

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100117\
 Data File : BM011806.D
 Acq On : 01 Oct 2017 13:31
 Operator : SJ/JU
 Sample : PB102762BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK62

Quant Time: Oct 02 02:26:38 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM092717.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 02 02:22:21 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	111199	20.00	ng/ul	0.00
18) Naphthalene-d8	10.69	136	506370	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.52	164	311000	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.26	188	752428	20.00	ng/ul	0.00
75) Chrysene-d12	21.43	240	878306	20.00	ng/ul	0.00
83) Perylene-d12	23.76	264	1013642	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	11073	4.44	ng/uL	0.00
5) Phenol-d5	7.04	99	216876	19.83	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.21	67	179599	22.64	ng/ul	0.00
9) 2-Chlorophenol-d4	7.41	132	185921	22.85	ng/ul	0.00
13) 4-Methylphenol-d8	8.58	113	182845	21.34	ng/ul	0.00
19) Nitrobenzene-d5	9.05	128	88538	22.42	ng/ul	0.00
22) 2-Nitrophenol-d4	9.77	143	99772	23.68	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.30	165	173492	21.58	ng/ul	0.00
29) 4-Chloroaniline-d4	10.82	131	261657	29.44	ng/ul	0.00
43) Dimethylphthalate-d6	13.93	166	743107	27.86	ng/ul	0.00
46) Acenaphthylene-d8	14.22	160	758822	23.52	ng/ul	0.00
51) 4-Nitrophenol-d4	14.71	143	93114	20.42	ng/ul	0.00
57) Fluorene-d10	15.52	176	573373	24.96	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.63	200	104163	20.44	ng/ul	0.00
70) Anthracene-d10	17.36	188	1010450	26.39	ng/ul	0.00
76) Pyrene-d10	19.65	212	1252031	29.87	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.61	264	1336700	27.54	ng/ul	0.00

Target Compounds Ovalue

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

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