

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP091019\
 Data File : BP000169.D
 Acq On : 10 Sep 2019 16:58
 Operator : HP/JU
 Sample : K4780-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0BM8

Quant Time: Sep 11 01:57:15 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP090519MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 10 02:43:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	363693	20.00	ng/ul	0.00
18) Naphthalene-d8	8.18	136	1407116	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.95	164	804387	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.45	188	1416640	20.00	ng/ul	0.00
78) Chrysene-d12	14.14	240	1105907	20.00	ng/ul	0.00
86) Perylene-d12	15.68	264	1124175	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.61	96	48076	4.69	ng/uL	0.00
5) Phenol-d5	6.50	99	217321	7.02	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.57	67	443912	27.40	ng/ul	0.00
9) 2-Chlorophenol-d4	6.66	132	587358	23.45	ng/ul	0.00
13) 4-Methylphenol-d8	7.26	113	370801	15.93	ng/ul	0.00
19) Nitrobenzene-d5	7.45	128	341757	31.99	ng/ul	0.00
22) 2-Nitrophenol-d4	7.78	143	346369	33.30	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	8.01	165	594070	26.91	ng/ul	0.00
29) 4-Chloroaniline-d4	8.23	131	121162	4.44	ng/ul	0.00
44) Dimethylphthalate-d6	9.63	166	1786810	30.08	ng/ul	0.00
47) Acenaphthylene-d8	9.79	160	2211084	30.12	ng/ul	0.00
52) 4-Nitrophenol-d4	10.02	143	71205	6.71	ng/ul	-0.01
58) Fluorene-d10	10.47	176	1507868	29.89	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.53	200	180887	27.78	ng/ul	0.00
71) Anthracene-d10	11.50	188	2258845	32.93	ng/ul	0.00
79) Pyrene-d10	12.90	212	2456647	37.35	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.59	264	1959225	33.62	ng/ul	0.00

Target Compounds					Ovalue	
45) Dimethylphthalate	9.65	163	202153	3.373	ng/ul	99

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

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