Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082915\
Data File : BF081269.D
Acq On : 29 Aug 2015 14:07
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 31 01:17:22 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 31 01:11:18 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	114	0.00
2		1,4-Dioxane	40.000	41.965	-4.9	116	0.00
3		Pyridine	40.000	37.884	5.3	96	0.00
4		n-Nitrosodimethylamine	40.000	42.330	-5.8	118	0.00
5	S	2-Fluorophenol	80.000	77.433	3.2	114	0.00
6		Aniline	40.000	40.338	-0.8	117	0.00
7	S	Phenol-d6	80.000	83.812	-4.8	112	0.00
8		2-Chlorophenol	40.000	39.790	0.5	104	0.00
9		Benzaldehyde	40.000	41.183	-3.0	109	0.00
10	C	Phenol	40.000	42.718	-6.8	109	0.00
11		bis(2-Chloroethyl)ether	40.000	41.555	-3.9	116	0.00
12		1,3-Dichlorobenzene	40.000	43.823	-9.6	124	0.00
13	C	1,4-Dichlorobenzene	40.000	43.704	-9.3	125	0.00
14		1,2-Dichlorobenzene	40.000	35.748	10.6	94	0.00
15		Benzyl Alcohol	40.000	43.335	-8.3	132	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.662	-4.2	114	0.00
17		2-Methylphenol	40.000	40.833	-2.1	112	0.00
18		Hexachloroethane	40.000	41.985	-5.0	115	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.796	-4.5	111	0.00
20		3+4-Methylphenols	40.000	39.174	2.1	112	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	123	0.00
22		Acetophenone	40.000	42.190	-5.5	121	0.00
23	S	Nitrobenzene-d5	80.000	75.241	5.9	116	0.00
24		Nitrobenzene	40.000	44.360	-10.9	129	0.00
25		Isophorone	40.000	42.141	-5.4	128	0.00
26	C	2-Nitrophenol	40.000	37.385	6.5	119	0.00
27		2,4-Dimethylphenol	40.000	36.328	9.2	101	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.506	-1.3	124	0.00
29	C	2,4-Dichlorophenol	40.000	43.086	-7.7	125	0.00
30		1,2,4-Trichlorobenzene	40.000	39.582	1.0	115	0.00
31		Naphthalene	40.000	37.165	7.1	112	0.00
32		Benzoic acid	40.000	48.333	-20.8	172	0.00
33		4-Chloroaniline	40.000	38.893	2.8	128	0.00
34	C	Hexachlorobutadiene	40.000	41.401	-3.5	125	0.00
35		Caprolactam	40.000	39.727	0.7	116	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.544	-6.4	128	0.00
37		2-Methylnaphthalene	40.000	38.388	4.0	113	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	137	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	35.162	12.1	119	0.00
40	P	Hexachlorocyclopentadiene	40.000	39.080	2.3	124	0.00
41	S	2,4,6-Tribromophenol	80.000	79.058	1.2	140	0.00
42	C	2,4,6-Trichlorophenol	40.000	36.408	9.0	124	0.00
43		2,4,5-Trichlorophenol	40.000	39.993	0.0	121	0.00
44	S	2-Fluorobiphenyl	80.000	69.991	12.5	132	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082915\
Data File : BF081269.D
Acq On : 29 Aug 2015 14:07
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 31 01:17:22 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 31 01:11:18 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	33.247	16.9	102	0.00
46		2-Chloronaphthalene	40.000	36.450	8.9	127	0.00
47		2-Nitroaniline	40.000	36.382	9.0	121	0.00
48		Acenaphthylene	40.000	36.713	8.2	133	0.00
49		Dimethylphthalate	40.000	35.127	12.2	111	0.00
50		2,6-Dinitrotoluene	40.000	38.434	3.9	120	0.00
51	C	Acenaphthene	40.000	34.856	12.9	121	0.00
52		3-Nitroaniline	40.000	32.424	18.9	113	0.00
53	P	2,4-Dinitrophenol	40.000	48.458	-21.1	168	0.00
54		Dibenzofuran	40.000	35.922	10.2	130	0.00
55	P	4-Nitrophenol	40.000	41.475	-3.7	148	0.00
56		2,4-Dinitrotoluene	40.000	37.184	7.0	123	0.00
57		Fluorene	40.000	32.819	18.0	118	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	37.892	5.3	117	0.00
59		Diethylphthalate	40.000	39.817	0.5	131	0.00
60		4-Chlorophenyl-phenylether	40.000	35.937	10.2	120	0.00
61		4-Nitroaniline	40.000	35.971	10.1	125	0.00
62		Azobenzene	40.000	33.990	15.0	107	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	130	0.00
64		4,6-Dinitro-2-methylphenol	40.000	44.171	-10.4	129	0.00
65	c	n-Nitrosodiphenylamine	40.000	35.245	11.9	110	0.00
66		4-Bromophenyl-phenylether	40.000	43.462	-8.7	144	0.00
67		Hexachlorobenzene	40.000	42.537	-6.3	128	0.00
68		Atrazine	40.000	36.178	9.6	118	0.00
69	C	Pentachlorophenol	40.000	41.681	-4.2	133	0.00
70		Phenanthrene	40.000	34.731	13.2	115	0.00
71		Anthracene	40.000	40.023	-0.1	126	0.00
72		Carbazole	40.000	39.792	0.5	115	0.00
73		Di-n-butylphthalate	40.000	40.460	-1.2	126	0.00
74	C	Fluoranthene	40.000	35.601	11.0	111	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	139	0.00
76		Benzidine	40.000	42.667	-6.7	120	0.00
77		Pyrene	40.000	37.644	5.9	138	0.00
78	S	Terphenyl-d14	80.000	78.019	2.5	129	0.00
79		Butylbenzylphthalate	40.000	42.847	-7.1	134	0.00
80		Benzo(a)anthracene	40.000	37.654	5.9	123	0.00
81		3,3'-Dichlorobenzidine	40.000	36.174	9.6	125	0.00
82		Chrysene	40.000	32.058	19.9	103	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	42.900	-7.2	142	0.00
84	c	Di-n-octyl phthalate	40.000	50.285	-25.7#	162	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	46.046	-15.1	152	0.00
86	I	Perylene-d12	20.000	20.000	0.0	152	0.00
87		Benzo(b)fluoranthene	40.000	41.561	-3.9	144	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082915\
Data File : BF081269.D
Acq On : 29 Aug 2015 14:07
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 31 01:17:22 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 31 01:11:18 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	31.682	20.8	131	0.00
89 C	Benzo(a)pyrene	40.000	38.371	4.1	144	0.00
90	Dibenzo(a,h)anthracene	40.000	41.571	-3.9	153	0.00
91	Benzo(g,h,i)perylene	40.000	39.256	1.9	147	0.00

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