Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA F\Data\BF031618\
Data File : BF103699.D
Acq On : 16 Mar 2018 16:32
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 16 17:15:14 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 16 13:24:19 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	102	0.00
2		1,4-Dioxane	0.647	0.650	-0.5	101	-0.01
3		Pyridine	1.781	1.798	-1.0	103	-0.01
4		n-Nitrosodimethylamine	0.657	0.649	1.2	101	-0.01
5	S	2-Fluorophenol	1.315	1.320	-0.4	102	0.00
6		Aniline	2.297	2.322	-1.1	102	0.00
7	S	Phenol-d6	1.609	1.625	-1.0	103	0.00
8		2-Chlorophenol	1.377	1.409	-2.3	103	0.00
9		Benzaldehyde	1.057	1.053	0.4	102	0.00
10	C	Phenol	1.709	1.722	-0.8	104	0.00
11		bis(2-Chloroethyl)ether	1.411	1.416	-0.4	103	0.00
12		1,3-Dichlorobenzene	1.569	1.559	0.6	102	0.00
13	C	1,4-Dichlorobenzene	1.576	1.535	2.6	100	0.00
14		1,2-Dichlorobenzene	1.463	1.463	0.0	103	0.00
15		Benzyl Alcohol	1.246	1.258	-1.0	102	0.00
16		2,2'-oxybis(1-Chloropropane	2.007	1.972	1.7	101	0.00
17		2-Methylphenol	1.227	1.248	-1.7	104	0.00
18		Hexachloroethane	0.555	0.579	-4.3	105	0.00
19	P	n-Nitroso-di-n-propylamine	1.031	1.008	2.2	101	0.00
20		3+4-Methylphenols	1.557	1.595	-2.4	104	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
22		Acetophenone	0.514	0.497	3.3	101	0.00
23	S	Nitrobenzene-d5	0.287	0.340	-18.5	116	0.00
24		Nitrobenzene	0.337	0.374	-11.0	110	0.00
25		Isophorone	0.664	0.651	2.0	103	0.00
26	C	2-Nitrophenol	0.110	0.169	-53.6#	155#	0.00
27		2,4-Dimethylphenol	0.284	0.294	-3.5	107	0.00
28		bis(2-Chloroethoxy)methane	0.402	0.398	1.0	103	0.00
29	C	2,4-Dichlorophenol	0.259	0.274	-5.8	106	0.00
30		1,2,4-Trichlorobenzene	0.300	0.300	0.0	104	0.00
31		Naphthalene	1.010	0.980	3.0	102	0.00
32		Benzoic acid	0.171	0.252	-47.4#	150	0.00
33		4-Chloroaniline	0.424	0.431	-1.7	104	0.00
34	C	Hexachlorobutadiene	0.165	0.162	1.8	102	0.00
35		Caprolactam	0.089	0.093	-4.5	107	0.00
36	C	4-Chloro-3-methylphenol	0.285	0.296	-3.9	105	0.00
37		2-Methylnaphthalene	0.641	0.630	1.7	102	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
39		1,2,4,5-Tetrachlorobenzene	0.571	0.565	1.1	104	0.00
40	P	Hexachlorocyclopentadiene	0.294	0.327	-11.2	110	0.00
41	S	2,4,6-Tribromophenol	0.172	0.201	-16.9	114	0.00
42	C	2,4,6-Trichlorophenol	0.365	0.397	-8.8	108	0.00
43		2,4,5-Trichlorophenol	0.412	0.453	-10.0	110	0.00
44	S	2-Fluorobiphenyl	1.254	1.183	5.7	103	0.00

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA F\Data\BF031618\
Data File : BF103699.D
Acq On : 16 Mar 2018 16:32
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 16 17:15:14 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 16 13:24:19 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.668	1.630	2.3	104	0.00
46		2-Chloronaphthalene	1.325	1.322	0.2	105	0.00
47		2-Nitroaniline	0.309	0.418	-35.3#	133	0.00
48		Acenaphthylene	2.054	2.024	1.5	105	0.00
49		Dimethylphthalate	1.526	1.544	-1.2	107	0.00
50		2,6-Dinitrotoluene	0.274	0.339	-23.7	122	0.00
51	C	Acenaphthene	1.220	1.240	-1.6	107	0.00
52		3-Nitroaniline	0.331	0.408	-23.3	121	0.00
53	P	2,4-Dinitrophenol	0.057	0.103	-80.7#	210#	0.00
54		Dibenzofuran	1.742	1.727	0.9	105	0.00
55	P	4-Nitrophenol	0.245	0.300	-22.4	121	0.00
56		2,4-Dinitrotoluene	0.336	0.447	-33.0#	130	0.00
57		Fluorene	1.352	1.340	0.9	105	0.00
58		2,3,4,6-Tetrachlorophenol	0.292	0.334	-14.4	112	0.00
59		Diethylphthalate	1.514	1.534	-1.3	107	0.00
60		4-Chlorophenyl-phenylether	0.618	0.616	0.3	107	0.00
61		4-Nitroaniline	0.319	0.392	-22.9	121	0.00
62		Azobenzene	1.540	1.523	1.1	104	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	107	0.00
64		4,6-Dinitro-2-methylphenol	0.065	0.107	-64.6#	178#	0.00
65	c	n-Nitrosodiphenylamine	0.681	0.672	1.3	106	0.00
66		4-Bromophenyl-phenylether	0.205	0.210	-2.4	108	0.00
67		Hexachlorobenzene	0.234	0.233	0.4	107	0.00
68		Atrazine	0.211	0.220	-4.3	111	0.00
69	C	Pentachlorophenol	0.135	0.153	-13.3	114	0.00
70		Phenanthrene	1.094	1.073	1.9	107	0.00
71		Anthracene	1.111	1.086	2.3	106	0.00
72		Carbazole	1.080	1.071	0.8	106	0.00
73		Di-n-butylphthalate	1.296	1.316	-1.5	107	0.00
74	C	Fluoranthene	1.127	1.103	2.1	105	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	105	0.00
76		Benzidine	0.853	0.758	11.1	91	0.00
77		Pyrene	1.469	1.442	1.8	105	0.00
78	S	Terphenyl-d14	0.862	0.835	3.1	106	0.00
79		Butylbenzylphthalate	0.651	0.707	-8.6	111	0.00
80		Benzo(a)anthracene	1.266	1.249	1.3	104	0.00
81		3,3'-Dichlorobenzidine	0.423	0.433	-2.4	101	0.00
82		Chrysene	1.207	1.173	2.8	103	0.00
83		Bis(2-ethylhexyl)phthalate	0.930	0.979	-5.3	106	0.00
84	c	Di-n-octyl phthalate	1.352	1.494	-10.5	106	0.00
85		Indeno(1,2,3-cd)pyrene	1.054	0.921	12.6	90	0.00
86	I	Perylene-d12	1.000	1.000	0.0	96	0.00
87		Benzo(b)fluoranthene	1.264	1.339	-5.9	103	0.00

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA F\Data\BF031618\
Data File : BF103699.D
Acq On : 16 Mar 2018 16:32
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 16 17:15:14 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 16 13:24:19 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.215	1.209	0.5	95	0.00
89 C	Benzo(a)pyrene	1.146	1.155	-0.8	96	0.00
90	Dibenzo(a,h)anthracene	0.992	0.927	6.6	90	0.00
91	Benzo(g,h,i)perylene	0.965	0.885	8.3	90	0.00

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