

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062516\
 Data File : BM005998.D
 Acq On : 25 Jun 2016 13:14
 Operator : UM/SJ
 Sample : SSTD16064
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD16064

Quant Time: Jun 26 04:35:01 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Jun 26 04:20:44 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	104987	20.00	ng/ul	0.00
7) Naphthalene-d8	10.44	136	603452	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.31	164	426621	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.06	188	1124085	20.00	ng/ul	0.00
29) Chrysene-d12	21.26	240	1258569	20.00	ng/ul	0.00
34) Perylene-d12	23.50	264	1643038	20.00	ng/ul	0.04

System Monitoring Compounds

2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
3) Phenol-d5	6.84	99	1858128	168.24	ng/ul	0.01
4) Bis-(2-Chloroethyl)ether-d	7.00	67	948792	154.42	ng/ul	0.00
5) 2-Chlorophenol-d4	0.00	132	0d	0.00	ng/ul	
6) 4-Methylphenol-d8	8.37	113	1552327	166.51	ng/ul	0.01
8) Nitrobenzene-d5	0.00	128	0d	0.00	ng/ul	
9) 2-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
10) 2,4-Dichlorophenol-d3	0.00	165	0d	0.00	ng/ul	
12) 4-Chloroaniline-d4	10.58	131	1029330	96.74	ng/ul	0.00
16) Dimethylphthalate-d6	0.00	166	0d	0.00	ng/ul	
17) Acenaphthylene-d8	0.00	160	0d	0.00	ng/ul	
20) 4-Nitrophenol-d4	14.55	143	1147038	147.23	ng/ul	0.04
21) Fluorene-d10	0.00	176	0d	0.00	ng/ul	
24) 4,6-Dinitro-2-methylphenol	15.45	200	978637	156.83	ng/ul	0.02
26) Anthracene-d10	0.00	188	0d	0.00	ng/ul	
30) Pyrene-d10	0.00	212	0d	0.00	ng/ul	
37) Benzo(a)pyrene-d12	0.00	264	0d	0.00	ng/ul	

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

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