

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050420\
 Data File : BG045231.D
 Acq On : 4 May 2020 12:25
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: May 04 13:05:56 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 04 13:03:14 2020
 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	78	0.00
2	1,4-Dioxane	0.538	0.601	-11.7	87	0.00
3	Pyridine	1.658	1.401	15.5	68	0.00
4	n-Nitrosodimethylamine	0.508	0.530	-4.3	85	0.00
5 S	2-Fluorophenol	1.211	1.170	3.4	77	0.00
6	Aniline	2.340	2.151	8.1	72	0.00
7 S	Phenol-d6	1.800	1.714	4.8	75	0.00
8	2-Chlorophenol	1.302	1.289	1.0	78	0.00
9	Benzaldehyde	1.067	1.055	1.1	81	0.00
10 C	Phenol	1.801	1.764	2.1	77	0.00
11	bis(2-Chloroethyl)ether	1.434	1.277	10.9	70	0.00
12	1,3-Dichlorobenzene	1.534	1.466	4.4	77	0.00
13 C	1,4-Dichlorobenzene	1.529	1.490	2.6	77	0.00
14	1,2-Dichlorobenzene	1.463	1.370	6.4	73	0.00
15	Benzyl Alcohol	1.509	1.308	13.3	68	0.00
16	2,2'-oxybis(1-Chloropropane	1.408	1.217	13.6	68	0.00
17	2-Methylphenol	1.390	1.166	16.1	67	0.00
18	Hexachloroethane	0.575	0.570	0.9	77	0.00
19 P	n-Nitroso-di-n-propylamine	1.273	1.089	14.5	67	0.00
20	3+4-Methylphenols	1.945	1.606	17.4	64	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	69	0.00
22	Acetophenone	0.541	0.561	-3.7	74	0.00
23 S	Nitrobenzene-d5	0.438	0.412	5.9	68	0.00
24	Nitrobenzene	0.449	0.449	0.0	71	0.00
25	Isophorone	0.898	0.821	8.6	64	0.00
26 C	2-Nitrophenol	0.194	0.192	1.0	70	0.00
27	2,4-Dimethylphenol	0.307	0.318	-3.6	74	0.00
28	bis(2-Chloroethoxy)methane	0.472	0.459	2.8	68	0.00
29 C	2,4-Dichlorophenol	0.355	0.345	2.8	69	0.00
30	1,2,4-Trichlorobenzene	0.401	0.405	-1.0	72	0.00
31	Naphthalene	1.094	1.098	-0.4	70	0.00
32	Benzoic acid	0.230	0.207	10.0	64	0.00
33	4-Chloroaniline	0.510	0.464	9.0	65	0.00
34 C	Hexachlorobutadiene	0.275	0.272	1.1	71	0.00
35	Caprolactam	0.141	0.129	8.5	64	0.00
36 C	4-Chloro-3-methylphenol	0.417	0.426	-2.2	73	0.00
37	2-Methylnaphthalene	0.849	0.842	0.8	70	0.00
38	1-Methylnaphthalene	0.802	0.779	2.9	68	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	71	0.00
40	1,2,4,5-Tetrachlorobenzene	0.644	0.638	0.9	72	0.00
41 P	Hexachlorocyclopentadiene	0.402	0.376	6.5	66	0.00
42 S	2,4,6-Tribromophenol	0.309	0.284	8.1	67	0.00
43 C	2,4,6-Trichlorophenol	0.454	0.429	5.5	68	0.00
44	2,4,5-Trichlorophenol	0.517	0.460	11.0	64	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050420\
 Data File : BG045231.D
 Acq On : 4 May 2020 12:25
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: May 04 13:05:56 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 04 13:03:14 2020
 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)
45 S	2-Fluorobiphenyl	1.364	1.277	6.4	67	0.00
46	1,1'-Biphenyl	1.520	1.468	3.4	70	0.00
47	2-Chloronaphthalene	1.162	1.081	7.0	68	0.00
48	2-Nitroaniline	0.382	0.370	3.1	72	0.00
49	Acenaphthylene	1.892	1.812	4.2	69	0.00
50	Dimethylphthalate	1.628	1.516	6.9	67	0.00
51	2,6-Dinitrotoluene	0.364	0.342	6.0	69	0.00
52 C	Acenaphthene	1.280	1.201	6.2	68	0.00
53	3-Nitroaniline	0.399	0.349	12.5	64	0.00
54 P	2,4-Dinitrophenol	0.217	0.178	18.0	61	0.00
55	Dibenzofuran	1.866	1.787	4.2	70	0.00
56 P	4-Nitrophenol	0.310	0.287	7.4	68	0.00
57	2,4-Dinitrotoluene	0.532	0.502	5.6	70	0.00
58	Fluorene	1.524	1.491	2.2	71	0.00
59	2,3,4,6-Tetrachlorophenol	0.460	0.448	2.6	72	0.00
60	Diethylphthalate	1.698	1.566	7.8	67	0.00
61	4-Chlorophenyl-phenylether	0.877	0.808	7.9	67	0.00
62	4-Nitroaniline	0.414	0.388	6.3	69	0.00
63	Azobenzene	1.565	1.539	1.7	72	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	73	0.00
65	4,6-Dinitro-2-methylphenol	0.131	0.126	3.8	74	0.00
66 c	n-Nitrosodiphenylamine	0.575	0.539	6.3	70	0.00
67	4-Bromophenyl-phenylether	0.258	0.230	10.9	67	0.00
68	Hexachlorobenzene	0.268	0.248	7.5	70	0.00
69	Atrazine	0.230	0.229	0.4	79	0.00
70 C	Pentachlorophenol	0.164	0.146	11.0	65	0.00
71	Phenanthrene	1.028	1.014	1.4	74	0.00
72	Anthracene	1.031	1.034	-0.3	75	0.00
73	Carbazole	1.046	0.974	6.9	72	0.00
74	Di-n-butylphthalate	1.227	1.160	5.5	72	0.00
75 C	Fluoranthene	1.305	1.287	1.4	77	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	72	0.00
77	Benzidine	0.644	0.563	12.6	65	0.00
78	Pyrene	1.379	1.424	-3.3	77	0.00
79 S	Terphenyl-d14	0.918	0.998	-8.7	76	0.00
80	Butylbenzylphthalate	0.572	0.582	-1.7	74	0.00
81	Benzo(a)anthracene	1.296	1.344	-3.7	76	0.00
82	3,3'-Dichlorobenzidine	0.516	0.543	-5.2	77	0.00
83	Chrysene	1.245	1.304	-4.7	77	0.00
84	Bis(2-ethylhexyl)phthalate	0.786	0.817	-3.9	76	0.00
85 c	Di-n-octyl phthalate	1.359	1.414	-4.0	76	0.00
86	Indeno(1,2,3-cd)pyrene	1.654	1.746	-5.6	79	0.00
87 I	Perylene-d12	1.000	1.000	0.0	77	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050420\
Data File : BG045231.D
Acq On : 4 May 2020 12:25
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: May 04 13:05:56 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 04 13:03:14 2020
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)
88	Benzo(b)fluoranthene	1.253	1.201	4.2	77	0.00
89	Benzo(k)fluoranthene	1.191	1.182	0.8	78	0.00
90 C	Benzo(a)pyrene	1.183	1.213	-2.5	81	0.00
91	Dibenzo(a,h)anthracene	1.192	1.198	-0.5	81	0.00
92	Benzo(q,h,i)perylene	1.195	1.194	0.1	80	0.00

(#) = Out of Range

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