Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG082416\
Data File : BG023891.D
Acq On : 24 Aug 2016 19:23
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 25 13:12:47 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG080116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 24 17:42:04 2016
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	74	0.00
2		1,4-Dioxane	0.506	0.482	4.7	77	0.00
3		Pyridine	1.664	1.510	9.3	70	0.00
4		n-Nitrosodimethylamine	0.617	0.696	-12.8	90	0.00
5	S	2-Fluorophenol	1.188	1.141	4.0	75	0.00
6		Aniline	2.674	2.190	18.1	64	-0.01
7	S	Phenol-d6	1.851	1.666	10.0	70	0.00
8		2-Chlorophenol	1.365	1.309	4.1	76	0.00
9		Benzaldehyde	1.096	1.198	-9.3	86	0.00
10	C	Phenol	1.980	1.794	9.4	71	0.00
11		bis(2-Chloroethyl)ether	1.579	1.448	8.3	73	0.00
12		1,3-Dichlorobenzene	1.563	1.560	0.2	78	0.00
13	C	1,4-Dichlorobenzene	1.600	1.571	1.8	79	0.00
14		1,2-Dichlorobenzene	1.569	1.487	5.2	76	0.00
15		Benzyl Alcohol	1.402	1.407	-0.4	77	0.00
16		2,2'-oxybis(1-Chloropropane	2.322	2.179	6.2	74	0.00
17		2-Methylphenol	1.327	1.215	8.4	72	0.00
18		Hexachloroethane	0.602	0.622	-3.3	83	0.00
19	P	n-Nitroso-di-n-propylamine	1.592	1.365	14.3	67	0.00
20		3+4-Methylphenols	1.871	1.691	9.6	70	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	66	-0.01
22		Acetophenone	0.548	0.539	1.6	69	0.00
23	S	Nitrobenzene-d5	0.402	0.419	-4.2	71	0.00
24		Nitrobenzene	0.445	0.492	-10.6	76	0.00
25		Isophorone	0.838	0.841	-0.4	69	0.00
26	C	2-Nitrophenol	0.188	0.199	-5.9	71	-0.01
27		2,4-Dimethylphenol	0.320	0.318	0.6	66	0.00
28		bis(2-Chloroethoxy)methane	0.504	0.445	11.7	61	0.00
29	C	2,4-Dichlorophenol	0.322	0.320	0.6	66	0.00
30		1,2,4-Trichlorobenzene	0.347	0.355	-2.3	71	0.00
31		Naphthalene	1.018	0.996	2.2	68	-0.01
32		Benzoic acid	0.265	0.242	8.7	57	0.00
33		4-Chloroaniline	0.487	0.412	15.4	58	-0.01
34	C	Hexachlorobutadiene	0.228	0.245	-7.5	76	0.00
35		Caprolactam	0.135	0.107	20.7	52	0.00
36	C	4-Chloro-3-methylphenol	0.423	0.383	9.5	62	0.00
37		2-Methylnaphthalene	0.774	0.716	7.5	64	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	58	0.00
39		1,2,4,5-Tetrachlorobenzene	0.725	0.774	-6.8	65	-0.01
40	P	Hexachlorocyclopentadiene	0.366	0.326	10.9	47#	0.00
41	S	2,4,6-Tribromophenol	0.293	0.239	18.4	49#	0.00
42	C	2,4,6-Trichlorophenol	0.432	0.431	0.2	59	0.00
43		2,4,5-Trichlorophenol	0.445	0.431	3.1	57	0.00
44	S	2-Fluorobiphenyl	1.186	1.262	-6.4	67	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG082416\
Data File : BG023891.D
Acq On : 24 Aug 2016 19:23
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 25 13:12:47 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG080116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 24 17:42:04 2016
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	1.529	1.567	-2.5	63	0.00
46		2-Chloronaphthalene	1.183	1.212	-2.5	62	0.00
47		2-Nitroaniline	0.491	0.479	2.4	57	0.00
48		Acenaphthylene	1.870	1.873	-0.2	61	0.00
49		Dimethylphthalate	1.534	1.558	-1.6	62	0.00
50		2,6-Dinitrotoluene	0.363	0.350	3.6	57	0.00
51	C	Acenaphthene	1.184	1.181	0.3	60	0.00
52		3-Nitroaniline	0.410	0.357	12.9	51	0.00
53	P	2,4-Dinitrophenol	0.233	0.236	-1.3	58	0.00
54		Dibenzofuran	1.720	1.665	3.2	60	0.00
55	P	4-Nitrophenol	0.322	0.254	21.1	45#	0.00
56		2,4-Dinitrotoluene	0.510	0.491	3.7	57	0.00
57		Fluorene	1.501	1.478	1.5	61	0.00
58		2,3,4,6-Tetrachlorophenol	0.408	0.364	10.8	53	0.00
59		Diethylphthalate	1.558	1.562	-0.3	61	0.00
60		4-Chlorophenyl-phenylether	0.794	0.758	4.5	58	0.00
61		4-Nitroaniline	0.436	0.385	11.7	52	0.00
62		Azobenzene	1.579	1.554	1.6	60	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	54	0.00
64		4,6-Dinitro-2-methylphenol	0.136	0.129	5.1	49#	0.00
65	c	n-Nitrosodiphenylamine	0.587	0.582	0.9	55	0.00
66		4-Bromophenyl-phenylether	0.233	0.232	0.4	55	0.00
67		Hexachlorobenzene	0.255	0.246	3.5	54	0.00
68		Atrazine	0.225	0.225	0.0	53	0.00
69	C	Pentachlorophenol	0.149	0.117	21.5#	37#	0.00
70		Phenanthrene	1.015	1.016	-0.1	60	0.00
71		Anthracene	1.021	1.041	-2.0	61	0.00
72		Carbazole	0.896	0.942	-5.1	59	0.00
73		Di-n-butylphthalate	1.021	1.137	-11.4	65	0.00
74	C	Fluoranthene	1.086	1.195	-10.0	65	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	69	0.00
76		Benzidine	0.556	0.571	-2.7	66	0.00
77		Pyrene	1.239	1.123	9.4	65	0.00
78	S	Terphenyl-d14	0.733	0.674	8.0	72	0.00
79		Butylbenzylphthalate	0.482	0.525	-8.9	76	-0.01
80		Benzo(a)anthracene	1.104	1.099	0.5	73	0.00
81		3,3'-Dichlorobenzidine	0.453	0.426	6.0	66	0.00
82		Chrysene	1.050	1.053	-0.3	75	0.00
83		Bis(2-ethylhexyl)phthalate	0.659	0.721	-9.4	78	-0.01
84	c	Di-n-octyl phthalate	1.126	1.206	-7.1	77	-0.01
85		Indeno(1,2,3-cd)pyrene	1.312	1.296	1.2	70	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	68	-0.02
87		Benzo(b)fluoranthene	1.125	1.102	2.0	70	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG082416\
Data File : BG023891.D
Acq On : 24 Aug 2016 19:23
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 25 13:12:47 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG080116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 24 17:42:04 2016
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	1.081	1.085	-0.4	72	-0.01
89 C	Benzo(a)pyrene	1.070	1.079	-0.8	72	-0.02
90	Dibenzo(a,h)anthracene	1.093	1.079	1.3	70	-0.02
91	Benzo(g,h,i)perylene	1.101	1.053	4.4	67	-0.02

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