Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051225\
Data File : BP024611.D
Acq On : 12 May 2025 21:25
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP051225

Quant Time: May 13 04:01:05 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue May 13 03:56:11 2025
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	113	0.00
2		1,4-Dioxane	0.502	0.481	4.2	109	0.00
3		Pyridine	1.176	1.205	-2.5	111	0.00
4		n-Nitrosodimethylamine	0.517	0.502	2.9	109	0.00
5	S	2-Fluorophenol	1.132	1.136	-0.4	111	0.00
6		Aniline	1.335	1.387	-3.9	112	0.00
7	S	Phenol-d6	1.489	1.533	-3.0	112	0.00
8		2-Chlorophenol	1.303	1.310	-0.5	112	0.00
9		Benzaldehyde	0.822	0.843	-2.6	113	0.00
10	C	Phenol	1.483	1.515	-2.2	113	0.00
11		bis(2-Chloroethyl)ether	1.204	1.192	1.0	111	0.00
12		1,3-Dichlorobenzene	1.434	1.422	0.8	112	0.00
13	C	1,4-Dichlorobenzene	1.460	1.436	1.6	110	0.00
14		1,2-Dichlorobenzene	1.413	1.374	2.8	111	0.00
15		Benzyl Alcohol	1.026	1.101	-7.3	114	0.00
16		2,2'-oxybis(1-Chloropropane	1.381	1.365	1.2	111	0.00
17		2-Methylphenol	1.020	1.055	-3.4	113	0.00
18		Hexachloroethane	0.554	0.540	2.5	110	0.00
19	P	n-Nitroso-di-n-propylamine	0.993	1.002	-0.9	111	0.00
20		3+4-Methylphenols	1.403	1.455	-3.7	113	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	111	0.00
22		Acetophenone	0.479	0.480	-0.2	112	0.00
23	S	Nitrobenzene-d5	0.348	0.350	-0.6	112	0.00
24		Nitrobenzene	0.338	0.342	-1.2	112	0.00
25		Isophorone	0.610	0.614	-0.7	112	0.00
26	C	2-Nitrophenol	0.173	0.184	-6.4	115	0.00
27		2,4-Dimethylphenol	0.208	0.214	-2.9	113	0.00
28		bis(2-Chloroethoxy)methane	0.396	0.396	0.0	112	0.00
29	C	2,4-Dichlorophenol	0.289	0.299	-3.5	112	0.00
30		1,2,4-Trichlorobenzene	0.321	0.315	1.9	112	0.00
31		Naphthalene	1.043	1.026	1.6	112	0.00
32		Benzoic acid	0.195	0.218	-11.8	121	0.01
33		4-Chloroaniline	0.351	0.370	-5.4	113	0.00
34	C	Hexachlorobutadiene	0.207	0.201	2.9	110	0.00
35		Caprolactam	0.101	0.109	-7.9	116	0.00
36	C	4-Chloro-3-methylphenol	0.353	0.362	-2.5	114	0.00
37		2-Methylnaphthalene	0.731	0.723	1.1	112	0.00
38		1-Methylnaphthalene	0.723	0.706	2.4	111	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
40		1,2,4,5-Tetrachlorobenzene	0.558	0.565	-1.3	112	0.00
41	P	Hexachlorocyclopentadiene	0.183	0.198	-8.2	110	0.00
42	S	2,4,6-Tribromophenol	0.328	0.341	-4.0	114	0.00
43	C	2,4,6-Trichlorophenol	0.375	0.398	-6.1	114	0.00
44		2,4,5-Trichlorophenol	0.418	0.437	-4.5	113	0.00
45	S	2-Fluorobiphenyl	1.337	1.342	-0.4	112	0.00
46		1,1'-Biphenyl	1.438	1.447	-0.6	111	0.00
47		2-Chloronaphthalene	1.092	1.107	-1.4	113	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.320	0.341	-6.6	114	0.00
49		Acenaphthylene	1.706	1.711	-0.3	111	0.00
50		Dimethylphthalate	1.492	1.518	-1.7	114	0.00
51		2,6-Dinitrotoluene	0.318	0.333	-4.7	114	0.00
52	C	Acenaphthene	1.087	1.075	1.1	110	0.00
53		3-Nitroaniline	0.307	0.332	-8.1	114	0.00
54	P	2,4-Dinitrophenol	0.175	0.191	-9.1	118	0.00
55		Dibenzofuran	1.770	1.732	2.1	109	0.00
56	P	4-Nitrophenol	0.252	0.273	-8.3	115	0.00
57		2,4-Dinitrotoluene	0.439	0.458	-4.3	112	0.00
58		Fluorene	1.375	1.389	-1.0	113	0.00
59		2,3,4,6-Tetrachlorophenol	0.394	0.408	-3.6	114	0.00
60		Diethylphthalate	1.545	1.563	-1.2	113	0.00
61		4-Chlorophenyl-phenylether	0.693	0.694	-0.1	113	0.00
62		4-Nitroaniline	0.306	0.340	-11.1	117	0.00
63		Azobenzene	1.318	1.338	-1.5	113	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	115	0.00
65		4,6-Dinitro-2-methylphenol	0.123	0.133	-8.1	115	0.00
66	c	n-Nitrosodiphenylamine	0.598	0.593	0.8	112	0.00
67		4-Bromophenyl-phenylether	0.236	0.232	1.7	112	0.00
68		Hexachlorobenzene	0.294	0.289	1.7	115	-0.01
69		Atrazine	0.202	0.203	-0.5	114	0.00
70	C	Pentachlorophenol	0.189	0.201	-6.3	118	0.00
71		Phenanthrene	1.079	1.059	1.9	113	0.00
72		Anthracene	1.054	1.034	1.9	112	0.00
73		Carbazole	0.988	1.009	-2.1	114	0.00
74		Di-n-butylphthalate	1.330	1.349	-1.4	114	0.00
75	C	Fluoranthene	1.315	1.310	0.4	113	-0.01
76	I	Chrysene-d12	1.000	1.000	0.0	112	0.00
77		Benzidine	0.234	0.212	9.4	87	-0.02
78		Pyrene	1.210	1.204	0.5	112	0.00
79	S	Terphenyl-d14	1.022	1.019	0.3	114	-0.01
80		Butylbenzylphthalate	0.540	0.557	-3.1	112	-0.01
81		Benzo(a)anthracene	1.218	1.215	0.2	113	0.00
82		3,3'-Dichlorobenzidine	0.291	0.306	-5.2	108	-0.01
83		Chrysene	1.180	1.157	1.9	111	0.00
84		Bis(2-ethylhexyl)phthalate	0.792	0.818	-3.3	114	0.00
85	c	Di-n-octyl phthalate	1.313	1.387	-5.6	113	0.00
86	I	Perylene-d12	1.000	1.000	0.0	110	0.00
87		Indeno(1,2,3-cd)pyrene	1.367	1.387	-1.5	110	0.00
88		Benzo(b)fluoranthene	1.154	1.168	-1.2	112	0.00
89		Benzo(k)fluoranthene	1.139	1.129	0.9	109	0.00
90	C	Benzo(a)pyrene	0.999	1.014	-1.5	111	0.00
91		Dibenzo(a,h)anthracene	1.121	1.151	-2.7	111	0.00
92		Benzo(g,h,i)perylene	1.159	1.175	-1.4	111	-0.01

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