

Data Path : Z:\HPCHEM1\BNA M\DATA\BM021216\
 Data File : BM004236.D
 Acq On : 12 Feb 2016 20:01
 Operator : SJ/IZ
 Sample : PB88316TB
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLEB14

Quant Time: Feb 12 22:33:04 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM021216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Feb 12 22:29:16 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	21363	20.00	ng/ul	0.00
18) Naphthalene-d8	10.43	136	93301	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.30	164	52629	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.06	188	111253	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	113314	20.00	ng/ul	0.00
86) Perylene-d12	23.50	264	113903	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.11	96	2438	5.75	ng/uL	0.00
5) Phenol-d5	6.82	99	49204	28.67	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	28101	30.32	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	44903	29.51	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	42275	28.92	ng/ul	0.00
19) Nitrobenzene-d5	8.80	128	21199	29.55	ng/ul	0.00
22) 2-Nitrophenol-d4	9.52	143	23972	29.89	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.06	165	39479	27.30	ng/ul	0.00
29) 4-Chloroaniline-d4	10.57	131	62722	36.43	ng/ul	0.00
44) Dimethylphthalate-d6	13.72	166	152492	34.82	ng/ul	0.00
47) Acenaphthylene-d8	14.00	160	167533	31.49	ng/ul	0.00
52) 4-Nitrophenol-d4	14.53	143	21957	30.99	ng/ul	0.00
58) Fluorene-d10	15.30	176	122766	33.26	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.44	200	20360	30.05	ng/ul	0.00
71) Anthracene-d10	17.16	188	213570	39.26	ng/ul	0.00
79) Pyrene-d10	19.46	212	231306	38.46	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.36	264	250157	41.16	ng/ul	0.00

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

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