

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040623\
 Data File : BM039333.D
 Acq On : 07 Apr 2023 01:32
 Operator : CG/JU
 Sample : 02092-10
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB975

Quant Time: Apr 07 02:43:01 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM040623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 06 23:58:10 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	5896	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.691	136	22867	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.532	164	14742	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.282	188	29319	0.400	ng/ul	0.00
17) Chrysene-d12	21.469	240	19632	0.400	ng/ul	0.00
23) Perylene-d12	23.857	264	21815	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	32242	2.954	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.286	152	9463	0.358	ng/ul	0.00
18) Fluoranthene-d10	19.310	212	26006	0.425	ng/ul	0.00
Target Compounds						
					Qvalue	
11) Acenaphthene	14.592	153	3273	0.073	ng/ul	96
12) Fluorene	15.582	166	1604	0.036	ng/ul#	93
15) Phenanthrene	17.324	178	3991	0.049	ng/ul#	77
16) Anthracene	17.413	178	2043	0.033	ng/ul#	58
19) Fluoranthene	19.343	202	8562	0.105	ng/ul#	90
20) Pyrene	19.705	202	6869	0.074	ng/ul#	79
21) Benzo(a)anthracene	21.454	228	2496	0.055	ng/ul#	56
22) Chrysene	21.507	228	1787	0.032	ng/ul#	51
24) Benzo(b)fluoranthene	23.129	252	2369	0.034	ng/ul#	1

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

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