

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
 Data File : BM044971.D
 Acq On : 06 Apr 2024 15:28
 Operator : MA/JU
 Sample : P2018-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Apr 08 01:50:59 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	69565	20.000	ng/ul	0.00
20) Naphthalene-d8	10.410	136	307082	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.280	164	193869	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.039	188	411097	20.000	ng/ul	0.00
79) Chrysene-d12	21.251	240	390804	20.000	ng/ul	0.00
88) Perylene-d12	23.474	264	467611	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	9649	5.179	ng/uL	0.00
4) Pyridine-d5	3.605	84	139635	27.563	ng/ul	0.00
7) Phenol-d5	6.810	99	184631	31.307	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67	116908	33.273	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	154970	33.541	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113	150748	32.657	ng/ul	0.00
21) Nitrobenzene-d5	8.793	128	76161	32.288	ng/ul	0.00
24) 2-Nitrophenol-d4	9.504	143	82422	33.106	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.039	165	141702	31.043	ng/ul	0.00
31) 4-Chloroaniline-d4	10.563	131	206503	28.415	ng/ul	0.00
46) Dimethylphthalate-d6	13.710	166	438306	32.471	ng/ul	0.00
49) Acenaphthylene-d8	13.975	160	530632	33.275	ng/ul	0.00
54) 4-Nitrophenol-d4	14.516	143	69244	24.579	ng/ul	0.00
60) Fluorene-d10	15.286	176	398602	33.122	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.422	200	54677	22.432	ng/ul	0.00
73) Anthracene-d10	17.139	188	627734	33.944	ng/ul	0.00
81) Pyrene-d10	19.445	212	746725	38.805	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.333	264	804526	35.269	ng/ul	0.00

Target Compounds Qvalue

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

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