



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 02:08 AM UTC

PDB ID : 3PU0 / pdb_00003pu0
Title : Crystal Structure of a vesicular stomatitis virus nucleocapsid-polyC complex
Authors : Luo, M.; Green, T.J.; Rowse, M.
Deposited on : 2010-12-03
Resolution : 3.09 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

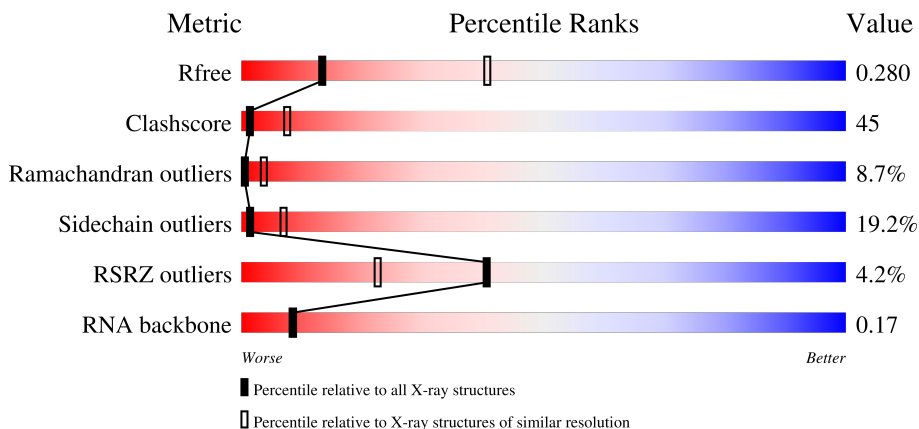
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.09 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.



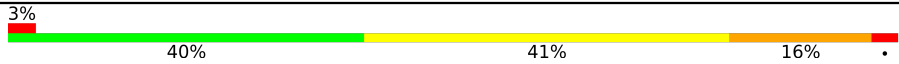

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)
RNA backbone	3983	1022 (3.32-2.88)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	421	 4% 39% 43% 16%
1	B	421	 4% 39% 41% 16%
1	C	421	 4% 39% 42% 15%
1	D	421	 6% 39% 43% 14%

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Mol	Chain	Length	Quality of chain
1	E	421	
2	R	45	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17437 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.

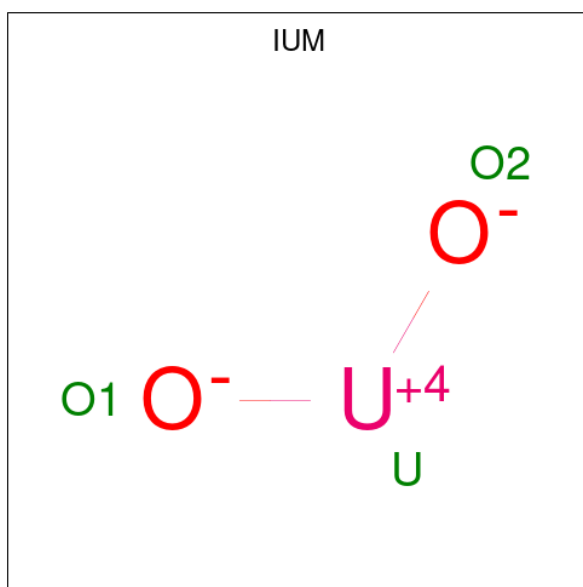
- Molecule 1 is a protein called Nucleoprotein.

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

- Molecule 2 is a RNA chain called RNA (45-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	R	45	900	405	135	315	45	0	0	0

- Molecule 3 is URANYL (VI) ION (CCD ID: IUM) (formula: O₂U).

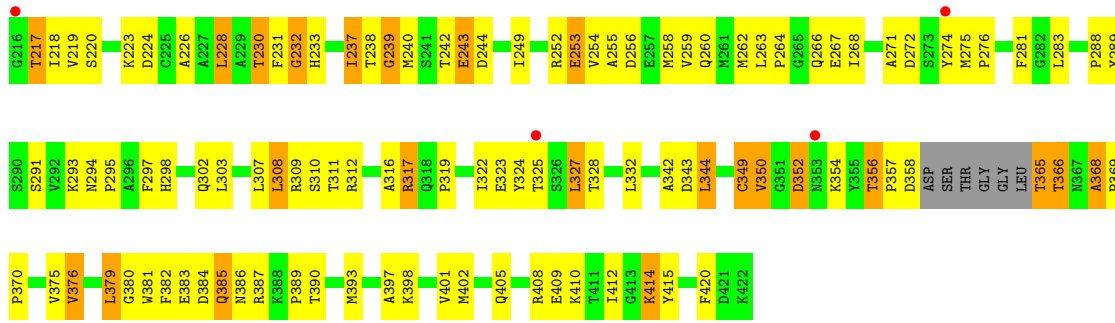


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total U 1 1	0	0
3	A	1	Total U 1 1	0	0
3	A	1	Total U 1 1	0	0
3	B	1	Total U 1 1	0	0
3	B	1	Total U 1 1	0	0
3	B	1	Total U 1 1	0	0
3	B	1	Total U 1 1	0	0
3	C	1	Total U 1 1	0	0
3	C	1	Total U 1 1	0	0
3	D	1	Total U 1 1	0	0
3	D	1	Total U 1 1	0	0
3	D	1	Total U 1 1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

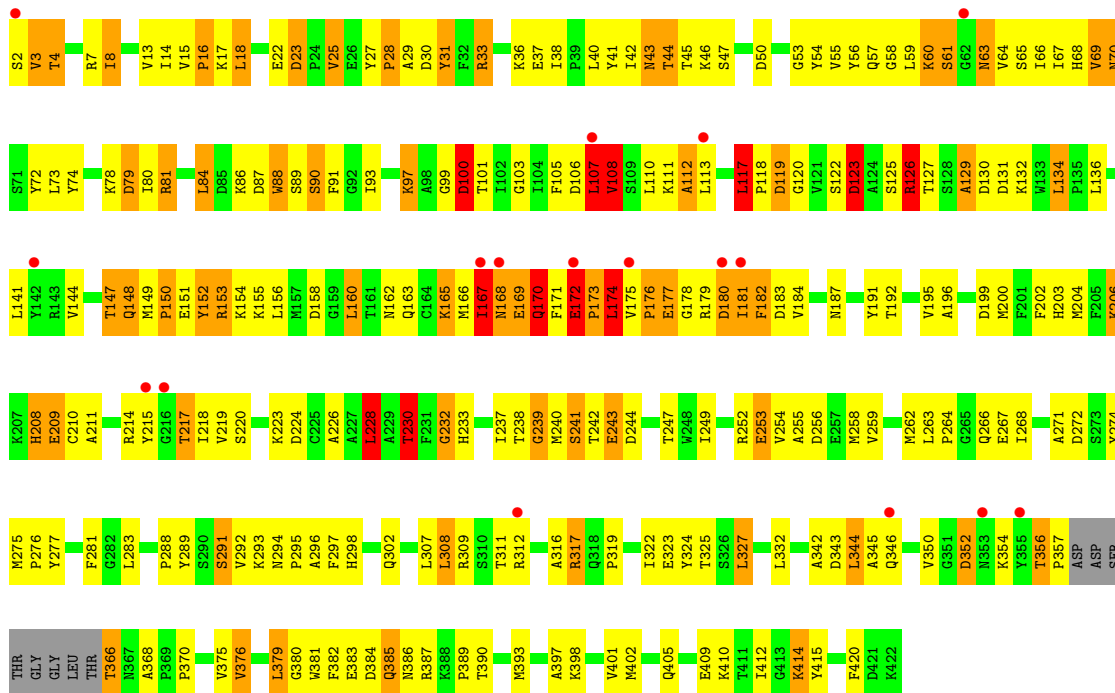
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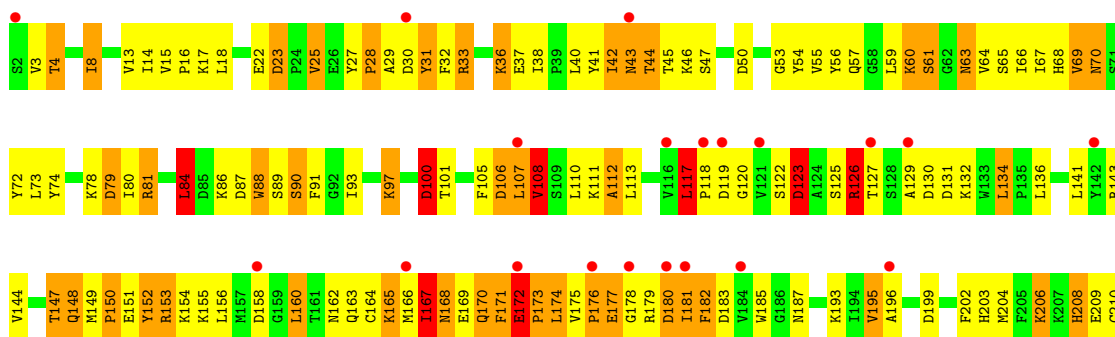
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	E	1	Total U 1 1	0	0
3	R	1	Total U 1 1	0	0
3	R	1	Total U 1 1	0	0
3	R	1	Total U 1 1	0	0
3	R	1	Total U 1 1	0	0
3	R	1	Total U 1 1	0	0

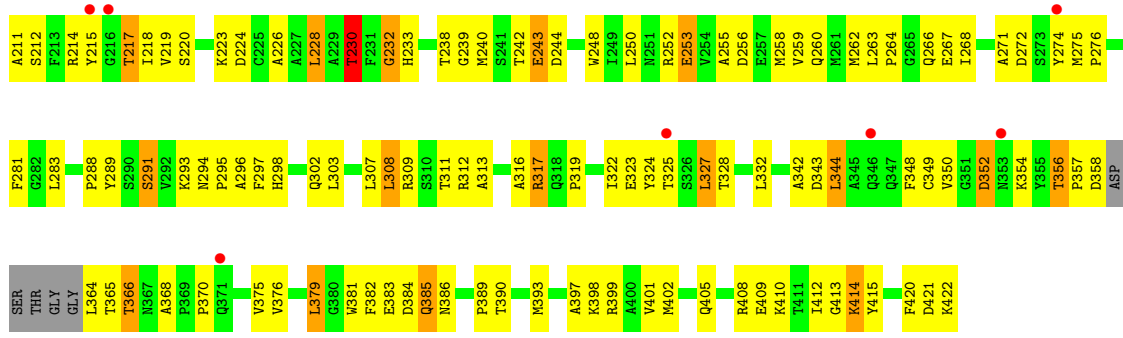


● Molecule 1: Nucleoprotein

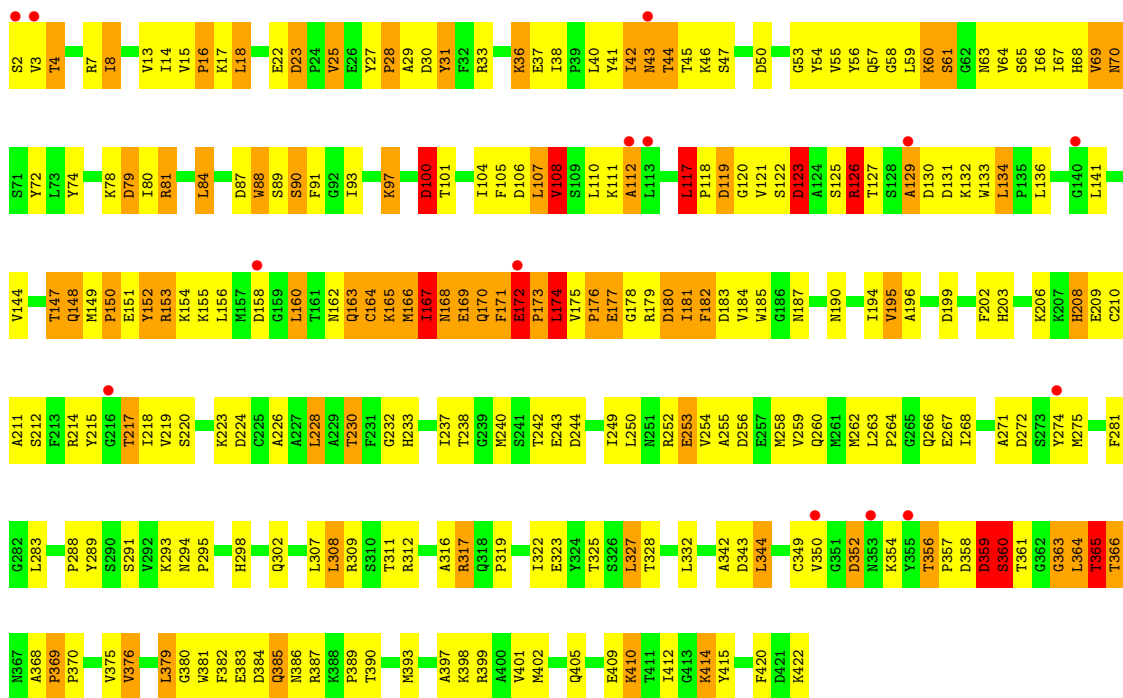


● Molecule 1: Nucleoprotein





• Molecule 1: Nucleoprotein



• Molecule 2: RNA (45-MER)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	164.65Å 235.55Å 75.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.75 – 3.09 49.75 – 3.09	Depositor EDS
% Data completeness (in resolution range)	78.5 (49.75-3.09) 78.5 (49.75-3.09)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.16 (at 3.07Å)	Xtrriage
Refinement program	PHENIX 1.6_289	Depositor
R, R_{free}	0.250 , 0.290 0.239 , 0.280	Depositor DCC
R_{free} test set	2000 reflections (4.34%)	wwPDB-VP
Wilson B-factor (Å ²)	73.2	Xtrriage
Anisotropy	0.879	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 57.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	17437	wwPDB-VP
Average B, all atoms (Å ²)	104.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.74% 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: 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	A	0.66	1/3403 (0.0%)	1.00	15/4607 (0.3%)
1	B	0.66	0/3365	0.97	13/4554 (0.3%)
1	C	0.65	0/3350	0.98	11/4533 (0.2%)
1	D	0.68	0/3373	0.99	12/4565 (0.3%)
1	E	0.65	0/3403	0.98	14/4607 (0.3%)
2	R	0.53	0/989	1.45	24/1526 (1.6%)
All	All	0.65	1/17883 (0.0%)	1.02	89/24392 (0.4%)

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	0	2
1	B	0	4
1	C	0	1
1	D	0	1
1	E	0	3
All	All	0	11

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	358	ASP	CA-C	5.74	1.55	1.52

The worst 5 of 89 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	R	15	C	N1-C1'-C2'	9.66	126.49	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	362	GLY	N-CA-C	9.56	124.56	110.60
1	D	168	ASN	N-CA-C	8.98	123.47	110.10
2	R	18	C	N1-C1'-C2'	-8.57	99.14	112.00
2	R	6	C	N1-C1'-C2'	7.06	122.60	112.00

There are no chirality outliers.

5 of 11 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	169	GLU	Peptide
1	A	78	LYS	Peptide
1	B	107	LEU	Peptide
1	B	167	ILE	Peptide
1	B	169	GLU	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	3327	0	3287	313	0
1	B	3290	0	3253	332	0
1	C	3275	0	3242	338	0
1	D	3298	0	3264	309	0
1	E	3327	0	3287	304	0
2	R	900	0	496	84	0
3	A	3	0	0	0	0
3	B	4	0	0	0	0
3	C	2	0	0	0	0
3	D	3	0	0	0	0
3	E	3	0	0	0	0
3	R	5	0	0	0	0
All	All	17437	0	16829	1542	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 45.

The worst 5 of 1542 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:E:163:GLN:O	1:E:167:ILE:HB	1.36	1.25
1:B:317:ARG:HD3	2:R:22:C:O2'	1.41	1.18
1:E:177:GLU:HA	1:E:181:ILE:HD11	1.31	1.12
1:C:317:ARG:NH2	2:R:14:C:H5'	1.64	1.11
1:A:317:ARG:NH2	2:R:32:C:H5'	1.64	1.11

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	419/421 (100%)	306 (73%)	82 (20%)	31 (7%)	1	5
1	B	411/421 (98%)	302 (74%)	73 (18%)	36 (9%)	0	3
1	C	409/421 (97%)	296 (72%)	73 (18%)	40 (10%)	0	3
1	D	412/421 (98%)	303 (74%)	75 (18%)	34 (8%)	0	4
1	E	419/421 (100%)	303 (72%)	76 (18%)	40 (10%)	0	3
All	All	2070/2105 (98%)	1510 (73%)	379 (18%)	181 (9%)	0	4

5 of 181 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	43	ASN
1	A	44	THR
1	A	60	LYS
1	A	61	SER
1	A	79	ASP

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	362/362 (100%)	292 (81%)	70 (19%)	1	7
1	B	358/362 (99%)	287 (80%)	71 (20%)	1	6
1	C	356/362 (98%)	290 (82%)	66 (18%)	1	8
1	D	359/362 (99%)	291 (81%)	68 (19%)	1	7
1	E	362/362 (100%)	292 (81%)	70 (19%)	1	7
All	All	1797/1810 (99%)	1452 (81%)	345 (19%)	1	7

5 of 345 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	153	ARG
1	E	70	ASN
1	D	172	GLU
1	D	325	THR
1	E	153	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 60 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	162	ASN
1	E	260	GLN
1	C	386	ASN
1	E	203	HIS
1	E	395	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	R	44/45 (97%)	18 (40%)	2 (4%)

5 of 18 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	R	3	C
2	R	5	C

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Mol	Chain	Res	Type
2	R	10	C
2	R	13	C
2	R	17	C

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	R	18	C
2	R	41	C

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 20 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	A	421/421 (100%)	0.33	17 (4%) 42 23	70, 102, 138, 158	0
1	B	415/421 (98%)	0.34	15 (3%) 46 26	71, 101, 137, 157	0
1	C	413/421 (98%)	0.33	17 (4%) 41 23	71, 101, 137, 158	0
1	D	416/421 (98%)	0.42	27 (6%) 25 13	71, 102, 137, 158	0
1	E	421/421 (100%)	0.30	14 (3%) 49 29	70, 102, 140, 163	0
2	R	45/45 (100%)	-0.20	0 100 100	97, 110, 120, 121	0
All	All	2131/2150 (99%)	0.33	90 (4%) 40 22	70, 102, 138, 163	0

The worst 5 of 90 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	2	SER	9.2
1	A	2	SER	7.2
1	A	356	THR	6.2
1	B	2	SER	6.1
1	A	353	ASN	6.0

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 oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
3	IUM	B	527	1/3	0.92	0.07	156,156,156,156	1
3	IUM	E	535	1/3	0.94	0.05	160,160,160,160	1
3	IUM	A	521	1/3	0.96	0.05	132,132,132,132	1
3	IUM	D	531	1/3	0.96	0.05	156,156,156,156	1
3	IUM	B	524	1/3	0.96	0.04	165,165,165,165	1
3	IUM	A	523	1/3	0.97	0.04	175,175,175,175	1
3	IUM	B	525	1/3	0.98	0.04	134,134,134,134	1
3	IUM	D	530	1/3	0.98	0.04	166,166,166,166	1
3	IUM	B	526	1/3	0.99	0.04	125,125,125,125	0
3	IUM	C	528	1/3	0.99	0.02	136,136,136,136	1
3	IUM	E	533	1/3	0.99	0.04	132,132,132,132	1
3	IUM	E	534	1/3	0.99	0.05	132,132,132,132	0
3	IUM	C	529	1/3	0.99	0.03	136,136,136,136	0
3	IUM	R	536	1/3	0.99	0.04	130,130,130,130	0
3	IUM	R	537	1/3	0.99	0.03	122,122,122,122	1
3	IUM	R	538	1/3	0.99	0.02	126,126,126,126	0
3	IUM	R	539	1/3	0.99	0.02	112,112,112,112	1
3	IUM	D	532	1/3	1.00	0.03	125,125,125,125	0
3	IUM	A	522	1/3	1.00	0.03	129,129,129,129	0
3	IUM	R	540	1/3	1.00	0.03	126,126,126,126	1

6.5 Other polymers [i](#)

There are no such residues in this entry.