

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030521\
 Data File : BG048339.D
 Acq On : 5 Mar 2021 15:36
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Mar 05 16:10:53 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG021021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 05 15:00:43 2021
 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	54	0.00
2	1,4-Dioxane	40.000	36.495	8.8	49	0.00
3	Pyridine	40.000	41.118	-2.8	51	0.00
4	n-Nitrosodimethylamine	40.000	52.046	-30.1#	71	0.00
5 S	2-Fluorophenol	80.000	79.020	1.2	53	0.00
6	Aniline	40.000	42.402	-6.0	58	0.00
7 S	Phenol-d6	80.000	84.721	-5.9	57	0.00
8	2-Chlorophenol	40.000	41.404	-3.5	55	0.00
9	Benzaldehyde	40.000	34.867	12.8	49	0.00
10 C	Phenol	40.000	43.404	-8.5	59	0.00
11	bis(2-Chloroethyl)ether	40.000	40.426	-1.1	54	0.00
12	1,3-Dichlorobenzene	40.000	39.005	2.5	54	0.00
13 C	1,4-Dichlorobenzene	40.000	40.950	-2.4	56	0.00
14	1,2-Dichlorobenzene	40.000	41.211	-3.0	57	0.00
15	Benzyl Alcohol	40.000	46.411	-16.0	60	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	43.789	-9.5	59	0.00
17	2-Methylphenol	40.000	42.724	-6.8	56	0.00
18	Hexachloroethane	40.000	41.006	-2.5	55	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	46.360	-15.9	62	0.00
20	3+4-Methylphenols	40.000	43.256	-8.1	57	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	57	0.00
22	Acetophenone	40.000	40.783	-2.0	60	0.00
23 S	Nitrobenzene-d5	80.000	85.252	-6.6	62	0.00
24	Nitrobenzene	40.000	43.753	-9.4	63	0.00
25	Isophorone	40.000	42.520	-6.3	60	0.00
26 C	2-Nitrophenol	40.000	40.864	-2.2	58	0.00
27	2,4-Dimethylphenol	40.000	40.014	-0.0	58	0.00
28	bis(2-Chloroethoxy)methane	40.000	39.128	2.2	57	0.00
29 C	2,4-Dichlorophenol	40.000	38.691	3.3	56	0.00
30	1,2,4-Trichlorobenzene	40.000	39.905	0.2	57	0.00
31	Naphthalene	40.000	39.757	0.6	58	0.00
32	Benzoic acid	40.000	27.812	30.5#	38	0.00
33	4-Chloroaniline	40.000	39.876	0.3	57	0.00
34 C	Hexachlorobutadiene	40.000	40.959	-2.4	59	0.00
35	Caprolactam	40.000	46.162	-15.4	64	0.00
36 C	4-Chloro-3-methylphenol	40.000	43.518	-8.8	63	0.00
37	2-Methylnaphthalene	40.000	39.998	0.0	58	0.00
38	1-Methylnaphthalene	40.000	40.465	-1.2	58	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	60	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	39.630	0.9	60	0.00
41 P	Hexachlorocyclopentadiene	40.000	36.924	7.7	53	0.00
42 S	2,4,6-Tribromophenol	80.000	80.852	-1.1	62	0.00
43 C	2,4,6-Trichlorophenol	40.000	38.898	2.8	59	0.00
44	2,4,5-Trichlorophenol	40.000	38.640	3.4	58	0.00
45 S	2-Fluorobiphenyl	80.000	79.709	0.4	61	0.00
46	1,1'-Biphenyl	40.000	39.039	2.4	59	0.00
47	2-Chloronaphthalene	40.000	39.273	1.8	59	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030521\
 Data File : BG048339.D
 Acq On : 5 Mar 2021 15:36
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Mar 05 16:10:53 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG021021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 05 15:00:43 2021
 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	47.347	-18.4	70	0.00
49	Acenaphthylene	40.000	39.490	1.3	60	0.00
50	Dimethylphthalate	40.000	40.375	-0.9	62	0.00
51	2,6-Dinitrotoluene	40.000	39.501	1.2	60	0.00
52 C	Acenaphthene	40.000	35.029	12.4	52	0.00
53	3-Nitroaniline	40.000	39.821	0.4	60	0.00
54 P	2,4-Dinitrophenol	40.000	32.489	18.8	46	0.00
55	Dibenzofuran	40.000	39.697	0.8	61	0.00
56 P	4-Nitrophenol	40.000	48.998	-22.5	72	0.00
57	2,4-Dinitrotoluene	40.000	40.468	-1.2	61	0.00
58	Fluorene	40.000	40.124	-0.3	61	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	35.762	10.6	54	0.00
60	Diethylphthalate	40.000	41.245	-3.1	62	0.00
61	4-Chlorophenyl-phenylether	40.000	40.594	-1.5	62	0.00
62	4-Nitroaniline	40.000	36.675	8.3	56	0.00
63	Azobenzene	40.000	44.069	-10.2	69	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	62	0.00
65	4,6-Dinitro-2-methylphenol	40.000	40.032	-0.1	61	0.00
66 c	n-Nitrosodiphenylamine	40.000	39.928	0.2	62	0.00
67	4-Bromophenyl-phenylether	40.000	39.378	1.6	61	0.00
68	Hexachlorobenzene	40.000	39.600	1.0	61	0.00
69	Atrazine	40.000	40.154	-0.4	62	0.00
70 C	Pentachlorophenol	40.000	42.300	-5.7	61	0.00
71	Phenanthrene	40.000	40.041	-0.1	64	0.00
72	Anthracene	40.000	40.115	-0.3	63	0.00
73	Carbazole	40.000	39.163	2.1	62	0.00
74	Di-n-butylphthalate	40.000	40.602	-1.5	64	0.00
75 C	Fluoranthene	40.000	39.866	0.3	63	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	64	0.00
77	Benzydine	40.000	38.027	4.9	60	0.00
78	Pyrene	40.000	39.289	1.8	64	0.00
79 S	Terphenyl-d14	80.000	78.193	2.3	66	0.00
80	Butylbenzylphthalate	40.000	39.939	0.2	63	0.00
81	Benzo(a)anthracene	40.000	39.262	1.8	64	0.00
82	3,3'-Dichlorobenzidine	40.000	37.489	6.3	60	0.00
83	Chrysene	40.000	38.949	2.6	63	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	39.335	1.7	63	0.00
85 c	Di-n-octyl phthalate	40.000	39.394	1.5	62	0.00
86 I	Perylene-d12	20.000	20.000	0.0	61	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	40.554	-1.4	62	0.00
88	Benzo(b)fluoranthene	40.000	40.690	-1.7	62	0.00
89	Benzo(k)fluoranthene	40.000	40.135	-0.3	61	0.00
90 C	Benzo(a)pyrene	40.000	40.533	-1.3	62	0.00
91	Dibenzo(a,h)anthracene	40.000	40.453	-1.1	61	0.00
92	Benzo(g,h,i)perylene	40.000	40.157	-0.4	62	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030521\
Data File : BG048339.D
Acq On : 5 Mar 2021 15:36
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Mar 05 16:10:53 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG021021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 05 15:00:43 2021
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