Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091420\
Data File : BP003269.D
Acq On : 14 Sep 2020 15:26
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Sep 14 17:12:21 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 01 14:37:05 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	77	-0.01
2		1,4-Dioxane	0.432	0.394	8.8	76	-0.01
3		Pyridine	1.138	1.079	5.2	77	-0.01
4		n-Nitrosodimethylamine	0.297	0.326	-9.8	88	-0.01
5	S	2-Fluorophenol	1.131	1.073	5.1	77	-0.01
6		Aniline	1.544	1.544	0.0	81	-0.01
7	S	Phenol-d6	1.418	1.382	2.5	80	0.00
8		2-Chlorophenol	1.261	1.226	2.8	80	-0.01
9		Benzaldehyde	0.661	0.671	-1.5	82	-0.01
10	C	Phenol	1.314	1.287	2.1	80	0.00
11		bis(2-Chloroethyl)ether	1.040	1.011	2.8	80	-0.01
12		1,3-Dichlorobenzene	1.533	1.467	4.3	80	-0.02
13	C	1,4-Dichlorobenzene	1.548	1.481	4.3	80	-0.01
14		1,2-Dichlorobenzene	1.468	1.400	4.6	80	-0.01
15		Benzyl Alcohol	0.799	0.873	-9.3	87	-0.01
16		2,2'-oxybis(1-Chloropropane	0.881	0.885	-0.5	82	-0.02
17		2-Methylphenol	0.936	0.941	-0.5	81	0.00
18		Hexachloroethane	0.506	0.507	-0.2	83	-0.02
19	P	n-Nitroso-di-n-propylamine	0.641	0.653	-1.9	82	-0.01
20		3+4-Methylphenols	1.261	1.258	0.2	81	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	77	-0.02
22		Acetophenone	0.445	0.429	3.6	80	-0.01
23	S	Nitrobenzene-d5	0.277	0.284	-2.5	84	0.00
24		Nitrobenzene	0.272	0.277	-1.8	84	0.00
25		Isophorone	0.489	0.505	-3.3	84	0.00
26	C	2-Nitrophenol	0.162	0.180	-11.1	87	-0.01
27		2,4-Dimethylphenol	0.244	0.243	0.4	81	0.00
28		bis(2-Chloroethoxy)methane	0.365	0.351	3.8	80	-0.01
29	C	2,4-Dichlorophenol	0.303	0.306	-1.0	83	0.00
30		1,2,4-Trichlorobenzene	0.361	0.348	3.6	82	-0.01
31		Naphthalene	1.028	0.976	5.1	80	-0.01
32		Benzoic acid	0.159	0.159	0.0	72	0.03
33		4-Chloroaniline	0.429	0.424	1.2	81	0.00
34	C	Hexachlorobutadiene	0.216	0.218	-0.9	86	-0.02
35		Caprolactam	0.091	0.099	-8.8	87	0.01
36	C	4-Chloro-3-methylphenol	0.283	0.293	-3.5	85	0.00
37		2-Methylnaphthalene	0.735	0.703	4.4	81	0.00
38		1-Methylnaphthalene	0.699	0.661	5.4	80	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	78	0.00
40		1,2,4,5-Tetrachlorobenzene	0.593	0.582	1.9	82	-0.01
41	P	Hexachlorocyclopentadiene	0.280	0.305	-8.9	84	-0.01
42	S	2,4,6-Tribromophenol	0.270	0.291	-7.8	90	0.00
43	C	2,4,6-Trichlorophenol	0.390	0.403	-3.3	84	0.00
44		2,4,5-Trichlorophenol	0.412	0.409	0.7	82	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.366	1.271	7.0	79	-0.01
46		1,1'-Biphenyl	1.500	1.435	4.3	81	0.00
47		2-Chloronaphthalene	1.228	1.179	4.0	81	-0.01
48		2-Nitroaniline	0.230	0.267	-16.1	92	0.00
49		Acenaphthylene	1.857	1.823	1.8	82	0.00
50		Dimethylphthalate	1.525	1.482	2.8	83	0.00
51		2,6-Dinitrotoluene	0.317	0.326	-2.8	84	0.00
52	C	Acenaphthene	1.141	1.084	5.0	80	0.00
53		3-Nitroaniline	0.359	0.375	-4.5	85	0.00
54	P	2,4-Dinitrophenol	0.142	0.198	-39.4#	109	0.00
55		Dibenzofuran	1.792	1.709	4.6	82	-0.01
56	P	4-Nitrophenol	0.258	0.247	4.3	74	0.01
57		2,4-Dinitrotoluene	0.426	0.460	-8.0	87	0.00
58		Fluorene	1.380	1.302	5.7	81	0.00
59		2,3,4,6-Tetrachlorophenol	0.373	0.366	1.9	81	0.00
60		Diethylphthalate	1.496	1.490	0.4	85	-0.01
61		4-Chlorophenyl-phenylether	0.713	0.677	5.0	82	-0.01
62		4-Nitroaniline	0.369	0.400	-8.4	89	0.00
63		Azobenzene	0.993	1.001	-0.8	85	-0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	79	0.00
65		4,6-Dinitro-2-methylphenol	0.096	0.108	-12.5	91	0.00
66	c	n-Nitrosodiphenylamine	0.585	0.569	2.7	83	-0.01
67		4-Bromophenyl-phenylether	0.221	0.220	0.5	85	-0.01
68		Hexachlorobenzene	0.263	0.263	0.0	87	0.00
69		Atrazine	0.197	0.202	-2.5	84	0.00
70	C	Pentachlorophenol	0.126	0.112	11.1	69	0.00
71		Phenanthrene	1.084	1.023	5.6	81	0.00
72		Anthracene	1.037	1.007	2.9	83	0.00
73		Carbazole	0.995	0.980	1.5	84	0.00
74		Di-n-butylphthalate	1.170	1.212	-3.6	86	-0.01
75	C	Fluoranthene	1.253	1.230	1.8	84	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	81	0.00
77		Benzidine	0.457	0.662	-44.9#	108	-0.01
78		Pyrene	1.298	1.208	6.9	83	0.00
79	S	Terphenyl-d14	0.909	0.876	3.6	83	0.00
80		Butylbenzylphthalate	0.534	0.571	-6.9	89	-0.01
81		Benzo(a)anthracene	1.253	1.209	3.5	85	0.00
82		3,3'-Dichlorobenzidine	0.461	0.483	-4.8	87	0.00
83		Chrysene	1.199	1.149	4.2	85	0.00
84		Bis(2-ethylhexyl)phthalate	0.772	0.756	2.1	82	-0.01
85	c	Di-n-octyl phthalate	1.323	1.425	-7.7	91	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	86	0.00
87		Indeno(1,2,3-cd)pyrene	1.491	1.409	5.5	91	0.00

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88	Benzo(b)fluoranthene	1.243	1.155	7.1	90	0.00
89	Benzo(k)fluoranthene	1.188	1.083	8.8	86	0.00
90 C	Benzo(a)pyrene	1.153	1.079	6.4	89	0.00
91	Dibenzo(a,h)anthracene	1.225	1.154	5.8	91	0.00
92	Benzo(g,h,i)perylene	1.230	1.182	3.9	93	0.00

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