

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092324\
 Data File : BP021988.D
 Acq On : 23 Sep 2024 18:52
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Sep 23 19:18:09 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 20 01:22:00 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	149	0.00
2	1,4-Dioxane	0.442	0.411	7.0	142	0.00
3	Pyridine	1.179	1.144	3.0	146	0.00
4	n-Nitrosodimethylamine	0.622	0.599	3.7	147	0.00
5 S	2-Fluorophenol	1.064	1.037	2.5	146	0.00
6	Aniline	1.214	1.239	-2.1	153#	0.00
7 S	Phenol-d6	1.382	1.411	-2.1	154#	0.00
8	2-Chlorophenol	1.138	1.146	-0.7	150	0.00
9	Benzaldehyde	0.722	0.788	-9.1	159#	0.00
10 C	Phenol	1.372	1.408	-2.6	155#	0.00
11	bis(2-Chloroethyl)ether	1.034	1.052	-1.7	155#	0.00
12	1,3-Dichlorobenzene	1.437	1.434	0.2	150#	0.00
13 C	1,4-Dichlorobenzene	1.457	1.444	0.9	149	0.00
14	1,2-Dichlorobenzene	1.398	1.386	0.9	148	0.00
15	Benzyl Alcohol	1.060	1.134	-7.0	159#	0.00
16	2,2'-oxybis(1-Chloropropane	1.102	1.174	-6.5	160#	0.00
17	2-Methylphenol	0.927	0.988	-6.6	155#	0.00
18	Hexachloroethane	0.536	0.540	-0.7	150	0.00
19 P	n-Nitroso-di-n-propylamine	0.859	0.950	-10.6	161#	0.00
20	3+4-Methylphenols	1.309	1.397	-6.7	158#	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	157#	0.00
22	Acetophenone	0.488	0.470	3.7	157#	0.00
23 S	Nitrobenzene-d5	0.378	0.379	-0.3	160#	0.00
24	Nitrobenzene	0.384	0.380	1.0	158#	0.00
25	Isophorone	0.623	0.632	-1.4	160#	0.00
26 C	2-Nitrophenol	0.168	0.166	1.2	154#	0.00
27	2,4-Dimethylphenol	0.194	0.195	-0.5	158#	0.00
28	bis(2-Chloroethoxy)methane	0.374	0.383	-2.4	163#	0.00
29 C	2,4-Dichlorophenol	0.330	0.333	-0.9	156#	0.00
30	1,2,4-Trichlorobenzene	0.411	0.396	3.6	154#	0.00
31	Naphthalene	0.998	0.989	0.9	159#	0.00
32	Benzoic acid	0.227	0.236	-4.0	157#	0.02
33	4-Chloroaniline	0.360	0.358	0.6	154#	0.00
34 C	Hexachlorobutadiene	0.321	0.308	4.0	156#	0.00
35	Caprolactam	0.098	0.102	-4.1	162#	0.00
36 C	4-Chloro-3-methylphenol	0.354	0.365	-3.1	162#	0.00
37	2-Methylnaphthalene	0.730	0.740	-1.4	163#	0.00
38	1-Methylnaphthalene	0.721	0.734	-1.8	163#	-0.01
39 I	Acenaphthene-d10	1.000	1.000	0.0	165#	0.00
40	1,2,4,5-Tetrachlorobenzene	0.700	0.684	2.3	163#	0.00
41 P	Hexachlorocyclopentadiene	0.260	0.249	4.2	162#	0.00
42 S	2,4,6-Tribromophenol	0.371	0.375	-1.1	171#	-0.02
43 C	2,4,6-Trichlorophenol	0.429	0.420	2.1	160#	0.00
44	2,4,5-Trichlorophenol	0.482	0.485	-0.6	167#	0.00
45 S	2-Fluorobiphenyl	1.370	1.323	3.4	164#	0.00
46	1,1'-Biphenyl	1.364	1.347	1.2	168#	-0.01
47	2-Chloronaphthalene	1.103	1.074	2.6	164#	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48	2-Nitroaniline	0.301	0.318	-5.6	166#	0.00
49	Acenaphthylene	1.556	1.555	0.1	164#	-0.01
50	Dimethylphthalate	1.451	1.408	3.0	162#	-0.01
51	2,6-Dinitrotoluene	0.323	0.324	-0.3	164#	0.00
52 C	Acenaphthene	1.013	1.009	0.4	167#	-0.02
53	3-Nitroaniline	0.288	0.289	-0.3	159#	0.00
54 P	2,4-Dinitrophenol	0.189	0.193	-2.1	164#	-0.01
55	Dibenzofuran	1.702	1.672	1.8	165#	-0.02
56 P	4-Nitrophenol	0.251	0.261	-4.0	161#	0.00
57	2,4-Dinitrotoluene	0.457	0.474	-3.7	167#	-0.01
58	Fluorene	1.421	1.392	2.0	165#	-0.01
59	2,3,4,6-Tetrachlorophenol	0.465	0.465	0.0	167#	0.00
60	Diethylphthalate	1.474	1.478	-0.3	170#	0.00
61	4-Chlorophenyl-phenylether	0.828	0.812	1.9	168#	-0.01
62	4-Nitroaniline	0.317	0.325	-2.5	163#	-0.01
63	Azobenzene	1.220	1.231	-0.9	168#	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	162#	-0.02
65	4,6-Dinitro-2-methylphenol	0.116	0.121	-4.3	165#	-0.02
66 c	n-Nitrosodiphenylamine	0.498	0.508	-2.0	165#	0.00
67	4-Bromophenyl-phenylether	0.233	0.241	-3.4	172#	0.00
68	Hexachlorobenzene	0.291	0.296	-1.7	169#	-0.01
69	Atrazine	0.178	0.148	16.9	150#	-0.01
70 C	Pentachlorophenol	0.196	0.201	-2.6	165#	-0.02
71	Phenanthrene	0.960	0.961	-0.1	165#	-0.01
72	Anthracene	0.929	0.947	-1.9	166#	0.00
73	Carbazole	0.898	0.913	-1.7	162#	0.00
74	Di-n-butylphthalate	1.054	1.088	-3.2	172#	0.00
75 C	Fluoranthene	1.340	1.340	0.0	163#	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	159#	0.00
77	Benzidine	0.244	0.385	-57.8#	151#	-0.01
78	Pyrene	1.210	1.219	-0.7	159#	0.00
79 S	Terphenyl-d14	1.065	1.070	-0.5	160#	0.00
80	Butylbenzylphthalate	0.432	0.447	-3.5	164#	0.00
81	Benzo(a)anthracene	1.285	1.264	1.6	159#	0.00
82	3,3'-Dichlorobenzidine	0.466	0.479	-2.8	153#	0.00
83	Chrysene	1.210	1.195	1.2	160#	-0.01
84	Bis(2-ethylhexyl)phthalate	0.633	0.651	-2.8	169#	0.00
85 c	Di-n-octyl phthalate	1.076	1.099	-2.1	167#	-0.02
86 I	Perylene-d12	1.000	1.000	0.0	152#	-0.02
87	Indeno(1,2,3-cd)pyrene	1.330	1.312	1.4	150#	-0.04
88	Benzo(b)fluoranthene	1.183	1.213	-2.5	154#	-0.02
89	Benzo(k)fluoranthene	1.119	1.116	0.3	155#	-0.02
90 C	Benzo(a)pyrene	1.019	1.030	-1.1	153#	0.00
91	Dibenzo(a,h)anthracene	1.098	1.076	2.0	151#	-0.04
92	Benzo(g,h,i)perylene	1.125	1.116	0.8	150	-0.02

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(#) = Out of Range

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