Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061319\
Data File : BG041230.D
Acq On : 13 Jun 2019 19:31
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 14 09:03:01 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Jun 09 23:08:34 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	73	0.00
2		1,4-Dioxane	0.503	0.434	13.7	63	0.00
3		Pyridine	1.489	1.314	11.8	64	0.00
4		n-Nitrosodimethylamine	0.691	0.707	-2.3	73	0.00
5	S	2-Fluorophenol	1.109	1.095	1.3	71	0.00
6		Aniline	2.286	2.219	2.9	71	0.00
7	S	Phenol-d6	1.713	1.648	3.8	70	0.00
8		2-Chlorophenol	1.322	1.283	3.0	71	0.00
9		Benzaldehyde	1.022	1.037	-1.5	73	0.00
10	C	Phenol	1.837	1.733	5.7	69	0.00
11		bis(2-Chloroethyl)ether	1.332	1.302	2.3	70	0.00
12		1,3-Dichlorobenzene	1.579	1.640	-3.9	75	0.00
13	C	1,4-Dichlorobenzene	1.599	1.673	-4.6	75	0.00
14		1,2-Dichlorobenzene	1.552	1.588	-2.3	74	0.00
15		Benzyl Alcohol	1.585	1.510	4.7	70	0.00
16		2,2'-oxybis(1-Chloropropane	2.177	2.117	2.8	71	0.00
17		2-Methylphenol	1.330	1.269	4.6	69	0.00
18		Hexachloroethane	0.616	0.659	-7.0	78	0.00
19	P	n-Nitroso-di-n-propylamine	1.318	1.291	2.0	71	0.00
20		3+4-Methylphenols	1.842	1.731	6.0	69	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	71	0.00
22		Acetophenone	0.561	0.567	-1.1	70	0.00
23	S	Nitrobenzene-d5	0.451	0.475	-5.3	74	0.00
24		Nitrobenzene	0.459	0.482	-5.0	73	0.00
25		Isophorone	0.857	0.847	1.2	68	0.00
26	C	2-Nitrophenol	0.189	0.201	-6.3	73	0.00
27		2,4-Dimethylphenol	0.313	0.307	1.9	68	0.00
28		bis(2-Chloroethoxy)methane	0.463	0.463	0.0	70	0.00
29	C	2,4-Dichlorophenol	0.397	0.398	-0.3	69	0.00
30		1,2,4-Trichlorobenzene	0.452	0.479	-6.0	73	0.00
31		Naphthalene	1.074	1.116	-3.9	72	0.00
32		Benzoic acid	0.223	0.216	3.1	67	0.00
33		4-Chloroaniline	0.498	0.488	2.0	68	0.00
34	C	Hexachlorobutadiene	0.346	0.376	-8.7	78	0.00
35		Caprolactam	0.131	0.115	12.2	61	0.00
36	C	4-Chloro-3-methylphenol	0.453	0.424	6.4	65	0.00
37		2-Methylnaphthalene	0.827	0.832	-0.6	70	0.00
38		1-Methylnaphthalene	0.779	0.788	-1.2	70	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	69	0.00
40		1,2,4,5-Tetrachlorobenzene	0.806	0.859	-6.6	72	0.00
41	P	Hexachlorocyclopentadiene	0.361	0.380	-5.3	72	0.00
42	S	2,4,6-Tribromophenol	0.297	0.284	4.4	65	0.00
43	C	2,4,6-Trichlorophenol	0.521	0.534	-2.5	68	0.00
44		2,4,5-Trichlorophenol	0.550	0.529	3.8	65	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061319\
Data File : BG041230.D
Acq On : 13 Jun 2019 19:31
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 14 09:03:01 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Jun 09 23:08:34 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.389	1.482	-6.7	71	0.00
46		1,1'-Biphenyl	1.480	1.527	-3.2	68	0.00
47		2-Chloronaphthalene	1.271	1.313	-3.3	69	0.00
48		2-Nitroaniline	0.456	0.460	-0.9	68	0.00
49		Acenaphthylene	1.913	1.905	0.4	66	0.00
50		Dimethylphthalate	1.693	1.648	2.7	65	0.00
51		2,6-Dinitrotoluene	0.365	0.362	0.8	67	0.00
52	C	Acenaphthene	1.159	1.134	2.2	66	0.00
53		3-Nitroaniline	0.365	0.347	4.9	62	0.00
54	P	2,4-Dinitrophenol	0.142	0.128	9.9	64	0.00
55		Dibenzofuran	1.950	1.933	0.9	65	0.00
56	P	4-Nitrophenol	0.250	0.217	13.2	57	0.02
57		2,4-Dinitrotoluene	0.516	0.496	3.9	64	0.00
58		Fluorene	1.553	1.478	4.8	64	0.00
59		2,3,4,6-Tetrachlorophenol	0.540	0.518	4.1	64	0.00
60		Diethylphthalate	1.728	1.657	4.1	64	0.00
61		4-Chlorophenyl-phenylether	1.009	0.976	3.3	66	0.00
62		4-Nitroaniline	0.386	0.362	6.2	63	0.00
63		Azobenzene	1.570	1.489	5.2	62	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	63	0.00
65		4,6-Dinitro-2-methylphenol	0.096	0.111	-15.6	72	0.00
66	c	n-Nitrosodiphenylamine	0.562	0.574	-2.1	63	0.00
67		4-Bromophenyl-phenylether	0.270	0.272	-0.7	62	0.00
68		Hexachlorobenzene	0.287	0.284	1.0	61	0.00
69		Atrazine	0.217	0.220	-1.4	58	0.00
70	C	Pentachlorophenol	0.139	0.143	-2.9	61	0.00
71		Phenanthrene	1.042	1.092	-4.8	64	0.00
72		Anthracene	1.027	1.075	-4.7	64	0.00
73		Carbazole	0.882	0.922	-4.5	64	0.00
74		Di-n-butylphthalate	1.096	1.138	-3.8	62	0.00
75	C	Fluoranthene	1.287	1.393	-8.2	66	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	65	0.00
77		Benzidine	0.477	0.341	28.5#	44#	0.00
78		Pyrene	1.300	1.372	-5.5	66	0.00
79	S	Terphenyl-d14	0.942	1.033	-9.7	67	0.00
80		Butylbenzylphthalate	0.506	0.523	-3.4	65	0.00
81		Benzo(a)anthracene	1.331	1.360	-2.2	65	0.00
82		3,3'-Dichlorobenzidine	0.468	0.448	4.3	62	0.00
83		Chrysene	1.274	1.321	-3.7	65	0.00
84		Bis(2-ethylhexyl)phthalate	0.712	0.712	0.0	63	0.00
85	c	Di-n-octyl phthalate	1.186	1.203	-1.4	64	0.00
86		Indeno(1,2,3-cd)pyrene	1.561	0.979	37.3#	41#	-0.01
87	I	Perylene-d12	1.000	1.000	0.0	58	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061319\
Data File : BG041230.D
Acq On : 13 Jun 2019 19:31
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 14 09:03:01 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Jun 09 23:08:34 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.213	1.245	-2.6	58	0.00
89	Benzo(k)fluoranthene	1.200	1.364	-13.7	64	0.00
90 C	Benzo(a)pyrene	1.157	1.230	-6.3	60	0.00
91	Dibenzo(a,h)anthracene	1.156	0.806	30.3#	39#	0.00
92	Benzo(g,h,i)perylene	1.124	0.637	43.3#	33#	-0.01

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