

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101719\
 Data File : BP000838.D
 Acq On : 17 Oct 2019 13:12
 Operator : JU
 Sample : K5346-14
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ET0N6

Quant Time: Oct 17 15:28:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 17 09:27:24 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	187928	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	717001	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	400138	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.30	188	747081	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	643688	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	727217	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.36	96	20691	4.57	ng/uL	-0.04
5) Phenol-d5	6.39	99	93889	6.48	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	228288	31.50	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	329360	26.97	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	188730	17.04	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	188869	34.65	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	203989	35.15	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	337603	30.39	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	51819	4.19	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	961511	34.33	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	1166434	33.85	ng/ul	0.00
52) 4-Nitrophenol-d4	9.93	143	21620	5.01	ng/ul	0.01
58) Fluorene-d10	10.32	176	815720	34.04	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	89845	25.25	ng/ul	0.00
71) Anthracene-d10	11.35	188	1178663	34.06	ng/ul	0.00
79) Pyrene-d10	12.74	212	1235797	33.95	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	1266203	35.27	ng/ul	0.00

Target Compounds Ovalue

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

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