Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF062215\
Data File : BF080104.D
Acq On : 22 Jun 2015 20:03
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 23 01:34:20 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 23 00:53:56 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	102	0.00
2		1,4-Dioxane	0.537	0.498	7.3	94	-0.01
3		Pyridine	1.428	1.300	9.0	92	0.00
4		n-Nitrosodimethylamine	0.679	0.667	1.8	94	-0.01
5	S	2-Fluorophenol	1.130	1.130	0.0	99	-0.01
6		Aniline	2.068	2.379	-15.0	111	0.00
7	S	Phenol-d6	1.503	1.584	-5.4	100	-0.01
8		2-Chlorophenol	1.306	1.334	-2.1	99	-0.01
9		Benzaldehyde	0.868	0.899	-3.6	101	0.00
10	C	Phenol	1.641	1.702	-3.7	101	0.00
11		bis(2-Chloroethyl)ether	1.302	1.342	-3.1	101	0.00
12		1,3-Dichlorobenzene	1.569	1.536	2.1	96	0.00
13	C	1,4-Dichlorobenzene	1.492	1.732	-16.1	115	0.00
14		1,2-Dichlorobenzene	1.452	1.590	-9.5	108	0.00
15		Benzyl Alcohol	1.229	1.238	-0.7	98	0.00
16		2,2'-oxybis(1-Chloropropane	1.932	2.363	-22.3	125	0.00
17		2-Methylphenol	1.115	1.188	-6.5	102	0.00
18		Hexachloroethane	0.497	0.540	-8.7	105	0.00
19	P	n-Nitroso-di-n-propylamine	1.044	1.259	-20.6	127	0.00
20		3+4-Methylphenols	1.440	1.652	-14.7	112	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	108	-0.01
22		Acetophenone	0.466	0.455	2.4	99	-0.01
23	S	Nitrobenzene-d5	0.344	0.366	-6.4	111	0.00
24		Nitrobenzene	0.355	0.400	-12.7	118	0.00
25		Isophorone	0.691	0.692	-0.1	105	0.00
26	C	2-Nitrophenol	0.148	0.184	-24.3#	130	0.00
27		2,4-Dimethylphenol	0.309	0.338	-9.4	120	0.00
28		bis(2-Chloroethoxy)methane	0.406	0.394	3.0	102	0.00
29	C	2,4-Dichlorophenol	0.265	0.313	-18.1	122	0.00
30		1,2,4-Trichlorobenzene	0.301	0.343	-14.0	121	0.00
31		Naphthalene	1.046	1.015	3.0	101	0.00
32		Benzoic acid	0.187	0.176	5.9	99	0.00
33		4-Chloroaniline	0.448	0.414	7.6	102	0.00
34	C	Hexachlorobutadiene	0.185	0.211	-14.1	127	0.00
35		Caprolactam	0.086	0.098	-14.0	114	0.00
36	C	4-Chloro-3-methylphenol	0.299	0.307	-2.7	105	0.00
37		2-Methylnaphthalene	0.676	0.724	-7.1	112	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	114	0.00
39		1,2,4,5-Tetrachlorobenzene	0.582	0.657	-12.9	124	0.00
40	P	Hexachlorocyclopentadiene	0.260	0.290	-11.5	120	0.00
41	S	2,4,6-Tribromophenol	0.196	0.222	-13.3	125	0.00
42	C	2,4,6-Trichlorophenol	0.396	0.446	-12.6	122	0.00
43		2,4,5-Trichlorophenol	0.456	0.453	0.7	107	0.00
44	S	2-Fluorobiphenyl	1.402	1.405	-0.2	112	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.627	1.729	-6.3	119	0.00
46		2-Chloronaphthalene	1.320	1.284	2.7	106	0.00
47		2-Nitroaniline	0.362	0.390	-7.7	111	0.00
48		Acenaphthylene	2.204	2.202	0.1	106	0.00
49		Dimethylphthalate	1.504	1.711	-13.8	129	0.00
50		2,6-Dinitrotoluene	0.301	0.325	-8.0	110	0.00
51	C	Acenaphthene	1.348	1.354	-0.4	110	0.00
52		3-Nitroaniline	0.348	0.375	-7.8	113	0.00
53	P	2,4-Dinitrophenol	0.106	0.138	-30.2#	139	0.00
54		Dibenzofuran	1.913	1.910	0.2	110	0.00
55	P	4-Nitrophenol	0.278	0.331	-19.1	135	0.00
56		2,4-Dinitrotoluene	0.383	0.445	-16.2	127	0.00
57		Fluorene	1.518	1.700	-12.0	124	0.00
58		2,3,4,6-Tetrachlorophenol	0.385	0.375	2.6	103	0.00
59		Diethylphthalate	1.514	1.515	-0.1	105	0.00
60		4-Chlorophenyl-phenylether	0.729	0.726	0.4	112	-0.01
61		4-Nitroaniline	0.359	0.358	0.3	105	0.00
62		Azobenzene	1.541	1.534	0.5	105	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	127	-0.02
64		4,6-Dinitro-2-methylphenol	0.094	0.096	-2.1	119	-0.01
65	c	n-Nitrosodiphenylamine	0.651	0.608	6.6	111	0.00
66		4-Bromophenyl-phenylether	0.213	0.221	-3.8	126	0.00
67		Hexachlorobenzene	0.236	0.239	-1.3	122	-0.01
68		Atrazine	0.206	0.218	-5.8	124	-0.01
69	C	Pentachlorophenol	0.134	0.123	8.2	116	-0.01
70		Phenanthrene	1.211	1.109	8.4	114	-0.02
71		Anthracene	1.166	1.165	0.1	125	-0.02
72		Carbazole	1.113	1.062	4.6	116	-0.02
73		Di-n-butylphthalate	1.286	1.131	12.1	113	-0.03
74	C	Fluoranthene	1.290	1.232	4.5	121	-0.08
75	I	Chrysene-d12	1.000	1.000	0.0	115	-0.17
76		Benzidine	0.559	0.600	-7.3	119	-0.08
77		Pyrene	1.231	1.210	1.7	112	-0.09
78	S	Terphenyl-d14	0.856	0.859	-0.4	116	-0.10
79		Butylbenzylphthalate	0.495	0.505	-2.0	112	-0.14
80		Benzo(a)anthracene	1.218	1.207	0.9	114	-0.17
81		3,3'-Dichlorobenzidine	0.420	0.405	3.6	109	-0.17
82		Chrysene	1.124	1.095	2.6	111	-0.17
83		Bis(2-ethylhexyl)phthalate	0.782	0.743	5.0	105	-0.18
84	c	Di-n-octyl phthalate	1.339	1.231	8.1	103	-0.25
85		Indeno(1,2,3-cd)pyrene	1.417	1.260	11.1	102	-0.27
86	I	Perylene-d12	1.000	1.000	0.0	110	-0.25
87		Benzo(b)fluoranthene	1.211	1.175	3.0	103	-0.25

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88	Benzo(k)fluoranthene	1.233	1.232	0.1	109	-0.25
89 C	Benzo(a)pyrene	1.150	1.119	2.7	106	-0.25
90	Dibenzo(a,h)anthracene	1.177	1.114	5.4	103	-0.27
91	Benzo(g,h,i)perylene	1.188	1.112	6.4	101	-0.27

(#) = Out of Range SPCC's out = 0 CCC's out = 1