

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111419\
 Data File : BM023726.D
 Acq On : 14 Nov 2019 20:53
 Operator : JU
 Sample : K5700-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BEGY1

Quant Time: Nov 15 06:33:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM11119MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 15 06:30:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.53	152	330000	20.00	ng/ul	0.00
18) Naphthalene-d8	10.30	136	1333206	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.19	164	849577	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.94	188	1804886	20.00	ng/ul	0.00
78) Chrysene-d12	21.14	240	1782674	20.00	ng/ul	0.00
86) Perylene-d12	23.30	264	2112065	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.06	96	45349	5.66	ng/uL	0.00
5) Phenol-d5	6.72	99	170011	5.65	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.88	67	418730	24.23	ng/ul	0.00
9) 2-Chlorophenol-d4	7.07	132	481773	22.11	ng/ul	0.00
13) 4-Methylphenol-d8	8.25	113	313964	13.08	ng/ul	0.00
19) Nitrobenzene-d5	8.68	128	291079	30.50	ng/ul	0.00
22) 2-Nitrophenol-d4	9.40	143	331226	33.02	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.94	165	597577	26.13	ng/ul	0.00
29) 4-Chloroaniline-d4	10.46	131	45314	1.80	ng/ul	0.00
44) Dimethylphthalate-d6	13.61	166	2033284	30.69	ng/ul	0.00
47) Acenaphthylene-d8	13.87	160	2280764	30.17	ng/ul	0.00
52) 4-Nitrophenol-d4	14.42	143	55530	4.91	ng/ul	0.00
58) Fluorene-d10	15.19	176	1749020	29.90	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.32	200	310833	29.77	ng/ul	0.00
71) Anthracene-d10	17.03	188	2937028	34.30	ng/ul	0.00
79) Pyrene-d10	19.34	212	3296952	35.47	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.16	264	3903447	35.12	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.75	94	40521	1.308	ng/ul 99
45) Dimethylphthalate	13.65	163	119409	1.839	ng/ul 99

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

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