

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012020\
 Data File : BG044098.D
 Acq On : 20 Jan 2020 15:06
 Operator : JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Jan 21 04:40:35 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 31 13:32:23 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	59	-0.03
2	1,4-Dioxane	0.475	0.499	-5.1	61	-0.04
3	Pyridine	1.403	1.462	-4.2	60	-0.03
4	n-Nitrosodimethylamine	0.625	0.600	4.0	57	-0.03
5 S	2-Fluorophenol	1.089	1.203	-10.5	63	-0.03
6	Aniline	2.000	2.065	-3.2	60	-0.03
7 S	Phenol-d6	1.523	1.688	-10.8	64	-0.02
8	2-Chlorophenol	1.279	1.346	-5.2	61	-0.03
9	Benzaldehyde	0.974	0.815	16.3	52	-0.03
10 C	Phenol	1.569	1.702	-8.5	63	-0.02
11	bis(2-Chloroethyl)ether	1.354	1.380	-1.9	59	-0.03
12	1,3-Dichlorobenzene	1.496	1.556	-4.0	61	-0.03
13 C	1,4-Dichlorobenzene	1.476	1.554	-5.3	61	-0.03
14	1,2-Dichlorobenzene	1.417	1.470	-3.7	61	-0.03
15	Benzyl Alcohol	1.202	1.218	-1.3	59	-0.03
16	2,2'-oxybis(1-Chloropropane	2.095	1.951	6.9	55	-0.03
17	2-Methylphenol	1.135	1.204	-6.1	62	-0.03
18	Hexachloroethane	0.575	0.594	-3.3	61	-0.03
19 P	n-Nitroso-di-n-propylamine	1.134	1.132	0.2	59	-0.03
20	3+4-Methylphenols	1.584	1.683	-6.2	62	-0.03
21 I	Naphthalene-d8	1.000	1.000	0.0	60	-0.03
22	Acetophenone	0.497	0.493	0.8	60	-0.03
23 S	Nitrobenzene-d5	0.374	0.361	3.5	60	-0.03
24	Nitrobenzene	0.370	0.357	3.5	59	-0.03
25	Isophorone	0.701	0.696	0.7	60	-0.03
26 C	2-Nitrophenol	0.175	0.191	-9.1	68	-0.03
27	2,4-Dimethylphenol	0.264	0.271	-2.7	64	-0.03
28	bis(2-Chloroethoxy)methane	0.468	0.462	1.3	60	-0.03
29 C	2,4-Dichlorophenol	0.315	0.332	-5.4	65	-0.02
30	1,2,4-Trichlorobenzene	0.353	0.354	-0.3	62	-0.03
31	Naphthalene	0.968	1.017	-5.1	63	-0.03
32	Benzoic acid	0.197	0.186	5.6	65	0.00
33	4-Chloroaniline	0.462	0.478	-3.5	62	-0.03
34 C	Hexachlorobutadiene	0.215	0.215	0.0	62	-0.03
35	Caprolactam	0.117	0.127	-8.5	66	-0.02
36 C	4-Chloro-3-methylphenol	0.335	0.375	-11.9	68	-0.02
37	2-Methylnaphthalene	0.729	0.777	-6.6	65	-0.03
38	1-Methylnaphthalene	0.691	0.736	-6.5	65	-0.03
39 I	Acenaphthene-d10	1.000	1.000	0.0	65	-0.03
40	1,2,4,5-Tetrachlorobenzene	0.586	0.571	2.6	64	-0.03
41 P	Hexachlorocyclopentadiene	0.304	0.279	8.2	60	-0.03
42 S	2,4,6-Tribromophenol	0.209	0.236	-12.9	72	-0.03
43 C	2,4,6-Trichlorophenol	0.403	0.415	-3.0	67	-0.03
44	2,4,5-Trichlorophenol	0.416	0.430	-3.4	67	-0.02

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	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
45 S	2-Fluorobiphenyl	1.104	1.201	-8.8	67	-0.03
46	1,1'-Biphenyl	1.434	1.474	-2.8	65	-0.03
47	2-Chloronaphthalene	1.159	1.176	-1.5	66	-0.03
48	2-Nitroaniline	0.373	0.377	-1.1	66	-0.02
49	Acenaphthylene	1.665	1.773	-6.5	68	-0.03
50	Dimethylphthalate	1.420	1.524	-7.3	68	-0.03
51	2,6-Dinitrotoluene	0.321	0.339	-5.6	70	-0.03
52 C	Acenaphthene	1.168	1.171	-0.3	65	-0.03
53	3-Nitroaniline	0.364	0.384	-5.5	68	-0.02
54 P	2,4-Dinitrophenol	0.145	0.185	-27.6#	84	-0.02
55	Dibenzofuran	1.608	1.758	-9.3	69	-0.02
56 P	4-Nitrophenol	0.279	0.283	-1.4	64	0.00
57	2,4-Dinitrotoluene	0.437	0.482	-10.3	71	-0.02
58	Fluorene	1.274	1.388	-8.9	69	-0.03
59	2,3,4,6-Tetrachlorophenol	0.358	0.376	-5.0	68	-0.02
60	Diethylphthalate	1.420	1.549	-9.1	69	-0.03
61	4-Chlorophenyl-phenylether	0.692	0.731	-5.6	69	-0.03
62	4-Nitroaniline	0.360	0.374	-3.9	66	-0.02
63	Azobenzene	1.317	1.382	-4.9	67	-0.03
64 I	Phenanthrene-d10	1.000	1.000	0.0	67	-0.02
65	4,6-Dinitro-2-methylphenol	0.097	0.120	-23.7	84	-0.02
66 c	n-Nitrosodiphenylamine	0.589	0.616	-4.6	71	-0.03
67	4-Bromophenyl-phenylether	0.225	0.229	-1.8	70	-0.03
68	Hexachlorobenzene	0.242	0.251	-3.7	71	-0.03
69	Atrazine	0.191	0.210	-9.9	70	-0.02
70 C	Pentachlorophenol	0.134	0.124	7.5	60	-0.02
71	Phenanthrene	0.959	1.034	-7.8	70	-0.03
72	Anthracene	0.948	1.021	-7.7	70	-0.03
73	Carbazole	0.876	0.936	-6.8	69	-0.02
74	Di-n-butylphthalate	1.118	1.185	-6.0	70	-0.03
75 C	Fluoranthene	1.097	1.161	-5.8	71	-0.02
76 I	Chrysene-d12	1.000	1.000	0.0	66	-0.03
77	Benzidine	0.503	0.500	0.6	58	-0.02
78	Pyrene	1.305	1.394	-6.8	71	-0.03
79 S	Terphenyl-d14	0.828	0.956	-15.5	74	-0.03
80	Butylbenzylphthalate	0.622	0.672	-8.0	70	-0.03
81	Benzo(a)anthracene	1.240	1.325	-6.9	69	-0.03
82	3,3'-Dichlorobenzidine	0.490	0.509	-3.9	67	-0.03
83	Chrysene	1.178	1.283	-8.9	70	-0.03
84	Bis(2-ethylhexyl)phthalate	0.858	0.916	-6.8	69	-0.03
85 c	Di-n-octyl phthalate	1.442	1.591	-10.3	70	-0.04
86	Indeno(1,2,3-cd)pyrene	1.601	1.644	-2.7	68	-0.07
87 I	Perylene-d12	1.000	1.000	0.0	66	-0.05

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88	Benzo(b)fluoranthene	1.182	1.211	-2.5	69	-0.04
89	Benzo(k)fluoranthene	1.124	1.162	-3.4	67	-0.04
90 C	Benzo(a)pyrene	1.116	1.146	-2.7	68	-0.04
91	Dibenzo(a,h)anthracene	1.121	1.146	-2.2	68	-0.07
92	Benzo(q,h,i)perylene	1.113	1.145	-2.9	69	-0.07

(#) = Out of Range

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