Full wwPDB X-ray Structure Validation Report (i)

Feb 19, 2016 – 06:44 PM GMT

PDB ID	:	3X04
Title	:	Crystal structure of PIP4KIIBETA complex with GMPPNP
Authors	:	Takeuchi, K.; Lo, Y.H.; Sumita, K.; Senda, M.; Terakawa, J.; Dimitoris, A.; Locasale, J.W.; Sasaki, M.; Yoshino, H.; Zhang, Y.; Kahoud, E.R.; Takano, T.; Yokota, T.; Emerling, B.; Asara, J.A.; Ishida, T.; Shimada, I.; Daikoku, T.; Cantley, L.C.; Senda, T.; Sasaki, A.T.
Deposited on	:	2014-10-09
Resolution	:	2.60 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 (i) symbol.

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

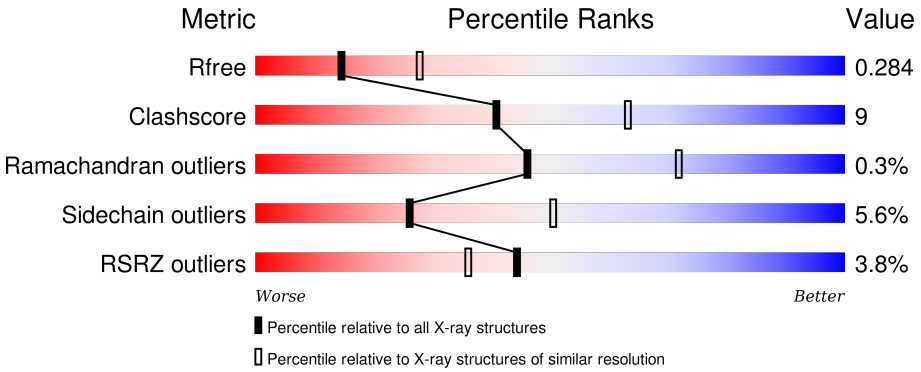
MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026982
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20026982

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 2.60 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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 <=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	393	% 65% 13% • 20%
1	B	393	5% 56% 18% • 24%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5172 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.

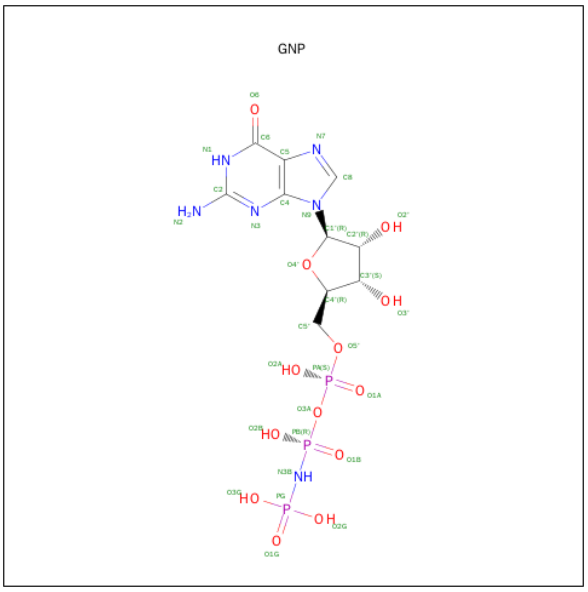
- Molecule 1 is a protein called Phosphatidylinositol 5-phosphate 4-kinase type-2 beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	314	Total	C	N	O	S	0	0	0
			2579	1643	440	482	14			
1	B	297	Total	C	N	O	S	0	0	0
			2456	1576	419	448	13			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	GLY	-	EXPRESSION TAG	UNP P78356
A	25	PRO	-	EXPRESSION TAG	UNP P78356
A	26	ASN	-	EXPRESSION TAG	UNP P78356
A	27	CYS	-	EXPRESSION TAG	UNP P78356
A	28	ALA	-	EXPRESSION TAG	UNP P78356
A	29	PRO	-	EXPRESSION TAG	UNP P78356
A	30	GLY	-	EXPRESSION TAG	UNP P78356
B	24	GLY	-	EXPRESSION TAG	UNP P78356
B	25	PRO	-	EXPRESSION TAG	UNP P78356
B	26	ASN	-	EXPRESSION TAG	UNP P78356
B	27	CYS	-	EXPRESSION TAG	UNP P78356
B	28	ALA	-	EXPRESSION TAG	UNP P78356
B	29	PRO	-	EXPRESSION TAG	UNP P78356
B	30	GLY	-	EXPRESSION TAG	UNP P78356

- Molecule 2 is PHOSPHOAMINOPHOSPHONIC ACID-GUANYLATE ESTER (three-letter code: GNP) (formula: $C_{10}H_{17}N_6O_{13}P_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C N O P 32 10 6 13 3	0	0
2	A	1	Total C N O P 32 10 6 13 3	0	0
2	A	1	Total C N O P 24 10 5 8 1	0	0
2	B	1	Total C N O P 32 10 6 13 3	0	0

- Molecule 3 is water.

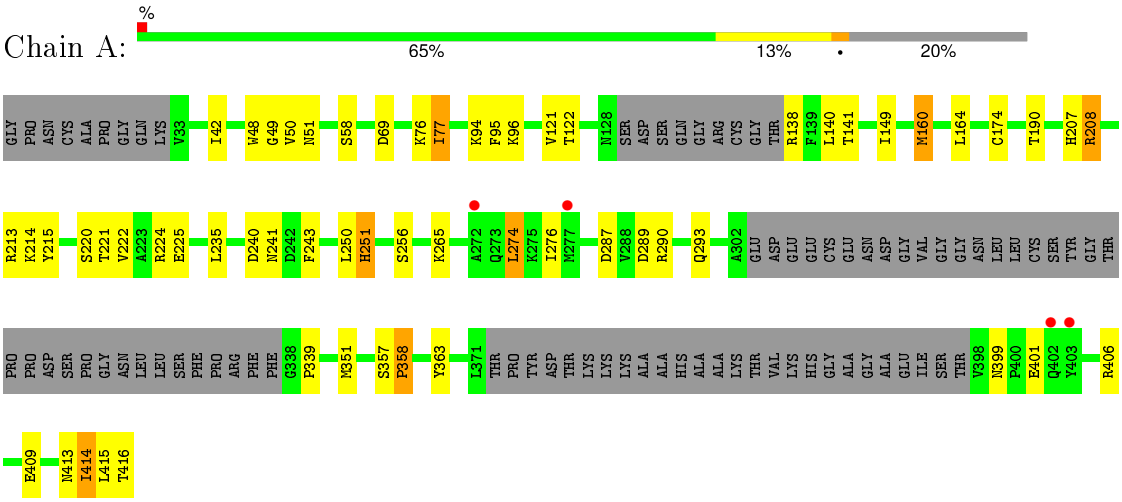
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	9	Total O 9 9	0	0
3	B	8	Total O 8 8	0	0

3 Residue-property plots (i)

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 (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.

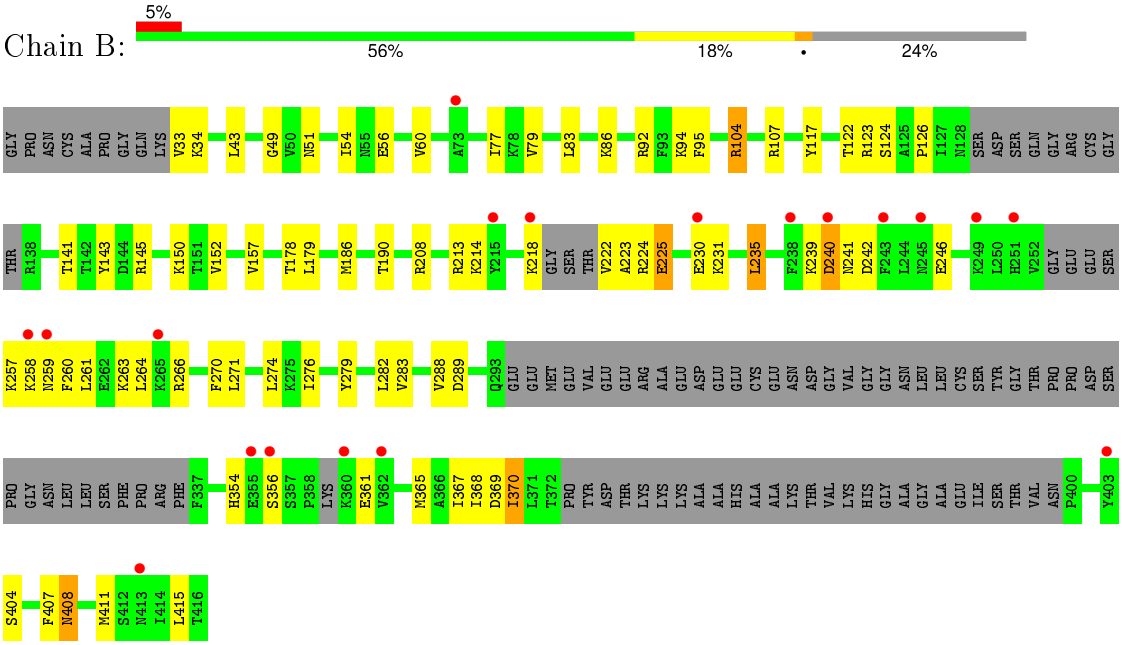
- Molecule 1: Phosphatidylinositol 5-phosphate 4-kinase type-2 beta

Chain A:



- Molecule 1: Phosphatidylinositol 5-phosphate 4-kinase type-2 beta

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	109.80Å 185.22Å 106.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	94.45 – 2.60 94.45 – 2.60	Depositor EDS
% Data completeness (in resolution range)	97.9 (94.45-2.60) 97.8 (94.45-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.02 (at 2.62Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.7.3_928)	Depositor
R, R_{free}	0.229 , 0.277 0.229 , 0.284	Depositor DCC
R_{free} test set	1652 reflections (5.27%)	DCC
Wilson B-factor (Å^2)	71.1	Xtriage
Anisotropy	0.152	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 67.6	EDS
Estimated twinning fraction	0.004 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.016 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 33029 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5172	wwPDB-VP
Average B, all atoms (Å^2)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: GNP

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	0.46	0/2632	0.57	0/3542
1	B	0.42	0/2507	0.58	0/3369
All	All	0.44	0/5139	0.58	0/6911

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2579	0	2548	27	0
1	B	2456	0	2444	65	0
2	A	88	0	38	1	0
2	B	32	0	13	2	0
3	A	9	0	0	0	0
3	B	8	0	0	0	0
All	All	5172	0	5043	88	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 9.

All (88) 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:B:224:ARG:HD3	1:B:239:LYS:NZ	1.55	1.21
1:B:222:VAL:HG12	1:B:223:ALA:N	1.56	1.17
1:B:222:VAL:CG1	1:B:223:ALA:H	1.55	1.13
1:B:224:ARG:HD3	1:B:239:LYS:HZ2	1.09	1.02
1:B:224:ARG:CD	1:B:239:LYS:HZ2	1.83	0.91
1:B:224:ARG:CD	1:B:239:LYS:NZ	2.35	0.90
1:B:222:VAL:HG12	1:B:223:ALA:H	0.70	0.83
1:B:33:VAL:HG12	1:B:33:VAL:O	1.85	0.74
1:B:157:VAL:HA	1:B:186:MET:HE1	1.68	0.73
1:A:48:TRP:NE1	1:B:56:GLU:OE2	2.25	0.70
1:A:224:ARG:NH1	1:A:240:ASP:OD1	2.23	0.69
1:B:94:LYS:HB2	1:B:190:THR:HB	1.76	0.68
1:B:257:LYS:HE2	1:B:261:LEU:HD11	1.76	0.68
1:B:283:VAL:HG22	1:B:365:MET:HG2	1.77	0.67
1:B:214:LYS:HG2	1:B:235:LEU:HD22	1.77	0.67
1:B:224:ARG:HD3	1:B:239:LYS:CE	2.26	0.65
1:A:96:LYS:NZ	2:A:502:GNP:O2A	2.29	0.65
1:B:141:THR:HG23	1:B:145:ARG:HA	1.82	0.62
1:B:271:LEU:HD21	1:B:370:ILE:HD12	1.84	0.60
1:B:282:LEU:HB2	1:B:368:ILE:HD13	1.85	0.59
1:A:94:LYS:HB2	1:A:190:THR:HB	1.85	0.58
1:B:225:GLU:HA	1:B:239:LYS:HB2	1.88	0.56
1:B:369:ASP:OD1	2:B:501:GNP:O1A	2.23	0.56
1:B:60:VAL:O	1:B:104:ARG:NH2	2.36	0.56
1:B:224:ARG:HD3	1:B:239:LYS:HZ3	1.59	0.54
1:B:224:ARG:HD3	1:B:239:LYS:CD	2.37	0.53
1:B:224:ARG:HB3	1:B:239:LYS:HD3	1.90	0.53
1:B:223:ALA:HB1	1:B:241:ASN:HD21	1.73	0.53
1:B:150:LYS:HD3	1:B:152:VAL:HG12	1.90	0.53
1:B:224:ARG:CD	1:B:239:LYS:HZ3	2.16	0.53
1:B:279:TYR:HB3	1:B:370:ILE:HG22	1.91	0.52
1:A:42:ILE:H	1:A:42:ILE:HD12	1.73	0.52
1:A:220:SER:HA	1:A:406:ARG:HD3	1.91	0.52
1:A:274:LEU:HB2	1:A:276:ILE:HG12	1.92	0.52
1:B:407:PHE:O	1:B:411:MET:HG2	2.10	0.51
1:A:256:SER:HB3	1:A:351:MET:CE	2.40	0.51
1:B:242:ASP:O	1:B:246:GLU:HB2	2.11	0.51
1:A:287:ASP:HB3	1:A:290:ARG:HB3	1.93	0.50
1:B:218:LYS:HG3	1:B:239:LYS:HZ1	1.77	0.50
1:B:240:ASP:N	1:B:240:ASP:OD1	2.33	0.50
1:B:141:THR:CG2	1:B:145:ARG:HA	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:276:ILE:HG21	1:B:370:ILE:O	2.12	0.50
1:B:354:HIS:CD2	1:B:356:SER:HB3	2.47	0.49
1:B:264:LEU:CD1	1:B:367:ILE:HD11	2.42	0.49
1:A:250:LEU:HD22	1:A:363:TYR:CE2	2.48	0.49
1:B:276:ILE:HD13	1:B:370:ILE:O	2.13	0.49
1:A:160:MET:HE1	1:A:164:LEU:HD13	1.94	0.48
1:B:213:ARG:NH2	1:B:246:GLU:OE1	2.45	0.48
1:B:369:ASP:OD2	2:B:501:GNP:O2A	2.32	0.48
1:A:357:SER:OG	1:A:358:PRO:HD2	2.14	0.47
1:A:50:VAL:HG21	1:A:121:VAL:HG11	1.96	0.47
1:B:222:VAL:CG1	1:B:223:ALA:N	2.30	0.47
1:A:69:ASP:HB3	1:B:83:LEU:HD13	1.95	0.47
1:B:288:VAL:HG23	1:B:361:GLU:HA	1.97	0.46
1:B:279:TYR:CB	1:B:370:ILE:HG22	2.46	0.46
1:B:178:THR:OG1	1:B:179:LEU:N	2.48	0.46
1:B:43:LEU:HD22	1:B:126:PRO:HG2	1.98	0.45
1:B:239:LYS:HB3	1:B:239:LYS:HE2	1.83	0.44
1:B:225:GLU:HB2	1:B:242:ASP:OD1	2.18	0.44
1:A:77:ILE:HG12	1:B:77:ILE:HD11	2.00	0.44
1:B:141:THR:HG21	1:B:145:ARG:HG2	2.00	0.43
1:A:51:ASN:HB2	1:A:122:THR:HG21	1.99	0.43
1:A:251:HIS:ND1	1:A:416:THR:O	2.51	0.43
1:B:266:ARG:HB3	1:B:266:ARG:HE	1.56	0.43
1:B:214:LYS:HE2	1:B:235:LEU:HD13	2.00	0.43
1:B:270:PHE:CZ	1:B:274:LEU:HD11	2.53	0.42
1:B:49:GLY:HA3	1:B:95:PHE:CE1	2.55	0.42
1:B:104:ARG:HA	1:B:107:ARG:NH1	2.35	0.42
1:B:54:ILE:HD11	1:B:117:TYR:CD1	2.54	0.42
1:A:207:HIS:NE2	1:A:208:ARG:HD3	2.35	0.42
1:A:243:PHE:CZ	1:A:414:ILE:HG12	2.55	0.42
1:A:409:GLU:O	1:A:413:ASN:ND2	2.44	0.42
1:B:124:SER:HB3	1:B:143:TYR:CG	2.55	0.41
1:A:290:ARG:O	1:A:293:GLN:HB3	2.20	0.41
1:A:213:ARG:HD3	1:A:215:TYR:OH	2.20	0.41
1:A:214:LYS:HG2	1:A:235:LEU:HG	2.02	0.41
1:B:51:ASN:HB2	1:B:122:THR:HG21	2.02	0.41
1:A:140:LEU:HB2	1:A:149:ILE:HB	2.03	0.41
1:A:225:GLU:HG2	1:A:241:ASN:HB2	2.03	0.41
1:B:257:LYS:HG3	1:B:415:LEU:HB2	2.01	0.41
1:B:279:TYR:HA	1:B:369:ASP:O	2.20	0.41
1:A:77:ILE:CG1	1:B:77:ILE:HD11	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:259:ASN:O	1:B:263:LYS:HB2	2.21	0.41
1:B:404:SER:O	1:B:408:ASN:HB2	2.20	0.41
1:B:79:VAL:O	1:B:92:ARG:HA	2.20	0.41
1:B:264:LEU:HD12	1:B:367:ILE:HD11	2.03	0.40
1:A:49:GLY:HA3	1:A:95:PHE:CE1	2.56	0.40
1:B:260:PHE:CD2	1:B:415:LEU:HD11	2.56	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	306/393 (78%)	296 (97%)	8 (3%)	2 (1%)	26 51
1	B	283/393 (72%)	273 (96%)	10 (4%)	0	100 100
All	All	589/786 (75%)	569 (97%)	18 (3%)	2 (0%)	46 72

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	358	PRO
1	A	339	PRO

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	290/352 (82%)	272 (94%)	18 (6%)	23	45
1	B	277/352 (79%)	263 (95%)	14 (5%)	29	55
All	All	567/704 (80%)	535 (94%)	32 (6%)	26	50

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

Mol	Chain	Res	Type
1	A	58	SER
1	A	76	LYS
1	A	77	ILE
1	A	138	ARG
1	A	141	THR
1	A	160	MET
1	A	174	CYS
1	A	208	ARG
1	A	221	THR
1	A	222	VAL
1	A	251	HIS
1	A	265	LYS
1	A	274	LEU
1	A	289	ASP
1	A	399	ASN
1	A	401	GLU
1	A	414	ILE
1	A	415	LEU
1	B	34	LYS
1	B	86	LYS
1	B	104	ARG
1	B	123	ARG
1	B	208	ARG
1	B	225	GLU
1	B	230	GLU
1	B	231	LYS
1	B	235	LEU
1	B	240	ASP
1	B	258	LYS
1	B	289	ASP
1	B	370	ILE
1	B	408	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	354	HIS
1	B	241	ASN
1	B	354	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

4 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	$\#\|Z\| > 2$
2	GNP	A	501	-	29,34,34	1.76	4 (13%)	29,54,54	2.75	8 (27%)
2	GNP	A	502	-	29,34,34	1.93	5 (17%)	29,54,54	2.59	7 (24%)
2	GNP	A	503	-	22,26,34	1.10	2 (9%)	25,40,54	2.57	4 (16%)
2	GNP	B	501	-	29,34,34	1.77	6 (20%)	29,54,54	1.91	6 (20%)

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
2	GNP	A	501	-	-	0/13/38/38	0/3/3/3
2	GNP	A	502	-	-	0/13/38/38	0/3/3/3
2	GNP	A	503	-	-	0/6/26/38	0/3/3/3
2	GNP	B	501	-	-	1/13/38/38	0/3/3/3

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	502	GNP	PB-O3A	-4.84	1.53	1.59
2	A	501	GNP	PB-O3A	-4.50	1.53	1.59
2	A	501	GNP	PB-O2B	-3.20	1.48	1.56
2	A	502	GNP	PB-O2B	-3.07	1.48	1.56
2	A	503	GNP	C8-N7	-2.12	1.30	1.34
2	B	501	GNP	PG-O1G	2.15	1.48	1.46
2	B	501	GNP	PB-O1B	2.19	1.48	1.46
2	A	502	GNP	PB-O1B	2.95	1.49	1.46
2	B	501	GNP	C5-C4	3.09	1.47	1.40
2	A	501	GNP	C6-N1	3.57	1.39	1.33
2	B	501	GNP	C6-C5	3.61	1.48	1.41
2	A	503	GNP	C6-N1	3.75	1.39	1.33
2	A	502	GNP	C6-N1	4.07	1.40	1.33
2	B	501	GNP	PB-N3B	4.42	1.75	1.63
2	B	501	GNP	PG-N3B	4.44	1.75	1.63
2	A	501	GNP	PG-O1G	4.68	1.51	1.46
2	A	502	GNP	PG-O1G	5.46	1.52	1.46

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	501	GNP	C5-C6-N1	-9.56	111.02	123.52
2	A	503	GNP	C5-C6-N1	-9.16	111.55	123.52
2	A	502	GNP	C5-C6-N1	-8.56	112.33	123.52
2	B	501	GNP	C5-C6-N1	-4.23	117.99	123.52
2	B	501	GNP	N3-C2-N1	-3.47	122.84	127.56
2	B	501	GNP	PA-O3A-PB	-3.36	120.51	132.71
2	A	501	GNP	O3G-PG-O1G	-3.27	104.96	113.58
2	B	501	GNP	C6-C5-C4	-3.16	117.25	120.86
2	A	501	GNP	PA-O3A-PB	-2.73	122.82	132.71
2	A	502	GNP	N3-C2-N1	-2.72	123.86	127.56
2	A	501	GNP	N3-C2-N1	-2.65	123.95	127.56
2	A	502	GNP	O3G-PG-O1G	-2.29	107.56	113.58
2	A	503	GNP	N3-C2-N1	-2.05	124.76	127.56
2	A	502	GNP	O2A-PA-O1A	2.01	123.01	112.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	501	GNP	C4'-O4'-C1'	2.02	111.78	109.64
2	A	502	GNP	O3G-PG-O2G	2.41	114.67	107.67
2	A	501	GNP	O5'-C5'-C4'	2.45	117.94	109.09
2	A	503	GNP	O2A-PA-O1A	2.46	118.65	110.63
2	A	501	GNP	O3G-PG-O2G	2.52	115.00	107.67
2	A	501	GNP	O2B-PB-O1B	3.12	116.16	110.02
2	B	501	GNP	C6-N1-C2	5.50	122.32	115.88
2	A	502	GNP	O2B-PB-O1B	5.66	121.17	110.02
2	A	502	GNP	C6-N1-C2	7.22	124.34	115.88
2	A	503	GNP	C6-N1-C2	7.40	124.55	115.88
2	A	501	GNP	C6-N1-C2	8.38	125.70	115.88

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	501	GNP	O1G-PG-N3B-PB

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	502	GNP	1	0
2	B	501	GNP	2	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, 95^{th} 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	314/393 (79%)	0.62	4 (1%) 79 75	40, 73, 126, 142	0
1	B	297/393 (75%)	0.82	19 (6%) 23 17	41, 80, 152, 174	0
All	All	611/786 (77%)	0.72	23 (3%) 44 36	40, 75, 142, 174	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	403	TYR	4.7
1	B	355	GLU	4.0
1	B	238	PHE	3.8
1	B	413	ASN	3.8
1	B	265	LYS	3.7
1	B	215	TYR	3.1
1	B	356	SER	3.1
1	A	402	GLN	2.8
1	B	259	ASN	2.7
1	B	258	LYS	2.7
1	B	251	HIS	2.6
1	A	272	ALA	2.6
1	B	249	LYS	2.5
1	B	240	ASP	2.4
1	B	245	ASN	2.4
1	B	73	ALA	2.4
1	B	362	VAL	2.3
1	B	243	PHE	2.2
1	B	403	TYR	2.1
1	A	277	MET	2.1
1	B	218	LYS	2.0
1	B	230	GLU	2.0
1	B	360	LYS	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, 95^{th} 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(Å^2)	Q<0.9
2	GNP	A	503	24/32	0.87	0.24	-0.45	85,96,99,104	0
2	GNP	A	502	32/32	0.90	0.20	-0.58	53,72,122,286	0
2	GNP	B	501	32/32	0.73	0.22	-0.61	89,121,197,333	0
2	GNP	A	501	32/32	0.87	0.21	-0.63	51,100,223,319	0

6.5 Other polymers (i)

There are no such residues in this entry.