

Data Path : Z:\HPCHEM1\BNA M\DATA\BM093017\
 Data File : BM011797.D
 Acq On : 30 Sep 2017 11:31
 Operator : SJ/JU
 Sample : PB102757BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK57

Quant Time: Oct 01 03:28:57 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM092717.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 30 06:38:23 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	252283	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.73	136	1111005	20.00	ng/ul	-0.01
35) Acenaphthene-d10	14.56	164	676883	20.00	ng/ul	-0.02
61) Phenanthrene-d10	17.30	188	1603586	20.00	ng/ul	-0.02
75) Chrysene-d12	21.47	240	1574721	20.00	ng/ul	-0.01
83) Perylene-d12	23.82	264	1254244	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.35	96	19352	3.42	ng/uL	-0.01
5) Phenol-d5	7.07	99	370616	14.94	ng/ul	-0.02
7) Bis-(2-Chloroethyl)ether-d	7.25	67	287643	15.98	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.45	132	306909	16.63	ng/ul	-0.02
13) 4-Methylphenol-d8	8.62	113	289314	14.89	ng/ul	-0.02
19) Nitrobenzene-d5	9.08	128	147874	17.06	ng/ul	-0.02
22) 2-Nitrophenol-d4	9.80	143	165831	17.94	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	10.34	165	287429	16.29	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.86	131	409558	21.00	ng/ul	-0.01
43) Dimethylphthalate-d6	13.96	166	1060531	18.27	ng/ul	-0.01
46) Acenaphthylene-d8	14.25	160	1180182	16.81	ng/ul	-0.02
51) 4-Nitrophenol-d4	14.75	143	137940	13.90	ng/ul	-0.02
57) Fluorene-d10	15.55	176	845667	16.92	ng/ul	-0.02
62) 4,6-Dinitro-2-methylphenol	15.66	200	95075	8.75	ng/ul	-0.02
70) Anthracene-d10	17.40	188	1378670	16.89	ng/ul	-0.01
76) Pyrene-d10	19.69	212	1590779	21.17	ng/ul	-0.01
87) Benzo(a)pyrene-d12	23.67	264	1093636	18.21	ng/ul	-0.02

Target Compounds	Ovalue

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

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