

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101519\
 Data File : BP000723.D
 Acq On : 14 Oct 2019 20:55
 Operator : JU
 Sample : K5097-13
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JLMFO

Quant Time: Oct 15 03:54:50 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 15 03:51:25 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	223503	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	846481	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	472963	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	865567	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	776046	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	867466	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.36	96	29982	5.56	ng/uL	0.00
5) Phenol-d5	6.39	99	507833	29.46	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	254807	29.56	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	422217	29.07	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	399306	30.32	ng/ul	0.00
19) Nitrobenzene-d5	7.31	128	200480	31.16	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	217330	31.72	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	384152	29.29	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	540358	37.05	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	1026383	31.00	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	1254345	30.80	ng/ul	0.00
52) 4-Nitrophenol-d4	9.91	143	129798	25.45	ng/ul	0.00
58) Fluorene-d10	10.32	176	857208	30.26	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	77760	18.86	ng/ul	0.00
71) Anthracene-d10	11.35	188	1252785	31.25	ng/ul	0.00
79) Pyrene-d10	12.74	212	1316857	30.01	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	1346345	31.44	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.40	94	66951	3.852	ng/ul 98
45) Dimethylphthalate	9.51	163	304497	9.230	ng/ul 99

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

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