Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF111317\
Data File : BF100497.D
Acq On : 13 Nov 2017 12:42
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 13 14:08:59 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 13 14:08:38 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	51	-0.01
2		1,4-Dioxane	0.608	0.631	-3.8	53	-0.02
3		Pyridine	1.659	1.790	-7.9	53	-0.02
4		n-Nitrosodimethylamine	0.805	0.880	-9.3	54	-0.02
5	S	2-Fluorophenol	1.248	1.202	3.7	50#	0.00
6		Aniline	2.174	1.963	9.7	46#	-0.01
7	S	Phenol-d6	1.539	1.450	5.8	48#	0.00
8		2-Chlorophenol	1.320	1.312	0.6	51	-0.01
9		Benzaldehyde	1.099	0.965	12.2	45#	0.00
10	C	Phenol	1.615	1.485	8.0	47#	0.00
11		bis(2-Chloroethyl)ether	1.357	1.242	8.5	47#	-0.01
12		1,3-Dichlorobenzene	1.522	1.543	-1.4	52	-0.01
13	C	1,4-Dichlorobenzene	1.540	1.567	-1.8	52	-0.01
14		1,2-Dichlorobenzene	1.424	1.475	-3.6	53	-0.01
15		Benzyl Alcohol	1.201	1.174	2.2	49#	-0.01
16		2,2'-oxybis(1-Chloropropane	2.170	1.843	15.1	43#	-0.01
17		2-Methylphenol	1.128	1.084	3.9	49#	-0.01
18		Hexachloroethane	0.508	0.534	-5.1	52	-0.01
19	P	n-Nitroso-di-n-propylamine	1.067	0.989	7.3	48#	-0.01
20		3+4-Methylphenols	1.427	1.375	3.6	49#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	48#	-0.01
22		Acetophenone	0.488	0.497	-1.8	50#	0.00
23	S	Nitrobenzene-d5	0.303	0.346	-14.2	52	-0.01
24		Nitrobenzene	0.311	0.348	-11.9	50#	-0.01
25		Isophorone	0.636	0.596	6.3	45#	-0.01
26	C	2-Nitrophenol	0.132	0.185	-40.2#	62	-0.01
27		2,4-Dimethylphenol	0.285	0.292	-2.5	48#	0.00
28		bis(2-Chloroethoxy)methane	0.416	0.404	2.9	47#	-0.01
29	C	2,4-Dichlorophenol	0.272	0.300	-10.3	51	0.00
30		1,2,4-Trichlorobenzene	0.294	0.327	-11.2	53	-0.01
31		Naphthalene	0.963	0.999	-3.7	50	0.00
32		Benzoic acid	0.186	0.255	-37.1#	63	0.00
33		4-Chloroaniline	0.401	0.410	-2.2	48#	-0.01
34	C	Hexachlorobutadiene	0.175	0.203	-16.0	55	-0.01
35		Caprolactam	0.093	0.085	8.6	43#	-0.01
36	C	4-Chloro-3-methylphenol	0.292	0.300	-2.7	48#	-0.01
37		2-Methylnaphthalene	0.640	0.678	-5.9	51	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	47#	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.663	0.766	-15.5	54	-0.01
40	P	Hexachlorocyclopentadiene	0.312	0.406	-30.1#	57	-0.01
41	S	2,4,6-Tribromophenol	0.204	0.243	-19.1	54	-0.01
42	C	2,4,6-Trichlorophenol	0.420	0.477	-13.6	51	-0.01
43		2,4,5-Trichlorophenol	0.436	0.501	-14.9	52	0.00
44	S	2-Fluorobiphenyl	1.313	1.479	-12.6	55	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.730	1.873	-8.3	52	-0.01
46		2-Chloronaphthalene	1.255	1.351	-7.6	51	-0.01
47		2-Nitroaniline	0.350	0.397	-13.4	50	-0.01
48		Acenaphthylene	2.080	2.180	-4.8	50#	-0.01
49		Dimethylphthalate	1.527	1.597	-4.6	50	-0.01
50		2,6-Dinitrotoluene	0.302	0.341	-12.9	51	-0.01
51	C	Acenaphthene	1.265	1.369	-8.2	52	-0.01
52		3-Nitroaniline	0.357	0.381	-6.7	49#	-0.01
53	P	2,4-Dinitrophenol	0.091	0.173	-90.1#	90	-0.01
54		Dibenzofuran	1.773	1.867	-5.3	50	-0.01
55	P	4-Nitrophenol	0.290	0.303	-4.5	47#	0.00
56		2,4-Dinitrotoluene	0.381	0.449	-17.8	52	0.00
57		Fluorene	1.299	1.438	-10.7	52	-0.01
58		2,3,4,6-Tetrachlorophenol	0.357	0.399	-11.8	51	-0.01
59		Diethylphthalate	1.528	1.560	-2.1	49#	-0.01
60		4-Chlorophenyl-phenylether	0.637	0.716	-12.4	53	-0.01
61		4-Nitroaniline	0.338	0.369	-9.2	49#	-0.01
62		Azobenzene	1.439	1.350	6.2	45#	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	47#	-0.01
64		4,6-Dinitro-2-methylphenol	0.087	0.139	-59.8#	70	0.00
65	c	n-Nitrosodiphenylamine	0.695	0.729	-4.9	50#	-0.01
66		4-Bromophenyl-phenylether	0.214	0.235	-9.8	52	-0.01
67		Hexachlorobenzene	0.224	0.250	-11.6	53	-0.01
68		Atrazine	0.211	0.221	-4.7	49#	-0.01
69	C	Pentachlorophenol	0.161	0.181	-12.4	52	0.00
70		Phenanthrene	1.053	1.083	-2.8	51	-0.01
71		Anthracene	1.068	1.103	-3.3	50	-0.01
72		Carbazole	0.986	0.983	0.3	48#	-0.01
73		Di-n-butylphthalate	1.154	1.224	-6.1	51	-0.02
74	C	Fluoranthene	1.116	1.143	-2.4	50	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	47#	-0.02
76		Benzidine	0.801	0.774	3.4	43#	-0.01
77		Pyrene	1.426	1.469	-3.0	49#	-0.01
78	S	Terphenyl-d14	0.870	0.925	-6.3	53	-0.01
79		Butylbenzylphthalate	0.581	0.640	-10.2	51	-0.01
80		Benzo(a)anthracene	1.204	1.271	-5.6	50	-0.02
81		3,3'-Dichlorobenzidine	0.432	0.469	-8.6	49#	-0.01
82		Chrysene	1.166	1.192	-2.2	49#	-0.01
83		Bis(2-ethylhexyl)phthalate	0.765	0.880	-15.0	52	-0.02
84	c	Di-n-octyl phthalate	1.314	1.428	-8.7	49#	-0.02
85		Indeno(1,2,3-cd)pyrene	0.943	0.845	10.4	44#	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	43#	-0.02
87		Benzo(b)fluoranthene	1.185	1.343	-13.3	49#	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.120	1.195	-6.7	43#	-0.01
89 C	Benzo(a)pyrene	1.050	1.119	-6.6	45#	-0.02
90	Dibenzo(a,h)anthracene	0.859	0.882	-2.7	45#	-0.03
91	Benzo(g,h,i)perylene	0.841	0.827	1.7	43#	-0.03

(#) = Out of Range SPCC's out = 0 CCC's out = 1