Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100325\
Data File : BP025913.D
Acq On : 03 Oct 2025 14:06
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Oct 03 14:28:48 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 11 16:45:04 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	153#	-0.02
2		1,4-Dioxane	0.520	0.505	2.9	142	0.00
3		Pyridine	1.432	1.406	1.8	137	0.00
4		n-Nitrosodimethylamine	0.676	0.583	13.8	124	0.00
5	S	2-Fluorophenol	1.240	1.313	-5.9	147	0.00
6		Aniline	2.262	2.202	2.7	134	-0.02
7	S	Phenol-d6	1.648	1.687	-2.4	139	0.00
8		2-Chlorophenol	1.360	1.404	-3.2	142	-0.02
9		Benzaldehyde	0.987	1.288	-30.5#	225#	-0.02
10	C	Phenol	1.737	1.759	-1.3	137	0.00
11		bis(2-Chloroethyl)ether	1.396	1.365	2.2	134	-0.02
12		1,3-Dichlorobenzene	1.548	1.570	-1.4	143	-0.02
13	C	1,4-Dichlorobenzene	1.576	1.581	-0.3	143	-0.02
14		1,2-Dichlorobenzene	1.531	1.513	1.2	140	-0.02
15		Benzyl Alcohol	1.361	1.320	3.0	134	-0.02
16		2,2'-oxybis(1-Chloropropane	2.282	1.956	14.3	120	-0.02
17		2-Methylphenol	1.171	1.177	-0.5	137	-0.01
18		Hexachloroethane	0.602	0.601	0.2	142	-0.02
19	P	n-Nitroso-di-n-propylamine	1.149	1.092	5.0	127	-0.02
20		3+4-Methylphenols	1.640	1.580	3.7	133	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	142	-0.02
22		Acetophenone	0.534	0.542	-1.5	131	-0.02
23	S	Nitrobenzene-d5	0.453	0.446	1.5	128	-0.02
24		Nitrobenzene	0.407	0.401	1.5	127	-0.02
25		Isophorone	0.797	0.774	2.9	128	-0.02
26	C	2-Nitrophenol	0.181	0.202	-11.6	142	-0.02
27		2,4-Dimethylphenol	0.317	0.323	-1.9	129	-0.02
28		bis(2-Chloroethoxy)methane	0.461	0.456	1.1	130	-0.03
29	C	2,4-Dichlorophenol	0.316	0.337	-6.6	137	-0.01
30		1,2,4-Trichlorobenzene	0.356	0.367	-3.1	138	-0.03
31		Naphthalene	1.078	1.066	1.1	132	-0.02
32		Benzoic acid	0.236	0.239	-1.3	130	0.01
33		4-Chloroaniline	0.449	0.465	-3.6	130	-0.02
34	C	Hexachlorobutadiene	0.229	0.238	-3.9	139	-0.03
35		Caprolactam	0.122	0.143	-17.2	154#	0.00
36	C	4-Chloro-3-methylphenol	0.377	0.378	-0.3	130	-0.01
37		2-Methylnaphthalene	0.693	0.687	0.9	131	-0.02
38		1-Methylnaphthalene	0.739	0.720	2.6	128	-0.02
39	I	Acenaphthene-d10	1.000	1.000	0.0	135	-0.03
40		1,2,4,5-Tetrachlorobenzene	0.606	0.655	-8.1	134	-0.02
41	P	Hexachlorocyclopentadiene	0.330	0.303	8.2	108	-0.04
42	S	2,4,6-Tribromophenol	0.249	0.284	-14.1	139	-0.04
43	C	2,4,6-Trichlorophenol	0.398	0.436	-9.5	132	-0.02
44		2,4,5-Trichlorophenol	0.443	0.484	-9.3	131	-0.01
45	S	2-Fluorobiphenyl	1.502	1.542	-2.7	131	-0.03
46		1,1'-Biphenyl	1.468	1.494	-1.8	127	-0.03
47		2-Chloronaphthalene	1.156	1.175	-1.6	128	-0.03

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.385	0.392	-1.8	121	-0.02
49		Acenaphthylene	1.915	1.936	-1.1	127	-0.04
50		Dimethylphthalate	1.527	1.508	1.2	123	-0.04
51		2,6-Dinitrotoluene	0.324	0.339	-4.6	128	-0.03
52	C	Acenaphthene	1.104	1.096	0.7	124	-0.04
53		3-Nitroaniline	0.340	0.363	-6.8	124	-0.02
54	P	2,4-Dinitrophenol	0.183	0.227	-24.0	147	-0.02
55		Dibenzofuran	1.800	1.777	1.3	125	-0.04
56	P	4-Nitrophenol	0.258	0.360	-39.5#	158#	0.00
57		2,4-Dinitrotoluene	0.460	0.489	-6.3	127	-0.03
58		Fluorene	1.431	1.413	1.3	123	-0.05
59		2,3,4,6-Tetrachlorophenol	0.399	0.426	-6.8	129	-0.04
60		Diethylphthalate	1.549	1.533	1.0	126	-0.04
61		4-Chlorophenyl-phenylether	0.721	0.735	-1.9	129	-0.04
62		4-Nitroaniline	0.349	0.383	-9.7	128	-0.03
63		Azobenzene	1.498	1.424	4.9	117	-0.04
64	I	Phenanthrene-d10	1.000	1.000	0.0	134	-0.04
65		4,6-Dinitro-2-methylphenol	0.133	0.151	-13.5	135	-0.02
66	c	n-Nitrosodiphenylamine	0.612	0.622	-1.6	126	-0.03
67		4-Bromophenyl-phenylether	0.220	0.230	-4.5	130	-0.05
68		Hexachlorobenzene	0.245	0.258	-5.3	133	-0.04
69		Atrazine	0.234	0.237	-1.3	122	-0.04
70	C	Pentachlorophenol	0.162	0.169	-4.3	127	-0.03
71		Phenanthrene	1.135	1.104	2.7	122	-0.04
72		Anthracene	1.151	1.138	1.1	124	-0.04
73		Carbazole	1.070	1.061	0.8	122	-0.04
74		Di-n-butylphthalate	1.313	1.329	-1.2	124	-0.04
75	C	Fluoranthene	1.361	1.336	1.8	124	-0.04
76	I	Chrysene-d12	1.000	1.000	0.0	131	-0.05
77		Benzidine	0.576	0.630	-9.4	126	-0.04
78		Pyrene	1.335	1.329	0.4	122	-0.04
79	S	Terphenyl-d14	1.112	1.133	-1.9	126	-0.04
80		Butylbenzylphthalate	0.585	0.594	-1.5	120	-0.05
81		Benzo(a)anthracene	1.369	1.359	0.7	123	-0.05
82		3,3'-Dichlorobenzidine	0.473	0.502	-6.1	127	-0.05
83		Chrysene	1.305	1.272	2.5	121	-0.05
84		Bis(2-ethylhexyl)phthalate	0.856	0.843	1.5	117	-0.05
85	c	Di-n-octyl phthalate	1.523	1.495	1.8	117	-0.06
86	I	Perylene-d12	1.000	1.000	0.0	136	-0.06
87		Indeno(1,2,3-cd)pyrene	1.454	1.488	-2.3	127	-0.08
88		Benzo(b)fluoranthene	1.223	1.194	2.4	122	-0.05
89		Benzo(k)fluoranthene	1.248	1.216	2.6	123	-0.05
90	C	Benzo(a)pyrene	1.198	1.178	1.7	123	-0.07
91		Dibenzo(a,h)anthracene	1.190	1.228	-3.2	129	-0.11
92		Benzo(g,h,i)perylene	1.170	1.195	-2.1	128	-0.08

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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