

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022616\
 Data File : BM004428.D
 Acq On : 26 Feb 2016 15:51
 Operator : SJ/UM
 Sample : H1564-10
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9R39

Quant Time: Feb 26 23:25:45 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM022216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Feb 26 23:20:20 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	22059	20.00	ng/ul	0.00
18) Naphthalene-d8	10.40	136	100752	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.28	164	63489	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	138182	20.00	ng/ul	0.00
78) Chrysene-d12	21.25	240	116908	20.00	ng/ul	0.00
86) Perylene-d12	23.47	264	109383	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	1661	3.97	ng/uL	0.00
5) Phenol-d5	6.80	99	35765	19.34	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.95	67	19524	20.75	ng/ul	0.00
9) 2-Chlorophenol-d4	7.15	132	33090	20.74	ng/ul	0.00
13) 4-Methylphenol-d8	8.34	113	31158	18.85	ng/ul	0.00
19) Nitrobenzene-d5	8.77	128	16192	21.12	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	18436	21.10	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	31883	20.19	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	46773	23.76	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	119870	22.01	ng/ul	0.00
47) Acenaphthylene-d8	13.97	160	144159	22.61	ng/ul	0.00
52) 4-Nitrophenol-d4	14.53	143	11103	12.69	ng/ul	0.01
58) Fluorene-d10	15.28	176	104894	22.65	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.43	200	11885	14.00	ng/ul	0.00
71) Anthracene-d10	17.13	188	156832	23.16	ng/ul	0.00
79) Pyrene-d10	19.44	212	157178	27.79	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.33	264	137239	23.91	ng/ul	0.00

Target Compounds					Ovalue
45) Dimethylphthalate	13.74	163	34876	6.14	ng/ul 99

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

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