Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF123114\
Data File : BF076716.D
Acq On : 31 Dec 2014 22:41
Operator : IZ / TP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 01 07:04:42 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121714.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 01 03:10:49 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	71	-0.20
2		1,4-Dioxane	0.504	0.514	-2.0	72	-0.28
3		Pyridine	1.269	1.454	-14.6	84	-0.40
4		n-Nitrosodimethylamine	0.614	0.706	-15.0	81	-0.38
5	S	2-Fluorophenol	1.180	1.278	-8.3	78	-0.28
6		Aniline	2.058	2.245	-9.1	77	-0.20
7	S	Phenol-d6	1.441	1.576	-9.4	77	-0.18
8		2-Chlorophenol	1.351	1.470	-8.8	76	-0.20
9		Benzaldehyde	1.096	1.079	1.6	69	-0.22
10	C	Phenol	1.749	1.851	-5.8	74	-0.18
11		bis(2-Chloroethyl)ether	1.259	1.378	-9.5	78	-0.19
12		1,3-Dichlorobenzene	1.568	1.742	-11.1	78	-0.21
13	C	1,4-Dichlorobenzene	1.589	1.727	-8.7	79	-0.20
14		1,2-Dichlorobenzene	1.506	1.599	-6.2	77	-0.20
15		Benzyl Alcohol	1.248	1.370	-9.8	76	-0.19
16		2,2'-oxybis(1-Chloropropane	1.823	1.906	-4.6	74	-0.18
17		2-Methylphenol	1.110	1.190	-7.2	74	-0.16
18		Hexachloroethane	0.568	0.630	-10.9	80	-0.20
19	P	n-Nitroso-di-n-propylamine	1.052	1.128	-7.2	79	-0.18
20		3+4-Methylphenols	1.499	1.581	-5.5	74	-0.16
21	I	Naphthalene-d8	1.000	1.000	0.0	74	-0.20
22		Acetophenone	0.479	0.472	1.5	73	-0.19
23	S	Nitrobenzene-d5	0.336	0.371	-10.4	82	-0.19
24		Nitrobenzene	0.349	0.381	-9.2	80	-0.19
25		Isophorone	0.648	0.688	-6.2	79	-0.19
26	C	2-Nitrophenol	0.170	0.171	-0.6	72	-0.20
27		2,4-Dimethylphenol	0.276	0.291	-5.4	78	-0.16
28		bis(2-Chloroethoxy)methane	0.380	0.398	-4.7	75	-0.18
29	C	2,4-Dichlorophenol	0.271	0.274	-1.1	76	-0.19
30		1,2,4-Trichlorobenzene	0.292	0.308	-5.5	77	-0.20
31		Naphthalene	0.991	0.976	1.5	74	-0.20
32		Benzoic acid	0.153	0.143	6.5	68	-0.16
33		4-Chloroaniline	0.404	0.415	-2.7	74	-0.19
34	C	Hexachlorobutadiene	0.170	0.184	-8.2	83	-0.19
35		Caprolactam	0.078	0.077	1.3	75	-0.20
36	C	4-Chloro-3-methylphenol	0.301	0.302	-0.3	71	-0.19
37		2-Methylnaphthalene	0.691	0.697	-0.9	74	-0.21
38	I	Acenaphthene-d10	1.000	1.000	0.0	78	-0.22
39		1,2,4,5-Tetrachlorobenzene	0.492	0.455	7.5	71	-0.21
40	P	Hexachlorocyclopentadiene	0.241	0.223	7.5	68	-0.21
41	S	2,4,6-Tribromophenol	0.169	0.167	1.2	76	-0.23
42	C	2,4,6-Trichlorophenol	0.357	0.362	-1.4	77	-0.20
43		2,4,5-Trichlorophenol	0.384	0.372	3.1	73	-0.20
44	S	2-Fluorobiphenyl	1.285	1.253	2.5	77	-0.21

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.486	1.397	6.0	70	-0.21
46		2-Chloronaphthalene	1.178	1.180	-0.2	78	-0.21
47		2-Nitroaniline	0.358	0.385	-7.5	79	-0.20
48		Acenaphthylene	2.004	1.973	1.5	76	-0.22
49		Dimethylphthalate	1.415	1.385	2.1	77	-0.20
50		2,6-Dinitrotoluene	0.290	0.277	4.5	72	-0.21
51	C	Acenaphthene	1.134	1.097	3.3	76	-0.22
52		3-Nitroaniline	0.329	0.337	-2.4	73	-0.21
53	P	2,4-Dinitrophenol	0.128	0.100	21.9	61	-0.21
54		Dibenzofuran	1.640	1.585	3.4	75	-0.22
55	P	4-Nitrophenol	0.252	0.263	-4.4	87	-0.19
56		2,4-Dinitrotoluene	0.386	0.411	-6.5	77	-0.21
57		Fluorene	1.378	1.393	-1.1	79	-0.22
58		2,3,4,6-Tetrachlorophenol	0.286	0.273	4.5	71	-0.22
59		Diethylphthalate	1.445	1.453	-0.6	78	-0.20
60		4-Chlorophenyl-phenylether	0.596	0.570	4.4	75	-0.22
61		4-Nitroaniline	0.346	0.349	-0.9	72	-0.22
62		Azobenzene	1.424	1.481	-4.0	82	-0.23
63	I	Phenanthrene-d10	1.000	1.000	0.0	73	-0.24
64		4,6-Dinitro-2-methylphenol	0.125	0.120	4.0	71	-0.21
65	c	n-Nitrosodiphenylamine	0.703	0.727	-3.4	75	-0.22
66		4-Bromophenyl-phenylether	0.195	0.212	-8.7	79	-0.22
67		Hexachlorobenzene	0.227	0.241	-6.2	78	-0.23
68		Atrazine	0.210	0.231	-10.0	78	-0.21
69	C	Pentachlorophenol	0.111	0.111	0.0	71	-0.23
70		Phenanthrene	1.207	1.240	-2.7	74	-0.24
71		Anthracene	1.184	1.214	-2.5	77	-0.24
72		Carbazole	1.148	1.195	-4.1	74	-0.24
73		Di-n-butylphthalate	1.410	1.469	-4.2	75	-0.22
74	C	Fluoranthene	1.232	1.325	-7.5	80	-0.26
75	I	Chrysene-d12	1.000	1.000	0.0	77	-0.29
76		Benzidine	0.833	0.769	7.7	68	-0.26
77		Pyrene	1.399	1.354	3.2	76	-0.28
78	S	Terphenyl-d14	0.907	0.863	4.9	75	-0.26
79		Butylbenzylphthalate	0.669	0.657	1.8	76	-0.26
80		Benzo(a)anthracene	1.245	1.233	1.0	78	-0.29
81		3,3'-Dichlorobenzidine	0.416	0.393	5.5	71	-0.28
82		Chrysene	1.140	1.170	-2.6	80	-0.29
83		Bis(2-ethylhexyl)phthalate	1.012	0.971	4.1	74	-0.24
84	c	Di-n-octyl phthalate	1.728	1.679	2.8	76	-0.28
85		Indeno(1,2,3-cd)pyrene	1.251	1.403	-12.2	92	-0.53#
86	I	Perylene-d12	1.000	1.000	0.0	87	-0.36
87		Benzo(b)fluoranthene	1.319	1.320	-0.1	86	-0.34

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.229	1.170	4.8	79	-0.35
89 C	Benzo(a)pyrene	1.177	1.185	-0.7	85	-0.36
90	Dibenzo(a,h)anthracene	1.167	1.223	-4.8	91	-0.53#
91	Benzo(g,h,i)perylene	1.135	1.197	-5.5	90	-0.60#

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