Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM080217\
Data File : BM011107.D
Acq On : 03 Aug 2017 02:59
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 03 03:52:44 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM072617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 26 17:29:06 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	56	-0.01
2		1,4-Dioxane	0.531	0.530	0.2	58	0.00
3		Pyridine	1.451	1.574	-8.5	61	-0.01
4		n-Nitrosodimethylamine	0.739	0.926	-25.3#	70	0.00
5	S	2-Fluorophenol	1.183	1.139	3.7	55	-0.01
6		Aniline	2.119	2.333	-10.1	62	-0.01
7	S	Phenol-d6	1.681	1.775	-5.6	59	-0.01
8		2-Chlorophenol	1.201	1.239	-3.2	59	-0.01
9		Benzaldehyde	1.088	1.194	-9.7	64	-0.01
10	C	Phenol	1.826	1.881	-3.0	58	-0.01
11		bis(2-Chloroethyl)ether	1.423	1.508	-6.0	60	-0.01
12		1,3-Dichlorobenzene	1.464	1.480	-1.1	58	-0.01
13	C	1,4-Dichlorobenzene	1.501	1.511	-0.7	57	-0.01
14		1,2-Dichlorobenzene	1.435	1.447	-0.8	58	-0.02
15		Benzyl Alcohol	1.613	1.972	-22.3	69	-0.01
16		2,2'-oxybis(1-Chloropropane	1.280	1.490	-16.4	66	0.00
17		2-Methylphenol	1.167	1.288	-10.4	62	-0.02
18		Hexachloroethane	0.654	0.713	-9.0	62	-0.02
19	P	n-Nitroso-di-n-propylamine	1.428	1.844	-29.1#	73	-0.01
20		3+4-Methylphenols	1.622	1.844	-13.7	64	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	65	-0.01
22		Acetophenone	0.600	0.574	4.3	65	-0.01
23	S	Nitrobenzene-d5	0.533	0.583	-9.4	72	-0.01
24		Nitrobenzene	0.552	0.600	-8.7	73	-0.01
25		Isophorone	0.888	0.981	-10.5	74	-0.02
26	C	2-Nitrophenol	0.176	0.174	1.1	64	-0.01
27		2,4-Dimethylphenol	0.309	0.298	3.6	64	-0.01
28		bis(2-Chloroethoxy)methane	0.496	0.508	-2.4	68	-0.01
29	C	2,4-Dichlorophenol	0.348	0.359	-3.2	69	-0.01
30		1,2,4-Trichlorobenzene	0.430	0.431	-0.2	68	-0.01
31		Naphthalene	1.006	1.019	-1.3	68	-0.02
32		Benzoic acid	0.249	0.273	-9.6	73	0.00
33		4-Chloroaniline	0.401	0.430	-7.2	72	-0.01
34	C	Hexachlorobutadiene	0.338	0.352	-4.1	69	-0.02
35		Caprolactam	0.099	0.104	-5.1	74	0.00
36	C	4-Chloro-3-methylphenol	0.449	0.533	-18.7	80	-0.01
37		2-Methylnaphthalene	0.739	0.815	-10.3	73	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	85	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.859	0.784	8.7	79	-0.02
40	P	Hexachlorocyclopentadiene	0.501	0.430	14.2	72	-0.02
41	S	2,4,6-Tribromophenol	0.321	0.379	-18.1	104	-0.01
42	C	2,4,6-Trichlorophenol	0.520	0.498	4.2	80	-0.01
43		2,4,5-Trichlorophenol	0.536	0.524	2.2	83	-0.01
44	S	2-Fluorobiphenyl	1.595	1.503	5.8	81	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.729	1.629	5.8	82	-0.01
46		2-Chloronaphthalene	1.274	1.188	6.8	81	-0.02
47		2-Nitroaniline	0.451	0.550	-22.0	103	-0.01
48		Acenaphthylene	1.902	1.935	-1.7	88	-0.01
49		Dimethylphthalate	1.711	1.773	-3.6	92	-0.01
50		2,6-Dinitrotoluene	0.346	0.373	-7.8	92	-0.01
51	C	Acenaphthene	1.177	1.220	-3.7	90	-0.01
52		3-Nitroaniline	0.300	0.342	-14.0	99	-0.01
53	P	2,4-Dinitrophenol	0.246	0.302	-22.8	108	-0.01
54		Dibenzofuran	1.908	2.039	-6.9	93	-0.01
55	P	4-Nitrophenol	0.264	0.300	-13.6	98	-0.01
56		2,4-Dinitrotoluene	0.486	0.579	-19.1	102	-0.01
57		Fluorene	1.661	1.889	-13.7	101	-0.01
58		2,3,4,6-Tetrachlorophenol	0.502	0.581	-15.7	102	-0.01
59		Diethylphthalate	1.747	1.984	-13.6	100	-0.01
60		4-Chlorophenyl-phenylether	1.003	1.114	-11.1	99	-0.01
61		4-Nitroaniline	0.319	0.359	-12.5	100	0.00
62		Azobenzene	1.870	2.375	-27.0#	111	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	108	-0.01
64		4,6-Dinitro-2-methylphenol	0.135	0.133	1.5	104	-0.01
65	c	n-Nitrosodiphenylamine	0.572	0.538	5.9	104	-0.01
66		4-Bromophenyl-phenylether	0.257	0.249	3.1	107	-0.01
67		Hexachlorobenzene	0.291	0.279	4.1	104	-0.01
68		Atrazine	0.229	0.235	-2.6	109	-0.01
69	C	Pentachlorophenol	0.184	0.196	-6.5	112	-0.01
70		Phenanthrene	1.047	1.086	-3.7	113	-0.01
71		Anthracene	1.044	1.091	-4.5	113	-0.01
72		Carbazole	0.913	0.991	-8.5	119	-0.01
73		Di-n-butylphthalate	1.112	1.243	-11.8	120	-0.01
74	C	Fluoranthene	1.298	1.507	-16.1	127	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	131	0.00
76		Benzidine	0.478	0.461	3.6	123	-0.01
77		Pyrene	1.188	1.150	3.2	128	0.00
78	S	Terphenyl-d14	1.010	1.012	-0.2	130	0.00
79		Butylbenzylphthalate	0.423	0.444	-5.0	133	-0.01
80		Benzo(a)anthracene	1.142	1.168	-2.3	136	0.00
81		3,3'-Dichlorobenzidine	0.448	0.453	-1.1	130	0.00
82		Chrysene	1.096	1.103	-0.6	135	0.00
83		Bis(2-ethylhexyl)phthalate	0.622	0.674	-8.4	140	-0.01
84	c	Di-n-octyl phthalate	1.009	1.068	-5.8	136	0.00
85		Indeno(1,2,3-cd)pyrene	1.135	0.880	22.5	105	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	118	-0.01
87		Benzo(b)fluoranthene	1.284	1.347	-4.9	124	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.230	1.320	-7.3	130	-0.01
89 C	Benzo(a)pyrene	1.162	1.201	-3.4	123	-0.01
90	Dibenzo(a,h)anthracene	1.077	0.934	13.3	106	-0.02
91	Benzo(g,h,i)perylene	1.057	0.901	14.8	103	-0.02

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