



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 04:48 PM EDT

PDB ID : 6SE7 / pdb_00006se7
Title : R600A mutant from Mycoplasma genitalium P110 Adhesin at 1.87 Angstroms resolution
Authors : Fita, I.; Aparicio, D.
Deposited on : 2019-07-29
Resolution : 1.87 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 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

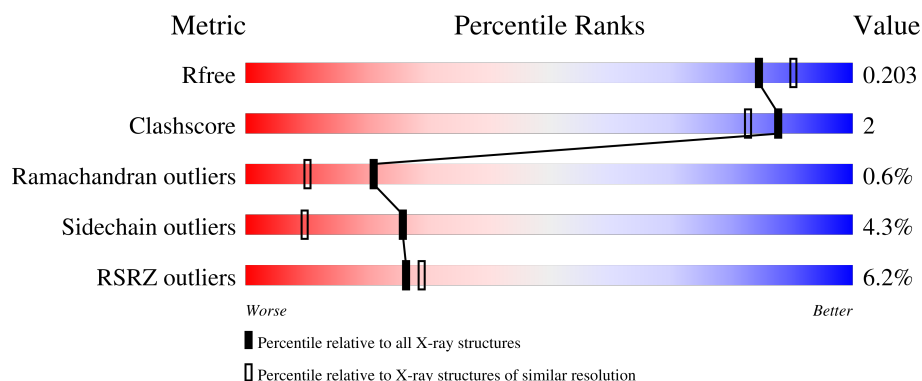
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.87 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	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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	923	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7515 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 Mgp-operon protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	885	Total	C	N	O	S	0	0	0
			6814	4276	1136	1396	6			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	939	HIS	-	expression tag	UNP P22747
A	940	HIS	-	expression tag	UNP P22747
A	941	HIS	-	expression tag	UNP P22747
A	942	HIS	-	expression tag	UNP P22747
A	943	HIS	-	expression tag	UNP P22747
A	944	HIS	-	expression tag	UNP P22747

- Molecule 2 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	K	0	0
			1	1		

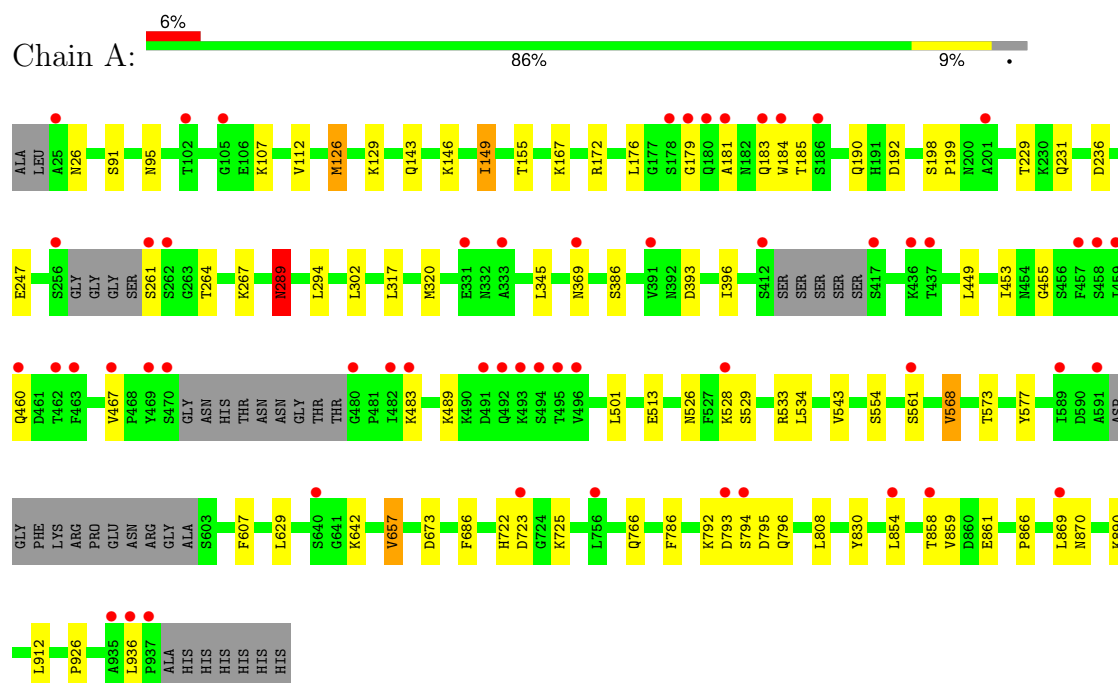
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	700	Total	O	0	0
			700	700		

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: Mgp-operon protein 3



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	109.39Å 150.70Å 176.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.36 – 1.87 27.36 – 1.87	Depositor EDS
% Data completeness (in resolution range)	61.5 (27.36-1.87) 61.5 (27.36-1.87)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 1.87Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.172 , 0.207 0.174 , 0.203	Depositor DCC
R_{free} test set	3693 reflections (3.06%)	wwPDB-VP
Wilson B-factor (Å ²)	40.2	Xtriage
Anisotropy	0.047	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 59.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7515	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% 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

5.1 Standard geometry

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

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	0.85	1/6954 (0.0%)	1.20	20/9459 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	126	MET	SD-CE	-13.08	1.46	1.79

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	766	GLN	CA-C-N	7.30	132.38	120.63
1	A	766	GLN	C-N-CA	7.30	132.38	120.63
1	A	95	ASN	CA-C-N	7.18	129.90	120.28
1	A	95	ASN	C-N-CA	7.18	129.90	120.28
1	A	722	HIS	CA-C-N	6.89	133.28	121.29
1	A	722	HIS	C-N-CA	6.89	133.28	121.29
1	A	236	ASP	CA-C-N	6.40	128.86	120.28
1	A	236	ASP	C-N-CA	6.40	128.86	120.28
1	A	155	THR	CA-C-N	6.28	129.55	120.38
1	A	155	THR	C-N-CA	6.28	129.55	120.38
1	A	607	PHE	CA-CB-CG	5.68	119.48	113.80
1	A	26	ASN	CA-CB-CG	5.57	118.17	112.60
1	A	795	ASP	N-CA-C	-5.56	107.13	114.31
1	A	289	ASN	CA-CB-CG	5.50	118.11	112.60
1	A	786	PHE	N-CA-C	5.45	118.33	109.06
1	A	526	ASN	CA-CB-CG	5.32	117.92	112.60
1	A	657	VAL	N-CA-CB	5.28	120.98	111.21
1	A	686	PHE	CA-CB-CG	5.26	119.06	113.80
1	A	673	ASP	CA-CB-CG	5.23	117.83	112.60
1	A	149	ILE	CB-CA-C	5.06	117.71	110.33

There are no chirality outliers.

There are no planarity outliers.

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	6814	0	6596	27	0
2	A	1	0	0	0	0
3	A	700	0	0	3	0
All	All	7515	0	6596	27	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 2.

All (27) 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:A:796:GLN:HB3	3:A:1321:HOH:O	1.67	0.94
1:A:126:MET:HE1	1:A:129:LYS:HD2	1.56	0.85
1:A:289:ASN:CG	3:A:1103:HOH:O	2.33	0.71
1:A:294:LEU:HD13	1:A:320:MET:HE2	1.73	0.69
1:A:167:LYS:HD2	1:A:513:GLU:HB3	1.76	0.67
1:A:554:SER:HB2	1:A:568:VAL:HG13	1.77	0.65
1:A:192:ASP:O	1:A:483:LYS:HD2	2.00	0.62
1:A:858:THR:HG23	1:A:861:GLU:H	1.72	0.55
1:A:723:ASP:HB3	1:A:725:LYS:H	1.71	0.55
1:A:126:MET:HE2	1:A:129:LYS:HB2	1.90	0.52
1:A:533:ARG:HD3	1:A:577:TYR:HE2	1.76	0.51
1:A:112:VAL:CG1	1:A:126:MET:HE3	2.41	0.50
1:A:112:VAL:HG11	1:A:126:MET:HE3	1.94	0.50
1:A:184:TRP:CD2	1:A:199:PRO:HB3	2.46	0.50
1:A:231:GLN:HG2	1:A:267:LYS:HB3	1.94	0.50
1:A:912:LEU:HD23	3:A:1140:HOH:O	2.15	0.47
1:A:176:LEU:HD13	1:A:183:GLN:HG2	1.97	0.47
1:A:294:LEU:CD1	1:A:320:MET:HE2	2.45	0.45
1:A:172:ARG:HD2	1:A:181:ALA:HA	1.99	0.45
1:A:859:VAL:HG11	1:A:890:LYS:HG3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:866:PRO:HA	1:A:869:LEU:CD2	2.49	0.42
1:A:172:ARG:HG3	1:A:176:LEU:HD12	2.01	0.42
1:A:345:LEU:HD22	1:A:573:THR:HG21	2.00	0.42
1:A:143:GLN:HA	1:A:146:LYS:HE3	2.03	0.41
1:A:501:LEU:C	1:A:501:LEU:HD12	2.45	0.41
1:A:396:ILE:HG22	1:A:629:LEU:HD21	2.03	0.40
1:A:830:TYR:CE1	1:A:926:PRO:HD3	2.56	0.40

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	875/923 (95%)	843 (96%)	27 (3%)	5 (1%)	21	10

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	794	SER
1	A	179	GLY
1	A	460	GLN
1	A	453	ILE
1	A	455	GLY

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	774/802 (96%)	741 (96%)	33 (4%)	26	10

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	91	SER
1	A	107	LYS
1	A	149	ILE
1	A	185	THR
1	A	190	GLN
1	A	198	SER
1	A	229	THR
1	A	247	GLU
1	A	261	SER
1	A	264	THR
1	A	289	ASN
1	A	302	LEU
1	A	317	LEU
1	A	369	ASN
1	A	386	SER
1	A	393	ASP
1	A	449	LEU
1	A	467	VAL
1	A	489	LYS
1	A	528	LYS
1	A	529	SER
1	A	534	LEU
1	A	543	VAL
1	A	561	SER
1	A	568	VAL
1	A	642	LYS
1	A	657	VAL
1	A	792	LYS
1	A	793	ASP
1	A	808	LEU
1	A	854	LEU
1	A	870	ASN
1	A	936	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	ASN
1	A	131	GLN
1	A	143	GLN
1	A	147	ASN
1	A	231	GLN
1	A	285	GLN
1	A	344	ASN
1	A	352	GLN
1	A	369	ASN
1	A	380	GLN
1	A	492	GLN
1	A	823	ASN
1	A	864	ASN
1	A	891	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](#)

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](#)

Of 1 ligands modelled in this entry, 1 is 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 ⓘ

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	885/923 (95%)	0.20	55 (6%) 26 29	26, 55, 107, 165	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	591	ALA	6.7
1	A	459	ILE	6.1
1	A	492	GLN	4.7
1	A	181	ALA	4.5
1	A	936	LEU	4.3
1	A	495	THR	4.0
1	A	589	ILE	4.0
1	A	482	ILE	3.6
1	A	256	SER	3.4
1	A	937	PRO	3.4
1	A	493	LYS	3.4
1	A	261	SER	3.4
1	A	437	THR	3.3
1	A	935	ALA	3.3
1	A	869	LEU	3.3
1	A	491	ASP	3.2
1	A	793	ASP	3.2
1	A	469	TYR	3.1
1	A	457	PHE	3.1
1	A	25	ALA	3.1
1	A	179	GLY	3.0
1	A	184	TRP	2.9
1	A	794	SER	2.9
1	A	480	GLY	2.8
1	A	494	SER	2.7
1	A	528	LYS	2.7
1	A	180	GLN	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	640	SER	2.6
1	A	183	GLN	2.6
1	A	470	SER	2.6
1	A	858	THR	2.6
1	A	369	ASN	2.5
1	A	417	SER	2.5
1	A	105	GLY	2.4
1	A	854	LEU	2.4
1	A	723	ASP	2.4
1	A	483	LYS	2.4
1	A	460	GLN	2.4
1	A	496	VAL	2.4
1	A	186	SER	2.4
1	A	333	ALA	2.3
1	A	436	LYS	2.3
1	A	262	SER	2.3
1	A	463	PHE	2.3
1	A	412	SER	2.3
1	A	462	THR	2.3
1	A	561	SER	2.2
1	A	102	THR	2.2
1	A	178	SER	2.2
1	A	391	VAL	2.2
1	A	467	VAL	2.2
1	A	331	GLU	2.1
1	A	756	LEU	2.1
1	A	458	SER	2.0
1	A	201	ALA	2.0

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](#)

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, 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	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	K	A	1001	1/1	1.00	0.01	32,32,32,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.