Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA F\Data\BF090717\
Data File : BF098312.D
Acq On : 7 Sep 2017 14:26
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 07 17:21:02 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 05 17:28:20 2017
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	-0.02
2		1,4-Dioxane	0.733	0.723	1.4	84	0.00
3		Pyridine	2.027	2.004	1.1	84	0.00
4		n-Nitrosodimethylamine	0.780	0.708	9.2	74	-0.01
5	S	2-Fluorophenol	1.315	1.305	0.8	85	-0.02
6		Aniline	2.510	2.385	5.0	81	-0.02
7	S	Phenol-d6	1.787	1.799	-0.7	86	-0.02
8		2-Chlorophenol	1.387	1.420	-2.4	85	-0.02
9		Benzaldehyde	1.132	1.005	11.2	74	-0.02
10	C	Phenol	2.089	2.040	2.3	79	-0.02
11		bis(2-Chloroethyl)ether	1.577	1.589	-0.8	85	-0.02
12		1,3-Dichlorobenzene	1.492	1.512	-1.3	86	-0.02
13	C	1,4-Dichlorobenzene	1.517	1.527	-0.7	86	-0.02
14		1,2-Dichlorobenzene	1.399	1.412	-0.9	86	-0.02
15		Benzyl Alcohol	1.338	1.207	9.8	75	-0.01
16		2,2'-oxybis(1-Chloropropane	2.122	1.981	6.6	79	-0.02
17		2-Methylphenol	1.222	1.226	-0.3	84	-0.01
18		Hexachloroethane	0.595	0.599	-0.7	84	-0.02
19	P	n-Nitroso-di-n-propylamine	1.223	1.127	7.8	78	-0.01
20		3+4-Methylphenols	1.493	1.476	1.1	84	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	85	-0.01
22		Acetophenone	0.528	0.503	4.7	83	-0.02
23	S	Nitrobenzene-d5	0.368	0.406	-10.3	90	-0.01
24		Nitrobenzene	0.410	0.441	-7.6	85	-0.01
25		Isophorone	0.766	0.749	2.2	83	-0.01
26	C	2-Nitrophenol	0.137	0.187	-36.5#	110	-0.02
27		2,4-Dimethylphenol	0.319	0.327	-2.5	88	-0.01
28		bis(2-Chloroethoxy)methane	0.503	0.506	-0.6	85	-0.02
29	C	2,4-Dichlorophenol	0.265	0.279	-5.3	87	-0.01
30		1,2,4-Trichlorobenzene	0.294	0.296	-0.7	86	-0.02
31		Naphthalene	1.038	1.037	0.1	86	-0.01
32		Benzoic acid	0.212	0.174	17.9	70	-0.02
33		4-Chloroaniline	0.438	0.445	-1.6	87	-0.01
34	C	Hexachlorobutadiene	0.172	0.177	-2.9	88	-0.02
35		Caprolactam	0.090	0.098	-8.9	88	-0.02
36	C	4-Chloro-3-methylphenol	0.341	0.346	-1.5	84	-0.02
37		2-Methylnaphthalene	0.637	0.640	-0.5	86	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	86	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.614	0.620	-1.0	88	-0.02
40	P	Hexachlorocyclopentadiene	0.295	0.278	5.8	77	-0.01
41	S	2,4,6-Tribromophenol	0.174	0.186	-6.9	89	-0.02
42	C	2,4,6-Trichlorophenol	0.412	0.411	0.2	84	-0.01
43		2,4,5-Trichlorophenol	0.409	0.424	-3.7	87	-0.02
44	S	2-Fluorobiphenyl	1.361	1.308	3.9	86	-0.02

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA F\Data\BF090717\
Data File : BF098312.D
Acq On : 7 Sep 2017 14:26
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 07 17:21:02 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 05 17:28:20 2017
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	1.824	1.811	0.7	87	-0.02
46		2-Chloronaphthalene	1.348	1.331	1.3	87	-0.02
47		2-Nitroaniline	0.452	0.515	-13.9	94	-0.02
48		Acenaphthylene	2.116	2.067	2.3	84	-0.01
49		Dimethylphthalate	1.462	1.502	-2.7	89	-0.02
50		2,6-Dinitrotoluene	0.292	0.321	-9.9	91	-0.02
51	C	Acenaphthene	1.335	1.256	5.9	82	-0.01
52		3-Nitroaniline	0.376	0.417	-10.9	94	-0.01
53	P	2,4-Dinitrophenol	0.106	0.132	-24.5	105	-0.02
54		Dibenzofuran	1.789	1.700	5.0	83	-0.01
55	P	4-Nitrophenol	0.300	0.246	18.0	68	-0.02
56		2,4-Dinitrotoluene	0.366	0.393	-7.4	88	-0.01
57		Fluorene	1.383	1.372	0.8	88	-0.02
58		2,3,4,6-Tetrachlorophenol	0.317	0.306	3.5	82	-0.01
59		Diethylphthalate	1.529	1.484	2.9	83	-0.02
60		4-Chlorophenyl-phenylether	0.642	0.633	1.4	86	-0.01
61		4-Nitroaniline	0.358	0.406	-13.4	95	-0.02
62		Azobenzene	1.816	1.677	7.7	80	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	86	-0.01
64		4,6-Dinitro-2-methylphenol	0.083	0.114	-37.3#	116	-0.01
65	c	n-Nitrosodiphenylamine	0.681	0.675	0.9	86	-0.01
66		4-Bromophenyl-phenylether	0.208	0.213	-2.4	88	-0.01
67		Hexachlorobenzene	0.212	0.218	-2.8	89	-0.02
68		Atrazine	0.181	0.172	5.0	83	-0.01
69	C	Pentachlorophenol	0.123	0.114	7.3	75	-0.02
70		Phenanthrene	1.066	1.023	4.0	84	-0.02
71		Anthracene	1.081	1.061	1.9	86	-0.02
72		Carbazole	1.026	1.000	2.5	86	-0.01
73		Di-n-butylphthalate	1.182	1.232	-4.2	89	-0.02
74	C	Fluoranthene	1.112	1.104	0.7	87	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	91	-0.02
76		Benzidine	0.656	0.663	-1.1	96	-0.01
77		Pyrene	1.636	1.570	4.0	87	-0.02
78	S	Terphenyl-d14	0.959	0.922	3.9	88	-0.02
79		Butylbenzylphthalate	0.627	0.669	-6.7	93	-0.01
80		Benzo(a)anthracene	1.199	1.128	5.9	86	-0.02
81		3,3'-Dichlorobenzidine	0.404	0.413	-2.2	88	-0.02
82		Chrysene	1.192	1.121	6.0	86	-0.02
83		Bis(2-ethylhexyl)phthalate	0.695	0.823	-18.4	99	-0.01
84	c	Di-n-octyl phthalate	1.209	1.494	-23.6#	105	-0.01
85		Indeno(1,2,3-cd)pyrene	0.877	0.935	-6.6	99	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	92	-0.02
87		Benzo(b)fluoranthene	1.261	1.218	3.4	88	-0.02

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA F\Data\BF090717\
Data File : BF098312.D
Acq On : 7 Sep 2017 14:26
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 07 17:21:02 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 05 17:28:20 2017
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	1.184	1.214	-2.5	92	-0.02
89 C	Benzo(a)pyrene	1.116	1.136	-1.8	91	-0.02
90	Dibenzo(a,h)anthracene	0.909	1.032	-13.5	101	-0.03
91	Benzo(g,h,i)perylene	0.948	1.047	-10.4	99	-0.03

(#) = Out of Range SPCC's out = 0 CCC's out = 2