

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070523\
 Data File : BM040664.D
 Acq On : 06 Jul 2023 06:20
 Operator : MA/JU
 Sample : PB153515BS
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS515

Quant Time: Jul 06 07:08:27 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 : Wed Jul 05 09:30:35 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.907	152	6347	0.400	ng/ul	0.00
4) Naphthalene-d8	10.712	136	23385	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.554	164	10391	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.307	188	17086	0.400	ng/ul	0.00
17) Chrysene-d12	21.491	240	11029	0.400	ng/ul	0.00
23) Perylene-d12	23.894	264	13907	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.312	96	5359	0.538	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.307	152	9737	0.297	ng/ul	0.00
18) Fluoranthene-d10	19.333	212	10968	0.310	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.346	88	13861	1.366	ng/ul	93
5) Naphthalene	10.762	128	18935	0.294	ng/ul	99
7) 2-Methylnaphthalene	12.378	142	10906	0.288	ng/ul	98
8) 1-Methylnaphthalene	12.598	142	10457	0.276	ng/ul	99
10) Acenaphthylene	14.276	152	14270	0.307	ng/ul	98
11) Acenaphthene	14.614	153	10968	0.280	ng/ul	98
12) Fluorene	15.604	166	11269	0.269	ng/ul	97
14) Pentachlorophenol	16.961	266	842	0.256	ng/ul	94
15) Phenanthrene	17.349	178	14669	0.276	ng/ul	97
16) Anthracene	17.442	178	12969	0.291	ng/ul	96
19) Fluoranthene	19.366	202	13776	0.294	ng/ul	100
20) Pyrene	19.728	202	14339	0.285	ng/ul	99
21) Benzo(a)anthracene	21.477	228	11849	0.318	ng/ul	97
22) Chrysene	21.526	228	12317	0.289	ng/ul	98
24) Benzo(b)fluoranthene	23.160	252	14289	0.265	ng/ul	85
25) Benzo(k)fluoranthene	23.210	252	14555	0.274	ng/ul#	86
26) Benzo(a)pyrene	23.786	252	13610	0.281	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	26.387	276	18617	0.276	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.397	278	14672	0.278	ng/ul	93
29) Benzo(g,h,i)perylene	27.148	276	15854	0.268	ng/ul	96

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

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