

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051116\
 Data File : BM005405.D
 Acq On : 12 May 2016 02:09
 Operator : UM/SJ
 Sample : H2846-02
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC8

Quant Time: May 12 05:32:03 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM050516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 12 02:20:05 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	67215	20.00	ng/ul	0.00
18) Naphthalene-d8	10.55	136	314873	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.40	164	172065	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.15	188	323896	20.00	ng/ul	0.00
75) Chrysene-d12	21.34	240	294947	20.00	ng/ul	0.00
83) Perylene-d12	23.61	264	290267	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.26	96	1377	0.96	ng/uL	0.00
5) Phenol-d5	6.93	99	33730	5.53	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.09	67	91506	26.31	ng/ul	0.00
9) 2-Chlorophenol-d4	7.29	132	101777	22.10	ng/ul	0.00
13) 4-Methylphenol-d8	8.47	113	67917	13.48	ng/ul	0.00
19) Nitrobenzene-d5	8.92	128	64795	28.82	ng/ul	0.00
22) 2-Nitrophenol-d4	9.63	143	77003	30.25	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.17	165	125958	26.63	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	5022	0.88	ng/ul	0.01
43) Dimethylphthalate-d6	13.81	166	412400	29.90	ng/ul	0.00
46) Acenaphthylene-d8	14.09	160	501216	30.98	ng/ul	0.00
51) 4-Nitrophenol-d4	14.63	143	10008	3.98	ng/ul	0.01
57) Fluorene-d10	15.40	176	338447	28.42	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.53	200	50488	27.71	ng/ul	0.00
70) Anthracene-d10	17.25	188	459542	32.10	ng/ul	0.00
76) Pyrene-d10	19.55	212	419223	30.79	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.46	264	395221	30.76	ng/ul	0.00

Target Compounds						Qvalue
6) Phenol	6.96	94	84815	13.46	ng/ul	100
16) 4-Methylphenol	8.53	108	24345	4.50	ng/ul	93
44) Dimethylphthalate	13.86	163	40823	2.96	ng/ul	99

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

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