Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF082018\
Data File : BF108277.D
Acq On : 20 Aug 2018 9:32
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 20 11:07:37 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 08 12:24:24 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	84	0.00
2		1,4-Dioxane	0.568	0.506	10.9	75	0.01
3		Pyridine	1.462	1.345	8.0	76	0.02
4		n-Nitrosodimethylamine	0.823	0.702	14.7	70	0.02
5	S	2-Fluorophenol	1.105	1.072	3.0	82	0.02
6		Aniline	1.997	1.976	1.1	84	0.01
7	S	Phenol-d6	1.468	1.437	2.1	83	0.02
8		2-Chlorophenol	1.244	1.235	0.7	83	0.00
9		Benzaldehyde	1.138	1.026	9.8	75	0.00
10	C	Phenol	1.518	1.495	1.5	84	0.02
11		bis(2-Chloroethyl)ether	1.228	1.201	2.2	82	0.00
12		1,3-Dichlorobenzene	1.439	1.408	2.2	83	0.00
13	C	1,4-Dichlorobenzene	1.445	1.412	2.3	81	0.00
14		1,2-Dichlorobenzene	1.344	1.315	2.2	83	0.00
15		Benzyl Alcohol	1.236	1.198	3.1	80	0.01
16		2,2'-oxybis(1-Chloropropane	2.529	2.209	12.7	73	0.00
17		2-Methylphenol	1.067	1.040	2.5	80	0.02
18		Hexachloroethane	0.515	0.516	-0.2	82	0.00
19	P	n-Nitroso-di-n-propylamine	0.993	0.928	6.5	78	0.00
20		3+4-Methylphenols	1.367	1.300	4.9	79	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	85	0.00
22		Acetophenone	0.468	0.441	5.8	80	0.00
23	S	Nitrobenzene-d5	0.358	0.348	2.8	81	0.01
24		Nitrobenzene	0.362	0.346	4.4	79	0.00
25		Isophorone	0.683	0.642	6.0	80	0.00
26	C	2-Nitrophenol	0.137	0.150	-9.5	91	0.00
27		2,4-Dimethylphenol	0.303	0.292	3.6	81	0.01
28		bis(2-Chloroethoxy)methane	0.399	0.382	4.3	82	0.00
29	C	2,4-Dichlorophenol	0.279	0.277	0.7	82	0.01
30		1,2,4-Trichlorobenzene	0.308	0.294	4.5	82	0.00
31		Naphthalene	0.910	0.882	3.1	83	0.00
32		Benzoic acid	0.161	0.139	13.7	71	0.03
33		4-Chloroaniline	0.396	0.388	2.0	82	0.00
34	C	Hexachlorobutadiene	0.195	0.191	2.1	83	0.00
35		Caprolactam	0.087	0.086	1.1	80	0.02
36	C	4-Chloro-3-methylphenol	0.301	0.290	3.7	80	0.01
37		2-Methylnaphthalene	0.644	0.615	4.5	81	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	83	0.00
39		1,2,4,5-Tetrachlorobenzene	0.614	0.611	0.5	82	0.00
40	P	Hexachlorocyclopentadiene	0.397	0.322	18.9	63	0.00
41	S	2,4,6-Tribromophenol	0.199	0.210	-5.5	83	0.00
42	C	2,4,6-Trichlorophenol	0.381	0.397	-4.2	83	0.00
43		2,4,5-Trichlorophenol	0.420	0.433	-3.1	81	0.01
44	S	2-Fluorobiphenyl	1.246	1.208	3.0	83	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF082018\
Data File : BF108277.D
Acq On : 20 Aug 2018 9:32
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 20 11:07:37 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 08 12:24:24 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	1.435	2.7	81	0.00
46		2-Chloronaphthalene	1.165	1.137	2.4	82	0.00
47		2-Nitroaniline	0.377	0.385	-2.1	80	0.00
48		Acenaphthylene	1.843	1.806	2.0	82	0.00
49		Dimethylphthalate	1.393	1.357	2.6	82	0.00
50		2,6-Dinitrotoluene	0.291	0.299	-2.7	82	0.00
51	C	Acenaphthene	1.129	1.106	2.0	82	0.00
52		3-Nitroaniline	0.319	0.325	-1.9	82	0.01
53	P	2,4-Dinitrophenol	0.092	0.110	-19.6	97	0.02
54		Dibenzofuran	1.635	1.591	2.7	82	0.00
55	P	4-Nitrophenol	0.241	0.235	2.5	79	0.02
56		2,4-Dinitrotoluene	0.375	0.401	-6.9	83	0.00
57		Fluorene	1.294	1.271	1.8	83	0.00
58		2,3,4,6-Tetrachlorophenol	0.351	0.364	-3.7	81	0.00
59		Diethylphthalate	1.349	1.301	3.6	80	0.00
60		4-Chlorophenyl-phenylether	0.674	0.655	2.8	81	0.00
61		4-Nitroaniline	0.315	0.337	-7.0	85	0.00
62		Azobenzene	1.393	1.318	5.4	78	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	82	0.00
64		4,6-Dinitro-2-methylphenol	0.069	0.075	-8.7	85	0.02
65	c	n-Nitrosodiphenylamine	0.591	0.580	1.9	80	0.00
66		4-Bromophenyl-phenylether	0.217	0.216	0.5	82	0.00
67		Hexachlorobenzene	0.229	0.226	1.3	81	0.00
68		Atrazine	0.198	0.196	1.0	80	0.00
69	C	Pentachlorophenol	0.109	0.104	4.6	75	0.01
70		Phenanthrene	0.943	0.921	2.3	82	0.00
71		Anthracene	0.967	0.945	2.3	81	0.00
72		Carbazole	0.859	0.835	2.8	80	0.00
73		Di-n-butylphthalate	0.973	0.981	-0.8	82	0.00
74	C	Fluoranthene	1.030	0.972	5.6	79	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	67	0.00
76		Benzidine	0.747	0.843	-12.9	72	0.00
77		Pyrene	1.266	1.462	-15.5	79	0.00
78	S	Terphenyl-d14	0.787	0.885	-12.5	78	0.00
79		Butylbenzylphthalate	0.503	0.583	-15.9	73	0.00
80		Benzo(a)anthracene	1.148	1.152	-0.3	68	0.00
81		3,3'-Dichlorobenzidine	0.411	0.401	2.4	62	0.00
82		Chrysene	1.052	1.047	0.5	67	0.01
83		Bis(2-ethylhexyl)phthalate	0.681	0.703	-3.2	67	0.00
84	c	Di-n-octyl phthalate	1.038	1.244	-19.8	76	0.00
85		Indeno(1,2,3-cd)pyrene	0.912	0.961	-5.4	72	0.03
86	I	Perylene-d12	1.000	1.000	0.0	65	0.01
87		Benzo(b)fluoranthene	1.195	1.129	5.5	58	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF082018\
Data File : BF108277.D
Acq On : 20 Aug 2018 9:32
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 20 11:07:37 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 08 12:24:24 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	1.089	0.9	66	0.00
89 C	Benzo(a)pyrene	1.070	1.059	1.0	63	0.01
90	Dibenzo(a,h)anthracene	0.885	0.991	-12.0	72	0.02
91	Benzo(g,h,i)perylene	0.872	1.007	-15.5	76	0.04

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