

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM050523\
 Data File : BM039843.D
 Acq On : 05 May 2023 13:19
 Operator : CG/JU
 Sample : PB152462BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS462

Quant Time: May 06 02:20:49 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 : Sat May 06 02:18:19 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.867	152	1462	0.400	ng/ul	0.02
4) Naphthalene-d8	10.617	136	6967	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.446	164	4381	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.208	188	8555	0.400	ng/ul	0.00
17) Chrysene-d12	21.398	240	9343	0.400	ng/ul	0.00
23) Perylene-d12	23.742	264	10203	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.176	96	1568	0.728	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.223	152	3053	0.342	ng/ul	0.00
18) Fluoranthene-d10	19.238	212	8402	0.334	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.210	88	3826	1.714	ng/ul#	40
5) Naphthalene	10.666	128	5514	0.315	ng/ul#	92
7) 2-Methylnaphthalene	12.300	142	3201	0.317	ng/ul	97
8) 1-Methylnaphthalene	12.498	142	3555	0.328	ng/ul	100
10) Acenaphthylene	14.173	152	6266	0.370	ng/ul	98
11) Acenaphthene	14.511	153	4656	0.350	ng/ul	100
12) Fluorene	15.515	166	4970	0.336	ng/ul	99
14) Pentachlorophenol	16.908	266	1048	0.562	ng/ul	99
15) Phenanthrene	17.250	178	7826	0.347	ng/ul	97
16) Anthracene	17.356	178	6487	0.335	ng/ul	98
19) Fluoranthene	19.266	202	10463	0.314	ng/ul	98
20) Pyrene	19.634	202	11934	0.332	ng/ul	99
21) Benzo(a)anthracene	21.386	228	9485	0.353	ng/ul	98
22) Chrysene	21.436	228	11302	0.372	ng/ul	98
24) Benzo(b)fluoranthene	23.029	252	11795	0.329	ng/ul	94
25) Benzo(k)fluoranthene	23.078	252	12203	0.334	ng/ul	94
26) Benzo(a)pyrene	23.646	252	11084	0.345	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.176	276	14038	0.345	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.189	278	11039	0.350	ng/ul	95
29) Benzo(g,h,i)perylene	26.927	276	11986	0.338	ng/ul	98

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

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