

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061221\
 Data File : BG048931.D
 Acq On : 13 Jun 2021 6:22
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 14 03:18:43 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 11 16:03:10 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	103	0.00
2	1,4-Dioxane	40.000	34.106	14.7	102	0.00
3	Pyridine	40.000	36.584	8.5	103	0.00
4	n-Nitrosodimethylamine	40.000	43.223	-8.1	117	0.00
5 S	2-Fluorophenol	80.000	73.597	8.0	107	0.00
6	Aniline	40.000	36.266	9.3	104	0.00
7 S	Phenol-d6	80.000	73.116	8.6	104	0.00
8	2-Chlorophenol	40.000	36.799	8.0	108	0.00
9	Benzaldehyde	40.000	42.712	-6.8	129	0.00
10 C	Phenol	40.000	36.301	9.2	106	0.00
11	bis(2-Chloroethyl)ether	40.000	35.463	11.3	102	0.00
12	1,3-Dichlorobenzene	40.000	36.006	10.0	103	0.00
13 C	1,4-Dichlorobenzene	40.000	36.373	9.1	106	0.00
14	1,2-Dichlorobenzene	40.000	36.225	9.4	107	0.00
15	Benzyl Alcohol	40.000	37.003	7.5	109	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	36.122	9.7	105	-0.01
17	2-Methylphenol	40.000	35.989	10.0	104	0.00
18	Hexachloroethane	40.000	36.225	9.4	104	-0.01
19 P	n-Nitroso-di-n-propylamine	40.000	35.378	11.6	104	0.00
20	3+4-Methylphenols	40.000	36.510	8.7	105	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	105	0.00
22	Acetophenone	40.000	34.531	13.7	105	0.00
23 S	Nitrobenzene-d5	80.000	75.582	5.5	113	0.00
24	Nitrobenzene	40.000	36.290	9.3	110	0.00
25	Isophorone	40.000	34.356	14.1	105	0.00
26 C	2-Nitrophenol	40.000	38.685	3.3	115	0.00
27	2,4-Dimethylphenol	40.000	35.451	11.4	108	0.00
28	bis(2-Chloroethoxy)methane	40.000	34.109	14.7	103	0.00
29 C	2,4-Dichlorophenol	40.000	35.891	10.3	108	0.00
30	1,2,4-Trichlorobenzene	40.000	35.302	11.7	108	0.00
31	Naphthalene	40.000	34.338	14.2	104	0.00
32	Benzoic acid	40.000	38.843	2.9	114	0.01
33	4-Chloroaniline	40.000	34.188	14.5	105	0.00
34 C	Hexachlorobutadiene	40.000	35.666	10.8	109	0.00
35	Caprolactam	40.000	37.896	5.3	115	0.00
36 C	4-Chloro-3-methylphenol	40.000	35.273	11.8	107	0.00
37	2-Methylnaphthalene	40.000	34.447	13.9	106	0.00
38	1-Methylnaphthalene	40.000	34.250	14.4	104	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	108	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	33.854	15.4	108	0.00
41 P	Hexachlorocyclopentadiene	40.000	34.892	12.8	109	0.00
42 S	2,4,6-Tribromophenol	80.000	68.039	15.0	106	0.00
43 C	2,4,6-Trichlorophenol	40.000	35.900	10.3	111	0.00
44	2,4,5-Trichlorophenol	40.000	37.907	5.2	118	0.00
45 S	2-Fluorobiphenyl	80.000	67.291	15.9	105	0.00
46	1,1'-Biphenyl	40.000	33.696	15.8	105	0.00
47	2-Chloronaphthalene	40.000	33.657	15.9	108	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG0611221\
 Data File : BG048931.D
 Acq On : 13 Jun 2021 6:22
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 14 03:18:43 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 11 16:03:10 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	40.101	-0.3	124	0.00
49	Acenaphthylene	40.000	33.667	15.8	106	0.00
50	Dimethylphthalate	40.000	34.202	14.5	108	0.00
51	2,6-Dinitrotoluene	40.000	38.136	4.7	117	0.00
52 C	Acenaphthene	40.000	34.084	14.8	106	0.00
53	3-Nitroaniline	40.000	37.604	6.0	116	0.00
54 P	2,4-Dinitrophenol	40.000	43.192	-8.0	132	0.00
55	Dibenzofuran	40.000	33.244	16.9	105	0.00
56 P	4-Nitrophenol	40.000	35.971	10.1	107	0.00
57	2,4-Dinitrotoluene	40.000	43.026	-7.6	119	0.00
58	Fluorene	40.000	33.146	17.1	105	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	34.823	12.9	107	0.00
60	Diethylphthalate	40.000	33.199	17.0	106	0.00
61	4-Chlorophenyl-phenylether	40.000	33.769	15.6	109	0.00
62	4-Nitroaniline	40.000	36.654	8.4	111	0.00
63	Azobenzene	40.000	32.800	18.0	105	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	106	0.00
65	4,6-Dinitro-2-methylphenol	40.000	45.340	-13.4	127	0.00
66 c	n-Nitrosodiphenylamine	40.000	33.344	16.6	104	0.00
67	4-Bromophenyl-phenylether	40.000	33.339	16.7	103	0.00
68	Hexachlorobenzene	40.000	33.885	15.3	106	0.00
69	Atrazine	40.000	37.548	6.1	108	0.00
70 C	Pentachlorophenol	40.000	37.012	7.5	112	0.00
71	Phenanthrene	40.000	33.586	16.0	105	0.00
72	Anthracene	40.000	33.429	16.4	106	0.00
73	Carbazole	40.000	34.030	14.9	106	0.00
74	Di-n-butylphthalate	40.000	34.878	12.8	110	0.00
75 C	Fluoranthene	40.000	34.509	13.7	110	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	110	-0.01
77	Benzydine	40.000	44.429	-11.1	123	0.00
78	Pyrene	40.000	33.779	15.6	108	0.00
79 S	Terphenyl-d14	80.000	68.009	15.0	109	0.00
80	Butylbenzylphthalate	40.000	35.762	10.6	114	0.00
81	Benzo(a)anthracene	40.000	34.155	14.6	109	0.00
82	3,3'-Dichlorobenzidine	40.000	38.259	4.4	120	0.00
83	Chrysene	40.000	34.305	14.2	110	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	35.365	11.6	114	0.00
85 c	Di-n-octyl phthalate	40.000	35.466	11.3	113	-0.02
86 I	Perylene-d12	20.000	20.000	0.0	112	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	36.817	8.0	115	-0.01
88	Benzo(b)fluoranthene	40.000	35.677	10.8	113	-0.01
89	Benzo(k)fluoranthene	40.000	34.777	13.1	109	0.00
90 C	Benzo(a)pyrene	40.000	35.404	11.5	112	-0.01
91	Dibenzo(a,h)anthracene	40.000	36.344	9.1	113	-0.02
92	Benzo(g,h,i)perylene	40.000	36.246	9.4	114	-0.02

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061221\
Data File : BG048931.D
Acq On : 13 Jun 2021 6:22
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Jun 14 03:18:43 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 11 16:03:10 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