

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061523\
 Data File : BM040265.D
 Acq On : 15 Jun 2023 18:39
 Operator : MA/JU
 Sample : PB153389BS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS389

Quant Time: Jun 15 23:43:31 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM061423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 15 00:31:16 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.974	152	8149	0.400	ng/ul	0.00
4) Naphthalene-d8	10.784	136	30283	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.609	164	15048	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.353	188	28278	0.400	ng/ul	0.00
17) Chrysene-d12	21.529	240	17839	0.400	ng/ul	0.00
23) Perylene-d12	23.955	264	16390	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.350	96	8195	0.814	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.367	152	14872	0.387	ng/ul	0.00
18) Fluoranthene-d10	19.375	212	21881	0.492	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.384	88	19009	1.799	ng/ul#	80
5) Naphthalene	10.833	128	25769	0.359	ng/ul	99
7) 2-Methylnaphthalene	12.439	142	15689	0.352	ng/ul	99
8) 1-Methylnaphthalene	12.659	142	15919	0.341	ng/ul	99
10) Acenaphthylene	14.332	152	19580	0.372	ng/ul	100
11) Acenaphthene	14.670	153	16723	0.371	ng/ul	98
12) Fluorene	15.655	166	17842	0.355	ng/ul	99
14) Pentachlorophenol	17.003	266	3828	0.663	ng/ul	99
15) Phenanthrene	17.396	178	26384	0.355	ng/ul	100
16) Anthracene	17.484	178	22181	0.352	ng/ul	99
19) Fluoranthene	19.407	202	26381	0.434	ng/ul	100
20) Pyrene	19.770	202	26981	0.427	ng/ul	99
21) Benzo(a)anthracene	21.512	228	20703	0.393	ng/ul	99
22) Chrysene	21.567	228	22129	0.398	ng/ul	99
24) Benzo(b)fluoranthene	23.213	252	22544	0.373	ng/ul	92
25) Benzo(k)fluoranthene	23.263	252	21148	0.372	ng/ul#	92
26) Benzo(a)pyrene	23.847	252	18042	0.364	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	26.471	276	23717	0.378	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.487	278	19040	0.378	ng/ul	96
29) Benzo(g,h,i)perylene	27.249	276	20277	0.387	ng/ul	96

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

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