

Data Path : Z:\HPCHEM1\BNA_M\Data\BM061216\
 Data File : BM005741.D
 Acq On : 12 Jun 2016 10:01
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02055

Quant Time: Jun 13 01:24:20 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jun 11 05:37:26 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	120	0.00
2 S	1,4-Dioxane-d8	0.461	0.386	16.3	104	0.00
3 S	Phenol-d5	1.726	1.657	4.0	115	0.00
4 S	Bis-(2-Chloroethyl)ether-d8	1.037	1.003	3.3	113	0.00
5 S	2-Chlorophenol-d4	1.395	1.413	-1.3	120	0.00
6 S	4-Methylphenol-d8	1.458	1.459	-0.1	119	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	121	0.00
8 S	Nitrobenzene-d5	0.144	0.143	0.7	117	0.00
9 S	2-Nitrophenol-d4	0.163	0.159	2.5	112	0.00
10 S	2,4-Dichlorophenol-d3	0.296	0.300	-1.4	121	0.00
11	Naphthalene	1.002	0.977	2.5	119	0.00
12 S	4-Chloroaniline-d4	0.357	0.417	-16.8	118	0.00
13	2-Methylnaphthalene	0.754	0.732	2.9	117	0.00
14	1-Methylnaphthalene	0.745	0.729	2.1	118	0.00
15 I	Acenaphthene-d10	1.000	1.000	0.0	116	0.00
16 S	Dimethylphthalate-d6	1.650	1.557	5.6	108	0.00
17 S	Acenaphthylene-d8	2.018	2.011	0.3	113	0.00
18	Acenaphthylene	2.066	2.032	1.6	113	0.00
19 C	Acenaphthene	1.344	1.330	1.0	114	0.00
20 S	4-Nitrophenol-d4	0.241	0.206	14.5	97	0.00
21 S	Fluorene-d10	1.397	1.365	2.3	111	0.00
22	Fluorene	1.534	1.511	1.5	111	0.00
23 I	Phenanthrene-d10	1.000	1.000	0.0	108	0.00
24 S	4,6-Dinitro-2-methylphenol-	0.103	0.063	38.8#	68	0.00
25	Phenanthrene	1.095	1.090	0.5	107	0.00
26 S	Anthracene-d10	0.942	0.933	1.0	106	0.00
27	Anthracene	1.111	1.103	0.7	105	0.00
28 C	Fluoranthene	1.286	1.161	9.7	96	0.00
29 I	Chrysene-d12	1.000	1.000	0.0	75	0.00
30 S	Pyrene-d10	0.961	1.142	-18.8	92	0.00
31	Pyrene	1.217	1.443	-18.6	91	0.00
32	Benzo(a)anthracene	1.155	1.149	0.5	75	0.00
33	Chrysene	1.139	1.113	2.3	73	0.00
34 I	Perylene-d12	1.000	1.000	0.0	73	0.00
35	Benzo(b)fluoranthene	1.188	1.150	3.2	73	0.00
36	Benzo(k)fluoranthene	1.201	1.148	4.4	69	0.00
37 S	Benzo(a)pyrene-d12	0.912	0.910	0.2	73	0.00
38 C	Benzo(a)pyrene	1.128	1.119	0.8	73	0.00
39	Indeno(1,2,3-cd)pyrene	1.101	1.283	-16.5	86	0.00
40	Dibenzo(a,h)anthracene	0.921	1.081	-17.4	87	0.00
41	Benzo(g,h,i)perylene	0.921	1.091	-18.5	90	0.00

Data Path : Z:\HPCHEM1\BNA_M\Data\BM061216\
Data File : BM005741.D
Acq On : 12 Jun 2016 10:01
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02055

Quant Time: Jun 13 01:24:20 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061016.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jun 11 05:37:26 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