



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 27, 2014 – 02:14 AM EDT

PDB ID : 3NQZ
Title : Crystal structure of the autoprocessed Vibriolysin MCP-02 with E369A mutation
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Deposited on : 2010-06-30
Resolution : 2.05 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>

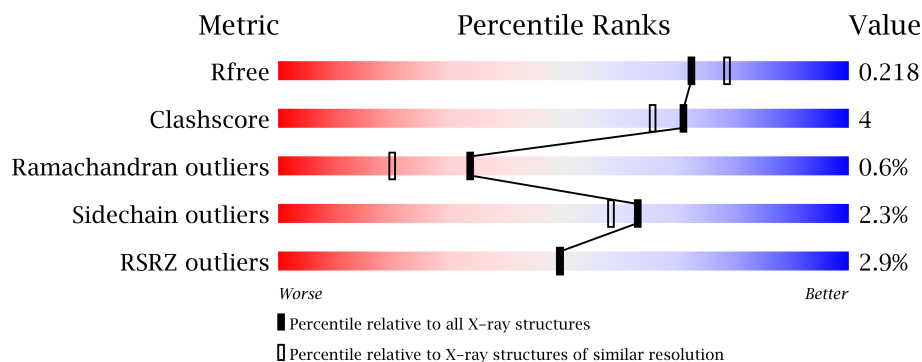
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1439
EDS : stable22978
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22978

1 Overall quality at a glance

The reported resolution of this entry is 2.05 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}	66092	1380 (2.06-2.02)
Clashscore	79885	1577 (2.06-2.02)
Ramachandran outliers	78287	1565 (2.06-2.02)
Sidechain outliers	78261	1565 (2.06-2.02)
RSRZ outliers	66119	1381 (2.06-2.02)

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.

Mol	Chain	Length	Quality of chain
1	A	180	
2	B	315	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4062 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 (with D amino acids) called Secreted metalloprotease Mcp02.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	175	Total	C	N	O	S	0	0	0
			1352	848	237	265	2			

- Molecule 2 is a protein (with D amino acids) called Secreted metalloprotease Mcp02.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	307	Total	C	N	O	S	0	0	0
			2326	1452	391	469	14			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	369	ALA	GLU	ENGINEERED MUTATION	UNP A1DRD5

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Ca	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	123	Total	O	0	0
			123	123		

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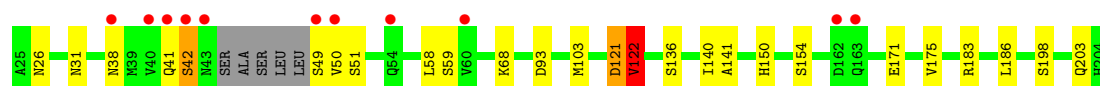
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	259	Total 259	O 259	0	0

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 ($RSRZ > 2$). 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: Secreted metalloprotease Mcp02

Chain A: 



- Molecule 2: Secreted metalloprotease Mcp02

Chain B: 



THR
ASP
ASP

4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	82.53Å 82.53Å 154.21Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	16.97 – 2.05 16.97 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.9 (16.97-2.05) 99.9 (16.97-2.05)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.40 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.179 , 0.220 0.182 , 0.218	Depositor DCC
R_{free} test set	1941 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	18.7	Xtriage
Anisotropy	0.040	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 36.6	EDS
Estimated twinning fraction	0.025 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 38780 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4062	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, CA

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	0.78	1/1372 (0.1%)	0.84	5/1856 (0.3%)
2	B	0.95	5/2385 (0.2%)	0.78	7/3241 (0.2%)
All	All	0.89	6/3757 (0.2%)	0.80	12/5097 (0.2%)

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	1
2	B	0	3
All	All	0	4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	237	CYS	CB-SG	-6.85	1.70	1.82
2	B	346	GLU	CD-OE1	-6.04	1.19	1.25
2	B	345	HIS	C-O	-5.69	1.12	1.23
2	B	305	LEU	C-N	5.52	1.46	1.34
2	B	384	LYS	CE-NZ	-5.28	1.35	1.49
1	A	58	LEU	C-N	-5.18	1.22	1.34

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	121	ASP	O-C-N	9.09	137.24	122.70
2	B	266	ASN	N-CA-C	7.35	130.85	111.00
1	A	59	SER	O-C-N	7.27	134.34	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	341	ASP	CB-CG-OD1	6.95	124.56	118.30
2	B	234	GLY	N-CA-C	-6.87	95.92	113.10
1	A	121	ASP	C-N-CA	6.59	138.17	121.70
2	B	305	LEU	C-N-CA	-5.87	107.03	121.70
1	A	122	VAL	N-CA-C	5.56	126.02	111.00
2	B	306	GLN	CA-C-N	-5.49	105.11	117.20
1	A	59	SER	CA-C-N	-5.37	105.39	117.20
2	B	267	THR	N-CA-CB	-5.36	100.12	110.30
2	B	471	ASP	CB-CG-OD1	5.08	122.87	118.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	121	ASP	Peptide
2	B	233	SER	Peptide
2	B	265	GLU	Peptide
2	B	306	GLN	Mainchain

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	1352	0	1352	19	0
2	B	2326	0	2119	17	0
3	B	1	0	0	0	0
4	B	1	0	0	0	0
5	A	123	0	0	5	0
5	B	259	0	0	1	0
All	All	4062	0	3471	32	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 4.

All (32) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:103:MET:HG2	5:A:471:HOH:O	1.54	1.06
1:A:41:GLN:O	1:A:42:SER:HB2	1.67	0.92
2:B:345:HIS:O	2:B:346:GLU:HB2	1.82	0.77
1:A:122:VAL:CG1	1:A:122:VAL:O	2.35	0.75
1:A:136:SER:HB2	5:A:226:HOH:O	1.88	0.73
1:A:68:LYS:HE2	5:A:251:HOH:O	1.89	0.72
2:B:345:HIS:O	2:B:346:GLU:CB	2.33	0.70
1:A:122:VAL:HG13	1:A:122:VAL:O	1.96	0.66
2:B:388:ASP:OD2	2:B:393:LYS:HE2	2.02	0.59
1:A:50:VAL:HG22	5:A:206:HOH:O	2.03	0.57
1:A:31:ASN:HD22	2:B:475:ASN:HD22	1.52	0.57
2:B:342:VAL:O	2:B:345:HIS:O	2.23	0.57
2:B:310:HIS:HD2	2:B:329:ASP:OD1	1.88	0.55
1:A:41:GLN:O	1:A:42:SER:CB	2.46	0.54
1:A:50:VAL:HG22	1:A:51:SER:H	1.73	0.53
1:A:154:SER:HA	1:A:171:GLU:O	2.12	0.50
2:B:388:ASP:OD2	2:B:393:LYS:NZ	2.44	0.50
2:B:388:ASP:OD2	2:B:393:LYS:CE	2.62	0.48
1:A:31:ASN:ND2	2:B:475:ASN:HD22	2.11	0.47
1:A:175:VAL:HG22	1:A:183:ARG:HG3	1.96	0.46
2:B:301:LEU:HD12	2:B:305:LEU:HD23	1.98	0.45
1:A:150:HIS:HD2	5:A:444:HOH:O	2.00	0.44
1:A:203:GLN:HG2	2:B:360:TYR:CG	2.53	0.44
1:A:38:ASN:O	1:A:42:SER:CB	2.66	0.44
2:B:240:ASN:ND2	2:B:245:ARG:HE	2.15	0.44
1:A:203:GLN:HG2	2:B:360:TYR:CD1	2.55	0.42
2:B:369:ALA:HB1	2:B:428:HIS:CD2	2.54	0.42
1:A:141:ALA:HB2	1:A:186:LEU:HD21	2.02	0.41
2:B:345:HIS:CG	2:B:372:SER:HB3	2.55	0.41
2:B:245:ARG:HB3	2:B:306:GLN:HG2	2.03	0.41
1:A:122:VAL:HG12	1:A:122:VAL:O	2.19	0.41
2:B:420:ASN:ND2	5:B:146:HOH:O	2.50	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

There are no protein chains in this entry.

5.3.2 Protein sidechains ⓘ

There are no protein chains in this entry.

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 ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - 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.

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	175/180 (97%)	0.07	11 (6%)	19 18	12, 24, 44, 65	0
2	B	307/315 (97%)	-0.65	3 (0%)	79 81	8, 14, 24, 38	0
All	All	482/495 (97%)	-0.39	14 (2%)	49 49	8, 16, 40, 65	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	50	VAL	6.1
1	A	42	SER	4.2
2	B	235	ASN	4.1
1	A	41	GLN	3.3
1	A	38	ASN	3.2
1	A	40	VAL	2.8
2	B	234	GLY	2.5
1	A	43	ASN	2.4
1	A	49	SER	2.4
1	A	60	VAL	2.3
1	A	54	GLN	2.3
1	A	163	GLN	2.2
1	A	162	ASP	2.2
2	B	256	SER	2.0

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 carbohydrates 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. 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	ZN	B	1	1/1	0.04	-2.06	16,16,16,16	0
4	CA	B	520	1/1	0.03	-2.99	19,19,19,19	0

6.5 Other polymers

There are no such residues in this entry.