

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121024\
 Data File : BM048924.D
 Acq On : 10 Dec 2024 15:33
 Operator : RC/JU
 Sample : PB165460BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS460

Quant Time: Dec 10 16:15:49 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	7.585	152	2987	0.400	ng/ul	0.00
4) Naphthalene-d8	10.348	136	8049	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.220	164	4408	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.964	188	9107	0.400	ng/ul	0.00
17) Chrysene-d12	21.160	240	7103	0.400	ng/ul	0.00
23) Perylene-d12	23.329	264	7347	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.159	96	2906	0.832	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.948	152	4460	0.435	ng/ul	0.00
18) Fluoranthene-d10	18.997	212	9355	0.441	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.193	88	8359	2.069	ng/ul#	81
5) Naphthalene	10.397	128	8547	0.414	ng/ul	99
7) 2-Methylnaphthalene	12.025	142	5089	0.404	ng/ul	98
8) 1-Methylnaphthalene	12.245	142	5283	0.412	ng/ul	98
10) Acenaphthylene	13.938	152	7539	0.405	ng/ul	99
11) Acenaphthene	14.280	153	5796	0.411	ng/ul	99
12) Fluorene	15.275	166	6515	0.416	ng/ul	100
14) Pentachlorophenol	16.618	266	1239	0.661	ng/ul	96
15) Phenanthrene	17.006	178	10339	0.414	ng/ul	99
16) Anthracene	17.099	178	8926	0.401	ng/ul	97
19) Fluoranthene	19.030	202	12111	0.411	ng/ul	100
20) Pyrene	19.392	202	12672	0.399	ng/ul	98
21) Benzo(a)anthracene	21.146	228	9641	0.386	ng/ul	99
22) Chrysene	21.195	228	12851	0.454	ng/ul	99
24) Benzo(b)fluoranthene	22.685	252	10914	0.430	ng/ul	99
25) Benzo(k)fluoranthene	22.726	252	13367	0.455	ng/ul	98
26) Benzo(a)pyrene	23.235	252	9688	0.413	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.480	276	13931	0.422	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.504	278	10606	0.413	ng/ul	96
29) Benzo(g,h,i)perylene	26.134	276	11710	0.441	ng/ul	98

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

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