



Full wwPDB X-ray Structure Validation Report

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PDB ID : 2F0Q
Title : Crystal structure of Staphylococcal nuclease mutant V66L/I92L
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Deposited on : 2005-11-13
Resolution : 1.95 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

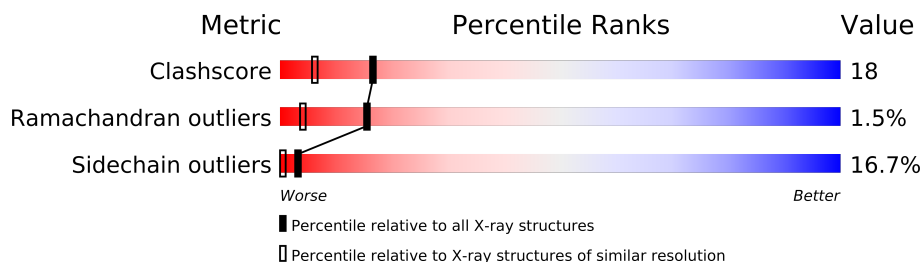
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.95 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	79885	1488 (1.96-1.96)
Ramachandran outliers	78287	1475 (1.96-1.96)
Sidechain outliers	78261	1475 (1.96-1.96)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	149	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1138 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Staphylococcal nuclease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	135	Total	C	N	O	S	0	1	0
			1089	695	193	197	4			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	66	LEU	VAL	ENGINEERED	UNP P00644
A	92	LEU	ILE	ENGINEERED	UNP P00644
A	124	HIS	LEU	VARIANT	UNP P00644

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	49	Total	O	0	0
			49	49		

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	48.15Å 48.14Å 63.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 1.95	Depositor
% Data completeness (in resolution range)	92.7 (8.00-1.95)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.194 , 0.270	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1138	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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	0.36	0/1115	1.07	1/1494 (0.1%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	35	ARG	NE-CZ-NH1	9.18	124.89	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1089	0	1129	40	0
2	A	49	0	0	2	0
All	All	1138	0	1129	40	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 18.

All (40) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:46:HIS:HB3	1:A:49:LYS:HB2	1.48	0.94

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:133:LYS:O	1:A:136:LYS:HD3	1.83	0.78
1:A:60:ALA:O	1:A:64:LYS:HD3	1.92	0.69
1:A:28:LYS:H	1:A:28:LYS:NZ	1.89	0.69
1:A:133:LYS:HD2	1:A:140:TRP:CZ2	2.33	0.62
1:A:46:HIS:CE1	1:A:48:LYS:HE2	2.36	0.60
1:A:74:VAL:HG12	1:A:92:LEU:CD1	2.33	0.58
1:A:43:GLU:O	1:A:52:GLU:HG3	2.04	0.58
1:A:46:HIS:CE1	1:A:48:LYS:HG3	2.40	0.57
1:A:133:LYS:HD2	1:A:140:TRP:CE2	2.39	0.56
1:A:46:HIS:HE1	1:A:48:LYS:HG3	1.71	0.55
1:A:38:LEU:HD23	2:A:219:HOH:O	2.07	0.55
1:A:60:ALA:HB1	1:A:64:LYS:NZ	2.22	0.54
1:A:75:GLU:HB3	1:A:91:TYR:HB2	1.91	0.53
1:A:74:VAL:HG12	1:A:92:LEU:HD11	1.90	0.52
1:A:53:LYS:O	1:A:54:TYR:HB2	2.10	0.51
1:A:46:HIS:CB	1:A:49:LYS:HD2	2.42	0.49
1:A:48:LYS:O	1:A:49:LYS:HG3	2.12	0.49
1:A:46:HIS:HB2	1:A:49:LYS:HD2	1.95	0.48
1:A:41:THR:HG22	1:A:109:ALA:HB2	1.96	0.48
1:A:116:LYS:HB3	1:A:116:LYS:HE3	1.31	0.48
1:A:107:GLY:O	1:A:139:ILE:HB	2.15	0.47
1:A:46:HIS:ND1	1:A:47:PRO:HD2	2.29	0.47
1:A:28:LYS:H	1:A:28:LYS:HZ1	1.58	0.46
1:A:80:GLN:HG2	2:A:232:HOH:O	2.16	0.45
1:A:53:LYS:HD3	1:A:54:TYR:CD2	2.52	0.44
1:A:53:LYS:HD3	1:A:54:TYR:CE2	2.51	0.44
1:A:127:LYS:HD2	1:A:127:LYS:HA	1.40	0.44
1:A:77:ASP:OD1	1:A:78:LYS:N	2.51	0.44
1:A:133:LYS:HG3	1:A:140:TRP:CD2	2.53	0.44
1:A:82:THR:HA	1:A:87:ARG:O	2.18	0.44
1:A:123:GLN:NE2	1:A:123:GLN:HA	2.33	0.43
1:A:46:HIS:ND1	1:A:48:LYS:O	2.52	0.42
1:A:23:VAL:HG21	1:A:92:LEU:HD21	2.02	0.42
1:A:92:LEU:HA	1:A:92:LEU:HD12	1.85	0.42
1:A:116:LYS:HA	1:A:117:PRO:HA	1.87	0.42
1:A:74:VAL:HG12	1:A:92:LEU:HD12	2.01	0.42
1:A:101:GLU:HG3	1:A:124:HIS:CD2	2.55	0.41
1:A:133:LYS:HD2	1:A:140:TRP:CH2	2.56	0.41
1:A:46:HIS:HA	1:A:47:PRO:HD3	1.87	0.40

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	134/149 (90%)	126 (94%)	6 (4%)	2 (2%)	15 4

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	49	LYS
1	A	138	ASN

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	115/125 (92%)	96 (84%)	19 (16%)	3 0

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	HIS
1	A	19	ASP
1	A	24	LYS
1	A	28	LYS
1	A	40	ASP
1	A	43	GLU
1	A	48	LYS
1	A	52	GLU
1	A	53	LYS
1	A	59	SER
1	A	84	LYS
1	A	92	LEU

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Mol	Chain	Res	Type
1	A	97	LYS
1	A	100	ASN
1	A	116	LYS
1	A	127	LYS
1	A	133	LYS
1	A	139	ILE
1	A	141	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	GLN
1	A	123	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.