

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100421\
 Data File : BG050412.D
 Acq On : 4 Oct 2021 13:25
 Operator : CG/JU
 Sample : M3878-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW2P3

Quant Time: Oct 04 14:08:24 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	8.271	152	45328	20.00	ng/ul	0.00
20) Naphthalene-d8	11.097	136	190618	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.887	164	125776	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.631	188	270131	20.00	ng/ul	0.00
79) Chrysene-d12	21.931	240	249918	20.00	ng/ul	-0.02
88) Perylene-d12	25.357	264	245811	20.00	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.641	96	5972	4.84	ng/uL	0.00
4) Pyridine-d5	4.070	84	19518	5.31	ng/ul	0.00
7) Phenol-d5	7.401	99	30633	7.74	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.583	67	93310	37.44	ng/ul	0.00
11) 2-Chlorophenol-d4	7.795	132	78147	28.16	ng/ul	0.00
15) 4-Methylphenol-d8	8.958	113	55774	18.61	ng/ul	0.00
21) Nitrobenzene-d5	9.440	128	53558	40.16	ng/ul	0.00
24) 2-Nitrophenol-d4	10.169	143	53346	37.27	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.703	165	88197	31.21	ng/ul	0.00
31) 4-Chloroaniline-d4	11.226	131	126307	31.29	ng/ul	0.00
46) Dimethylphthalate-d6	14.282	166	301244	35.00	ng/ul	-0.01
49) Acenaphthylene-d8	14.587	160	388136	35.91	ng/ul	0.00
54) 4-Nitrophenol-d4	15.051	143	10499	6.33	ng/ul	0.00
60) Fluorene-d10	15.874	176	272865	35.18	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.980	200	35768	26.55	ng/ul	0.00
73) Anthracene-d10	17.730	188	432115	37.87	ng/ul	0.00
81) Pyrene-d10	20.004	212	501907	37.77	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.128	264	443413	35.87	ng/ul	-0.01

Target Compounds Qvalue

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

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