Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121015\
Data File : BF083684.D
Acq On : 10 Dec 2015 19:01
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 17 11:22:12 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 17 09:56:43 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	81	-0.09
2		1,4-Dioxane	0.651	0.687	-5.5	89	-0.09
3		Pyridine	1.583	1.584	-0.1	82	-0.09
4		n-Nitrosodimethylamine	0.833	0.728	12.6	66	-0.08
5	S	2-Fluorophenol	1.319	1.366	-3.6	80	-0.08
6		Aniline	2.171	2.155	0.7	75	-0.08
7	S	Phenol-d6	1.762	1.780	-1.0	79	-0.09
8		2-Chlorophenol	1.442	1.467	-1.7	81	-0.09
9		Benzaldehyde	0.936	0.990	-5.8	83	-0.08
10	C	Phenol	2.134	2.173	-1.8	79	-0.08
11		bis(2-Chloroethyl)ether	1.570	1.604	-2.2	81	-0.08
12		1,3-Dichlorobenzene	1.607	1.636	-1.8	81	-0.08
13	C	1,4-Dichlorobenzene	1.649	1.648	0.1	81	-0.09
14		1,2-Dichlorobenzene	1.530	1.549	-1.2	80	-0.09
15		Benzyl Alcohol	1.254	1.237	1.4	74	-0.08
16		2,2'-oxybis(1-Chloropropane	2.903	2.834	2.4	77	-0.08
17		2-Methylphenol	1.173	1.171	0.2	77	-0.08
18		Hexachloroethane	0.582	0.579	0.5	78	-0.10
19	P	n-Nitroso-di-n-propylamine	1.230	1.212	1.5	79	-0.09
20		3+4-Methylphenols	1.545	1.456	5.8	76	-0.08
21	I	Naphthalene-d8	1.000	1.000	0.0	76	-0.10
22		Acetophenone	0.561	0.571	-1.8	76	-0.08
23	S	Nitrobenzene-d5	0.429	0.445	-3.7	79	-0.09
24		Nitrobenzene	0.474	0.490	-3.4	80	-0.09
25		Isophorone	0.832	0.870	-4.6	79	-0.09
26	C	2-Nitrophenol	0.188	0.203	-8.0	78	-0.09
27		2,4-Dimethylphenol	0.329	0.355	-7.9	81	-0.09
28		bis(2-Chloroethoxy)methane	0.462	0.480	-3.9	78	-0.09
29	C	2,4-Dichlorophenol	0.299	0.312	-4.3	77	-0.08
30		1,2,4-Trichlorobenzene	0.331	0.340	-2.7	78	-0.09
31		Naphthalene	1.059	1.108	-4.6	81	-0.09
32		Benzoic acid	0.185	0.162	12.4	62	-0.07
33		4-Chloroaniline	0.386	0.347	10.1	66	-0.07
34	C	Hexachlorobutadiene	0.193	0.203	-5.2	80	-0.10
35		Caprolactam	0.090	0.100	-11.1	80	-0.09
36	C	4-Chloro-3-methylphenol	0.336	0.365	-8.6	81	-0.08
37		2-Methylnaphthalene	0.665	0.710	-6.8	82	-0.09
38	I	Acenaphthene-d10	1.000	1.000	0.0	78	-0.09
39		1,2,4,5-Tetrachlorobenzene	0.657	0.678	-3.2	80	-0.09
40	P	Hexachlorocyclopentadiene	0.089	0.034#	61.8#	31#	-0.10
41	S	2,4,6-Tribromophenol	0.178	0.189	-6.2	81	-0.10
42	C	2,4,6-Trichlorophenol	0.391	0.388	0.8	75	-0.08
43		2,4,5-Trichlorophenol	0.387	0.369	4.7	71	-0.07
44	S	2-Fluorobiphenyl	1.421	1.443	-1.5	81	-0.09

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.738	1.817	-4.5	82	-0.09
46		2-Chloronaphthalene	1.312	1.369	-4.3	81	-0.09
47		2-Nitroaniline	0.473	0.499	-5.5	80	-0.08
48		Acenaphthylene	2.164	2.234	-3.2	81	-0.09
49		Dimethylphthalate	1.592	1.610	-1.1	80	-0.09
50		2,6-Dinitrotoluene	0.329	0.344	-4.6	80	-0.09
51	C	Acenaphthene	1.239	1.283	-3.6	81	-0.09
52		3-Nitroaniline	0.349	0.369	-5.7	78	-0.09
53	P	2,4-Dinitrophenol	0.140	0.155	-10.7	82	-0.05
54		Dibenzofuran	1.718	1.802	-4.9	81	-0.09
55	P	4-Nitrophenol	0.154	0.122	20.8	61	-0.03
56		2,4-Dinitrotoluene	0.436	0.476	-9.2	80	-0.08
57		Fluorene	1.500	1.550	-3.3	82	-0.09
58		2,3,4,6-Tetrachlorophenol	0.286	0.305	-6.6	76	-0.08
59		Diethylphthalate	1.548	1.564	-1.0	80	-0.09
60		4-Chlorophenyl-phenylether	0.714	0.743	-4.1	80	-0.10
61		4-Nitroaniline	0.317	0.356	-12.3	82	-0.08
62		Azobenzene	1.704	1.869	-9.7	84	-0.09
63	I	Phenanthrene-d10	1.000	1.000	0.0	79	-0.09
64		4,6-Dinitro-2-methylphenol	0.120	0.133	-10.8	77	-0.07
65	c	n-Nitrosodiphenylamine	0.660	0.679	-2.9	81	-0.09
66		4-Bromophenyl-phenylether	0.223	0.234	-4.9	82	-0.09
67		Hexachlorobenzene	0.237	0.244	-3.0	81	-0.09
68		Atrazine	0.196	0.205	-4.6	79	-0.10
69	C	Pentachlorophenol	0.052	0.030	42.3#	44#	-0.06
70		Phenanthrene	1.149	1.206	-5.0	83	-0.09
71		Anthracene	1.190	1.272	-6.9	84	-0.09
72		Carbazole	1.021	1.119	-9.6	87	-0.09
73		Di-n-butylphthalate	1.284	1.356	-5.6	82	-0.10
74	C	Fluoranthene	1.176	1.276	-8.5	85	-0.09
75	I	Chrysene-d12	1.000	1.000	0.0	85	-0.08
76		Benzidine	0.622	0.611	1.8	78	-0.08
77		Pyrene	1.440	1.466	-1.8	86	-0.09
78	S	Terphenyl-d14	0.850	0.860	-1.2	88	-0.10
79		Butylbenzylphthalate	0.639	0.667	-4.4	86	-0.09
80		Benzo(a)anthracene	1.171	1.187	-1.4	87	-0.08
81		3,3'-Dichlorobenzidine	0.395	0.442	-11.9	91	-0.08
82		Chrysene	1.174	1.256	-7.0	91	-0.08
83		Bis(2-ethylhexyl)phthalate	0.887	0.914	-3.0	84	-0.08
84	c	Di-n-octyl phthalate	1.355	1.447	-6.8	89	-0.08
85		Indeno(1,2,3-cd)pyrene	0.792	0.878	-10.9	100	-0.09
86	I	Perylene-d12	1.000	1.000	0.0	92	-0.06
87		Benzo(b)fluoranthene	1.139	1.269	-11.4	104	-0.07

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.255	1.243	1.0	89	-0.07
89 C	Benzo(a)pyrene	1.111	1.169	-5.2	98	-0.07
90	Dibenzo(a,h)anthracene	0.850	0.912	-7.3	103	-0.10
91	Benzo(g,h,i)perylene	0.828	0.817	1.3	96	-0.09

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