Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082118\
Data File : BF108343.D
Acq On : 21 Aug 2018 17:54
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 21 23:44:19 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 21 15:20:21 2018
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	79	0.00
2		1,4-Dioxane	0.568	0.498	12.3	69	0.04
3		Pyridine	1.462	1.298	11.2	69	0.02
4		n-Nitrosodimethylamine	0.823	0.715	13.1	67	0.02
5	S	2-Fluorophenol	1.105	1.046	5.3	74	0.00
6		Aniline	1.997	1.909	4.4	76	0.00
7	S	Phenol-d6	1.468	1.407	4.2	76	0.00
8		2-Chlorophenol	1.244	1.195	3.9	75	0.00
9		Benzaldehyde	1.138	1.013	11.0	69	0.00
10	C	Phenol	1.518	1.460	3.8	77	0.00
11		bis(2-Chloroethyl)ether	1.228	1.153	6.1	73	0.00
12		1,3-Dichlorobenzene	1.439	1.355	5.8	74	0.00
13	C	1,4-Dichlorobenzene	1.445	1.367	5.4	74	0.00
14		1,2-Dichlorobenzene	1.344	1.264	6.0	74	0.00
15		Benzyl Alcohol	1.236	1.138	7.9	71	0.00
16		2,2'-oxybis(1-Chloropropane	2.529	2.174	14.0	67	0.00
17		2-Methylphenol	1.067	1.027	3.7	74	0.00
18		Hexachloroethane	0.515	0.493	4.3	74	0.00
19	P	n-Nitroso-di-n-propylamine	0.993	0.885	10.9	70	0.00
20		3+4-Methylphenols	1.367	1.271	7.0	72	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	78	0.00
22		Acetophenone	0.468	0.429	8.3	71	0.00
23	S	Nitrobenzene-d5	0.358	0.345	3.6	74	0.00
24		Nitrobenzene	0.362	0.341	5.8	71	0.00
25		Isophorone	0.683	0.634	7.2	72	0.00
26	C	2-Nitrophenol	0.137	0.152	-10.9	84	0.00
27		2,4-Dimethylphenol	0.303	0.288	5.0	73	0.00
28		bis(2-Chloroethoxy)methane	0.399	0.372	6.8	72	0.00
29	C	2,4-Dichlorophenol	0.279	0.272	2.5	74	0.00
30		1,2,4-Trichlorobenzene	0.308	0.292	5.2	74	0.00
31		Naphthalene	0.910	0.863	5.2	74	0.00
32		Benzoic acid	0.161	0.113	29.8#	53	0.02
33		4-Chloroaniline	0.396	0.375	5.3	73	0.00
34	C	Hexachlorobutadiene	0.195	0.186	4.6	74	0.00
35		Caprolactam	0.087	0.085	2.3	72	0.00
36	C	4-Chloro-3-methylphenol	0.301	0.284	5.6	72	0.00
37		2-Methylnaphthalene	0.644	0.597	7.3	72	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	75	0.00
39		1,2,4,5-Tetrachlorobenzene	0.614	0.602	2.0	73	0.00
40	P	Hexachlorocyclopentadiene	0.397	0.381	4.0	67	0.00
41	S	2,4,6-Tribromophenol	0.199	0.206	-3.5	73	0.00
42	C	2,4,6-Trichlorophenol	0.381	0.395	-3.7	74	0.00
43		2,4,5-Trichlorophenol	0.420	0.424	-1.0	72	0.00
44	S	2-Fluorobiphenyl	1.246	1.218	2.2	75	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082118\
Data File : BF108343.D
Acq On : 21 Aug 2018 17:54
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 21 23:44:19 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 21 15:20:21 2018
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		1,1'-Biphenyl	1.475	1.413	4.2	72	0.00
46		2-Chloronaphthalene	1.165	1.116	4.2	72	0.00
47		2-Nitroaniline	0.377	0.381	-1.1	71	0.00
48		Acenaphthylene	1.843	1.764	4.3	72	0.00
49		Dimethylphthalate	1.393	1.318	5.4	71	0.00
50		2,6-Dinitrotoluene	0.291	0.294	-1.0	72	0.00
51	C	Acenaphthene	1.129	1.082	4.2	72	0.00
52		3-Nitroaniline	0.319	0.317	0.6	71	0.00
53	P	2,4-Dinitrophenol	0.092	0.092	0.0	73	0.00
54		Dibenzofuran	1.635	1.558	4.7	72	0.00
55	P	4-Nitrophenol	0.241	0.217	10.0	65	0.00
56		2,4-Dinitrotoluene	0.375	0.387	-3.2	72	0.00
57		Fluorene	1.294	1.240	4.2	73	0.00
58		2,3,4,6-Tetrachlorophenol	0.351	0.346	1.4	69	0.00
59		Diethylphthalate	1.349	1.265	6.2	70	0.00
60		4-Chlorophenyl-phenylether	0.674	0.635	5.8	70	0.00
61		4-Nitroaniline	0.315	0.310	1.6	70	0.01
62		Azobenzene	1.393	1.291	7.3	69	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	73	0.00
64		4,6-Dinitro-2-methylphenol	0.069	0.071	-2.9	72	0.00
65	c	n-Nitrosodiphenylamine	0.591	0.571	3.4	71	0.00
66		4-Bromophenyl-phenylether	0.217	0.211	2.8	71	0.00
67		Hexachlorobenzene	0.229	0.219	4.4	70	0.00
68		Atrazine	0.198	0.191	3.5	69	0.00
69	C	Pentachlorophenol	0.109	0.096	11.9	62	0.00
70		Phenanthrene	0.943	0.891	5.5	71	0.00
71		Anthracene	0.967	0.925	4.3	71	0.00
72		Carbazole	0.859	0.786	8.5	67	0.00
73		Di-n-butylphthalate	0.973	0.966	0.7	72	0.00
74	C	Fluoranthene	1.030	0.944	8.3	68	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	56	0.00
76		Benzidine	0.747	0.696	6.8	50#	0.00
77		Pyrene	1.266	1.486	-17.4	67	0.00
78	S	Terphenyl-d14	0.787	0.925	-17.5	68	0.00
79		Butylbenzylphthalate	0.503	0.588	-16.9	62	0.00
80		Benzo(a)anthracene	1.148	1.115	2.9	55	0.00
81		3,3'-Dichlorobenzidine	0.411	0.335	18.5	43#	0.00
82		Chrysene	1.052	0.999	5.0	54	0.00
83		Bis(2-ethylhexyl)phthalate	0.681	0.760	-11.6	60	0.00
84	c	Di-n-octyl phthalate	1.038	1.070	-3.1	54	0.00
85		Indeno(1,2,3-cd)pyrene	0.912	1.012	-11.0	63	0.00
86	I	Perylene-d12	1.000	1.000	0.0	54	0.00
87		Benzo(b)fluoranthene	1.195	1.124	5.9	48#	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082118\
Data File : BF108343.D
Acq On : 21 Aug 2018 17:54
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 21 23:44:19 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 21 15:20:21 2018
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(k)fluoranthene	1.099	0.957	12.9	48#	0.00
89 C	Benzo(a)pyrene	1.070	1.015	5.1	50	0.00
90	Dibenzo(a,h)anthracene	0.885	1.063	-20.1	65	0.00
91	Benzo(g,h,i)perylene	0.872	1.073	-23.1	67	0.01

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