Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
Data File : BF088538.D
Acq On : 30 Jun 2016 19:31
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 01 01:54:41 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 30 18:01:55 2016
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	96	0.00
2		1,4-Dioxane	40.000	39.092	2.3	97	0.00
3		Pyridine	40.000	38.804	3.0	97	0.00
4		n-Nitrosodimethylamine	40.000	38.776	3.1	97	0.00
5	S	2-Fluorophenol	80.000	82.899	-3.6	99	0.00
6		Aniline	40.000	40.383	-1.0	97	0.00
7	S	Phenol-d6	80.000	77.219	3.5	97	0.00
8		2-Chlorophenol	40.000	38.052	4.9	98	0.00
9		Benzaldehyde	40.000	36.822	7.9	95	0.00
10	C	Phenol	40.000	40.515	-1.3	97	0.00
11		bis(2-Chloroethyl)ether	40.000	38.543	3.6	99	0.00
12		1,3-Dichlorobenzene	40.000	36.821	7.9	97	0.00
13	C	1,4-Dichlorobenzene	40.000	39.747	0.6	102	0.00
14		1,2-Dichlorobenzene	40.000	37.566	6.1	101	0.00
15		Benzyl Alcohol	40.000	40.874	-2.2	96	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.125	-0.3	98	0.00
17		2-Methylphenol	40.000	40.652	-1.6	98	0.00
18		Hexachloroethane	40.000	40.689	-1.7	100	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	35.511	11.2	99	0.00
20		3+4-Methylphenols	40.000	37.341	6.6	96	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	98	0.00
22		Acetophenone	40.000	38.864	2.8	98	0.00
23	S	Nitrobenzene-d5	80.000	76.298	4.6	99	0.00
24		Nitrobenzene	40.000	38.614	3.5	100	0.00
25		Isophorone	40.000	37.543	6.1	97	0.00
26	C	2-Nitrophenol	40.000	37.175	7.1	99	0.00
27		2,4-Dimethylphenol	40.000	39.406	1.5	96	0.00
28		bis(2-Chloroethoxy)methane	40.000	36.518	8.7	98	0.00
29	C	2,4-Dichlorophenol	40.000	36.733	8.2	96	0.00
30		1,2,4-Trichlorobenzene	40.000	39.034	2.4	97	0.00
31		Naphthalene	40.000	39.682	0.8	97	0.00
32		Benzoic acid	40.000	38.920	2.7	97	0.00
33		4-Chloroaniline	40.000	38.221	4.4	95	0.00
34	C	Hexachlorobutadiene	40.000	40.501	-1.3	100	0.00
35		Caprolactam	40.000	40.494	-1.2	97	0.01
36	C	4-Chloro-3-methylphenol	40.000	38.322	4.2	96	0.00
37		2-Methylnaphthalene	40.000	34.785	13.0	97	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	98	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	41.559	-3.9	99	0.00
40	P	Hexachlorocyclopentadiene	40.000	39.243	1.9	97	0.00
41	S	2,4,6-Tribromophenol	80.000	84.050	-5.1	99	0.00
42	C	2,4,6-Trichlorophenol	40.000	36.499	8.8	98	0.00
43		2,4,5-Trichlorophenol	40.000	41.366	-3.4	98	0.00
44	S	2-Fluorobiphenyl	80.000	72.736	9.1	99	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
Data File : BF088538.D
Acq On : 30 Jun 2016 19:31
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 01 01:54:41 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 30 18:01:55 2016
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	40.896	-2.2	98	0.00
46		2-Chloronaphthalene	40.000	41.286	-3.2	99	0.00
47		2-Nitroaniline	40.000	43.714	-9.3	97	0.00
48		Acenaphthylene	40.000	39.368	1.6	96	0.00
49		Dimethylphthalate	40.000	36.887	7.8	98	0.00
50		2,6-Dinitrotoluene	40.000	37.088	7.3	98	0.00
51	C	Acenaphthene	40.000	37.688	5.8	97	0.00
52		3-Nitroaniline	40.000	44.674	-11.7	97	0.00
53	P	2,4-Dinitrophenol	40.000	39.136	2.2	98	0.00
54		Dibenzofuran	40.000	39.087	2.3	97	0.00
55	P	4-Nitrophenol	40.000	42.888	-7.2	96	0.00
56		2,4-Dinitrotoluene	40.000	39.255	1.9	101	0.00
57		Fluorene	40.000	41.402	-3.5	98	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	41.195	-3.0	98	0.00
59		Diethylphthalate	40.000	40.015	-0.0	98	-0.01
60		4-Chlorophenyl-phenylether	40.000	36.222	9.4	97	0.00
61		4-Nitroaniline	40.000	36.192	9.5	96	0.00
62		Azobenzene	40.000	38.062	4.8	98	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	99	0.00
64		4,6-Dinitro-2-methylphenol	40.000	42.445	-6.1	95	0.00
65	c	n-Nitrosodiphenylamine	40.000	41.199	-3.0	98	0.00
66		4-Bromophenyl-phenylether	40.000	38.232	4.4	97	0.00
67		Hexachlorobenzene	40.000	40.447	-1.1	100	0.00
68		Atrazine	40.000	40.804	-2.0	97	0.00
69	C	Pentachlorophenol	40.000	40.890	-2.2	95	0.00
70		Phenanthrene	40.000	37.224	6.9	97	0.00
71		Anthracene	40.000	39.548	1.1	97	0.00
72		Carbazole	40.000	34.812	13.0	96	0.00
73		Di-n-butylphthalate	40.000	36.193	9.5	94	0.00
74	C	Fluoranthene	40.000	39.981	0.0	94	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	94	0.00
76		Benzidine	40.000	39.362	1.6	90	0.00
77		Pyrene	40.000	42.691	-6.7	95	0.00
78	S	Terphenyl-d14	80.000	80.129	-0.2	96	0.00
79		Butylbenzylphthalate	40.000	38.428	3.9	91	0.00
80		Benzo(a)anthracene	40.000	38.153	4.6	89	0.00
81		3,3'-Dichlorobenzidine	40.000	36.156	9.6	88	0.00
82		Chrysene	40.000	36.221	9.4	86	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	40.627	-1.6	90	0.00
84	c	Di-n-octyl phthalate	40.000	38.749	3.1	87	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	37.657	5.9	92	0.00
86	I	Perylene-d12	20.000	20.000	0.0	86	0.00
87		Benzo(b)fluoranthene	40.000	39.364	1.6	89	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
Data File : BF088538.D
Acq On : 30 Jun 2016 19:31
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 01 01:54:41 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 30 18:01:55 2016
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	38.994	2.5	84	0.00
89 C	Benzo(a)pyrene	40.000	40.265	-0.7	86	-0.01
90	Dibenzo(a,h)anthracene	40.000	41.548	-3.9	91	0.00
91	Benzo(g,h,i)perylene	40.000	42.327	-5.8	93	0.00

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