

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111624\
 Data File : BM048736.D
 Acq On : 16 Nov 2024 11:35
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4086

Quant Time: Nov 17 22:31:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM110624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Nov 17 22:31:02 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.679	152	2746	0.400	ng/ul	0.00
4) Naphthalene-d8	10.443	136	9502	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.286	164	5285	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.045	188	9710	0.400	ng/ul	0.00
17) Chrysene-d12	21.228	240	8208	0.400	ng/ul	0.00
23) Perylene-d12	23.435	264	8231	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	1723	0.520	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.048	152	4823	0.393	ng/ul	0.00
18) Fluoranthene-d10	19.073	212	8794	0.382	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.127	88	1896	0.498	ng/ul#	72
5) Naphthalene	10.492	128	10328	0.434	ng/ul	99
7) 2-Methylnaphthalene	12.120	142	5990	0.400	ng/ul	95
8) 1-Methylnaphthalene	12.323	142	6640	0.433	ng/ul	97
10) Acenaphthylene	14.013	152	9147	0.456	ng/ul	99
11) Acenaphthene	14.350	153	7005	0.436	ng/ul	99
12) Fluorene	15.354	166	7281	0.417	ng/ul	98
14) Pentachlorophenol	16.720	266	987	0.277	ng/ul	99
15) Phenanthrene	17.087	178	9991	0.392	ng/ul	99
16) Anthracene	17.193	178	9061	0.382	ng/ul	99
19) Fluoranthene	19.101	202	12111	0.398	ng/ul	99
20) Pyrene	19.463	202	12929	0.392	ng/ul	99
21) Benzo(a)anthracene	21.217	228	9303	0.346	ng/ul	98
22) Chrysene	21.264	228	13376	0.415	ng/ul	99
24) Benzo(b)fluoranthene	22.771	252	10581	0.413	ng/ul	98
25) Benzo(k)fluoranthene	22.815	252	13190	0.439	ng/ul	99
26) Benzo(a)pyrene	23.341	252	10622	0.424	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.662	276	15179	0.434	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.676	278	11576	0.425	ng/ul	98
29) Benzo(g,h,i)perylene	26.340	276	12874	0.439	ng/ul	99

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

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