Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061821\
Data File : BP005970.D
Acq On : 18 Jun 2021 14:14
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Jun 18 14:50:25 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 16 11:35:30 2021
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	83	0.00
2		1,4-Dioxane	0.563	0.570	-1.2	88	0.00
3		Pyridine	1.323	1.521	-15.0	95	0.00
4		n-Nitrosodimethylamine	0.615	0.607	1.3	86	0.00
5	S	2-Fluorophenol	1.246	1.342	-7.7	93	0.00
6		Aniline	1.804	1.975	-9.5	95	0.00
7	S	Phenol-d6	1.616	1.760	-8.9	94	0.00
8		2-Chlorophenol	1.429	1.562	-9.3	96	0.00
9		Benzaldehyde	0.689	0.802	-16.4	92	0.00
10	C	Phenol	1.614	1.764	-9.3	96	0.00
11		bis(2-Chloroethyl)ether	1.223	1.356	-10.9	97	0.00
12		1,3-Dichlorobenzene	1.598	1.731	-8.3	95	0.00
13	C	1,4-Dichlorobenzene	1.630	1.755	-7.7	95	0.00
14		1,2-Dichlorobenzene	1.558	1.686	-8.2	95	0.00
15		Benzyl Alcohol	1.217	1.294	-6.3	91	0.00
16		2,2'-oxybis(1-Chloropropane	1.130	1.178	-4.2	91	-0.01
17		2-Methylphenol	1.100	1.225	-11.4	97	0.00
18		Hexachloroethane	0.589	0.632	-7.3	93	0.00
19	P	n-Nitroso-di-n-propylamine	0.994	1.091	-9.8	95	0.00
20		3+4-Methylphenols	1.485	1.649	-11.0	96	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	83	0.00
22		Acetophenone	0.526	0.561	-6.7	93	0.00
23	S	Nitrobenzene-d5	0.408	0.422	-3.4	90	0.00
24		Nitrobenzene	0.424	0.436	-2.8	92	0.00
25		Isophorone	0.753	0.791	-5.0	93	0.00
26	C	2-Nitrophenol	0.202	0.217	-7.4	95	0.00
27		2,4-Dimethylphenol	0.301	0.317	-5.3	95	0.00
28		bis(2-Chloroethoxy)methane	0.393	0.426	-8.4	99	0.00
29	C	2,4-Dichlorophenol	0.362	0.375	-3.6	93	0.00
30		1,2,4-Trichlorobenzene	0.407	0.418	-2.7	94	0.00
31		Naphthalene	1.157	1.220	-5.4	96	0.00
32		Benzoic acid	0.193	0.206	-6.7	82	0.00
33		4-Chloroaniline	0.467	0.498	-6.6	96	0.00
34	C	Hexachlorobutadiene	0.275	0.283	-2.9	95	0.00
35		Caprolactam	0.104	0.114	-9.6	92	0.00
36	C	4-Chloro-3-methylphenol	0.367	0.394	-7.4	97	0.00
37		2-Methylnaphthalene	0.824	0.863	-4.7	96	0.00
38		1-Methylnaphthalene	0.776	0.821	-5.8	96	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	82	0.00
40		1,2,4,5-Tetrachlorobenzene	0.704	0.731	-3.8	90	0.00
41	P	Hexachlorocyclopentadiene	0.297	0.286	3.7	80	0.00
42	S	2,4,6-Tribromophenol	0.343	0.368	-7.3	93	0.00
43	C	2,4,6-Trichlorophenol	0.483	0.493	-2.1	89	0.00
44		2,4,5-Trichlorophenol	0.520	0.531	-2.1	89	0.00
45	S	2-Fluorobiphenyl	1.523	1.589	-4.3	92	0.00
46		1,1'-Biphenyl	1.613	1.694	-5.0	91	0.00
47		2-Chloronaphthalene	1.317	1.378	-4.6	95	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061821\
Data File : BP005970.D
Acq On : 18 Jun 2021 14:14
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Jun 18 14:50:25 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 16 11:35:30 2021
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	0.346	0.368	-6.4	93	0.00
49		Acenaphthylene	2.020	2.136	-5.7	94	0.00
50		Dimethylphthalate	1.644	1.721	-4.7	94	0.00
51		2,6-Dinitrotoluene	0.374	0.390	-4.3	92	0.00
52	C	Acenaphthene	1.227	1.281	-4.4	94	0.00
53		3-Nitroaniline	0.362	0.393	-8.6	95	0.00
54	P	2,4-Dinitrophenol	0.191	0.218	-14.1	92	0.00
55		Dibenzofuran	2.018	2.067	-2.4	92	0.00
56	P	4-Nitrophenol	0.294	0.282	4.1	82	0.00
57		2,4-Dinitrotoluene	0.500	0.531	-6.2	92	0.00
58		Fluorene	1.572	1.626	-3.4	93	0.00
59		2,3,4,6-Tetrachlorophenol	0.458	0.470	-2.6	90	0.00
60		Diethylphthalate	1.650	1.706	-3.4	92	0.00
61		4-Chlorophenyl-phenylether	0.815	0.839	-2.9	93	0.00
62		4-Nitroaniline	0.373	0.405	-8.6	95	0.00
63		Azobenzene	1.493	1.516	-1.5	90	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	79	0.00
65		4,6-Dinitro-2-methylphenol	0.147	0.136	7.5	78	0.00
66	c	n-Nitrosodiphenylamine	0.663	0.706	-6.5	93	0.00
67		4-Bromophenyl-phenylether	0.257	0.273	-6.2	92	0.00
68		Hexachlorobenzene	0.308	0.325	-5.5	92	0.00
69		Atrazine	0.206	0.186	9.7	67	0.00
70	C	Pentachlorophenol	0.171	0.176	-2.9	81	0.00
71		Phenanthrene	1.201	1.263	-5.2	91	-0.01
72		Anthracene	1.182	1.257	-6.3	92	0.00
73		Carbazole	1.135	1.223	-7.8	93	0.00
74		Di-n-butylphthalate	1.315	1.422	-8.1	93	0.00
75	C	Fluoranthene	1.459	1.568	-7.5	94	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	89	0.00
77		Benzidine	0.416	0.385	7.5	72	0.00
78		Pyrene	1.396	1.359	2.7	95	0.00
79	S	Terphenyl-d14	1.068	1.083	-1.4	98	0.00
80		Butylbenzylphthalate	0.574	0.601	-4.7	100	0.00
81		Benzo(a)anthracene	1.444	1.481	-2.6	101	0.00
82		3,3'-Dichlorobenzidine	0.499	0.526	-5.4	105	0.00
83		Chrysene	1.398	1.449	-3.6	102	0.00
84		Bis(2-ethylhexyl)phthalate	0.842	0.877	-4.2	101	0.00
85	c	Di-n-octyl phthalate	1.507	1.582	-5.0	101	0.00
86	I	Perylene-d12	1.000	1.000	0.0	87	0.00
87		Indeno(1,2,3-cd)pyrene	1.710	1.865	-9.1	104	-0.01
88		Benzo(b)fluoranthene	1.370	1.481	-8.1	102	-0.01
89		Benzo(k)fluoranthene	1.328	1.480	-11.4	106	0.00
90	C	Benzo(a)pyrene	1.288	1.406	-9.2	104	0.00
91		Dibenzo(a,h)anthracene	1.472	1.602	-8.8	104	-0.01
92		Benzo(g,h,i)perylene	1.454	1.553	-6.8	102	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061821\
Data File : BP005970.D
Acq On : 18 Jun 2021 14:14
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jun 18 14:50:25 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 16 11:35:30 2021
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