

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
 Data File : BG050347.D
 Acq On : 1 Oct 2021 7:31
 Operator : CG/JU
 Sample : M3873-16
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW7Y5

Quant Time: Oct 01 08:14:13 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG093021.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 30 16:27:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.273	152	40553	20.000	ng/u1	0.00
20) Naphthalene-d8	11.099	136	167219	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.889	164	122404	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.633	188	296091	20.000	ng/u1	0.00
79) Chrysene-d12	21.933	240	286076	20.000	ng/u1	-0.02
88) Perylene-d12	25.359	264	275375	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.631	96	7640	6.922	ng/uL	0.00
4) Pyridine-d5	4.066	84	35512	10.797	ng/u1	0.00
7) Phenol-d5	7.397	99	39741	11.216	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.585	67	90403	40.542	ng/u1	0.00
11) 2-Chlorophenol-d4	7.797	132	79204	31.904	ng/u1	0.00
15) 4-Methylphenol-d8	8.954	113	62874	23.445	ng/u1	0.00
21) Nitrobenzene-d5	9.436	128	50733	43.369	ng/u1	0.00
24) 2-Nitrophenol-d4	10.171	143	50417	40.158	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.699	165	88524	35.707	ng/u1	0.00
31) 4-Chloroaniline-d4	11.222	131	119653	33.785	ng/u1	0.00
46) Dimethylphthalate-d6	14.284	166	358716	42.821	ng/u1	0.00
49) Acenaphthylene-d8	14.583	160	417432	39.679	ng/u1	0.00
54) 4-Nitrophenol-d4	15.041	143	19707	12.214	ng/u1	-0.02
60) Fluorene-d10	15.876	176	323065	42.805	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.976	200	60290	40.824	ng/u1	0.00
73) Anthracene-d10	17.732	188	562672	44.983	ng/u1	0.00
81) Pyrene-d10	20.006	212	675141	44.383	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.124	264	588461	42.498	ng/u1	-0.02

Target Compounds	Qvalue

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

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