

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102220\
 Data File : BG047034.D
 Acq On : 22 Oct 2020 16:24
 Operator : CG/JU
 Sample : L4483-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0FH1

Quant Time: Oct 22 17:08:07 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 22 05:48:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	51582	20.00	ng/ul	0.00
18) Naphthalene-d8	10.87	136	205910	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.69	164	146909	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.43	188	365328	20.00	ng/ul	0.00
78) Chrysene-d12	21.70	240	398694	20.00	ng/ul	0.00
86) Perylene-d12	24.93	264	401433	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.48	96	6510	5.09	ng/uL	0.00
5) Phenol-d5	7.21	99	28859	6.45	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.38	67	81312	31.58	ng/ul	0.00
9) 2-Chlorophenol-d4	7.59	132	86588	24.87	ng/ul	0.00
13) 4-Methylphenol-d8	8.76	113	53876	15.08	ng/ul	0.00
19) Nitrobenzene-d5	9.22	128	54995	31.86	ng/ul	0.00
22) 2-Nitrophenol-d4	9.95	143	63861	31.67	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.48	165	109280	27.78	ng/ul	0.00
29) 4-Chloroaniline-d4	11.00	131	12153	3.62	ng/ul	0.00
44) Dimethylphthalate-d6	14.09	166	413516	34.23	ng/ul	0.00
47) Acenaphthylene-d8	14.39	160	486835	34.06	ng/ul	0.00
52) 4-Nitrophenol-d4	14.87	143	10498	5.52	ng/ul	0.00
58) Fluorene-d10	15.68	176	382428	34.10	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	87981	34.92	ng/ul	0.00
71) Anthracene-d10	17.53	188	654451	36.96	ng/ul	0.00
79) Pyrene-d10	19.82	212	851097	37.42	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.72	264	821882	37.24	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	7.24	94	8302	1.900 ng/ul	87
20) Nitrobenzene	9.26	77	69769	14.971 ng/ul	99
45) Dimethylphthalate	14.13	163	38005	3.135 ng/ul#	93

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

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