W O R L D W I D E
wwPDB
P R O T E I N   D A T A   B A N K

wwPDB X-ray Structure Validation Summary Report (i)

Nov 15, 2023 – 02:57 PM JST

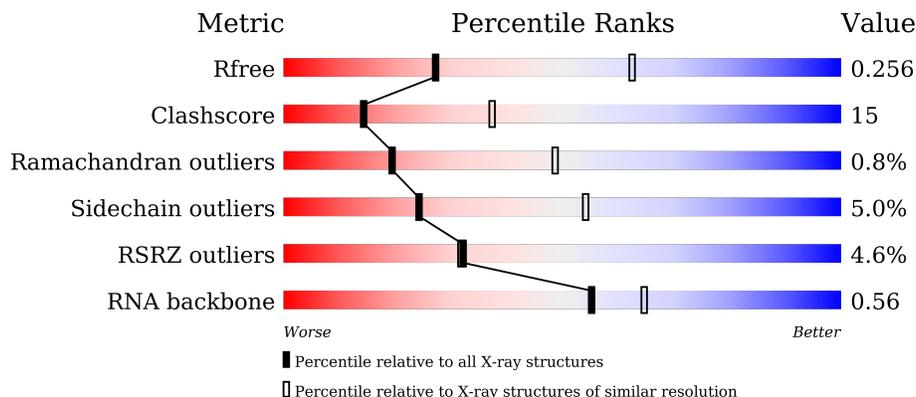|		
PDB ID	:	6JE3
Title	:	Crystal structure of Nme2Cas9 in complex with sgRNA and target DNA (AG-GCCC PAM) with 5 nt overhang
Authors	:	Sun, W.; Yang, J.; Cheng, Z.; Liu, C.; Wang, K.; Huang, X.; Wang, Y.
Deposited on	:	2019-02-03
Resolution	:	2.93 Å(reported)

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 2.93 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	Percentile Ranks	Value
Rfree		0.256
Clashscore		15
Ramachandran outliers		0.8%
Sidechain outliers		5.0%
RSRZ outliers		4.6%
RNA backbone		0.56

Worse ___ *Better*

■ Percentile relative to all X-ray structures
▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2969 (2.98-2.90)
Clashscore	141614	3218 (2.98-2.90)
Ramachandran outliers	138981	3122 (2.98-2.90)
Sidechain outliers	138945	3124 (2.98-2.90)
RSRZ outliers	127900	2902 (2.98-2.90)
RNA backbone	3102	1060 (3.20-2.68)

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 <=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	1083	4% 53% 24% • 21%
2	B	135	4% 32% 43% 11% • 13%
3	C	35	40% 60%
4	D	16	31% 50% 19%

2 Entry composition ⓘ

There are 6 unique types of molecules in this entry. The entry contains 9806 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 CRISPR-associated endonuclease Cas9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	858	Total	C	N	O	S	0	0	0
			6327	3994	1154	1159	20			

- Molecule 2 is a RNA chain called sgRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	117	Total	C	N	O	P	0	0	0
			2472	1106	424	825	117			

- Molecule 3 is a DNA chain called target DNA strand.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	35	Total	C	N	O	P	0	0	0
			720	344	139	203	34			

- Molecule 4 is a DNA chain called non-target DNA strand.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	13	Total	C	N	O	P	0	0	0
			268	127	50	78	13			

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).

EDO

O2

OH

C1

C2

HO

O1
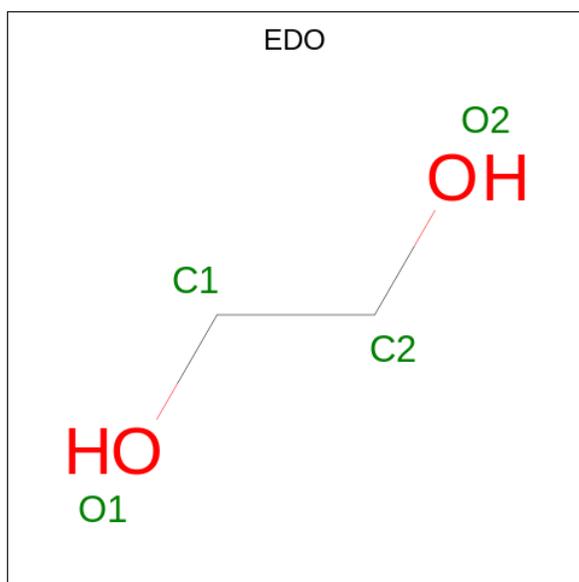
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total 4	C 2	O 2	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	7	Total 7	O 7	0	0
6	B	5	Total 5	O 5	0	0
6	C	3	Total 3	O 3	0	0

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: CRISPR-associated endonuclease Cas9

A906 D911 N912 P913 F914 Y915 K916 A924 V925 R926 V927 L936 K939 N948 M951 V952 R953 V954 D955 V956 F957 C958 K959 V960 D961 K962 K965 N966 Q967 Y968 F969 I970 A975 W976 Q977 V978 N981 I982 L983 P984 D985 I986 D987 Y991 R992 I993 C1000 F1001

H1004 K1005 Y1006 D1007 L1008 I1009 A1010 F1011 Q1012 S1017 E1020 F1021 A1022 C1027 G1032 R1033 F1034 Y1035 L1036 A1037 W1038 H1039 K1044 E1045 F1048 R1049 I1050 Q1053 N1054 L1055 I1058 Q1062 V1063 L1066 G1067 R1071 P1072 C1073 R1078 P1079 P1080 V1081 R1082

● Molecule 2: sgRNA

Chain B:

4%

32%   43%   11% · 13%

G1 G2 U3 C4 U9 G10 U16 A17 A18 C19 U20 U21 C24 G25 U26 U27 G28 U29 A30 G31 C32 U33 C34 C35 C36 U37 U38 V39 C40 U41 C42 G43 G47 A48 A51 C52 C53 U56 A60 C61 A65 G68 C69 C70 G71 U72 C73 U74 G75 A76 A77

A78 A79 G80 A81 U82 G83 U84 G85 C86 C87 G88 C89 A90 A91 C92 G93 C94 U95 C96 U97 G98 C99 C100 C101 C102 U U A A A G C U C C C U G G C U U U A A120 G121 G122 G123 G124 C128 G129 U130 U131 U132 A133 U134 C

● Molecule 3: target DNA strand

Chain C:

40%   60%

T1 A2 T5 G6 A14 A15 A16 G17 T18 T19 A20 A21 A22 T23 A24 G25 C26 A27 G28 G32 A33 C34 C35

● Molecule 4: non-target DNA strand

Chain D:

31%   50%   19%

DA DA DA T4 G5 A6 G7 C10 C11 A12 T15 A16
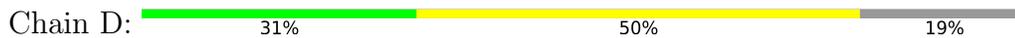
4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	166.39Å 165.97Å 178.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.60 – 2.93 41.60 – 2.93	Depositor EDS
% Data completeness (in resolution range)	75.3 (41.60-2.93) 75.3 (41.60-2.93)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.35 (at 2.95Å)	Xtriage
Refinement program	PHENIX (1.14_3247: ???)	Depositor
R, R_{free}	0.227 , 0.256 0.227 , 0.256	Depositor DCC
R_{free} test set	1906 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å^2)	49.1	Xtriage
Anisotropy	0.548	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.25 , 54.3	EDS
L-test for twinning[2]	$< |L| > = 0.46, < L^2 > = 0.29$	Xtriage
Estimated twinning fraction	0.038 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.89	EDS
Total number of atoms	9806	wwPDB-VP
Average B, all atoms (Å^2)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.10% 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, 0.333 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: EDO

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	0.30	0/6437	0.56	2/8721 (0.0%)
2	B	0.32	0/2757	0.91	4/4287 (0.1%)
3	C	0.61	0/810	0.93	1/1250 (0.1%)
4	D	0.66	0/300	0.97	0/461
All	All	0.35	0/10304	0.73	7/14719 (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
1	A	0	5

There are no bond length outliers.

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	725	LEU	CA-CB-CG	6.63	130.54	115.30
2	B	70	C	C5-C6-N1	6.12	124.06	121.00
1	A	791	PRO	N-CA-CB	5.78	110.23	103.30
2	B	87	C	C2-N1-C1'	5.33	124.66	118.80
3	C	17	DG	O4'-C4'-C3'	-5.30	102.38	104.50

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1010	ALA	Peptide

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Group
1	A	512	SER	Peptide
1	A	814	LEU	Peptide
1	A	820	ALA	Peptide
1	A	821	VAL	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	A	6327	0	5883	213	0
2	B	2472	0	1256	48	0
3	C	720	0	392	17	0
4	D	268	0	147	5	0
5	A	4	0	6	0	0
6	A	7	0	0	0	0
6	B	5	0	0	0	0
6	C	3	0	0	0	0
All	All	9806	0	7684	269	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 15.

The worst 5 of 269 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:687:LEU:HD22	1:A:693:ARG:HA	1.35	1.08
1:A:305:LEU:HD23	1:A:323:LEU:HD13	1.37	1.02
1:A:358:ARG:NH1	1:A:368:LYS:HE3	1.77	0.99
1:A:289:LEU:O	1:A:331:PHE:N	2.01	0.94
1:A:687:LEU:HD12	1:A:687:LEU:O	1.67	0.93

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	
1	A	840/1083 (78%)	758 (90%)	75 (9%)	7 (1%)	19	49

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	816	SER
1	A	327	GLU
1	A	700	GLY
1	A	815	SER
1	A	821	VAL

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	576/935 (62%)	547 (95%)	29 (5%)	24	54

5 of 29 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	725	LEU
1	A	1049	ARG
1	A	812	GLU
1	A	992	ARG
1	A	797	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	722	HIS
1	A	966	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	115/135 (85%)	29 (25%)	6 (5%)

5 of 29 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	24	C
2	B	43	G
2	B	47	G
2	B	48	A
2	B	65	A

5 of 6 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	82	U
2	B	93	G
2	B	97	U
2	B	73	C
2	B	68	G

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

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

1 ligand is 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	$\#|Z| > 2$
5	EDO	A	1101	-	3,3,3	0.47	0	2,2,2	0.35	0

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
5	EDO	A	1101	-	-	0/1/1/1	-

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 ⓘ

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, 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	858/1083 (79%)	0.11	42 (4%)	29	29	12, 70, 136, 166	0
2	B	117/135 (86%)	0.03	5 (4%)	35	34	12, 53, 161, 207	0
3	C	35/35 (100%)	-0.55	0	100	100	17, 39, 75, 79	0
4	D	13/16 (81%)	-0.36	0	100	100	32, 43, 74, 110	0
All	All	1023/1269 (80%)	0.08	47 (4%)	32	32	12, 65, 141, 207	0

The worst 5 of 47 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	730	VAL	6.0
1	A	731	ALA	5.8
1	A	704	ASN	5.6
1	A	38	LEU	4.5
1	A	691	GLY	4.4

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

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

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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
5	EDO	A	1101	4/4	0.95	0.12	33,40,54,56	0

6.5 Other polymers ⓘ

There are no such residues in this entry.