Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF060515\
Data File : BF079745.D
Acq On : 5 Jun 2015 23:14
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF060515

Quant Time: Jun 08 07:36:57 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 08 04:31:55 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	105	0.00
2		1,4-Dioxane	0.528	0.523	0.9	105	0.00
3		Pyridine	1.477	1.559	-5.6	101	0.00
4		n-Nitrosodimethylamine	0.641	0.637	0.6	113	0.00
5	S	2-Fluorophenol	1.193	1.243	-4.2	109	0.00
6		Aniline	2.237	2.257	-0.9	106	0.00
7	S	Phenol-d6	1.541	1.555	-0.9	106	0.00
8		2-Chlorophenol	1.292	1.283	0.7	105	0.00
9		Benzaldehyde	0.998	0.989	0.9	104	0.00
10	C	Phenol	1.678	1.759	-4.8	108	0.00
11		bis(2-Chloroethyl)ether	1.301	1.386	-6.5	113	0.00
12		1,3-Dichlorobenzene	1.617	1.572	2.8	103	0.00
13	C	1,4-Dichlorobenzene	1.716	1.749	-1.9	107	0.00
14		1,2-Dichlorobenzene	1.603	1.611	-0.5	107	0.00
15		Benzyl Alcohol	1.274	1.274	0.0	107	0.00
16		2,2'-oxybis(1-Chloropropane	1.968	1.979	-0.6	108	0.00
17		2-Methylphenol	1.205	1.230	-2.1	110	0.00
18		Hexachloroethane	0.499	0.506	-1.4	110	-0.01
19	P	n-Nitroso-di-n-propylamine	1.069	1.102	-3.1	112	0.00
20		3+4-Methylphenols	1.535	1.540	-0.3	105	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
22		Acetophenone	0.477	0.512	-7.3	109	0.00
23	S	Nitrobenzene-d5	0.326	0.347	-6.4	106	0.00
24		Nitrobenzene	0.325	0.347	-6.8	109	0.00
25		Isophorone	0.665	0.691	-3.9	109	0.00
26	C	2-Nitrophenol	0.137	0.145	-5.8	110	0.00
27		2,4-Dimethylphenol	0.308	0.336	-9.1	113	0.00
28		bis(2-Chloroethoxy)methane	0.411	0.407	1.0	104	0.00
29	C	2,4-Dichlorophenol	0.276	0.284	-2.9	106	0.00
30		1,2,4-Trichlorobenzene	0.335	0.357	-6.6	110	0.00
31		Naphthalene	1.049	1.060	-1.0	107	0.00
32		Benzoic acid	0.113	0.129	-14.2	109	0.00
33		4-Chloroaniline	0.533	0.567	-6.4	108	0.00
34	C	Hexachlorobutadiene	0.204	0.203	0.5	104	0.00
35		Caprolactam	0.084	0.092	-9.5	109	0.00
36	C	4-Chloro-3-methylphenol	0.309	0.320	-3.6	106	0.00
37		2-Methylnaphthalene	0.667	0.704	-5.5	110	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
39		1,2,4,5-Tetrachlorobenzene	0.623	0.620	0.5	111	0.00
40	P	Hexachlorocyclopentadiene	0.268	0.275	-2.6	113	0.00
41	S	2,4,6-Tribromophenol	0.191	0.180	5.8	98	0.00
42	C	2,4,6-Trichlorophenol	0.414	0.421	-1.7	112	0.00
43		2,4,5-Trichlorophenol	0.436	0.414	5.0	105	-0.01
44	S	2-Fluorobiphenyl	1.476	1.348	8.7	103	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.740	1.710	1.7	111	0.00
46		2-Chloronaphthalene	1.421	1.319	7.2	107	0.00
47		2-Nitroaniline	0.298	0.279	6.4	108	0.00
48		Acenaphthylene	2.349	2.257	3.9	108	0.00
49		Dimethylphthalate	1.541	1.439	6.6	107	0.00
50		2,6-Dinitrotoluene	0.247	0.239	3.2	110	-0.01
51	C	Acenaphthene	1.330	1.301	2.2	111	0.00
52		3-Nitroaniline	0.325	0.305	6.2	106	-0.01
53	P	2,4-Dinitrophenol	0.062	0.060	3.2	109	0.00
54		Dibenzofuran	1.884	1.911	-1.4	112	0.00
55	P	4-Nitrophenol	0.243	0.255	-4.9	117	0.00
56		2,4-Dinitrotoluene	0.293	0.336	-14.7	126	0.00
57		Fluorene	1.630	1.503	7.8	106	-0.01
58		2,3,4,6-Tetrachlorophenol	0.353	0.363	-2.8	107	0.00
59		Diethylphthalate	1.507	1.412	6.3	109	-0.01
60		4-Chlorophenyl-phenylether	0.739	0.710	3.9	107	0.00
61		4-Nitroaniline	0.323	0.308	4.6	110	-0.01
62		Azobenzene	1.522	1.526	-0.3	110	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	108	-0.01
64		4,6-Dinitro-2-methylphenol	0.050	0.046	8.0	111	-0.01
65	c	n-Nitrosodiphenylamine	0.668	0.622	6.9	103	0.00
66		4-Bromophenyl-phenylether	0.217	0.203	6.5	100	-0.01
67		Hexachlorobenzene	0.244	0.241	1.2	113	0.00
68		Atrazine	0.201	0.213	-6.0	119	0.00
69	C	Pentachlorophenol	0.104	0.112	-7.7	111	-0.01
70		Phenanthrene	1.171	1.192	-1.8	113	-0.02
71		Anthracene	1.155	1.177	-1.9	111	-0.01
72		Carbazole	1.097	1.056	3.7	108	-0.02
73		Di-n-butylphthalate	1.244	1.268	-1.9	115	-0.03
74	C	Fluoranthene	1.282	1.303	-1.6	111	-0.08
75	I	Chrysene-d12	1.000	1.000	0.0	104	-0.23
76		Benzidine	0.631	0.648	-2.7	112	-0.09
77		Pyrene	1.226	1.281	-4.5	110	-0.10
78	S	Terphenyl-d14	0.873	0.872	0.1	106	-0.11
79		Butylbenzylphthalate	0.489	0.522	-6.7	112	-0.16
80		Benzo(a)anthracene	1.224	1.215	0.7	106	-0.22
81		3,3'-Dichlorobenzidine	0.416	0.447	-7.5	112	-0.22
82		Chrysene	1.132	1.150	-1.6	109	-0.23
83		Bis(2-ethylhexyl)phthalate	0.772	0.794	-2.8	110	-0.23
84	c	Di-n-octyl phthalate	1.360	1.368	-0.6	109	-0.31
85		Indeno(1,2,3-cd)pyrene	1.310	1.358	-3.7	112	-0.34
86	I	Perylene-d12	1.000	1.000	0.0	107	-0.32
87		Benzo(b)fluoranthene	1.242	1.348	-8.5	123	-0.31

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.192	1.124	5.7	100	-0.31
89 C	Benzo(a)pyrene	1.147	1.155	-0.7	110	-0.31
90	Dibenzo(a,h)anthracene	1.107	1.098	0.8	110	-0.35
91	Benzo(g,h,i)perylene	1.144	1.139	0.4	111	-0.35

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