Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM080718\
Data File : BM016224.D
Acq On : 07 Aug 2018 19:40
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 08 02:21:20 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 07 16:15:57 2018
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	114	0.02
2		1,4-Dioxane	0.564	0.535	5.1	103	0.00
3		Pyridine	1.359	1.377	-1.3	110	0.00
4		n-Nitrosodimethylamine	1.060	1.220	-15.1	123	0.00
5	S	2-Fluorophenol	1.204	1.245	-3.4	112	0.02
6		Aniline	1.810	1.906	-5.3	115	0.02
7	S	Phenol-d6	1.572	1.685	-7.2	116	0.02
8		2-Chlorophenol	1.431	1.457	-1.8	111	0.02
9		Benzaldehyde	1.105	1.085	1.8	106	0.02
10	C	Phenol	1.572	1.654	-5.2	114	0.01
11		bis(2-Chloroethyl)ether	1.242	1.266	-1.9	114	0.02
12		1,3-Dichlorobenzene	1.675	1.607	4.1	108	0.02
13	C	1,4-Dichlorobenzene	1.699	1.647	3.1	108	0.02
14		1,2-Dichlorobenzene	1.662	1.609	3.2	111	0.02
15		Benzyl Alcohol	1.252	1.442	-15.2	126	0.02
16		2,2'-oxybis(1-Chloropropane	1.901	1.769	6.9	105	0.04
17		2-Methylphenol	1.075	1.165	-8.4	116	0.02
18		Hexachloroethane	0.734	0.760	-3.5	112	0.02
19	P	n-Nitroso-di-n-propylamine	1.149	1.245	-8.4	114	0.02
20		3+4-Methylphenols	1.546	1.639	-6.0	116	0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	116	0.03
22		Acetophenone	0.504	0.515	-2.2	116	0.02
23	S	Nitrobenzene-d5	0.430	0.468	-8.8	121	0.02
24		Nitrobenzene	0.433	0.443	-2.3	112	0.02
25		Isophorone	0.588	0.635	-8.0	119	0.02
26	C	2-Nitrophenol	0.181	0.190	-5.0	121	0.02
27		2,4-Dimethylphenol	0.252	0.257	-2.0	114	0.02
28		bis(2-Chloroethoxy)methane	0.382	0.374	2.1	107	0.02
29	C	2,4-Dichlorophenol	0.299	0.318	-6.4	115	0.02
30		1,2,4-Trichlorobenzene	0.363	0.345	5.0	109	0.02
31		Naphthalene	1.012	0.968	4.3	110	0.02
32		Benzoic acid	0.192	0.217	-13.0	132	0.04
33		4-Chloroaniline	0.413	0.437	-5.8	115	0.02
34	C	Hexachlorobutadiene	0.252	0.251	0.4	113	0.03
35		Caprolactam	0.083	0.103	-24.1	137	0.04
36	C	4-Chloro-3-methylphenol	0.329	0.370	-12.5	122	0.02
37		2-Methylnaphthalene	0.678	0.683	-0.7	112	0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	120	0.02
39		1,2,4,5-Tetrachlorobenzene	0.681	0.681	0.0	120	0.02
40	P	Hexachlorocyclopentadiene	0.278	0.327	-17.6	138	0.02
41	S	2,4,6-Tribromophenol	0.254	0.273	-7.5	130	0.02
42	C	2,4,6-Trichlorophenol	0.426	0.459	-7.7	123	0.02
43		2,4,5-Trichlorophenol	0.439	0.466	-6.2	123	0.01
44	S	2-Fluorobiphenyl	1.644	1.747	-6.3	128	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM080718\
Data File : BM016224.D
Acq On : 07 Aug 2018 19:40
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 08 02:21:20 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 07 16:15:57 2018
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	1.805	1.771	1.9	116	0.02
46		2-Chloronaphthalene	1.404	1.379	1.8	116	0.02
47		2-Nitroaniline	0.464	0.525	-13.1	131	0.02
48		Acenaphthylene	2.055	2.103	-2.3	120	0.02
49		Dimethylphthalate	1.750	1.826	-4.3	124	0.02
50		2,6-Dinitrotoluene	0.352	0.383	-8.8	125	0.02
51	C	Acenaphthene	1.264	1.256	0.6	120	0.02
52		3-Nitroaniline	0.380	0.416	-9.5	128	0.02
53	P	2,4-Dinitrophenol	0.131	0.213	-62.6#	193#	0.01
54		Dibenzofuran	2.083	2.144	-2.9	123	0.02
55	P	4-Nitrophenol	0.273	0.316	-15.8	136	0.00
56		2,4-Dinitrotoluene	0.504	0.577	-14.5	135	0.02
57		Fluorene	1.548	1.632	-5.4	125	0.02
58		2,3,4,6-Tetrachlorophenol	0.378	0.425	-12.4	128	0.02
59		Diethylphthalate	1.887	2.118	-12.2	132	0.02
60		4-Chlorophenyl-phenylether	0.862	0.906	-5.1	125	0.02
61		4-Nitroaniline	0.365	0.420	-15.1	137	0.02
62		Azobenzene	1.987	2.231	-12.3	128	0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	130	0.02
64		4,6-Dinitro-2-methylphenol	0.121	0.134	-10.7	141	0.02
65	c	n-Nitrosodiphenylamine	0.659	0.653	0.9	126	0.02
66		4-Bromophenyl-phenylether	0.226	0.226	0.0	127	0.02
67		Hexachlorobenzene	0.249	0.245	1.6	125	0.02
68		Atrazine	0.236	0.249	-5.5	132	0.02
69	C	Pentachlorophenol	0.154	0.163	-5.8	136	0.02
70		Phenanthrene	1.120	1.092	2.5	126	0.02
71		Anthracene	1.088	1.093	-0.5	129	0.02
72		Carbazole	1.127	1.178	-4.5	131	0.02
73		Di-n-butylphthalate	1.405	1.614	-14.9	145	0.02
74	C	Fluoranthene	1.249	1.348	-7.9	139	0.02
75	I	Chrysene-d12	1.000	1.000	0.0	164#	0.02
76		Benzidine	0.559	0.438	21.6	122	0.01
77		Pyrene	1.199	1.053	12.2	139	0.02
78	S	Terphenyl-d14	0.955	1.027	-7.5	173#	0.02
79		Butylbenzylphthalate	0.609	0.621	-2.0	160#	0.02
80		Benzo(a)anthracene	1.232	1.196	2.9	157#	0.02
81		3,3'-Dichlorobenzidine	0.496	0.497	-0.2	159#	0.02
82		Chrysene	1.182	1.125	4.8	155#	0.02
83		Bis(2-ethylhexyl)phthalate	0.882	0.931	-5.6	166#	0.02
84	c	Di-n-octyl phthalate	1.483	1.608	-8.4	171#	0.02
85		Indeno(1,2,3-cd)pyrene	1.402	1.466	-4.6	172#	0.05
86	I	Perylene-d12	1.000	1.000	0.0	168#	0.04
87		Benzo(b)fluoranthene	1.178	1.181	-0.3	169#	0.03

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM080718\
Data File : BM016224.D
Acq On : 07 Aug 2018 19:40
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 08 02:21:20 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 07 16:15:57 2018
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	1.193	1.145	4.0	158#	0.03
89 C	Benzo(a)pyrene	1.132	1.120	1.1	165#	0.03
90	Dibenzo(a,h)anthracene	1.173	1.230	-4.9	174#	0.05
91	Benzo(g,h,i)perylene	1.109	1.051	5.2	159#	0.05

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