Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122816\
Data File : BF092014.D
Acq On : 29 Dec 2016 00:21
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 29 02:45:31 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 28 17:32:10 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	111	0.00
2		1,4-Dioxane	0.590	0.526	10.8	98	0.01
3		Pyridine	2.058	1.664	19.1	87	0.01
4		n-Nitrosodimethylamine	0.776	0.673	13.3	92	0.01
5	S	2-Fluorophenol	1.198	1.289	-7.6	124	0.00
6		Aniline	1.899	1.672	12.0	98	0.00
7	S	Phenol-d6	1.462	1.285	12.1	96	0.00
8		2-Chlorophenol	1.362	1.393	-2.3	111	0.00
9		Benzaldehyde	0.904	0.864	4.4	113	0.00
10	C	Phenol	1.666	1.737	-4.3	106	0.00
11		bis(2-Chloroethyl)ether	1.326	1.380	-4.1	107	0.00
12		1,3-Dichlorobenzene	1.455	1.464	-0.6	106	0.00
13	C	1,4-Dichlorobenzene	1.480	1.535	-3.7	113	0.00
14		1,2-Dichlorobenzene	1.285	1.271	1.1	112	0.01
15		Benzyl Alcohol	1.136	1.112	2.1	107	0.01
16		2,2'-oxybis(1-Chloropropane	1.975	2.213	-12.1	124	-0.01
17		2-Methylphenol	1.096	1.208	-10.2	122	0.00
18		Hexachloroethane	0.549	0.604	-10.0	116	0.00
19	P	n-Nitroso-di-n-propylamine	0.973	1.096	-12.6	118	0.00
20		3+4-Methylphenols	1.307	1.527	-16.8	123	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
22		Acetophenone	0.493	0.589	-19.5	126	0.00
23	S	Nitrobenzene-d5	0.360	0.388	-7.8	120	0.00
24		Nitrobenzene	0.383	0.437	-14.1	110	0.01
25		Isophorone	0.664	0.709	-6.8	114	0.00
26	C	2-Nitrophenol	0.189	0.198	-4.8	115	0.00
27		2,4-Dimethylphenol	0.313	0.295	5.8	93	0.00
28		bis(2-Chloroethoxy)methane	0.402	0.392	2.5	101	0.00
29	C	2,4-Dichlorophenol	0.265	0.266	-0.4	104	0.00
30		1,2,4-Trichlorobenzene	0.291	0.295	-1.4	103	0.00
31		Naphthalene	0.915	0.885	3.3	94	0.00
32		Benzoic acid	0.232	0.229	1.3	99	0.00
33		4-Chloroaniline	0.413	0.370	10.4	94	0.00
34	C	Hexachlorobutadiene	0.153	0.172	-12.4	117	0.00
35		Caprolactam	0.087	0.066	24.1	79	0.00
36	C	4-Chloro-3-methylphenol	0.305	0.245	19.7	80	0.00
37		2-Methylnaphthalene	0.628	0.562	10.5	105	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	95	0.00
39		1,2,4,5-Tetrachlorobenzene	0.505	0.480	5.0	81	0.00
40	P	Hexachlorocyclopentadiene	0.199	0.153	23.1	63	0.00
41	S	2,4,6-Tribromophenol	0.166	0.161	3.0	95	0.00
42	C	2,4,6-Trichlorophenol	0.353	0.361	-2.3	88	0.00
43		2,4,5-Trichlorophenol	0.364	0.350	3.8	87	0.00
44	S	2-Fluorobiphenyl	1.042	1.087	-4.3	102	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.369	1.505	-9.9	92	0.01
46		2-Chloronaphthalene	1.084	1.095	-1.0	91	0.00
47		2-Nitroaniline	0.386	0.433	-12.2	94	0.00
48		Acenaphthylene	1.668	1.709	-2.5	96	0.01
49		Dimethylphthalate	1.279	1.310	-2.4	92	0.00
50		2,6-Dinitrotoluene	0.280	0.258	7.9	86	0.00
51	C	Acenaphthene	1.131	1.151	-1.8	97	0.00
52		3-Nitroaniline	0.346	0.290	16.2	69	0.00
53	P	2,4-Dinitrophenol	0.156	0.085	45.5#	45#	0.00
54		Dibenzofuran	1.527	1.436	6.0	96	0.00
55	P	4-Nitrophenol	0.235	0.203	13.6	74	0.00
56		2,4-Dinitrotoluene	0.351	0.309	12.0	85	0.00
57		Fluorene	1.111	1.118	-0.6	91	0.01
58		2,3,4,6-Tetrachlorophenol	0.288	0.299	-3.8	91	0.00
59		Diethylphthalate	1.242	1.351	-8.8	94	0.01
60		4-Chlorophenyl-phenylether	0.486	0.490	-0.8	95	0.00
61		4-Nitroaniline	0.359	0.310	13.6	73	0.00
62		Azobenzene	1.368	1.448	-5.8	101	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	86	0.00
64		4,6-Dinitro-2-methylphenol	0.121	0.100	17.4	67	0.00
65	c	n-Nitrosodiphenylamine	0.593	0.676	-14.0	99	0.00
66		4-Bromophenyl-phenylether	0.208	0.240	-15.4	100	0.00
67		Hexachlorobenzene	0.222	0.223	-0.5	84	0.00
68		Atrazine	0.191	0.196	-2.6	89	0.01
69	C	Pentachlorophenol	0.109	0.105	3.7	74	0.00
70		Phenanthrene	1.084	0.997	8.0	76	0.00
71		Anthracene	1.086	1.145	-5.4	101	0.00
72		Carbazole	1.005	0.962	4.3	90	0.00
73		Di-n-butylphthalate	1.174	1.176	-0.2	93	0.00
74	C	Fluoranthene	1.082	1.008	6.8	90	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	58	-0.01
76		Benzidine	0.580	0.564	2.8	60	0.00
77		Pyrene	1.467	1.983	-35.2#	84	0.00
78	S	Terphenyl-d14	0.825	1.142	-38.4#	88	0.00
79		Butylbenzylphthalate	0.667	0.945	-41.7#	77	0.00
80		Benzo(a)anthracene	1.129	1.297	-14.9	67	0.00
81		3,3'-Dichlorobenzidine	0.428	0.550	-28.5#	79	0.00
82		Chrysene	1.132	1.432	-26.5#	81	0.00
83		Bis(2-ethylhexyl)phthalate	0.786	1.124	-43.0#	84	0.00
84	c	Di-n-octyl phthalate	1.456	1.827	-25.5#	74	0.00
85		Indeno(1,2,3-cd)pyrene	0.981	1.620	-65.1#	99	0.00
86	I	Perylene-d12	1.000	1.000	0.0	83	0.00
87		Benzo(b)fluoranthene	1.245	1.121	10.0	70	0.00

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88	Benzo(k)fluoranthene	1.062	1.059	0.3	86	0.00
89 C	Benzo(a)pyrene	1.105	1.093	1.1	81	0.00
90	Dibenzo(a,h)anthracene	0.908	1.113	-22.6	100	0.00
91	Benzo(g,h,i)perylene	0.945	1.152	-21.9	100	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1