Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG041715\
Data File : BG016506.D
Acq On : 18 Apr 2015 4:16
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 20 12:22:05 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 14 14:00:56 2015
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	122	-0.01
2		1,4-Dioxane	0.576	0.521	9.5	114	-0.02
3		Pyridine	1.722	1.548	10.1	110	-0.02
4		n-Nitrosodimethylamine	0.512	0.467	8.8	111	-0.02
5	S	2-Fluorophenol	1.267	1.240	2.1	121	-0.01
6		Aniline	2.717	2.425	10.7	109	-0.01
7	S	Phenol-d6	1.902	1.785	6.2	114	-0.01
8		2-Chlorophenol	1.522	1.509	0.9	121	-0.01
9		Benzaldehyde	1.114	0.891	20.0	102	-0.02
10	C	Phenol	2.035	1.883	7.5	112	0.00
11		bis(2-Chloroethyl)ether	1.623	1.425	12.2	109	-0.02
12		1,3-Dichlorobenzene	1.537	1.526	0.7	124	-0.01
13	C	1,4-Dichlorobenzene	1.594	1.566	1.8	125	-0.02
14		1,2-Dichlorobenzene	1.525	1.477	3.1	121	-0.01
15		Benzyl Alcohol	1.326	1.160	12.5	104	-0.01
16		2,2'-oxybis(1-Chloropropane	1.828	1.426	22.0	96	-0.01
17		2-Methylphenol	1.365	1.256	8.0	111	0.00
18		Hexachloroethane	0.610	0.578	5.2	119	-0.02
19	P	n-Nitroso-di-n-propylamine	1.363	1.116	18.1	97	-0.02
20		3+4-Methylphenols	1.891	1.736	8.2	109	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	106	-0.02
22		Acetophenone	0.464	0.447	3.7	105	-0.01
23	S	Nitrobenzene-d5	0.345	0.334	3.2	104	-0.02
24		Nitrobenzene	0.369	0.347	6.0	102	-0.01
25		Isophorone	0.767	0.681	11.2	95	-0.01
26	C	2-Nitrophenol	0.172	0.198	-15.1	117	-0.01
27		2,4-Dimethylphenol	0.321	0.327	-1.9	109	-0.01
28		bis(2-Chloroethoxy)methane	0.438	0.402	8.2	100	-0.02
29	C	2,4-Dichlorophenol	0.254	0.279	-9.8	115	-0.01
30		1,2,4-Trichlorobenzene	0.265	0.282	-6.4	117	-0.02
31		Naphthalene	1.042	1.035	0.7	109	-0.01
32		Benzoic acid	0.182	0.135	25.8#	77	0.00
33		4-Chloroaniline	0.489	0.486	0.6	107	-0.01
34	C	Hexachlorobutadiene	0.116	0.124	-6.9	118	-0.02
35		Caprolactam	0.160	0.159	0.6	100	0.00
36	C	4-Chloro-3-methylphenol	0.335	0.335	0.0	105	0.00
37		2-Methylnaphthalene	0.750	0.734	2.1	107	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	99	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.440	0.461	-4.8	109	-0.01
40	P	Hexachlorocyclopentadiene	0.202	0.163	19.3	83	-0.01
41	S	2,4,6-Tribromophenol	0.135	0.144	-6.7	105	-0.01
42	C	2,4,6-Trichlorophenol	0.327	0.360	-10.1	110	-0.01
43		2,4,5-Trichlorophenol	0.350	0.383	-9.4	111	0.00
44	S	2-Fluorobiphenyl	1.175	1.176	-0.1	105	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.534	1.534	0.0	105	-0.01
46		2-Chloronaphthalene	1.169	1.188	-1.6	107	-0.01
47		2-Nitroaniline	0.377	0.381	-1.1	99	-0.01
48		Acenaphthylene	2.078	2.072	0.3	103	-0.02
49		Dimethylphthalate	1.466	1.449	1.2	102	-0.01
50		2,6-Dinitrotoluene	0.355	0.374	-5.4	106	-0.01
51	C	Acenaphthene	1.242	1.231	0.9	105	-0.01
52		3-Nitroaniline	0.453	0.475	-4.9	104	0.00
53	P	2,4-Dinitrophenol	0.167	0.107	35.9#	63	0.00
54		Dibenzofuran	1.723	1.716	0.4	105	-0.01
55	P	4-Nitrophenol	0.374	0.340	9.1	91	0.00
56		2,4-Dinitrotoluene	0.483	0.522	-8.1	108	0.00
57		Fluorene	1.520	1.522	-0.1	104	-0.01
58		2,3,4,6-Tetrachlorophenol	0.270	0.292	-8.1	110	0.00
59		Diethylphthalate	1.560	1.521	2.5	100	-0.01
60		4-Chlorophenyl-phenylether	0.622	0.603	3.1	101	-0.01
61		4-Nitroaniline	0.514	0.532	-3.5	103	0.00
62		Azobenzene	1.661	1.489	10.4	94	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	-0.02
64		4,6-Dinitro-2-methylphenol	0.119	0.118	0.8	93	0.00
65	c	n-Nitrosodiphenylamine	0.676	0.677	-0.1	104	-0.01
66		4-Bromophenyl-phenylether	0.160	0.160	0.0	104	-0.02
67		Hexachlorobenzene	0.166	0.162	2.4	102	-0.01
68		Atrazine	0.188	0.192	-2.1	105	0.00
69	C	Pentachlorophenol	0.101	0.080	20.8#	80	0.00
70		Phenanthrene	1.125	1.103	2.0	105	-0.01
71		Anthracene	1.115	1.111	0.4	105	-0.01
72		Carbazole	1.119	1.116	0.3	105	-0.01
73		Di-n-butylphthalate	1.369	1.369	0.0	103	-0.01
74	C	Fluoranthene	1.159	1.134	2.2	103	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	104	-0.01
76		Benzidine	0.848	0.681	19.7	83	0.00
77		Pyrene	1.393	1.354	2.8	104	-0.01
78	S	Terphenyl-d14	0.855	0.806	5.7	102	-0.01
79		Butylbenzylphthalate	0.684	0.745	-8.9	111	-0.01
80		Benzo(a)anthracene	1.182	1.172	0.8	108	0.00
81		3,3'-Dichlorobenzidine	0.389	0.399	-2.6	107	0.00
82		Chrysene	1.141	1.121	1.8	107	0.00
83		Bis(2-ethylhexyl)phthalate	0.926	1.023	-10.5	113	-0.01
84	c	Di-n-octyl phthalate	1.509	1.708	-13.2	113	-0.01
85		Indeno(1,2,3-cd)pyrene	1.186	1.297	-9.4	117	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	108	-0.01
87		Benzo(b)fluoranthene	1.171	1.173	-0.2	112	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.146	1.132	1.2	114	0.00
89 C	Benzo(a)pyrene	1.098	1.112	-1.3	113	-0.01
90	Dibenzo(a,h)anthracene	1.068	1.101	-3.1	116	-0.02
91	Benzo(g,h,i)perylene	1.067	1.116	-4.6	118	-0.02

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