

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\
 Data File : BG050821.D
 Acq On : 2 Nov 2021 17:40
 Operator : CG/JU
 Sample : M4364-12
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG322

Quant Time: Nov 02 18:37:09 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	8.234	152	31219	20.000	ng/u1	0.00
20) Naphthalene-d8	11.054	136	147839	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.856	164	105243	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.600	188	250163	20.000	ng/u1	0.00
79) Chrysene-d12	21.895	240	226063	20.000	ng/u1	-0.01
88) Perylene-d12	25.297	264	211670	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.587	96	3805	3.934	ng/uL	0.00
4) Pyridine-d5	4.021	84	19197	6.634	ng/u1	0.00
7) Phenol-d5	7.376	99	19732	5.925	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.547	67	55635	25.860	ng/u1	0.00
11) 2-Chlorophenol-d4	7.758	132	47132	20.420	ng/u1	0.00
15) 4-Methylphenol-d8	8.928	113	34389	13.116	ng/u1	0.00
21) Nitrobenzene-d5	9.403	128	31773	25.290	ng/u1	0.00
24) 2-Nitrophenol-d4	10.132	143	33531	24.003	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.673	165	51530	21.898	ng/u1	0.00
31) 4-Chloroaniline-d4	11.190	131	77668	21.795	ng/u1	0.00
46) Dimethylphthalate-d6	14.245	166	219112	27.213	ng/u1	0.00
49) Acenaphthylene-d8	14.550	160	252134	25.134	ng/u1	0.00
54) 4-Nitrophenol-d4	15.050	143	7494	5.133	ng/u1	0.00
60) Fluorene-d10	15.843	176	191425	26.838	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.955	200	37774	24.906	ng/u1	0.00
73) Anthracene-d10	17.700	188	351202	29.694	ng/u1	0.00
81) Pyrene-d10	19.973	212	417873	28.619	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.062	264	345596	29.534	ng/u1	-0.01

Target Compounds Qvalue

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

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