Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF101817\
Data File : BF099607.D
Acq On : 18 Oct 2017 10:21
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 18 15:26:19 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 18 15:18:49 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	64	0.00
2		1,4-Dioxane	0.600	0.602	-0.3	63	0.00
3		Pyridine	1.624	1.477	9.1	60	0.00
4		n-Nitrosodimethylamine	0.688	0.580	15.7	54	0.00
5	S	2-Fluorophenol	1.256	1.345	-7.1	68	0.00
6		Aniline	2.092	2.072	1.0	64	0.00
7	S	Phenol-d6	1.550	1.666	-7.5	69	0.00
8		2-Chlorophenol	1.423	1.552	-9.1	71	0.00
9		Benzaldehyde	0.899	0.784	12.8	57	0.00
10	C	Phenol	1.733	1.865	-7.6	70	0.00
11		bis(2-Chloroethyl)ether	1.348	1.422	-5.5	69	0.00
12		1,3-Dichlorobenzene	1.490	1.628	-9.3	71	0.00
13	C	1,4-Dichlorobenzene	1.516	1.664	-9.8	72	0.00
14		1,2-Dichlorobenzene	1.401	1.512	-7.9	70	0.00
15		Benzyl Alcohol	1.137	1.029	9.5	58	0.00
16		2,2'-oxybis(1-Chloropropane	2.099	1.786	14.9	56	0.00
17		2-Methylphenol	1.164	1.214	-4.3	68	0.00
18		Hexachloroethane	0.545	0.560	-2.8	67	0.00
19	P	n-Nitroso-di-n-propylamine	0.990	0.914	7.7	62	0.00
20		3+4-Methylphenols	1.473	1.459	1.0	64	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	67	0.00
22		Acetophenone	0.485	0.460	5.2	65	0.00
23	S	Nitrobenzene-d5	0.312	0.318	-1.9	68	0.00
24		Nitrobenzene	0.354	0.349	1.4	66	0.00
25		Isophorone	0.645	0.630	2.3	67	0.00
26	C	2-Nitrophenol	0.170	0.204	-20.0	77	0.00
27		2,4-Dimethylphenol	0.321	0.310	3.4	66	0.00
28		bis(2-Chloroethoxy)methane	0.402	0.412	-2.5	71	0.00
29	C	2,4-Dichlorophenol	0.276	0.298	-8.0	73	0.00
30		1,2,4-Trichlorobenzene	0.276	0.293	-6.2	72	0.00
31		Naphthalene	0.939	0.973	-3.6	71	0.00
32		Benzoic acid	0.264	0.270	-2.3	66	0.00
33		4-Chloroaniline	0.392	0.415	-5.9	72	0.00
34	C	Hexachlorobutadiene	0.142	0.150	-5.6	73	0.00
35		Caprolactam	0.088	0.097	-10.2	76	0.00
36	C	4-Chloro-3-methylphenol	0.310	0.322	-3.9	70	0.00
37		2-Methylnaphthalene	0.611	0.644	-5.4	72	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	69	0.00
39		1,2,4,5-Tetrachlorobenzene	0.596	0.640	-7.4	75	0.00
40	P	Hexachlorocyclopentadiene	0.273	0.249	8.8	61	0.00
41	S	2,4,6-Tribromophenol	0.164	0.179	-9.1	73	0.00
42	C	2,4,6-Trichlorophenol	0.423	0.433	-2.4	71	0.00
43		2,4,5-Trichlorophenol	0.422	0.441	-4.5	71	0.00
44	S	2-Fluorobiphenyl	1.198	1.270	-6.0	73	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF101817\
Data File : BF099607.D
Acq On : 18 Oct 2017 10:21
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 18 15:26:19 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 18 15:18:49 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	1.719	1.808	-5.2	74	0.00
46		2-Chloronaphthalene	1.291	1.349	-4.5	73	0.00
47		2-Nitroaniline	0.395	0.387	2.0	66	0.00
48		Acenaphthylene	1.976	2.021	-2.3	72	0.00
49		Dimethylphthalate	1.509	1.631	-8.1	76	0.00
50		2,6-Dinitrotoluene	0.329	0.369	-12.2	75	0.00
51	C	Acenaphthene	1.251	1.223	2.2	69	0.00
52		3-Nitroaniline	0.383	0.421	-9.9	73	0.00
53	P	2,4-Dinitrophenol	0.148	0.146	1.4	65	0.00
54		Dibenzofuran	1.666	1.656	0.6	70	0.00
55	P	4-Nitrophenol	0.324	0.363	-12.0	75	0.00
56		2,4-Dinitrotoluene	0.426	0.442	-3.8	69	0.00
57		Fluorene	1.339	1.444	-7.8	76	0.00
58		2,3,4,6-Tetrachlorophenol	0.291	0.300	-3.1	71	0.00
59		Diethylphthalate	1.475	1.519	-3.0	72	0.00
60		4-Chlorophenyl-phenylether	0.602	0.645	-7.1	75	0.00
61		4-Nitroaniline	0.370	0.397	-7.3	72	0.00
62		Azobenzene	1.356	1.269	6.4	66	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	70	0.00
64		4,6-Dinitro-2-methylphenol	0.122	0.138	-13.1	78	0.00
65	c	n-Nitrosodiphenylamine	0.693	0.709	-2.3	74	0.00
66		4-Bromophenyl-phenylether	0.196	0.207	-5.6	76	0.00
67		Hexachlorobenzene	0.209	0.224	-7.2	77	0.00
68		Atrazine	0.208	0.209	-0.5	71	0.00
69	C	Pentachlorophenol	0.130	0.116	10.8	60	0.00
70		Phenanthrene	1.071	1.100	-2.7	75	0.00
71		Anthracene	1.095	1.121	-2.4	74	0.00
72		Carbazole	1.073	1.082	-0.8	72	0.00
73		Di-n-butylphthalate	1.291	1.310	-1.5	72	0.00
74	C	Fluoranthene	1.084	1.113	-2.7	75	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	67	0.00
76		Benzidine	0.740	0.630	14.9	58	0.00
77		Pyrene	1.525	1.635	-7.2	73	0.00
78	S	Terphenyl-d14	0.879	0.946	-7.6	73	0.00
79		Butylbenzylphthalate	0.717	0.787	-9.8	73	0.00
80		Benzo(a)anthracene	1.211	1.269	-4.8	72	0.00
81		3,3'-Dichlorobenzidine	0.444	0.416	6.3	63	0.00
82		Chrysene	1.153	1.217	-5.6	73	0.00
83		Bis(2-ethylhexyl)phthalate	0.987	1.037	-5.1	71	0.00
84	c	Di-n-octyl phthalate	1.589	1.752	-10.3	76	0.00
85		Indeno(1,2,3-cd)pyrene	0.922	0.875	5.1	70	0.00
86	I	Perylene-d12	1.000	1.000	0.0	62	0.00
87		Benzo(b)fluoranthene	1.230	1.362	-10.7	67	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF101817\
Data File : BF099607.D
Acq On : 18 Oct 2017 10:21
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 18 15:26:19 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 18 15:18:49 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	1.215	1.334	-9.8	69	0.00
89 C	Benzo(a)pyrene	1.129	1.210	-7.2	67	0.00
90	Dibenzo(a,h)anthracene	0.903	0.990	-9.6	70	0.00
91	Benzo(g,h,i)perylene	0.893	1.047	-17.2	74	0.00

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