

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062516\
 Data File : BM006009.D
 Acq On : 25 Jun 2016 19:50
 Operator : UM/SJ
 Sample : H3670-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9Y75

Quant Time: Jun 26 04:42:28 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Jun 26 04:39:40 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	172337	20.00	ng/ul	0.00
7) Naphthalene-d8	10.44	136	969067	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.30	164	727355	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.05	188	1982124	20.00	ng/ul	0.00
29) Chrysene-d12	21.24	240	2581201	20.00	ng/ul	0.00
34) Perylene-d12	23.46	264	2856089	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.18	96	4715	1.22	ng/uL	0.00
3) Phenol-d5	6.83	99	166406	9.18	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.99	67	125048	12.40	ng/ul	0.00
5) 2-Chlorophenol-d4	7.19	132	134852	10.68	ng/ul	0.00
6) 4-Methylphenol-d8	8.36	113	118549	7.75	ng/ul	0.00
8) Nitrobenzene-d5	8.81	128	79237	10.74	ng/ul	0.00
9) 2-Nitrophenol-d4	9.53	143	84232	10.26	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.06	165	156588	9.94	ng/ul	0.00
12) 4-Chloroaniline-d4	10.57	131	227621	13.32	ng/ul	0.00
16) Dimethylphthalate-d6	13.72	166	714383	11.47	ng/ul	0.00
17) Acenaphthylene-d8	13.99	160	801941	11.10	ng/ul	0.00
20) 4-Nitrophenol-d4	14.50	143	118960	8.96	ng/ul	0.00
21) Fluorene-d10	15.30	176	631801	11.88	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.42	200	90742	8.25	ng/ul	0.00
26) Anthracene-d10	17.15	188	1082569	11.89	ng/ul	0.00
30) Pyrene-d10	19.45	212	1398152	12.37	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.32	264	1600741	12.41	ng/ul	0.00

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

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