



wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 2BHV / pdb_00002bhv
Title : Structure of ComB10 of the Com Type IV secretion system of *Helicobacter pylori*
Authors : Terradot, L.; Oomen, C.; Bayliss, R.; Leonard, G.; Waksman, G.
Deposited on : 2005-01-18
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 ⓘ 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

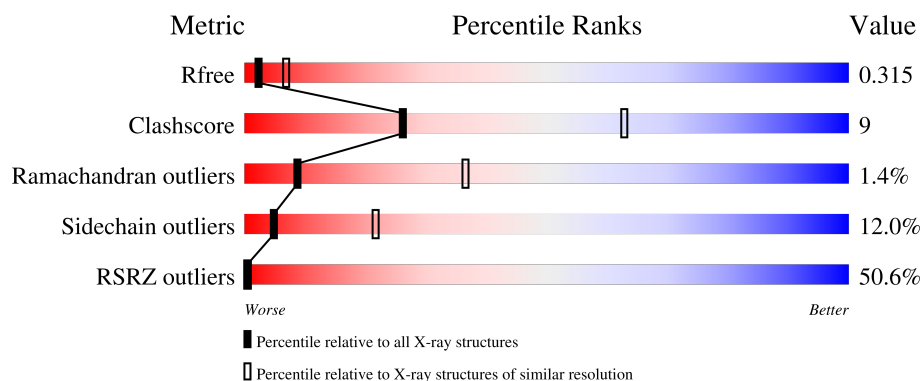
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (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 $\leq 5\%$. 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	246	77% 55% 17% • • 22%
1	B	246	72% 57% 17% • • 22%
1	C	246	5% 59% 14% • 24%
1	D	246	4% 54% 18% • 23%
1	E	246	5% 53% 21% • • 23%

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Mol	Chain	Length	Quality of chain
1	F	246	71% 56% 19% • 24%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 8759 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 COMB10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	191	Total	C	N	O	S	0	0	0
			1468	945	253	263	7			
1	B	192	Total	C	N	O	S	0	0	0
			1466	944	254	261	7			
1	C	188	Total	C	N	O	S	0	0	0
			1457	941	250	259	7			
1	D	189	Total	C	N	O	S	0	0	0
			1464	944	251	262	7			
1	E	190	Total	C	N	O	S	0	0	0
			1447	931	251	258	7			
1	F	188	Total	C	N	O	S	0	0	0
			1457	939	250	261	7			





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.73Å 139.41Å 168.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.00 20.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	98.3 (20.00-3.00) 97.9 (20.00-3.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 3.00Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.257 , 0.293 0.291 , 0.315	Depositor DCC
R_{free} test set	1674 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	74.0	Xtriage
Anisotropy	0.346	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 73.2	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	0.70	EDS
Total number of atoms	8759	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.58% 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	0.86	5/1490 (0.3%)	0.87	2/2016 (0.1%)
1	B	0.76	1/1487 (0.1%)	0.92	4/2012 (0.2%)
1	C	0.77	1/1480 (0.1%)	0.85	1/2002 (0.0%)
1	D	0.68	0/1487	0.84	1/2011 (0.0%)
1	E	1.32	11/1468 (0.7%)	0.92	4/1986 (0.2%)
1	F	0.64	0/1480	0.84	1/2002 (0.0%)
All	All	0.87	18/8892 (0.2%)	0.88	13/12029 (0.1%)

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	A	0	1
1	B	0	1
All	All	0	2

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	319	MET	SD-CE	22.23	2.35	1.79
1	E	271	ARG	CZ-NH1	19.96	1.60	1.32
1	E	274	GLN	CD-OE1	19.22	1.60	1.23
1	E	274	GLN	CD-NE2	12.66	1.59	1.33
1	C	311	GLY	N-CA	12.31	1.65	1.45

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	271	ARG	NE-CZ-NH2	-7.51	112.44	119.20
1	B	321	GLN	N-CA-C	7.32	120.25	110.88
1	F	252	ASN	N-CA-C	6.16	117.83	110.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	252	ASN	N-CA-C	5.67	118.07	110.35
1	E	252	ASN	N-CA-C	5.65	118.24	109.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	228	MET	Peptide
1	B	226	ASN	Peptide

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	1468	0	1517	39	3
1	B	1466	0	1519	29	0
1	C	1457	0	1513	26	2
1	D	1464	0	1515	35	0
1	E	1447	0	1493	34	0
1	F	1457	0	1506	29	1
All	All	8759	0	9063	167	3

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 9.

The worst 5 of 167 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:175:ASP:OD2	1:E:208:LYS:HE3	1.30	1.26
1:A:226:ASN:HB3	1:A:228:MET:HG3	1.21	1.19
1:E:319:MET:SD	1:E:319:MET:CE	2.35	1.15
1:A:319:MET:HE3	1:A:319:MET:HA	1.45	0.94
1:A:319:MET:HE1	1:D:197:GLN:HE22	1.30	0.93

All (3) 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)
1:A:208:LYS:CE	1:F:175:ASP:OD2[3_555]	1.95	0.25
1:A:232:ARG:NH1	1:C:366:ARG:CD[4_455]	2.06	0.14
1:A:234:ASP:OD1	1:C:366:ARG:NE[4_455]	2.09	0.11

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	185/246 (75%)	166 (90%)	15 (8%)	4 (2%)	5	26
1	B	186/246 (76%)	172 (92%)	8 (4%)	6 (3%)	3	18
1	C	182/246 (74%)	172 (94%)	10 (6%)	0	100	100
1	D	183/246 (74%)	171 (93%)	10 (6%)	2 (1%)	11	43
1	E	184/246 (75%)	166 (90%)	16 (9%)	2 (1%)	11	43
1	F	182/246 (74%)	171 (94%)	10 (6%)	1 (0%)	24	60
All	All	1102/1476 (75%)	1018 (92%)	69 (6%)	15 (1%)	9	36

5 of 15 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	228	MET
1	B	254	LYS
1	A	313	TYR
1	B	229	GLY
1	B	274	GLN

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	159/209 (76%)	143 (90%)	16 (10%)	7	29
1	B	158/209 (76%)	135 (85%)	23 (15%)	3	15
1	C	159/209 (76%)	142 (89%)	17 (11%)	6	26
1	D	160/209 (77%)	137 (86%)	23 (14%)	3	15
1	E	155/209 (74%)	136 (88%)	19 (12%)	4	21
1	F	159/209 (76%)	143 (90%)	16 (10%)	7	29
All	All	950/1254 (76%)	836 (88%)	114 (12%)	5	22

5 of 114 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	189	GLN
1	F	352	VAL
1	D	330	VAL
1	F	351	ARG
1	F	189	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 14 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	272	ASN
1	D	189	GLN
1	F	262	ASN
1	E	226	ASN
1	F	189	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 [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	191/246 (77%)	6.13	190 (99%) 0 0	33, 47, 56, 68	0
1	B	192/246 (78%)	4.20	177 (92%) 0 0	28, 47, 57, 68	0
1	C	188/246 (76%)	0.77	12 (6%) 25 13	32, 47, 54, 67	0
1	D	189/246 (76%)	0.74	10 (5%) 32 16	34, 47, 57, 62	0
1	E	190/246 (77%)	0.82	12 (6%) 26 13	25, 47, 56, 63	0
1	F	188/246 (76%)	4.35	175 (93%) 0 0	35, 47, 54, 67	0
All	All	1138/1476 (77%)	2.84	576 (50%) 0 0	25, 47, 56, 68	0

The worst 5 of 576 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	230	GLU	15.3
1	A	235	ILE	15.0
1	A	276	TYR	13.9
1	A	288	GLY	13.7
1	F	376	LYS	12.3

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.