Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF102518\
Data File : BF110163.D
Acq On : 25 Oct 2018 16:27
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 25 18:08:08 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 23 15:06:18 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	86	-0.01
2		1,4-Dioxane	0.643	0.612	4.8	83	-0.02
3		Pyridine	1.433	1.400	2.3	81	-0.02
4		n-Nitrosodimethylamine	0.600	0.586	2.3	82	-0.02
5	S	2-Fluorophenol	1.228	1.201	2.2	84	-0.01
6		Aniline	2.046	2.002	2.2	84	-0.01
7	S	Phenol-d6	1.571	1.509	3.9	83	-0.01
8		2-Chlorophenol	1.416	1.398	1.3	84	-0.02
9		Benzaldehyde	0.954	0.878	8.0	80	-0.01
10	C	Phenol	1.679	1.636	2.6	84	-0.01
11		bis(2-Chloroethyl)ether	1.381	1.332	3.5	84	-0.01
12		1,3-Dichlorobenzene	1.558	1.515	2.8	84	-0.01
13	C	1,4-Dichlorobenzene	1.576	1.545	2.0	85	-0.01
14		1,2-Dichlorobenzene	1.463	1.430	2.3	84	-0.02
15		Benzyl Alcohol	1.208	1.221	-1.1	85	-0.02
16		2,2'-oxybis(1-Chloropropane	2.209	2.159	2.3	84	-0.02
17		2-Methylphenol	1.128	1.109	1.7	84	-0.01
18		Hexachloroethane	0.577	0.570	1.2	86	-0.02
19	P	n-Nitroso-di-n-propylamine	1.008	0.972	3.6	84	-0.01
20		3+4-Methylphenols	1.459	1.428	2.1	84	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	86	-0.02
22		Acetophenone	0.461	0.440	4.6	83	-0.01
23	S	Nitrobenzene-d5	0.337	0.335	0.6	84	-0.02
24		Nitrobenzene	0.348	0.346	0.6	85	-0.02
25		Isophorone	0.575	0.549	4.5	82	-0.02
26	C	2-Nitrophenol	0.166	0.175	-5.4	87	-0.01
27		2,4-Dimethylphenol	0.275	0.267	2.9	84	-0.02
28		bis(2-Chloroethoxy)methane	0.390	0.375	3.8	84	-0.02
29	C	2,4-Dichlorophenol	0.277	0.272	1.8	83	-0.01
30		1,2,4-Trichlorobenzene	0.311	0.307	1.3	85	-0.01
31		Naphthalene	0.922	0.893	3.1	85	-0.02
32		Benzoic acid	0.212	0.213	-0.5	84	0.01
33		4-Chloroaniline	0.415	0.395	4.8	83	-0.01
34	C	Hexachlorobutadiene	0.173	0.171	1.2	87	-0.01
35		Caprolactam	0.097	0.092	5.2	79	0.00
36	C	4-Chloro-3-methylphenol	0.300	0.289	3.7	82	0.00
37		2-Methylnaphthalene	0.610	0.584	4.3	83	-0.02
38		1-Methylnaphthalene	0.589	0.561	4.8	83	-0.02
39	I	Acenaphthene-d10	1.000	1.000	0.0	87	-0.01
40		1,2,4,5-Tetrachlorobenzene	0.623	0.596	4.3	84	-0.01
41	P	Hexachlorocyclopentadiene	0.319	0.310	2.8	79	-0.01
42	S	2,4,6-Tribromophenol	0.198	0.191	3.5	84	-0.01
43	C	2,4,6-Trichlorophenol	0.402	0.393	2.2	83	-0.01
44		2,4,5-Trichlorophenol	0.425	0.408	4.0	83	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF102518\
Data File : BF110163.D
Acq On : 25 Oct 2018 16:27
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 25 18:08:08 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 23 15:06:18 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	S	2-Fluorobiphenyl	1.274	1.239	2.7	85	-0.01
46		1,1'-Biphenyl	1.809	1.722	4.8	85	-0.01
47		2-Chloronaphthalene	1.327	1.276	3.8	84	-0.02
48		2-Nitroaniline	0.402	0.396	1.5	84	-0.01
49		Acenaphthylene	1.916	1.815	5.3	83	-0.02
50		Dimethylphthalate	1.476	1.394	5.6	83	-0.01
51		2,6-Dinitrotoluene	0.318	0.321	-0.9	84	-0.01
52	C	Acenaphthene	1.268	1.221	3.7	85	-0.02
53		3-Nitroaniline	0.378	0.367	2.9	81	0.00
54	P	2,4-Dinitrophenol	0.112	0.115	-2.7	88	-0.01
55		Dibenzofuran	1.775	1.672	5.8	82	-0.01
56	P	4-Nitrophenol	0.280	0.278	0.7	80	0.00
57		2,4-Dinitrotoluene	0.404	0.413	-2.2	84	-0.01
58		Fluorene	1.321	1.232	6.7	82	-0.01
59		2,3,4,6-Tetrachlorophenol	0.346	0.334	3.5	82	-0.01
60		Diethylphthalate	1.441	1.372	4.8	83	-0.01
61		4-Chlorophenyl-phenylether	0.697	0.659	5.5	84	-0.01
62		4-Nitroaniline	0.376	0.366	2.7	83	-0.01
63		Azobenzene	1.431	1.344	6.1	83	-0.02
64	I	Phenanthrene-d10	1.000	1.000	0.0	85	-0.01
65		4,6-Dinitro-2-methylphenol	0.091	0.091	0.0	82	-0.01
66	c	n-Nitrosodiphenylamine	0.656	0.637	2.9	83	-0.02
67		4-Bromophenyl-phenylether	0.217	0.211	2.8	83	-0.01
68		Hexachlorobenzene	0.230	0.223	3.0	83	-0.02
69		Atrazine	0.207	0.199	3.9	81	-0.01
70	C	Pentachlorophenol	0.134	0.140	-4.5	80	-0.01
71		Phenanthrene	1.046	0.986	5.7	82	-0.02
72		Anthracene	1.049	1.007	4.0	83	-0.02
73		Carbazole	1.024	0.975	4.8	82	-0.01
74		Di-n-butylphthalate	1.157	1.113	3.8	82	-0.02
75	C	Fluoranthene	1.062	0.985	7.3	80	-0.01
76	I	Chrysene-d12	1.000	1.000	0.0	78	-0.01
77		Benzidine	0.681	0.653	4.1	84	-0.01
78		Pyrene	1.434	1.456	-1.5	80	-0.02
79	S	Terphenyl-d14	0.962	0.968	-0.6	81	-0.01
80		Butylbenzylphthalate	0.622	0.621	0.2	77	-0.01
81		Benzo(a)anthracene	1.186	1.140	3.9	75	-0.01
82		3,3'-Dichlorobenzidine	0.413	0.392	5.1	73	-0.01
83		Chrysene	1.127	1.055	6.4	74	-0.01
84		Bis(2-ethylhexyl)phthalate	0.841	0.789	6.2	72	-0.02
85	c	Di-n-octyl phthalate	1.308	1.162	11.2	69	-0.02
86		Indeno(1,2,3-cd)pyrene	0.935	0.833	10.9	77	-0.02
87	I	Perylene-d12	1.000	1.000	0.0	71	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF102518\
Data File : BF110163.D
Acq On : 25 Oct 2018 16:27
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 25 18:08:08 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 23 15:06:18 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(b)fluoranthene	1.155	1.122	2.9	68	-0.01
89	Benzo(k)fluoranthene	1.135	1.151	-1.4	72	-0.02
90 C	Benzo(a)pyrene	1.063	1.043	1.9	70	-0.02
91	Dibenzo(a,h)anthracene	0.917	0.958	-4.5	76	-0.02
92	Benzo(g,h,i)perylene	0.904	0.963	-6.5	78	-0.02

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