Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101420\
Data File : BP003613.D
Acq On : 14 Oct 2020 17:57
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : ICVBP101420

Quant Time: Oct 14 18:49:22 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 14 17:03:23 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	112	0.00
2		1,4-Dioxane	0.498	0.463	7.0	108	0.00
3		Pyridine	1.452	1.407	3.1	111	0.00
4		n-Nitrosodimethylamine	0.624	0.597	4.3	110	0.00
5	S	2-Fluorophenol	1.163	1.144	1.6	111	0.00
6		Aniline	1.983	1.919	3.2	111	0.00
7	S	Phenol-d6	1.695	1.684	0.6	113	0.00
8		2-Chlorophenol	1.183	1.164	1.6	112	0.00
9		Benzaldehyde	1.058	0.882	16.6	110	0.00
10	C	Phenol	1.637	1.594	2.6	112	0.00
11		bis(2-Chloroethyl)ether	1.314	1.254	4.6	109	0.00
12		1,3-Dichlorobenzene	1.538	1.463	4.9	110	0.00
13	C	1,4-Dichlorobenzene	1.543	1.480	4.1	109	0.00
14		1,2-Dichlorobenzene	1.480	1.399	5.5	110	0.00
15		Benzyl Alcohol	1.361	1.352	0.7	113	0.00
16		2,2'-oxybis(1-Chloropropane	1.345	1.291	4.0	111	0.00
17		2-Methylphenol	1.109	1.084	2.3	111	0.00
18		Hexachloroethane	0.600	0.578	3.7	111	0.00
19	P	n-Nitroso-di-n-propylamine	1.130	1.095	3.1	113	0.00
20		3+4-Methylphenols	1.476	1.471	0.3	113	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	111	0.00
22		Acetophenone	0.590	0.567	3.9	111	0.00
23	S	Nitrobenzene-d5	0.431	0.449	-4.2	116	0.00
24		Nitrobenzene	0.425	0.434	-2.1	116	0.00
25		Isophorone	0.805	0.781	3.0	111	0.00
26	C	2-Nitrophenol	0.128	0.151	-18.0	124	0.00
27		2,4-Dimethylphenol	0.265	0.262	1.1	114	0.00
28		bis(2-Chloroethoxy)methane	0.495	0.480	3.0	113	0.00
29	C	2,4-Dichlorophenol	0.331	0.339	-2.4	116	0.00
30		1,2,4-Trichlorobenzene	0.439	0.419	4.6	111	0.00
31		Naphthalene	1.072	1.031	3.8	110	0.00
32		Benzoic acid	0.123	0.151	-22.8	128	0.01
33		4-Chloroaniline	0.463	0.447	3.5	111	0.00
34	C	Hexachlorobutadiene	0.326	0.310	4.9	109	0.00
35		Caprolactam	0.105	0.106	-1.0	113	0.01
36	C	4-Chloro-3-methylphenol	0.378	0.379	-0.3	114	0.00
37		2-Methylnaphthalene	0.798	0.763	4.4	111	0.00
38		1-Methylnaphthalene	0.750	0.722	3.7	111	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	111	0.00
40		1,2,4,5-Tetrachlorobenzene	0.758	0.731	3.6	111	0.00
41	P	Hexachlorocyclopentadiene	0.423	0.423	0.0	112	0.00
42	S	2,4,6-Tribromophenol	0.267	0.278	-4.1	119	0.01
43	C	2,4,6-Trichlorophenol	0.420	0.435	-3.6	118	0.00
44		2,4,5-Trichlorophenol	0.442	0.451	-2.0	115	0.00
45	S	2-Fluorobiphenyl	1.524	1.457	4.4	110	0.00
46		1,1'-Biphenyl	1.545	1.481	4.1	111	0.00
47		2-Chloronaphthalene	1.287	1.234	4.1	111	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.320	0.364	-13.7	122	0.00
49		Acenaphthylene	1.874	1.811	3.4	111	0.00
50		Dimethylphthalate	1.642	1.569	4.4	111	0.00
51		2,6-Dinitrotoluene	0.296	0.315	-6.4	117	0.00
52	C	Acenaphthene	1.221	1.184	3.0	112	0.00
53		3-Nitroaniline	0.298	0.319	-7.0	117	0.01
54	P	2,4-Dinitrophenol	0.116	0.132	-13.8	135	0.00
55		Dibenzofuran	1.898	1.812	4.5	111	0.00
56	P	4-Nitrophenol	0.214	0.232	-8.4	121	0.00
57		2,4-Dinitrotoluene	0.383	0.424	-10.7	117	0.00
58		Fluorene	1.535	1.472	4.1	111	0.00
59		2,3,4,6-Tetrachlorophenol	0.406	0.432	-6.4	120	0.00
60		Diethylphthalate	1.614	1.547	4.2	111	0.00
61		4-Chlorophenyl-phenylether	0.910	0.877	3.6	112	0.00
62		4-Nitroaniline	0.332	0.356	-7.2	117	0.01
63		Azobenzene	1.553	1.478	4.8	110	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	110	0.01
65		4,6-Dinitro-2-methylphenol	0.077	0.086	-11.7	126	0.00
66	c	n-Nitrosodiphenylamine	0.597	0.574	3.9	110	0.00
67		4-Bromophenyl-phenylether	0.263	0.255	3.0	112	0.00
68		Hexachlorobenzene	0.300	0.288	4.0	110	0.00
69		Atrazine	0.233	0.227	2.6	110	0.00
70	C	Pentachlorophenol	0.136	0.145	-6.6	120	0.00
71		Phenanthrene	1.111	1.065	4.1	109	0.00
72		Anthracene	1.084	1.037	4.3	109	0.00
73		Carbazole	0.979	0.935	4.5	109	0.00
74		Di-n-butylphthalate	1.192	1.162	2.5	109	0.00
75	C	Fluoranthene	1.375	1.313	4.5	109	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	109	0.00
77		Benzidine	0.469	0.536	-14.3	116	0.00
78		Pyrene	1.329	1.297	2.4	110	0.00
79	S	Terphenyl-d14	1.105	1.084	1.9	110	0.00
80		Butylbenzylphthalate	0.464	0.485	-4.5	113	0.00
81		Benzo(a)anthracene	1.345	1.297	3.6	109	0.00
82		3,3'-Dichlorobenzidine	0.475	0.474	0.2	110	0.00
83		Chrysene	1.272	1.218	4.2	108	0.00
84		Bis(2-ethylhexyl)phthalate	0.705	0.697	1.1	109	0.00
85	c	Di-n-octyl phthalate	1.138	1.154	-1.4	111	0.00
86	I	Perylene-d12	1.000	1.000	0.0	108	0.01
87		Indeno(1,2,3-cd)pyrene	1.518	1.492	1.7	109	0.01
88		Benzo(b)fluoranthene	1.390	1.327	4.5	106	0.00
89		Benzo(k)fluoranthene	1.305	1.302	0.2	111	0.01
90	C	Benzo(a)pyrene	1.252	1.229	1.8	109	0.01
91		Dibenzo(a,h)anthracene	1.266	1.245	1.7	109	0.01
92		Benzo(g,h,i)perylene	1.195	1.174	1.8	109	0.02

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