Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG042318\
Data File : BG033915.D
Acq On : 23 Apr 2018 11:24
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 24 03:15:36 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	146	0.00
2		1,4-Dioxane	0.529	0.486	8.1	139	0.00
3		Pyridine	1.532	1.438	6.1	128	0.00
4		n-Nitrosodimethylamine	0.698	0.611	12.5	128	0.00
5	S	2-Fluorophenol	1.253	1.157	7.7	133	0.00
6		Aniline	2.442	2.241	8.2	127	0.00
7	S	Phenol-d6	1.827	1.660	9.1	128	0.00
8		2-Chlorophenol	1.404	1.292	8.0	131	0.00
9		Benzaldehyde	0.996	0.824	17.3	121	0.00
10	C	Phenol	1.872	1.673	10.6	127	0.00
11		bis(2-Chloroethyl)ether	1.432	1.332	7.0	130	-0.01
12		1,3-Dichlorobenzene	1.616	1.507	6.7	131	0.00
13	C	1,4-Dichlorobenzene	1.637	1.531	6.5	133	-0.01
14		1,2-Dichlorobenzene	1.592	1.467	7.9	131	0.00
15		Benzyl Alcohol	1.593	1.403	11.9	124	0.00
16		2,2'-oxybis(1-Chloropropane	2.792	2.408	13.8	121	-0.01
17		2-Methylphenol	1.387	1.217	12.3	125	0.00
18		Hexachloroethane	0.598	0.551	7.9	133	0.00
19	P	n-Nitroso-di-n-propylamine	1.557	1.313	15.7	118	0.00
20		3+4-Methylphenols	2.015	1.743	13.5	122	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	132	-0.01
22		Acetophenone	0.544	0.509	6.4	124	0.00
23	S	Nitrobenzene-d5	0.351	0.360	-2.6	134	0.00
24		Nitrobenzene	0.383	0.376	1.8	126	0.00
25		Isophorone	0.790	0.703	11.0	116	-0.01
26	C	2-Nitrophenol	0.151	0.171	-13.2	149	0.00
27		2,4-Dimethylphenol	0.283	0.266	6.0	121	0.00
28		bis(2-Chloroethoxy)methane	0.418	0.390	6.7	122	-0.01
29	C	2,4-Dichlorophenol	0.316	0.303	4.1	124	0.00
30		1,2,4-Trichlorobenzene	0.353	0.344	2.5	128	0.00
31		Naphthalene	1.028	0.965	6.1	123	0.00
32		Benzoic acid	0.262	0.284	-8.4	138	0.00
33		4-Chloroaniline	0.483	0.444	8.1	119	0.00
34	C	Hexachlorobutadiene	0.211	0.219	-3.8	129	0.00
35		Caprolactam	0.134	0.122	9.0	120	0.00
36	C	4-Chloro-3-methylphenol	0.385	0.339	11.9	115	0.00
37		2-Methylnaphthalene	0.787	0.728	7.5	121	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	120	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.664	0.656	1.2	120	-0.01
40	P	Hexachlorocyclopentadiene	0.365	0.392	-7.4	127	0.00
41	S	2,4,6-Tribromophenol	0.233	0.238	-2.1	118	-0.01
42	C	2,4,6-Trichlorophenol	0.405	0.412	-1.7	123	0.00
43		2,4,5-Trichlorophenol	0.466	0.472	-1.3	122	0.00
44	S	2-Fluorobiphenyl	1.361	1.305	4.1	116	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.619	1.549	4.3	116	0.00
46		2-Chloronaphthalene	1.228	1.179	4.0	117	-0.01
47		2-Nitroaniline	0.383	0.396	-3.4	123	-0.01
48		Acenaphthylene	2.041	1.919	6.0	114	0.00
49		Dimethylphthalate	1.759	1.585	9.9	109	-0.01
50		2,6-Dinitrotoluene	0.333	0.341	-2.4	121	-0.01
51	C	Acenaphthene	1.282	1.187	7.4	109	0.00
52		3-Nitroaniline	0.362	0.380	-5.0	119	0.00
53	P	2,4-Dinitrophenol	0.119	0.148	-24.4	151#	-0.01
54		Dibenzofuran	1.900	1.746	8.1	112	0.00
55	P	4-Nitrophenol	0.284	0.295	-3.9	117	0.00
56		2,4-Dinitrotoluene	0.443	0.477	-7.7	128	0.00
57		Fluorene	1.645	1.496	9.1	110	0.00
58		2,3,4,6-Tetrachlorophenol	0.426	0.416	2.3	112	0.00
59		Diethylphthalate	1.860	1.639	11.9	107	-0.01
60		4-Chlorophenyl-phenylether	0.860	0.799	7.1	114	-0.01
61		4-Nitroaniline	0.384	0.393	-2.3	118	0.00
62		Azobenzene	1.658	1.445	12.8	106	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	115	-0.01
64		4,6-Dinitro-2-methylphenol	0.088	0.109	-23.9	138	0.00
65	c	n-Nitrosodiphenylamine	0.581	0.549	5.5	110	0.00
66		4-Bromophenyl-phenylether	0.220	0.215	2.3	112	-0.01
67		Hexachlorobenzene	0.234	0.226	3.4	110	-0.01
68		Atrazine	0.228	0.211	7.5	108	-0.01
69	C	Pentachlorophenol	0.161	0.162	-0.6	114	-0.01
70		Phenanthrene	1.073	0.994	7.4	107	-0.01
71		Anthracene	1.073	1.007	6.2	107	-0.01
72		Carbazole	1.035	0.946	8.6	107	-0.01
73		Di-n-butylphthalate	1.288	1.210	6.1	109	-0.01
74	C	Fluoranthene	1.298	1.181	9.0	106	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	115	-0.01
76		Benzidine	0.540	0.470	13.0	101	-0.01
77		Pyrene	1.312	1.223	6.8	106	-0.01
78	S	Terphenyl-d14	0.890	0.826	7.2	106	-0.01
79		Butylbenzylphthalate	0.576	0.537	6.8	106	-0.02
80		Benzo(a)anthracene	1.261	1.178	6.6	107	-0.02
81		3,3'-Dichlorobenzidine	0.470	0.453	3.6	109	-0.01
82		Chrysene	1.204	1.118	7.1	106	-0.01
83		Bis(2-ethylhexyl)phthalate	0.814	0.769	5.5	107	-0.02
84	c	Di-n-octyl phthalate	1.291	1.229	4.8	107	-0.03
85		Indeno(1,2,3-cd)pyrene	1.371	1.340	2.3	110	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	118	-0.02
87		Benzo(b)fluoranthene	1.220	1.133	7.1	109	-0.02

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88	Benzo(k)fluoranthene	1.212	1.110	8.4	108	-0.02
89 C	Benzo(a)pyrene	1.169	1.092	6.6	110	-0.03
90	Dibenzo(a,h)anthracene	1.131	1.061	6.2	110	-0.04
91	Benzo(g,h,i)perylene	1.113	1.048	5.8	110	-0.04

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