU:HD21	2.42	0.50
1:A:32:TYR:C	1:A:32:TYR:CD1	2.90	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:67[B]:ARG:HD2	2:B:68:THR:N	2.28	0.49
1:A:73:ARG:NH1	1:A:104:LYS:CE	2.76	0.49
1:A:23:LEU:HD22	1:A:156:PHE:CG	2.48	0.48
2:B:63:PRO:O	2:B:64:ASN:HB2	2.14	0.48
2:B:64:ASN:O	2:B:65:LYS:CB	2.58	0.48
2:B:127:GLN:NE2	2:B:129:ASP:OD2	2.46	0.48
1:A:41:ARG:HD3	1:A:54:ASP:CG	2.39	0.48
2:B:121:LEU:O	2:B:123:GLY:N	2.37	0.48
2:B:67[B]:ARG:NH2	2:B:69:VAL:CG2	2.76	0.48
2:B:67[A]:ARG:NH1	2:B:67[A]:ARG:CG	2.38	0.48
2:B:119:ALA:C	2:B:121:LEU:H	2.22	0.48
2:B:61:PHE:HB2	2:B:127:GLN:HB2	1.95	0.47
2:B:101:LEU:CD1	2:B:125:GLU:OE2	2.62	0.47
2:B:112:LEU:N	2:B:112:LEU:CD2	2.73	0.47
2:B:84:LYS:O	2:B:88:VAL:HG22	2.14	0.46
2:B:67[B]:ARG:NH2	2:B:69:VAL:HG21	2.30	0.46
2:B:83:MET:HG2	2:B:93:PRO:HG3	1.97	0.46
2:B:99:PHE:HZ	2:B:129:ASP:CB	2.14	0.46
1:A:99:GLN:CD	1:A:102:ARG:NH2	2.74	0.46
1:A:148:THR:HG22	1:A:150:GLN:N	2.15	0.46
2:B:100:ARG:NH2	2:B:116:THR:HG22	2.29	0.45
2:B:65:LYS:C	2:B:67[A]:ARG:HH12	2.23	0.45
2:B:103:HIS:HB3	2:B:106:LYS:NZ	2.31	0.45
2:B:100:ARG:HE	2:B:112:LEU:CD2	2.30	0.45
2:B:62:LEU:N	2:B:66:GLN:O	2.31	0.45
2:B:92:GLN:HB2	2:B:95:CYS:HB2	1.98	0.45
1:A:24:ILE:HG22	1:A:25:GLN:HB2	1.99	0.44
2:B:56:ASN:N	2:B:56:ASN:HD22	2.14	0.44
1:A:73:ARG:HA	1:A:73:ARG:HD2	1.81	0.44
1:A:37:GLU:OE1	1:A:56:LEU:HD21	2.18	0.44
1:A:73:ARG:HH11	1:A:104:LYS:CE	2.30	0.44
1:A:97:ARG:O	1:A:101:LYS:HG3	2.19	0.43
2:B:103:HIS:CB	2:B:106:LYS:NZ	2.80	0.43
1:A:28:PHE:CD1	1:A:147:LYS:HA	2.53	0.43
1:A:25:GLN:CG	1:A:25:GLN:NE2	2.68	0.43
1:A:21:ILE:HD13	1:A:21:ILE:N	2.33	0.42
2:B:103:HIS:HB2	2:B:106:LYS:HZ3	1.85	0.42
2:B:78[B]:LEU:HD21	2:B:98:VAL:HG11	2.01	0.42
2:B:56:ASN:HA	2:B:122:ILE:CG2	2.48	0.42
1:A:59:ALA:O	1:A:60:GLY:C	2.62	0.42
1:A:63:GLU:OE2	1:A:64:TYR:HD2	2.03	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:MET:HE3	1:A:111:MET:HB2	1.89	0.42
2:B:56:ASN:CA	2:B:122:ILE:HG23	2.48	0.42
2:B:74:ASN:HD22	2:B:74:ASN:HA	1.71	0.42
2:B:64:ASN:O	2:B:65:LYS:HD3	2.20	0.41
1:A:36:ILE:O	1:A:67:MET:HE2	2.21	0.41
2:B:78[B]:LEU:CD2	2:B:114:TRP:HA	2.51	0.41
1:A:68:ARG:HB2	1:A:72:MET:HE3	2.01	0.41
1:A:120:LEU:HD23	1:A:120:LEU:HA	1.84	0.41
2:B:59:ARG:HG2	2:B:69:VAL:HG22	2.02	0.41
2:B:100:ARG:HG2	2:B:126:LEU:HD22	2.02	0.40
2:B:92:GLN:HA	2:B:93:PRO:HD2	1.77	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:539:HOH:O	5:A:541:HOH:O[18_444]	1.16	1.04

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	163/166 (98%)	148 (91%)	10 (6%)	5 (3%)	3	0
2	B	76/81 (94%)	59 (78%)	11 (14%)	6 (8%)	1	0
All	All	239/247 (97%)	207 (87%)	21 (9%)	11 (5%)	2	0

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	60	GLY
1	A	66	ALA

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Mol	Chain	Res	Type
1	A	72	MET
2	B	65	LYS
2	B	105	HIS
2	B	120	SER
2	B	122	ILE
1	A	11	ALA
1	A	70	GLN
2	B	117	ASP
2	B	123	GLY

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	143/144 (99%)	120 (84%)	23 (16%)	2	1
2	B	70/73 (96%)	46 (66%)	24 (34%)	0	0
All	All	213/217 (98%)	166 (78%)	47 (22%)	1	0

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	3	GLU
1	A	19	LEU
1	A	21	ILE
1	A	24	ILE
1	A	30	ASP
1	A	31	GLU
1	A	45	VAL
1	A	53	LEU
1	A	56	LEU
1	A	61	GLN
1	A	63	GLU
1	A	68	ARG
1	A	72	MET

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Mol	Chain	Res	Type
1	A	73	ARG
1	A	85	ASN
1	A	97	ARG
1	A	106	SER
1	A	133	LEU
1	A	136	SER
1	A	144	THR
1	A	148	THR
1	A	159	LEU
2	B	56	ASN
2	B	67[A]	ARG
2	B	67[B]	ARG
2	B	73	ARG
2	B	74	ASN
2	B	78[A]	LEU
2	B	78[B]	LEU
2	B	82	LEU
2	B	88	VAL
2	B	94	GLU
2	B	98	VAL
2	B	99	PHE
2	B	102	LEU
2	B	103	HIS
2	B	104	GLU
2	B	105	HIS
2	B	106	LYS
2	B	112	LEU
2	B	115	ASN
2	B	117	ASP
2	B	120	SER
2	B	121	LEU
2	B	129	ASP
2	B	131	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	HIS
1	A	85	ASN
1	A	95	GLN
1	A	99	GLN
1	A	131	GLN

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Mol	Chain	Res	Type
1	A	150	GLN
1	A	165	GLN
2	B	56	ASN
2	B	74	ASN
2	B	103	HIS
2	B	115	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	A	170	-	29,30,30	1.08	2 (6%)	45,47,47	1.69	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	A	170	-	-	4/16/32/32	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	170	GDP	C5-C4	3.30	1.47	1.38
3	A	170	GDP	C6-N1	-2.46	1.34	1.38

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	170	GDP	C5-C4-N3	-4.89	120.61	128.39
3	A	170	GDP	N9-C4-N3	4.01	133.97	125.95
3	A	170	GDP	C2-N3-C4	3.34	118.06	112.30
3	A	170	GDP	C6-C5-N7	3.30	136.30	130.29
3	A	170	GDP	O3'-C3'-C2'	-2.92	102.46	111.82
3	A	170	GDP	C4-C5-N7	-2.40	106.87	110.67
3	A	170	GDP	O3A-PA-O1A	-2.17	104.17	110.70
3	A	170	GDP	C2'-C1'-N9	-2.03	107.59	113.25

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	170	GDP	PA-O3A-PB-O2B
3	A	170	GDP	PA-O3A-PB-O1B
3	A	170	GDP	C5'-O5'-PA-O3A
3	A	170	GDP	C5'-O5'-PA-O1A

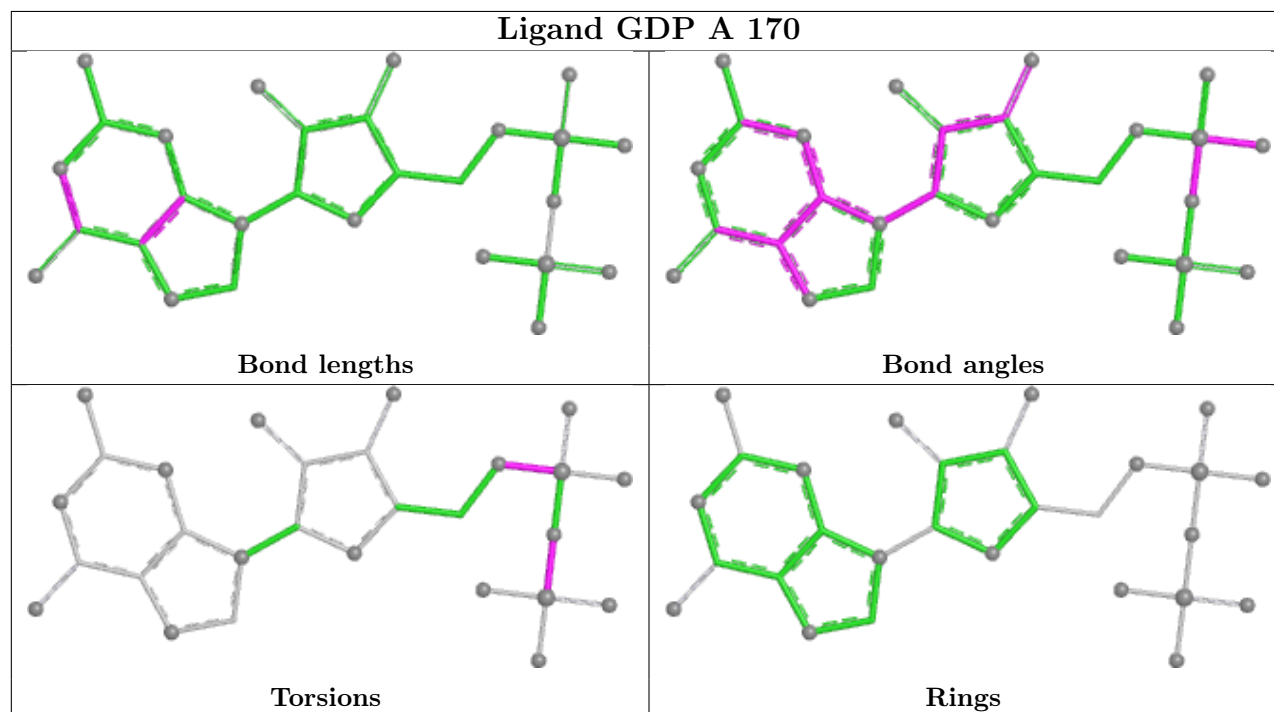
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	170	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, 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	165/166 (99%)	1.63	59 (35%) 1 1	18, 29, 64, 82	0
2	B	76/81 (93%)	2.42	43 (56%) 0 0	25, 52, 61, 65	2 (2%)
All	All	241/247 (97%)	1.88	102 (42%) 0 1	18, 35, 63, 82	2 (0%)

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	66	ALA	7.9
2	B	106	LYS	7.5
2	B	102	LEU	5.8
2	B	122	ILE	5.6
2	B	67[A]	ARG	5.4
2	B	128	VAL	4.9
2	B	105	HIS	4.6
1	A	121	ALA	4.6
1	A	64	TYR	4.5
2	B	131	LEU	4.5
1	A	84	ILE	4.4
1	A	52	LEU	4.2
2	B	78[A]	LEU	4.1
1	A	125	VAL	4.1
2	B	103	HIS	4.0
1	A	120	LEU	3.9
1	A	1	MET	3.9
1	A	28	PHE	3.9
1	A	68	ARG	3.9
2	B	116	THR	3.8
1	A	106	SER	3.8
1	A	105	ASP	3.8
2	B	65	LYS	3.8
1	A	67	MET	3.7

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Mol	Chain	Res	Type	RSRZ
2	B	114	TRP	3.6
1	A	34	PRO	3.6
1	A	130	ALA	3.6
1	A	107	ASP	3.5
1	A	63	GLU	3.5
1	A	60	GLY	3.5
1	A	29	VAL	3.5
1	A	148	THR	3.5
1	A	83	ALA	3.4
1	A	18	ALA	3.4
1	A	72	MET	3.4
1	A	59	ALA	3.4
1	A	32	TYR	3.3
1	A	27	HIS	3.3
1	A	71	TYR	3.3
1	A	122	ALA	3.3
2	B	100	ARG	3.3
2	B	101	LEU	3.3
2	B	108	LYS	3.2
1	A	61	GLN	3.2
1	A	70	GLN	3.2
2	B	61	PHE	3.2
1	A	118	CYS	3.2
2	B	58	ILE	3.1
1	A	147	LYS	3.1
2	B	123	GLY	3.1
2	B	76	MET	3.1
2	B	62	LEU	3.0
2	B	121	LEU	3.0
2	B	60	VAL	3.0
2	B	91	LEU	3.0
1	A	108	ASP	3.0
1	A	25	GLN	2.9
1	A	74	THR	2.9
1	A	75	GLY	2.9
2	B	107	GLY	2.9
1	A	30	ASP	2.9
2	B	117	ASP	2.9
1	A	36	ILE	2.9
1	A	124	THR	2.9
1	A	127	SER	2.8
2	B	98	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	62	GLU	2.8
1	A	129	GLN	2.6
1	A	102	ARG	2.6
2	B	64	ASN	2.6
1	A	78	PHE	2.5
2	B	69	VAL	2.5
2	B	92	GLN	2.5
1	A	37	GLU	2.5
1	A	69	ASP	2.4
1	A	123	ARG	2.4
2	B	72	VAL	2.4
1	A	144	THR	2.4
2	B	86	LEU	2.4
2	B	129	ASP	2.4
1	A	31	GLU	2.3
2	B	130	PHE	2.3
2	B	124	GLU	2.3
2	B	74	ASN	2.3
1	A	132	ASP	2.3
1	A	151	GLY	2.2
2	B	99	PHE	2.2
2	B	81	CYS	2.2
2	B	79	HIS	2.2
1	A	40	TYR	2.2
2	B	88	VAL	2.2
1	A	3	GLU	2.2
1	A	65	SER	2.1
1	A	115	GLY	2.1
1	A	85	ASN	2.1
2	B	89	ARG	2.1
1	A	38	ASP	2.1
1	A	119	ASP	2.0
2	B	127	GLN	2.0
2	B	73	ARG	2.0
1	A	117	LYS	2.0
2	B	126	LEU	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 [i](#)

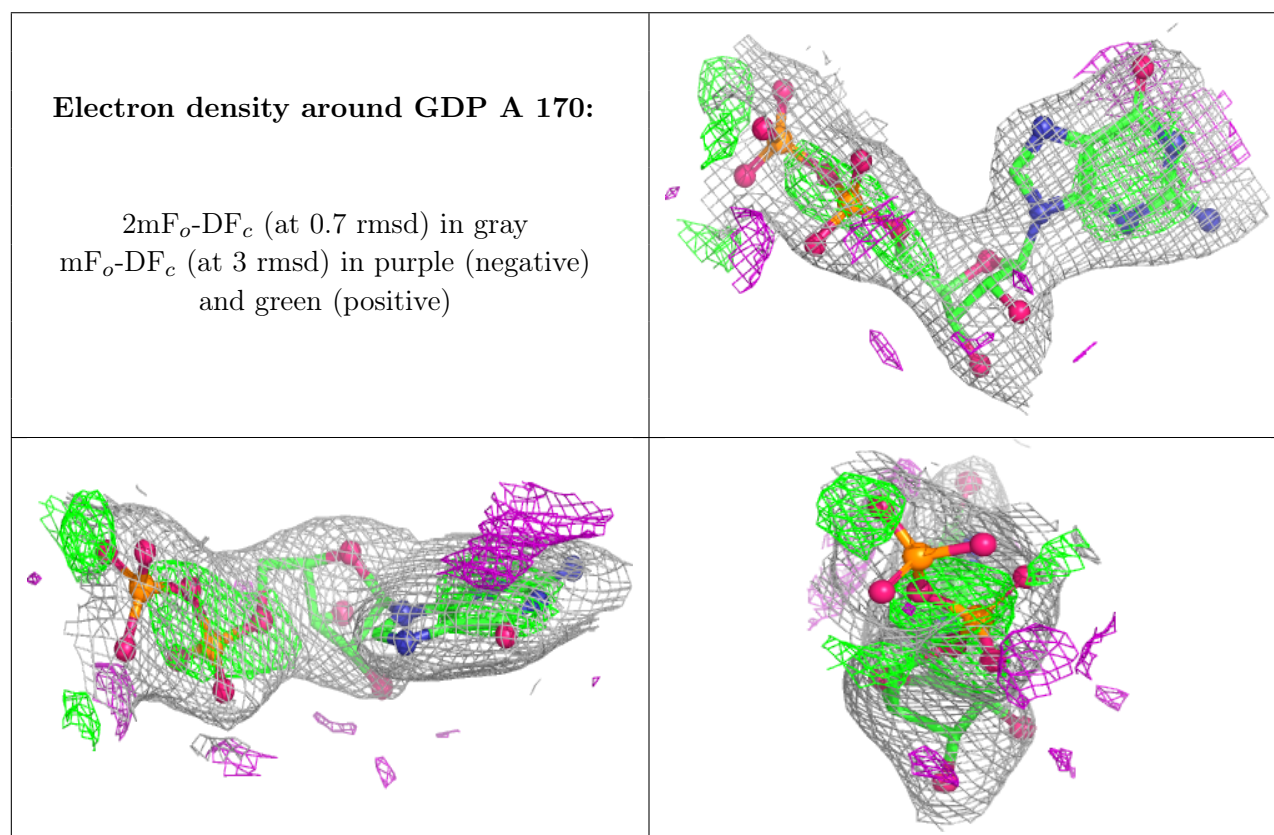
There are no oligosaccharides in this entry.

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($\text{\AA}^2$)	Q<0.9
4	MG	A	171	1/1	0.89	0.12	39,39,39,39	0
3	GDP	A	170	28/28	0.91	0.14	17,24,29,31	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.