Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP070620\
Data File : BP002646.D
Acq On : 06 Jul 2020 16:17
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Jul 06 16:47:15 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 29 16:51:57 2020
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	274#	0.00
2		1,4-Dioxane	0.506	0.465	8.1	260#	0.00
3		Pyridine	1.501	1.344	10.5	248#	0.00
4		n-Nitrosodimethylamine	0.637	0.533	16.3	231#	0.00
5	S	2-Fluorophenol	1.181	1.089	7.8	259#	0.00
6		Aniline	2.071	1.886	8.9	253#	0.00
7	S	Phenol-d6	1.649	1.486	9.9	251#	0.00
8		2-Chlorophenol	1.312	1.219	7.1	261#	0.00
9		Benzaldehyde	0.857	0.810	5.5	263#	0.00
10	C	Phenol	1.663	1.520	8.6	256#	0.00
11		bis(2-Chloroethyl)ether	1.365	1.275	6.6	264#	0.00
12		1,3-Dichlorobenzene	1.514	1.446	4.5	270#	0.00
13	C	1,4-Dichlorobenzene	1.543	1.450	6.0	269#	0.00
14		1,2-Dichlorobenzene	1.485	1.356	8.7	261#	0.00
15		Benzyl Alcohol	1.239	1.029	16.9	226#	0.00
16		2,2'-oxybis(1-Chloropropane	1.901	1.724	9.3	256#	0.00
17		2-Methylphenol	1.244	1.074	13.7	239#	0.00
18		Hexachloroethane	0.557	0.516	7.4	261#	0.00
19	P	n-Nitroso-di-n-propylamine	1.082	0.843	22.1	218#	0.00
20		3+4-Methylphenols	1.678	1.414	15.7	230#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	240#	0.00
22		Acetophenone	0.506	0.463	8.5	225#	0.00
23	S	Nitrobenzene-d5	0.388	0.357	8.0	224#	0.00
24		Nitrobenzene	0.411	0.384	6.6	230#	0.00
25		Isophorone	0.778	0.687	11.7	217#	0.00
26	C	2-Nitrophenol	0.187	0.187	0.0	239#	0.00
27		2,4-Dimethylphenol	0.284	0.273	3.9	236#	0.00
28		bis(2-Chloroethoxy)methane	0.447	0.428	4.3	237#	0.00
29	C	2,4-Dichlorophenol	0.327	0.321	1.8	239#	0.00
30		1,2,4-Trichlorobenzene	0.368	0.389	-5.7	261#	0.00
31		Naphthalene	1.059	1.009	4.7	238#	0.00
32		Benzoic acid	0.190	0.156	17.9	185#	0.00
33		4-Chloroaniline	0.463	0.427	7.8	224#	0.00
34	C	Hexachlorobutadiene	0.225	0.240	-6.7	264#	0.00
35		Caprolactam	0.110	0.089	19.1	198#	0.00
36	C	4-Chloro-3-methylphenol	0.359	0.299	16.7	203#	0.00
37		2-Methylnaphthalene	0.772	0.701	9.2	225#	0.00
38		1-Methylnaphthalene	0.727	0.648	10.9	221#	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	216#	0.00
40		1,2,4,5-Tetrachlorobenzene	0.625	0.674	-7.8	242#	0.00
41	P	Hexachlorocyclopentadiene	0.263	0.378	-43.7#	278#	0.00
42	S	2,4,6-Tribromophenol	0.243	0.267	-9.9	245#	0.00
43	C	2,4,6-Trichlorophenol	0.423	0.444	-5.0	227#	0.00
44		2,4,5-Trichlorophenol	0.486	0.486	0.0	213#	0.00

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Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.359	1.240	8.8	206#	0.00
46		1,1'-Biphenyl	1.513	1.430	5.5	214#	0.00
47		2-Chloronaphthalene	1.231	1.188	3.5	217#	0.00
48		2-Nitroaniline	0.393	0.338	14.0	186#	0.00
49		Acenaphthylene	1.938	1.796	7.3	205#	0.00
50		Dimethylphthalate	1.571	1.395	11.2	199#	0.00
51		2,6-Dinitrotoluene	0.339	0.316	6.8	207#	0.00
52	C	Acenaphthene	1.143	1.064	6.9	210#	0.00
53		3-Nitroaniline	0.369	0.335	9.2	197#	0.00
54	P	2,4-Dinitrophenol	0.179	0.158	11.7	181#	0.00
55		Dibenzofuran	1.853	1.699	8.3	207#	0.00
56	P	4-Nitrophenol	0.258	0.229	11.2	179#	0.01
57		2,4-Dinitrotoluene	0.456	0.416	8.8	197#	0.00
58		Fluorene	1.411	1.225	13.2	198#	0.00
59		2,3,4,6-Tetrachlorophenol	0.393	0.386	1.8	216#	0.00
60		Diethylphthalate	1.562	1.298	16.9	186#	0.00
61		4-Chlorophenyl-phenylether	0.757	0.697	7.9	210#	0.00
62		4-Nitroaniline	0.369	0.323	12.5	186#	0.00
63		Azobenzene	1.510	1.276	15.5	189#	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	197#	0.00
65		4,6-Dinitro-2-methylphenol	0.115	0.120	-4.3	186#	0.00
66	c	n-Nitrosodiphenylamine	0.584	0.583	0.2	204#	0.00
67		4-Bromophenyl-phenylether	0.225	0.249	-10.7	224#	0.00
68		Hexachlorobenzene	0.241	0.264	-9.5	223#	0.00
69		Atrazine	0.187	0.174	7.0	181#	0.00
70	C	Pentachlorophenol	0.120	0.134	-11.7	201#	0.00
71		Phenanthrene	1.076	1.021	5.1	192#	0.00
72		Anthracene	1.071	1.040	2.9	196#	0.00
73		Carbazole	1.035	0.964	6.9	188#	0.00
74		Di-n-butylphthalate	1.219	1.099	9.8	178#	0.00
75	C	Fluoranthene	1.303	1.203	7.7	186#	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	202#	0.00
77		Benzidine	0.511	0.559	-9.4	196#	0.00
78		Pyrene	1.361	1.275	6.3	192#	0.00
79	S	Terphenyl-d14	0.927	0.826	10.9	189#	0.00
80		Butylbenzylphthalate	0.594	0.518	12.8	176#	0.00
81		Benzo(a)anthracene	1.308	1.253	4.2	200#	0.00
82		3,3'-Dichlorobenzidine	0.479	0.487	-1.7	202#	0.00
83		Chrysene	1.254	1.183	5.7	198#	0.00
84		Bis(2-ethylhexyl)phthalate	0.853	0.718	15.8	172#	0.00
85	c	Di-n-octyl phthalate	1.507	1.329	11.8	179#	0.00
86	I	Perylene-d12	1.000	1.000	0.0	231#	0.00
87		Indeno(1,2,3-cd)pyrene	1.440	1.491	-3.5	239#	0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.276	1.187	7.0	213#	0.00
89	Benzo(k)fluoranthene	1.191	1.139	4.4	225#	0.00
90 C	Benzo(a)pyrene	1.171	1.135	3.1	225#	0.00
91	Dibenzo(a,h)anthracene	1.177	1.207	-2.5	236#	0.01
92	Benzo(g,h,i)perylene	1.177	1.275	-8.3	248#	0.00

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