Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061725\
Data File : BP024971.D
Acq On : 17 Jun 2025 11:54
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jun 17 12:29:21 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 06 16:20:27 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	102	0.00
2		1,4-Dioxane	0.527	0.515	2.3	106	0.00
3		Pyridine	1.268	1.291	-1.8	107	0.00
4		n-Nitrosodimethylamine	0.518	0.498	3.9	101	0.00
5	S	2-Fluorophenol	1.198	1.218	-1.7	106	0.00
6		Aniline	2.023	2.018	0.2	104	0.00
7	S	Phenol-d6	1.585	1.549	2.3	102	0.00
8		2-Chlorophenol	1.358	1.337	1.5	103	0.00
9		Benzaldehyde	0.914	0.896	2.0	104	0.00
10	C	Phenol	1.634	1.587	2.9	102	0.00
11		bis(2-Chloroethyl)ether	1.285	1.257	2.2	104	0.00
12		1,3-Dichlorobenzene	1.515	1.465	3.3	104	0.00
13	C	1,4-Dichlorobenzene	1.529	1.480	3.2	105	0.00
14		1,2-Dichlorobenzene	1.501	1.426	5.0	103	0.00
15		Benzyl Alcohol	1.217	1.191	2.1	104	0.00
16		2,2'-oxybis(1-Chloropropane	1.682	1.622	3.6	104	0.00
17		2-Methylphenol	1.141	1.087	4.7	100	0.00
18		Hexachloroethane	0.576	0.553	4.0	103	0.00
19	P	n-Nitroso-di-n-propylamine	1.078	1.017	5.7	99	0.00
20		3+4-Methylphenols	1.557	1.466	5.8	100	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
22		Acetophenone	0.505	0.484	4.2	97	0.00
23	S	Nitrobenzene-d5	0.412	0.406	1.5	99	0.00
24		Nitrobenzene	0.366	0.358	2.2	98	0.00
25		Isophorone	0.713	0.708	0.7	101	0.00
26	C	2-Nitrophenol	0.180	0.185	-2.8	101	0.00
27		2,4-Dimethylphenol	0.309	0.303	1.9	99	0.00
28		bis(2-Chloroethoxy)methane	0.426	0.421	1.2	100	0.00
29	C	2,4-Dichlorophenol	0.296	0.295	0.3	100	0.01
30		1,2,4-Trichlorobenzene	0.334	0.317	5.1	99	0.00
31		Naphthalene	1.025	0.992	3.2	99	0.00
32		Benzoic acid	0.209	0.200	4.3	97	0.02
33		4-Chloroaniline	0.429	0.430	-0.2	100	0.00
34	C	Hexachlorobutadiene	0.201	0.196	2.5	100	0.00
35		Caprolactam	0.109	0.114	-4.6	102	0.02
36	C	4-Chloro-3-methylphenol	0.342	0.338	1.2	98	0.00
37		2-Methylnaphthalene	0.650	0.637	2.0	99	0.00
38		1-Methylnaphthalene	0.695	0.672	3.3	99	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
40		1,2,4,5-Tetrachlorobenzene	0.568	0.541	4.8	98	0.00
41	P	Hexachlorocyclopentadiene	0.346	0.303	12.4	88	0.00
42	S	2,4,6-Tribromophenol	0.277	0.261	5.8	95	0.00
43	C	2,4,6-Trichlorophenol	0.381	0.370	2.9	96	0.01
44		2,4,5-Trichlorophenol	0.408	0.403	1.2	97	0.02
45	S	2-Fluorobiphenyl	1.485	1.409	5.1	99	0.00
46		1,1'-Biphenyl	1.453	1.397	3.9	98	0.00
47		2-Chloronaphthalene	1.117	1.073	3.9	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061725\
Data File : BP024971.D
Acq On : 17 Jun 2025 11:54
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jun 17 12:29:21 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 06 16:20:27 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)
48		2-Nitroaniline	0.343	0.351	-2.3	99	0.01
49		Acenaphthylene	1.863	1.794	3.7	99	0.00
50		Dimethylphthalate	1.475	1.392	5.6	97	0.00
51		2,6-Dinitrotoluene	0.318	0.310	2.5	97	0.00
52	C	Acenaphthene	1.067	1.005	5.8	96	0.00
53		3-Nitroaniline	0.330	0.337	-2.1	98	0.00
54	P	2,4-Dinitrophenol	0.178	0.171	3.9	93	0.02
55		Dibenzofuran	1.713	1.600	6.6	96	0.00
56	P	4-Nitrophenol	0.239	0.271	-13.4	107	0.04
57		2,4-Dinitrotoluene	0.445	0.440	1.1	98	0.02
58		Fluorene	1.384	1.314	5.1	98	0.00
59		2,3,4,6-Tetrachlorophenol	0.365	0.343	6.0	95	0.00
60		Diethylphthalate	1.470	1.423	3.2	100	0.00
61		4-Chlorophenyl-phenylether	0.677	0.626	7.5	96	0.00
62		4-Nitroaniline	0.299	0.334	-11.7	105	0.00
63		Azobenzene	1.349	1.319	2.2	99	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	97	0.00
65		4,6-Dinitro-2-methylphenol	0.131	0.132	-0.8	98	0.01
66	c	n-Nitrosodiphenylamine	0.620	0.603	2.7	98	0.00
67		4-Bromophenyl-phenylether	0.226	0.217	4.0	98	0.00
68		Hexachlorobenzene	0.274	0.259	5.5	96	0.00
69		Atrazine	0.225	0.222	1.3	99	0.00
70	C	Pentachlorophenol	0.142	0.145	-2.1	101	0.00
71		Phenanthrene	1.105	1.049	5.1	96	0.00
72		Anthracene	1.119	1.083	3.2	98	0.00
73		Carbazole	1.038	1.044	-0.6	101	0.00
74		Di-n-butylphthalate	1.285	1.306	-1.6	99	-0.02
75	C	Fluoranthene	1.281	1.223	4.5	97	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	94	0.02
77		Benzidine	0.619	0.666	-7.6	96	0.00
78		Pyrene	1.249	1.218	2.5	97	0.00
79	S	Terphenyl-d14	1.116	1.069	4.2	92	0.00
80		Butylbenzylphthalate	0.572	0.601	-5.1	101	0.00
81		Benzo(a)anthracene	1.279	1.246	2.6	96	0.02
82		3,3'-Dichlorobenzidine	0.508	0.512	-0.8	98	0.00
83		Chrysene	1.212	1.173	3.2	96	0.02
84		Bis(2-ethylhexyl)phthalate	0.820	0.844	-2.9	99	0.00
85	c	Di-n-octyl phthalate	1.445	1.533	-6.1	104	0.00
86	I	Perylene-d12	1.000	1.000	0.0	98	0.05
87		Indeno(1,2,3-cd)pyrene	1.459	1.412	3.2	98	0.07
88		Benzo(b)fluoranthene	1.145	1.087	5.1	95	0.03
89		Benzo(k)fluoranthene	1.165	1.092	6.3	97	0.03
90	C	Benzo(a)pyrene	1.118	1.081	3.3	100	0.04
91		Dibenzo(a,h)anthracene	1.188	1.150	3.2	99	0.08
92		Benzo(g,h,i)perylene	1.179	1.126	4.5	98	0.09

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061725\
Data File : BP024971.D
Acq On : 17 Jun 2025 11:54
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jun 17 12:29:21 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 06 16:20:27 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)

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