

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101519\
 Data File : BP000724.D
 Acq On : 14 Oct 2019 21:19
 Operator : JU
 Sample : K5097-14
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JLMF1

Quant Time: Oct 15 03:55:02 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	261258	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	964852	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	514089	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.30	188	900471	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	787966	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	881452	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.36	96	36233	5.75	ng/uL	0.00
5) Phenol-d5	6.39	99	586779	29.12	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	299994	29.78	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	503129	29.63	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	463928	30.13	ng/ul	0.00
19) Nitrobenzene-d5	7.31	128	234247	31.94	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	250092	32.02	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	435589	29.14	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	592508	35.64	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	1103265	30.66	ng/ul	0.00
47) Acenaphthylene-d8	9.65	160	1364975	30.83	ng/ul	0.00
52) 4-Nitrophenol-d4	9.91	143	142804	25.76	ng/ul	0.00
58) Fluorene-d10	10.32	176	923711	30.00	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	88668	20.68	ng/ul	0.00
71) Anthracene-d10	11.35	188	1305985	31.31	ng/ul	0.00
79) Pyrene-d10	12.74	212	1350919	30.32	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	1380967	31.73	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.40	94	76020	3.742	ng/ul 98
45) Dimethylphthalate	9.51	163	340984	9.509	ng/ul 99
84) Bis(2-ethylhexyl)phthalate	13.96	149	67045	2.208	ng/ul 99

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

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