Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092625\
Data File : BP025869.D
Acq On : 26 Sep 2025 14:06
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Sep 26 14:40:45 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	122	-0.02
2		1,4-Dioxane	0.520	0.487	6.3	109	0.00
3		Pyridine	1.432	1.317	8.0	102	0.00
4		n-Nitrosodimethylamine	0.676	0.532	21.3	90	0.00
5	S	2-Fluorophenol	1.240	1.246	-0.5	112	0.00
6		Aniline	2.262	2.090	7.6	102	-0.01
7	S	Phenol-d6	1.648	1.571	4.7	103	0.00
8		2-Chlorophenol	1.360	1.357	0.2	110	-0.01
9		Benzaldehyde	0.987	1.197	-21.3	167#	-0.02
10	C	Phenol	1.737	1.629	6.2	101	0.00
11		bis(2-Chloroethyl)ether	1.396	1.305	6.5	102	-0.01
12		1,3-Dichlorobenzene	1.548	1.551	-0.2	113	-0.01
13	C	1,4-Dichlorobenzene	1.576	1.550	1.6	112	-0.01
14		1,2-Dichlorobenzene	1.531	1.511	1.3	111	-0.02
15		Benzyl Alcohol	1.361	1.249	8.2	101	-0.01
16		2,2'-oxybis(1-Chloropropane	2.282	1.825	20.0	89	-0.02
17		2-Methylphenol	1.171	1.109	5.3	102	0.00
18		Hexachloroethane	0.602	0.593	1.5	112	-0.02
19	P	n-Nitroso-di-n-propylamine	1.149	1.055	8.2	97	-0.02
20		3+4-Methylphenols	1.640	1.494	8.9	100	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	111	-0.01
22		Acetophenone	0.534	0.540	-1.1	101	-0.01
23	S	Nitrobenzene-d5	0.453	0.440	2.9	98	-0.01
24		Nitrobenzene	0.407	0.388	4.7	96	-0.02
25		Isophorone	0.797	0.768	3.6	99	-0.01
26	C	2-Nitrophenol	0.181	0.201	-11.0	109	0.00
27		2,4-Dimethylphenol	0.317	0.317	0.0	99	0.00
28		bis(2-Chloroethoxy)methane	0.461	0.459	0.4	102	-0.02
29	C	2,4-Dichlorophenol	0.316	0.335	-6.0	106	0.00
30		1,2,4-Trichlorobenzene	0.356	0.380	-6.7	111	-0.01
31		Naphthalene	1.078	1.064	1.3	103	-0.02
32		Benzoic acid	0.236	0.246	-4.2	104	0.01
33		4-Chloroaniline	0.449	0.460	-2.4	100	-0.01
34	C	Hexachlorobutadiene	0.229	0.259	-13.1	118	-0.02
35		Caprolactam	0.122	0.143	-17.2	119	0.01
36	C	4-Chloro-3-methylphenol	0.377	0.376	0.3	100	0.01
37		2-Methylnaphthalene	0.693	0.690	0.4	102	-0.01
38		1-Methylnaphthalene	0.739	0.737	0.3	102	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	111	-0.02
40		1,2,4,5-Tetrachlorobenzene	0.606	0.653	-7.8	110	0.00
41	P	Hexachlorocyclopentadiene	0.330	0.378	-14.5	110	-0.02
42	S	2,4,6-Tribromophenol	0.249	0.308	-23.7	124	-0.01
43	C	2,4,6-Trichlorophenol	0.398	0.429	-7.8	107	0.00
44		2,4,5-Trichlorophenol	0.443	0.470	-6.1	104	0.01
45	S	2-Fluorobiphenyl	1.502	1.532	-2.0	106	-0.02
46		1,1'-Biphenyl	1.468	1.485	-1.2	104	0.00
47		2-Chloronaphthalene	1.156	1.156	0.0	103	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.385	0.356	7.5	90	0.00
49		Acenaphthylene	1.915	1.902	0.7	102	-0.02
50		Dimethylphthalate	1.527	1.557	-2.0	104	-0.01
51		2,6-Dinitrotoluene	0.324	0.333	-2.8	103	0.00
52	C	Acenaphthene	1.104	1.086	1.6	101	-0.02
53		3-Nitroaniline	0.340	0.335	1.5	94	0.00
54	P	2,4-Dinitrophenol	0.183	0.219	-19.7	116	0.00
55		Dibenzofuran	1.800	1.793	0.4	103	-0.01
56	P	4-Nitrophenol	0.258	0.311	-20.5	112	0.05
57		2,4-Dinitrotoluene	0.460	0.491	-6.7	105	0.00
58		Fluorene	1.431	1.445	-1.0	104	-0.01
59		2,3,4,6-Tetrachlorophenol	0.399	0.444	-11.3	110	0.00
60		Diethylphthalate	1.549	1.580	-2.0	107	-0.01
61		4-Chlorophenyl-phenylether	0.721	0.768	-6.5	111	-0.01
62		4-Nitroaniline	0.349	0.360	-3.2	99	0.00
63		Azobenzene	1.498	1.379	7.9	93	-0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	114	0.00
65		4,6-Dinitro-2-methylphenol	0.133	0.153	-15.0	116	0.00
66	c	n-Nitrosodiphenylamine	0.612	0.604	1.3	104	0.00
67		4-Bromophenyl-phenylether	0.220	0.238	-8.2	114	-0.01
68		Hexachlorobenzene	0.245	0.273	-11.4	119	0.00
69		Atrazine	0.234	0.248	-6.0	109	0.00
70	C	Pentachlorophenol	0.162	0.173	-6.8	110	0.00
71		Phenanthrene	1.135	1.096	3.4	103	0.00
72		Anthracene	1.151	1.138	1.1	105	0.00
73		Carbazole	1.070	1.064	0.6	103	0.00
74		Di-n-butylphthalate	1.313	1.357	-3.4	108	0.00
75	C	Fluoranthene	1.361	1.376	-1.1	108	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	118	-0.01
77		Benzidine	0.576	0.590	-2.4	106	0.00
78		Pyrene	1.335	1.283	3.9	106	0.00
79	S	Terphenyl-d14	1.112	1.145	-3.0	114	-0.01
80		Butylbenzylphthalate	0.585	0.598	-2.2	109	-0.02
81		Benzo(a)anthracene	1.369	1.379	-0.7	112	0.00
82		3,3'-Dichlorobenzidine	0.473	0.505	-6.8	115	0.00
83		Chrysene	1.305	1.290	1.1	110	-0.01
84		Bis(2-ethylhexyl)phthalate	0.856	0.877	-2.5	109	-0.02
85	c	Di-n-octyl phthalate	1.523	1.564	-2.7	110	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	121	0.00
87		Indeno(1,2,3-cd)pyrene	1.454	1.493	-2.7	113	0.00
88		Benzo(b)fluoranthene	1.223	1.218	0.4	110	0.00
89		Benzo(k)fluoranthene	1.248	1.251	-0.2	113	0.00
90	C	Benzo(a)pyrene	1.198	1.204	-0.5	111	0.00
91		Dibenzo(a,h)anthracene	1.190	1.234	-3.7	115	-0.01
92		Benzo(g,h,i)perylene	1.170	1.207	-3.2	115	0.01

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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