Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 11:20 am BST

PDB ID	:	4SGA
Title	:	STRUCTURES OF PRODUCT AND INHIBITOR COMPLEXES OF STREPTOMYCES GRISEUS PROTEASE A AT 1.8 ANGSTROMS RESOLUTION. A MODEL FOR SERINE PROTEASE CATALYSIS
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Deposited on	:	1990-05-29
Resolution	:	1.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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 following versions of software and data (see references ⓘ) were used in the production of this report:

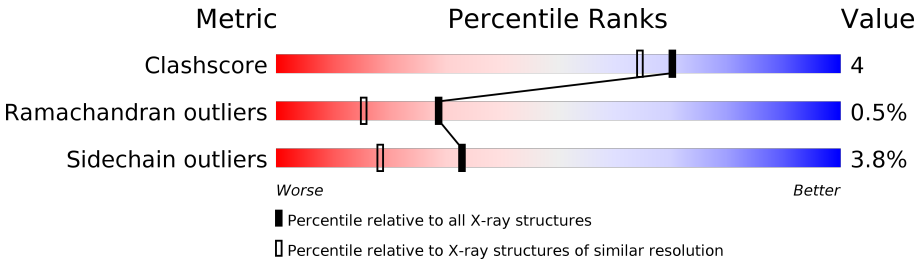
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 1.80 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	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)

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 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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	E	181	85% 14% .
2	P	5	80% 20%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1477 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 PROTEINASE A (SGPA).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	181	Total	C	N	O	S	0	0	0
			1259	768	224	262	5			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	192A	GLN	GLU	CONFLICT	UNP P00776

- Molecule 2 is a protein called TETRAPEPTIDE ACE-PRO-ALA-PRO-PHE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	P	5	Total	C	N	O	0	0	0
			34	24	4	6			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	181	Total	O	0	0
			181	181		
3	P	3	Total	O	0	0
			3	3		

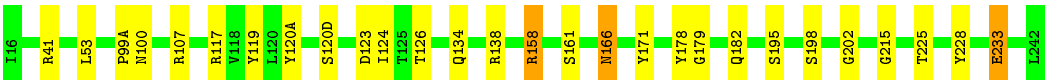
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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. 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. 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: PROTEINASE A (SGPA)

Chain E:



- Molecule 2: TETRAPEPTIDE ACE-PRO-ALA-PRO-PHE

Chain P:

4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 42	Depositor
Cell constants a, b, c, α, β, γ	55.20Å 55.20Å 54.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.80	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-1.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.116 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1477	wwPDB-VP
Average B, all atoms (Å^2)	14.0	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: ACE

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	E	1.18	2/1280 (0.2%)	1.77	15/1742 (0.9%)
2	P	1.48	0/34	1.47	1/46 (2.2%)
All	All	1.19	2/1314 (0.2%)	1.77	16/1788 (0.9%)

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	E	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	202	GLY	N-CA	5.77	1.54	1.46
1	E	215	GLY	N-CA	5.20	1.53	1.46

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	138	ARG	NE-CZ-NH1	21.51	131.06	120.30
1	E	41	ARG	NE-CZ-NH1	12.88	126.74	120.30
1	E	107	ARG	NE-CZ-NH1	12.72	126.66	120.30
1	E	138	ARG	NE-CZ-NH2	-11.15	114.72	120.30
1	E	107	ARG	NE-CZ-NH2	-8.23	116.19	120.30
1	E	41	ARG	NE-CZ-NH2	-7.67	116.46	120.30
1	E	123	ASP	CB-CG-OD1	7.35	124.92	118.30
1	E	138	ARG	CD-NE-CZ	7.07	133.50	123.60
1	E	117	ARG	NE-CZ-NH2	6.94	123.77	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	178	TYR	CB-CG-CD2	-6.20	117.28	121.00
1	E	171	TYR	CB-CG-CD1	-5.47	117.72	121.00
1	E	158	ARG	NE-CZ-NH2	-5.39	117.60	120.30
1	E	233	GLU	CG-CD-OE1	5.20	128.70	118.30
1	E	195	SER	O-C-N	5.11	131.89	123.20
2	P	5	ACE	C-N-CD	5.11	139.14	128.40
1	E	123	ASP	CB-CG-OD2	-5.00	113.80	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	158	ARG	Sidechain

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	E	1259	0	1188	9	0
2	P	34	0	31	0	0
3	E	181	0	0	3	0
3	P	3	0	0	0	0
All	All	1477	0	1219	9	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 4.

All (9) 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:E:182:GLN:HE21	1:E:225:THR:HG23	1.74	0.51
1:E:166:ASN:HA	1:E:179:GLY:HA2	1.95	0.48
1:E:126:THR:HG22	3:E:335:HOH:O	2.13	0.47
1:E:119:TYR:CE2	1:E:120(A):TYR:HA	2.51	0.45
1:E:182:GLN:HE21	1:E:225:THR:CG2	2.29	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:124:ILE:O	3:E:284:HOH:O	2.21	0.45
1:E:228:TYR:N	1:E:228:TYR:CD2	2.84	0.44
1:E:134:GLN:O	1:E:161:SER:HA	2.20	0.42
1:E:233:GLU:HB2	3:E:270:HOH:O	2.20	0.42

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	E	179/181 (99%)	174 (97%)	4 (2%)	1 (1%)	25 12
2	P	3/5 (60%)	3 (100%)	0	0	100 100
All	All	182/186 (98%)	177 (97%)	4 (2%)	1 (0%)	29 15

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	100	ASN

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	E	129/130 (99%)	124 (96%)	5 (4%)	32 17

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	P	3/3 (100%)	3 (100%)	0	100 100
All	All	132/133 (99%)	127 (96%)	5 (4%)	33 18

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	53	LEU
1	E	99(A)	PRO
1	E	120(D)	SER
1	E	166	ASN
1	E	198	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	122	GLN
1	E	166	ASN
1	E	182	GLN

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

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 is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.