Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060821\
Data File : BF124184.D
Acq On : 8 Jun 2021 17:19
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 08 18:08:13 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 07 13:32:21 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	126	0.00
2		1,4-Dioxane	0.581	0.562	3.3	125	0.00
3		Pyridine	1.489	1.371	7.9	118	0.00
4		n-Nitrosodimethylamine	0.885	0.814	8.0	118	0.02
5	S	2-Fluorophenol	1.329	1.160	12.7	112	0.00
6		Aniline	2.040	1.825	10.5	119	0.00
7	S	Phenol-d6	1.786	1.607	10.0	118	0.00
8		2-Chlorophenol	1.478	1.334	9.7	115	0.00
9		Benzaldehyde	0.794	0.749	5.7	128	0.00
10	C	Phenol	1.930	1.737	10.0	118	0.00
11		bis(2-Chloroethyl)ether	1.407	1.279	9.1	116	0.00
12		1,3-Dichlorobenzene	1.728	1.599	7.5	120	0.00
13	C	1,4-Dichlorobenzene	1.734	1.619	6.6	121	0.00
14		1,2-Dichlorobenzene	1.658	1.547	6.7	121	0.00
15		Benzyl Alcohol	1.610	1.496	7.1	123	0.00
16		2,2'-oxybis(1-Chloropropane	2.195	1.847	15.9	112	0.00
17		2-Methylphenol	1.202	1.109	7.7	120	0.00
18		Hexachloroethane	0.683	0.645	5.6	122	0.00
19	P	n-Nitroso-di-n-propylamine	1.261	1.168	7.4	126	0.00
20		3+4-Methylphenols	1.592	1.423	10.6	117	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	126	0.00
22		Acetophenone	0.570	0.545	4.4	124	0.00
23	S	Nitrobenzene-d5	0.504	0.462	8.3	121	0.00
24		Nitrobenzene	0.506	0.456	9.9	116	0.00
25		Isophorone	0.909	0.829	8.8	121	0.00
26	C	2-Nitrophenol	0.226	0.213	5.8	123	0.00
27		2,4-Dimethylphenol	0.319	0.296	7.2	122	0.00
28		bis(2-Chloroethoxy)methane	0.450	0.420	6.7	127	0.00
29	C	2,4-Dichlorophenol	0.369	0.349	5.4	125	0.00
30		1,2,4-Trichlorobenzene	0.413	0.386	6.5	125	0.00
31		Naphthalene	1.196	1.088	9.0	120	0.00
32		Benzoic acid	0.196	0.177	9.7	126	0.01
33		4-Chloroaniline	0.445	0.418	6.1	124	0.00
34	C	Hexachlorobutadiene	0.292	0.268	8.2	120	0.00
35		Caprolactam	0.101	0.092	8.9	122	0.00
36	C	4-Chloro-3-methylphenol	0.403	0.381	5.5	130	0.00
37		2-Methylnaphthalene	0.839	0.797	5.0	124	0.00
38		1-Methylnaphthalene	0.784	0.724	7.7	121	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	122	0.00
40		1,2,4,5-Tetrachlorobenzene	0.709	0.669	5.6	126	0.00
41	P	Hexachlorocyclopentadiene	0.314	0.312	0.6	125	0.00
42	S	2,4,6-Tribromophenol	0.283	0.260	8.1	121	0.00
43	C	2,4,6-Trichlorophenol	0.462	0.454	1.7	126	0.00
44		2,4,5-Trichlorophenol	0.496	0.461	7.1	121	0.00
45	S	2-Fluorobiphenyl	1.589	1.520	4.3	125	0.00
46		1,1'-Biphenyl	1.688	1.611	4.6	127	0.00
47		2-Chloronaphthalene	1.374	1.288	6.3	126	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060821\
Data File : BF124184.D
Acq On : 8 Jun 2021 17:19
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 08 18:08:13 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 07 13:32:21 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	0.496	0.465	6.2	123	0.00
49		Acenaphthylene	1.998	1.797	10.1	118	0.00
50		Dimethylphthalate	1.655	1.524	7.9	126	0.00
51		2,6-Dinitrotoluene	0.379	0.336	11.3	121	0.00
52	C	Acenaphthene	1.305	1.178	9.7	119	0.00
53		3-Nitroaniline	0.355	0.322	9.3	121	0.00
54	P	2,4-Dinitrophenol	0.160	0.173	-8.1	128	0.00
55		Dibenzofuran	2.042	1.845	9.6	120	0.00
56	P	4-Nitrophenol	0.254	0.223	12.2	111	0.01
57		2,4-Dinitrotoluene	0.500	0.470	6.0	119	0.00
58		Fluorene	1.567	1.434	8.5	119	0.00
59		2,3,4,6-Tetrachlorophenol	0.402	0.360	10.4	121	0.00
60		Diethylphthalate	1.670	1.507	9.8	119	0.00
61		4-Chlorophenyl-phenylether	0.801	0.713	11.0	120	0.00
62		4-Nitroaniline	0.357	0.311	12.9	113	0.00
63		Azobenzene	1.741	1.530	12.1	119	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	122	0.00
65		4,6-Dinitro-2-methylphenol	0.161	0.136	15.5	111	0.00
66	c	n-Nitrosodiphenylamine	0.760	0.664	12.6	115	0.00
67		4-Bromophenyl-phenylether	0.275	0.248	9.8	124	0.00
68		Hexachlorobenzene	0.311	0.282	9.3	120	0.00
69		Atrazine	0.215	0.212	1.4	118	0.00
70	C	Pentachlorophenol	0.146	0.167	-14.4	146	0.00
71		Phenanthrene	1.249	1.159	7.2	121	0.00
72		Anthracene	1.258	1.168	7.2	126	0.00
73		Carbazole	1.096	0.997	9.0	120	0.00
74		Di-n-butylphthalate	1.409	1.368	2.9	130	0.00
75	C	Fluoranthene	1.311	1.168	10.9	117	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	114	0.00
77		Benzidine	0.429	0.435	-1.4	101	0.00
78		Pyrene	1.921	1.851	3.6	116	0.00
79	S	Terphenyl-d14	1.457	1.445	0.8	118	0.00
80		Butylbenzylphthalate	0.779	0.806	-3.5	125	0.00
81		Benzo(a)anthracene	1.457	1.345	7.7	113	0.00
82		3,3'-Dichlorobenzidine	0.427	0.423	0.9	121	0.00
83		Chrysene	1.406	1.295	7.9	112	0.00
84		Bis(2-ethylhexyl)phthalate	0.927	1.009	-8.8	131	0.00
85	c	Di-n-octyl phthalate	1.237	1.383	-11.8	137	0.00
86	I	Perylene-d12	1.000	1.000	0.0	121	0.00
87		Indeno(1,2,3-cd)pyrene	1.448	1.385	4.4	114	0.01
88		Benzo(b)fluoranthene	1.554	1.481	4.7	128	0.00
89		Benzo(k)fluoranthene	1.471	1.310	10.9	116	0.00
90	C	Benzo(a)pyrene	1.349	1.253	7.1	122	0.00
91		Dibenzo(a,h)anthracene	1.252	1.207	3.6	117	0.00
92		Benzo(g,h,i)perylene	1.235	1.121	9.2	108	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060821\
Data File : BF124184.D
Acq On : 8 Jun 2021 17:19
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jun 08 18:08:13 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 07 13:32:21 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