



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 6, 2017 – 09:01 PM EDT

PDB ID : 5TZD
Title : Structure of the WT *S. venezulae* BldD-(CTD-c-di-GMP)₂ assembly intermediate
Authors : Schumacher, M.
Deposited on : unknown
Resolution : 1.75 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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-20029824

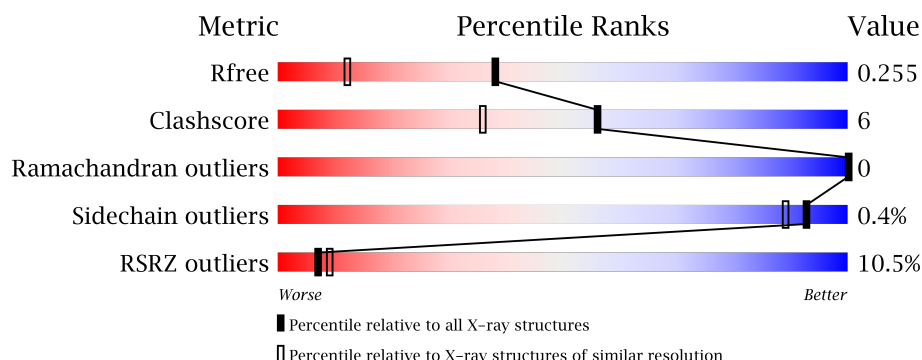
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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}	100719	2694 (1.76-1.72)
Clashscore	112137	2854 (1.76-1.72)
Ramachandran outliers	110173	2824 (1.76-1.72)
Sidechain outliers	110143	2824 (1.76-1.72)
RSRZ outliers	101464	2705 (1.76-1.72)

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	91	<div> <div>14%</div> <div> <div></div> <div>64%</div> <div>13%</div> <div>•</div> <div>22%</div> </div> </div>
1	B	91	<div> <div>14%</div> <div> <div></div> <div>74%</div> <div>9%</div> <div>18%</div> </div> </div>
1	C	91	<div> <div>4%</div> <div> <div></div> <div>75%</div> <div>8%</div> <div>18%</div> </div> </div>
1	J	91	<div> <div>•%</div> <div> <div></div> <div>76%</div> <div>7%</div> <div>18%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5288 atoms, of which 2460 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.

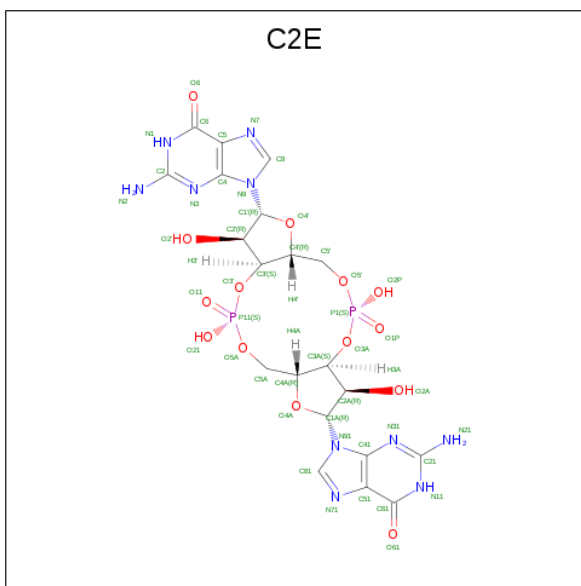
- Molecule 1 is a protein called DNA-binding protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	J	75	Total	C	H	N	O	S	0	0	0
			1192	370	600	107	114	1			
1	B	75	Total	C	H	N	O	S	0	0	0
			1192	370	600	107	114	1			
1	C	75	Total	C	H	N	O	S	0	0	0
			1180	364	595	107	113	1			
1	A	71	Total	C	H	N	O	S	0	0	0
			1136	354	573	100	108	1			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
J	76	GLY	-	expression tag	UNP F2RCL8
J	77	SER	-	expression tag	UNP F2RCL8
J	78	HIS	-	expression tag	UNP F2RCL8
J	79	MET	-	expression tag	UNP F2RCL8
J	92	MET	LEU	engineered mutation	UNP F2RCL8
B	76	GLY	-	expression tag	UNP F2RCL8
B	77	SER	-	expression tag	UNP F2RCL8
B	78	HIS	-	expression tag	UNP F2RCL8
B	79	MET	-	expression tag	UNP F2RCL8
B	92	MET	LEU	engineered mutation	UNP F2RCL8
C	76	GLY	-	expression tag	UNP F2RCL8
C	77	SER	-	expression tag	UNP F2RCL8
C	78	HIS	-	expression tag	UNP F2RCL8
C	79	MET	-	expression tag	UNP F2RCL8
C	92	MET	LEU	engineered mutation	UNP F2RCL8
A	76	GLY	-	expression tag	UNP F2RCL8
A	77	SER	-	expression tag	UNP F2RCL8
A	78	HIS	-	expression tag	UNP F2RCL8
A	79	MET	-	expression tag	UNP F2RCL8
A	92	MET	LEU	engineered mutation	UNP F2RCL8

- Molecule 2 is 9,9'-[(2R,3R,3aS,5S,7aR,9R,10R,10aS,12S,14aR)-3,5,10,12-tetrahydroxy-5,12-dioxidoctahydro-2H,7H-difuro[3,2-d:3',2'-j][1,3,7,9,2,8]tetraoxadiphosphacyclododecine-2,9-diyl]bis(2-amino-1,9-dihydro-6H-purin-6-one) (three-letter code: C2E) (formula: $C_{20}H_{24}N_{10}O_{14}P_2$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	J	1	Total 69	C 20	H 23	N 10	O 14	P 2	0	0
2	J	1	Total 69	C 20	H 23	N 10	O 14	P 2	0	0
2	C	1	Total 69	C 20	H 23	N 10	O 14	P 2	0	0
2	C	1	Total 69	C 20	H 23	N 10	O 14	P 2	0	0

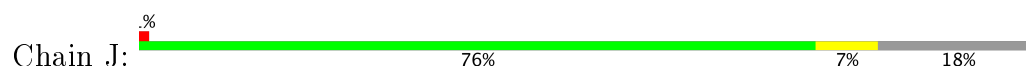
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	J	130	Total	O	0	0
			130	130		
3	B	52	Total	O	0	0
			52	52		
3	C	75	Total	O	0	0
			75	75		
3	A	55	Total	O	0	0
			55	55		

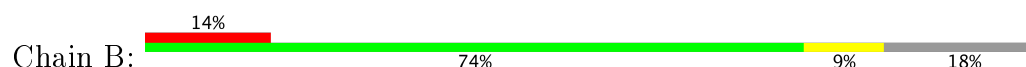
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

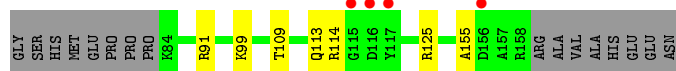
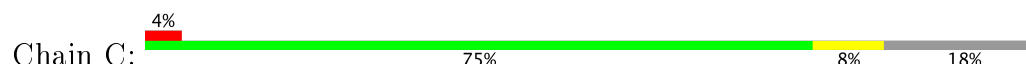
- Molecule 1: DNA-binding protein



- Molecule 1: DNA-binding protein



- Molecule 1: DNA-binding protein



- Molecule 1: DNA-binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	108.30 Å 108.30 Å 49.66 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.89 – 1.75 43.89 – 1.75	Depositor EDS
% Data completeness (in resolution range)	94.2 (43.89-1.75) 94.2 (43.89-1.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.20 (at 1.75 Å)	Xtriage
Refinement program	PHENIX 1.6.4_486	Depositor
R, R_{free}	0.207 , 0.240 0.220 , 0.255	Depositor DCC
R_{free} test set	1994 reflections (6.70%)	DCC
Wilson B-factor (Å ²)	18.8	Xtriage
Anisotropy	0.016	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 51.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.038 for h,-h-k,-l	Xtriage
Reported twinning fraction	0.038 for h,-h-k,-l	Depositor
Outliers	0 of 31775 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5288	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.42% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: C2E

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	0.33	0/571	0.49	0/774
1	B	0.37	0/600	0.53	0/813
1	C	0.53	0/592	0.66	0/802
1	J	0.50	0/600	0.67	0/813
All	All	0.44	0/2363	0.59	0/3202

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 [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	563	573	572	10	1
1	B	592	600	599	7	1
1	C	585	595	592	6	0
1	J	592	600	599	6	0
2	C	92	46	43	3	0
2	J	92	46	42	3	0
3	A	55	0	0	1	1
3	B	52	0	0	0	1
3	C	75	0	0	3	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	J	130	0	0	3	1
All	All	2828	2460	2447	29	3

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 6.

All (29) 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:C:99:LYS:NZ	3:C:301:HOH:O	2.03	0.90
1:B:85:LEU:HD23	1:B:85:LEU:N	2.08	0.68
1:C:114:ARG:NH2	2:C:201:C2E:O61	2.27	0.67
1:J:99:LYS:NZ	3:J:302:HOH:O	2.26	0.64
1:A:146:GLN:HG3	1:A:150:TRP:CZ2	2.36	0.60
2:C:201:C2E:H8	1:A:125:ARG:HG2	1.83	0.60
1:C:125:ARG:HG2	2:C:202:C2E:H8	1.85	0.57
1:C:155:ALA:HB3	3:C:357:HOH:O	2.05	0.56
1:J:126:GLN:NE2	3:J:306:HOH:O	2.40	0.53
1:B:85:LEU:HD13	1:B:129:LEU:HD13	1.92	0.52
1:J:114:ARG:NH2	2:J:202:C2E:O61	2.42	0.52
1:C:109:THR:O	1:C:113:GLN:HG3	2.10	0.51
1:J:99:LYS:NZ	3:J:301:HOH:O	2.17	0.51
1:A:105:ARG:O	1:A:109:THR:HG23	2.11	0.49
1:A:146:GLN:CG	1:A:150:TRP:CZ2	2.96	0.49
1:A:146:GLN:HG3	1:A:150:TRP:CE2	2.48	0.48
1:B:132:LEU:HB3	1:B:143:LEU:HD21	1.94	0.48
1:A:103:LEU:HD22	1:A:147:LEU:CD2	2.45	0.47
1:A:127:ASP:HB2	3:A:226:HOH:O	2.15	0.46
1:A:91:ARG:O	1:A:91:ARG:HD3	2.17	0.45
2:J:202:C2E:O6	1:B:125:ARG:NH2	2.44	0.45
1:B:85:LEU:CD1	1:B:129:LEU:HD13	2.48	0.44
1:C:91:ARG:NH2	3:C:302:HOH:O	2.06	0.44
1:B:95:VAL:CG1	1:B:100:ALA:HB2	2.49	0.42
1:B:95:VAL:HG11	1:B:100:ALA:HB2	2.03	0.41
1:A:95:VAL:HG13	1:A:96:PRO:HD2	2.03	0.41
1:J:95:VAL:HG13	1:J:96:PRO:HD2	2.03	0.41
1:J:125:ARG:HG2	2:J:201:C2E:H8	2.03	0.41
1:A:95:VAL:HG11	1:A:100:ALA:HB2	2.02	0.41

All (3) 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)
3:B:251:HOH:O	3:A:250:HOH:O[3_454]	2.08	0.12
1:B:119:GLY:O	1:A:117:TYR:HH[3_454]	1.53	0.07
3:J:392:HOH:O	3:C:349:HOH:O[6_555]	2.14	0.06

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	69/91 (76%)	67 (97%)	2 (3%)	0	100	100
1	B	73/91 (80%)	73 (100%)	0	0	100	100
1	C	73/91 (80%)	72 (99%)	1 (1%)	0	100	100
1	J	73/91 (80%)	72 (99%)	1 (1%)	0	100	100
All	All	288/364 (79%)	284 (99%)	4 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	62/77 (80%)	61 (98%)	1 (2%)	68	49
1	B	64/77 (83%)	64 (100%)	0	100	100
1	C	63/77 (82%)	63 (100%)	0	100	100
1	J	64/77 (83%)	64 (100%)	0	100	100
All	All	253/308 (82%)	252 (100%)	1 (0%)	93	89

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	91	ARG

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

Mol	Chain	Res	Type
1	B	113	GLN
1	C	118	ASN
1	C	126	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no 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	C2E	C	201	-	44,52,52	2.95	12 (27%)	44,82,82	2.32	13 (29%)
2	C2E	C	202	-	44,52,52	2.78	11 (25%)	44,82,82	2.18	9 (20%)
2	C2E	J	201	-	44,52,52	2.78	12 (27%)	44,82,82	2.18	9 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	C2E	J	202	-	44,52,52	2.95	13 (29%)	44,82,82	2.31	13 (29%)

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	C2E	C	201	-	-	0/22/62/62	0/6/7/7
2	C2E	C	202	-	-	0/22/62/62	0/6/7/7
2	C2E	J	201	-	-	0/22/62/62	0/6/7/7
2	C2E	J	202	-	-	0/22/62/62	0/6/7/7

All (48) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	J	201	C2E	C2'-C1'	-6.57	1.43	1.53
2	C	202	C2E	C2'-C1'	-6.57	1.43	1.53
2	J	201	C2E	C8-N7	-5.82	1.23	1.34
2	C	202	C2E	C8-N7	-5.81	1.23	1.34
2	C	201	C2E	C2'-C1'	-5.71	1.44	1.53
2	J	202	C2E	C2'-C1'	-5.67	1.44	1.53
2	J	202	C2E	C81-N71	-5.60	1.24	1.34
2	C	201	C2E	C81-N71	-5.59	1.24	1.34
2	C	201	C2E	C8-N7	-5.37	1.24	1.34
2	J	202	C2E	C8-N7	-5.36	1.24	1.34
2	C	202	C2E	C2'-C3'	-4.79	1.42	1.53
2	J	201	C2E	C2'-C3'	-4.77	1.42	1.53
2	J	201	C2E	C81-N71	-4.32	1.26	1.34
2	C	201	C2E	C2'-C3'	-4.29	1.43	1.53
2	C	202	C2E	C81-N71	-4.29	1.26	1.34
2	J	202	C2E	C2'-C3'	-4.28	1.43	1.53
2	J	202	C2E	O4'-C4'	-3.24	1.37	1.45
2	C	201	C2E	O4'-C4'	-3.20	1.37	1.45
2	J	202	C2E	O2A-C2A	-2.62	1.36	1.43
2	C	201	C2E	O2A-C2A	-2.58	1.37	1.43
2	J	201	C2E	O2A-C2A	-2.37	1.37	1.43
2	C	202	C2E	O2A-C2A	-2.35	1.37	1.43
2	C	201	C2E	C61-N11	-2.28	1.33	1.36
2	J	202	C2E	C61-N11	-2.26	1.33	1.36
2	J	201	C2E	O5'-C5'	-2.08	1.36	1.44
2	C	202	C2E	O5'-C5'	-2.06	1.36	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	J	201	C2E	C21-N11	-2.01	1.33	1.36
2	J	202	C2E	C2-N3	2.05	1.36	1.33
2	J	202	C2E	C5-C4	2.52	1.46	1.40
2	C	201	C2E	C5-C4	2.57	1.46	1.40
2	J	201	C2E	C5-C4	2.60	1.46	1.40
2	C	202	C2E	C5-C4	2.61	1.46	1.40
2	J	202	C2E	O61-C61	2.71	1.31	1.24
2	C	201	C2E	O61-C61	2.73	1.31	1.24
2	C	202	C2E	O61-C61	2.93	1.31	1.24
2	J	201	C2E	O61-C61	2.97	1.31	1.24
2	C	202	C2E	O6-C6	3.55	1.33	1.24
2	J	201	C2E	O6-C6	3.55	1.33	1.24
2	J	202	C2E	O6-C6	4.60	1.36	1.24
2	C	201	C2E	O6-C6	4.65	1.36	1.24
2	J	201	C2E	C21-N21	7.15	1.42	1.32
2	C	202	C2E	C21-N21	7.16	1.42	1.32
2	C	201	C2E	C21-N21	8.51	1.44	1.32
2	J	202	C2E	C21-N21	8.57	1.44	1.32
2	C	202	C2E	C2-N2	9.98	1.46	1.32
2	J	201	C2E	C2-N2	10.02	1.46	1.32
2	C	201	C2E	C2-N2	10.03	1.46	1.32
2	J	202	C2E	C2-N2	10.07	1.46	1.32

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	202	C2E	C41-C51-N71	-6.57	103.06	109.41
2	C	201	C2E	C41-C51-N71	-6.57	103.06	109.41
2	J	201	C2E	C41-C51-N71	-6.46	103.17	109.41
2	C	202	C2E	C41-C51-N71	-6.45	103.17	109.41
2	J	202	C2E	C4A-O4A-C1A	-6.27	103.09	109.77
2	C	201	C2E	C4A-O4A-C1A	-6.24	103.12	109.77
2	C	202	C2E	C2'-C3'-C4'	-5.76	92.92	103.23
2	J	201	C2E	C2'-C3'-C4'	-5.76	92.93	103.23
2	C	202	C2E	N21-C21-N31	-5.14	113.95	120.26
2	J	201	C2E	N21-C21-N31	-5.14	113.95	120.26
2	C	201	C2E	C4'-O4'-C1'	-4.74	104.72	109.77
2	C	201	C2E	C4-C5-N7	-4.68	104.89	109.41
2	J	202	C2E	C4'-O4'-C1'	-4.65	104.82	109.77
2	J	202	C2E	C4-C5-N7	-4.63	104.94	109.41
2	C	202	C2E	C1'-N9-C4	-3.48	120.63	126.64
2	J	201	C2E	C1'-N9-C4	-3.44	120.69	126.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	202	C2E	C4-C5-N7	-3.18	106.33	109.41
2	J	201	C2E	C4-C5-N7	-3.15	106.37	109.41
2	J	202	C2E	C1'-N9-C4	-2.73	121.92	126.64
2	C	201	C2E	C1'-N9-C4	-2.71	121.95	126.64
2	C	201	C2E	C1A-N91-C41	-2.60	122.15	126.64
2	C	201	C2E	C2A-C3A-C4A	-2.54	98.68	103.23
2	J	202	C2E	C1A-N91-C41	-2.53	122.27	126.64
2	J	202	C2E	C2A-C3A-C4A	-2.49	98.77	103.23
2	C	201	C2E	N21-C21-N31	-2.17	117.59	120.26
2	J	202	C2E	N21-C21-N31	-2.15	117.62	120.26
2	J	202	C2E	C3'-C2'-C1'	2.01	104.46	99.95
2	C	201	C2E	N21-C21-N11	2.01	119.98	117.84
2	C	202	C2E	N11-C21-N31	2.31	124.72	121.79
2	J	201	C2E	N11-C21-N31	2.32	124.74	121.79
2	J	202	C2E	C5'-C4'-C3'	2.32	122.18	114.30
2	C	201	C2E	C5'-C4'-C3'	2.33	122.19	114.30
2	C	201	C2E	C5A-C4A-C3A	2.48	122.70	114.30
2	J	202	C2E	C5A-C4A-C3A	2.51	122.82	114.30
2	J	202	C2E	O4'-C4'-C3'	2.92	111.25	104.81
2	C	201	C2E	O4'-C4'-C3'	2.92	111.25	104.81
2	J	201	C2E	N21-C21-N11	3.24	121.30	117.84
2	C	202	C2E	N21-C21-N11	3.27	121.32	117.84
2	C	201	C2E	O4A-C4A-C3A	4.45	114.64	104.81
2	J	202	C2E	O4A-C4A-C3A	4.45	114.64	104.81
2	J	201	C2E	O2'-C2'-C3'	4.54	124.08	111.18
2	C	202	C2E	O2'-C2'-C3'	4.54	124.09	111.18
2	J	201	C2E	C3'-C2'-C1'	4.57	110.22	99.95
2	C	202	C2E	C3'-C2'-C1'	4.59	110.26	99.95

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	201	C2E	2	0
2	C	202	C2E	1	0
2	J	201	C2E	1	0
2	J	202	C2E	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, 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	71/91 (78%)	1.11	13 (18%)  	19, 34, 65, 84	0
1	B	75/91 (82%)	0.99	13 (17%)  	16, 33, 64, 100	0
1	C	75/91 (82%)	0.56	4 (5%)  	10, 21, 54, 76	0
1	J	75/91 (82%)	0.49	1 (1%)  	8, 15, 38, 66	0
All	All	296/364 (81%)	0.78	31 (10%)  	8, 27, 62, 100	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	156	ASP	4.5
1	C	115	GLY	3.9
1	A	148	ILE	3.9
1	B	157	ALA	3.4
1	A	154	ASP	3.3
1	B	116	ASP	3.3
1	A	153	LEU	3.2
1	A	137	ASP	3.2
1	A	99	LYS	3.1
1	B	154	ASP	3.1
1	B	84	LYS	2.8
1	A	98	GLU	2.8
1	A	100	ALA	2.6
1	B	158	ARG	2.6
1	A	150	TRP	2.6
1	J	97	GLN	2.5
1	A	147	LEU	2.5
1	B	121	VAL	2.4
1	B	97	GLN	2.4
1	B	150	TRP	2.3
1	A	142	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	95	VAL	2.2
1	A	121	VAL	2.2
1	B	155	ALA	2.1
1	C	156	ASP	2.1
1	B	85	LEU	2.1
1	A	87	LEU	2.1
1	B	120	LYS	2.0
1	C	117	TYR	2.0
1	C	116	ASP	2.0
1	A	143	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(Å ²)	Q<0.9
2	C2E	J	201	46/46	0.94	0.13	0.19	13,20,29,32	0
2	C2E	J	202	46/46	0.92	0.15	0.17	11,20,27,30	0
2	C2E	C	202	46/46	0.93	0.13	0.04	15,23,31,32	0
2	C2E	C	201	46/46	0.92	0.13	-0.37	16,22,28,35	0

6.5 Other polymers [i](#)

There are no such residues in this entry.