Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030717\
Data File : BF093418.D
Acq On : 7 Mar 2017 14:15
Operator : SJ/MA
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF030717

Quant Time: Mar 07 15:56:52 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 07 15:55:22 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	107	0.00
2		1,4-Dioxane	0.662	0.692	-4.5	109	0.00
3		Pyridine	1.688	1.860	-10.2	108	0.00
4		n-Nitrosodimethylamine	0.806	0.787	2.4	111	0.00
5	S	2-Fluorophenol	1.202	1.183	1.6	106	0.00
6		Aniline	2.080	1.953	6.1	107	0.00
7	S	Phenol-d6	1.515	1.478	2.4	109	0.00
8		2-Chlorophenol	1.346	1.311	2.6	106	0.00
9		Benzaldehyde	1.017	0.940	7.6	110	0.00
10	C	Phenol	1.857	1.759	5.3	111	0.00
11		bis(2-Chloroethyl)ether	1.501	1.435	4.4	107	0.00
12		1,3-Dichlorobenzene	1.541	1.464	5.0	106	0.00
13	C	1,4-Dichlorobenzene	1.578	1.533	2.9	108	0.00
14		1,2-Dichlorobenzene	1.397	1.360	2.6	107	0.00
15		Benzyl Alcohol	1.080	1.021	5.5	109	0.01
16		2,2'-oxybis(1-Chloropropane	2.493	2.371	4.9	103	0.00
17		2-Methylphenol	1.139	1.056	7.3	104	0.00
18		Hexachloroethane	0.532	0.500	6.0	104	0.00
19	P	n-Nitroso-di-n-propylamine	1.042	0.881	15.5	101	0.00
20		3+4-Methylphenols	1.477	1.401	5.1	103	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	103	-0.01
22		Acetophenone	0.481	0.469	2.5	103	0.00
23	S	Nitrobenzene-d5	0.360	0.345	4.2	100	0.00
24		Nitrobenzene	0.405	0.406	-0.2	100	0.00
25		Isophorone	0.727	0.707	2.8	103	0.00
26	C	2-Nitrophenol	0.185	0.186	-0.5	99	0.00
27		2,4-Dimethylphenol	0.301	0.297	1.3	94	0.00
28		bis(2-Chloroethoxy)methane	0.459	0.421	8.3	99	0.00
29	C	2,4-Dichlorophenol	0.283	0.285	-0.7	104	0.00
30		1,2,4-Trichlorobenzene	0.301	0.306	-1.7	104	0.00
31		Naphthalene	1.002	0.926	7.6	100	-0.01
32		Benzoic acid	0.156	0.143	8.3	99	0.00
33		4-Chloroaniline	0.407	0.405	0.5	103	0.00
34	C	Hexachlorobutadiene	0.155	0.149	3.9	102	0.00
35		Caprolactam	0.097	0.102	-5.2	104	0.00
36	C	4-Chloro-3-methylphenol	0.302	0.308	-2.0	104	0.00
37		2-Methylnaphthalene	0.638	0.650	-1.9	106	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	125	0.00
39		1,2,4,5-Tetrachlorobenzene	0.546	0.477	12.6	100	0.00
40	P	Hexachlorocyclopentadiene	0.098	0.095	3.1	109	0.00
41	S	2,4,6-Tribromophenol	0.130	0.118	9.2	105	0.00
42	C	2,4,6-Trichlorophenol	0.341	0.311	8.8	105	0.00
43		2,4,5-Trichlorophenol	0.366	0.350	4.4	105	0.00
44	S	2-Fluorobiphenyl	1.075	1.010	6.0	107	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.457	1.236	15.2	101	0.00
46		2-Chloronaphthalene	1.115	1.070	4.0	104	0.00
47		2-Nitroaniline	0.394	0.381	3.3	109	0.00
48		Acenaphthylene	1.839	1.723	6.3	107	0.00
49		Dimethylphthalate	1.326	1.197	9.7	110	0.00
50		2,6-Dinitrotoluene	0.291	0.261	10.3	115	0.00
51	C	Acenaphthene	1.088	1.027	5.6	107	0.00
52		3-Nitroaniline	0.350	0.304	13.1	112	0.00
53	P	2,4-Dinitrophenol	0.132	0.153	-15.9	117	0.00
54		Dibenzofuran	1.361	1.335	1.9	108	0.00
55	P	4-Nitrophenol	0.170	0.167	1.8	105	0.01
56		2,4-Dinitrotoluene	0.357	0.303	15.1	107	0.00
57		Fluorene	1.154	1.139	1.3	106	0.00
58		2,3,4,6-Tetrachlorophenol	0.258	0.258	0.0	105	0.00
59		Diethylphthalate	1.267	1.133	10.6	104	0.00
60		4-Chlorophenyl-phenylether	0.517	0.521	-0.8	107	0.00
61		4-Nitroaniline	0.357	0.307	14.0	109	0.00
62		Azobenzene	1.440	1.424	1.1	118	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	113	0.00
64		4,6-Dinitro-2-methylphenol	0.130	0.131	-0.8	112	0.01
65	c	n-Nitrosodiphenylamine	0.668	0.679	-1.6	110	0.00
66		4-Bromophenyl-phenylether	0.211	0.202	4.3	110	0.00
67		Hexachlorobenzene	0.202	0.207	-2.5	112	0.00
68		Atrazine	0.195	0.190	2.6	109	0.00
69	C	Pentachlorophenol	0.070	0.070	0.0	110	0.00
70		Phenanthrene	1.142	1.105	3.2	110	0.00
71		Anthracene	1.155	1.005	13.0	107	0.00
72		Carbazole	1.085	0.981	9.6	108	0.00
73		Di-n-butylphthalate	1.255	1.143	8.9	108	0.00
74	C	Fluoranthene	1.103	1.085	1.6	113	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	107	0.01
76		Benzidine	0.658	0.642	2.4	106	0.00
77		Pyrene	1.455	1.467	-0.8	112	0.01
78	S	Terphenyl-d14	0.769	0.783	-1.8	100	0.00
79		Butylbenzylphthalate	0.612	0.641	-4.7	106	0.00
80		Benzo(a)anthracene	1.181	1.213	-2.7	110	0.01
81		3,3'-Dichlorobenzidine	0.414	0.402	2.9	100	0.00
82		Chrysene	1.093	1.211	-10.8	105	0.00
83		Bis(2-ethylhexyl)phthalate	0.853	0.915	-7.3	108	0.00
84	c	Di-n-octyl phthalate	1.422	1.456	-2.4	105	0.00
85		Indeno(1,2,3-cd)pyrene	0.915	0.898	1.9	105	0.01
86	I	Perylene-d12	1.000	1.000	0.0	106	0.01
87		Benzo(b)fluoranthene	1.218	1.186	2.6	94	0.00

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88	Benzo(k)fluoranthene	1.175	1.039	11.6	112	0.00
89 C	Benzo(a)pyrene	1.113	1.066	4.2	101	0.01
90	Dibenzo(a,h)anthracene	0.921	0.862	6.4	105	0.01
91	Benzo(g,h,i)perylene	0.945	0.897	5.1	106	0.00

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