

Data Path : Z:\HPCHEM1\BNA_M\Data\BM062916\
 Data File : BM006117.D
 Acq On : 29 Jun 2016 09:06
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02050

Quant Time: Jun 29 09:37:21 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062816.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 29 00:57:09 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	134	0.00
2 S	1,4-Dioxane-d8	0.431	0.347	19.5	135	0.00
3 S	Phenol-d5	1.634	1.378	15.7	137	0.00
4 S	Bis-(2-Chloroethyl)ether-d8	0.963	0.840	12.8	136	0.00
5 S	2-Chlorophenol-d4	1.254	1.029	17.9	135	0.00
6 S	4-Methylphenol-d8	1.315	1.119	14.9	137	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	136	0.00
8 S	Nitrobenzene-d5	0.135	0.109	19.3	138	0.00
9 S	2-Nitrophenol-d4	0.154	0.121	21.4#	132	0.00
10 S	2,4-Dichlorophenol-d3	0.281	0.225	19.9	135	0.00
11	Naphthalene	0.876	0.714	18.5	138	0.00
12 S	4-Chloroaniline-d4	0.222	0.182	18.0	123	0.00
13	2-Methylnaphthalene	0.665	0.549	17.4	139	0.00
14	1-Methylnaphthalene	0.632	0.516	18.4	138	0.00
15 I	Acenaphthene-d10	1.000	1.000	0.0	136	0.00
16 S	Dimethylphthalate-d6	1.434	1.146	20.1#	136	0.00
17 S	Acenaphthylene-d8	1.740	1.429	17.9	138	0.00
18	Acenaphthylene	1.822	1.501	17.6	136	0.00
19 C	Acenaphthene	1.162	0.944	18.8	134	0.00
20 S	4-Nitrophenol-d4	0.272	0.220	19.1	129	0.00
21 S	Fluorene-d10	1.223	1.001	18.2	138	0.00
22	Fluorene	1.323	1.089	17.7	135	0.00
23 I	Phenanthrene-d10	1.000	1.000	0.0	135	0.00
24 S	4,6-Dinitro-2-methylphenol-	0.071	0.075	-5.6	213#	0.00
25	Phenanthrene	0.946	0.769	18.7	135	0.00
26 S	Anthracene-d10	0.802	0.659	17.8	134	0.00
27	Anthracene	0.973	0.798	18.0	134	0.00
28 C	Fluoranthene	1.159	0.940	18.9	131	0.00
29 I	Chrysene-d12	1.000	1.000	0.0	118	0.00
30 S	Pyrene-d10	0.805	0.694	13.8	129	0.00
31	Pyrene	1.055	0.912	13.6	129	0.00
32	Benzo(a)anthracene	1.040	0.855	17.8	119	0.00
33	Chrysene	0.957	0.794	17.0	119	0.00
34 I	Perylene-d12	1.000	1.000	0.0	98	0.00
35	Benzo(b)fluoranthene	1.038	0.885	14.7	106	0.00
36	Benzo(k)fluoranthene	0.974	0.830	14.8	103	0.00
37 S	Benzo(a)pyrene-d12	0.796	0.648	18.6	100	0.00
38 C	Benzo(a)pyrene	0.981	0.804	18.0	99	0.00
39	Indeno(1,2,3-cd)pyrene	1.019	0.795	22.0#	90	0.00
40	Dibenzo(a,h)anthracene	0.843	0.662	21.5#	90	0.00
41	Benzo(g,h,i)perylene	0.853	0.671	21.3#	91	0.00

Data Path : Z:\HPCHEM1\BNA_M\Data\BM062916\
Data File : BM006117.D
Acq On : 29 Jun 2016 09:06
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02050

Quant Time: Jun 29 09:37:21 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062816.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 29 00:57:09 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