Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG110419\
Data File : BG043498.D
Acq On : 5 Nov 2019 2:50
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 05 04:54:55 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 04 15:22:15 2019
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	135	0.00
2		1,4-Dioxane	0.425	0.402	5.4	122	0.00
3		Pyridine	1.073	1.067	0.6	115	0.00
4		n-Nitrosodimethylamine	0.541	0.510	5.7	122	0.00
5	S	2-Fluorophenol	1.091	1.085	0.5	129	0.00
6		Aniline	1.809	1.699	6.1	119	0.00
7	S	Phenol-d6	1.570	1.516	3.4	125	0.00
8		2-Chlorophenol	1.205	1.175	2.5	126	0.00
9		Benzaldehyde	0.772	0.801	-3.8	139	0.00
10	C	Phenol	1.435	1.368	4.7	121	0.00
11		bis(2-Chloroethyl)ether	1.109	1.023	7.8	118	0.00
12		1,3-Dichlorobenzene	1.510	1.500	0.7	130	0.00
13	C	1,4-Dichlorobenzene	1.527	1.476	3.3	125	0.00
14		1,2-Dichlorobenzene	1.491	1.416	5.0	123	0.00
15		Benzyl Alcohol	1.103	1.065	3.4	121	0.00
16		2,2'-oxybis(1-Chloropropane	1.613	1.420	12.0	111	0.00
17		2-Methylphenol	1.046	0.991	5.3	125	0.00
18		Hexachloroethane	0.551	0.525	4.7	124	0.00
19	P	n-Nitroso-di-n-propylamine	1.015	0.888	12.5	114	0.00
20		3+4-Methylphenols	1.434	1.324	7.7	120	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	124	0.00
22		Acetophenone	0.512	0.503	1.8	116	0.00
23	S	Nitrobenzene-d5	0.388	0.377	2.8	117	0.00
24		Nitrobenzene	0.370	0.364	1.6	117	0.00
25		Isophorone	0.709	0.650	8.3	109	0.00
26	C	2-Nitrophenol	0.178	0.181	-1.7	124	0.00
27		2,4-Dimethylphenol	0.256	0.257	-0.4	120	0.00
28		bis(2-Chloroethoxy)methane	0.391	0.366	6.4	108	0.00
29	C	2,4-Dichlorophenol	0.328	0.341	-4.0	122	0.00
30		1,2,4-Trichlorobenzene	0.413	0.422	-2.2	123	0.00
31		Naphthalene	1.015	1.008	0.7	120	0.00
32		Benzoic acid	0.179	0.172	3.9	128	0.00
33		4-Chloroaniline	0.416	0.417	-0.2	116	0.00
34	C	Hexachlorobutadiene	0.305	0.306	-0.3	122	0.00
35		Caprolactam	0.113	0.109	3.5	112	0.02
36	C	4-Chloro-3-methylphenol	0.357	0.347	2.8	115	0.00
37		2-Methylnaphthalene	0.779	0.751	3.6	115	0.00
38		1-Methylnaphthalene	0.741	0.716	3.4	115	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	118	0.00
40		1,2,4,5-Tetrachlorobenzene	0.740	0.759	-2.6	116	0.00
41	P	Hexachlorocyclopentadiene	0.389	0.131	66.3#	38#	0.00
42	S	2,4,6-Tribromophenol	0.381	0.363	4.7	106	0.00
43	C	2,4,6-Trichlorophenol	0.436	0.437	-0.2	111	0.00
44		2,4,5-Trichlorophenol	0.441	0.431	2.3	109	0.00

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Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.447	1.431	1.1	111	0.00
46		1,1'-Biphenyl	1.521	1.512	0.6	111	0.00
47		2-Chloronaphthalene	1.158	1.164	-0.5	114	0.00
48		2-Nitroaniline	0.346	0.339	2.0	107	0.00
49		Acenaphthylene	1.802	1.804	-0.1	112	0.00
50		Dimethylphthalate	1.549	1.435	7.4	105	0.00
51		2,6-Dinitrotoluene	0.337	0.327	3.0	108	0.00
52	C	Acenaphthene	1.099	1.078	1.9	111	0.00
53		3-Nitroaniline	0.325	0.321	1.2	110	0.00
54	P	2,4-Dinitrophenol	0.178	0.073	59.0#	46#	0.00
55		Dibenzofuran	1.782	1.730	2.9	109	0.00
56	P	4-Nitrophenol	0.218	0.202	7.3	102	0.00
57		2,4-Dinitrotoluene	0.481	0.462	4.0	109	0.00
58		Fluorene	1.494	1.439	3.7	107	0.00
59		2,3,4,6-Tetrachlorophenol	0.468	0.442	5.6	101	0.00
60		Diethylphthalate	1.557	1.427	8.3	103	0.00
61		4-Chlorophenyl-phenylether	0.872	0.831	4.7	108	0.00
62		4-Nitroaniline	0.336	0.348	-3.6	114	0.00
63		Azobenzene	1.260	1.181	6.3	105	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	117	0.00
65		4,6-Dinitro-2-methylphenol	0.118	0.058	50.8#	56	0.00
66	c	n-Nitrosodiphenylamine	0.565	0.540	4.4	106	0.00
67		4-Bromophenyl-phenylether	0.269	0.256	4.8	107	0.00
68		Hexachlorobenzene	0.309	0.293	5.2	107	0.00
69		Atrazine	0.196	0.174	11.2	101	0.00
70	C	Pentachlorophenol	0.141	0.123	12.8	94	0.00
71		Phenanthrene	1.024	0.977	4.6	106	0.00
72		Anthracene	1.025	0.986	3.8	105	0.00
73		Carbazole	0.928	0.894	3.7	107	0.00
74		Di-n-butylphthalate	1.126	1.083	3.8	106	0.00
75	C	Fluoranthene	1.335	1.240	7.1	101	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	112	0.00
77		Benzidine	0.403	0.428	-6.2	106	0.00
78		Pyrene	1.238	1.211	2.2	104	0.00
79	S	Terphenyl-d14	1.066	1.040	2.4	103	0.00
80		Butylbenzylphthalate	0.469	0.478	-1.9	109	0.00
81		Benzo(a)anthracene	1.253	1.261	-0.6	109	0.00
82		3,3'-Dichlorobenzidine	0.485	0.516	-6.4	113	0.00
83		Chrysene	1.245	1.241	0.3	106	0.00
84		Bis(2-ethylhexyl)phthalate	0.666	0.680	-2.1	110	0.00
85	c	Di-n-octyl phthalate	1.130	1.158	-2.5	111	0.00
86		Indeno(1,2,3-cd)pyrene	1.562	1.516	2.9	105	0.00
87	I	Perylene-d12	1.000	1.000	0.0	117	0.00

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88	Benzo(b)fluoranthene	1.133	1.125	0.7	111	0.00
89	Benzo(k)fluoranthene	1.140	1.098	3.7	107	0.00
90 C	Benzo(a)pyrene	1.082	1.072	0.9	111	0.00
91	Dibenzo(a,h)anthracene	1.106	1.041	5.9	105	0.00
92	Benzo(g,h,i)perylene	1.091	0.921	15.6	94	0.02

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