Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF071717\
Data File : BF096810.D
Acq On : 17 Jul 2017 10:48
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 18 16:41:11 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF071317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 18 13:33:22 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	116	0.01
2		1,4-Dioxane	0.684	0.557	18.6	89	0.02
3		Pyridine	1.438	1.366	5.0	95	0.02
4		n-Nitrosodimethylamine	0.804	0.758	5.7	100	0.02
5	S	2-Fluorophenol	1.330	1.218	8.4	100	0.01
6		Aniline	1.966	1.929	1.9	119	0.01
7	S	Phenol-d6	1.596	1.581	0.9	120	0.01
8		2-Chlorophenol	1.435	1.400	2.4	116	0.01
9		Benzaldehyde	0.923	0.881	4.6	105	0.01
10	C	Phenol	1.804	1.750	3.0	119	0.00
11		bis(2-Chloroethyl)ether	1.357	1.333	1.8	119	0.00
12		1,3-Dichlorobenzene	1.525	1.502	1.5	116	0.01
13	C	1,4-Dichlorobenzene	1.545	1.548	-0.2	118	0.01
14		1,2-Dichlorobenzene	1.405	1.433	-2.0	118	0.01
15		Benzyl Alcohol	1.045	1.019	2.5	113	0.01
16		2,2'-oxybis(1-Chloropropane	2.800	2.940	-5.0	128	0.01
17		2-Methylphenol	1.116	1.147	-2.8	123	0.00
18		Hexachloroethane	0.548	0.526	4.0	115	0.01
19	P	n-Nitroso-di-n-propylamine	0.962	0.911	5.3	116	0.01
20		3+4-Methylphenols	1.354	1.310	3.2	117	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	116	0.00
22		Acetophenone	0.447	0.454	-1.6	123	0.01
23	S	Nitrobenzene-d5	0.323	0.314	2.8	114	0.01
24		Nitrobenzene	0.334	0.321	3.9	114	0.01
25		Isophorone	0.601	0.626	-4.2	125	0.01
26	C	2-Nitrophenol	0.186	0.187	-0.5	115	0.01
27		2,4-Dimethylphenol	0.305	0.315	-3.3	121	0.00
28		bis(2-Chloroethoxy)methane	0.406	0.425	-4.7	127	0.01
29	C	2,4-Dichlorophenol	0.277	0.267	3.6	113	0.01
30		1,2,4-Trichlorobenzene	0.263	0.258	1.9	114	0.01
31		Naphthalene	0.947	0.961	-1.5	119	0.00
32		Benzoic acid	0.195	0.241	-23.6	146	0.00
33		4-Chloroaniline	0.375	0.383	-2.1	118	0.00
34	C	Hexachlorobutadiene	0.140	0.140	0.0	115	0.00
35		Caprolactam	0.089	0.082	7.9	113	0.00
36	C	4-Chloro-3-methylphenol	0.297	0.281	5.4	111	0.01
37		2-Methylnaphthalene	0.604	0.578	4.3	111	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	115	0.00
39		1,2,4,5-Tetrachlorobenzene	0.541	0.538	0.6	112	0.00
40	P	Hexachlorocyclopentadiene	0.180	0.182	-1.1	109	0.00
41	S	2,4,6-Tribromophenol	0.171	0.176	-2.9	114	0.00
42	C	2,4,6-Trichlorophenol	0.399	0.396	0.8	110	0.00
43		2,4,5-Trichlorophenol	0.402	0.399	0.7	113	0.00
44	S	2-Fluorobiphenyl	1.266	1.214	4.1	110	0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.715	1.688	1.6	111	0.00
46		2-Chloronaphthalene	1.263	1.259	0.3	112	0.00
47		2-Nitroaniline	0.434	0.435	-0.2	115	0.00
48		Acenaphthylene	2.012	1.973	1.9	112	0.01
49		Dimethylphthalate	1.506	1.487	1.3	114	0.00
50		2,6-Dinitrotoluene	0.326	0.328	-0.6	113	0.00
51	C	Acenaphthene	1.189	1.137	4.4	112	0.00
52		3-Nitroaniline	0.386	0.400	-3.6	119	0.00
53	P	2,4-Dinitrophenol	0.137	0.168	-22.6	138	0.00
54		Dibenzofuran	1.666	1.590	4.6	110	0.00
55	P	4-Nitrophenol	0.282	0.284	-0.7	115	0.01
56		2,4-Dinitrotoluene	0.419	0.418	0.2	114	0.00
57		Fluorene	1.231	1.221	0.8	111	0.00
58		2,3,4,6-Tetrachlorophenol	0.279	0.288	-3.2	114	0.00
59		Diethylphthalate	1.466	1.425	2.8	114	0.00
60		4-Chlorophenyl-phenylether	0.518	0.522	-0.8	112	0.00
61		4-Nitroaniline	0.363	0.388	-6.9	122	0.00
62		Azobenzene	1.293	1.280	1.0	116	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	122	0.01
64		4,6-Dinitro-2-methylphenol	0.126	0.132	-4.8	129	0.00
65	c	n-Nitrosodiphenylamine	0.715	0.650	9.1	114	0.01
66		4-Bromophenyl-phenylether	0.204	0.201	1.5	122	0.01
67		Hexachlorobenzene	0.227	0.217	4.4	117	0.01
68		Atrazine	0.204	0.201	1.5	122	0.00
69	C	Pentachlorophenol	0.110	0.116	-5.5	119	0.00
70		Phenanthrene	1.087	1.042	4.1	120	0.00
71		Anthracene	1.100	1.012	8.0	114	0.00
72		Carbazole	1.065	1.009	5.3	119	0.00
73		Di-n-butylphthalate	1.326	1.267	4.4	118	0.00
74	C	Fluoranthene	1.041	0.997	4.2	123	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	151#	0.00
76		Benzidine	0.655	0.780	-19.1	175#	0.00
77		Pyrene	1.815	1.596	12.1	131	0.00
78	S	Terphenyl-d14	1.153	0.971	15.8	128	0.00
79		Butylbenzylphthalate	0.862	0.764	11.4	134	0.00
80		Benzo(a)anthracene	1.246	1.237	0.7	148	0.00
81		3,3'-Dichlorobenzidine	0.448	0.486	-8.5	157#	0.00
82		Chrysene	1.227	1.239	-1.0	152#	0.00
83		Bis(2-ethylhexyl)phthalate	1.050	1.003	4.5	141	0.00
84	c	Di-n-octyl phthalate	1.736	1.744	-0.5	154#	0.00
85		Indeno(1,2,3-cd)pyrene	1.113	0.906	18.6	118	0.01
86	I	Perylene-d12	1.000	1.000	0.0	151#	0.00
87		Benzo(b)fluoranthene	1.206	1.182	2.0	150	0.00

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88	Benzo(k)fluoranthene	1.142	1.155	-1.1	152#	0.00
89 C	Benzo(a)pyrene	1.106	1.101	0.5	147	0.00
90	Dibenzo(a,h)anthracene	1.018	0.852	16.3	123	0.01
91	Benzo(g,h,i)perylene	1.090	0.912	16.3	121	0.00

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