

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111124\
 Data File : BP023007.D
 Acq On : 11 Nov 2024 10:12
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Nov 11 11:07:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 07 23:36:11 2024
 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	97	-0.01
2	1,4-Dioxane	0.421	0.434	-3.1	100	0.00
3	Pyridine	1.155	1.176	-1.8	98	0.00
4	n-Nitrosodimethylamine	0.557	0.647	-16.2	112	0.00
5 S	2-Fluorophenol	1.054	1.074	-1.9	97	0.00
6	Aniline	1.225	1.236	-0.9	97	0.00
7 S	Phenol-d6	1.470	1.457	0.9	96	-0.01
8	2-Chlorophenol	1.105	1.105	0.0	96	-0.01
9	Benzaldehyde	0.849	0.823	3.1	94	-0.01
10 C	Phenol	1.456	1.460	-0.3	96	-0.01
11	bis(2-Chloroethyl)ether	1.100	1.101	-0.1	97	-0.01
12	1,3-Dichlorobenzene	1.410	1.437	-1.9	100	-0.01
13 C	1,4-Dichlorobenzene	1.440	1.456	-1.1	100	-0.02
14	1,2-Dichlorobenzene	1.392	1.398	-0.4	99	-0.02
15	Benzyl Alcohol	1.113	1.125	-1.1	98	-0.01
16	2,2'-oxybis(1-Chloropropane	1.193	1.199	-0.5	99	-0.01
17	2-Methylphenol	0.988	0.997	-0.9	96	-0.01
18	Hexachloroethane	0.533	0.556	-4.3	103	-0.02
19 P	n-Nitroso-di-n-propylamine	1.006	1.030	-2.4	99	-0.01
20	3+4-Methylphenols	1.381	1.418	-2.7	99	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	103	-0.01
22	Acetophenone	0.532	0.504	5.3	99	0.00
23 S	Nitrobenzene-d5	0.423	0.423	0.0	102	-0.01
24	Nitrobenzene	0.430	0.427	0.7	101	-0.01
25	Isophorone	0.717	0.688	4.0	99	-0.01
26 C	2-Nitrophenol	0.177	0.173	2.3	100	-0.01
27	2,4-Dimethylphenol	0.202	0.195	3.5	100	0.00
28	bis(2-Chloroethoxy)methane	0.407	0.398	2.2	101	-0.01
29 C	2,4-Dichlorophenol	0.349	0.357	-2.3	103	-0.01
30	1,2,4-Trichlorobenzene	0.446	0.450	-0.9	105	-0.02
31	Naphthalene	1.002	0.978	2.4	101	-0.01
32	Benzoic acid	0.242	0.242	0.0	98	0.00
33	4-Chloroaniline	0.357	0.364	-2.0	103	-0.01
34 C	Hexachlorobutadiene	0.327	0.338	-3.4	106	-0.01
35	Caprolactam	0.104	0.106	-1.9	103	0.00
36 C	4-Chloro-3-methylphenol	0.389	0.407	-4.6	105	0.00
37	2-Methylnaphthalene	0.751	0.778	-3.6	107	-0.01
38	1-Methylnaphthalene	0.745	0.768	-3.1	105	-0.02
39 I	Acenaphthene-d10	1.000	1.000	0.0	114	-0.01
40	1,2,4,5-Tetrachlorobenzene	0.677	0.644	4.9	110	-0.01
41 P	Hexachlorocyclopentadiene	0.190	0.197	-3.7	111	-0.01
42 S	2,4,6-Tribromophenol	0.388	0.415	-7.0	123	-0.02
43 C	2,4,6-Trichlorophenol	0.431	0.415	3.7	108	-0.02
44	2,4,5-Trichlorophenol	0.490	0.469	4.3	107	-0.01
45 S	2-Fluorobiphenyl	1.400	1.325	5.4	110	-0.01
46	1,1'-Biphenyl	1.330	1.263	5.0	108	-0.01
47	2-Chloronaphthalene	1.103	1.058	4.1	111	-0.01

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48	2-Nitroaniline	0.321	0.332	-3.4	115	-0.01
49	Acenaphthylene	1.541	1.510	2.0	112	-0.02
50	Dimethylphthalate	1.464	1.414	3.4	111	-0.01
51	2,6-Dinitrotoluene	0.333	0.326	2.1	112	0.00
52 C	Acenaphthene	0.991	0.969	2.2	111	-0.02
53	3-Nitroaniline	0.271	0.280	-3.3	115	0.00
54 P	2,4-Dinitrophenol	0.207	0.212	-2.4	111	-0.02
55	Dibenzofuran	1.740	1.789	-2.8	120	-0.02
56 P	4-Nitrophenol	0.222	0.238	-7.2	115	0.00
57	2,4-Dinitrotoluene	0.481	0.499	-3.7	116	-0.02
58	Fluorene	1.446	1.471	-1.7	117	-0.01
59	2,3,4,6-Tetrachlorophenol	0.473	0.487	-3.0	116	0.00
60	Diethylphthalate	1.456	1.495	-2.7	118	-0.02
61	4-Chlorophenyl-phenylether	0.852	0.876	-2.8	120	-0.02
62	4-Nitroaniline	0.278	0.305	-9.7	118	-0.02
63	Azobenzene	1.252	1.364	-8.9	123	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	125	0.00
65	4,6-Dinitro-2-methylphenol	0.122	0.124	-1.6	120	-0.02
66 c	n-Nitrosodiphenylamine	0.494	0.476	3.6	120	-0.01
67	4-Bromophenyl-phenylether	0.240	0.232	3.3	121	-0.02
68	Hexachlorobenzene	0.307	0.298	2.9	124	-0.01
69	Atrazine	0.167	0.172	-3.0	117	-0.02
70 C	Pentachlorophenol	0.196	0.197	-0.5	121	0.00
71	Phenanthrene	0.968	0.956	1.2	124	0.00
72	Anthracene	0.929	0.938	-1.0	126	-0.02
73	Carbazole	0.887	0.893	-0.7	126	0.00
74	Di-n-butylphthalate	0.989	1.024	-3.5	128	-0.01
75 C	Fluoranthene	1.343	1.345	-0.1	123	-0.01
76 I	Chrysene-d12	1.000	1.000	0.0	126	-0.01
77	Benzidine	0.395	0.428	-8.4	114	-0.02
78	Pyrene	1.184	1.188	-0.3	128	0.00
79 S	Terphenyl-d14	1.065	1.078	-1.2	130	-0.01
80	Butylbenzylphthalate	0.404	0.414	-2.5	127	0.00
81	Benzo(a)anthracene	1.245	1.231	1.1	128	-0.01
82	3,3'-Dichlorobenzidine	0.504	0.504	0.0	128	0.00
83	Chrysene	1.166	1.177	-0.9	129	0.00
84	Bis(2-ethylhexyl)phthalate	0.579	0.595	-2.8	132	-0.02
85 c	Di-n-octyl phthalate	0.968	1.012	-4.5	131	-0.02
86 I	Perylene-d12	1.000	1.000	0.0	127	-0.02
87	Indeno(1,2,3-cd)pyrene	1.349	1.354	-0.4	129	-0.02
88	Benzo(b)fluoranthene	1.175	1.185	-0.9	125	-0.02
89	Benzo(k)fluoranthene	1.110	1.130	-1.8	133	-0.02
90 C	Benzo(a)pyrene	1.005	1.012	-0.7	128	0.00
91	Dibenzo(a,h)anthracene	1.106	1.117	-1.0	132	-0.01
92	Benzo(g,h,i)perylene	1.133	1.111	1.9	125	-0.02

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(#) = Out of Range

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