

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM073016\
 Data File : BM006741.D
 Acq On : 29 Jul 2016 18:00
 Operator : UM/SJ
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02058

Quant Time: Jul 30 01:38:11 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	87	0.00
2 S	1,4-Dioxane-d8	0.495	0.511	-3.2	89	0.00
3 S	Phenol-d5	1.702	1.688	0.8	85	0.00
4 S	Bis-(2-Chloroethyl)ether-d8	1.054	1.096	-4.0	88	0.00
5 S	2-Chlorophenol-d4	1.360	1.362	-0.1	87	0.00
6 S	4-Methylphenol-d8	1.320	1.351	-2.3	87	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	88	0.00
8 S	Nitrobenzene-d5	0.151	0.149	1.3	89	0.00
9 S	2-Nitrophenol-d4	0.161	0.161	0.0	89	0.00
10 S	2,4-Dichlorophenol-d3	0.304	0.304	0.0	88	0.00
11	Naphthalene	1.031	1.008	2.2	87	0.00
12 S	4-Chloroaniline-d4	0.335	0.363	-8.4	78	0.00
13	2-Methylnaphthalene	0.757	0.751	0.8	88	0.00
14	1-Methylnaphthalene	0.722	0.727	-0.7	89	0.00
15 I	Acenaphthene-d10	1.000	1.000	0.0	90	0.00
16 S	Dimethylphthalate-d6	1.568	1.595	-1.7	91	0.00
17 S	Acenaphthylene-d8	2.002	2.031	-1.4	91	0.00
18	Acenaphthylene	2.144	2.131	0.6	88	0.00
19 C	Acenaphthene	1.378	1.367	0.8	88	0.00
20 S	4-Nitrophenol-d4	0.272	0.258	5.1	86	0.00
21 S	Fluorene-d10	1.402	1.412	-0.7	90	0.00
22	Fluorene	1.552	1.592	-2.6	89	0.00
23 I	Phenanthrene-d10	1.000	1.000	0.0	90	0.00
24 S	4,6-Dinitro-2-methylphenol-	0.104	0.101	2.9	88	0.00
25	Phenanthrene	1.125	1.116	0.8	88	0.00
26 S	Anthracene-d10	0.950	0.959	-0.9	89	0.00
27	Anthracene	1.158	1.155	0.3	88	0.00
28 C	Fluoranthene	1.321	1.335	-1.1	88	0.00
29 I	Chrysene-d12	1.000	1.000	0.0	88	0.00
30 S	Pyrene-d10	0.911	0.907	0.4	88	0.00
31	Pyrene	1.211	1.205	0.5	87	0.00
32	Benzo(a)anthracene	1.212	1.216	-0.3	87	0.00
33	Chrysene	1.118	1.120	-0.2	87	0.00
34 I	Perylene-d12	1.000	1.000	0.0	88	0.00
35	Benzo(b)fluoranthene	1.199	1.236	-3.1	90	0.00
36	Benzo(k)fluoranthene	1.159	1.111	4.1	82	0.00
37 S	Benzo(a)pyrene-d12	0.920	0.925	-0.5	87	0.00
38 C	Benzo(a)pyrene	1.167	1.175	-0.7	87	0.00
39	Indeno(1,2,3-cd)pyrene	1.351	1.357	-0.4	87	0.00
40	Dibenzo(a,h)anthracene	1.120	1.139	-1.7	87	0.00
41	Benzo(g,h,i)perylene	1.158	1.158	0.0	87	0.00

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(#) = Out of Range

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