Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF123114\
Data File : BF076714.D
Acq On : 31 Dec 2014 21:15
Operator : IZ / TP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SSTDCCC040EC

Quant Time: Jan 01 07:00:18 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121714.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 31 23:59:08 2014
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	75	-0.20
2		1,4-Dioxane	0.504	0.430	14.7	64	-0.28
3		Pyridine	1.269	1.284	-1.2	78	-0.40
4		n-Nitrosodimethylamine	0.614	0.628	-2.3	76	-0.38
5	S	2-Fluorophenol	1.180	1.206	-2.2	77	-0.28
6		Aniline	2.058	2.115	-2.8	77	-0.20
7	S	Phenol-d6	1.441	1.538	-6.7	79	-0.18
8		2-Chlorophenol	1.351	1.371	-1.5	75	-0.20
9		Benzaldehyde	1.096	0.984	10.2	66	-0.22
10	C	Phenol	1.749	1.817	-3.9	77	-0.18
11		bis(2-Chloroethyl)ether	1.259	1.350	-7.2	80	-0.19
12		1,3-Dichlorobenzene	1.568	1.613	-2.9	76	-0.21
13	C	1,4-Dichlorobenzene	1.589	1.687	-6.2	81	-0.20
14		1,2-Dichlorobenzene	1.506	1.545	-2.6	78	-0.21
15		Benzyl Alcohol	1.248	1.310	-5.0	76	-0.19
16		2,2'-oxybis(1-Chloropropane	1.823	1.765	3.2	71	-0.18
17		2-Methylphenol	1.110	1.183	-6.6	78	-0.16
18		Hexachloroethane	0.568	0.598	-5.3	80	-0.20
19	P	n-Nitroso-di-n-propylamine	1.052	1.099	-4.5	81	-0.18
20		3+4-Methylphenols	1.499	1.515	-1.1	75	-0.16
21	I	Naphthalene-d8	1.000	1.000	0.0	75	-0.20
22		Acetophenone	0.479	0.473	1.3	74	-0.19
23	S	Nitrobenzene-d5	0.336	0.365	-8.6	81	-0.19
24		Nitrobenzene	0.349	0.371	-6.3	79	-0.19
25		Isophorone	0.648	0.656	-1.2	76	-0.19
26	C	2-Nitrophenol	0.170	0.171	-0.6	72	-0.20
27		2,4-Dimethylphenol	0.276	0.285	-3.3	77	-0.16
28		bis(2-Chloroethoxy)methane	0.380	0.412	-8.4	78	-0.18
29	C	2,4-Dichlorophenol	0.271	0.276	-1.8	77	-0.19
30		1,2,4-Trichlorobenzene	0.292	0.296	-1.4	75	-0.20
31		Naphthalene	0.991	0.985	0.6	75	-0.20
32		Benzoic acid	0.153	0.153	0.0	74	-0.15
33		4-Chloroaniline	0.404	0.403	0.2	72	-0.19
34	C	Hexachlorobutadiene	0.170	0.179	-5.3	81	-0.19
35		Caprolactam	0.078	0.072	7.7	70	-0.20
36	C	4-Chloro-3-methylphenol	0.301	0.289	4.0	68	-0.19
37		2-Methylnaphthalene	0.691	0.687	0.6	73	-0.21
38	I	Acenaphthene-d10	1.000	1.000	0.0	74	-0.22
39		1,2,4,5-Tetrachlorobenzene	0.492	0.484	1.6	71	-0.21
40	P	Hexachlorocyclopentadiene	0.241	0.254	-5.4	73	-0.21
41	S	2,4,6-Tribromophenol	0.169	0.175	-3.6	75	-0.23
42	C	2,4,6-Trichlorophenol	0.357	0.385	-7.8	77	-0.20
43		2,4,5-Trichlorophenol	0.384	0.383	0.3	71	-0.20
44	S	2-Fluorobiphenyl	1.285	1.312	-2.1	76	-0.21

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.486	1.469	1.1	70	-0.21
46		2-Chloronaphthalene	1.178	1.225	-4.0	77	-0.21
47		2-Nitroaniline	0.358	0.389	-8.7	76	-0.20
48		Acenaphthylene	2.004	2.099	-4.7	77	-0.22
49		Dimethylphthalate	1.415	1.448	-2.3	76	-0.20
50		2,6-Dinitrotoluene	0.290	0.300	-3.4	74	-0.21
51	C	Acenaphthene	1.134	1.134	0.0	74	-0.22
52		3-Nitroaniline	0.329	0.332	-0.9	68	-0.21
53	P	2,4-Dinitrophenol	0.128	0.114	10.9	66	-0.21
54		Dibenzofuran	1.640	1.707	-4.1	77	-0.22
55	P	4-Nitrophenol	0.252	0.270	-7.1	85	-0.19
56		2,4-Dinitrotoluene	0.386	0.414	-7.3	74	-0.21
57		Fluorene	1.378	1.443	-4.7	77	-0.22
58		2,3,4,6-Tetrachlorophenol	0.286	0.291	-1.7	72	-0.22
59		Diethylphthalate	1.445	1.486	-2.8	76	-0.20
60		4-Chlorophenyl-phenylether	0.596	0.624	-4.7	78	-0.22
61		4-Nitroaniline	0.346	0.358	-3.5	70	-0.22
62		Azobenzene	1.424	1.509	-6.0	79	-0.22
63	I	Phenanthrene-d10	1.000	1.000	0.0	71	-0.24
64		4,6-Dinitro-2-methylphenol	0.125	0.122	2.4	71	-0.21
65	c	n-Nitrosodiphenylamine	0.703	0.726	-3.3	74	-0.22
66		4-Bromophenyl-phenylether	0.195	0.210	-7.7	76	-0.22
67		Hexachlorobenzene	0.227	0.238	-4.8	75	-0.23
68		Atrazine	0.210	0.226	-7.6	75	-0.21
69	C	Pentachlorophenol	0.111	0.107	3.6	68	-0.23
70		Phenanthrene	1.207	1.212	-0.4	71	-0.24
71		Anthracene	1.184	1.218	-2.9	75	-0.24
72		Carbazole	1.148	1.131	1.5	69	-0.24
73		Di-n-butylphthalate	1.410	1.490	-5.7	75	-0.22
74	C	Fluoranthene	1.232	1.330	-8.0	79	-0.26
75	I	Chrysene-d12	1.000	1.000	0.0	72	-0.29
76		Benzidine	0.833	0.780	6.4	65	-0.26
77		Pyrene	1.399	1.341	4.1	71	-0.28
78	S	Terphenyl-d14	0.907	0.864	4.7	70	-0.26
79		Butylbenzylphthalate	0.669	0.685	-2.4	74	-0.26
80		Benzo(a)anthracene	1.245	1.267	-1.8	75	-0.29
81		3,3'-Dichlorobenzidine	0.416	0.418	-0.5	70	-0.28
82		Chrysene	1.140	1.200	-5.3	77	-0.29
83		Bis(2-ethylhexyl)phthalate	1.012	1.009	0.3	72	-0.24
84	c	Di-n-octyl phthalate	1.728	1.728	0.0	73	-0.29
85		Indeno(1,2,3-cd)pyrene	1.251	1.447	-15.7	89	-0.59#
86	I	Perylene-d12	1.000	1.000	0.0	84	-0.38
87		Benzo(b)fluoranthene	1.319	1.484	-12.5	94	-0.36

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.229	1.028	16.4	67	-0.36
89 C	Benzo(a)pyrene	1.177	1.186	-0.8	83	-0.38
90	Dibenzo(a,h)anthracene	1.167	1.239	-6.2	89	-0.58#
91	Benzo(g,h,i)perylene	1.135	1.213	-6.9	89	-0.66#

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