

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011025\
 Data File : BG063951.D
 Acq On : 10 Jan 2025 15:25
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Jan 10 17:06:42 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG010325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 03 23:17:54 2025
 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)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	124	-0.03
2	1,4-Dioxane	0.660	0.658	0.3	126	-0.04
3	Pyridine	1.778	1.779	-0.1	120	-0.04
4	n-Nitrosodimethylamine	0.680	0.663	2.5	117	-0.04
5 S	2-Fluorophenol	1.255	1.323	-5.4	129	-0.03
6	Aniline	1.577	1.635	-3.7	123	-0.03
7 S	Phenol-d6	1.816	1.866	-2.8	124	-0.03
8	2-Chlorophenol	1.220	1.238	-1.5	124	-0.03
9	Benzaldehyde	1.160	1.177	-1.5	124	-0.03
10 C	Phenol	1.891	1.935	-2.3	123	-0.02
11	bis(2-Chloroethyl)ether	1.482	1.487	-0.3	122	-0.03
12	1,3-Dichlorobenzene	1.471	1.465	0.4	121	-0.03
13 C	1,4-Dichlorobenzene	1.470	1.528	-3.9	124	-0.03
14	1,2-Dichlorobenzene	1.447	1.481	-2.3	126	-0.03
15	Benzyl Alcohol	1.303	1.292	0.8	116	-0.02
16	2,2'-oxybis(1-Chloropropane	2.134	2.157	-1.1	122	-0.03
17	2-Methylphenol	1.227	1.288	-5.0	127	-0.03
18	Hexachloroethane	0.570	0.575	-0.9	127	-0.02
19 P	n-Nitroso-di-n-propylamine	1.326	1.282	3.3	114	-0.02
20	3+4-Methylphenols	1.651	1.696	-2.7	120	-0.02
21 I	Naphthalene-d8	1.000	1.000	0.0	119	-0.02
22	Acetophenone	0.601	0.587	2.3	119	-0.03
23 S	Nitrobenzene-d5	0.508	0.499	1.8	119	-0.02
24	Nitrobenzene	0.545	0.538	1.3	119	-0.02
25	Isophorone	0.895	0.883	1.3	119	-0.03
26 C	2-Nitrophenol	0.173	0.179	-3.5	121	-0.03
27	2,4-Dimethylphenol	0.221	0.223	-0.9	122	-0.02
28	bis(2-Chloroethoxy)methane	0.534	0.527	1.3	118	-0.02
29 C	2,4-Dichlorophenol	0.330	0.339	-2.7	125	-0.02
30	1,2,4-Trichlorobenzene	0.399	0.399	0.0	122	-0.02
31	Naphthalene	1.084	1.100	-1.5	125	-0.02
32	Benzoic acid	0.148	0.151	-2.0	122	-0.02
33	4-Chloroaniline	0.352	0.375	-6.5	125	-0.02
34 C	Hexachlorobutadiene	0.273	0.260	4.8	117	-0.03
35	Caprolactam	0.109	0.116	-6.4	127	-0.02
36 C	4-Chloro-3-methylphenol	0.366	0.374	-2.2	123	-0.03
37	2-Methylnaphthalene	0.762	0.765	-0.4	122	-0.02
38	1-Methylnaphthalene	0.758	0.757	0.1	123	-0.02
39 I	Acenaphthene-d10	1.000	1.000	0.0	115	-0.02
40	1,2,4,5-Tetrachlorobenzene	0.671	0.660	1.6	118	-0.02
41 P	Hexachlorocyclopentadiene	0.079	0.075	5.1	101	-0.02
42 S	2,4,6-Tribromophenol	0.241	0.228	5.4	112	-0.02
43 C	2,4,6-Trichlorophenol	0.395	0.402	-1.8	118	-0.02
44	2,4,5-Trichlorophenol	0.444	0.460	-3.6	120	-0.02
45 S	2-Fluorobiphenyl	1.381	1.385	-0.3	120	-0.02
46	1,1'-Biphenyl	1.460	1.491	-2.1	123	-0.02
47	2-Chloronaphthalene	1.200	1.219	-1.6	120	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48	2-Nitroaniline	0.402	0.412	-2.5	119	-0.02
49	Acenaphthylene	1.778	1.804	-1.5	122	-0.02
50	Dimethylphthalate	1.461	1.460	0.1	120	-0.02
51	2,6-Dinitrotoluene	0.325	0.341	-4.9	122	-0.02
52 C	Acenaphthene	1.087	1.105	-1.7	122	-0.02
53	3-Nitroaniline	0.299	0.318	-6.4	123	-0.02
54 P	2,4-Dinitrophenol	0.140	0.130	7.1	101	-0.02
55	Dibenzofuran	1.887	1.858	1.5	119	-0.02
56 P	4-Nitrophenol	0.205	0.217	-5.9	114	-0.02
57	2,4-Dinitrotoluene	0.457	0.465	-1.8	122	-0.02
58	Fluorene	1.466	1.433	2.3	119	-0.02
59	2,3,4,6-Tetrachlorophenol	0.391	0.366	6.4	108	-0.02
60	Diethylphthalate	1.440	1.423	1.2	121	-0.02
61	4-Chlorophenyl-phenylether	0.795	0.766	3.6	115	-0.02
62	4-Nitroaniline	0.310	0.325	-4.8	118	-0.02
63	Azobenzene	1.821	1.764	3.1	116	-0.02
64 I	Phenanthrene-d10	1.000	1.000	0.0	111	-0.02
65	4,6-Dinitro-2-methylphenol	0.104	0.109	-4.8	108	-0.02
66 c	n-Nitrosodiphenylamine	0.521	0.574	-10.2	120	-0.02
67	4-Bromophenyl-phenylether	0.221	0.234	-5.9	116	-0.02
68	Hexachlorobenzene	0.252	0.259	-2.8	114	-0.02
69	Atrazine	0.167	0.148	11.4	104	-0.02
70 C	Pentachlorophenol	0.105	0.104	1.0	100	-0.02
71	Phenanthrene	1.040	1.089	-4.7	115	-0.02
72	Anthracene	1.013	1.077	-6.3	116	-0.02
73	Carbazole	0.935	1.000	-7.0	117	-0.02
74	Di-n-butylphthalate	1.042	1.093	-4.9	114	-0.02
75 C	Fluoranthene	1.313	1.357	-3.4	113	-0.02
76 I	Chrysene-d12	1.000	1.000	0.0	106	-0.02
77	Benzydine	0.280	0.305	-8.9	103	-0.02
78	Pyrene	1.372	1.394	-1.6	112	-0.02
79 S	Terphenyl-d14	1.007	0.975	3.2	108	-0.02
80	Butylbenzylphthalate	0.423	0.419	0.9	109	-0.02
81	Benzo(a)anthracene	1.338	1.321	1.3	110	-0.02
82	3,3'-Dichlorobenzidine	0.416	0.423	-1.7	109	-0.02
83	Chrysene	1.299	1.288	0.8	111	-0.02
84	Bis(2-ethylhexyl)phthalate	0.647	0.642	0.8	108	-0.02
85 c	Di-n-octyl phthalate	1.127	1.051	6.7	101	-0.02
86 I	Perylene-d12	1.000	1.000	0.0	106	-0.04
87	Indeno(1,2,3-cd)pyrene	1.389	1.453	-4.6	112	-0.07
88	Benzo(b)fluoranthene	1.250	1.306	-4.5	112	-0.03
89	Benzo(k)fluoranthene	1.255	1.224	2.5	103	-0.03
90 C	Benzo(a)pyrene	1.110	1.148	-3.4	111	-0.04
91	Dibenzo(a,h)anthracene	1.149	1.178	-2.5	109	-0.06
92	Benzo(g,h,i)perylene	1.182	1.236	-4.6	111	-0.06

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(#) = Out of Range

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