

Data Path : Z:\HPCHEM1\BNA M\DATA\BM021516\
 Data File : BM004268.D
 Acq On : 15 Feb 2016 20:29
 Operator : SJ/UM
 Sample : H1416-12
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9QE7

Quant Time: Feb 16 01:26:19 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM021216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 16 01:21:35 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	29992	20.00	ng/ul	0.00
18) Naphthalene-d8	10.43	136	139439	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.30	164	75902	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.06	188	149534	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	134097	20.00	ng/ul	0.00
86) Perylene-d12	23.50	264	128551	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.11	96	949	1.59	ng/uL	0.00
5) Phenol-d5	6.83	99	22852	9.49	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	42747	32.85	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	62786	29.39	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	42837	20.88	ng/ul	0.00
19) Nitrobenzene-d5	8.80	128	35356	32.98	ng/ul	0.00
22) 2-Nitrophenol-d4	9.52	143	40126	33.48	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.06	165	65589	30.35	ng/ul	0.00
29) 4-Chloroaniline-d4	10.58	131	7795	3.03	ng/ul	0.00
44) Dimethylphthalate-d6	13.72	166	208763	33.05	ng/ul	0.00
47) Acenaphthylene-d8	13.99	160	256098	33.37	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	7378	7.22	ng/ul	0.01
58) Fluorene-d10	15.30	176	174502	32.78	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.44	200	27464	30.16	ng/ul	0.00
71) Anthracene-d10	17.16	188	246898	33.77	ng/ul	0.00
79) Pyrene-d10	19.46	212	235832	33.13	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.36	264	226943	33.08	ng/ul	0.00

Target Compounds Ovalue

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

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