

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM073016\
 Data File : BM006779.D
 Acq On : 30 Jul 2016 19:14
 Operator : UM/SJ
 Sample : SSTDC0020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02061

Quant Time: Aug 01 04:26:20 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	116	0.00
2 S	1,4-Dioxane-d8	0.495	0.502	-1.4	116	0.00
3 S	Phenol-d5	1.702	1.757	-3.2	118	0.00
4 S	Bis-(2-Chloroethyl)ether-d8	1.054	1.099	-4.3	117	0.00
5 S	2-Chlorophenol-d4	1.360	1.363	-0.2	116	0.00
6 S	4-Methylphenol-d8	1.320	1.344	-1.8	116	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	116	0.00
8 S	Nitrobenzene-d5	0.151	0.152	-0.7	118	0.00
9 S	2-Nitrophenol-d4	0.161	0.164	-1.9	119	0.00
10 S	2,4-Dichlorophenol-d3	0.304	0.307	-1.0	117	0.00
11	Naphthalene	1.031	1.019	1.2	116	0.00
12 S	4-Chloroaniline-d4	0.335	0.370	-10.4	105	0.00
13	2-Methylnaphthalene	0.757	0.745	1.6	114	0.00
14	1-Methylnaphthalene	0.722	0.706	2.2	114	0.00
15 I	Acenaphthene-d10	1.000	1.000	0.0	115	0.00
16 S	Dimethylphthalate-d6	1.568	1.555	0.8	113	0.00
17 S	Acenaphthylene-d8	2.002	2.022	-1.0	115	0.00
18	Acenaphthylene	2.144	2.109	1.6	112	0.00
19 C	Acenaphthene	1.378	1.346	2.3	111	0.00
20 S	4-Nitrophenol-d4	0.272	0.264	2.9	112	0.00
21 S	Fluorene-d10	1.402	1.390	0.9	114	0.00
22	Fluorene	1.552	1.556	-0.3	112	0.00
23 I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
24 S	4,6-Dinitro-2-methylphenol-	0.104	0.107	-2.9	115	0.00
25	Phenanthrene	1.125	1.117	0.7	109	0.00
26 S	Anthracene-d10	0.950	0.954	-0.4	109	0.00
27	Anthracene	1.158	1.158	0.0	109	0.00
28 C	Fluoranthene	1.321	1.312	0.7	107	0.00
29 I	Chrysene-d12	1.000	1.000	0.0	105	0.00
30 S	Pyrene-d10	0.911	0.930	-2.1	108	0.00
31	Pyrene	1.211	1.215	-0.3	105	0.00
32	Benzo(a)anthracene	1.212	1.219	-0.6	104	0.00
33	Chrysene	1.118	1.107	1.0	103	0.00
34 I	Perylene-d12	1.000	1.000	0.0	103	0.00
35	Benzo(b)fluoranthene	1.199	1.235	-3.0	105	0.00
36	Benzo(k)fluoranthene	1.159	1.125	2.9	97	0.00
37 S	Benzo(a)pyrene-d12	0.920	0.927	-0.8	102	-0.01
38 C	Benzo(a)pyrene	1.167	1.172	-0.4	101	0.00
39	Indeno(1,2,3-cd)pyrene	1.351	1.359	-0.6	101	0.00
40	Dibenzo(a,h)anthracene	1.120	1.132	-1.1	101	-0.01
41	Benzo(g,h,i)perylene	1.158	1.157	0.1	102	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM073016\
Data File : BM006779.D
Acq On : 30 Jul 2016 19:14
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02061

Quant Time: Aug 01 04:26:20 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)
----------	-------	------	-----------	------------

(#) = Out of Range

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