Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061615\
Data File : BF079989.D
Acq On : 16 Jun 2015 14:14
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 17 09:51:32 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 15 15:44:31 2015
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	144	0.00
2		1,4-Dioxane	0.496	0.521	-5.0	148	0.00
3		Pyridine	1.272	1.242	2.4	150#	0.00
4		n-Nitrosodimethylamine	0.627	0.679	-8.3	170#	0.00
5	S	2-Fluorophenol	1.091	1.161	-6.4	150	0.00
6		Aniline	2.151	2.325	-8.1	147	0.00
7	S	Phenol-d6	1.486	1.557	-4.8	140	0.00
8		2-Chlorophenol	1.326	1.401	-5.7	145	0.00
9		Benzaldehyde	0.854	0.797	6.7	142	0.00
10	C	Phenol	1.634	1.591	2.6	130	0.00
11		bis(2-Chloroethyl)ether	1.315	1.230	6.5	124	0.00
12		1,3-Dichlorobenzene	1.546	1.556	-0.6	139	0.00
13	C	1,4-Dichlorobenzene	1.606	1.725	-7.4	150#	0.00
14		1,2-Dichlorobenzene	1.427	1.495	-4.8	146	0.00
15		Benzyl Alcohol	1.191	1.353	-13.6	155#	0.00
16		2,2'-oxybis(1-Chloropropane	1.948	1.964	-0.8	138	0.00
17		2-Methylphenol	1.158	1.124	2.9	133	0.00
18		Hexachloroethane	0.497	0.550	-10.7	153#	0.00
19	P	n-Nitroso-di-n-propylamine	1.062	1.034	2.6	132	0.00
20		3+4-Methylphenols	1.480	1.519	-2.6	138	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	134	0.00
22		Acetophenone	0.472	0.521	-10.4	142	0.00
23	S	Nitrobenzene-d5	0.303	0.393	-29.7#	169#	0.00
24		Nitrobenzene	0.322	0.365	-13.4	147	0.00
25		Isophorone	0.705	0.689	2.3	125	0.00
26	C	2-Nitrophenol	0.100	0.157	-57.0#	199#	0.00
27		2,4-Dimethylphenol	0.316	0.311	1.6	133	0.00
28		bis(2-Chloroethoxy)methane	0.419	0.452	-7.9	140	0.00
29	C	2,4-Dichlorophenol	0.265	0.291	-9.8	146	0.00
30		1,2,4-Trichlorobenzene	0.337	0.359	-6.5	142	0.00
31		Naphthalene	1.093	1.088	0.5	131	0.00
32		Benzoic acid	0.107	0.137	-28.0#	159#	0.00
33		4-Chloroaniline	0.424	0.561	-32.3#	170#	0.00
34	C	Hexachlorobutadiene	0.187	0.200	-7.0	146	0.00
35		Caprolactam	0.084	0.091	-8.3	126	0.00
36	C	4-Chloro-3-methylphenol	0.302	0.301	0.3	124	0.00
37		2-Methylnaphthalene	0.739	0.765	-3.5	137	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	126	0.00
39		1,2,4,5-Tetrachlorobenzene	0.670	0.706	-5.4	134	0.00
40	P	Hexachlorocyclopentadiene	0.211	0.296	-40.3#	160#	0.00
41	S	2,4,6-Tribromophenol	0.182	0.209	-14.8	128	0.00
42	C	2,4,6-Trichlorophenol	0.413	0.484	-17.2	141	0.00
43		2,4,5-Trichlorophenol	0.456	0.469	-2.9	124	0.00
44	S	2-Fluorobiphenyl	1.428	1.430	-0.1	131	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061615\
Data File : BF079989.D
Acq On : 16 Jun 2015 14:14
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 17 09:51:32 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 15 15:44:31 2015
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	1.619	1.699	-4.9	134	0.01
46		2-Chloronaphthalene	1.378	1.401	-1.7	130	0.00
47		2-Nitroaniline	0.256	0.355	-38.7#	161#	0.00
48		Acenaphthylene	2.326	2.442	-5.0	131	0.00
49		Dimethylphthalate	1.602	1.556	2.9	123	0.01
50		2,6-Dinitrotoluene	0.250	0.304	-21.6	132	0.00
51	C	Acenaphthene	1.322	1.374	-3.9	132	0.00
52		3-Nitroaniline	0.319	0.356	-11.6	128	0.00
53	P	2,4-Dinitrophenol	0.037	0.076	-105.4#	245#	0.01
54		Dibenzofuran	1.932	1.966	-1.8	127	0.00
55	P	4-Nitrophenol	0.215	0.278	-29.3#	143	0.00
56		2,4-Dinitrotoluene	0.314	0.444	-41.4#	149	0.00
57		Fluorene	1.684	1.754	-4.2	126	0.00
58		2,3,4,6-Tetrachlorophenol	0.345	0.386	-11.9	129	0.00
59		Diethylphthalate	1.602	1.517	5.3	115	0.00
60		4-Chlorophenyl-phenylether	0.752	0.736	2.1	123	0.00
61		4-Nitroaniline	0.326	0.374	-14.7	124	0.00
62		Azobenzene	1.542	1.591	-3.2	127	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	122	0.02
64		4,6-Dinitro-2-methylphenol	0.038	0.074	-94.7#	226#	0.00
65	c	n-Nitrosodiphenylamine	0.646	0.647	-0.2	119	0.00
66		4-Bromophenyl-phenylether	0.224	0.210	6.3	110	0.00
67		Hexachlorobenzene	0.244	0.239	2.0	114	0.01
68		Atrazine	0.199	0.211	-6.0	118	0.02
69	C	Pentachlorophenol	0.095	0.107	-12.6	132	0.01
70		Phenanthrene	1.192	1.144	4.0	112	0.02
71		Anthracene	1.153	1.166	-1.1	119	0.02
72		Carbazole	1.119	1.107	1.1	113	0.03
73		Di-n-butylphthalate	1.219	1.189	2.5	112	0.05
74	C	Fluoranthene	1.318	1.271	3.6	111	0.09
75	I	Chrysene-d12	1.000	1.000	0.0	111	0.24
76		Benzidine	0.554	0.588	-6.1	120	0.10
77		Pyrene	1.256	1.212	3.5	110	0.10
78	S	Terphenyl-d14	0.867	0.859	0.9	112	0.13
79		Butylbenzylphthalate	0.424	0.482	-13.7	118	0.18
80		Benzo(a)anthracene	1.270	1.204	5.2	110	0.24
81		3,3'-Dichlorobenzidine	0.406	0.408	-0.5	108	0.24
82		Chrysene	1.229	1.170	4.8	113	0.24
83		Bis(2-ethylhexyl)phthalate	0.679	0.724	-6.6	111	0.25
84	c	Di-n-octyl phthalate	1.107	1.266	-14.4	116	0.35
85		Indeno(1,2,3-cd)pyrene	1.330	1.390	-4.5	118	0.42
86	I	Perylene-d12	1.000	1.000	0.0	115	0.39
87		Benzo(b)fluoranthene	1.225	1.203	1.8	116	0.38

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061615\
Data File : BF079989.D
Acq On : 16 Jun 2015 14:14
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 17 09:51:32 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 15 15:44:31 2015
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	1.250	1.256	-0.5	110	0.38
89 C	Benzo(a)pyrene	1.165	1.141	2.1	112	0.39
90	Dibenzo(a,h)anthracene	1.146	1.137	0.8	118	0.42
91	Benzo(g,h,i)perylene	1.185	1.153	2.7	116	0.43

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