

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP091019\
 Data File : BP000167.D
 Acq On : 10 Sep 2019 16:09
 Operator : HP/JU
 Sample : K4751-10
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Sep 11 01:57:00 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	328174	20.00	ng/ul	0.00
18) Naphthalene-d8	8.18	136	1275507	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.95	164	711322	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.45	188	1311553	20.00	ng/ul	0.00
78) Chrysene-d12	14.13	240	1021011	20.00	ng/ul	-0.01
86) Perylene-d12	15.67	264	1039783	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	2.61	96	83	0.01	ng/uL	0.00
5) Phenol-d5	6.51	99	15	0.00	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.57	67	43	0.00	ng/ul	0.00
9) 2-Chlorophenol-d4	6.66	132	17	0.00	ng/ul	0.00
13) 4-Methylphenol-d8	7.26	113	21	0.00	ng/ul	0.00
19) Nitrobenzene-d5	7.45	128	137	0.01	ng/ul	0.00
22) 2-Nitrophenol-d4	7.77	143	13	0.00	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	8.12	165	32	0.00	ng/ul	0.10
29) 4-Chloroaniline-d4	8.23	131	153	0.01	ng/ul	0.00
44) Dimethylphthalate-d6	9.63	166	206	0.00	ng/ul	0.00
47) Acenaphthylene-d8	9.95	160	326411	5.03	ng/ul	0.15
52) 4-Nitrophenol-d4	10.03	143	40	0.00	ng/ul	0.00
58) Fluorene-d10	10.46	176	1420	0.03	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.52	200	12	0.00	ng/ul	-0.01
71) Anthracene-d10	11.50	188	7585	0.12	ng/ul	-0.01
79) Pyrene-d10	12.89	212	120	0.00	ng/ul	-0.01
90) Benzo(a)pyrene-d12	15.67	264	1039471	19.29	ng/ul	0.09

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

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