Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011017\
Data File : BF092174.D
Acq On : 10 Jan 2017 11:45
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 10 14:45:14 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 06 15:43:51 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	95	-0.01
2		1,4-Dioxane	0.590	0.522	11.5	83	0.02
3		Pyridine	2.058	1.663	19.2	75	0.01
4		n-Nitrosodimethylamine	0.776	0.689	11.2	81	0.03
5	S	2-Fluorophenol	1.198	1.080	9.8	89	0.00
6		Aniline	1.899	2.042	-7.5	102	-0.02
7	S	Phenol-d6	1.462	1.500	-2.6	96	0.00
8		2-Chlorophenol	1.362	1.394	-2.3	95	-0.01
9		Benzaldehyde	0.904	0.871	3.7	98	-0.01
10	C	Phenol	1.666	1.967	-18.1	103	0.00
11		bis(2-Chloroethyl)ether	1.326	1.431	-7.9	95	-0.01
12		1,3-Dichlorobenzene	1.455	1.469	-1.0	91	-0.02
13	C	1,4-Dichlorobenzene	1.480	1.405	5.1	89	-0.02
14		1,2-Dichlorobenzene	1.285	1.332	-3.7	100	-0.01
15		Benzyl Alcohol	1.136	1.048	7.7	87	-0.01
16		2,2'-oxybis(1-Chloropropane	1.975	2.059	-4.3	99	-0.02
17		2-Methylphenol	1.096	1.117	-1.9	97	0.00
18		Hexachloroethane	0.549	0.551	-0.4	91	-0.02
19	P	n-Nitroso-di-n-propylamine	0.973	0.980	-0.7	91	-0.01
20		3+4-Methylphenols	1.307	1.512	-15.7	104	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	97	-0.01
22		Acetophenone	0.493	0.496	-0.6	96	-0.01
23	S	Nitrobenzene-d5	0.360	0.367	-1.9	103	-0.01
24		Nitrobenzene	0.383	0.369	3.7	84	-0.01
25		Isophorone	0.664	0.668	-0.6	98	-0.01
26	C	2-Nitrophenol	0.189	0.206	-9.0	108	-0.01
27		2,4-Dimethylphenol	0.313	0.298	4.8	85	-0.01
28		bis(2-Chloroethoxy)methane	0.402	0.393	2.2	92	-0.01
29	C	2,4-Dichlorophenol	0.265	0.266	-0.4	95	-0.01
30		1,2,4-Trichlorobenzene	0.291	0.285	2.1	90	-0.01
31		Naphthalene	0.915	0.980	-7.1	94	-0.01
32		Benzoic acid	0.232	0.202	12.9	79	0.03
33		4-Chloroaniline	0.413	0.414	-0.2	95	-0.01
34	C	Hexachlorobutadiene	0.153	0.155	-1.3	95	-0.02
35		Caprolactam	0.087	0.087	0.0	95	0.01
36	C	4-Chloro-3-methylphenol	0.305	0.320	-4.9	94	0.01
37		2-Methylnaphthalene	0.628	0.562	10.5	95	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	98	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.505	0.491	2.8	85	-0.01
40	P	Hexachlorocyclopentadiene	0.199	0.179	10.1	76	-0.01
41	S	2,4,6-Tribromophenol	0.166	0.176	-6.0	107	0.00
42	C	2,4,6-Trichlorophenol	0.353	0.361	-2.3	90	0.00
43		2,4,5-Trichlorophenol	0.364	0.345	5.2	89	0.00
44	S	2-Fluorobiphenyl	1.042	1.140	-9.4	110	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.369	1.336	2.4	84	-0.01
46		2-Chloronaphthalene	1.084	1.055	2.7	90	-0.01
47		2-Nitroaniline	0.386	0.373	3.4	83	-0.01
48		Acenaphthylene	1.668	1.567	6.1	91	-0.01
49		Dimethylphthalate	1.279	1.385	-8.3	100	0.00
50		2,6-Dinitrotoluene	0.280	0.290	-3.6	100	0.00
51	C	Acenaphthene	1.131	0.989	12.6	86	-0.01
52		3-Nitroaniline	0.346	0.408	-17.9	100	0.00
53	P	2,4-Dinitrophenol	0.156	0.179	-14.7	98	0.00
54		Dibenzofuran	1.527	1.342	12.1	92	-0.01
55	P	4-Nitrophenol	0.235	0.237	-0.9	89	0.01
56		2,4-Dinitrotoluene	0.351	0.320	8.8	91	-0.01
57		Fluorene	1.111	1.037	6.7	87	-0.01
58		2,3,4,6-Tetrachlorophenol	0.288	0.288	0.0	90	0.00
59		Diethylphthalate	1.242	1.218	1.9	87	-0.01
60		4-Chlorophenyl-phenylether	0.486	0.513	-5.6	103	0.00
61		4-Nitroaniline	0.359	0.398	-10.9	96	0.00
62		Azobenzene	1.368	1.447	-5.8	104	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	96	-0.01
64		4,6-Dinitro-2-methylphenol	0.121	0.153	-26.4#	113	0.01
65	c	n-Nitrosodiphenylamine	0.593	0.656	-10.6	106	0.00
66		4-Bromophenyl-phenylether	0.208	0.201	3.4	93	-0.01
67		Hexachlorobenzene	0.222	0.209	5.9	88	0.00
68		Atrazine	0.191	0.151	20.9	76	-0.01
69	C	Pentachlorophenol	0.109	0.102	6.4	79	0.00
70		Phenanthrene	1.084	1.129	-4.2	96	0.00
71		Anthracene	1.086	0.952	12.3	93	-0.01
72		Carbazole	1.005	0.942	6.3	98	-0.01
73		Di-n-butylphthalate	1.174	1.250	-6.5	110	0.00
74	C	Fluoranthene	1.082	1.003	7.3	99	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	94	0.00
76		Benzidine	0.580	0.591	-1.9	101	-0.01
77		Pyrene	1.467	1.394	5.0	96	0.00
78	S	Terphenyl-d14	0.825	0.820	0.6	103	-0.01
79		Butylbenzylphthalate	0.667	0.763	-14.4	100	-0.01
80		Benzo(a)anthracene	1.129	1.262	-11.8	106	0.00
81		3,3'-Dichlorobenzidine	0.428	0.553	-29.2#	129	0.00
82		Chrysene	1.132	1.340	-18.4	123	0.01
83		Bis(2-ethylhexyl)phthalate	0.786	0.870	-10.7	105	-0.01
84	c	Di-n-octyl phthalate	1.456	1.270	12.8	83	0.00
85		Indeno(1,2,3-cd)pyrene	0.981	0.868	11.5	85	0.01
86	I	Perylene-d12	1.000	1.000	0.0	89	0.01
87		Benzo(b)fluoranthene	1.245	1.188	4.6	79	0.00

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88	Benzo(k)fluoranthene	1.062	1.197	-12.7	103	0.01
89 C	Benzo(a)pyrene	1.105	1.137	-2.9	89	0.00
90	Dibenzo(a,h)anthracene	0.908	0.908	0.0	87	0.01
91	Benzo(g,h,i)perylene	0.945	0.927	1.9	85	0.02

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