

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100919\
 Data File : BG043128.D
 Acq On : 10 Oct 2019 9:04
 Operator : JU
 Sample : K5097-11
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JLME7

Manual Integrations
 APPROVED

Quant Time: Oct 10 12:01:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 10 11:52:02 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	49236	20.00	ng/ul	0.00
18) Naphthalene-d8	10.86	136	207796	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.68	164	158007	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.42	188	380092	20.00	ng/ul	0.00
77) Chrysene-d12	21.69	240	414643	20.00	ng/ul	0.00
85) Perylene-d12	24.92	264	490576	20.00	ng/ul	0.01

mohammad

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.50	96	3285	3.23	ng/uL	0.00
5) Phenol-d5	7.22	99	57838	15.89	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.38	67	36671	18.05	ng/ul	0.00
9) 2-Chlorophenol-d4	7.59	132	49630	17.01	ng/ul	0.00
13) 4-Methylphenol-d8	8.77	113	50648	16.14	ng/ul	0.00
19) Nitrobenzene-d5	9.22	128	26970	18.44	ng/ul	0.00
22) 2-Nitrophenol-d4	9.94	143	31834	19.09	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.49	165	56690	16.25	ng/ul	0.00
29) 4-Chloroaniline-d4	11.01	131	61911	17.04	ng/ul	0.01
43) Dimethylphthalate-d6	14.08	166	242158	19.06	ng/ul	0.00
46) Acenaphthylene-d8	14.37	160	264902	19.17	ng/ul	0.00
51) 4-Nitrophenol-d4	14.91	143	11937m	7.09	ng/ul	0.02
57) Fluorene-d10	15.67	176	205048	19.01	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.78	200	26505	10.87	ng/ul	0.00
70) Anthracene-d10	17.52	188	361994	20.16	ng/ul	0.00
78) Pyrene-d10	19.80	212	450978	22.06	ng/ul	0.00
89) Benzo(a)pyrene-d12	24.69	264	507914	20.95	ng/ul	0.00

10/12/2019 7:32:22 AM

Target Compounds

					Ovalue
6) Phenol	7.25	94	9605	2.585	ng/ul 98
44) Dimethylphthalate	14.13	163	79233	6.458	ng/ul 99

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

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