Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF033115\
Data File : BF078153.D
Acq On : 1 Apr 2015 11:41
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 01 13:18:53 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 31 14:58:34 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	94	0.00
2		1,4-Dioxane	0.520	0.459	11.7	83	-0.02
3		Pyridine	1.398	1.349	3.5	93	-0.06
4		n-Nitrosodimethylamine	0.609	0.550	9.7	79	-0.03
5	S	2-Fluorophenol	1.146	1.163	-1.5	99	-0.01
6		Aniline	2.025	2.045	-1.0	98	0.00
7	S	Phenol-d6	1.416	1.493	-5.4	100	0.00
8		2-Chlorophenol	1.364	1.381	-1.2	100	0.00
9		Benzaldehyde	0.580	0.705	-21.6	116	0.00
10	C	Phenol	1.673	1.766	-5.6	103	-0.01
11		bis(2-Chloroethyl)ether	1.253	1.250	0.2	95	0.00
12		1,3-Dichlorobenzene	1.517	1.523	-0.4	99	0.00
13	C	1,4-Dichlorobenzene	1.576	1.566	0.6	99	0.00
14		1,2-Dichlorobenzene	1.445	1.427	1.2	98	0.00
15		Benzyl Alcohol	1.105	1.127	-2.0	97	0.00
16		2,2'-oxybis(1-Chloropropane	1.689	1.682	0.4	97	0.00
17		2-Methylphenol	1.110	1.095	1.4	94	-0.01
18		Hexachloroethane	0.517	0.534	-3.3	104	0.00
19	P	n-Nitroso-di-n-propylamine	0.979	0.990	-1.1	99	0.00
20		3+4-Methylphenols	1.457	1.524	-4.6	103	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
22		Acetophenone	0.443	0.452	-2.0	99	0.00
23	S	Nitrobenzene-d5	0.305	0.323	-5.9	104	0.00
24		Nitrobenzene	0.329	0.350	-6.4	107	0.00
25		Isophorone	0.616	0.624	-1.3	97	0.00
26	C	2-Nitrophenol	0.161	0.161	0.0	90	0.00
27		2,4-Dimethylphenol	0.293	0.297	-1.4	96	0.00
28		bis(2-Chloroethoxy)methane	0.374	0.378	-1.1	99	0.00
29	C	2,4-Dichlorophenol	0.279	0.276	1.1	94	0.00
30		1,2,4-Trichlorobenzene	0.313	0.305	2.6	94	0.00
31		Naphthalene	1.027	1.032	-0.5	98	0.00
32		Benzoic acid	0.141	0.151	-7.1	93	0.01
33		4-Chloroaniline	0.417	0.402	3.6	91	-0.01
34	C	Hexachlorobutadiene	0.177	0.178	-0.6	98	0.00
35		Caprolactam	0.082	0.078	4.9	90	0.02
36	C	4-Chloro-3-methylphenol	0.281	0.270	3.9	94	0.00
37		2-Methylnaphthalene	0.690	0.670	2.9	92	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	92	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.597	0.600	-0.5	93	0.00
40	P	Hexachlorocyclopentadiene	0.254	0.207	18.5	69	0.00
41	S	2,4,6-Tribromophenol	0.191	0.181	5.2	86	-0.01
42	C	2,4,6-Trichlorophenol	0.413	0.400	3.1	86	0.00
43		2,4,5-Trichlorophenol	0.446	0.421	5.6	88	0.00
44	S	2-Fluorobiphenyl	1.361	1.377	-1.2	97	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.599	1.611	-0.8	92	0.00
46		2-Chloronaphthalene	1.299	1.284	1.2	93	0.00
47		2-Nitroaniline	0.323	0.345	-6.8	93	0.00
48		Acenaphthylene	2.161	2.134	1.2	93	0.00
49		Dimethylphthalate	1.483	1.486	-0.2	94	0.00
50		2,6-Dinitrotoluene	0.307	0.325	-5.9	97	0.00
51	C	Acenaphthene	1.239	1.258	-1.5	94	0.00
52		3-Nitroaniline	0.357	0.346	3.1	87	-0.01
53	P	2,4-Dinitrophenol	0.093	0.054	41.9#	53	-0.01
54		Dibenzofuran	1.837	1.824	0.7	92	0.00
55	P	4-Nitrophenol	0.265	0.243	8.3	79	-0.01
56		2,4-Dinitrotoluene	0.401	0.413	-3.0	95	0.00
57		Fluorene	1.526	1.501	1.6	90	0.00
58		2,3,4,6-Tetrachlorophenol	0.344	0.316	8.1	84	0.00
59		Diethylphthalate	1.445	1.431	1.0	91	0.00
60		4-Chlorophenyl-phenylether	0.696	0.669	3.9	90	0.00
61		4-Nitroaniline	0.356	0.349	2.0	89	0.00
62		Azobenzene	1.381	1.527	-10.6	94	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	91	0.00
64		4,6-Dinitro-2-methylphenol	0.078	0.061	21.8	62	0.00
65	c	n-Nitrosodiphenylamine	0.666	0.680	-2.1	90	0.00
66		4-Bromophenyl-phenylether	0.213	0.204	4.2	84	0.00
67		Hexachlorobenzene	0.235	0.220	6.4	84	-0.01
68		Atrazine	0.187	0.180	3.7	83	-0.01
69	C	Pentachlorophenol	0.104	0.089	14.4	70	0.00
70		Phenanthrene	1.148	1.109	3.4	87	-0.01
71		Anthracene	1.134	1.153	-1.7	95	0.00
72		Carbazole	1.079	1.032	4.4	90	-0.01
73		Di-n-butylphthalate	1.210	1.234	-2.0	92	-0.01
74	C	Fluoranthene	1.266	1.190	6.0	80	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	82	-0.09
76		Benzidine	0.546	0.476	12.8	77	-0.01
77		Pyrene	1.258	1.257	0.1	76	-0.01
78	S	Terphenyl-d14	0.819	0.802	2.1	77	-0.02
79		Butylbenzylphthalate	0.496	0.560	-12.9	92	-0.03
80		Benzo(a)anthracene	1.186	1.172	1.2	85	-0.08
81		3,3'-Dichlorobenzidine	0.391	0.408	-4.3	91	-0.08
82		Chrysene	1.066	1.089	-2.2	85	-0.03
83		Bis(2-ethylhexyl)phthalate	0.751	0.830	-10.5	93	-0.10
84	c	Di-n-octyl phthalate	1.284	1.425	-11.0	94	-0.19
85		Indeno(1,2,3-cd)pyrene	1.331	1.344	-1.0	90	-0.37
86	I	Perylene-d12	1.000	1.000	0.0	86	-0.29
87		Benzo(b)fluoranthene	1.306	1.229	5.9	82	-0.23

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88	Benzo(k)fluoranthene	1.020	1.093	-7.2	97	-0.23
89 C	Benzo(a)pyrene	1.089	1.102	-1.2	89	-0.26
90	Dibenzo(a,h)anthracene	1.081	1.099	-1.7	89	-0.37
91	Benzo(g,h,i)perylene	1.100	1.120	-1.8	88	-0.37

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