Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022122\
Data File : BF126973.D
Acq On : 21 Feb 2022 15:00
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Feb 21 17:01:54 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 21 13:16:57 2022
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	80	0.00
2		1,4-Dioxane	0.592	0.588	0.7	82	-0.01
3		Pyridine	1.553	1.517	2.3	79	-0.01
4		n-Nitrosodimethylamine	0.761	0.753	1.1	78	-0.01
5	S	2-Fluorophenol	1.294	1.256	2.9	79	0.00
6		Aniline	2.069	1.987	4.0	76	0.00
7	S	Phenol-d6	1.746	1.680	3.8	78	0.00
8		2-Chlorophenol	1.388	1.360	2.0	79	0.00
9		Benzaldehyde	1.063	0.902	15.1	76	0.00
10	C	Phenol	1.764	1.696	3.9	79	0.00
11		bis(2-Chloroethyl)ether	1.384	1.282	7.4	75	0.00
12		1,3-Dichlorobenzene	1.498	1.447	3.4	79	0.00
13	C	1,4-Dichlorobenzene	1.518	1.494	1.6	81	0.00
14		1,2-Dichlorobenzene	1.448	1.398	3.5	79	0.00
15		Benzyl Alcohol	1.471	1.443	1.9	78	0.00
16		2,2'-oxybis(1-Chloropropane	2.179	1.931	11.4	71	0.00
17		2-Methylphenol	1.201	1.155	3.8	77	0.00
18		Hexachloroethane	0.582	0.576	1.0	81	0.00
19	P	n-Nitroso-di-n-propylamine	1.125	1.050	6.7	76	0.00
20		3+4-Methylphenols	1.560	1.504	3.6	77	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	75	0.00
22		Acetophenone	0.527	0.518	1.7	77	0.00
23	S	Nitrobenzene-d5	0.443	0.441	0.5	78	0.00
24		Nitrobenzene	0.419	0.411	1.9	76	0.00
25		Isophorone	0.733	0.696	5.0	74	0.00
26	C	2-Nitrophenol	0.178	0.180	-1.1	76	0.00
27		2,4-Dimethylphenol	0.293	0.287	2.0	77	0.00
28		bis(2-Chloroethoxy)methane	0.445	0.430	3.4	76	0.00
29	C	2,4-Dichlorophenol	0.275	0.271	1.5	76	0.00
30		1,2,4-Trichlorobenzene	0.308	0.308	0.0	79	0.00
31		Naphthalene	1.044	1.033	1.1	78	0.00
32		Benzoic acid	0.208	0.196	5.8	67	0.00
33		4-Chloroaniline	0.411	0.406	1.2	77	0.00
34	C	Hexachlorobutadiene	0.180	0.181	-0.6	79	0.00
35		Caprolactam	0.096	0.096	0.0	76	0.00
36	C	4-Chloro-3-methylphenol	0.352	0.353	-0.3	76	0.00
37		2-Methylnaphthalene	0.677	0.668	1.3	77	0.00
38		1-Methylnaphthalene	0.649	0.646	0.5	78	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	75	0.00
40		1,2,4,5-Tetrachlorobenzene	0.526	0.533	-1.3	80	0.00
41	P	Hexachlorocyclopentadiene	0.286	0.291	-1.7	76	0.00
42	S	2,4,6-Tribromophenol	0.230	0.242	-5.2	80	0.00
43	C	2,4,6-Trichlorophenol	0.386	0.387	-0.3	78	0.00
44		2,4,5-Trichlorophenol	0.406	0.400	1.5	76	0.00
45	S	2-Fluorobiphenyl	1.359	1.350	0.7	80	0.00
46		1,1'-Biphenyl	1.585	1.544	2.6	77	0.00
47		2-Chloronaphthalene	1.175	1.153	1.9	77	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022122\
Data File : BF126973.D
Acq On : 21 Feb 2022 15:00
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Feb 21 17:01:54 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 21 13:16:57 2022
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	0.464	0.456	1.7	76	0.00
49		Acenaphthylene	1.909	1.886	1.2	77	0.00
50		Dimethylphthalate	1.430	1.396	2.4	76	0.00
51		2,6-Dinitrotoluene	0.291	0.293	-0.7	77	0.00
52	C	Acenaphthene	1.241	1.245	-0.3	78	0.00
53		3-Nitroaniline	0.356	0.357	-0.3	77	0.00
54	P	2,4-Dinitrophenol	0.154	0.174	-13.0	81	0.00
55		Dibenzofuran	1.707	1.700	0.4	78	0.00
56	P	4-Nitrophenol	0.263	0.279	-6.1	79	0.00
57		2,4-Dinitrotoluene	0.393	0.399	-1.5	77	0.00
58		Fluorene	1.330	1.324	0.5	79	0.00
59		2,3,4,6-Tetrachlorophenol	0.322	0.332	-3.1	79	0.00
60		Diethylphthalate	1.494	1.498	-0.3	78	0.00
61		4-Chlorophenyl-phenylether	0.627	0.616	1.8	77	0.00
62		4-Nitroaniline	0.351	0.360	-2.6	77	0.00
63		Azobenzene	1.663	1.616	2.8	76	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	77	0.00
65		4,6-Dinitro-2-methylphenol	0.118	0.129	-9.3	80	0.00
66	c	n-Nitrosodiphenylamine	0.656	0.638	2.7	78	0.00
67		4-Bromophenyl-phenylether	0.223	0.225	-0.9	80	0.00
68		Hexachlorobenzene	0.241	0.238	1.2	80	0.00
69		Atrazine	0.203	0.204	-0.5	81	0.00
70	C	Pentachlorophenol	0.131	0.141	-7.6	79	0.00
71		Phenanthrene	1.119	1.101	1.6	80	0.00
72		Anthracene	1.114	1.107	0.6	80	0.00
73		Carbazole	1.022	1.036	-1.4	82	0.00
74		Di-n-butylphthalate	1.314	1.299	1.1	79	0.00
75	C	Fluoranthene	1.063	1.110	-4.4	85	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	94	0.00
77		Benzidine	0.544	0.590	-8.5	100	0.00
78		Pyrene	1.738	1.578	9.2	86	0.00
79	S	Terphenyl-d14	1.361	1.261	7.3	90	0.00
80		Butylbenzylphthalate	0.823	0.766	6.9	87	0.00
81		Benzo(a)anthracene	1.348	1.327	1.6	95	0.00
82		3,3'-Dichlorobenzidine	0.425	0.438	-3.1	97	0.00
83		Chrysene	1.268	1.254	1.1	95	0.00
84		Bis(2-ethylhexyl)phthalate	1.080	1.020	5.6	90	0.00
85	c	Di-n-octyl phthalate	1.555	1.512	2.8	95	0.00
86	I	Perylene-d12	1.000	1.000	0.0	91	0.00
87		Indeno(1,2,3-cd)pyrene	1.311	1.228	6.3	82	0.00
88		Benzo(b)fluoranthene	1.343	1.366	-1.7	95	0.00
89		Benzo(k)fluoranthene	1.296	1.327	-2.4	93	0.00
90	C	Benzo(a)pyrene	1.049	1.039	1.0	90	0.00
91		Dibenzo(a,h)anthracene	1.085	1.001	7.7	81	0.00
92		Benzo(g,h,i)perylene	1.069	0.996	6.8	81	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022122\
Data File : BF126973.D
Acq On : 21 Feb 2022 15:00
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Feb 21 17:01:54 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 21 13:16:57 2022
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