

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101919\
 Data File : BP000918.D
 Acq On : 19 Oct 2019 16:39
 Operator : JU
 Sample : K5321-20
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E0AB7

Quant Time: Oct 19 22:46:51 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	127222	20.00	ng/ul	0.00
18) Naphthalene-d8	8.02	136	478350	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.79	164	269101	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.28	188	518541	20.00	ng/ul	0.00
78) Chrysene-d12	13.96	240	504373	20.00	ng/ul	0.00
86) Perylene-d12	15.43	264	580587	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.33	96	17263	5.63	ng/uL	0.00
5) Phenol-d5	6.39	99	64870	6.61	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.42	67	129222	26.34	ng/ul	0.00
9) 2-Chlorophenol-d4	6.51	132	194854	23.57	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	117970	15.73	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	110250	30.32	ng/ul	0.00
22) 2-Nitrophenol-d4	7.62	143	120656	31.16	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	199945	26.98	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	2304	0.28	ng/ul	0.00
44) Dimethylphthalate-d6	9.48	166	587812	31.20	ng/ul	0.00
47) Acenaphthylene-d8	9.63	160	674763	29.12	ng/ul	0.00
52) 4-Nitrophenol-d4	9.92	143	18798	6.48	ng/ul	0.00
58) Fluorene-d10	10.30	176	502501	31.18	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	63050	25.53	ng/ul	0.00
71) Anthracene-d10	11.34	188	816321	33.99	ng/ul	0.00
79) Pyrene-d10	12.72	212	904394	31.71	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	981211	34.23	ng/ul	0.00

Target Compounds Ovalue

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

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