Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG062618\
Data File : BG035427.D
Acq On : 27 Jun 2018 2:41
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 27 08:25:47 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	153#	-0.02
2		1,4-Dioxane	0.565	0.534	5.5	146	-0.02
3		Pyridine	1.526	1.491	2.3	140	-0.02
4		n-Nitrosodimethylamine	0.655	0.574	12.4	127	-0.02
5	S	2-Fluorophenol	1.185	1.163	1.9	147	-0.02
6		Aniline	2.261	2.210	2.3	147	-0.02
7	S	Phenol-d6	1.749	1.795	-2.6	153#	-0.01
8		2-Chlorophenol	1.242	1.206	2.9	144	-0.02
9		Benzaldehyde	1.347	1.133	15.9	122	-0.02
10	C	Phenol	1.810	1.758	2.9	146	-0.01
11		bis(2-Chloroethyl)ether	1.405	1.343	4.4	144	-0.02
12		1,3-Dichlorobenzene	1.482	1.470	0.8	154#	-0.02
13	C	1,4-Dichlorobenzene	1.530	1.494	2.4	148	-0.02
14		1,2-Dichlorobenzene	1.437	1.446	-0.6	151#	-0.03
15		Benzyl Alcohol	1.657	1.543	6.9	139	-0.02
16		2,2'-oxybis(1-Chloropropane	1.398	1.180	15.6	129	-0.03
17		2-Methylphenol	1.193	1.198	-0.4	147	-0.02
18		Hexachloroethane	0.548	0.558	-1.8	150#	-0.03
19	P	n-Nitroso-di-n-propylamine	1.486	1.307	12.0	131	-0.03
20		3+4-Methylphenols	1.711	1.706	0.3	149	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	157#	-0.03
22		Acetophenone	0.620	0.574	7.4	144	-0.02
23	S	Nitrobenzene-d5	0.421	0.447	-6.2	158#	-0.02
24		Nitrobenzene	0.431	0.437	-1.4	152#	-0.02
25		Isophorone	0.888	0.796	10.4	139	-0.03
26	C	2-Nitrophenol	0.149	0.185	-24.2#	194#	-0.02
27		2,4-Dimethylphenol	0.312	0.280	10.3	141	-0.02
28		bis(2-Chloroethoxy)methane	0.477	0.445	6.7	144	-0.03
29	C	2,4-Dichlorophenol	0.340	0.351	-3.2	158#	-0.02
30		1,2,4-Trichlorobenzene	0.407	0.403	1.0	150	-0.02
31		Naphthalene	1.039	0.998	3.9	150	-0.02
32		Benzoic acid	0.209	0.218	-4.3	162#	0.00
33		4-Chloroaniline	0.451	0.447	0.9	152#	-0.02
34	C	Hexachlorobutadiene	0.317	0.315	0.6	154#	-0.03
35		Caprolactam	0.126	0.128	-1.6	160#	-0.01
36	C	4-Chloro-3-methylphenol	0.423	0.440	-4.0	160#	-0.01
37		2-Methylnaphthalene	0.788	0.786	0.3	152#	-0.03
38	I	Acenaphthene-d10	1.000	1.000	0.0	170#	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.745	0.738	0.9	163#	-0.02
40	P	Hexachlorocyclopentadiene	0.467	0.417	10.7	144	-0.03
41	S	2,4,6-Tribromophenol	0.256	0.287	-12.1	187#	-0.02
42	C	2,4,6-Trichlorophenol	0.446	0.471	-5.6	172#	-0.02
43		2,4,5-Trichlorophenol	0.490	0.521	-6.3	179#	-0.01
44	S	2-Fluorobiphenyl	1.538	1.425	7.3	156#	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG062618\
Data File : BG035427.D
Acq On : 27 Jun 2018 2:41
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 27 08:25:47 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.608	1.509	6.2	157#	-0.02
46		2-Chloronaphthalene	1.216	1.137	6.5	159#	-0.03
47		2-Nitroaniline	0.359	0.390	-8.6	182#	-0.02
48		Acenaphthylene	2.039	1.945	4.6	160#	-0.02
49		Dimethylphthalate	1.744	1.606	7.9	156#	-0.03
50		2,6-Dinitrotoluene	0.318	0.355	-11.6	183#	-0.02
51	C	Acenaphthene	1.332	1.228	7.8	153#	-0.02
52		3-Nitroaniline	0.312	0.355	-13.8	180#	-0.01
53	P	2,4-Dinitrophenol	0.129	0.166	-28.7#	210#	-0.02
54		Dibenzofuran	1.995	1.915	4.0	160#	-0.02
55	P	4-Nitrophenol	0.263	0.247	6.1	148	0.00
56		2,4-Dinitrotoluene	0.423	0.504	-19.1	195#	-0.02
57		Fluorene	1.667	1.598	4.1	159#	-0.02
58		2,3,4,6-Tetrachlorophenol	0.476	0.508	-6.7	177#	-0.02
59		Diethylphthalate	1.779	1.629	8.4	155#	-0.03
60		4-Chlorophenyl-phenylether	0.951	0.958	-0.7	171#	-0.03
61		4-Nitroaniline	0.335	0.368	-9.9	175#	-0.02
62		Azobenzene	1.744	1.528	12.4	148	-0.03
63	I	Phenanthrene-d10	1.000	1.000	0.0	174#	-0.02
64		4,6-Dinitro-2-methylphenol	0.091	0.115	-26.4#	216#	-0.02
65	c	n-Nitrosodiphenylamine	0.601	0.565	6.0	159#	-0.02
66		4-Bromophenyl-phenylether	0.241	0.245	-1.7	175#	-0.03
67		Hexachlorobenzene	0.245	0.250	-2.0	177#	-0.02
68		Atrazine	0.256	0.235	8.2	156#	-0.03
69	C	Pentachlorophenol	0.165	0.165	0.0	168#	-0.02
70		Phenanthrene	1.016	0.959	5.6	161#	-0.02
71		Anthracene	1.018	0.956	6.1	159#	-0.02
72		Carbazole	0.944	0.875	7.3	157#	-0.02
73		Di-n-butylphthalate	1.125	1.018	9.5	153#	-0.04
74	C	Fluoranthene	1.322	1.226	7.3	157#	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	178#	-0.04
76		Benzidine	0.744	0.595	20.0	139	-0.02
77		Pyrene	1.318	1.174	10.9	157#	-0.02
78	S	Terphenyl-d14	0.955	0.864	9.5	158#	-0.02
79		Butylbenzylphthalate	0.479	0.439	8.4	156#	-0.04
80		Benzo(a)anthracene	1.278	1.188	7.0	165#	-0.03
81		3,3'-Dichlorobenzidine	0.481	0.435	9.6	155#	-0.04
82		Chrysene	1.170	1.088	7.0	164#	-0.04
83		Bis(2-ethylhexyl)phthalate	0.676	0.594	12.1	152#	-0.05
84	c	Di-n-octyl phthalate	1.161	1.010	13.0	150	-0.07
85		Indeno(1,2,3-cd)pyrene	1.370	1.289	5.9	164#	-0.09
86	I	Perylene-d12	1.000	1.000	0.0	174#	-0.06
87		Benzo(b)fluoranthene	1.232	1.181	4.1	162#	-0.05

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG062618\
Data File : BG035427.D
Acq On : 27 Jun 2018 2:41
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 27 08:25:47 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.205	1.114	7.6	164#	-0.05
89 C	Benzo(a)pyrene	1.159	1.094	5.6	162#	-0.06
90	Dibenzo(a,h)anthracene	1.144	1.082	5.4	164#	-0.10
91	Benzo(g,h,i)perylene	1.120	1.074	4.1	165#	-0.10

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