

Data Path : Z:\HPCHEM1\BNA M\DATA\BM060216\
 Data File : BM005655.D
 Acq On : 02 Jun 2016 09:19
 Operator : UM/SJ
 Sample : SSTD16039
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD16039

Quant Time: Jun 02 16:43:19 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM060216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 02 16:26:07 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	105016	20.00	ng/ul	0.00
7) Naphthalene-d8	10.56	136	467912	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.40	164	227894	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.15	188	415318	20.00	ng/ul	0.00
29) Chrysene-d12	21.33	240	344317	20.00	ng/ul	0.00
34) Perylene-d12	23.59	264	369154	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
3) Phenol-d5	6.95	99	1449608	156.33	ng/ul	0.01
4) Bis-(2-Chloroethyl)ether-d	7.10	67	837695	169.61	ng/ul	0.01
5) 2-Chlorophenol-d4	0.00	132	0d	0.00	ng/ul	
6) 4-Methylphenol-d8	8.50	113	1133429	137.22	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	10.70	131	866782	177.77	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	14.65	143	398287	128.85	ng/ul	0.03
21) Fluorene-d10	0.00	176	0d	0.00	ng/ul	
24) 4,6-Dinitro-2-methylphenol	15.54	200	358540	173.95	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	Ovalue
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

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