

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062516\
 Data File : BM006003.D
 Acq On : 25 Jun 2016 16:14
 Operator : UM/SJ
 Sample : H3669-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9Y62

Quant Time: Jun 26 04:41:32 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	209114	20.00	ng/ul	0.00
7) Naphthalene-d8	10.44	136	1098945	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.30	164	768824	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.05	188	1989612	20.00	ng/ul	0.00
29) Chrysene-d12	21.24	240	2591003	20.00	ng/ul	0.00
34) Perylene-d12	23.47	264	2885760	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.17	96	9455	2.02	ng/uL	0.00
3) Phenol-d5	6.83	99	376155	17.10	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.99	67	279462	22.83	ng/ul	0.00
5) 2-Chlorophenol-d4	7.19	132	312931	20.42	ng/ul	0.00
6) 4-Methylphenol-d8	8.36	113	246800	13.29	ng/ul	0.00
8) Nitrobenzene-d5	8.81	128	186141	22.25	ng/ul	0.00
9) 2-Nitrophenol-d4	9.53	143	212356	22.81	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.06	165	365544	20.47	ng/ul	0.00
12) 4-Chloroaniline-d4	10.57	131	518110	26.74	ng/ul	0.00
16) Dimethylphthalate-d6	13.72	166	1486736	22.59	ng/ul	0.00
17) Acenaphthylene-d8	13.99	160	1764111	23.11	ng/ul	0.00
20) 4-Nitrophenol-d4	14.50	143	287881	20.50	ng/ul	0.00
21) Fluorene-d10	15.30	176	1295309	23.05	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.42	200	241416	21.86	ng/ul	0.00
26) Anthracene-d10	17.15	188	2129001	23.30	ng/ul	0.00
30) Pyrene-d10	19.45	212	2648744	23.34	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.32	264	3003437	23.05	ng/ul	0.00

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

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