Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040925\
Data File : BP024243.D
Acq On : 09 Apr 2025 20:24
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Apr 10 01:01:24 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 13 05:55:58 2025
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	75	-0.03
2		1,4-Dioxane	0.497	0.507	-2.0	81	-0.02
3		Pyridine	1.358	1.378	-1.5	79	-0.02
4		n-Nitrosodimethylamine	0.454	0.456	-0.4	79	-0.02
5	S	2-Fluorophenol	1.253	1.314	-4.9	81	-0.02
6		Aniline	1.527	1.587	-3.9	80	-0.03
7	S	Phenol-d6	1.675	1.750	-4.5	80	-0.02
8		2-Chlorophenol	1.353	1.406	-3.9	81	-0.03
9		Benzaldehyde	0.809	0.906	-12.0	87	-0.03
10	C	Phenol	1.691	1.751	-3.5	81	-0.02
11		bis(2-Chloroethyl)ether	1.334	1.331	0.2	78	-0.02
12		1,3-Dichlorobenzene	1.459	1.490	-2.1	80	-0.03
13	C	1,4-Dichlorobenzene	1.490	1.501	-0.7	80	-0.04
14		1,2-Dichlorobenzene	1.430	1.444	-1.0	80	-0.03
15		Benzyl Alcohol	1.061	1.160	-9.3	83	-0.03
16		2,2'-oxybis(1-Chloropropane	1.361	1.377	-1.2	80	-0.04
17		2-Methylphenol	1.061	1.128	-6.3	81	-0.03
18		Hexachloroethane	0.529	0.539	-1.9	80	-0.03
19	P	n-Nitroso-di-n-propylamine	0.972	1.022	-5.1	80	-0.03
20		3+4-Methylphenols	1.457	1.572	-7.9	82	-0.03
21	I	Naphthalene-d8	1.000	1.000	0.0	76	-0.03
22		Acetophenone	0.506	0.520	-2.8	80	-0.03
23	S	Nitrobenzene-d5	0.354	0.369	-4.2	80	-0.03
24		Nitrobenzene	0.350	0.364	-4.0	81	-0.03
25		Isophorone	0.608	0.663	-9.0	84	-0.03
26	C	2-Nitrophenol	0.167	0.188	-12.6	84	-0.03
27		2,4-Dimethylphenol	0.215	0.231	-7.4	82	-0.03
28		bis(2-Chloroethoxy)methane	0.427	0.435	-1.9	79	-0.03
29	C	2,4-Dichlorophenol	0.291	0.316	-8.6	81	-0.03
30		1,2,4-Trichlorobenzene	0.321	0.325	-1.2	79	-0.03
31		Naphthalene	1.060	1.091	-2.9	80	-0.03
32		Benzoic acid	0.209	0.280	-34.0#	97	-0.03
33		4-Chloroaniline	0.379	0.413	-9.0	83	-0.03
34	C	Hexachlorobutadiene	0.185	0.188	-1.6	79	-0.04
35		Caprolactam	0.095	0.118	-24.2	92	-0.03
36	C	4-Chloro-3-methylphenol	0.313	0.361	-15.3	87	-0.03
37		2-Methylnaphthalene	0.716	0.753	-5.2	81	-0.02
38		1-Methylnaphthalene	0.697	0.733	-5.2	83	-0.03
39	I	Acenaphthene-d10	1.000	1.000	0.0	80	-0.02
40		1,2,4,5-Tetrachlorobenzene	0.578	0.567	1.9	81	-0.03
41	P	Hexachlorocyclopentadiene	0.208	0.176	15.4	66	-0.02
42	S	2,4,6-Tribromophenol	0.252	0.290	-15.1	92	-0.02
43	C	2,4,6-Trichlorophenol	0.367	0.392	-6.8	84	-0.03
44		2,4,5-Trichlorophenol	0.403	0.432	-7.2	85	-0.02
45	S	2-Fluorobiphenyl	1.356	1.326	2.2	81	-0.02
46		1,1'-Biphenyl	1.503	1.493	0.7	82	-0.02
47		2-Chloronaphthalene	1.130	1.137	-0.6	83	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.289	0.334	-15.6	92	-0.03
49		Acenaphthylene	1.710	1.796	-5.0	87	-0.03
50		Dimethylphthalate	1.399	1.465	-4.7	86	-0.02
51		2,6-Dinitrotoluene	0.294	0.318	-8.2	87	-0.02
52	C	Acenaphthene	1.093	1.121	-2.6	85	-0.02
53		3-Nitroaniline	0.318	0.354	-11.3	88	-0.02
54	P	2,4-Dinitrophenol	0.156	0.152	2.6	76	-0.03
55		Dibenzofuran	1.770	1.800	-1.7	84	-0.02
56	P	4-Nitrophenol	0.266	0.321	-20.7	93	-0.02
57		2,4-Dinitrotoluene	0.389	0.444	-14.1	90	-0.02
58		Fluorene	1.379	1.424	-3.3	85	-0.02
59		2,3,4,6-Tetrachlorophenol	0.356	0.392	-10.1	88	-0.02
60		Diethylphthalate	1.394	1.481	-6.2	87	-0.02
61		4-Chlorophenyl-phenylether	0.676	0.684	-1.2	83	-0.02
62		4-Nitroaniline	0.334	0.384	-15.0	90	-0.02
63		Azobenzene	1.314	1.403	-6.8	87	-0.02
64	I	Phenanthrene-d10	1.000	1.000	0.0	85	-0.02
65		4,6-Dinitro-2-methylphenol	0.119	0.112	5.9	79	-0.02
66	c	n-Nitrosodiphenylamine	0.607	0.621	-2.3	88	-0.02
67		4-Bromophenyl-phenylether	0.220	0.225	-2.3	88	-0.02
68		Hexachlorobenzene	0.260	0.266	-2.3	89	-0.02
69		Atrazine	0.147	0.162	-10.2	101	-0.02
70	C	Pentachlorophenol	0.168	0.199	-18.5	95	-0.02
71		Phenanthrene	1.093	1.106	-1.2	87	-0.02
72		Anthracene	1.064	1.103	-3.7	88	-0.02
73		Carbazole	1.026	1.094	-6.6	89	-0.02
74		Di-n-butylphthalate	1.218	1.336	-9.7	90	-0.02
75	C	Fluoranthene	1.265	1.324	-4.7	89	-0.02
76	I	Chrysene-d12	1.000	1.000	0.0	77	-0.03
77		Benzidine	0.219	0.350	-59.8#	105	-0.02
78		Pyrene	1.243	1.354	-8.9	87	-0.02
79	S	Terphenyl-d14	0.969	1.033	-6.6	84	-0.02
80		Butylbenzylphthalate	0.513	0.603	-17.5	89	-0.02
81		Benzo(a)anthracene	1.243	1.312	-5.6	84	-0.02
82		3,3'-Dichlorobenzidine	0.417	0.517	-24.0	93	-0.02
83		Chrysene	1.195	1.232	-3.1	82	-0.03
84		Bis(2-ethylhexyl)phthalate	0.774	0.887	-14.6	87	-0.02
85	c	Di-n-octyl phthalate	1.240	1.532	-23.5#	93	-0.04
86	I	Perylene-d12	1.000	1.000	0.0	87	-0.08
87		Indeno(1,2,3-cd)pyrene	1.396	1.453	-4.1	92	-0.11
88		Benzo(b)fluoranthene	1.206	1.190	1.3	87	-0.05
89		Benzo(k)fluoranthene	1.180	1.144	3.1	88	-0.05
90	C	Benzo(a)pyrene	1.049	1.077	-2.7	91	-0.06
91		Dibenzo(a,h)anthracene	1.161	1.199	-3.3	92	-0.12
92		Benzo(g,h,i)perylene	1.180	1.217	-3.1	92	-0.12

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

(#) = Out of Range SPCC's out = 0 CCC's out = 1