Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG092618\
Data File : BG036981.D
Acq On : 26 Sep 2018 14:39
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 26 15:37:20 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 24 10:41:23 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	121	0.00
2		1,4-Dioxane	0.477	0.580	-21.6	137	0.00
3		Pyridine	1.378	1.530	-11.0	127	0.00
4		n-Nitrosodimethylamine	0.612	0.733	-19.8	141	0.00
5	S	2-Fluorophenol	1.043	1.196	-14.7	134	0.00
6		Aniline	2.094	1.983	5.3	112	0.00
7	S	Phenol-d6	1.613	1.638	-1.5	118	0.00
8		2-Chlorophenol	1.211	1.263	-4.3	121	0.00
9		Benzaldehyde	1.066	1.045	2.0	116	0.00
10	C	Phenol	1.665	1.713	-2.9	120	0.00
11		bis(2-Chloroethyl)ether	1.540	1.462	5.1	118	0.00
12		1,3-Dichlorobenzene	1.485	1.519	-2.3	120	0.00
13	C	1,4-Dichlorobenzene	1.519	1.500	1.3	118	0.00
14		1,2-Dichlorobenzene	1.472	1.432	2.7	118	0.00
15		Benzyl Alcohol	1.309	1.231	6.0	111	0.00
16		2,2'-oxybis(1-Chloropropane	2.957	2.904	1.8	116	0.00
17		2-Methylphenol	1.160	1.110	4.3	114	0.00
18		Hexachloroethane	0.559	0.589	-5.4	124	0.00
19	P	n-Nitroso-di-n-propylamine	1.342	1.182	11.9	105	0.00
20		3+4-Methylphenols	1.614	1.518	5.9	110	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
22		Acetophenone	0.470	0.484	-3.0	108	0.00
23	S	Nitrobenzene-d5	0.373	0.405	-8.6	110	0.00
24		Nitrobenzene	0.399	0.422	-5.8	108	0.00
25		Isophorone	0.682	0.690	-1.2	104	0.00
26	C	2-Nitrophenol	0.159	0.182	-14.5	114	0.00
27		2,4-Dimethylphenol	0.248	0.247	0.4	101	0.00
28		bis(2-Chloroethoxy)methane	0.421	0.418	0.7	101	0.00
29	C	2,4-Dichlorophenol	0.287	0.310	-8.0	107	0.00
30		1,2,4-Trichlorobenzene	0.359	0.373	-3.9	106	0.00
31		Naphthalene	0.952	0.954	-0.2	104	0.00
32		Benzoic acid	0.175	0.175	0.0	102	0.00
33		4-Chloroaniline	0.433	0.407	6.0	97	0.00
34	C	Hexachlorobutadiene	0.248	0.265	-6.9	109	0.00
35		Caprolactam	0.118	0.117	0.8	99	0.00
36	C	4-Chloro-3-methylphenol	0.343	0.359	-4.7	102	0.00
37		2-Methylnaphthalene	0.725	0.707	2.5	102	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	96	0.00
39		1,2,4,5-Tetrachlorobenzene	0.698	0.731	-4.7	100	0.00
40	P	Hexachlorocyclopentadiene	0.359	0.351	2.2	89	0.00
41	S	2,4,6-Tribromophenol	0.239	0.258	-7.9	100	0.00
42	C	2,4,6-Trichlorophenol	0.387	0.438	-13.2	105	0.00
43		2,4,5-Trichlorophenol	0.413	0.477	-15.5	105	0.00
44	S	2-Fluorobiphenyl	1.343	1.382	-2.9	97	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.532	1.568	-2.3	98	0.00
46		2-Chloronaphthalene	1.205	1.224	-1.6	98	0.00
47		2-Nitroaniline	0.417	0.463	-11.0	102	0.00
48		Acenaphthylene	1.888	1.921	-1.7	97	0.00
49		Dimethylphthalate	1.610	1.576	2.1	93	0.00
50		2,6-Dinitrotoluene	0.351	0.353	-0.6	95	0.00
51	C	Acenaphthene	1.141	1.139	0.2	96	0.00
52		3-Nitroaniline	0.370	0.391	-5.7	98	0.00
53	P	2,4-Dinitrophenol	0.136	0.127	6.6	86	0.00
54		Dibenzofuran	1.930	1.961	-1.6	97	0.00
55	P	4-Nitrophenol	0.236	0.244	-3.4	94	0.00
56		2,4-Dinitrotoluene	0.490	0.510	-4.1	98	0.00
57		Fluorene	1.442	1.457	-1.0	97	0.00
58		2,3,4,6-Tetrachlorophenol	0.419	0.451	-7.6	98	0.00
59		Diethylphthalate	1.624	1.608	1.0	95	0.00
60		4-Chlorophenyl-phenylether	0.913	0.916	-0.3	95	0.00
61		4-Nitroaniline	0.389	0.399	-2.6	96	0.00
62		Azobenzene	1.560	1.653	-6.0	101	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	96	0.00
64		4,6-Dinitro-2-methylphenol	0.105	0.106	-1.0	96	0.00
65	c	n-Nitrosodiphenylamine	0.552	0.563	-2.0	96	0.00
66		4-Bromophenyl-phenylether	0.227	0.228	-0.4	94	0.00
67		Hexachlorobenzene	0.234	0.236	-0.9	95	0.00
68		Atrazine	0.224	0.225	-0.4	94	0.00
69	C	Pentachlorophenol	0.148	0.158	-6.8	98	0.00
70		Phenanthrene	0.964	0.980	-1.7	97	0.00
71		Anthracene	0.953	0.997	-4.6	100	0.00
72		Carbazole	0.929	0.971	-4.5	99	0.00
73		Di-n-butylphthalate	1.032	1.077	-4.4	99	0.00
74	C	Fluoranthene	1.160	1.204	-3.8	99	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
76		Benzidine	0.605	0.631	-4.3	109	0.00
77		Pyrene	1.200	1.153	3.9	99	0.00
78	S	Terphenyl-d14	0.761	0.728	4.3	100	0.00
79		Butylbenzylphthalate	0.478	0.481	-0.6	104	0.00
80		Benzo(a)anthracene	1.167	1.156	0.9	105	0.00
81		3,3'-Dichlorobenzidine	0.444	0.443	0.2	100	0.00
82		Chrysene	1.128	1.110	1.6	103	0.00
83		Bis(2-ethylhexyl)phthalate	0.668	0.672	-0.6	106	0.00
84	c	Di-n-octyl phthalate	1.100	1.118	-1.6	105	-0.02
85		Indeno(1,2,3-cd)pyrene	1.211	1.256	-3.7	105	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	104	-0.02
87		Benzo(b)fluoranthene	1.066	1.090	-2.3	104	0.00

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88	Benzo(k)fluoranthene	1.069	1.058	1.0	104	-0.02
89 C	Benzo(a)pyrene	1.030	1.036	-0.6	103	-0.02
90	Dibenzo(a,h)anthracene	0.966	0.988	-2.3	105	-0.02
91	Benzo(g,h,i)perylene	0.969	1.003	-3.5	104	-0.02

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