Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF102017\
Data File : BF099693.D
Acq On : 20 Oct 2017 11:37
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 21 23:57:21 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 20 13:11:43 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	0.590	1.7	62	0.00
3		Pyridine	1.624	1.524	6.2	62	0.00
4		n-Nitrosodimethylamine	0.688	0.551	19.9	51	0.00
5	S	2-Fluorophenol	1.256	1.312	-4.5	67	0.00
6		Aniline	2.092	1.979	5.4	61	0.00
7	S	Phenol-d6	1.550	1.561	-0.7	64	0.00
8		2-Chlorophenol	1.423	1.480	-4.0	68	0.00
9		Benzaldehyde	0.899	0.707	21.4	52	0.00
10	C	Phenol	1.733	1.807	-4.3	68	0.00
11		bis(2-Chloroethyl)ether	1.348	1.333	1.1	64	0.00
12		1,3-Dichlorobenzene	1.490	1.569	-5.3	69	0.00
13	C	1,4-Dichlorobenzene	1.516	1.590	-4.9	69	0.00
14		1,2-Dichlorobenzene	1.401	1.446	-3.2	67	0.00
15		Benzyl Alcohol	1.137	0.937	17.6	53	0.00
16		2,2'-oxybis(1-Chloropropane	2.099	1.676	20.2	53	0.00
17		2-Methylphenol	1.164	1.128	3.1	63	0.00
18		Hexachloroethane	0.545	0.538	1.3	65	0.00
19	P	n-Nitroso-di-n-propylamine	0.990	0.833	15.9	57	0.00
20		3+4-Methylphenols	1.473	1.343	8.8	59	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	63	0.00
22		Acetophenone	0.485	0.442	8.9	60	0.00
23	S	Nitrobenzene-d5	0.312	0.302	3.2	61	0.00
24		Nitrobenzene	0.354	0.336	5.1	60	0.00
25		Isophorone	0.645	0.603	6.5	61	0.00
26	C	2-Nitrophenol	0.170	0.195	-14.7	70	0.00
27		2,4-Dimethylphenol	0.321	0.301	6.2	61	0.00
28		bis(2-Chloroethoxy)methane	0.402	0.395	1.7	64	0.00
29	C	2,4-Dichlorophenol	0.276	0.289	-4.7	67	0.00
30		1,2,4-Trichlorobenzene	0.276	0.293	-6.2	68	0.00
31		Naphthalene	0.939	0.961	-2.3	67	0.00
32		Benzoic acid	0.264	0.248	6.1	58	0.00
33		4-Chloroaniline	0.392	0.398	-1.5	65	0.00
34	C	Hexachlorobutadiene	0.142	0.147	-3.5	68	0.00
35		Caprolactam	0.088	0.089	-1.1	66	0.00
36	C	4-Chloro-3-methylphenol	0.310	0.302	2.6	62	0.00
37		2-Methylnaphthalene	0.611	0.617	-1.0	66	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	63	0.00
39		1,2,4,5-Tetrachlorobenzene	0.596	0.643	-7.9	70	0.00
40	P	Hexachlorocyclopentadiene	0.273	0.251	8.1	57	0.00
41	S	2,4,6-Tribromophenol	0.164	0.173	-5.5	65	0.00
42	C	2,4,6-Trichlorophenol	0.423	0.421	0.5	64	0.00
43		2,4,5-Trichlorophenol	0.422	0.427	-1.2	63	0.00
44	S	2-Fluorobiphenyl	1.198	1.275	-6.4	67	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF102017\
Data File : BF099693.D
Acq On : 20 Oct 2017 11:37
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 21 23:57:21 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 20 13:11:43 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	1.779	-3.5	67	0.00
46		2-Chloronaphthalene	1.291	1.327	-2.8	66	0.00
47		2-Nitroaniline	0.395	0.363	8.1	57	0.00
48		Acenaphthylene	1.976	1.959	0.9	64	0.00
49		Dimethylphthalate	1.509	1.557	-3.2	67	0.00
50		2,6-Dinitrotoluene	0.329	0.352	-7.0	66	0.00
51	C	Acenaphthene	1.251	1.179	5.8	61	0.00
52		3-Nitroaniline	0.383	0.405	-5.7	64	0.00
53	P	2,4-Dinitrophenol	0.148	0.135	8.8	55	0.00
54		Dibenzofuran	1.666	1.614	3.1	62	0.00
55	P	4-Nitrophenol	0.324	0.365	-12.7	69	0.00
56		2,4-Dinitrotoluene	0.426	0.421	1.2	61	0.00
57		Fluorene	1.339	1.402	-4.7	68	0.00
58		2,3,4,6-Tetrachlorophenol	0.291	0.289	0.7	63	0.00
59		Diethylphthalate	1.475	1.453	1.5	64	0.00
60		4-Chlorophenyl-phenylether	0.602	0.629	-4.5	68	0.00
61		4-Nitroaniline	0.370	0.370	0.0	62	0.00
62		Azobenzene	1.356	1.206	11.1	57	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	62	0.00
64		4,6-Dinitro-2-methylphenol	0.122	0.137	-12.3	70	0.00
65	c	n-Nitrosodiphenylamine	0.693	0.707	-2.0	66	0.00
66		4-Bromophenyl-phenylether	0.196	0.209	-6.6	68	0.00
67		Hexachlorobenzene	0.209	0.221	-5.7	68	0.00
68		Atrazine	0.208	0.209	-0.5	64	0.00
69	C	Pentachlorophenol	0.130	0.114	12.3	53	0.00
70		Phenanthrene	1.071	1.099	-2.6	67	0.00
71		Anthracene	1.095	1.128	-3.0	66	0.00
72		Carbazole	1.073	1.076	-0.3	64	0.00
73		Di-n-butylphthalate	1.291	1.304	-1.0	64	0.00
74	C	Fluoranthene	1.084	1.105	-1.9	66	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	60	0.00
76		Benzidine	0.740	0.605	18.2	50	0.00
77		Pyrene	1.525	1.616	-6.0	65	0.00
78	S	Terphenyl-d14	0.879	0.948	-7.8	66	0.00
79		Butylbenzylphthalate	0.717	0.778	-8.5	65	0.00
80		Benzo(a)anthracene	1.211	1.244	-2.7	64	0.00
81		3,3'-Dichlorobenzidine	0.444	0.411	7.4	56	0.00
82		Chrysene	1.153	1.171	-1.6	63	0.00
83		Bis(2-ethylhexyl)phthalate	0.987	1.060	-7.4	65	0.00
84	c	Di-n-octyl phthalate	1.589	1.734	-9.1	68	0.00
85		Indeno(1,2,3-cd)pyrene	0.922	0.906	1.7	65	0.00
86	I	Perylene-d12	1.000	1.000	0.0	57	0.00
87		Benzo(b)fluoranthene	1.230	1.220	0.8	56	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF102017\
Data File : BF099693.D
Acq On : 20 Oct 2017 11:37
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 21 23:57:21 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 20 13:11:43 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	1.268	-4.4	60	0.00
89 C	Benzo(a)pyrene	1.129	1.134	-0.4	58	0.00
90	Dibenzo(a,h)anthracene	0.903	1.000	-10.7	65	0.00
91	Benzo(g,h,i)perylene	0.893	1.060	-18.7	70	0.00

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