

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071423\
 Data File : BM040899.D
 Acq On : 15 Jul 2023 04:20
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4179

Quant Time: Jul 15 05:22:17 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 15 03:58:02 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	3431	0.400	ng/ul	0.00
4) Naphthalene-d8	10.658	136	11256	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.504	164	4669	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.257	188	7740	0.400	ng/ul	0.00
17) Chrysene-d12	21.448	240	5598	0.400	ng/ul	0.00
23) Perylene-d12	23.828	264	5899	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.284	96	2202	0.409	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.253	152	5950	0.377	ng/ul	0.00
18) Fluoranthene-d10	19.287	212	7135	0.398	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.318	88	2317	0.422	ng/ul#	90
5) Naphthalene	10.708	128	12321	0.397	ng/ul	98
7) 2-Methylnaphthalene	12.330	142	6785	0.372	ng/ul	99
8) 1-Methylnaphthalene	12.544	142	6810	0.374	ng/ul	98
10) Acenaphthylene	14.227	152	7847	0.376	ng/ul	98
11) Acenaphthene	14.569	153	6752	0.383	ng/ul	98
12) Fluorene	15.559	166	6630	0.353	ng/ul	99
14) Pentachlorophenol	16.919	266	648	0.435	ng/ul	90
15) Phenanthrene	17.299	178	9190	0.382	ng/ul	99
16) Anthracene	17.392	178	7377	0.365	ng/ul	98
19) Fluoranthene	19.320	202	9048	0.381	ng/ul	98
20) Pyrene	19.682	202	9493	0.372	ng/ul	99
21) Benzo(a)anthracene	21.434	228	7374	0.389	ng/ul	100
22) Chrysene	21.483	228	8173	0.378	ng/ul	99
24) Benzo(b)fluoranthene	23.100	252	8607	0.377	ng/ul	98
25) Benzo(k)fluoranthene	23.149	252	8375	0.372	ng/ul	96
26) Benzo(a)pyrene	23.719	252	7610	0.370	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.280	276	11010	0.384	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.297	278	8813	0.393	ng/ul	99
29) Benzo(g,h,i)perylene	27.038	276	9575	0.381	ng/ul	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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