

Data Path : Z:\HPCHEM1\BNA_M\Data\BM110115\
 Data File : BM003019.D
 Acq On : 01 Nov 2015 08:21
 Operator : UM/IZ
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02061

Quant Time: Nov 02 01:01:47 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM103015.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 30 18:02:20 2015
 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	100	0.00
2	1,4-Dioxane	0.463	0.424	8.4	99	0.00
3 S	1,4-Dioxane-d8	0.424	0.389	8.3	97	0.00
4	Benzaldehyde	0.812	0.883	-8.7	98	0.00
5 S	Phenol-d5	1.581	1.536	2.8	98	0.00
6	Phenol	1.627	1.609	1.1	98	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	0.924	0.896	3.0	97	0.00
8	Bis(2-Chloroethyl)ether	1.298	1.272	2.0	98	0.00
9 S	2-Chlorophenol-d4	1.340	1.275	4.9	98	0.00
10	2-Chlorophenol	1.384	1.328	4.0	99	0.00
11	2-Methylphenol	1.269	1.245	1.9	98	0.00
12	2,2'-oxybis(1-Chloropropane	1.597	1.561	2.3	96	0.00
13 S	4-Methylphenol-d8	1.275	1.260	1.2	98	0.00
14	Acetophenone	1.936	1.982	-2.4	97	0.00
15 P	N-Nitroso-di-n-propylamine	0.959	0.951	0.8	96	0.00
16	4-Methylphenol	1.371	1.371	0.0	98	0.00
17	Hexachloroethane	0.526	0.498	5.3	99	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
19 S	Nitrobenzene-d5	0.122	0.120	1.6	100	0.00
20	Nitrobenzene	0.296	0.292	1.4	100	0.00
21	Isophorone	0.599	0.584	2.5	96	0.00
22 S	2-Nitrophenol-d4	0.124	0.124	0.0	104	0.00
23 C	2-Nitrophenol	0.144	0.146	-1.4	102	0.00
24	2,4-Dimethylphenol	0.335	0.325	3.0	98	0.00
25	Bis(2-Chloroethoxy)methane	0.387	0.375	3.1	98	0.00
26 S	2,4-Dichlorophenol-d3	0.277	0.270	2.5	98	0.00
27 C	2,4-Dichlorophenol	0.284	0.278	2.1	98	0.00
28	Naphthalene	1.004	0.962	4.2	99	0.00
29 S	4-Chloroaniline-d4	0.359	0.380	-5.8	97	0.00
30	4-Chloroaniline	0.376	0.403	-7.2	97	0.00
31 C	Hexachlorobutadiene	0.176	0.169	4.0	101	0.00
32	Caprolactam	0.080	0.076	5.0	90	0.00
33 C	4-Chloro-3-methylphenol	0.301	0.293	2.7	95	0.00
34	2-Methylnaphthalene	0.723	0.695	3.9	97	0.00
35	1-Methylnaphthalene	0.709	0.688	3.0	98	0.00
36 I	Acenaphthene-d10	1.000	1.000	0.0	96	0.00
37	1,2,4,5-Tetrachlorobenzene	0.640	0.613	4.2	99	0.00
38	Hexachlorocyclopentadiene	0.377	0.345	8.5	99	0.00
39 C	2,4,6-Trichlorophenol	0.391	0.380	2.8	97	0.00
40	2,4,5-Trichlorophenol	0.420	0.414	1.4	98	0.00
41	1,1'-Biphenyl	1.685	1.614	4.2	97	0.00
42	2-Chloronaphthalene	1.254	1.207	3.7	97	0.00
43	2-Nitroaniline	0.252	0.257	-2.0	97	0.00
44 S	Dimethylphthalate-d6	1.505	1.450	3.7	93	0.00

Data Path : Z:\HPCHEM1\BNA_M\Data\BM110115\
 Data File : BM003019.D
 Acq On : 01 Nov 2015 08:21
 Operator : UM/IZ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02061

Quant Time: Nov 02 01:01:47 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM103015.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 30 18:02:20 2015
 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)
45	Dimethylphthalate	1.533	1.472	4.0	93	0.00
46	2,6-Dinitrotoluene	0.251	0.257	-2.4	96	0.00
47 S	Acenaphthylene-d8	1.907	1.846	3.2	95	0.00
48	Acenaphthylene	2.036	1.985	2.5	97	0.00
49	3-Nitroaniline	0.265	0.268	-1.1	96	0.00
50 C	Acenaphthene	1.371	1.309	4.5	97	0.00
51	2,4-Dinitrophenol	0.139	0.128	7.9	100	0.00
52 S	4-Nitrophenol-d4	0.254	0.236	7.1	92	0.00
53	4-Nitrophenol	0.193	0.183	5.2	93	0.00
54	Dibenzofuran	1.917	1.844	3.8	96	0.00
55	2,4-Dinitrotoluene	0.353	0.363	-2.8	94	0.00
56	2,3,4,6-Tetrachlorophenol	0.347	0.340	2.0	94	0.00
57	Diethylphthalate	1.512	1.457	3.6	91	0.00
58 S	Fluorene-d10	1.301	1.254	3.6	95	0.00
59	Fluorene	1.490	1.439	3.4	94	0.00
60	4-Chlorophenyl-phenylether	0.700	0.681	2.7	95	0.00
61	4-Nitroaniline	0.284	0.258	9.2	91	0.00
62 I	Phenanthrene-d10	1.000	1.000	0.0	91	0.00
63 S	4,6-Dinitro-2-methylphenol-	0.093	0.087	6.5	96	0.00
64	4,6-Dinitro-2-methylphenol	0.100	0.095	5.0	94	0.00
65	N-Nitrosodiphenylamine	0.601	0.581	3.3	92	0.00
66	4-Bromophenyl-phenylether	0.195	0.190	2.6	93	0.00
67	Hexachlorobenzene	0.236	0.225	4.7	93	0.00
68	Atrazine	0.207	0.200	3.4	88	0.00
69 C	Pentachlorophenol	0.132	0.124	6.1	90	0.00
70	Phenanthrene	1.133	1.071	5.5	91	0.00
71 S	Anthracene-d10	0.887	0.848	4.4	90	0.00
72	Anthracene	1.106	1.065	3.7	91	0.00
73	1,2,3,4-Tetrachlorobenzene	0.303	0.291	4.0	98	0.00
74	Pentachlorobenzene	0.287	0.276	3.8	96	0.00
75	Carbazole	0.945	0.920	2.6	88	0.00
76	Di-n-butylphthalate	1.088	1.055	3.0	85	0.00
77 C	Fluoranthene	1.145	1.111	3.0	86	0.00
78 I	Chrysene-d12	1.000	1.000	0.0	90	0.00
79 S	Pyrene-d10	0.985	0.953	3.2	88	0.00
80	Pyrene	1.308	1.275	2.5	88	0.00
81	Butylbenzylphthalate	0.386	0.396	-2.6	90	0.00
82	3,3'-Dichlorobenzidine	0.280	0.276	1.4	93	0.00
83	Benzo(a)anthracene	1.098	1.058	3.6	88	0.00
84	Bis(2-ethylhexyl)phthalate	0.604	0.611	-1.2	87	0.00
85	Chrysene	1.087	1.032	5.1	89	0.00
86 I	Perylene-d12	1.000	1.000	0.0	95	0.00
87	Di-n-octyl phthalate	1.030	0.940	8.7	92	0.00

Data Path : Z:\HPCHEM1\BNA_M\Data\BM110115\
Data File : BM003019.D
Acq On : 01 Nov 2015 08:21
Operator : UM/IZ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02061

Quant Time: Nov 02 01:01:47 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM103015.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 30 18:02:20 2015
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)
88	Benzo(b)fluoranthene	1.152	1.125	2.3	96	0.00
89	Benzo(k)fluoranthene	1.075	1.047	2.6	92	0.00
90 S	Benzo(a)pyrene-d12	0.950	0.928	2.3	95	0.00
91 C	Benzo(a)pyrene	1.091	1.057	3.1	95	0.00
92	Indeno(1,2,3-cd)pyrene	1.200	1.170	2.5	99	0.00
93	Dibenzo(a,h)anthracene	0.945	0.937	0.8	99	0.00
94	Benzo(g,h,i)perylene	1.079	1.041	3.5	101	0.00

(#) = Out of Range

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