Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG093016\
Data File : BG024177.D
Acq On : 1 Oct 2016 2:23
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 01 04:08:55 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG092316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 23 15:35:28 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	105	0.00
2		1,4-Dioxane	0.442	0.520	-17.6	111	-0.01
3		Pyridine	1.562	1.635	-4.7	102	-0.01
4		n-Nitrosodimethylamine	0.792	0.730	7.8	95	0.00
5	S	2-Fluorophenol	1.154	1.170	-1.4	98	0.00
6		Aniline	2.728	2.398	12.1	92	0.00
7	S	Phenol-d6	1.942	1.785	8.1	95	0.00
8		2-Chlorophenol	1.382	1.216	12.0	92	-0.01
9		Benzaldehyde	1.058	1.079	-2.0	106	0.00
10	C	Phenol	2.042	1.915	6.2	97	0.00
11		bis(2-Chloroethyl)ether	1.610	1.438	10.7	91	-0.01
12		1,3-Dichlorobenzene	1.541	1.410	8.5	95	0.00
13	C	1,4-Dichlorobenzene	1.573	1.469	6.6	97	-0.01
14		1,2-Dichlorobenzene	1.531	1.427	6.8	95	-0.01
15		Benzyl Alcohol	1.880	1.591	15.4	90	0.00
16		2,2'-oxybis(1-Chloropropane	2.223	1.938	12.8	90	-0.02
17		2-Methylphenol	1.358	1.252	7.8	96	0.00
18		Hexachloroethane	0.663	0.627	5.4	97	-0.01
19	P	n-Nitroso-di-n-propylamine	1.908	1.589	16.7	88	-0.01
20		3+4-Methylphenols	1.972	1.807	8.4	97	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	105	-0.01
22		Acetophenone	0.577	0.538	6.8	95	-0.01
23	S	Nitrobenzene-d5	0.507	0.467	7.9	94	-0.01
24		Nitrobenzene	0.544	0.512	5.9	96	0.00
25		Isophorone	1.035	0.935	9.7	95	-0.01
26	C	2-Nitrophenol	0.193	0.173	10.4	91	0.00
27		2,4-Dimethylphenol	0.323	0.299	7.4	94	-0.01
28		bis(2-Chloroethoxy)methane	0.503	0.464	7.8	96	-0.01
29	C	2,4-Dichlorophenol	0.336	0.320	4.8	94	0.00
30		1,2,4-Trichlorobenzene	0.359	0.323	10.0	91	-0.01
31		Naphthalene	1.013	0.957	5.5	98	-0.01
32		Benzoic acid	0.241	0.190	21.2	79	-0.01
33		4-Chloroaniline	0.452	0.415	8.2	95	0.00
34	C	Hexachlorobutadiene	0.283	0.240	15.2	87	-0.01
35		Caprolactam	0.137	0.113	17.5	90	-0.01
36	C	4-Chloro-3-methylphenol	0.472	0.439	7.0	93	-0.01
37		2-Methylnaphthalene	0.776	0.697	10.2	93	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	106	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.584	0.533	8.7	91	-0.02
40	P	Hexachlorocyclopentadiene	0.220	0.199	9.5	84	-0.01
41	S	2,4,6-Tribromophenol	0.287	0.223	22.3	81	-0.01
42	C	2,4,6-Trichlorophenol	0.378	0.339	10.3	88	-0.01
43		2,4,5-Trichlorophenol	0.389	0.340	12.6	86	0.00
44	S	2-Fluorobiphenyl	1.190	1.110	6.7	95	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.368	1.297	5.2	96	-0.01
46		2-Chloronaphthalene	1.021	0.957	6.3	95	-0.02
47		2-Nitroaniline	0.475	0.421	11.4	89	-0.01
48		Acenaphthylene	1.734	1.630	6.0	98	-0.02
49		Dimethylphthalate	1.534	1.398	8.9	94	-0.01
50		2,6-Dinitrotoluene	0.322	0.294	8.7	92	0.00
51	C	Acenaphthene	1.051	0.987	6.1	96	-0.02
52		3-Nitroaniline	0.322	0.309	4.0	101	-0.01
53	P	2,4-Dinitrophenol	0.190	0.142	25.3#	75	0.00
54		Dibenzofuran	1.637	1.508	7.9	95	-0.01
55	P	4-Nitrophenol	0.225	0.188	16.4	86	0.00
56		2,4-Dinitrotoluene	0.490	0.444	9.4	95	-0.01
57		Fluorene	1.425	1.304	8.5	96	-0.01
58		2,3,4,6-Tetrachlorophenol	0.375	0.324	13.6	86	-0.01
59		Diethylphthalate	1.652	1.462	11.5	95	-0.01
60		4-Chlorophenyl-phenylether	0.763	0.663	13.1	92	-0.01
61		4-Nitroaniline	0.333	0.292	12.3	93	-0.01
62		Azobenzene	1.679	1.561	7.0	97	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	-0.01
64		4,6-Dinitro-2-methylphenol	0.139	0.115	17.3	79	-0.01
65	c	n-Nitrosodiphenylamine	0.547	0.538	1.6	94	-0.01
66		4-Bromophenyl-phenylether	0.232	0.206	11.2	87	-0.01
67		Hexachlorobenzene	0.261	0.224	14.2	84	-0.01
68		Atrazine	0.242	0.204	15.7	83	-0.01
69	C	Pentachlorophenol	0.124	0.096	22.6#	76	0.00
70		Phenanthrene	1.019	0.975	4.3	95	-0.02
71		Anthracene	1.046	0.978	6.5	93	-0.01
72		Carbazole	0.974	0.919	5.6	95	-0.01
73		Di-n-butylphthalate	1.275	1.109	13.0	89	-0.02
74	C	Fluoranthene	1.330	1.155	13.2	91	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	128	-0.02
76		Benzidine	0.566	0.395	30.2#	81	-0.01
77		Pyrene	1.397	1.099	21.3	92	-0.01
78	S	Terphenyl-d14	0.977	0.737	24.6	90	-0.01
79		Butylbenzylphthalate	0.514	0.450	12.5	108	-0.01
80		Benzo(a)anthracene	1.139	1.043	8.4	114	-0.02
81		3,3'-Dichlorobenzidine	0.408	0.396	2.9	118	-0.02
82		Chrysene	1.070	0.999	6.6	117	-0.01
83		Bis(2-ethylhexyl)phthalate	0.614	0.644	-4.9	125	-0.02
84	c	Di-n-octyl phthalate	1.060	1.133	-6.9	123	-0.02
85		Indeno(1,2,3-cd)pyrene	1.193	1.258	-5.4	122	-0.07
86	I	Perylene-d12	1.000	1.000	0.0	136	-0.03
87		Benzo(b)fluoranthene	1.164	1.064	8.6	120	-0.03

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88	Benzo(k)fluoranthene	1.156	1.101	4.8	125	-0.03
89 C	Benzo(a)pyrene	1.125	1.049	6.8	123	-0.03
90	Dibenzo(a,h)anthracene	1.123	1.032	8.1	123	-0.07
91	Benzo(g,h,i)perylene	1.132	1.057	6.6	125	-0.06

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