Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG061818\
Data File : BG035212.D
Acq On : 18 Jun 2018 10:43
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 18 16:35:41 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG060718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 08 17:16:28 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	121	-0.01
2		1,4-Dioxane	0.565	0.558	1.2	120	0.00
3		Pyridine	1.526	1.601	-4.9	119	0.00
4		n-Nitrosodimethylamine	0.655	0.548	16.3	95	-0.01
5	S	2-Fluorophenol	1.185	1.176	0.8	117	-0.01
6		Aniline	2.261	2.228	1.5	117	-0.01
7	S	Phenol-d6	1.749	1.754	-0.3	118	0.00
8		2-Chlorophenol	1.242	1.229	1.0	116	-0.01
9		Benzaldehyde	1.347	1.184	12.1	101	-0.01
10	C	Phenol	1.810	1.829	-1.0	120	-0.01
11		bis(2-Chloroethyl)ether	1.405	1.330	5.3	113	-0.01
12		1,3-Dichlorobenzene	1.482	1.448	2.3	120	-0.01
13	C	1,4-Dichlorobenzene	1.530	1.480	3.3	115	-0.01
14		1,2-Dichlorobenzene	1.437	1.396	2.9	115	-0.02
15		Benzyl Alcohol	1.657	1.543	6.9	110	-0.02
16		2,2'-oxybis(1-Chloropropane	1.398	1.258	10.0	109	-0.03
17		2-Methylphenol	1.193	1.197	-0.3	116	-0.01
18		Hexachloroethane	0.548	0.509	7.1	108	-0.01
19	P	n-Nitroso-di-n-propylamine	1.486	1.367	8.0	108	-0.02
20		3+4-Methylphenols	1.711	1.685	1.5	116	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	124	-0.02
22		Acetophenone	0.620	0.577	6.9	115	-0.02
23	S	Nitrobenzene-d5	0.421	0.434	-3.1	121	-0.02
24		Nitrobenzene	0.431	0.443	-2.8	122	-0.02
25		Isophorone	0.888	0.812	8.6	112	-0.02
26	C	2-Nitrophenol	0.149	0.187	-25.5#	155#	-0.02
27		2,4-Dimethylphenol	0.312	0.292	6.4	116	-0.02
28		bis(2-Chloroethoxy)methane	0.477	0.461	3.4	118	-0.02
29	C	2,4-Dichlorophenol	0.340	0.343	-0.9	122	-0.01
30		1,2,4-Trichlorobenzene	0.407	0.398	2.2	117	-0.01
31		Naphthalene	1.039	0.990	4.7	117	-0.01
32		Benzoic acid	0.209	0.256	-22.5	151#	-0.02
33		4-Chloroaniline	0.451	0.443	1.8	119	-0.01
34	C	Hexachlorobutadiene	0.317	0.303	4.4	118	-0.02
35		Caprolactam	0.126	0.166	-31.7#	164#	-0.02
36	C	4-Chloro-3-methylphenol	0.423	0.454	-7.3	130	0.00
37		2-Methylnaphthalene	0.788	0.770	2.3	118	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	134	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.745	0.722	3.1	126	-0.01
40	P	Hexachlorocyclopentadiene	0.467	0.425	9.0	116	-0.02
41	S	2,4,6-Tribromophenol	0.256	0.332	-29.7#	170#	-0.01
42	C	2,4,6-Trichlorophenol	0.446	0.475	-6.5	137	-0.01
43		2,4,5-Trichlorophenol	0.490	0.543	-10.8	147	-0.01
44	S	2-Fluorobiphenyl	1.538	1.416	7.9	122	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.608	1.519	5.5	124	-0.01
46		2-Chloronaphthalene	1.216	1.158	4.8	128	-0.02
47		2-Nitroaniline	0.359	0.438	-22.0	161#	-0.01
48		Acenaphthylene	2.039	1.978	3.0	129	-0.01
49		Dimethylphthalate	1.744	1.808	-3.7	139	-0.02
50		2,6-Dinitrotoluene	0.318	0.402	-26.4#	163#	-0.01
51	C	Acenaphthene	1.332	1.380	-3.6	136	-0.01
52		3-Nitroaniline	0.312	0.446	-42.9#	179#	0.00
53	P	2,4-Dinitrophenol	0.129	0.223	-72.9#	223#	-0.01
54		Dibenzofuran	1.995	2.020	-1.3	133	-0.01
55	P	4-Nitrophenol	0.263	0.404	-53.6#	191#	0.00
56		2,4-Dinitrotoluene	0.423	0.645	-52.5#	197#	-0.01
57		Fluorene	1.667	1.772	-6.3	139	-0.01
58		2,3,4,6-Tetrachlorophenol	0.476	0.584	-22.7	160#	-0.01
59		Diethylphthalate	1.779	1.988	-11.7	149	-0.03
60		4-Chlorophenyl-phenylether	0.951	1.003	-5.5	141	-0.03
61		4-Nitroaniline	0.335	0.526	-57.0#	198#	-0.01
62		Azobenzene	1.744	1.758	-0.8	134	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	180#	-0.01
64		4,6-Dinitro-2-methylphenol	0.091	0.123	-35.2#	240#	-0.01
65	c	n-Nitrosodiphenylamine	0.601	0.508	15.5	148	-0.01
66		4-Bromophenyl-phenylether	0.241	0.206	14.5	153#	-0.02
67		Hexachlorobenzene	0.245	0.219	10.6	160#	-0.01
68		Atrazine	0.256	0.278	-8.6	190#	-0.02
69	C	Pentachlorophenol	0.165	0.176	-6.7	185#	-0.01
70		Phenanthrene	1.016	0.944	7.1	164#	-0.02
71		Anthracene	1.018	0.968	4.9	166#	-0.01
72		Carbazole	0.944	1.024	-8.5	190#	-0.01
73		Di-n-butylphthalate	1.125	1.238	-10.0	193#	-0.03
74	C	Fluoranthene	1.322	1.476	-11.6	195#	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	176#	-0.02
76		Benzidine	0.744	0.575	22.7	132	-0.01
77		Pyrene	1.318	1.480	-12.3	195#	-0.01
78	S	Terphenyl-d14	0.955	1.067	-11.7	192#	-0.01
79		Butylbenzylphthalate	0.479	0.487	-1.7	170#	-0.03
80		Benzo(a)anthracene	1.278	1.210	5.3	165#	-0.02
81		3,3'-Dichlorobenzidine	0.481	0.438	8.9	154#	-0.02
82		Chrysene	1.170	1.100	6.0	164#	-0.02
83		Bis(2-ethylhexyl)phthalate	0.676	0.590	12.7	149	-0.04
84	c	Di-n-octyl phthalate	1.161	1.055	9.1	154#	-0.05
85		Indeno(1,2,3-cd)pyrene	1.370	1.287	6.1	161#	-0.06
86	I	Perylene-d12	1.000	1.000	0.0	170#	-0.04
87		Benzo(b)fluoranthene	1.232	1.163	5.6	156#	-0.03

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88	Benzo(k)fluoranthene	1.205	1.129	6.3	162#	-0.03
89 C	Benzo(a)pyrene	1.159	1.080	6.8	157#	-0.04
90	Dibenzo(a,h)anthracene	1.144	1.075	6.0	159#	-0.07
91	Benzo(g,h,i)perylene	1.120	1.064	5.0	160#	-0.05

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