

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062716\
 Data File : BM006081.D
 Acq On : 27 Jun 2016 21:57
 Operator : UM/SJ
 Sample : H3770-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SVOC-GPCBLANK

Quant Time: Jun 28 02:14:27 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 28 02:11:41 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	148609	20.00	ng/ul	0.00
7) Naphthalene-d8	10.44	136	797917	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.31	164	562003	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.06	188	1430670	20.00	ng/ul	0.00
29) Chrysene-d12	21.25	240	1735272	20.00	ng/ul	0.00
34) Perylene-d12	23.47	264	1527320	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
3) Phenol-d5	0.00	99	0	0.00	ng/ul	
4) Bis-(2-Chloroethyl)ether-d	7.07	67	217	0.02	ng/ul	0.07
5) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
6) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
8) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul	
9) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
10) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul	
12) 4-Chloroaniline-d4	10.45	131	404	0.03	ng/ul	-0.13
16) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
17) Acenaphthylene-d8	14.31	160	267774	4.80	ng/ul	0.31
20) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
21) Fluorene-d10	15.30	176	3781	0.09	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
26) Anthracene-d10	17.06	188	1430670	21.77	ng/ul	-0.10
30) Pyrene-d10	19.46	212	1521	0.02	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.47	264	1527320	22.15	ng/ul	0.14

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

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