

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092316\
 Data File : BM007530.D
 Acq On : 23 Sep 2016 19:14
 Operator : UM/SJ
 Sample : PB93536BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK36

Quant Time: Sep 24 02:24:27 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM092316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 24 02:24:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	216348	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	1024832	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.42	164	727134	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.16	188	1946827	20.00	ng/ul	0.00
75) Chrysene-d12	21.34	240	2901836	20.00	ng/ul	0.00
83) Perylene-d12	23.61	264	3107866	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.30	96	17863	4.06	ng/uL	0.00
5) Phenol-d5	6.96	99	352096	18.33	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.12	67	215049	20.83	ng/ul	0.00
9) 2-Chlorophenol-d4	7.32	132	281482	20.18	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	306939	18.73	ng/ul	0.00
19) Nitrobenzene-d5	8.94	128	152371	20.80	ng/ul	0.00
22) 2-Nitrophenol-d4	9.66	143	176970	21.26	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.19	165	311270	18.77	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	441779	25.37	ng/ul	0.00
43) Dimethylphthalate-d6	13.83	166	1309590	24.76	ng/ul	0.00
46) Acenaphthylene-d8	14.11	160	1438913	23.44	ng/ul	0.00
51) 4-Nitrophenol-d4	14.63	143	215046	23.03	ng/ul	0.00
57) Fluorene-d10	15.41	176	1106220	23.85	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.54	200	206490	17.38	ng/ul	0.00
70) Anthracene-d10	17.26	188	1934498	21.62	ng/ul	0.00
76) Pyrene-d10	19.55	212	2518136	21.87	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.46	264	2977994	21.66	ng/ul	0.00

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

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