Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF092518\
Data File : BF109282.D
Acq On : 25 Sep 2018 13:43
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 25 15:23:26 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Sep 22 13:32:18 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	106	0.00
2		1,4-Dioxane	0.559	0.533	4.7	102	0.00
3		Pyridine	1.439	1.464	-1.7	103	0.01
4		n-Nitrosodimethylamine	0.613	0.582	5.1	100	0.00
5	S	2-Fluorophenol	1.214	1.204	0.8	108	0.00
6		Aniline	1.982	1.940	2.1	106	0.00
7	S	Phenol-d6	1.549	1.535	0.9	108	0.00
8		2-Chlorophenol	1.350	1.356	-0.4	109	0.00
9		Benzaldehyde	0.995	0.953	4.2	106	0.00
10	C	Phenol	1.755	1.700	3.1	106	0.00
11		bis(2-Chloroethyl)ether	1.373	1.337	2.6	107	0.00
12		1,3-Dichlorobenzene	1.555	1.559	-0.3	109	0.00
13	C	1,4-Dichlorobenzene	1.566	1.544	1.4	108	0.00
14		1,2-Dichlorobenzene	1.445	1.456	-0.8	110	0.00
15		Benzyl Alcohol	1.186	1.218	-2.7	110	0.00
16		2,2'-oxybis(1-Chloropropane	1.948	1.869	4.1	105	0.00
17		2-Methylphenol	1.093	1.088	0.5	110	0.00
18		Hexachloroethane	0.552	0.560	-1.4	110	0.00
19	P	n-Nitroso-di-n-propylamine	1.015	0.998	1.7	109	-0.01
20		3+4-Methylphenols	1.319	1.325	-0.5	112	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	111	0.00
22		Acetophenone	0.462	0.446	3.5	110	0.00
23	S	Nitrobenzene-d5	0.339	0.356	-5.0	113	0.00
24		Nitrobenzene	0.360	0.369	-2.5	113	0.00
25		Isophorone	0.610	0.587	3.8	107	0.00
26	C	2-Nitrophenol	0.142	0.174	-22.5#	129	0.00
27		2,4-Dimethylphenol	0.266	0.268	-0.8	110	0.00
28		bis(2-Chloroethoxy)methane	0.404	0.396	2.0	111	0.00
29	C	2,4-Dichlorophenol	0.279	0.287	-2.9	112	0.00
30		1,2,4-Trichlorobenzene	0.312	0.312	0.0	112	0.00
31		Naphthalene	0.935	0.918	1.8	110	0.00
32		Benzoic acid	0.163	0.177	-8.6	115	-0.04
33		4-Chloroaniline	0.408	0.406	0.5	112	0.00
34	C	Hexachlorobutadiene	0.188	0.191	-1.6	113	0.00
35		Caprolactam	0.089	0.091	-2.2	110	-0.02
36	C	4-Chloro-3-methylphenol	0.294	0.304	-3.4	114	0.00
37		2-Methylnaphthalene	0.602	0.598	0.7	112	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	114	0.00
39		1,2,4,5-Tetrachlorobenzene	0.637	0.627	1.6	114	0.00
40	P	Hexachlorocyclopentadiene	0.278	0.265	4.7	105	0.00
41	S	2,4,6-Tribromophenol	0.197	0.217	-10.2	124	0.00
42	C	2,4,6-Trichlorophenol	0.409	0.425	-3.9	116	0.00
43		2,4,5-Trichlorophenol	0.431	0.453	-5.1	116	0.00
44	S	2-Fluorobiphenyl	1.326	1.311	1.1	114	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.630	1.592	2.3	113	0.00
46		2-Chloronaphthalene	1.302	1.271	2.4	114	0.00
47		2-Nitroaniline	0.369	0.393	-6.5	117	0.00
48		Acenaphthylene	1.964	1.908	2.9	112	0.00
49		Dimethylphthalate	1.455	1.423	2.2	114	0.00
50		2,6-Dinitrotoluene	0.288	0.319	-10.8	118	0.00
51	C	Acenaphthene	1.187	1.124	5.3	109	0.00
52		3-Nitroaniline	0.334	0.368	-10.2	119	0.00
53	P	2,4-Dinitrophenol	0.082	0.102	-24.4	135	0.00
54		Dibenzofuran	1.766	1.717	2.8	114	0.00
55	P	4-Nitrophenol	0.251	0.261	-4.0	116	0.00
56		2,4-Dinitrotoluene	0.365	0.418	-14.5	126	0.00
57		Fluorene	1.260	1.222	3.0	116	0.00
58		2,3,4,6-Tetrachlorophenol	0.342	0.367	-7.3	120	0.00
59		Diethylphthalate	1.473	1.420	3.6	111	0.00
60		4-Chlorophenyl-phenylether	0.658	0.645	2.0	118	0.00
61		4-Nitroaniline	0.325	0.354	-8.9	118	-0.01
62		Azobenzene	1.530	1.427	6.7	108	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	114	0.00
64		4,6-Dinitro-2-methylphenol	0.072	0.089	-23.6	136	0.00
65	c	n-Nitrosodiphenylamine	0.661	0.645	2.4	114	0.00
66		4-Bromophenyl-phenylether	0.222	0.219	1.4	114	0.00
67		Hexachlorobenzene	0.236	0.237	-0.4	118	0.00
68		Atrazine	0.203	0.204	-0.5	115	0.00
69	C	Pentachlorophenol	0.141	0.146	-3.5	114	0.00
70		Phenanthrene	0.999	0.981	1.8	115	0.00
71		Anthracene	1.013	0.982	3.1	112	0.00
72		Carbazole	1.008	0.978	3.0	114	0.00
73		Di-n-butylphthalate	1.160	1.117	3.7	111	0.00
74	C	Fluoranthene	1.067	1.025	3.9	112	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	108	0.00
76		Benzidine	0.721	0.934	-29.5#	135	0.00
77		Pyrene	1.433	1.483	-3.5	111	0.00
78	S	Terphenyl-d14	0.815	0.829	-1.7	112	0.00
79		Butylbenzylphthalate	0.610	0.650	-6.6	112	0.00
80		Benzo(a)anthracene	1.205	1.115	7.5	100	0.00
81		3,3'-Dichlorobenzidine	0.458	0.444	3.1	100	0.00
82		Chrysene	1.194	1.142	4.4	103	0.00
83		Bis(2-ethylhexyl)phthalate	0.769	0.760	1.2	104	0.00
84	c	Di-n-octyl phthalate	1.315	1.247	5.2	97	0.00
85		Indeno(1,2,3-cd)pyrene	0.968	0.872	9.9	102	0.02
86	I	Perylene-d12	1.000	1.000	0.0	87	0.01
87		Benzo(b)fluoranthene	1.099	1.118	-1.7	89	0.00

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88	Benzo(k)fluoranthene	1.043	1.058	-1.4	90	0.00
89 C	Benzo(a)pyrene	0.990	0.993	-0.3	88	0.00
90	Dibenzo(a,h)anthracene	0.812	0.893	-10.0	101	0.02
91	Benzo(g,h,i)perylene	0.798	0.920	-15.3	109	0.02

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