

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062716\
 Data File : BM006083.D
 Acq On : 27 Jun 2016 23:09
 Operator : UM/SJ
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02040

Quant Time: Jun 28 02:14:38 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 28 02:11:41 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	70	0.00
2 S	1,4-Dioxane-d8	0.448	0.405	9.6	62	0.00
3 S	Phenol-d5	2.104	2.063	1.9	70	0.00
4 S	Bis-(2-Chloroethyl)ether-d8	1.170	1.103	5.7	65	0.00
5 S	2-Chlorophenol-d4	1.465	1.449	1.1	69	0.00
6 S	4-Methylphenol-d8	1.776	1.692	4.7	67	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	72	0.00
8 S	Nitrobenzene-d5	0.152	0.146	3.9	68	0.00
9 S	2-Nitrophenol-d4	0.169	0.156	7.7	65	0.00
10 S	2,4-Dichlorophenol-d3	0.325	0.337	-3.7	73	0.00
11	Naphthalene	0.999	1.024	-2.5	72	0.00
12 S	4-Chloroaniline-d4	0.353	0.443	-25.5#	72	0.00
13	2-Methylnaphthalene	0.804	0.828	-3.0	72	0.00
14	1-Methylnaphthalene	0.770	0.798	-3.6	72	0.00
15 I	Acenaphthene-d10	1.000	1.000	0.0	74	0.00
16 S	Dimethylphthalate-d6	1.712	1.528	10.7	66	0.00
17 S	Acenaphthylene-d8	1.986	2.050	-3.2	75	0.00
18	Acenaphthylene	2.094	2.166	-3.4	75	0.00
19 C	Acenaphthene	1.330	1.380	-3.8	76	0.00
20 S	4-Nitrophenol-d4	0.365	0.384	-5.2	73	-0.01
21 S	Fluorene-d10	1.462	1.552	-6.2	77	0.00
22	Fluorene	1.602	1.742	-8.7	78	0.00
23 I	Phenanthrene-d10	1.000	1.000	0.0	82	0.00
24 S	4,6-Dinitro-2-methylphenol-	0.111	0.076	31.5#	54	0.00
25	Phenanthrene	1.076	1.124	-4.5	84	0.00
26 S	Anthracene-d10	0.919	0.958	-4.2	84	0.00
27	Anthracene	1.111	1.173	-5.6	85	0.00
28 C	Fluoranthene	1.378	1.614	-17.1	93	0.00
29 I	Chrysene-d12	1.000	1.000	0.0	106	0.00
30 S	Pyrene-d10	0.876	0.823	6.1	97	0.00
31	Pyrene	1.145	1.079	5.8	96	0.00
32	Benzo(a)anthracene	1.188	1.196	-0.7	103	0.00
33	Chrysene	1.085	1.118	-3.0	106	0.00
34 I	Perylene-d12	1.000	1.000	0.0	98	0.00
35	Benzo(b)fluoranthene	1.162	1.203	-3.5	99	0.00
36	Benzo(k)fluoranthene	1.081	1.175	-8.7	102	0.00
37 S	Benzo(a)pyrene-d12	0.903	0.933	-3.3	98	0.00
38 C	Benzo(a)pyrene	1.114	1.142	-2.5	98	0.00
39	Indeno(1,2,3-cd)pyrene	1.328	1.245	6.2	88	0.00
40	Dibenzo(a,h)anthracene	1.104	1.043	5.5	88	0.00
41	Benzo(g,h,i)perylene	1.149	1.055	8.2	86	0.02

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(#) = Out of Range

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