

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110424\
 Data File : BM048425.D
 Acq On : 04 Nov 2024 14:09
 Operator : RC/JU
 Sample : PB164424BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS424

Quant Time: Nov 04 14:54:39 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM103124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 31 14:58:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.729	152	4194	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.518	136	13287	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.359	164	7110	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.116	188	14013	0.400	ng/ul	0.00
17) Chrysene-d12	21.295	240	10548	0.400	ng/ul	0.00
23) Perylene-d12	23.536	264	10036	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.168	96	4146	0.805	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.130	152	6376	0.377	ng/ul	-0.01
18) Fluoranthene-d10	19.137	212	12984	0.485	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.202	88	10663	1.908	ng/ul#	83
5) Naphthalene	10.568	128	12759	0.367	ng/ul	99
7) 2-Methylnaphthalene	12.207	142	7228	0.350	ng/ul	99
8) 1-Methylnaphthalene	12.410	142	8069	0.366	ng/ul	97
10) Acenaphthylene	14.086	152	10153	0.369	ng/ul	99
11) Acenaphthene	14.423	153	8302	0.377	ng/ul	99
12) Fluorene	15.427	166	8808	0.360	ng/ul	99
14) Pentachlorophenol	16.782	266	1992	0.511	ng/ul	92
15) Phenanthrene	17.158	178	12756	0.366	ng/ul	99
16) Anthracene	17.264	178	11308	0.347	ng/ul	99
19) Fluoranthene	19.169	202	16562	0.455	ng/ul	99
20) Pyrene	19.532	202	17606	0.451	ng/ul	99
21) Benzo(a)anthracene	21.286	228	11049	0.350	ng/ul	100
22) Chrysene	21.327	228	19212	0.442	ng/ul	99
24) Benzo(b)fluoranthene	22.861	252	12899	0.376	ng/ul	97
25) Benzo(k)fluoranthene	22.905	252	16910	0.416	ng/ul	96
26) Benzo(a)pyrene	23.451	252	12761	0.396	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	25.829	276	16418	0.365	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.859	278	12659	0.360	ng/ul	94
29) Benzo(g,h,i)perylene	26.510	276	14556	0.383	ng/ul	99

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

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