Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF043015\
Data File : BF078857.D
Acq On : 30 Apr 2015 18:56
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF043015

Quant Time: Apr 30 19:32:41 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 30 19:07:47 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	104	0.00
2		1,4-Dioxane	0.505	0.496	1.8	103	0.00
3		Pyridine	1.478	1.331	9.9	100	0.01
4		n-Nitrosodimethylamine	0.633	0.625	1.3	106	0.00
5	S	2-Fluorophenol	1.162	1.119	3.7	103	0.00
6		Aniline	2.168	2.110	2.7	103	0.00
7	S	Phenol-d6	1.536	1.409	8.3	102	-0.01
8		2-Chlorophenol	1.318	1.237	6.1	106	-0.01
9		Benzaldehyde	1.035	0.992	4.2	104	0.00
10	C	Phenol	1.782	1.732	2.8	104	0.00
11		bis(2-Chloroethyl)ether	1.306	1.203	7.9	105	0.00
12		1,3-Dichlorobenzene	1.481	1.391	6.1	105	-0.01
13	C	1,4-Dichlorobenzene	1.524	1.452	4.7	105	0.00
14		1,2-Dichlorobenzene	1.419	1.359	4.2	105	0.00
15		Benzyl Alcohol	1.140	1.058	7.2	103	-0.01
16		2,2'-oxybis(1-Chloropropane	1.998	1.891	5.4	106	0.00
17		2-Methylphenol	1.060	0.990	6.6	103	0.00
18		Hexachloroethane	0.525	0.505	3.8	103	0.00
19	P	n-Nitroso-di-n-propylamine	1.037	1.035	0.2	104	0.00
20		3+4-Methylphenols	1.413	1.341	5.1	102	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
22		Acetophenone	0.452	0.413	8.6	102	-0.01
23	S	Nitrobenzene-d5	0.348	0.328	5.7	103	-0.01
24		Nitrobenzene	0.345	0.322	6.7	105	-0.01
25		Isophorone	0.645	0.622	3.6	104	0.00
26	C	2-Nitrophenol	0.143	0.145	-1.4	105	0.00
27		2,4-Dimethylphenol	0.286	0.277	3.1	103	0.00
28		bis(2-Chloroethoxy)methane	0.398	0.393	1.3	104	0.00
29	C	2,4-Dichlorophenol	0.254	0.241	5.1	102	-0.01
30		1,2,4-Trichlorobenzene	0.279	0.266	4.7	104	0.00
31		Naphthalene	0.953	0.881	7.6	102	-0.01
32		Benzoic acid	0.164	0.152	7.3	92	-0.03
33		4-Chloroaniline	0.403	0.386	4.2	107	-0.01
34	C	Hexachlorobutadiene	0.161	0.154	4.3	106	0.00
35		Caprolactam	0.082	0.083	-1.2	108	-0.02
36	C	4-Chloro-3-methylphenol	0.267	0.272	-1.9	103	0.00
37		2-Methylnaphthalene	0.660	0.632	4.2	104	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
39		1,2,4,5-Tetrachlorobenzene	0.563	0.542	3.7	106	0.00
40	P	Hexachlorocyclopentadiene	0.283	0.273	3.5	103	-0.01
41	S	2,4,6-Tribromophenol	0.216	0.212	1.9	101	-0.01
42	C	2,4,6-Trichlorophenol	0.387	0.388	-0.3	105	0.00
43		2,4,5-Trichlorophenol	0.392	0.389	0.8	105	0.00
44	S	2-Fluorobiphenyl	1.436	1.391	3.1	104	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.590	1.519	4.5	106	0.00
46		2-Chloronaphthalene	1.240	1.160	6.5	105	-0.01
47		2-Nitroaniline	0.359	0.341	5.0	99	-0.01
48		Acenaphthylene	2.036	1.940	4.7	104	-0.01
49		Dimethylphthalate	1.468	1.365	7.0	104	-0.01
50		2,6-Dinitrotoluene	0.290	0.280	3.4	103	0.00
51	C	Acenaphthene	1.214	1.152	5.1	102	-0.01
52		3-Nitroaniline	0.362	0.347	4.1	103	-0.01
53	P	2,4-Dinitrophenol	0.085	0.074	12.9	94	-0.01
54		Dibenzofuran	1.754	1.675	4.5	103	0.00
55	P	4-Nitrophenol	0.286	0.259	9.4	98	-0.01
56		2,4-Dinitrotoluene	0.383	0.380	0.8	101	0.00
57		Fluorene	1.440	1.345	6.6	104	-0.01
58		2,3,4,6-Tetrachlorophenol	0.320	0.312	2.5	102	0.00
59		Diethylphthalate	1.454	1.413	2.8	106	-0.01
60		4-Chlorophenyl-phenylether	0.639	0.616	3.6	105	0.00
61		4-Nitroaniline	0.362	0.340	6.1	97	-0.01
62		Azobenzene	1.433	1.396	2.6	103	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
64		4,6-Dinitro-2-methylphenol	0.080	0.073	8.8	96	0.00
65	c	n-Nitrosodiphenylamine	0.666	0.651	2.3	102	0.00
66		4-Bromophenyl-phenylether	0.208	0.200	3.8	105	0.00
67		Hexachlorobenzene	0.238	0.224	5.9	104	-0.01
68		Atrazine	0.194	0.183	5.7	103	-0.01
69	C	Pentachlorophenol	0.120	0.118	1.7	102	0.00
70		Phenanthrene	1.119	1.071	4.3	102	0.00
71		Anthracene	1.098	1.035	5.7	101	-0.01
72		Carbazole	1.046	1.005	3.9	102	0.00
73		Di-n-butylphthalate	1.255	1.235	1.6	101	-0.01
74	C	Fluoranthene	1.166	1.125	3.5	103	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	97	-0.01
76		Benzidine	0.539	0.546	-1.3	96	0.00
77		Pyrene	1.152	1.151	0.1	103	0.00
78	S	Terphenyl-d14	0.835	0.830	0.6	103	0.00
79		Butylbenzylphthalate	0.504	0.510	-1.2	96	0.00
80		Benzo(a)anthracene	1.095	1.065	2.7	99	-0.01
81		3,3'-Dichlorobenzidine	0.390	0.394	-1.0	99	0.00
82		Chrysene	1.053	1.003	4.7	101	-0.01
83		Bis(2-ethylhexyl)phthalate	0.763	0.806	-5.6	101	0.00
84	c	Di-n-octyl phthalate	1.344	1.332	0.9	101	0.00
85		Indeno(1,2,3-cd)pyrene	1.342	1.323	1.4	101	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	101	-0.01
87		Benzo(b)fluoranthene	1.095	1.015	7.3	96	-0.01

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88	Benzo(k)fluoranthene	1.060	1.090	-2.8	106	-0.01
89 C	Benzo(a)pyrene	1.008	0.980	2.8	101	-0.01
90	Dibenzo(a,h)anthracene	1.071	1.045	2.4	101	-0.02
91	Benzo(g,h,i)perylene	1.074	1.035	3.6	101	-0.02

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