

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM073016\
 Data File : BM006764.D
 Acq On : 30 Jul 2016 09:25
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02060

Quant Time: Aug 01 04:24:28 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 30 01:37:26 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	96	0.00
2 S	1,4-Dioxane-d8	0.495	0.507	-2.4	98	0.00
3 S	Phenol-d5	1.702	1.735	-1.9	97	0.00
4 S	Bis-(2-Chloroethyl)ether-d8	1.054	1.110	-5.3	99	0.00
5 S	2-Chlorophenol-d4	1.360	1.376	-1.2	98	0.00
6 S	4-Methylphenol-d8	1.320	1.366	-3.5	98	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
8 S	Nitrobenzene-d5	0.151	0.152	-0.7	101	0.00
9 S	2-Nitrophenol-d4	0.161	0.165	-2.5	101	0.00
10 S	2,4-Dichlorophenol-d3	0.304	0.308	-1.3	100	0.00
11	Naphthalene	1.031	1.017	1.4	98	0.00
12 S	4-Chloroaniline-d4	0.335	0.369	-10.1	88	0.00
13	2-Methylnaphthalene	0.757	0.756	0.1	98	0.00
14	1-Methylnaphthalene	0.722	0.723	-0.1	99	0.00
15 I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
16 S	Dimethylphthalate-d6	1.568	1.577	-0.6	101	0.00
17 S	Acenaphthylene-d8	2.002	2.037	-1.7	102	0.00
18	Acenaphthylene	2.144	2.130	0.7	99	0.00
19 C	Acenaphthene	1.378	1.355	1.7	98	0.00
20 S	4-Nitrophenol-d4	0.272	0.264	2.9	99	0.00
21 S	Fluorene-d10	1.402	1.404	-0.1	101	0.00
22	Fluorene	1.552	1.565	-0.8	99	0.00
23 I	Phenanthrene-d10	1.000	1.000	0.0	100	0.00
24 S	4,6-Dinitro-2-methylphenol-	0.104	0.104	0.0	101	0.00
25	Phenanthrene	1.125	1.120	0.4	98	0.00
26 S	Anthracene-d10	0.950	0.959	-0.9	98	0.00
27	Anthracene	1.158	1.162	-0.3	98	0.00
28 C	Fluoranthene	1.321	1.342	-1.6	98	0.00
29 I	Chrysene-d12	1.000	1.000	0.0	99	0.00
30 S	Pyrene-d10	0.911	0.901	1.1	99	0.00
31	Pyrene	1.211	1.195	1.3	97	0.00
32	Benzo(a)anthracene	1.212	1.198	1.2	96	0.00
33	Chrysene	1.118	1.112	0.5	98	0.00
34 I	Perylene-d12	1.000	1.000	0.0	100	0.00
35	Benzo(b)fluoranthene	1.199	1.219	-1.7	101	0.00
36	Benzo(k)fluoranthene	1.159	1.119	3.5	94	0.00
37 S	Benzo(a)pyrene-d12	0.920	0.920	0.0	99	-0.01
38 C	Benzo(a)pyrene	1.167	1.163	0.3	98	0.00
39	Indeno(1,2,3-cd)pyrene	1.351	1.353	-0.1	98	-0.01
40	Dibenzo(a,h)anthracene	1.120	1.128	-0.7	98	-0.01
41	Benzo(g,h,i)perylene	1.158	1.156	0.2	99	-0.01

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(#) = Out of Range

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