Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF080218\
Data File : BF107906.D
Acq On : 2 Aug 2018 13:46
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 02 14:41:20 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 30 17:48:53 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	94	0.00
2		1,4-Dioxane	0.515	0.493	4.3	85	-0.01
3		Pyridine	1.258	1.082	14.0	78	0.01
4		n-Nitrosodimethylamine	0.599	0.318	46.9#	52	0.01
5	S	2-Fluorophenol	1.177	1.105	6.1	82	0.01
6		Aniline	1.613	1.503	6.8	82	0.00
7	S	Phenol-d6	1.551	1.425	8.1	89	0.00
8		2-Chlorophenol	1.478	1.377	6.8	89	0.00
9		Benzaldehyde	0.936	0.904	3.4	92	0.00
10	C	Phenol	1.850	1.784	3.6	105	0.00
11		bis(2-Chloroethyl)ether	1.645	1.643	0.1	96	0.00
12		1,3-Dichlorobenzene	1.550	1.444	6.8	87	-0.01
13	C	1,4-Dichlorobenzene	1.726	1.700	1.5	90	0.00
14		1,2-Dichlorobenzene	1.602	1.588	0.9	96	0.00
15		Benzyl Alcohol	0.964	0.902	6.4	84	0.00
16		2,2'-oxybis(1-Chloropropane	2.365	2.235	5.5	90	-0.01
17		2-Methylphenol	1.091	1.048	3.9	87	0.00
18		Hexachloroethane	0.620	0.574	7.4	90	0.00
19	P	n-Nitroso-di-n-propylamine	0.984	0.983	0.1	92	0.00
20		3+4-Methylphenols	1.340	1.254	6.4	80	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	93	0.00
22		Acetophenone	0.481	0.454	5.6	91	0.00
23	S	Nitrobenzene-d5	0.347	0.332	4.3	92	0.00
24		Nitrobenzene	0.364	0.339	6.9	89	0.00
25		Isophorone	0.570	0.548	3.9	88	0.00
26	C	2-Nitrophenol	0.183	0.186	-1.6	90	0.00
27		2,4-Dimethylphenol	0.245	0.232	5.3	89	0.00
28		bis(2-Chloroethoxy)methane	0.376	0.371	1.3	88	0.00
29	C	2,4-Dichlorophenol	0.277	0.273	1.4	89	0.00
30		1,2,4-Trichlorobenzene	0.316	0.303	4.1	88	0.00
31		Naphthalene	0.930	0.919	1.2	88	0.00
32		Benzoic acid	0.157	0.144	8.3	76	0.00
33		4-Chloroaniline	0.397	0.381	4.0	88	0.00
34	C	Hexachlorobutadiene	0.195	0.178	8.7	86	-0.01
35		Caprolactam	0.083	0.072	13.3	79	-0.01
36	C	4-Chloro-3-methylphenol	0.300	0.285	5.0	84	0.00
37		2-Methylnaphthalene	0.584	0.570	2.4	86	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	90	0.00
39		1,2,4,5-Tetrachlorobenzene	0.682	0.677	0.7	86	-0.01
40	P	Hexachlorocyclopentadiene	0.243	0.221	9.1	76	0.00
41	S	2,4,6-Tribromophenol	0.272	0.248	8.8	80	0.00
42	C	2,4,6-Trichlorophenol	0.377	0.379	-0.5	83	0.00
43		2,4,5-Trichlorophenol	0.454	0.455	-0.2	83	0.00
44	S	2-Fluorobiphenyl	1.435	1.404	2.2	89	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF080218\
Data File : BF107906.D
Acq On : 2 Aug 2018 13:46
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 02 14:41:20 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 30 17:48:53 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	1.820	1.811	0.5	87	-0.01
46		2-Chloronaphthalene	1.373	1.353	1.5	87	0.00
47		2-Nitroaniline	0.446	0.429	3.8	83	0.00
48		Acenaphthylene	2.128	1.964	7.7	85	-0.01
49		Dimethylphthalate	1.502	1.399	6.9	82	0.00
50		2,6-Dinitrotoluene	0.326	0.329	-0.9	90	0.00
51	C	Acenaphthene	1.236	1.146	7.3	87	-0.01
52		3-Nitroaniline	0.411	0.387	5.8	83	0.00
53	P	2,4-Dinitrophenol	0.140	0.144	-2.9	85	0.01
54		Dibenzofuran	1.911	1.783	6.7	84	0.00
55	P	4-Nitrophenol	0.150	0.128	14.7	67	0.03
56		2,4-Dinitrotoluene	0.429	0.397	7.5	81	0.00
57		Fluorene	1.485	1.365	8.1	84	0.00
58		2,3,4,6-Tetrachlorophenol	0.417	0.432	-3.6	87	0.00
59		Diethylphthalate	1.615	1.474	8.7	83	0.00
60		4-Chlorophenyl-phenylether	0.809	0.743	8.2	84	0.00
61		4-Nitroaniline	0.365	0.344	5.8	84	0.00
62		Azobenzene	1.504	1.414	6.0	86	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	87	0.00
64		4,6-Dinitro-2-methylphenol	0.109	0.100	8.3	77	0.00
65	c	n-Nitrosodiphenylamine	0.659	0.617	6.4	83	-0.01
66		4-Bromophenyl-phenylether	0.233	0.230	1.3	89	0.00
67		Hexachlorobenzene	0.258	0.235	8.9	80	-0.01
68		Atrazine	0.195	0.178	8.7	79	0.00
69	C	Pentachlorophenol	0.127	0.121	4.7	78	0.00
70		Phenanthrene	0.950	0.883	7.1	81	0.00
71		Anthracene	1.095	1.056	3.6	86	0.00
72		Carbazole	1.082	1.045	3.4	84	0.00
73		Di-n-butylphthalate	1.181	1.071	9.3	80	0.00
74	C	Fluoranthene	1.116	1.002	10.2	80	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	72	0.00
76		Benzidine	0.647	0.540	16.5	57	0.00
77		Pyrene	1.440	1.555	-8.0	79	0.00
78	S	Terphenyl-d14	0.922	1.010	-9.5	82	0.00
79		Butylbenzylphthalate	0.631	0.615	2.5	69	0.00
80		Benzo(a)anthracene	1.133	1.009	10.9	63	0.00
81		3,3'-Dichlorobenzidine	0.439	0.415	5.5	68	0.00
82		Chrysene	1.222	1.205	1.4	72	-0.01
83		Bis(2-ethylhexyl)phthalate	0.793	0.731	7.8	67	-0.01
84	c	Di-n-octyl phthalate	1.273	1.113	12.6	62	-0.01
85		Indeno(1,2,3-cd)pyrene	0.929	1.048	-12.8	87	0.00
86	I	Perylene-d12	1.000	1.000	0.0	73	0.00
87		Benzo(b)fluoranthene	0.986	0.854	13.4	64	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF080218\
Data File : BF107906.D
Acq On : 2 Aug 2018 13:46
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 02 14:41:20 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 30 17:48:53 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	1.211	1.085	10.4	64	0.00
89 C	Benzo(a)pyrene	1.018	0.922	9.4	68	-0.01
90	Dibenzo(a,h)anthracene	0.888	1.023	-15.2	88	0.00
91	Benzo(g,h,i)perylene	0.859	1.022	-19.0	91	0.00

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