

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG032719\
 Data File : BG040182.D
 Acq On : 27 Mar 2019 21:44
 Operator : JU/SJ
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ICVBG032719

Quant Time: Mar 28 06:35:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG032719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 28 05:47:00 2019
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	107	0.00
2	1,4-Dioxane	0.566	0.569	-0.5	106	0.00
3	Pyridine	1.523	1.573	-3.3	102	0.00
4	n-Nitrosodimethylamine	0.705	0.717	-1.7	107	0.00
5 S	2-Fluorophenol	1.203	1.224	-1.7	105	0.00
6	Aniline	2.160	2.175	-0.7	103	0.00
7 S	Phenol-d6	1.663	1.689	-1.6	105	0.00
8	2-Chlorophenol	1.408	1.428	-1.4	105	0.00
9	Benzaldehyde	1.140	1.192	-4.6	104	0.00
10 C	Phenol	1.805	1.844	-2.2	106	0.00
11	bis(2-Chloroethyl)ether	1.445	1.464	-1.3	104	0.00
12	1,3-Dichlorobenzene	1.531	1.549	-1.2	104	0.00
13 C	1,4-Dichlorobenzene	1.525	1.579	-3.5	107	0.00
14	1,2-Dichlorobenzene	1.458	1.508	-3.4	107	0.00
15	Benzyl Alcohol	1.339	1.350	-0.8	104	0.00
16	2,2'-oxybis(1-Chloropropane	2.774	2.750	0.9	102	0.00
17	2-Methylphenol	1.249	1.239	0.8	102	0.00
18	Hexachloroethane	0.580	0.587	-1.2	105	0.00
19 P	n-Nitroso-di-n-propylamine	1.192	1.188	0.3	104	0.00
20	3+4-Methylphenols	1.692	1.696	-0.2	103	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
22	Acetophenone	0.499	0.511	-2.4	104	0.00
23 S	Nitrobenzene-d5	0.376	0.388	-3.2	103	0.00
24	Nitrobenzene	0.390	0.404	-3.6	104	0.00
25	Isophorone	0.774	0.795	-2.7	103	0.00
26 C	2-Nitrophenol	0.185	0.195	-5.4	106	0.00
27	2,4-Dimethylphenol	0.293	0.303	-3.4	104	0.00
28	bis(2-Chloroethoxy)methane	0.458	0.474	-3.5	102	0.00
29 C	2,4-Dichlorophenol	0.318	0.331	-4.1	103	0.00
30	1,2,4-Trichlorobenzene	0.350	0.364	-4.0	105	0.00
31	Naphthalene	1.025	1.075	-4.9	104	0.00
32	Benzoic acid	0.177	0.178	-0.6	108	0.00
33	4-Chloroaniline	0.437	0.453	-3.7	102	0.00
34 C	Hexachlorobutadiene	0.224	0.235	-4.9	106	0.00
35	Caprolactam	0.126	0.126	0.0	99	0.00
36 C	4-Chloro-3-methylphenol	0.366	0.375	-2.5	102	0.00
37	2-Methylnaphthalene	0.758	0.786	-3.7	104	0.00
38	1-Methylnaphthalene	0.735	0.756	-2.9	103	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	102	0.00
40	1,2,4,5-Tetrachlorobenzene	0.636	0.661	-3.9	104	0.00
41 P	Hexachlorocyclopentadiene	0.349	0.353	-1.1	104	0.00
42 S	2,4,6-Tribromophenol	0.216	0.231	-6.9	107	0.00
43 C	2,4,6-Trichlorophenol	0.410	0.429	-4.6	105	0.00
44	2,4,5-Trichlorophenol	0.447	0.456	-2.0	104	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG032719\
 Data File : BG040182.D
 Acq On : 27 Mar 2019 21:44
 Operator : JU/SJ
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ICVBG032719

Quant Time: Mar 28 06:35:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG032719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 28 05:47:00 2019
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
45 S	2-Fluorobiphenyl	1.350	1.415	-4.8	104	0.00
46	1,1'-Biphenyl	1.580	1.652	-4.6	105	0.00
47	2-Chloronaphthalene	1.212	1.256	-3.6	104	0.00
48	2-Nitroaniline	0.409	0.421	-2.9	102	0.00
49	Acenaphthylene	2.055	2.148	-4.5	104	0.00
50	Dimethylphthalate	1.565	1.624	-3.8	104	0.00
51	2,6-Dinitrotoluene	0.351	0.370	-5.4	106	0.00
52 C	Acenaphthene	1.178	1.213	-3.0	104	0.00
53	3-Nitroaniline	0.372	0.384	-3.2	103	0.00
54 P	2,4-Dinitrophenol	0.187	0.196	-4.8	112	0.00
55	Dibenzofuran	1.892	1.987	-5.0	105	0.00
56 P	4-Nitrophenol	0.300	0.313	-4.3	106	0.00
57	2,4-Dinitrotoluene	0.488	0.513	-5.1	105	0.00
58	Fluorene	1.488	1.543	-3.7	104	0.00
59	2,3,4,6-Tetrachlorophenol	0.405	0.417	-3.0	102	0.00
60	Diethylphthalate	1.602	1.664	-3.9	104	0.00
61	4-Chlorophenyl-phenylether	0.795	0.822	-3.4	103	0.00
62	4-Nitroaniline	0.399	0.412	-3.3	104	0.00
63	Azobenzene	1.511	1.547	-2.4	103	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	105	0.00
65	4,6-Dinitro-2-methylphenol	0.112	0.115	-2.7	111	0.00
66 c	n-Nitrosodiphenylamine	0.585	0.594	-1.5	103	0.00
67	4-Bromophenyl-phenylether	0.225	0.234	-4.0	106	0.00
68	Hexachlorobenzene	0.226	0.230	-1.8	105	0.00
69	Atrazine	0.218	0.221	-1.4	103	0.00
70 C	Pentachlorophenol	0.131	0.131	0.0	105	0.00
71	Phenanthrene	1.051	1.085	-3.2	105	0.00
72	Anthracene	1.052	1.089	-3.5	105	0.00
73	Carbazole	0.996	1.025	-2.9	104	0.00
74	Di-n-butylphthalate	1.180	1.216	-3.1	104	0.00
75 C	Fluoranthene	1.245	1.291	-3.7	105	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	104	0.00
77	Benzidine	0.722	0.692	4.2	96	0.00
78	Pyrene	1.315	1.357	-3.2	105	0.00
79 S	Terphenyl-d14	0.889	0.904	-1.7	104	0.00
80	Butylbenzylphthalate	0.563	0.580	-3.0	105	0.00
81	Benzo(a)anthracene	1.232	1.258	-2.1	104	0.00
82	3,3'-Dichlorobenzidine	0.469	0.470	-0.2	101	0.00
83	Chrysene	1.184	1.199	-1.3	103	0.00
84	Bis(2-ethylhexyl)phthalate	0.805	0.814	-1.1	104	0.00
85 c	Di-n-octyl phthalate	1.371	1.387	-1.2	102	0.00
86	Indeno(1,2,3-cd)pyrene	1.448	1.489	-2.8	106	0.00
87 I	Perylene-d12	1.000	1.000	0.0	105	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG032719\
Data File : BG040182.D
Acq On : 27 Mar 2019 21:44
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ICVBG032719

Quant Time: Mar 28 06:35:49 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG032719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 28 05:47:00 2019
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.224	1.223	0.1	102	0.00
89	Benzo(k)fluoranthene	1.126	1.155	-2.6	106	0.00
90 C	Benzo(a)pyrene	1.149	1.158	-0.8	103	0.00
91	Dibenzo(a,h)anthracene	1.067	1.097	-2.8	106	0.00
92	Benzo(g,h,i)perylene	1.097	1.120	-2.1	105	-0.01

(#) = Out of Range

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