Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040517\
Data File : BM009547.D
Acq On : 05 Apr 2017 01:39
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Apr 05 02:20:44 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 31 12:10:26 2017
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	81	-0.02
2		1,4-Dioxane	0.603	0.530	12.1	74	-0.01
3		Pyridine	1.705	1.702	0.2	78	-0.01
4		n-Nitrosodimethylamine	0.908	0.942	-3.7	84	-0.01
5	S	2-Fluorophenol	1.200	1.225	-2.1	82	-0.02
6		Aniline	2.224	2.329	-4.7	82	-0.02
7	S	Phenol-d6	1.636	1.759	-7.5	83	-0.02
8		2-Chlorophenol	1.217	1.318	-8.3	83	-0.02
9		Benzaldehyde	1.288	1.158	10.1	72	-0.02
10	C	Phenol	1.773	1.871	-5.5	83	-0.01
11		bis(2-Chloroethyl)ether	1.453	1.543	-6.2	84	-0.02
12		1,3-Dichlorobenzene	1.457	1.452	0.3	79	-0.02
13	C	1,4-Dichlorobenzene	1.505	1.532	-1.8	81	-0.02
14		1,2-Dichlorobenzene	1.461	1.491	-2.1	80	-0.02
15		Benzyl Alcohol	1.430	1.530	-7.0	85	-0.02
16		2,2'-oxybis(1-Chloropropane	2.526	2.585	-2.3	82	-0.02
17		2-Methylphenol	1.160	1.281	-10.4	85	-0.02
18		Hexachloroethane	0.616	0.656	-6.5	83	-0.02
19	P	n-Nitroso-di-n-propylamine	1.319	1.449	-9.9	87	-0.02
20		3+4-Methylphenols	1.577	1.711	-8.5	84	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	87	-0.02
22		Acetophenone	0.564	0.567	-0.5	84	-0.02
23	S	Nitrobenzene-d5	0.533	0.543	-1.9	86	-0.02
24		Nitrobenzene	0.523	0.530	-1.3	85	-0.02
25		Isophorone	0.814	0.863	-6.0	88	-0.02
26	C	2-Nitrophenol	0.158	0.166	-5.1	86	-0.02
27		2,4-Dimethylphenol	0.288	0.304	-5.6	89	-0.02
28		bis(2-Chloroethoxy)methane	0.504	0.508	-0.8	86	-0.02
29	C	2,4-Dichlorophenol	0.301	0.320	-6.3	88	-0.02
30		1,2,4-Trichlorobenzene	0.374	0.368	1.6	83	-0.02
31		Naphthalene	1.062	1.070	-0.8	85	-0.02
32		Benzoic acid	0.196	0.173	11.7	75	0.00
33		4-Chloroaniline	0.409	0.429	-4.9	87	-0.02
34	C	Hexachlorobutadiene	0.256	0.243	5.1	81	-0.02
35		Caprolactam	0.092	0.104	-13.0	94	-0.01
36	C	4-Chloro-3-methylphenol	0.341	0.382	-12.0	92	-0.02
37		2-Methylnaphthalene	0.733	0.762	-4.0	88	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	95	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.755	0.723	4.2	87	-0.02
40	P	Hexachlorocyclopentadiene	0.433	0.378	12.7	78	-0.02
41	S	2,4,6-Tribromophenol	0.248	0.275	-10.9	99	-0.02
42	C	2,4,6-Trichlorophenol	0.432	0.426	1.4	87	-0.02
43		2,4,5-Trichlorophenol	0.482	0.480	0.4	89	-0.02
44	S	2-Fluorobiphenyl	1.660	1.623	2.2	89	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040517\
Data File : BM009547.D
Acq On : 05 Apr 2017 01:39
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Apr 05 02:20:44 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 31 12:10:26 2017
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	1.665	1.664	0.1	91	-0.02
46		2-Chloronaphthalene	1.293	1.268	1.9	90	-0.02
47		2-Nitroaniline	0.476	0.540	-13.4	101	-0.02
48		Acenaphthylene	2.106	2.162	-2.7	92	-0.02
49		Dimethylphthalate	1.536	1.576	-2.6	95	-0.01
50		2,6-Dinitrotoluene	0.331	0.351	-6.0	93	-0.02
51	C	Acenaphthene	1.271	1.293	-1.7	93	-0.02
52		3-Nitroaniline	0.326	0.365	-12.0	98	-0.01
53	P	2,4-Dinitrophenol	0.195	0.195	0.0	94	-0.01
54		Dibenzofuran	1.879	1.925	-2.4	95	-0.02
55	P	4-Nitrophenol	0.269	0.297	-10.4	99	-0.01
56		2,4-Dinitrotoluene	0.448	0.500	-11.6	102	-0.01
57		Fluorene	1.689	1.765	-4.5	96	-0.01
58		2,3,4,6-Tetrachlorophenol	0.407	0.433	-6.4	97	-0.02
59		Diethylphthalate	1.558	1.657	-6.4	98	-0.02
60		4-Chlorophenyl-phenylether	0.882	0.892	-1.1	94	-0.02
61		4-Nitroaniline	0.354	0.401	-13.3	106	0.00
62		Azobenzene	1.819	1.997	-9.8	100	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	-0.01
64		4,6-Dinitro-2-methylphenol	0.126	0.127	-0.8	100	-0.01
65	c	n-Nitrosodiphenylamine	0.608	0.610	-0.3	96	-0.02
66		4-Bromophenyl-phenylether	0.232	0.230	0.9	98	-0.01
67		Hexachlorobenzene	0.254	0.252	0.8	99	-0.01
68		Atrazine	0.228	0.228	0.0	98	-0.01
69	C	Pentachlorophenol	0.155	0.150	3.2	97	-0.02
70		Phenanthrene	1.144	1.158	-1.2	101	-0.02
71		Anthracene	1.158	1.198	-3.5	101	-0.02
72		Carbazole	1.109	1.211	-9.2	109	-0.01
73		Di-n-butylphthalate	1.133	1.244	-9.8	107	-0.02
74	C	Fluoranthene	1.355	1.463	-8.0	110	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	115	-0.01
76		Benzidine	0.596	0.482	19.1	89	-0.02
77		Pyrene	1.134	1.144	-0.9	111	-0.02
78	S	Terphenyl-d14	0.957	0.974	-1.8	111	-0.01
79		Butylbenzylphthalate	0.408	0.443	-8.6	117	-0.01
80		Benzo(a)anthracene	1.168	1.201	-2.8	117	0.00
81		3,3'-Dichlorobenzidine	0.437	0.460	-5.3	118	-0.01
82		Chrysene	1.115	1.137	-2.0	116	-0.01
83		Bis(2-ethylhexyl)phthalate	0.610	0.655	-7.4	117	-0.01
84	c	Di-n-octyl phthalate	1.014	1.120	-10.5	121	-0.02
85		Indeno(1,2,3-cd)pyrene	1.216	1.274	-4.8	121	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	120	-0.02
87		Benzo(b)fluoranthene	1.268	1.288	-1.6	118	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040517\
Data File : BM009547.D
Acq On : 05 Apr 2017 01:39
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Apr 05 02:20:44 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 31 12:10:26 2017
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	1.219	1.266	-3.9	121	-0.02
89 C	Benzo(a)pyrene	1.174	1.220	-3.9	121	-0.02
90	Dibenzo(a,h)anthracene	1.150	1.201	-4.4	122	-0.03
91	Benzo(g,h,i)perylene	1.121	1.154	-2.9	120	-0.03

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