

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061521\
 Data File : BG048974.D
 Acq On : 15 Jun 2021 12:15
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 15 12:50:44 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	87	0.00
2	1,4-Dioxane	40.000	36.082	9.8	90	0.00
3	Pyridine	40.000	39.516	1.2	93	0.00
4	n-Nitrosodimethylamine	40.000	45.073	-12.7	103	0.00
5 S	2-Fluorophenol	80.000	76.007	5.0	93	0.00
6	Aniline	40.000	37.079	7.3	90	0.00
7 S	Phenol-d6	80.000	75.367	5.8	90	0.00
8	2-Chlorophenol	40.000	36.791	8.0	91	0.00
9	Benzaldehyde	40.000	34.018	15.0	87	0.00
10 C	Phenol	40.000	37.021	7.4	91	0.00
11	bis(2-Chloroethyl)ether	40.000	35.239	11.9	85	0.00
12	1,3-Dichlorobenzene	40.000	37.529	6.2	90	0.00
13 C	1,4-Dichlorobenzene	40.000	37.909	5.2	93	0.00
14	1,2-Dichlorobenzene	40.000	37.334	6.7	92	0.00
15	Benzyl Alcohol	40.000	36.092	9.8	89	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	36.752	8.1	90	0.00
17	2-Methylphenol	40.000	37.211	7.0	90	0.00
18	Hexachloroethane	40.000	36.646	8.4	89	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	35.440	11.4	88	0.00
20	3+4-Methylphenols	40.000	36.561	8.6	88	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	89	0.00
22	Acetophenone	40.000	35.044	12.4	91	0.00
23 S	Nitrobenzene-d5	80.000	71.472	10.7	91	0.00
24	Nitrobenzene	40.000	35.818	10.5	92	0.00
25	Isophorone	40.000	33.903	15.2	88	0.00
26 C	2-Nitrophenol	40.000	34.386	14.0	87	0.00
27	2,4-Dimethylphenol	40.000	33.571	16.1	87	0.00
28	bis(2-Chloroethoxy)methane	40.000	33.558	16.1	86	0.00
29 C	2,4-Dichlorophenol	40.000	34.550	13.6	88	0.00
30	1,2,4-Trichlorobenzene	40.000	35.035	12.4	91	0.00
31	Naphthalene	40.000	34.384	14.0	88	0.00
32	Benzoic acid	40.000	34.940	12.7	87	0.00
33	4-Chloroaniline	40.000	33.444	16.4	87	0.00
34 C	Hexachlorobutadiene	40.000	35.153	12.1	91	0.00
35	Caprolactam	40.000	39.350	1.6	101	0.00
36 C	4-Chloro-3-methylphenol	40.000	33.634	15.9	87	0.00
37	2-Methylnaphthalene	40.000	33.641	15.9	88	0.00
38	1-Methylnaphthalene	40.000	34.244	14.4	88	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	86	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	35.029	12.4	89	0.00
41 P	Hexachlorocyclopentadiene	40.000	34.384	14.0	86	0.00
42 S	2,4,6-Tribromophenol	80.000	68.160	14.8	85	0.00
43 C	2,4,6-Trichlorophenol	40.000	36.701	8.2	91	0.00
44	2,4,5-Trichlorophenol	40.000	39.422	1.4	98	0.00
45 S	2-Fluorobiphenyl	80.000	70.358	12.1	88	0.00
46	1,1'-Biphenyl	40.000	34.718	13.2	87	0.00
47	2-Chloronaphthalene	40.000	35.448	11.4	90	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061521\
 Data File : BG048974.D
 Acq On : 15 Jun 2021 12:15
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 15 12:50:44 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	36.253	9.4	87	0.00
49	Acenaphthylene	40.000	34.893	12.8	88	0.00
50	Dimethylphthalate	40.000	34.809	13.0	87	0.00
51	2,6-Dinitrotoluene	40.000	35.528	11.2	87	0.00
52 C	Acenaphthene	40.000	34.661	13.3	86	0.00
53	3-Nitroaniline	40.000	36.394	9.0	89	0.00
54 P	2,4-Dinitrophenol	40.000	36.845	7.9	90	0.00
55	Dibenzofuran	40.000	34.552	13.6	87	0.00
56 P	4-Nitrophenol	40.000	36.135	9.7	86	0.00
57	2,4-Dinitrotoluene	40.000	40.874	-2.2	90	0.00
58	Fluorene	40.000	34.640	13.4	88	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	33.947	15.1	83	0.00
60	Diethylphthalate	40.000	33.819	15.5	86	0.00
61	4-Chlorophenyl-phenylether	40.000	34.608	13.5	89	0.00
62	4-Nitroaniline	40.000	38.759	3.1	94	0.00
63	Azobenzene	40.000	33.660	15.9	86	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	87	0.00
65	4,6-Dinitro-2-methylphenol	40.000	37.939	5.2	87	0.00
66 c	n-Nitrosodiphenylamine	40.000	33.836	15.4	86	0.00
67	4-Bromophenyl-phenylether	40.000	33.240	16.9	84	0.00
68	Hexachlorobenzene	40.000	33.675	15.8	87	0.00
69	Atrazine	40.000	35.782	10.5	84	0.00
70 C	Pentachlorophenol	40.000	33.358	16.6	83	0.00
71	Phenanthrene	40.000	33.596	16.0	86	0.00
72	Anthracene	40.000	33.870	15.3	88	0.00
73	Carbazole	40.000	34.281	14.3	88	0.00
74	Di-n-butylphthalate	40.000	34.950	12.6	91	0.00
75 C	Fluoranthene	40.000	35.256	11.9	92	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	92	0.00
77	Benzydine	40.000	42.658	-6.6	99	0.00
78	Pyrene	40.000	33.521	16.2	90	0.00
79 S	Terphenyl-d14	80.000	71.579	10.5	96	0.00
80	Butylbenzylphthalate	40.000	34.571	13.6	92	0.00
81	Benzo(a)anthracene	40.000	34.106	14.7	91	0.00
82	3,3'-Dichlorobenzidine	40.000	37.820	5.4	100	0.00
83	Chrysene	40.000	34.126	14.7	92	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	35.677	10.8	97	0.00
85 c	Di-n-octyl phthalate	40.000	36.186	9.5	97	0.01
86 I	Perylene-d12	20.000	20.000	0.0	96	0.01
87	Indeno(1,2,3-cd)pyrene	40.000	37.383	6.5	100	0.01
88	Benzo(b)fluoranthene	40.000	36.086	9.8	98	0.01
89	Benzo(k)fluoranthene	40.000	34.763	13.1	92	0.01
90 C	Benzo(a)pyrene	40.000	36.004	10.0	97	0.01
91	Dibenzo(a,h)anthracene	40.000	37.549	6.1	99	0.02
92	Benzo(g,h,i)perylene	40.000	36.834	7.9	98	0.02

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061521\
Data File : BG048974.D
Acq On : 15 Jun 2021 12:15
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Jun 15 12:50:44 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