

Data Path : Z:\HPCHEM1\BNA_M\Data\BM060616\
 Data File : BM005688.D
 Acq On : 07 Jun 2016 02:25
 Operator : UM/SJ
 Sample : SSTDC0020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02047

Quant Time: Jun 07 04:29:32 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM060216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 07 04:25:20 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	128	0.00
2 S	1,4-Dioxane-d8	0.476	0.450	5.5	124	0.00
3 S	Phenol-d5	1.758	1.682	4.3	122	0.01
4 S	Bis-(2-Chloroethyl)ether-d8	1.070	1.049	2.0	122	0.00
5 S	2-Chlorophenol-d4	1.451	1.411	2.8	125	0.01
6 S	4-Methylphenol-d8	1.391	1.262	9.3	116	0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	112	0.00
8 S	Nitrobenzene-d5	0.154	0.161	-4.5	118	0.01
9 S	2-Nitrophenol-d4	0.164	0.161	1.8	109	0.00
10 S	2,4-Dichlorophenol-d3	0.292	0.286	2.1	111	0.01
11	Naphthalene	1.017	1.007	1.0	112	0.01
12 S	4-Chloroaniline-d4	0.347	0.381	-9.8	110	0.01
13	2-Methylnaphthalene	0.707	0.649	8.2	104	0.00
14	1-Methylnaphthalene	0.695	0.630	9.4	103	0.01
15 I	Acenaphthene-d10	1.000	1.000	0.0	96	0.01
16 S	Dimethylphthalate-d6	1.538	1.543	-0.3	97	0.00
17 S	Acenaphthylene-d8	2.000	2.043	-2.2	98	0.01
18	Acenaphthylene	2.069	2.083	-0.7	96	0.01
19 C	Acenaphthene	1.343	1.350	-0.5	97	0.01
20 S	4-Nitrophenol-d4	0.220	0.211	4.1	94	0.02
21 S	Fluorene-d10	1.321	1.297	1.8	94	0.01
22	Fluorene	1.453	1.436	1.2	93	0.00
23 I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
24 S	4,6-Dinitro-2-methylphenol-	0.094	0.020	78.7#	22	0.02
25	Phenanthrene	1.104	1.094	0.9	94	0.01
26 S	Anthracene-d10	0.950	0.954	-0.4	95	0.00
27	Anthracene	1.122	1.128	-0.5	95	0.00
28 C	Fluoranthene	1.128	1.214	-7.6	101	0.00
29 I	Chrysene-d12	1.000	1.000	0.0	103	0.01
30 S	Pyrene-d10	1.005	0.959	4.6	99	0.00
31	Pyrene	1.262	1.200	4.9	99	0.00
32	Benzo(a)anthracene	1.173	1.179	-0.5	103	0.00
33	Chrysene	1.108	1.110	-0.2	105	0.01
34 I	Perylene-d12	1.000	1.000	0.0	112	0.01
35	Benzo(b)fluoranthene	1.206	1.197	0.7	113	0.01
36	Benzo(k)fluoranthene	1.132	1.079	4.7	106	0.01
37 S	Benzo(a)pyrene-d12	0.920	0.908	1.3	111	0.02
38 C	Benzo(a)pyrene	1.161	1.142	1.6	110	0.01
39	Indeno(1,2,3-cd)pyrene	1.374	1.360	1.0	110	0.03
40	Dibenzo(a,h)anthracene	1.153	1.137	1.4	109	0.02
41	Benzo(g,h,i)perylene	1.152	1.144	0.7	113	0.04

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(#) = Out of Range

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