Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070720\
Data File : BF120775.D
Acq On : 7 Jul 2020 9:19
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 07 10:57:49 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 02 18:51:52 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	98	0.00
2		1,4-Dioxane	0.592	0.604	-2.0	101	0.00
3		Pyridine	1.621	1.613	0.5	97	0.00
4		n-Nitrosodimethylamine	0.811	0.833	-2.7	102	0.00
5	S	2-Fluorophenol	1.296	1.304	-0.6	98	0.00
6		Aniline	2.125	2.079	2.2	96	0.00
7	S	Phenol-d6	1.655	1.625	1.8	97	0.00
8		2-Chlorophenol	1.379	1.368	0.8	95	0.00
9		Benzaldehyde	0.984	0.861	12.5	95	0.00
10	C	Phenol	1.702	1.678	1.4	97	0.00
11		bis(2-Chloroethyl)ether	1.385	1.362	1.7	97	0.00
12		1,3-Dichlorobenzene	1.518	1.518	0.0	97	0.00
13	C	1,4-Dichlorobenzene	1.521	1.528	-0.5	99	0.00
14		1,2-Dichlorobenzene	1.447	1.448	-0.1	98	0.00
15		Benzyl Alcohol	1.243	1.225	1.4	96	0.00
16		2,2'-oxybis(1-Chloropropane	2.160	2.114	2.1	96	0.00
17		2-Methylphenol	1.234	1.200	2.8	96	0.00
18		Hexachloroethane	0.572	0.572	0.0	98	0.00
19	P	n-Nitroso-di-n-propylamine	0.996	0.968	2.8	97	0.00
20		3+4-Methylphenols	1.520	1.526	-0.4	97	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	94	0.00
22		Acetophenone	0.497	0.525	-5.6	99	0.00
23	S	Nitrobenzene-d5	0.345	0.386	-11.9	101	0.00
24		Nitrobenzene	0.388	0.422	-8.8	99	0.00
25		Isophorone	0.744	0.740	0.5	94	0.00
26	C	2-Nitrophenol	0.142	0.152	-7.0	94	0.00
27		2,4-Dimethylphenol	0.297	0.301	-1.3	95	0.00
28		bis(2-Chloroethoxy)methane	0.441	0.451	-2.3	96	0.00
29	C	2,4-Dichlorophenol	0.297	0.300	-1.0	94	0.00
30		1,2,4-Trichlorobenzene	0.333	0.341	-2.4	95	0.00
31		Naphthalene	1.069	1.092	-2.2	96	0.00
32		Benzoic acid	0.161	0.173	-7.5	81	0.00
33		4-Chloroaniline	0.454	0.453	0.2	94	0.00
34	C	Hexachlorobutadiene	0.200	0.210	-5.0	96	0.00
35		Caprolactam	0.086	0.085	1.2	92	0.00
36	C	4-Chloro-3-methylphenol	0.309	0.308	0.3	92	0.00
37		2-Methylnaphthalene	0.695	0.699	-0.6	93	0.00
38		1-Methylnaphthalene	0.651	0.659	-1.2	94	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	92	0.00
40		1,2,4,5-Tetrachlorobenzene	0.604	0.652	-7.9	95	0.00
41	P	Hexachlorocyclopentadiene	0.357	0.373	-4.5	88	0.00
42	S	2,4,6-Tribromophenol	0.201	0.213	-6.0	90	0.00
43	C	2,4,6-Trichlorophenol	0.412	0.412	0.0	88	0.00
44		2,4,5-Trichlorophenol	0.457	0.480	-5.0	92	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070720\
Data File : BF120775.D
Acq On : 7 Jul 2020 9:19
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 07 10:57:49 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 02 18:51:52 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	1.322	1.413	-6.9	96	0.00
46		1,1'-Biphenyl	1.613	1.713	-6.2	94	0.00
47		2-Chloronaphthalene	1.292	1.356	-5.0	94	0.00
48		2-Nitroaniline	0.346	0.431	-24.6	105	0.00
49		Acenaphthylene	2.026	2.094	-3.4	94	0.00
50		Dimethylphthalate	1.459	1.485	-1.8	92	0.00
51		2,6-Dinitrotoluene	0.264	0.297	-12.5	96	0.00
52	C	Acenaphthene	1.257	1.320	-5.0	94	0.00
53		3-Nitroaniline	0.326	0.377	-15.6	100	0.00
54	P	2,4-Dinitrophenol	0.087	0.106	-21.8	114	0.00
55		Dibenzofuran	1.812	1.885	-4.0	93	0.00
56	P	4-Nitrophenol	0.239	0.264	-10.5	93	0.00
57		2,4-Dinitrotoluene	0.316	0.381	-20.6	99	0.00
58		Fluorene	1.349	1.449	-7.4	96	0.00
59		2,3,4,6-Tetrachlorophenol	0.340	0.344	-1.2	89	0.00
60		Diethylphthalate	1.520	1.550	-2.0	92	0.00
61		4-Chlorophenyl-phenylether	0.678	0.731	-7.8	96	0.00
62		4-Nitroaniline	0.310	0.361	-16.5	100	0.00
63		Azobenzene	1.555	1.626	-4.6	96	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	89	0.00
65		4,6-Dinitro-2-methylphenol	0.064	0.072	-12.5	89	0.00
66	c	n-Nitrosodiphenylamine	0.670	0.711	-6.1	92	0.00
67		4-Bromophenyl-phenylether	0.232	0.244	-5.2	90	0.00
68		Hexachlorobenzene	0.243	0.256	-5.3	89	0.00
69		Atrazine	0.173	0.173	0.0	85	0.00
70	C	Pentachlorophenol	0.138	0.136	1.4	82	0.00
71		Phenanthrene	1.085	1.158	-6.7	93	0.00
72		Anthracene	1.097	1.175	-7.1	93	0.00
73		Carbazole	1.054	1.117	-6.0	93	0.00
74		Di-n-butylphthalate	1.305	1.362	-4.4	90	0.00
75	C	Fluoranthene	1.156	1.202	-4.0	91	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	92	0.00
77		Benzidine	0.565	0.522	7.6	78	0.00
78		Pyrene	1.597	1.628	-1.9	92	0.00
79	S	Terphenyl-d14	1.022	1.093	-6.9	99	0.00
80		Butylbenzylphthalate	0.690	0.672	2.6	86	0.00
81		Benzo(a)anthracene	1.282	1.385	-8.0	96	0.00
82		3,3'-Dichlorobenzidine	0.433	0.423	2.3	85	0.00
83		Chrysene	1.308	1.355	-3.6	93	0.00
84		Bis(2-ethylhexyl)phthalate	0.904	0.956	-5.8	95	0.00
85	c	Di-n-octyl phthalate	1.600	1.526	4.6	84	0.00
86	I	Perylene-d12	1.000	1.000	0.0	87	0.00
87		Indeno(1,2,3-cd)pyrene	1.283	1.374	-7.1	88	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070720\
Data File : BF120775.D
Acq On : 7 Jul 2020 9:19
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 07 10:57:49 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 02 18:51:52 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	1.315	1.378	-4.8	86	0.00
89	Benzo(k)fluoranthene	1.279	1.424	-11.3	91	0.00
90 C	Benzo(a)pyrene	1.173	1.243	-6.0	87	0.00
91	Dibenzo(a,h)anthracene	1.069	1.160	-8.5	89	0.00
92	Benzo(g,h,i)perylene	0.977	1.054	-7.9	90	0.00

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