

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021916\
 Data File : BM004283.D
 Acq On : 19 Feb 2016 12:55
 Operator : SJ/UM
 Sample : PB88413BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK13

Quant Time: Feb 19 23:33:51 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM021216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Feb 19 23:29:57 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	19709	20.00	ng/ul	0.00
18) Naphthalene-d8	10.42	136	97504	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.29	164	65169	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.04	188	160825	20.00	ng/ul	0.00
78) Chrysene-d12	21.26	240	193793	20.00	ng/ul	0.00
86) Perylene-d12	23.49	264	183153	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.09	96	2304	5.89	ng/uL	0.00
5) Phenol-d5	6.83	99	50740	32.05	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.97	67	28946	33.85	ng/ul	0.00
9) 2-Chlorophenol-d4	7.17	132	46230	32.93	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	48691	36.11	ng/ul	0.00
19) Nitrobenzene-d5	8.79	128	24765	33.03	ng/ul	0.00
22) 2-Nitrophenol-d4	9.51	143	29156	34.79	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.06	165	47368	31.34	ng/ul	0.00
29) 4-Chloroaniline-d4	10.57	131	75382	41.90	ng/ul	0.00
44) Dimethylphthalate-d6	13.71	166	212787	39.24	ng/ul	0.00
47) Acenaphthylene-d8	13.99	160	224341	34.05	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	31958	36.43	ng/ul	0.00
58) Fluorene-d10	15.29	176	167443	36.63	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.44	200	26158	26.71	ng/ul	0.00
71) Anthracene-d10	17.14	188	283393	36.04	ng/ul	0.00
79) Pyrene-d10	19.46	212	342259	33.27	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.34	264	364640	37.31	ng/ul	0.00

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

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