Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
Data File : BF083006.D
Acq On : 18 Nov 2015 12:58
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 19 01:44:52 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 18 13:32:32 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	157#	0.00
2		1,4-Dioxane	0.555	0.565	-1.8	157#	0.00
3		Pyridine	1.609	1.768	-9.9	167#	0.00
4		n-Nitrosodimethylamine	0.798	0.783	1.9	158#	0.00
5	S	2-Fluorophenol	1.285	1.251	2.6	151#	0.00
6		Aniline	2.399	2.214	7.7	153#	0.00
7	S	Phenol-d6	1.709	1.540	9.9	146	0.00
8		2-Chlorophenol	1.425	1.264	11.3	141	0.00
9		Benzaldehyde	0.944	0.850	10.0	134	0.00
10	C	Phenol	1.939	1.807	6.8	139	0.00
11		bis(2-Chloroethyl)ether	1.450	1.348	7.0	139	0.00
12		1,3-Dichlorobenzene	1.607	1.466	8.8	142	0.00
13	C	1,4-Dichlorobenzene	1.672	1.437	14.1	149	0.00
14		1,2-Dichlorobenzene	1.521	1.508	0.9	156#	0.00
15		Benzyl Alcohol	1.501	1.232	17.9	134	0.00
16		2,2'-oxybis(1-Chloropropane	2.127	2.059	3.2	156#	0.00
17		2-Methylphenol	1.207	1.106	8.4	143	0.00
18		Hexachloroethane	0.598	0.585	2.2	164#	0.00
19	P	n-Nitroso-di-n-propylamine	1.256	1.148	8.6	147	0.00
20		3+4-Methylphenols	1.569	1.466	6.6	164#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	157#	0.00
22		Acetophenone	0.573	0.571	0.3	171#	0.00
23	S	Nitrobenzene-d5	0.460	0.417	9.3	144	0.00
24		Nitrobenzene	0.470	0.441	6.2	153#	0.00
25		Isophorone	0.850	0.823	3.2	158#	0.00
26	C	2-Nitrophenol	0.196	0.190	3.1	139	0.00
27		2,4-Dimethylphenol	0.353	0.330	6.5	154#	0.00
28		bis(2-Chloroethoxy)methane	0.480	0.493	-2.7	157#	0.00
29	C	2,4-Dichlorophenol	0.319	0.312	2.2	158#	0.00
30		1,2,4-Trichlorobenzene	0.358	0.347	3.1	154#	0.00
31		Naphthalene	1.103	0.968	12.2	144	0.00
32		Benzoic acid	0.241	0.227	5.8	145	0.00
33		4-Chloroaniline	0.476	0.402	15.5	133	0.00
34	C	Hexachlorobutadiene	0.224	0.193	13.8	138	0.00
35		Caprolactam	0.100	0.098	2.0	146	0.00
36	C	4-Chloro-3-methylphenol	0.375	0.384	-2.4	166#	0.00
37		2-Methylnaphthalene	0.714	0.696	2.5	160#	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	163#	0.00
39		1,2,4,5-Tetrachlorobenzene	0.657	0.609	7.3	139	0.00
40	P	Hexachlorocyclopentadiene	0.353	0.339	4.0	148	0.00
41	S	2,4,6-Tribromophenol	0.229	0.192	16.2	143	0.00
42	C	2,4,6-Trichlorophenol	0.491	0.460	6.3	158#	0.00
43		2,4,5-Trichlorophenol	0.485	0.468	3.5	155#	0.00
44	S	2-Fluorobiphenyl	1.395	1.092	21.7	136	0.00

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Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.716	1.764	-2.8	155#	0.00
46		2-Chloronaphthalene	1.339	1.327	0.9	152#	0.00
47		2-Nitroaniline	0.497	0.440	11.5	131	0.00
48		Acenaphthylene	2.098	1.989	5.2	159#	0.00
49		Dimethylphthalate	1.639	1.374	16.2	150#	0.00
50		2,6-Dinitrotoluene	0.350	0.365	-4.3	190#	0.00
51	C	Acenaphthene	1.330	1.305	1.9	166#	0.00
52		3-Nitroaniline	0.385	0.386	-0.3	146	0.00
53	P	2,4-Dinitrophenol	0.192	0.171	10.9	125	0.00
54		Dibenzofuran	1.793	1.741	2.9	165#	0.00
55	P	4-Nitrophenol	0.295	0.313	-6.1	174#	0.00
56		2,4-Dinitrotoluene	0.474	0.462	2.5	175#	0.00
57		Fluorene	1.452	1.434	1.2	164#	0.00
58		2,3,4,6-Tetrachlorophenol	0.396	0.342	13.6	151#	0.00
59		Diethylphthalate	1.600	1.523	4.8	157#	0.00
60		4-Chlorophenyl-phenylether	0.726	0.555	23.6	130	0.00
61		4-Nitroaniline	0.387	0.437	-12.9	191#	0.00
62		Azobenzene	1.719	1.531	10.9	146	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	154#	0.00
64		4,6-Dinitro-2-methylphenol	0.142	0.150	-5.6	167#	0.00
65	c	n-Nitrosodiphenylamine	0.723	0.644	10.9	137	0.00
66		4-Bromophenyl-phenylether	0.241	0.231	4.1	163#	0.00
67		Hexachlorobenzene	0.269	0.251	6.7	158#	0.00
68		Atrazine	0.223	0.221	0.9	162#	0.00
69	C	Pentachlorophenol	0.163	0.141	13.5	122	0.00
70		Phenanthrene	1.161	0.939	19.1	125	0.00
71		Anthracene	1.223	1.152	5.8	157#	0.00
72		Carbazole	1.071	0.918	14.3	130	0.00
73		Di-n-butylphthalate	1.334	1.102	17.4	130	0.00
74	C	Fluoranthene	1.246	1.174	5.8	144	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	124	0.00
76		Benzidine	0.670	0.673	-0.4	119	0.00
77		Pyrene	1.373	1.582	-15.2	149	0.00
78	S	Terphenyl-d14	0.843	0.777	7.8	113	0.00
79		Butylbenzylphthalate	0.607	0.755	-24.4	165#	0.00
80		Benzo(a)anthracene	1.251	1.398	-11.8	147	0.00
81		3,3'-Dichlorobenzidine	0.445	0.500	-12.4	140	0.00
82		Chrysene	1.146	1.298	-13.3	153#	0.00
83		Bis(2-ethylhexyl)phthalate	0.831	0.905	-8.9	143	0.00
84	c	Di-n-octyl phthalate	1.385	1.433	-3.5	130	0.00
85		Indeno(1,2,3-cd)pyrene	0.987	0.964	2.3	128	0.00
86	I	Perylene-d12	1.000	1.000	0.0	124	0.00
87		Benzo(b)fluoranthene	1.284	1.277	0.5	137	0.00

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88	Benzo(k)fluoranthene	1.139	0.978	14.1	102	0.00
89 C	Benzo(a)pyrene	1.112	1.049	5.7	120	0.00
90	Dibenzo(a,h)anthracene	0.942	0.940	0.2	126	0.00
91	Benzo(g,h,i)perylene	0.902	0.969	-7.4	139	0.00

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