Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF020315\
Data File : BF077161.D
Acq On : 3 Feb 2015 12:14
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 04 12:40:40 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 03 23:50:37 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	111	0.00
2		1,4-Dioxane	0.521	0.463	11.1	104	0.00
3		Pyridine	1.348	1.365	-1.3	94	-0.02
4		n-Nitrosodimethylamine	0.668	0.683	-2.2	113	0.00
5	S	2-Fluorophenol	1.216	1.156	4.9	101	-0.02
6		Aniline	2.124	2.062	2.9	101	0.00
7	S	Phenol-d6	1.535	1.474	4.0	102	0.00
8		2-Chlorophenol	1.402	1.381	1.5	103	0.00
9		Benzaldehyde	0.894	0.838	6.3	119	0.00
10	C	Phenol	1.629	1.517	6.9	95	-0.02
11		bis(2-Chloroethyl)ether	1.275	1.265	0.8	107	0.00
12		1,3-Dichlorobenzene	1.595	1.609	-0.9	109	0.00
13	C	1,4-Dichlorobenzene	1.692	1.657	2.1	106	0.00
14		1,2-Dichlorobenzene	1.559	1.572	-0.8	107	0.00
15		Benzyl Alcohol	1.242	1.267	-2.0	105	0.00
16		2,2'-oxybis(1-Chloropropane	1.714	1.769	-3.2	111	0.00
17		2-Methylphenol	1.146	1.101	3.9	100	-0.02
18		Hexachloroethane	0.588	0.612	-4.1	110	0.00
19	P	n-Nitroso-di-n-propylamine	1.074	1.074	0.0	104	0.00
20		3+4-Methylphenols	1.517	1.374	9.4	96	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	113	0.00
22		Acetophenone	0.490	0.494	-0.8	107	0.00
23	S	Nitrobenzene-d5	0.361	0.356	1.4	104	0.00
24		Nitrobenzene	0.368	0.354	3.8	100	0.00
25		Isophorone	0.657	0.648	1.4	103	0.00
26	C	2-Nitrophenol	0.172	0.158	8.1	91	0.00
27		2,4-Dimethylphenol	0.284	0.282	0.7	102	0.00
28		bis(2-Chloroethoxy)methane	0.374	0.381	-1.9	104	0.00
29	C	2,4-Dichlorophenol	0.265	0.245	7.5	95	0.00
30		1,2,4-Trichlorobenzene	0.294	0.294	0.0	102	0.00
31		Naphthalene	1.030	1.006	2.3	102	0.00
32		Benzoic acid	0.158	0.129	18.4	85	0.00
33		4-Chloroaniline	0.416	0.400	3.8	102	0.00
34	C	Hexachlorobutadiene	0.175	0.179	-2.3	106	0.00
35		Caprolactam	0.078	0.072	7.7	93	0.00
36	C	4-Chloro-3-methylphenol	0.303	0.296	2.3	102	0.00
37		2-Methylnaphthalene	0.703	0.700	0.4	106	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	109	0.00
39		1,2,4,5-Tetrachlorobenzene	0.494	0.486	1.6	105	0.00
40	P	Hexachlorocyclopentadiene	0.202	0.244	-20.8	107	0.00
41	S	2,4,6-Tribromophenol	0.162	0.159	1.9	99	0.00
42	C	2,4,6-Trichlorophenol	0.360	0.340	5.6	96	0.00
43		2,4,5-Trichlorophenol	0.367	0.316	13.9	87	0.00
44	S	2-Fluorobiphenyl	1.314	1.344	-2.3	107	0.00

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Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 04 12:40:40 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 03 23:50:37 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)
45		1,1'-Biphenyl	1.483	1.522	-2.6	105	0.00
46		2-Chloronaphthalene	1.220	1.225	-0.4	106	0.00
47		2-Nitroaniline	0.370	0.360	2.7	98	0.00
48		Acenaphthylene	2.058	2.052	0.3	104	0.00
49		Dimethylphthalate	1.418	1.451	-2.3	109	0.00
50		2,6-Dinitrotoluene	0.283	0.299	-5.7	104	0.00
51	C	Acenaphthene	1.168	1.155	1.1	104	0.00
52		3-Nitroaniline	0.339	0.326	3.8	95	0.00
53	P	2,4-Dinitrophenol	0.095	0.094	1.1	96	0.00
54		Dibenzofuran	1.701	1.686	0.9	105	0.00
55	P	4-Nitrophenol	0.222	0.177	20.3	81	0.00
56		2,4-Dinitrotoluene	0.384	0.408	-6.2	104	0.00
57		Fluorene	1.441	1.435	0.4	105	0.00
58		2,3,4,6-Tetrachlorophenol	0.268	0.234	12.7	88	0.00
59		Diethylphthalate	1.433	1.517	-5.9	111	0.00
60		4-Chlorophenyl-phenylether	0.590	0.589	0.2	106	0.00
61		4-Nitroaniline	0.328	0.325	0.9	98	0.00
62		Azobenzene	1.494	1.552	-3.9	109	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	112	0.00
64		4,6-Dinitro-2-methylphenol	0.104	0.112	-7.7	95	0.00
65	c	n-Nitrosodiphenylamine	0.713	0.737	-3.4	107	0.00
66		4-Bromophenyl-phenylether	0.203	0.208	-2.5	104	0.00
67		Hexachlorobenzene	0.214	0.222	-3.7	111	0.00
68		Atrazine	0.216	0.223	-3.2	101	0.00
69	C	Pentachlorophenol	0.082	0.069	15.9	85	0.00
70		Phenanthrene	1.247	1.241	0.5	106	0.00
71		Anthracene	1.216	1.240	-2.0	109	0.00
72		Carbazole	1.180	1.213	-2.8	108	0.00
73		Di-n-butylphthalate	1.365	1.515	-11.0	115	0.00
74	C	Fluoranthene	1.278	1.345	-5.2	109	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	118	0.00
76		Benzidine	0.640	0.683	-6.7	126	0.00
77		Pyrene	1.371	1.432	-4.4	112	0.00
78	S	Terphenyl-d14	0.869	0.909	-4.6	114	0.00
79		Butylbenzylphthalate	0.621	0.711	-14.5	119	0.00
80		Benzo(a)anthracene	1.255	1.280	-2.0	111	0.00
81		3,3'-Dichlorobenzidine	0.399	0.412	-3.3	113	0.00
82		Chrysene	1.144	1.159	-1.3	110	0.00
83		Bis(2-ethylhexyl)phthalate	0.859	1.014	-18.0	129	0.00
84	c	Di-n-octyl phthalate	1.491	1.773	-18.9	128	0.00
85		Indeno(1,2,3-cd)pyrene	1.213	1.241	-2.3	110	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	116	-0.02
87		Benzo(b)fluoranthene	1.347	1.435	-6.5	129	-0.02

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 03 23:50:37 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)
88	Benzo(k)fluoranthene	1.177	1.111	5.6	96	0.00
89 C	Benzo(a)pyrene	1.188	1.195	-0.6	110	0.00
90	Dibenzo(a,h)anthracene	1.090	1.136	-4.2	114	-0.03
91	Benzo(g,h,i)perylene	1.084	1.120	-3.3	113	-0.02

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