

	Glucose, mM					C3		D3						
	0'	15'	30'	60'	120'			0'	15'	30'	60'	120'		
Nº1C3		3.9	14.7	11.7	8.3	5.8	Nº1D3		5.2	15.5	16.8	12.9	9.6	C3-D3 C8-D8 C18-D18
Nº2C3		3.9	9.8	10.1	7.4	5.0	Nº2D3		4.9	13.2	15.9	13.6	11.2	
Nº3C3		4.3	14.5	13.2	8.3	5.6	Nº3D3		6.1	17.5	19.7	15.8	12.5	
Nº4C3		3.9	14.3	12.0	7.2	4.9	Nº4D3		5.3	18.2	17.5	13.0	9.0	
Nº5C3		4.5	13.9	10.8	8.7	6.3	Nº5D3		5.6	14.0	18.0	15.8	12.2	
Nº6C3		4.0	9.1	8.8	7.5	5.4	Nº6D3		4.9	16.6	15.5	11.9	9.7	
Nº7C3		3.6	12.3	10.6	8.2	4.8								
Nº8C3		3.8	13.7	12.0	9.0	6.2								
means		4.0	12.8	11.2	8.1	5.5			5.3	15.8	17.2	13.8	10.7	
SD		0.3	2.2	1.4	0.6	0.6			0.5	2.0	1.5	1.6	1.5	

	Glucose, mM					C8		D8						
	0'	15'	30'	60'	120'			0'	15'	30'	60'	120'		
Nº1C8		3.7	13.6	13.8	9.0	6.7	Nº1D8		5.7	18.7	19.8	15.6	13.3	
Nº2C8		4.0	10.5	12.5	8.9	4.5	Nº2D8		6.4	19.5	17.7	12.4	10.4	
Nº3C8		4.8	16.8	13.3	7.4	5.5	Nº3D8		7.5	13.8	14.3	10.9	8.9	
Nº4C8		3.8	11.9	9.8	6.9	6.0	Nº4D8		6.0	21.6	19.0	16.1	12.6	
Nº5C8		5.1	12.6	13.3	8.3	6.3	Nº5D8		7.4	17.0	18.1	14.0	11.7	
Nº6C8		3.7	15.4	14.0	7.7	4.9	Nº6D8		4.9	12.9	13.6	10.5	9.0	
Nº7C8		3.8	9.0	9.1	7.2	4.1	Nº7D8		6.6	14.2	14.3	11.8	10.8	
Nº8C8		4.0	16.9	12.1	8.1	5.9	Nº8D8		6.6	19.4	22.2	18.1	14.1	
Nº9C8		3.3	9.8	9.0	6.3	5.7	Nº9D8		5.1	17.1	16.9	14.9	11.0	
Nº10C8		4.5	14.7	11.1	9.0	6.4	Nº10D8		7.3	20.0	19.7	17.3	13.7	
means		4.1	13.1	11.8	7.9	5.6			6.4	17.4	17.6	14.2	11.6	
SD		0.6	2.8	1.9	0.9	0.9			0.9	2.9	2.8	2.7	1.9	

	Glucose, mM					C18		D18						
	0'	15'	30'	60'	120'			0'	15'	30'	60'	120'		
Nº1C18		4.6	15.1	16.2	9.8	7.9	Nº1D18		5.6	17.9	15.8	13.4	12.0	
Nº2C18		3.4	13.6	12.7	7.0	4.8	Nº2D18		5.6	14.7	13.9	9.7	7.9	
Nº3C18		5.0	15.8	14.4	9.4	6.9	Nº3D18		6.4	16.8	16.7	14.9	12.8	
Nº4C18		3.6	11.9	12.1	9.6	6.6	Nº4D18		6.8	18.3	16.5	14.6	10.1	
Nº5C18		4.2	14.4	12.0	7.7	5.2	Nº5D18		5.9	13.7	12.5	11.4	9.8	
Nº6C18		3.5	11.7	11.8	8.5	4.8	Nº6D18		5.3	12.0	11.4	9.9	8.6	
Nº7C18		4.0	14.6	15.3	10.3	7.8	Nº7D18		6.4	16.1	15.8	14.0	12.3	
means		4.0	13.9	13.5	8.9	6.3			6.0	15.6	14.7	12.6	10.5	
SD		0.6	1.6	1.8	1.2	1.4			0.5	2.3	2.1	2.2	1.9	

0'	15'	30'	60'	120'
0.000246297	0.018979	1.48989E-05	0.000145002	0.000155169
7.78615E-06	0.003778	6.59743E-05	2.07762E-05	5.78714E-07
3.58998E-05	0.120952	0.28648807	0.003622318	0.000607154