Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM061920\
Data File : BM026482.D
Acq On : 20 Jun 2020 01:07
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jun 20 06:49:05 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 19 10:20:20 2020
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	73	0.00
2		1,4-Dioxane	0.510	0.544	-6.7	72	0.00
3		Pyridine	1.498	1.488	0.7	68	0.00
4		n-Nitrosodimethylamine	0.742	0.829	-11.7	75	0.00
5	S	2-Fluorophenol	1.131	1.158	-2.4	70	0.00
6		Aniline	2.149	2.315	-7.7	73	0.00
7	S	Phenol-d6	1.638	1.738	-6.1	73	0.00
8		2-Chlorophenol	1.305	1.374	-5.3	73	0.00
9		Benzaldehyde	1.044	0.872	16.5	73	0.00
10	C	Phenol	1.744	1.837	-5.3	73	0.00
11		bis(2-Chloroethyl)ether	1.405	1.497	-6.5	73	0.00
12		1,3-Dichlorobenzene	1.445	1.538	-6.4	73	0.00
13	C	1,4-Dichlorobenzene	1.471	1.559	-6.0	73	0.00
14		1,2-Dichlorobenzene	1.437	1.538	-7.0	73	0.00
15		Benzyl Alcohol	1.391	1.496	-7.5	74	0.00
16		2,2'-oxybis(1-Chloropropane	2.624	2.932	-11.7	75	0.00
17		2-Methylphenol	1.275	1.374	-7.8	74	0.00
18		Hexachloroethane	0.557	0.614	-10.2	74	0.00
19	P	n-Nitroso-di-n-propylamine	1.276	1.436	-12.5	77	0.00
20		3+4-Methylphenols	1.738	1.882	-8.3	74	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	74	0.00
22		Acetophenone	0.514	0.555	-8.0	76	0.00
23	S	Nitrobenzene-d5	0.407	0.438	-7.6	75	0.00
24		Nitrobenzene	0.426	0.455	-6.8	76	0.00
25		Isophorone	0.765	0.842	-10.1	77	0.00
26	C	2-Nitrophenol	0.169	0.181	-7.1	75	0.00
27		2,4-Dimethylphenol	0.266	0.284	-6.8	75	0.00
28		bis(2-Chloroethoxy)methane	0.434	0.460	-6.0	75	0.00
29	C	2,4-Dichlorophenol	0.292	0.318	-8.9	76	0.00
30		1,2,4-Trichlorobenzene	0.321	0.346	-7.8	76	0.00
31		Naphthalene	1.008	1.057	-4.9	74	0.00
32		Benzoic acid	0.198	0.215	-8.6	74	0.00
33		4-Chloroaniline	0.440	0.465	-5.7	75	0.00
34	C	Hexachlorobutadiene	0.202	0.225	-11.4	78	0.00
35		Caprolactam	0.103	0.116	-12.6	80	0.00
36	C	4-Chloro-3-methylphenol	0.356	0.386	-8.4	77	0.00
37		2-Methylnaphthalene	0.718	0.771	-7.4	76	0.00
38		1-Methylnaphthalene	0.680	0.736	-8.2	77	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	78	0.00
40		1,2,4,5-Tetrachlorobenzene	0.569	0.609	-7.0	78	0.00
41	P	Hexachlorocyclopentadiene	0.335	0.320	4.5	69	0.00
42	S	2,4,6-Tribromophenol	0.242	0.272	-12.4	85	0.00
43	C	2,4,6-Trichlorophenol	0.385	0.416	-8.1	79	0.00
44		2,4,5-Trichlorophenol	0.447	0.483	-8.1	79	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM061920\
Data File : BM026482.D
Acq On : 20 Jun 2020 01:07
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jun 20 06:49:05 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 19 10:20:20 2020
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	S	2-Fluorobiphenyl	1.318	1.424	-8.0	79	0.00
46		1,1'-Biphenyl	1.402	1.491	-6.3	78	0.00
47		2-Chloronaphthalene	1.139	1.215	-6.7	79	0.00
48		2-Nitroaniline	0.451	0.489	-8.4	79	0.00
49		Acenaphthylene	1.822	1.948	-6.9	79	0.00
50		Dimethylphthalate	1.463	1.586	-8.4	81	0.00
51		2,6-Dinitrotoluene	0.321	0.349	-8.7	80	0.00
52	C	Acenaphthene	1.094	1.171	-7.0	79	0.00
53		3-Nitroaniline	0.357	0.377	-5.6	79	0.00
54	P	2,4-Dinitrophenol	0.195	0.211	-8.2	81	0.00
55		Dibenzofuran	1.762	1.888	-7.2	79	0.00
56	P	4-Nitrophenol	0.283	0.288	-1.8	77	0.00
57		2,4-Dinitrotoluene	0.446	0.504	-13.0	85	0.00
58		Fluorene	1.431	1.561	-9.1	81	0.00
59		2,3,4,6-Tetrachlorophenol	0.383	0.418	-9.1	81	0.00
60		Diethylphthalate	1.490	1.649	-10.7	83	0.00
61		4-Chlorophenyl-phenylether	0.759	0.841	-10.8	82	0.00
62		4-Nitroaniline	0.381	0.398	-4.5	79	0.00
63		Azobenzene	1.630	1.817	-11.5	82	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	83	0.00
65		4,6-Dinitro-2-methylphenol	0.122	0.134	-9.8	87	0.00
66	c	n-Nitrosodiphenylamine	0.579	0.611	-5.5	82	0.00
67		4-Bromophenyl-phenylether	0.220	0.239	-8.6	84	0.00
68		Hexachlorobenzene	0.230	0.247	-7.4	84	0.00
69		Atrazine	0.173	0.187	-8.1	80	0.00
70	C	Pentachlorophenol	0.138	0.146	-5.8	82	0.00
71		Phenanthrene	1.042	1.124	-7.9	85	0.00
72		Anthracene	1.039	1.126	-8.4	85	0.00
73		Carbazole	0.986	1.072	-8.7	87	0.00
74		Di-n-butylphthalate	1.162	1.340	-15.3	93	0.00
75	C	Fluoranthene	1.194	1.358	-13.7	94	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	111	0.00
77		Benzidine	0.415	0.370	10.8	87	0.00
78		Pyrene	1.321	1.276	3.4	95	0.00
79	S	Terphenyl-d14	1.030	1.055	-2.4	102	0.00
80		Butylbenzylphthalate	0.534	0.582	-9.0	115	0.00
81		Benzo(a)anthracene	1.200	1.273	-6.1	112	0.00
82		3,3'-Dichlorobenzidine	0.371	0.422	-13.7	118	0.00
83		Chrysene	1.144	1.236	-8.0	112	0.00
84		Bis(2-ethylhexyl)phthalate	0.762	0.899	-18.0	128	0.00
85	c	Di-n-octyl phthalate	1.175	1.491	-26.9#	141	0.00
86	I	Perylene-d12	1.000	1.000	0.0	111	0.00
87		Indeno(1,2,3-cd)pyrene	1.157	1.176	-1.6	99	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM061920\
Data File : BM026482.D
Acq On : 20 Jun 2020 01:07
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jun 20 06:49:05 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 19 10:20:20 2020
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(b)fluoranthene	1.203	1.316	-9.4	116	0.00
89	Benzo(k)fluoranthene	1.169	1.290	-10.4	117	0.00
90 C	Benzo(a)pyrene	1.069	1.156	-8.1	112	-0.01
91	Dibenzo(a,h)anthracene	0.966	0.990	-2.5	99	-0.01
92	Benzo(g,h,i)perylene	0.919	0.909	1.1	94	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1