

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011616\
 Data File : BM004059.D
 Acq On : 15 Jan 2016 22:01
 Operator : SJ/UM
 Sample : H1100-16
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BC9D0

Quant Time: Jan 16 03:07:37 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jan 16 03:00:58 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	34086	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	167471	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.32	164	97784	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	213104	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	185719	20.00	ng/ul	0.00
86) Perylene-d12	23.52	264	167389	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.13	96	3180	4.17	ng/uL	0.00
5) Phenol-d5	6.85	99	83333	29.45	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.00	67	48066	30.05	ng/ul	0.00
9) 2-Chlorophenol-d4	7.20	132	70171	30.17	ng/ul	0.00
13) 4-Methylphenol-d8	8.37	113	72841	31.76	ng/ul	0.00
19) Nitrobenzene-d5	8.82	128	37144	29.44	ng/ul	0.00
22) 2-Nitrophenol-d4	9.54	143	41020	29.67	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.08	165	70236	28.59	ng/ul	0.00
29) 4-Chloroaniline-d4	10.59	131	107950	33.28	ng/ul	0.00
44) Dimethylphthalate-d6	13.73	166	239467	31.19	ng/ul	0.00
47) Acenaphthylene-d8	14.01	160	293553	31.39	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	34681	25.15	ng/ul	0.00
58) Fluorene-d10	15.32	176	208978	32.09	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	22883	19.04	ng/ul	0.00
71) Anthracene-d10	17.17	188	315519	32.01	ng/ul	0.00
79) Pyrene-d10	19.47	212	329571	34.26	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	282960	33.85	ng/ul	0.00

Target Compounds					Qvalue
45) Dimethylphthalate	13.78	163	131152	16.76 ng/ul	99

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

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