

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041921\
 Data File : BG048760.D
 Acq On : 19 Apr 2021 10:35
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Apr 19 11:25:39 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG041421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 16 15:59:00 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	98	0.00
2	1,4-Dioxane	40.000	39.877	0.3	96	0.00
3	Pyridine	40.000	43.042	-7.6	100	0.00
4	n-Nitrosodimethylamine	40.000	41.491	-3.7	98	0.00
5 S	2-Fluorophenol	80.000	81.582	-2.0	101	0.00
6	Aniline	40.000	41.337	-3.3	97	0.01
7 S	Phenol-d6	80.000	82.445	-3.1	100	0.01
8	2-Chlorophenol	40.000	40.851	-2.1	101	0.00
9	Benzaldehyde	40.000	37.268	6.8	94	0.00
10 C	Phenol	40.000	40.978	-2.4	101	0.00
11	bis(2-Chloroethyl)ether	40.000	40.864	-2.2	102	0.01
12	1,3-Dichlorobenzene	40.000	39.086	2.3	98	0.00
13 C	1,4-Dichlorobenzene	40.000	40.086	-0.2	98	0.00
14	1,2-Dichlorobenzene	40.000	39.884	0.3	96	0.00
15	Benzyl Alcohol	40.000	41.165	-2.9	97	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	41.920	-4.8	100	0.01
17	2-Methylphenol	40.000	40.499	-1.2	97	0.01
18	Hexachloroethane	40.000	39.755	0.6	96	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	41.642	-4.1	102	0.00
20	3+4-Methylphenols	40.000	41.463	-3.7	99	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	98	0.00
22	Acetophenone	40.000	40.373	-0.9	103	0.00
23 S	Nitrobenzene-d5	80.000	79.966	0.0	99	0.00
24	Nitrobenzene	40.000	39.356	1.6	97	0.00
25	Isophorone	40.000	40.573	-1.4	100	0.00
26 C	2-Nitrophenol	40.000	40.250	-0.6	98	0.01
27	2,4-Dimethylphenol	40.000	39.825	0.4	100	0.00
28	bis(2-Chloroethoxy)methane	40.000	41.168	-2.9	102	0.00
29 C	2,4-Dichlorophenol	40.000	39.398	1.5	97	0.01
30	1,2,4-Trichlorobenzene	40.000	39.108	2.2	100	0.00
31	Naphthalene	40.000	39.916	0.2	98	0.01
32	Benzoic acid	40.000	21.784	45.5#	53	0.00
33	4-Chloroaniline	40.000	39.221	1.9	95	0.01
34 C	Hexachlorobutadiene	40.000	38.182	4.5	91	0.01
35	Caprolactam	40.000	39.683	0.8	99	0.00
36 C	4-Chloro-3-methylphenol	40.000	39.760	0.6	101	0.01
37	2-Methylnaphthalene	40.000	39.099	2.3	98	0.00
38	1-Methylnaphthalene	40.000	39.203	2.0	98	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	96	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	41.305	-3.3	96	0.00
41 P	Hexachlorocyclopentadiene	40.000	41.180	-2.9	100	0.00
42 S	2,4,6-Tribromophenol	80.000	77.631	3.0	91	0.00
43 C	2,4,6-Trichlorophenol	40.000	40.128	-0.3	92	0.01
44	2,4,5-Trichlorophenol	40.000	38.430	3.9	89	0.00
45 S	2-Fluorobiphenyl	80.000	83.145	-3.9	98	0.00
46	1,1'-Biphenyl	40.000	41.452	-3.6	97	0.00
47	2-Chloronaphthalene	40.000	41.406	-3.5	98	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041921\
 Data File : BG048760.D
 Acq On : 19 Apr 2021 10:35
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Apr 19 11:25:39 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG041421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 16 15:59:00 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	42.792	-7.0	94	0.00
49	Acenaphthylene	40.000	41.131	-2.8	97	0.00
50	Dimethylphthalate	40.000	40.659	-1.6	95	0.00
51	2,6-Dinitrotoluene	40.000	39.822	0.4	91	0.00
52 C	Acenaphthene	40.000	40.991	-2.5	94	0.00
53	3-Nitroaniline	40.000	42.482	-6.2	102	0.00
54 P	2,4-Dinitrophenol	40.000	47.058	-17.6	112	0.00
55	Dibenzofuran	40.000	40.994	-2.5	95	0.00
56 P	4-Nitrophenol	40.000	38.186	4.5	86	0.01
57	2,4-Dinitrotoluene	40.000	40.451	-1.1	94	0.00
58	Fluorene	40.000	40.124	-0.3	94	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	37.192	7.0	85	0.00
60	Diethylphthalate	40.000	40.606	-1.5	96	0.00
61	4-Chlorophenyl-phenylether	40.000	40.884	-2.2	97	0.00
62	4-Nitroaniline	40.000	40.735	-1.8	99	0.00
63	Azobenzene	40.000	41.740	-4.4	99	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	95	0.00
65	4,6-Dinitro-2-methylphenol	40.000	37.980	5.1	85	0.00
66 c	n-Nitrosodiphenylamine	40.000	40.537	-1.3	97	0.00
67	4-Bromophenyl-phenylether	40.000	37.817	5.5	96	0.00
68	Hexachlorobenzene	40.000	39.745	0.6	95	0.00
69	Atrazine	40.000	40.908	-2.3	100	0.00
70 C	Pentachlorophenol	40.000	36.171	9.6	88	0.00
71	Phenanthrene	40.000	40.125	-0.3	97	0.00
72	Anthracene	40.000	40.510	-1.3	98	0.00
73	Carbazole	40.000	39.052	2.4	95	0.00
74	Di-n-butylphthalate	40.000	40.472	-1.2	99	0.00
75 C	Fluoranthene	40.000	38.922	2.7	95	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	99	0.00
77	Benzydine	40.000	44.080	-10.2	89	0.00
78	Pyrene	40.000	40.324	-0.8	97	0.00
79 S	Terphenyl-d14	80.000	76.853	3.9	95	0.00
80	Butylbenzylphthalate	40.000	39.788	0.5	97	0.00
81	Benzo(a)anthracene	40.000	39.171	2.1	96	0.00
82	3,3'-Dichlorobenzidine	40.000	37.151	7.1	91	0.00
83	Chrysene	40.000	38.714	3.2	94	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	39.619	1.0	96	0.00
85 c	Di-n-octyl phthalate	40.000	39.739	0.7	96	0.00
86 I	Perylene-d12	20.000	20.000	0.0	94	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	39.730	0.7	94	0.00
88	Benzo(b)fluoranthene	40.000	39.756	0.6	93	0.00
89	Benzo(k)fluoranthene	40.000	40.481	-1.2	96	0.00
90 C	Benzo(a)pyrene	40.000	40.680	-1.7	96	0.00
91	Dibenzo(a,h)anthracene	40.000	39.931	0.2	96	0.00
92	Benzo(g,h,i)perylene	40.000	38.363	4.1	92	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041921\
Data File : BG048760.D
Acq On : 19 Apr 2021 10:35
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Apr 19 11:25:39 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG041421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 16 15:59:00 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