

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062816\
 Data File : BM006098.D
 Acq On : 28 Jun 2016 16:33
 Operator : UM/SJ
 Sample : SSTD16047
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD16047

Quant Time: Jun 28 17:21:17 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062816.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 28 15:37:16 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	153418	20.00	ng/ul	0.00
7) Naphthalene-d8	10.44	136	731200	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.31	164	456903	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.06	188	1107387	20.00	ng/ul	0.00
29) Chrysene-d12	21.26	240	1019986	20.00	ng/ul	0.01
34) Perylene-d12	23.47	264	1148257	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
3) Phenol-d5	6.84	99	2160239	171.42	ng/ul	0.01
4) Bis-(2-Chloroethyl)ether-d	7.00	67	1228854	163.84	ng/ul	0.00
5) 2-Chlorophenol-d4	0.00	132	0d	0.00	ng/ul	
6) 4-Methylphenol-d8	8.38	113	1734676	162.93	ng/ul	0.02
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	827109	68.37	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	1044827	250.56	ng/ul	0.04
21) Fluorene-d10	0.00	176	0d	0.00	ng/ul	
24) 4,6-Dinitro-2-methylphenol	15.45	200	854078	172.39	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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