Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062515\
Data File : BF080159.D
Acq On : 25 Jun 2015 11:20
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 26 01:03:21 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 26 01:01:50 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	71	-0.03
2		1,4-Dioxane	40.000	40.341	-0.9	71	-0.06
3		Pyridine	40.000	40.862	-2.2	72	-0.03
4		n-Nitrosodimethylamine	40.000	39.228	1.9	65	-0.05
5	S	2-Fluorophenol	80.000	88.110	-10.1	76	-0.03
6		Aniline	40.000	46.096	-15.2	78	-0.02
7	S	Phenol-d6	80.000	87.059	-8.8	72	-0.02
8		2-Chlorophenol	40.000	42.022	-5.1	71	-0.03
9		Benzaldehyde	40.000	36.219	9.5	62	-0.02
10	C	Phenol	40.000	44.339	-10.8	75	-0.02
11		bis(2-Chloroethyl)ether	40.000	40.272	-0.7	69	-0.02
12		1,3-Dichlorobenzene	40.000	40.254	-0.6	69	-0.03
13	C	1,4-Dichlorobenzene	40.000	48.798	-22.0#	84	-0.02
14		1,2-Dichlorobenzene	40.000	41.594	-4.0	71	-0.02
15		Benzyl Alcohol	40.000	39.788	0.5	67	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	47.390	-18.5	85	-0.02
17		2-Methylphenol	40.000	43.910	-9.8	73	-0.02
18		Hexachloroethane	40.000	44.455	-11.1	75	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	44.236	-10.6	81	-0.02
20		3+4-Methylphenols	40.000	45.068	-12.7	76	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	72	-0.03
22		Acetophenone	40.000	40.548	-1.4	68	-0.03
23	S	Nitrobenzene-d5	80.000	89.426	-11.8	77	-0.02
24		Nitrobenzene	40.000	44.393	-11.0	77	-0.02
25		Isophorone	40.000	38.297	4.3	67	-0.02
26	C	2-Nitrophenol	40.000	51.638	-29.1#	90	-0.02
27		2,4-Dimethylphenol	40.000	45.000	-12.5	82	-0.02
28		bis(2-Chloroethoxy)methane	40.000	37.221	6.9	65	-0.02
29	C	2,4-Dichlorophenol	40.000	48.446	-21.1#	83	-0.02
30		1,2,4-Trichlorobenzene	40.000	46.638	-16.6	83	-0.02
31		Naphthalene	40.000	39.453	1.4	68	-0.02
32		Benzoic acid	40.000	41.195	-3.0	72	-0.02
33		4-Chloroaniline	40.000	37.703	5.7	69	-0.02
34	C	Hexachlorobutadiene	40.000	43.436	-8.6	80	-0.02
35		Caprolactam	40.000	43.246	-8.1	72	-0.03
36	C	4-Chloro-3-methylphenol	40.000	41.018	-2.5	70	-0.02
37		2-Methylnaphthalene	40.000	43.932	-9.8	76	-0.02
38	I	Acenaphthene-d10	20.000	20.000	0.0	73	-0.03
39		1,2,4,5-Tetrachlorobenzene	40.000	46.057	-15.1	82	-0.02
40	P	Hexachlorocyclopentadiene	40.000	39.770	0.6	73	-0.02
41	S	2,4,6-Tribromophenol	80.000	90.816	-13.5	80	-0.02
42	C	2,4,6-Trichlorophenol	40.000	47.134	-17.8	82	-0.02
43		2,4,5-Trichlorophenol	40.000	38.967	2.6	67	-0.03
44	S	2-Fluorobiphenyl	80.000	77.906	2.6	70	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062515\
Data File : BF080159.D
Acq On : 25 Jun 2015 11:20
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 26 01:03:21 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 26 01:01:50 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	40.134	-0.3	72	-0.02
46		2-Chloronaphthalene	40.000	39.372	1.6	69	-0.02
47		2-Nitroaniline	40.000	42.287	-5.7	70	-0.02
48		Acenaphthylene	40.000	42.821	-7.1	73	-0.02
49		Dimethylphthalate	40.000	43.259	-8.1	79	-0.02
50		2,6-Dinitrotoluene	40.000	41.157	-2.9	68	-0.02
51	C	Acenaphthene	40.000	40.330	-0.8	71	-0.03
52		3-Nitroaniline	40.000	43.756	-9.4	74	-0.02
53	P	2,4-Dinitrophenol	40.000	52.764	-31.9#	100	-0.02
54		Dibenzofuran	40.000	38.668	3.3	68	-0.03
55	P	4-Nitrophenol	40.000	44.714	-11.8	81	-0.02
56		2,4-Dinitrotoluene	40.000	46.344	-15.9	81	-0.02
57		Fluorene	40.000	43.436	-8.6	77	-0.02
58		2,3,4,6-Tetrachlorophenol	40.000	39.619	1.0	67	-0.02
59		Diethylphthalate	40.000	40.621	-1.6	69	-0.02
60		4-Chlorophenyl-phenylether	40.000	39.271	1.8	71	-0.03
61		4-Nitroaniline	40.000	41.799	-4.5	71	-0.02
62		Azobenzene	40.000	38.337	4.2	65	-0.03
63	I	Phenanthrene-d10	20.000	20.000	0.0	82	-0.05
64		4,6-Dinitro-2-methylphenol	40.000	40.373	-0.9	76	-0.02
65	c	n-Nitrosodiphenylamine	40.000	36.338	9.2	70	-0.02
66		4-Bromophenyl-phenylether	40.000	40.584	-1.5	79	-0.02
67		Hexachlorobenzene	40.000	39.640	0.9	77	-0.03
68		Atrazine	40.000	42.083	-5.2	79	-0.03
69	C	Pentachlorophenol	40.000	36.888	7.8	75	-0.03
70		Phenanthrene	40.000	37.444	6.4	75	-0.05
71		Anthracene	40.000	39.431	1.4	80	-0.05
72		Carbazole	40.000	38.435	3.9	75	-0.05
73		Di-n-butylphthalate	40.000	36.516	8.7	76	-0.07
74	C	Fluoranthene	40.000	38.364	4.1	78	-0.10
75	I	Chrysene-d12	20.000	20.000	0.0	76	-0.17
76		Benzidine	40.000	41.679	-4.2	77	-0.10
77		Pyrene	40.000	39.381	1.5	74	-0.11
78	S	Terphenyl-d14	80.000	77.422	3.2	74	-0.13
79		Butylbenzylphthalate	40.000	41.709	-4.3	76	-0.15
80		Benzo(a)anthracene	40.000	38.789	3.0	74	-0.17
81		3,3'-Dichlorobenzidine	40.000	39.967	0.1	75	-0.17
82		Chrysene	40.000	39.662	0.8	75	-0.17
83		Bis(2-ethylhexyl)phthalate	40.000	38.716	3.2	71	-0.17
84	c	Di-n-octyl phthalate	40.000	38.738	3.2	72	-0.21
85		Indeno(1,2,3-cd)pyrene	40.000	38.783	3.0	74	-0.24
86	I	Perylene-d12	20.000	20.000	0.0	77	-0.22
87		Benzo(b)fluoranthene	40.000	41.045	-2.6	76	-0.22

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062515\
Data File : BF080159.D
Acq On : 25 Jun 2015 11:20
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 26 01:03:21 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 26 01:01:50 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	36.475	8.8	70	-0.22
89 C	Benzo(a)pyrene	40.000	40.630	-1.6	77	-0.22
90	Dibenzo(a,h)anthracene	40.000	38.824	2.9	74	-0.24
91	Benzo(g,h,i)perylene	40.000	38.792	3.0	73	-0.24

(#) = Out of Range SPCC's out = 0 CCC's out = 3