

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110522\
 Data File : BG055467.D
 Acq On : 5 Nov 2022 12:28
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Nov 07 00:52:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 07 00:51:08 2022
 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	101	0.00
2	1,4-Dioxane	40.000	35.458	11.4	90	0.00
3	Pyridine	40.000	36.711	8.2	93	0.00
4	n-Nitrosodimethylamine	40.000	38.690	3.3	98	0.00
5 S	2-Fluorophenol	80.000	75.945	5.1	97	0.00
6	Aniline	40.000	37.206	7.0	98	0.00
7 S	Phenol-d6	80.000	76.993	3.8	99	0.00
8	2-Chlorophenol	40.000	37.515	6.2	99	0.00
9	Benzaldehyde	40.000	36.627	8.4	94	0.00
10 C	Phenol	40.000	38.074	4.8	99	0.00
11	bis(2-Chloroethyl)ether	40.000	36.766	8.1	97	0.00
12	1,3-Dichlorobenzene	40.000	37.215	7.0	98	0.00
13 C	1,4-Dichlorobenzene	40.000	37.371	6.6	98	0.00
14	1,2-Dichlorobenzene	40.000	37.646	5.9	100	0.00
15	Benzyl Alcohol	40.000	38.637	3.4	101	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	34.349	14.1	92	0.00
17	2-Methylphenol	40.000	38.006	5.0	100	0.00
18	Hexachloroethane	40.000	37.012	7.5	99	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	36.249	9.4	98	0.00
20	3+4-Methylphenols	40.000	38.244	4.4	102	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	104	0.00
22	Acetophenone	40.000	37.516	6.2	101	0.00
23 S	Nitrobenzene-d5	80.000	75.258	5.9	100	0.00
24	Nitrobenzene	40.000	37.261	6.8	99	0.00
25	Isophorone	40.000	37.022	7.4	101	0.00
26 C	2-Nitrophenol	40.000	38.444	3.9	103	0.00
27	2,4-Dimethylphenol	40.000	38.314	4.2	105	0.00
28	bis(2-Chloroethoxy)methane	40.000	36.638	8.4	101	0.00
29 C	2,4-Dichlorophenol	40.000	39.328	1.7	107	0.00
30	1,2,4-Trichlorobenzene	40.000	38.493	3.8	104	0.00
31	Naphthalene	40.000	37.986	5.0	103	0.00
32	Benzoic acid	40.000	39.087	2.3	112	0.00
33	4-Chloroaniline	40.000	39.001	2.5	105	0.00
34 C	Hexachlorobutadiene	40.000	38.104	4.7	104	0.00
35	Caprolactam	40.000	39.010	2.5	106	0.00
36 C	4-Chloro-3-methylphenol	40.000	39.266	1.8	106	0.00
37	2-Methylnaphthalene	40.000	38.443	3.9	106	0.00
38	1-Methylnaphthalene	40.000	38.281	4.3	106	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	108	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	38.266	4.3	107	0.00
41 P	Hexachlorocyclopentadiene	40.000	36.456	8.9	106	0.00
42 S	2,4,6-Tribromophenol	80.000	82.977	-3.7	114	0.00
43 C	2,4,6-Trichlorophenol	40.000	39.578	1.1	110	0.00
44	2,4,5-Trichlorophenol	40.000	39.405	1.5	109	0.00
45 S	2-Fluorobiphenyl	80.000	75.399	5.8	107	0.00
46	1,1'-Biphenyl	40.000	37.897	5.3	107	0.00
47	2-Chloronaphthalene	40.000	38.557	3.6	108	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110522\
 Data File : BG055467.D
 Acq On : 5 Nov 2022 12:28
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Nov 07 00:52:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 07 00:51:08 2022
 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	Amount	Calc.	%Dev	Area%	Dev(min)
48	2-Nitroaniline	40.000	38.659	3.4	106	0.00
49	Acenaphthylene	40.000	38.780	3.0	109	0.00
50	Dimethylphthalate	40.000	38.903	2.7	110	0.00
51	2,6-Dinitrotoluene	40.000	40.153	-0.4	112	0.00
52 C	Acenaphthene	40.000	38.490	3.8	108	0.00
53	3-Nitroaniline	40.000	40.112	-0.3	109	0.00
54 P	2,4-Dinitrophenol	40.000	37.566	6.1	108	0.00
55	Dibenzofuran	40.000	39.020	2.4	109	0.00
56 P	4-Nitrophenol	40.000	40.450	-1.1	105	0.00
57	2,4-Dinitrotoluene	40.000	40.796	-2.0	111	0.00
58	Fluorene	40.000	39.232	1.9	111	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	39.970	0.1	111	0.00
60	Diethylphthalate	40.000	38.567	3.6	109	0.00
61	4-Chlorophenyl-phenylether	40.000	39.010	2.5	111	0.00
62	4-Nitroaniline	40.000	39.931	0.2	109	0.00
63	Azobenzene	40.000	37.659	5.9	107	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	111	0.00
65	4,6-Dinitro-2-methylphenol	40.000	41.459	-3.6	111	0.00
66 c	n-Nitrosodiphenylamine	40.000	38.046	4.9	112	0.00
67	4-Bromophenyl-phenylether	40.000	38.287	4.3	112	0.00
68	Hexachlorobenzene	40.000	38.957	2.6	114	0.00
69	Atrazine	40.000	37.256	6.9	108	0.00
70 C	Pentachlorophenol	40.000	37.378	6.6	110	0.00
71	Phenanthrene	40.000	38.081	4.8	109	0.00
72	Anthracene	40.000	38.141	4.6	110	0.00
73	Carbazole	40.000	37.926	5.2	108	0.00
74	Di-n-butylphthalate	40.000	37.678	5.8	107	0.00
75 C	Fluoranthene	40.000	37.878	5.3	109	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	111	0.00
77	Benzydine	40.000	38.265	4.3	99	0.00
78	Pyrene	40.000	37.827	5.4	109	0.00
79 S	Terphenyl-d14	80.000	74.064	7.4	108	0.00
80	Butylbenzylphthalate	40.000	38.373	4.1	110	0.00
81	Benzo(a)anthracene	40.000	38.715	3.2	111	0.00
82	3,3'-Dichlorobenzidine	40.000	39.890	0.3	113	0.00
83	Chrysene	40.000	38.524	3.7	112	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	38.056	4.9	109	0.00
85 c	Di-n-octyl phthalate	40.000	38.009	5.0	109	0.00
86 I	Perylene-d12	20.000	20.000	0.0	111	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	38.208	4.5	112	0.00
88	Benzo(b)fluoranthene	40.000	39.188	2.0	112	0.00
89	Benzo(k)fluoranthene	40.000	38.103	4.7	112	0.00
90 C	Benzo(a)pyrene	40.000	38.549	3.6	113	0.00
91	Dibenzo(a,h)anthracene	40.000	38.303	4.2	112	0.00
92	Benzo(g,h,i)perylene	40.000	38.191	4.5	111	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110522\
Data File : BG055467.D
Acq On : 5 Nov 2022 12:28
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Nov 07 00:52:37 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 07 00:51:08 2022
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	Amount	Calc.	%Dev	Area%	Dev(min)
----------	--------	-------	------	-------	----------

(#) = Out of Range SPCC's out = 0 CCC's out = 0