

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102016\
 Data File : BG024462.D
 Acq On : 20 Oct 2016 14:36
 Operator : UM/SJ
 Sample : SSTD16012
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTD16012

Manual Integrations
 APPROVED

sohil
 10/21/2016 6:56:58 PM

Quant Time: Oct 21 08:17:37 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102016-MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 21 07:54:54 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.28	152	142756	20.00	ng/ul	0.00
7) Naphthalene-d8	11.11	136	633907	20.00	ng/ul	0.00
13) Acenaphthene-d10	14.90	164	491708	20.00	ng/ul	0.00
18) Phenanthrene-d10	17.64	188	968552	20.00	ng/ul	0.00
23) Chrysene-d12	21.95	240	1138889	20.00	ng/ul	0.01
25) Perylene-d12	25.38	264	1154830	20.00	ng/ul	0.01

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	7.42	99	2033694	183.86	ng/ul	0.01
4) Bis-(2-Chloroethyl)ether-d	7.60	67	1249905	190.66	ng/ul	0.01
5) 2-Chlorophenol-d4	0.00	132	0d	0.00	ng/ul	
6) 4-Methylphenol-d8	8.98	113	1732632	199.32	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	
11) 4-Chloroaniline-d4	11.25	131	1483263	135.91	ng/ul	0.00
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	15.09	143	940020m	166.66	ng/ul	-0.03
17) Fluorene-d10	0.00	176	0d	0.00	ng/ul	
19) 4,6-Dinitro-2-methylphenol	16.01	200	1140616	194.34	ng/ul	0.02
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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