



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:11 PM UTC

PDB ID : 8G0K / pdb_00008g0k
Title : Bacteroides multispecies type VI secretion system Ntox15 domain effector and immunity Tde1/Tdi1
Authors : Bosch, D.E.; Mougous, J.D.
Deposited on : 2023-01-31
Resolution : 2.20 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

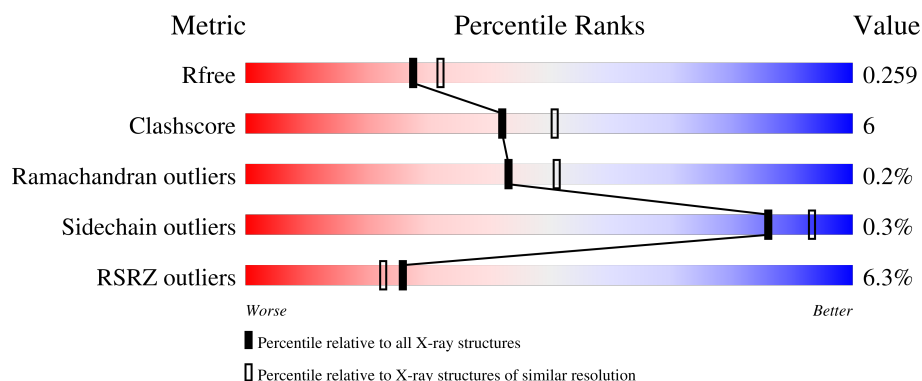
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	195	6% 62% 7% 31%
1	C	195	9% 42% 16% 43%
2	B	195	2% 87% 11% ..
2	D	195	4% 88% 11% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5607 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 Ntox15 domain-containing protein.

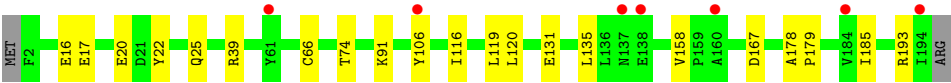
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	135	Total	C	N	O	S	0	0	0
			1132	728	191	207	6			
1	C	112	Total	C	N	O	S	0	0	0
			945	605	162	174	4			

- Molecule 2 is a protein called DUF1851 domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	193	Total	C	N	O	S	0	0	0
			1590	1046	245	290	9			
2	D	193	Total	C	N	O	S	0	0	0
			1590	1046	245	290	9			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	83	Total	O	0	0
			83	83		
3	B	113	Total	O	0	0
			113	113		
3	C	66	Total	O	0	0
			66	66		
3	D	88	Total	O	0	0
			88	88		



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	109.77Å 112.27Å 149.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.79 – 2.20 34.79 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.2 (34.79-2.20) 99.3 (34.79-2.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.20	Depositor
R, R_{free}	0.207 , 0.258 0.210 , 0.259	Depositor DCC
R_{free} test set	1997 reflections (4.24%)	wwPDB-VP
Wilson B-factor (Å ²)	34.2	Xtriage
Anisotropy	0.630	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 30.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.185 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5607	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.64% 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	0.55	0/1146	0.61	0/1528
1	C	0.48	0/959	0.66	2/1277 (0.2%)
2	B	0.58	0/1627	0.75	0/2205
2	D	0.54	0/1627	0.69	0/2205
All	All	0.54	0/5359	0.68	2/7215 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	268	ALA	CA-C-N	6.39	130.50	121.61
1	C	268	ALA	C-N-CA	6.39	130.50	121.61

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	1132	0	1164	14	0
1	C	945	0	953	27	0
2	B	1590	0	1587	16	0
2	D	1590	0	1587	16	0
3	A	83	0	0	1	0
3	B	113	0	0	0	0
3	C	66	0	0	4	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	88	0	0	2	0
All	All	5607	0	5291	66	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 6.

All (66) 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:C:173:LEU:HD12	1:C:342:ILE:HD13	1.59	0.82
1:A:167:VAL:HG22	1:A:304:ARG:HG2	1.67	0.73
1:A:185:ILE:HD12	1:A:185:ILE:H	1.55	0.70
1:C:173:LEU:HB2	1:C:340:LEU:HD21	1.72	0.70
1:A:167:VAL:CG2	1:A:304:ARG:HG2	2.27	0.64
2:B:32:GLU:H	2:B:32:GLU:CD	2.06	0.64
1:C:218:ARG:NH1	2:D:193:ARG:HD2	2.14	0.62
1:C:214:LYS:NZ	1:C:309:GLU:OE1	2.32	0.62
1:A:304:ARG:HB2	2:B:186:GLN:O	2.04	0.57
2:B:106:TYR:OH	2:B:167:ASP:OD2	2.21	0.57
1:C:193:MET:HG2	2:D:158:VAL:HG11	1.88	0.56
1:A:185:ILE:O	1:A:189:LYS:HG3	2.07	0.55
2:B:59:TYR:O	2:B:63:ILE:HG12	2.07	0.55
1:C:195:GLN:HB2	1:C:336:LEU:HD22	1.89	0.55
1:A:342:ILE:O	1:A:343:VAL:HB	2.06	0.55
2:D:16:GLU:HG3	2:D:17:GLU:N	2.23	0.53
1:C:224:LYS:HG3	2:D:193:ARG:NH2	2.24	0.53
1:C:190:GLN:HA	1:C:193:MET:HE3	1.91	0.53
1:C:206:TRP:HE3	1:C:207:LEU:HG	1.74	0.53
1:A:218:ARG:NH2	2:B:193:ARG:HD2	2.26	0.51
1:C:222:SER:N	3:C:407:HOH:O	2.43	0.51
1:A:175:LYS:HG3	1:A:343:VAL:HG22	1.92	0.50
1:C:195:GLN:C	1:C:197:GLY:H	2.20	0.50
1:C:247:ASP:OD2	1:C:247:ASP:N	2.42	0.50
1:A:216:TYR:OH	3:A:401:HOH:O	2.20	0.49
2:D:20:GLU:OE1	2:D:39:ARG:HD3	2.12	0.49
1:A:218:ARG:NH2	1:A:227:GLU:OE2	2.45	0.49
2:D:116:ILE:HG23	2:D:120:LEU:HD12	1.95	0.48
2:D:66:CYS:HB3	2:D:185:ILE:HD13	1.94	0.47
2:B:184:VAL:HG12	2:B:188:MET:HE3	1.95	0.47
1:C:184:GLU:O	1:C:188:ARG:HG3	2.15	0.47
2:D:131:GLU:O	2:D:135:LEU:HD23	2.14	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:14:LYS:NZ	2:B:17:GLU:OE1	2.43	0.47
2:D:74:THR:HG23	3:D:262:HOH:O	2.14	0.46
2:B:16:GLU:HG3	2:B:17:GLU:O	2.15	0.46
1:C:173:LEU:HD12	1:C:342:ILE:CD1	2.38	0.46
1:C:262:GLU:HB3	3:C:409:HOH:O	2.15	0.46
2:D:106:TYR:OH	2:D:167:ASP:OD2	2.34	0.46
2:D:22:TYR:O	2:D:25:GLN:HG2	2.16	0.46
2:B:63:ILE:HG13	2:B:77:PRO:HG3	1.98	0.45
2:B:178:ALA:HB3	2:B:179:PRO:HD3	1.98	0.45
1:C:206:TRP:N	3:C:408:HOH:O	2.50	0.45
1:C:167:VAL:HG23	1:C:308:ILE:HD11	1.99	0.45
2:B:63:ILE:HD11	2:B:86:ILE:CD1	2.47	0.44
2:B:185:ILE:HA	2:B:188:MET:HG2	1.99	0.44
1:C:186:GLU:HB3	2:D:179:PRO:HG2	1.99	0.44
1:A:203:ILE:HD13	1:A:312:LEU:HD11	1.98	0.44
2:D:119:LEU:HB2	3:D:255:HOH:O	2.17	0.44
2:B:127:LYS:HD2	2:B:127:LYS:HA	1.70	0.44
2:D:178:ALA:HB3	2:D:179:PRO:HD3	1.99	0.43
2:B:116:ILE:HG23	2:B:120:LEU:HD12	1.99	0.43
1:C:179:LYS:HD2	1:C:180:ASP:N	2.33	0.43
1:C:218:ARG:HH12	2:D:193:ARG:HD2	1.83	0.43
1:A:185:ILE:H	1:A:185:ILE:CD1	2.30	0.42
1:A:237:LYS:HD2	1:A:237:LYS:HA	1.79	0.42
1:C:224:LYS:NZ	1:C:228:ASP:OD2	2.52	0.42
1:C:206:TRP:CE3	1:C:207:LEU:HG	2.52	0.42
2:B:3:GLU:CD	2:B:3:GLU:H	2.28	0.41
1:C:173:LEU:CB	1:C:340:LEU:HD21	2.44	0.41
1:C:184:GLU:N	3:C:402:HOH:O	2.35	0.41
1:C:338:GLU:CD	1:C:339:ASN:H	2.28	0.41
1:C:175:LYS:O	1:C:178:GLN:HG2	2.21	0.41
2:D:91:LYS:HA	2:D:91:LYS:HD3	1.99	0.41
2:B:43:ILE:HD12	2:B:43:ILE:N	2.35	0.40
1:A:314:LYS:HA	1:A:314:LYS:HD2	1.81	0.40
1:C:175:LYS:HD2	1:C:343:VAL:HG22	2.03	0.40

There are no symmetry-related clashes.

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	127/195 (65%)	122 (96%)	5 (4%)	0	100	100
1	C	104/195 (53%)	96 (92%)	8 (8%)	0	100	100
2	B	191/195 (98%)	187 (98%)	3 (2%)	1 (0%)	24	27
2	D	191/195 (98%)	188 (98%)	3 (2%)	0	100	100
All	All	613/780 (79%)	593 (97%)	19 (3%)	1 (0%)	43	51

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	189	GLY

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	125/178 (70%)	124 (99%)	1 (1%)	73	85
1	C	102/178 (57%)	102 (100%)	0	100	100
2	B	176/178 (99%)	175 (99%)	1 (1%)	78	89
2	D	176/178 (99%)	176 (100%)	0	100	100
All	All	579/712 (81%)	577 (100%)	2 (0%)	86	93

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	304	ARG
2	B	32	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	11	ASN
2	B	48	ASN
2	B	140	ASN
2	D	25	GLN
2	D	48	ASN

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

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	135/195 (69%)	0.64	12 (8%) 15 13	26, 45, 64, 73	0
1	C	112/195 (57%)	0.91	18 (16%) 4 3	27, 59, 74, 78	0
2	B	193/195 (98%)	0.22	3 (1%) 70 67	25, 33, 48, 58	0
2	D	193/195 (98%)	0.34	7 (3%) 46 43	26, 37, 58, 64	0
All	All	633/780 (81%)	0.47	40 (6%) 26 23	25, 39, 67, 78	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	336	LEU	4.9
1	A	270	PRO	4.8
1	A	330	ILE	3.9
1	C	343	VAL	3.7
1	A	329	LYS	3.5
1	A	162	LEU	3.5
1	A	343	VAL	3.4
1	C	166	PHE	3.3
1	C	221	LYS	3.3
1	C	197	GLY	3.1
1	C	337	VAL	3.1
1	C	206	TRP	3.1
2	D	138	GLU	3.0
1	C	182	PHE	2.7
1	A	316	ILE	2.7
1	C	342	ILE	2.7
2	D	106	TYR	2.6
1	C	270	PRO	2.6
2	D	194	ILE	2.6
2	B	138	GLU	2.5
1	A	262	GLU	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	191	LEU	2.4
2	B	32	GLU	2.3
1	C	217	GLY	2.2
1	C	179	LYS	2.2
1	C	308	ILE	2.2
1	A	204	PHE	2.2
2	D	160	ALA	2.2
1	A	165	ILE	2.2
1	A	218	ARG	2.1
1	A	163	SER	2.1
1	C	339	ASN	2.1
2	D	61	TYR	2.1
2	D	137	ASN	2.1
1	A	314	LYS	2.1
2	B	169	ILE	2.1
2	D	184	VAL	2.1
1	C	181	PRO	2.0
1	C	303	GLY	2.0
1	C	185	ILE	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.