

Data Path : Z:\HPCHEM1\BNA G\Data\BG020918\
 Data File : BG032631.D
 Acq On : 9 Feb 2018 14:03
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Feb 09 23:30:49 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 18:57:08 2018
 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	137	-0.02
2	1,4-Dioxane	40.000	43.242	-8.1	162	-0.01
3	Pyridine	40.000	41.775	-4.4	140	-0.01
4	n-Nitrosodimethylamine	40.000	41.268	-3.2	133	-0.01
5 S	2-Fluorophenol	80.000	73.058	8.7	122	-0.01
6	Aniline	40.000	38.597	3.5	126	-0.02
7 S	Phenol-d6	80.000	81.979	-2.5	139	-0.01
8	2-Chlorophenol	40.000	37.570	6.1	125	-0.02
9	Benzaldehyde	40.000	28.559	28.6#	98	-0.01
10 C	Phenol	40.000	40.430	-1.1	134	-0.02
11	bis(2-Chloroethyl)ether	40.000	44.414	-11.0	150	-0.02
12	1,3-Dichlorobenzene	40.000	36.912	7.7	128	-0.02
13 C	1,4-Dichlorobenzene	40.000	38.354	4.1	132	-0.02
14	1,2-Dichlorobenzene	40.000	38.197	4.5	129	-0.02
15	Benzyl Alcohol	40.000	37.186	7.0	122	-0.02
16	2,2'-oxybis(1-Chloropropane)	40.000	46.709	-16.8	157	-0.02
17	2-Methylphenol	40.000	39.149	2.1	130	-0.02
18	Hexachloroethane	40.000	38.095	4.8	133	-0.02
19 P	n-Nitroso-di-n-propylamine	40.000	42.426	-6.1	139	-0.01
20	3+4-Methylphenols	40.000	39.792	0.5	128	-0.02
21 I	Naphthalene-d8	20.000	20.000	0.0	134	-0.02
22	Acetophenone	40.000	36.972	7.6	126	-0.02
23 S	Nitrobenzene-d5	80.000	81.314	-1.6	135	-0.02
24	Nitrobenzene	40.000	40.976	-2.4	139	-0.02
25	Isophorone	40.000	42.755	-6.9	145	-0.02
26 C	2-Nitrophenol	40.000	37.278	6.8	120	-0.03
27	2,4-Dimethylphenol	40.000	35.583	11.0	119	-0.02
28	bis(2-Chloroethoxy)methane	40.000	39.844	0.4	134	-0.02
29 C	2,4-Dichlorophenol	40.000	37.030	7.4	125	-0.02
30	1,2,4-Trichlorobenzene	40.000	37.219	7.0	129	-0.03
31	Naphthalene	40.000	37.835	5.4	129	-0.02
32	Benzoic acid	40.000	38.504	3.7	133	-0.03
33	4-Chloroaniline	40.000	36.260	9.4	124	-0.02
34 C	Hexachlorobutadiene	40.000	38.543	3.6	132	-0.03
35	Caprolactam	40.000	36.939	7.7	122	-0.02
36 C	4-Chloro-3-methylphenol	40.000	38.433	3.9	132	-0.03
37	2-Methylnaphthalene	40.000	37.715	5.7	128	-0.03
38 I	Acenaphthene-d10	20.000	20.000	0.0	139	-0.02
39	1,2,4,5-Tetrachlorobenzene	40.000	37.336	6.7	134	-0.02
40 P	Hexachlorocyclopentadiene	40.000	36.512	8.7	129	-0.02
41 S	2,4,6-Tribromophenol	80.000	57.148	28.6#	102	-0.01
42 C	2,4,6-Trichlorophenol	40.000	34.806	13.0	126	-0.02
43	2,4,5-Trichlorophenol	40.000	36.842	7.9	131	-0.03
44 S	2-Fluorobiphenyl	80.000	71.634	10.5	129	-0.02

Data Path : Z:\HPCHEM1\BNA G\Data\BG020918\
 Data File : BG032631.D
 Acq On : 9 Feb 2018 14:03
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Feb 09 23:30:49 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 18:57:08 2018
 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)
45	1,1'-Biphenyl	40.000	36.879	7.8	132	-0.02
46	2-Chloronaphthalene	40.000	35.956	10.1	129	-0.02
47	2-Nitroaniline	40.000	36.950	7.6	131	-0.02
48	Acenaphthylene	40.000	36.399	9.0	131	-0.02
49	Dimethylphthalate	40.000	34.778	13.1	128	-0.01
50	2,6-Dinitrotoluene	40.000	34.810	13.0	127	-0.01
51 C	Acenaphthene	40.000	35.153	12.1	125	-0.02
52	3-Nitroaniline	40.000	32.053	19.9	114	-0.01
53 P	2,4-Dinitrophenol	40.000	25.501	36.2#	90	-0.01
54	Dibenzofuran	40.000	35.879	10.3	131	-0.02
55 P	4-Nitrophenol	40.000	24.207	39.5#	87	-0.01
56	2,4-Dinitrotoluene	40.000	32.698	18.3	117	-0.01
57	Fluorene	40.000	34.561	13.6	128	-0.02
58	2,3,4,6-Tetrachlorophenol	40.000	32.464	18.8	115	-0.02
59	Diethylphthalate	40.000	33.493	16.3	124	-0.01
60	4-Chlorophenyl-phenylether	40.000	36.481	8.8	132	-0.01
61	4-Nitroaniline	40.000	28.843	27.9#	107	-0.01
62	Azobenzene	40.000	37.029	7.4	137	-0.01
63 I	Phenanthrene-d10	20.000	20.000	0.0	125	-0.01
64	4,6-Dinitro-2-methylphenol	40.000	34.094	14.8	105	-0.01
65 c	n-Nitrosodiphenylamine	40.000	38.953	2.6	125	-0.01
66	4-Bromophenyl-phenylether	40.000	38.635	3.4	123	-0.01
67	Hexachlorobenzene	40.000	36.914	7.7	119	-0.01
68	Atrazine	40.000	24.421	38.9#	76	0.00
69 C	Pentachlorophenol	40.000	32.217	19.5	103	-0.01
70	Phenanthrene	40.000	35.291	11.8	114	0.00
71	Anthracene	40.000	36.295	9.3	115	-0.01
72	Carbazole	40.000	32.527	18.7	104	0.00
73	Di-n-butylphthalate	40.000	32.860	17.9	104	0.00
74 C	Fluoranthene	40.000	32.601	18.5	107	0.00
75 I	Chrysene-d12	20.000	20.000	0.0	121	0.00
76	Benzidine	40.000	13.126	67.2#	36	0.00
77	Pyrene	40.000	35.318	11.7	109	0.00
78 S	Terphenyl-d14	80.000	68.661	14.2	107	0.00
79	Butylbenzylphthalate	40.000	38.172	4.6	114	0.00
80	Benzo(a)anthracene	40.000	37.001	7.5	114	0.00
81	3,3'-Dichlorobenzidine	40.000	33.475	16.3	101	0.00
82	Chrysene	40.000	37.118	7.2	115	0.00
83	Bis(2-ethylhexyl)phthalate	40.000	38.467	3.8	115	0.00
84 c	Di-n-octyl phthalate	40.000	39.118	2.2	117	0.00
85	Indeno(1,2,3-cd)pyrene	40.000	35.865	10.3	109	0.00
86 I	Perylene-d12	20.000	20.000	0.0	118	0.00
87	Benzo(b)fluoranthene	40.000	36.549	8.6	110	0.00

Data Path : Z:\HPCHEM1\BNA G\Data\BG020918\
Data File : BG032631.D
Acq On : 9 Feb 2018 14:03
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Feb 09 23:30:49 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 06 18:57:08 2018
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)
88	Benzo(k)fluoranthene	40.000	38.534	3.7	117	0.00
89 C	Benzo(a)pyrene	40.000	37.622	5.9	114	0.00
90	Dibenzo(a,h)anthracene	40.000	35.797	10.5	107	0.00
91	Benzo(a,h,i)perylene	40.000	36.105	9.7	108	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0