

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP091019\
 Data File : BP000171.D
 Acq On : 10 Sep 2019 17:47
 Operator : HP/JU
 Sample : K4780-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0BN6

Quant Time: Sep 11 01:57:31 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	415852	20.00	ng/ul	0.00
18) Naphthalene-d8	8.18	136	1602942	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.95	164	927519	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.45	188	1627715	20.00	ng/ul	0.00
78) Chrysene-d12	14.14	240	1283682	20.00	ng/ul	0.00
86) Perylene-d12	15.68	264	1279772	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.61	96	51456	4.39	ng/uL	0.00
5) Phenol-d5	6.50	99	227710	6.43	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.57	67	433445	23.40	ng/ul	0.00
9) 2-Chlorophenol-d4	6.66	132	583204	20.36	ng/ul	0.00
13) 4-Methylphenol-d8	7.26	113	385733	14.50	ng/ul	0.00
19) Nitrobenzene-d5	7.45	128	329695	27.09	ng/ul	0.00
22) 2-Nitrophenol-d4	7.78	143	336500	28.40	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	8.01	165	586303	23.31	ng/ul	0.00
29) 4-Chloroaniline-d4	8.23	131	133844	4.30	ng/ul	0.00
44) Dimethylphthalate-d6	9.63	166	1744087	25.47	ng/ul	0.00
47) Acenaphthylene-d8	9.79	160	2050780	24.23	ng/ul	0.00
52) 4-Nitrophenol-d4	10.02	143	85921	7.02	ng/ul	-0.01
58) Fluorene-d10	10.47	176	1408053	24.21	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.53	200	201224	26.90	ng/ul	0.00
71) Anthracene-d10	11.50	188	2290096	29.05	ng/ul	0.00
79) Pyrene-d10	12.90	212	2456576	32.18	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.59	264	2255734	34.01	ng/ul	0.00

Target Compounds					Ovalue
45) Dimethylphthalate	9.65	163	339216	4.908	ng/ul 100

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

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