

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\
 Data File : BG050869.D
 Acq On : 4 Nov 2021 5:51
 Operator : CG/JU
 Sample : M4412-18
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG372

Quant Time: Nov 07 08:25:14 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 02 14:49:05 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.230	152	31069	20.000	ng/u1	0.00
20) Naphthalene-d8	11.056	136	140394	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.851	164	99471	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.595	188	233247	20.000	ng/u1	-0.01
79) Chrysene-d12	21.890	240	205333	20.000	ng/u1	-0.02
88) Perylene-d12	25.292	264	200048	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.582	96	4850	5.038	ng/uL	0.00
4) Pyridine-d5	4.017	84	20344	7.064	ng/u1	0.00
7) Phenol-d5	7.366	99	22629	6.827	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.542	67	63884	29.837	ng/u1	0.00
11) 2-Chlorophenol-d4	7.754	132	52998	23.072	ng/u1	-0.01
15) 4-Methylphenol-d8	8.923	113	40637	15.574	ng/u1	-0.01
21) Nitrobenzene-d5	9.399	128	37218	31.195	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.127	143	38304	28.874	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.668	165	59695	26.713	ng/u1	0.00
31) 4-Chloroaniline-d4	11.185	131	84878	25.081	ng/u1	0.00
46) Dimethylphthalate-d6	14.246	166	262902	34.546	ng/u1	0.00
49) Acenaphthylene-d8	14.546	160	306116	32.286	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.039	143	10595	7.678	ng/u1	0.00
60) Fluorene-d10	15.838	176	230335	34.166	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.950	200	41073	29.045	ng/u1	-0.01
73) Anthracene-d10	17.695	188	403444	36.585	ng/u1	-0.01
81) Pyrene-d10	19.969	212	463644	34.960	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.057	264	392621	35.502	ng/u1	-0.02

Target Compounds	Qvalue

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

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