

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101919\
 Data File : BP000913.D
 Acq On : 19 Oct 2019 14:39
 Operator : JU
 Sample : K5321-15
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E0AA9

Quant Time: Oct 19 22:34:56 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	146385	20.00	ng/ul	0.00
18) Naphthalene-d8	8.02	136	547230	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.79	164	305929	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.28	188	585662	20.00	ng/ul	0.00
78) Chrysene-d12	13.96	240	541885	20.00	ng/ul	0.00
86) Perylene-d12	15.43	264	617112	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.34	96	17606	4.99	ng/uL	0.00
5) Phenol-d5	6.39	99	64499	5.71	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.42	67	139210	24.66	ng/ul	0.00
9) 2-Chlorophenol-d4	6.51	132	207344	21.79	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	124496	14.43	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	117785	28.32	ng/ul	0.00
22) 2-Nitrophenol-d4	7.62	143	129582	29.25	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	215601	25.43	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	32670	3.47	ng/ul	0.00
44) Dimethylphthalate-d6	9.48	166	628818	29.36	ng/ul	0.00
47) Acenaphthylene-d8	9.63	160	751444	28.52	ng/ul	0.00
52) 4-Nitrophenol-d4	9.92	143	17805	5.40	ng/ul	0.00
58) Fluorene-d10	10.31	176	533672	29.12	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	63830	22.89	ng/ul	0.00
71) Anthracene-d10	11.34	188	829304	30.57	ng/ul	0.00
79) Pyrene-d10	12.72	212	913964	29.83	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	982464	32.25	ng/ul	0.00

Target Compounds					Ovalue	
45) Dimethylphthalate	9.50	163	146881	6.883	ng/ul	99

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

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