Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020818\
Data File : BF102845.D
Acq On : 8 Feb 2018 10:16
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 08 15:36:05 2018
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 05 12:44:16 2018
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	91	0.00
2		1,4-Dioxane	40.000	42.945	-7.4	96	-0.01
3		Pyridine	40.000	44.404	-11.0	98	-0.01
4		n-Nitrosodimethylamine	40.000	44.816	-12.0	99	-0.01
5	S	2-Fluorophenol	80.000	80.236	-0.3	92	0.00
6		Aniline	40.000	39.983	0.0	91	0.00
7	S	Phenol-d6	80.000	79.491	0.6	91	0.00
8		2-Chlorophenol	40.000	39.577	1.1	90	0.00
9		Benzaldehyde	40.000	38.335	4.2	87	0.00
10	C	Phenol	40.000	40.306	-0.8	94	0.00
11		bis(2-Chloroethyl)ether	40.000	39.673	0.8	91	0.00
12		1,3-Dichlorobenzene	40.000	38.713	3.2	89	0.00
13	C	1,4-Dichlorobenzene	40.000	38.641	3.4	89	0.00
14		1,2-Dichlorobenzene	40.000	38.869	2.8	89	0.00
15		Benzyl Alcohol	40.000	41.033	-2.6	92	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	36.535	8.7	82	0.00
17		2-Methylphenol	40.000	39.852	0.4	91	0.00
18		Hexachloroethane	40.000	42.185	-5.5	94	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.017	5.0	89	0.00
20		3+4-Methylphenols	40.000	38.832	2.9	87	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	93	0.00
22		Acetophenone	40.000	36.831	7.9	88	0.00
23	S	Nitrobenzene-d5	80.000	93.679	-17.1	104	0.00
24		Nitrobenzene	40.000	43.315	-8.3	98	0.00
25		Isophorone	40.000	38.670	3.3	92	0.00
26	C	2-Nitrophenol	40.000	50.754	-26.9#	135	0.00
27		2,4-Dimethylphenol	40.000	36.593	8.5	84	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.755	3.1	91	0.00
29	C	2,4-Dichlorophenol	40.000	39.870	0.3	90	0.00
30		1,2,4-Trichlorobenzene	40.000	38.013	5.0	90	0.00
31		Naphthalene	40.000	37.505	6.2	88	0.00
32		Benzoic acid	40.000	33.514	16.2	76	0.00
33		4-Chloroaniline	40.000	38.489	3.8	90	0.00
34	C	Hexachlorobutadiene	40.000	37.912	5.2	89	0.00
35		Caprolactam	40.000	40.948	-2.4	93	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.557	1.1	91	0.00
37		2-Methylnaphthalene	40.000	36.410	9.0	86	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	92	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	37.263	6.8	87	0.00
40	P	Hexachlorocyclopentadiene	40.000	39.132	2.2	84	0.00
41	S	2,4,6-Tribromophenol	80.000	87.119	-8.9	95	0.00
42	C	2,4,6-Trichlorophenol	40.000	40.390	-1.0	90	0.00
43		2,4,5-Trichlorophenol	40.000	40.256	-0.6	88	0.00
44	S	2-Fluorobiphenyl	80.000	74.847	6.4	86	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020818\
Data File : BF102845.D
Acq On : 8 Feb 2018 10:16
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 08 15:36:05 2018
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 05 12:44:16 2018
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	36.661	8.3	87	0.00
46		2-Chloronaphthalene	40.000	37.594	6.0	88	0.00
47		2-Nitroaniline	40.000	45.513	-13.8	108	0.00
48		Acenaphthylene	40.000	37.518	6.2	88	0.00
49		Dimethylphthalate	40.000	37.697	5.8	89	0.00
50		2,6-Dinitrotoluene	40.000	44.301	-10.8	98	0.00
51	C	Acenaphthene	40.000	39.100	2.2	93	0.00
52		3-Nitroaniline	40.000	44.898	-12.2	100	0.00
53	P	2,4-Dinitrophenol	40.000	54.503	-36.3#	172	0.00
54		Dibenzofuran	40.000	38.474	3.8	91	0.00
55	P	4-Nitrophenol	40.000	41.625	-4.1	99	0.00
56		2,4-Dinitrotoluene	40.000	45.332	-13.3	107	0.00
57		Fluorene	40.000	40.072	-0.2	92	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	42.230	-5.6	93	0.00
59		Diethylphthalate	40.000	38.615	3.5	91	0.00
60		4-Chlorophenyl-phenylether	40.000	40.011	-0.0	93	0.00
61		4-Nitroaniline	40.000	47.014	-17.5	107	0.00
62		Azobenzene	40.000	39.228	1.9	92	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	95	0.00
64		4,6-Dinitro-2-methylphenol	40.000	52.739	-31.8#	150	0.00
65	c	n-Nitrosodiphenylamine	40.000	37.809	5.5	90	0.00
66		4-Bromophenyl-phenylether	40.000	37.906	5.2	91	0.00
67		Hexachlorobenzene	40.000	37.812	5.5	91	0.00
68		Atrazine	40.000	38.398	4.0	89	0.00
69	C	Pentachlorophenol	40.000	37.286	6.8	84	0.00
70		Phenanthrene	40.000	36.608	8.5	88	0.00
71		Anthracene	40.000	37.150	7.1	90	0.00
72		Carbazole	40.000	37.219	7.0	90	0.00
73		Di-n-butylphthalate	40.000	38.317	4.2	90	0.00
74	C	Fluoranthene	40.000	36.915	7.7	90	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	97	0.00
76		Benzidine	40.000	45.379	-13.4	102	0.00
77		Pyrene	40.000	37.904	5.2	92	0.00
78	S	Terphenyl-d14	80.000	72.814	9.0	90	0.00
79		Butylbenzylphthalate	40.000	42.270	-5.7	101	0.00
80		Benzo(a)anthracene	40.000	39.394	1.5	97	0.00
81		3,3'-Dichlorobenzidine	40.000	39.249	1.9	90	0.00
82		Chrysene	40.000	37.408	6.5	92	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	44.403	-11.0	105	0.00
84	c	Di-n-octyl phthalate	40.000	43.390	-8.5	101	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	32.905	17.7	86	0.02
86	I	Perylene-d12	20.000	20.000	0.0	92	0.00
87		Benzo(b)fluoranthene	40.000	40.991	-2.5	92	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020818\
Data File : BF102845.D
Acq On : 8 Feb 2018 10:16
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 08 15:36:05 2018
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 05 12:44:16 2018
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	37.455	6.4	85	0.00
89 C	Benzo(a)pyrene	40.000	38.355	4.1	87	0.01
90	Dibenzo(a,h)anthracene	40.000	36.506	8.7	85	0.01
91	Benzo(g,h,i)perylene	40.000	37.113	7.2	87	0.02

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