Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF010318\
Data File : BF101855.D
Acq On : 3 Jan 2018 12:59
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 04 01:14:06 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 29 13:21:45 2017
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	-0.01
2		1,4-Dioxane	0.690	0.623	9.7	64	-0.01
3		Pyridine	1.917	1.640	14.4	60	-0.01
4		n-Nitrosodimethylamine	0.883	0.721	18.3	57	-0.02
5	S	2-Fluorophenol	1.343	1.392	-3.6	74	-0.01
6		Aniline	2.322	2.094	9.8	65	-0.02
7	S	Phenol-d6	1.671	1.562	6.5	69	-0.01
8		2-Chlorophenol	1.412	1.385	1.9	72	-0.01
9		Benzaldehyde	1.234	0.994	19.4	58	0.00
10	C	Phenol	1.905	1.726	9.4	66	-0.02
11		bis(2-Chloroethyl)ether	1.494	1.382	7.5	66	-0.01
12		1,3-Dichlorobenzene	1.594	1.565	1.8	71	-0.01
13	C	1,4-Dichlorobenzene	1.628	1.623	0.3	73	-0.01
14		1,2-Dichlorobenzene	1.465	1.431	2.3	71	-0.01
15		Benzyl Alcohol	1.150	1.127	2.0	72	-0.01
16		2,2'-oxybis(1-Chloropropane	2.286	2.110	7.7	65	-0.02
17		2-Methylphenol	1.299	1.191	8.3	69	-0.01
18		Hexachloroethane	0.566	0.542	4.2	69	-0.02
19	P	n-Nitroso-di-n-propylamine	1.097	1.060	3.4	74	-0.02
20		3+4-Methylphenols	1.377	1.579	-14.7	83	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	71	-0.02
22		Acetophenone	0.456	0.470	-3.1	74	-0.01
23	S	Nitrobenzene-d5	0.359	0.347	3.3	68	-0.02
24		Nitrobenzene	0.398	0.381	4.3	69	-0.02
25		Isophorone	0.721	0.655	9.2	66	-0.01
26	C	2-Nitrophenol	0.171	0.188	-9.9	75	-0.01
27		2,4-Dimethylphenol	0.290	0.283	2.4	71	-0.01
28		bis(2-Chloroethoxy)methane	0.458	0.435	5.0	69	-0.01
29	C	2,4-Dichlorophenol	0.287	0.293	-2.1	73	-0.01
30		1,2,4-Trichlorobenzene	0.303	0.289	4.6	69	-0.01
31		Naphthalene	1.007	0.967	4.0	70	-0.01
32		Benzoic acid	0.173	0.202	-16.8	85	-0.01
33		4-Chloroaniline	0.438	0.424	3.2	71	-0.01
34	C	Hexachlorobutadiene	0.175	0.172	1.7	71	-0.02
35		Caprolactam	0.100	0.095	5.0	68	-0.02
36	C	4-Chloro-3-methylphenol	0.315	0.315	0.0	71	-0.02
37		2-Methylnaphthalene	0.656	0.633	3.5	71	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	71	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.675	0.664	1.6	72	-0.02
40	P	Hexachlorocyclopentadiene	0.088	0.079	10.2	67	-0.02
41	S	2,4,6-Tribromophenol	0.193	0.201	-4.1	75	-0.02
42	C	2,4,6-Trichlorophenol	0.407	0.398	2.2	69	-0.02
43		2,4,5-Trichlorophenol	0.445	0.382	14.2	61	-0.01
44	S	2-Fluorobiphenyl	1.289	1.253	2.8	71	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.774	1.708	3.7	72	-0.02
46		2-Chloronaphthalene	1.357	1.286	5.2	71	-0.02
47		2-Nitroaniline	0.469	0.471	-0.4	72	-0.02
48		Acenaphthylene	2.144	2.031	5.3	71	-0.02
49		Dimethylphthalate	1.609	1.528	5.0	71	-0.02
50		2,6-Dinitrotoluene	0.327	0.319	2.4	69	-0.02
51	C	Acenaphthene	1.288	1.219	5.4	71	-0.02
52		3-Nitroaniline	0.408	0.419	-2.7	73	-0.02
53	P	2,4-Dinitrophenol	0.087	0.134	-54.0#	108	0.02
54		Dibenzofuran	1.637	1.749	-6.8	78	-0.02
55	P	4-Nitrophenol	0.178	0.199	-11.8	77	-0.01
56		2,4-Dinitrotoluene	0.368	0.459	-24.7	89	-0.02
57		Fluorene	1.385	1.419	-2.5	74	-0.02
58		2,3,4,6-Tetrachlorophenol	0.320	0.323	-0.9	72	-0.02
59		Diethylphthalate	1.593	1.546	3.0	74	-0.01
60		4-Chlorophenyl-phenylether	0.664	0.694	-4.5	75	-0.02
61		4-Nitroaniline	0.410	0.405	1.2	71	-0.02
62		Azobenzene	1.650	1.505	8.8	69	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	73	-0.02
64		4,6-Dinitro-2-methylphenol	0.102	0.114	-11.8	78	-0.01
65	c	n-Nitrosodiphenylamine	0.738	0.690	6.5	71	-0.02
66		4-Bromophenyl-phenylether	0.225	0.222	1.3	74	-0.02
67		Hexachlorobenzene	0.238	0.232	2.5	73	-0.02
68		Atrazine	0.220	0.204	7.3	69	-0.02
69	C	Pentachlorophenol	0.079	0.069	12.7	64	-0.02
70		Phenanthrene	1.112	1.067	4.0	74	-0.02
71		Anthracene	1.136	1.101	3.1	75	-0.02
72		Carbazole	1.064	1.051	1.2	76	-0.02
73		Di-n-butylphthalate	1.291	1.274	1.3	76	-0.02
74	C	Fluoranthene	1.167	1.141	2.2	76	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	71	-0.02
76		Benzidine	0.845	0.690	18.3	58	-0.02
77		Pyrene	1.501	1.533	-2.1	75	-0.02
78	S	Terphenyl-d14	0.830	0.831	-0.1	76	-0.02
79		Butylbenzylphthalate	0.679	0.700	-3.1	74	-0.02
80		Benzo(a)anthracene	1.321	1.250	5.4	69	-0.02
81		3,3'-Dichlorobenzidine	0.431	0.419	2.8	70	-0.02
82		Chrysene	1.247	1.221	2.1	72	-0.02
83		Bis(2-ethylhexyl)phthalate	0.860	0.853	0.8	72	-0.02
84	c	Di-n-octyl phthalate	1.534	1.505	1.9	72	-0.02
85		Indeno(1,2,3-cd)pyrene	1.110	0.974	12.3	66	-0.04
86	I	Perylene-d12	1.000	1.000	0.0	66	-0.02
87		Benzo(b)fluoranthene	1.219	1.101	9.7	60	-0.03

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.185	1.244	-5.0	74	-0.02
89 C	Benzo(a)pyrene	1.124	1.061	5.6	65	-0.02
90	Dibenzo(a,h)anthracene	0.965	0.910	5.7	66	-0.04
91	Benzo(g,h,i)perylene	0.955	0.918	3.9	67	-0.04

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