

Data Path : Z:\HPCHEM1\BNA_B\DATA\BB101916\
 Data File : BB058651.D
 Acq On : 19 Oct 2016 14:01
 Operator : UM/SJ
 Sample : SSTD16095
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_B
 ClientSampleId :
 SSTD16095

Quant Time: Oct 19 14:42:46 2016
 Quant Method : Z:\HPCHEM1\BNA_B\METHOD\SOM02.2-EPA-BB101916-MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 17 17:52:13 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.70	152	550136	20.00	ng/ul	0.04
7) Naphthalene-d8	11.66	136	2200077	20.00	ng/ul	0.04
13) Acenaphthene-d10	15.61	164	1288465	20.00	ng/ul	0.04
18) Phenanthrene-d10	18.46	188	1989205	20.00	ng/ul	0.04
23) Chrysene-d12	22.85	240	2343069	20.00	ng/ul	0.06
25) Perylene-d12	25.93	264	1623629	20.00	ng/ul	0.08

System Monitoring Compounds

2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
3) Phenol-d5	7.89	99	7025277	175.85	ng/ul	0.08
4) Bis-(2-Chloroethyl)ether-d	7.98	67	4542298	169.74	ng/ul	0.04
5) 2-Chlorophenol-d4	0.00	132	0d	0.00	ng/ul	
6) 4-Methylphenol-d8	9.52	113	5768778	174.40	ng/ul	0.08
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	
11) 4-Chloroaniline-d4	11.81	131	5863433	140.39	ng/ul	0.08
14) Dimethylphthalate-d6	0.00	166	0d	0.00	ng/ul	
15) Acenaphthylene-d8	0.00	160	0d	0.00	ng/ul	
16) 4-Nitrophenol-d4	15.92	143	3733641	189.76	ng/ul	0.15
17) Fluorene-d10	0.00	176	0d	0.00	ng/ul	
19) 4,6-Dinitro-2-methylphenol	16.84	200	4241493	200.65	ng/ul	0.11
20) Anthracene-d10	0.00	188	0d	0.00	ng/ul	
24) Pyrene-d10	0.00	212	0d	0.00	ng/ul	
26) 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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