Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG071116\
Data File : BG023020.D
Acq On : 12 Jul 2016 3:58
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 12 04:31:49 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG070816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 07 16:51:07 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	85	0.00
2		1,4-Dioxane	0.475	0.459	3.4	86	0.00
3		Pyridine	1.627	1.550	4.7	82	0.00
4		n-Nitrosodimethylamine	0.698	0.689	1.3	83	0.00
5	S	2-Fluorophenol	1.223	1.209	1.1	86	0.00
6		Aniline	2.654	2.676	-0.8	85	0.00
7	S	Phenol-d6	1.915	2.058	-7.5	90	0.00
8		2-Chlorophenol	1.301	1.369	-5.2	90	0.00
9		Benzaldehyde	1.280	1.267	1.0	86	0.00
10	C	Phenol	1.971	2.083	-5.7	89	0.00
11		bis(2-Chloroethyl)ether	1.562	1.533	1.9	84	0.00
12		1,3-Dichlorobenzene	1.593	1.587	0.4	88	0.00
13	C	1,4-Dichlorobenzene	1.595	1.623	-1.8	88	0.00
14		1,2-Dichlorobenzene	1.585	1.585	0.0	89	0.00
15		Benzyl Alcohol	1.754	1.923	-9.6	91	0.00
16		2,2'-oxybis(1-Chloropropane	1.908	1.982	-3.9	89	0.00
17		2-Methylphenol	1.283	1.349	-5.1	89	0.00
18		Hexachloroethane	0.673	0.658	2.2	90	0.00
19	P	n-Nitroso-di-n-propylamine	1.727	1.767	-2.3	85	0.00
20		3+4-Methylphenols	1.789	1.972	-10.2	89	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	88	0.00
22		Acetophenone	0.600	0.593	1.2	87	0.00
23	S	Nitrobenzene-d5	0.467	0.468	-0.2	89	0.00
24		Nitrobenzene	0.496	0.488	1.6	89	0.00
25		Isophorone	0.939	0.892	5.0	81	0.00
26	C	2-Nitrophenol	0.195	0.195	0.0	83	0.00
27		2,4-Dimethylphenol	0.337	0.333	1.2	88	0.00
28		bis(2-Chloroethoxy)methane	0.524	0.511	2.5	85	0.00
29	C	2,4-Dichlorophenol	0.359	0.365	-1.7	88	0.00
30		1,2,4-Trichlorobenzene	0.411	0.410	0.2	89	0.00
31		Naphthalene	1.018	1.041	-2.3	91	0.00
32		Benzoic acid	0.306	0.287	6.2	78	0.00
33		4-Chloroaniline	0.498	0.493	1.0	87	0.00
34	C	Hexachlorobutadiene	0.334	0.316	5.4	86	0.00
35		Caprolactam	0.148	0.130	12.2	77	0.00
36	C	4-Chloro-3-methylphenol	0.491	0.500	-1.8	88	0.00
37		2-Methylnaphthalene	0.805	0.827	-2.7	89	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	85	0.00
39		1,2,4,5-Tetrachlorobenzene	0.862	0.850	1.4	85	0.00
40	P	Hexachlorocyclopentadiene	0.496	0.531	-7.1	92	0.00
41	S	2,4,6-Tribromophenol	0.359	0.298	17.0	73	0.00
42	C	2,4,6-Trichlorophenol	0.495	0.487	1.6	83	0.00
43		2,4,5-Trichlorophenol	0.509	0.493	3.1	85	0.00
44	S	2-Fluorobiphenyl	1.270	1.267	0.2	88	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG071116\
Data File : BG023020.D
Acq On : 12 Jul 2016 3:58
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 12 04:31:49 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG070816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 07 16:51:07 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	1.501	1.532	-2.1	88	0.00
46		2-Chloronaphthalene	1.217	1.227	-0.8	87	0.00
47		2-Nitroaniline	0.519	0.519	0.0	87	0.00
48		Acenaphthylene	1.826	1.798	1.5	85	0.00
49		Dimethylphthalate	1.634	1.515	7.3	82	0.00
50		2,6-Dinitrotoluene	0.367	0.338	7.9	80	0.00
51	C	Acenaphthene	1.193	1.175	1.5	86	0.00
52		3-Nitroaniline	0.366	0.346	5.5	81	0.00
53	P	2,4-Dinitrophenol	0.252	0.296	-17.5	97	0.00
54		Dibenzofuran	1.786	1.702	4.7	84	0.00
55	P	4-Nitrophenol	0.307	0.260	15.3	73	0.00
56		2,4-Dinitrotoluene	0.535	0.488	8.8	80	0.00
57		Fluorene	1.537	1.436	6.6	83	0.00
58		2,3,4,6-Tetrachlorophenol	0.509	0.439	13.8	74	0.00
59		Diethylphthalate	1.662	1.516	8.8	80	0.00
60		4-Chlorophenyl-phenylether	0.936	0.851	9.1	78	0.00
61		4-Nitroaniline	0.396	0.348	12.1	77	0.00
62		Azobenzene	1.590	1.539	3.2	85	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	77	0.00
64		4,6-Dinitro-2-methylphenol	0.149	0.131	12.1	65	0.00
65	c	n-Nitrosodiphenylamine	0.576	0.606	-5.2	80	0.00
66		4-Bromophenyl-phenylether	0.260	0.255	1.9	74	0.00
67		Hexachlorobenzene	0.283	0.277	2.1	75	0.00
68		Atrazine	0.256	0.248	3.1	76	0.00
69	C	Pentachlorophenol	0.183	0.158	13.7	63	0.00
70		Phenanthrene	0.980	0.989	-0.9	79	0.00
71		Anthracene	0.992	1.018	-2.6	80	0.00
72		Carbazole	0.858	0.857	0.1	78	0.00
73		Di-n-butylphthalate	0.954	1.013	-6.2	80	0.00
74	C	Fluoranthene	1.203	1.163	3.3	79	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	74	0.03
76		Benzidine	0.573	0.640	-11.7	81	0.00
77		Pyrene	1.124	1.136	-1.1	80	0.00
78	S	Terphenyl-d14	0.646	0.706	-9.3	82	0.00
79		Butylbenzylphthalate	0.445	0.463	-4.0	78	0.01
80		Benzo(a)anthracene	1.119	1.126	-0.6	77	0.03
81		3,3'-Dichlorobenzidine	0.491	0.484	1.4	72	0.03
82		Chrysene	1.050	1.068	-1.7	78	0.03
83		Bis(2-ethylhexyl)phthalate	0.618	0.647	-4.7	80	0.03
84	c	Di-n-octyl phthalate	1.031	1.091	-5.8	79	0.04
85		Indeno(1,2,3-cd)pyrene	1.338	1.221	8.7	70	0.12
86	I	Perylene-d12	1.000	1.000	0.0	73	0.07
87		Benzo(b)fluoranthene	1.154	1.154	0.0	75	0.06

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG071116\
Data File : BG023020.D
Acq On : 12 Jul 2016 3:58
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 12 04:31:49 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG070816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 07 16:51:07 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	1.095	1.102	-0.6	72	0.06
89 C	Benzo(a)pyrene	1.106	1.105	0.1	74	0.07
90	Dibenzo(a,h)anthracene	1.098	1.057	3.7	71	0.12
91	Benzo(g,h,i)perylene	1.099	1.030	6.3	69	0.14

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