

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
 Data File : BM044969.D
 Acq On : 06 Apr 2024 14:16
 Operator : MA/JU
 Sample : P2018-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Apr 08 01:50:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 08 01:45:36 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	71725	20.000	ng/ul	0.00
20) Naphthalene-d8	10.410	136	311046	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.280	164	195574	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.045	188	415969	20.000	ng/ul	0.00
79) Chrysene-d12	21.250	240	406305	20.000	ng/ul	0.00
88) Perylene-d12	23.474	264	495175	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	7776	4.048	ng/uL	0.00
4) Pyridine-d5	3.616	84	62282	11.924	ng/ul	0.00
7) Phenol-d5	6.810	99	125454	20.632	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67	80697	22.275	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	106031	22.258	ng/ul	0.00
15) 4-Methylphenol-d8	8.339	113	99325	20.869	ng/ul	0.00
21) Nitrobenzene-d5	8.792	128	50764	21.247	ng/ul	0.00
24) 2-Nitrophenol-d4	9.510	143	52782	20.930	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.039	165	95088	20.566	ng/ul	0.00
31) 4-Chloroaniline-d4	10.569	131	128937	17.515	ng/ul	0.00
46) Dimethylphthalate-d6	13.710	166	308782	22.676	ng/ul	0.00
49) Acenaphthylene-d8	13.974	160	361439	22.468	ng/ul	0.00
54) 4-Nitrophenol-d4	14.521	143	36136	12.715	ng/ul	0.00
60) Fluorene-d10	15.286	176	283833	23.380	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.427	200	34107	13.829	ng/ul	0.00
73) Anthracene-d10	17.139	188	437881	23.400	ng/ul	0.00
81) Pyrene-d10	19.445	212	546672	27.325	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.333	264	607507	25.149	ng/ul	0.00

Target Compounds Qvalue

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

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