

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121024\
 Data File : BM048930.D
 Acq On : 10 Dec 2024 19:09
 Operator : RC/JU
 Sample : PB165472BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS472

Quant Time: Dec 10 22:50:18 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	7.590	152	3420	0.400	ng/ul	0.00
4) Naphthalene-d8	10.352	136	9157	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.218	164	4764	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.962	188	9677	0.400	ng/ul	0.00
17) Chrysene-d12	21.158	240	7592	0.400	ng/ul	0.00
23) Perylene-d12	23.326	264	7815	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.159	96	3023	0.756	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.947	152	4538	0.389	ng/ul	0.00
18) Fluoranthene-d10	18.996	212	9434	0.416	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.193	88	8251	1.784	ng/ul#	81
5) Naphthalene	10.396	128	8796	0.375	ng/ul	99
7) 2-Methylnaphthalene	12.024	142	5077	0.354	ng/ul	100
8) 1-Methylnaphthalene	12.244	142	5278	0.361	ng/ul	99
10) Acenaphthylene	13.936	152	7264	0.361	ng/ul	99
11) Acenaphthene	14.279	153	5659	0.371	ng/ul	98
12) Fluorene	15.273	166	6193	0.366	ng/ul	99
14) Pentachlorophenol	16.621	266	1115	0.559	ng/ul	94
15) Phenanthrene	17.005	178	9982	0.376	ng/ul	99
16) Anthracene	17.098	178	8552	0.362	ng/ul	98
19) Fluoranthene	19.029	202	11995	0.381	ng/ul	99
20) Pyrene	19.391	202	12529	0.369	ng/ul	99
21) Benzo(a)anthracene	21.143	228	9055	0.339	ng/ul	97
22) Chrysene	21.193	228	12346	0.408	ng/ul	99
24) Benzo(b)fluoranthene	22.680	252	10879	0.403	ng/ul	99
25) Benzo(k)fluoranthene	22.724	252	12534	0.401	ng/ul	98
26) Benzo(a)pyrene	23.230	252	9838	0.394	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	25.471	276	13305	0.379	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.497	278	10073	0.369	ng/ul	97
29) Benzo(g,h,i)perylene	26.131	276	11127	0.394	ng/ul	98

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

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