

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101919\
 Data File : BP000909.D
 Acq On : 19 Oct 2019 13:03
 Operator : JU
 Sample : K5321-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E0AA7

Quant Time: Oct 19 22:24:03 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	136131	20.00	ng/ul	0.00
18) Naphthalene-d8	8.02	136	520402	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.79	164	299326	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.28	188	571017	20.00	ng/ul	0.00
78) Chrysene-d12	13.96	240	521781	20.00	ng/ul	0.00
86) Perylene-d12	15.43	264	580942	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.34	96	15646	4.77	ng/uL	0.00
5) Phenol-d5	6.39	99	68656	6.54	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.42	67	149315	28.44	ng/ul	0.00
9) 2-Chlorophenol-d4	6.51	132	222138	25.11	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	133158	16.60	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	128942	32.60	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	145496	34.54	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	235347	29.19	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	27371	3.05	ng/ul	0.00
44) Dimethylphthalate-d6	9.48	166	720863	34.40	ng/ul	0.00
47) Acenaphthylene-d8	9.63	160	843312	32.72	ng/ul	0.00
52) 4-Nitrophenol-d4	9.92	143	20807	6.45	ng/ul	0.00
58) Fluorene-d10	10.31	176	617482	34.44	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	72229	26.56	ng/ul	0.00
71) Anthracene-d10	11.34	188	947936	35.84	ng/ul	0.00
79) Pyrene-d10	12.73	212	1027511	34.82	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	1052644	36.70	ng/ul	0.00

Target Compounds					Ovalue	
45) Dimethylphthalate	9.50	163	38431	1.841	ng/ul	98

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

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