

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110423\
 Data File : BG059650.D
 Acq On : 5 Nov 2023 10:57
 Operator : MA/JU
 Sample : 05016-11
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A49P2

Quant Time: Nov 05 23:59:12 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	22706	20.000	ng/u1	0.00
20) Naphthalene-d8	11.204	136	110974	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.976	164	85305	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.732	188	192164	20.000	ng/u1	# 0.00
79) Chrysene-d12	22.062	240	222604	20.000	ng/u1	# 0.00
88) Perylene-d12	25.658	264	251388	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.648	96	2461	4.062	ng/uL	0.00
4) Pyridine-d5	4.083	84	18400	10.328	ng/u1	0.00
7) Phenol-d5	7.462	99	20993	9.043	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.667	67	36611	29.497	ng/u1	0.00
11) 2-Chlorophenol-d4	7.861	132	46489	29.487	ng/u1	0.00
15) 4-Methylphenol-d8	9.030	113	37886	19.941	ng/u1	0.00
21) Nitrobenzene-d5	9.553	128	30906	38.082	ng/u1	0.00
24) 2-Nitrophenol-d4	10.270	143	38819	39.635	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.793	165	63851	36.694	ng/u1	0.00
31) 4-Chloroaniline-d4	11.345	131	83433	30.470	ng/u1	0.00
46) Dimethylphthalate-d6	14.371	166	276952	40.453	ng/u1	0.00
49) Acenaphthylene-d8	14.683	160	302839	37.096	ng/u1	0.00
54) 4-Nitrophenol-d4	15.147	143	11515	9.526	ng/u1	0.00
60) Fluorene-d10	15.969	176	238262	36.966	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	16.069	200	46459	38.197	ng/u1	0.00
73) Anthracene-d10	17.832	188	390265	38.803	ng/u1	0.00
81) Pyrene-d10	20.100	212	542424	38.077	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.411	264	573573	38.976	ng/u1	0.00

Target Compounds	Qvalue

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

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