

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM050823\
 Data File : BM039883.D
 Acq On : 09 May 2023 00:25
 Operator : CG/JU
 Sample : PB152658BS
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS658

Quant Time: May 09 02:37:27 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM042723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 09 02:35:28 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	2030	0.400	ng/ul	0.03
4) Naphthalene-d8	10.599	136	9917	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.440	164	6308	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.199	188	12167	0.400	ng/ul	0.00
17) Chrysene-d12	21.394	240	9794	0.400	ng/ul	0.00
23) Perylene-d12	23.741	264	8346	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.176	96	2240	0.749	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.205	152	4856	0.382	ng/ul	0.02
18) Fluoranthene-d10	19.233	212	11350	0.431	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.210	88	5597	1.806	ng/ul#	37
5) Naphthalene	10.649	128	8660	0.348	ng/ul	97
7) 2-Methylnaphthalene	12.277	142	5313	0.370	ng/ul	99
8) 1-Methylnaphthalene	12.480	142	5667	0.368	ng/ul	100
10) Acenaphthylene	14.163	152	9513	0.390	ng/ul	99
11) Acenaphthene	14.500	153	7161	0.373	ng/ul	98
12) Fluorene	15.504	166	7826	0.368	ng/ul	100
14) Pentachlorophenol	16.886	266	1582	0.597	ng/ul	99
15) Phenanthrene	17.241	178	11823	0.369	ng/ul	99
16) Anthracene	17.346	178	10030	0.364	ng/ul	99
19) Fluoranthene	19.265	202	14457	0.414	ng/ul	99
20) Pyrene	19.628	202	15798	0.419	ng/ul	99
21) Benzo(a)anthracene	21.379	228	11187	0.397	ng/ul	98
22) Chrysene	21.432	228	12615	0.396	ng/ul	99
24) Benzo(b)fluoranthene	23.028	252	10780	0.368	ng/ul	95
25) Benzo(k)fluoranthene	23.074	252	11081	0.371	ng/ul	95
26) Benzo(a)pyrene	23.642	252	9665	0.367	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.181	276	11987	0.360	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.185	278	9318	0.361	ng/ul	96
29) Benzo(g,h,i)perylene	26.922	276	10296	0.355	ng/ul	98

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

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