Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121720\
Data File : BG047767.D
Acq On : 17 Dec 2020 15:12
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG121720

Quant Time: Dec 17 15:57:16 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG121720.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 17 15:55:16 2020
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	100	0.00
2		1,4-Dioxane	0.497	0.416	16.3	106	0.00
3		Pyridine	1.282	1.158	9.7	103	0.00
4		n-Nitrosodimethylamine	1.014	0.940	7.3	106	0.00
5	S	2-Fluorophenol	1.156	1.028	11.1	106	0.00
6		Aniline	1.875	1.681	10.3	107	0.00
7	S	Phenol-d6	1.690	1.508	10.8	103	0.00
8		2-Chlorophenol	1.178	1.051	10.8	104	0.00
9		Benzaldehyde	1.108	0.935	15.6	106	0.00
10	C	Phenol	1.532	1.380	9.9	103	0.00
11		bis(2-Chloroethyl)ether	1.092	0.959	12.2	106	0.00
12		1,3-Dichlorobenzene	1.589	1.393	12.3	103	0.00
13	C	1,4-Dichlorobenzene	1.587	1.378	13.2	103	0.00
14		1,2-Dichlorobenzene	1.493	1.291	13.5	104	0.00
15		Benzyl Alcohol	1.497	1.409	5.9	107	0.00
16		2,2'-oxybis(1-Chloropropane	1.085	0.924	14.8	103	0.00
17		2-Methylphenol	1.081	0.929	14.1	103	0.00
18		Hexachloroethane	0.622	0.562	9.6	104	0.00
19	P	n-Nitroso-di-n-propylamine	1.176	1.030	12.4	103	0.00
20		3+4-Methylphenols	1.481	1.312	11.4	105	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
22		Acetophenone	0.572	0.519	9.3	108	0.00
23	S	Nitrobenzene-d5	0.544	0.480	11.8	104	0.00
24		Nitrobenzene	0.513	0.453	11.7	105	0.00
25		Isophorone	0.820	0.738	10.0	107	0.00
26	C	2-Nitrophenol	0.203	0.183	9.9	105	0.00
27		2,4-Dimethylphenol	0.292	0.249	14.7	104	0.00
28		bis(2-Chloroethoxy)methane	0.424	0.373	12.0	109	0.00
29	C	2,4-Dichlorophenol	0.381	0.332	12.9	104	0.00
30		1,2,4-Trichlorobenzene	0.488	0.441	9.6	103	0.00
31		Naphthalene	1.099	0.980	10.8	104	0.00
32		Benzoic acid	0.153	0.083	45.8#	117	0.00
33		4-Chloroaniline	0.467	0.411	12.0	101	0.00
34	C	Hexachlorobutadiene	0.431	0.384	10.9	106	0.00
35		Caprolactam	0.121	0.108	10.7	109	0.00
36	C	4-Chloro-3-methylphenol	0.416	0.374	10.1	105	0.00
37		2-Methylnaphthalene	0.853	0.772	9.5	107	0.00
38		1-Methylnaphthalene	0.806	0.715	11.3	105	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	112	0.00
40		1,2,4,5-Tetrachlorobenzene	0.821	0.700	14.7	106	0.00
41	P	Hexachlorocyclopentadiene	0.444	0.409	7.9	103	0.00
42	S	2,4,6-Tribromophenol	0.341	0.310	9.1	110	0.00
43	C	2,4,6-Trichlorophenol	0.524	0.464	11.5	109	0.00
44		2,4,5-Trichlorophenol	0.519	0.463	10.8	109	0.00
45	S	2-Fluorobiphenyl	1.546	1.337	13.5	106	0.00
46		1,1'-Biphenyl	1.526	1.321	13.4	107	0.00
47		2-Chloronaphthalene	1.241	1.068	13.9	103	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.459	0.426	7.2	113	0.00
49		Acenaphthylene	1.842	1.621	12.0	107	0.00
50		Dimethylphthalate	1.724	1.492	13.5	106	0.00
51		2,6-Dinitrotoluene	0.350	0.312	10.9	110	0.00
52	C	Acenaphthene	1.299	1.164	10.4	107	0.00
53		3-Nitroaniline	0.347	0.308	11.2	107	0.00
54	P	2,4-Dinitrophenol	0.225	0.210	6.7	111	0.00
55		Dibenzofuran	1.883	1.633	13.3	108	0.00
56	P	4-Nitrophenol	0.250	0.241	3.6	115	0.00
57		2,4-Dinitrotoluene	0.521	0.452	13.2	109	0.00
58		Fluorene	1.600	1.366	14.6	107	0.00
59		2,3,4,6-Tetrachlorophenol	0.530	0.471	11.1	109	0.00
60		Diethylphthalate	1.662	1.436	13.6	107	0.00
61		4-Chlorophenyl-phenylether	1.008	0.887	12.0	109	0.00
62		4-Nitroaniline	0.382	0.334	12.6	106	0.00
63		Azobenzene	1.550	1.356	12.5	108	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
65		4,6-Dinitro-2-methylphenol	0.128	0.116	9.4	107	0.00
66	c	n-Nitrosodiphenylamine	0.572	0.495	13.5	104	0.00
67		4-Bromophenyl-phenylether	0.269	0.233	13.4	103	0.00
68		Hexachlorobenzene	0.292	0.263	9.9	108	0.00
69		Atrazine	0.240	0.216	10.0	109	0.00
70	C	Pentachlorophenol	0.133	0.149	-12.0	114	0.00
71		Phenanthrene	1.078	0.937	13.1	107	0.00
72		Anthracene	1.045	0.898	14.1	106	0.00
73		Carbazole	0.915	0.786	14.1	108	0.00
74		Di-n-butylphthalate	1.079	0.943	12.6	106	0.00
75	C	Fluoranthene	1.446	1.247	13.8	107	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	107	0.00
77		Benzidine	0.538	0.477	11.3	98	0.00
78		Pyrene	1.433	1.227	14.4	106	0.00
79	S	Terphenyl-d14	1.214	1.040	14.3	105	0.00
80		Butylbenzylphthalate	0.491	0.440	10.4	109	0.00
81		Benzo(a)anthracene	1.437	1.255	12.7	106	0.00
82		3,3'-Dichlorobenzidine	0.506	0.440	13.0	104	0.00
83		Chrysene	1.357	1.150	15.3	104	0.00
84		Bis(2-ethylhexyl)phthalate	0.677	0.598	11.7	107	0.00
85	c	Di-n-octyl phthalate	1.148	1.011	11.9	107	0.00
86	I	Perylene-d12	1.000	1.000	0.0	106	0.00
87		Indeno(1,2,3-cd)pyrene	1.519	1.330	12.4	104	0.02
88		Benzo(b)fluoranthene	1.379	1.223	11.3	105	0.00
89		Benzo(k)fluoranthene	1.337	1.174	12.2	107	0.00
90	C	Benzo(a)pyrene	1.283	1.124	12.4	105	0.00
91		Dibenzo(a,h)anthracene	1.230	1.068	13.2	104	0.00
92		Benzo(g,h,i)perylene	1.203	1.036	13.9	103	0.00

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