Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 10:58 PM GMT

PDB ID	:	1W2Y
Title	:	THE CRYSTAL STRUCTURE OF A COMPLEX OF CAMPYLOBACTER JEJUNI DUTPASE WITH SUBSTRATE ANALOGUE DUPNHP
Authors	:	Moroz, O.V.; Harkiolaki, M.; Galperin, M.Y.; Vagin, A.A.; Gonzalez-Pacanowska, D.; Wilson, K.S.
Deposited on	:	2004-07-09
Resolution	:	1.65 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:

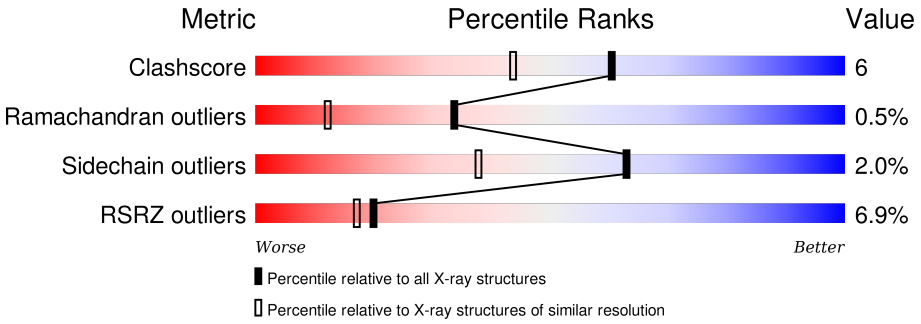
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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))
Clashscore	102246	1323 (1.66-1.66)
Ramachandran outliers	100387	1295 (1.66-1.66)
Sidechain outliers	100360	1295 (1.66-1.66)
RSRZ outliers	91569	1227 (1.66-1.66)

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	229	5% 82% 12%
1	B	229	8% 80% 10% 9%

2 Entry composition

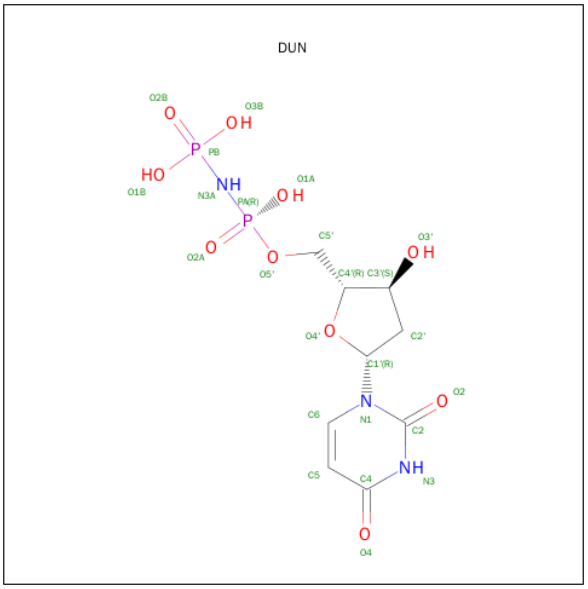
There are 4 unique types of molecules in this entry. The entry contains 4067 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDE HYDROLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	226	Total	C	N	O	S	83	4	0
			1895	1220	306	360	9			
1	B	209	Total	C	N	O	S	59	2	0
			1735	1121	285	320	9			

- Molecule 2 is 2'-DEOXYURIDINE 5'-ALPHA,BETA-IMIDO-DIPHOSPHATE (three-letter code: DUN) (formula: $C_9H_{15}N_3O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			24	9	3	10	2		
2	B	1	Total	C	N	O	P	0	0
			24	9	3	10	2		

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

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

- Molecule 4 is water.

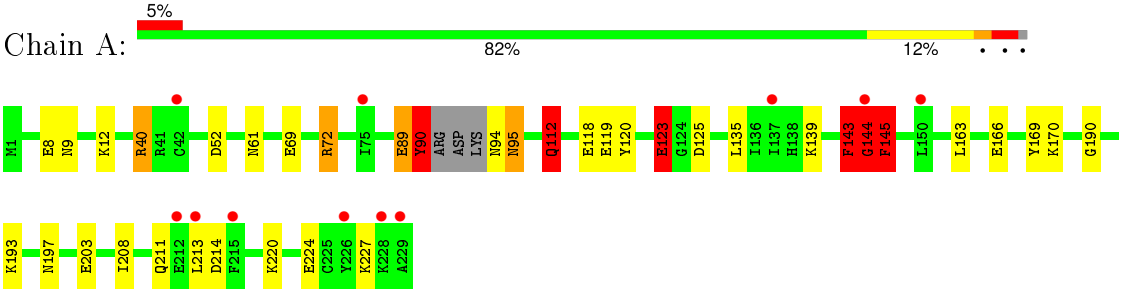
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	195	Total O 195 195	0	0
4	B	188	Total O 188 188	0	0

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.

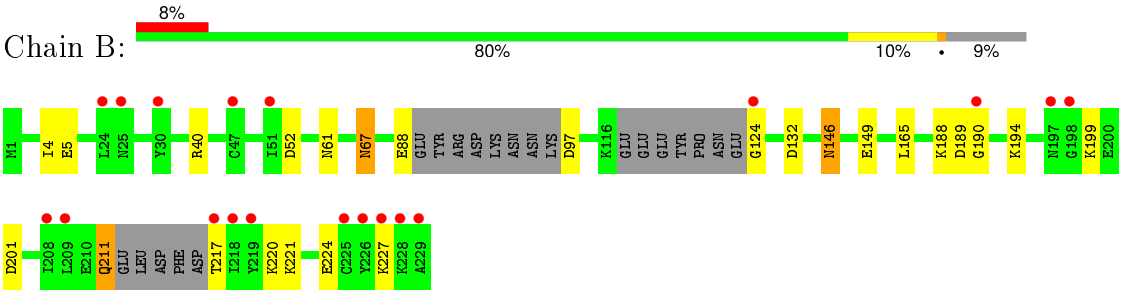
- Molecule 1: DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDE HYDROLASE

Chain A:



- Molecule 1: DEOXYURIDINE 5'-TRIPHOSPHATE NUCLEOTIDE HYDROLASE

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	66.96Å 70.63Å 92.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	55.90 – 1.65 48.59 – 1.65	Depositor EDS
% Data completeness (in resolution range)	99.0 (55.90-1.65) 99.0 (48.59-1.65)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.26 (at 1.65Å)	Xtriage
Refinement program	REFMAC 5.2.0003	Depositor
R, R_{free}	0.151 , 0.194 0.237 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å^2)	25.0	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 54.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.35$	Xtriage
Outliers	1 of 52878 reflections (0.002%)	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	4067	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.77% 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: MG, DUN

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	1.41	18/1953 (0.9%)	1.56	24/2633 (0.9%)
1	B	1.85	8/1777 (0.5%)	1.14	13/2391 (0.5%)
All	All	1.64	26/3730 (0.7%)	1.38	37/5024 (0.7%)

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	2	5
1	B	1	0
All	All	3	5

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	211	GLN	CB-CG	-41.82	0.39	1.52
1	B	224	GLU	CB-CG	32.54	2.13	1.52
1	B	146	ASN	CG-OD1	29.28	1.88	1.24
1	B	146	ASN	CG-ND2	-28.83	0.60	1.32
1	A	90	TYR	CB-CG	-20.79	1.20	1.51
1	A	143	PHE	C-N	-17.70	1.01	1.33
1	A	143	PHE	CA-C	-16.14	1.10	1.52
1	A	144	GLY	C-N	-15.52	0.98	1.34
1	A	8	GLU	CB-CG	-14.52	1.24	1.52
1	B	227	LYS	CA-CB	-13.44	1.24	1.53
1	A	94	ASN	CA-CB	-12.12	1.21	1.53
1	B	220	LYS	CA-CB	11.74	1.79	1.53
1	A	227	LYS	CA-CB	-11.23	1.29	1.53
1	B	149	GLU	CA-CB	-11.21	1.29	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	123	GLU	CG-CD	10.66	1.68	1.51
1	A	95	ASN	CA-CB	-10.25	1.26	1.53
1	A	145	PHE	N-CA	9.78	1.66	1.46
1	A	197	ASN	CA-CB	-9.55	1.28	1.53
1	A	224	GLU	CG-CD	9.17	1.65	1.51
1	A	112	GLN	CG-CD	8.91	1.71	1.51
1	A	220	LYS	CG-CD	-7.75	1.26	1.52
1	A	166	GLU	CD-OE1	7.25	1.33	1.25
1	A	118	GLU	CD-OE2	6.79	1.33	1.25
1	A	69	GLU	CB-CG	-6.79	1.39	1.52
1	B	221	LYS	CA-CB	6.46	1.68	1.53
1	A	89	GLU	CA-CB	-5.36	1.42	1.53

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	143	PHE	O-C-N	-50.35	37.60	123.20
1	A	143	PHE	CA-C-N	23.65	163.50	116.20
1	B	146	ASN	CB-CG-OD1	-19.89	81.83	121.60
1	B	146	ASN	CB-CG-ND2	16.51	156.32	116.70
1	B	221	LYS	N-CA-CB	-16.38	81.12	110.60
1	A	143	PHE	N-CA-C	12.41	144.50	111.00
1	A	143	PHE	C-N-CA	-11.87	97.38	122.30
1	A	224	GLU	CB-CG-CD	-11.80	82.35	114.20
1	B	227	LYS	CB-CA-C	11.34	133.08	110.40
1	B	220	LYS	CB-CA-C	10.05	130.50	110.40
1	B	224	GLU	CA-CB-CG	-8.76	94.14	113.40
1	A	40[A]	ARG	NE-CZ-NH2	-8.74	115.93	120.30
1	A	40[B]	ARG	NE-CZ-NH2	-8.74	115.93	120.30
1	A	119	GLU	CA-CB-CG	8.56	132.23	113.40
1	B	199	LYS	CD-CE-NZ	7.95	129.98	111.70
1	A	90	TYR	CA-CB-CG	7.47	127.59	113.40
1	A	94	ASN	N-CA-CB	7.45	124.01	110.60
1	B	52	ASP	CB-CG-OD1	7.16	124.74	118.30
1	A	220	LYS	CB-CG-CD	6.87	129.46	111.60
1	A	120	TYR	CB-CA-C	-6.84	96.72	110.40
1	B	132	ASP	CB-CG-OD1	6.78	124.40	118.30
1	A	227	LYS	CB-CA-C	6.66	123.72	110.40
1	A	52	ASP	CB-CG-OD1	6.34	124.01	118.30
1	A	227	LYS	N-CA-CB	6.33	121.99	110.60
1	B	201	ASP	CB-CG-OD1	6.33	124.00	118.30
1	A	125	ASP	CB-CG-OD2	6.27	123.95	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	220	LYS	N-CA-CB	-6.23	99.39	110.60
1	A	94	ASN	CB-CA-C	6.04	122.48	110.40
1	A	193	LYS	CD-CE-NZ	5.87	125.20	111.70
1	A	123	GLU	CG-CD-OE1	-5.68	106.94	118.30
1	A	214	ASP	CB-CG-OD2	5.62	123.35	118.30
1	A	143	PHE	CB-CA-C	-5.27	99.86	110.40
1	A	72[A]	ARG	CA-CB-CG	5.24	124.93	113.40
1	A	72[B]	ARG	CA-CB-CG	5.24	124.93	113.40
1	B	149	GLU	CB-CA-C	5.16	120.71	110.40
1	B	211	GLN	CA-CB-CG	-5.14	102.09	113.40
1	A	12	LYS	CD-CE-NZ	-5.08	100.01	111.70

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	94	ASN	CA
1	A	227	LYS	CA
1	B	227	LYS	CA

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	123	GLU	Sidechain
1	A	143	PHE	Mainchain,Peptide
1	A	144	GLY	Mainchain,Peptide

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	1895	0	1865	26	0
1	B	1735	0	1736	17	0
2	A	24	0	12	0	0
2	B	24	0	12	0	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
4	A	195	0	0	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	188	0	0	11	0
All	All	4067	0	3625	39	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 6.

All (39) 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:217:THR:N	4:B:2178:HOH:O	1.96	0.97
1:B:189:ASP:HB2	4:B:2170:HOH:O	1.67	0.92
1:A:9:ASN:HB2	4:A:2011:HOH:O	1.85	0.77
1:A:144:GLY:O	4:A:2147:HOH:O	2.01	0.77
1:A:61:ASN:ND2	1:B:190:GLY:HA2	2.00	0.76
1:A:61:ASN:HB3	4:A:2070:HOH:O	1.93	0.69
1:B:88:GLU:C	4:B:2109:HOH:O	2.30	0.68
1:A:89:GLU:O	1:A:90:TYR:HB2	1.93	0.67
1:A:9:ASN:CB	4:A:2011:HOH:O	2.43	0.66
1:A:190:GLY:HA2	1:B:61:ASN:ND2	2.11	0.66
1:A:203:GLU:OE2	4:A:2182:HOH:O	2.14	0.64
1:B:61:ASN:ND2	4:B:2082:HOH:O	2.30	0.61
1:B:124:GLY:HA2	4:B:2126:HOH:O	2.00	0.61
1:A:61:ASN:HD21	1:B:190:GLY:HA2	1.66	0.60
1:A:190:GLY:HA2	1:B:61:ASN:HD21	1.67	0.59
1:A:144:GLY:CA	1:A:145:PHE:CB	2.86	0.53
1:A:144:GLY:CA	1:A:145:PHE:HB2	2.39	0.53
1:A:170:LYS:HD2	1:A:213:LEU:HD22	1.90	0.53
1:A:61:ASN:ND2	4:A:2071:HOH:O	2.38	0.52
1:A:9:ASN:ND2	1:A:169:TYR:OH	2.42	0.52
1:A:144:GLY:HA3	1:A:145:PHE:CB	2.40	0.52
1:A:208:ILE:O	1:A:211:GLN:HG2	2.12	0.50
1:B:124:GLY:CA	4:B:2126:HOH:O	2.57	0.50
4:A:2066:HOH:O	1:B:194:LYS:HE3	2.13	0.49
1:A:61:ASN:CB	4:A:2070:HOH:O	2.57	0.48
1:B:5:GLU:OE1	4:B:2006:HOH:O	2.20	0.48
1:B:67:ASN:C	1:B:67:ASN:HD22	2.17	0.47
1:B:40[A]:ARG:NH2	4:B:2047:HOH:O	2.44	0.46
1:A:72[B]:ARG:HD2	1:A:163:LEU:HG	1.96	0.46
1:B:97:ASP:N	4:B:2112:HOH:O	2.49	0.45
1:B:4:ILE:HD11	4:B:2118:HOH:O	2.16	0.45
1:A:112:GLN:HG3	1:A:112:GLN:H	1.61	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:LYS:NZ	1:A:213:LEU:HD13	2.33	0.43
1:A:40[A]:ARG:CZ	4:A:2044:HOH:O	2.67	0.42
1:A:40[B]:ARG:HD3	1:A:145:PHE:CD1	2.55	0.41
1:A:135:LEU:HG	1:A:139:LYS:HE2	2.03	0.41
1:A:123:GLU:HG3	4:A:2079:HOH:O	2.20	0.41
1:A:40[B]:ARG:NE	1:A:145:PHE:CD2	2.90	0.40
1:B:40[A]:ARG:NE	4:B:2047:HOH:O	2.48	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	226/229 (99%)	223 (99%)	1 (0%)	2 (1%)	21 4
1	B	203/229 (89%)	202 (100%)	1 (0%)	0	100 100
All	All	429/458 (94%)	425 (99%)	2 (0%)	2 (0%)	34 12

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	143	PHE
1	A	145	PHE

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	210/209 (100%)	207 (99%)	3 (1%)	74 53
1	B	191/209 (91%)	186 (97%)	5 (3%)	54 25
All	All	401/418 (96%)	393 (98%)	8 (2%)	63 38

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90	TYR
1	A	95	ASN
1	A	112	GLN
1	B	67	ASN
1	B	146	ASN
1	B	165	LEU
1	B	188	LYS
1	B	211	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	61	ASN
1	B	61	ASN
1	B	67	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 carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 8 ligands modelled in this entry, 6 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	$\#\|Z\| > 2$
2	DUN	A	1230	3	20,25,25	1.44	3 (15%)	26,38,38	2.56	6 (23%)
2	DUN	B	1230	3	20,25,25	2.31	4 (20%)	26,38,38	2.81	6 (23%)

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	DUN	A	1230	3	-	0/9/28/28	0/2/2/2
2	DUN	B	1230	3	-	0/9/28/28	0/2/2/2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1230	DUN	PB-O1B	-2.80	1.48	1.56
2	B	1230	DUN	PA-O1A	-2.12	1.50	1.56
2	A	1230	DUN	PA-N3A	2.19	1.69	1.63
2	A	1230	DUN	PB-O2B	2.59	1.49	1.46
2	B	1230	DUN	PB-O2B	3.91	1.50	1.46
2	A	1230	DUN	PA-O2A	4.33	1.51	1.46
2	B	1230	DUN	PA-O2A	8.16	1.55	1.46

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	1230	DUN	O2B-PB-N3A	-7.14	100.95	111.90
2	B	1230	DUN	O2A-PA-N3A	-5.65	103.24	111.90
2	A	1230	DUN	O2A-PA-N3A	-3.53	106.48	111.90
2	A	1230	DUN	O2B-PB-N3A	-3.21	106.98	111.90
2	A	1230	DUN	O5'-PA-O2A	-2.63	102.41	113.31
2	B	1230	DUN	O5'-PA-O2A	-2.43	103.24	113.31
2	A	1230	DUN	C5-C4-N3	-2.25	117.35	123.12
2	B	1230	DUN	O4'-C1'-N1	2.51	112.06	107.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1230	DUN	O1A-PA-O2A	4.05	118.46	110.00
2	B	1230	DUN	O1A-PA-O2A	4.46	119.32	110.00
2	B	1230	DUN	C4-N3-C2	8.74	122.80	114.14
2	A	1230	DUN	C4-N3-C2	10.11	124.15	114.14

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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	226/229 (98%)	0.57	11 (4%) 33 31	27, 33, 46, 59	23 (10%)
1	B	209/229 (91%)	0.71	19 (9%) 11 9	27, 34, 43, 52	15 (7%)
All	All	435/458 (94%)	0.63	30 (6%) 20 17	27, 33, 45, 59	38 (8%)

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	213	LEU	10.3
1	B	229	ALA	6.7
1	B	218	ILE	4.8
1	B	197	ASN	4.7
1	B	226	TYR	4.5
1	A	229	ALA	4.3
1	B	227	LYS	4.3
1	B	217	THR	3.7
1	B	198	GLY	3.6
1	B	228	LYS	3.4
1	A	212	GLU	3.4
1	B	208	ILE	3.4
1	B	225	CYS	3.3
1	B	219	TYR	3.3
1	B	209	LEU	3.0
1	A	144	GLY	2.8
1	B	24	LEU	2.7
1	A	137	ILE	2.6
1	B	124	GLY	2.6
1	A	75	ILE	2.4
1	A	228	LYS	2.3
1	B	30	TYR	2.2
1	B	25	ASN	2.2
1	A	226	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	51	ILE	2.1
1	B	47	CYS	2.1
1	A	150	LEU	2.1
1	A	42	CYS	2.0
1	A	215	PHE	2.0
1	B	190	GLY	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	DUN	B	1230	24/24	0.97	0.06	-1.21	18,22,25,32	0
3	MG	A	1233	1/1	0.94	0.08	-1.90	17,17,17,17	1
3	MG	B	1233	1/1	0.93	0.05	-2.65	16,16,16,16	1
2	DUN	A	1230	24/24	0.99	0.04	-3.15	15,18,21,30	0
3	MG	B	1232	1/1	0.99	0.05	-4.69	19,19,19,19	0
3	MG	A	1232	1/1	0.99	0.03	-6.27	18,18,18,18	0
3	MG	A	1231	1/1	0.97	0.06	-	17,17,17,17	0
3	MG	B	1231	1/1	0.99	0.03	-	19,19,19,19	0

6.5 Other polymers (i)

There are no such residues in this entry.