

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP091019\
 Data File : BP000174.D
 Acq On : 10 Sep 2019 19:00
 Operator : HP/JU
 Sample : K4780-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0BM5

Quant Time: Sep 11 01:58:16 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	435311	20.00	ng/ul	0.00
18) Naphthalene-d8	8.18	136	1658917	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.95	164	937404	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.45	188	1634382	20.00	ng/ul	0.00
78) Chrysene-d12	14.14	240	1275504	20.00	ng/ul	0.00
86) Perylene-d12	15.69	264	1314124	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.61	96	50594	4.12	ng/uL	0.00
5) Phenol-d5	6.50	99	227785	6.14	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.57	67	468934	24.19	ng/ul	0.00
9) 2-Chlorophenol-d4	6.66	132	612020	20.41	ng/ul	0.00
13) 4-Methylphenol-d8	7.26	113	394451	14.16	ng/ul	0.00
19) Nitrobenzene-d5	7.45	128	366522	29.10	ng/ul	0.00
22) 2-Nitrophenol-d4	7.78	143	381073	31.07	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	8.01	165	625448	24.03	ng/ul	0.00
29) 4-Chloroaniline-d4	8.23	131	174238	5.41	ng/ul	0.00
44) Dimethylphthalate-d6	9.63	166	1938688	28.01	ng/ul	0.00
47) Acenaphthylene-d8	9.79	160	2301476	26.91	ng/ul	0.00
52) 4-Nitrophenol-d4	10.02	143	72119	5.83	ng/ul	-0.01
58) Fluorene-d10	10.47	176	1584767	26.96	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.53	200	192579	25.64	ng/ul	0.00
71) Anthracene-d10	11.50	188	2302759	29.09	ng/ul	0.00
79) Pyrene-d10	12.90	212	2296181	30.27	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.59	264	1991377	29.24	ng/ul	0.00

Target Compounds					Ovalue	
45) Dimethylphthalate	9.65	163	232036	3.322	ng/ul	99

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

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