Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG042318\
Data File : BG033933.D
Acq On : 23 Apr 2018 23:32
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 24 03:24:15 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 19 13:56:22 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	152#	0.00
2		1,4-Dioxane	0.529	0.450	14.9	134	0.00
3		Pyridine	1.532	1.429	6.7	131	0.00
4		n-Nitrosodimethylamine	0.698	0.639	8.5	139	0.00
5	S	2-Fluorophenol	1.253	1.173	6.4	139	0.00
6		Aniline	2.442	2.233	8.6	131	0.00
7	S	Phenol-d6	1.827	1.733	5.1	138	0.00
8		2-Chlorophenol	1.404	1.308	6.8	137	-0.01
9		Benzaldehyde	0.996	0.879	11.7	134	0.00
10	C	Phenol	1.872	1.731	7.5	136	0.00
11		bis(2-Chloroethyl)ether	1.432	1.349	5.8	136	-0.01
12		1,3-Dichlorobenzene	1.616	1.537	4.9	138	-0.01
13	C	1,4-Dichlorobenzene	1.637	1.564	4.5	140	-0.01
14		1,2-Dichlorobenzene	1.592	1.516	4.8	140	0.00
15		Benzyl Alcohol	1.593	1.431	10.2	131	0.00
16		2,2'-oxybis(1-Chloropropane	2.792	2.392	14.3	124	-0.02
17		2-Methylphenol	1.387	1.274	8.1	135	0.00
18		Hexachloroethane	0.598	0.567	5.2	141	0.00
19	P	n-Nitroso-di-n-propylamine	1.557	1.354	13.0	125	0.00
20		3+4-Methylphenols	2.015	1.840	8.7	134	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	140	-0.01
22		Acetophenone	0.544	0.511	6.1	131	0.00
23	S	Nitrobenzene-d5	0.351	0.359	-2.3	142	0.00
24		Nitrobenzene	0.383	0.379	1.0	134	0.00
25		Isophorone	0.790	0.717	9.2	126	-0.01
26	C	2-Nitrophenol	0.151	0.175	-15.9	162#	-0.01
27		2,4-Dimethylphenol	0.283	0.275	2.8	133	0.00
28		bis(2-Chloroethoxy)methane	0.418	0.399	4.5	132	-0.01
29	C	2,4-Dichlorophenol	0.316	0.319	-0.9	138	0.00
30		1,2,4-Trichlorobenzene	0.353	0.351	0.6	138	0.00
31		Naphthalene	1.028	0.967	5.9	131	0.00
32		Benzoic acid	0.262	0.272	-3.8	140	0.00
33		4-Chloroaniline	0.483	0.459	5.0	130	0.00
34	C	Hexachlorobutadiene	0.211	0.218	-3.3	136	-0.01
35		Caprolactam	0.134	0.137	-2.2	142	0.00
36	C	4-Chloro-3-methylphenol	0.385	0.362	6.0	130	-0.01
37		2-Methylnaphthalene	0.787	0.743	5.6	131	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	135	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.664	0.646	2.7	133	-0.01
40	P	Hexachlorocyclopentadiene	0.365	0.330	9.6	120	-0.01
41	S	2,4,6-Tribromophenol	0.233	0.249	-6.9	139	-0.01
42	C	2,4,6-Trichlorophenol	0.405	0.419	-3.5	141	0.00
43		2,4,5-Trichlorophenol	0.466	0.491	-5.4	143	0.00
44	S	2-Fluorobiphenyl	1.361	1.281	5.9	128	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.619	1.542	4.8	130	0.00
46		2-Chloronaphthalene	1.228	1.174	4.4	131	-0.01
47		2-Nitroaniline	0.383	0.403	-5.2	142	-0.01
48		Acenaphthylene	2.041	1.902	6.8	127	0.00
49		Dimethylphthalate	1.759	1.631	7.3	127	-0.01
50		2,6-Dinitrotoluene	0.333	0.353	-6.0	141	-0.01
51	C	Acenaphthene	1.282	1.196	6.7	124	-0.01
52		3-Nitroaniline	0.362	0.392	-8.3	139	0.00
53	P	2,4-Dinitrophenol	0.119	0.101	15.1	116	-0.01
54		Dibenzofuran	1.900	1.775	6.6	128	-0.01
55	P	4-Nitrophenol	0.284	0.292	-2.8	131	0.00
56		2,4-Dinitrotoluene	0.443	0.492	-11.1	149	0.00
57		Fluorene	1.645	1.529	7.1	127	-0.01
58		2,3,4,6-Tetrachlorophenol	0.426	0.432	-1.4	131	-0.01
59		Diethylphthalate	1.860	1.684	9.5	124	-0.01
60		4-Chlorophenyl-phenylether	0.860	0.819	4.8	132	-0.01
61		4-Nitroaniline	0.384	0.403	-4.9	137	0.00
62		Azobenzene	1.658	1.460	11.9	120	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	135	-0.01
64		4,6-Dinitro-2-methylphenol	0.088	0.093	-5.7	139	0.00
65	c	n-Nitrosodiphenylamine	0.581	0.555	4.5	130	-0.01
66		4-Bromophenyl-phenylether	0.220	0.213	3.2	131	-0.02
67		Hexachlorobenzene	0.234	0.227	3.0	129	-0.01
68		Atrazine	0.228	0.215	5.7	128	-0.02
69	C	Pentachlorophenol	0.161	0.152	5.6	125	-0.01
70		Phenanthrene	1.073	0.993	7.5	125	-0.02
71		Anthracene	1.073	1.000	6.8	125	-0.01
72		Carbazole	1.035	0.941	9.1	124	-0.01
73		Di-n-butylphthalate	1.288	1.150	10.7	121	-0.02
74	C	Fluoranthene	1.298	1.188	8.5	124	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	139	-0.02
76		Benzidine	0.540	0.471	12.8	122	-0.01
77		Pyrene	1.312	1.195	8.9	125	-0.01
78	S	Terphenyl-d14	0.890	0.797	10.4	124	-0.01
79		Butylbenzylphthalate	0.576	0.533	7.5	127	-0.02
80		Benzo(a)anthracene	1.261	1.176	6.7	129	-0.02
81		3,3'-Dichlorobenzidine	0.470	0.460	2.1	134	-0.02
82		Chrysene	1.204	1.127	6.4	130	-0.01
83		Bis(2-ethylhexyl)phthalate	0.814	0.755	7.2	127	-0.03
84	c	Di-n-octyl phthalate	1.291	1.203	6.8	126	-0.04
85		Indeno(1,2,3-cd)pyrene	1.371	1.360	0.8	135	-0.06
86	I	Perylene-d12	1.000	1.000	0.0	143	-0.03
87		Benzo(b)fluoranthene	1.220	1.164	4.6	136	-0.03

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88	Benzo(k)fluoranthene	1.212	1.107	8.7	130	-0.03
89 C	Benzo(a)pyrene	1.169	1.091	6.7	132	-0.04
90	Dibenzo(a,h)anthracene	1.131	1.076	4.9	135	-0.07
91	Benzo(g,h,i)perylene	1.113	1.062	4.6	135	-0.07

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