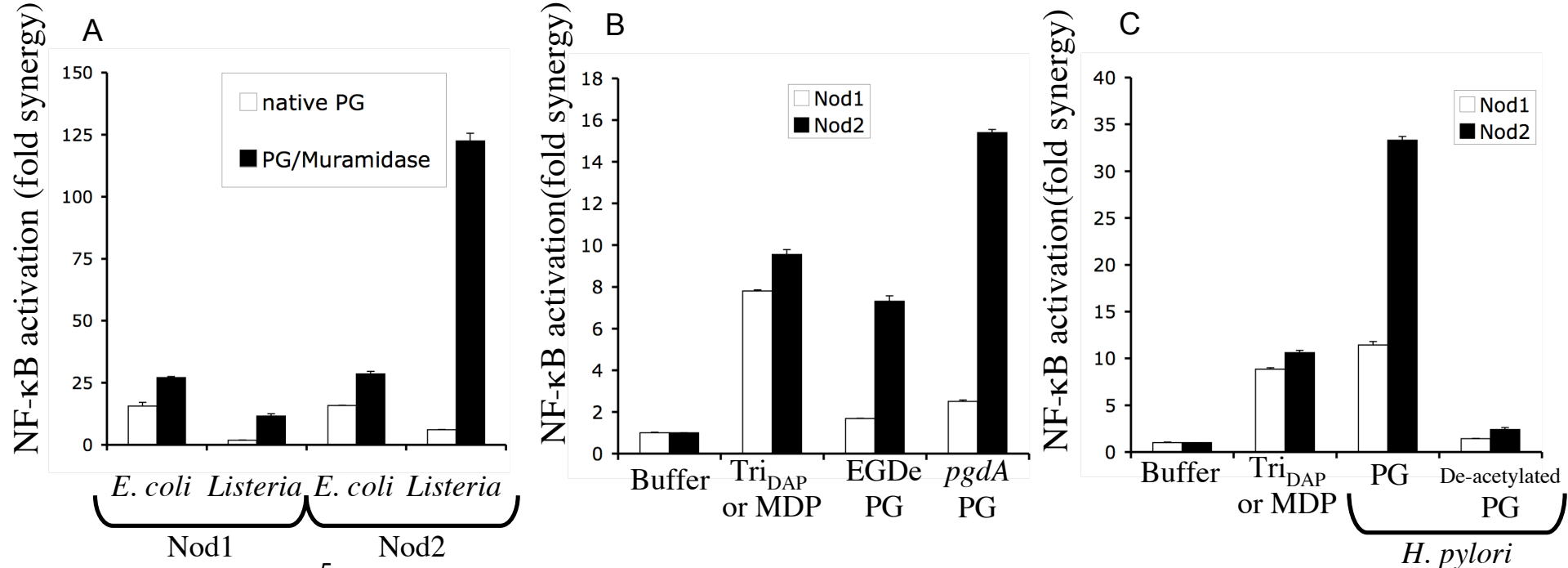

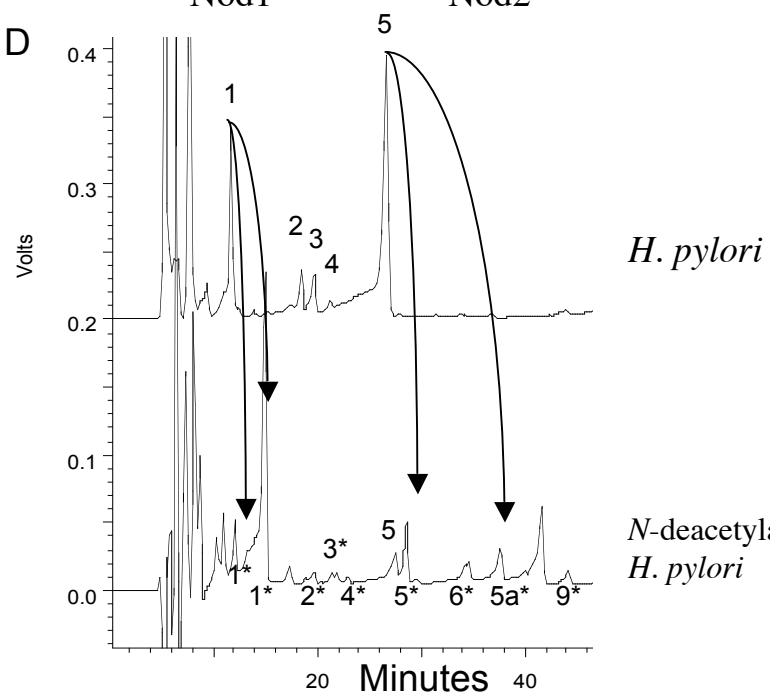

Peak no.	[M+Na]$^{+,a}$	Proposed structure[b]
1	893,3479	Glc*N*Ac-Mur*N*Ac-Ala-Glu-*m*Dap
2	964,4356	Glc*N*Ac-Mur*N*Ac-Ala-Glu-*m*Dap-Ala
3	1021,4159	Glc*N*Ac-Mur*N*Ac-Ala-Glu-*m*Dap-Ala-Gly
4	721,2981	Glc*N*Ac-Mur*N*Ac-Ala-Glu
5	1035,4398	Glc*N*Ac-Mur*N*Ac-Ala-Glu-*m*Dap-Ala-Ala
1*	851,3619	Glc*N*-Mur*N*Ac-Ala-Glu-*m*Dap
1a*	851,3745	Glc*N*-Mur*N*Ac-Ala-Glu-*m*Dap
2*	922,3697	Glc*N*-Mur*N*Ac-Ala-Glu-*m*Dap-Ala
3*	979,4075	Glc*N*-Mur*N*Ac-Ala-Glu-*m*Dap-Ala-Gly
4*	679,2736	Glc*N*-Mur*N*Ac-Ala-Glu
5*	993,4572	Glc*N*-Mur*N*Ac-Ala-Glu-*m*Dap-Ala-Ala
5a*	993,4518	Glc*N*-Mur*N*Ac-Ala-Glu-*m*Dap-Ala-Ala
6*	902,3679	Glc*N*-[(an)Mur*N*Ac]-Ala-Glu-*m*Dap-Ala
9*	973,4416	Glc*N*-[(an)Mur*N*Ac]-Ala-Glu-*m*Dap-Ala-Ala

[a] Molecular masses were measured by MALDI-TOF. De*N*-acetylation of the N-acetylgluycosamine residue was confirmed by MALDI-PSD.

[b] Abbreviations; Glc*N*Ac, *N*-acetylglucosamine; Mur*N*Ac, *N*-acetylmuramic acid; GlcN, glucosamine; (an)Mur*N*Ac, (1-6)anhydro-*N*-acetylmuramic acid; Ala, alanine; Glu, glutamic acid; *m*Dap, *meso*-diaminopimelic acid; Gly, glycine