

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022625\
 Data File : BF141775.D
 Acq On : 26 Feb 2025 10:45
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Feb 26 11:15:26 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF020625.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 06 16:58:59 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	96	-0.01
2	1,4-Dioxane	0.513	0.531	-3.5	99	-0.02
3	Pyridine	1.222	1.317	-7.8	100	-0.02
4	n-Nitrosodimethylamine	0.809	0.726	10.3	85	-0.04
5 S	2-Fluorophenol	1.276	1.366	-7.1	102	0.00
6	Aniline	1.513	1.528	-1.0	95	-0.02
7 S	Phenol-d6	1.612	1.663	-3.2	99	-0.01
8	2-Chlorophenol	1.378	1.467	-6.5	102	-0.01
9	Benzaldehyde	0.857	0.907	-5.8	106	0.00
10 C	Phenol	1.678	1.722	-2.6	98	-0.01
11	bis(2-Chloroethyl)ether	1.261	1.285	-1.9	97	-0.01
12	1,3-Dichlorobenzene	1.455	1.552	-6.7	103	-0.01
13 C	1,4-Dichlorobenzene	1.487	1.554	-4.5	101	-0.01
14	1,2-Dichlorobenzene	1.380	1.467	-6.3	102	-0.01
15	Benzyl Alcohol	1.236	1.239	-0.2	94	0.00
16	2,2'-oxybis(1-Chloropropane	2.158	2.011	6.8	89	-0.01
17	2-Methylphenol	1.079	1.099	-1.9	98	0.00
18	Hexachloroethane	0.550	0.572	-4.0	99	-0.01
19 P	n-Nitroso-di-n-propylamine	0.964	0.977	-1.3	97	-0.02
20	3+4-Methylphenols	1.371	1.432	-4.4	100	-0.01
21 I	Naphthalene-d8	1.000	1.000	0.0	98	-0.01
22	Acetophenone	0.498	0.499	-0.2	99	-0.02
23 S	Nitrobenzene-d5	0.383	0.375	2.1	96	-0.01
24	Nitrobenzene	0.374	0.361	3.5	95	-0.01
25	Isophorone	0.611	0.607	0.7	98	-0.01
26 C	2-Nitrophenol	0.186	0.196	-5.4	100	0.00
27	2,4-Dimethylphenol	0.217	0.221	-1.8	98	0.00
28	bis(2-Chloroethoxy)methane	0.392	0.404	-3.1	101	-0.01
29 C	2,4-Dichlorophenol	0.294	0.303	-3.1	101	0.00
30	1,2,4-Trichlorobenzene	0.320	0.326	-1.9	101	0.00
31	Naphthalene	1.041	1.050	-0.9	99	-0.01
32	Benzoic acid	0.226	0.188	16.8	80	-0.02
33	4-Chloroaniline	0.363	0.381	-5.0	103	-0.02
34 C	Hexachlorobutadiene	0.192	0.196	-2.1	100	-0.01
35	Caprolactam	0.089	0.090	-1.1	95	-0.03
36 C	4-Chloro-3-methylphenol	0.325	0.327	-0.6	98	-0.01
37	2-Methylnaphthalene	0.677	0.692	-2.2	102	-0.01
38	1-Methylnaphthalene	0.656	0.662	-0.9	99	-0.01
39 I	Acenaphthene-d10	1.000	1.000	0.0	97	-0.02
40	1,2,4,5-Tetrachlorobenzene	0.579	0.600	-3.6	100	0.00
41 P	Hexachlorocyclopentadiene	0.192	0.186	3.1	89	-0.01
42 S	2,4,6-Tribromophenol	0.201	0.205	-2.0	100	-0.02
43 C	2,4,6-Trichlorophenol	0.374	0.378	-1.1	98	0.00
44	2,4,5-Trichlorophenol	0.405	0.404	0.2	96	0.00
45 S	2-Fluorobiphenyl	1.298	1.311	-1.0	101	-0.01
46	1,1'-Biphenyl	1.551	1.592	-2.6	100	-0.02
47	2-Chloronaphthalene	1.147	1.179	-2.8	101	-0.01

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48	2-Nitroaniline	0.385	0.371	3.6	93	-0.01
49	Acenaphthylene	1.705	1.716	-0.6	99	-0.01
50	Dimethylphthalate	1.358	1.404	-3.4	102	-0.02
51	2,6-Dinitrotoluene	0.297	0.311	-4.7	101	-0.01
52 C	Acenaphthene	1.147	1.148	-0.1	97	-0.02
53	3-Nitroaniline	0.319	0.316	0.9	96	-0.01
54 P	2,4-Dinitrophenol	0.176	0.165	6.2	91	0.00
55	Dibenzofuran	1.683	1.683	0.0	98	-0.02
56 P	4-Nitrophenol	0.237	0.220	7.2	86	0.00
57	2,4-Dinitrotoluene	0.395	0.392	0.8	96	-0.01
58	Fluorene	1.301	1.325	-1.8	100	-0.01
59	2,3,4,6-Tetrachlorophenol	0.349	0.345	1.1	96	-0.01
60	Diethylphthalate	1.336	1.362	-1.9	100	-0.02
61	4-Chlorophenyl-phenylether	0.626	0.660	-5.4	103	-0.02
62	4-Nitroaniline	0.324	0.308	4.9	90	-0.02
63	Azobenzene	1.328	1.308	1.5	96	-0.02
64 I	Phenanthrene-d10	1.000	1.000	0.0	94	-0.01
65	4,6-Dinitro-2-methylphenol	0.139	0.138	0.7	91	-0.01
66 c	n-Nitrosodiphenylamine	0.640	0.669	-4.5	100	-0.02
67	4-Bromophenyl-phenylether	0.224	0.239	-6.7	102	-0.02
68	Hexachlorobenzene	0.230	0.248	-7.8	103	-0.01
69	Atrazine	0.192	0.176	8.3	89	-0.02
70 C	Pentachlorophenol	0.147	0.154	-4.8	93	-0.01
71	Phenanthrene	1.101	1.138	-3.4	98	-0.02
72	Anthracene	1.107	1.135	-2.5	98	-0.02
73	Carbazole	0.996	0.991	0.5	94	-0.01
74	Di-n-butylphthalate	1.150	1.240	-7.8	102	-0.02
75 C	Fluoranthene	1.167	1.148	1.6	93	-0.02
76 I	Chrysene-d12	1.000	1.000	0.0	84	-0.02
77	Benzidine	0.441	0.446	-1.1	71	-0.01
78	Pyrene	1.703	1.867	-9.6	90	-0.02
79 S	Terphenyl-d14	1.185	1.316	-11.1	93	-0.02
80	Butylbenzylphthalate	0.590	0.688	-16.6	95	-0.02
81	Benzo(a)anthracene	1.329	1.363	-2.6	85	-0.02
82	3,3'-Dichlorobenzidine	0.381	0.427	-12.1	95	-0.02
83	Chrysene	1.228	1.285	-4.6	90	-0.02
84	Bis(2-ethylhexyl)phthalate	0.665	0.883	-32.8#	108	-0.02
85 c	Di-n-octyl phthalate	0.949	1.339	-41.1#	117	-0.04
86 I	Perylene-d12	1.000	1.000	0.0	104	-0.02
87	Indeno(1,2,3-cd)pyrene	1.236	1.406	-13.8	113	-0.04
88	Benzo(b)fluoranthene	1.343	1.304	2.9	101	-0.03
89	Benzo(k)fluoranthene	1.147	1.152	-0.4	104	-0.04
90 C	Benzo(a)pyrene	1.087	1.110	-2.1	106	-0.04
91	Dibenzo(a,h)anthracene	1.001	1.141	-14.0	112	-0.05
92	Benzo(g,h,i)perylene	1.048	1.186	-13.2	112	-0.05

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(#) = Out of Range

SPCC's out = 0 CCC's out = 1