Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071315\
Data File : BF080445.D
Acq On : 14 Jul 2015 00:25
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 14 08:31:19 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 14 00:52:37 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	66	-0.02
2		1,4-Dioxane	40.000	42.757	-6.9	68	-0.07
3		Pyridine	40.000	40.611	-1.5	67	-0.10
4		n-Nitrosodimethylamine	40.000	41.766	-4.4	67	-0.06
5	S	2-Fluorophenol	80.000	86.388	-8.0	68	-0.02
6		Aniline	40.000	43.529	-8.8	72	-0.01
7	S	Phenol-d6	80.000	77.667	2.9	63	-0.02
8		2-Chlorophenol	40.000	38.702	3.2	60	-0.01
9		Benzaldehyde	40.000	38.911	2.7	63	-0.01
10	C	Phenol	40.000	41.309	-3.3	68	-0.01
11		bis(2-Chloroethyl)ether	40.000	37.395	6.5	61	-0.01
12		1,3-Dichlorobenzene	40.000	42.937	-7.3	72	-0.01
13	C	1,4-Dichlorobenzene	40.000	41.548	-3.9	68	-0.01
14		1,2-Dichlorobenzene	40.000	39.847	0.4	66	-0.01
15		Benzyl Alcohol	40.000	35.924	10.2	58	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	46.647	-16.6	76	-0.01
17		2-Methylphenol	40.000	35.432	11.4	54	-0.01
18		Hexachloroethane	40.000	40.598	-1.5	64	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	37.784	5.5	63	-0.02
20		3+4-Methylphenols	40.000	41.814	-4.5	65	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	66	-0.01
22		Acetophenone	40.000	40.147	-0.4	61	-0.02
23	S	Nitrobenzene-d5	80.000	87.293	-9.1	71	-0.01
24		Nitrobenzene	40.000	38.451	3.9	57	-0.01
25		Isophorone	40.000	44.771	-11.9	72	-0.02
26	C	2-Nitrophenol	40.000	36.047	9.9	55	-0.01
27		2,4-Dimethylphenol	40.000	43.794	-9.5	70	-0.01
28		bis(2-Chloroethoxy)methane	40.000	44.167	-10.4	69	-0.02
29	C	2,4-Dichlorophenol	40.000	44.138	-10.3	68	-0.02
30		1,2,4-Trichlorobenzene	40.000	38.187	4.5	60	-0.01
31		Naphthalene	40.000	41.643	-4.1	69	-0.01
32		Benzoic acid	40.000	34.956	12.6	57	-0.03
33		4-Chloroaniline	40.000	44.315	-10.8	67	-0.01
34	C	Hexachlorobutadiene	40.000	40.315	-0.8	65	-0.01
35		Caprolactam	40.000	34.943	12.6	58	-0.04
36	C	4-Chloro-3-methylphenol	40.000	44.118	-10.3	68	-0.02
37		2-Methylnaphthalene	40.000	37.162	7.1	59	-0.02
38	I	Acenaphthene-d10	20.000	20.000	0.0	63	-0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	40.709	-1.8	62	-0.01
40	P	Hexachlorocyclopentadiene	40.000	38.262	4.3	58	-0.01
41	S	2,4,6-Tribromophenol	80.000	65.180	18.5	46	-0.02
42	C	2,4,6-Trichlorophenol	40.000	41.554	-3.9	60	-0.02
43		2,4,5-Trichlorophenol	40.000	44.253	-10.6	65	-0.01
44	S	2-Fluorobiphenyl	80.000	83.275	-4.1	65	-0.01

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Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 14 08:31:19 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 14 00:52:37 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	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	39.978	0.1	61	-0.01
46		2-Chloronaphthalene	40.000	38.736	3.2	58	-0.02
47		2-Nitroaniline	40.000	32.967	17.6	50	-0.01
48		Acenaphthylene	40.000	42.700	-6.8	65	-0.01
49		Dimethylphthalate	40.000	40.338	-0.8	62	-0.02
50		2,6-Dinitrotoluene	40.000	34.014	15.0	51	-0.01
51	C	Acenaphthene	40.000	40.683	-1.7	62	-0.01
52		3-Nitroaniline	40.000	35.288	11.8	52	-0.01
53	P	2,4-Dinitrophenol	40.000	38.582	3.5	60	-0.02
54		Dibenzofuran	40.000	42.117	-5.3	65	-0.01
55	P	4-Nitrophenol	40.000	35.820	10.4	56	-0.01
56		2,4-Dinitrotoluene	40.000	42.976	-7.4	65	-0.01
57		Fluorene	40.000	40.568	-1.4	64	-0.01
58		2,3,4,6-Tetrachlorophenol	40.000	37.941	5.1	54	-0.01
59		Diethylphthalate	40.000	37.274	6.8	59	-0.01
60		4-Chlorophenyl-phenylether	40.000	36.738	8.2	56	-0.01
61		4-Nitroaniline	40.000	35.569	11.1	55	-0.02
62		Azobenzene	40.000	37.012	7.5	56	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	51	-0.01
64		4,6-Dinitro-2-methylphenol	40.000	43.355	-8.4	55	-0.01
65	c	n-Nitrosodiphenylamine	40.000	48.186	-20.5#	59	-0.02
66		4-Bromophenyl-phenylether	40.000	43.902	-9.8	55	-0.01
67		Hexachlorobenzene	40.000	43.669	-9.2	55	-0.02
68		Atrazine	40.000	43.998	-10.0	49	-0.02
69	C	Pentachlorophenol	40.000	40.562	-1.4	48	-0.01
70		Phenanthrene	40.000	41.873	-4.7	54	-0.01
71		Anthracene	40.000	47.637	-19.1	59	-0.02
72		Carbazole	40.000	47.657	-19.1	59	-0.02
73		Di-n-butylphthalate	40.000	48.879	-22.2	55	-0.02
74	C	Fluoranthene	40.000	41.969	-4.9	53	-0.02
75	I	Chrysene-d12	20.000	20.000	0.0	55	-0.06
76		Benzidine	40.000	52.273	-30.7#	70	-0.03
77		Pyrene	40.000	41.624	-4.1	58	-0.02
78	S	Terphenyl-d14	80.000	81.482	-1.9	56	-0.03
79		Butylbenzylphthalate	40.000	37.326	6.7	48	-0.03
80		Benzo(a)anthracene	40.000	39.994	0.0	54	-0.06
81		3,3'-Dichlorobenzidine	40.000	42.472	-6.2	55	-0.05
82		Chrysene	40.000	40.456	-1.1	54	-0.06
83		Bis(2-ethylhexyl)phthalate	40.000	36.106	9.7	48	-0.06
84	c	Di-n-octyl phthalate	40.000	34.155	14.6	45	-0.07
85		Indeno(1,2,3-cd)pyrene	40.000	67.773	-69.4#	94	-0.11
86	I	Perylene-d12	20.000	20.000	0.0	66	-0.08
87		Benzo(b)fluoranthene	40.000	35.914	10.2	60	-0.08

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Instrument : BNA_F
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 14 00:52:37 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	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	35.893	10.3	60	-0.08
89 C	Benzo(a)pyrene	40.000	38.065	4.8	64	-0.09
90	Dibenzo(a,h)anthracene	40.000	53.167	-32.9#	91	-0.11
91	Benzo(g,h,i)perylene	40.000	59.535	-48.8#	101	-0.10

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