

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101024\
 Data File : BM048005.D
 Acq On : 11 Oct 2024 09:18
 Operator : RC/JU
 Sample : PB163729BS
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS729

Quant Time: Oct 11 09:50:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM092524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 10 18:01:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.796	152	7670	0.400	ng/ul	0.00
4) Naphthalene-d8	10.583	136	23007	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.422	164	13161	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.165	188	21915	0.400	ng/ul	0.00
17) Chrysene-d12	21.333	240	17639	0.400	ng/ul	0.00
23) Perylene-d12	23.598	264	17201	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.273	96	6767	0.687	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.178	152	8912	0.290	ng/ul	0.00
18) Fluoranthene-d10	19.187	212	16982	0.314	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.307	88	19043	1.615	ng/ul#	83
5) Naphthalene	10.633	128	18355	0.293	ng/ul	99
7) 2-Methylnaphthalene	12.255	142	10245	0.272	ng/ul	99
8) 1-Methylnaphthalene	12.469	142	14056	0.356	ng/ul	100
10) Acenaphthylene	14.145	152	16081	0.251	ng/ul	99
11) Acenaphthene	14.487	153	12832	0.292	ng/ul	98
12) Fluorene	15.477	166	12400	0.250	ng/ul	99
14) Pentachlorophenol	16.823	266	3747	0.569	ng/ul	100
15) Phenanthrene	17.207	178	18233	0.280	ng/ul	100
16) Anthracene	17.305	178	12376	0.214	ng/ul	98
19) Fluoranthene	19.219	202	21985	0.296	ng/ul	96
20) Pyrene	19.577	202	23792	0.301	ng/ul	99
21) Benzo(a)anthracene	21.318	228	16823	0.226	ng/ul	99
22) Chrysene	21.368	228	22208	0.286	ng/ul	99
24) Benzo(b)fluoranthene	22.914	252	19360	0.259	ng/ul	98
25) Benzo(k)fluoranthene	22.958	252	23013	0.306	ng/ul	97
26) Benzo(a)pyrene	23.502	252	17632	0.299	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.913	276	23826	0.254	ng/ul#	93
28) Dibenzo(a,h)anthracene	25.933	278	17035	0.240	ng/ul	94
29) Benzo(g,h,i)perylene	26.608	276	21510	0.285	ng/ul	98

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

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