

Supplemental Table S3

Quantitative Estimation of Target Metabolites (Total Amount)

Metabolite	KEGG ID	Concentration (pmol/10 ⁶ cells)										Comparative Analysis	
		DMSO 48hr			FK866 48hr			DMSO 48hr		FK866 48hr		FK866 48hr vs DMSO 48hr	
		MGG4-1	MGG4-2	MGG4-3	MGG4FK-1	MGG4FK-2	MGG4FK-3	Mean	S.D.	Mean	S.D.	Ratio [†]	p-value [‡]
2-Phosphoglyceric acid	C00631	4.6	2.3	3.0	5.2	4.3	3.8	3.3	1.2	4.4	0.7	1.3	0.239
3-Phosphoglyceric acid	C00197	16	19	20	24	21	23	18	2.0	23	1.7	1.3	0.039
ATP	C00002	4,833	4,411	4,790	1,060	1,078	1,061	4,678	232	1,066	10	0.2	0.001
Dihydroxyacetone phosphate	C00111	84	87	99	2,360	2,709	2,660	90	8.0	2,576	189	29	0.002
Fructose 1,6-diphosphate	C00354	121	105	123	9,597	9,872	9,716	116	9.8	9,728	138	84	6.3E-05
Fructose 6-phosphate	C05345,C00085	7.2	7.8	10	46	44	45	8.5	1.7	45	1.0	5.3	3.9E-05
Glucose 6-phosphate	C00668,C01172,C00092	27	22	25	161	163	154	25	2.5	159	4.5	6.5	1.8E-05
Glyceraldehyde 3-phosphate	C00118,C00661	28	28	35	424	483	442	31	3.6	449	30	15	0.002
Lactic acid	C00186,C00256,C01432	6,228	5,330	5,430	2,304	3,110	2,823	5,663	493	2,746	409	0.5	0.002
Phosphoenolpyruvic acid	C00074	17	14	14	15	13	9.7	15	1.4	12	2.6	0.8	0.225
Pyruvic acid	C00022	74	61	72	44	53	52	69	7.0	50	5.0	0.7	0.022

ID consists of analysis mode and number. 'A' showed anion mode.

N.D. (Not Detected): The target peak or metabolite was below detection limits.

N.A. (Not Available): The calculation was impossible because of insufficiency of the data.