

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110723\
 Data File : BM042631.D
 Acq On : 08 Nov 2023 03:09
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4126

Quant Time: Nov 08 03:40:14 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM110623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 07 02:23:03 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.868	152	1007	0.400	ng/ul	0.00
4) Naphthalene-d8	10.678	136	2394	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.511	164	1431	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.268	188	2858	0.400	ng/ul	0.00
17) Chrysene-d12	21.449	240	2084	0.400	ng/ul	0.00
23) Perylene-d12	23.826	264	2562	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.206	96	628	0.472	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.267	152	1341	0.401	ng/ul	0.00
18) Fluoranthene-d10	19.295	212	2973	0.412	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.244	88	728	0.460	ng/ul#	76
5) Naphthalene	10.728	128	3192	0.413	ng/ul	100
7) 2-Methylnaphthalene	12.344	142	1983	0.400	ng/ul	94
8) 1-Methylnaphthalene	12.559	142	2035	0.405	ng/ul	99
10) Acenaphthylene	14.243	152	3829	0.391	ng/ul	99
11) Acenaphthene	14.576	153	2687	0.404	ng/ul	99
12) Fluorene	15.566	166	3018	0.396	ng/ul	99
14) Pentachlorophenol	16.913	266	537	0.368	ng/ul	98
15) Phenanthrene	17.310	178	4516	0.408	ng/ul	100
16) Anthracene	17.403	178	4062	0.405	ng/ul	99
19) Fluoranthene	19.323	202	5245	0.403	ng/ul	99
20) Pyrene	19.690	202	5671	0.396	ng/ul	98
21) Benzo(a)anthracene	21.432	228	3502	0.378	ng/ul	100
22) Chrysene	21.485	228	3824	0.399	ng/ul	100
24) Benzo(b)fluoranthene	23.092	252	4183	0.404	ng/ul	97
25) Benzo(k)fluoranthene	23.145	252	4217	0.389	ng/ul	96
26) Benzo(a)pyrene	23.724	252	4344	0.405	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.319	276	5344	0.384	ng/ul	94
28) Dibenzo(a,h)anthracene	26.332	278	3888	0.391	ng/ul	94
29) Benzo(g,h,i)perylene	27.087	276	4880	0.377	ng/ul	99

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

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