

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM063023\
 Data File : BM040493.D
 Acq On : 01 Jul 2023 04:25
 Operator : MA/JU
 Sample : PB153636BS
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS636

Quant Time: Jul 01 04:55:53 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 01 03:00:05 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.936	152	7714	0.400	ng/ul	0.00
4) Naphthalene-d8	10.745	136	29497	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.577	164	14351	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.328	188	25242	0.400	ng/ul	0.00
17) Chrysene-d12	21.506	240	14255	0.400	ng/ul	0.00
23) Perylene-d12	23.920	264	14547	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.325	96	5075	0.419	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.334	152	11160	0.270	ng/ul	0.00
18) Fluoranthene-d10	19.352	212	14364	0.314	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.359	88	13753	1.115	ng/ul	89
5) Naphthalene	10.795	128	21811	0.268	ng/ul	99
7) 2-Methylnaphthalene	12.406	142	13051	0.273	ng/ul	99
8) 1-Methylnaphthalene	12.626	142	12777	0.268	ng/ul	99
10) Acenaphthylene	14.300	152	18537	0.289	ng/ul	99
11) Acenaphthene	14.642	153	13813	0.255	ng/ul	99
12) Fluorene	15.627	166	14622	0.253	ng/ul	100
14) Pentachlorophenol	16.982	266	967	0.199	ng/ul	93
15) Phenanthrene	17.370	178	19699	0.251	ng/ul	99
16) Anthracene	17.463	178	18030	0.274	ng/ul	98
19) Fluoranthene	19.384	202	19170	0.317	ng/ul	98
20) Pyrene	19.747	202	19911	0.307	ng/ul	100
21) Benzo(a)anthracene	21.491	228	15063	0.312	ng/ul	96
22) Chrysene	21.544	228	15035	0.273	ng/ul	97
24) Benzo(b)fluoranthene	23.184	252	14960	0.266	ng/ul	84
25) Benzo(k)fluoranthene	23.230	252	14783	0.266	ng/ul#	82
26) Benzo(a)pyrene	23.812	252	13203	0.260	ng/ul#	81
27) Indeno(1,2,3-cd)pyrene	26.424	276	18036	0.255	ng/ul	93
28) Dibenzo(a,h)anthracene	26.441	278	14353	0.260	ng/ul	90
29) Benzo(g,h,i)perylene	27.192	276	15037	0.243	ng/ul	94

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

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