

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP091019\
 Data File : BP000166.D
 Acq On : 10 Sep 2019 15:45
 Operator : HP/JU
 Sample : K4751-05
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Sep 11 01:56:52 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	341898	20.00	ng/ul	0.00
18) Naphthalene-d8	8.18	136	1327744	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.95	164	737084	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.45	188	1345125	20.00	ng/ul	0.00
78) Chrysene-d12	14.14	240	1029251	20.00	ng/ul	0.00
86) Perylene-d12	15.68	264	1044541	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.59	96	67	0.01	ng/uL	-0.01
5) Phenol-d5	6.50	99	59	0.00	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.57	67	119	0.01	ng/ul	0.00
9) 2-Chlorophenol-d4	6.65	132	60	0.00	ng/ul	-0.01
13) 4-Methylphenol-d8	7.26	113	6	0.00	ng/ul	0.00
19) Nitrobenzene-d5	7.47	128	11	0.00	ng/ul	0.02
22) 2-Nitrophenol-d4	7.78	143	68	0.01	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	8.01	165	25	0.00	ng/ul	0.00
29) 4-Chloroaniline-d4	8.23	131	880	0.03	ng/ul	0.00
44) Dimethylphthalate-d6	9.63	166	136	0.00	ng/ul	0.00
47) Acenaphthylene-d8	9.95	160	333206	4.95	ng/ul	0.15
52) 4-Nitrophenol-d4	10.03	143	24	0.00	ng/ul	0.00
58) Fluorene-d10	10.46	176	1446	0.03	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.52	200	11	0.00	ng/ul	0.00
71) Anthracene-d10	11.45	188	1345125	20.65	ng/ul	-0.06
79) Pyrene-d10	12.90	212	112	0.00	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.68	264	1043878	19.28	ng/ul	0.09

Target Compounds Ovalue

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

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