Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062221\
Data File : BP006008.D
Acq On : 22 Jun 2021 16:51
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Jun 22 17:52:56 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 22 17:52:38 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	74	0.00
2		1,4-Dioxane	0.563	0.587	-4.3	82	0.00
3		Pyridine	1.323	1.503	-13.6	84	0.00
4		n-Nitrosodimethylamine	0.615	0.593	3.6	75	0.00
5	S	2-Fluorophenol	1.246	1.401	-12.4	87	0.00
6		Aniline	1.804	2.051	-13.7	89	0.00
7	S	Phenol-d6	1.616	1.839	-13.8	88	0.00
8		2-Chlorophenol	1.429	1.598	-11.8	88	0.00
9		Benzaldehyde	0.689	0.868	-26.0#	89	0.00
10	C	Phenol	1.614	1.827	-13.2	89	0.00
11		bis(2-Chloroethyl)ether	1.223	1.415	-15.7	91	0.00
12		1,3-Dichlorobenzene	1.598	1.762	-10.3	87	0.00
13	C	1,4-Dichlorobenzene	1.630	1.792	-9.9	87	0.00
14		1,2-Dichlorobenzene	1.558	1.730	-11.0	88	0.00
15		Benzyl Alcohol	1.217	1.321	-8.5	83	0.00
16		2,2'-oxybis(1-Chloropropane	1.130	1.251	-10.7	86	0.00
17		2-Methylphenol	1.100	1.283	-16.6	91	0.00
18		Hexachloroethane	0.589	0.637	-8.1	84	0.00
19	P	n-Nitroso-di-n-propylamine	0.994	1.144	-15.1	90	0.00
20		3+4-Methylphenols	1.485	1.708	-15.0	89	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	74	0.00
22		Acetophenone	0.526	0.618	-17.5	91	0.00
23	S	Nitrobenzene-d5	0.408	0.455	-11.5	87	0.00
24		Nitrobenzene	0.424	0.458	-8.0	86	0.00
25		Isophorone	0.753	0.847	-12.5	89	0.00
26	C	2-Nitrophenol	0.202	0.230	-13.9	90	0.00
27		2,4-Dimethylphenol	0.301	0.335	-11.3	89	0.00
28		bis(2-Chloroethoxy)methane	0.393	0.454	-15.5	93	0.00
29	C	2,4-Dichlorophenol	0.362	0.399	-10.2	88	0.00
30		1,2,4-Trichlorobenzene	0.407	0.441	-8.4	88	0.00
31		Naphthalene	1.157	1.291	-11.6	90	0.00
32		Benzoic acid	0.193	0.229	-18.7	80	0.00
33		4-Chloroaniline	0.467	0.522	-11.8	89	0.00
34	C	Hexachlorobutadiene	0.275	0.300	-9.1	89	0.00
35		Caprolactam	0.104	0.126	-21.2	90	0.00
36	C	4-Chloro-3-methylphenol	0.367	0.423	-15.3	92	0.00
37		2-Methylnaphthalene	0.824	0.931	-13.0	92	0.00
38		1-Methylnaphthalene	0.776	0.874	-12.6	90	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	75	0.00
40		1,2,4,5-Tetrachlorobenzene	0.704	0.784	-11.4	89	0.00
41	P	Hexachlorocyclopentadiene	0.297	0.280	5.7	71	0.00
42	S	2,4,6-Tribromophenol	0.343	0.399	-16.3	92	0.00
43	C	2,4,6-Trichlorophenol	0.483	0.511	-5.8	85	0.00
44		2,4,5-Trichlorophenol	0.520	0.553	-6.3	85	0.00
45	S	2-Fluorobiphenyl	1.523	1.699	-11.6	90	0.00
46		1,1'-Biphenyl	1.613	1.819	-12.8	89	0.00
47		2-Chloronaphthalene	1.317	1.430	-8.6	90	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.346	0.377	-9.0	87	0.00
49		Acenaphthylene	2.020	2.228	-10.3	89	0.00
50		Dimethylphthalate	1.644	1.818	-10.6	91	0.00
51		2,6-Dinitrotoluene	0.374	0.412	-10.2	89	0.00
52	C	Acenaphthene	1.227	1.350	-10.0	91	0.00
53		3-Nitroaniline	0.362	0.406	-12.2	90	0.00
54	P	2,4-Dinitrophenol	0.191	0.229	-19.9	88	0.00
55		Dibenzofuran	2.018	2.178	-7.9	89	0.00
56	P	4-Nitrophenol	0.294	0.287	2.4	76	0.00
57		2,4-Dinitrotoluene	0.500	0.567	-13.4	90	0.00
58		Fluorene	1.572	1.733	-10.2	91	0.00
59		2,3,4,6-Tetrachlorophenol	0.458	0.484	-5.7	85	0.00
60		Diethylphthalate	1.650	1.829	-10.8	91	0.00
61		4-Chlorophenyl-phenylether	0.815	0.891	-9.3	90	0.00
62		4-Nitroaniline	0.373	0.417	-11.8	90	0.00
63		Azobenzene	1.493	1.616	-8.2	88	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	71	0.00
65		4,6-Dinitro-2-methylphenol	0.147	0.149	-1.4	77	0.00
66	c	n-Nitrosodiphenylamine	0.663	0.759	-14.5	89	0.00
67		4-Bromophenyl-phenylether	0.257	0.299	-16.3	90	0.00
68		Hexachlorobenzene	0.308	0.353	-14.6	90	0.00
69		Atrazine	0.206	0.208	-1.0	67	0.00
70	C	Pentachlorophenol	0.171	0.188	-9.9	78	0.00
71		Phenanthrene	1.201	1.384	-15.2	89	0.00
72		Anthracene	1.182	1.369	-15.8	89	0.00
73		Carbazole	1.135	1.311	-15.5	89	0.00
74		Di-n-butylphthalate	1.315	1.561	-18.7	91	0.00
75	C	Fluoranthene	1.459	1.686	-15.6	90	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	77	0.00
77		Benzidine	0.416	0.497	-19.5	81	0.00
78		Pyrene	1.396	1.507	-8.0	91	0.00
79	S	Terphenyl-d14	1.068	1.237	-15.8	96	0.00
80		Butylbenzylphthalate	0.574	0.654	-13.9	94	0.00
81		Benzo(a)anthracene	1.444	1.619	-12.1	95	0.00
82		3,3'-Dichlorobenzidine	0.499	0.540	-8.2	93	0.00
83		Chrysene	1.398	1.581	-13.1	96	0.00
84		Bis(2-ethylhexyl)phthalate	0.842	0.973	-15.6	96	0.00
85	c	Di-n-octyl phthalate	1.507	1.715	-13.8	95	0.00
86	I	Perylene-d12	1.000	1.000	0.0	76	0.00
87		Indeno(1,2,3-cd)pyrene	1.710	1.909	-11.6	93	0.00
88		Benzo(b)fluoranthene	1.370	1.570	-14.6	95	0.00
89		Benzo(k)fluoranthene	1.328	1.532	-15.4	96	0.00
90	C	Benzo(a)pyrene	1.288	1.470	-14.1	95	0.00
91		Dibenzo(a,h)anthracene	1.472	1.647	-11.9	93	0.00
92		Benzo(g,h,i)perylene	1.454	1.588	-9.2	91	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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