

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061516\
 Data File : BM005806.D
 Acq On : 15 Jun 2016 04:07
 Operator : UM/SJ
 Sample : SSTDC0020EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02041

Quant Time: Jun 15 05:53:29 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 15 01:42:36 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	103	0.00
2 S	1,4-Dioxane-d8	0.432	0.387	10.4	93	0.00
3 S	Phenol-d5	1.643	1.637	0.4	102	0.00
4 S	Bis-(2-Chloroethyl)ether-d8	0.978	0.988	-1.0	101	0.00
5 S	2-Chlorophenol-d4	1.369	1.395	-1.9	105	0.00
6 S	4-Methylphenol-d8	1.388	1.452	-4.6	105	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
8 S	Nitrobenzene-d5	0.142	0.148	-4.2	106	0.00
9 S	2-Nitrophenol-d4	0.157	0.168	-7.0	108	0.00
10 S	2,4-Dichlorophenol-d3	0.301	0.308	-2.3	106	0.00
11	Naphthalene	0.997	0.987	1.0	106	0.00
12 S	4-Chloroaniline-d4	0.331	0.375	-13.3	106	0.00
13	2-Methylnaphthalene	0.717	0.718	-0.1	107	0.00
14	1-Methylnaphthalene	0.724	0.707	2.3	107	0.00
15 I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
16 S	Dimethylphthalate-d6	1.588	1.614	-1.6	107	0.00
17 S	Acenaphthylene-d8	2.017	2.033	-0.8	107	0.00
18	Acenaphthylene	2.034	2.049	-0.7	109	0.00
19 C	Acenaphthene	1.332	1.333	-0.1	109	0.00
20 S	4-Nitrophenol-d4	0.183	0.167	8.7	116	0.00
21 S	Fluorene-d10	1.322	1.275	3.6	108	0.00
22	Fluorene	1.431	1.394	2.6	108	0.00
23 I	Phenanthrene-d10	1.000	1.000	0.0	105	0.00
24 S	4,6-Dinitro-2-methylphenol-	0.089	0.086	3.4	111	0.00
25	Phenanthrene	1.097	1.091	0.5	106	0.00
26 S	Anthracene-d10	0.928	0.922	0.6	106	0.00
27	Anthracene	1.086	1.093	-0.6	106	0.00
28 C	Fluoranthene	1.168	1.198	-2.6	106	0.00
29 I	Chrysene-d12	1.000	1.000	0.0	106	0.00
30 S	Pyrene-d10	0.973	0.935	3.9	107	0.00
31	Pyrene	1.228	1.178	4.1	106	0.00
32	Benzo(a)anthracene	1.157	1.154	0.3	106	0.00
33	Chrysene	1.134	1.096	3.4	104	0.00
34 I	Perylene-d12	1.000	1.000	0.0	103	0.00
35	Benzo(b)fluoranthene	1.195	1.208	-1.1	104	0.00
36	Benzo(k)fluoranthene	1.106	1.074	2.9	102	0.00
37 S	Benzo(a)pyrene-d12	0.924	0.920	0.4	104	0.00
38 C	Benzo(a)pyrene	1.141	1.133	0.7	104	0.00
39	Indeno(1,2,3-cd)pyrene	1.309	1.347	-2.9	106	0.00
40	Dibenzo(a,h)anthracene	1.095	1.120	-2.3	105	0.00
41	Benzo(g,h,i)perylene	1.134	1.145	-1.0	105	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061516\
Data File : BM005806.D
Acq On : 15 Jun 2016 04:07
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02041

Quant Time: Jun 15 05:53:29 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061516.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 15 01:42:36 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