Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062018\

Evaluate Continuing Calibration Report

Data File : BG035278.D
Acq On : 21 Jun 2018 00:07
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 21 05:53:53 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG060718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 08 17:16:28 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	192	-0.02
2		1,4-Dioxane	40.000	38.964	2.6	188	0.00
3		Pyridine	40.000	39.681	0.8	179	0.00
4		n-Nitrosodimethylamine	40.000	33.094	17.3	150	0.00
5	S	2-Fluorophenol	80.000	78.760	1.5	185	0.00
6		Aniline	40.000	37.469	6.3	177	-0.02
7	S	Phenol-d6	80.000	77.819	2.7	181	0.00
8		2-Chlorophenol	40.000	38.101	4.7	178	-0.02
9		Benzaldehyde	40.000	32.052	19.9	146	-0.02
10	C	Phenol	40.000	37.597	6.0	177	-0.02
11		bis(2-Chloroethyl)ether	40.000	37.360	6.6	177	-0.02
12		1,3-Dichlorobenzene	40.000	38.373	4.1	187	-0.02
13	C	1,4-Dichlorobenzene	40.000	38.255	4.4	182	-0.02
14		1,2-Dichlorobenzene	40.000	37.935	5.2	179	-0.02
15		Benzyl Alcohol	40.000	35.374	11.6	166	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	34.664	13.3	166	-0.02
17		2-Methylphenol	40.000	39.271	1.8	180	0.00
18		Hexachloroethane	40.000	37.012	7.5	171	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	35.405	11.5	166	-0.02
20		3+4-Methylphenols	40.000	38.167	4.6	179	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	198	-0.02
22		Acetophenone	40.000	36.277	9.3	178	-0.02
23	S	Nitrobenzene-d5	80.000	80.977	-1.2	190	-0.02
24		Nitrobenzene	40.000	38.516	3.7	182	-0.02
25		Isophorone	40.000	34.397	14.0	168	-0.02
26	C	2-Nitrophenol	40.000	48.187	-20.5#	236	-0.02
27		2,4-Dimethylphenol	40.000	35.716	10.7	177	-0.02
28		bis(2-Chloroethoxy)methane	40.000	36.788	8.0	180	-0.02
29	C	2,4-Dichlorophenol	40.000	39.195	2.0	189	-0.02
30		1,2,4-Trichlorobenzene	40.000	38.396	4.0	183	-0.02
31		Naphthalene	40.000	36.902	7.7	181	-0.02
32		Benzoic acid	40.000	41.522	-3.8	203	0.00
33		4-Chloroaniline	40.000	37.153	7.1	179	-0.02
34	C	Hexachlorobutadiene	40.000	38.234	4.4	187	-0.02
35		Caprolactam	40.000	37.007	7.5	183	0.00
36	C	4-Chloro-3-methylphenol	40.000	38.356	4.1	186	0.00
37		2-Methylnaphthalene	40.000	37.505	6.2	181	-0.02
38	I	Acenaphthene-d10	20.000	20.000	0.0	198	-0.02
39		1,2,4,5-Tetrachlorobenzene	40.000	40.429	-1.1	194	-0.02
40	P	Hexachlorocyclopentadiene	40.000	39.497	1.3	185	-0.03
41	S	2,4,6-Tribromophenol	80.000	82.892	-3.6	201	-0.02
42	C	2,4,6-Trichlorophenol	40.000	41.459	-3.6	196	-0.02
43		2,4,5-Trichlorophenol	40.000	41.334	-3.3	203	-0.02
44	S	2-Fluorobiphenyl	80.000	73.725	7.8	181	-0.02

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062018\

Evaluate Continuing Calibration Report

Data File : BG035278.D
Acq On : 21 Jun 2018 00:07
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 21 05:53:53 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG060718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 08 17:16:28 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	37.613	6.0	183	-0.02
46		2-Chloronaphthalene	40.000	38.529	3.7	191	-0.02
47		2-Nitroaniline	40.000	41.894	-4.7	204	-0.02
48		Acenaphthylene	40.000	37.313	6.7	182	-0.02
49		Dimethylphthalate	40.000	36.328	9.2	179	-0.02
50		2,6-Dinitrotoluene	40.000	43.317	-8.3	206	-0.02
51	C	Acenaphthene	40.000	38.604	3.5	187	-0.02
52		3-Nitroaniline	40.000	44.064	-10.2	203	0.00
53	P	2,4-Dinitrophenol	40.000	53.900	-34.8#	256	-0.02
54		Dibenzofuran	40.000	36.948	7.6	179	-0.02
55	P	4-Nitrophenol	40.000	38.836	2.9	179	0.00
56		2,4-Dinitrotoluene	40.000	45.028	-12.6	214	-0.02
57		Fluorene	40.000	36.397	9.0	176	-0.02
58		2,3,4,6-Tetrachlorophenol	40.000	41.344	-3.4	199	-0.02
59		Diethylphthalate	40.000	35.194	12.0	174	-0.03
60		4-Chlorophenyl-phenylether	40.000	38.298	4.3	189	-0.03
61		4-Nitroaniline	40.000	41.343	-3.4	192	-0.02
62		Azobenzene	40.000	33.498	16.3	165	-0.02
63	I	Phenanthrene-d10	20.000	20.000	0.0	197	-0.02
64		4,6-Dinitro-2-methylphenol	40.000	51.575	-28.9#	250	-0.02
65	c	n-Nitrosodiphenylamine	40.000	37.038	7.4	176	-0.02
66		4-Bromophenyl-phenylether	40.000	39.498	1.3	193	-0.02
67		Hexachlorobenzene	40.000	39.739	0.7	195	-0.02
68		Atrazine	40.000	33.435	16.4	160	-0.03
69	C	Pentachlorophenol	40.000	40.807	-2.0	193	-0.02
70		Phenanthrene	40.000	36.835	7.9	177	-0.02
71		Anthracene	40.000	36.925	7.7	176	-0.02
72		Carbazole	40.000	35.915	10.2	172	-0.02
73		Di-n-butylphthalate	40.000	34.888	12.8	167	-0.03
74	C	Fluoranthene	40.000	36.341	9.1	174	-0.02
75	I	Chrysene-d12	20.000	20.000	0.0	197	-0.02
76		Benzidine	40.000	19.274	51.8#	93	-0.02
77		Pyrene	40.000	35.358	11.6	173	-0.02
78	S	Terphenyl-d14	80.000	70.529	11.8	171	-0.02
79		Butylbenzylphthalate	40.000	35.861	10.3	169	-0.03
80		Benzo(a)anthracene	40.000	36.356	9.1	178	-0.02
81		3,3'-Dichlorobenzidine	40.000	36.283	9.3	172	-0.03
82		Chrysene	40.000	36.910	7.7	181	-0.03
83		Bis(2-ethylhexyl)phthalate	40.000	34.912	12.7	167	-0.04
84	c	Di-n-octyl phthalate	40.000	34.823	12.9	166	-0.06
85		Indeno(1,2,3-cd)pyrene	40.000	37.627	5.9	182	-0.07
86	I	Perylene-d12	20.000	20.000	0.0	195	-0.04
87		Benzo(b)fluoranthene	40.000	36.460	8.8	173	-0.03

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062018\

Evaluate Continuing Calibration Report

Data File : BG035278.D
Acq On : 21 Jun 2018 00:07
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 21 05:53:53 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG060718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 08 17:16:28 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	37.453	6.4	186	-0.03
89 C	Benzo(a)pyrene	40.000	37.117	7.2	179	-0.04
90	Dibenzo(a,h)anthracene	40.000	37.216	7.0	181	-0.07
91	Benzo(g,h,i)perylene	40.000	37.449	6.4	181	-0.07

(#) = Out of Range SPCC's out = 0 CCC's out = 1