Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061917\
Data File : BF095999.D
Acq On : 19 Jun 2017 11:13
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 20 02:18:19 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 16 17:12:03 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	75	0.01
2		1,4-Dioxane	0.597	0.724	-21.3	87	0.02
3		Pyridine	1.403	1.271	9.4	65	0.03
4		n-Nitrosodimethylamine	0.534	0.546	-2.2	75	0.03
5	S	2-Fluorophenol	1.255	1.308	-4.2	71	0.00
6		Aniline	1.966	2.012	-2.3	72	0.02
7	S	Phenol-d6	1.503	1.526	-1.5	70	0.00
8		2-Chlorophenol	1.335	1.375	-3.0	71	0.02
9		Benzaldehyde	1.014	0.933	8.0	65	0.02
10	C	Phenol	1.559	1.640	-5.2	71	0.00
11		bis(2-Chloroethyl)ether	1.235	1.347	-9.1	75	0.02
12		1,3-Dichlorobenzene	1.511	1.538	-1.8	75	0.01
13	C	1,4-Dichlorobenzene	1.532	1.576	-2.9	75	0.01
14		1,2-Dichlorobenzene	1.412	1.478	-4.7	76	0.02
15		Benzyl Alcohol	1.031	1.067	-3.5	73	0.01
16		2,2'-oxybis(1-Chloropropane	1.735	1.825	-5.2	76	0.00
17		2-Methylphenol	1.120	1.185	-5.8	76	0.00
18		Hexachloroethane	0.506	0.530	-4.7	76	0.01
19	P	n-Nitroso-di-n-propylamine	0.952	1.024	-7.6	77	0.02
20		3+4-Methylphenols	1.422	1.553	-9.2	78	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	76	0.01
22		Acetophenone	0.468	0.476	-1.7	76	0.02
23	S	Nitrobenzene-d5	0.322	0.324	-0.6	75	0.02
24		Nitrobenzene	0.344	0.343	0.3	75	0.01
25		Isophorone	0.601	0.604	-0.5	76	0.02
26	C	2-Nitrophenol	0.180	0.185	-2.8	75	0.02
27		2,4-Dimethylphenol	0.280	0.287	-2.5	77	0.01
28		bis(2-Chloroethoxy)methane	0.396	0.404	-2.0	75	0.01
29	C	2,4-Dichlorophenol	0.269	0.272	-1.1	75	0.00
30		1,2,4-Trichlorobenzene	0.280	0.281	-0.4	76	0.01
31		Naphthalene	1.004	0.998	0.6	76	0.01
32		Benzoic acid	0.161	0.156	3.1	70	0.02
33		4-Chloroaniline	0.392	0.407	-3.8	79	0.01
34	C	Hexachlorobutadiene	0.145	0.145	0.0	76	0.01
35		Caprolactam	0.080	0.082	-2.5	74	0.02
36	C	4-Chloro-3-methylphenol	0.282	0.288	-2.1	76	0.01
37		2-Methylnaphthalene	0.678	0.683	-0.7	77	0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	81	0.00
39		1,2,4,5-Tetrachlorobenzene	0.599	0.617	-3.0	78	0.01
40	P	Hexachlorocyclopentadiene	0.138	0.174	-26.1#	92	0.01
41	S	2,4,6-Tribromophenol	0.177	0.178	-0.6	79	0.00
42	C	2,4,6-Trichlorophenol	0.385	0.395	-2.6	77	0.01
43		2,4,5-Trichlorophenol	0.409	0.387	5.4	68	0.00
44	S	2-Fluorobiphenyl	1.437	1.411	1.8	77	0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.648	1.654	-0.4	77	0.01
46		2-Chloronaphthalene	1.288	1.282	0.5	76	0.01
47		2-Nitroaniline	0.377	0.381	-1.1	72	0.01
48		Acenaphthylene	2.202	2.097	4.8	71	0.01
49		Dimethylphthalate	1.519	1.472	3.1	74	0.01
50		2,6-Dinitrotoluene	0.331	0.315	4.8	69	0.01
51	C	Acenaphthene	1.272	1.244	2.2	79	0.01
52		3-Nitroaniline	0.386	0.389	-0.8	80	0.01
53	P	2,4-Dinitrophenol	0.160	0.152	5.0	74	0.01
54		Dibenzofuran	1.790	1.772	1.0	80	0.01
55	P	4-Nitrophenol	0.201	0.173	13.9	64	0.00
56		2,4-Dinitrotoluene	0.447	0.460	-2.9	80	0.01
57		Fluorene	1.558	1.529	1.9	79	0.01
58		2,3,4,6-Tetrachlorophenol	0.280	0.274	2.1	75	0.01
59		Diethylphthalate	1.461	1.443	1.2	79	0.02
60		4-Chlorophenyl-phenylether	0.677	0.678	-0.1	80	0.01
61		4-Nitroaniline	0.360	0.350	2.8	76	0.01
62		Azobenzene	1.324	1.302	1.7	78	0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	77	0.01
64		4,6-Dinitro-2-methylphenol	0.131	0.125	4.6	70	0.01
65	c	n-Nitrosodiphenylamine	0.719	0.721	-0.3	79	0.01
66		4-Bromophenyl-phenylether	0.208	0.210	-1.0	77	0.01
67		Hexachlorobenzene	0.205	0.215	-4.9	80	0.00
68		Atrazine	0.200	0.207	-3.5	81	0.00
69	C	Pentachlorophenol	0.067	0.068	-1.5	78	0.00
70		Phenanthrene	1.154	1.144	0.9	78	0.01
71		Anthracene	1.205	1.195	0.8	78	0.01
72		Carbazole	1.174	1.212	-3.2	79	0.01
73		Di-n-butylphthalate	1.249	1.309	-4.8	79	0.00
74	C	Fluoranthene	1.142	1.172	-2.6	79	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	78	0.00
76		Benzidine	0.768	0.890	-15.9	88	0.01
77		Pyrene	1.492	1.518	-1.7	79	0.01
78	S	Terphenyl-d14	0.806	0.823	-2.1	80	0.00
79		Butylbenzylphthalate	0.652	0.681	-4.4	79	0.01
80		Benzo(a)anthracene	1.163	1.163	0.0	77	0.00
81		3,3'-Dichlorobenzidine	0.413	0.419	-1.5	77	0.00
82		Chrysene	1.162	1.172	-0.9	78	0.00
83		Bis(2-ethylhexyl)phthalate	0.919	0.951	-3.5	78	0.01
84	c	Di-n-octyl phthalate	1.515	1.534	-1.3	77	0.00
85		Indeno(1,2,3-cd)pyrene	0.994	0.881	11.4	73	0.00
86	I	Perylene-d12	1.000	1.000	0.0	73	0.00
87		Benzo(b)fluoranthene	1.252	1.303	-4.1	72	0.00

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88	Benzo(k)fluoranthene	1.197	1.228	-2.6	77	0.00
89 C	Benzo(a)pyrene	1.151	1.161	-0.9	74	0.00
90	Dibenzo(a,h)anthracene	0.976	0.918	5.9	73	0.00
91	Benzo(g,h,i)perylene	0.995	0.928	6.7	72	0.00

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