



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 2, 2016 – 10:47 AM EDT

PDB ID : 5JCF
Title : Crystal structure of chicken MDA5 with 5'p 10-mer dsRNA and ADP-Mg2+ at 2.6 Å resolution (orthorhombic form).
Authors : Cusack, S.; Uchikawa, E.
Deposited on : 2016-04-15
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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027674
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-20027674

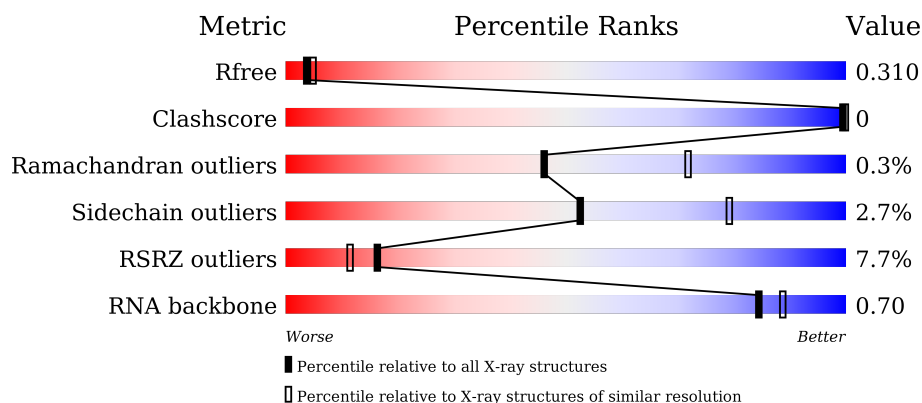
1 Overall quality at a glance

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)
RNA backbone	2183	1022 (3.00-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	701	<div> <div>8%</div> <div>91%</div> <div>6%</div> </div>
1	B	701	<div> <div>7%</div> <div>91%</div> <div>6%</div> </div>
2	C	10	<div> <div>90%</div> <div>10%</div> </div>
2	D	10	<div> <div>90%</div> <div>10%</div> </div>

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Mol	Chain	Length	Quality of chain
2	X	10	 90%10%
3	Y	11	 9%100%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 11870 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 Melanoma differentiation associated protein-5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	660	Total	C	N	O	S	0	0	0
			5386	3394	964	996	32			
1	B	660	Total	C	N	O	S	0	0	0
			5385	3393	964	996	32			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	294	GLY	-	expression tag	UNP D9N195
A	295	ALA	-	expression tag	UNP D9N195
A	296	MET	-	expression tag	UNP D9N195
A	297	GLY	-	expression tag	UNP D9N195
A	436	GLN	GLU	engineered mutation	UNP D9N195
B	294	GLY	-	expression tag	UNP D9N195
B	295	ALA	-	expression tag	UNP D9N195
B	296	MET	-	expression tag	UNP D9N195
B	297	GLY	-	expression tag	UNP D9N195
B	436	GLN	GLU	engineered mutation	UNP D9N195

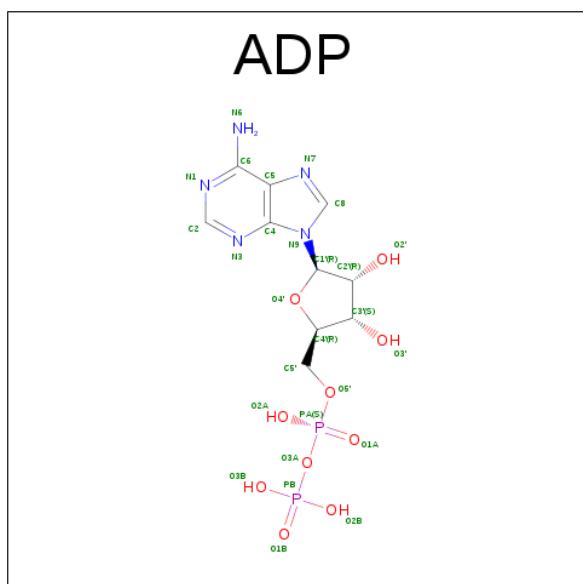
- Molecule 2 is a RNA chain called RNA (5'-R(P*GP*GP*UP*AP*CP*GP*UP*AP*CP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	X	10	Total	C	N	O	P	0	0	0
			214	95	38	71	10			
2	C	10	Total	C	N	O	P	0	0	0
			214	95	38	71	10			
2	D	10	Total	C	N	O	P	0	0	0
			214	95	38	71	10			

- Molecule 3 is a RNA chain called RNA (5'-R(P*AP*GP*GP*UP*AP*CP*GP*UP*AP*CP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	Y	11	Total	C	N	O	P	0	0	0
			231	105	43	73	10			

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $\text{C}_{10}\text{H}_{15}\text{N}_5\text{O}_{10}\text{P}_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 27	C 10	N 5	O 10	P 2	0	0
4	B	1	Total 27	C 10	N 5	O 10	P 2	0	0

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Mg 1 1	0	0
5	A	1	Total Mg 1 1	0	0

- Molecule 6 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total Zn 1 1	0	0
6	A	1	Total Zn 1 1	0	0

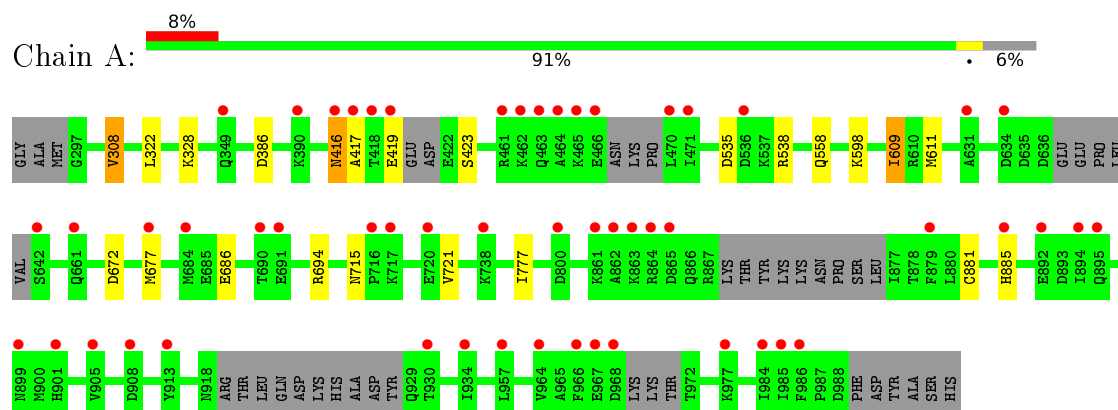
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	74	Total 74	O 74	0	0
7	X	5	Total 5	O 5	0	0
7	Y	7	Total 7	O 7	0	0
7	B	70	Total 70	O 70	0	0
7	C	5	Total 5	O 5	0	0
7	D	7	Total 7	O 7	0	0

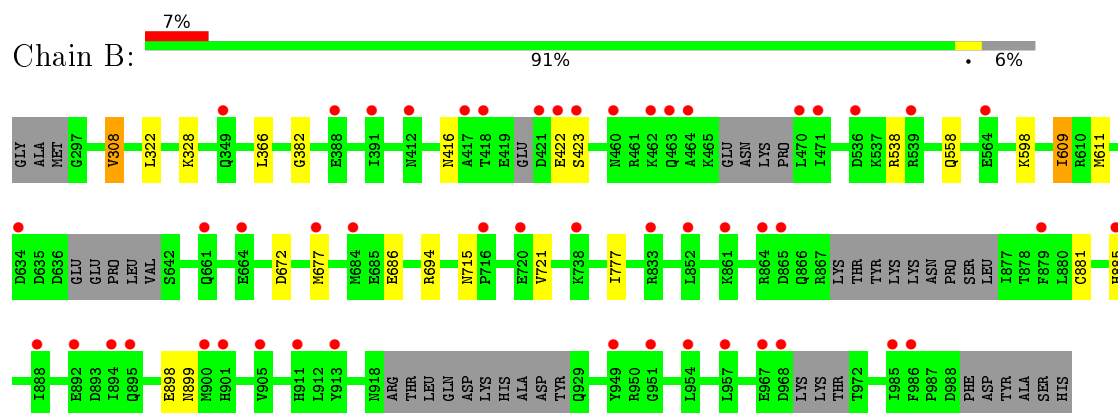
3 Residue-property plots

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 ($RSRZ > 2$). 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.

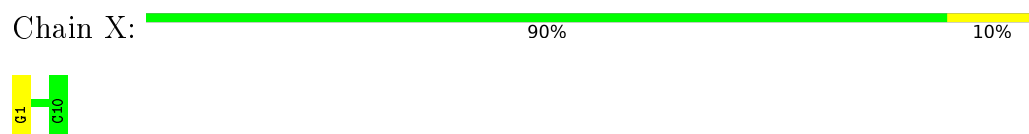
- Molecule 1: Melanoma differentiation associated protein-5



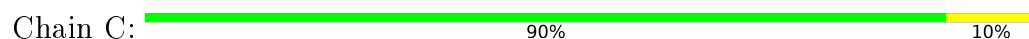
- Molecule 1: Melanoma differentiation associated protein-5



- Molecule 2: RNA (5'-R(P*GP*GP*UP*AP*CP*GP*UP*AP*CP*C)-3')




- Molecule 2: RNA (5'-R(P*GP*GP*UP*AP*CP*GP*UP*AP*CP*C)-3')





- Molecule 2: RNA (5'-R(P*GP*GP*UP*AP*CP*GP*UP*AP*CP*C)-3')

Chain D:  90% 10%



- Molecule 3: RNA (5'-R(P*AP*GP*GP*UP*AP*CP*GP*UP*AP*CP*C)-3')

Chain Y:  9% 100%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.92Å 132.47Å 139.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.60 47.95 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.2 (50.00-2.60) 98.2 (47.95-2.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.277 , 0.312 0.279 , 0.310	Depositor DCC
R_{free} test set	2901 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	51.3	Xtriage
Anisotropy	0.783	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 32.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11870	wwPDB-VP
Average B, all atoms (Å ²)	66.0	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 60.90 % 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 1.4199e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: ZN, MG, ADP

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	0.37	0/5471	0.58	0/7338
1	B	0.36	0/5470	0.58	0/7337
2	C	0.68	1/238 (0.4%)	0.69	0/367
2	D	0.70	1/238 (0.4%)	0.73	0/367
2	X	0.68	1/238 (0.4%)	0.69	0/367
3	Y	0.22	0/258	0.69	0/401
All	All	0.39	3/11913 (0.0%)	0.59	0/16177

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	G	OP3-P	-10.18	1.49	1.61
2	C	1	G	OP3-P	-9.94	1.49	1.61
2	X	1	G	OP3-P	-9.86	1.49	1.61

There are no bond angle outliers.

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	5386	0	5463	4	0
1	B	5385	0	5461	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	214	0	109	0	0
2	D	214	0	109	0	0
2	X	214	0	109	0	0
3	Y	231	0	118	0	0
4	A	27	0	12	0	0
4	B	27	0	12	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	A	74	0	0	0	0
7	B	70	0	0	0	0
7	C	5	0	0	0	0
7	D	7	0	0	0	0
7	X	5	0	0	0	0
7	Y	7	0	0	0	0
All	All	11870	0	11393	8	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 0.

All (8) 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:538:ARG:NH1	1:A:672:ASP:O	2.30	0.64
1:B:538:ARG:NH1	1:B:672:ASP:O	2.30	0.64
1:B:308:VAL:HG21	1:B:322:LEU:HD21	1.99	0.45
1:A:308:VAL:HG21	1:A:322:LEU:HD21	1.99	0.45
1:B:686:GLU:HG2	1:B:777:ILE:HB	2.00	0.44
1:A:686:GLU:HG2	1:A:777:ILE:HB	2.00	0.42
1:A:416:ASN:O	1:A:417:ALA:HB3	2.20	0.42
1:B:366:LEU:HD11	1:B:382:GLY:HA3	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	646/701 (92%)	619 (96%)	25 (4%)	2 (0%)	46	72
1	B	646/701 (92%)	619 (96%)	25 (4%)	2 (0%)	46	72
All	All	1292/1402 (92%)	1238 (96%)	50 (4%)	4 (0%)	46	72

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	416	ASN
1	B	416	ASN
1	A	609	ILE
1	B	609	ILE

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	593/630 (94%)	577 (97%)	16 (3%)	52	79
1	B	593/630 (94%)	577 (97%)	16 (3%)	52	79
All	All	1186/1260 (94%)	1154 (97%)	32 (3%)	52	79

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

Mol	Chain	Res	Type
1	A	308	VAL
1	A	328	LYS

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Mol	Chain	Res	Type
1	A	386	ASP
1	A	419	GLU
1	A	423	SER
1	A	535	ASP
1	A	558	GLN
1	A	598	LYS
1	A	609	ILE
1	A	611	MET
1	A	677	MET
1	A	694	ARG
1	A	715	ASN
1	A	721	VAL
1	A	881	CYS
1	A	885	HIS
1	B	308	VAL
1	B	328	LYS
1	B	422	GLU
1	B	423	SER
1	B	558	GLN
1	B	598	LYS
1	B	609	ILE
1	B	611	MET
1	B	677	MET
1	B	694	ARG
1	B	715	ASN
1	B	721	VAL
1	B	881	CYS
1	B	885	HIS
1	B	898	GLU
1	B	899	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	9/10 (90%)	0	0
2	D	9/10 (90%)	0	0
2	X	9/10 (90%)	0	0
3	Y	10/11 (90%)	0	0

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
All	All	37/41 (90%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

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](#)

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	$\# Z > 2$
4	ADP	A	1001	5	24,29,29	1.07	2 (8%)	23,45,45	1.95	2 (8%)
4	ADP	B	1001	5	24,29,29	1.05	1 (4%)	23,45,45	2.00	3 (13%)

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	ADP	A	1001	5	-	0/12/32/32	0/3/3/3
4	ADP	B	1001	5	-	0/12/32/32	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1001	ADP	C2-N3	2.02	1.35	1.32
4	B	1001	ADP	C5-C4	3.24	1.47	1.40
4	A	1001	ADP	C5-C4	3.26	1.47	1.40

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1001	ADP	N3-C2-N1	-7.85	122.70	128.87
4	A	1001	ADP	N3-C2-N1	-7.70	122.82	128.87
4	B	1001	ADP	C2-N1-C6	2.02	122.38	118.77
4	A	1001	ADP	C4'-O4'-C1'	2.22	112.00	109.64
4	B	1001	ADP	C4'-O4'-C1'	2.39	112.18	109.64

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	660/701 (94%)	0.48	54 (8%) 14 10	41, 63, 102, 142	0
1	B	660/701 (94%)	0.42	50 (7%) 17 12	41, 63, 103, 128	0
2	C	10/10 (100%)	-0.06	0 100 100	41, 51, 70, 75	0
2	D	10/10 (100%)	-0.13	0 100 100	42, 46, 71, 84	0
2	X	10/10 (100%)	-0.20	0 100 100	43, 51, 67, 74	0
3	Y	11/11 (100%)	0.40	1 (9%) 11 7	41, 46, 86, 106	0
All	All	1361/1443 (94%)	0.44	105 (7%) 16 11	41, 63, 102, 142	0

All (105) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	968	ASP	6.1
1	A	417	ALA	5.8
1	B	716	PRO	5.8
1	B	905	VAL	5.1
3	Y	0	A	5.1
1	A	463	GLN	5.1
1	B	470	LEU	5.0
1	A	716	PRO	4.9
1	B	463	GLN	4.9
1	A	862	ALA	4.8
1	A	466	GLU	4.7
1	A	905	VAL	4.5
1	B	967	GLU	4.4
1	B	418	THR	4.4
1	B	349	GLN	4.4
1	B	892	GLU	4.4
1	A	892	GLU	4.3
1	A	677	MET	4.1
1	A	631	ALA	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	894	ILE	4.1
1	B	986	PHE	4.1
1	B	423	SER	4.0
1	B	894	ILE	4.0
1	B	536	ASP	3.9
1	A	464	ALA	3.9
1	B	634	ASP	3.9
1	A	968	ASP	3.8
1	A	416	ASN	3.7
1	A	899	ASN	3.7
1	A	536	ASP	3.7
1	A	967	GLU	3.6
1	B	985	ILE	3.6
1	A	985	ILE	3.6
1	A	462	LYS	3.4
1	A	986	PHE	3.4
1	B	417	ALA	3.4
1	A	895	GLN	3.4
1	A	690	THR	3.4
1	B	462	LYS	3.4
1	B	421	ASP	3.4
1	A	465	LYS	3.4
1	B	539	ARG	3.3
1	A	418	THR	3.3
1	B	895	GLN	3.2
1	B	677	MET	3.1
1	A	930	THR	3.0
1	A	349	GLN	3.0
1	A	885	HIS	3.0
1	B	471	ILE	2.9
1	A	691	GLU	2.9
1	A	800	ASP	2.9
1	B	885	HIS	2.8
1	A	957	LEU	2.8
1	A	634	ASP	2.8
1	B	879	PHE	2.7
1	A	863	LYS	2.7
1	A	661	GLN	2.7
1	B	684	MET	2.7
1	A	964	VAL	2.6
1	B	422	GLU	2.6
1	A	738	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	391	ILE	2.6
1	B	951	GLY	2.6
1	A	390	LYS	2.5
1	A	966	PHE	2.5
1	A	419	GLU	2.5
1	A	934	ILE	2.5
1	A	865	ASP	2.4
1	A	879	PHE	2.4
1	B	738	LYS	2.4
1	B	661	GLN	2.4
1	A	861	LYS	2.4
1	A	864	ARG	2.4
1	A	977	LYS	2.3
1	B	564	GLU	2.3
1	B	388	GLU	2.3
1	B	852	LEU	2.3
1	A	471	ILE	2.3
1	A	984	ILE	2.3
1	B	720	GLU	2.3
1	A	461	ARG	2.3
1	B	464	ALA	2.3
1	A	470	LEU	2.3
1	A	642	SER	2.3
1	B	957	LEU	2.3
1	B	949	TYR	2.3
1	A	684	MET	2.2
1	B	460	ASN	2.2
1	B	664	GLU	2.2
1	A	717	LYS	2.2
1	B	861	LYS	2.2
1	B	954	LEU	2.2
1	B	911	HIS	2.2
1	B	913	TYR	2.1
1	B	901	HIS	2.1
1	A	913	TYR	2.1
1	A	901	HIS	2.1
1	B	888	ILE	2.1
1	B	833	ARG	2.1
1	B	865	ASP	2.1
1	A	720	GLU	2.1
1	A	908	ASP	2.1
1	B	864	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	900	MET	2.0
1	B	412	ASN	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(Å ²)	Q<0.9
4	ADP	B	1001	27/27	0.95	0.16	-1.03	58,62,65,66	0
4	ADP	A	1001	27/27	0.94	0.14	-1.26	64,68,71,71	0
6	ZN	A	1003	1/1	0.98	0.09	-1.44	63,63,63,63	0
6	ZN	B	1003	1/1	0.99	0.11	-2.06	65,65,65,65	0
5	MG	A	1002	1/1	0.96	0.21	-	42,42,42,42	0
5	MG	B	1002	1/1	0.91	0.25	-	41,41,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.