Evaluate Continuing Calibration Report

Data Path : U:\HPCHEM1\BNA G\Data\BG122617\
Data File : BG031756.D
Acq On : 26 Dec 2017 11:36
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 26 12:09:22 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 21 14:09:19 2017
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	82	-0.03
2		1,4-Dioxane	0.407	0.417	-2.5	88	-0.02
3		Pyridine	1.088	1.175	-8.0	85	-0.02
4		n-Nitrosodimethylamine	0.439	0.443	-0.9	82	-0.02
5	S	2-Fluorophenol	1.098	1.074	2.2	81	-0.02
6		Aniline	1.834	1.757	4.2	77	-0.02
7	S	Phenol-d6	1.426	1.411	1.1	79	-0.02
8		2-Chlorophenol	1.274	1.264	0.8	81	-0.03
9		Benzaldehyde	0.858	0.784	8.6	74	-0.03
10	C	Phenol	1.439	1.380	4.1	78	-0.02
11		bis(2-Chloroethyl)ether	1.195	1.079	9.7	74	-0.02
12		1,3-Dichlorobenzene	1.582	1.543	2.5	81	-0.03
13	C	1,4-Dichlorobenzene	1.616	1.552	4.0	80	-0.03
14		1,2-Dichlorobenzene	1.531	1.476	3.6	80	-0.02
15		Benzyl Alcohol	1.070	1.054	1.5	78	-0.03
16		2,2'-oxybis(1-Chloropropane	0.973	0.802	17.6	68	-0.02
17		2-Methylphenol	1.030	0.997	3.2	78	-0.02
18		Hexachloroethane	0.489	0.466	4.7	81	-0.03
19	P	n-Nitroso-di-n-propylamine	0.974	0.909	6.7	77	-0.03
20		3+4-Methylphenols	1.464	1.428	2.5	78	-0.03
21	I	Naphthalene-d8	1.000	1.000	0.0	80	-0.03
22		Acetophenone	0.465	0.445	4.3	77	-0.03
23	S	Nitrobenzene-d5	0.265	0.270	-1.9	79	-0.03
24		Nitrobenzene	0.273	0.275	-0.7	77	-0.03
25		Isophorone	0.570	0.540	5.3	75	-0.03
26	C	2-Nitrophenol	0.143	0.146	-2.1	80	-0.03
27		2,4-Dimethylphenol	0.301	0.291	3.3	79	-0.02
28		bis(2-Chloroethoxy)methane	0.383	0.354	7.6	74	-0.03
29	C	2,4-Dichlorophenol	0.354	0.355	-0.3	79	-0.03
30		1,2,4-Trichlorobenzene	0.432	0.414	4.2	77	-0.03
31		Naphthalene	1.009	0.993	1.6	79	-0.03
32		Benzoic acid	0.186	0.196	-5.4	79	-0.03
33		4-Chloroaniline	0.449	0.434	3.3	77	-0.02
34	C	Hexachlorobutadiene	0.296	0.289	2.4	79	-0.03
35		Caprolactam	0.107	0.110	-2.8	79	-0.02
36	C	4-Chloro-3-methylphenol	0.327	0.330	-0.9	80	-0.03
37		2-Methylnaphthalene	0.818	0.793	3.1	78	-0.03
38	I	Acenaphthene-d10	1.000	1.000	0.0	80	-0.03
39		1,2,4,5-Tetrachlorobenzene	0.842	0.804	4.5	77	-0.03
40	P	Hexachlorocyclopentadiene	0.444	0.451	-1.6	79	-0.03
41	S	2,4,6-Tribromophenol	0.305	0.339	-11.1	88	-0.03
42	C	2,4,6-Trichlorophenol	0.486	0.483	0.6	79	-0.02
43		2,4,5-Trichlorophenol	0.538	0.546	-1.5	81	-0.03
44	S	2-Fluorobiphenyl	1.593	1.541	3.3	79	-0.03

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.708	1.652	3.3	78	-0.03
46		2-Chloronaphthalene	1.303	1.258	3.5	78	-0.03
47		2-Nitroaniline	0.225	0.228	-1.3	80	-0.02
48		Acenaphthylene	2.084	2.046	1.8	80	-0.03
49		Dimethylphthalate	1.762	1.731	1.8	80	-0.03
50		2,6-Dinitrotoluene	0.324	0.331	-2.2	80	-0.03
51	C	Acenaphthene	1.365	1.349	1.2	80	-0.03
52		3-Nitroaniline	0.316	0.327	-3.5	81	-0.03
53	P	2,4-Dinitrophenol	0.115	0.112	2.6	80	-0.03
54		Dibenzofuran	2.093	2.055	1.8	80	-0.03
55	P	4-Nitrophenol	0.257	0.251	2.3	75	-0.02
56		2,4-Dinitrotoluene	0.417	0.438	-5.0	81	-0.03
57		Fluorene	1.700	1.675	1.5	81	-0.03
58		2,3,4,6-Tetrachlorophenol	0.525	0.536	-2.1	82	-0.03
59		Diethylphthalate	1.717	1.678	2.3	79	-0.03
60		4-Chlorophenyl-phenylether	1.040	1.014	2.5	79	-0.03
61		4-Nitroaniline	0.309	0.340	-10.0	82	-0.03
62		Azobenzene	1.101	1.033	6.2	76	-0.03
63	I	Phenanthrene-d10	1.000	1.000	0.0	82	-0.03
64		4,6-Dinitro-2-methylphenol	0.086	0.092	-7.0	87	-0.03
65	c	n-Nitrosodiphenylamine	0.664	0.643	3.2	80	-0.03
66		4-Bromophenyl-phenylether	0.289	0.283	2.1	81	-0.03
67		Hexachlorobenzene	0.296	0.303	-2.4	85	-0.03
68		Atrazine	0.258	0.244	5.4	78	-0.03
69	C	Pentachlorophenol	0.203	0.210	-3.4	81	-0.03
70		Phenanthrene	1.132	1.091	3.6	80	-0.03
71		Anthracene	1.142	1.098	3.9	80	-0.03
72		Carbazole	1.011	0.968	4.3	79	-0.03
73		Di-n-butylphthalate	1.123	1.075	4.3	79	-0.04
74	C	Fluoranthene	1.389	1.347	3.0	81	-0.03
75	I	Chrysene-d12	1.000	1.000	0.0	87	-0.05
76		Benzidine	0.617	0.567	8.1	78	-0.03
77		Pyrene	1.278	1.192	6.7	82	-0.03
78	S	Terphenyl-d14	0.875	0.835	4.6	84	-0.03
79		Butylbenzylphthalate	0.429	0.425	0.9	85	-0.05
80		Benzo(a)anthracene	1.272	1.225	3.7	85	-0.05
81		3,3'-Dichlorobenzidine	0.493	0.489	0.8	86	-0.05
82		Chrysene	1.204	1.139	5.4	83	-0.05
83		Bis(2-ethylhexyl)phthalate	0.621	0.611	1.6	85	-0.06
84	c	Di-n-octyl phthalate	1.046	1.038	0.8	85	-0.09
85		Indeno(1,2,3-cd)pyrene	1.544	1.556	-0.8	86	-0.13
86	I	Perylene-d12	1.000	1.000	0.0	89	-0.08
87		Benzo(b)fluoranthene	1.241	1.173	5.5	86	-0.07

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.242	1.175	5.4	84	-0.07
89 C	Benzo(a)pyrene	1.189	1.125	5.4	85	-0.07
90	Dibenzo(a,h)anthracene	1.228	1.184	3.6	86	-0.15
91	Benzo(g,h,i)perylene	1.450	1.404	3.2	86	-0.13

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