Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG070218\
Data File : BG035551.D
Acq On : 2 Jul 2018 15:04
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 02 15:39:03 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	91	0.00
2		1,4-Dioxane	0.565	0.586	-3.7	95	0.00
3		Pyridine	1.526	1.592	-4.3	89	0.00
4		n-Nitrosodimethylamine	0.655	0.609	7.0	80	0.00
5	S	2-Fluorophenol	1.185	1.190	-0.4	89	0.00
6		Aniline	2.261	2.206	2.4	87	0.00
7	S	Phenol-d6	1.749	1.787	-2.2	90	0.00
8		2-Chlorophenol	1.242	1.244	-0.2	89	0.00
9		Benzaldehyde	1.347	1.256	6.8	81	0.00
10	C	Phenol	1.810	1.825	-0.8	90	0.00
11		bis(2-Chloroethyl)ether	1.405	1.405	0.0	90	0.00
12		1,3-Dichlorobenzene	1.482	1.494	-0.8	93	0.00
13	C	1,4-Dichlorobenzene	1.530	1.532	-0.1	90	0.00
14		1,2-Dichlorobenzene	1.437	1.454	-1.2	91	-0.02
15		Benzyl Alcohol	1.657	1.557	6.0	84	0.00
16		2,2'-oxybis(1-Chloropropane	1.398	1.307	6.5	85	-0.02
17		2-Methylphenol	1.193	1.209	-1.3	88	0.00
18		Hexachloroethane	0.548	0.541	1.3	87	-0.02
19	P	n-Nitroso-di-n-propylamine	1.486	1.412	5.0	84	-0.02
20		3+4-Methylphenols	1.711	1.696	0.9	88	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	91	-0.02
22		Acetophenone	0.620	0.614	1.0	89	-0.02
23	S	Nitrobenzene-d5	0.421	0.474	-12.6	97	0.00
24		Nitrobenzene	0.431	0.464	-7.7	94	0.00
25		Isophorone	0.888	0.828	6.8	84	-0.02
26	C	2-Nitrophenol	0.149	0.182	-22.1#	110	0.00
27		2,4-Dimethylphenol	0.312	0.300	3.8	87	0.00
28		bis(2-Chloroethoxy)methane	0.477	0.471	1.3	88	-0.02
29	C	2,4-Dichlorophenol	0.340	0.362	-6.5	94	0.00
30		1,2,4-Trichlorobenzene	0.407	0.425	-4.4	91	-0.02
31		Naphthalene	1.039	1.073	-3.3	93	0.00
32		Benzoic acid	0.209	0.171	18.2	74	0.00
33		4-Chloroaniline	0.451	0.458	-1.6	90	0.00
34	C	Hexachlorobutadiene	0.317	0.340	-7.3	97	-0.02
35		Caprolactam	0.126	0.129	-2.4	93	0.00
36	C	4-Chloro-3-methylphenol	0.423	0.448	-5.9	94	0.00
37		2-Methylnaphthalene	0.788	0.818	-3.8	92	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
39		1,2,4,5-Tetrachlorobenzene	0.745	0.792	-6.3	99	0.00
40	P	Hexachlorocyclopentadiene	0.467	0.388	16.9	76	-0.02
41	S	2,4,6-Tribromophenol	0.256	0.301	-17.6	111	0.00
42	C	2,4,6-Trichlorophenol	0.446	0.481	-7.8	100	0.00
43		2,4,5-Trichlorophenol	0.490	0.520	-6.1	102	0.00
44	S	2-Fluorobiphenyl	1.538	1.533	0.3	95	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.608	1.602	0.4	94	-0.02
46		2-Chloronaphthalene	1.216	1.218	-0.2	97	-0.02
47		2-Nitroaniline	0.359	0.413	-15.0	110	0.00
48		Acenaphthylene	2.039	2.031	0.4	95	0.00
49		Dimethylphthalate	1.744	1.694	2.9	93	-0.02
50		2,6-Dinitrotoluene	0.318	0.370	-16.4	108	0.00
51	C	Acenaphthene	1.332	1.279	4.0	91	0.00
52		3-Nitroaniline	0.312	0.365	-17.0	105	0.00
53	P	2,4-Dinitrophenol	0.129	0.131	-1.6	95	0.00
54		Dibenzofuran	1.995	2.038	-2.2	97	-0.02
55	P	4-Nitrophenol	0.263	0.279	-6.1	95	0.00
56		2,4-Dinitrotoluene	0.423	0.531	-25.5#	117	0.00
57		Fluorene	1.667	1.723	-3.4	98	0.00
58		2,3,4,6-Tetrachlorophenol	0.476	0.509	-6.9	101	0.00
59		Diethylphthalate	1.779	1.722	3.2	93	-0.02
60		4-Chlorophenyl-phenylether	0.951	1.005	-5.7	102	-0.02
61		4-Nitroaniline	0.335	0.399	-19.1	108	0.00
62		Azobenzene	1.744	1.690	3.1	93	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	105	-0.02
64		4,6-Dinitro-2-methylphenol	0.091	0.108	-18.7	122	0.00
65	c	n-Nitrosodiphenylamine	0.601	0.571	5.0	97	-0.02
66		4-Bromophenyl-phenylether	0.241	0.245	-1.7	106	-0.02
67		Hexachlorobenzene	0.245	0.253	-3.3	108	0.00
68		Atrazine	0.256	0.246	3.9	98	-0.02
69	C	Pentachlorophenol	0.165	0.148	10.3	91	0.00
70		Phenanthrene	1.016	1.015	0.1	103	-0.02
71		Anthracene	1.018	1.016	0.2	102	-0.02
72		Carbazole	0.944	0.951	-0.7	103	0.00
73		Di-n-butylphthalate	1.125	1.071	4.8	97	-0.03
74	C	Fluoranthene	1.322	1.351	-2.2	104	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	110	-0.02
76		Benzidine	0.744	0.683	8.2	98	-0.02
77		Pyrene	1.318	1.269	3.7	105	-0.02
78	S	Terphenyl-d14	0.955	0.941	1.5	106	-0.02
79		Butylbenzylphthalate	0.479	0.458	4.4	100	-0.03
80		Benzo(a)anthracene	1.278	1.288	-0.8	110	-0.02
81		3,3'-Dichlorobenzidine	0.481	0.466	3.1	102	-0.02
82		Chrysene	1.170	1.187	-1.5	111	-0.02
83		Bis(2-ethylhexyl)phthalate	0.676	0.613	9.3	97	-0.05
84	c	Di-n-octyl phthalate	1.161	1.038	10.6	95	-0.06
85		Indeno(1,2,3-cd)pyrene	1.370	1.376	-0.4	108	-0.06
86	I	Perylene-d12	1.000	1.000	0.0	110	-0.04
87		Benzo(b)fluoranthene	1.232	1.211	1.7	105	-0.03

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88	Benzo(k)fluoranthene	1.205	1.181	2.0	110	-0.03
89 C	Benzo(a)pyrene	1.159	1.139	1.7	107	-0.04
90	Dibenzo(a,h)anthracene	1.144	1.141	0.3	109	-0.07
91	Benzo(g,h,i)perylene	1.120	1.110	0.9	108	-0.06

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