

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM070516\
 Data File : BM006263.D
 Acq On : 05 Jul 2016 18:20
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02038

Quant Time: Jul 05 18:51:21 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM070216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 05 17:52:15 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	159#	0.00
2	1,4-Dioxane	0.512	0.518	-1.2	166#	0.00
3 S	1,4-Dioxane-d8	0.468	0.465	0.6	156#	0.00
4	Benzaldehyde	1.083	1.281	-18.3	157#	0.00
5 S	Phenol-d5	1.873	1.903	-1.6	158#	0.00
6	Phenol	1.946	2.005	-3.0	160#	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.105	1.168	-5.7	161#	0.00
8	Bis(2-Chloroethyl)ether	1.438	1.553	-8.0	163#	0.00
9 S	2-Chlorophenol-d4	1.433	1.443	-0.7	160#	0.00
10	2-Chlorophenol	1.484	1.495	-0.7	160#	0.00
11	2-Methylphenol	1.491	1.510	-1.3	156#	0.00
12	2,2'-oxybis(1-Chloropropane	2.147	2.234	-4.1	160#	0.00
13 S	4-Methylphenol-d8	1.516	1.524	-0.5	155#	0.00
14	Acetophenone	2.310	2.383	-3.2	153#	0.00
15 P	N-Nitroso-di-n-propylamine	1.319	1.329	-0.8	156#	0.00
16	4-Methylphenol	1.618	1.655	-2.3	156#	0.00
17	Hexachloroethane	0.608	0.605	0.5	157#	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	155#	0.00
19 S	Nitrobenzene-d5	0.159	0.161	-1.3	156#	0.00
20	Nitrobenzene	0.415	0.415	0.0	154#	0.00
21	Isophorone	0.790	0.785	0.6	152#	0.00
22 S	2-Nitrophenol-d4	0.181	0.173	4.4	146	0.00
23 C	2-Nitrophenol	0.195	0.193	1.0	153#	0.00
24	2,4-Dimethylphenol	0.411	0.404	1.7	150#	0.00
25	Bis(2-Chloroethoxy)methane	0.472	0.481	-1.9	157#	0.00
26 S	2,4-Dichlorophenol-d3	0.327	0.320	2.1	153#	0.00
27 C	2,4-Dichlorophenol	0.324	0.319	1.5	152#	0.00
28	Naphthalene	1.029	1.039	-1.0	157#	0.00
29 S	4-Chloroaniline-d4	0.340	0.414	-21.8#	179#	0.00
30	4-Chloroaniline	0.353	0.439	-24.4#	183#	0.00
31 C	Hexachlorobutadiene	0.202	0.193	4.5	148	0.00
32	Caprolactam	0.128	0.111	13.3	133	0.00
33 C	4-Chloro-3-methylphenol	0.395	0.375	5.1	147	0.00
34	2-Methylnaphthalene	0.781	0.772	1.2	151#	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	144	0.00
36	1,2,4,5-Tetrachlorobenzene	0.665	0.691	-3.9	147	0.00
37	Hexachlorocyclopentadiene	0.378	0.320	15.3	120	0.00
38 C	2,4,6-Trichlorophenol	0.470	0.462	1.7	140	0.00
39	2,4,5-Trichlorophenol	0.491	0.483	1.6	136	0.00
40	1,1'-Biphenyl	1.649	1.734	-5.2	149	0.00
41	2-Chloronaphthalene	1.285	1.339	-4.2	149	0.00
42	2-Nitroaniline	0.491	0.476	3.1	136	0.00
43 S	Dimethylphthalate-d6	1.669	1.631	2.3	140	0.00
44	Dimethylphthalate	1.678	1.643	2.1	140	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM070516\
 Data File : BM006263.D
 Acq On : 05 Jul 2016 18:20
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02038

Quant Time: Jul 05 18:51:21 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM070216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 05 17:52:15 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)
45	2,6-Dinitrotoluene	0.364	0.347	4.7	136	0.00
46 S	Acenaphthylene-d8	2.067	2.095	-1.4	144	0.00
47	Acenaphthylene	2.169	2.202	-1.5	144	0.00
48	3-Nitroaniline	0.332	0.377	-13.6	153#	0.00
49 C	Acenaphthene	1.378	1.387	-0.7	144	0.00
50	2,4-Dinitrophenol	0.202	0.145	28.2#	110	0.00
51 S	4-Nitrophenol-d4	0.316	0.289	8.5	126	0.00
52	4-Nitrophenol	0.318	0.264	17.0	115	0.00
53	Dibenzofuran	1.967	1.969	-0.1	141	0.00
54	2,4-Dinitrotoluene	0.514	0.487	5.3	132	0.00
55	2,3,4,6-Tetrachlorophenol	0.442	0.409	7.5	130	0.00
56	Diethylphthalate	1.751	1.692	3.4	137	0.00
57 S	Fluorene-d10	1.439	1.402	2.6	139	0.00
58	Fluorene	1.577	1.529	3.0	135	0.00
59	4-Chlorophenyl-phenylether	0.774	0.748	3.4	136	0.00
60	4-Nitroaniline	0.392	0.400	-2.0	131	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	127	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.117	0.109	6.8	114	0.00
63	4,6-Dinitro-2-methylphenol	0.126	0.119	5.6	115	0.00
64	N-Nitrosodiphenylamine	0.604	0.653	-8.1	136	0.00
65	4-Bromophenyl-phenylether	0.224	0.232	-3.6	130	0.00
66	Hexachlorobenzene	0.247	0.250	-1.2	126	0.00
67	Atrazine	0.223	0.234	-4.9	123	0.00
68 C	Pentachlorophenol	0.131	0.120	8.4	117	0.00
69	Phenanthrene	1.116	1.136	-1.8	128	0.00
70 S	Anthracene-d10	0.954	0.963	-0.9	127	0.00
71	Anthracene	1.152	1.165	-1.1	126	0.00
72	Carbazole	0.966	1.015	-5.1	122	0.00
73	Di-n-butylphthalate	1.327	1.305	1.7	121	0.00
74 C	Fluoranthene	1.242	1.209	2.7	110	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	91	0.00
76 S	Pyrene-d10	0.954	1.140	-19.5	109	0.00
77	Pyrene	1.254	1.488	-18.7	107	0.00
78	Butylbenzylphthalate	0.573	0.641	-11.9	102	0.00
79	3,3'-Dichlorobenzidine	0.370	0.417	-12.7	90	0.00
80	Benzo(a)anthracene	1.227	1.261	-2.8	92	0.00
81	Bis(2-ethylhexyl)phthalate	0.769	0.851	-10.7	98	0.00
82	Chrysene	1.120	1.157	-3.3	93	0.00
83 I	Perylene-d12	1.000	1.000	0.0	81	0.00
84	Di-n-octyl phthalate	1.188	1.492	-25.6#	90	0.00
85	Benzo(b)fluoranthene	1.184	1.273	-7.5	87	0.00
86	Benzo(k)fluoranthene	1.102	1.163	-5.5	82	0.00
87 S	Benzo(a)pyrene-d12	0.929	0.955	-2.8	83	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM070516\
Data File : BM006263.D
Acq On : 05 Jul 2016 18:20
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02038

Quant Time: Jul 05 18:51:21 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM070216.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jul 05 17:52:15 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)
88 C	Benzo(a)pyrene	1.146	1.200	-4.7	83	0.00
89	Indeno(1,2,3-cd)pyrene	1.311	1.348	-2.8	83	0.00
90	Dibenzo(a,h)anthracene	1.081	1.124	-4.0	82	0.00
91	Benzo(g,h,i)perylene	1.131	1.158	-2.4	83	0.00

(#) = Out of Range

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