

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM032916\
 Data File : BM004818.D
 Acq On : 29 Mar 2016 13:22
 Operator : UM/SJ
 Sample : H1939-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9SF9

Quant Time: Mar 30 03:15:33 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM031516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 30 02:30:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	39952	20.00	ng/ul	0.00
18) Naphthalene-d8	10.62	136	201692	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.46	164	133760	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.20	188	334144	20.00	ng/ul	0.00
78) Chrysene-d12	21.39	240	448966	20.00	ng/ul	0.00
86) Perylene-d12	23.69	264	482996	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.32	96	1093	0.94	ng/uL	0.00
5) Phenol-d5	6.99	99	24331	7.41	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.16	67	58641	31.52	ng/ul	0.00
9) 2-Chlorophenol-d4	7.36	132	61561	23.29	ng/ul	0.00
13) 4-Methylphenol-d8	8.52	113	46182	17.45	ng/ul	0.00
19) Nitrobenzene-d5	8.98	128	37978	27.08	ng/ul	0.00
22) 2-Nitrophenol-d4	9.70	143	40600	26.80	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.23	165	73518	24.74	ng/ul	0.00
29) 4-Chloroaniline-d4	10.76	131	2110	0.59	ng/ul	0.01
44) Dimethylphthalate-d6	13.87	166	307111	29.37	ng/ul	0.00
47) Acenaphthylene-d8	14.16	160	365807	28.53	ng/ul	0.00
52) 4-Nitrophenol-d4	14.65	143	11093	5.49	ng/ul	0.00
58) Fluorene-d10	15.46	176	277865	30.49	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.57	200	41973	23.90	ng/ul	0.00
71) Anthracene-d10	17.30	188	462722	30.08	ng/ul	0.00
79) Pyrene-d10	19.60	212	570243	27.82	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.54	264	710137	32.92	ng/ul	0.00

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

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