Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091522\
Data File : BP011706.D
Acq On : 15 Sep 2022 10:23
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Sep 15 15:52:09 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 14 19:12:50 2022
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	98	0.00
2		1,4-Dioxane	0.484	0.492	-1.7	104	0.00
3		Pyridine	1.395	1.438	-3.1	103	0.00
4		n-Nitrosodimethylamine	0.535	0.538	-0.6	101	0.00
5	S	2-Fluorophenol	1.099	1.126	-2.5	103	0.00
6		Aniline	1.804	1.805	-0.1	98	0.00
7	S	Phenol-d6	1.465	1.487	-1.5	100	0.00
8		2-Chlorophenol	1.303	1.307	-0.3	99	0.00
9		Benzaldehyde	0.938	0.928	1.1	102	0.00
10	C	Phenol	1.647	1.653	-0.4	99	0.00
11		bis(2-Chloroethyl)ether	1.312	1.303	0.7	99	0.00
12		1,3-Dichlorobenzene	1.445	1.434	0.8	99	0.00
13	C	1,4-Dichlorobenzene	1.469	1.453	1.1	99	0.00
14		1,2-Dichlorobenzene	1.401	1.382	1.4	99	0.00
15		Benzyl Alcohol	1.046	1.043	0.3	98	0.00
16		2,2'-oxybis(1-Chloropropane	1.753	1.762	-0.5	100	0.00
17		2-Methylphenol	1.070	1.067	0.3	98	0.00
18		Hexachloroethane	0.508	0.510	-0.4	100	0.00
19	P	n-Nitroso-di-n-propylamine	0.948	0.936	1.3	96	0.00
20		3+4-Methylphenols	1.469	1.457	0.8	97	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	97	0.00
22		Acetophenone	0.469	0.468	0.2	97	0.00
23	S	Nitrobenzene-d5	0.337	0.343	-1.8	98	0.00
24		Nitrobenzene	0.351	0.354	-0.9	97	0.00
25		Isophorone	0.597	0.597	0.0	95	0.00
26	C	2-Nitrophenol	0.144	0.142	1.4	92	0.00
27		2,4-Dimethylphenol	0.249	0.253	-1.6	96	0.00
28		bis(2-Chloroethoxy)methane	0.381	0.378	0.8	96	0.00
29	C	2,4-Dichlorophenol	0.276	0.279	-1.1	96	0.00
30		1,2,4-Trichlorobenzene	0.306	0.301	1.6	96	0.00
31		Naphthalene	1.038	1.030	0.8	97	0.00
32		Benzoic acid	0.187	0.168	10.2	86	0.00
33		4-Chloroaniline	0.407	0.405	0.5	95	0.00
34	C	Hexachlorobutadiene	0.168	0.167	0.6	97	0.00
35		Caprolactam	0.090	0.083	7.8	89	0.00
36	C	4-Chloro-3-methylphenol	0.297	0.297	0.0	95	0.00
37		2-Methylnaphthalene	0.695	0.681	2.0	95	0.00
38		1-Methylnaphthalene	0.679	0.664	2.2	95	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	94	0.00
40		1,2,4,5-Tetrachlorobenzene	0.528	0.522	1.1	94	0.00
41	P	Hexachlorocyclopentadiene	0.263	0.267	-1.5	94	0.00
42	S	2,4,6-Tribromophenol	0.156	0.155	0.6	91	0.00
43	C	2,4,6-Trichlorophenol	0.359	0.367	-2.2	94	0.00
44		2,4,5-Trichlorophenol	0.401	0.405	-1.0	94	0.00
45	S	2-Fluorobiphenyl	1.287	1.294	-0.5	95	0.00
46		1,1'-Biphenyl	1.455	1.455	0.0	95	0.00
47		2-Chloronaphthalene	1.145	1.143	0.2	95	0.00

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Instrument :
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LabSampleId :
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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.321	0.336	-4.7	93	0.00
49		Acenaphthylene	1.771	1.803	-1.8	94	0.00
50		Dimethylphthalate	1.395	1.378	1.2	93	0.00
51		2,6-Dinitrotoluene	0.283	0.288	-1.8	92	0.00
52	C	Acenaphthene	1.178	1.169	0.8	93	0.00
53		3-Nitroaniline	0.326	0.338	-3.7	93	0.00
54	P	2,4-Dinitrophenol	0.138	0.129	6.5	89	0.00
55		Dibenzofuran	1.697	1.687	0.6	94	0.00
56	P	4-Nitrophenol	0.275	0.275	0.0	93	0.00
57		2,4-Dinitrotoluene	0.366	0.377	-3.0	91	0.00
58		Fluorene	1.376	1.388	-0.9	94	0.00
59		2,3,4,6-Tetrachlorophenol	0.330	0.324	1.8	92	0.00
60		Diethylphthalate	1.389	1.400	-0.8	94	0.00
61		4-Chlorophenyl-phenylether	0.636	0.634	0.3	94	0.00
62		4-Nitroaniline	0.328	0.348	-6.1	94	0.00
63		Azobenzene	1.335	1.386	-3.8	95	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	93	0.00
65		4,6-Dinitro-2-methylphenol	0.098	0.093	5.1	89	0.00
66	c	n-Nitrosodiphenylamine	0.595	0.602	-1.2	94	0.00
67		4-Bromophenyl-phenylether	0.187	0.186	0.5	92	0.00
68		Hexachlorobenzene	0.202	0.201	0.5	94	0.00
69		Atrazine	0.182	0.183	-0.5	92	0.00
70	C	Pentachlorophenol	0.128	0.127	0.8	93	0.00
71		Phenanthrene	1.088	1.090	-0.2	94	0.00
72		Anthracene	1.029	1.053	-2.3	94	0.00
73		Carbazole	0.971	0.990	-2.0	93	0.00
74		Di-n-butylphthalate	1.170	1.214	-3.8	93	0.00
75	C	Fluoranthene	1.174	1.183	-0.8	93	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	97	0.00
77		Benzidine	0.476	0.550	-15.5	96	0.00
78		Pyrene	1.349	1.332	1.3	95	0.00
79	S	Terphenyl-d14	0.920	0.923	-0.3	96	0.00
80		Butylbenzylphthalate	0.565	0.576	-1.9	94	0.00
81		Benzo(a)anthracene	1.289	1.283	0.5	96	0.00
82		3,3'-Dichlorobenzidine	0.392	0.408	-4.1	96	0.00
83		Chrysene	1.232	1.229	0.2	96	0.00
84		Bis(2-ethylhexyl)phthalate	0.813	0.843	-3.7	93	0.00
85	c	Di-n-octyl phthalate	1.395	1.410	-1.1	95	0.00
86	I	Perylene-d12	1.000	1.000	0.0	98	0.00
87		Indeno(1,2,3-cd)pyrene	1.441	1.508	-4.6	101	0.00
88		Benzo(b)fluoranthene	1.237	1.268	-2.5	99	0.00
89		Benzo(k)fluoranthene	1.247	1.246	0.1	97	0.00
90	C	Benzo(a)pyrene	1.022	1.042	-2.0	97	0.00
91		Dibenzo(a,h)anthracene	1.196	1.255	-4.9	101	0.00
92		Benzo(g,h,i)perylene	1.166	1.207	-3.5	100	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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