

Table S1: Comparison of the binding energies (eV) for C_{1s}, O_{1s} and Cl_{2p}

	C 1s	O 1s	Cl 2p
HOCH ₂ CH ₂ CH ₂ Cl/Si(100) ^a			
HO-C ¹ H ₂ C ² H ₂ C ³ H ₂ -Cl	286.5 for C ¹ , C ³ 285.4 for C ²	533.1	201.9(2p _{1/2}) 200.3(2p _{3/2})
Si-O-C ¹ H ₂ C ² H ₂ C ³ H ₂ -Cl	286.6 for C ¹ , C ³ 285.5 for C ²	532.2	202.0(2p _{1/2}) 200.4(2p _{3/2})
Si-O-C ¹ H ₂ C ² H ₂ C ³ H ₂ C ³ H ₂ C ² H ₂ C ¹ H ₂ -O-Si	286.5 for C ¹ 285.5 for C ² , C ³	532.3	
CH ₂ Cl(CH ₃) ₂ COH/Ni(100) ^b			
CH ₂ Cl(CH ₃) ₂ COH		532.8	202.0(2p _{1/2}) 200.2(2p _{3/2})
-CH ₂ (CH ₃) ₂ COH		532.8	
-CH ₂ (CH ₃) ₂ CO-		531.7	
O-		530.2	
Cl-			200.1(2p _{1/2}) 198.4(2p _{3/2})
ICH ₂ CH ₂ OH/Ni(100) ^c			
ICH ₂ CH ₂ OH		532.8	
ICH ₂ CH ₂ O-		531.8	
O-		530.8	
CO/Ni(111) ^d			
Hollow	285.24	530.86	
Bridge	285.32	531.01	
On-top	285.96	532.15	
C ₂ H ₄ /Ni(111) ^e			
C ₂ H ₄	284.0		
C ₂ H ₂	283.5		

a – e: References 26, 9, 6, 27 and 28.