

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121024\
 Data File : BM048925.D
 Acq On : 10 Dec 2024 16:09
 Operator : RC/JU
 Sample : PB165458BS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS458

Quant Time: Dec 10 17:08:11 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM120924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 09 17:52:26 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.586	152	3096	0.400	ng/ul	0.00
4) Naphthalene-d8	10.352	136	8306	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.219	164	4528	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.967	188	9306	0.400	ng/ul	0.00
17) Chrysene-d12	21.161	240	7336	0.400	ng/ul	0.00
23) Perylene-d12	23.329	264	7554	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.160	96	2515	0.695	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.953	152	3774	0.356	ng/ul	0.00
18) Fluoranthene-d10	19.001	212	7928	0.362	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.193	88	7336	1.752	ng/ul#	81
5) Naphthalene	10.402	128	7312	0.343	ng/ul	97
7) 2-Methylnaphthalene	12.024	142	4250	0.327	ng/ul	99
8) 1-Methylnaphthalene	12.244	142	4423	0.334	ng/ul	100
10) Acenaphthylene	13.936	152	6360	0.333	ng/ul	99
11) Acenaphthene	14.283	153	4899	0.338	ng/ul	98
12) Fluorene	15.278	166	5479	0.341	ng/ul	98
14) Pentachlorophenol	16.621	266	1081	0.564	ng/ul	94
15) Phenanthrene	17.005	178	8658	0.339	ng/ul	98
16) Anthracene	17.102	178	7531	0.331	ng/ul	98
19) Fluoranthene	19.029	202	10259	0.337	ng/ul	99
20) Pyrene	19.391	202	10679	0.326	ng/ul	97
21) Benzo(a)anthracene	21.146	228	8156	0.316	ng/ul	99
22) Chrysene	21.196	228	10821	0.370	ng/ul	99
24) Benzo(b)fluoranthene	22.683	252	9379	0.359	ng/ul	98
25) Benzo(k)fluoranthene	22.727	252	11103	0.368	ng/ul	98
26) Benzo(a)pyrene	23.233	252	8073	0.334	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.484	276	11746	0.346	ng/ul	100
28) Dibenzo(a,h)anthracene	25.497	278	8924	0.338	ng/ul	96
29) Benzo(g,h,i)perylene	26.135	276	10068	0.369	ng/ul	98

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

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