Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG061418\
Data File : BG035137.D
Acq On : 14 Jun 2018 14:39
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 15 15:38:56 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	180#	0.00
2		1,4-Dioxane	0.565	0.534	5.5	171#	0.00
3		Pyridine	1.526	1.583	-3.7	175#	0.00
4		n-Nitrosodimethylamine	0.655	0.628	4.1	163#	-0.01
5	S	2-Fluorophenol	1.185	1.208	-1.9	179#	0.00
6		Aniline	2.261	2.330	-3.1	182#	0.00
7	S	Phenol-d6	1.749	1.879	-7.4	188#	0.00
8		2-Chlorophenol	1.242	1.277	-2.8	180#	0.00
9		Benzaldehyde	1.347	1.273	5.5	161#	0.00
10	C	Phenol	1.810	1.910	-5.5	186#	0.00
11		bis(2-Chloroethyl)ether	1.405	1.408	-0.2	178#	0.00
12		1,3-Dichlorobenzene	1.482	1.518	-2.4	187#	0.00
13	C	1,4-Dichlorobenzene	1.530	1.560	-2.0	181#	0.00
14		1,2-Dichlorobenzene	1.437	1.501	-4.5	185#	-0.01
15		Benzyl Alcohol	1.657	1.699	-2.5	180#	-0.01
16		2,2'-oxybis(1-Chloropropane	1.398	1.153	17.5	148	-0.01
17		2-Methylphenol	1.193	1.282	-7.5	185#	0.00
18		Hexachloroethane	0.548	0.566	-3.3	179#	-0.01
19	P	n-Nitroso-di-n-propylamine	1.486	1.480	0.4	175#	0.00
20		3+4-Methylphenols	1.711	1.861	-8.8	191#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	188#	-0.01
22		Acetophenone	0.620	0.609	1.8	183#	-0.01
23	S	Nitrobenzene-d5	0.421	0.471	-11.9	199#	0.00
24		Nitrobenzene	0.431	0.467	-8.4	195#	0.00
25		Isophorone	0.888	0.861	3.0	180#	0.00
26	C	2-Nitrophenol	0.149	0.197	-32.2#	247#	-0.01
27		2,4-Dimethylphenol	0.312	0.307	1.6	185#	0.00
28		bis(2-Chloroethoxy)methane	0.477	0.477	0.0	185#	-0.01
29	C	2,4-Dichlorophenol	0.340	0.376	-10.6	202#	0.00
30		1,2,4-Trichlorobenzene	0.407	0.419	-2.9	186#	-0.01
31		Naphthalene	1.039	1.048	-0.9	188#	0.00
32		Benzoic acid	0.209	0.224	-7.2	200#	0.02
33		4-Chloroaniline	0.451	0.474	-5.1	193#	0.00
34	C	Hexachlorobutadiene	0.317	0.325	-2.5	191#	-0.01
35		Caprolactam	0.126	0.139	-10.3	208#	0.01
36	C	4-Chloro-3-methylphenol	0.423	0.454	-7.3	198#	0.00
37		2-Methylnaphthalene	0.788	0.817	-3.7	190#	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	199#	0.00
39		1,2,4,5-Tetrachlorobenzene	0.745	0.786	-5.5	204#	0.00
40	P	Hexachlorocyclopentadiene	0.467	0.435	6.9	175#	-0.01
41	S	2,4,6-Tribromophenol	0.256	0.289	-12.9	220#	0.00
42	C	2,4,6-Trichlorophenol	0.446	0.500	-12.1	214#	0.00
43		2,4,5-Trichlorophenol	0.490	0.541	-10.4	218#	0.00
44	S	2-Fluorobiphenyl	1.538	1.499	2.5	192#	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.608	1.610	-0.1	196#	0.00
46		2-Chloronaphthalene	1.216	1.214	0.2	199#	-0.01
47		2-Nitroaniline	0.359	0.430	-19.8	235#	0.00
48		Acenaphthylene	2.039	2.057	-0.9	198#	0.00
49		Dimethylphthalate	1.744	1.717	1.5	195#	-0.01
50		2,6-Dinitrotoluene	0.318	0.378	-18.9	228#	0.00
51	C	Acenaphthene	1.332	1.353	-1.6	198#	0.00
52		3-Nitroaniline	0.312	0.380	-21.8	226#	0.00
53	P	2,4-Dinitrophenol	0.129	0.133	-3.1	197#	0.00
54		Dibenzofuran	1.995	2.014	-1.0	197#	0.00
55	P	4-Nitrophenol	0.263	0.282	-7.2	198#	0.00
56		2,4-Dinitrotoluene	0.423	0.534	-26.2#	241#	0.00
57		Fluorene	1.667	1.686	-1.1	197#	0.00
58		2,3,4,6-Tetrachlorophenol	0.476	0.524	-10.1	214#	0.00
59		Diethylphthalate	1.779	1.729	2.8	193#	-0.01
60		4-Chlorophenyl-phenylether	0.951	0.998	-4.9	208#	-0.02
61		4-Nitroaniline	0.335	0.391	-16.7	218#	0.00
62		Azobenzene	1.744	1.656	5.0	188#	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	197#	0.00
64		4,6-Dinitro-2-methylphenol	0.091	0.108	-18.7	229#	0.00
65	c	n-Nitrosodiphenylamine	0.601	0.618	-2.8	196#	0.00
66		4-Bromophenyl-phenylether	0.241	0.263	-9.1	213#	-0.01
67		Hexachlorobenzene	0.245	0.262	-6.9	210#	0.00
68		Atrazine	0.256	0.267	-4.3	199#	-0.01
69	C	Pentachlorophenol	0.165	0.176	-6.7	201#	0.00
70		Phenanthrene	1.016	1.040	-2.4	197#	-0.01
71		Anthracene	1.018	1.047	-2.8	196#	0.00
72		Carbazole	0.944	0.953	-1.0	193#	0.00
73		Di-n-butylphthalate	1.125	1.111	1.2	189#	-0.02
74	C	Fluoranthene	1.322	1.334	-0.9	192#	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	202#	-0.01
76		Benzidine	0.744	0.674	9.4	178#	-0.01
77		Pyrene	1.318	1.270	3.6	193#	0.00
78	S	Terphenyl-d14	0.955	0.914	4.3	190#	0.00
79		Butylbenzylphthalate	0.479	0.480	-0.2	193#	-0.01
80		Benzo(a)anthracene	1.278	1.277	0.1	201#	0.00
81		3,3'-Dichlorobenzidine	0.481	0.474	1.5	192#	-0.01
82		Chrysene	1.170	1.185	-1.3	203#	-0.01
83		Bis(2-ethylhexyl)phthalate	0.676	0.656	3.0	190#	-0.02
84	c	Di-n-octyl phthalate	1.161	1.137	2.1	191#	-0.03
85		Indeno(1,2,3-cd)pyrene	1.370	1.355	1.1	196#	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	199#	-0.01
87		Benzo(b)fluoranthene	1.232	1.277	-3.7	201#	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.205	1.212	-0.6	204#	-0.01
89 C	Benzo(a)pyrene	1.159	1.195	-3.1	203#	-0.02
90	Dibenzo(a,h)anthracene	1.144	1.127	1.5	195#	-0.03
91	Benzo(g,h,i)perylene	1.120	1.105	1.3	195#	-0.03

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