

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102220\
 Data File : BG047035.D
 Acq On : 22 Oct 2020 17:04
 Operator : CG/JU
 Sample : L4483-05DL 5X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0FH3DL

Quant Time: Oct 22 18:09:11 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	8.07	152	62272	20.00	ng/ul	0.00
18) Naphthalene-d8	10.88	136	220392	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.69	164	158422	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.43	188	388483	20.00	ng/ul	0.00
78) Chrysene-d12	21.70	240	447030	20.00	ng/ul	0.00
86) Perylene-d12	24.93	264	446139	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
5) Phenol-d5	7.22	99	7800	1.44	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.38	67	20630	6.64	ng/ul	0.00
9) 2-Chlorophenol-d4	7.59	132	23320	5.55	ng/ul	0.00
13) 4-Methylphenol-d8	8.76	113	13643	3.16	ng/ul	0.00
19) Nitrobenzene-d5	9.22	128	13586	7.35	ng/ul	0.00
22) 2-Nitrophenol-d4	9.95	143	15367	7.12	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.49	165	26991	6.41	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
44) Dimethylphthalate-d6	14.09	166	96695	7.42	ng/ul	0.00
47) Acenaphthylene-d8	14.39	160	120378	7.81	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
58) Fluorene-d10	15.68	176	91858	7.59	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	16952	6.33	ng/ul	0.00
71) Anthracene-d10	17.53	188	159385	8.46	ng/ul	0.00
79) Pyrene-d10	19.81	212	203270	7.97	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.70	264	193480	7.89	ng/ul	-0.01

Target Compounds

					Ovalue
20) Nitrobenzene	9.27	77	117827	23.622	ng/ul 98
37) 1,2,4,5-Tetrachlorobenzene	12.89	216	63073	10.222	ng/ul 94
73) 1,2,3,4-Tetrachlorobenzene	13.49	216	173782	26.629	ng/uL 98

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

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