Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG062518\
Data File : BG035371.D
Acq On : 25 Jun 2018 10:28
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 25 11:09:18 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	147	-0.02
2		1,4-Dioxane	0.565	0.547	3.2	143	-0.01
3		Pyridine	1.526	1.481	2.9	133	-0.01
4		n-Nitrosodimethylamine	0.655	0.570	13.0	120	-0.02
5	S	2-Fluorophenol	1.185	1.160	2.1	140	-0.02
6		Aniline	2.261	2.128	5.9	136	-0.02
7	S	Phenol-d6	1.749	1.718	1.8	140	-0.01
8		2-Chlorophenol	1.242	1.174	5.5	135	-0.02
9		Benzaldehyde	1.347	1.185	12.0	122	-0.02
10	C	Phenol	1.810	1.720	5.0	137	-0.01
11		bis(2-Chloroethyl)ether	1.405	1.299	7.5	134	-0.02
12		1,3-Dichlorobenzene	1.482	1.416	4.5	142	-0.02
13	C	1,4-Dichlorobenzene	1.530	1.455	4.9	138	-0.02
14		1,2-Dichlorobenzene	1.437	1.378	4.1	138	-0.02
15		Benzyl Alcohol	1.657	1.435	13.4	124	-0.02
16		2,2'-oxybis(1-Chloropropane	1.398	1.153	17.5	121	-0.03
17		2-Methylphenol	1.193	1.155	3.2	136	-0.02
18		Hexachloroethane	0.548	0.541	1.3	140	-0.02
19	P	n-Nitroso-di-n-propylamine	1.486	1.252	15.7	121	-0.02
20		3+4-Methylphenols	1.711	1.617	5.5	135	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	147	-0.02
22		Acetophenone	0.620	0.563	9.2	132	-0.02
23	S	Nitrobenzene-d5	0.421	0.423	-0.5	140	-0.02
24		Nitrobenzene	0.431	0.424	1.6	138	-0.02
25		Isophorone	0.888	0.754	15.1	123	-0.02
26	C	2-Nitrophenol	0.149	0.169	-13.4	166#	-0.02
27		2,4-Dimethylphenol	0.312	0.280	10.3	132	-0.02
28		bis(2-Chloroethoxy)methane	0.477	0.428	10.3	130	-0.02
29	C	2,4-Dichlorophenol	0.340	0.335	1.5	140	-0.02
30		1,2,4-Trichlorobenzene	0.407	0.398	2.2	138	-0.02
31		Naphthalene	1.039	0.951	8.5	133	-0.02
32		Benzoic acid	0.209	0.209	0.0	145	0.00
33		4-Chloroaniline	0.451	0.417	7.5	132	-0.02
34	C	Hexachlorobutadiene	0.317	0.310	2.2	142	-0.03
35		Caprolactam	0.126	0.117	7.1	136	-0.02
36	C	4-Chloro-3-methylphenol	0.423	0.412	2.6	140	-0.01
37		2-Methylnaphthalene	0.788	0.728	7.6	132	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	146	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.745	0.744	0.1	142	-0.02
40	P	Hexachlorocyclopentadiene	0.467	0.419	10.3	124	-0.03
41	S	2,4,6-Tribromophenol	0.256	0.274	-7.0	153#	-0.02
42	C	2,4,6-Trichlorophenol	0.446	0.472	-5.8	148	-0.02
43		2,4,5-Trichlorophenol	0.490	0.515	-5.1	152#	-0.02
44	S	2-Fluorobiphenyl	1.538	1.433	6.8	135	-0.03

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.608	1.501	6.7	134	-0.02
46		2-Chloronaphthalene	1.216	1.151	5.3	138	-0.02
47		2-Nitroaniline	0.359	0.384	-7.0	154#	-0.02
48		Acenaphthylene	2.039	1.892	7.2	134	-0.02
49		Dimethylphthalate	1.744	1.583	9.2	132	-0.03
50		2,6-Dinitrotoluene	0.318	0.347	-9.1	154#	-0.02
51	C	Acenaphthene	1.332	1.292	3.0	139	-0.02
52		3-Nitroaniline	0.312	0.351	-12.5	153#	-0.01
53	P	2,4-Dinitrophenol	0.129	0.152	-17.8	165#	-0.02
54		Dibenzofuran	1.995	1.876	6.0	135	-0.02
55	P	4-Nitrophenol	0.263	0.250	4.9	129	0.00
56		2,4-Dinitrotoluene	0.423	0.499	-18.0	166#	-0.02
57		Fluorene	1.667	1.570	5.8	135	-0.02
58		2,3,4,6-Tetrachlorophenol	0.476	0.501	-5.3	150	-0.02
59		Diethylphthalate	1.779	1.619	9.0	133	-0.03
60		4-Chlorophenyl-phenylether	0.951	0.936	1.6	143	-0.03
61		4-Nitroaniline	0.335	0.366	-9.3	150#	-0.02
62		Azobenzene	1.744	1.483	15.0	123	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	152#	-0.02
64		4,6-Dinitro-2-methylphenol	0.091	0.108	-18.7	177#	-0.02
65	c	n-Nitrosodiphenylamine	0.601	0.550	8.5	134	-0.02
66		4-Bromophenyl-phenylether	0.241	0.233	3.3	145	-0.03
67		Hexachlorobenzene	0.245	0.235	4.1	145	-0.02
68		Atrazine	0.256	0.233	9.0	134	-0.03
69	C	Pentachlorophenol	0.165	0.163	1.2	144	-0.02
70		Phenanthrene	1.016	0.959	5.6	140	-0.02
71		Anthracene	1.018	0.950	6.7	137	-0.02
72		Carbazole	0.944	0.890	5.7	139	-0.02
73		Di-n-butylphthalate	1.125	0.993	11.7	130	-0.04
74	C	Fluoranthene	1.322	1.265	4.3	141	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	165#	-0.03
76		Benzidine	0.744	0.447	39.9#	97	-0.02
77		Pyrene	1.318	1.148	12.9	143	-0.02
78	S	Terphenyl-d14	0.955	0.828	13.3	140	-0.02
79		Butylbenzylphthalate	0.479	0.400	16.5	131	-0.03
80		Benzo(a)anthracene	1.278	1.141	10.7	147	-0.03
81		3,3'-Dichlorobenzidine	0.481	0.426	11.4	141	-0.03
82		Chrysene	1.170	1.063	9.1	149	-0.03
83		Bis(2-ethylhexyl)phthalate	0.676	0.558	17.5	132	-0.05
84	c	Di-n-octyl phthalate	1.161	0.948	18.3	130	-0.07
85		Indeno(1,2,3-cd)pyrene	1.370	1.255	8.4	148	-0.09
86	I	Perylene-d12	1.000	1.000	0.0	160#	-0.05
87		Benzo(b)fluoranthene	1.232	1.124	8.8	142	-0.05

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88	Benzo(k)fluoranthene	1.205	1.117	7.3	151#	-0.05
89 C	Benzo(a)pyrene	1.159	1.058	8.7	144	-0.05
90	Dibenzo(a,h)anthracene	1.144	1.057	7.6	147	-0.09
91	Benzo(g,h,i)perylene	1.120	1.039	7.2	147	-0.09

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