

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\
 Data File : BG050859.D
 Acq On : 3 Nov 2021 21:02
 Operator : CG/JU
 Sample : M4412-10
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG366

Quant Time: Nov 07 07:50:17 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	8.229	152	27438	20.000	ng/u1	0.00
20) Naphthalene-d8	11.055	136	120599	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.850	164	86502	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.594	188	202296	20.000	ng/u1	-0.01
79) Chrysene-d12	21.889	240	178665	20.000	ng/u1	-0.02
88) Perylene-d12	25.291	264	176765	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.587	96	1850	2.176	ng/uL	0.00
4) Pyridine-d5	4.028	84	7938	3.121	ng/u1	0.01
7) Phenol-d5	7.371	99	18942	6.471	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.541	67	53372	28.226	ng/u1	0.00
11) 2-Chlorophenol-d4	7.753	132	43938	21.659	ng/u1	-0.01
15) 4-Methylphenol-d8	8.922	113	34061	14.781	ng/u1	-0.01
21) Nitrobenzene-d5	9.398	128	30916	30.166	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.126	143	31632	27.758	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.667	165	49909	26.000	ng/u1	0.00
31) 4-Chloroaniline-d4	11.190	131	25624	8.815	ng/u1	0.00
46) Dimethylphthalate-d6	14.245	166	234315	35.406	ng/u1	0.00
49) Acenaphthylene-d8	14.551	160	258572	31.360	ng/u1	0.00
54) 4-Nitrophenol-d4	15.044	143	8037	6.698	ng/u1	0.00
60) Fluorene-d10	15.837	176	198995	33.943	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.955	200	36353	29.640	ng/u1	0.00
73) Anthracene-d10	17.694	188	348122	36.398	ng/u1	-0.01
81) Pyrene-d10	19.974	212	419794	36.378	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.056	264	340553	34.850	ng/u1	-0.02

Target Compounds	Qvalue

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

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