



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:25 PM GMT

PDB ID : 1K8R
Title : Crystal structure of Ras-Bry2RBD complex
Authors : Scheffzek, K.; Gruenewald, P.; Wohlgemuth, S.; Kabsch, W.; Tu, H.; Wigler, M.; Wittinghofer, A.; Herrmann, C.
Deposited on : 2001-10-25
Resolution : 3.00 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

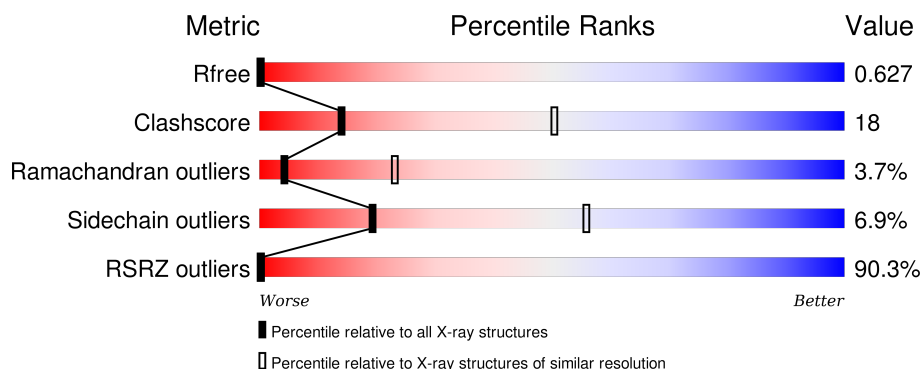
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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}	91344	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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. 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	<div> <div>86%</div> <div>63% 32% 5%</div> </div>
2	B	110	<div> <div>75%</div> <div>39% 30% 5% 25%</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
4	GNP	A	167	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2012 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 Transforming protein P21/H-RAS-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	166	Total	C	N	O	S	0	0	0
			1319	822	228	263	6			

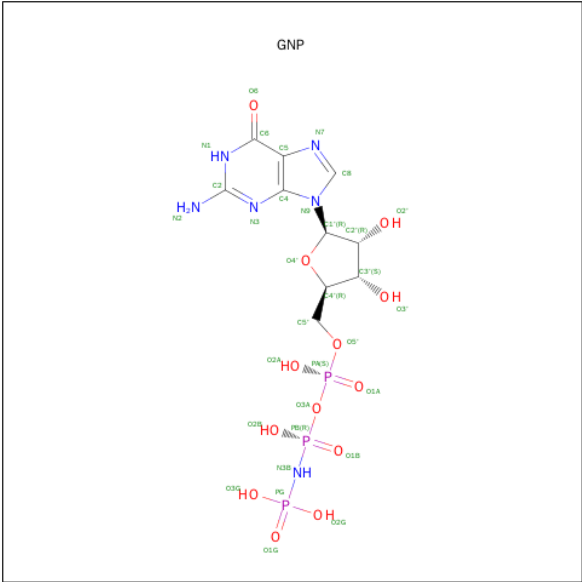
- Molecule 2 is a protein called Protein kinase byr2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	82	Total	C	N	O	S	0	0	0
			660	416	119	121	4			

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 4 is PHOSPHOAMINOPHOSPHONIC ACID-GUANYLATE ESTER (three-letter code: GNP) (formula: C₁₀H₁₇N₆O₁₃P₃).

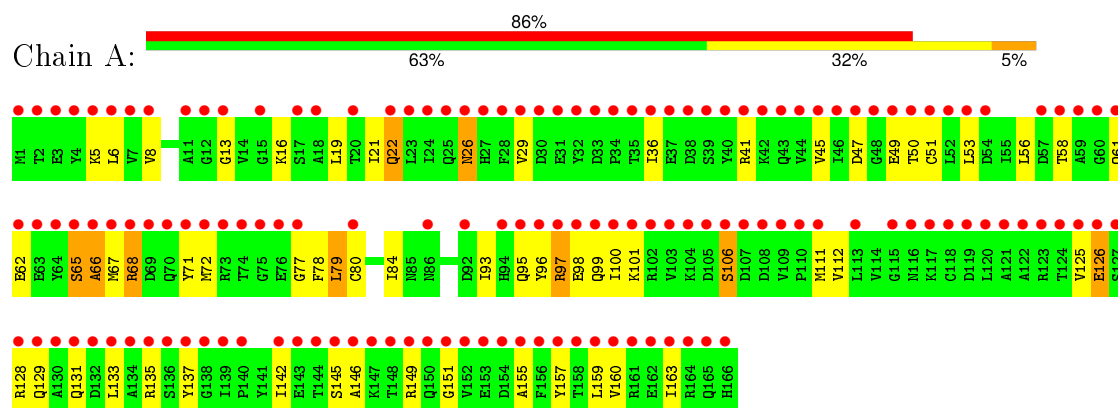


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			32	10	6	13	3		

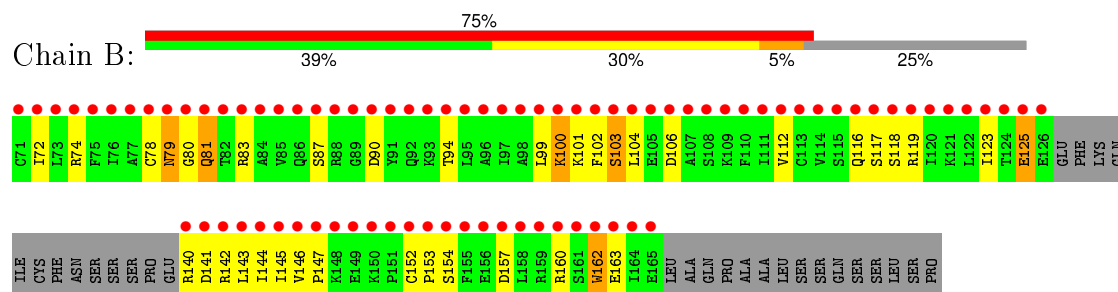
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ($\text{RSRZ} > 2$). 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: Transforming protein P21/H-RAS-1



• Molecule 2: Protein kinase byr2



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	139.03Å 139.03Å 112.80Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.00 – 3.00 60.20 – 1.73	Depositor EDS
% Data completeness (in resolution range)	99.9 (25.00-3.00) 18.6 (60.20-1.73)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.46 (at 1.73Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.235 , 0.305 0.564 , 0.627	Depositor DCC
R_{free} test set	176 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	4.4	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , -1.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	5 of 8167 reflections (0.061%)	Xtriage
F_o, F_c correlation	0.08	EDS
Total number of atoms	2012	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 29.89 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.4336e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: GNP, MG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.44	0/1338	0.67	1/1807 (0.1%)
2	B	0.41	0/668	0.66	0/894
All	All	0.43	0/2006	0.67	1/2701 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	79	LEU	CA-CB-CG	6.15	129.44	115.30

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	1319	0	1289	49	0
2	B	660	0	680	31	0
3	A	1	0	0	0	0
4	A	32	0	13	1	0
All	All	2012	0	1982	73	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 18.

All (73) 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:41:ARG:HG2	2:B:81:GLN:HE21	1.38	0.89
1:A:41:ARG:HH21	2:B:80:GLY:HA3	1.47	0.79
1:A:6:LEU:HG	1:A:159:LEU:HD12	1.64	0.77
1:A:8:VAL:HG22	1:A:79:LEU:HD21	1.68	0.75
2:B:146:VAL:HG22	2:B:147:PRO:HD2	1.68	0.74
1:A:68:ARG:HA	1:A:71:TYR:CZ	2.23	0.74
1:A:112:VAL:HG23	1:A:159:LEU:HD23	1.71	0.70
1:A:41:ARG:HG2	2:B:81:GLN:NE2	2.11	0.66
1:A:68:ARG:NH1	1:A:99:GLN:HG2	2.14	0.62
1:A:41:ARG:H	2:B:81:GLN:HE22	1.46	0.61
2:B:146:VAL:CG2	2:B:147:PRO:HD2	2.32	0.59
2:B:160:ARG:O	2:B:163:GLU:HG2	2.02	0.59
2:B:144:ILE:HG21	2:B:153:PRO:HB3	1.85	0.59
1:A:157:TYR:O	1:A:160:VAL:HG22	2.03	0.58
1:A:68:ARG:HH22	1:A:99:GLN:HB3	1.68	0.57
1:A:101:LYS:HB2	1:A:106:SER:O	2.04	0.57
2:B:154:SER:HB3	2:B:157:ASP:OD2	2.04	0.57
2:B:162:TRP:HD1	2:B:163:GLU:N	2.02	0.56
1:A:26:ASN:C	1:A:26:ASN:HD22	2.07	0.56
2:B:162:TRP:CD1	2:B:163:GLU:N	2.74	0.56
2:B:78:CYS:O	2:B:79:ASN:HB2	2.08	0.54
2:B:102:PHE:O	2:B:103:SER:HB2	2.08	0.54
1:A:41:ARG:N	2:B:81:GLN:HE22	2.06	0.53
1:A:21:ILE:HD12	1:A:29:VAL:HG21	1.90	0.53
1:A:80:CYS:HB3	1:A:93:ILE:HD12	1.91	0.53
1:A:97:ARG:HG2	1:A:97:ARG:HH21	1.73	0.53
1:A:41:ARG:HH21	2:B:80:GLY:CA	2.21	0.52
1:A:84:ILE:HA	1:A:125:VAL:CG1	2.40	0.52
2:B:125:GLU:OE1	2:B:125:GLU:HA	2.11	0.51
2:B:74:ARG:HD2	2:B:141:ASP:O	2.11	0.51
2:B:72:ILE:HG23	2:B:72:ILE:O	2.10	0.51
2:B:142:ARG:O	2:B:143:LEU:HD23	2.11	0.51
1:A:135:ARG:HG2	1:A:135:ARG:HH21	1.77	0.50
2:B:99:LEU:HD22	2:B:104:LEU:HB2	1.95	0.49
1:A:45:VAL:O	1:A:45:VAL:HG13	2.13	0.48
1:A:22:GLN:O	1:A:26:ASN:HA	2.13	0.48
1:A:16:LYS:HE3	1:A:58:THR:HG22	1.95	0.48
1:A:77:GLY:HA3	1:A:163:ILE:HD11	1.96	0.48
1:A:72:MET:HE2	1:A:100:ILE:HA	1.97	0.47
1:A:5:LYS:O	1:A:6:LEU:HD12	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:GLU:N	1:A:129:GLN:OE1	2.40	0.47
1:A:56:LEU:CD2	1:A:71:TYR:HB2	2.45	0.47
1:A:97:ARG:CG	1:A:97:ARG:HH21	2.28	0.46
1:A:128:ARG:HD2	1:A:131:GLN:OE1	2.16	0.46
1:A:13:GLY:H	4:A:167:GNP:HNB3	1.63	0.46
1:A:41:ARG:HA	1:A:53:LEU:O	2.15	0.46
1:A:19:LEU:HD23	1:A:146:ALA:HB2	1.98	0.46
2:B:90:ASP:C	2:B:90:ASP:OD2	2.55	0.46
1:A:71:TYR:CE1	1:A:72:MET:HG3	2.52	0.45
1:A:96:TYR:O	1:A:100:ILE:HG13	2.17	0.44
2:B:116:GLN:O	2:B:118:SER:N	2.49	0.44
2:B:118:SER:O	2:B:119:ARG:HB3	2.18	0.44
1:A:128:ARG:NH2	1:A:131:GLN:OE1	2.51	0.44
1:A:6:LEU:HG	1:A:159:LEU:CD1	2.41	0.44
1:A:8:VAL:HG22	1:A:79:LEU:CD2	2.42	0.44
2:B:112:VAL:O	2:B:123:ILE:HG22	2.18	0.44
1:A:53:LEU:O	1:A:53:LEU:HD12	2.17	0.44
2:B:81:GLN:H	2:B:81:GLN:HG2	1.54	0.44
1:A:68:ARG:HA	1:A:71:TYR:CE2	2.53	0.43
2:B:83:ARG:HD2	2:B:101:LYS:HE3	2.01	0.43
2:B:99:LEU:HD21	2:B:145:ILE:HG21	2.00	0.43
1:A:5:LYS:C	1:A:6:LEU:HD12	2.39	0.43
1:A:145:SER:O	1:A:149:ARG:N	2.51	0.43
2:B:100:LYS:O	2:B:100:LYS:HD3	2.18	0.43
1:A:142:ILE:HD12	1:A:155:ALA:HA	2.00	0.42
2:B:87:SER:HB2	2:B:94:THR:OG1	2.19	0.42
1:A:78:PHE:O	1:A:111:MET:HA	2.19	0.42
1:A:133:LEU:HG	1:A:137:TYR:CE2	2.55	0.42
1:A:36:ILE:HG21	2:B:72:ILE:HD11	2.01	0.41
1:A:98:GLU:CD	1:A:101:LYS:HE2	2.41	0.41
1:A:47:ASP:C	1:A:49:GLU:H	2.23	0.41
1:A:65:SER:O	1:A:66:ALA:C	2.58	0.41
2:B:152:CYS:HA	2:B:153:PRO:HD2	1.91	0.41

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	164/166 (99%)	145 (88%)	15 (9%)	4 (2%)	7	35
2	B	78/110 (71%)	60 (77%)	13 (17%)	5 (6%)	2	9
All	All	242/276 (88%)	205 (85%)	28 (12%)	9 (4%)	4	23

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	66	ALA
2	B	106	ASP
1	A	65	SER
2	B	81	GLN
2	B	117	SER
2	B	79	ASN
1	A	67	MET
2	B	103	SER
1	A	151	GLY

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	143/144 (99%)	132 (92%)	11 (8%)	16	50
2	B	74/100 (74%)	70 (95%)	4 (5%)	27	66
All	All	217/244 (89%)	202 (93%)	15 (7%)	19	56

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	26	ASN
1	A	50	THR
1	A	51	CYS
1	A	61	GLN
1	A	62	GLU
1	A	68	ARG
1	A	95	GLN
1	A	97	ARG
1	A	106	SER
1	A	126	GLU
2	B	100	LYS
2	B	125	GLU
2	B	140	ARG
2	B	162	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	GLN
1	A	26	ASN
1	A	70	GLN
1	A	95	GLN
2	B	81	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

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$
4	GNP	A	167	3	28,34,34	1.72	6 (21%)	33,54,54	2.36	8 (24%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GNP	A	167	3	-	1/12/38/38	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	167	GNP	PB-O2B	-4.16	1.45	1.56
4	A	167	GNP	PG-O2G	-4.09	1.45	1.56
4	A	167	GNP	C8-N7	-3.37	1.28	1.34
4	A	167	GNP	C6-C5	2.32	1.45	1.41
4	A	167	GNP	C4-N3	2.55	1.39	1.35
4	A	167	GNP	C6-N1	3.25	1.39	1.33

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	167	GNP	C5-C6-N1	-7.53	113.30	123.59
4	A	167	GNP	O3G-PG-O1G	-4.11	102.56	113.49
4	A	167	GNP	N3-C2-N1	-3.70	121.82	127.44
4	A	167	GNP	C2'-C1'-N9	-2.99	109.72	114.29
4	A	167	GNP	C6-C5-C4	-2.25	118.20	120.90
4	A	167	GNP	O1B-PB-N3B	2.12	115.15	111.90
4	A	167	GNP	O1G-PG-N3B	2.48	115.69	111.90
4	A	167	GNP	C6-N1-C2	7.56	126.43	115.94

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	167	GNP	O1B-PB-N3B-PG

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	167	GNP	1	0

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	166/166 (100%)	3.90	142 (85%)  	14, 44, 76, 82	0
2	B	82/110 (74%)	5.42	82 (100%)  	35, 63, 79, 86	0
All	All	248/276 (89%)	4.41	224 (90%)  	14, 51, 77, 86	0

All (224) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	118	SER	9.4
2	B	162	TRP	8.4
1	A	108	ASP	8.4
1	A	71	TYR	8.3
1	A	2	THR	8.3
2	B	113	CYS	8.1
2	B	87	SER	7.9
1	A	103	VAL	7.8
1	A	74	THR	7.8
2	B	124	THR	7.7
1	A	166	HIS	7.6
2	B	117	SER	7.5
2	B	141	ASP	7.4
2	B	120	ILE	7.4
2	B	90	ASP	7.3
2	B	154	SER	7.3
2	B	115	SER	7.2
1	A	1	MET	7.1
2	B	116	GLN	7.1
2	B	151	PRO	7.1
2	B	112	VAL	7.1
1	A	65	SER	7.0
2	B	153	PRO	7.0
2	B	104	LEU	6.9

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Mol	Chain	Res	Type	RSRZ
2	B	161	SER	6.8
2	B	125	GLU	6.7
1	A	44	VAL	6.7
2	B	106	ASP	6.7
1	A	64	TYR	6.7
2	B	89	GLY	6.7
2	B	122	LEU	6.6
1	A	105	ASP	6.6
1	A	107	ASP	6.5
2	B	114	VAL	6.4
1	A	61	GLN	6.4
2	B	107	ALA	6.4
2	B	123	ILE	6.3
2	B	147	PRO	6.3
2	B	94	THR	6.3
2	B	95	LEU	6.3
2	B	71	CYS	6.3
1	A	165	GLN	6.3
1	A	69	ASP	6.2
1	A	43	GLN	6.2
2	B	91	TYR	6.1
1	A	72	MET	6.1
1	A	106	SER	6.1
1	A	67	MET	6.0
2	B	126	GLU	6.0
1	A	27	HIS	5.9
2	B	105	GLU	5.9
2	B	103	SER	5.8
1	A	46	ILE	5.8
2	B	86	GLN	5.8
1	A	70	GLN	5.8
2	B	152	CYS	5.8
2	B	165	GLU	5.8
1	A	47	ASP	5.8
2	B	164	ILE	5.7
1	A	50	THR	5.7
2	B	121	LYS	5.7
2	B	119	ARG	5.7
1	A	99	GLN	5.7
2	B	150	LYS	5.6
2	B	108	SER	5.6
1	A	124	THR	5.6

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Mol	Chain	Res	Type	RSRZ
1	A	63	GLU	5.5
2	B	163	GLU	5.5
2	B	73	LEU	5.4
1	A	51	CYS	5.4
2	B	155	PHE	5.4
1	A	32	TYR	5.4
1	A	5	LYS	5.4
2	B	157	ASP	5.4
1	A	75	GLY	5.3
2	B	72	ILE	5.3
2	B	80	GLY	5.3
1	A	45	VAL	5.3
1	A	73	ARG	5.3
1	A	31	GLU	5.2
1	A	66	ALA	5.2
1	A	60	GLY	5.1
1	A	148	THR	5.1
1	A	3	GLU	5.1
2	B	158	LEU	5.1
1	A	68	ARG	5.1
1	A	150	GLN	5.1
1	A	154	ASP	5.1
1	A	94	HIS	5.0
1	A	52	LEU	5.0
2	B	97	ILE	5.0
1	A	62	GLU	5.0
1	A	157	TYR	5.0
1	A	100	ILE	4.9
2	B	92	GLN	4.9
1	A	136	SER	4.9
1	A	48	GLY	4.9
2	B	88	ARG	4.9
1	A	25	GLN	4.8
2	B	110	PHE	4.8
1	A	109	VAL	4.8
1	A	30	ASP	4.8
2	B	149	GLU	4.7
1	A	135	ARG	4.7
2	B	111	ILE	4.7
2	B	77	ALA	4.6
2	B	93	LYS	4.6
1	A	121	ALA	4.6

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Mol	Chain	Res	Type	RSRZ
2	B	79	ASN	4.6
1	A	95	GLN	4.6
1	A	153	GLU	4.6
1	A	76	GLU	4.6
1	A	102	ARG	4.6
1	A	137	TYR	4.5
2	B	96	ALA	4.5
1	A	151	GLY	4.5
2	B	78	CYS	4.5
2	B	140	ARG	4.5
1	A	163	ILE	4.5
2	B	109	LYS	4.4
1	A	4	TYR	4.4
1	A	26	ASN	4.4
1	A	35	THR	4.3
1	A	164	ARG	4.3
1	A	54	ASP	4.3
1	A	104	LYS	4.3
1	A	110	PRO	4.3
1	A	138	GLY	4.2
2	B	148	LYS	4.2
1	A	149	ARG	4.1
1	A	49	GLU	4.1
1	A	162	GLU	4.1
1	A	118	CYS	4.1
1	A	101	LYS	4.1
1	A	42	LYS	4.1
1	A	161	ARG	4.1
1	A	98	GLU	4.0
2	B	159	ARG	4.0
2	B	100	LYS	4.0
2	B	146	VAL	4.0
2	B	144	ILE	4.0
1	A	132	ASP	4.0
2	B	98	ALA	4.0
1	A	119	ASP	4.0
1	A	120	LEU	4.0
1	A	29	VAL	4.0
1	A	34	PRO	4.0
1	A	96	TYR	3.9
1	A	158	THR	3.9
2	B	143	LEU	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	28	PHE	3.8
1	A	134	ALA	3.8
2	B	142	ARG	3.8
2	B	145	ILE	3.8
2	B	85	VAL	3.8
1	A	59	ALA	3.8
1	A	122	ALA	3.7
1	A	127	SER	3.7
1	A	128	ARG	3.7
2	B	160	ARG	3.6
2	B	156	GLU	3.6
1	A	57	ASP	3.6
1	A	77	GLY	3.5
2	B	84	ALA	3.5
2	B	74	ARG	3.5
1	A	24	ILE	3.5
2	B	99	LEU	3.5
1	A	145	SER	3.5
1	A	111	MET	3.5
1	A	17	SER	3.4
2	B	76	ILE	3.4
1	A	40	TYR	3.4
1	A	33	ASP	3.4
2	B	82	THR	3.4
1	A	159	LEU	3.4
1	A	37	GLU	3.3
1	A	58	THR	3.3
1	A	155	ALA	3.3
1	A	41	ARG	3.3
1	A	140	PRO	3.3
1	A	133	LEU	3.3
1	A	97	ARG	3.2
1	A	131	GLN	3.2
1	A	126	GLU	3.2
1	A	116	ASN	3.2
1	A	39	SER	3.2
1	A	147	LYS	3.1
2	B	101	LYS	3.1
1	A	123	ARG	3.1
2	B	81	GLN	3.1
1	A	92	ASP	3.1
1	A	156	PHE	3.0

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Mol	Chain	Res	Type	RSRZ
2	B	102	PHE	3.0
1	A	129	GLN	3.0
1	A	115	GLY	2.9
1	A	38	ASP	2.9
1	A	53	LEU	2.8
1	A	18	ALA	2.8
1	A	139	ILE	2.8
1	A	6	LEU	2.8
1	A	160	VAL	2.7
1	A	7	VAL	2.7
1	A	15	GLY	2.7
2	B	75	PHE	2.7
1	A	146	ALA	2.7
1	A	23	LEU	2.6
1	A	143	GLU	2.6
1	A	22	GLN	2.6
1	A	11	ALA	2.6
1	A	130	ALA	2.5
1	A	80	CYS	2.5
1	A	125	VAL	2.4
1	A	144	THR	2.4
1	A	152	VAL	2.4
1	A	12	GLY	2.4
1	A	142	ILE	2.3
1	A	20	THR	2.3
1	A	36	ILE	2.2
1	A	13	GLY	2.2
1	A	86	ASN	2.1
1	A	8	VAL	2.1
1	A	117	LYS	2.0
2	B	83	ARG	2.0
1	A	113	LEU	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	LLDF	B-factors(\AA^2)	Q<0.9
4	GNP	A	167	32/32	0.58	0.42	-1.19	20,23,36,37	0
3	MG	A	168	1/1	0.80	0.32	-2.66	41,41,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.