

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061016\
 Data File : BM005719.D
 Acq On : 11 Jun 2016 00:40
 Operator : UM/SJ
 Sample : SSTD16034
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD16034

Quant Time: Jun 11 05:34:32 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jun 11 05:17:22 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	78066	20.00	ng/ul	0.00
7) Naphthalene-d8	10.54	136	391315	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.39	164	227121	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.14	188	516815	20.00	ng/ul	0.00
29) Chrysene-d12	21.33	240	438983	20.00	ng/ul	0.01
34) Perylene-d12	23.57	264	402058	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	6.94	99	1100876	161.18	ng/ul	0.01
4) Bis-(2-Chloroethyl)ether-d	7.09	67	600781	144.74	ng/ul	0.01
5) 2-Chlorophenol-d4	0.00	132	0d	0.00	ng/ul	
6) 4-Methylphenol-d8	8.49	113	912400	166.08	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	10.68	131	657960	94.65	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	14.66	143	444877	173.23	ng/ul	0.04
21) Fluorene-d10	0.00	176	0d	0.00	ng/ul	
24) 4,6-Dinitro-2-methylphenol	15.54	200	439357	175.15	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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