

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081219\
 Data File : BG042169.D
 Acq On : 13 Aug 2019 7:31
 Operator : HP/JU
 Sample : K4215-22
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ETOC6

Quant Time: Aug 13 13:48:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 13 04:44:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	83652	20.00	ng/ul	0.00
18) Naphthalene-d8	10.85	136	349492	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.68	164	237790	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.44	188	509863	20.00	ng/ul	0.00
77) Chrysene-d12	21.71	240	521087	20.00	ng/ul	0.00
85) Perylene-d12	24.98	264	554417	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.41	96	8504	4.45	ng/uL	0.00
5) Phenol-d5	7.18	99	182653	27.82	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.34	67	95268	25.95	ng/ul	0.00
9) 2-Chlorophenol-d4	7.55	132	161745	28.82	ng/ul	0.00
13) 4-Methylphenol-d8	8.74	113	159072	28.86	ng/ul	0.00
19) Nitrobenzene-d5	9.19	128	84258	32.03	ng/ul	0.00
22) 2-Nitrophenol-d4	9.92	143	101284	33.16	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.47	165	180891	29.33	ng/ul	0.00
29) 4-Chloroaniline-d4	10.98	131	224109	32.75	ng/ul	0.00
43) Dimethylphthalate-d6	14.08	166	554826	30.08	ng/ul	0.00
46) Acenaphthylene-d8	14.38	160	666878	31.66	ng/ul	0.00
51) 4-Nitrophenol-d4	14.89	143	42687	13.78	ng/ul	0.00
57) Fluorene-d10	15.67	176	508915	31.02	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.79	200	39338	12.03	ng/ul	0.00
70) Anthracene-d10	17.53	188	774602	31.67	ng/ul	0.00
78) Pyrene-d10	19.82	212	871551	29.97	ng/ul	0.00
89) Benzo(a)pyrene-d12	24.76	264	946407	32.68	ng/ul	0.00

Target Compounds

					Ovalue	
6) Phenol	7.21	94	15045	2.220	ng/ul#	89
44) Dimethylphthalate	14.13	163	189800	10.365	ng/ul	99

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

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