Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100225\
Data File : BG064482.D
Acq On : 2 Oct 2025 18:09
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ICVBG100225

Quant Time: Oct 02 18:56:29 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 02 16:40:42 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	91	0.00
2		1,4-Dioxane	0.427	0.392	8.2	81	0.00
3		Pyridine	1.149	1.017	11.5	75	0.00
4		n-Nitrosodimethylamine	0.833	0.813	2.4	86	0.00
5	S	2-Fluorophenol	1.081	1.018	5.8	78	0.00
6		Aniline	1.832	1.883	-2.8	85	0.00
7	S	Phenol-d6	1.452	1.407	3.1	79	0.01
8		2-Chlorophenol	1.327	1.231	7.2	77	0.00
9		Benzaldehyde	1.289	1.040	19.3	62	0.00
10	C	Phenol	1.532	1.471	4.0	80	0.00
11		bis(2-Chloroethyl)ether	1.038	0.969	6.6	78	0.00
12		1,3-Dichlorobenzene	1.452	1.361	6.3	78	0.00
13	C	1,4-Dichlorobenzene	1.457	1.452	0.3	84	0.00
14		1,2-Dichlorobenzene	1.436	1.381	3.8	82	0.00
15		Benzyl Alcohol	1.337	1.363	-1.9	86	0.00
16		2,2'-oxybis(1-Chloropropane	1.672	1.597	4.5	81	0.01
17		2-Methylphenol	1.109	1.069	3.6	80	0.00
18		Hexachloroethane	0.575	0.545	5.2	82	0.00
19	P	n-Nitroso-di-n-propylamine	0.971	0.995	-2.5	83	0.01
20		3+4-Methylphenols	1.548	1.505	2.8	81	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
22		Acetophenone	0.484	0.483	0.2	86	0.00
23	S	Nitrobenzene-d5	0.372	0.339	8.9	79	0.01
24		Nitrobenzene	0.371	0.345	7.0	80	0.00
25		Isophorone	0.674	0.660	2.1	84	0.01
26	C	2-Nitrophenol	0.185	0.163	11.9	78	0.00
27		2,4-Dimethylphenol	0.282	0.296	-5.0	91	0.00
28		bis(2-Chloroethoxy)methane	0.343	0.331	3.5	81	0.00
29	C	2,4-Dichlorophenol	0.326	0.317	2.8	85	0.00
30		1,2,4-Trichlorobenzene	0.356	0.344	3.4	84	0.00
31		Naphthalene	1.030	0.991	3.8	82	0.00
32		Benzoic acid	0.216	0.218	-0.9	87	0.05
33		4-Chloroaniline	0.395	0.423	-7.1	93	0.00
34	C	Hexachlorobutadiene	0.304	0.278	8.6	81	0.00
35		Caprolactam	0.112	0.108	3.6	86	0.02
36	C	4-Chloro-3-methylphenol	0.397	0.381	4.0	83	0.00
37		2-Methylnaphthalene	0.791	0.696	12.0	76	0.00
38		1-Methylnaphthalene	0.788	0.728	7.6	81	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	95	0.00
40		1,2,4,5-Tetrachlorobenzene	0.646	0.622	3.7	84	0.00
41	P	Hexachlorocyclopentadiene	0.338	0.426	-26.0#	112	0.00
42	S	2,4,6-Tribromophenol	0.363	0.348	4.1	87	0.00
43	C	2,4,6-Trichlorophenol	0.413	0.403	2.4	86	0.00
44		2,4,5-Trichlorophenol	0.461	0.436	5.4	85	0.00
45	S	2-Fluorobiphenyl	1.357	1.264	6.9	84	0.00
46		1,1'-Biphenyl	1.438	1.393	3.1	87	0.00
47		2-Chloronaphthalene	1.074	1.013	5.7	84	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.367	0.359	2.2	86	0.00
49		Acenaphthylene	1.594	1.747	-9.6	98	0.00
50		Dimethylphthalate	1.566	1.484	5.2	87	0.00
51		2,6-Dinitrotoluene	0.326	0.316	3.1	89	0.00
52	C	Acenaphthene	1.108	1.015	8.4	83	0.00
53		3-Nitroaniline	0.301	0.306	-1.7	91	0.00
54	P	2,4-Dinitrophenol	0.237	0.198	16.5	91	0.00
55		Dibenzofuran	1.829	1.744	4.6	87	0.00
56	P	4-Nitrophenol	0.237	0.235	0.8	86	0.00
57		2,4-Dinitrotoluene	0.499	0.491	1.6	89	0.00
58		Fluorene	1.520	1.500	1.3	89	0.00
59		2,3,4,6-Tetrachlorophenol	0.470	0.471	-0.2	89	0.00
60		Diethylphthalate	1.680	1.610	4.2	89	0.00
61		4-Chlorophenyl-phenylether	0.830	0.756	8.9	83	0.00
62		4-Nitroaniline	0.314	0.325	-3.5	93	0.00
63		Azobenzene	1.301	1.254	3.6	89	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	100	0.00
65		4,6-Dinitro-2-methylphenol	0.139	0.132	5.0	91	0.00
66	c	n-Nitrosodiphenylamine	0.534	0.500	6.4	88	0.00
67		4-Bromophenyl-phenylether	0.235	0.217	7.7	85	0.00
68		Hexachlorobenzene	0.269	0.248	7.8	86	0.00
69		Atrazine	0.238	0.250	-5.0	100	0.00
70	C	Pentachlorophenol	0.157	0.142	9.6	84	0.00
71		Phenanthrene	1.046	0.970	7.3	87	0.00
72		Anthracene	1.049	1.001	4.6	89	0.00
73		Carbazole	0.907	0.900	0.8	91	0.00
74		Di-n-butylphthalate	1.107	1.075	2.9	90	0.00
75	C	Fluoranthene	1.250	1.210	3.2	91	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
77		Benzidine	0.498	0.545	-9.4	85	0.00
78		Pyrene	1.298	1.196	7.9	88	0.00
79	S	Terphenyl-d14	1.057	1.005	4.9	91	0.00
80		Butylbenzylphthalate	0.460	0.432	6.1	89	0.00
81		Benzo(a)anthracene	1.277	1.209	5.3	91	0.00
82		3,3'-Dichlorobenzidine	0.432	0.458	-6.0	102	0.00
83		Chrysene	1.211	1.103	8.9	87	0.00
84		Bis(2-ethylhexyl)phthalate	0.649	0.602	7.2	89	0.00
85	c	Di-n-octyl phthalate	1.091	1.002	8.2	88	0.00
86	I	Perylene-d12	1.000	1.000	0.0	101	0.00
87		Indeno(1,2,3-cd)pyrene	1.392	1.316	5.5	89	0.00
88		Benzo(b)fluoranthene	1.138	1.064	6.5	89	0.00
89		Benzo(k)fluoranthene	1.146	1.136	0.9	93	0.00
90	C	Benzo(a)pyrene	1.049	1.035	1.3	93	0.00
91		Dibenzo(a,h)anthracene	1.113	1.074	3.5	91	0.01
92		Benzo(g,h,i)perylene	1.082	1.068	1.3	93	0.02

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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