Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF010915\
Data File : BF076796.D
Acq On : 9 Jan 2015 21:05
Operator : IZ / TP
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF010915

Quant Time: Jan 12 11:20:13 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 12 10:43:48 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	116	0.00
2		1,4-Dioxane	0.511	0.524	-2.5	111	0.00
3		Pyridine	1.360	1.439	-5.8	137	-0.01
4		n-Nitrosodimethylamine	0.684	0.733	-7.2	113	0.00
5	S	2-Fluorophenol	1.199	1.208	-0.8	109	-0.01
6		Aniline	2.083	2.061	1.1	106	0.00
7	S	Phenol-d6	1.509	1.532	-1.5	110	-0.01
8		2-Chlorophenol	1.417	1.409	0.6	111	0.00
9		Benzaldehyde	0.896	0.937	-4.6	138	0.00
10	C	Phenol	1.644	1.600	2.7	106	0.00
11		bis(2-Chloroethyl)ether	1.317	1.331	-1.1	112	0.00
12		1,3-Dichlorobenzene	1.574	1.664	-5.7	119	0.00
13	C	1,4-Dichlorobenzene	1.650	1.625	1.5	109	0.00
14		1,2-Dichlorobenzene	1.538	1.595	-3.7	114	0.00
15		Benzyl Alcohol	1.255	1.299	-3.5	111	0.00
16		2,2'-oxybis(1-Chloropropane	1.770	1.713	3.2	109	0.00
17		2-Methylphenol	1.135	1.174	-3.4	114	0.00
18		Hexachloroethane	0.597	0.610	-2.2	112	0.00
19	P	n-Nitroso-di-n-propylamine	1.072	1.118	-4.3	114	0.00
20		3+4-Methylphenols	1.499	1.523	-1.6	112	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	116	0.00
22		Acetophenone	0.491	0.497	-1.2	111	0.00
23	S	Nitrobenzene-d5	0.360	0.369	-2.5	111	0.00
24		Nitrobenzene	0.364	0.379	-4.1	111	0.00
25		Isophorone	0.655	0.675	-3.1	113	0.01
26	C	2-Nitrophenol	0.174	0.184	-5.7	114	0.00
27		2,4-Dimethylphenol	0.282	0.287	-1.8	110	0.00
28		bis(2-Chloroethoxy)methane	0.384	0.403	-4.9	117	0.00
29	C	2,4-Dichlorophenol	0.269	0.278	-3.3	108	0.00
30		1,2,4-Trichlorobenzene	0.303	0.303	0.0	109	0.00
31		Naphthalene	1.026	1.041	-1.5	112	0.00
32		Benzoic acid	0.138	0.131	5.1	102	0.01
33		4-Chloroaniline	0.402	0.414	-3.0	112	0.00
34	C	Hexachlorobutadiene	0.173	0.188	-8.7	119	0.00
35		Caprolactam	0.080	0.077	3.8	106	0.00
36	C	4-Chloro-3-methylphenol	0.298	0.313	-5.0	111	0.00
37		2-Methylnaphthalene	0.720	0.749	-4.0	116	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	121	0.00
39		1,2,4,5-Tetrachlorobenzene	0.497	0.500	-0.6	115	0.00
40	P	Hexachlorocyclopentadiene	0.247	0.262	-6.1	122	0.00
41	S	2,4,6-Tribromophenol	0.167	0.171	-2.4	115	-0.01
42	C	2,4,6-Trichlorophenol	0.361	0.361	0.0	110	0.00
43		2,4,5-Trichlorophenol	0.379	0.396	-4.5	115	-0.01
44	S	2-Fluorobiphenyl	1.359	1.332	2.0	112	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.538	1.526	0.8	114	0.00
46		2-Chloronaphthalene	1.236	1.237	-0.1	112	0.00
47		2-Nitroaniline	0.368	0.374	-1.6	114	0.00
48		Acenaphthylene	2.094	2.070	1.1	114	0.00
49		Dimethylphthalate	1.454	1.468	-1.0	114	0.01
50		2,6-Dinitrotoluene	0.308	0.287	6.8	108	0.00
51	C	Acenaphthene	1.183	1.152	2.6	111	0.01
52		3-Nitroaniline	0.352	0.338	4.0	113	0.00
53	P	2,4-Dinitrophenol	0.075	0.077	-2.7	126	0.00
54		Dibenzofuran	1.731	1.753	-1.3	117	0.00
55	P	4-Nitrophenol	0.227	0.218	4.0	113	-0.01
56		2,4-Dinitrotoluene	0.394	0.410	-4.1	113	0.00
57		Fluorene	1.445	1.467	-1.5	116	0.00
58		2,3,4,6-Tetrachlorophenol	0.274	0.280	-2.2	119	0.00
59		Diethylphthalate	1.497	1.510	-0.9	112	0.00
60		4-Chlorophenyl-phenylether	0.620	0.609	1.8	114	0.00
61		4-Nitroaniline	0.326	0.347	-6.4	114	0.00
62		Azobenzene	1.535	1.551	-1.0	114	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	121	0.00
64		4,6-Dinitro-2-methylphenol	0.113	0.118	-4.4	117	0.00
65	c	n-Nitrosodiphenylamine	0.726	0.712	1.9	111	0.00
66		4-Bromophenyl-phenylether	0.208	0.216	-3.8	114	0.00
67		Hexachlorobenzene	0.217	0.218	-0.5	115	0.00
68		Atrazine	0.220	0.221	-0.5	112	0.00
69	C	Pentachlorophenol	0.073	0.078	-6.8	119	0.00
70		Phenanthrene	1.246	1.227	1.5	113	0.00
71		Anthracene	1.211	1.216	-0.4	113	0.00
72		Carbazole	1.187	1.171	1.3	112	0.00
73		Di-n-butylphthalate	1.410	1.448	-2.7	115	0.00
74	C	Fluoranthene	1.277	1.283	-0.5	117	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	118	0.00
76		Benzidine	0.626	0.676	-8.0	141	0.00
77		Pyrene	1.439	1.394	3.1	112	0.00
78	S	Terphenyl-d14	0.911	0.912	-0.1	117	0.00
79		Butylbenzylphthalate	0.668	0.700	-4.8	117	0.00
80		Benzo(a)anthracene	1.274	1.300	-2.0	117	0.00
81		3,3'-Dichlorobenzidine	0.407	0.410	-0.7	118	0.00
82		Chrysene	1.153	1.096	4.9	109	0.00
83		Bis(2-ethylhexyl)phthalate	0.926	0.979	-5.7	124	0.00
84	c	Di-n-octyl phthalate	1.562	1.700	-8.8	125	0.00
85		Indeno(1,2,3-cd)pyrene	1.141	1.165	-2.1	116	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	125	-0.01
87		Benzo(b)fluoranthene	1.385	1.348	2.7	105	0.00

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88	Benzo(k)fluoranthene	1.211	1.134	6.4	118	-0.01
89 C	Benzo(a)pyrene	1.190	1.165	2.1	111	-0.01
90	Dibenzo(a,h)anthracene	1.085	1.059	2.4	113	-0.03
91	Benzo(g,h,i)perylene	1.107	1.072	3.2	113	-0.02

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