Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA N\Data\BN041818\
Data File : BN000759.D
Acq On : 19 Apr 2018 02:35
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Apr 19 05:59:24 2018
Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN041818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 18 16:58:01 2018
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	72	0.00
2		1,4-Dioxane	0.618	0.627	-1.5	74	0.00
3		Pyridine	1.627	1.725	-6.0	70	0.00
4		n-Nitrosodimethylamine	0.447	0.465	-4.0	69	0.00
5	S	2-Fluorophenol	1.346	1.407	-4.5	72	0.00
6		Aniline	2.434	2.452	-0.7	70	0.00
7	S	Phenol-d6	1.835	1.876	-2.2	70	0.00
8		2-Chlorophenol	1.392	1.419	-1.9	70	0.00
9		Benzaldehyde	0.846	0.858	-1.4	70	0.00
10	C	Phenol	1.934	1.958	-1.2	70	0.00
11		bis(2-Chloroethyl)ether	1.582	1.578	0.3	70	0.00
12		1,3-Dichlorobenzene	1.617	1.661	-2.7	73	0.00
13	C	1,4-Dichlorobenzene	1.642	1.683	-2.5	72	0.00
14		1,2-Dichlorobenzene	1.567	1.587	-1.3	72	0.00
15		Benzyl Alcohol	1.262	1.291	-2.3	68	0.00
16		2,2'-oxybis(1-Chloropropane	1.680	1.649	1.8	69	0.00
17		2-Methylphenol	1.300	1.320	-1.5	70	0.00
18		Hexachloroethane	0.605	0.623	-3.0	71	0.00
19	P	n-Nitroso-di-n-propylamine	1.141	1.145	-0.4	67	0.00
20		3+4-Methylphenols	1.787	1.804	-1.0	68	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	69	0.00
22		Acetophenone	0.573	0.583	-1.7	68	0.00
23	S	Nitrobenzene-d5	0.384	0.402	-4.7	67	0.00
24		Nitrobenzene	0.411	0.425	-3.4	67	0.00
25		Isophorone	0.695	0.711	-2.3	66	0.00
26	C	2-Nitrophenol	0.169	0.171	-1.2	68	0.00
27		2,4-Dimethylphenol	0.275	0.282	-2.5	68	0.00
28		bis(2-Chloroethoxy)methane	0.453	0.455	-0.4	66	0.00
29	C	2,4-Dichlorophenol	0.271	0.283	-4.4	69	0.00
30		1,2,4-Trichlorobenzene	0.327	0.335	-2.4	70	0.00
31		Naphthalene	1.075	1.095	-1.9	71	0.00
32		Benzoic acid	0.238	0.230	3.4	62	-0.02
33		4-Chloroaniline	0.439	0.452	-3.0	68	0.00
34	C	Hexachlorobutadiene	0.177	0.184	-4.0	73	0.00
35		Caprolactam	0.094	0.097	-3.2	67	-0.01
36	C	4-Chloro-3-methylphenol	0.320	0.336	-5.0	68	-0.01
37		2-Methylnaphthalene	0.704	0.715	-1.6	70	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	69	0.00
39		1,2,4,5-Tetrachlorobenzene	0.620	0.641	-3.4	71	0.00
40	P	Hexachlorocyclopentadiene	0.364	0.381	-4.7	72	0.00
41	S	2,4,6-Tribromophenol	0.201	0.221	-10.0	71	0.00
42	C	2,4,6-Trichlorophenol	0.376	0.400	-6.4	70	0.00
43		2,4,5-Trichlorophenol	0.435	0.464	-6.7	71	0.00
44	S	2-Fluorobiphenyl	1.449	1.478	-2.0	72	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.790	1.819	-1.6	69	0.00
46		2-Chloronaphthalene	1.342	1.369	-2.0	71	0.00
47		2-Nitroaniline	0.389	0.405	-4.1	69	0.00
48		Acenaphthylene	2.128	2.215	-4.1	70	0.00
49		Dimethylphthalate	1.598	1.645	-2.9	70	-0.01
50		2,6-Dinitrotoluene	0.339	0.348	-2.7	68	0.00
51	C	Acenaphthene	1.297	1.307	-0.8	69	0.00
52		3-Nitroaniline	0.385	0.405	-5.2	69	0.00
53	P	2,4-Dinitrophenol	0.154	0.165	-7.1	71	0.00
54		Dibenzofuran	1.888	1.909	-1.1	68	0.00
55	P	4-Nitrophenol	0.286	0.309	-8.0	72	0.00
56		2,4-Dinitrotoluene	0.443	0.468	-5.6	70	0.00
57		Fluorene	1.493	1.536	-2.9	69	0.00
58		2,3,4,6-Tetrachlorophenol	0.334	0.354	-6.0	69	0.00
59		Diethylphthalate	1.584	1.648	-4.0	68	0.00
60		4-Chlorophenyl-phenylether	0.698	0.710	-1.7	69	0.00
61		4-Nitroaniline	0.379	0.410	-8.2	71	0.00
62		Azobenzene	1.696	1.755	-3.5	67	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	75	0.00
64		4,6-Dinitro-2-methylphenol	0.124	0.129	-4.0	72	0.00
65	c	n-Nitrosodiphenylamine	0.671	0.674	-0.4	69	0.00
66		4-Bromophenyl-phenylether	0.203	0.206	-1.5	71	0.00
67		Hexachlorobenzene	0.235	0.236	-0.4	71	0.00
68		Atrazine	0.182	0.201	-10.4	72	0.00
69	C	Pentachlorophenol	0.152	0.157	-3.3	72	0.00
70		Phenanthrene	1.158	1.165	-0.6	73	0.00
71		Anthracene	1.134	1.159	-2.2	74	0.00
72		Carbazole	1.066	1.127	-5.7	77	0.00
73		Di-n-butylphthalate	1.215	1.322	-8.8	77	0.00
74	C	Fluoranthene	1.167	1.257	-7.7	80	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	89	0.00
76		Benzidine	0.542	0.611	-12.7	98	0.00
77		Pyrene	1.441	1.393	3.3	79	0.00
78	S	Terphenyl-d14	0.928	0.899	3.1	81	0.00
79		Butylbenzylphthalate	0.607	0.614	-1.2	85	0.00
80		Benzo(a)anthracene	1.260	1.299	-3.1	91	0.00
81		3,3'-Dichlorobenzidine	0.397	0.438	-10.3	95	0.00
82		Chrysene	1.220	1.243	-1.9	90	0.00
83		Bis(2-ethylhexyl)phthalate	0.927	0.949	-2.4	88	0.00
84	c	Di-n-octyl phthalate	1.326	1.431	-7.9	93	0.00
85		Indeno(1,2,3-cd)pyrene	1.049	1.233	-17.5	111	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	107	0.00
87		Benzo(b)fluoranthene	1.215	1.205	0.8	104	0.00

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88	Benzo(k)fluoranthene	1.210	1.225	-1.2	105	0.00
89 C	Benzo(a)pyrene	1.109	1.152	-3.9	105	0.00
90	Dibenzo(a,h)anthracene	1.093	1.181	-8.1	111	0.00
91	Benzo(g,h,i)perylene	1.067	1.166	-9.3	111	-0.01

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