



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 06:12 PM UTC

PDB ID : 7JTN / pdb_00007jtn
Title : Human Complement Factor B Inhibited by a Slow Off-Rate Modified Aptamer of 29 Bases
Authors : Xu, X.; Geisbrecht, B.V.
Deposited on : 2020-08-18
Resolution : 3.10 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

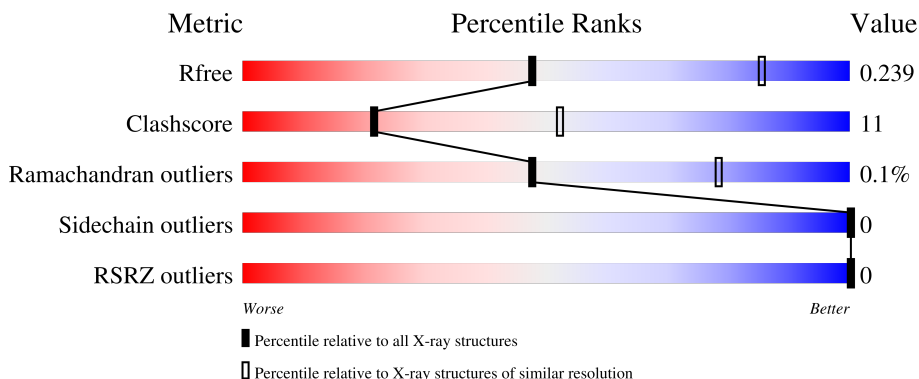
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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)

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	764	 75% 16% 8%
1	C	764	 74% 18% 8%
2	B	30	 67% 30%
2	D	30	 10% 57% 33%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12542 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 Complement factor B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	702	Total	C	N	O	S	0	0	0
			5540	3488	961	1058	33			
1	C	702	Total	C	N	O	S	0	0	0
			5540	3488	961	1058	33			

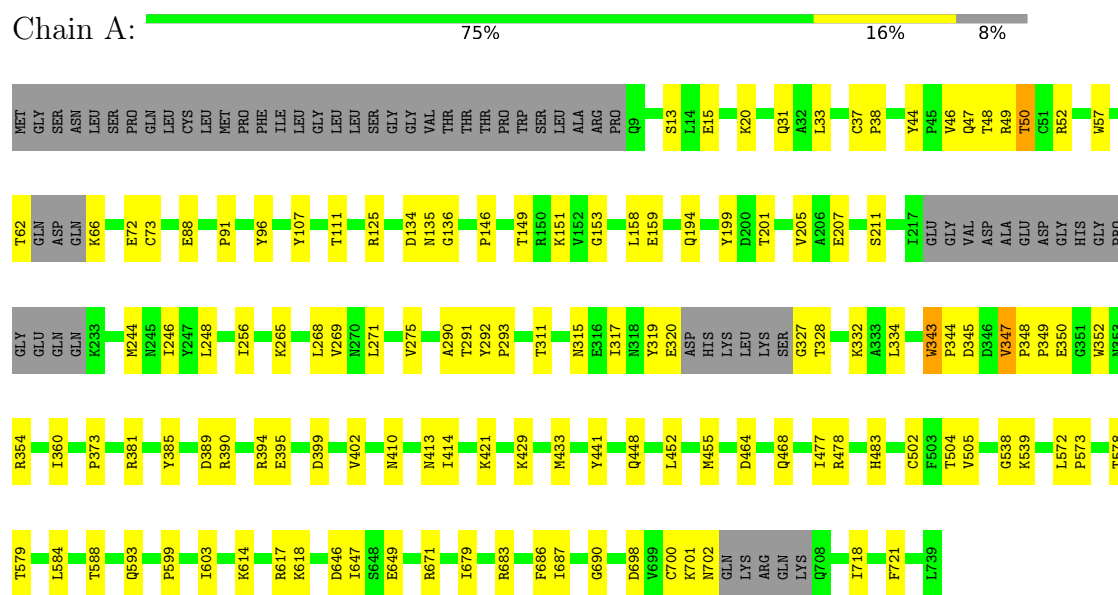
- Molecule 2 is a DNA chain called DNA (30-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	30	Total	C	N	O	P	0	0	0
			731	381	127	193	30			
2	D	30	Total	C	N	O	P	0	0	0
			731	381	127	193	30			

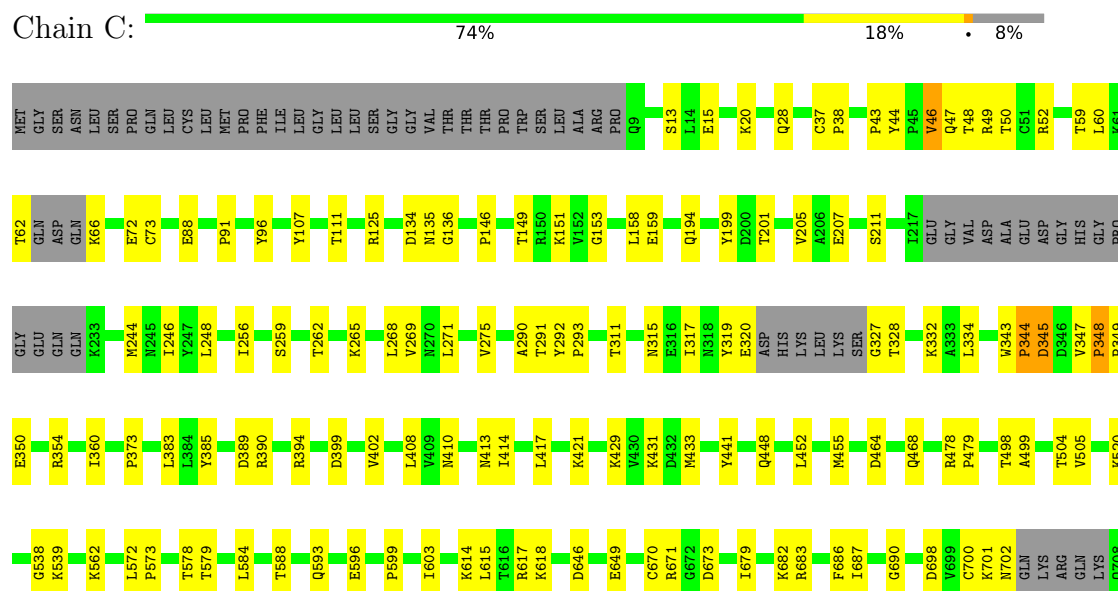
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Complement factor B

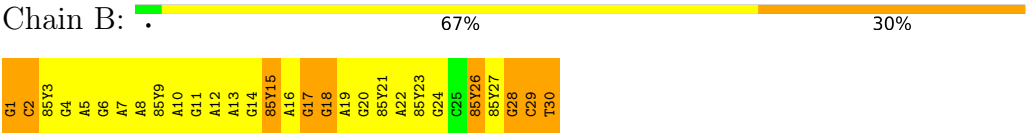


• Molecule 1: Complement factor B

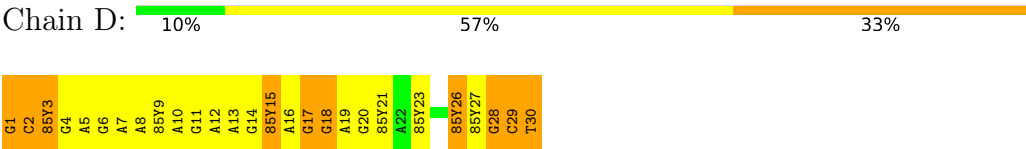




• Molecule 2: DNA (30-MER)



• Molecule 2: DNA (30-MER)



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.13Å 144.41Å 87.19Å 90.00° 109.46° 90.00°	Depositor
Resolution (Å)	42.90 – 3.10 42.90 – 3.10	Depositor EDS
% Data completeness (in resolution range)	98.7 (42.90-3.10) 98.7 (42.90-3.10)	Depositor EDS
R_{merge}	0.23	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 3.12Å)	Xtriage
Refinement program	PHENIX v1.17	Depositor
R, R_{free}	0.208 , 0.241 0.207 , 0.239	Depositor DCC
R_{free} test set	1994 reflections (5.40%)	wwPDB-VP
Wilson B-factor (Å ²)	62.6	Xtriage
Anisotropy	0.304	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 59.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.438 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12542	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.93% 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: OMC, OMG, T3P, 85Y

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	0.65	7/5664 (0.1%)	0.61	12/7667 (0.2%)
1	C	0.66	10/5664 (0.2%)	0.60	8/7667 (0.1%)
2	B	0.71	0/358	0.96	0/544
2	D	0.75	0/358	0.99	0/544
All	All	0.66	17/12044 (0.1%)	0.64	20/16422 (0.1%)

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	343	TRP	C-N	-14.39	1.20	1.33
1	C	38	PRO	N-CD	-12.82	1.29	1.47
1	A	38	PRO	N-CD	-12.79	1.29	1.47
1	C	348	PRO	C-N	8.94	1.44	1.33
1	C	345	ASP	CA-C	-6.43	1.45	1.52

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	343	TRP	O-C-N	-16.51	105.31	120.99
1	A	50	THR	CB-CA-C	-11.06	89.44	110.24
1	C	50	THR	CB-CA-C	-11.05	89.47	110.24
1	A	343	TRP	CA-C-N	9.82	132.78	120.89
1	A	343	TRP	C-N-CA	9.82	132.78	120.89

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	5540	0	5406	102	0
1	C	5540	0	5406	87	0
2	B	731	0	265	34	0
2	D	731	0	265	33	0
All	All	12542	0	11342	253	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 11.

The worst 5 of 253 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:343:TRP:CZ2	1:A:352:TRP:HB2	1.82	1.13
1:A:343:TRP:CZ2	1:A:352:TRP:CB	2.33	1.11
1:A:343:TRP:HZ2	1:A:352:TRP:HB2	1.05	1.10
1:A:343:TRP:CE3	1:A:349:PRO:HD2	1.87	1.09
1:A:343:TRP:CE3	1:A:349:PRO:CD	2.39	1.04

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	692/764 (91%)	655 (95%)	37 (5%)	0	100 100
1	C	692/764 (91%)	652 (94%)	39 (6%)	1 (0%)	48 78

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1384/1528 (91%)	1307 (94%)	76 (6%)	1 (0%)	48 78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	479	PRO

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	613/665 (92%)	613 (100%)	0	100	100
1	C	613/665 (92%)	613 (100%)	0	100	100
All	All	1226/1330 (92%)	1226 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	459	HIS
1	C	468	GLN
1	C	668	ASN
1	A	468	GLN
1	A	77	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

30 non-standard protein/DNA/RNA residues 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	85Y	B	15	2	33,36,37	4.54	13 (39%)	46,51,54	3.01	15 (32%)
2	85Y	B	23	2	33,36,37	4.50	15 (45%)	46,51,54	3.24	16 (34%)
2	T3P	D	30	2	18,21,22	4.94	12 (66%)	26,30,33	2.64	10 (38%)
2	85Y	D	3	2	33,36,37	4.58	15 (45%)	46,51,54	3.12	17 (36%)
2	OMC	B	2	2	19,22,23	3.31	8 (42%)	25,31,34	1.52	3 (12%)
2	85Y	D	27	2	33,36,37	4.93	16 (48%)	46,51,54	3.64	17 (36%)
2	OMG	D	4	2	23,26,27	2.49	9 (39%)	32,38,41	2.42	12 (37%)
2	85Y	D	23	2	33,36,37	4.46	14 (42%)	46,51,54	3.42	14 (30%)
2	85Y	B	9	2	33,36,37	4.53	16 (48%)	46,51,54	2.70	13 (28%)
2	85Y	B	21	2	33,36,37	4.64	16 (48%)	46,51,54	2.94	15 (32%)
2	85Y	D	15	2	33,36,37	4.61	14 (42%)	46,51,54	3.10	15 (32%)
2	OMG	D	28	2	23,26,27	2.49	8 (34%)	32,38,41	2.19	11 (34%)
2	OMG	B	4	2	23,26,27	2.43	9 (39%)	32,38,41	2.49	14 (43%)
2	OMG	B	17	2	23,26,27	2.55	8 (34%)	32,38,41	2.11	11 (34%)
2	OMG	B	28	2	23,26,27	2.57	7 (30%)	32,38,41	2.25	11 (34%)
2	85Y	D	9	2	33,36,37	4.47	16 (48%)	46,51,54	2.77	11 (23%)
2	OMG	D	1	2	23,26,27	2.67	7 (30%)	32,38,41	3.20	16 (50%)
2	85Y	B	26	2	33,36,37	4.56	15 (45%)	46,51,54	3.46	14 (30%)
2	85Y	B	3	2	33,36,37	4.55	14 (42%)	46,51,54	3.38	16 (34%)
2	OMG	B	1	2	23,26,27	2.62	8 (34%)	32,38,41	2.90	16 (50%)
2	OMG	D	17	2	23,26,27	2.51	9 (39%)	32,38,41	2.10	12 (37%)
2	OMG	D	18	2	23,26,27	2.26	9 (39%)	32,38,41	1.83	11 (34%)
2	85Y	B	27	2	33,36,37	4.76	16 (48%)	46,51,54	3.72	18 (39%)
2	85Y	D	21	2	33,36,37	4.71	16 (48%)	46,51,54	3.10	16 (34%)
2	OMC	D	29	2	19,22,23	4.06	11 (57%)	25,31,34	1.93	6 (24%)
2	OMG	B	18	2	23,26,27	2.34	9 (39%)	32,38,41	1.85	10 (31%)
2	OMC	B	29	2	19,22,23	4.44	13 (68%)	25,31,34	0.91	1 (4%)
2	OMC	D	2	2	19,22,23	3.25	8 (42%)	25,31,34	2.08	6 (24%)
2	T3P	B	30	2	18,21,22	5.27	13 (72%)	26,30,33	2.36	10 (38%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	85Y	D	26	2	33,36,37	4.62	16 (48%)	46,51,54	3.52	15 (32%)

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	85Y	B	15	2	-	4/16/30/31	0/4/4/4
2	85Y	B	23	2	-	1/16/30/31	0/4/4/4
2	T3P	D	30	2	-	2/8/22/23	0/2/2/2
2	85Y	D	3	2	-	0/16/30/31	0/4/4/4
2	OMC	B	2	2	-	4/9/27/28	0/2/2/2
2	85Y	D	27	2	-	2/16/30/31	0/4/4/4
2	OMG	D	4	2	-	2/9/27/28	0/3/3/3
2	85Y	D	23	2	-	1/16/30/31	0/4/4/4
2	85Y	B	9	2	-	4/16/30/31	0/4/4/4
2	85Y	B	21	2	-	4/16/30/31	0/4/4/4
2	85Y	D	15	2	-	2/16/30/31	0/4/4/4
2	OMG	D	28	2	-	1/9/27/28	0/3/3/3
2	OMG	B	4	2	-	3/9/27/28	0/3/3/3
2	OMG	B	17	2	-	2/9/27/28	0/3/3/3
2	OMG	B	28	2	-	2/9/27/28	0/3/3/3
2	85Y	D	9	2	-	4/16/30/31	0/4/4/4
2	OMG	D	1	2	-	4/9/27/28	0/3/3/3
2	85Y	B	26	2	-	2/16/30/31	0/4/4/4
2	85Y	B	3	2	-	2/16/30/31	0/4/4/4
2	OMG	B	1	2	-	3/9/27/28	0/3/3/3
2	OMG	D	17	2	-	2/9/27/28	0/3/3/3
2	OMG	D	18	2	-	1/9/27/28	0/3/3/3
2	85Y	B	27	2	-	2/16/30/31	0/4/4/4
2	85Y	D	21	2	-	4/16/30/31	0/4/4/4
2	OMC	D	29	2	-	4/9/27/28	0/2/2/2
2	OMG	B	18	2	-	3/9/27/28	0/3/3/3
2	OMC	B	29	2	-	3/9/27/28	0/2/2/2
2	OMC	D	2	2	-	4/9/27/28	0/2/2/2
2	T3P	B	30	2	-	4/8/22/23	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	85Y	D	26	2	-	2/16/30/31	0/4/4/4

The worst 5 of 360 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	30	T3P	C2'-C3'	-14.99	1.21	1.52
2	B	30	T3P	C2'-C3'	-14.18	1.22	1.52
2	D	27	85Y	C6-N1	13.53	1.61	1.38
2	B	27	85Y	C6-N1	12.54	1.59	1.38
2	B	21	85Y	C6-N1	12.06	1.58	1.38

The worst 5 of 372 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	3	85Y	C24-C23-N22	-13.41	84.80	113.07
2	D	3	85Y	C24-C23-N22	-13.26	85.12	113.07
2	B	23	85Y	C24-C23-N22	11.29	136.88	113.07
2	D	15	85Y	C23-N22-C20	11.08	137.90	122.07
2	D	23	85Y	C24-C23-N22	10.66	135.54	113.07

There are no chirality outliers.

5 of 78 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1	OMG	C1'-C2'-O2'-CM2
2	D	1	OMG	C1'-C2'-O2'-CM2
2	B	2	OMC	C1'-C2'-O2'-CM2
2	B	2	OMC	C3'-C4'-C5'-O5'
2	B	2	OMC	C4'-C5'-O5'-P

There are no ring outliers.

19 monomers are involved in 36 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	15	85Y	2	0
2	D	30	T3P	1	0
2	D	3	85Y	1	0
2	B	2	OMC	5	0
2	D	15	85Y	3	0
2	D	28	OMG	3	0
2	B	17	OMG	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	28	OMG	2	0
2	D	1	OMG	1	0
2	B	26	85Y	1	0
2	B	1	OMG	3	0
2	D	17	OMG	2	0
2	D	18	OMG	3	0
2	D	29	OMC	3	0
2	B	18	OMG	3	0
2	B	29	OMC	4	0
2	D	2	OMC	3	0
2	B	30	T3P	1	0
2	D	26	85Y	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	702/764 (91%)	-1.11	0 100 100	26, 56, 104, 174	0
1	C	702/764 (91%)	-1.13	0 100 100	25, 56, 100, 164	0
2	B	15/30 (50%)	-1.36	0 100 100	38, 46, 54, 54	0
2	D	15/30 (50%)	-1.47	0 100 100	40, 47, 52, 56	0
All	All	1434/1588 (90%)	-1.13	0 100 100	25, 56, 102, 174	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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
2	T3P	D	30	20/21	0.94	0.10	97,127,142,156	0
2	T3P	B	30	20/21	0.95	0.10	113,147,168,172	0
2	OMC	D	29	21/22	0.96	0.06	62,106,122,137	0
2	OMG	B	1	24/25	0.96	0.08	89,106,129,134	0
2	OMG	D	1	24/25	0.96	0.08	69,95,130,140	0
2	OMC	B	2	21/22	0.97	0.05	71,87,95,116	0
2	OMG	B	28	24/25	0.97	0.07	69,90,108,125	0
2	OMC	B	29	21/22	0.97	0.06	60,110,132,142	0
2	OMG	D	28	24/25	0.98	0.05	59,84,111,126	0
2	85Y	D	9	33/34	0.99	0.05	37,51,92,97	0
2	85Y	B	15	33/34	0.99	0.05	32,53,78,85	0
2	85Y	D	15	33/34	0.99	0.05	37,52,75,85	0
2	OMG	B	17	24/25	0.99	0.05	49,71,92,93	0
2	OMG	D	17	24/25	0.99	0.05	47,67,93,94	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	OMG	B	18	24/25	0.99	0.05	33,49,56,60	0
2	OMG	D	18	24/25	0.99	0.05	38,48,62,67	0
2	85Y	B	21	33/34	0.99	0.04	34,46,58,66	0
2	85Y	D	21	33/34	0.99	0.04	29,46,58,61	0
2	85Y	B	23	33/34	0.99	0.05	37,48,58,62	0
2	85Y	D	23	33/34	0.99	0.05	37,47,56,59	0
2	85Y	B	26	33/34	0.99	0.04	27,45,64,89	0
2	85Y	D	26	33/34	0.99	0.04	31,48,66,86	0
2	85Y	B	27	33/34	0.99	0.05	29,47,72,82	0
2	85Y	D	27	33/34	0.99	0.06	36,51,73,82	0
2	OMC	D	2	21/22	0.99	0.04	60,80,94,96	0
2	85Y	B	3	33/34	0.99	0.06	42,64,84,92	0
2	85Y	D	3	33/34	0.99	0.06	42,67,82,89	0
2	OMG	B	4	24/25	0.99	0.04	42,50,63,78	0
2	OMG	D	4	24/25	0.99	0.05	36,48,64,74	0
2	85Y	B	9	33/34	0.99	0.05	36,52,95,106	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.