



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 10:57 AM UTC

PDB ID : 9CGO / pdb_00009cgo
Title : Tylosin thioesterase domain (TylG5 TE)
Authors : Smith, J.L.; Choudhary, V.
Deposited on : 2024-06-30
Resolution : 1.93 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

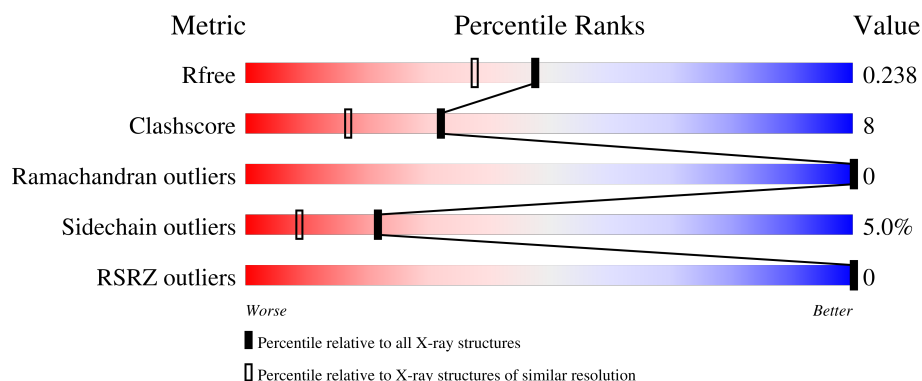
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.93 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	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	281	
1	B	281	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4209 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 Tylactone synthase module 7.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	Se	0	0	0
			1986	1264	361	350	1	10			
1	B	273	Total	C	N	O	S	Se	0	0	0
			2013	1279	366	357	1	10			

- Molecule 2 is water.

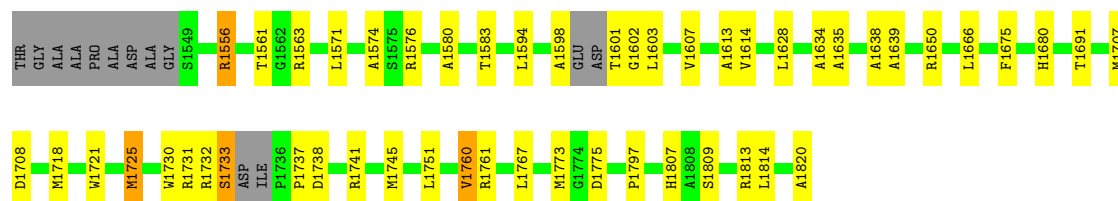
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	109	Total	O	0	0
			109	109		
2	B	101	Total	O	0	0
			101	101		

3 Residue-property plots

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.

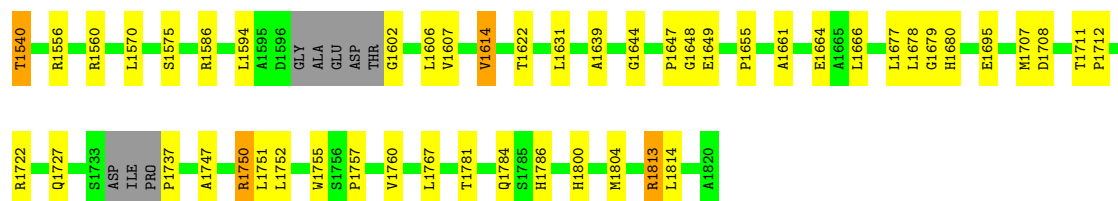
• Molecule 1: Tylactone synthase module 7

Chain A: 



• Molecule 1: Tylactone synthase module 7

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	56.88Å 56.88Å 292.47Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.26 – 1.93 49.26 – 1.93	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.26-1.93) 100.0 (49.26-1.93)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.97 (at 1.94Å)	Xtriage
Refinement program	PHENIX 1.21_5207	Depositor
R, R_{free}	0.176 , 0.210 (Not available) , 0.238	Depositor DCC
R_{free} test set	2042 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	28.6	Xtriage
Anisotropy	0.409	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 40.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.327 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4209	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.22% 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

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	0.33	0/2039	0.57	0/2775
1	B	0.33	0/2066	0.56	0/2812
All	All	0.33	0/4105	0.57	0/5587

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

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	1986	0	1928	33	0
1	B	2013	0	1950	33	0
2	A	109	0	0	6	0
2	B	101	0	0	11	1
All	All	4209	0	3878	63	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 8.

All (63) 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:1614:VAL:HG23	1:A:1741:ARG:HB3	1.65	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1614:VAL:HG21	1:A:1737:PRO:HB3	1.64	0.78
1:A:1732:ARG:NH1	2:A:1901:HOH:O	2.20	0.74
1:B:1711:THR:OG1	2:B:1901:HOH:O	1.75	0.74
1:B:1695:GLU:OE2	2:B:1902:HOH:O	2.09	0.70
1:B:1586:ARG:NH1	2:B:1907:HOH:O	2.27	0.67
1:A:1721:TRP:HB3	1:A:1773:MSE:HE1	1.77	0.66
1:A:1603:LEU:HD21	1:A:1820:ALA:HB2	1.83	0.60
1:A:1580:ALA:H	1:B:1540:THR:HA	1.66	0.60
1:B:1647:PRO:O	2:B:1903:HOH:O	2.16	0.59
1:A:1601:THR:HB	1:A:1603:LEU:H	1.66	0.59
1:B:1678:LEU:HD11	1:B:1707:MSE:HE3	1.86	0.57
1:A:1576:ARG:NH2	2:A:1909:HOH:O	2.38	0.56
1:A:1718:MSE:SE	1:A:1725:MSE:SE	3.24	0.56
1:A:1594:LEU:HD11	1:A:1639:ALA:HB2	1.89	0.55
1:A:1601:THR:N	1:A:1602:GLY:HA2	2.23	0.54
1:A:1556:ARG:HD2	2:A:1917:HOH:O	2.09	0.52
1:A:1730:TRP:CD1	1:A:1737:PRO:HG2	2.45	0.52
1:B:1722:ARG:NH2	2:B:1912:HOH:O	2.42	0.51
1:B:1712:PRO:HG2	1:B:1752:LEU:HB3	1.92	0.50
1:B:1607:VAL:HB	1:B:1677:LEU:HD23	1.92	0.50
1:B:1644:GLY:HA2	1:B:1649:GLU:HB2	1.93	0.50
1:A:1731:ARG:NH1	2:A:1911:HOH:O	2.42	0.49
1:B:1614:VAL:HG11	1:B:1737:PRO:HB3	1.95	0.49
1:B:1664:GLU:OE2	2:B:1904:HOH:O	2.18	0.49
1:A:1571:LEU:HD12	1:A:1745:MSE:HE2	1.95	0.48
1:A:1809:SER:O	1:A:1813:ARG:HG3	2.13	0.48
1:A:1580:ALA:O	1:B:1540:THR:HA	2.13	0.48
1:A:1580:ALA:O	1:A:1650:ARG:NH2	2.46	0.48
1:B:1707:MSE:HG2	1:B:1767:LEU:HD12	1.96	0.48
1:A:1767:LEU:HD21	1:A:1814:LEU:HD22	1.95	0.47
1:B:1655:PRO:HG3	1:B:1751:LEU:HD22	1.94	0.47
1:B:1755:TRP:HZ3	2:B:1901:HOH:O	1.97	0.47
1:B:1556:ARG:O	1:B:1560:ARG:HG3	2.15	0.47
1:A:1561:THR:OG1	1:A:1563:ARG:HD3	2.16	0.46
1:A:1775:ASP:OD2	2:A:1902:HOH:O	2.21	0.46
1:A:1613:ALA:HA	1:A:1745:MSE:HG3	1.98	0.46
1:A:1680:HIS:HA	1:A:1707:MSE:HB2	1.98	0.46
1:B:1586:ARG:HD2	1:B:1661:ALA:HB2	1.99	0.45
1:A:1598:ALA:HB3	1:A:1634:ALA:N	2.31	0.45
1:B:1594:LEU:HD11	1:B:1639:ALA:HB2	1.99	0.45
1:B:1707:MSE:SE	1:B:1804:MSE:SE	3.35	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1607:VAL:HA	1:A:1638:ALA:O	2.16	0.44
1:B:1647:PRO:HA	1:B:1648:GLY:HA2	1.60	0.44
1:B:1602:GLY:N	2:B:1921:HOH:O	2.51	0.44
1:B:1680:HIS:NE2	1:B:1800:HIS:O	2.51	0.43
1:A:1666:LEU:HD21	1:A:1675:PHE:CE2	2.53	0.43
1:B:1575:SER:HB2	1:B:1747:ALA:HB2	2.01	0.42
1:B:1678:LEU:HD23	1:B:1679:GLY:N	2.33	0.42
1:A:1603:LEU:HB3	1:A:1635:ALA:HB2	2.02	0.42
1:A:1797:PRO:HB2	1:A:1807:HIS:CG	2.54	0.42
1:A:1730:TRP:O	1:A:1733:SER:HB2	2.19	0.41
1:A:1761:ARG:NH1	2:A:1918:HOH:O	2.53	0.41
1:A:1574:ALA:HA	1:B:1570:LEU:HD21	2.03	0.41
1:B:1607:VAL:HG21	1:B:1666:LEU:HD11	2.02	0.41
1:B:1781:THR:HB	1:B:1784:GLN:HG3	2.02	0.41
1:A:1738:ASP:OD1	1:A:1738:ASP:N	2.48	0.41
1:B:1678:LEU:HD11	1:B:1707:MSE:CE	2.50	0.41
1:B:1760:VAL:O	2:B:1905:HOH:O	2.22	0.41
1:B:1813:ARG:HD2	1:B:1813:ARG:HA	1.54	0.41
1:B:1750:ARG:HD2	2:B:1942:HOH:O	2.20	0.40
1:A:1691:THR:HG21	1:A:1760:VAL:HG11	2.04	0.40
1:B:1757:PRO:HD3	2:B:1901:HOH:O	2.22	0.40

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)
2:B:1909:HOH:O	2:B:1984:HOH:O[6_554]	2.16	0.04

5.3 Torsion angles

5.3.1 Protein backbone

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	33/281 (12%)	32 (97%)	1 (3%)	0	100	100
1	B	219/281 (78%)	208 (95%)	11 (5%)	0	100	100
All	All	252/562 (45%)	240 (95%)	12 (5%)	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	188/185 (102%)	180 (96%)	8 (4%)	26	13
1	B	189/185 (102%)	178 (94%)	11 (6%)	18	6
All	All	377/370 (102%)	358 (95%)	19 (5%)	22	9

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1556	ARG
1	A	1583	THR
1	A	1628	LEU
1	A	1708	ASP
1	A	1725	MSE
1	A	1733	SER
1	A	1751	LEU
1	A	1760	VAL
1	B	1540	THR
1	B	1606	LEU
1	B	1614	VAL
1	B	1622	THR
1	B	1631	LEU
1	B	1708	ASP
1	B	1727	GLN
1	B	1750	ARG
1	B	1786	HIS
1	B	1813	ARG

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Mol	Chain	Res	Type
1	B	1814	LEU

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

Mol	Chain	Res	Type
1	B	1693	HIS
1	B	1697	ASN
1	B	1786	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	258/281 (91%)	-1.26	0 100 100	15, 32, 49, 71	0
1	B	263/281 (93%)	-1.29	0 100 100	17, 31, 48, 58	0
All	All	521/562 (92%)	-1.27	0 100 100	15, 31, 49, 71	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.