Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF070517\
Data File : BF096472.D
Acq On : 5 Jul 2017 9:46
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 05 13:15:08 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF070117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 03 01:43:12 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	78	0.00
2		1,4-Dioxane	0.643	0.592	7.9	69	0.00
3		Pyridine	1.763	1.479	16.1	66	0.00
4		n-Nitrosodimethylamine	0.870	0.789	9.3	70	0.00
5	S	2-Fluorophenol	1.355	1.261	6.9	74	0.00
6		Aniline	2.153	2.029	5.8	74	0.00
7	S	Phenol-d6	1.666	1.580	5.2	75	0.00
8		2-Chlorophenol	1.445	1.533	-6.1	84	0.00
9		Benzaldehyde	1.043	0.930	10.8	70	0.00
10	C	Phenol	1.898	1.742	8.2	72	0.00
11		bis(2-Chloroethyl)ether	1.467	1.423	3.0	75	0.00
12		1,3-Dichlorobenzene	1.515	1.550	-2.3	82	0.00
13	C	1,4-Dichlorobenzene	1.514	1.553	-2.6	80	0.00
14		1,2-Dichlorobenzene	1.390	1.431	-2.9	82	0.00
15		Benzyl Alcohol	1.114	1.095	1.7	77	0.00
16		2,2'-oxybis(1-Chloropropane	2.982	2.842	4.7	74	0.00
17		2-Methylphenol	1.152	1.160	-0.7	79	0.00
18		Hexachloroethane	0.549	0.526	4.2	74	0.00
19	P	n-Nitroso-di-n-propylamine	0.992	0.899	9.4	71	0.00
20		3+4-Methylphenols	1.284	1.276	0.6	77	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	77	0.00
22		Acetophenone	0.437	0.418	4.3	76	0.00
23	S	Nitrobenzene-d5	0.333	0.319	4.2	73	0.00
24		Nitrobenzene	0.349	0.332	4.9	73	0.00
25		Isophorone	0.645	0.595	7.8	70	0.00
26	C	2-Nitrophenol	0.185	0.191	-3.2	77	0.00
27		2,4-Dimethylphenol	0.315	0.306	2.9	75	0.00
28		bis(2-Chloroethoxy)methane	0.426	0.397	6.8	71	0.00
29	C	2,4-Dichlorophenol	0.271	0.281	-3.7	80	0.00
30		1,2,4-Trichlorobenzene	0.256	0.262	-2.3	79	0.00
31		Naphthalene	0.944	0.940	0.4	77	0.00
32		Benzoic acid	0.264	0.244	7.6	69	0.00
33		4-Chloroaniline	0.392	0.398	-1.5	78	0.00
34	C	Hexachlorobutadiene	0.131	0.136	-3.8	80	0.00
35		Caprolactam	0.094	0.090	4.3	76	0.00
36	C	4-Chloro-3-methylphenol	0.300	0.301	-0.3	79	0.00
37		2-Methylnaphthalene	0.601	0.615	-2.3	80	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	79	0.00
39		1,2,4,5-Tetrachlorobenzene	0.519	0.548	-5.6	84	0.00
40	P	Hexachlorocyclopentadiene	0.253	0.264	-4.3	78	0.00
41	S	2,4,6-Tribromophenol	0.156	0.176	-12.8	91	0.00
42	C	2,4,6-Trichlorophenol	0.411	0.454	-10.5	87	0.00
43		2,4,5-Trichlorophenol	0.406	0.451	-11.1	88	0.00
44	S	2-Fluorobiphenyl	1.170	1.269	-8.5	85	0.00

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.714	1.824	-6.4	86	0.00
46		2-Chloronaphthalene	1.277	1.354	-6.0	85	0.00
47		2-Nitroaniline	0.465	0.448	3.7	75	0.00
48		Acenaphthylene	2.024	2.158	-6.6	86	0.00
49		Dimethylphthalate	1.493	1.533	-2.7	83	0.00
50		2,6-Dinitrotoluene	0.330	0.369	-11.8	87	0.00
51	C	Acenaphthene	1.247	1.166	6.5	75	0.00
52		3-Nitroaniline	0.412	0.457	-10.9	89	0.00
53	P	2,4-Dinitrophenol	0.175	0.177	-1.1	76	0.00
54		Dibenzofuran	1.647	1.651	-0.2	80	0.00
55	P	4-Nitrophenol	0.341	0.322	5.6	73	0.00
56		2,4-Dinitrotoluene	0.418	0.442	-5.7	81	0.00
57		Fluorene	1.163	1.240	-6.6	84	0.00
58		2,3,4,6-Tetrachlorophenol	0.281	0.292	-3.9	82	0.00
59		Diethylphthalate	1.411	1.410	0.1	80	0.00
60		4-Chlorophenyl-phenylether	0.506	0.524	-3.6	83	0.00
61		4-Nitroaniline	0.374	0.394	-5.3	83	0.00
62		Azobenzene	1.365	1.260	7.7	73	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	82	0.00
64		4,6-Dinitro-2-methylphenol	0.138	0.144	-4.3	80	0.00
65	c	n-Nitrosodiphenylamine	0.702	0.704	-0.3	83	0.00
66		4-Bromophenyl-phenylether	0.195	0.203	-4.1	85	0.00
67		Hexachlorobenzene	0.213	0.224	-5.2	86	0.00
68		Atrazine	0.200	0.212	-6.0	87	0.00
69	C	Pentachlorophenol	0.126	0.132	-4.8	80	0.00
70		Phenanthrene	1.081	1.071	0.9	83	0.00
71		Anthracene	1.102	1.086	1.5	82	0.00
72		Carbazole	1.108	1.080	2.5	81	0.00
73		Di-n-butylphthalate	1.342	1.304	2.8	80	0.00
74	C	Fluoranthene	1.060	1.076	-1.5	84	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	89	0.00
76		Benzidine	0.960	0.903	5.9	83	0.00
77		Pyrene	1.605	1.519	5.4	85	0.00
78	S	Terphenyl-d14	0.936	0.907	3.1	90	0.00
79		Butylbenzylphthalate	0.813	0.758	6.8	82	0.00
80		Benzo(a)anthracene	1.158	1.139	1.6	90	0.00
81		3,3'-Dichlorobenzidine	0.476	0.499	-4.8	92	0.00
82		Chrysene	1.213	1.188	2.1	88	0.00
83		Bis(2-ethylhexyl)phthalate	0.907	0.869	4.2	87	0.00
84	c	Di-n-octyl phthalate	1.786	1.757	1.6	88	0.00
85		Indeno(1,2,3-cd)pyrene	0.957	1.042	-8.9	102	0.00
86	I	Perylene-d12	1.000	1.000	0.0	102	0.00
87		Benzo(b)fluoranthene	1.253	1.141	8.9	91	0.00

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88	Benzo(k)fluoranthene	1.121	1.157	-3.2	110	0.00
89 C	Benzo(a)pyrene	1.119	1.101	1.6	101	0.00
90	Dibenzo(a,h)anthracene	0.880	0.861	2.2	103	0.00
91	Benzo(g,h,i)perylene	0.912	0.869	4.7	100	0.00

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