

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\
 Data File : BG050862.D
 Acq On : 4 Nov 2021 00:26
 Operator : CG/JU
 Sample : M4412-22
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG376

Quant Time: Nov 07 07:57:30 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	29686	20.000	ng/u1	0.00
20) Naphthalene-d8	11.056	136	136899	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.851	164	95885	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.595	188	223116	20.000	ng/u1	-0.01
79) Chrysene-d12	21.890	240	195450	20.000	ng/u1	-0.02
88) Perylene-d12	25.292	264	191979	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.588	96	3953	4.298	ng/uL	0.00
4) Pyridine-d5	4.017	84	18900	6.869	ng/u1	0.00
7) Phenol-d5	7.372	99	22539	7.117	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	62155	30.382	ng/u1	0.00
11) 2-Chlorophenol-d4	7.754	132	52493	23.917	ng/u1	-0.01
15) 4-Methylphenol-d8	8.929	113	38829	15.574	ng/u1	0.00
21) Nitrobenzene-d5	9.399	128	36681	31.530	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.127	143	38374	29.665	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.668	165	60110	27.586	ng/u1	0.00
31) 4-Chloroaniline-d4	11.185	131	79078	23.964	ng/u1	0.00
46) Dimethylphthalate-d6	14.246	166	265669	36.215	ng/u1	0.00
49) Acenaphthylene-d8	14.551	160	304544	33.322	ng/u1	0.00
54) 4-Nitrophenol-d4	15.045	143	9597	7.215	ng/u1	0.00
60) Fluorene-d10	15.838	176	230468	35.465	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.956	200	40122	29.661	ng/u1	0.00
73) Anthracene-d10	17.695	188	395365	37.480	ng/u1	-0.01
81) Pyrene-d10	19.969	212	460122	36.449	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.057	264	381581	35.954	ng/u1	-0.02

Target Compounds	Qvalue

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

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