

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120420\
 Data File : BG047604.D
 Acq On : 5 Dec 2020 13:52
 Operator : CG/JU
 Sample : L4864-13
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0BH5

Quant Time: Dec 07 05:37:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 07 05:06:33 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.25	152	27986	20.00	ng/ul	-0.02
18) Naphthalene-d8	11.08	136	111881	20.00	ng/ul	-0.02
36) Acenaphthene-d10	14.86	164	91353	20.00	ng/ul	-0.02
62) Phenanthrene-d10	17.60	188	231705	20.00	ng/ul	-0.02
78) Chrysene-d12	21.88	240	220784	20.00	ng/ul	-0.02
86) Perylene-d12	25.26	264	231313	20.00	ng/ul	-0.03

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.58	96	4090	5.31	ng/uL	-0.03
5) Phenol-d5	7.39	99	101333	36.11	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	7.56	67	54432	34.00	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.77	132	79618	42.30	ng/ul	-0.02
13) 4-Methylphenol-d8	8.94	113	82067	38.05	ng/ul	-0.02
19) Nitrobenzene-d5	9.42	128	40967	44.33	ng/ul	-0.02
22) 2-Nitrophenol-d4	10.15	143	50047	49.63	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	10.69	165	98825	42.81	ng/ul	-0.02
29) 4-Chloroaniline-d4	11.20	131	115566	49.50	ng/ul	-0.02
44) Dimethylphthalate-d6	14.26	166	352729	44.98	ng/ul	-0.02
47) Acenaphthylene-d8	14.56	160	399968	44.13	ng/ul	-0.02
52) 4-Nitrophenol-d4	15.03	143	53104	44.38	ng/ul	0.00
58) Fluorene-d10	15.85	176	327535	45.59	ng/ul	-0.02
63) 4,6-Dinitro-2-methylphenol	15.95	200	67296	51.03	ng/ul	-0.02
71) Anthracene-d10	17.70	188	536622	46.69	ng/ul	-0.02
79) Pyrene-d10	19.96	212	600128	46.71	ng/ul	-0.01
90) Benzo(a)pyrene-d12	25.04	264	617526	46.38	ng/ul	-0.02

Target Compounds

					Ovalue	
6) Phenol	7.41	94	6988	2.599	ng/ul#	78
45) Dimethylphthalate	14.30	163	92483	11.561	ng/ul	97

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

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