

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN061918\
 Data File : BN001876.D
 Acq On : 19 Jun 2018 14:51
 Operator : JU/SJ
 Sample : SSTDICV020
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV87

Quant Time: Jun 19 15:27:51 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 19 14:50:02 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)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	107	0.00
2	1,4-Dioxane	0.474	0.457	3.6	105	0.00
3 S	1,4-Dioxane-d8	0.460	0.454	1.3	105	0.00
4	Benzaldehyde	1.105	1.149	-4.0	106	0.00
5 S	Phenol-d5	1.810	1.763	2.6	105	0.00
6	Phenol	1.825	1.805	1.1	105	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.117	1.121	-0.4	108	0.00
8	Bis(2-Chloroethyl)ether	1.415	1.419	-0.3	107	0.00
9 S	2-Chlorophenol-d4	1.435	1.415	1.4	105	0.00
10	2-Chlorophenol	1.453	1.427	1.8	107	0.00
11	2-Methylphenol	1.396	1.365	2.2	105	0.00
12	2,2'-oxybis(1-Chloropropane	2.220	2.234	-0.6	106	0.00
13 S	4-Methylphenol-d8	1.450	1.417	2.3	105	0.00
14	Acetophenone	2.227	2.243	-0.7	105	0.00
15 P	N-Nitroso-di-n-propylamine	1.148	1.173	-2.2	108	0.00
16	4-Methylphenol	1.521	1.517	0.3	105	0.00
17	Hexachloroethane	0.573	0.565	1.4	107	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
19 S	Nitrobenzene-d5	0.149	0.154	-3.4	110	0.00
20	Nitrobenzene	0.365	0.368	-0.8	109	0.00
21	Isophorone	0.716	0.727	-1.5	107	0.00
22 S	2-Nitrophenol-d4	0.154	0.164	-6.5	113	0.00
23 C	2-Nitrophenol	0.167	0.178	-6.6	111	0.00
24	2,4-Dimethylphenol	0.377	0.377	0.0	106	0.00
25	Bis(2-Chloroethoxy)methane	0.451	0.449	0.4	107	0.00
26 S	2,4-Dichlorophenol-d3	0.328	0.327	0.3	105	0.00
27 C	2,4-Dichlorophenol	0.319	0.316	0.9	105	0.00
28	Naphthalene	1.061	1.035	2.5	104	0.00
29 S	4-Chloroaniline-d4	0.340	0.323	5.0	103	0.00
30	4-Chloroaniline	0.336	0.321	4.5	103	0.00
31 C	Hexachlorobutadiene	0.195	0.195	0.0	108	0.00
32	Caprolactam	0.111	0.104	6.3	101	0.00
33 C	4-Chloro-3-methylphenol	0.351	0.342	2.6	102	0.00
34	2-Methylnaphthalene	0.770	0.756	1.8	105	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	104	0.00
36	1,2,4,5-Tetrachlorobenzene	0.650	0.658	-1.2	106	0.00
37	Hexachlorocyclopentadiene	0.396	0.414	-4.5	111	0.00
38 C	2,4,6-Trichlorophenol	0.424	0.432	-1.9	105	0.00
39	2,4,5-Trichlorophenol	0.451	0.455	-0.9	103	0.00
40	1,1'-Biphenyl	1.661	1.665	-0.2	104	0.00
41	2-Chloronaphthalene	1.297	1.296	0.1	104	0.00
42	2-Nitroaniline	0.362	0.375	-3.6	103	0.00
43 S	Dimethylphthalate-d6	1.689	1.663	1.5	102	0.00
44	Dimethylphthalate	1.660	1.625	2.1	101	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN061918\
 Data File : BN001876.D
 Acq On : 19 Jun 2018 14:51
 Operator : JU/SJ
 Sample : SSTDICV020
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV87

Quant Time: Jun 19 15:27:51 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 19 14:50:02 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	0.313	0.325	-3.8	103	0.00
46 S	Acenaphthylene-d8	2.085	2.081	0.2	103	0.00
47	Acenaphthylene	2.013	2.028	-0.7	103	0.00
48	3-Nitroaniline	0.301	0.282	6.3	98	0.00
49 C	Acenaphthene	1.396	1.375	1.5	102	0.00
50	2,4-Dinitrophenol	0.161	0.140	13.0	109	0.00
51 S	4-Nitrophenol-d4	0.311	0.286	8.0	98	0.00
52	4-Nitrophenol	0.231	0.214	7.4	97	0.00
53	Dibenzofuran	2.003	1.959	2.2	102	0.00
54	2,4-Dinitrotoluene	0.466	0.472	-1.3	101	0.00
55	2,3,4,6-Tetrachlorophenol	0.398	0.392	1.5	101	0.00
56	Diethylphthalate	1.682	1.639	2.6	101	0.00
57 S	Fluorene-d10	1.480	1.432	3.2	101	0.00
58	Fluorene	1.598	1.538	3.8	100	0.00
59	4-Chlorophenyl-phenylether	0.801	0.786	1.9	103	0.00
60	4-Nitroaniline	0.372	0.349	6.2	97	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	97	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.109	0.106	2.8	102	0.00
63	4,6-Dinitro-2-methylphenol	0.116	0.117	-0.9	104	0.00
64	N-Nitrosodiphenylamine	0.598	0.622	-4.0	100	0.00
65	4-Bromophenyl-phenylether	0.217	0.226	-4.1	100	0.00
66	Hexachlorobenzene	0.245	0.245	0.0	97	0.00
67	Atrazine	0.218	0.220	-0.9	96	0.00
68 C	Pentachlorophenol	0.135	0.128	5.2	96	0.00
69	Phenanthrene	1.127	1.117	0.9	96	0.00
70 S	Anthracene-d10	0.966	0.975	-0.9	97	0.00
71	Anthracene	1.140	1.144	-0.4	96	0.00
72	Carbazole	0.955	0.992	-3.9	93	0.00
73	Di-n-butylphthalate	1.206	1.234	-2.3	95	0.00
74 C	Fluoranthene	1.184	1.247	-5.3	91	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	81	0.00
76 S	Pyrene-d10	1.008	1.148	-13.9	89	0.00
77	Pyrene	1.252	1.434	-14.5	89	0.00
78	Butylbenzylphthalate	0.534	0.582	-9.0	85	0.00
79	3,3'-Dichlorobenzidine	0.380	0.348	8.4	75	0.00
80	Benzo(a)anthracene	1.264	1.267	-0.2	80	0.00
81	Bis(2-ethylhexyl)phthalate	0.796	0.826	-3.8	81	0.00
82	Chrysene	1.176	1.169	0.6	80	0.00
83 I	Perylene-d12	1.000	1.000	0.0	73	0.00
84	Di-n-octyl phthalate	1.326	1.468	-10.7	71	0.00
85	Benzo(b)fluoranthene	1.263	1.328	-5.1	75	0.00
86	Benzo(k)fluoranthene	1.225	1.209	1.3	70	0.00
87 S	Benzo(a)pyrene-d12	1.088	1.088	0.0	72	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN061918\
Data File : BN001876.D
Acq On : 19 Jun 2018 14:51
Operator : JU/SJ
Sample : SSTDICV020
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SICV87

Quant Time: Jun 19 15:27:51 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jun 19 14:50:02 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	1.180	1.187	-0.6	73	0.00
89	Indeno(1,2,3-cd)pyrene	1.256	1.258	-0.2	77	0.00
90	Dibenzo(a,h)anthracene	1.045	1.060	-1.4	78	0.00
91	Benzo(g,h,i)perylene	1.036	1.040	-0.4	78	0.00

(#) = Out of Range

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