Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG082818\
Data File : BG036461.D
Acq On : 28 Aug 2018 17:01
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 28 17:46:01 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 23 12:07:02 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	123	0.00
2		1,4-Dioxane	0.588	0.522	11.2	114	0.00
3		Pyridine	1.652	1.520	8.0	112	0.00
4		n-Nitrosodimethylamine	0.736	0.615	16.4	103	0.00
5	S	2-Fluorophenol	1.261	1.255	0.5	123	0.00
6		Aniline	2.291	2.075	9.4	112	0.00
7	S	Phenol-d6	1.805	1.789	0.9	122	0.00
8		2-Chlorophenol	1.367	1.311	4.1	118	0.00
9		Benzaldehyde	1.243	1.076	13.4	114	0.00
10	C	Phenol	1.889	1.867	1.2	122	0.00
11		bis(2-Chloroethyl)ether	1.498	1.361	9.1	114	0.00
12		1,3-Dichlorobenzene	1.561	1.450	7.1	117	0.00
13	C	1,4-Dichlorobenzene	1.576	1.473	6.5	117	0.00
14		1,2-Dichlorobenzene	1.505	1.380	8.3	116	0.00
15		Benzyl Alcohol	1.455	1.353	7.0	113	0.00
16		2,2'-oxybis(1-Chloropropane	3.020	2.504	17.1	104	0.00
17		2-Methylphenol	1.254	1.199	4.4	119	0.00
18		Hexachloroethane	0.565	0.522	7.6	114	0.00
19	P	n-Nitroso-di-n-propylamine	1.349	1.148	14.9	107	0.00
20		3+4-Methylphenols	1.751	1.684	3.8	119	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	126	0.00
22		Acetophenone	0.525	0.468	10.9	112	-0.01
23	S	Nitrobenzene-d5	0.369	0.368	0.3	122	0.00
24		Nitrobenzene	0.397	0.375	5.5	115	0.00
25		Isophorone	0.732	0.626	14.5	107	-0.01
26	C	2-Nitrophenol	0.158	0.164	-3.8	130	0.00
27		2,4-Dimethylphenol	0.292	0.271	7.2	117	0.00
28		bis(2-Chloroethoxy)methane	0.449	0.400	10.9	112	0.00
29	C	2,4-Dichlorophenol	0.320	0.333	-4.1	128	0.00
30		1,2,4-Trichlorobenzene	0.374	0.355	5.1	121	0.00
31		Naphthalene	1.000	0.919	8.1	115	0.00
32		Benzoic acid	0.202	0.202	0.0	123	0.00
33		4-Chloroaniline	0.458	0.426	7.0	115	-0.01
34	C	Hexachlorobutadiene	0.251	0.243	3.2	120	-0.01
35		Caprolactam	0.127	0.118	7.1	112	0.00
36	C	4-Chloro-3-methylphenol	0.371	0.359	3.2	119	0.00
37		2-Methylnaphthalene	0.732	0.675	7.8	115	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	122	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.708	0.690	2.5	119	-0.01
40	P	Hexachlorocyclopentadiene	0.368	0.328	10.9	107	-0.01
41	S	2,4,6-Tribromophenol	0.239	0.256	-7.1	124	0.00
42	C	2,4,6-Trichlorophenol	0.431	0.451	-4.6	126	0.00
43		2,4,5-Trichlorophenol	0.460	0.484	-5.2	125	0.00
44	S	2-Fluorobiphenyl	1.401	1.377	1.7	122	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.660	1.528	8.0	115	-0.01
46		2-Chloronaphthalene	1.267	1.198	5.4	116	-0.01
47		2-Nitroaniline	0.398	0.403	-1.3	117	0.00
48		Acenaphthylene	1.981	1.841	7.1	114	0.00
49		Dimethylphthalate	1.633	1.513	7.3	113	-0.01
50		2,6-Dinitrotoluene	0.336	0.341	-1.5	122	0.00
51	C	Acenaphthene	1.269	1.188	6.4	115	0.00
52		3-Nitroaniline	0.362	0.383	-5.8	120	0.00
53	P	2,4-Dinitrophenol	0.127	0.145	-14.2	137	0.00
54		Dibenzofuran	1.987	1.880	5.4	117	0.00
55	P	4-Nitrophenol	0.296	0.292	1.4	114	0.00
56		2,4-Dinitrotoluene	0.438	0.462	-5.5	126	0.00
57		Fluorene	1.486	1.383	6.9	113	-0.01
58		2,3,4,6-Tetrachlorophenol	0.432	0.458	-6.0	126	0.00
59		Diethylphthalate	1.646	1.510	8.3	111	-0.01
60		4-Chlorophenyl-phenylether	0.917	0.878	4.3	119	-0.01
61		4-Nitroaniline	0.379	0.392	-3.4	118	0.00
62		Azobenzene	1.675	1.463	12.7	107	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	122	0.00
64		4,6-Dinitro-2-methylphenol	0.088	0.085	3.4	119	0.00
65	c	n-Nitrosodiphenylamine	0.594	0.544	8.4	115	-0.01
66		4-Bromophenyl-phenylether	0.234	0.226	3.4	119	-0.01
67		Hexachlorobenzene	0.239	0.230	3.8	118	0.00
68		Atrazine	0.223	0.204	8.5	115	-0.01
69	C	Pentachlorophenol	0.163	0.162	0.6	114	0.00
70		Phenanthrene	1.016	0.940	7.5	114	-0.01
71		Anthracene	1.013	0.928	8.4	113	0.00
72		Carbazole	1.012	0.935	7.6	113	0.00
73		Di-n-butylphthalate	1.127	1.032	8.4	110	-0.01
74	C	Fluoranthene	1.239	1.169	5.6	114	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	125	0.00
76		Benzidine	0.638	0.524	17.9	110	-0.01
77		Pyrene	1.230	1.123	8.7	113	0.00
78	S	Terphenyl-d14	0.856	0.814	4.9	120	-0.01
79		Butylbenzylphthalate	0.489	0.459	6.1	113	-0.01
80		Benzo(a)anthracene	1.212	1.114	8.1	117	-0.01
81		3,3'-Dichlorobenzidine	0.483	0.448	7.2	113	0.00
82		Chrysene	1.148	1.052	8.4	115	0.00
83		Bis(2-ethylhexyl)phthalate	0.685	0.621	9.3	111	-0.03
84	c	Di-n-octyl phthalate	1.144	1.049	8.3	111	-0.03
85		Indeno(1,2,3-cd)pyrene	1.334	1.259	5.6	118	-0.04
86	I	Perylene-d12	1.000	1.000	0.0	125	-0.03
87		Benzo(b)fluoranthene	1.111	1.033	7.0	115	-0.02

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88	Benzo(k)fluoranthene	1.085	1.018	6.2	117	-0.02
89 C	Benzo(a)pyrene	1.067	0.995	6.7	116	-0.02
90	Dibenzo(a,h)anthracene	1.030	0.960	6.8	116	-0.06
91	Benzo(g,h,i)perylene	1.035	0.970	6.3	116	-0.04

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