

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110423\
 Data File : BG059658.D
 Acq On : 5 Nov 2023 17:02
 Operator : MA/JU
 Sample : 05013-15
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4919

Quant Time: Nov 06 00:00:39 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	8.347	152	25195	20.000	ng/u1	# 0.00
20) Naphthalene-d8	11.197	136	123796	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.980	164	89568	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.730	188	215136	20.000	ng/u1	# 0.00
79) Chrysene-d12	22.066	240	249089	20.000	ng/u1	# 0.00
88) Perylene-d12	25.668	264	292885	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.647	96	2143	3.188	ng/uL	0.00
4) Pyridine-d5	4.081	84	18109	9.161	ng/u1	0.00
7) Phenol-d5	7.460	99	14370	5.579	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.665	67	31228	22.674	ng/u1	0.00
11) 2-Chlorophenol-d4	7.865	132	38903	22.237	ng/u1	0.00
15) 4-Methylphenol-d8	9.029	113	30908	14.661	ng/u1	-0.01
21) Nitrobenzene-d5	9.552	128	26290	29.039	ng/u1	0.00
24) 2-Nitrophenol-d4	10.262	143	31334	28.679	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.791	165	56068	28.884	ng/u1	0.00
31) 4-Chloroaniline-d4	11.338	131	69011	22.593	ng/u1	0.00
46) Dimethylphthalate-d6	14.369	166	248103	34.515	ng/u1	0.00
49) Acenaphthylene-d8	14.681	160	262285	30.599	ng/u1	0.00
54) 4-Nitrophenol-d4	15.145	143	11216	8.837	ng/u1	-0.01
60) Fluorene-d10	15.962	176	206278	30.481	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	16.073	200	42587	31.275	ng/u1	0.00
73) Anthracene-d10	17.830	188	354438	31.478	ng/u1	0.00
81) Pyrene-d10	20.104	212	528389	33.148	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.415	264	571010	33.304	ng/u1	0.00

Target Compounds	Qvalue

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

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