Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF111017\
Data File : BF100427.D
Acq On : 11 Nov 2017 7:57
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 13 08:31:39 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 10 15:44:40 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	73	0.00
2		1,4-Dioxane	0.608	0.631	-3.8	76	-0.02
3		Pyridine	1.659	1.714	-3.3	74	-0.01
4		n-Nitrosodimethylamine	0.805	0.826	-2.6	73	-0.02
5	S	2-Fluorophenol	1.248	1.272	-1.9	76	0.00
6		Aniline	2.174	2.153	1.0	72	0.00
7	S	Phenol-d6	1.539	1.561	-1.4	75	0.00
8		2-Chlorophenol	1.320	1.362	-3.2	76	0.00
9		Benzaldehyde	1.099	1.105	-0.5	74	0.00
10	C	Phenol	1.615	1.624	-0.6	75	0.00
11		bis(2-Chloroethyl)ether	1.357	1.398	-3.0	76	0.00
12		1,3-Dichlorobenzene	1.522	1.530	-0.5	74	0.00
13	C	1,4-Dichlorobenzene	1.540	1.568	-1.8	75	0.00
14		1,2-Dichlorobenzene	1.424	1.431	-0.5	74	0.00
15		Benzyl Alcohol	1.201	1.181	1.7	71	0.00
16		2,2'-oxybis(1-Chloropropane	2.170	1.797	17.2	61	0.00
17		2-Methylphenol	1.128	1.120	0.7	73	0.00
18		Hexachloroethane	0.508	0.387	23.8	55	0.00
19	P	n-Nitroso-di-n-propylamine	1.067	0.994	6.8	69	0.00
20		3+4-Methylphenols	1.427	1.366	4.3	70	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	69	0.00
22		Acetophenone	0.488	0.484	0.8	70	0.00
23	S	Nitrobenzene-d5	0.303	0.359	-18.5	78	0.00
24		Nitrobenzene	0.311	0.374	-20.3	76	0.00
25		Isophorone	0.636	0.630	0.9	68	0.00
26	C	2-Nitrophenol	0.132	0.152	-15.2	73	0.00
27		2,4-Dimethylphenol	0.285	0.283	0.7	67	0.00
28		bis(2-Chloroethoxy)methane	0.416	0.437	-5.0	73	0.00
29	C	2,4-Dichlorophenol	0.272	0.280	-2.9	68	0.00
30		1,2,4-Trichlorobenzene	0.294	0.302	-2.7	71	0.00
31		Naphthalene	0.963	0.974	-1.1	70	0.00
32		Benzoic acid	0.186	0.183	1.6	64	-0.03
33		4-Chloroaniline	0.401	0.406	-1.2	69	0.00
34	C	Hexachlorobutadiene	0.175	0.180	-2.9	70	0.00
35		Caprolactam	0.093	0.084	9.7	61	0.00
36	C	4-Chloro-3-methylphenol	0.292	0.279	4.5	64	0.00
37		2-Methylnaphthalene	0.640	0.619	3.3	67	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	59	0.00
39		1,2,4,5-Tetrachlorobenzene	0.663	0.766	-15.5	69	0.00
40	P	Hexachlorocyclopentadiene	0.312	0.041#	86.9#	7#	0.00
41	S	2,4,6-Tribromophenol	0.204	0.190	6.9	53	0.00
42	C	2,4,6-Trichlorophenol	0.420	0.450	-7.1	61	0.00
43		2,4,5-Trichlorophenol	0.436	0.459	-5.3	61	0.00
44	S	2-Fluorobiphenyl	1.313	1.427	-8.7	67	0.00

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Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

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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	1.730	1.871	-8.2	66	0.00
46		2-Chloronaphthalene	1.255	1.354	-7.9	64	0.00
47		2-Nitroaniline	0.350	0.450	-28.6#	72	0.00
48		Acenaphthylene	2.080	2.100	-1.0	60	0.00
49		Dimethylphthalate	1.527	1.653	-8.3	66	0.00
50		2,6-Dinitrotoluene	0.302	0.322	-6.6	61	0.00
51	C	Acenaphthene	1.265	1.231	2.7	59	0.00
52		3-Nitroaniline	0.357	0.400	-12.0	64	0.00
53	P	2,4-Dinitrophenol	0.091	0.008#	91.2#	5#	0.00
54		Dibenzofuran	1.773	1.767	0.3	60	0.00
55	P	4-Nitrophenol	0.290	0.233	19.7	46#	0.00
56		2,4-Dinitrotoluene	0.381	0.381	0.0	55	0.00
57		Fluorene	1.299	1.273	2.0	58	0.00
58		2,3,4,6-Tetrachlorophenol	0.357	0.318	10.9	51	0.00
59		Diethylphthalate	1.528	1.589	-4.0	63	0.00
60		4-Chlorophenyl-phenylether	0.637	0.668	-4.9	62	0.00
61		4-Nitroaniline	0.338	0.348	-3.0	58	0.00
62		Azobenzene	1.439	1.373	4.6	58	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	49#	0.00
64		4,6-Dinitro-2-methylphenol	0.087	0.015	82.8#	8#	0.00
65	c	n-Nitrosodiphenylamine	0.695	0.814	-17.1	58	0.00
66		4-Bromophenyl-phenylether	0.214	0.249	-16.4	57	0.00
67		Hexachlorobenzene	0.224	0.248	-10.7	54	0.00
68		Atrazine	0.211	0.232	-10.0	53	0.00
69	C	Pentachlorophenol	0.161	0.148	8.1	44#	0.00
70		Phenanthrene	1.053	1.080	-2.6	52	0.00
71		Anthracene	1.068	1.081	-1.2	51	0.00
72		Carbazole	0.986	0.955	3.1	49#	0.00
73		Di-n-butylphthalate	1.154	1.376	-19.2	59	0.00
74	C	Fluoranthene	1.116	1.063	4.7	48#	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	47#	0.00
76		Benzidine	0.801	0.684	14.6	37#	0.00
77		Pyrene	1.426	1.450	-1.7	48#	0.00
78	S	Terphenyl-d14	0.870	0.905	-4.0	51	0.00
79		Butylbenzylphthalate	0.581	0.653	-12.4	51	0.00
80		Benzo(a)anthracene	1.204	1.265	-5.1	49#	0.00
81		3,3'-Dichlorobenzidine	0.432	0.479	-10.9	49#	0.00
82		Chrysene	1.166	1.193	-2.3	49#	0.00
83		Bis(2-ethylhexyl)phthalate	0.765	0.922	-20.5	54	0.00
84	c	Di-n-octyl phthalate	1.314	1.563	-18.9	53	0.00
85		Indeno(1,2,3-cd)pyrene	0.943	0.867	8.1	44#	0.00
86	I	Perylene-d12	1.000	1.000	0.0	42#	0.00
87		Benzo(b)fluoranthene	1.185	1.335	-12.7	48#	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.120	1.147	-2.4	41#	0.00
89 C	Benzo(a)pyrene	1.050	1.090	-3.8	43#	0.00
90	Dibenzo(a,h)anthracene	0.859	0.921	-7.2	46#	0.00
91	Benzo(g,h,i)perylene	0.841	0.860	-2.3	44#	0.00

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