

Table S1. Target compounds and selected instrumental parameters for quantification of each compound by UPLC-(ESI+)MS/MS.

Compound	Precursor > product ion (m/z)	Cone voltage (V)	Collision energy (eV)	Internal Standard used
CML	<u>205.2 > 84.2</u>	20	20	CML-d2
	<i>205.2 > 130.2</i>	20	12	
CEL	<u>219.2 > 84.2</u>	20	20	CEL-d4
	<i>219.2 > 130.2</i>	20	10	
MG-H1	<u>229.2 > 114.1</u>	20	13	MG-H1-d3
	<i>229.2 > 166.2</i>	20	13	
Pentosidine	<u>379.4 > 135.1</u>	20	40	MG-H1-d3
	<i>379.4 > 187.1</i>	20	35	

Note: the underlined precursor > product ion transition was used for quantification, the *italic* precursor > product ion transition was used for confirmation.

Table S2. Instrumental detection and quantitation limits (pg on column) and method detection and quantitation limits (pg on column)

Compound	IDL (pg on column)	IQL (pg on column)	MDL (µg/ml)	MLQ (µg/ml)
CML	0.5	1.5	0.001	0.003
CEL	1.0	3.5	0.01	0.03
MG-H1	1.7	5.7	0.03	0.1
Pentosidine	6.4	21	1.2	4