

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080916\
 Data File : BM006995.D
 Acq On : 09 Aug 2016 23:08
 Operator : UM/SJ
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02040

Quant Time: Aug 10 18:45:52 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM080916.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 10 01:30:54 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	70	0.00
2 S	1,4-Dioxane-d8	0.439	0.453	-3.2	69	0.00
3 S	Phenol-d5	1.780	1.840	-3.4	70	0.00
4 S	Bis-(2-Chloroethyl)ether-d8	0.995	1.038	-4.3	69	0.00
5 S	2-Chlorophenol-d4	1.350	1.393	-3.2	71	0.00
6 S	4-Methylphenol-d8	1.507	1.560	-3.5	70	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	71	0.00
8 S	Nitrobenzene-d5	0.153	0.158	-3.3	71	0.00
9 S	2-Nitrophenol-d4	0.175	0.180	-2.9	71	0.00
10 S	2,4-Dichlorophenol-d3	0.342	0.352	-2.9	71	0.00
11	Naphthalene	1.003	1.035	-3.2	71	0.00
12 S	4-Chloroaniline-d4	0.399	0.448	-12.3	72	0.00
13	2-Methylnaphthalene	0.794	0.838	-5.5	72	0.00
14	1-Methylnaphthalene	0.762	0.797	-4.6	72	0.00
15 I	Acenaphthene-d10	1.000	1.000	0.0	73	0.00
16 S	Dimethylphthalate-d6	1.709	1.799	-5.3	74	0.00
17 S	Acenaphthylene-d8	2.002	2.095	-4.6	73	0.00
18	Acenaphthylene	2.069	2.175	-5.1	74	0.00
19 C	Acenaphthene	1.338	1.403	-4.9	73	0.00
20 S	4-Nitrophenol-d4	0.307	0.342	-11.4	79	0.00
21 S	Fluorene-d10	1.453	1.542	-6.1	75	0.00
22	Fluorene	1.588	1.696	-6.8	74	0.00
23 I	Phenanthrene-d10	1.000	1.000	0.0	76	0.00
24 S	4,6-Dinitro-2-methylphenol-	0.108	0.111	-2.8	80	0.00
25	Phenanthrene	1.072	1.122	-4.7	76	0.00
26 S	Anthracene-d10	0.932	0.973	-4.4	76	0.00
27	Anthracene	1.107	1.172	-5.9	78	0.00
28 C	Fluoranthene	1.350	1.468	-8.7	80	0.00
29 I	Chrysene-d12	1.000	1.000	0.0	85	0.00
30 S	Pyrene-d10	0.909	0.890	2.1	80	0.00
31	Pyrene	1.160	1.146	1.2	80	0.00
32	Benzo(a)anthracene	1.178	1.229	-4.3	87	0.00
33	Chrysene	1.069	1.130	-5.7	87	0.00
34 I	Perylene-d12	1.000	1.000	0.0	101	0.00
35	Benzo(b)fluoranthene	1.242	1.246	-0.3	98	0.00
36	Benzo(k)fluoranthene	1.174	1.172	0.2	96	0.00
37 S	Benzo(a)pyrene-d12	0.919	0.950	-3.4	101	0.00
38 C	Benzo(a)pyrene	1.143	1.185	-3.7	101	0.00
39	Indeno(1,2,3-cd)pyrene	1.133	1.316	-16.2	116	0.00
40	Dibenzo(a,h)anthracene	0.938	1.097	-17.0	116	0.00
41	Benzo(g,h,i)perylene	0.932	1.086	-16.5	117	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080916\
Data File : BM006995.D
Acq On : 09 Aug 2016 23:08
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02040

Quant Time: Aug 10 18:45:52 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM080916.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 10 01:30:54 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