Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM081617\
Data File : BM011259.D
Acq On : 16 Aug 2017 18:07
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 17 03:53:11 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	48#	-0.01
2		1,4-Dioxane	0.531	0.639	-20.3	61	-0.01
3		Pyridine	1.451	1.752	-20.7	58	-0.01
4		n-Nitrosodimethylamine	0.739	0.992	-34.2#	65	-0.01
5	S	2-Fluorophenol	1.183	1.311	-10.8	55	-0.01
6		Aniline	2.119	2.299	-8.5	53	-0.02
7	S	Phenol-d6	1.681	1.926	-14.6	55	-0.01
8		2-Chlorophenol	1.201	1.254	-4.4	51	-0.01
9		Benzaldehyde	1.088	1.160	-6.6	54	-0.01
10	C	Phenol	1.826	2.113	-15.7	56	-0.01
11		bis(2-Chloroethyl)ether	1.423	1.589	-11.7	55	-0.01
12		1,3-Dichlorobenzene	1.464	1.544	-5.5	53	-0.02
13	C	1,4-Dichlorobenzene	1.501	1.579	-5.2	51	-0.01
14		1,2-Dichlorobenzene	1.435	1.434	0.1	50#	-0.01
15		Benzyl Alcohol	1.613	1.886	-16.9	57	-0.01
16		2,2'-oxybis(1-Chloropropane	1.280	1.587	-24.0	61	0.00
17		2-Methylphenol	1.167	1.329	-13.9	55	-0.01
18		Hexachloroethane	0.654	0.723	-10.6	55	-0.01
19	P	n-Nitroso-di-n-propylamine	1.428	1.888	-32.2#	64	-0.01
20		3+4-Methylphenols	1.622	1.801	-11.0	54	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	45#	-0.01
22		Acetophenone	0.600	0.692	-15.3	54	-0.02
23	S	Nitrobenzene-d5	0.533	0.702	-31.7#	60	-0.02
24		Nitrobenzene	0.552	0.723	-31.0#	61	-0.02
25		Isophorone	0.888	1.151	-29.6#	61	-0.02
26	C	2-Nitrophenol	0.176	0.192	-9.1	49#	-0.02
27		2,4-Dimethylphenol	0.309	0.347	-12.3	52	-0.01
28		bis(2-Chloroethoxy)methane	0.496	0.602	-21.4	56	-0.01
29	C	2,4-Dichlorophenol	0.348	0.388	-11.5	51	-0.02
30		1,2,4-Trichlorobenzene	0.430	0.487	-13.3	53	-0.01
31		Naphthalene	1.006	1.061	-5.5	49#	-0.02
32		Benzoic acid	0.249	0.268	-7.6	50#	-0.02
33		4-Chloroaniline	0.401	0.396	1.2	46#	-0.01
34	C	Hexachlorobutadiene	0.338	0.397	-17.5	54	-0.02
35		Caprolactam	0.099	0.125	-26.3#	62	-0.01
36	C	4-Chloro-3-methylphenol	0.449	0.547	-21.8#	57	-0.02
37		2-Methylnaphthalene	0.739	0.816	-10.4	51	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	50#	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.859	0.930	-8.3	55	-0.02
40	P	Hexachlorocyclopentadiene	0.501	0.527	-5.2	52	-0.02
41	S	2,4,6-Tribromophenol	0.321	0.353	-10.0	57	-0.02
42	C	2,4,6-Trichlorophenol	0.520	0.575	-10.6	54	-0.01
43		2,4,5-Trichlorophenol	0.536	0.600	-11.9	56	-0.02
44	S	2-Fluorobiphenyl	1.595	1.695	-6.3	54	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.729	1.819	-5.2	54	-0.01
46		2-Chloronaphthalene	1.274	1.348	-5.8	54	-0.01
47		2-Nitroaniline	0.451	0.659	-46.1#	73	-0.02
48		Acenaphthylene	1.902	2.102	-10.5	56	-0.02
49		Dimethylphthalate	1.711	1.864	-8.9	57	-0.01
50		2,6-Dinitrotoluene	0.346	0.391	-13.0	57	-0.01
51	C	Acenaphthene	1.177	1.296	-10.1	57	-0.01
52		3-Nitroaniline	0.300	0.303	-1.0	52	-0.02
53	P	2,4-Dinitrophenol	0.246	0.297	-20.7	63	-0.02
54		Dibenzofuran	1.908	2.120	-11.1	57	-0.01
55	P	4-Nitrophenol	0.264	0.279	-5.7	54	-0.01
56		2,4-Dinitrotoluene	0.486	0.580	-19.3	60	-0.01
57		Fluorene	1.661	1.882	-13.3	59	-0.02
58		2,3,4,6-Tetrachlorophenol	0.502	0.577	-14.9	60	-0.01
59		Diethylphthalate	1.747	2.054	-17.6	61	-0.01
60		4-Chlorophenyl-phenylether	1.003	1.107	-10.4	58	-0.01
61		4-Nitroaniline	0.319	0.294	7.8	48#	-0.02
62		Azobenzene	1.870	2.539	-35.8#	70	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	57	-0.01
64		4,6-Dinitro-2-methylphenol	0.135	0.156	-15.6	65	-0.01
65	c	n-Nitrosodiphenylamine	0.572	0.590	-3.1	60	-0.01
66		4-Bromophenyl-phenylether	0.257	0.263	-2.3	60	-0.01
67		Hexachlorobenzene	0.291	0.295	-1.4	59	-0.02
68		Atrazine	0.229	0.235	-2.6	58	0.00
69	C	Pentachlorophenol	0.184	0.198	-7.6	60	-0.01
70		Phenanthrene	1.047	1.126	-7.5	62	-0.01
71		Anthracene	1.044	1.128	-8.0	62	-0.01
72		Carbazole	0.913	1.018	-11.5	65	-0.01
73		Di-n-butylphthalate	1.112	1.279	-15.0	65	-0.01
74	C	Fluoranthene	1.298	1.523	-17.3	68	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	68	-0.01
76		Benzidine	0.478	0.126	73.6#	18#	0.00
77		Pyrene	1.188	1.212	-2.0	71	-0.02
78	S	Terphenyl-d14	1.010	1.026	-1.6	69	-0.01
79		Butylbenzylphthalate	0.423	0.450	-6.4	71	-0.01
80		Benzo(a)anthracene	1.142	1.220	-6.8	74	-0.01
81		3,3'-Dichlorobenzidine	0.448	0.443	1.1	67	-0.01
82		Chrysene	1.096	1.163	-6.1	74	-0.01
83		Bis(2-ethylhexyl)phthalate	0.622	0.650	-4.5	71	-0.01
84	c	Di-n-octyl phthalate	1.009	1.087	-7.7	72	0.00
85		Indeno(1,2,3-cd)pyrene	1.135	1.178	-3.8	74	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	69	-0.02
87		Benzo(b)fluoranthene	1.284	1.338	-4.2	72	-0.02

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88	Benzo(k)fluoranthene	1.230	1.338	-8.8	77	-0.02
89 C	Benzo(a)pyrene	1.162	1.242	-6.9	75	-0.02
90	Dibenzo(a,h)anthracene	1.077	1.123	-4.3	75	-0.03
91	Benzo(g,h,i)perylene	1.057	1.111	-5.1	75	-0.03

(#) = Out of Range SPCC's out = 0 CCC's out = 1