Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062916\
Data File : BF088510.D
Acq On : 29 Jun 2016 23:57
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 30 02:27:40 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF062916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 30 02:07:38 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	92	0.00
2		1,4-Dioxane	0.660	0.641	2.9	93	-0.01
3		Pyridine	1.920	1.891	1.5	93	-0.01
4		n-Nitrosodimethylamine	0.756	0.775	-2.5	94	-0.01
5	S	2-Fluorophenol	1.292	1.327	-2.7	92	0.00
6		Aniline	2.412	2.463	-2.1	90	0.00
7	S	Phenol-d6	1.652	1.735	-5.0	94	0.00
8		2-Chlorophenol	1.418	1.438	-1.4	95	0.00
9		Benzaldehyde	0.663	0.739	-11.5	93	0.00
10	C	Phenol	1.846	2.128	-15.3	91	0.00
11		bis(2-Chloroethyl)ether	1.496	1.415	5.4	96	0.00
12		1,3-Dichlorobenzene	1.494	1.533	-2.6	92	0.00
13	C	1,4-Dichlorobenzene	1.523	1.599	-5.0	94	0.00
14		1,2-Dichlorobenzene	1.399	1.389	0.7	97	0.00
15		Benzyl Alcohol	1.366	1.388	-1.6	89	-0.01
16		2,2'-oxybis(1-Chloropropane	2.173	2.050	5.7	90	0.00
17		2-Methylphenol	1.206	1.192	1.2	89	0.00
18		Hexachloroethane	0.587	0.587	0.0	94	0.00
19	P	n-Nitroso-di-n-propylamine	1.124	1.123	0.1	94	0.00
20		3+4-Methylphenols	1.332	1.388	-4.2	96	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	91	0.00
22		Acetophenone	0.482	0.451	6.4	90	0.00
23	S	Nitrobenzene-d5	0.381	0.407	-6.8	91	0.00
24		Nitrobenzene	0.409	0.385	5.9	93	0.00
25		Isophorone	0.769	0.722	6.1	85	-0.01
26	C	2-Nitrophenol	0.133	0.169	-27.1#	99	0.00
27		2,4-Dimethylphenol	0.345	0.322	6.7	89	0.00
28		bis(2-Chloroethoxy)methane	0.474	0.463	2.3	93	0.00
29	C	2,4-Dichlorophenol	0.278	0.291	-4.7	91	0.00
30		1,2,4-Trichlorobenzene	0.306	0.302	1.3	91	0.00
31		Naphthalene	1.015	1.013	0.2	90	-0.01
32		Benzoic acid	0.145	0.117	19.3	90	-0.01
33		4-Chloroaniline	0.453	0.437	3.5	86	0.00
34	C	Hexachlorobutadiene	0.158	0.150	5.1	89	0.00
35		Caprolactam	0.085	0.091	-7.1	97	0.00
36	C	4-Chloro-3-methylphenol	0.309	0.282	8.7	81	-0.01
37		2-Methylnaphthalene	0.624	0.674	-8.0	91	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	86	0.00
39		1,2,4,5-Tetrachlorobenzene	0.601	0.551	8.3	90	0.00
40	P	Hexachlorocyclopentadiene	0.320	0.336	-5.0	85	0.00
41	S	2,4,6-Tribromophenol	0.153	0.152	0.7	86	0.00
42	C	2,4,6-Trichlorophenol	0.399	0.431	-8.0	90	0.00
43		2,4,5-Trichlorophenol	0.399	0.452	-13.3	94	0.00
44	S	2-Fluorobiphenyl	1.363	1.392	-2.1	93	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062916\
Data File : BF088510.D
Acq On : 29 Jun 2016 23:57
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 30 02:27:40 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF062916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 30 02:07:38 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.779	1.559	12.4	82	0.00
46		2-Chloronaphthalene	1.391	1.259	9.5	86	0.00
47		2-Nitroaniline	0.380	0.391	-2.9	90	0.00
48		Acenaphthylene	2.111	2.128	-0.8	83	-0.01
49		Dimethylphthalate	1.522	1.582	-3.9	95	0.00
50		2,6-Dinitrotoluene	0.301	0.350	-16.3	86	0.00
51	C	Acenaphthene	1.243	1.291	-3.9	85	0.00
52		3-Nitroaniline	0.342	0.337	1.5	88	0.00
53	P	2,4-Dinitrophenol	0.048	0.072	-50.0#	93	0.00
54		Dibenzofuran	1.834	1.827	0.4	84	-0.01
55	P	4-Nitrophenol	0.234	0.227	3.0	83	0.00
56		2,4-Dinitrotoluene	0.376	0.404	-7.4	79	0.00
57		Fluorene	1.398	1.160	17.0	79	0.00
58		2,3,4,6-Tetrachlorophenol	0.300	0.301	-0.3	82	0.00
59		Diethylphthalate	1.430	1.492	-4.3	86	0.00
60		4-Chlorophenyl-phenylether	0.594	0.576	3.0	89	0.00
61		4-Nitroaniline	0.290	0.277	4.5	72	-0.01
62		Azobenzene	1.861	1.976	-6.2	86	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	81	0.00
64		4,6-Dinitro-2-methylphenol	0.051	0.079	-54.9#	100	0.00
65	c	n-Nitrosodiphenylamine	0.716	0.668	6.7	84	0.00
66		4-Bromophenyl-phenylether	0.216	0.216	0.0	81	0.00
67		Hexachlorobenzene	0.220	0.224	-1.8	91	0.00
68		Atrazine	0.207	0.177	14.5	72	0.00
69	C	Pentachlorophenol	0.099	0.111	-12.1	82	0.00
70		Phenanthrene	1.090	1.102	-1.1	88	0.00
71		Anthracene	1.203	1.074	10.7	78	-0.01
72		Carbazole	0.934	1.003	-7.4	83	0.00
73		Di-n-butylphthalate	1.199	1.252	-4.4	82	0.00
74	C	Fluoranthene	0.945	0.922	2.4	80	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	89	0.00
76		Benzidine	0.546	0.648	-18.7	87	0.00
77		Pyrene	1.725	1.799	-4.3	88	0.00
78	S	Terphenyl-d14	1.004	0.918	8.6	77	0.00
79		Butylbenzylphthalate	0.720	0.770	-6.9	84	0.00
80		Benzo(a)anthracene	1.350	1.419	-5.1	89	0.00
81		3,3'-Dichlorobenzidine	0.447	0.491	-9.8	92	0.00
82		Chrysene	1.253	1.397	-11.5	92	0.00
83		Bis(2-ethylhexyl)phthalate	1.001	1.022	-2.1	79	-0.01
84	c	Di-n-octyl phthalate	1.712	1.853	-8.2	90	0.00
85		Indeno(1,2,3-cd)pyrene	1.634	1.804	-10.4	101	0.00
86	I	Perylene-d12	1.000	1.000	0.0	98	0.00
87		Benzo(b)fluoranthene	1.158	1.331	-14.9	103	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062916\
Data File : BF088510.D
Acq On : 29 Jun 2016 23:57
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 30 02:27:40 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF062916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 30 02:07:38 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.037	0.863	16.8	93	0.00
89 C	Benzo(a)pyrene	1.109	1.100	0.8	97	0.00
90	Dibenzo(a,h)anthracene	1.125	1.165	-3.6	100	-0.01
91	Benzo(g,h,i)perylene	1.178	1.194	-1.4	98	0.00

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