

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122324\
 Data File : BM049189.D
 Acq On : 24 Dec 2024 17:42
 Operator : RC/JU
 Sample : PB165780BS
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS780

Quant Time: Dec 24 21:41:40 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM121824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 18 15:33:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.552	152	5638	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.315	136	17522	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.188	164	10418	0.400	ng/ul	-0.01
13) Phenanthrene-d10	16.939	188	21645	0.400	ng/ul	-0.01
17) Chrysene-d12	21.137	240	18090	0.400	ng/ul	0.00
23) Perylene-d12	23.294	264	19638	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	3388	0.521	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.915	152	7172	0.305	ng/ul	0.00
18) Fluoranthene-d10	18.974	212	15851	0.309	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.176	88	9472	1.246	ng/ul#	77
5) Naphthalene	10.365	128	12514	0.285	ng/ul	99
7) 2-Methylnaphthalene	11.987	142	7819	0.281	ng/ul	99
8) 1-Methylnaphthalene	12.207	142	7997	0.281	ng/ul	100
10) Acenaphthylene	13.905	152	11628	0.269	ng/ul	99
11) Acenaphthene	14.252	153	8625	0.274	ng/ul	99
12) Fluorene	15.247	166	9768	0.275	ng/ul	99
14) Pentachlorophenol	16.597	266	2952	0.493	ng/ul	96
15) Phenanthrene	16.981	178	15369	0.272	ng/ul	100
16) Anthracene	17.074	178	14117	0.272	ng/ul	98
19) Fluoranthene	19.002	202	18707	0.286	ng/ul	99
20) Pyrene	19.365	202	19806	0.284	ng/ul	99
21) Benzo(a)anthracene	21.122	228	16476	0.262	ng/ul	99
22) Chrysene	21.172	228	18874	0.279	ng/ul	99
24) Benzo(b)fluoranthene	22.651	252	17530	0.279	ng/ul	98
25) Benzo(k)fluoranthene	22.694	252	19216	0.279	ng/ul	98
26) Benzo(a)pyrene	23.197	252	16466	0.274	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.427	276	22244	0.261	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.443	278	17314	0.262	ng/ul	99
29) Benzo(g,h,i)perylene	26.074	276	18285	0.267	ng/ul	100

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

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