Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM072417\
Data File : BM010959.D
Acq On : 24 Jul 2017 12:05
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jul 24 17:33:24 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM071217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 12 14:21:16 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	99	-0.03
2		1,4-Dioxane	0.533	0.486	8.8	90	-0.02
3		Pyridine	1.456	1.211	16.8	79	-0.02
4		n-Nitrosodimethylamine	0.744	0.682	8.3	91	-0.02
5	S	2-Fluorophenol	1.207	1.192	1.2	97	-0.02
6		Aniline	2.085	1.868	10.4	88	-0.02
7	S	Phenol-d6	1.664	1.767	-6.2	105	-0.02
8		2-Chlorophenol	1.195	1.251	-4.7	102	-0.03
9		Benzaldehyde	1.156	1.157	-0.1	98	-0.02
10	C	Phenol	1.760	1.885	-7.1	105	-0.02
11		bis(2-Chloroethyl)ether	1.355	1.416	-4.5	104	-0.02
12		1,3-Dichlorobenzene	1.477	1.457	1.4	99	-0.03
13	C	1,4-Dichlorobenzene	1.525	1.497	1.8	99	-0.02
14		1,2-Dichlorobenzene	1.453	1.407	3.2	98	-0.03
15		Benzyl Alcohol	1.576	1.515	3.9	96	-0.02
16		2,2'-oxybis(1-Chloropropane	1.175	1.272	-8.3	107	-0.02
17		2-Methylphenol	1.138	1.227	-7.8	105	-0.02
18		Hexachloroethane	0.639	0.614	3.9	97	-0.03
19	P	n-Nitroso-di-n-propylamine	1.318	1.463	-11.0	112	-0.03
20		3+4-Methylphenols	1.581	1.729	-9.4	107	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	106	-0.03
22		Acetophenone	0.602	0.593	1.5	106	-0.02
23	S	Nitrobenzene-d5	0.520	0.524	-0.8	108	-0.02
24		Nitrobenzene	0.533	0.536	-0.6	108	-0.02
25		Isophorone	0.852	0.907	-6.5	113	-0.02
26	C	2-Nitrophenol	0.177	0.180	-1.7	110	-0.03
27		2,4-Dimethylphenol	0.317	0.322	-1.6	108	-0.03
28		bis(2-Chloroethoxy)methane	0.478	0.490	-2.5	108	-0.03
29	C	2,4-Dichlorophenol	0.358	0.351	2.0	105	-0.02
30		1,2,4-Trichlorobenzene	0.437	0.412	5.7	101	-0.03
31		Naphthalene	1.017	0.997	2.0	104	-0.03
32		Benzoic acid	0.248	0.270	-8.9	111	-0.02
33		4-Chloroaniline	0.410	0.342	16.6	87	-0.02
34	C	Hexachlorobutadiene	0.339	0.312	8.0	98	-0.03
35		Caprolactam	0.095	0.120	-26.3#	128	-0.02
36	C	4-Chloro-3-methylphenol	0.433	0.486	-12.2	118	-0.02
37		2-Methylnaphthalene	0.730	0.741	-1.5	109	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	119	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.872	0.783	10.2	108	-0.03
40	P	Hexachlorocyclopentadiene	0.500	0.411	17.8	94	-0.03
41	S	2,4,6-Tribromophenol	0.306	0.327	-6.9	127	-0.02
42	C	2,4,6-Trichlorophenol	0.537	0.510	5.0	115	-0.02
43		2,4,5-Trichlorophenol	0.530	0.518	2.3	113	-0.02
44	S	2-Fluorobiphenyl	1.608	1.479	8.0	112	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM072417\
Data File : BM010959.D
Acq On : 24 Jul 2017 12:05
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jul 24 17:33:24 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM071217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 12 14:21:16 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	1.742	1.623	6.8	112	-0.02
46		2-Chloronaphthalene	1.278	1.186	7.2	112	-0.02
47		2-Nitroaniline	0.420	0.471	-12.1	130	-0.02
48		Acenaphthylene	1.900	1.877	1.2	117	-0.02
49		Dimethylphthalate	1.701	1.711	-0.6	122	-0.02
50		2,6-Dinitrotoluene	0.334	0.365	-9.3	129	-0.02
51	C	Acenaphthene	1.161	1.129	2.8	118	-0.02
52		3-Nitroaniline	0.303	0.294	3.0	114	-0.02
53	P	2,4-Dinitrophenol	0.231	0.254	-10.0	129	-0.02
54		Dibenzofuran	1.878	1.842	1.9	118	-0.02
55	P	4-Nitrophenol	0.263	0.302	-14.8	134	-0.02
56		2,4-Dinitrotoluene	0.463	0.539	-16.4	134	-0.02
57		Fluorene	1.615	1.647	-2.0	123	-0.02
58		2,3,4,6-Tetrachlorophenol	0.492	0.502	-2.0	120	-0.02
59		Diethylphthalate	1.680	1.762	-4.9	125	-0.02
60		4-Chlorophenyl-phenylether	0.992	0.957	3.5	119	-0.02
61		4-Nitroaniline	0.317	0.341	-7.6	123	-0.02
62		Azobenzene	1.706	1.845	-8.1	128	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	135	-0.02
64		4,6-Dinitro-2-methylphenol	0.132	0.138	-4.5	140	-0.02
65	c	n-Nitrosodiphenylamine	0.572	0.538	5.9	127	-0.02
66		4-Bromophenyl-phenylether	0.253	0.234	7.5	123	-0.02
67		Hexachlorobenzene	0.290	0.270	6.9	128	-0.02
68		Atrazine	0.237	0.226	4.6	127	-0.02
69	C	Pentachlorophenol	0.186	0.192	-3.2	135	-0.02
70		Phenanthrene	1.040	1.028	1.2	134	-0.02
71		Anthracene	1.033	1.022	1.1	133	-0.02
72		Carbazole	0.910	0.941	-3.4	139	-0.02
73		Di-n-butylphthalate	1.076	1.163	-8.1	141	-0.02
74	C	Fluoranthene	1.289	1.400	-8.6	146	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	152#	-0.01
76		Benzidine	0.535	0.130	75.7#	34#	-0.01
77		Pyrene	1.159	1.121	3.3	147	-0.02
78	S	Terphenyl-d14	1.034	0.976	5.6	146	-0.02
79		Butylbenzylphthalate	0.406	0.442	-8.9	161#	-0.01
80		Benzo(a)anthracene	1.159	1.118	3.5	147	-0.01
81		3,3'-Dichlorobenzidine	0.456	0.322	29.4#	105	-0.02
82		Chrysene	1.104	1.064	3.6	147	-0.01
83		Bis(2-ethylhexyl)phthalate	0.591	0.651	-10.2	158#	-0.01
84	c	Di-n-octyl phthalate	0.969	1.106	-14.1	167#	-0.02
85		Indeno(1,2,3-cd)pyrene	1.266	0.929	26.6#	112	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	136	-0.02
87		Benzo(b)fluoranthene	1.270	1.347	-6.1	143	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM072417\
Data File : BM010959.D
Acq On : 24 Jul 2017 12:05
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jul 24 17:33:24 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM071217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 12 14:21:16 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	1.213	1.241	-2.3	139	-0.02
89 C	Benzo(a)pyrene	1.156	1.148	0.7	134	-0.02
90	Dibenzo(a,h)anthracene	1.122	0.882	21.4	106	-0.03
91	Benzo(g,h,i)perylene	1.103	0.907	17.8	111	-0.03

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