

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
 Data File : BM036747.D
 Acq On : 26 Sep 2022 16:12
 Operator : CG/JU
 Sample : N4748-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Sep 26 23:14:18 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 26 23:10:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.004	152	394631	20.000	ng/ul	0.00
20) Naphthalene-d8	10.816	136	1728806	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.627	164	1074370	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.368	188	2066897	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240	1509294	20.000	ng/ul	0.00
88) Perylene-d12	23.968	264	1306936	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	50591	4.806	ng/uL	0.00
4) Pyridine-d5	3.816	84	227173	7.399	ng/ul	0.00
7) Phenol-d5	7.157	99	223288	6.332	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	667636	31.173	ng/ul	0.00
11) 2-Chlorophenol-d4	7.534	132	634044	24.972	ng/ul	0.00
15) 4-Methylphenol-d8	8.710	113	425266	16.320	ng/ul	0.00
21) Nitrobenzene-d5	9.169	128	419803	35.876	ng/ul	0.00
24) 2-Nitrophenol-d4	9.892	143	381573	36.405	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.433	165	685748	28.231	ng/ul	0.00
31) 4-Chloroaniline-d4	10.951	131	902331	22.343	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	2562723	35.367	ng/ul	0.00
49) Acenaphthylene-d8	14.327	160	2925571	34.038	ng/ul	0.00
54) 4-Nitrophenol-d4	14.821	143	63304	4.504	ng/ul	0.00
60) Fluorene-d10	15.616	176	2276631	35.060	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.727	200	234389	28.662	ng/ul	0.00
73) Anthracene-d10	17.462	188	3437984	36.379	ng/ul	0.00
81) Pyrene-d10	19.751	212	3587707	48.318	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.809	264	2409889	37.742	ng/ul	0.00

Target Compounds	Qvalue

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

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