

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060616\
 Data File : BM005674.D
 Acq On : 06 Jun 2016 17:44
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02046

Quant Time: Jun 07 04:26:50 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	109	0.00
2 S	1,4-Dioxane-d8	0.476	0.468	1.7	109	0.00
3 S	Phenol-d5	1.758	1.721	2.1	106	0.00
4 S	Bis-(2-Chloroethyl)ether-d8	1.070	1.072	-0.2	107	0.00
5 S	2-Chlorophenol-d4	1.451	1.434	1.2	109	0.00
6 S	4-Methylphenol-d8	1.391	1.311	5.8	103	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
8 S	Nitrobenzene-d5	0.154	0.155	-0.6	104	0.00
9 S	2-Nitrophenol-d4	0.164	0.162	1.2	101	0.00
10 S	2,4-Dichlorophenol-d3	0.292	0.285	2.4	101	0.00
11	Naphthalene	1.017	1.012	0.5	103	0.00
12 S	4-Chloroaniline-d4	0.347	0.375	-8.1	100	0.00
13	2-Methylnaphthalene	0.707	0.676	4.4	99	0.00
14	1-Methylnaphthalene	0.695	0.661	4.9	99	0.00
15 I	Acenaphthene-d10	1.000	1.000	0.0	95	0.00
16 S	Dimethylphthalate-d6	1.538	1.494	2.9	94	0.00
17 S	Acenaphthylene-d8	2.000	2.008	-0.4	96	0.00
18	Acenaphthylene	2.069	2.084	-0.7	95	0.00
19 C	Acenaphthene	1.343	1.352	-0.7	96	0.00
20 S	4-Nitrophenol-d4	0.220	0.207	5.9	92	0.00
21 S	Fluorene-d10	1.321	1.290	2.3	94	0.00
22	Fluorene	1.453	1.445	0.6	94	0.00
23 I	Phenanthrene-d10	1.000	1.000	0.0	93	0.00
24 S	4,6-Dinitro-2-methylphenol-	0.094	0.082	12.8	88	0.00
25	Phenanthrene	1.104	1.101	0.3	93	0.00
26 S	Anthracene-d10	0.950	0.958	-0.8	93	0.00
27	Anthracene	1.122	1.122	0.0	92	0.00
28 C	Fluoranthene	1.128	1.145	-1.5	93	0.00
29 I	Chrysene-d12	1.000	1.000	0.0	94	0.00
30 S	Pyrene-d10	1.005	0.973	3.2	92	0.00
31	Pyrene	1.262	1.228	2.7	92	0.00
32	Benzo(a)anthracene	1.173	1.162	0.9	93	0.00
33	Chrysene	1.108	1.100	0.7	94	0.00
34 I	Perylene-d12	1.000	1.000	0.0	96	-0.01
35	Benzo(b)fluoranthene	1.206	1.206	0.0	98	0.00
36	Benzo(k)fluoranthene	1.132	1.095	3.3	93	0.00
37 S	Benzo(a)pyrene-d12	0.920	0.905	1.6	96	0.00
38 C	Benzo(a)pyrene	1.161	1.145	1.4	95	0.00
39	Indeno(1,2,3-cd)pyrene	1.374	1.386	-0.9	97	0.00
40	Dibenzo(a,h)anthracene	1.153	1.175	-1.9	97	0.00
41	Benzo(g,h,i)perylene	1.152	1.151	0.1	98	0.00

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(#) = Out of Range

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