Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090325\
Data File : BF143632.D
Acq On : 03 Sep 2025 18:08
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF090325

Quant Time: Sep 03 18:50:39 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF090325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 03 18:35:49 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	105	0.00
2		1,4-Dioxane	0.533	0.536	-0.6	103	0.00
3		Pyridine	1.411	1.414	-0.2	102	0.00
4		n-Nitrosodimethylamine	0.586	0.589	-0.5	103	0.00
5	S	2-Fluorophenol	1.129	1.159	-2.7	105	0.00
6		Aniline	1.950	1.969	-1.0	104	0.00
7	S	Phenol-d6	1.401	1.426	-1.8	105	0.00
8		2-Chlorophenol	1.169	1.277	-9.2	108	0.00
9		Benzaldehyde	1.041	1.138	-9.3	104	0.00
10	C	Phenol	1.536	1.574	-2.5	104	0.00
11		bis(2-Chloroethyl)ether	1.203	1.207	-0.3	105	0.00
12		1,3-Dichlorobenzene	1.443	1.442	0.1	104	0.00
13	C	1,4-Dichlorobenzene	1.447	1.457	-0.7	106	0.00
14		1,2-Dichlorobenzene	1.370	1.372	-0.1	103	0.00
15		Benzyl Alcohol	1.065	1.106	-3.8	106	0.00
16		2,2'-oxybis(1-Chloropropane	1.894	1.893	0.1	104	0.00
17		2-Methylphenol	1.022	1.049	-2.6	105	0.00
18		Hexachloroethane	0.451	0.488	-8.2	109	0.00
19	P	n-Nitroso-di-n-propylamine	0.892	0.920	-3.1	105	0.00
20		3+4-Methylphenols	1.287	1.334	-3.7	104	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
22		Acetophenone	0.452	0.451	0.2	103	0.00
23	S	Nitrobenzene-d5	0.285	0.337	-18.2	113	0.00
24		Nitrobenzene	0.304	0.352	-15.8	110	0.00
25		Isophorone	0.641	0.656	-2.3	107	0.00
26	C	2-Nitrophenol	0.112	0.161	-43.8#	137	0.00
27		2,4-Dimethylphenol	0.279	0.296	-6.1	107	0.00
28		bis(2-Chloroethoxy)methane	0.392	0.392	0.0	104	0.00
29	C	2,4-Dichlorophenol	0.274	0.298	-8.8	107	0.00
30		1,2,4-Trichlorobenzene	0.299	0.305	-2.0	104	0.00
31		Naphthalene	0.978	0.991	-1.3	106	0.00
32		Benzoic acid	0.145	0.192	-32.4#	139	0.00
33		4-Chloroaniline	0.339	0.356	-5.0	109	0.00
34	C	Hexachlorobutadiene	0.194	0.200	-3.1	106	0.00
35		Caprolactam	0.073	0.086	-17.8	116	0.00
36	C	4-Chloro-3-methylphenol	0.287	0.307	-7.0	108	0.00
37		2-Methylnaphthalene	0.662	0.677	-2.3	106	0.00
38		1-Methylnaphthalene	0.633	0.647	-2.2	107	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
40		1,2,4,5-Tetrachlorobenzene	0.553	0.555	-0.4	107	0.00
41	P	Hexachlorocyclopentadiene	0.268	0.294	-9.7	112	0.00
42	S	2,4,6-Tribromophenol	0.167	0.201	-20.4	116	0.00
43	C	2,4,6-Trichlorophenol	0.341	0.377	-10.6	113	0.00
44		2,4,5-Trichlorophenol	0.352	0.394	-11.9	111	0.00
45	S	2-Fluorobiphenyl	1.252	1.227	2.0	106	0.00
46		1,1'-Biphenyl	1.422	1.402	1.4	105	0.00
47		2-Chloronaphthalene	1.113	1.111	0.2	107	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.261	0.324	-24.1	119	0.00
49		Acenaphthylene	1.604	1.608	-0.2	106	0.00
50		Dimethylphthalate	1.266	1.325	-4.7	110	0.00
51		2,6-Dinitrotoluene	0.193	0.256	-32.6#	125	0.00
52	C	Acenaphthene	1.073	1.082	-0.8	108	0.00
53		3-Nitroaniline	0.223	0.285	-27.8#	120	0.00
54	P	2,4-Dinitrophenol	0.074	0.104	-40.5#	158#	0.00
55		Dibenzofuran	1.559	1.576	-1.1	108	0.00
56	P	4-Nitrophenol	0.188	0.231	-22.9	123	0.00
57		2,4-Dinitrotoluene	0.260	0.349	-34.2#	126	0.00
58		Fluorene	1.206	1.198	0.7	107	0.00
59		2,3,4,6-Tetrachlorophenol	0.267	0.321	-20.2	115	0.00
60		Diethylphthalate	1.202	1.266	-5.3	109	0.00
61		4-Chlorophenyl-phenylether	0.602	0.603	-0.2	107	0.00
62		4-Nitroaniline	0.219	0.274	-25.1#	117	0.00
63		Azobenzene	1.133	1.158	-2.2	108	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
65		4,6-Dinitro-2-methylphenol	0.070	0.096	-37.1#	147	0.00
66	c	n-Nitrosodiphenylamine	0.646	0.638	1.2	107	0.00
67		4-Bromophenyl-phenylether	0.232	0.236	-1.7	110	0.00
68		Hexachlorobenzene	0.244	0.245	-0.4	107	0.00
69		Atrazine	0.193	0.214	-10.9	112	0.00
70	C	Pentachlorophenol	0.137	0.160	-16.8	121	0.00
71		Phenanthrene	1.081	1.068	1.2	108	0.00
72		Anthracene	1.085	1.088	-0.3	108	0.00
73		Carbazole	0.914	0.942	-3.1	110	0.00
74		Di-n-butylphthalate	1.027	1.151	-12.1	113	0.00
75	C	Fluoranthene	0.993	1.018	-2.5	109	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	107	0.00
77		Benzidine	0.458	0.497	-8.5	108	0.00
78		Pyrene	1.650	1.721	-4.3	108	0.00
79	S	Terphenyl-d14	1.200	1.269	-5.7	111	0.00
80		Butylbenzylphthalate	0.452	0.579	-28.1#	118	0.00
81		Benzo(a)anthracene	1.295	1.275	1.5	107	0.00
82		3,3'-Dichlorobenzidine	0.353	0.382	-8.2	112	0.00
83		Chrysene	1.164	1.217	-4.6	107	0.00
84		Bis(2-ethylhexyl)phthalate	0.648	0.772	-19.1	115	0.00
85	c	Di-n-octyl phthalate	1.187	1.321	-11.3	115	0.00
86	I	Perylene-d12	1.000	1.000	0.0	105	0.00
87		Indeno(1,2,3-cd)pyrene	1.394	1.444	-3.6	104	0.00
88		Benzo(b)fluoranthene	1.205	1.257	-4.3	102	0.00
89		Benzo(k)fluoranthene	1.042	1.041	0.1	106	0.00
90	C	Benzo(a)pyrene	1.036	1.067	-3.0	104	0.00
91		Dibenzo(a,h)anthracene	1.147	1.184	-3.2	104	0.00
92		Benzo(g,h,i)perylene	1.105	1.122	-1.5	104	0.00

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(#) = Out of Range SPCC's out = 0 CCC's out = 1