Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM081717\
Data File : BM011286.D
Acq On : 18 Aug 2017 03:48
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 18 10:41:14 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	36#	-0.02
2		1,4-Dioxane	0.531	0.680	-28.1#	47#	-0.01
3		Pyridine	1.451	1.752	-20.7	43#	-0.01
4		n-Nitrosodimethylamine	0.739	1.149	-55.5#	55	-0.02
5	S	2-Fluorophenol	1.183	1.293	-9.3	40#	-0.02
6		Aniline	2.119	2.516	-18.7	43#	-0.02
7	S	Phenol-d6	1.681	1.992	-18.5	42#	-0.02
8		2-Chlorophenol	1.201	1.297	-8.0	39#	-0.01
9		Benzaldehyde	1.088	1.270	-16.7	43#	-0.02
10	C	Phenol	1.826	2.203	-20.6#	43#	-0.01
11		bis(2-Chloroethyl)ether	1.423	1.611	-13.2	41#	-0.02
12		1,3-Dichlorobenzene	1.464	1.588	-8.5	40#	-0.02
13	C	1,4-Dichlorobenzene	1.501	1.607	-7.1	38#	-0.01
14		1,2-Dichlorobenzene	1.435	1.570	-9.4	40#	-0.02
15		Benzyl Alcohol	1.613	2.088	-29.4#	47#	-0.02
16		2,2'-oxybis(1-Chloropropane	1.280	1.706	-33.3#	49#	-0.01
17		2-Methylphenol	1.167	1.314	-12.6	40#	-0.02
18		Hexachloroethane	0.654	0.795	-21.6	44#	-0.02
19	P	n-Nitroso-di-n-propylamine	1.428	2.056	-44.0#	52	-0.01
20		3+4-Methylphenols	1.622	1.861	-14.7	41#	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	36#	-0.02
22		Acetophenone	0.600	0.681	-13.5	42#	-0.02
23	S	Nitrobenzene-d5	0.533	0.721	-35.3#	49#	-0.02
24		Nitrobenzene	0.552	0.763	-38.2#	50	-0.02
25		Isophorone	0.888	1.179	-32.8#	49#	-0.02
26	C	2-Nitrophenol	0.176	0.191	-8.5	39#	-0.02
27		2,4-Dimethylphenol	0.309	0.341	-10.4	40#	-0.02
28		bis(2-Chloroethoxy)methane	0.496	0.602	-21.4	44#	-0.02
29	C	2,4-Dichlorophenol	0.348	0.391	-12.4	41#	-0.02
30		1,2,4-Trichlorobenzene	0.430	0.492	-14.4	42#	-0.02
31		Naphthalene	1.006	1.108	-10.1	41#	-0.02
32		Benzoic acid	0.249	0.253	-1.6	37#	-0.03
33		4-Chloroaniline	0.401	0.421	-5.0	38#	-0.02
34	C	Hexachlorobutadiene	0.338	0.421	-24.6#	45#	-0.02
35		Caprolactam	0.099	0.116	-17.2	45#	-0.02
36	C	4-Chloro-3-methylphenol	0.449	0.572	-27.4#	47#	-0.02
37		2-Methylnaphthalene	0.739	0.841	-13.8	41#	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	43#	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.859	0.933	-8.6	47#	-0.02
40	P	Hexachlorocyclopentadiene	0.501	0.504	-0.6	43#	-0.02
41	S	2,4,6-Tribromophenol	0.321	0.365	-13.7	50	-0.02
42	C	2,4,6-Trichlorophenol	0.520	0.571	-9.8	46#	-0.02
43		2,4,5-Trichlorophenol	0.536	0.575	-7.3	46#	-0.02
44	S	2-Fluorobiphenyl	1.595	1.662	-4.2	45#	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.729	1.787	-3.4	45#	-0.02
46		2-Chloronaphthalene	1.274	1.343	-5.4	46#	-0.02
47		2-Nitroaniline	0.451	0.643	-42.6#	61	-0.02
48		Acenaphthylene	1.902	2.022	-6.3	46#	-0.02
49		Dimethylphthalate	1.711	1.849	-8.1	48#	-0.02
50		2,6-Dinitrotoluene	0.346	0.391	-13.0	49#	-0.01
51	C	Acenaphthene	1.177	1.281	-8.8	48#	-0.02
52		3-Nitroaniline	0.300	0.319	-6.3	47#	-0.02
53	P	2,4-Dinitrophenol	0.246	0.284	-15.4	51	-0.02
54		Dibenzofuran	1.908	2.125	-11.4	49#	-0.02
55	P	4-Nitrophenol	0.264	0.221	16.3	37#	-0.02
56		2,4-Dinitrotoluene	0.486	0.563	-15.8	50	-0.02
57		Fluorene	1.661	1.897	-14.2	51	-0.02
58		2,3,4,6-Tetrachlorophenol	0.502	0.591	-17.7	52	-0.02
59		Diethylphthalate	1.747	1.989	-13.9	51	-0.02
60		4-Chlorophenyl-phenylether	1.003	1.178	-17.4	53	-0.02
61		4-Nitroaniline	0.319	0.295	7.5	41#	-0.02
62		Azobenzene	1.870	2.698	-44.3#	64	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	51	-0.02
64		4,6-Dinitro-2-methylphenol	0.135	0.146	-8.1	54	-0.01
65	c	n-Nitrosodiphenylamine	0.572	0.587	-2.6	54	-0.01
66		4-Bromophenyl-phenylether	0.257	0.259	-0.8	53	-0.02
67		Hexachlorobenzene	0.291	0.296	-1.7	52	-0.02
68		Atrazine	0.229	0.229	0.0	50	-0.01
69	C	Pentachlorophenol	0.184	0.189	-2.7	51	-0.02
70		Phenanthrene	1.047	1.121	-7.1	55	-0.01
71		Anthracene	1.044	1.143	-9.5	56	-0.02
72		Carbazole	0.913	1.011	-10.7	58	-0.02
73		Di-n-butylphthalate	1.112	1.201	-8.0	55	-0.02
74	C	Fluoranthene	1.298	1.532	-18.0	61	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	63	-0.02
76		Benzidine	0.478	0.227	52.5#	29#	-0.02
77		Pyrene	1.188	1.185	0.3	64	-0.02
78	S	Terphenyl-d14	1.010	1.020	-1.0	63	-0.02
79		Butylbenzylphthalate	0.423	0.418	1.2	61	-0.01
80		Benzo(a)anthracene	1.142	1.228	-7.5	69	-0.01
81		3,3'-Dichlorobenzidine	0.448	0.457	-2.0	63	-0.02
82		Chrysene	1.096	1.163	-6.1	69	-0.02
83		Bis(2-ethylhexyl)phthalate	0.622	0.621	0.2	62	-0.02
84	c	Di-n-octyl phthalate	1.009	1.045	-3.6	64	-0.01
85		Indeno(1,2,3-cd)pyrene	1.135	1.208	-6.4	70	-0.04
86	I	Perylene-d12	1.000	1.000	0.0	65	-0.02
87		Benzo(b)fluoranthene	1.284	1.346	-4.8	68	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.230	1.364	-10.9	74	-0.02
89 C	Benzo(a)pyrene	1.162	1.268	-9.1	72	-0.02
90	Dibenzo(a,h)anthracene	1.077	1.130	-4.9	71	-0.04
91	Benzo(g,h,i)perylene	1.057	1.120	-6.0	71	-0.04

(#) = Out of Range SPCC's out = 0 CCC's out = 3