Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM081718\
Data File : BM016443.D
Acq On : 17 Aug 2018 18:37
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 17 19:19:19 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM081718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 17 16:27:00 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	111	0.00
2		1,4-Dioxane	0.596	0.565	5.2	114	0.00
3		Pyridine	1.200	1.261	-5.1	121	0.00
4		n-Nitrosodimethylamine	1.382	1.415	-2.4	119	0.00
5	S	2-Fluorophenol	1.190	1.244	-4.5	117	0.00
6		Aniline	1.655	1.738	-5.0	116	0.00
7	S	Phenol-d6	1.600	1.638	-2.4	118	0.00
8		2-Chlorophenol	1.467	1.451	1.1	112	0.00
9		Benzaldehyde	1.052	1.084	-3.0	111	0.00
10	C	Phenol	1.549	1.566	-1.1	115	0.00
11		bis(2-Chloroethyl)ether	1.176	1.184	-0.7	113	0.00
12		1,3-Dichlorobenzene	1.682	1.629	3.2	108	0.00
13	C	1,4-Dichlorobenzene	1.700	1.656	2.6	106	0.00
14		1,2-Dichlorobenzene	1.670	1.678	-0.5	108	0.00
15		Benzyl Alcohol	1.452	1.477	-1.7	116	0.00
16		2,2'-oxybis(1-Chloropropane	1.481	1.457	1.6	111	0.00
17		2-Methylphenol	1.120	1.124	-0.4	113	0.00
18		Hexachloroethane	0.875	0.889	-1.6	111	0.00
19	P	n-Nitroso-di-n-propylamine	1.214	1.309	-7.8	117	0.00
20		3+4-Methylphenols	1.539	1.649	-7.1	118	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	126	0.00
22		Acetophenone	0.536	0.525	2.1	116	0.00
23	S	Nitrobenzene-d5	0.480	0.501	-4.4	123	0.00
24		Nitrobenzene	0.463	0.460	0.6	118	0.00
25		Isophorone	0.641	0.636	0.8	118	0.00
26	C	2-Nitrophenol	0.195	0.186	4.6	111	0.00
27		2,4-Dimethylphenol	0.260	0.267	-2.7	119	0.00
28		bis(2-Chloroethoxy)methane	0.371	0.370	0.3	121	0.00
29	C	2,4-Dichlorophenol	0.352	0.342	2.8	121	0.00
30		1,2,4-Trichlorobenzene	0.416	0.400	3.8	119	0.00
31		Naphthalene	1.014	0.995	1.9	123	0.00
32		Benzoic acid	0.211	0.214	-1.4	118	0.02
33		4-Chloroaniline	0.435	0.436	-0.2	126	0.00
34	C	Hexachlorobutadiene	0.359	0.347	3.3	121	0.00
35		Caprolactam	0.090	0.100	-11.1	139	0.00
36	C	4-Chloro-3-methylphenol	0.401	0.397	1.0	127	0.00
37		2-Methylnaphthalene	0.752	0.747	0.7	126	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	132	0.00
39		1,2,4,5-Tetrachlorobenzene	0.819	0.770	6.0	128	0.00
40	P	Hexachlorocyclopentadiene	0.254	0.271	-6.7	146	0.00
41	S	2,4,6-Tribromophenol	0.302	0.309	-2.3	138	0.00
42	C	2,4,6-Trichlorophenol	0.512	0.495	3.3	133	0.00
43		2,4,5-Trichlorophenol	0.499	0.502	-0.6	136	0.00
44	S	2-Fluorobiphenyl	1.923	1.960	-1.9	139	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM081718\
Data File : BM016443.D
Acq On : 17 Aug 2018 18:37
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 17 19:19:19 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM081718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 17 16:27:00 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	1.851	1.746	5.7	127	0.00
46		2-Chloronaphthalene	1.444	1.366	5.4	127	0.00
47		2-Nitroaniline	0.501	0.507	-1.2	135	0.00
48		Acenaphthylene	2.172	2.162	0.5	134	0.00
49		Dimethylphthalate	1.961	1.931	1.5	134	0.00
50		2,6-Dinitrotoluene	0.372	0.389	-4.6	138	0.00
51	C	Acenaphthene	1.310	1.274	2.7	130	0.00
52		3-Nitroaniline	0.367	0.385	-4.9	139	0.00
53	P	2,4-Dinitrophenol	0.162	0.190	-17.3	145	0.00
54		Dibenzofuran	2.309	2.348	-1.7	133	0.00
55	P	4-Nitrophenol	0.240	0.262	-9.2	146	0.00
56		2,4-Dinitrotoluene	0.622	0.655	-5.3	143	0.00
57		Fluorene	1.764	1.821	-3.2	136	0.00
58		2,3,4,6-Tetrachlorophenol	0.468	0.493	-5.3	137	0.00
59		Diethylphthalate	2.226	2.246	-0.9	134	0.00
60		4-Chlorophenyl-phenylether	1.116	1.113	0.3	131	0.00
61		4-Nitroaniline	0.355	0.382	-7.6	142	0.00
62		Azobenzene	2.567	2.551	0.6	131	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	147	0.00
64		4,6-Dinitro-2-methylphenol	0.139	0.139	0.0	141	0.00
65	c	n-Nitrosodiphenylamine	0.677	0.641	5.3	135	0.00
66		4-Bromophenyl-phenylether	0.269	0.256	4.8	135	0.00
67		Hexachlorobenzene	0.264	0.258	2.3	138	0.00
68		Atrazine	0.248	0.251	-1.2	141	0.00
69	C	Pentachlorophenol	0.151	0.152	-0.7	145	0.00
70		Phenanthrene	1.133	1.100	2.9	141	0.00
71		Anthracene	1.125	1.112	1.2	142	0.00
72		Carbazole	1.089	1.136	-4.3	150	0.00
73		Di-n-butylphthalate	1.563	1.658	-6.1	152#	0.00
74	C	Fluoranthene	1.369	1.638	-19.6	178#	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	208#	0.00
76		Benzidine	0.440	0.471	-7.0	237#	0.00
77		Pyrene	1.320	1.130	14.4	174#	0.00
78	S	Terphenyl-d14	1.323	1.241	6.2	201#	0.00
79		Butylbenzylphthalate	0.619	0.589	4.8	191#	0.00
80		Benzo(a)anthracene	1.294	1.265	2.2	204#	0.00
81		3,3'-Dichlorobenzidine	0.461	0.488	-5.9	216#	0.00
82		Chrysene	1.216	1.183	2.7	204#	0.00
83		Bis(2-ethylhexyl)phthalate	0.882	0.897	-1.7	210#	0.00
84	c	Di-n-octyl phthalate	1.324	1.529	-15.5	241#	0.00
85		Indeno(1,2,3-cd)pyrene	0.995	1.387	-39.4#	296#	0.02
86	I	Perylene-d12	1.000	1.000	0.0	253#	0.00
87		Benzo(b)fluoranthene	1.325	1.221	7.8	238#	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM081718\
Data File : BM016443.D
Acq On : 17 Aug 2018 18:37
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 17 19:19:19 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM081718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 17 16:27:00 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	1.324	1.248	5.7	244#	0.00
89 C	Benzo(a)pyrene	1.199	1.179	1.7	252#	0.00
90	Dibenzo(a,h)anthracene	1.075	1.295	-20.5	307#	0.01
91	Benzo(g,h,i)perylene	0.967	1.060	-9.6	276#	0.01

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