Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050715\
Data File : BF078991.D
Acq On : 7 May 2015 12:37
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 08 00:59:43 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 08 00:46:40 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	64	-0.03
2		1,4-Dioxane	40.000	44.144	-10.4	72	-0.05
3		Pyridine	40.000	44.829	-12.1	77	-0.05
4		n-Nitrosodimethylamine	40.000	42.481	-6.2	70	-0.06
5	S	2-Fluorophenol	80.000	84.684	-5.9	70	-0.03
6		Aniline	40.000	43.166	-7.9	71	-0.03
7	S	Phenol-d6	80.000	83.619	-4.5	72	-0.03
8		2-Chlorophenol	40.000	42.563	-6.4	74	-0.03
9		Benzaldehyde	40.000	41.469	-3.7	69	-0.03
10	C	Phenol	40.000	43.552	-8.9	72	-0.02
11		bis(2-Chloroethyl)ether	40.000	43.163	-7.9	76	-0.03
12		1,3-Dichlorobenzene	40.000	40.994	-2.5	71	-0.03
13	C	1,4-Dichlorobenzene	40.000	41.246	-3.1	70	-0.03
14		1,2-Dichlorobenzene	40.000	41.124	-2.8	69	-0.03
15		Benzyl Alcohol	40.000	44.022	-10.1	75	-0.03
16		2,2'-oxybis(1-Chloropropane	40.000	40.399	-1.0	70	-0.02
17		2-Methylphenol	40.000	40.837	-2.1	70	-0.03
18		Hexachloroethane	40.000	41.644	-4.1	69	-0.03
19	P	n-Nitroso-di-n-propylamine	40.000	41.604	-4.0	67	-0.03
20		3+4-Methylphenols	40.000	42.904	-7.3	71	-0.03
21	I	Naphthalene-d8	20.000	20.000	0.0	65	-0.03
22		Acetophenone	40.000	41.291	-3.2	72	-0.03
23	S	Nitrobenzene-d5	80.000	89.290	-11.6	76	-0.03
24		Nitrobenzene	40.000	42.654	-6.6	75	-0.03
25		Isophorone	40.000	42.519	-6.3	72	-0.03
26	C	2-Nitrophenol	40.000	46.899	-17.2	76	-0.03
27		2,4-Dimethylphenol	40.000	40.384	-1.0	67	-0.03
28		bis(2-Chloroethoxy)methane	40.000	40.672	-1.7	67	-0.03
29	C	2,4-Dichlorophenol	40.000	44.287	-10.7	75	-0.03
30		1,2,4-Trichlorobenzene	40.000	40.430	-1.1	69	-0.03
31		Naphthalene	40.000	41.484	-3.7	72	-0.03
32		Benzoic acid	40.000	39.638	0.9	65	-0.06
33		4-Chloroaniline	40.000	44.071	-10.2	77	-0.03
34	C	Hexachlorobutadiene	40.000	40.079	-0.2	70	-0.03
35		Caprolactam	40.000	44.079	-10.2	73	-0.06
36	C	4-Chloro-3-methylphenol	40.000	42.810	-7.0	68	-0.02
37		2-Methylnaphthalene	40.000	41.731	-4.3	71	-0.03
38	I	Acenaphthene-d10	20.000	20.000	0.0	66	-0.03
39		1,2,4,5-Tetrachlorobenzene	40.000	40.640	-1.6	72	-0.03
40	P	Hexachlorocyclopentadiene	40.000	35.357	11.6	60	-0.03
41	S	2,4,6-Tribromophenol	80.000	84.755	-5.9	70	-0.03
42	C	2,4,6-Trichlorophenol	40.000	41.866	-4.7	70	-0.03
43		2,4,5-Trichlorophenol	40.000	43.513	-8.8	74	-0.03
44	S	2-Fluorobiphenyl	80.000	81.952	-2.4	70	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050715\
Data File : BF078991.D
Acq On : 7 May 2015 12:37
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 08 00:59:43 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 08 00:46:40 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	40.521	-1.3	72	-0.03
46		2-Chloronaphthalene	40.000	40.504	-1.3	72	-0.03
47		2-Nitroaniline	40.000	45.028	-12.6	75	-0.03
48		Acenaphthylene	40.000	41.844	-4.6	73	-0.05
49		Dimethylphthalate	40.000	39.478	1.3	71	-0.03
50		2,6-Dinitrotoluene	40.000	43.658	-9.1	74	-0.03
51	C	Acenaphthene	40.000	38.488	3.8	66	-0.05
52		3-Nitroaniline	40.000	44.088	-10.2	75	-0.03
53	P	2,4-Dinitrophenol	40.000	45.251	-13.1	78	-0.03
54		Dibenzofuran	40.000	40.261	-0.7	70	-0.03
55	P	4-Nitrophenol	40.000	45.285	-13.2	78	-0.02
56		2,4-Dinitrotoluene	40.000	41.494	-3.7	68	-0.03
57		Fluorene	40.000	40.407	-1.0	72	-0.05
58		2,3,4,6-Tetrachlorophenol	40.000	39.925	0.2	67	-0.03
59		Diethylphthalate	40.000	39.413	1.5	69	-0.05
60		4-Chlorophenyl-phenylether	40.000	38.960	2.6	67	-0.03
61		4-Nitroaniline	40.000	48.679	-21.7	81	-0.03
62		Azobenzene	40.000	39.350	1.6	67	-0.03
63	I	Phenanthrene-d10	20.000	20.000	0.0	67	-0.03
64		4,6-Dinitro-2-methylphenol	40.000	43.317	-8.3	74	-0.03
65	c	n-Nitrosodiphenylamine	40.000	41.659	-4.1	71	-0.03
66		4-Bromophenyl-phenylether	40.000	39.343	1.6	70	-0.03
67		Hexachlorobenzene	40.000	39.923	0.2	72	-0.03
68		Atrazine	40.000	42.637	-6.6	76	-0.03
69	C	Pentachlorophenol	40.000	34.060	14.8	58	-0.03
70		Phenanthrene	40.000	40.616	-1.5	71	-0.03
71		Anthracene	40.000	41.213	-3.0	72	-0.03
72		Carbazole	40.000	41.854	-4.6	73	-0.02
73		Di-n-butylphthalate	40.000	40.883	-2.2	68	-0.03
74	C	Fluoranthene	40.000	42.059	-5.1	73	-0.03
75	I	Chrysene-d12	20.000	20.000	0.0	66	-0.02
76		Benzidine	40.000	43.588	-9.0	70	-0.03
77		Pyrene	40.000	39.827	0.4	70	-0.03
78	S	Terphenyl-d14	80.000	78.584	1.8	69	-0.03
79		Butylbenzylphthalate	40.000	42.019	-5.0	68	-0.02
80		Benzo(a)anthracene	40.000	40.676	-1.7	71	-0.02
81		3,3'-Dichlorobenzidine	40.000	44.321	-10.8	74	-0.01
82		Chrysene	40.000	39.352	1.6	71	-0.02
83		Bis(2-ethylhexyl)phthalate	40.000	39.681	0.8	65	-0.01
84	c	Di-n-octyl phthalate	40.000	40.143	-0.4	70	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	43.135	-7.8	75	0.01
86	I	Perylene-d12	20.000	20.000	0.0	72	0.02
87		Benzo(b)fluoranthene	40.000	39.884	0.3	74	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050715\
Data File : BF078991.D
Acq On : 7 May 2015 12:37
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 08 00:59:43 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 08 00:46:40 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	40.853	-2.1	75	0.01
89 C	Benzo(a)pyrene	40.000	41.006	-2.5	76	0.02
90	Dibenzo(a,h)anthracene	40.000	40.625	-1.6	75	0.01
91	Benzo(g,h,i)perylene	40.000	41.088	-2.7	77	0.00

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