

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021216\
 Data File : BM004237.D
 Acq On : 12 Feb 2016 20:37
 Operator : SJ/IZ
 Sample : PB88346BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK46

Quant Time: Feb 12 22:33:12 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	21499	20.00	ng/ul	0.00
18) Naphthalene-d8	10.43	136	94144	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.30	164	53255	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.06	188	112342	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	113900	20.00	ng/ul	0.00
86) Perylene-d12	23.50	264	111927	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.11	96	2883	6.76	ng/uL	0.00
5) Phenol-d5	6.82	99	57917	33.54	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	33014	35.40	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	52492	34.27	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	50421	34.28	ng/ul	0.00
19) Nitrobenzene-d5	8.80	128	25343	35.01	ng/ul	0.00
22) 2-Nitrophenol-d4	9.52	143	27993	34.60	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.06	165	47148	32.31	ng/ul	0.00
29) 4-Chloroaniline-d4	10.57	131	72843	41.93	ng/ul	0.00
44) Dimethylphthalate-d6	13.72	166	164257	37.06	ng/ul	0.00
47) Acenaphthylene-d8	14.00	160	193972	36.03	ng/ul	0.00
52) 4-Nitrophenol-d4	14.53	143	22353	31.18	ng/ul	0.00
58) Fluorene-d10	15.30	176	136053	36.42	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.44	200	18967	27.72	ng/ul	0.00
71) Anthracene-d10	17.16	188	209299	38.11	ng/ul	0.00
79) Pyrene-d10	19.46	212	218369	36.12	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.36	264	227639	38.11	ng/ul	0.00

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

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