

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110723\
 Data File : BM042647.D
 Acq On : 08 Nov 2023 13:46
 Operator : MA/JU
 Sample : PB156699BS
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS699

Quant Time: Nov 08 22:54:33 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	863	0.400	ng/ul	0.00
4) Naphthalene-d8	10.678	136	1964	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.511	164	1075	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.268	188	2066	0.400	ng/ul	0.00
17) Chrysene-d12	21.450	240	1580	0.400	ng/ul	0.00
23) Perylene-d12	23.832	264	1889	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.202	96	999	0.876	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.267	152	1074	0.392	ng/ul	0.00
18) Fluoranthene-d10	19.295	212	2251	0.411	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.235	88	3227	2.381	ng/ul#	58
5) Naphthalene	10.728	128	2653	0.418	ng/ul	98
7) 2-Methylnaphthalene	12.344	142	1575	0.387	ng/ul	94
8) 1-Methylnaphthalene	12.559	142	1553	0.377	ng/ul	98
10) Acenaphthylene	14.243	152	2936	0.399	ng/ul	98
11) Acenaphthene	14.576	153	2074	0.416	ng/ul	99
12) Fluorene	15.566	166	2312	0.404	ng/ul	99
14) Pentachlorophenol	16.905	266	946	0.898	ng/ul	96
15) Phenanthrene	17.306	178	3328	0.415	ng/ul	98
16) Anthracene	17.403	178	2887	0.398	ng/ul	96
19) Fluoranthene	19.323	202	3899	0.395	ng/ul	98
20) Pyrene	19.690	202	4282	0.394	ng/ul	97
21) Benzo(a)anthracene	21.432	228	2713	0.386	ng/ul	98
22) Chrysene	21.485	228	3125	0.430	ng/ul	99
24) Benzo(b)fluoranthene	23.095	252	3279	0.430	ng/ul	93
25) Benzo(k)fluoranthene	23.142	252	3513	0.439	ng/ul#	94
26) Benzo(a)pyrene	23.724	252	3661	0.463	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.315	276	4546	0.443	ng/ul#	93
28) Dibenzo(a,h)anthracene	26.332	278	3289	0.449	ng/ul	95
29) Benzo(g,h,i)perylene	27.087	276	4184	0.438	ng/ul	98

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

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