

wwPDB X-ray Structure Validation Summary Report (i)

Oct 10, 2021 – 01:29 PM EDT

PDB ID : 3GDU

Title: Crystal structure of DegS H198P/D320A mutant modified by DFP and in

complex with YRF peptide

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Deposited on : 2009-02-24

Resolution : 3.00 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
A user guide is available at

https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity : 4.02b-467

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.23.2

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Refmac : 5.8.0158

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

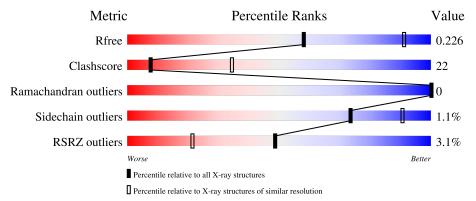
Validation Pipeline (wwPDB-VP) : 2.23.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries},{\rm resolution\ range}({\rm \AA})) \end{array}$
R_{free}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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 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	Λ	340	2%					
1	A	340	47%	37%	• 13%			
1	В	340	55%	27%	18%			
1	С	340	61%	26%	14%			
2	D	3	67%	33	%			
			67%					
2	Е	3	67%	339				

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Mol	Chain	Length	Quality of chain
2	F	3	100%



2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 6530 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 DegS protease.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace		
1	Λ	295	Total	C N C	О	Р	S	0	0	0	
1	A	290	2171	1355	378	432	1	5	U	U	U
1	D	279	Total	С	N	О	Р	S	0	0	0
1	Б	219	2069	1293	361	410	1	4	U	U	0
1	C	293	Total	С	N	О	Р	S	0	0	0
1		∠95	2185	1366	388	425	1	5	U		U

There are 39 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	MET	-	expression tag	UNP P0AEE3
A	17	ARG	-	expression tag	UNP P0AEE3
A	18	GLY	-	expression tag	UNP P0AEE3
A	19	SER	-	expression tag	UNP P0AEE3
A	20	HIS	-	expression tag	UNP P0AEE3
A	21	HIS	-	expression tag	UNP P0AEE3
A	22	HIS	-	expression tag	UNP P0AEE3
A	23	HIS	-	expression tag	UNP P0AEE3
A	24	HIS	-	expression tag	UNP P0AEE3
A	25	HIS	-	expression tag	UNP P0AEE3
A	26	GLY	-	expression tag	UNP P0AEE3
A	198	PRO	HIS	engineered mutation	UNP P0AEE3
A	320	ALA	ASP	engineered mutation	UNP P0AEE3
В	16	MET	-	expression tag	UNP P0AEE3
В	17	ARG	-	expression tag	UNP P0AEE3
В	18	GLY	-	expression tag	UNP P0AEE3
В	19	SER	-	expression tag	UNP P0AEE3
В	20	HIS	-	expression tag	UNP P0AEE3
В	21	HIS	-	expression tag	UNP P0AEE3
В	22	HIS	-	expression tag	UNP P0AEE3
В	23	HIS	-	expression tag	UNP P0AEE3
В	24	HIS	-	expression tag	UNP P0AEE3
В	25	HIS		expression tag	UNP P0AEE3

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Chain	Residue	Modelled	Actual	Comment	Reference
В	26	GLY	-	expression tag	UNP P0AEE3
В	198	PRO	HIS engineered mutation		UNP P0AEE3
В	320	ALA	ASP	engineered mutation	UNP P0AEE3
С	16	MET	-	expression tag	UNP P0AEE3
С	17	ARG	-	expression tag	UNP P0AEE3
С	18	GLY	-	expression tag	UNP P0AEE3
С	19	SER	-	expression tag	UNP P0AEE3
С	20	HIS	-	expression tag	UNP P0AEE3
С	21	HIS	-	expression tag	UNP P0AEE3
С	22	HIS	-	expression tag	UNP P0AEE3
С	23	HIS	-	expression tag	UNP P0AEE3
С	24	HIS	-	expression tag	UNP P0AEE3
С	25	HIS	-	expression tag	UNP P0AEE3
С	26	GLY	-	expression tag	UNP P0AEE3
С	198	PRO	HIS	engineered mutation	UNP P0AEE3
С	320	ALA	ASP	engineered mutation	UNP P0AEE3

• Molecule 2 is a protein called YRF peptide.

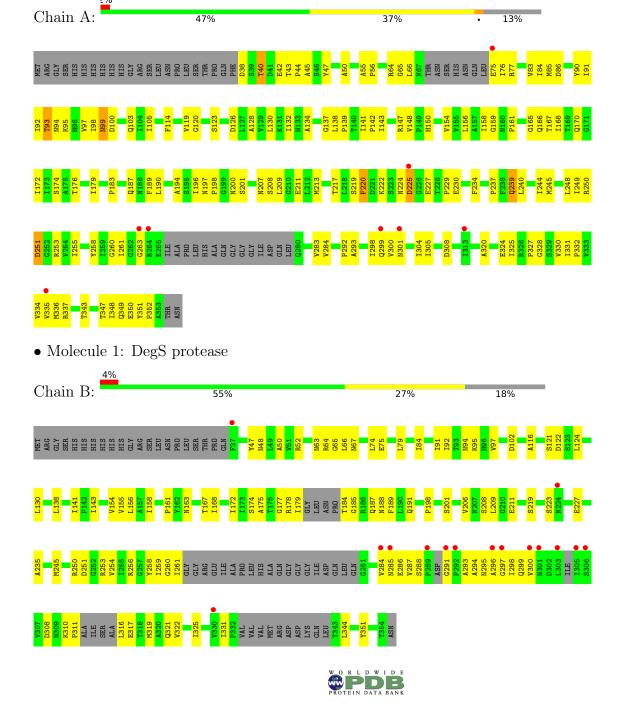
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
2	D	3	Total C N O 35 24 6 5	0	0	0
2	E	3	Total C N O 35 24 6 5	0	0	0
2	F	3	Total C N O 35 24 6 5	0	0	0

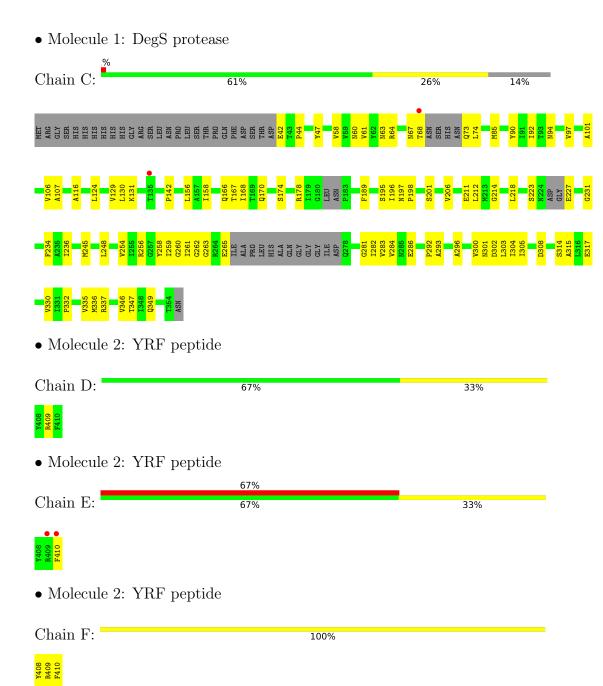


3 Residue-property plots (i)

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: DegS protease







4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants	117.57Å 171.28Å 111.69Å	Donogitor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.15 - 3.00	Depositor
Resolution (A)	36.15 - 2.93	EDS
% Data completeness	92.9 (36.15-3.00)	Depositor
(in resolution range)	92.9 (36.15-2.93)	EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	2.51 (at 2.95Å)	Xtriage
Refinement program	PHENIX	Depositor
D D.	0.209 , 0.231	Depositor
R, R_{free}	0.203 , 0.226	DCC
R_{free} test set	1117 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	75.8	Xtriage
Anisotropy	0.414	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.27, 56.3	EDS
L-test for twinning ²	$ < L >=0.48, < L^2>=0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6530	wwPDB-VP
Average B, all atoms (Å ²)	113.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.04% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities estimated from amplitudes.

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: MIS

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	
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	A	0.22	0/2183	0.43	0/2974
1	В	0.23	0/2079	0.43	0/2828
1	С	0.21	0/2195	0.41	0/2980
2	D	0.27	0/36	0.26	0/45
2	Е	0.24	0/36	0.25	0/45
2	F	0.24	0/36	0.27	0/45
All	All	0.22	0/6565	0.42	0/8917

There are no bond length outliers.

There are no bond angle outliers.

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	2171	0	2180	132	0
1	В	2069	0	2075	94	0
1	С	2185	0	2233	76	0
2	D	35	0	30	1	0
2	Ε	35	0	30	2	0
2	F	35	0	30	3	0
All	All	6530	0	6578	287	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 22.

The worst 5 of 287 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} \text{Interatomic} \\ \text{distance (Å)} \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
1:A:305:ILE:HD11	1:A:336:MET:HG2	1.40	1.04
1:A:292:PRO:HG3	1:A:349:GLN:HG2	1.42	0.96
1:B:64:ARG:HH21	1:B:74:LEU:HD13	1.31	0.94
1:A:179:ILE:HD12	1:A:183:PRO:HA	1.57	0.85
1:A:224:ASN:O	1:A:227:GLU:HG2	1.78	0.84

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	Favoured Allowed		Perce	ntiles
1	A	288/340~(85%)	274 (95%)	14 (5%)	0	100	100
1	В	$266/340 \ (78\%)$	253 (95%)	13 (5%)	0	100	100
1	С	282/340 (83%)	277 (98%)	5 (2%)	0	100	100
2	D	1/3 (33%)	0	1 (100%)	0	100	100
2	E	1/3 (33%)	0	1 (100%)	0	100	100
2	F	1/3 (33%)	1 (100%)	0	0	100	100
All	All	839/1029 (82%)	805 (96%)	34 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	232/275~(84%)	225 (97%)	7 (3%)	41	75	
1	В	223/275~(81%)	223 (100%)	0	100	100	
1	C	235/275~(86%)	234 (100%)	1 (0%)	91	97	
2	D	3/3 (100%)	3 (100%)	0	100	100	
2	E	3/3 (100%)	3 (100%)	0	100	100	
2	F	3/3 (100%)	3 (100%)	0	100	100	
All	All	699/834 (84%)	691 (99%)	8 (1%)	73	90	

5 of 8 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	С	64	ARG
1	A	251	ASP
1	A	225	ASP
1	A	220	PHE
1	A	239	GLN

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

Mol	Chain	Res	Type
1	С	295	ASN
1	С	321	GLN
1	A	295	ASN
1	A	321	GLN
1	A	345	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

3 non-standard protein/DNA/RNA residues are 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	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Chain Res	Link	Bo	Bond lengths			Bond angles		
Mol Type	nes		Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2													
1	MIS	В	201	1	11,12,13	0.65	0	10,16,18	0.87	0												
1	MIS	С	201	1	11,12,13	0.65	0	10,16,18	0.58	0												
1	MIS	A	201	1	11,12,13	0.65	0	10,16,18	0.57	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
1	MIS	В	201	1	-	5/10/13/15	-
1	MIS	С	201	1	-	6/10/13/15	-
1	MIS	A	201	1	-	6/10/13/15	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 17 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	201	MIS	CB-OG-P-O2P
1	A	201	MIS	C1-O3P-P-O1P
1	В	201	MIS	CB-OG-P-O1P
1	В	201	MIS	CB-OG-P-O2P
1	В	201	MIS	CB-OG-P-O3P

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	В	201	MIS	1	0
1	С	201	MIS	2	0
1	A	201	MIS	4	0



5.5 Carbohydrates (i)

There are no monosaccharides 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, 95^{th} 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>	$\#\mathrm{RSRZ}{>}2$	$\mathrm{OWAB}(\mathrm{\AA}^2)$	Q < 0.9
1	A	294/340~(86%)	0.01	8 (2%) 54 26	66, 133, 179, 197	0
1	В	278/340 (81%)	0.07	15 (5%) 25 9	51, 81, 206, 234	0
1	С	292/340 (85%)	-0.20	2 (0%) 87 69	60, 100, 151, 184	0
2	D	3/3 (100%)	0.73	0 100 100	175, 175, 181, 199	0
2	E	3/3 (100%)	2.65	2 (66%) 0 0	205, 205, 212, 219	0
2	F	3/3 (100%)	-0.24	0 100 100	117, 117, 152, 155	0
All	All	873/1029 (84%)	-0.03	27 (3%) 49 21	51, 105, 181, 234	0

The worst 5 of 27 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	297	GLY	5.3
1	В	291	GLY	4.8
1	В	301	ASN	4.4
1	A	263	GLY	4.1
1	В	224	ASN	3.8

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

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, 95^{th} 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
1	MIS	A	201	13/14	0.97	0.16	71,86,98,105	0
1	MIS	В	201	13/14	0.98	0.19	48,49,66,66	0
1	MIS	С	201	13/14	0.98	0.16	57,64,78,83	0



6.3 Carbohydrates (i)

There are no monosaccharides 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.

