Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122921\
Data File : BP008599.D
Acq On : 29 Dec 2021 18:53
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Dec 30 00:44:18 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP122821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 29 13:29:13 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	133	0.00
2		1,4-Dioxane	0.443	0.304	31.4#	93	0.00
3		Pyridine	1.207	0.806	33.2#	86	0.00
4		n-Nitrosodimethylamine	0.390	0.250	35.9#	85	0.00
5	S	2-Fluorophenol	1.096	0.780	28.8#	93	0.00
6		Aniline	1.773	1.055	40.5#	77	0.00
7	S	Phenol-d6	1.525	0.939	38.4#	79	0.00
8		2-Chlorophenol	1.291	0.830	35.7#	84	0.00
9		Benzaldehyde	0.868	0.579	33.3#	85	0.00
10	C	Phenol	1.540	0.952	38.2#	80	0.00
11		bis(2-Chloroethyl)ether	1.242	0.733	41.0#	77	0.00
12		1,3-Dichlorobenzene	1.511	0.946	37.4#	83	0.00
13	C	1,4-Dichlorobenzene	1.537	0.962	37.4#	83	0.00
14		1,2-Dichlorobenzene	1.501	0.919	38.8#	81	0.00
15		Benzyl Alcohol	1.092	0.626	42.7#	73	0.00
16		2,2'-oxybis(1-Chloropropane	1.413	0.788	44.2#	73	0.00
17		2-Methylphenol	1.071	0.626	41.5#	75	0.00
18		Hexachloroethane	0.501	0.309	38.3#	80	0.00
19	P	n-Nitroso-di-n-propylamine	0.974	0.513	47.3#	68	0.00
20		3+4-Methylphenols	1.505	0.872	42.1#	74	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	117	0.00
22		Acetophenone	0.480	0.297	38.1#	71	0.00
23	S	Nitrobenzene-d5	0.336	0.214	36.3#	72	0.00
24		Nitrobenzene	0.320	0.203	36.6#	73	0.00
25		Isophorone	0.628	0.359	42.8#	64	0.00
26	C	2-Nitrophenol	0.152	0.105	30.9#	75	0.00
27		2,4-Dimethylphenol	0.274	0.170	38.0#	70	0.00
28		bis(2-Chloroethoxy)methane	0.435	0.254	41.6#	67	0.00
29	C	2,4-Dichlorophenol	0.332	0.216	34.9#	73	0.00
30		1,2,4-Trichlorobenzene	0.392	0.247	37.0#	73	0.00
31		Naphthalene	1.038	0.643	38.1#	72	0.00
32		Benzoic acid	0.173	0.170	1.7	108	-0.15
33		4-Chloroaniline	0.445	0.262	41.1#	67	0.00
34	C	Hexachlorobutadiene	0.238	0.151	36.6#	73	0.00
35		Caprolactam	0.109	0.056	48.6#	56	-0.02
36	C	4-Chloro-3-methylphenol	0.332	0.190	42.8#	64	0.00
37		2-Methylnaphthalene	0.781	0.453	42.0#	66	0.00
38		1-Methylnaphthalene	0.764	0.435	43.1#	65	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
40		1,2,4,5-Tetrachlorobenzene	0.630	0.425	32.5#	66	0.00
41	P	Hexachlorocyclopentadiene	0.335	0.201	40.0#	57	0.00
42	S	2,4,6-Tribromophenol	0.271	0.158	41.7#	58	0.00
43	C	2,4,6-Trichlorophenol	0.410	0.280	31.7#	65	0.00
44		2,4,5-Trichlorophenol	0.453	0.306	32.5#	64	0.00
45	S	2-Fluorobiphenyl	1.380	0.925	33.0#	66	0.00
46		1,1'-Biphenyl	1.479	0.952	35.6#	64	0.00
47		2-Chloronaphthalene	1.159	0.752	35.1#	64	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.254	0.156	38.6#	57	0.00
49		Acenaphthylene	1.781	1.118	37.2#	62	0.00
50		Dimethylphthalate	1.540	0.882	42.7#	57	-0.01
51		2,6-Dinitrotoluene	0.316	0.185	41.5#	56	0.00
52	C	Acenaphthene	1.165	0.673	42.2#	57	0.00
53		3-Nitroaniline	0.329	0.190	42.2#	55	-0.01
54	P	2,4-Dinitrophenol	0.141	0.004#	97.2#	2#	0.00
55		Dibenzofuran	1.868	1.126	39.7#	60	0.00
56	P	4-Nitrophenol	0.259	0.134	48.3#	48#	0.00
57		2,4-Dinitrotoluene	0.422	0.229	45.7#	51	-0.01
58		Fluorene	1.456	0.860	40.9#	59	0.00
59		2,3,4,6-Tetrachlorophenol	0.406	0.238	41.4#	56	0.00
60		Diethylphthalate	1.506	0.823	45.4#	54	0.00
61		4-Chlorophenyl-phenylether	0.811	0.474	41.6#	58	0.00
62		4-Nitroaniline	0.334	0.186	44.3#	53	-0.01
63		Azobenzene	1.215	0.693	43.0#	56	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	89	0.00
65		4,6-Dinitro-2-methylphenol	0.101	0.013	87.1#	10#	0.00
66	c	n-Nitrosodiphenylamine	0.604	0.390	35.4#	56	0.00
67		4-Bromophenyl-phenylether	0.219	0.135	38.4#	58	0.00
68		Hexachlorobenzene	0.269	0.167	37.9#	55	0.00
69		Atrazine	0.226	0.135	40.3#	51	0.00
70	C	Pentachlorophenol	0.146	0.108	26.0#	60	0.00
71		Phenanthrene	1.078	0.674	37.5#	54	0.00
72		Anthracene	1.051	0.662	37.0#	54	0.00
73		Carbazole	1.031	0.631	38.8#	53	0.00
74		Di-n-butylphthalate	1.128	0.698	38.1#	50	0.00
75	C	Fluoranthene	1.240	0.760	38.7#	51	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	88	-0.01
77		Benzidine	0.601	0.420	30.1#	57	0.00
78		Pyrene	1.301	0.794	39.0#	52	0.00
79	S	Terphenyl-d14	0.980	0.650	33.7#	64	0.00
80		Butylbenzylphthalate	0.519	0.310	40.3#	49#	0.00
81		Benzo(a)anthracene	1.297	0.814	37.2#	54	0.00
82		3,3'-Dichlorobenzidine	0.477	0.315	34.0#	54	0.00
83		Chrysene	1.213	0.773	36.3#	55	0.00
84		Bis(2-ethylhexyl)phthalate	0.784	0.490	37.5#	52	0.00
85	c	Di-n-octyl phthalate	1.213	0.819	32.5#	53	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	108	-0.01
87		Indeno(1,2,3-cd)pyrene	1.343	0.927	31.0#	77	-0.02
88		Benzo(b)fluoranthene	1.286	0.730	43.2#	58	-0.01
89		Benzo(k)fluoranthene	1.279	0.729	43.0#	60	-0.01
90	C	Benzo(a)pyrene	1.037	0.619	40.3#	62	-0.01
91		Dibenzo(a,h)anthracene	1.133	0.775	31.6#	75	-0.02
92		Benzo(g,h,i)perylene	1.060	0.745	29.7#	80	-0.02

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(#) = Out of Range SPCC's out = 1 CCC's out = 13