Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121120\
Data File : BP004228.D
Acq On : 11 Dec 2020 14:48
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : ICVBP121120

Quant Time: Dec 15 10:35:15 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP121120.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 15 10:27:21 2020
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	111	0.00
2		1,4-Dioxane	0.579	0.576	0.5	109	0.00
3		Pyridine	1.569	1.648	-5.0	112	0.00
4		n-Nitrosodimethylamine	0.507	0.536	-5.7	111	0.00
5	S	2-Fluorophenol	1.221	1.311	-7.4	114	0.00
6		Aniline	1.981	2.153	-8.7	114	0.01
7	S	Phenol-d6	1.718	1.871	-8.9	114	0.01
8		2-Chlorophenol	1.303	1.397	-7.2	114	0.01
9		Benzaldehyde	1.071	0.998	6.8	112	0.01
10	C	Phenol	1.698	1.827	-7.6	113	0.01
11		bis(2-Chloroethyl)ether	1.412	1.468	-4.0	111	0.01
12		1,3-Dichlorobenzene	1.679	1.721	-2.5	112	0.01
13	C	1,4-Dichlorobenzene	1.690	1.744	-3.2	113	0.00
14		1,2-Dichlorobenzene	1.601	1.637	-2.2	111	0.01
15		Benzyl Alcohol	1.068	1.223	-14.5	117	0.01
16		2,2'-oxybis(1-Chloropropane	1.801	1.841	-2.2	110	0.01
17		2-Methylphenol	1.083	1.189	-9.8	114	0.01
18		Hexachloroethane	0.612	0.630	-2.9	112	0.01
19	P	n-Nitroso-di-n-propylamine	1.016	1.097	-8.0	113	0.01
20		3+4-Methylphenols	1.439	1.611	-12.0	114	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	113	0.00
22		Acetophenone	0.540	0.558	-3.3	112	0.01
23	S	Nitrobenzene-d5	0.421	0.439	-4.3	113	0.01
24		Nitrobenzene	0.414	0.433	-4.6	113	0.01
25		Isophorone	0.692	0.751	-8.5	114	0.01
26	C	2-Nitrophenol	0.150	0.179	-19.3	121	0.01
27		2,4-Dimethylphenol	0.261	0.285	-9.2	114	0.01
28		bis(2-Chloroethoxy)methane	0.492	0.515	-4.7	113	0.01
29	C	2,4-Dichlorophenol	0.307	0.347	-13.0	116	0.01
30		1,2,4-Trichlorobenzene	0.410	0.424	-3.4	113	0.01
31		Naphthalene	1.137	1.164	-2.4	112	0.01
32		Benzoic acid	0.152	0.178	-17.1	123	0.02
33		4-Chloroaniline	0.459	0.494	-7.6	113	0.01
34	C	Hexachlorobutadiene	0.257	0.265	-3.1	113	0.01
35		Caprolactam	0.101	0.113	-11.9	119	0.02
36	C	4-Chloro-3-methylphenol	0.308	0.348	-13.0	116	0.01
37		2-Methylnaphthalene	0.794	0.821	-3.4	112	0.00
38		1-Methylnaphthalene	0.754	0.778	-3.2	112	0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	113	0.01
40		1,2,4,5-Tetrachlorobenzene	0.723	0.750	-3.7	113	0.01
41	P	Hexachlorocyclopentadiene	0.363	0.394	-8.5	115	0.01
42	S	2,4,6-Tribromophenol	0.267	0.290	-8.6	117	0.01
43	C	2,4,6-Trichlorophenol	0.402	0.469	-16.7	118	0.01
44		2,4,5-Trichlorophenol	0.435	0.492	-13.1	116	0.01
45	S	2-Fluorobiphenyl	1.578	1.607	-1.8	112	0.00
46		1,1'-Biphenyl	1.690	1.729	-2.3	113	0.01
47		2-Chloronaphthalene	1.387	1.426	-2.8	113	0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.356	0.408	-14.6	116	0.01
49		Acenaphthylene	2.040	2.138	-4.8	113	0.01
50		Dimethylphthalate	1.646	1.734	-5.3	114	0.01
51		2,6-Dinitrotoluene	0.346	0.377	-9.0	114	0.01
52	C	Acenaphthene	1.273	1.318	-3.5	113	0.01
53		3-Nitroaniline	0.379	0.426	-12.4	115	0.01
54	P	2,4-Dinitrophenol	0.140	0.160	-14.3	124	0.01
55		Dibenzofuran	2.014	2.054	-2.0	112	0.01
56	P	4-Nitrophenol	0.284	0.315	-10.9	118	0.01
57		2,4-Dinitrotoluene	0.455	0.517	-13.6	115	0.01
58		Fluorene	1.581	1.627	-2.9	112	0.01
59		2,3,4,6-Tetrachlorophenol	0.400	0.455	-13.7	117	0.01
60		Diethylphthalate	1.579	1.688	-6.9	114	0.01
61		4-Chlorophenyl-phenylether	0.860	0.888	-3.3	113	0.00
62		4-Nitroaniline	0.395	0.445	-12.7	114	0.00
63		Azobenzene	1.503	1.581	-5.2	112	0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	113	0.01
65		4,6-Dinitro-2-methylphenol	0.091	0.105	-15.4	124	0.01
66	c	n-Nitrosodiphenylamine	0.630	0.668	-6.0	114	0.01
67		4-Bromophenyl-phenylether	0.249	0.263	-5.6	114	0.01
68		Hexachlorobenzene	0.297	0.308	-3.7	114	0.01
69		Atrazine	0.204	0.225	-10.3	113	0.01
70	C	Pentachlorophenol	0.141	0.160	-13.5	121	0.00
71		Phenanthrene	1.203	1.233	-2.5	113	0.01
72		Anthracene	1.136	1.189	-4.7	112	0.01
73		Carbazole	1.052	1.116	-6.1	112	0.01
74		Di-n-butylphthalate	1.141	1.302	-14.1	115	0.01
75	C	Fluoranthene	1.381	1.449	-4.9	112	0.01
76	I	Chrysene-d12	1.000	1.000	0.0	111	0.01
77		Benzidine	0.346	0.490	-41.6#	115	0.00
78		Pyrene	1.363	1.451	-6.5	111	0.01
79	S	Terphenyl-d14	1.044	1.074	-2.9	110	0.01
80		Butylbenzylphthalate	0.437	0.549	-25.6#	120	0.00
81		Benzo(a)anthracene	1.373	1.436	-4.6	111	0.02
82		3,3'-Dichlorobenzidine	0.389	0.475	-22.1	114	0.01
83		Chrysene	1.323	1.372	-3.7	111	0.01
84		Bis(2-ethylhexyl)phthalate	0.672	0.820	-22.0	115	0.01
85	c	Di-n-octyl phthalate	1.044	1.328	-27.2#	118	0.01
86	I	Perylene-d12	1.000	1.000	0.0	113	0.02
87		Indeno(1,2,3-cd)pyrene	1.529	1.638	-7.1	114	0.02
88		Benzo(b)fluoranthene	1.318	1.361	-3.3	109	0.01
89		Benzo(k)fluoranthene	1.309	1.374	-5.0	116	0.01
90	C	Benzo(a)pyrene	1.201	1.290	-7.4	113	0.02
91		Dibenzo(a,h)anthracene	1.257	1.335	-6.2	113	0.03
92		Benzo(g,h,i)perylene	1.227	1.300	-5.9	113	0.03

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Compound	AvgRF	CCRF	%Dev Area%	Dev(min)

(#) = Out of Range SPCC's out = 0 CCC's out = 1