Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112222\
Data File : BP012700.D
Acq On : 22 Nov 2022 14:01
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP112222

Quant Time: Nov 23 00:48:27 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 23 00:38:14 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	104	0.00
2		1,4-Dioxane	0.389	0.356	8.5	102	0.00
3		Pyridine	1.236	1.163	5.9	103	0.00
4		n-Nitrosodimethylamine	0.541	0.495	8.5	101	0.00
5	S	2-Fluorophenol	0.923	0.920	0.3	107	0.00
6		Aniline	1.674	1.650	1.4	103	0.00
7	S	Phenol-d6	1.324	1.337	-1.0	106	0.00
8		2-Chlorophenol	1.191	1.188	0.3	106	0.00
9		Benzaldehyde	0.925	0.804	13.1	105	0.00
10	C	Phenol	1.509	1.501	0.5	105	0.00
11		bis(2-Chloroethyl)ether	1.317	1.247	5.3	103	0.00
12		1,3-Dichlorobenzene	1.480	1.386	6.4	103	0.00
13	C	1,4-Dichlorobenzene	1.520	1.420	6.6	103	0.00
14		1,2-Dichlorobenzene	1.477	1.386	6.2	103	0.00
15		Benzyl Alcohol	0.941	0.995	-5.7	107	0.00
16		2,2'-oxybis(1-Chloropropane	1.952	1.814	7.1	101	0.00
17		2-Methylphenol	1.030	1.064	-3.3	105	0.00
18		Hexachloroethane	0.555	0.513	7.6	103	0.00
19	P	n-Nitroso-di-n-propylamine	0.935	0.968	-3.5	103	0.00
20		3+4-Methylphenols	1.387	1.464	-5.6	107	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
22		Acetophenone	0.488	0.461	5.5	103	0.00
23	S	Nitrobenzene-d5	0.359	0.342	4.7	102	0.00
24		Nitrobenzene	0.367	0.350	4.6	101	0.00
25		Isophorone	0.658	0.641	2.6	104	0.00
26	C	2-Nitrophenol	0.096	0.132	-37.5#	117	0.00
27		2,4-Dimethylphenol	0.225	0.239	-6.2	108	0.00
28		bis(2-Chloroethoxy)methane	0.389	0.377	3.1	102	0.00
29	C	2,4-Dichlorophenol	0.271	0.305	-12.5	105	0.00
30		1,2,4-Trichlorobenzene	0.359	0.344	4.2	103	0.00
31		Naphthalene	1.085	1.021	5.9	102	0.00
32		Benzoic acid	0.117	0.131	-12.0	124	0.00
33		4-Chloroaniline	0.396	0.400	-1.0	104	0.00
34	C	Hexachlorobutadiene	0.225	0.214	4.9	104	0.00
35		Caprolactam	0.080	0.088	-10.0	117	0.00
36	C	4-Chloro-3-methylphenol	0.301	0.313	-4.0	106	0.00
37		2-Methylnaphthalene	0.773	0.743	3.9	102	0.00
38		1-Methylnaphthalene	0.759	0.727	4.2	102	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
40		1,2,4,5-Tetrachlorobenzene	0.637	0.603	5.3	103	0.00
41	P	Hexachlorocyclopentadiene	0.303	0.284	6.3	99	0.00
42	S	2,4,6-Tribromophenol	0.225	0.230	-2.2	112	0.00
43	C	2,4,6-Trichlorophenol	0.369	0.388	-5.1	106	0.00
44		2,4,5-Trichlorophenol	0.412	0.444	-7.8	106	0.00
45	S	2-Fluorobiphenyl	1.404	1.311	6.6	102	0.00
46		1,1'-Biphenyl	1.513	1.425	5.8	103	0.00
47		2-Chloronaphthalene	1.198	1.122	6.3	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112222\
Data File : BP012700.D
Acq On : 22 Nov 2022 14:01
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP112222

Quant Time: Nov 23 00:48:27 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 23 00:38:14 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	0.263	0.279	-6.1	105	0.00
49		Acenaphthylene	1.903	1.787	6.1	103	0.00
50		Dimethylphthalate	1.472	1.430	2.9	104	0.00
51		2,6-Dinitrotoluene	0.309	0.304	1.6	103	0.00
52	C	Acenaphthene	1.206	1.147	4.9	104	0.00
53		3-Nitroaniline	0.314	0.321	-2.2	104	0.00
54	P	2,4-Dinitrophenol	0.120	0.122	-1.7	112	0.00
55		Dibenzofuran	1.876	1.769	5.7	103	0.00
56	P	4-Nitrophenol	0.265	0.264	0.4	107	0.00
57		2,4-Dinitrotoluene	0.408	0.414	-1.5	104	0.00
58		Fluorene	1.542	1.449	6.0	102	0.00
59		2,3,4,6-Tetrachlorophenol	0.403	0.411	-2.0	106	0.00
60		Diethylphthalate	1.392	1.396	-0.3	105	0.00
61		4-Chlorophenyl-phenylether	0.766	0.726	5.2	103	0.00
62		4-Nitroaniline	0.301	0.320	-6.3	104	0.00
63		Azobenzene	1.467	1.378	6.1	103	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
65		4,6-Dinitro-2-methylphenol	0.081	0.085	-4.9	114	0.00
66	c	n-Nitrosodiphenylamine	0.559	0.543	2.9	104	0.00
67		4-Bromophenyl-phenylether	0.205	0.205	0.0	108	0.00
68		Hexachlorobenzene	0.253	0.238	5.9	104	0.00
69		Atrazine	0.150	0.169	-12.7	109	0.00
70	C	Pentachlorophenol	0.146	0.151	-3.4	108	0.00
71		Phenanthrene	1.131	1.071	5.3	103	0.00
72		Anthracene	1.075	1.022	4.9	103	0.00
73		Carbazole	0.971	0.929	4.3	103	0.00
74		Di-n-butylphthalate	0.850	0.981	-15.4	114	0.00
75	C	Fluoranthene	1.335	1.284	3.8	103	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
77		Benzidine	0.118	0.125	-5.9	117	0.00
78		Pyrene	1.320	1.282	2.9	103	0.00
79	S	Terphenyl-d14	0.941	0.929	1.3	102	0.00
80		Butylbenzylphthalate	0.132	0.159	-20.5	126	0.00
81		Benzo(a)anthracene	1.352	1.295	4.2	103	0.00
82		3,3'-Dichlorobenzidine	0.286	0.297	-3.8	108	0.00
83		Chrysene	1.294	1.242	4.0	103	0.00
84		Bis(2-ethylhexyl)phthalate	0.249	0.324	-30.1#	129	0.00
85	c	Di-n-octyl phthalate	0.581	0.622	-7.1	127	0.00
86	I	Perylene-d12	1.000	1.000	0.0	99	0.00
87		Indeno(1,2,3-cd)pyrene	1.401	1.284	8.4	94	0.00
88		Benzo(b)fluoranthene	1.338	1.353	-1.1	105	0.00
89		Benzo(k)fluoranthene	1.372	1.312	4.4	97	0.00
90	C	Benzo(a)pyrene	1.067	1.045	2.1	99	0.00
91		Dibenzo(a,h)anthracene	1.194	1.096	8.2	95	0.00
92		Benzo(g,h,i)perylene	1.078	0.972	9.8	94	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112222\
Data File : BP012700.D
Acq On : 22 Nov 2022 14:01
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP112222

Quant Time: Nov 23 00:48:27 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 23 00:38:14 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 = 1