

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031616\
 Data File : BM004634.D
 Acq On : 16 Mar 2016 13:50
 Operator : UM/SJ
 Sample : H1783-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AA1

Quant Time: Mar 17 01:35:49 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM031516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 17 00:56:57 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	42352	20.00	ng/ul	0.03
18) Naphthalene-d8	10.66	136	270531	20.00	ng/ul	0.02
36) Acenaphthene-d10	14.47	164	144095	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.22	188	325779	20.00	ng/ul	0.00
78) Chrysene-d12	21.40	240	369691	20.00	ng/ul	0.00
86) Perylene-d12	23.70	264	356136	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.34	96	2128	1.73	ng/uL	0.00
5) Phenol-d5	7.00	99	36877	10.60	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.20	67	102954	52.19	ng/ul	0.02
9) 2-Chlorophenol-d4	7.39	132	113355	40.46	ng/ul	0.01
13) 4-Methylphenol-d8	8.55	113	71491	25.48	ng/ul	0.00
19) Nitrobenzene-d5	9.01	128	65588	34.86	ng/ul	0.01
22) 2-Nitrophenol-d4	9.73	143	68907	33.91	ng/ul	0.01
26) 2,4-Dichlorophenol-d3	10.26	165	123686	31.03	ng/ul	0.00
29) 4-Chloroaniline-d4	10.78	131	7803	1.64	ng/ul	0.01
44) Dimethylphthalate-d6	13.89	166	420988	37.37	ng/ul	0.00
47) Acenaphthylene-d8	14.17	160	521294	37.74	ng/ul	0.00
52) 4-Nitrophenol-d4	14.66	143	11774	5.41	ng/ul	0.00
58) Fluorene-d10	15.47	176	370081	37.69	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.58	200	50191	29.32	ng/ul	0.00
71) Anthracene-d10	17.32	188	581756	38.79	ng/ul	0.00
79) Pyrene-d10	19.61	212	675906	40.05	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.56	264	696103	43.76	ng/ul	0.00

Target Compounds

						Qvalue
20) Nitrobenzene	9.05	77	7422	1.60	ng/ul	97
23) 2-Nitrophenol	9.76	139	6395	2.84	ng/ul	92
28) Naphthalene	10.70	128	64140	4.55	ng/ul	99
37) 1,2,4,5-Tetrachlorobenzene	12.66	216	105769	22.98	ng/ul	98
73) 1,2,3,4-Tetrachlorobenzene	13.27	216	424874	94.79	ng/uL	100
74) Pentachlorobenzene	14.79	250	13304	3.08	ng/uL	98

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

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