Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081324\
Data File : BP021489.D
Acq On : 13 Aug 2024 19:03
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP081324

Quant Time: Aug 14 00:10:09 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 14 00:00:30 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	132	0.00
2		1,4-Dioxane	0.639	0.606	5.2	125	0.00
3		Pyridine	1.783	1.803	-1.1	133	0.00
4		n-Nitrosodimethylamine	0.536	0.539	-0.6	134	0.00
5	S	2-Fluorophenol	1.347	1.361	-1.0	134	0.00
6		Aniline	1.964	2.043	-4.0	137	0.00
7	S	Phenol-d6	1.965	2.109	-7.3	141	0.00
8		2-Chlorophenol	1.245	1.316	-5.7	140	0.00
9		Benzaldehyde	1.258	1.272	-1.1	141	0.00
10	C	Phenol	2.035	2.167	-6.5	140	0.00
11		bis(2-Chloroethyl)ether	1.722	1.788	-3.8	139	0.00
12		1,3-Dichlorobenzene	1.479	1.466	0.9	134	0.00
13	C	1,4-Dichlorobenzene	1.493	1.477	1.1	133	0.00
14		1,2-Dichlorobenzene	1.439	1.434	0.3	135	0.00
15		Benzyl Alcohol	1.410	1.567	-11.1	147	0.00
16		2,2'-oxybis(1-Chloropropane	1.585	1.660	-4.7	142	0.00
17		2-Methylphenol	1.286	1.405	-9.3	144	0.00
18		Hexachloroethane	0.599	0.610	-1.8	136	0.00
19	P	n-Nitroso-di-n-propylamine	1.357	1.469	-8.3	145	0.00
20		3+4-Methylphenols	1.734	1.945	-12.2	147	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	145	0.00
22		Acetophenone	0.614	0.613	0.2	145	0.00
23	S	Nitrobenzene-d5	0.387	0.433	-11.9	153#	0.00
24		Nitrobenzene	0.421	0.455	-8.1	149	0.00
25		Isophorone	0.915	0.931	-1.7	148	0.00
26	C	2-Nitrophenol	0.086	0.125	-45.3#	193#	0.00
27		2,4-Dimethylphenol	0.226	0.233	-3.1	150	0.00
28		bis(2-Chloroethoxy)methane	0.559	0.565	-1.1	149	0.00
29	C	2,4-Dichlorophenol	0.267	0.290	-8.6	154#	0.00
30		1,2,4-Trichlorobenzene	0.342	0.335	2.0	143	0.00
31		Naphthalene	1.045	1.035	1.0	144	0.00
32		Benzoic acid	0.137	0.187	-36.5#	210#	0.02
33		4-Chloroaniline	0.392	0.399	-1.8	146	0.00
34	C	Hexachlorobutadiene	0.215	0.211	1.9	142	0.00
35		Caprolactam	0.111	0.116	-4.5	148	0.00
36	C	4-Chloro-3-methylphenol	0.370	0.398	-7.6	151#	0.00
37		2-Methylnaphthalene	0.713	0.720	-1.0	148	0.00
38		1-Methylnaphthalene	0.703	0.706	-0.4	147	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	150	0.00
40		1,2,4,5-Tetrachlorobenzene	0.586	0.590	-0.7	149	0.01
41	P	Hexachlorocyclopentadiene	0.186	0.201	-8.1	157#	0.00
42	S	2,4,6-Tribromophenol	0.223	0.232	-4.0	147	0.00
43	C	2,4,6-Trichlorophenol	0.326	0.360	-10.4	158#	0.00
44		2,4,5-Trichlorophenol	0.362	0.401	-10.8	157#	0.00
45	S	2-Fluorobiphenyl	1.310	1.288	1.7	146	0.00
46		1,1'-Biphenyl	1.415	1.399	1.1	147	0.00
47		2-Chloronaphthalene	1.102	1.093	0.8	147	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.256	0.328	-28.1#	172#	0.00
49		Acenaphthylene	1.619	1.600	1.2	146	0.01
50		Dimethylphthalate	1.339	1.310	2.2	145	0.00
51		2,6-Dinitrotoluene	0.232	0.276	-19.0	162#	0.00
52	C	Acenaphthene	1.063	1.046	1.6	146	0.00
53		3-Nitroaniline	0.228	0.264	-15.8	156#	0.00
54	P	2,4-Dinitrophenol	0.069	0.081	-17.4	184#	0.00
55		Dibenzofuran	1.635	1.575	3.7	143	0.00
56	P	4-Nitrophenol	0.195	0.205	-5.1	152#	-0.01
57		2,4-Dinitrotoluene	0.284	0.344	-21.1	161#	0.00
58		Fluorene	1.343	1.278	4.8	141	0.00
59		2,3,4,6-Tetrachlorophenol	0.307	0.325	-5.9	148	0.00
60		Diethylphthalate	1.374	1.297	5.6	139	-0.01
61		4-Chlorophenyl-phenylether	0.694	0.673	3.0	145	0.00
62		4-Nitroaniline	0.241	0.257	-6.6	141	0.00
63		Azobenzene	1.689	1.598	5.4	140	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	133	0.00
65		4,6-Dinitro-2-methylphenol	0.054	0.068	-25.9#	172#	0.00
66	c	n-Nitrosodiphenylamine	0.536	0.564	-5.2	140	0.00
67		4-Bromophenyl-phenylether	0.212	0.224	-5.7	144	0.00
68		Hexachlorobenzene	0.250	0.257	-2.8	136	0.00
69		Atrazine	0.184	0.179	2.7	125	0.00
70	C	Pentachlorophenol	0.137	0.152	-10.9	148	0.00
71		Phenanthrene	0.962	0.951	1.1	131	0.00
72		Anthracene	0.943	0.935	0.8	131	0.00
73		Carbazole	0.909	0.851	6.4	122	0.00
74		Di-n-butylphthalate	1.116	1.075	3.7	125	0.00
75	C	Fluoranthene	1.180	1.071	9.2	119	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	106	0.02
77		Benzidine	0.420	0.354	15.7	79	0.00
78		Pyrene	1.307	1.407	-7.7	116	0.01
79	S	Terphenyl-d14	1.029	1.109	-7.8	117	0.01
80		Butylbenzylphthalate	0.437	0.510	-16.7	115	0.00
81		Benzo(a)anthracene	1.256	1.261	-0.4	106	0.01
82		3,3'-Dichlorobenzidine	0.403	0.412	-2.2	102	0.00
83		Chrysene	1.185	1.178	0.6	105	0.01
84		Bis(2-ethylhexyl)phthalate	0.742	0.805	-8.5	109	0.01
85	c	Di-n-octyl phthalate	1.199	1.270	-5.9	108	0.01
86	I	Perylene-d12	1.000	1.000	0.0	102	0.02
87		Indeno(1,2,3-cd)pyrene	1.296	1.332	-2.8	102	0.00
88		Benzo(b)fluoranthene	1.149	1.160	-1.0	102	0.01
89		Benzo(k)fluoranthene	1.121	1.120	0.1	100	0.01
90	C	Benzo(a)pyrene	0.997	1.013	-1.6	102	0.02
91		Dibenzo(a,h)anthracene	1.077	1.109	-3.0	103	0.02
92		Benzo(g,h,i)perylene	1.080	1.121	-3.8	104	0.03

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