

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112423\
 Data File : BM043001.D
 Acq On : 25 Nov 2023 04:49
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4134

Quant Time: Nov 25 06:01:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM112423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 24 15:22:31 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.880	152	698	0.400	ng/ul	0.00
4) Naphthalene-d8	10.656	136	2262	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.470	164	1295	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.243	188	2588	0.400	ng/ul	0.00
17) Chrysene-d12	21.423	240	1722	0.400	ng/ul	0.00
23) Perylene-d12	23.779	264	2289	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.168	96	479	0.447	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.262	152	1196	0.395	ng/ul	0.01
18) Fluoranthene-d10	19.267	212	2373	0.378	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.206	88	475	0.413	ng/ul#	91
5) Naphthalene	10.705	128	2438	0.397	ng/ul	96
7) 2-Methylnaphthalene	12.328	142	1416	0.391	ng/ul	99
8) 1-Methylnaphthalene	12.526	142	1576	0.390	ng/ul	100
10) Acenaphthylene	14.206	152	2601	0.394	ng/ul	100
11) Acenaphthene	14.534	153	2046	0.409	ng/ul	98
12) Fluorene	15.543	166	2024	0.394	ng/ul	96
14) Pentachlorophenol	16.913	266	201	0.287	ng/ul	93
15) Phenanthrene	17.285	178	2845	0.385	ng/ul	99
16) Anthracene	17.399	178	2740	0.392	ng/ul	98
19) Fluoranthene	19.295	202	3703	0.379	ng/ul	97
20) Pyrene	19.662	202	3933	0.379	ng/ul	100
21) Benzo(a)anthracene	21.414	228	2226	0.398	ng/ul	98
22) Chrysene	21.458	228	4171	0.401	ng/ul	99
24) Benzo(b)fluoranthene	23.048	252	3165	0.410	ng/ul	96
25) Benzo(k)fluoranthene	23.098	252	4294	0.404	ng/ul	97
26) Benzo(a)pyrene	23.680	252	3286	0.390	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.245	276	4569	0.400	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.258	278	3493	0.396	ng/ul	95
29) Benzo(g,h,i)perylene	27.003	276	4536	0.409	ng/ul	97

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

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