

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
 Data File : BG050888.D
 Acq On : 8 Nov 2021 20:20
 Operator : CG/JU
 Sample : M4445-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG378

Quant Time: Nov 09 01:03:05 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	37689	20.000	ng/u1	0.00
20) Naphthalene-d8	11.055	136	162374	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.851	164	109844	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.594	188	246518	20.000	ng/u1	-0.01
79) Chrysene-d12	21.889	240	216066	20.000	ng/u1	-0.02
88) Perylene-d12	25.291	264	210149	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.593	96	5170	4.427	ng/uL	0.00
4) Pyridine-d5	4.028	84	8123	2.325	ng/u1	0.01
7) Phenol-d5	7.371	99	22515	5.600	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	65766	25.321	ng/u1	0.00
11) 2-Chlorophenol-d4	7.759	132	54809	19.670	ng/u1	0.00
15) 4-Methylphenol-d8	8.928	113	40014	12.641	ng/u1	0.00
21) Nitrobenzene-d5	9.398	128	37859	27.437	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.127	143	38161	24.872	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.667	165	61029	23.613	ng/u1	0.00
31) 4-Chloroaniline-d4	11.184	131	56021	14.313	ng/u1	0.00
46) Dimethylphthalate-d6	14.245	166	250595	29.819	ng/u1	0.00
49) Acenaphthylene-d8	14.551	160	291188	27.811	ng/u1	0.00
54) 4-Nitrophenol-d4	15.044	143	7251	4.759	ng/u1	0.00
60) Fluorene-d10	15.838	176	220670	29.642	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.955	200	36589	24.481	ng/u1	0.00
73) Anthracene-d10	17.694	188	365103	31.326	ng/u1	-0.01
81) Pyrene-d10	19.968	212	425827	30.513	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.050	264	354897	30.549	ng/u1	-0.02

Target Compounds Qvalue

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

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