Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG110518\
Data File : BG037825.D
Acq On : 6 Nov 2018 5:38
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 06 06:29:50 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 18 14:06:42 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	96	0.00
2		1,4-Dioxane	0.607	0.574	5.4	94	0.00
3		Pyridine	1.529	1.413	7.6	90	0.00
4		n-Nitrosodimethylamine	0.545	0.517	5.1	93	0.00
5	S	2-Fluorophenol	1.196	1.142	4.5	92	0.00
6		Aniline	2.207	2.070	6.2	91	0.00
7	S	Phenol-d6	1.809	1.677	7.3	89	0.00
8		2-Chlorophenol	1.399	1.337	4.4	92	-0.01
9		Benzaldehyde	1.132	0.921	18.6	79	-0.01
10	C	Phenol	1.909	1.752	8.2	89	0.00
11		bis(2-Chloroethyl)ether	1.450	1.368	5.7	92	-0.01
12		1,3-Dichlorobenzene	1.558	1.536	1.4	97	-0.01
13	C	1,4-Dichlorobenzene	1.594	1.528	4.1	94	0.00
14		1,2-Dichlorobenzene	1.533	1.449	5.5	92	0.00
15		Benzyl Alcohol	1.337	1.226	8.3	87	0.00
16		2,2'-oxybis(1-Chloropropane	2.594	2.438	6.0	92	-0.01
17		2-Methylphenol	1.262	1.157	8.3	87	0.00
18		Hexachloroethane	0.543	0.549	-1.1	98	-0.01
19	P	n-Nitroso-di-n-propylamine	1.246	1.113	10.7	84	-0.01
20		3+4-Methylphenols	1.804	1.657	8.1	88	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	93	0.00
22		Acetophenone	0.502	0.467	7.0	87	0.00
23	S	Nitrobenzene-d5	0.357	0.354	0.8	92	0.00
24		Nitrobenzene	0.371	0.368	0.8	91	-0.01
25		Isophorone	0.652	0.610	6.4	85	-0.01
26	C	2-Nitrophenol	0.167	0.187	-12.0	101	-0.01
27		2,4-Dimethylphenol	0.298	0.282	5.4	87	0.00
28		bis(2-Chloroethoxy)methane	0.427	0.412	3.5	88	-0.01
29	C	2,4-Dichlorophenol	0.332	0.326	1.8	90	0.00
30		1,2,4-Trichlorobenzene	0.361	0.364	-0.8	95	-0.01
31		Naphthalene	0.982	0.967	1.5	93	0.00
32		Benzoic acid	0.194	0.169	12.9	79	-0.03
33		4-Chloroaniline	0.462	0.440	4.8	87	-0.01
34	C	Hexachlorobutadiene	0.214	0.223	-4.2	95	-0.01
35		Caprolactam	0.133	0.126	5.3	85	0.00
36	C	4-Chloro-3-methylphenol	0.375	0.353	5.9	87	0.00
37		2-Methylnaphthalene	0.716	0.703	1.8	90	-0.01
38		1-Methylnaphthalene	0.695	0.674	3.0	91	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	90	0.00
40		1,2,4,5-Tetrachlorobenzene	0.645	0.676	-4.8	95	-0.01
41	P	Hexachlorocyclopentadiene	0.308	0.323	-4.9	92	0.00
42	S	2,4,6-Tribromophenol	0.218	0.215	1.4	84	0.00
43	C	2,4,6-Trichlorophenol	0.431	0.438	-1.6	88	0.00
44		2,4,5-Trichlorophenol	0.469	0.462	1.5	86	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.326	1.331	-0.4	92	-0.01
46		1,1'-Biphenyl	1.656	1.665	-0.5	91	-0.01
47		2-Chloronaphthalene	1.253	1.268	-1.2	92	0.00
48		2-Nitroaniline	0.380	0.389	-2.4	90	0.00
49		Acenaphthylene	1.885	1.900	-0.8	90	0.00
50		Dimethylphthalate	1.547	1.522	1.6	88	0.00
51		2,6-Dinitrotoluene	0.342	0.346	-1.2	88	0.00
52	C	Acenaphthene	1.161	1.145	1.4	90	0.00
53		3-Nitroaniline	0.380	0.380	0.0	86	0.00
54	P	2,4-Dinitrophenol	0.100	0.062	38.0#	56	0.00
55		Dibenzofuran	1.923	1.876	2.4	89	-0.01
56	P	4-Nitrophenol	0.281	0.226	19.6	69	0.00
57		2,4-Dinitrotoluene	0.449	0.473	-5.3	90	0.00
58		Fluorene	1.440	1.400	2.8	89	-0.01
59		2,3,4,6-Tetrachlorophenol	0.402	0.381	5.2	84	-0.01
60		Diethylphthalate	1.588	1.547	2.6	87	-0.01
61		4-Chlorophenyl-phenylether	0.840	0.824	1.9	89	-0.01
62		4-Nitroaniline	0.378	0.373	1.3	85	0.00
63		Azobenzene	1.516	1.446	4.6	87	-0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	86	0.00
65		4,6-Dinitro-2-methylphenol	0.093	0.085	8.6	79	0.00
66	c	n-Nitrosodiphenylamine	0.602	0.603	-0.2	86	-0.01
67		4-Bromophenyl-phenylether	0.220	0.221	-0.5	87	0.00
68		Hexachlorobenzene	0.228	0.224	1.8	85	0.00
69		Atrazine	0.218	0.204	6.4	79	-0.01
70	C	Pentachlorophenol	0.152	0.127	16.4	70	0.00
71		Phenanthrene	1.016	0.999	1.7	86	0.00
72		Anthracene	0.998	0.990	0.8	86	0.00
73		Carbazole	0.998	0.993	0.5	86	0.00
74		Di-n-butylphthalate	1.110	1.119	-0.8	85	-0.01
75	C	Fluoranthene	1.153	1.140	1.1	86	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	87	0.00
77		Benzidine	0.680	0.463	31.9#	57	0.00
78		Pyrene	1.307	1.312	-0.4	87	0.00
79	S	Terphenyl-d14	0.936	0.943	-0.7	87	-0.01
80		Butylbenzylphthalate	0.535	0.572	-6.9	89	-0.01
81		Benzo(a)anthracene	1.183	1.199	-1.4	88	0.00
82		3,3'-Dichlorobenzidine	0.459	0.443	3.5	81	0.00
83		Chrysene	1.138	1.152	-1.2	88	0.00
84		Bis(2-ethylhexyl)phthalate	0.750	0.806	-7.5	89	-0.02
85	c	Di-n-octyl phthalate	1.223	1.316	-7.6	89	-0.03
86		Indeno(1,2,3-cd)pyrene	1.220	0.989	18.9	69	-0.04
87	I	Perylene-d12	1.000	1.000	0.0	85	-0.02

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88	Benzo(b)fluoranthene	1.096	1.110	-1.3	85	-0.01
89	Benzo(k)fluoranthene	1.074	1.157	-7.7	91	-0.01
90 C	Benzo(a)pyrene	1.042	1.043	-0.1	84	-0.02
91	Dibenzo(a,h)anthracene	0.961	0.776	19.3	67	-0.03
92	Benzo(g,h,i)perylene	0.972	0.836	14.0	72	-0.04

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