Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM081717\
Data File : BM011267.D
Acq On : 17 Aug 2017 15:08
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 18 00:58:38 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM072617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 16 02:25:04 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	40#	-0.01
2		1,4-Dioxane	0.531	0.682	-28.4#	53	-0.01
3		Pyridine	1.451	1.886	-30.0#	52	-0.01
4		n-Nitrosodimethylamine	0.739	1.073	-45.2#	58	-0.02
5	S	2-Fluorophenol	1.183	1.334	-12.8	46#	-0.01
6		Aniline	2.119	2.477	-16.9	47#	-0.02
7	S	Phenol-d6	1.681	1.961	-16.7	46#	-0.01
8		2-Chlorophenol	1.201	1.283	-6.8	43#	-0.01
9		Benzaldehyde	1.088	1.254	-15.3	48#	-0.02
10	C	Phenol	1.826	2.165	-18.6	48#	-0.01
11		bis(2-Chloroethyl)ether	1.423	1.641	-15.3	47#	-0.01
12		1,3-Dichlorobenzene	1.464	1.582	-8.1	45#	-0.02
13	C	1,4-Dichlorobenzene	1.501	1.663	-10.8	45#	-0.01
14		1,2-Dichlorobenzene	1.435	1.502	-4.7	43#	-0.02
15		Benzyl Alcohol	1.613	2.030	-25.9#	51	-0.01
16		2,2'-oxybis(1-Chloropropane	1.280	1.621	-26.6#	52	0.00
17		2-Methylphenol	1.167	1.295	-11.0	45#	-0.02
18		Hexachloroethane	0.654	0.777	-18.8	48#	-0.01
19	P	n-Nitroso-di-n-propylamine	1.428	1.971	-38.0#	56	-0.01
20		3+4-Methylphenols	1.622	1.814	-11.8	45#	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	39#	-0.02
22		Acetophenone	0.600	0.674	-12.3	45#	-0.02
23	S	Nitrobenzene-d5	0.533	0.708	-32.8#	52	-0.02
24		Nitrobenzene	0.552	0.757	-37.1#	54	-0.02
25		Isophorone	0.888	1.145	-28.9#	52	-0.02
26	C	2-Nitrophenol	0.176	0.197	-11.9	43#	-0.02
27		2,4-Dimethylphenol	0.309	0.346	-12.0	44#	-0.01
28		bis(2-Chloroethoxy)methane	0.496	0.586	-18.1	47#	-0.01
29	C	2,4-Dichlorophenol	0.348	0.380	-9.2	43#	-0.02
30		1,2,4-Trichlorobenzene	0.430	0.480	-11.6	45#	-0.02
31		Naphthalene	1.006	1.076	-7.0	43#	-0.02
32		Benzoic acid	0.249	0.244	2.0	39#	-0.03
33		4-Chloroaniline	0.401	0.411	-2.5	41#	-0.01
34	C	Hexachlorobutadiene	0.338	0.413	-22.2#	48#	-0.02
35		Caprolactam	0.099	0.118	-19.2	50	-0.02
36	C	4-Chloro-3-methylphenol	0.449	0.538	-19.8	48#	-0.02
37		2-Methylnaphthalene	0.739	0.817	-10.6	44#	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	45#	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.859	0.925	-7.7	49#	-0.02
40	P	Hexachlorocyclopentadiene	0.501	0.498	0.6	44#	-0.02
41	S	2,4,6-Tribromophenol	0.321	0.339	-5.6	49#	-0.02
42	C	2,4,6-Trichlorophenol	0.520	0.560	-7.7	47#	-0.01
43		2,4,5-Trichlorophenol	0.536	0.559	-4.3	47#	-0.02
44	S	2-Fluorobiphenyl	1.595	1.658	-3.9	47#	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.729	1.752	-1.3	46#	-0.01
46		2-Chloronaphthalene	1.274	1.299	-2.0	46#	-0.02
47		2-Nitroaniline	0.451	0.623	-38.1#	61	-0.02
48		Acenaphthylene	1.902	1.986	-4.4	47#	-0.02
49		Dimethylphthalate	1.711	1.770	-3.4	48#	-0.02
50		2,6-Dinitrotoluene	0.346	0.371	-7.2	48#	-0.01
51	C	Acenaphthene	1.177	1.254	-6.5	49#	-0.01
52		3-Nitroaniline	0.300	0.303	-1.0	46#	-0.02
53	P	2,4-Dinitrophenol	0.246	0.271	-10.2	51	-0.02
54		Dibenzofuran	1.908	2.044	-7.1	49#	-0.02
55	P	4-Nitrophenol	0.264	0.223	15.5	38#	0.00
56		2,4-Dinitrotoluene	0.486	0.540	-11.1	50#	-0.02
57		Fluorene	1.661	1.782	-7.3	50#	-0.02
58		2,3,4,6-Tetrachlorophenol	0.502	0.578	-15.1	53	-0.01
59		Diethylphthalate	1.747	1.916	-9.7	51	-0.01
60		4-Chlorophenyl-phenylether	1.003	1.109	-10.6	52	-0.02
61		4-Nitroaniline	0.319	0.269	15.7	39#	-0.02
62		Azobenzene	1.870	2.513	-34.4#	62	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	49#	-0.01
64		4,6-Dinitro-2-methylphenol	0.135	0.152	-12.6	54	-0.01
65	c	n-Nitrosodiphenylamine	0.572	0.603	-5.4	52	-0.01
66		4-Bromophenyl-phenylether	0.257	0.272	-5.8	52	-0.01
67		Hexachlorobenzene	0.291	0.299	-2.7	50	-0.02
68		Atrazine	0.229	0.239	-4.4	50#	-0.01
69	C	Pentachlorophenol	0.184	0.200	-8.7	52	-0.01
70		Phenanthrene	1.047	1.180	-12.7	55	-0.01
71		Anthracene	1.044	1.153	-10.4	54	-0.02
72		Carbazole	0.913	1.040	-13.9	57	-0.02
73		Di-n-butylphthalate	1.112	1.250	-12.4	54	-0.02
74	C	Fluoranthene	1.298	1.565	-20.6#	60	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	58	-0.02
76		Benzidine	0.478	0.281	41.2#	33#	-0.01
77		Pyrene	1.188	1.216	-2.4	61	-0.02
78	S	Terphenyl-d14	1.010	1.049	-3.9	60	-0.01
79		Butylbenzylphthalate	0.423	0.454	-7.3	61	-0.01
80		Benzo(a)anthracene	1.142	1.252	-9.6	65	-0.01
81		3,3'-Dichlorobenzidine	0.448	0.481	-7.4	62	-0.02
82		Chrysene	1.096	1.206	-10.0	66	-0.01
83		Bis(2-ethylhexyl)phthalate	0.622	0.656	-5.5	61	-0.02
84	c	Di-n-octyl phthalate	1.009	1.116	-10.6	63	-0.01
85		Indeno(1,2,3-cd)pyrene	1.135	1.246	-9.8	67	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	61	-0.02
87		Benzo(b)fluoranthene	1.284	1.405	-9.4	67	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.230	1.287	-4.6	65	-0.02
89 C	Benzo(a)pyrene	1.162	1.265	-8.9	67	-0.02
90	Dibenzo(a,h)anthracene	1.077	1.131	-5.0	67	-0.03
91	Benzo(g,h,i)perylene	1.057	1.135	-7.4	67	-0.03

(#) = Out of Range SPCC's out = 0 CCC's out = 2