

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022124\
 Data File : BM044247.D
 Acq On : 22 Feb 2024 19:22
 Operator : MA/JU
 Sample : PB159095BS
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS095

Quant Time: Feb 22 23:35:41 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM021524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 15 13:40:24 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.896	152	4177	0.400	ng/ul	0.00
4) Naphthalene-d8	10.699	136	12076	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.543	164	5183	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.301	188	10723	0.400	ng/ul	0.00
17) Chrysene-d12	21.500	240	9268	0.400	ng/ul	0.00
23) Perylene-d12	23.897	264	10540	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	5427	0.917	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.294	152	7542	0.446	ng/ul	0.00
18) Fluoranthene-d10	19.331	212	14378	0.520	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.382	88	17203	2.910	ng/ul#	73
5) Naphthalene	10.748	128	17220	0.480	ng/ul	99
7) 2-Methylnaphthalene	12.365	142	9919	0.448	ng/ul	98
8) 1-Methylnaphthalene	12.585	142	9724	0.434	ng/ul	99
10) Acenaphthylene	14.265	152	14334	0.503	ng/ul	99
11) Acenaphthene	14.607	153	10136	0.491	ng/ul	98
12) Fluorene	15.602	166	10554	0.453	ng/ul	99
14) Pentachlorophenol	16.946	266	4030	1.050	ng/ul	98
15) Phenanthrene	17.343	178	16878	0.497	ng/ul	99
16) Anthracene	17.436	178	17551	0.522	ng/ul	99
19) Fluoranthene	19.359	202	21903	0.518	ng/ul	97
20) Pyrene	19.726	202	23218	0.523	ng/ul	100
21) Benzo(a)anthracene	21.485	228	19513	0.460	ng/ul	99
22) Chrysene	21.538	228	24406	0.533	ng/ul	99
24) Benzo(b)fluoranthene	23.166	252	21897	0.503	ng/ul	98
25) Benzo(k)fluoranthene	23.218	252	26033	0.551	ng/ul	99
26) Benzo(a)pyrene	23.791	252	20826	0.515	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.386	276	25284	0.491	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.433	278	19636	0.484	ng/ul	97
29) Benzo(g,h,i)perylene	27.138	276	23000	0.505	ng/ul	98

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

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