Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF111414\
Data File : BF075572.D
Acq On : 16 Nov 2014 7:57
Operator : TP / JJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 17 04:22:14 2014
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF111214.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 17 04:16:34 2014
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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	72	-0.02
2		1,4-Dioxane	40.000	42.541	-6.4	77	-0.07
3		Pyridine	40.000	38.322	4.2	69	-0.16
4		n-Nitrosodimethylamine	40.000	41.307	-3.3	78	-0.10
5	S	2-Fluorophenol	80.000	82.683	-3.4	76	-0.01
6		Aniline	40.000	41.894	-4.7	77	-0.01
7	S	Phenol-d6	80.000	84.624	-5.8	76	0.00
8		2-Chlorophenol	40.000	42.250	-5.6	77	-0.01
9		Benzaldehyde	40.000	40.577	-1.4	74	-0.01
10	C	Phenol	40.000	43.114	-7.8	76	0.00
11		bis(2-Chloroethyl)ether	40.000	39.482	1.3	72	-0.01
12		1,3-Dichlorobenzene	40.000	40.699	-1.7	75	-0.02
13	C	1,4-Dichlorobenzene	40.000	42.075	-5.2	76	-0.01
14		1,2-Dichlorobenzene	40.000	42.146	-5.4	77	-0.01
15		Benzyl Alcohol	40.000	42.148	-5.4	73	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.781	0.5	73	-0.01
17		2-Methylphenol	40.000	40.487	-1.2	72	-0.01
18		Hexachloroethane	40.000	41.068	-2.7	75	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	41.857	-4.6	76	-0.01
20		3+4-Methylphenols	40.000	41.220	-3.0	75	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	70	-0.02
22		Acetophenone	40.000	41.444	-3.6	74	-0.01
23	S	Nitrobenzene-d5	80.000	88.192	-10.2	76	-0.01
24		Nitrobenzene	40.000	42.618	-6.5	75	-0.01
25		Isophorone	40.000	40.843	-2.1	73	-0.01
26	C	2-Nitrophenol	40.000	45.069	-12.7	80	-0.01
27		2,4-Dimethylphenol	40.000	39.756	0.6	71	-0.01
28		bis(2-Chloroethoxy)methane	40.000	39.420	1.4	69	-0.01
29	C	2,4-Dichlorophenol	40.000	42.426	-6.1	75	-0.01
30		1,2,4-Trichlorobenzene	40.000	39.022	2.4	70	-0.02
31		Naphthalene	40.000	42.249	-5.6	78	-0.01
32		Benzoic acid	40.000	44.062	-10.2	71	0.06
33		4-Chloroaniline	40.000	41.002	-2.5	72	-0.01
34	C	Hexachlorobutadiene	40.000	40.254	-0.6	71	-0.02
35		Caprolactam	40.000	42.526	-6.3	77	0.02
36	C	4-Chloro-3-methylphenol	40.000	42.805	-7.0	76	0.00
37		2-Methylnaphthalene	40.000	40.177	-0.4	73	-0.01
38	I	Acenaphthene-d10	20.000	20.000	0.0	69	-0.02
39		1,2,4,5-Tetrachlorobenzene	40.000	41.263	-3.2	70	-0.01
40	P	Hexachlorocyclopentadiene	40.000	38.955	2.6	67	-0.01
41	S	2,4,6-Tribromophenol	80.000	79.896	0.1	67	-0.01
42	C	2,4,6-Trichlorophenol	40.000	41.444	-3.6	70	-0.01
43		2,4,5-Trichlorophenol	40.000	42.416	-6.0	73	0.00
44	S	2-Fluorobiphenyl	80.000	79.518	0.6	69	-0.01

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Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	42.253	-5.6	74	-0.01
46		2-Chloronaphthalene	40.000	41.627	-4.1	71	-0.01
47		2-Nitroaniline	40.000	45.499	-13.7	78	-0.01
48		Acenaphthylene	40.000	41.928	-4.8	72	-0.01
49		Dimethylphthalate	40.000	39.072	2.3	66	-0.02
50		2,6-Dinitrotoluene	40.000	43.325	-8.3	75	-0.01
51	C	Acenaphthene	40.000	39.880	0.3	69	-0.01
52		3-Nitroaniline	40.000	43.554	-8.9	73	-0.01
53	P	2,4-Dinitrophenol	40.000	43.712	-9.3	75	0.00
54		Dibenzofuran	40.000	41.182	-3.0	72	-0.01
55	P	4-Nitrophenol	40.000	46.972	-17.4	79	0.01
56		2,4-Dinitrotoluene	40.000	45.713	-14.3	71	-0.01
57		Fluorene	40.000	40.914	-2.3	70	-0.02
58		2,3,4,6-Tetrachlorophenol	40.000	40.689	-1.7	69	-0.01
59		Diethylphthalate	40.000	39.005	2.5	65	-0.02
60		4-Chlorophenyl-phenylether	40.000	39.058	2.4	69	-0.02
61		4-Nitroaniline	40.000	42.346	-5.9	72	-0.01
62		Azobenzene	40.000	39.688	0.8	69	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	70	-0.02
64		4,6-Dinitro-2-methylphenol	40.000	45.630	-14.1	76	-0.01
65	c	n-Nitrosodiphenylamine	40.000	38.954	2.6	66	-0.02
66		4-Bromophenyl-phenylether	40.000	41.033	-2.6	71	-0.02
67		Hexachlorobenzene	40.000	40.040	-0.1	69	-0.01
68		Atrazine	40.000	38.294	4.3	69	-0.01
69	C	Pentachlorophenol	40.000	33.200	17.0	56	-0.02
70		Phenanthrene	40.000	41.533	-3.8	72	-0.01
71		Anthracene	40.000	40.883	-2.2	72	-0.01
72		Carbazole	40.000	40.044	-0.1	73	-0.02
73		Di-n-butylphthalate	40.000	38.656	3.4	68	-0.01
74	C	Fluoranthene	40.000	41.836	-4.6	74	-0.01
75	I	Chrysene-d12	20.000	20.000	0.0	71	-0.06
76		Benzidine	40.000	4.922	87.7#	8	-0.01
77		Pyrene	40.000	41.143	-2.9	70	-0.02
78	S	Terphenyl-d14	80.000	78.313	2.1	69	-0.02
79		Butylbenzylphthalate	40.000	38.761	3.1	66	-0.05
80		Benzo(a)anthracene	40.000	40.438	-1.1	71	-0.06
81		3,3'-Dichlorobenzidine	40.000	37.201	7.0	63	-0.07
82		Chrysene	40.000	41.128	-2.8	73	-0.06
83		Bis(2-ethylhexyl)phthalate	40.000	36.905	7.7	64	-0.07
84	c	Di-n-octyl phthalate	40.000	36.008	10.0	65	-0.11
85		Indeno(1,2,3-cd)pyrene	40.000	42.885	-7.2	74	-0.17
86	I	Perylene-d12	20.000	20.000	0.0	75	-0.15
87		Benzo(b)fluoranthene	40.000	39.502	1.2	78	-0.14

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	39.204	2.0	69	-0.14
89 C	Benzo(a)pyrene	40.000	40.072	-0.2	74	-0.15
90	Dibenzo(a,h)anthracene	40.000	40.668	-1.7	75	-0.17
91	Benzo(g,h,i)perylene	40.000	41.218	-3.0	76	-0.16

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