Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072325\
Data File : BP025219.D
Acq On : 23 Jul 2025 11:22
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jul 23 12:04:55 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072125.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 22 02:55:21 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	65	0.00
2		1,4-Dioxane	0.444	0.442	0.5	67	0.00
3		Pyridine	1.188	1.182	0.5	66	0.00
4		n-Nitrosodimethylamine	0.471	0.467	0.8	65	0.00
5	S	2-Fluorophenol	1.184	1.179	0.4	66	0.00
6		Aniline	1.916	1.952	-1.9	66	0.00
7	S	Phenol-d6	1.466	1.476	-0.7	65	0.00
8		2-Chlorophenol	1.343	1.350	-0.5	66	0.00
9		Benzaldehyde	0.994	0.962	3.2	56	0.00
10	C	Phenol	1.520	1.537	-1.1	66	0.00
11		bis(2-Chloroethyl)ether	1.166	1.175	-0.8	66	0.00
12		1,3-Dichlorobenzene	1.510	1.485	1.7	65	0.00
13	C	1,4-Dichlorobenzene	1.525	1.514	0.7	65	0.00
14		1,2-Dichlorobenzene	1.467	1.463	0.3	66	0.00
15		Benzyl Alcohol	1.057	1.081	-2.3	66	0.00
16		2,2'-oxybis(1-Chloropropane	1.491	1.471	1.3	64	0.00
17		2-Methylphenol	1.054	1.073	-1.8	66	0.00
18		Hexachloroethane	0.524	0.519	1.0	65	0.00
19	P	n-Nitroso-di-n-propylamine	0.882	0.913	-3.5	67	0.00
20		3+4-Methylphenols	1.438	1.469	-2.2	66	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	65	0.00
22		Acetophenone	0.465	0.463	0.4	67	0.00
23	S	Nitrobenzene-d5	0.360	0.360	0.0	66	0.00
24		Nitrobenzene	0.321	0.322	-0.3	66	0.00
25		Isophorone	0.633	0.629	0.6	65	0.00
26	C	2-Nitrophenol	0.184	0.188	-2.2	67	0.00
27		2,4-Dimethylphenol	0.308	0.308	0.0	66	0.00
28		bis(2-Chloroethoxy)methane	0.396	0.392	1.0	66	0.00
29	C	2,4-Dichlorophenol	0.317	0.318	-0.3	66	0.01
30		1,2,4-Trichlorobenzene	0.346	0.336	2.9	65	0.00
31		Naphthalene	1.023	1.020	0.3	66	0.00
32		Benzoic acid	0.235	0.226	3.8	62	0.00
33		4-Chloroaniline	0.450	0.456	-1.3	68	0.00
34	C	Hexachlorobutadiene	0.208	0.203	2.4	65	0.00
35		Caprolactam	0.105	0.106	-1.0	68	0.00
36	C	4-Chloro-3-methylphenol	0.317	0.321	-1.3	68	0.00
37		2-Methylnaphthalene	0.665	0.659	0.9	66	0.00
38		1-Methylnaphthalene	0.698	0.698	0.0	66	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	66	0.00
40		1,2,4,5-Tetrachlorobenzene	0.599	0.599	0.0	67	0.00
41	P	Hexachlorocyclopentadiene	0.358	0.377	-5.3	70	0.00
42	S	2,4,6-Tribromophenol	0.295	0.297	-0.7	68	-0.01
43	C	2,4,6-Trichlorophenol	0.410	0.416	-1.5	67	0.00
44		2,4,5-Trichlorophenol	0.451	0.455	-0.9	66	0.00
45	S	2-Fluorobiphenyl	1.419	1.437	-1.3	67	0.00
46		1,1'-Biphenyl	1.450	1.460	-0.7	67	0.00
47		2-Chloronaphthalene	1.132	1.135	-0.3	67	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.289	0.300	-3.8	68	-0.01
49		Acenaphthylene	1.831	1.855	-1.3	67	-0.01
50		Dimethylphthalate	1.430	1.428	0.1	68	0.00
51		2,6-Dinitrotoluene	0.311	0.315	-1.3	67	0.00
52	C	Acenaphthene	1.059	1.059	0.0	67	0.00
53		3-Nitroaniline	0.350	0.351	-0.3	67	0.00
54	P	2,4-Dinitrophenol	0.184	0.191	-3.8	69	0.00
55		Dibenzofuran	1.710	1.713	-0.2	67	0.00
56	P	4-Nitrophenol	0.300	0.303	-1.0	69	0.00
57		2,4-Dinitrotoluene	0.435	0.448	-3.0	68	0.00
58		Fluorene	1.358	1.377	-1.4	68	0.00
59		2,3,4,6-Tetrachlorophenol	0.399	0.404	-1.3	68	0.00
60		Diethylphthalate	1.410	1.410	0.0	67	0.00
61		4-Chlorophenyl-phenylether	0.686	0.685	0.1	67	0.00
62		4-Nitroaniline	0.356	0.368	-3.4	70	0.00
63		Azobenzene	1.149	1.151	-0.2	67	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	69	0.00
65		4,6-Dinitro-2-methylphenol	0.129	0.133	-3.1	70	0.00
66	c	n-Nitrosodiphenylamine	0.607	0.594	2.1	68	0.00
67		4-Bromophenyl-phenylether	0.233	0.229	1.7	68	0.00
68		Hexachlorobenzene	0.277	0.268	3.2	66	0.00
69		Atrazine	0.225	0.221	1.8	69	-0.01
70	C	Pentachlorophenol	0.188	0.185	1.6	67	0.00
71		Phenanthrene	1.079	1.056	2.1	68	0.00
72		Anthracene	1.087	1.071	1.5	68	-0.01
73		Carbazole	1.037	1.015	2.1	68	0.00
74		Di-n-butylphthalate	1.197	1.238	-3.4	71	0.00
75	C	Fluoranthene	1.270	1.247	1.8	69	0.02
76	I	Chrysene-d12	1.000	1.000	0.0	70	0.01
77		Benzidine	0.702	0.782	-11.4	64	0.01
78		Pyrene	1.233	1.269	-2.9	70	0.01
79	S	Terphenyl-d14	1.013	1.053	-3.9	82	0.00
80		Butylbenzylphthalate	0.558	0.561	-0.5	69	0.00
81		Benzo(a)anthracene	1.263	1.259	0.3	70	0.00
82		3,3'-Dichlorobenzidine	0.524	0.522	0.4	71	0.02
83		Chrysene	1.179	1.202	-2.0	72	0.01
84		Bis(2-ethylhexyl)phthalate	0.798	0.794	0.5	69	0.01
85	c	Di-n-octyl phthalate	1.378	1.420	-3.0	71	0.02
86	I	Perylene-d12	1.000	1.000	0.0	71	0.01
87		Indeno(1,2,3-cd)pyrene	1.462	1.474	-0.8	74	0.04
88		Benzo(b)fluoranthene	1.145	1.121	2.1	70	0.02
89		Benzo(k)fluoranthene	1.130	1.116	1.2	70	0.01
90	C	Benzo(a)pyrene	1.117	1.103	1.3	71	0.02
91		Dibenzo(a,h)anthracene	1.192	1.207	-1.3	73	0.04
92		Benzo(g,h,i)perylene	1.172	1.178	-0.5	73	0.03

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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