

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101919\
 Data File : BP000911.D
 Acq On : 19 Oct 2019 13:51
 Operator : JU
 Sample : K5321-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E0AB4

Quant Time: Oct 19 22:30:31 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	147288	20.00	ng/ul	0.00
18) Naphthalene-d8	8.02	136	570277	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.79	164	335161	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	658155	20.00	ng/ul	0.00
78) Chrysene-d12	13.96	240	621572	20.00	ng/ul	0.00
86) Perylene-d12	15.43	264	664832	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.34	96	17169	4.84	ng/uL	0.00
5) Phenol-d5	6.39	99	64035	5.64	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.42	67	155359	27.35	ng/ul	0.00
9) 2-Chlorophenol-d4	6.51	132	222341	23.23	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	131783	15.18	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	135509	31.26	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	150323	32.56	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.87	165	247021	27.96	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	30238	3.08	ng/ul	0.00
44) Dimethylphthalate-d6	9.48	166	755830	32.21	ng/ul	0.00
47) Acenaphthylene-d8	9.63	160	908422	31.47	ng/ul	0.00
52) 4-Nitrophenol-d4	9.92	143	19370	5.36	ng/ul	0.00
58) Fluorene-d10	10.31	176	648433	32.30	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	80867	25.80	ng/ul	0.00
71) Anthracene-d10	11.34	188	1044219	34.25	ng/ul	0.00
79) Pyrene-d10	12.73	212	1128542	32.11	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	1150502	35.05	ng/ul	0.00
Target Compounds						
6) Phenol	6.40	94	12563	1.097	ng/ul	93
45) Dimethylphthalate	9.50	163	83261	3.562	ng/ul	100

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

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