Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031821\
Data File : BF123379.D
Acq On : 18 Mar 2021 11:17
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 18 12:34:35 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 18 02:38:45 2021
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	95	0.00
2		1,4-Dioxane	0.655	0.663	-1.2	100	0.00
3		Pyridine	1.724	1.734	-0.6	97	-0.01
4		n-Nitrosodimethylamine	0.765	0.759	0.8	94	-0.01
5	S	2-Fluorophenol	1.408	1.314	6.7	98	0.00
6		Aniline	2.155	1.912	11.3	98	0.00
7	S	Phenol-d6	1.829	1.589	13.1	97	0.00
8		2-Chlorophenol	1.520	1.393	8.4	95	0.00
9		Benzaldehyde	0.948	0.814	14.1	99	0.00
10	C	Phenol	1.904	1.703	10.6	98	0.00
11		bis(2-Chloroethyl)ether	1.455	1.407	3.3	95	0.00
12		1,3-Dichlorobenzene	1.508	1.436	4.8	95	0.00
13	C	1,4-Dichlorobenzene	1.575	1.426	9.5	95	0.00
14		1,2-Dichlorobenzene	1.530	1.314	14.1	96	0.00
15		Benzyl Alcohol	1.172	1.008	14.0	97	0.00
16		2,2'-oxybis(1-Chloropropane	2.402	2.353	2.0	97	0.00
17		2-Methylphenol	1.203	1.115	7.3	96	0.00
18		Hexachloroethane	0.546	0.530	2.9	96	0.00
19	P	n-Nitroso-di-n-propylamine	1.004	0.849	15.4	95	0.00
20		3+4-Methylphenols	1.452	1.261	13.2	97	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
22		Acetophenone	0.478	0.406	15.1	94	0.00
23	S	Nitrobenzene-d5	0.341	0.328	3.8	95	0.00
24		Nitrobenzene	0.377	0.362	4.0	95	0.00
25		Isophorone	0.707	0.686	3.0	95	0.00
26	C	2-Nitrophenol	0.204	0.200	2.0	94	0.00
27		2,4-Dimethylphenol	0.295	0.285	3.4	96	0.00
28		bis(2-Chloroethoxy)methane	0.410	0.399	2.7	96	0.00
29	C	2,4-Dichlorophenol	0.288	0.280	2.8	95	0.00
30		1,2,4-Trichlorobenzene	0.280	0.264	5.7	95	0.00
31		Naphthalene	1.048	0.949	9.4	96	0.00
32		Benzoic acid	0.218	0.228	-4.6	93	0.00
33		4-Chloroaniline	0.431	0.423	1.9	95	0.00
34	C	Hexachlorobutadiene	0.137	0.130	5.1	96	0.00
35		Caprolactam	0.100	0.101	-1.0	94	0.00
36	C	4-Chloro-3-methylphenol	0.302	0.294	2.6	95	0.00
37		2-Methylnaphthalene	0.691	0.623	9.8	95	0.00
38		1-Methylnaphthalene	0.650	0.584	10.2	94	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	95	0.00
40		1,2,4,5-Tetrachlorobenzene	0.470	0.441	6.2	95	0.00
41	P	Hexachlorocyclopentadiene	0.161	0.158	1.9	87	0.00
42	S	2,4,6-Tribromophenol	0.140	0.134	4.3	93	0.00
43	C	2,4,6-Trichlorophenol	0.376	0.365	2.9	94	0.00
44		2,4,5-Trichlorophenol	0.406	0.387	4.7	91	0.00
45	S	2-Fluorobiphenyl	1.099	1.015	7.6	95	0.00
46		1,1'-Biphenyl	1.631	1.396	14.4	95	0.00
47		2-Chloronaphthalene	1.293	1.176	9.0	96	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031821\
Data File : BF123379.D
Acq On : 18 Mar 2021 11:17
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 18 12:34:35 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 18 02:38:45 2021
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)
48		2-Nitroaniline	0.423	0.411	2.8	92	0.00
49		Acenaphthylene	1.994	1.792	10.1	94	0.00
50		Dimethylphthalate	1.414	1.347	4.7	94	0.00
51		2,6-Dinitrotoluene	0.337	0.330	2.1	93	0.00
52	C	Acenaphthene	1.172	1.037	11.5	93	0.00
53		3-Nitroaniline	0.422	0.418	0.9	94	0.00
54	P	2,4-Dinitrophenol	0.168	0.172	-2.4	88	0.00
55		Dibenzofuran	1.764	1.539	12.8	94	0.00
56	P	4-Nitrophenol	0.282	0.299	-6.0	90	0.00
57		2,4-Dinitrotoluene	0.429	0.424	1.2	94	0.00
58		Fluorene	1.316	1.134	13.8	94	0.00
59		2,3,4,6-Tetrachlorophenol	0.287	0.284	1.0	93	0.00
60		Diethylphthalate	1.400	1.274	9.0	94	0.00
61		4-Chlorophenyl-phenylether	0.534	0.468	12.4	94	0.00
62		4-Nitroaniline	0.419	0.424	-1.2	94	0.00
63		Azobenzene	1.497	1.380	7.8	96	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	94	0.00
65		4,6-Dinitro-2-methylphenol	0.141	0.144	-2.1	92	0.00
66	c	n-Nitrosodiphenylamine	0.715	0.653	8.7	94	0.00
67		4-Bromophenyl-phenylether	0.190	0.179	5.8	95	0.00
68		Hexachlorobenzene	0.198	0.185	6.6	95	0.00
69		Atrazine	0.191	0.184	3.7	95	0.00
70	C	Pentachlorophenol	0.107	0.110	-2.8	91	0.00
71		Phenanthrene	1.221	1.025	16.1	96	0.00
72		Anthracene	1.228	1.036	15.6	96	0.00
73		Carbazole	1.250	1.069	14.5	94	0.00
74		Di-n-butylphthalate	1.420	1.244	12.4	98	0.00
75	C	Fluoranthene	1.134	1.020	10.1	95	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	101	0.00
77		Benzidine	0.726	0.745	-2.6	110	0.00
78		Pyrene	1.835	1.655	9.8	98	0.00
79	S	Terphenyl-d14	1.027	0.843	17.9	97	0.00
80		Butylbenzylphthalate	0.865	0.835	3.5	100	0.00
81		Benzo(a)anthracene	1.318	1.202	8.8	100	0.00
82		3,3'-Dichlorobenzidine	0.450	0.434	3.6	101	0.00
83		Chrysene	1.361	1.249	8.2	100	0.00
84		Bis(2-ethylhexyl)phthalate	1.042	0.979	6.0	100	0.00
85	c	Di-n-octyl phthalate	1.773	1.764	0.5	105	0.00
86	I	Perylene-d12	1.000	1.000	0.0	107	0.00
87		Indeno(1,2,3-cd)pyrene	1.340	1.159	13.5	101	0.00
88		Benzo(b)fluoranthene	1.437	1.341	6.7	99	0.00
89		Benzo(k)fluoranthene	1.287	1.270	1.3	113	0.00
90	C	Benzo(a)pyrene	1.268	1.203	5.1	106	0.00
91		Dibenzo(a,h)anthracene	1.107	0.959	13.4	102	0.00
92		Benzo(g,h,i)perylene	1.177	1.008	14.4	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031821\
Data File : BF123379.D
Acq On : 18 Mar 2021 11:17
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Mar 18 12:34:35 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 18 02:38:45 2021
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)

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