

Data Path : Z:\HPCHEM1\BNA M\DATA\BM060916\
 Data File : BM005707.D
 Acq On : 09 Jun 2016 10:05
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02050

Quant Time: Jun 10 01:19:28 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM060216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 08 01:24:32 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	85	0.00
2 S	1,4-Dioxane-d8	0.476	0.419	12.0	76	0.00
3 S	Phenol-d5	1.758	1.780	-1.3	85	0.00
4 S	Bis-(2-Chloroethyl)ether-d8	1.070	1.053	1.6	82	0.00
5 S	2-Chlorophenol-d4	1.451	1.446	0.3	85	0.00
6 S	4-Methylphenol-d8	1.391	1.505	-8.2	92	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	93	0.00
8 S	Nitrobenzene-d5	0.154	0.152	1.3	93	0.00
9 S	2-Nitrophenol-d4	0.164	0.170	-3.7	96	0.00
10 S	2,4-Dichlorophenol-d3	0.292	0.304	-4.1	98	0.00
11	Naphthalene	1.017	0.987	2.9	91	0.00
12 S	4-Chloroaniline-d4	0.347	0.424	-22.2#	102	0.00
13	2-Methylnaphthalene	0.707	0.735	-4.0	98	0.00
14	1-Methylnaphthalene	0.695	0.730	-5.0	99	0.00
15 I	Acenaphthene-d10	1.000	1.000	0.0	109	0.00
16 S	Dimethylphthalate-d6	1.538	1.616	-5.1	116	0.00
17 S	Acenaphthylene-d8	2.000	1.997	0.1	109	0.00
18	Acenaphthylene	2.069	2.043	1.3	107	0.00
19 C	Acenaphthene	1.343	1.330	1.0	109	0.00
20 S	4-Nitrophenol-d4	0.220	0.242	-10.0	123	0.00
21 S	Fluorene-d10	1.321	1.394	-5.5	116	0.00
22	Fluorene	1.453	1.521	-4.7	113	0.00
23 I	Phenanthrene-d10	1.000	1.000	0.0	129	0.00
24 S	4,6-Dinitro-2-methylphenol-	0.094	0.077	18.1	115	0.00
25	Phenanthrene	1.104	1.098	0.5	128	0.00
26 S	Anthracene-d10	0.950	0.941	0.9	127	0.00
27	Anthracene	1.122	1.122	0.0	128	0.00
28 C	Fluoranthene	1.128	1.246	-10.5	141	0.00
29 I	Chrysene-d12	1.000	1.000	0.0	140	0.00
30 S	Pyrene-d10	1.005	0.972	3.3	137	0.00
31	Pyrene	1.262	1.240	1.7	139	0.00
32	Benzo(a)anthracene	1.173	1.168	0.4	139	0.00
33	Chrysene	1.108	1.124	-1.4	144	0.00
34 I	Perylene-d12	1.000	1.000	0.0	133	0.00
35	Benzo(b)fluoranthene	1.206	1.218	-1.0	138	0.00
36	Benzo(k)fluoranthene	1.132	1.180	-4.2	138	0.00
37 S	Benzo(a)pyrene-d12	0.920	0.924	-0.4	135	0.00
38 C	Benzo(a)pyrene	1.161	1.138	2.0	131	0.00
39	Indeno(1,2,3-cd)pyrene	1.374	1.104	19.7	107	0.00
40	Dibenzo(a,h)anthracene	1.153	0.926	19.7	106	0.00
41	Benzo(g,h,i)perylene	1.152	0.902	21.7#	107	0.00

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(#) = Out of Range

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