

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040116\
 Data File : BM004857.D
 Acq On : 02 Apr 2016 01:28
 Operator : UM/SJ
 Sample : H2004-05
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AF8

Quant Time: Apr 02 02:15:38 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM040116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Apr 02 00:53:13 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	108095	20.00	ng/ul	0.00
18) Naphthalene-d8	10.62	136	499167	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.46	164	313856	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.20	188	736709	20.00	ng/ul	0.00
78) Chrysene-d12	21.39	240	871911	20.00	ng/ul	0.00
86) Perylene-d12	23.69	264	920063	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.32	96	1197	0.45	ng/uL	0.00
5) Phenol-d5	6.99	99	25044	2.61	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.16	67	78133	15.01	ng/ul	0.00
9) 2-Chlorophenol-d4	7.36	132	81683	11.20	ng/ul	0.00
13) 4-Methylphenol-d8	8.52	113	52864	6.55	ng/ul	0.00
19) Nitrobenzene-d5	8.98	128	54099	15.27	ng/ul	0.00
22) 2-Nitrophenol-d4	9.70	143	61637	15.65	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.23	165	99536	13.56	ng/ul	0.00
29) 4-Chloroaniline-d4	10.76	131	1848	0.21	ng/ul	0.00
44) Dimethylphthalate-d6	13.87	166	397508	16.05	ng/ul	0.00
47) Acenaphthylene-d8	14.15	160	471082	15.90	ng/ul	0.00
52) 4-Nitrophenol-d4	14.66	143	9443	1.93	ng/ul	0.00
58) Fluorene-d10	15.45	176	347433	16.23	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.57	200	60531	13.71	ng/ul	0.00
71) Anthracene-d10	17.30	188	542420	16.46	ng/ul	0.00
79) Pyrene-d10	19.60	212	631475	16.68	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.54	264	739230	18.13	ng/ul	0.00

Target Compounds	Ovalue

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

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