Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080221\
Data File : BP006464.D
Acq On : 03 Aug 2021 00:03
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Aug 03 05:27:59 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 03 05:26:21 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	133	0.00
2		1,4-Dioxane	0.568	0.549	3.3	129	0.00
3		Pyridine	1.641	1.627	0.9	130	0.00
4		n-Nitrosodimethylamine	0.620	0.660	-6.5	140	0.00
5	S	2-Fluorophenol	1.282	1.291	-0.7	133	0.00
6		Aniline	1.979	2.008	-1.5	132	0.00
7	S	Phenol-d6	1.811	1.841	-1.7	132	0.00
8		2-Chlorophenol	1.387	1.398	-0.8	133	0.00
9		Benzaldehyde	1.135	1.095	3.5	137	0.00
10	C	Phenol	1.824	1.848	-1.3	132	0.00
11		bis(2-Chloroethyl)ether	1.396	1.410	-1.0	134	0.00
12		1,3-Dichlorobenzene	1.464	1.451	0.9	133	0.00
13	C	1,4-Dichlorobenzene	1.481	1.475	0.4	133	0.00
14		1,2-Dichlorobenzene	1.423	1.410	0.9	134	0.00
15		Benzyl Alcohol	1.356	1.432	-5.6	137	0.00
16		2,2'-oxybis(1-Chloropropane	1.648	1.674	-1.6	136	0.00
17		2-Methylphenol	1.141	1.190	-4.3	134	0.00
18		Hexachloroethane	0.580	0.602	-3.8	139	0.00
19	P	n-Nitroso-di-n-propylamine	1.201	1.273	-6.0	136	0.00
20		3+4-Methylphenols	1.557	1.625	-4.4	135	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	133	0.00
22		Acetophenone	0.527	0.529	-0.4	134	0.00
23	S	Nitrobenzene-d5	0.424	0.437	-3.1	136	0.00
24		Nitrobenzene	0.440	0.453	-3.0	135	0.00
25		Isophorone	0.757	0.788	-4.1	135	0.00
26	C	2-Nitrophenol	0.158	0.166	-5.1	136	0.00
27		2,4-Dimethylphenol	0.270	0.280	-3.7	133	0.00
28		bis(2-Chloroethoxy)methane	0.414	0.417	-0.7	134	0.00
29	C	2,4-Dichlorophenol	0.273	0.285	-4.4	136	0.00
30		1,2,4-Trichlorobenzene	0.296	0.298	-0.7	138	0.00
31		Naphthalene	1.065	1.066	-0.1	135	0.00
32		Benzoic acid	0.205	0.215	-4.9	139	0.02
33		4-Chloroaniline	0.426	0.437	-2.6	132	0.00
34	C	Hexachlorobutadiene	0.173	0.177	-2.3	140	0.00
35		Caprolactam	0.100	0.102	-2.0	125	0.00
36	C	4-Chloro-3-methylphenol	0.340	0.362	-6.5	134	0.00
37		2-Methylnaphthalene	0.715	0.728	-1.8	134	0.00
38		1-Methylnaphthalene	0.679	0.693	-2.1	135	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	132	0.00
40		1,2,4,5-Tetrachlorobenzene	0.521	0.509	2.3	134	0.00
41	P	Hexachlorocyclopentadiene	0.276	0.287	-4.0	139	0.00
42	S	2,4,6-Tribromophenol	0.205	0.212	-3.4	131	0.00
43	C	2,4,6-Trichlorophenol	0.350	0.358	-2.3	133	0.00
44		2,4,5-Trichlorophenol	0.387	0.394	-1.8	132	0.00
45	S	2-Fluorobiphenyl	1.334	1.301	2.5	132	0.00
46		1,1'-Biphenyl	1.512	1.470	2.8	132	0.00
47		2-Chloronaphthalene	1.164	1.130	2.9	132	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080221\
Data File : BP006464.D
Acq On : 03 Aug 2021 00:03
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Aug 03 05:27:59 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 03 05:26:21 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	0.390	0.422	-8.2	133	0.00
49		Acenaphthylene	1.850	1.855	-0.3	130	0.00
50		Dimethylphthalate	1.438	1.431	0.5	129	0.00
51		2,6-Dinitrotoluene	0.313	0.322	-2.9	128	0.00
52	C	Acenaphthene	1.141	1.127	1.2	131	0.00
53		3-Nitroaniline	0.347	0.364	-4.9	127	0.00
54	P	2,4-Dinitrophenol	0.170	0.172	-1.2	129	0.00
55		Dibenzofuran	1.787	1.754	1.8	130	0.00
56	P	4-Nitrophenol	0.315	0.318	-1.0	124	0.00
57		2,4-Dinitrotoluene	0.418	0.446	-6.7	128	0.00
58		Fluorene	1.419	1.422	-0.2	130	0.00
59		2,3,4,6-Tetrachlorophenol	0.325	0.333	-2.5	131	0.00
60		Diethylphthalate	1.495	1.536	-2.7	131	0.00
61		4-Chlorophenyl-phenylether	0.645	0.644	0.2	132	0.00
62		4-Nitroaniline	0.367	0.389	-6.0	124	0.00
63		Azobenzene	1.726	1.842	-6.7	133	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	127	0.00
65		4,6-Dinitro-2-methylphenol	0.120	0.119	0.8	127	0.00
66	c	n-Nitrosodiphenylamine	0.619	0.623	-0.6	130	0.00
67		4-Bromophenyl-phenylether	0.192	0.193	-0.5	131	0.00
68		Hexachlorobenzene	0.216	0.215	0.5	131	0.00
69		Atrazine	0.191	0.200	-4.7	127	0.00
70	C	Pentachlorophenol	0.134	0.143	-6.7	130	0.00
71		Phenanthrene	1.122	1.098	2.1	126	0.00
72		Anthracene	1.075	1.084	-0.8	126	0.00
73		Carbazole	1.092	1.098	-0.5	124	0.00
74		Di-n-butylphthalate	1.233	1.328	-7.7	130	0.00
75	C	Fluoranthene	1.248	1.243	0.4	122	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	117	0.00
77		Benzidine	0.492	0.535	-8.7	113	0.00
78		Pyrene	1.319	1.333	-1.1	121	0.00
79	S	Terphenyl-d14	0.993	0.993	0.0	123	0.00
80		Butylbenzylphthalate	0.570	0.624	-9.5	124	0.00
81		Benzo(a)anthracene	1.288	1.284	0.3	118	0.00
82		3,3'-Dichlorobenzidine	0.341	0.355	-4.1	118	0.00
83		Chrysene	1.255	1.244	0.9	118	0.00
84		Bis(2-ethylhexyl)phthalate	0.880	0.964	-9.5	127	0.00
85	c	Di-n-octyl phthalate	1.490	1.605	-7.7	128	0.00
86	I	Perylene-d12	1.000	1.000	0.0	116	0.00
87		Indeno(1,2,3-cd)pyrene	1.435	1.463	-2.0	117	0.00
88		Benzo(b)fluoranthene	1.284	1.291	-0.5	118	0.00
89		Benzo(k)fluoranthene	1.245	1.261	-1.3	117	0.00
90	C	Benzo(a)pyrene	1.149	1.178	-2.5	117	0.00
91		Dibenzo(a,h)anthracene	1.243	1.261	-1.4	117	0.00
92		Benzo(g,h,i)perylene	1.191	1.205	-1.2	116	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080221\
Data File : BP006464.D
Acq On : 03 Aug 2021 00:03
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Aug 03 05:27:59 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 03 05:26:21 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