

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080916\
 Data File : BM006994.D
 Acq On : 09 Aug 2016 22:32
 Operator : UM/SJ
 Sample : SSTD16039
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD16039

Quant Time: Aug 10 01:28:24 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM080916.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 10 01:21:42 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	296298	20.00	ng/ul	0.00
7) Naphthalene-d8	10.62	136	1401983	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.46	164	933325	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.20	188	2300838	20.00	ng/ul	0.00
29) Chrysene-d12	21.39	240	2360891	20.00	ng/ul	0.01
34) Perylene-d12	23.67	264	2165905	20.00	ng/ul	0.01

System Monitoring Compounds

2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
3) Phenol-d5	7.00	99	4183066	148.70	ng/ul	0.02
4) Bis-(2-Chloroethyl)ether-d	7.16	67	2206710	128.63	ng/ul	0.00
5) 2-Chlorophenol-d4	0.00	132	0d	0.00	ng/ul	
6) 4-Methylphenol-d8	8.54	113	3545362	150.57	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	10.76	131	3714379	131.26	ng/ul	0.02
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	14.68	143	2192216	170.38	ng/ul	0.04
21) Fluorene-d10	0.00	176	0d	0.00	ng/ul	
24) 4,6-Dinitro-2-methylphenol	15.59	200	2156996	204.68	ng/ul	0.03
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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