Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM071817\
Data File : BM010877.D
Acq On : 18 Jul 2017 19:05
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jul 19 01:01:00 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM071217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 12 14:21:16 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	129	-0.01
2		1,4-Dioxane	0.533	0.458	14.1	110	0.00
3		Pyridine	1.456	1.394	4.3	119	0.00
4		n-Nitrosodimethylamine	0.744	0.691	7.1	120	0.00
5	S	2-Fluorophenol	1.207	1.219	-1.0	130	0.00
6		Aniline	2.085	2.213	-6.1	137	0.00
7	S	Phenol-d6	1.664	1.791	-7.6	139	0.00
8		2-Chlorophenol	1.195	1.189	0.5	126	0.00
9		Benzaldehyde	1.156	1.174	-1.6	130	0.00
10	C	Phenol	1.760	1.794	-1.9	131	0.00
11		bis(2-Chloroethyl)ether	1.355	1.447	-6.8	139	0.00
12		1,3-Dichlorobenzene	1.477	1.455	1.5	129	0.00
13	C	1,4-Dichlorobenzene	1.525	1.507	1.2	130	0.00
14		1,2-Dichlorobenzene	1.453	1.436	1.2	130	0.00
15		Benzyl Alcohol	1.576	1.609	-2.1	133	0.00
16		2,2'-oxybis(1-Chloropropane	1.175	1.274	-8.4	139	0.00
17		2-Methylphenol	1.138	1.276	-12.1	142	0.00
18		Hexachloroethane	0.639	0.651	-1.9	134	-0.01
19	P	n-Nitroso-di-n-propylamine	1.318	1.472	-11.7	147	0.00
20		3+4-Methylphenols	1.581	1.783	-12.8	144	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	140	0.00
22		Acetophenone	0.602	0.591	1.8	141	0.00
23	S	Nitrobenzene-d5	0.520	0.512	1.5	139	0.00
24		Nitrobenzene	0.533	0.531	0.4	142	0.00
25		Isophorone	0.852	0.898	-5.4	148	0.00
26	C	2-Nitrophenol	0.177	0.179	-1.1	144	0.00
27		2,4-Dimethylphenol	0.317	0.304	4.1	135	0.00
28		bis(2-Chloroethoxy)methane	0.478	0.505	-5.6	148	0.00
29	C	2,4-Dichlorophenol	0.358	0.349	2.5	138	0.00
30		1,2,4-Trichlorobenzene	0.437	0.422	3.4	138	-0.01
31		Naphthalene	1.017	1.051	-3.3	146	0.00
32		Benzoic acid	0.248	0.240	3.2	131	0.02
33		4-Chloroaniline	0.410	0.432	-5.4	146	0.00
34	C	Hexachlorobutadiene	0.339	0.320	5.6	133	-0.01
35		Caprolactam	0.095	0.107	-12.6	152#	0.01
36	C	4-Chloro-3-methylphenol	0.433	0.442	-2.1	142	0.00
37		2-Methylnaphthalene	0.730	0.807	-10.5	158#	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	147	0.00
39		1,2,4,5-Tetrachlorobenzene	0.872	0.840	3.7	142	0.00
40	P	Hexachlorocyclopentadiene	0.500	0.514	-2.8	145	0.00
41	S	2,4,6-Tribromophenol	0.306	0.333	-8.8	159#	0.00
42	C	2,4,6-Trichlorophenol	0.537	0.515	4.1	143	0.00
43		2,4,5-Trichlorophenol	0.530	0.550	-3.8	147	0.00
44	S	2-Fluorobiphenyl	1.608	1.646	-2.4	153#	0.00

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Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.742	1.610	7.6	137	0.00
46		2-Chloronaphthalene	1.278	1.300	-1.7	152#	0.00
47		2-Nitroaniline	0.420	0.469	-11.7	159#	0.00
48		Acenaphthylene	1.900	2.081	-9.5	160#	0.00
49		Dimethylphthalate	1.701	1.653	2.8	145	0.00
50		2,6-Dinitrotoluene	0.334	0.369	-10.5	161#	0.00
51	C	Acenaphthene	1.161	1.264	-8.9	162#	0.00
52		3-Nitroaniline	0.303	0.325	-7.3	156#	0.00
53	P	2,4-Dinitrophenol	0.231	0.235	-1.7	147	0.00
54		Dibenzofuran	1.878	1.937	-3.1	153#	0.00
55	P	4-Nitrophenol	0.263	0.253	3.8	138	0.00
56		2,4-Dinitrotoluene	0.463	0.506	-9.3	155#	0.00
57		Fluorene	1.615	1.755	-8.7	161#	0.00
58		2,3,4,6-Tetrachlorophenol	0.492	0.519	-5.5	152#	0.00
59		Diethylphthalate	1.680	1.694	-0.8	148	0.00
60		4-Chlorophenyl-phenylether	0.992	1.016	-2.4	155#	0.00
61		4-Nitroaniline	0.317	0.336	-6.0	150	0.00
62		Azobenzene	1.706	1.827	-7.1	156#	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	145	0.00
64		4,6-Dinitro-2-methylphenol	0.132	0.140	-6.1	152#	0.00
65	c	n-Nitrosodiphenylamine	0.572	0.606	-5.9	153#	0.00
66		4-Bromophenyl-phenylether	0.253	0.273	-7.9	154#	0.00
67		Hexachlorobenzene	0.290	0.290	0.0	147	0.00
68		Atrazine	0.237	0.219	7.6	131	0.00
69	C	Pentachlorophenol	0.186	0.180	3.2	135	0.00
70		Phenanthrene	1.040	1.119	-7.6	156#	0.00
71		Anthracene	1.033	1.136	-10.0	158#	0.00
72		Carbazole	0.910	1.014	-11.4	161#	0.00
73		Di-n-butylphthalate	1.076	1.091	-1.4	142	0.00
74	C	Fluoranthene	1.289	1.324	-2.7	148	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	121	0.00
76		Benzidine	0.535	0.445	16.8	93	0.00
77		Pyrene	1.159	1.384	-19.4	144	0.00
78	S	Terphenyl-d14	1.034	1.005	2.8	119	0.00
79		Butylbenzylphthalate	0.406	0.446	-9.9	129	0.00
80		Benzo(a)anthracene	1.159	1.255	-8.3	131	0.00
81		3,3'-Dichlorobenzidine	0.456	0.458	-0.4	118	0.00
82		Chrysene	1.104	1.167	-5.7	128	0.00
83		Bis(2-ethylhexyl)phthalate	0.591	0.617	-4.4	119	0.00
84	c	Di-n-octyl phthalate	0.969	0.965	0.4	116	0.00
85		Indeno(1,2,3-cd)pyrene	1.266	1.137	10.2	109	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	109	0.00
87		Benzo(b)fluoranthene	1.270	1.415	-11.4	121	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.213	1.319	-8.7	119	0.00
89 C	Benzo(a)pyrene	1.156	1.264	-9.3	119	0.00
90	Dibenzo(a,h)anthracene	1.122	1.166	-3.9	113	-0.01
91	Benzo(g,h,i)perylene	1.103	1.145	-3.8	113	-0.01

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