Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020316\
Data File : BF084802.D
Acq On : 3 Feb 2016 23:27
Operator : SJ/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 04 04:01:32 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 03 14:16:01 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	93	0.00
2		1,4-Dioxane	0.562	0.566	-0.7	96	0.00
3		Pyridine	1.465	1.474	-0.6	99	0.00
4		n-Nitrosodimethylamine	0.758	0.769	-1.5	94	0.00
5	S	2-Fluorophenol	1.284	1.381	-7.6	107	0.00
6		Aniline	1.948	2.011	-3.2	98	0.00
7	S	Phenol-d6	1.592	1.551	2.6	98	0.00
8		2-Chlorophenol	1.453	1.555	-7.0	99	0.00
9		Benzaldehyde	0.940	0.942	-0.2	92	-0.01
10	C	Phenol	1.783	1.671	6.3	100	0.00
11		bis(2-Chloroethyl)ether	1.391	1.432	-2.9	105	0.00
12		1,3-Dichlorobenzene	1.536	1.555	-1.2	100	0.00
13	C	1,4-Dichlorobenzene	1.535	1.644	-7.1	105	0.00
14		1,2-Dichlorobenzene	1.430	1.361	4.8	87	0.00
15		Benzyl Alcohol	1.099	1.101	-0.2	97	0.00
16		2,2'-oxybis(1-Chloropropane	2.339	2.304	1.5	98	0.00
17		2-Methylphenol	1.149	1.218	-6.0	95	0.00
18		Hexachloroethane	0.591	0.615	-4.1	96	0.00
19	P	n-Nitroso-di-n-propylamine	0.998	0.958	4.0	96	0.00
20		3+4-Methylphenols	1.427	1.409	1.3	94	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	94	0.00
22		Acetophenone	0.527	0.518	1.7	94	0.00
23	S	Nitrobenzene-d5	0.355	0.387	-9.0	101	0.00
24		Nitrobenzene	0.380	0.350	7.9	96	0.00
25		Isophorone	0.690	0.699	-1.3	94	0.00
26	C	2-Nitrophenol	0.188	0.199	-5.9	101	0.00
27		2,4-Dimethylphenol	0.326	0.305	6.4	94	-0.01
28		bis(2-Chloroethoxy)methane	0.410	0.396	3.4	94	0.00
29	C	2,4-Dichlorophenol	0.279	0.286	-2.5	93	0.00
30		1,2,4-Trichlorobenzene	0.315	0.309	1.9	95	0.00
31		Naphthalene	1.003	0.963	4.0	96	0.00
32		Benzoic acid	0.209	0.237	-13.4	106	0.00
33		4-Chloroaniline	0.415	0.399	3.9	99	0.00
34	C	Hexachlorobutadiene	0.182	0.193	-6.0	98	0.00
35		Caprolactam	0.086	0.093	-8.1	90	0.00
36	C	4-Chloro-3-methylphenol	0.318	0.300	5.7	95	0.00
37		2-Methylnaphthalene	0.628	0.662	-5.4	95	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	87	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.635	0.611	3.8	92	0.00
40	P	Hexachlorocyclopentadiene	0.223	0.273	-22.4	106	0.00
41	S	2,4,6-Tribromophenol	0.209	0.217	-3.8	87	0.00
42	C	2,4,6-Trichlorophenol	0.409	0.413	-1.0	87	0.00
43		2,4,5-Trichlorophenol	0.416	0.436	-4.8	95	0.00
44	S	2-Fluorobiphenyl	1.321	1.366	-3.4	102	0.00

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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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	1.786	1.839	-3.0	90	0.00
46		2-Chloronaphthalene	1.316	1.267	3.7	87	0.00
47		2-Nitroaniline	0.417	0.386	7.4	92	0.00
48		Acenaphthylene	1.996	1.935	3.1	85	-0.01
49		Dimethylphthalate	1.523	1.572	-3.2	89	0.00
50		2,6-Dinitrotoluene	0.333	0.386	-15.9	96	0.00
51	C	Acenaphthene	1.299	1.268	2.4	86	-0.01
52		3-Nitroaniline	0.374	0.345	7.8	84	-0.01
53	P	2,4-Dinitrophenol	0.133	0.154	-15.8	95	0.00
54		Dibenzofuran	1.587	1.586	0.1	87	-0.01
55	P	4-Nitrophenol	0.275	0.278	-1.1	81	0.00
56		2,4-Dinitrotoluene	0.424	0.473	-11.6	96	0.00
57		Fluorene	1.326	1.242	6.3	80	0.00
58		2,3,4,6-Tetrachlorophenol	0.317	0.330	-4.1	102	0.00
59		Diethylphthalate	1.487	1.482	0.3	90	-0.01
60		4-Chlorophenyl-phenylether	0.621	0.665	-7.1	100	0.00
61		4-Nitroaniline	0.364	0.376	-3.3	93	0.00
62		Azobenzene	1.430	1.564	-9.4	97	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	93	0.00
64		4,6-Dinitro-2-methylphenol	0.123	0.133	-8.1	95	0.00
65	c	n-Nitrosodiphenylamine	0.635	0.667	-5.0	91	0.00
66		4-Bromophenyl-phenylether	0.221	0.229	-3.6	90	0.00
67		Hexachlorobenzene	0.241	0.247	-2.5	100	0.00
68		Atrazine	0.201	0.188	6.5	93	0.00
69	C	Pentachlorophenol	0.113	0.117	-3.5	92	0.00
70		Phenanthrene	1.109	1.073	3.2	97	0.00
71		Anthracene	1.118	1.127	-0.8	88	0.00
72		Carbazole	0.975	0.952	2.4	84	-0.01
73		Di-n-butylphthalate	1.241	1.174	5.4	92	0.00
74	C	Fluoranthene	1.123	1.103	1.8	88	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	89	0.00
76		Benzidine	0.597	0.694	-16.2	87	0.00
77		Pyrene	1.531	1.471	3.9	82	0.00
78	S	Terphenyl-d14	0.883	0.970	-9.9	103	0.00
79		Butylbenzylphthalate	0.695	0.718	-3.3	87	0.00
80		Benzo(a)anthracene	1.241	1.239	0.2	89	0.00
81		3,3'-Dichlorobenzidine	0.440	0.431	2.0	79	-0.01
82		Chrysene	1.204	1.111	7.7	80	0.00
83		Bis(2-ethylhexyl)phthalate	0.864	0.932	-7.9	94	0.00
84	c	Di-n-octyl phthalate	1.426	1.415	0.8	86	0.00
85		Indeno(1,2,3-cd)pyrene	0.947	0.949	-0.2	91	0.00
86	I	Perylene-d12	1.000	1.000	0.0	84	0.00
87		Benzo(b)fluoranthene	1.344	1.237	8.0	77	0.00

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88	Benzo(k)fluoranthene	1.111	1.159	-4.3	91	0.00
89 C	Benzo(a)pyrene	1.145	1.105	3.5	81	0.00
90	Dibenzo(a,h)anthracene	0.964	1.012	-5.0	90	0.00
91	Benzo(g,h,i)perylene	0.971	1.035	-6.6	91	0.00

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