Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122122\
Data File : BF131746.D
Acq On : 21 Dec 2022 15:03
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Dec 22 05:20:31 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 21 03:12:37 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	86	0.00
2		1,4-Dioxane	0.555	0.569	-2.5	88	0.00
3		Pyridine	1.583	1.619	-2.3	87	0.01
4		n-Nitrosodimethylamine	0.716	0.734	-2.5	87	0.00
5	S	2-Fluorophenol	1.188	1.197	-0.8	87	0.01
6		Aniline	1.974	1.915	3.0	83	0.01
7	S	Phenol-d6	1.599	1.573	1.6	86	0.01
8		2-Chlorophenol	1.276	1.253	1.8	85	0.01
9		Benzaldehyde	1.073	0.940	12.4	89	0.01
10	C	Phenol	1.909	1.890	1.0	86	0.00
11		bis(2-Chloroethyl)ether	1.366	1.356	0.7	85	0.01
12		1,3-Dichlorobenzene	1.442	1.420	1.5	85	0.01
13	C	1,4-Dichlorobenzene	1.474	1.457	1.2	86	0.01
14		1,2-Dichlorobenzene	1.380	1.352	2.0	87	0.01
15		Benzyl Alcohol	1.428	1.379	3.4	84	0.00
16		2,2'-oxybis(1-Chloropropane	1.742	1.705	2.1	86	0.01
17		2-Methylphenol	1.204	1.166	3.2	84	0.00
18		Hexachloroethane	0.584	0.593	-1.5	89	0.01
19	P	n-Nitroso-di-n-propylamine	1.153	1.095	5.0	83	0.01
20		3+4-Methylphenols	1.550	1.499	3.3	84	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	84	0.01
22		Acetophenone	0.584	0.581	0.5	84	0.00
23	S	Nitrobenzene-d5	0.493	0.495	-0.4	83	0.01
24		Nitrobenzene	0.501	0.503	-0.4	83	0.01
25		Isophorone	0.837	0.840	-0.4	85	0.00
26	C	2-Nitrophenol	0.196	0.200	-2.0	85	0.01
27		2,4-Dimethylphenol	0.279	0.280	-0.4	85	0.01
28		bis(2-Chloroethoxy)methane	0.444	0.451	-1.6	85	0.00
29	C	2,4-Dichlorophenol	0.337	0.344	-2.1	86	0.01
30		1,2,4-Trichlorobenzene	0.396	0.402	-1.5	84	0.01
31		Naphthalene	1.092	1.097	-0.5	85	0.01
32		Benzoic acid	0.223	0.233	-4.5	82	0.00
33		4-Chloroaniline	0.426	0.421	1.2	83	0.01
34	C	Hexachlorobutadiene	0.266	0.270	-1.5	85	0.01
35		Caprolactam	0.101	0.101	0.0	85	0.01
36	C	4-Chloro-3-methylphenol	0.384	0.391	-1.8	85	0.01
37		2-Methylnaphthalene	0.745	0.748	-0.4	85	0.01
38		1-Methylnaphthalene	0.722	0.720	0.3	85	0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	87	0.01
40		1,2,4,5-Tetrachlorobenzene	0.684	0.705	-3.1	85	0.01
41	P	Hexachlorocyclopentadiene	0.310	0.325	-4.8	82	0.01
42	S	2,4,6-Tribromophenol	0.219	0.227	-3.7	88	0.01
43	C	2,4,6-Trichlorophenol	0.447	0.463	-3.6	86	0.01
44		2,4,5-Trichlorophenol	0.484	0.513	-6.0	87	0.01
45	S	2-Fluorobiphenyl	1.428	1.467	-2.7	87	0.01
46		1,1'-Biphenyl	1.504	1.537	-2.2	85	0.01
47		2-Chloronaphthalene	1.232	1.280	-3.9	86	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.444	0.467	-5.2	88	0.00
49		Acenaphthylene	1.848	1.913	-3.5	86	0.00
50		Dimethylphthalate	1.505	1.545	-2.7	88	0.01
51		2,6-Dinitrotoluene	0.324	0.346	-6.8	89	0.01
52	C	Acenaphthene	1.132	1.172	-3.5	87	0.01
53		3-Nitroaniline	0.331	0.336	-1.5	85	0.01
54	P	2,4-Dinitrophenol	0.194	0.203	-4.6	86	0.00
55		Dibenzofuran	1.736	1.787	-2.9	87	0.01
56	P	4-Nitrophenol	0.237	0.244	-3.0	84	0.01
57		2,4-Dinitrotoluene	0.435	0.465	-6.9	91	0.00
58		Fluorene	1.403	1.456	-3.8	88	0.01
59		2,3,4,6-Tetrachlorophenol	0.404	0.408	-1.0	84	0.01
60		Diethylphthalate	1.450	1.524	-5.1	91	0.01
61		4-Chlorophenyl-phenylether	0.746	0.774	-3.8	89	0.01
62		4-Nitroaniline	0.333	0.345	-3.6	88	0.01
63		Azobenzene	1.592	1.667	-4.7	89	0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	90	0.01
65		4,6-Dinitro-2-methylphenol	0.137	0.146	-6.6	90	0.01
66	c	n-Nitrosodiphenylamine	0.620	0.621	-0.2	89	0.01
67		4-Bromophenyl-phenylether	0.236	0.240	-1.7	89	0.01
68		Hexachlorobenzene	0.260	0.261	-0.4	88	0.01
69		Atrazine	0.207	0.216	-4.3	92	0.01
70	C	Pentachlorophenol	0.138	0.146	-5.8	90	0.01
71		Phenanthrene	1.091	1.078	1.2	89	0.01
72		Anthracene	1.068	1.068	0.0	90	0.01
73		Carbazole	0.924	0.931	-0.8	92	0.01
74		Di-n-butylphthalate	1.150	1.204	-4.7	98	0.01
75	C	Fluoranthene	1.215	1.208	0.6	92	0.02
76	I	Chrysene-d12	1.000	1.000	0.0	97	0.01
77		Benzidine	0.367	0.348	5.2	89	0.01
78		Pyrene	1.818	1.994	-9.7	93	0.01
79	S	Terphenyl-d14	1.283	1.437	-12.0	96	0.01
80		Butylbenzylphthalate	0.641	0.716	-11.7	100	0.02
81		Benzo(a)anthracene	1.440	1.546	-7.4	98	0.01
82		3,3'-Dichlorobenzidine	0.402	0.438	-9.0	105	0.01
83		Chrysene	1.360	1.442	-6.0	97	0.02
84		Bis(2-ethylhexyl)phthalate	0.765	0.848	-10.8	103	0.02
85	c	Di-n-octyl phthalate	1.068	1.256	-17.6	111	0.02
86	I	Perylene-d12	1.000	1.000	0.0	97	0.02
87		Indeno(1,2,3-cd)pyrene	1.343	1.574	-17.2	103	0.03
88		Benzo(b)fluoranthene	1.415	1.357	4.1	97	0.02
89		Benzo(k)fluoranthene	1.381	1.343	2.8	97	0.02
90	C	Benzo(a)pyrene	1.121	1.102	1.7	95	0.02
91		Dibenzo(a,h)anthracene	1.110	1.309	-17.9	104	0.04
92		Benzo(g,h,i)perylene	1.097	1.287	-17.3	103	0.03

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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