Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112122\
Data File : BP012689.D
Acq On : 21 Nov 2022 16:49
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP112122

Quant Time: Nov 21 23:59:53 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 21 23:51:17 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	113	0.00
2		1,4-Dioxane	0.406	0.380	6.4	111	0.00
3		Pyridine	1.267	1.223	3.5	111	0.00
4		n-Nitrosodimethylamine	0.553	0.520	6.0	108	0.00
5	S	2-Fluorophenol	0.987	0.983	0.4	114	0.00
6		Aniline	1.730	1.759	-1.7	114	0.00
7	S	Phenol-d6	1.370	1.384	-1.0	114	0.00
8		2-Chlorophenol	1.248	1.261	-1.0	114	0.00
9		Benzaldehyde	0.844	0.748	11.4	114	0.00
10	C	Phenol	1.554	1.566	-0.8	114	0.00
11		bis(2-Chloroethyl)ether	1.340	1.290	3.7	111	0.00
12		1,3-Dichlorobenzene	1.517	1.463	3.6	112	0.00
13	C	1,4-Dichlorobenzene	1.562	1.498	4.1	112	0.00
14		1,2-Dichlorobenzene	1.514	1.453	4.0	112	0.00
15		Benzyl Alcohol	0.961	1.029	-7.1	117	0.00
16		2,2'-oxybis(1-Chloropropane	1.972	1.871	5.1	109	0.00
17		2-Methylphenol	1.073	1.112	-3.6	115	0.00
18		Hexachloroethane	0.567	0.537	5.3	111	0.00
19	P	n-Nitroso-di-n-propylamine	0.944	0.995	-5.4	113	0.00
20		3+4-Methylphenols	1.451	1.520	-4.8	115	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	111	0.00
22		Acetophenone	0.494	0.479	3.0	112	0.00
23	S	Nitrobenzene-d5	0.357	0.354	0.8	111	0.00
24		Nitrobenzene	0.369	0.361	2.2	110	0.00
25		Isophorone	0.690	0.673	2.5	112	0.00
26	C	2-Nitrophenol	0.120	0.140	-16.7	125	0.00
27		2,4-Dimethylphenol	0.237	0.252	-6.3	116	0.00
28		bis(2-Chloroethoxy)methane	0.391	0.392	-0.3	112	0.00
29	C	2,4-Dichlorophenol	0.301	0.318	-5.6	114	0.00
30		1,2,4-Trichlorobenzene	0.370	0.360	2.7	113	0.00
31		Naphthalene	1.117	1.074	3.8	111	0.00
32		Benzoic acid	0.130	0.137	-5.4	127	0.00
33		4-Chloroaniline	0.393	0.420	-6.9	114	0.00
34	C	Hexachlorobutadiene	0.232	0.225	3.0	113	0.00
35		Caprolactam	0.089	0.096	-7.9	120	0.00
36	C	4-Chloro-3-methylphenol	0.315	0.324	-2.9	112	0.00
37		2-Methylnaphthalene	0.793	0.771	2.8	110	0.00
38		1-Methylnaphthalene	0.787	0.758	3.7	111	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
40		1,2,4,5-Tetrachlorobenzene	0.637	0.632	0.8	112	0.00
41	P	Hexachlorocyclopentadiene	0.298	0.297	0.3	108	0.00
42	S	2,4,6-Tribromophenol	0.220	0.238	-8.2	126	0.00
43	C	2,4,6-Trichlorophenol	0.388	0.417	-7.5	114	0.00
44		2,4,5-Trichlorophenol	0.451	0.475	-5.3	113	0.00
45	S	2-Fluorobiphenyl	1.417	1.379	2.7	109	0.00
46		1,1'-Biphenyl	1.527	1.495	2.1	110	0.00
47		2-Chloronaphthalene	1.227	1.200	2.2	110	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.260	0.292	-12.3	115	0.00
49		Acenaphthylene	1.962	1.907	2.8	109	0.00
50		Dimethylphthalate	1.543	1.531	0.8	111	0.00
51		2,6-Dinitrotoluene	0.312	0.323	-3.5	111	0.00
52	C	Acenaphthene	1.247	1.220	2.2	110	0.00
53		3-Nitroaniline	0.320	0.344	-7.5	112	0.00
54	P	2,4-Dinitrophenol	0.127	0.133	-4.7	121	0.00
55		Dibenzofuran	1.929	1.869	3.1	110	0.00
56	P	4-Nitrophenol	0.276	0.285	-3.3	114	0.00
57		2,4-Dinitrotoluene	0.414	0.437	-5.6	111	0.00
58		Fluorene	1.577	1.518	3.7	109	0.00
59		2,3,4,6-Tetrachlorophenol	0.414	0.436	-5.3	113	0.00
60		Diethylphthalate	1.479	1.509	-2.0	112	0.00
61		4-Chlorophenyl-phenylether	0.778	0.763	1.9	112	0.00
62		4-Nitroaniline	0.298	0.339	-13.8	113	0.00
63		Azobenzene	1.510	1.458	3.4	109	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	109	0.00
65		4,6-Dinitro-2-methylphenol	0.085	0.089	-4.7	119	0.00
66	c	n-Nitrosodiphenylamine	0.578	0.574	0.7	110	0.00
67		4-Bromophenyl-phenylether	0.204	0.211	-3.4	120	0.00
68		Hexachlorobenzene	0.258	0.254	1.6	112	0.00
69		Atrazine	0.140	0.173	-23.6	116	0.00
70	C	Pentachlorophenol	0.154	0.162	-5.2	115	0.00
71		Phenanthrene	1.166	1.127	3.3	108	0.00
72		Anthracene	1.110	1.090	1.8	109	0.00
73		Carbazole	0.996	0.997	-0.1	110	0.00
74		Di-n-butylphthalate	0.947	1.083	-14.4	118	0.00
75	C	Fluoranthene	1.384	1.361	1.7	108	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	107	0.00
77		Benzidine	0.122	0.135	-10.7	147	0.00
78		Pyrene	1.362	1.340	1.6	108	0.00
79	S	Terphenyl-d14	0.950	0.954	-0.4	107	0.00
80		Butylbenzylphthalate	0.152	0.186	-22.4	136	0.00
81		Benzo(a)anthracene	1.396	1.376	1.4	107	0.00
82		3,3'-Dichlorobenzidine	0.278	0.303	-9.0	119	0.00
83		Chrysene	1.339	1.304	2.6	107	0.00
84		Bis(2-ethylhexyl)phthalate	0.292	0.381	-30.5#	134	0.00
85	c	Di-n-octyl phthalate	0.679	0.747	-10.0	134	0.00
86	I	Perylene-d12	1.000	1.000	0.0	106	0.00
87		Indeno(1,2,3-cd)pyrene	1.481	1.441	2.7	100	0.01
88		Benzo(b)fluoranthene	1.363	1.386	-1.7	107	0.00
89		Benzo(k)fluoranthene	1.393	1.361	2.3	106	0.00
90	C	Benzo(a)pyrene	1.099	1.106	-0.6	106	0.00
91		Dibenzo(a,h)anthracene	1.249	1.219	2.4	101	0.01
92		Benzo(g,h,i)perylene	1.157	1.102	4.8	99	0.02

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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