

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
 Data File : BG062850.D
 Acq On : 15 Sep 2024 17:34
 Operator : JU/RC
 Sample : P3952-20
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0B41

Quant Time: Sep 15 23:04:44 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 10 01:28:17 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.902	152	35070	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	143202	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.529	164	126621	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.373	188	592968	20.000	ng/ul	0.10
79) Chrysene-d12	21.521	240	427359	20.000	ng/ul	0.00
88) Perylene-d12	24.582	264	473723	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.360	96	3222	4.153	ng/uL	0.00
4) Pyridine-d5	3.771	84	187	0.076	ng/ul	0.00
7) Phenol-d5	7.067	99	34940	11.362	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.232	67	51678	28.029	ng/ul	0.00
11) 2-Chlorophenol-d4	7.432	132	67030	29.924	ng/ul	0.00
15) 4-Methylphenol-d8	8.601	113	61237	23.896	ng/ul	0.00
21) Nitrobenzene-d5	9.053	128	37489	36.629	ng/ul	0.00
24) 2-Nitrophenol-d4	9.776	143	43358	38.736	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.317	165	86358	35.427	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	7249	2.163	ng/ul	0.00
46) Dimethylphthalate-d6	13.930	166	329278	33.770	ng/ul	0.00
49) Acenaphthylene-d8	14.218	160	334501	30.889	ng/ul	0.00
54) 4-Nitrophenol-d4	14.729	143	19182	11.580	ng/ul	0.00
60) Fluorene-d10	15.522	176	303604	34.553	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.734	200	468	0.137	ng/ul	0.09
73) Anthracene-d10	17.373	188	592337	21.167	ng/ul	0.00
81) Pyrene-d10	19.665	212	830641	34.540	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.377	264	922710	36.888	ng/ul	0.00

Target Compounds Qvalue

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

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