Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121421\
Data File : BP008398.D
Acq On : 14 Dec 2021 17:24
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP121421

Quant Time: Dec 14 18:31:17 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP121421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 14 17:59:47 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	100	0.00
2		1,4-Dioxane	0.541	0.495	8.5	94	0.00
3		Pyridine	1.527	1.481	3.0	95	0.00
4		n-Nitrosodimethylamine	0.602	0.590	2.0	97	0.00
5	S	2-Fluorophenol	1.266	1.232	2.7	97	0.00
6		Aniline	1.969	2.059	-4.6	99	0.00
7	S	Phenol-d6	1.701	1.736	-2.1	97	0.00
8		2-Chlorophenol	1.291	1.311	-1.5	98	0.00
9		Benzaldehyde	1.152	0.968	16.0	98	0.00
10	C	Phenol	1.727	1.754	-1.6	97	0.00
11		bis(2-Chloroethyl)ether	1.441	1.401	2.8	98	0.00
12		1,3-Dichlorobenzene	1.493	1.477	1.1	97	0.00
13	C	1,4-Dichlorobenzene	1.519	1.508	0.7	97	0.00
14		1,2-Dichlorobenzene	1.473	1.477	-0.3	98	0.00
15		Benzyl Alcohol	1.368	1.467	-7.2	98	0.00
16		2,2'-oxybis(1-Chloropropane	2.039	2.010	1.4	99	0.00
17		2-Methylphenol	1.143	1.190	-4.1	98	0.00
18		Hexachloroethane	0.564	0.567	-0.5	99	0.00
19	P	n-Nitroso-di-n-propylamine	1.262	1.264	-0.2	99	0.00
20		3+4-Methylphenols	1.574	1.674	-6.4	100	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
22		Acetophenone	0.588	0.606	-3.1	100	0.00
23	S	Nitrobenzene-d5	0.472	0.483	-2.3	99	0.00
24		Nitrobenzene	0.443	0.456	-2.9	99	0.00
25		Isophorone	0.790	0.841	-6.5	102	0.00
26	C	2-Nitrophenol	0.196	0.208	-6.1	101	0.00
27		2,4-Dimethylphenol	0.281	0.294	-4.6	100	0.00
28		bis(2-Chloroethoxy)methane	0.488	0.517	-5.9	100	0.00
29	C	2,4-Dichlorophenol	0.345	0.361	-4.6	99	0.00
30		1,2,4-Trichlorobenzene	0.409	0.415	-1.5	100	0.00
31		Naphthalene	1.049	1.078	-2.8	101	0.00
32		Benzoic acid	0.206	0.247	-19.9	104	0.00
33		4-Chloroaniline	0.420	0.460	-9.5	104	0.00
34	C	Hexachlorobutadiene	0.299	0.306	-2.3	100	0.00
35		Caprolactam	0.065	0.076	-16.9	110	0.00
36	C	4-Chloro-3-methylphenol	0.386	0.424	-9.8	104	0.00
37		2-Methylnaphthalene	0.723	0.778	-7.6	103	0.00
38		1-Methylnaphthalene	0.704	0.758	-7.7	103	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
40		1,2,4,5-Tetrachlorobenzene	0.764	0.762	0.3	104	0.00
41	P	Hexachlorocyclopentadiene	0.162	0.179	-10.5	104	0.00
42	S	2,4,6-Tribromophenol	0.294	0.328	-11.6	115	0.00
43	C	2,4,6-Trichlorophenol	0.480	0.498	-3.8	104	0.00
44		2,4,5-Trichlorophenol	0.501	0.531	-6.0	108	0.00
45	S	2-Fluorobiphenyl	1.573	1.594	-1.3	105	0.00
46		1,1'-Biphenyl	1.545	1.576	-2.0	105	0.00
47		2-Chloronaphthalene	1.232	1.249	-1.4	105	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.415	0.444	-7.0	109	0.00
49		Acenaphthylene	1.812	1.860	-2.6	106	0.00
50		Dimethylphthalate	1.514	1.640	-8.3	112	0.00
51		2,6-Dinitrotoluene	0.316	0.347	-9.8	114	0.00
52	C	Acenaphthene	1.078	1.104	-2.4	105	0.00
53		3-Nitroaniline	0.298	0.329	-10.4	113	0.00
54	P	2,4-Dinitrophenol	0.165	0.195	-18.2	117	0.00
55		Dibenzofuran	1.833	1.892	-3.2	108	0.00
56	P	4-Nitrophenol	0.173	0.201	-16.2	114	0.00
57		2,4-Dinitrotoluene	0.422	0.468	-10.9	115	0.00
58		Fluorene	1.421	1.494	-5.1	110	0.00
59		2,3,4,6-Tetrachlorophenol	0.428	0.463	-8.2	112	0.00
60		Diethylphthalate	1.437	1.562	-8.7	115	0.00
61		4-Chlorophenyl-phenylether	0.812	0.858	-5.7	110	0.00
62		4-Nitroaniline	0.268	0.298	-11.2	115	0.00
63		Azobenzene	1.457	1.540	-5.7	109	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	115	0.00
65		4,6-Dinitro-2-methylphenol	0.131	0.151	-15.3	120	0.00
66	c	n-Nitrosodiphenylamine	0.630	0.644	-2.2	112	0.00
67		4-Bromophenyl-phenylether	0.260	0.272	-4.6	115	0.00
68		Hexachlorobenzene	0.287	0.297	-3.5	115	0.00
69		Atrazine	0.139	0.157	-12.9	121	0.00
70	C	Pentachlorophenol	0.128	0.146	-14.1	119	0.00
71		Phenanthrene	1.083	1.116	-3.0	117	0.00
72		Anthracene	1.069	1.113	-4.1	117	0.00
73		Carbazole	0.978	1.019	-4.2	119	0.00
74		Di-n-butylphthalate	1.122	1.232	-9.8	126	0.00
75	C	Fluoranthene	1.223	1.266	-3.5	122	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	115	0.00
77		Benzidine	0.203	0.240	-18.2	122	0.00
78		Pyrene	1.431	1.522	-6.4	122	0.00
79	S	Terphenyl-d14	1.149	1.321	-15.0	126	0.00
80		Butylbenzylphthalate	0.521	0.576	-10.6	125	0.00
81		Benzo(a)anthracene	1.340	1.386	-3.4	116	0.00
82		3,3'-Dichlorobenzidine	0.617	0.612	0.8	113	0.00
83		Chrysene	1.280	1.314	-2.7	117	0.00
84		Bis(2-ethylhexyl)phthalate	0.734	0.769	-4.8	118	0.00
85	c	Di-n-octyl phthalate	1.265	1.244	1.7	111	0.00
86	I	Perylene-d12	1.000	1.000	0.0	104	0.00
87		Indeno(1,2,3-cd)pyrene	1.574	1.585	-0.7	101	0.00
88		Benzo(b)fluoranthene	1.203	1.258	-4.6	107	0.00
89		Benzo(k)fluoranthene	1.213	1.245	-2.6	105	0.00
90	C	Benzo(a)pyrene	1.044	1.064	-1.9	104	0.00
91		Dibenzo(a,h)anthracene	1.318	1.319	-0.1	101	0.00
92		Benzo(g,h,i)perylene	1.305	1.308	-0.2	101	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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