

Data Path : Z:\HPCHEM1\BNA M\DATA\BM050616\
 Data File : BM005297.D
 Acq On : 07 May 2016 07:36
 Operator : UM/SJ
 Sample : H2743-06
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H4005

Quant Time: May 08 00:33:07 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM050516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat May 07 07:05:19 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	58454	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	284001	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.41	164	194656	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.16	188	467935	20.00	ng/ul	0.00
75) Chrysene-d12	21.35	240	548945	20.00	ng/ul	0.00
83) Perylene-d12	23.63	264	510992	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	6.96	99	3371	0.64	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.10	67	13508	4.47	ng/ul	0.00
9) 2-Chlorophenol-d4	7.31	132	13927	3.48	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	7909	1.80	ng/ul	0.00
19) Nitrobenzene-d5	8.93	128	8104	4.00	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	8238	3.59	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.19	165	17766	4.16	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
43) Dimethylphthalate-d6	13.82	166	93826	6.01	ng/ul	0.00
46) Acenaphthylene-d8	14.10	160	105988	5.79	ng/ul	0.00
51) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
57) Fluorene-d10	15.40	176	94453	7.01	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.53	200	8186	3.11	ng/ul	0.00
70) Anthracene-d10	17.26	188	162842	7.87	ng/ul	0.00
76) Pyrene-d10	19.56	212	194708	7.68	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.48	264	172468	7.62	ng/ul	0.00

Target Compounds	Ovalue

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

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