Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110116\
Data File : BF090948.D
Acq On : 1 Nov 2016 14:05
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 02 11:34:49 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 01 11:53:50 2016
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	122	-0.01
2		1,4-Dioxane	0.585	0.700	-19.7	141	0.01
3		Pyridine	1.586	1.639	-3.3	118	0.00
4		n-Nitrosodimethylamine	0.447	0.585	-30.9#	162#	0.01
5	S	2-Fluorophenol	1.144	1.260	-10.1	135	0.00
6		Aniline	1.763	1.919	-8.8	125	0.00
7	S	Phenol-d6	1.543	1.610	-4.3	121	-0.01
8		2-Chlorophenol	1.411	1.365	3.3	109	-0.01
9		Benzaldehyde	0.797	0.881	-10.5	134	-0.01
10	C	Phenol	1.701	1.776	-4.4	129	0.00
11		bis(2-Chloroethyl)ether	1.407	1.495	-6.3	117	0.00
12		1,3-Dichlorobenzene	1.443	1.499	-3.9	119	-0.01
13	C	1,4-Dichlorobenzene	1.518	1.468	3.3	109	0.00
14		1,2-Dichlorobenzene	1.455	1.355	6.9	113	-0.01
15		Benzyl Alcohol	1.014	1.117	-10.2	121	-0.01
16		2,2'-oxybis(1-Chloropropane	1.400	1.953	-39.5#	147	-0.01
17		2-Methylphenol	1.073	1.152	-7.4	120	0.00
18		Hexachloroethane	0.552	0.585	-6.0	129	-0.01
19	P	n-Nitroso-di-n-propylamine	0.892	1.076	-20.6	130	0.00
20		3+4-Methylphenols	1.243	1.400	-12.6	121	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	130	-0.01
22		Acetophenone	0.411	0.422	-2.7	125	0.00
23	S	Nitrobenzene-d5	0.306	0.340	-11.1	132	-0.01
24		Nitrobenzene	0.323	0.392	-21.4	159#	-0.01
25		Isophorone	0.620	0.654	-5.5	136	-0.01
26	C	2-Nitrophenol	0.173	0.172	0.6	104	0.00
27		2,4-Dimethylphenol	0.280	0.303	-8.2	119	0.00
28		bis(2-Chloroethoxy)methane	0.353	0.404	-14.4	131	0.00
29	C	2,4-Dichlorophenol	0.260	0.246	5.4	112	-0.01
30		1,2,4-Trichlorobenzene	0.300	0.279	7.0	114	-0.01
31		Naphthalene	0.935	0.929	0.6	135	-0.01
32		Benzoic acid	0.203	0.214	-5.4	126	0.02
33		4-Chloroaniline	0.378	0.390	-3.2	115	0.00
34	C	Hexachlorobutadiene	0.176	0.144	18.2	94	-0.01
35		Caprolactam	0.081	0.083	-2.5	115	0.01
36	C	4-Chloro-3-methylphenol	0.282	0.306	-8.5	130	0.00
37		2-Methylnaphthalene	0.604	0.536	11.3	102	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	107	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.549	0.568	-3.5	118	-0.01
40	P	Hexachlorocyclopentadiene	0.252	0.243	3.6	106	-0.01
41	S	2,4,6-Tribromophenol	0.206	0.193	6.3	104	0.00
42	C	2,4,6-Trichlorophenol	0.354	0.361	-2.0	115	-0.01
43		2,4,5-Trichlorophenol	0.365	0.375	-2.7	104	0.00
44	S	2-Fluorobiphenyl	1.139	1.124	1.3	109	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.441	1.411	2.1	105	-0.01
46		2-Chloronaphthalene	1.096	1.099	-0.3	106	-0.01
47		2-Nitroaniline	0.309	0.462	-49.5#	164#	0.00
48		Acenaphthylene	1.659	1.584	4.5	98	-0.01
49		Dimethylphthalate	1.336	1.204	9.9	99	0.00
50		2,6-Dinitrotoluene	0.295	0.276	6.4	92	-0.01
51	C	Acenaphthene	1.143	1.098	3.9	104	-0.01
52		3-Nitroaniline	0.314	0.315	-0.3	104	0.00
53	P	2,4-Dinitrophenol	0.155	0.152	1.9	104	0.00
54		Dibenzofuran	1.475	1.336	9.4	97	-0.01
55	P	4-Nitrophenol	0.192	0.217	-13.0	114	0.00
56		2,4-Dinitrotoluene	0.367	0.408	-11.2	110	0.00
57		Fluorene	1.208	1.070	11.4	92	-0.01
58		2,3,4,6-Tetrachlorophenol	0.326	0.314	3.7	111	-0.01
59		Diethylphthalate	1.327	1.349	-1.7	117	0.00
60		4-Chlorophenyl-phenylether	0.590	0.577	2.2	115	0.00
61		4-Nitroaniline	0.311	0.342	-10.0	116	0.00
62		Azobenzene	1.305	1.530	-17.2	146	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	116	0.00
64		4,6-Dinitro-2-methylphenol	0.133	0.128	3.8	99	0.00
65	c	n-Nitrosodiphenylamine	0.613	0.581	5.2	109	-0.01
66		4-Bromophenyl-phenylether	0.219	0.183	16.4	97	-0.01
67		Hexachlorobenzene	0.259	0.220	15.1	91	-0.01
68		Atrazine	0.190	0.171	10.0	123	0.00
69	C	Pentachlorophenol	0.139	0.114	18.0	90	-0.01
70		Phenanthrene	1.018	0.936	8.1	102	-0.01
71		Anthracene	1.072	1.058	1.3	121	0.00
72		Carbazole	0.947	0.869	8.2	106	-0.01
73		Di-n-butylphthalate	1.221	1.212	0.7	119	0.00
74	C	Fluoranthene	1.115	1.091	2.2	111	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
76		Benzidine	0.437	0.479	-9.6	112	0.00
77		Pyrene	1.266	1.379	-8.9	109	0.00
78	S	Terphenyl-d14	0.794	0.839	-5.7	104	0.00
79		Butylbenzylphthalate	0.574	0.601	-4.7	112	-0.01
80		Benzo(a)anthracene	1.061	1.098	-3.5	110	-0.01
81		3,3'-Dichlorobenzidine	0.412	0.444	-7.8	103	0.00
82		Chrysene	1.074	1.111	-3.4	110	0.00
83		Bis(2-ethylhexyl)phthalate	0.748	0.847	-13.2	115	0.00
84	c	Di-n-octyl phthalate	1.256	1.359	-8.2	115	-0.01
85		Indeno(1,2,3-cd)pyrene	0.951	0.829	12.8	99	0.00
86	I	Perylene-d12	1.000	1.000	0.0	96	0.00
87		Benzo(b)fluoranthene	1.345	1.618	-20.3	101	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.063	0.996	6.3	93	0.00
89 C	Benzo(a)pyrene	1.156	1.174	-1.6	92	0.00
90	Dibenzo(a,h)anthracene	0.978	1.009	-3.2	98	0.00
91	Benzo(g,h,i)perylene	0.921	0.911	1.1	96	0.00

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