Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061716\
Data File : BF088091.D
Acq On : 17 Jun 2016 13:07
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 17 17:11:17 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 17 16:09:13 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	98	0.00
2		1,4-Dioxane	40.000	41.816	-4.5	105	0.00
3		Pyridine	40.000	45.602	-14.0	109	0.00
4		n-Nitrosodimethylamine	40.000	37.807	5.5	93	0.00
5	S	2-Fluorophenol	80.000	71.498	10.6	89	0.00
6		Aniline	40.000	33.585	16.0	83	0.00
7	S	Phenol-d6	80.000	76.596	4.3	98	0.00
8		2-Chlorophenol	40.000	39.316	1.7	99	0.00
9		Benzaldehyde	40.000	31.163	22.1	86	0.00
10	C	Phenol	40.000	43.759	-9.4	112	0.00
11		bis(2-Chloroethyl)ether	40.000	41.872	-4.7	110	0.00
12		1,3-Dichlorobenzene	40.000	38.774	3.1	95	0.00
13	C	1,4-Dichlorobenzene	40.000	39.843	0.4	102	0.00
14		1,2-Dichlorobenzene	40.000	35.972	10.1	87	0.00
15		Benzyl Alcohol	40.000	32.708	18.2	77	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.276	4.3	93	0.00
17		2-Methylphenol	40.000	37.048	7.4	88	0.00
18		Hexachloroethane	40.000	40.032	-0.1	97	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	33.772	15.6	90	0.00
20		3+4-Methylphenols	40.000	34.719	13.2	92	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	104	0.00
22		Acetophenone	40.000	36.623	8.4	97	0.00
23	S	Nitrobenzene-d5	80.000	76.026	5.0	97	0.00
24		Nitrobenzene	40.000	34.389	14.0	97	0.00
25		Isophorone	40.000	38.055	4.9	96	0.00
26	C	2-Nitrophenol	40.000	43.265	-8.2	98	0.00
27		2,4-Dimethylphenol	40.000	35.580	11.1	89	0.00
28		bis(2-Chloroethoxy)methane	40.000	35.774	10.6	90	0.00
29	C	2,4-Dichlorophenol	40.000	36.709	8.2	93	0.00
30		1,2,4-Trichlorobenzene	40.000	35.148	12.1	94	0.00
31		Naphthalene	40.000	36.014	10.0	98	0.00
32		Benzoic acid	40.000	30.651	23.4	68	0.00
33		4-Chloroaniline	40.000	38.393	4.0	93	0.00
34	C	Hexachlorobutadiene	40.000	34.497	13.8	86	0.00
35		Caprolactam	40.000	36.864	7.8	87	0.00
36	C	4-Chloro-3-methylphenol	40.000	37.364	6.6	99	0.00
37		2-Methylnaphthalene	40.000	33.750	15.6	83	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	91	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	40.839	-2.1	98	0.00
40	P	Hexachlorocyclopentadiene	40.000	24.659	38.4#	56	0.00
41	S	2,4,6-Tribromophenol	80.000	58.375	27.0#	65	0.00
42	C	2,4,6-Trichlorophenol	40.000	36.648	8.4	81	0.00
43		2,4,5-Trichlorophenol	40.000	38.279	4.3	90	0.00
44	S	2-Fluorobiphenyl	80.000	78.193	2.3	91	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061716\
Data File : BF088091.D
Acq On : 17 Jun 2016 13:07
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 17 17:11:17 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 17 16:09:13 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	36.389	9.0	86	0.00
46		2-Chloronaphthalene	40.000	35.500	11.3	87	0.00
47		2-Nitroaniline	40.000	39.979	0.1	84	0.00
48		Acenaphthylene	40.000	31.844	20.4	74	0.00
49		Dimethylphthalate	40.000	38.634	3.4	97	0.00
50		2,6-Dinitrotoluene	40.000	41.194	-3.0	104	0.00
51	C	Acenaphthene	40.000	31.257	21.9#	71	0.00
52		3-Nitroaniline	40.000	34.802	13.0	76	0.00
53	P	2,4-Dinitrophenol	40.000	25.648	35.9#	49	0.00
54		Dibenzofuran	40.000	31.377	21.6	70	0.00
55	P	4-Nitrophenol	40.000	28.670	28.3#	57	0.00
56		2,4-Dinitrotoluene	40.000	36.848	7.9	91	0.00
57		Fluorene	40.000	36.344	9.1	82	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	38.230	4.4	82	0.00
59		Diethylphthalate	40.000	37.663	5.8	90	0.00
60		4-Chlorophenyl-phenylether	40.000	34.482	13.8	86	0.00
61		4-Nitroaniline	40.000	39.579	1.1	82	0.00
62		Azobenzene	40.000	37.355	6.6	84	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	93	0.00
64		4,6-Dinitro-2-methylphenol	40.000	33.676	15.8	68	0.00
65	c	n-Nitrosodiphenylamine	40.000	35.313	11.7	81	0.00
66		4-Bromophenyl-phenylether	40.000	34.487	13.8	77	0.00
67		Hexachlorobenzene	40.000	38.770	3.1	98	0.00
68		Atrazine	40.000	42.176	-5.4	90	0.00
69	C	Pentachlorophenol	40.000	32.523	18.7	67	0.00
70		Phenanthrene	40.000	39.480	1.3	103	0.00
71		Anthracene	40.000	34.155	14.6	76	0.00
72		Carbazole	40.000	35.906	10.2	92	0.00
73		Di-n-butylphthalate	40.000	35.545	11.1	82	0.00
74	C	Fluoranthene	40.000	32.733	18.2	75	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	77	0.00
76		Benzidine	40.000	45.035	-12.6	86	0.00
77		Pyrene	40.000	40.396	-1.0	79	0.00
78	S	Terphenyl-d14	80.000	70.914	11.4	71	0.00
79		Butylbenzylphthalate	40.000	46.421	-16.1	86	0.00
80		Benzo(a)anthracene	40.000	36.029	9.9	76	0.00
81		3,3'-Dichlorobenzidine	40.000	46.441	-16.1	84	0.00
82		Chrysene	40.000	40.862	-2.2	84	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	43.313	-8.3	86	0.00
84	c	Di-n-octyl phthalate	40.000	45.356	-13.4	88	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	43.068	-7.7	83	0.00
86	I	Perylene-d12	20.000	20.000	0.0	89	0.00
87		Benzo(b)fluoranthene	40.000	32.365	19.1	67	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061716\
Data File : BF088091.D
Acq On : 17 Jun 2016 13:07
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 17 17:11:17 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 17 16:09:13 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	39.799	0.5	108	0.00
89 C	Benzo(a)pyrene	40.000	38.339	4.2	83	0.00
90	Dibenzo(a,h)anthracene	40.000	37.507	6.2	83	0.00
91	Benzo(g,h,i)perylene	40.000	39.420	1.4	86	0.00

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