



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 14, 2026 – 05:48 PM UTC

PDB ID : 5NGD / pdb_00005ngd
Title : Crystal structure of the V325G mutant of the bacteriophage G20C portal protein
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Deposited on : 2017-03-17
Resolution : 1.90 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
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

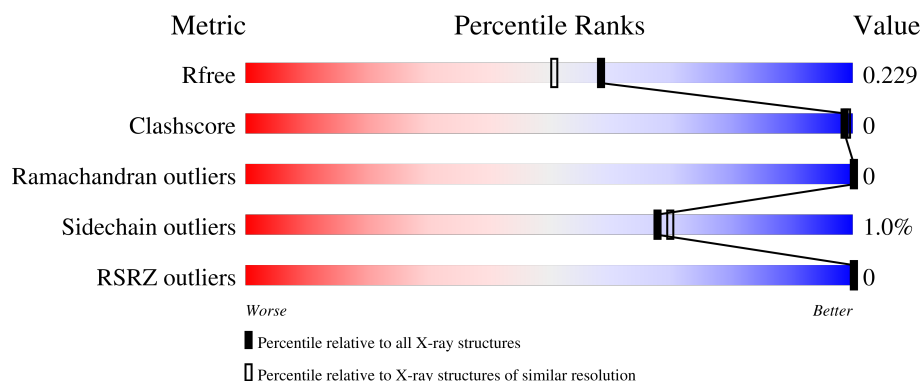
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

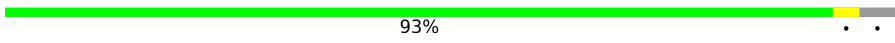
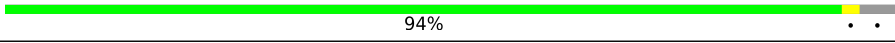
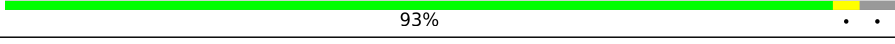
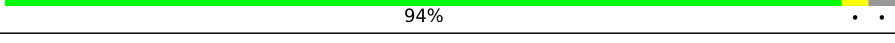
The reported resolution of this entry is 1.90 Å.

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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	417	 93%
1	B	417	 94%
1	C	417	 93%
1	D	417	 94%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13061 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 Portal protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	402	Total	C	N	O	S	0	2	0
			3081	1992	511	568	10			
1	B	401	Total	C	N	O	S	0	1	0
			3072	1985	511	566	10			
1	C	401	Total	C	N	O	S	0	2	0
			3093	1999	511	573	10			
1	D	403	Total	C	N	O	S	0	2	0
			3112	2013	516	573	10			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	GLY	-	expression tag	UNP A7XXR3
A	23	PRO	-	expression tag	UNP A7XXR3
A	24	ALA	-	expression tag	UNP A7XXR3
A	325	GLY	VAL	engineered mutation	UNP A7XXR3
B	22	GLY	-	expression tag	UNP A7XXR3
B	23	PRO	-	expression tag	UNP A7XXR3
B	24	ALA	-	expression tag	UNP A7XXR3
B	325	GLY	VAL	engineered mutation	UNP A7XXR3
C	22	GLY	-	expression tag	UNP A7XXR3
C	23	PRO	-	expression tag	UNP A7XXR3
C	24	ALA	-	expression tag	UNP A7XXR3
C	325	GLY	VAL	engineered mutation	UNP A7XXR3
D	22	GLY	-	expression tag	UNP A7XXR3
D	23	PRO	-	expression tag	UNP A7XXR3
D	24	ALA	-	expression tag	UNP A7XXR3
D	325	GLY	VAL	engineered mutation	UNP A7XXR3

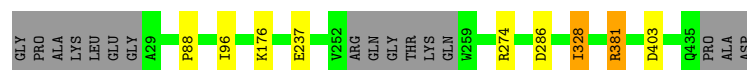
- Molecule 2 is water.

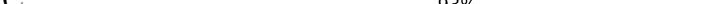
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	159	Total 159	O 159	0	0
2	B	172	Total 172	O 172	0	0
2	C	184	Total 184	O 184	0	0
2	D	188	Total 188	O 188	0	0

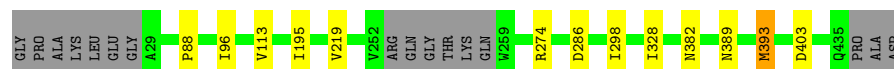
- Molecule 1: Portal protein

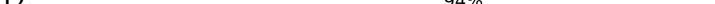


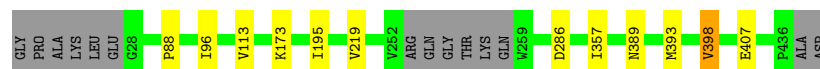
Chain B: 94% . .



Chain C:  93% . .



Chain D:  94%



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	223.23Å 223.23Å 115.82Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	111.62 – 1.90 111.62 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.8 (111.62-1.90) 97.8 (111.62-1.90)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.206 , 0.225 0.213 , 0.229	Depositor DCC
R_{free} test set	1623 reflections (0.93%)	wwPDB-VP
Wilson B-factor (Å ²)	22.1	Xtriage
Anisotropy	0.242	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 28.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.174 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13061	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.98% 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](#)

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	1.11	2/3157 (0.1%)	1.03	5/4301 (0.1%)
1	B	1.12	0/3144	1.03	3/4284 (0.1%)
1	C	1.12	1/3168 (0.0%)	1.03	4/4312 (0.1%)
1	D	1.13	2/3188 (0.1%)	1.03	2/4336 (0.0%)
All	All	1.12	5/12657 (0.0%)	1.03	14/17233 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	219	VAL	CA-C	7.12	1.59	1.52
1	A	219	VAL	CA-C	6.54	1.58	1.52
1	C	219	VAL	CA-C	5.48	1.57	1.52
1	A	210	THR	CA-C	-5.14	1.46	1.52
1	D	357	ILE	CA-C	5.09	1.57	1.52

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	393	MET	CG-SD-CE	6.46	115.11	100.90
1	B	381	ARG	CG-CD-NE	-5.48	99.94	112.00
1	A	403	ASP	CA-C-N	5.32	123.59	120.24
1	A	403	ASP	C-N-CA	5.32	123.59	120.24
1	C	403	ASP	CA-C-N	5.16	124.46	120.33

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	3081	0	3056	4	1
1	B	3072	0	3042	5	1
1	C	3093	0	3081	5	0
1	D	3112	0	3120	2	0
2	A	159	0	0	2	0
2	B	172	0	0	0	0
2	C	184	0	0	0	0
2	D	188	0	0	0	0
All	All	13061	0	12299	12	1

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 0.

The worst 5 of 12 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:B:381:ARG:HH11	1:C:382:ASN:HD22	1.51	0.59
1:B:328:ILE:HD13	1:C:328:ILE:HG12	1.89	0.53
1:A:436:PRO:O	2:A:501:HOH:O	2.18	0.52
1:B:88:PRO:HG3	1:B:96:ILE:HD13	1.96	0.48
1:B:237:GLU:HG3	1:C:274:ARG:CZ	2.44	0.47

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:ALA:O	1:B:176:LYS:NZ[8_664]	2.18	0.02

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	400/417 (96%)	392 (98%)	8 (2%)	0	100	100
1	B	398/417 (95%)	390 (98%)	8 (2%)	0	100	100
1	C	399/417 (96%)	392 (98%)	7 (2%)	0	100	100
1	D	401/417 (96%)	393 (98%)	8 (2%)	0	100	100
All	All	1598/1668 (96%)	1567 (98%)	31 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	321/344 (93%)	318 (99%)	3 (1%)	70	73
1	B	319/344 (93%)	317 (99%)	2 (1%)	78	81
1	C	325/344 (94%)	322 (99%)	3 (1%)	70	73
1	D	328/344 (95%)	323 (98%)	5 (2%)	57	56
All	All	1293/1376 (94%)	1280 (99%)	13 (1%)	68	70

5 of 13 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	298	ILE
1	D	173	LYS
1	D	407	GLU
1	D	286	ASP
1	D	398	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	275	HIS
1	D	60	HIS
1	D	396	ASN

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Mol	Chain	Res	Type
1	D	275	HIS
1	B	60	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	402/417 (96%)	-1.38	0 100 100	13, 28, 53, 76	2 (0%)
1	B	401/417 (96%)	-1.39	0 100 100	12, 27, 50, 74	1 (0%)
1	C	401/417 (96%)	-1.42	0 100 100	13, 26, 45, 66	2 (0%)
1	D	403/417 (96%)	-1.44	0 100 100	12, 26, 43, 63	2 (0%)
All	All	1607/1668 (96%)	-1.41	0 100 100	12, 27, 48, 76	7 (0%)

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

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.