

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110423\
 Data File : BG059649.D
 Acq On : 5 Nov 2023 10:16
 Operator : MA/JU
 Sample : 05016-10
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A49P1

Quant Time: Nov 05 23:59:00 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110323.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Nov 04 02:46:50 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.355	152	22531	20.000	ng/u1	0.00
20) Naphthalene-d8	11.198	136	112415	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.976	164	83681	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.732	188	192398	20.000	ng/u1	# 0.00
79) Chrysene-d12	22.062	240	208544	20.000	ng/u1	0.00
88) Perylene-d12	25.670	264	240281	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.648	96	3870	6.438	ng/uL	0.00
4) Pyridine-d5	4.077	84	17798	10.068	ng/u1	0.00
7) Phenol-d5	7.462	99	23309	10.119	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.661	67	40261	32.689	ng/u1	0.00
11) 2-Chlorophenol-d4	7.867	132	47687	30.481	ng/u1	0.00
15) 4-Methylphenol-d8	9.036	113	40207	21.327	ng/u1	0.00
21) Nitrobenzene-d5	9.542	128	33067	40.222	ng/u1	0.00
24) 2-Nitrophenol-d4	10.264	143	38459	38.764	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.793	165	68398	38.804	ng/u1	0.00
31) 4-Chloroaniline-d4	11.339	131	70983	25.591	ng/u1	0.00
46) Dimethylphthalate-d6	14.371	166	314449	46.822	ng/u1	0.00
49) Acenaphthylene-d8	14.683	160	330884	41.318	ng/u1	0.00
54) 4-Nitrophenol-d4	15.147	143	14508	12.235	ng/u1	0.00
60) Fluorene-d10	15.969	176	274960	43.488	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	16.069	200	55616	45.670	ng/u1	0.00
73) Anthracene-d10	17.832	188	448827	44.572	ng/u1	0.00
81) Pyrene-d10	20.100	212	603279	45.205	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.417	264	646870	45.988	ng/u1	0.00

Target Compounds	Qvalue

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

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