Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF081815\
Data File : BF081043.D
Acq On : 18 Aug 2015 15:06
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 19 02:01:01 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 17 17:57:25 2015
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	91	-0.02
2		1,4-Dioxane	0.627	0.140	77.7#	20#	-0.06
3		Pyridine	1.769	1.895	-7.1	86	-0.09
4		n-Nitrosodimethylamine	0.985	0.974	1.1	88	-0.05
5	S	2-Fluorophenol	1.330	1.267	4.7	89	-0.02
6		Aniline	2.523	2.668	-5.7	98	-0.01
7	S	Phenol-d6	1.735	1.771	-2.1	87	-0.01
8		2-Chlorophenol	1.455	1.559	-7.1	90	-0.01
9		Benzaldehyde	1.118	1.200	-7.3	91	-0.01
10	C	Phenol	2.049	2.020	1.4	80	0.00
11		bis(2-Chloroethyl)ether	1.572	1.716	-9.2	97	-0.01
12		1,3-Dichlorobenzene	1.564	1.692	-8.2	97	-0.01
13	C	1,4-Dichlorobenzene	1.549	1.707	-10.2	101	-0.01
14		1,2-Dichlorobenzene	1.449	1.429	1.4	82	-0.02
15		Benzyl Alcohol	1.404	1.675	-19.3	116	-0.01
16		2,2'-oxybis(1-Chloropropane	3.089	3.196	-3.5	90	-0.01
17		2-Methylphenol	1.249	1.248	0.1	87	-0.01
18		Hexachloroethane	0.626	0.661	-5.6	92	-0.02
19	P	n-Nitroso-di-n-propylamine	1.202	1.253	-4.2	88	-0.01
20		3+4-Methylphenols	1.585	1.569	1.0	90	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	100	-0.01
22		Acetophenone	0.525	0.486	7.4	87	-0.01
23	S	Nitrobenzene-d5	0.438	0.432	1.4	99	-0.01
24		Nitrobenzene	0.458	0.460	-0.4	95	-0.01
25		Isophorone	0.816	0.848	-3.9	103	0.00
26	C	2-Nitrophenol	0.190	0.211	-11.1	115	-0.01
27		2,4-Dimethylphenol	0.337	0.310	8.0	83	-0.01
28		bis(2-Chloroethoxy)methane	0.482	0.502	-4.1	104	-0.01
29	C	2,4-Dichlorophenol	0.284	0.284	0.0	94	-0.01
30		1,2,4-Trichlorobenzene	0.315	0.304	3.5	91	-0.02
31		Naphthalene	1.115	1.169	-4.8	103	-0.01
32		Benzoic acid	0.189	0.225	-19.0	138	0.05
33		4-Chloroaniline	0.470	0.529	-12.6	120	-0.01
34	C	Hexachlorobutadiene	0.178	0.194	-9.0	107	-0.01
35		Caprolactam	0.099	0.100	-1.0	96	0.02
36	C	4-Chloro-3-methylphenol	0.338	0.371	-9.8	107	0.00
37		2-Methylnaphthalene	0.704	0.669	5.0	91	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	108	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.645	0.662	-2.6	110	-0.01
40	P	Hexachlorocyclopentadiene	0.334	0.355	-6.3	107	-0.01
41	S	2,4,6-Tribromophenol	0.173	0.178	-2.9	116	-0.02
42	C	2,4,6-Trichlorophenol	0.456	0.505	-10.7	120	-0.01
43		2,4,5-Trichlorophenol	0.456	0.429	5.9	91	-0.01
44	S	2-Fluorobiphenyl	1.526	1.402	8.1	110	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.762	1.894	-7.5	105	-0.01
46		2-Chloronaphthalene	1.415	1.415	0.0	111	-0.01
47		2-Nitroaniline	0.579	0.537	7.3	98	-0.01
48		Acenaphthylene	2.532	2.458	2.9	111	-0.01
49		Dimethylphthalate	1.647	1.571	4.6	95	-0.01
50		2,6-Dinitrotoluene	0.354	0.372	-5.1	104	-0.01
51	C	Acenaphthene	1.516	1.583	-4.4	115	-0.01
52		3-Nitroaniline	0.443	0.426	3.8	106	-0.01
53	P	2,4-Dinitrophenol	0.167	0.208	-24.6	128	-0.01
54		Dibenzofuran	2.031	2.025	0.3	114	-0.01
55	P	4-Nitrophenol	0.311	0.328	-5.5	119	-0.01
56		2,4-Dinitrotoluene	0.469	0.506	-7.9	113	-0.01
57		Fluorene	1.562	1.415	9.4	103	-0.02
58		2,3,4,6-Tetrachlorophenol	0.353	0.377	-6.8	105	-0.07
59		Diethylphthalate	1.743	1.900	-9.0	114	-0.01
60		4-Chlorophenyl-phenylether	0.661	0.667	-0.9	107	-0.02
61		4-Nitroaniline	0.452	0.489	-8.2	119	0.00
62		Azobenzene	2.009	2.122	-5.6	106	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	112	-0.02
64		4,6-Dinitro-2-methylphenol	0.119	0.132	-10.9	112	-0.01
65	c	n-Nitrosodiphenylamine	0.714	0.772	-8.1	117	-0.01
66		4-Bromophenyl-phenylether	0.196	0.215	-9.7	126	-0.01
67		Hexachlorobenzene	0.199	0.212	-6.5	111	-0.01
68		Atrazine	0.199	0.213	-7.0	120	-0.01
69	C	Pentachlorophenol	0.119	0.142	-19.3	131	-0.01
70		Phenanthrene	1.185	1.212	-2.3	117	-0.01
71		Anthracene	1.153	1.163	-0.9	110	-0.01
72		Carbazole	1.110	1.234	-11.2	111	-0.01
73		Di-n-butylphthalate	1.344	1.495	-11.2	120	-0.02
74	C	Fluoranthene	1.279	1.387	-8.4	117	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	129	-0.01
76		Benzidine	0.756	0.847	-12.0	117	-0.02
77		Pyrene	1.626	1.416	12.9	119	-0.02
78	S	Terphenyl-d14	0.933	0.902	3.3	119	-0.01
79		Butylbenzylphthalate	0.699	0.835	-19.5	139	-0.02
80		Benzo(a)anthracene	1.341	1.366	-1.9	124	-0.01
81		3,3'-Dichlorobenzidine	0.434	0.479	-10.4	142	-0.02
82		Chrysene	1.209	1.219	-0.8	120	-0.01
83		Bis(2-ethylhexyl)phthalate	0.895	1.101	-23.0	151#	-0.02
84	c	Di-n-octyl phthalate	1.361	1.965	-44.4#	172#	-0.01
85		Indeno(1,2,3-cd)pyrene	0.849	1.128	-32.9#	163#	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	156#	-0.02
87		Benzo(b)fluoranthene	1.415	1.502	-6.1	151#	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.318	1.069	18.9	137	-0.02
89 C	Benzo(a)pyrene	1.233	1.286	-4.3	161#	-0.01
90	Dibenzo(a,h)anthracene	0.960	1.032	-7.5	163#	-0.02
91	Benzo(g,h,i)perylene	0.974	0.998	-2.5	157#	-0.01

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