Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092521\
Data File : BP007182.D
Acq On : 25 Sep 2021 12:27
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Sep 26 02:12:52 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 23 11:17:25 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	76	-0.01
2		1,4-Dioxane	0.483	0.441	8.7	72	-0.01
3		Pyridine	1.322	1.221	7.6	71	-0.01
4		n-Nitrosodimethylamine	0.453	0.424	6.4	72	-0.01
5	S	2-Fluorophenol	1.195	1.128	5.6	74	0.00
6		Aniline	1.800	1.675	6.9	71	-0.01
7	S	Phenol-d6	1.633	1.526	6.6	71	-0.01
8		2-Chlorophenol	1.299	1.238	4.7	74	-0.01
9		Benzaldehyde	0.920	0.837	9.0	69	-0.01
10	C	Phenol	1.574	1.453	7.7	71	0.00
11		bis(2-Chloroethyl)ether	1.243	1.131	9.0	70	-0.01
12		1,3-Dichlorobenzene	1.465	1.402	4.3	75	-0.01
13	C	1,4-Dichlorobenzene	1.493	1.436	3.8	75	-0.01
14		1,2-Dichlorobenzene	1.447	1.385	4.3	75	-0.01
15		Benzyl Alcohol	1.046	1.020	2.5	74	0.00
16		2,2'-oxybis(1-Chloropropane	1.423	1.279	10.1	70	0.00
17		2-Methylphenol	1.061	1.012	4.6	73	-0.01
18		Hexachloroethane	0.501	0.487	2.8	76	-0.02
19	P	n-Nitroso-di-n-propylamine	0.888	0.819	7.8	71	-0.01
20		3+4-Methylphenols	1.468	1.391	5.2	72	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	73	-0.01
22		Acetophenone	0.482	0.462	4.1	71	-0.01
23	S	Nitrobenzene-d5	0.343	0.335	2.3	72	-0.01
24		Nitrobenzene	0.327	0.318	2.8	71	-0.01
25		Isophorone	0.599	0.581	3.0	71	-0.01
26	C	2-Nitrophenol	0.171	0.181	-5.8	77	-0.01
27		2,4-Dimethylphenol	0.269	0.262	2.6	71	0.00
28		bis(2-Chloroethoxy)methane	0.422	0.394	6.6	69	-0.01
29	C	2,4-Dichlorophenol	0.314	0.318	-1.3	74	-0.01
30		1,2,4-Trichlorobenzene	0.357	0.359	-0.6	75	-0.01
31		Naphthalene	1.021	0.992	2.8	72	-0.01
32		Benzoic acid	0.207	0.219	-5.8	74	0.00
33		4-Chloroaniline	0.422	0.416	1.4	72	-0.01
34	C	Hexachlorobutadiene	0.214	0.225	-5.1	79	-0.02
35		Caprolactam	0.096	0.104	-8.3	76	0.00
36	C	4-Chloro-3-methylphenol	0.319	0.319	0.0	74	0.00
37		2-Methylnaphthalene	0.715	0.694	2.9	73	0.00
38		1-Methylnaphthalene	0.698	0.676	3.2	72	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	74	-0.01
40		1,2,4,5-Tetrachlorobenzene	0.612	0.616	-0.7	76	-0.01
41	P	Hexachlorocyclopentadiene	0.339	0.318	6.2	69	0.00
42	S	2,4,6-Tribromophenol	0.259	0.289	-11.6	83	-0.01
43	C	2,4,6-Trichlorophenol	0.418	0.428	-2.4	76	-0.01
44		2,4,5-Trichlorophenol	0.450	0.451	-0.2	74	0.00
45	S	2-Fluorobiphenyl	1.396	1.350	3.3	74	0.00
46		1,1'-Biphenyl	1.488	1.421	4.5	73	-0.01
47		2-Chloronaphthalene	1.152	1.113	3.4	74	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.281	0.288	-2.5	75	0.00
49		Acenaphthylene	1.775	1.742	1.9	73	0.00
50		Dimethylphthalate	1.438	1.412	1.8	75	0.00
51		2,6-Dinitrotoluene	0.294	0.303	-3.1	76	-0.01
52	C	Acenaphthene	1.075	1.029	4.3	72	0.00
53		3-Nitroaniline	0.318	0.330	-3.8	75	0.00
54	P	2,4-Dinitrophenol	0.169	0.184	-8.9	76	0.00
55		Dibenzofuran	1.794	1.734	3.3	74	0.00
56	P	4-Nitrophenol	0.258	0.272	-5.4	73	0.00
57		2,4-Dinitrotoluene	0.388	0.415	-7.0	78	0.00
58		Fluorene	1.386	1.346	2.9	74	-0.01
59		2,3,4,6-Tetrachlorophenol	0.395	0.409	-3.5	78	0.00
60		Diethylphthalate	1.394	1.383	0.8	75	0.00
61		4-Chlorophenyl-phenylether	0.732	0.723	1.2	75	-0.01
62		4-Nitroaniline	0.320	0.341	-6.6	76	0.00
63		Azobenzene	1.208	1.161	3.9	72	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	76	0.00
65		4,6-Dinitro-2-methylphenol	0.120	0.132	-10.0	79	0.00
66	c	n-Nitrosodiphenylamine	0.595	0.573	3.7	75	0.00
67		4-Bromophenyl-phenylether	0.219	0.219	0.0	78	-0.01
68		Hexachlorobenzene	0.242	0.248	-2.5	80	0.00
69		Atrazine	0.210	0.214	-1.9	77	-0.01
70	C	Pentachlorophenol	0.150	0.152	-1.3	76	0.00
71		Phenanthrene	1.072	1.039	3.1	76	0.00
72		Anthracene	1.049	1.031	1.7	76	-0.01
73		Carbazole	1.028	1.016	1.2	75	-0.01
74		Di-n-butylphthalate	1.139	1.144	-0.4	76	0.00
75	C	Fluoranthene	1.236	1.242	-0.5	78	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	78	0.00
77		Benzidine	0.615	0.565	8.1	72	-0.01
78		Pyrene	1.312	1.264	3.7	77	0.00
79	S	Terphenyl-d14	1.044	1.012	3.1	77	-0.01
80		Butylbenzylphthalate	0.525	0.532	-1.3	79	-0.01
81		Benzo(a)anthracene	1.298	1.251	3.6	77	-0.01
82		3,3'-Dichlorobenzidine	0.446	0.456	-2.2	79	0.00
83		Chrysene	1.241	1.187	4.4	76	-0.01
84		Bis(2-ethylhexyl)phthalate	0.755	0.740	2.0	76	-0.01
85	c	Di-n-octyl phthalate	1.286	1.286	0.0	78	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	81	-0.01
87		Indeno(1,2,3-cd)pyrene	1.437	1.406	2.2	81	-0.02
88		Benzo(b)fluoranthene	1.195	1.144	4.3	78	-0.01
89		Benzo(k)fluoranthene	1.210	1.127	6.9	78	-0.02
90	C	Benzo(a)pyrene	1.015	0.984	3.1	80	-0.02
91		Dibenzo(a,h)anthracene	1.205	1.171	2.8	80	-0.02
92		Benzo(g,h,i)perylene	1.175	1.160	1.3	82	-0.01

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