

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP102119\
 Data File : BP000939.D
 Acq On : 21 Oct 2019 18:31
 Operator : JU
 Sample : K5321-10
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E0AA6

Quant Time: Oct 21 18:53:08 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 19 22:14:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	157447	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	601149	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.79	164	349657	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	680049	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	657495	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	746679	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.33	96	11900	3.14	ng/uL	0.00
5) Phenol-d5	6.39	99	44295	3.65	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.42	67	110884	18.26	ng/ul	0.00
9) 2-Chlorophenol-d4	6.51	132	163195	15.95	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	96184	10.37	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	94665	20.72	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	104431	21.46	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	174459	18.73	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	13950	1.35	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	539567	22.04	ng/ul	0.00
47) Acenaphthylene-d8	9.63	160	616419	20.47	ng/ul	0.00
52) 4-Nitrophenol-d4	9.93	143	14008	3.72	ng/ul	0.01
58) Fluorene-d10	10.32	176	462725	22.09	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	58059	17.93	ng/ul	0.00
71) Anthracene-d10	11.35	188	729728	23.17	ng/ul	0.00
79) Pyrene-d10	12.73	212	828168	22.27	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	914961	24.82	ng/ul	0.00

Target Compounds

				Ovalue	
84) Bis(2-ethylhexyl)phthalate	13.95	149	54625	2.156	ng/ul 99

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

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