Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM033117\
Data File : BM009487.D
Acq On : 31 Mar 2017 23:09
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Apr 01 02:14:31 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	85	-0.02
2		1,4-Dioxane	0.603	0.537	10.9	79	-0.01
3		Pyridine	1.705	1.681	1.4	81	-0.01
4		n-Nitrosodimethylamine	0.908	0.907	0.1	84	-0.01
5	S	2-Fluorophenol	1.200	1.194	0.5	84	-0.01
6		Aniline	2.224	2.311	-3.9	85	-0.01
7	S	Phenol-d6	1.636	1.710	-4.5	85	-0.02
8		2-Chlorophenol	1.217	1.245	-2.3	82	-0.01
9		Benzaldehyde	1.288	1.181	8.3	76	-0.01
10	C	Phenol	1.773	1.841	-3.8	85	-0.01
11		bis(2-Chloroethyl)ether	1.453	1.455	-0.1	83	-0.01
12		1,3-Dichlorobenzene	1.457	1.444	0.9	82	-0.01
13	C	1,4-Dichlorobenzene	1.505	1.486	1.3	82	-0.02
14		1,2-Dichlorobenzene	1.461	1.442	1.3	81	-0.02
15		Benzyl Alcohol	1.430	1.489	-4.1	87	-0.02
16		2,2'-oxybis(1-Chloropropane	2.526	2.523	0.1	84	-0.02
17		2-Methylphenol	1.160	1.238	-6.7	86	-0.02
18		Hexachloroethane	0.616	0.635	-3.1	84	-0.02
19	P	n-Nitroso-di-n-propylamine	1.319	1.419	-7.6	89	-0.02
20		3+4-Methylphenols	1.577	1.668	-5.8	86	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	87	-0.02
22		Acetophenone	0.564	0.572	-1.4	85	-0.02
23	S	Nitrobenzene-d5	0.533	0.546	-2.4	87	-0.01
24		Nitrobenzene	0.523	0.533	-1.9	86	-0.01
25		Isophorone	0.814	0.864	-6.1	88	-0.01
26	C	2-Nitrophenol	0.158	0.168	-6.3	87	-0.01
27		2,4-Dimethylphenol	0.288	0.301	-4.5	89	-0.02
28		bis(2-Chloroethoxy)methane	0.504	0.511	-1.4	87	-0.01
29	C	2,4-Dichlorophenol	0.301	0.314	-4.3	87	-0.02
30		1,2,4-Trichlorobenzene	0.374	0.372	0.5	84	-0.02
31		Naphthalene	1.062	1.080	-1.7	86	-0.02
32		Benzoic acid	0.196	0.189	3.6	82	0.00
33		4-Chloroaniline	0.409	0.438	-7.1	89	-0.01
34	C	Hexachlorobutadiene	0.256	0.254	0.8	85	-0.02
35		Caprolactam	0.092	0.106	-15.2	97	-0.01
36	C	4-Chloro-3-methylphenol	0.341	0.368	-7.9	89	-0.02
37		2-Methylnaphthalene	0.733	0.771	-5.2	89	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	92	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.755	0.746	1.2	88	-0.01
40	P	Hexachlorocyclopentadiene	0.433	0.398	8.1	80	-0.01
41	S	2,4,6-Tribromophenol	0.248	0.271	-9.3	95	-0.01
42	C	2,4,6-Trichlorophenol	0.432	0.451	-4.4	90	-0.01
43		2,4,5-Trichlorophenol	0.482	0.499	-3.5	90	-0.01
44	S	2-Fluorobiphenyl	1.660	1.674	-0.8	90	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM033117\
Data File : BM009487.D
Acq On : 31 Mar 2017 23:09
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Apr 01 02:14:31 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	1.689	-1.4	90	-0.01
46		2-Chloronaphthalene	1.293	1.299	-0.5	90	-0.02
47		2-Nitroaniline	0.476	0.537	-12.8	98	-0.01
48		Acenaphthylene	2.106	2.216	-5.2	93	-0.01
49		Dimethylphthalate	1.536	1.597	-4.0	94	-0.01
50		2,6-Dinitrotoluene	0.331	0.361	-9.1	94	-0.01
51	C	Acenaphthene	1.271	1.322	-4.0	93	-0.01
52		3-Nitroaniline	0.326	0.371	-13.8	97	0.00
53	P	2,4-Dinitrophenol	0.195	0.188	3.6	89	0.00
54		Dibenzofuran	1.879	1.956	-4.1	94	-0.01
55	P	4-Nitrophenol	0.269	0.291	-8.2	95	0.00
56		2,4-Dinitrotoluene	0.448	0.503	-12.3	100	-0.01
57		Fluorene	1.689	1.759	-4.1	94	-0.01
58		2,3,4,6-Tetrachlorophenol	0.407	0.439	-7.9	96	-0.01
59		Diethylphthalate	1.558	1.679	-7.8	97	-0.01
60		4-Chlorophenyl-phenylether	0.882	0.904	-2.5	93	-0.01
61		4-Nitroaniline	0.354	0.413	-16.7	106	0.00
62		Azobenzene	1.819	1.991	-9.5	97	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	100	-0.01
64		4,6-Dinitro-2-methylphenol	0.126	0.126	0.0	98	-0.01
65	c	n-Nitrosodiphenylamine	0.608	0.609	-0.2	95	-0.01
66		4-Bromophenyl-phenylether	0.232	0.233	-0.4	99	-0.01
67		Hexachlorobenzene	0.254	0.247	2.8	96	-0.01
68		Atrazine	0.228	0.238	-4.4	101	-0.01
69	C	Pentachlorophenol	0.155	0.152	1.9	98	-0.01
70		Phenanthrene	1.144	1.154	-0.9	100	-0.02
71		Anthracene	1.158	1.195	-3.2	100	-0.01
72		Carbazole	1.109	1.174	-5.9	105	0.00
73		Di-n-butylphthalate	1.133	1.221	-7.8	104	-0.01
74	C	Fluoranthene	1.355	1.445	-6.6	108	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	112	0.00
76		Benzidine	0.596	0.563	5.5	101	-0.01
77		Pyrene	1.134	1.131	0.3	106	-0.01
78	S	Terphenyl-d14	0.957	0.964	-0.7	107	0.00
79		Butylbenzylphthalate	0.408	0.447	-9.6	115	0.00
80		Benzo(a)anthracene	1.168	1.203	-3.0	114	0.00
81		3,3'-Dichlorobenzidine	0.437	0.471	-7.8	117	0.00
82		Chrysene	1.115	1.131	-1.4	112	0.00
83		Bis(2-ethylhexyl)phthalate	0.610	0.657	-7.7	114	0.00
84	c	Di-n-octyl phthalate	1.014	1.128	-11.2	118	-0.01
85		Indeno(1,2,3-cd)pyrene	1.216	1.288	-5.9	119	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	118	-0.01
87		Benzo(b)fluoranthene	1.268	1.305	-2.9	117	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM033117\
Data File : BM009487.D
Acq On : 31 Mar 2017 23:09
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Apr 01 02:14:31 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	1.252	-2.7	117	-0.01
89 C	Benzo(a)pyrene	1.174	1.216	-3.6	118	-0.02
90	Dibenzo(a,h)anthracene	1.150	1.192	-3.7	119	-0.02
91	Benzo(g,h,i)perylene	1.121	1.154	-2.9	118	-0.02

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