

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040424\
 Data File : BM044950.D
 Acq On : 05 Apr 2024 15:32
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4063

Quant Time: Apr 05 16:01:18 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM040424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 03 23:34:03 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.644	152	2083	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.430	136	5961	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.297	164	3207	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.056	188	7531	0.400	ng/ul	-0.01
17) Chrysene-d12	21.263	240	7375	0.400	ng/ul	0.00
23) Perylene-d12	23.490	264	9619	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	1119	0.409	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.030	152	3232	0.400	ng/ul	-0.02
18) Fluoranthene-d10	19.094	212	7772	0.417	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.251	88	1066	0.406	ng/ul	95
5) Naphthalene	10.479	128	6448	0.408	ng/ul	98
7) 2-Methylnaphthalene	12.107	142	3791	0.401	ng/ul	100
8) 1-Methylnaphthalene	12.321	142	4150	0.402	ng/ul	100
10) Acenaphthylene	14.015	152	6459	0.404	ng/ul	99
11) Acenaphthene	14.358	153	4642	0.405	ng/ul	99
12) Fluorene	15.361	166	5148	0.401	ng/ul	99
14) Pentachlorophenol	16.722	266	743	0.421	ng/ul	98
15) Phenanthrene	17.098	178	8392	0.396	ng/ul	100
16) Anthracene	17.199	178	8279	0.396	ng/ul	99
19) Fluoranthene	19.122	202	10750	0.408	ng/ul	99
20) Pyrene	19.489	202	11411	0.410	ng/ul	100
21) Benzo(a)anthracene	21.249	228	10820	0.421	ng/ul	99
22) Chrysene	21.298	228	12295	0.384	ng/ul	100
24) Benzo(b)fluoranthene	22.821	252	13467	0.388	ng/ul	98
25) Benzo(k)fluoranthene	22.865	252	14632	0.376	ng/ul	99
26) Benzo(a)pyrene	23.394	252	13070	0.394	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.736	276	17668	0.389	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.753	278	13985	0.388	ng/ul	99
29) Benzo(g,h,i)perylene	26.413	276	14779	0.377	ng/ul	98

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

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