Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030822\
Data File : BP009297.D
Acq On : 08 Mar 2022 10:45
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Mar 09 02:59:36 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Mar 05 01:55:59 2022
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	125	0.00
2		1,4-Dioxane	0.583	0.507	13.0	121	0.00
3		Pyridine	1.243	1.128	9.3	122	0.00
4		n-Nitrosodimethylamine	0.518	0.500	3.5	127	0.00
5	S	2-Fluorophenol	1.370	1.246	9.1	125	0.00
6		Aniline	2.205	2.007	9.0	124	0.00
7	S	Phenol-d6	1.741	1.585	9.0	125	0.00
8		2-Chlorophenol	1.404	1.307	6.9	126	0.00
9		Benzaldehyde	1.109	1.008	9.1	122	0.00
10	C	Phenol	1.803	1.633	9.4	124	0.00
11		bis(2-Chloroethyl)ether	1.395	1.274	8.7	123	0.00
12		1,3-Dichlorobenzene	1.654	1.478	10.6	125	0.00
13	C	1,4-Dichlorobenzene	1.678	1.505	10.3	125	0.00
14		1,2-Dichlorobenzene	1.603	1.437	10.4	124	0.00
15		Benzyl Alcohol	1.228	1.146	6.7	126	0.00
16		2,2'-oxybis(1-Chloropropane	1.811	1.541	14.9	118	0.00
17		2-Methylphenol	1.167	1.075	7.9	125	0.00
18		Hexachloroethane	0.584	0.536	8.2	127	0.00
19	P	n-Nitroso-di-n-propylamine	1.050	0.985	6.2	126	0.00
20		3+4-Methylphenols	1.564	1.445	7.6	125	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	124	0.00
22		Acetophenone	0.535	0.500	6.5	125	0.00
23	S	Nitrobenzene-d5	0.363	0.357	1.7	127	0.00
24		Nitrobenzene	0.385	0.369	4.2	126	0.00
25		Isophorone	0.679	0.652	4.0	126	0.00
26	C	2-Nitrophenol	0.149	0.168	-12.8	141	0.00
27		2,4-Dimethylphenol	0.328	0.303	7.6	123	0.00
28		bis(2-Chloroethoxy)methane	0.444	0.418	5.9	125	0.00
29	C	2,4-Dichlorophenol	0.316	0.307	2.8	126	0.00
30		1,2,4-Trichlorobenzene	0.370	0.347	6.2	128	0.00
31		Naphthalene	1.108	1.016	8.3	124	0.00
32		Benzoic acid	0.167	0.203	-21.6	155#	0.02
33		4-Chloroaniline	0.453	0.426	6.0	125	0.00
34	C	Hexachlorobutadiene	0.225	0.217	3.6	131	0.00
35		Caprolactam	0.093	0.096	-3.2	129	0.00
36	C	4-Chloro-3-methylphenol	0.330	0.313	5.2	125	0.00
37		2-Methylnaphthalene	0.787	0.727	7.6	124	0.00
38		1-Methylnaphthalene	0.725	0.667	8.0	125	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	125	0.00
40		1,2,4,5-Tetrachlorobenzene	0.674	0.651	3.4	130	0.00
41	P	Hexachlorocyclopentadiene	0.427	0.420	1.6	131	0.00
42	S	2,4,6-Tribromophenol	0.242	0.265	-9.5	142	0.00
43	C	2,4,6-Trichlorophenol	0.405	0.410	-1.2	131	0.00
44		2,4,5-Trichlorophenol	0.444	0.439	1.1	128	0.00
45	S	2-Fluorobiphenyl	1.454	1.361	6.4	127	0.00
46		1,1'-Biphenyl	1.633	1.511	7.5	125	0.00
47		2-Chloronaphthalene	1.237	1.157	6.5	126	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030822\
Data File : BP009297.D
Acq On : 08 Mar 2022 10:45
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Mar 09 02:59:36 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Mar 05 01:55:59 2022
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	0.284	0.310	-9.2	135	0.00
49		Acenaphthylene	1.911	1.817	4.9	126	0.00
50		Dimethylphthalate	1.475	1.395	5.4	127	0.00
51		2,6-Dinitrotoluene	0.291	0.301	-3.4	130	0.00
52	C	Acenaphthene	1.225	1.157	5.6	126	0.00
53		3-Nitroaniline	0.309	0.316	-2.3	129	0.00
54	P	2,4-Dinitrophenol	0.125	0.152	-21.6	150#	0.00
55		Dibenzofuran	1.855	1.704	8.1	125	0.00
56	P	4-Nitrophenol	0.257	0.263	-2.3	127	0.00
57		2,4-Dinitrotoluene	0.373	0.403	-8.0	132	0.00
58		Fluorene	1.453	1.348	7.2	127	0.00
59		2,3,4,6-Tetrachlorophenol	0.374	0.375	-0.3	132	0.00
60		Diethylphthalate	1.436	1.386	3.5	128	0.00
61		4-Chlorophenyl-phenylether	0.740	0.699	5.5	129	0.00
62		4-Nitroaniline	0.300	0.316	-5.3	131	0.00
63		Azobenzene	1.411	1.337	5.2	125	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	126	0.00
65		4,6-Dinitro-2-methylphenol	0.088	0.104	-18.2	145	0.00
66	c	n-Nitrosodiphenylamine	0.628	0.581	7.5	124	0.00
67		4-Bromophenyl-phenylether	0.216	0.219	-1.4	136	0.00
68		Hexachlorobenzene	0.253	0.248	2.0	132	0.00
69		Atrazine	0.221	0.223	-0.9	130	0.00
70	C	Pentachlorophenol	0.152	0.160	-5.3	133	0.00
71		Phenanthrene	1.149	1.050	8.6	124	0.00
72		Anthracene	1.144	1.068	6.6	125	0.00
73		Carbazole	1.042	0.965	7.4	122	0.00
74		Di-n-butylphthalate	1.154	1.167	-1.1	129	0.00
75	C	Fluoranthene	1.327	1.230	7.3	122	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	125	0.00
77		Benzidine	0.698	0.758	-8.6	128	0.00
78		Pyrene	1.350	1.239	8.2	122	0.00
79	S	Terphenyl-d14	1.060	0.992	6.4	126	0.00
80		Butylbenzylphthalate	0.487	0.519	-6.6	132	0.00
81		Benzo(a)anthracene	1.356	1.243	8.3	123	0.00
82		3,3'-Dichlorobenzidine	0.498	0.502	-0.8	127	0.00
83		Chrysene	1.316	1.208	8.2	124	0.00
84		Bis(2-ethylhexyl)phthalate	0.780	0.799	-2.4	129	0.00
85	c	Di-n-octyl phthalate	1.276	1.391	-9.0	134	0.00
86	I	Perylene-d12	1.000	1.000	0.0	126	0.00
87		Indeno(1,2,3-cd)pyrene	1.554	1.521	2.1	131	0.00
88		Benzo(b)fluoranthene	1.350	1.273	5.7	124	0.00
89		Benzo(k)fluoranthene	1.290	1.204	6.7	127	0.00
90	C	Benzo(a)pyrene	1.281	1.222	4.6	126	0.00
91		Dibenzo(a,h)anthracene	1.318	1.294	1.8	130	0.00
92		Benzo(g,h,i)perylene	1.302	1.280	1.7	131	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030822\
Data File : BP009297.D
Acq On : 08 Mar 2022 10:45
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Mar 09 02:59:36 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Mar 05 01:55:59 2022
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