

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050616\
 Data File : BM005296.D
 Acq On : 07 May 2016 07:00
 Operator : UM/SJ
 Sample : H2743-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H4003

Quant Time: May 08 00:33:01 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	63426	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	314230	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.41	164	214185	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.16	188	512784	20.00	ng/ul	0.00
75) Chrysene-d12	21.35	240	589151	20.00	ng/ul	0.00
83) Perylene-d12	23.62	264	532753	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	3448	0.60	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.11	67	13426	4.09	ng/ul	0.00
9) 2-Chlorophenol-d4	7.31	132	12886	2.97	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	7644	1.61	ng/ul	0.00
19) Nitrobenzene-d5	8.93	128	7398	3.30	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	8366	3.29	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.19	165	17014	3.60	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
43) Dimethylphthalate-d6	13.82	166	92864	5.41	ng/ul	0.00
46) Acenaphthylene-d8	14.10	160	104050	5.17	ng/ul	0.00
51) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
57) Fluorene-d10	15.40	176	91617	6.18	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.53	200	8020	2.78	ng/ul	0.00
70) Anthracene-d10	17.26	188	153927	6.79	ng/ul	0.00
76) Pyrene-d10	19.56	212	186542	6.86	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.47	264	158999	6.74	ng/ul	0.00

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

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