Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100121\
Data File : BM032304.D
Acq On : 30 Sep 2021 20:41
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM100121

Quant Time: Oct 01 10:41:47 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 01 10:39:43 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	80	0.00
2		1,4-Dioxane	0.495	0.456	7.9	78	0.00
3		Pyridine	1.425	1.338	6.1	77	0.00
4		n-Nitrosodimethylamine	0.581	0.531	8.6	75	0.00
5	S	2-Fluorophenol	1.184	1.136	4.1	79	0.00
6		Aniline	1.931	1.871	3.1	78	0.00
7	S	Phenol-d6	1.719	1.658	3.5	78	0.00
8		2-Chlorophenol	1.320	1.290	2.3	80	0.00
9		Benzaldehyde	0.995	0.913	8.2	80	0.00
10	C	Phenol	1.656	1.598	3.5	78	0.00
11		bis(2-Chloroethyl)ether	1.320	1.251	5.2	78	0.00
12		1,3-Dichlorobenzene	1.464	1.402	4.2	80	0.00
13	C	1,4-Dichlorobenzene	1.491	1.430	4.1	80	0.00
14		1,2-Dichlorobenzene	1.450	1.391	4.1	80	0.00
15		Benzyl Alcohol	1.184	1.208	-2.0	81	0.00
16		2,2'-oxybis(1-Chloropropane	1.756	1.594	9.2	74	0.00
17		2-Methylphenol	1.116	1.117	-0.1	80	0.00
18		Hexachloroethane	0.496	0.496	0.0	80	0.00
19	P	n-Nitroso-di-n-propylamine	0.971	0.970	0.1	78	0.00
20		3+4-Methylphenols	1.536	1.548	-0.8	80	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	81	0.00
22		Acetophenone	0.490	0.470	4.1	79	0.00
23	S	Nitrobenzene-d5	0.356	0.350	1.7	79	0.00
24		Nitrobenzene	0.346	0.336	2.9	79	0.00
25		Isophorone	0.609	0.622	-2.1	80	0.00
26	C	2-Nitrophenol	0.155	0.168	-8.4	84	0.00
27		2,4-Dimethylphenol	0.273	0.272	0.4	80	0.00
28		bis(2-Chloroethoxy)methane	0.442	0.421	4.8	78	0.00
29	C	2,4-Dichlorophenol	0.298	0.303	-1.7	82	0.00
30		1,2,4-Trichlorobenzene	0.331	0.324	2.1	83	0.00
31		Naphthalene	1.025	0.984	4.0	80	0.00
32		Benzoic acid	0.195	0.221	-13.3	86	0.00
33		4-Chloroaniline	0.436	0.432	0.9	80	0.00
34	C	Hexachlorobutadiene	0.193	0.191	1.0	84	0.00
35		Caprolactam	0.109	0.112	-2.8	83	0.00
36	C	4-Chloro-3-methylphenol	0.333	0.343	-3.0	83	0.00
37		2-Methylnaphthalene	0.707	0.695	1.7	82	0.00
38		1-Methylnaphthalene	0.692	0.679	1.9	81	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	84	0.00
40		1,2,4,5-Tetrachlorobenzene	0.559	0.529	5.4	83	0.00
41	P	Hexachlorocyclopentadiene	0.245	0.254	-3.7	84	0.00
42	S	2,4,6-Tribromophenol	0.217	0.232	-6.9	89	0.00
43	C	2,4,6-Trichlorophenol	0.388	0.392	-1.0	85	0.00
44		2,4,5-Trichlorophenol	0.420	0.422	-0.5	85	0.00
45	S	2-Fluorobiphenyl	1.237	1.195	3.4	83	0.00
46		1,1'-Biphenyl	1.468	1.394	5.0	83	0.00
47		2-Chloronaphthalene	1.119	1.067	4.6	83	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.309	0.328	-6.1	83	0.00
49		Acenaphthylene	1.744	1.730	0.8	83	0.00
50		Dimethylphthalate	1.399	1.393	0.4	85	0.00
51		2,6-Dinitrotoluene	0.279	0.297	-6.5	86	0.00
52	C	Acenaphthene	1.164	1.138	2.2	84	0.00
53		3-Nitroaniline	0.326	0.349	-7.1	85	0.00
54	P	2,4-Dinitrophenol	0.163	0.179	-9.8	87	0.00
55		Dibenzofuran	1.769	1.705	3.6	84	0.00
56	P	4-Nitrophenol	0.277	0.301	-8.7	87	0.00
57		2,4-Dinitrotoluene	0.371	0.415	-11.9	88	0.00
58		Fluorene	1.323	1.245	5.9	83	0.00
59		2,3,4,6-Tetrachlorophenol	0.366	0.375	-2.5	86	0.00
60		Diethylphthalate	1.383	1.404	-1.5	86	0.00
61		4-Chlorophenyl-phenylether	0.669	0.643	3.9	85	0.00
62		4-Nitroaniline	0.333	0.362	-8.7	85	0.00
63		Azobenzene	1.336	1.328	0.6	81	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	87	0.00
65		4,6-Dinitro-2-methylphenol	0.115	0.122	-6.1	88	0.00
66	c	n-Nitrosodiphenylamine	0.585	0.553	5.5	84	0.00
67		4-Bromophenyl-phenylether	0.200	0.195	2.5	87	0.00
68		Hexachlorobenzene	0.221	0.216	2.3	89	0.00
69		Atrazine	0.196	0.199	-1.5	89	0.00
70	C	Pentachlorophenol	0.137	0.148	-8.0	90	0.00
71		Phenanthrene	1.056	0.995	5.8	85	0.00
72		Anthracene	1.024	0.994	2.9	86	0.00
73		Carbazole	1.045	1.009	3.4	85	0.00
74		Di-n-butylphthalate	1.062	1.120	-5.5	89	0.00
75	C	Fluoranthene	1.196	1.164	2.7	87	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	87	0.00
77		Benzidine	0.494	0.493	0.2	86	0.00
78		Pyrene	1.246	1.216	2.4	87	0.00
79	S	Terphenyl-d14	0.864	0.807	6.6	90	0.00
80		Butylbenzylphthalate	0.476	0.533	-12.0	92	0.00
81		Benzo(a)anthracene	1.230	1.203	2.2	88	0.00
82		3,3'-Dichlorobenzidine	0.395	0.408	-3.3	88	0.00
83		Chrysene	1.199	1.153	3.8	88	0.00
84		Bis(2-ethylhexyl)phthalate	0.643	0.711	-10.6	90	0.00
85	c	Di-n-octyl phthalate	1.155	1.318	-14.1	90	0.00
86	I	Perylene-d12	1.000	1.000	0.0	89	0.00
87		Indeno(1,2,3-cd)pyrene	1.334	1.277	4.3	88	0.00
88		Benzo(b)fluoranthene	1.125	1.127	-0.2	93	0.00
89		Benzo(k)fluoranthene	1.143	1.028	10.1	84	0.00
90	C	Benzo(a)pyrene	0.960	0.949	1.1	89	0.00
91		Dibenzo(a,h)anthracene	1.106	1.057	4.4	88	0.00
92		Benzo(g,h,i)perylene	1.146	1.107	3.4	87	0.00

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