Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010319\
Data File : BG038907.D
Acq On : 3 Jan 2019 17:25
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 03 18:12:31 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 12 15:26:27 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	106	-0.01
2		1,4-Dioxane	0.525	0.510	2.9	102	-0.01
3		Pyridine	1.445	1.211	16.2	74	0.00
4		n-Nitrosodimethylamine	0.708	0.603	14.8	83	-0.01
5	S	2-Fluorophenol	1.128	1.118	0.9	105	0.00
6		Aniline	1.976	1.894	4.1	101	-0.01
7	S	Phenol-d6	1.548	1.553	-0.3	107	0.00
8		2-Chlorophenol	1.305	1.279	2.0	104	0.00
9		Benzaldehyde	1.004	0.882	12.2	101	-0.01
10	C	Phenol	1.645	1.621	1.5	105	0.00
11		bis(2-Chloroethyl)ether	1.309	1.171	10.5	97	-0.01
12		1,3-Dichlorobenzene	1.543	1.535	0.5	107	-0.01
13	C	1,4-Dichlorobenzene	1.580	1.509	4.5	103	-0.01
14		1,2-Dichlorobenzene	1.490	1.463	1.8	107	-0.02
15		Benzyl Alcohol	1.208	1.228	-1.7	103	0.00
16		2,2'-oxybis(1-Chloropropane	2.999	2.361	21.3	85	0.00
17		2-Methylphenol	1.102	1.116	-1.3	105	0.00
18		Hexachloroethane	0.577	0.539	6.6	101	-0.02
19	P	n-Nitroso-di-n-propylamine	1.087	1.011	7.0	98	-0.01
20		3+4-Methylphenols	1.522	1.578	-3.7	108	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	107	-0.02
22		Acetophenone	0.500	0.492	1.6	105	-0.02
23	S	Nitrobenzene-d5	0.391	0.366	6.4	100	-0.01
24		Nitrobenzene	0.396	0.365	7.8	98	-0.01
25		Isophorone	0.703	0.615	12.5	80	-0.01
26	C	2-Nitrophenol	0.203	0.200	1.5	105	-0.01
27		2,4-Dimethylphenol	0.291	0.279	4.1	102	0.00
28		bis(2-Chloroethoxy)methane	0.397	0.378	4.8	100	-0.02
29	C	2,4-Dichlorophenol	0.323	0.364	-12.7	115	0.00
30		1,2,4-Trichlorobenzene	0.390	0.413	-5.9	114	-0.01
31		Naphthalene	0.985	0.984	0.1	107	-0.02
32		Benzoic acid	0.154	0.136	11.7	86	-0.04
33		4-Chloroaniline	0.428	0.443	-3.5	107	-0.01
34	C	Hexachlorobutadiene	0.264	0.286	-8.3	113	-0.02
35		Caprolactam	0.134	0.135	-0.7	113	0.00
36	C	4-Chloro-3-methylphenol	0.369	0.390	-5.7	114	0.00
37		2-Methylnaphthalene	0.703	0.726	-3.3	111	-0.02
38		1-Methylnaphthalene	0.682	0.698	-2.3	110	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	119	-0.01
40		1,2,4,5-Tetrachlorobenzene	0.748	0.750	-0.3	118	-0.01
41	P	Hexachlorocyclopentadiene	0.411	0.435	-5.8	122	-0.02
42	S	2,4,6-Tribromophenol	0.276	0.317	-14.9	133	-0.01
43	C	2,4,6-Trichlorophenol	0.460	0.493	-7.2	120	0.00
44		2,4,5-Trichlorophenol	0.487	0.529	-8.6	123	0.00

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Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.458	1.425	2.3	116	-0.01
46		1,1'-Biphenyl	1.677	1.600	4.6	112	-0.01
47		2-Chloronaphthalene	1.295	1.255	3.1	115	-0.02
48		2-Nitroaniline	0.413	0.382	7.5	107	0.00
49		Acenaphthylene	1.936	1.860	3.9	112	-0.01
50		Dimethylphthalate	1.633	1.586	2.9	114	0.00
51		2,6-Dinitrotoluene	0.370	0.361	2.4	115	-0.01
52	C	Acenaphthene	1.158	1.117	3.5	114	-0.01
53		3-Nitroaniline	0.377	0.372	1.3	114	0.00
54	P	2,4-Dinitrophenol	0.193	0.154	20.2	91	0.00
55		Dibenzofuran	1.985	1.958	1.4	118	0.00
56	P	4-Nitrophenol	0.286	0.256	10.5	100	0.00
57		2,4-Dinitrotoluene	0.521	0.509	2.3	112	0.00
58		Fluorene	1.470	1.477	-0.5	118	-0.01
59		2,3,4,6-Tetrachlorophenol	0.438	0.479	-9.4	124	0.00
60		Diethylphthalate	1.793	1.684	6.1	112	0.00
61		4-Chlorophenyl-phenylether	0.917	0.931	-1.5	119	-0.01
62		4-Nitroaniline	0.388	0.385	0.8	112	0.00
63		Azobenzene	1.498	1.350	9.9	106	-0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	124	0.00
65		4,6-Dinitro-2-methylphenol	0.143	0.109	23.8	93	0.00
66	c	n-Nitrosodiphenylamine	0.588	0.570	3.1	119	-0.01
67		4-Bromophenyl-phenylether	0.244	0.249	-2.0	124	-0.01
68		Hexachlorobenzene	0.253	0.259	-2.4	123	0.00
69		Atrazine	0.218	0.204	6.4	112	-0.01
70	C	Pentachlorophenol	0.150	0.137	8.7	109	0.00
71		Phenanthrene	1.037	1.000	3.6	120	-0.01
72		Anthracene	1.024	0.991	3.2	118	0.00
73		Carbazole	1.013	0.972	4.0	117	0.00
74		Di-n-butylphthalate	1.162	1.084	6.7	113	-0.01
75	C	Fluoranthene	1.283	1.216	5.2	119	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	121	0.00
77		Benzidine	0.591	0.531	10.2	90	-0.01
78		Pyrene	1.321	1.262	4.5	118	0.00
79	S	Terphenyl-d14	0.919	0.894	2.7	118	0.00
80		Butylbenzylphthalate	0.516	0.495	4.1	113	-0.01
81		Benzo(a)anthracene	1.219	1.192	2.2	116	0.00
82		3,3'-Dichlorobenzidine	0.483	0.482	0.2	115	0.00
83		Chrysene	1.174	1.113	5.2	114	-0.01
84		Bis(2-ethylhexyl)phthalate	0.732	0.691	5.6	109	-0.01
85	c	Di-n-octyl phthalate	1.226	1.144	6.7	107	-0.02
86		Indeno(1,2,3-cd)pyrene	1.380	1.333	3.4	112	-0.02
87	I	Perylene-d12	1.000	1.000	0.0	118	-0.01

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Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.098	1.070	2.6	112	0.00
89	Benzo(k)fluoranthene	1.093	1.052	3.8	111	-0.01
90 C	Benzo(a)pyrene	1.059	1.034	2.4	113	-0.02
91	Dibenzo(a,h)anthracene	1.055	1.021	3.2	111	-0.02
92	Benzo(g,h,i)perylene	1.039	1.014	2.4	112	-0.03

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