Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061419\
Data File : BG041254.D
Acq On : 14 Jun 2019 15:31
Operator : HP/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG061419

Quant Time: Jun 14 16:11:22 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 14 15:26:06 2019
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	169	-0.02
2		1,4-Dioxane	40.000	38.949	2.6	156	-0.02
3		Pyridine	40.000	40.186	-0.5	162	-0.02
4		n-Nitrosodimethylamine	40.000	41.880	-4.7	167	-0.02
5	S	2-Fluorophenol	80.000	80.184	-0.2	161	-0.01
6		Aniline	40.000	41.831	-4.6	168	-0.01
7	S	Phenol-d6	80.000	84.120	-5.2	168	0.00
8		2-Chlorophenol	40.000	40.588	-1.5	167	-0.02
9		Benzaldehyde	40.000	42.456	-6.1	181	-0.01
10	C	Phenol	40.000	40.990	-2.5	163	0.00
11		bis(2-Chloroethyl)ether	40.000	41.132	-2.8	169	-0.01
12		1,3-Dichlorobenzene	40.000	40.739	-1.8	166	-0.01
13	C	1,4-Dichlorobenzene	40.000	40.904	-2.3	167	-0.01
14		1,2-Dichlorobenzene	40.000	40.588	-1.5	161	-0.02
15		Benzyl Alcohol	40.000	41.842	-4.6	167	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.636	-4.1	168	-0.01
17		2-Methylphenol	40.000	41.164	-2.9	164	0.00
18		Hexachloroethane	40.000	41.064	-2.7	165	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	42.492	-6.2	172	-0.01
20		3+4-Methylphenols	40.000	41.499	-3.7	166	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	168	-0.01
22		Acetophenone	40.000	40.775	-1.9	168	-0.01
23	S	Nitrobenzene-d5	80.000	80.638	-0.8	165	0.00
24		Nitrobenzene	40.000	40.800	-2.0	166	-0.01
25		Isophorone	40.000	40.940	-2.3	167	0.00
26	C	2-Nitrophenol	40.000	41.134	-2.8	166	-0.02
27		2,4-Dimethylphenol	40.000	40.660	-1.6	165	-0.01
28		bis(2-Chloroethoxy)methane	40.000	41.267	-3.2	170	-0.01
29	C	2,4-Dichlorophenol	40.000	40.320	-0.8	167	-0.01
30		1,2,4-Trichlorobenzene	40.000	41.079	-2.7	171	-0.01
31		Naphthalene	40.000	40.800	-2.0	167	-0.01
32		Benzoic acid	40.000	53.694	-34.2#	399	0.01
33		4-Chloroaniline	40.000	41.407	-3.5	170	-0.01
34	C	Hexachlorobutadiene	40.000	40.478	-1.2	166	-0.01
35		Caprolactam	40.000	42.303	-5.8	173	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.830	-4.6	169	0.00
37		2-Methylnaphthalene	40.000	40.696	-1.7	164	0.00
38		1-Methylnaphthalene	40.000	40.671	-1.7	164	-0.01
39	I	Acenaphthene-d10	20.000	20.000	0.0	167	-0.01
40		1,2,4,5-Tetrachlorobenzene	40.000	39.898	0.3	165	-0.01
41	P	Hexachlorocyclopentadiene	40.000	44.129	-10.3	198	-0.01
42	S	2,4,6-Tribromophenol	80.000	85.582	-7.0	173	0.00
43	C	2,4,6-Trichlorophenol	40.000	41.631	-4.1	169	0.00
44		2,4,5-Trichlorophenol	40.000	41.599	-4.0	168	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061419\
Data File : BG041254.D
Acq On : 14 Jun 2019 15:31
Operator : HP/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG061419

Quant Time: Jun 14 16:11:22 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 14 15:26:06 2019
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	S	2-Fluorobiphenyl	80.000	78.295	2.1	161	0.00
46		1,1'-Biphenyl	40.000	40.501	-1.3	168	0.00
47		2-Chloronaphthalene	40.000	39.891	0.3	162	-0.01
48		2-Nitroaniline	40.000	40.039	-0.1	162	0.00
49		Acenaphthylene	40.000	39.580	1.1	161	0.00
50		Dimethylphthalate	40.000	40.596	-1.5	164	0.00
51		2,6-Dinitrotoluene	40.000	40.887	-2.2	170	0.00
52	C	Acenaphthene	40.000	39.908	0.2	164	-0.01
53		3-Nitroaniline	40.000	42.372	-5.9	173	0.00
54	P	2,4-Dinitrophenol	40.000	55.962	-39.9#	382	0.00
55		Dibenzofuran	40.000	40.128	-0.3	163	0.00
56	P	4-Nitrophenol	40.000	42.830	-7.1	196	0.00
57		2,4-Dinitrotoluene	40.000	40.957	-2.4	167	0.00
58		Fluorene	40.000	40.358	-0.9	164	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	43.900	-9.7	171	0.00
60		Diethylphthalate	40.000	41.146	-2.9	168	-0.01
61		4-Chlorophenyl-phenylether	40.000	40.647	-1.6	168	-0.01
62		4-Nitroaniline	40.000	42.038	-5.1	175	0.00
63		Azobenzene	40.000	40.395	-1.0	164	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	166	-0.01
65		4,6-Dinitro-2-methylphenol	40.000	49.921	-24.8	249	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.448	-1.1	167	0.00
67		4-Bromophenyl-phenylether	40.000	41.297	-3.2	172	0.00
68		Hexachlorobenzene	40.000	40.136	-0.3	168	-0.01
69		Atrazine	40.000	42.365	-5.9	166	0.00
70	C	Pentachlorophenol	40.000	45.976	-14.9	216	0.00
71		Phenanthrene	40.000	39.268	1.8	161	0.00
72		Anthracene	40.000	39.739	0.7	161	0.00
73		Carbazole	40.000	39.854	0.4	164	0.00
74		Di-n-butylphthalate	40.000	40.159	-0.4	164	-0.01
75	C	Fluoranthene	40.000	39.464	1.3	160	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	167	0.00
77		Benzidine	40.000	46.468	-16.2	190	0.00
78		Pyrene	40.000	40.016	-0.0	161	-0.01
79	S	Terphenyl-d14	80.000	78.807	1.5	156	0.00
80		Butylbenzylphthalate	40.000	40.731	-1.8	166	-0.01
81		Benzo(a)anthracene	40.000	41.121	-2.8	164	0.00
82		3,3'-Dichlorobenzidine	40.000	44.477	-11.2	178	0.00
83		Chrysene	40.000	41.307	-3.3	163	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	40.397	-1.0	163	-0.01
85	c	Di-n-octyl phthalate	40.000	40.919	-2.3	163	-0.02
86		Indeno(1,2,3-cd)pyrene	40.000	44.430	-11.1	182	-0.03
87	I	Perylene-d12	20.000	20.000	0.0	159	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061419\
Data File : BG041254.D
Acq On : 14 Jun 2019 15:31
Operator : HP/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG061419

Quant Time: Jun 14 16:11:22 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 14 15:26:06 2019
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(b)fluoranthene	40.000	41.139	-2.8	171	0.00
89	Benzo(k)fluoranthene	40.000	39.528	1.2	170	-0.02
90 C	Benzo(a)pyrene	40.000	41.333	-3.3	168	-0.01
91	Dibenzo(a,h)anthracene	40.000	43.712	-9.3	180	-0.02
92	Benzo(g,h,i)perylene	40.000	47.511	-18.8	215	-0.04

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