

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070921\
 Data File : BG049250.D
 Acq On : 9 Jul 2021 23:29
 Operator : CG/JU
 Sample : M2878-02
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGDZ1

Manual Integrations
 APPROVED

mohammad
 7/12/2021 12:33:45 PM

Quant Time: Jul 10 00:07:04 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 09 22:53:59 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.079	152	28226	20.000	ng/u1	0.00
20) Naphthalene-d8	10.893	136	112022	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.718	164	84819	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.468	188	198137	20.000	ng/u1	0.00
79) Chrysene-d12	21.751	240	217211	20.000	ng/u1	0.00
88) Perylene-d12	25.053	264	216781	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.472	96	139	0.232	ng/uL	0.00
4) Pyridine-d5	3.919	84	3417m	1.937	ng/u1	0.03
7) Phenol-d5	7.227	99	7420	3.022	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.397	67	4943	3.471	ng/u1	0.00
11) 2-Chlorophenol-d4	7.603	132	6113	3.369	ng/u1	0.00
15) 4-Methylphenol-d8	8.778	113	7044	3.581	ng/u1	0.00
21) Nitrobenzene-d5	9.242	128	3025	3.519	ng/u1	0.00
24) 2-Nitrophenol-d4	9.971	143	3728	3.778	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.517	165	6518	3.385	ng/u1	0.00
31) 4-Chloroaniline-d4	11.028	131	8748	3.486	ng/u1	0.00
46) Dimethylphthalate-d6	14.119	166	27261	4.179	ng/u1	0.00
49) Acenaphthylene-d8	14.413	160	30652	4.005	ng/u1	0.00
54) 4-Nitrophenol-d4	14.918	143	2604m	2.412	ng/u1	0.00
60) Fluorene-d10	15.711	176	22846	4.137	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.823	200	3552m	2.820	ng/u1	0.00
73) Anthracene-d10	17.568	188	40905	4.534	ng/u1	0.00
81) Pyrene-d10	19.853	212	51605	4.635	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.818	264	47160	4.185	ng/u1	0.00

Target Compounds Qvalue

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

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