



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 07:18 PM GMT

PDB ID : 1EZL  
Title : CRYSTAL STRUCTURE OF THE DISULPHIDE BOND-DEFICIENT  
AZURIN MUTANT C3A/C26A: HOW IMPORTANT IS THE S-S BOND  
FOR FOLDING AND STABILITY?  
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Deposited on : 2000-05-11  
Resolution : 2.00 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

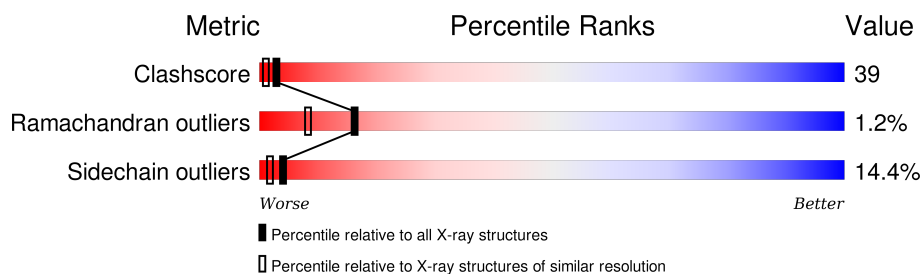
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	128	
1	B	128	
1	C	128	
1	D	128	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4267 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 AZURIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	128	Total	C	N	O	S	0	0	0
			972	607	164	194	7			
1	B	128	Total	C	N	O	S	0	0	0
			972	607	164	194	7			
1	C	128	Total	C	N	O	S	0	0	0
			972	607	164	194	7			
1	D	128	Total	C	N	O	S	0	0	0
			972	607	164	194	7			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	3	ALA	CYS	ENGINEERED	UNP P00282
A	26	ALA	CYS	ENGINEERED	UNP P00282
B	133	ALA	CYS	ENGINEERED	UNP P00282
B	156	ALA	CYS	ENGINEERED	UNP P00282
C	263	ALA	CYS	ENGINEERED	UNP P00282
C	286	ALA	CYS	ENGINEERED	UNP P00282
D	393	ALA	CYS	ENGINEERED	UNP P00282
D	416	ALA	CYS	ENGINEERED	UNP P00282

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cu	0	0
			1	1		
2	A	1	Total	Cu	0	0
			1	1		
2	D	1	Total	Cu	0	0
			1	1		
2	C	1	Total	Cu	0	0
			1	1		

- Molecule 3 is water.

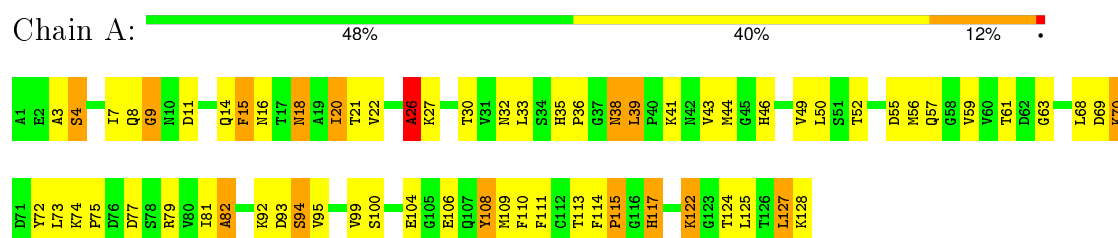
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	93	Total 93	O 93	0	0
3	B	97	Total 97	O 97	0	0
3	C	97	Total 97	O 97	0	0
3	D	88	Total 88	O 88	0	0

### 3 Residue-property plots

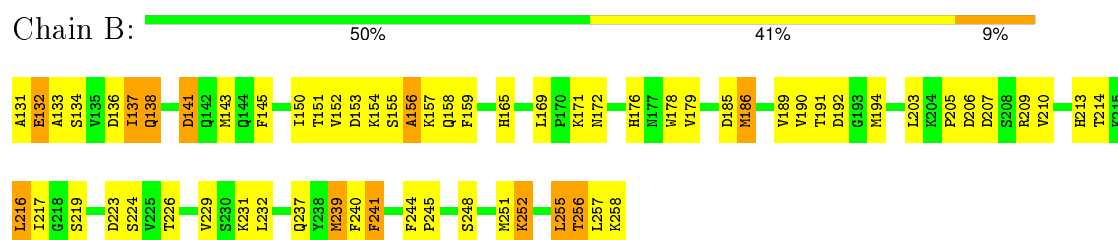
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ( $\text{RSRZ} > 2$ ). 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.

Note EDS was not executed.

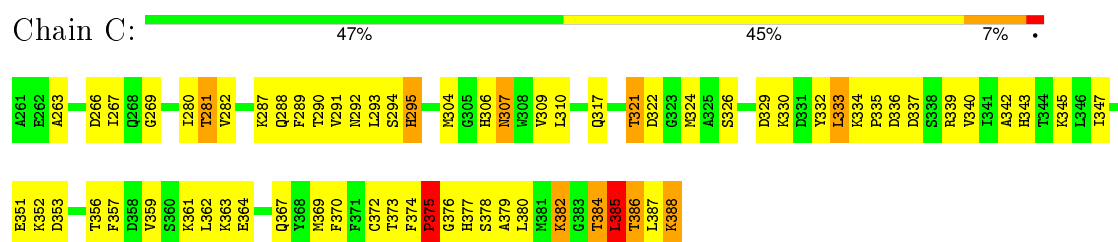
#### • Molecule 1: AZURIN



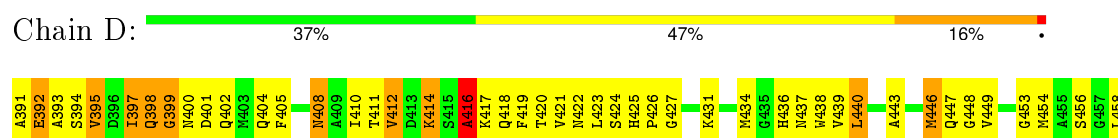
#### • Molecule 1: AZURIN



#### • Molecule 1: AZURIN



#### • Molecule 1: AZURIN



D459	K460	D461	Y462	L463	K464	D467	S468	K469	V470	I471	A472	H473	L476	S479	K482	T486	F487	D488	V489	S490	K491	L492	K493	E494	G495	E496	Q497	Y498	N499	F500	F501	G502	P505	G506	M511	K512	L515	T516	L517	K518
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## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	57.40 Å 80.40 Å 110.30 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.00)	Depositor
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.189 , 0.256	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4267	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

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

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	1.07	1/989 (0.1%)	1.13	4/1333 (0.3%)
1	B	1.06	2/989 (0.2%)	1.18	3/1333 (0.2%)
1	C	1.05	1/989 (0.1%)	1.23	4/1333 (0.3%)
1	D	1.22	3/989 (0.3%)	1.24	5/1333 (0.4%)
All	All	1.10	7/3956 (0.2%)	1.20	16/5332 (0.3%)

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	C	0	1
1	D	0	1
All	All	0	4

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	416	ALA	CA-CB	15.89	1.85	1.52
1	A	26	ALA	CA-CB	9.29	1.72	1.52
1	B	239	MET	CG-SD	-6.10	1.65	1.81
1	C	263	ALA	CA-CB	5.42	1.63	1.52
1	D	438	TRP	CB-CG	5.35	1.59	1.50

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	385	LEU	N-CA-C	-7.84	89.84	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	333	LEU	CA-CB-CG	6.80	130.93	115.30
1	D	416	ALA	CB-CA-C	6.78	120.27	110.10
1	D	399	GLY	N-CA-C	-6.66	96.44	113.10
1	A	9	GLY	N-CA-C	-6.27	97.43	113.10

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	108	TYR	Sidechain
1	A	26	ALA	Mainchain
1	C	332	TYR	Sidechain
1	D	416	ALA	Mainchain

## 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	972	0	954	61	0
1	B	972	0	951	58	0
1	C	972	0	951	72	0
1	D	972	0	951	120	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	93	0	0	8	0
3	B	97	0	0	6	0
3	C	97	0	0	14	0
3	D	88	0	0	13	0
All	All	4267	0	3807	303	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 39.

The worst 5 of 303 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:D:416:ALA:CB	1:D:416:ALA:CA	1.85	1.49
1:D:434:MET:SD	1:D:434:MET:CE	2.01	1.48
1:D:391:ALA:HB1	1:D:410:ILE:HG21	1.34	1.09
1:D:391:ALA:N	1:D:419:PHE:HB3	1.74	1.01
1:D:412:VAL:HG13	1:D:419:PHE:CD2	1.96	1.00

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	126/128 (98%)	118 (94%)	6 (5%)	2 (2%)	12	5
1	B	126/128 (98%)	112 (89%)	13 (10%)	1 (1%)	24	15
1	C	126/128 (98%)	112 (89%)	12 (10%)	2 (2%)	12	5
1	D	126/128 (98%)	111 (88%)	14 (11%)	1 (1%)	24	15
All	All	504/512 (98%)	453 (90%)	45 (9%)	6 (1%)	16	8

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	132	GLU
1	C	386	THR
1	A	82	ALA
1	D	414	LYS
1	C	375	PRO

### 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	108/108 (100%)	89 (82%)	19 (18%)	2	1
1	B	108/108 (100%)	93 (86%)	15 (14%)	4	2
1	C	108/108 (100%)	97 (90%)	11 (10%)	9	5
1	D	108/108 (100%)	91 (84%)	17 (16%)	3	1
All	All	432/432 (100%)	370 (86%)	62 (14%)	4	2

5 of 62 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	231	LYS
1	C	281	THR
1	D	496	GLU
1	B	248	SER
1	C	295	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	272	GLN
1	D	497	GLN
1	D	418	GLN
1	B	172	ASN
1	C	307	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules 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 4 ligands modelled in this entry, 4 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.

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 ⓘ

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.