

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101519\
 Data File : BP000731.D
 Acq On : 15 Oct 2019 00:08
 Operator : JU
 Sample : K5096-07RX
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JLMH9RX

Quant Time: Oct 15 04:42:33 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	197397	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	749970	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	413675	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	770228	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	682490	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	777600	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.36	96	22411	4.71	ng/uL	0.00
5) Phenol-d5	6.39	99	87312	5.73	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	175415	23.04	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	254644	19.85	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	156327	13.44	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	140132	24.58	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	147829	24.35	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	262589	22.60	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	38884	3.01	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	790650	27.30	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	894378	25.11	ng/ul	0.00
52) 4-Nitrophenol-d4	9.92	143	24702	5.54	ng/ul	0.00
58) Fluorene-d10	10.32	176	661687	26.71	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	72131	19.67	ng/ul	0.00
71) Anthracene-d10	11.35	188	1023116	28.68	ng/ul	0.00
79) Pyrene-d10	12.73	212	1099896	28.50	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	1091925	28.44	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	6.40	94	25961	1.691	ng/ul 96
45) Dimethylphthalate	9.51	163	166673	5.776	ng/ul 99

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

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