

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102016\
 Data File : BG024481.D
 Acq On : 21 Oct 2016 3:02
 Operator : UM/SJ
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02015

Quant Time: Oct 21 18:03:58 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102016-MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 21 17:27:43 2016
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	100	0.00
2 S	1,4-Dioxane-d8	0.374	0.363	2.9	103	0.00
3 S	Phenol-d5	1.819	1.839	-1.1	104	0.00
4 S	Bis-(2-Chloroethyl)ether-d8	1.124	1.099	2.2	96	0.00
5 S	2-Chlorophenol-d4	1.280	1.258	1.7	99	0.00
6 S	4-Methylphenol-d8	1.517	1.506	0.7	101	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
8 S	Nitrobenzene-d5	0.150	0.146	2.7	103	0.00
9 S	2-Nitrophenol-d4	0.174	0.171	1.7	97	0.00
10 S	2,4-Dichlorophenol-d3	0.353	0.353	0.0	100	0.00
11 S	4-Chloroaniline-d4	0.367	0.393	-7.1	103	0.00
12	1-Methylnaphthalene	0.743	0.748	-0.7	101	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
14 S	Dimethylphthalate-d6	1.349	1.336	1.0	103	0.00
15 S	Acenaphthylene-d8	1.605	1.583	1.4	99	0.00
16 S	4-Nitrophenol-d4	0.231	0.221	4.3	104	0.00
17 S	Fluorene-d10	1.285	1.274	0.9	106	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
19 S	4,6-Dinitro-2-methylphenol-	0.134	0.127	5.2	105	0.00
20 S	Anthracene-d10	0.889	0.914	-2.8	103	0.00
21	1,2,3,4-Tetrachlorobenzene	0.305	0.314	-3.0	101	0.00
22	Pentachlorobenzene	0.320	0.318	0.6	101	0.00
23 I	Chrysene-d12	1.000	1.000	0.0	101	0.00
24 S	Pyrene-d10	0.869	0.894	-2.9	101	0.00
25 I	Perylene-d12	1.000	1.000	0.0	102	0.00
26 S	Benzo(a)pyrene-d12	0.875	0.873	0.2	102	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0