

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120524\
 Data File : BM048871.D
 Acq On : 06 Dec 2024 16:56
 Operator : RC/JU
 Sample : PB165412BS
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS412

Quant Time: Dec 06 16:24:35 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM120324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 06 00:05:00 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.594	152	4017	0.400	ng/ul	0.00
4) Naphthalene-d8	10.359	136	11392	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.229	164	6442	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.981	188	11975	0.400	ng/ul	0.00
17) Chrysene-d12	21.175	240	7362	0.400	ng/ul	0.00
23) Perylene-d12	23.349	264	7463	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	3269	0.637	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.959	152	4717	0.344	ng/ul	0.00
18) Fluoranthene-d10	19.011	212	7905	0.348	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.198	88	9138	1.505	ng/ul#	81
5) Naphthalene	10.409	128	9091	0.316	ng/ul	99
7) 2-Methylnaphthalene	12.036	142	5430	0.326	ng/ul	97
8) 1-Methylnaphthalene	12.256	142	5662	0.322	ng/ul	100
10) Acenaphthylene	13.947	152	8243	0.292	ng/ul	99
11) Acenaphthene	14.294	153	6260	0.307	ng/ul	98
12) Fluorene	15.289	166	6588	0.312	ng/ul	98
14) Pentachlorophenol	16.635	266	978	0.944	ng/ul#	8
15) Phenanthrene	17.019	178	9591	0.316	ng/ul	99
16) Anthracene	17.112	178	8184	0.303	ng/ul	98
19) Fluoranthene	19.039	202	10385	0.307	ng/ul	98
20) Pyrene	19.407	202	10664	0.303	ng/ul	99
21) Benzo(a)anthracene	21.160	228	6773	0.326	ng/ul	99
22) Chrysene	21.210	228	9756	0.297	ng/ul	100
24) Benzo(b)fluoranthene	22.703	252	7845	0.374	ng/ul	99
25) Benzo(k)fluoranthene	22.747	252	9720	0.331	ng/ul	98
26) Benzo(a)pyrene	23.253	252	7161	0.333	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.514	276	9556	0.336	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.531	278	7246	0.341	ng/ul	98
29) Benzo(g,h,i)perylene	26.168	276	8178	0.332	ng/ul	97

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

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