

	H ₁	H ₂	H ₃	H _{4/5}	C ₁	C ₂	C ₃	C ₄	C ₅	C ₆	N ₁	N ₂	N ₃
H ₁	-5800												
H ₂	1.0	-4989											
H ₃	0.0	0.0	-2700										
H _{4/5}	0.0	-0.7	7.3	-2205									
C ₁	0.0	0.0	-5.0	3.5	-30443								
C ₂	0.0	0.0	146.6	-4.3	26.80	-9458							
C ₃	1.5	0.6	-4.5	132.5	-1.0	34.8	-4589						
C ₄	6.5	10.7	5.0	-6.9	2.4	-2.6	52.8	-22562					
C ₅	-5.4	198.2	0.0	0.0	0.0	1.5	4.7	72.9	-20657				
C ₆	218.5	-6.8	0.0	0.0	0.0	0.0	0.0	2.9	-1.6	-23633			
N ₁	-6.5	-3.4	0.0	0.0	0.0	0.0	0.0	-7.4	-0.8	-13.4	-13342		
N ₂	-5.9	-4.6	0.0	0.0	0.0	0.0	0.0	0.0	-11.0	-14.7	0.0	-12351	
N ₃	0.0	0.0	0.9	2.7	0.0	-5.8	0.0	0.0	0.0	0.0	0.0	0.0	-2762
T ₂ (ms)	180	80	140	NA	180	125	80	65	135	105	50	90	250
T ₂ (ms)	295	285	190	65	260	255	215	170	190	155	140	215	395

