Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF042816\
Data File : BF086470.D
Acq On : 28 Apr 2016 17:08
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 29 01:03:28 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 27 14:20:52 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	140	-0.01
2		1,4-Dioxane	0.609	0.593	2.6	144	-0.01
3		Pyridine	1.442	1.433	0.6	146	-0.01
4		n-Nitrosodimethylamine	0.822	0.951	-15.7	151#	0.00
5	S	2-Fluorophenol	1.160	1.262	-8.8	157#	0.00
6		Aniline	1.958	2.053	-4.9	139	0.00
7	S	Phenol-d6	1.618	1.576	2.6	142	0.00
8		2-Chlorophenol	1.444	1.382	4.3	136	-0.01
9		Benzaldehyde	0.941	0.879	6.6	139	-0.01
10	C	Phenol	1.805	1.679	7.0	145	0.00
11		bis(2-Chloroethyl)ether	1.561	1.613	-3.3	153#	0.00
12		1,3-Dichlorobenzene	1.553	1.503	3.2	144	-0.01
13	C	1,4-Dichlorobenzene	1.603	1.544	3.7	147	0.00
14		1,2-Dichlorobenzene	1.470	1.246	15.2	118	-0.01
15		Benzyl Alcohol	1.024	1.061	-3.6	135	-0.01
16		2,2'-oxybis(1-Chloropropane	3.124	2.948	5.6	135	-0.01
17		2-Methylphenol	1.167	1.190	-2.0	142	-0.01
18		Hexachloroethane	0.599	0.576	3.8	141	-0.01
19	P	n-Nitroso-di-n-propylamine	1.070	1.141	-6.6	163#	0.00
20		3+4-Methylphenols	1.333	1.198	10.1	125	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	147	-0.01
22		Acetophenone	0.439	0.423	3.6	143	0.00
23	S	Nitrobenzene-d5	0.371	0.367	1.1	145	-0.01
24		Nitrobenzene	0.382	0.392	-2.6	157#	0.00
25		Isophorone	0.714	0.711	0.4	149	-0.01
26	C	2-Nitrophenol	0.176	0.170	3.4	134	-0.01
27		2,4-Dimethylphenol	0.308	0.284	7.8	131	-0.01
28		bis(2-Chloroethoxy)methane	0.389	0.409	-5.1	158#	0.00
29	C	2,4-Dichlorophenol	0.260	0.273	-5.0	154#	0.00
30		1,2,4-Trichlorobenzene	0.302	0.289	4.3	145	0.00
31		Naphthalene	1.018	0.934	8.3	150	0.00
32		Benzoic acid	0.160	0.206	-28.7#	166#	0.00
33		4-Chloroaniline	0.381	0.363	4.7	135	-0.01
34	C	Hexachlorobutadiene	0.169	0.150	11.2	132	-0.01
35		Caprolactam	0.087	0.084	3.4	135	0.00
36	C	4-Chloro-3-methylphenol	0.298	0.284	4.7	142	0.00
37		2-Methylnaphthalene	0.601	0.523	13.0	125	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	141	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.572	0.561	1.9	152#	0.00
40	P	Hexachlorocyclopentadiene	0.284	0.278	2.1	143	0.00
41	S	2,4,6-Tribromophenol	0.211	0.198	6.2	129	-0.01
42	C	2,4,6-Trichlorophenol	0.362	0.389	-7.5	155#	0.00
43		2,4,5-Trichlorophenol	0.394	0.362	8.1	130	-0.01
44	S	2-Fluorobiphenyl	1.182	1.057	10.6	134	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.672	1.455	13.0	130	-0.01
46		2-Chloronaphthalene	1.271	1.128	11.3	131	-0.01
47		2-Nitroaniline	0.408	0.472	-15.7	159#	0.00
48		Acenaphthylene	1.987	1.686	15.1	126	-0.01
49		Dimethylphthalate	1.506	1.378	8.5	144	0.00
50		2,6-Dinitrotoluene	0.311	0.288	7.4	125	-0.01
51	C	Acenaphthene	1.276	1.039	18.6	120	-0.01
52		3-Nitroaniline	0.354	0.380	-7.3	156#	0.00
53	P	2,4-Dinitrophenol	0.139	0.164	-18.0	158#	0.00
54		Dibenzofuran	1.595	1.336	16.2	122	-0.01
55	P	4-Nitrophenol	0.232	0.237	-2.2	144	0.00
56		2,4-Dinitrotoluene	0.397	0.358	9.8	119	-0.01
57		Fluorene	1.385	1.064	23.2	118	-0.01
58		2,3,4,6-Tetrachlorophenol	0.308	0.316	-2.6	145	0.00
59		Diethylphthalate	1.425	1.433	-0.6	162#	0.00
60		4-Chlorophenyl-phenylether	0.629	0.578	8.1	152#	0.00
61		4-Nitroaniline	0.346	0.339	2.0	137	0.00
62		Azobenzene	1.561	1.525	2.3	147	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	149	0.00
64		4,6-Dinitro-2-methylphenol	0.117	0.121	-3.4	147	0.00
65	c	n-Nitrosodiphenylamine	0.625	0.572	8.5	131	-0.01
66		4-Bromophenyl-phenylether	0.206	0.200	2.9	135	-0.01
67		Hexachlorobenzene	0.239	0.234	2.1	147	0.00
68		Atrazine	0.188	0.200	-6.4	156#	0.00
69	C	Pentachlorophenol	0.126	0.134	-6.3	147	0.00
70		Phenanthrene	1.050	0.933	11.1	131	-0.01
71		Anthracene	1.120	1.051	6.2	144	-0.01
72		Carbazole	0.993	0.869	12.5	127	-0.01
73		Di-n-butylphthalate	1.125	0.970	13.8	121	-0.01
74	C	Fluoranthene	1.059	0.858	19.0	121	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	112	-0.01
76		Benzidine	0.687	0.526	23.4	82	-0.01
77		Pyrene	1.441	1.510	-4.8	115	-0.01
78	S	Terphenyl-d14	0.908	0.913	-0.6	119	-0.01
79		Butylbenzylphthalate	0.624	0.642	-2.9	115	0.00
80		Benzo(a)anthracene	1.066	0.973	8.7	109	-0.01
81		3,3'-Dichlorobenzidine	0.713	0.656	8.0	104	-0.01
82		Chrysene	1.159	1.075	7.2	108	-0.01
83		Bis(2-ethylhexyl)phthalate	0.775	0.686	11.5	99	-0.01
84	c	Di-n-octyl phthalate	1.213	1.154	4.9	104	-0.01
85		Indeno(1,2,3-cd)pyrene	0.859	1.197	-39.3#	151#	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	134	-0.01
87		Benzo(b)fluoranthene	1.177	1.136	3.5	127	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.227	0.948	22.7	113	-0.01
89 C	Benzo(a)pyrene	1.111	1.033	7.0	128	-0.01
90	Dibenzo(a,h)anthracene	0.909	1.076	-18.4	149	-0.01
91	Benzo(g,h,i)perylene	0.911	1.101	-20.9	153#	-0.01

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