Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110715\
Data File : BF082760.D
Acq On : 6 Nov 2015 14:38
Operator : UM/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF110715

Quant Time: Nov 06 17:41:10 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 06 14:39:58 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	123	0.00
2		1,4-Dioxane	0.555	0.523	5.8	113	-0.01
3		Pyridine	1.609	1.612	-0.2	119	-0.01
4		n-Nitrosodimethylamine	0.798	0.770	3.5	121	0.01
5	S	2-Fluorophenol	1.285	1.240	3.5	117	0.00
6		Aniline	2.399	2.211	7.8	120	0.00
7	S	Phenol-d6	1.709	1.539	9.9	114	0.01
8		2-Chlorophenol	1.425	1.325	7.0	115	0.00
9		Benzaldehyde	0.944	0.904	4.2	111	0.00
10	C	Phenol	1.939	1.733	10.6	105	0.01
11		bis(2-Chloroethyl)ether	1.450	1.351	6.8	109	0.00
12		1,3-Dichlorobenzene	1.607	1.600	0.4	121	0.00
13	C	1,4-Dichlorobenzene	1.672	1.470	12.1	119	0.00
14		1,2-Dichlorobenzene	1.521	1.391	8.5	112	0.00
15		Benzyl Alcohol	1.501	1.511	-0.7	128	0.01
16		2,2'-oxybis(1-Chloropropane	2.127	1.964	7.7	117	0.00
17		2-Methylphenol	1.207	1.185	1.8	120	0.01
18		Hexachloroethane	0.598	0.547	8.5	120	0.00
19	P	n-Nitroso-di-n-propylamine	1.256	1.101	12.3	110	0.01
20		3+4-Methylphenols	1.569	1.357	13.5	118	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	123	0.00
22		Acetophenone	0.573	0.506	11.7	119	0.00
23	S	Nitrobenzene-d5	0.460	0.454	1.3	123	0.01
24		Nitrobenzene	0.470	0.402	14.5	110	0.00
25		Isophorone	0.850	0.804	5.4	121	0.01
26	C	2-Nitrophenol	0.196	0.202	-3.1	116	0.00
27		2,4-Dimethylphenol	0.353	0.338	4.2	124	0.01
28		bis(2-Chloroethoxy)methane	0.480	0.444	7.5	111	0.00
29	C	2,4-Dichlorophenol	0.319	0.287	10.0	114	0.00
30		1,2,4-Trichlorobenzene	0.358	0.341	4.7	118	0.00
31		Naphthalene	1.103	1.014	8.1	118	0.01
32		Benzoic acid	0.241	0.241	0.0	121	0.03
33		4-Chloroaniline	0.476	0.452	5.0	118	0.01
34	C	Hexachlorobutadiene	0.224	0.209	6.7	117	0.00
35		Caprolactam	0.100	0.092	8.0	108	0.02
36	C	4-Chloro-3-methylphenol	0.375	0.325	13.3	110	0.00
37		2-Methylnaphthalene	0.714	0.626	12.3	113	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	119	0.00
39		1,2,4,5-Tetrachlorobenzene	0.657	0.680	-3.5	114	0.00
40	P	Hexachlorocyclopentadiene	0.353	0.371	-5.1	119	0.00
41	S	2,4,6-Tribromophenol	0.229	0.209	8.7	114	0.00
42	C	2,4,6-Trichlorophenol	0.491	0.474	3.5	120	0.00
43		2,4,5-Trichlorophenol	0.485	0.492	-1.4	120	0.01
44	S	2-Fluorobiphenyl	1.395	1.250	10.4	115	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.716	1.743	-1.6	112	0.00
46		2-Chloronaphthalene	1.339	1.354	-1.1	114	0.00
47		2-Nitroaniline	0.497	0.540	-8.7	118	0.01
48		Acenaphthylene	2.098	1.944	7.3	114	0.00
49		Dimethylphthalate	1.639	1.714	-4.6	138	0.01
50		2,6-Dinitrotoluene	0.350	0.336	4.0	128	0.00
51	C	Acenaphthene	1.330	1.236	7.1	115	0.00
52		3-Nitroaniline	0.385	0.370	3.9	103	0.01
53	P	2,4-Dinitrophenol	0.192	0.210	-9.4	113	0.01
54		Dibenzofuran	1.793	1.623	9.5	113	0.00
55	P	4-Nitrophenol	0.295	0.321	-8.8	131	0.01
56		2,4-Dinitrotoluene	0.474	0.514	-8.4	143	0.01
57		Fluorene	1.452	1.329	8.5	112	0.00
58		2,3,4,6-Tetrachlorophenol	0.396	0.388	2.0	126	0.01
59		Diethylphthalate	1.600	1.671	-4.4	127	0.01
60		4-Chlorophenyl-phenylether	0.726	0.727	-0.1	125	0.01
61		4-Nitroaniline	0.387	0.407	-5.2	131	0.01
62		Azobenzene	1.719	1.772	-3.1	124	0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	126	0.00
64		4,6-Dinitro-2-methylphenol	0.142	0.140	1.4	127	0.00
65	c	n-Nitrosodiphenylamine	0.723	0.626	13.4	109	0.00
66		4-Bromophenyl-phenylether	0.241	0.211	12.4	121	0.00
67		Hexachlorobenzene	0.269	0.236	12.3	121	0.00
68		Atrazine	0.223	0.183	17.9	110	0.00
69	C	Pentachlorophenol	0.163	0.169	-3.7	119	0.00
70		Phenanthrene	1.161	1.009	13.1	109	0.00
71		Anthracene	1.223	1.158	5.3	129	0.01
72		Carbazole	1.071	1.015	5.2	117	0.00
73		Di-n-butylphthalate	1.334	1.272	4.6	123	0.00
74	C	Fluoranthene	1.246	1.183	5.1	118	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	108	0.01
76		Benzidine	0.670	0.640	4.5	98	0.00
77		Pyrene	1.373	1.515	-10.3	123	0.01
78	S	Terphenyl-d14	0.843	0.878	-4.2	111	0.00
79		Butylbenzylphthalate	0.607	0.690	-13.7	131	0.01
80		Benzo(a)anthracene	1.251	1.299	-3.8	119	0.00
81		3,3'-Dichlorobenzidine	0.445	0.435	2.2	106	0.01
82		Chrysene	1.146	1.282	-11.9	131	0.01
83		Bis(2-ethylhexyl)phthalate	0.831	0.944	-13.6	129	0.01
84	c	Di-n-octyl phthalate	1.385	1.467	-5.9	116	0.01
85		Indeno(1,2,3-cd)pyrene	0.987	1.091	-10.5	126	0.02
86	I	Perylene-d12	1.000	1.000	0.0	128	0.02
87		Benzo(b)fluoranthene	1.284	1.179	8.2	130	0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.139	1.074	5.7	115	0.02
89 C	Benzo(a)pyrene	1.112	1.062	4.5	125	0.02
90	Dibenzo(a,h)anthracene	0.942	0.894	5.1	124	0.03
91	Benzo(g,h,i)perylene	0.902	0.850	5.8	126	0.03

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