Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091924\
Data File : BP021956.D
Acq On : 19 Sep 2024 18:25
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP091924

Quant Time: Sep 20 01:25:53 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 20 01:22:00 2024
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	116	0.00
2		1,4-Dioxane	0.442	0.414	6.3	111	0.00
3		Pyridine	1.179	1.110	5.9	110	0.00
4		n-Nitrosodimethylamine	0.622	0.577	7.2	110	0.00
5	S	2-Fluorophenol	1.064	1.031	3.1	113	0.00
6		Aniline	1.214	1.190	2.0	114	0.00
7	S	Phenol-d6	1.382	1.350	2.3	115	0.00
8		2-Chlorophenol	1.138	1.113	2.2	113	0.00
9		Benzaldehyde	0.722	0.735	-1.8	116	0.00
10	C	Phenol	1.372	1.331	3.0	114	0.00
11		bis(2-Chloroethyl)ether	1.034	1.020	1.4	117	0.00
12		1,3-Dichlorobenzene	1.437	1.412	1.7	115	0.00
13	C	1,4-Dichlorobenzene	1.457	1.422	2.4	114	0.00
14		1,2-Dichlorobenzene	1.398	1.338	4.3	111	0.00
15		Benzyl Alcohol	1.060	1.067	-0.7	116	0.00
16		2,2'-oxybis(1-Chloropropane	1.102	1.103	-0.1	117	0.00
17		2-Methylphenol	0.927	0.936	-1.0	114	0.00
18		Hexachloroethane	0.536	0.523	2.4	113	0.00
19	P	n-Nitroso-di-n-propylamine	0.859	0.882	-2.7	116	0.00
20		3+4-Methylphenols	1.309	1.306	0.2	115	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	111	0.00
22		Acetophenone	0.488	0.480	1.6	114	0.00
23	S	Nitrobenzene-d5	0.378	0.380	-0.5	114	0.00
24		Nitrobenzene	0.384	0.385	-0.3	113	0.00
25		Isophorone	0.623	0.647	-3.9	116	0.00
26	C	2-Nitrophenol	0.168	0.173	-3.0	115	0.00
27		2,4-Dimethylphenol	0.194	0.200	-3.1	116	0.00
28		bis(2-Chloroethoxy)methane	0.374	0.376	-0.5	114	0.00
29	C	2,4-Dichlorophenol	0.330	0.335	-1.5	111	0.00
30		1,2,4-Trichlorobenzene	0.411	0.405	1.5	112	0.00
31		Naphthalene	0.998	0.989	0.9	113	0.00
32		Benzoic acid	0.227	0.236	-4.0	112	0.00
33		4-Chloroaniline	0.360	0.361	-0.3	111	0.00
34	C	Hexachlorobutadiene	0.321	0.319	0.6	114	0.01
35		Caprolactam	0.098	0.100	-2.0	112	0.00
36	C	4-Chloro-3-methylphenol	0.354	0.359	-1.4	113	0.00
37		2-Methylnaphthalene	0.730	0.732	-0.3	114	0.00
38		1-Methylnaphthalene	0.721	0.713	1.1	113	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	115	0.01
40		1,2,4,5-Tetrachlorobenzene	0.700	0.674	3.7	113	0.00
41	P	Hexachlorocyclopentadiene	0.260	0.255	1.9	116	0.00
42	S	2,4,6-Tribromophenol	0.371	0.380	-2.4	121	-0.01
43	C	2,4,6-Trichlorophenol	0.429	0.432	-0.7	115	0.00
44		2,4,5-Trichlorophenol	0.482	0.483	-0.2	116	0.00
45	S	2-Fluorobiphenyl	1.370	1.333	2.7	116	0.00
46		1,1'-Biphenyl	1.364	1.327	2.7	116	0.00
47		2-Chloronaphthalene	1.103	1.088	1.4	116	0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.301	0.312	-3.7	114	0.00
49		Acenaphthylene	1.556	1.530	1.7	113	0.00
50		Dimethylphthalate	1.451	1.445	0.4	116	0.00
51		2,6-Dinitrotoluene	0.323	0.328	-1.5	116	0.01
52	C	Acenaphthene	1.013	0.989	2.4	115	0.00
53		3-Nitroaniline	0.288	0.292	-1.4	112	0.00
54	P	2,4-Dinitrophenol	0.189	0.198	-4.8	118	0.00
55		Dibenzofuran	1.702	1.674	1.6	116	0.00
56	P	4-Nitrophenol	0.251	0.258	-2.8	111	0.01
57		2,4-Dinitrotoluene	0.457	0.471	-3.1	116	0.00
58		Fluorene	1.421	1.393	2.0	115	0.00
59		2,3,4,6-Tetrachlorophenol	0.465	0.463	0.4	116	0.00
60		Diethylphthalate	1.474	1.488	-0.9	119	0.00
61		4-Chlorophenyl-phenylether	0.828	0.817	1.3	118	0.00
62		4-Nitroaniline	0.317	0.321	-1.3	112	0.00
63		Azobenzene	1.220	1.200	1.6	115	0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	115	-0.01
65		4,6-Dinitro-2-methylphenol	0.116	0.122	-5.2	118	-0.02
66	c	n-Nitrosodiphenylamine	0.498	0.500	-0.4	115	0.01
67		4-Bromophenyl-phenylether	0.233	0.236	-1.3	120	0.00
68		Hexachlorobenzene	0.291	0.292	-0.3	119	0.00
69		Atrazine	0.178	0.160	10.1	115	0.00
70	C	Pentachlorophenol	0.196	0.198	-1.0	115	0.00
71		Phenanthrene	0.960	0.938	2.3	115	-0.01
72		Anthracene	0.929	0.916	1.4	114	0.00
73		Carbazole	0.898	0.897	0.1	113	0.00
74		Di-n-butylphthalate	1.054	1.111	-5.4	125	0.00
75	C	Fluoranthene	1.340	1.362	-1.6	118	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	114	0.00
77		Benzidine	0.244	0.349	-43.0#	99	0.00
78		Pyrene	1.210	1.215	-0.4	114	0.00
79	S	Terphenyl-d14	1.065	1.087	-2.1	117	0.00
80		Butylbenzylphthalate	0.432	0.474	-9.7	125	0.00
81		Benzo(a)anthracene	1.285	1.301	-1.2	118	0.00
82		3,3'-Dichlorobenzidine	0.466	0.495	-6.2	114	0.00
83		Chrysene	1.210	1.197	1.1	115	0.00
84		Bis(2-ethylhexyl)phthalate	0.633	0.706	-11.5	132	0.00
85	c	Di-n-octyl phthalate	1.076	1.208	-12.3	132	0.00
86	I	Perylene-d12	1.000	1.000	0.0	115	0.00
87		Indeno(1,2,3-cd)pyrene	1.330	1.309	1.6	114	-0.03
88		Benzo(b)fluoranthene	1.183	1.166	1.4	112	0.00
89		Benzo(k)fluoranthene	1.119	1.114	0.4	117	0.00
90	C	Benzo(a)pyrene	1.019	1.031	-1.2	116	0.00
91		Dibenzo(a,h)anthracene	1.098	1.076	2.0	115	0.00
92		Benzo(g,h,i)perylene	1.125	1.114	1.0	114	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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