Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062921\
Data File : BF124488.D
Acq On : 29 Jun 2021 12:15
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 29 12:49:45 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF062321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 29 12:49:33 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	89	0.00
2		1,4-Dioxane	0.547	0.553	-1.1	88	0.00
3		Pyridine	1.199	1.119	6.7	77	0.00
4		n-Nitrosodimethylamine	0.647	0.596	7.9	81	0.00
5	S	2-Fluorophenol	1.272	1.252	1.6	85	0.00
6		Aniline	1.855	1.807	2.6	86	0.00
7	S	Phenol-d6	1.660	1.669	-0.5	87	0.00
8		2-Chlorophenol	1.413	1.431	-1.3	89	0.00
9		Benzaldehyde	0.773	0.839	-8.5	90	0.00
10	C	Phenol	1.795	1.715	4.5	88	0.00
11		bis(2-Chloroethyl)ether	1.298	1.240	4.5	86	0.00
12		1,3-Dichlorobenzene	1.717	1.663	3.1	86	0.00
13	C	1,4-Dichlorobenzene	1.741	1.675	3.8	87	0.00
14		1,2-Dichlorobenzene	1.639	1.631	0.5	89	0.00
15		Benzyl Alcohol	1.424	1.364	4.2	87	0.00
16		2,2'-oxybis(1-Chloropropane	1.666	1.518	8.9	82	0.00
17		2-Methylphenol	1.131	1.110	1.9	89	0.00
18		Hexachloroethane	0.634	0.634	0.0	88	0.00
19	P	n-Nitroso-di-n-propylamine	1.117	1.098	1.7	90	0.00
20		3+4-Methylphenols	1.476	1.428	3.3	85	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
22		Acetophenone	0.557	0.555	0.4	88	0.00
23	S	Nitrobenzene-d5	0.482	0.459	4.8	86	0.00
24		Nitrobenzene	0.482	0.470	2.5	94	0.00
25		Isophorone	0.829	0.794	4.2	90	0.00
26	C	2-Nitrophenol	0.234	0.220	6.0	88	0.00
27		2,4-Dimethylphenol	0.312	0.303	2.9	90	0.00
28		bis(2-Chloroethoxy)methane	0.420	0.411	2.1	89	0.00
29	C	2,4-Dichlorophenol	0.371	0.352	5.1	88	0.00
30		1,2,4-Trichlorobenzene	0.421	0.408	3.1	88	0.00
31		Naphthalene	1.177	1.147	2.5	91	0.00
32		Benzoic acid	0.162	0.143	11.7	79	0.00
33		4-Chloroaniline	0.442	0.440	0.5	90	0.00
34	C	Hexachlorobutadiene	0.281	0.265	5.7	91	0.00
35		Caprolactam	0.097	0.093	4.1	86	0.00
36	C	4-Chloro-3-methylphenol	0.375	0.359	4.3	91	0.00
37		2-Methylnaphthalene	0.833	0.803	3.6	89	0.00
38		1-Methylnaphthalene	0.774	0.727	6.1	88	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	94	0.00
40		1,2,4,5-Tetrachlorobenzene	0.719	0.712	1.0	88	0.00
41	P	Hexachlorocyclopentadiene	0.056	0.032#	42.9#	90	0.00
42	S	2,4,6-Tribromophenol	0.239	0.239	0.0	92	0.00
43	C	2,4,6-Trichlorophenol	0.451	0.432	4.2	89	0.00
44		2,4,5-Trichlorophenol	0.476	0.423	11.1	81	0.00
45	S	2-Fluorobiphenyl	1.633	1.633	0.0	91	0.00
46		1,1'-Biphenyl	1.700	1.727	-1.6	91	0.00
47		2-Chloronaphthalene	1.390	1.365	1.8	93	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062921\
Data File : BF124488.D
Acq On : 29 Jun 2021 12:15
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 29 12:49:45 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF062321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 29 12:49:33 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	0.457	0.414	9.4	80	0.00
49		Acenaphthylene	2.010	1.972	1.9	92	0.00
50		Dimethylphthalate	1.577	1.556	1.3	92	0.00
51		2,6-Dinitrotoluene	0.364	0.350	3.8	90	0.00
52	C	Acenaphthene	1.276	1.253	1.8	92	0.00
53		3-Nitroaniline	0.322	0.335	-4.0	96	0.00
54	P	2,4-Dinitrophenol	0.109	0.096	11.9	78	0.00
55		Dibenzofuran	2.034	1.975	2.9	91	0.00
56	P	4-Nitrophenol	0.169	0.127	24.9	70	0.00
57		2,4-Dinitrotoluene	0.483	0.484	-0.2	93	0.00
58		Fluorene	1.563	1.530	2.1	93	0.00
59		2,3,4,6-Tetrachlorophenol	0.348	0.337	3.2	88	0.00
60		Diethylphthalate	1.569	1.572	-0.2	97	0.00
61		4-Chlorophenyl-phenylether	0.775	0.741	4.4	90	0.00
62		4-Nitroaniline	0.310	0.294	5.2	90	0.00
63		Azobenzene	1.562	1.483	5.1	91	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	92	0.00
65		4,6-Dinitro-2-methylphenol	0.143	0.136	4.9	86	0.00
66	c	n-Nitrosodiphenylamine	0.767	0.735	4.2	91	0.00
67		4-Bromophenyl-phenylether	0.269	0.252	6.3	90	0.00
68		Hexachlorobenzene	0.292	0.283	3.1	95	0.00
69		Atrazine	0.198	0.216	-9.1	95	0.00
70	C	Pentachlorophenol	0.061	0.056	8.2	82	0.00
71		Phenanthrene	1.261	1.207	4.3	92	0.00
72		Anthracene	1.256	1.223	2.6	92	0.00
73		Carbazole	1.095	1.047	4.4	93	0.00
74		Di-n-butylphthalate	1.304	1.332	-2.1	100	0.00
75	C	Fluoranthene	1.269	1.201	5.4	94	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	93	0.00
77		Benzidine	0.401	0.530	-32.2#	125	0.00
78		Pyrene	1.826	1.774	2.8	97	0.00
79	S	Terphenyl-d14	1.335	1.368	-2.5	98	0.00
80		Butylbenzylphthalate	0.725	0.745	-2.8	104	0.00
81		Benzo(a)anthracene	1.442	1.460	-1.2	99	0.00
82		3,3'-Dichlorobenzidine	0.462	0.462	0.0	102	0.00
83		Chrysene	1.429	1.407	1.5	97	0.00
84		Bis(2-ethylhexyl)phthalate	0.971	1.041	-7.2	106	0.00
85	c	Di-n-octyl phthalate	1.513	1.709	-13.0	111	0.00
86	I	Perylene-d12	1.000	1.000	0.0	94	0.00
87		Indeno(1,2,3-cd)pyrene	1.577	1.596	-1.2	95	0.00
88		Benzo(b)fluoranthene	1.516	1.387	8.5	91	0.00
89		Benzo(k)fluoranthene	1.387	1.393	-0.4	103	0.00
90	C	Benzo(a)pyrene	1.347	1.284	4.7	98	0.00
91		Dibenzo(a,h)anthracene	1.359	1.401	-3.1	97	0.00
92		Benzo(g,h,i)perylene	1.337	1.320	1.3	93	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062921\
Data File : BF124488.D
Acq On : 29 Jun 2021 12:15
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jun 29 12:49:45 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF062321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 29 12:49:33 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 = 1 CCC's out = 0