

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101719\
 Data File : BP000834.D
 Acq On : 17 Oct 2019 11:36
 Operator : JU
 Sample : K5346-10
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ET0N4

Quant Time: Oct 17 14:58:53 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	211410	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	790739	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	422394	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.30	188	753160	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	634270	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	694235	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	2.38	96	26888	5.28	ng/uL	-0.02
5) Phenol-d5	6.39	99	80493	4.94	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	197816	24.26	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	281861	20.51	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	159596	12.81	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	161907	26.94	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	176976	27.65	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.89	165	294610	24.05	ng/ul	0.00
29) 4-Chloroaniline-d4	8.10	131	42237	3.10	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	837155	28.31	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	985273	27.09	ng/ul	0.00
52) 4-Nitrophenol-d4	9.93	143	21391	4.70	ng/ul	0.01
58) Fluorene-d10	10.32	176	705734	27.90	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	87680	24.45	ng/ul	0.00
71) Anthracene-d10	11.35	188	1118564	32.06	ng/ul	0.00
79) Pyrene-d10	12.74	212	1158116	32.29	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	1123027	32.77	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
45) Dimethylphthalate	9.51	163	87524	2.971	ng/ul	100

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

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