

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112423\
 Data File : BM042998.D
 Acq On : 25 Nov 2023 03:00
 Operator : MA/JU
 Sample : PB157190BS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS190

Quant Time: Nov 25 03:37:31 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	7.906	152	600	0.400	ng/ul	0.03
4) Naphthalene-d8	10.662	136	2174	0.400	ng/ul	0.01
9) Acenaphthene-d10	14.470	164	1366	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.238	188	2952	0.400	ng/ul	0.00
17) Chrysene-d12	21.420	240	1886	0.400	ng/ul	0.00
23) Perylene-d12	23.779	264	2183	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.168	96	711	0.772	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.267	152	1019	0.350	ng/ul	0.02
18) Fluoranthene-d10	19.267	212	2437	0.355	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.202	88	2025	2.048	ng/ul#	86
5) Naphthalene	10.717	128	1871	0.317	ng/ul	96
7) 2-Methylnaphthalene	12.350	142	1087	0.312	ng/ul	97
8) 1-Methylnaphthalene	12.526	142	1165	0.300	ng/ul	98
10) Acenaphthylene	14.206	152	2170	0.311	ng/ul	100
11) Acenaphthene	14.534	153	1676	0.317	ng/ul	99
12) Fluorene	15.543	166	1776	0.327	ng/ul#	96
14) Pentachlorophenol	16.892	266	522	0.653	ng/ul	97
15) Phenanthrene	17.281	178	2704	0.321	ng/ul	98
16) Anthracene	17.395	178	2591	0.325	ng/ul	97
19) Fluoranthene	19.295	202	3549	0.332	ng/ul	100
20) Pyrene	19.662	202	3784	0.333	ng/ul	99
21) Benzo(a)anthracene	21.414	228	2110	0.345	ng/ul	99
22) Chrysene	21.458	228	3851	0.338	ng/ul	99
24) Benzo(b)fluoranthene	23.051	252	2768	0.376	ng/ul	99
25) Benzo(k)fluoranthene	23.098	252	3672	0.362	ng/ul	99
26) Benzo(a)pyrene	23.683	252	2716	0.338	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.248	276	3433	0.316	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.275	278	2623	0.312	ng/ul	98
29) Benzo(g,h,i)perylene	27.010	276	3396	0.321	ng/ul	98

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

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