Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF103015\
Data File : BF082586.D
Acq On : 30 Oct 2015 22:01
Operator : UM/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF103015

Quant Time: Oct 30 23:50:38 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF103015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Oct 31 00:17:39 2015
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	98	0.00
2		1,4-Dioxane	0.614	1.081	-76.1#	178#	0.00
3		Pyridine	1.800	3.299	-83.3#	176#	0.00
4		n-Nitrosodimethylamine	1.013	1.843	-81.9#	184#	0.01
5	S	2-Fluorophenol	1.328	2.154	-62.2#	153#	0.00
6		Aniline	2.456	4.237	-72.5#	169#	0.01
7	S	Phenol-d6	1.764	2.952	-67.3#	173#	0.01
8		2-Chlorophenol	1.437	2.285	-59.0#	162#	0.00
9		Benzaldehyde	1.204	1.945	-61.5#	171#	0.01
10	C	Phenol	1.999	3.162	-58.2#	156#	0.01
11		bis(2-Chloroethyl)ether	1.475	2.357	-59.8#	180#	0.01
12		1,3-Dichlorobenzene	1.638	3.066	-87.2#	200#	0.01
13	C	1,4-Dichlorobenzene	1.625	2.746	-69.0#	171#	0.00
14		1,2-Dichlorobenzene	1.508	2.365	-56.8#	148	0.00
15		Benzyl Alcohol	1.645	3.132	-90.4#	182#	0.01
16		2,2'-oxybis(1-Chloropropane	2.344	4.151	-77.1#	179#	0.01
17		2-Methylphenol	1.222	2.201	-80.1#	175#	0.00
18		Hexachloroethane	0.634	1.078	-70.0#	168#	0.00
19	P	n-Nitroso-di-n-propylamine	1.337	2.364	-76.8#	185#	0.02
20		3+4-Methylphenols	1.588	2.826	-78.0#	192#	0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	104	0.01
22		Acetophenone	0.589	1.051	-78.4#	170#	0.00
23	S	Nitrobenzene-d5	0.490	0.854	-74.3#	178#	0.01
24		Nitrobenzene	0.495	0.925	-86.9#	179#	0.01
25		Isophorone	0.901	1.582	-75.6#	183#	0.01
26	C	2-Nitrophenol	0.181	0.354	-95.6#	203#	0.01
27		2,4-Dimethylphenol	0.353	0.680	-92.6#	192#	0.01
28		bis(2-Chloroethoxy)methane	0.493	0.759	-54.0#	160#	0.00
29	C	2,4-Dichlorophenol	0.322	0.526	-63.4#	168#	0.00
30		1,2,4-Trichlorobenzene	0.357	0.651	-82.4#	195#	0.01
31		Naphthalene	1.082	1.931	-78.5#	199#	0.01
32		Benzoic acid	0.256	0.491	-91.8#	189#	0.05
33		4-Chloroaniline	0.462	0.864	-87.0#	211#	0.01
34	C	Hexachlorobutadiene	0.227	0.431	-89.9#	197#	0.01
35		Caprolactam	0.099	0.170	-71.7#	181#	0.03
36	C	4-Chloro-3-methylphenol	0.379	0.625	-64.9#	182#	0.01
37		2-Methylnaphthalene	0.701	1.094	-56.1#	150	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
39		1,2,4,5-Tetrachlorobenzene	0.667	1.157	-73.5#	164#	0.00
40	P	Hexachlorocyclopentadiene	0.420	0.807	-92.1#	181#	0.00
41	S	2,4,6-Tribromophenol	0.219	0.471	-115.1#	218#	0.01
42	C	2,4,6-Trichlorophenol	0.482	0.902	-87.1#	166#	0.00
43		2,4,5-Trichlorophenol	0.480	0.922	-92.1#	203#	0.01
44	S	2-Fluorobiphenyl	1.409	2.707	-92.1#	188#	0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.689	2.654	-57.1#	140	0.00
46		2-Chloronaphthalene	1.319	2.226	-68.8#	155#	0.00
47		2-Nitroaniline	0.513	0.976	-90.3#	194#	0.01
48		Acenaphthylene	2.110	3.987	-89.0#	183#	0.01
49		Dimethylphthalate	1.633	2.542	-55.7#	163#	0.01
50		2,6-Dinitrotoluene	0.339	0.723	-113.3#	187#	0.01
51	C	Acenaphthene	1.339	2.657	-98.4#	202#	0.01
52		3-Nitroaniline	0.377	0.705	-87.0#	178#	0.01
53	P	2,4-Dinitrophenol	0.152	0.341	-124.3#	223#	0.01
54		Dibenzofuran	1.818	3.542	-94.8#	205#	0.01
55	P	4-Nitrophenol	0.283	0.607	-114.5#	173#	0.01
56		2,4-Dinitrotoluene	0.458	0.911	-98.9#	175#	0.01
57		Fluorene	1.466	2.873	-96.0#	201#	0.01
58		2,3,4,6-Tetrachlorophenol	0.406	0.803	-97.8#	203#	0.01
59		Diethylphthalate	1.663	2.839	-70.7#	162#	0.00
60		4-Chlorophenyl-phenylether	0.713	1.279	-79.4#	166#	0.00
61		4-Nitroaniline	0.359	0.708	-97.2#	192#	0.02
62		Azobenzene	1.875	3.375	-80.0#	166#	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	94	0.00
64		4,6-Dinitro-2-methylphenol	0.121	0.272	-124.8#	182#	0.01
65	c	n-Nitrosodiphenylamine	0.688	1.303	-89.4#	170#	0.01
66		4-Bromophenyl-phenylether	0.227	0.448	-97.4#	193#	0.00
67		Hexachlorobenzene	0.251	0.502	-100.0#	204#	0.01
68		Atrazine	0.225	0.434	-92.9#	198#	0.01
69	C	Pentachlorophenol	0.166	0.317	-91.0#	161#	0.00
70		Phenanthrene	1.128	2.156	-91.1#	193#	0.01
71		Anthracene	1.136	1.782	-56.9#	136	0.00
72		Carbazole	0.989	1.634	-65.2#	142	0.00
73		Di-n-butylphthalate	1.259	2.049	-62.7#	144	0.00
74	C	Fluoranthene	1.153	2.101	-82.2#	179#	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	83	0.01
76		Benzidine	0.902	1.808	-100.4#	145	0.00
77		Pyrene	1.481	3.292	-122.3#	180#	0.01
78	S	Terphenyl-d14	0.914	2.045	-123.7#	194#	0.01
79		Butylbenzylphthalate	0.649	1.334	-105.5#	153#	0.00
80		Benzo(a)anthracene	1.270	2.583	-103.4#	161#	0.01
81		3,3'-Dichlorobenzidine	0.438	0.891	-103.4#	165#	0.01
82		Chrysene	1.152	2.271	-97.1#	136	0.00
83		Bis(2-ethylhexyl)phthalate	0.841	1.895	-125.3#	163#	0.00
84	c	Di-n-octyl phthalate	1.312	2.840	-116.5#	158#	-0.01
85		Indeno(1,2,3-cd)pyrene	0.950	2.181	-129.6#	180#	0.00
86	I	Perylene-d12	1.000	1.000	0.0	86	-0.01
87		Benzo(b)fluoranthene	1.379	2.177	-57.9#	147	-0.01

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	Compound	AvgRF	CCRF	%Dev Area% Dev(min)
88	Benzo(k)fluoranthene	1.051	2.264	-115.4# 162# -0.01
89 C	Benzo(a)pyrene	1.086	2.084	-91.9# 161# 0.00
90	Dibenzo(a,h)anthracene	1.013	2.120	-109.3# 180# 0.00
91	Benzo(g,h,i)perylene	0.982	2.059	-109.7# 183# 0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 13