



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 02:18 AM GMT

PDB ID : 2GIC
Title : Crystal Structure of a vesicular stomatitis virus nucleocapsid-RNA complex
Authors : Green, T.J.; Zhang, X.; Wertz, G.W.; Luo, M.
Deposited on : 2006-03-28
Resolution : 2.92 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

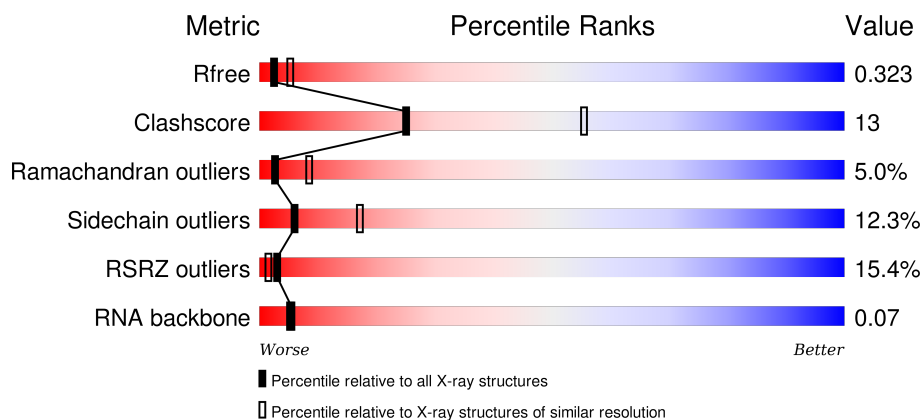
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.92 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	1643 (2.94-2.90)
Clashscore	102246	1871 (2.94-2.90)
Ramachandran outliers	100387	1824 (2.94-2.90)
Sidechain outliers	100360	1826 (2.94-2.90)
RSRZ outliers	91569	1650 (2.94-2.90)
RNA backbone	2183	1004 (3.30-2.54)

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	R	45	<div> <div>38%</div> <div> <div>9%</div> <div>29%</div> <div>51%</div> <div>11%</div> </div> </div>
2	A	422	<div> <div>15%</div> <div>70%</div> <div>24%</div> <div>5%</div> </div>
2	B	422	<div> <div>12%</div> <div>67%</div> <div>25%</div> <div>7%</div> <div>.</div> </div>
2	C	422	<div> <div>12%</div> <div>68%</div> <div>23%</div> <div>6%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
2	D	422	<div><div></div><div>13%</div><div>70%</div><div>24%</div><div>5%</div></div>
2	E	422	<div><div></div><div>22%</div><div>68%</div><div>26%</div><div>5%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17432 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.

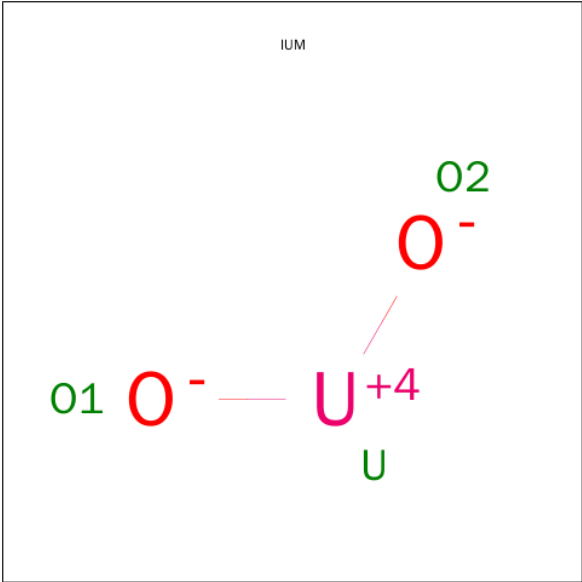
- Molecule 1 is a RNA chain called 45-MER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	R	45	Total	C	N	O	P	0	0	0
			900	405	90	360	45			

- Molecule 2 is a protein called Nucleocapsid protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	421	Total	C	N	O	S	0	0	0
			3327	2118	558	633	18			
2	B	415	Total	C	N	O	S	0	0	0
			3290	2097	552	623	18			
2	C	413	Total	C	N	O	S	0	0	0
			3275	2089	550	618	18			
2	D	416	Total	C	N	O	S	0	0	0
			3298	2103	553	624	18			
2	E	421	Total	C	N	O	S	0	0	0
			3327	2118	558	633	18			

- Molecule 3 is URANYL (VI) ION (three-letter code: IUM) (formula: O₂U).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	U	0	0
			1	1		
3	A	1	Total	U	0	0
			1	1		
3	B	1	Total	U	0	0
			1	1		
3	B	1	Total	U	0	0
			1	1		
3	C	1	Total	U	0	0
			1	1		
3	C	1	Total	U	0	0
			1	1		
3	D	1	Total	U	0	0
			1	1		
3	D	1	Total	U	0	0
			1	1		
3	E	1	Total	U	0	0
			1	1		
3	E	1	Total	U	0	0
			1	1		
3	A	1	Total	U	0	0
			1	1		
3	E	1	Total	U	0	0
			1	1		
3	A	1	Total	U	0	0
			1	1		
3	B	1	Total	U	0	0
			1	1		

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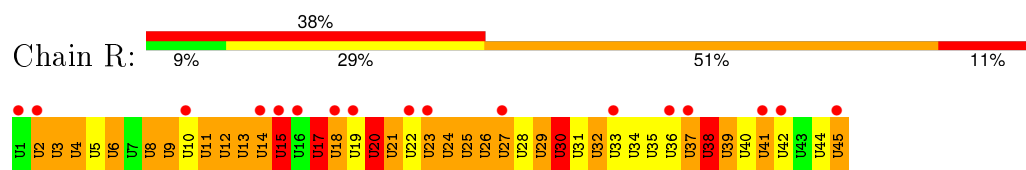
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	1	Total	U	0	0
			1	1		

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.

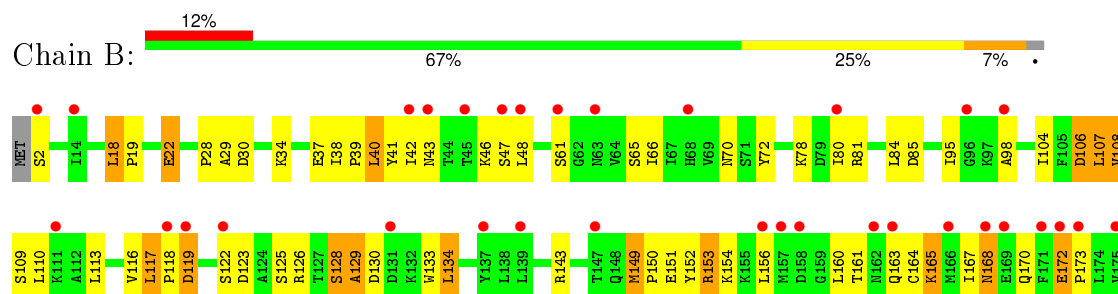
- Molecule 1: 45-MER

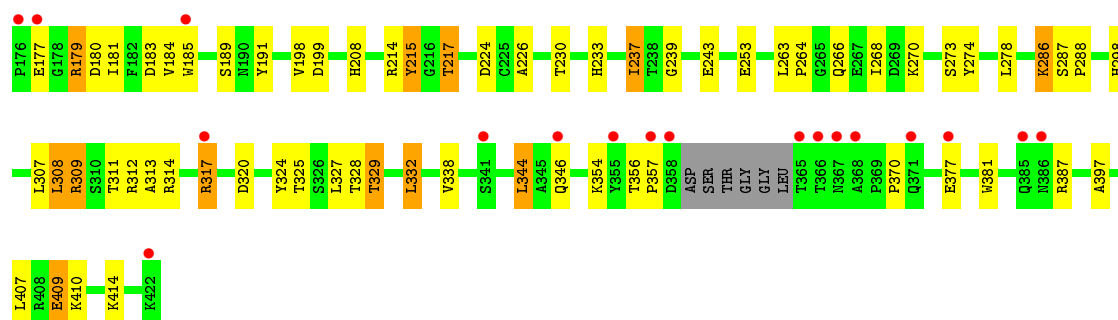


- Molecule 2: Nucleocapsid protein

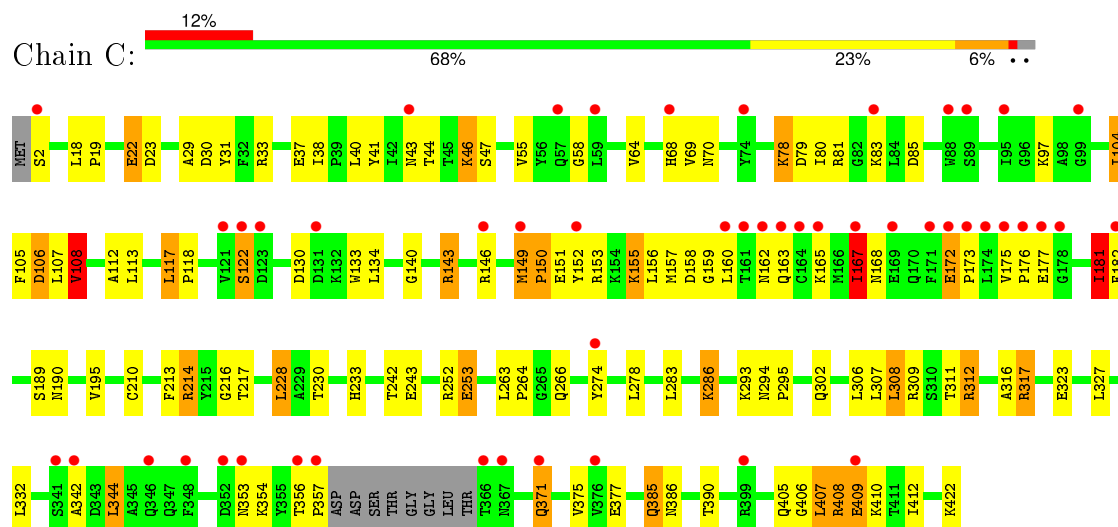


- Molecule 2: Nucleocapsid protein

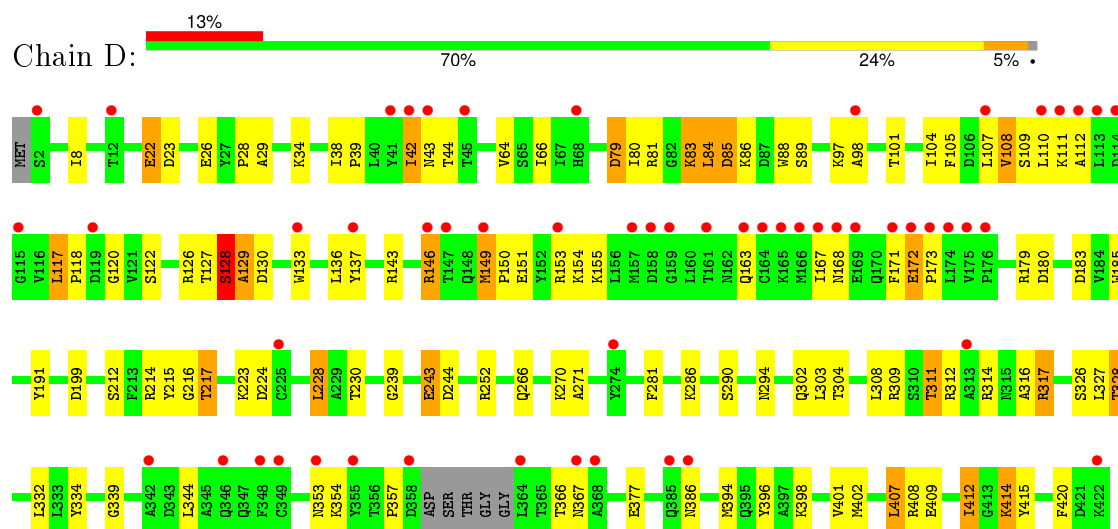




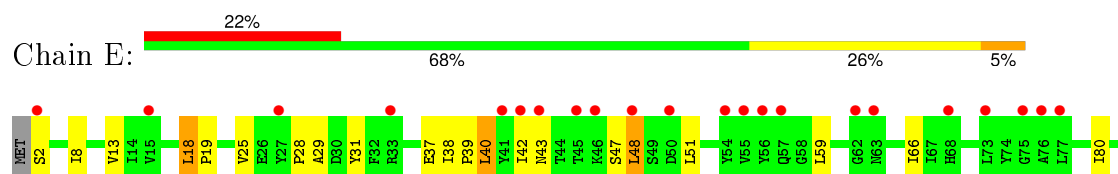
• Molecule 2: Nucleocapsid protein

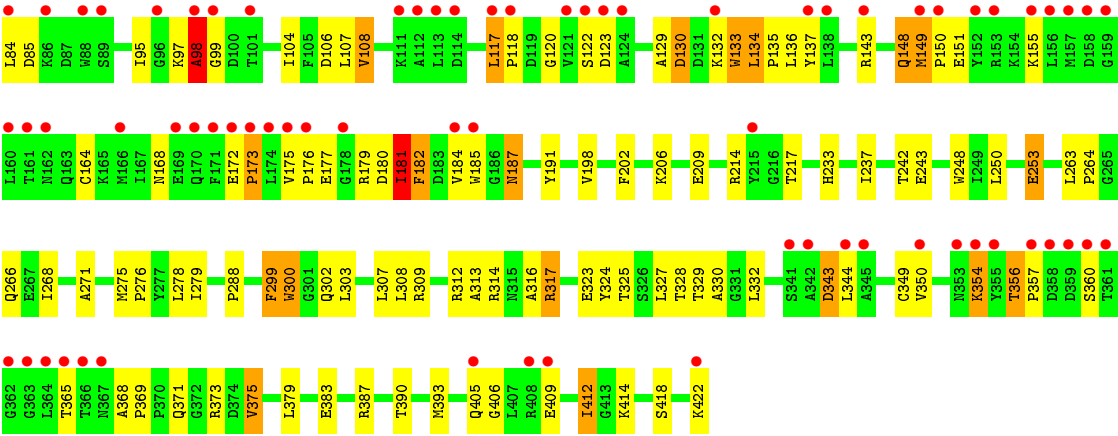


• Molecule 2: Nucleocapsid protein



• Molecule 2: Nucleocapsid protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	166.16 Å 236.32 Å 75.65 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.86 – 2.92 44.84 – 2.92	Depositor EDS
% Data completeness (in resolution range)	91.9 (44.86-2.92) 90.5 (44.84-2.92)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.95 (at 2.90 Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.255 , 0.306 0.270 , 0.323	Depositor DCC
R_{free} test set	3262 reflections (5.79%)	DCC
Wilson B-factor (Å ²)	73.1	Xtriage
Anisotropy	0.676	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 33.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	4 of 64651 reflections (0.006%)	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	17432	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.35% 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, 0.375 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: IUM

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	R	0.93	0/989	1.56	10/1526 (0.7%)
2	A	0.39	0/3403	0.54	0/4607
2	B	0.40	0/3365	0.55	0/4554
2	C	0.39	0/3350	0.56	0/4533
2	D	0.39	0/3373	0.56	0/4565
2	E	0.40	0/3403	0.54	0/4607
All	All	0.44	0/17883	0.66	10/24392 (0.0%)

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
2	B	0	1
2	E	0	1
All	All	0	2

There are no bond length outliers.

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	R	20	U	P-O3'-C3'	8.96	130.45	119.70
1	R	9	U	O4'-C1'-N1	6.62	113.50	108.20
1	R	9	U	O4'-C4'-C3'	-6.12	97.88	104.00
1	R	34	U	O4'-C1'-N1	5.87	112.90	108.20
1	R	15	U	C2-N1-C1'	5.75	124.61	117.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	106	ASP	Peptide
2	E	98	ALA	Peptide

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	R	900	0	451	62	0
2	A	3327	0	3287	87	0
2	B	3290	0	3253	105	0
2	C	3275	0	3242	81	0
2	D	3298	0	3264	75	0
2	E	3327	0	3287	92	0
3	A	4	0	0	0	0
3	B	3	0	0	0	0
3	C	3	0	0	0	0
3	D	2	0	0	0	0
3	E	3	0	0	0	0
All	All	17432	0	16784	435	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 13.

The worst 5 of 435 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)
2:A:143:ARG:HD2	2:A:155:LYS:CE	1.44	1.45
2:A:143:ARG:CD	2:A:155:LYS:HE2	1.52	1.36
2:B:109:SER:O	2:B:110:LEU:HD23	1.31	1.22
2:D:107:LEU:O	2:D:108:VAL:HG22	1.42	1.18
2:B:107:LEU:N	2:B:107:LEU:HD23	1.56	1.16

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	
2	A	419/422 (99%)	354 (84%)	46 (11%)	19 (4%)	3	11
2	B	411/422 (97%)	344 (84%)	45 (11%)	22 (5%)	2	7
2	C	409/422 (97%)	340 (83%)	49 (12%)	20 (5%)	3	9
2	D	412/422 (98%)	343 (83%)	48 (12%)	21 (5%)	2	8
2	E	419/422 (99%)	349 (83%)	48 (12%)	22 (5%)	2	7
All	All	2070/2110 (98%)	1730 (84%)	236 (11%)	104 (5%)	3	8

5 of 104 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	47	SER
2	A	98	ALA
2	A	113	LEU
2	A	172	GLU
2	B	98	ALA

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	
2	A	362/363 (100%)	323 (89%)	39 (11%)	8	23
2	B	358/363 (99%)	311 (87%)	47 (13%)	5	15
2	C	356/363 (98%)	303 (85%)	53 (15%)	4	11
2	D	359/363 (99%)	315 (88%)	44 (12%)	6	17

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	E	362/363 (100%)	324 (90%)	38 (10%)	8	25
All	All	1797/1815 (99%)	1576 (88%)	221 (12%)	6	17

5 of 221 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	C	134	LEU
2	C	308	LEU
2	E	242	THR
2	C	152	TYR
2	C	189	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 37 such sidechains are listed below:

Mol	Chain	Res	Type
2	C	68	HIS
2	D	43	ASN
2	E	315	ASN
2	C	70	ASN
2	C	168	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	R	44/45 (97%)	35 (79%)	8 (18%)

5 of 35 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	R	2	U
1	R	3	U
1	R	4	U
1	R	5	U
1	R	6	U

5 of 8 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	R	12	U

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Mol	Chain	Res	Type
1	R	39	U
1	R	20	U
1	R	6	U
1	R	14	U

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 15 ligands modelled in this entry, 15 are modelled with single atom - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	R	45/45 (100%)	1.88	17 (37%) 0 0	29, 44, 53, 58	0
2	A	421/422 (99%)	0.96	63 (14%) 3 2	23, 27, 35, 40	0
2	B	415/422 (98%)	0.89	51 (12%) 5 3	23, 27, 33, 42	0
2	C	413/422 (97%)	0.84	50 (12%) 6 3	23, 27, 33, 39	0
2	D	416/422 (98%)	0.94	55 (13%) 4 2	23, 27, 34, 43	0
2	E	421/422 (99%)	1.18	92 (21%) 1 1	23, 27, 36, 45	0
All	All	2131/2155 (98%)	0.98	328 (15%) 3 1	23, 27, 37, 58	0

The worst 5 of 328 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	2	SER	13.4
2	A	2	SER	11.6
2	A	361	THR	9.8
2	E	101	THR	8.1
2	D	166	MET	8.1

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
3	IUM	B	423	1/3	0.93	0.10	-1.35	129,129,129,129	0
3	IUM	A	424	1/3	0.99	0.14	-1.72	21,21,21,21	0
3	IUM	C	424	1/3	0.99	0.14	-1.97	22,22,22,22	0
3	IUM	E	424	1/3	0.99	0.14	-2.65	21,21,21,21	0
3	IUM	E	423	1/3	0.94	0.06	-3.12	107,107,107,107	0
3	IUM	B	424	1/3	0.99	0.12	-3.37	15,15,15,15	0
3	IUM	D	424	1/3	0.99	0.12	-5.04	17,17,17,17	0
3	IUM	A	423	1/3	0.95	0.09	-	92,92,92,92	0
3	IUM	C	423	1/3	0.97	0.11	-	54,54,54,54	0
3	IUM	B	425	1/3	0.96	0.04	-	184,184,184,184	0
3	IUM	E	425	1/3	0.79	0.07	-	192,192,192,192	0
3	IUM	A	425	1/3	0.97	0.03	-	129,129,129,129	0
3	IUM	C	425	1/3	0.94	0.05	-	123,123,123,123	0
3	IUM	A	426	1/3	0.95	0.14	-	114,114,114,114	0
3	IUM	D	423	1/3	0.85	0.09	-	120,120,120,120	0

6.5 Other polymers [i](#)

There are no such residues in this entry.