

Data Path : Z:\HPCHEM1\BNA M\Data\BM021918\
 Data File : BM014297.D
 Acq On : 19 Feb 2018 16:30
 Operator : SJ/JU
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02045

Quant Time: Feb 20 01:11:25 2018
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Feb 17 06:25:12 2018
 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	74	0.00
2	1,4-Dioxane	0.519	0.588	-13.3	88	0.00
3 S	1,4-Dioxane-d8	0.467	0.564	-20.8#	96	0.00
4	Benzaldehyde	0.689	0.794	-15.2	74	0.00
5 S	Phenol-d5	1.690	1.620	4.1	77	0.00
6	Phenol	1.774	1.710	3.6	78	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.023	1.018	0.5	76	0.00
8	Bis(2-Chloroethyl)ether	1.287	1.263	1.9	75	0.00
9 S	2-Chlorophenol-d4	1.296	1.349	-4.1	78	0.00
10	2-Chlorophenol	1.323	1.311	0.9	79	0.00
11	2-Methylphenol	1.357	1.314	3.2	79	0.00
12	2,2'-oxybis(1-Chloropropane	1.299	1.315	-1.2	81	0.00
13 S	4-Methylphenol-d8	1.479	1.472	0.5	80	0.00
14	Acetophenone	2.552	2.530	0.9	83	0.00
15 P	N-Nitroso-di-n-propylamine	1.279	1.366	-6.8	85	0.00
16	4-Methylphenol	1.478	1.422	3.8	78	0.00
17	Hexachloroethane	0.689	0.742	-7.7	83	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	80	0.00
19 S	Nitrobenzene-d5	0.148	0.137	7.4	75	0.00
20	Nitrobenzene	0.437	0.417	4.6	79	0.00
21	Isophorone	0.748	0.749	-0.1	82	0.00
22 S	2-Nitrophenol-d4	0.153	0.149	2.6	75	0.00
23 C	2-Nitrophenol	0.166	0.172	-3.6	83	0.00
24	2,4-Dimethylphenol	0.406	0.419	-3.2	86	0.00
25	Bis(2-Chloroethoxy)methane	0.394	0.387	1.8	79	0.00
26 S	2,4-Dichlorophenol-d3	0.317	0.299	5.7	75	0.00
27 C	2,4-Dichlorophenol	0.301	0.294	2.3	83	0.00
28	Naphthalene	1.027	1.020	0.7	82	0.00
29 S	4-Chloroaniline-d4	0.312	0.355	-13.8	86	0.00
30	4-Chloroaniline	0.318	0.361	-13.5	86	0.00
31 C	Hexachlorobutadiene	0.239	0.257	-7.5	90	0.00
32	Caprolactam	0.096	0.095	1.0	81	0.00
33 C	4-Chloro-3-methylphenol	0.363	0.374	-3.0	84	0.00
34	2-Methylnaphthalene	0.782	0.796	-1.8	83	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	85	0.00
36	1,2,4,5-Tetrachlorobenzene	0.642	0.628	2.2	86	0.00
37	Hexachlorocyclopentadiene	0.349	0.301	13.8	86	0.00
38 C	2,4,6-Trichlorophenol	0.401	0.391	2.5	86	0.00
39	2,4,5-Trichlorophenol	0.409	0.389	4.9	78	0.00
40	1,1'-Biphenyl	1.541	1.542	-0.1	86	0.00
41	2-Chloronaphthalene	1.231	1.199	2.6	84	0.00
42	2-Nitroaniline	0.416	0.417	-0.2	85	0.00
43 S	Dimethylphthalate-d6	1.652	1.621	1.9	85	0.00
44	Dimethylphthalate	1.629	1.602	1.7	87	0.00

Data Path : Z:\HPCHEM1\BNA M\Data\BM021918\
 Data File : BM014297.D
 Acq On : 19 Feb 2018 16:30
 Operator : SJ/JU
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02045

Quant Time: Feb 20 01:11:25 2018
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Feb 17 06:25:12 2018
 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	0.309	0.328	-6.1	91	0.00
46 S	Acenaphthylene-d8	2.037	2.042	-0.2	87	0.00
47	Acenaphthylene	1.925	1.907	0.9	85	0.00
48	3-Nitroaniline	0.262	0.240	8.4	89	0.00
49 C	Acenaphthene	1.355	1.397	-3.1	91	0.00
50	2,4-Dinitrophenol	0.166	0.129	22.3#	91	0.00
51 S	4-Nitrophenol-d4	0.228	0.184	19.3	86	0.00
52	4-Nitrophenol	0.299	0.258	13.7	85	0.00
53	Dibenzofuran	2.032	2.074	-2.1	89	0.00
54	2,4-Dinitrotoluene	0.496	0.518	-4.4	89	0.00
55	2,3,4,6-Tetrachlorophenol	0.414	0.421	-1.7	88	0.00
56	Diethylphthalate	1.815	1.907	-5.1	91	0.00
57 S	Fluorene-d10	1.480	1.528	-3.2	90	0.00
58	Fluorene	1.755	1.781	-1.5	89	0.00
59	4-Chlorophenyl-phenylether	0.918	0.946	-3.1	91	0.00
60	4-Nitroaniline	0.294	0.230	21.8#	82	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	89	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.120	0.107	10.8	92	0.00
63	4,6-Dinitro-2-methylphenol	0.122	0.110	9.8	91	0.00
64	N-Nitrosodiphenylamine	0.575	0.565	1.7	90	0.00
65	4-Bromophenyl-phenylether	0.223	0.223	0.0	94	0.00
66	Hexachlorobenzene	0.235	0.235	0.0	95	0.00
67	Atrazine	0.227	0.228	-0.4	95	0.00
68 C	Pentachlorophenol	0.123	0.109	11.4	95	0.00
69	Phenanthrene	1.152	1.148	0.3	93	0.00
70 S	Anthracene-d10	0.965	0.981	-1.7	93	0.00
71	Anthracene	1.159	1.161	-0.2	91	0.00
72	Carbazole	0.974	0.946	2.9	93	0.00
73	Di-n-butylphthalate	1.257	1.307	-4.0	93	0.00
74 C	Fluoranthene	1.377	1.405	-2.0	95	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	95	0.00
76 S	Pyrene-d10	1.032	0.998	3.3	95	0.00
77	Pyrene	1.329	1.280	3.7	95	0.00
78	Butylbenzylphthalate	0.550	0.529	3.8	93	0.00
79	3,3'-Dichlorobenzidine	0.352	0.289	17.9	93	0.00
80	Benzo(a)anthracene	1.248	1.214	2.7	96	0.00
81	Bis(2-ethylhexyl)phthalate	0.787	0.796	-1.1	98	0.00
82	Chrysene	1.164	1.131	2.8	94	0.00
83 I	Perylene-d12	1.000	1.000	0.0	101	0.00
84	Di-n-octyl phthalate	1.716	1.519	11.5	100	0.00
85	Benzo(b)fluoranthene	1.339	1.287	3.9	98	0.00
86	Benzo(k)fluoranthene	1.273	1.307	-2.7	107	0.00
87 S	Benzo(a)pyrene-d12	0.973	0.974	-0.1	103	0.00

Data Path : Z:\HPCHEM1\BNA M\Data\BM021918\
Data File : BM014297.D
Acq On : 19 Feb 2018 16:30
Operator : SJ/JU
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02045

Quant Time: Feb 20 01:11:25 2018
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Feb 17 06:25:12 2018
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	1.197	1.191	0.5	105	0.00
89	Indeno(1,2,3-cd)pyrene	1.161	1.126	3.0	103	0.00
90	Dibenzo(a,h)anthracene	0.965	0.930	3.6	103	0.00
91	Benzo(a,h,i)perylene	0.941	0.869	7.7	100	0.01

(#) = Out of Range

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