Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF082018\
Data File : BF108290.D
Acq On : 20 Aug 2018 15:19
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 20 16:46:09 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 08 12:24:24 2018
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	95	0.01
2		1,4-Dioxane	0.568	0.495	12.9	82	0.05
3		Pyridine	1.462	1.320	9.7	84	0.04
4		n-Nitrosodimethylamine	0.823	0.690	16.2	78	0.04
5	S	2-Fluorophenol	1.105	1.050	5.0	90	0.01
6		Aniline	1.997	1.929	3.4	93	0.02
7	S	Phenol-d6	1.468	1.401	4.6	91	0.01
8		2-Chlorophenol	1.244	1.209	2.8	91	0.01
9		Benzaldehyde	1.138	1.009	11.3	83	0.02
10	C	Phenol	1.518	1.464	3.6	92	0.02
11		bis(2-Chloroethyl)ether	1.228	1.157	5.8	89	0.02
12		1,3-Dichlorobenzene	1.439	1.383	3.9	91	0.01
13	C	1,4-Dichlorobenzene	1.445	1.383	4.3	90	0.01
14		1,2-Dichlorobenzene	1.344	1.275	5.1	91	0.01
15		Benzyl Alcohol	1.236	1.059	14.3	79	0.02
16		2,2'-oxybis(1-Chloropropane	2.529	2.151	14.9	80	0.01
17		2-Methylphenol	1.067	1.078	-1.0	93	0.01
18		Hexachloroethane	0.515	0.502	2.5	90	0.01
19	P	n-Nitroso-di-n-propylamine	0.993	0.896	9.8	85	0.02
20		3+4-Methylphenols	1.367	1.276	6.7	88	0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	93	0.02
22		Acetophenone	0.468	0.437	6.6	87	0.02
23	S	Nitrobenzene-d5	0.358	0.349	2.5	89	0.02
24		Nitrobenzene	0.362	0.347	4.1	87	0.02
25		Isophorone	0.683	0.649	5.0	88	0.02
26	C	2-Nitrophenol	0.137	0.155	-13.1	102	0.01
27		2,4-Dimethylphenol	0.303	0.292	3.6	89	0.02
28		bis(2-Chloroethoxy)methane	0.399	0.383	4.0	89	0.01
29	C	2,4-Dichlorophenol	0.279	0.279	0.0	90	0.02
30		1,2,4-Trichlorobenzene	0.308	0.299	2.9	91	0.01
31		Naphthalene	0.910	0.871	4.3	89	0.01
32		Benzoic acid	0.161	0.135	16.1	75	0.04
33		4-Chloroaniline	0.396	0.382	3.5	88	0.01
34	C	Hexachlorobutadiene	0.195	0.191	2.1	91	0.01
35		Caprolactam	0.087	0.087	0.0	88	0.02
36	C	4-Chloro-3-methylphenol	0.301	0.291	3.3	88	0.02
37		2-Methylnaphthalene	0.644	0.610	5.3	88	0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	90	0.01
39		1,2,4,5-Tetrachlorobenzene	0.614	0.609	0.8	89	0.01
40	P	Hexachlorocyclopentadiene	0.397	0.339	14.6	72	0.02
41	S	2,4,6-Tribromophenol	0.199	0.215	-8.0	91	0.02
42	C	2,4,6-Trichlorophenol	0.381	0.397	-4.2	90	0.02
43		2,4,5-Trichlorophenol	0.420	0.447	-6.4	91	0.02
44	S	2-Fluorobiphenyl	1.246	1.201	3.6	89	0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.475	1.432	2.9	88	0.02
46		2-Chloronaphthalene	1.165	1.138	2.3	88	0.02
47		2-Nitroaniline	0.377	0.390	-3.4	88	0.02
48		Acenaphthylene	1.843	1.783	3.3	88	0.01
49		Dimethylphthalate	1.393	1.351	3.0	88	0.02
50		2,6-Dinitrotoluene	0.291	0.301	-3.4	89	0.02
51	C	Acenaphthene	1.129	1.111	1.6	89	0.01
52		3-Nitroaniline	0.319	0.326	-2.2	89	0.02
53	P	2,4-Dinitrophenol	0.092	0.110	-19.6	105	0.02
54		Dibenzofuran	1.635	1.565	4.3	87	0.01
55	P	4-Nitrophenol	0.241	0.238	1.2	86	0.02
56		2,4-Dinitrotoluene	0.375	0.403	-7.5	90	0.02
57		Fluorene	1.294	1.255	3.0	89	0.02
58		2,3,4,6-Tetrachlorophenol	0.351	0.378	-7.7	91	0.02
59		Diethylphthalate	1.349	1.278	5.3	85	0.02
60		4-Chlorophenyl-phenylether	0.674	0.645	4.3	86	0.01
61		4-Nitroaniline	0.315	0.334	-6.0	91	0.02
62		Azobenzene	1.393	1.298	6.8	83	0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	88	0.01
64		4,6-Dinitro-2-methylphenol	0.069	0.080	-15.9	96	0.02
65	c	n-Nitrosodiphenylamine	0.591	0.580	1.9	86	0.02
66		4-Bromophenyl-phenylether	0.217	0.218	-0.5	88	0.01
67		Hexachlorobenzene	0.229	0.227	0.9	87	0.01
68		Atrazine	0.198	0.198	0.0	86	0.02
69	C	Pentachlorophenol	0.109	0.109	0.0	84	0.01
70		Phenanthrene	0.943	0.908	3.7	87	0.02
71		Anthracene	0.967	0.942	2.6	86	0.02
72		Carbazole	0.859	0.832	3.1	86	0.02
73		Di-n-butylphthalate	0.973	0.959	1.4	86	0.02
74	C	Fluoranthene	1.030	0.993	3.6	86	0.02
75	I	Chrysene-d12	1.000	1.000	0.0	80	0.02
76		Benzidine	0.747	0.772	-3.3	79	0.02
77		Pyrene	1.266	1.327	-4.8	85	0.02
78	S	Terphenyl-d14	0.787	0.819	-4.1	86	0.02
79		Butylbenzylphthalate	0.503	0.560	-11.3	84	0.01
80		Benzo(a)anthracene	1.148	1.146	0.2	80	0.02
81		3,3'-Dichlorobenzidine	0.411	0.392	4.6	72	0.02
82		Chrysene	1.052	1.051	0.1	81	0.02
83		Bis(2-ethylhexyl)phthalate	0.681	0.687	-0.9	78	0.02
84	c	Di-n-octyl phthalate	1.038	1.054	-1.5	76	0.02
85		Indeno(1,2,3-cd)pyrene	0.912	0.809	11.3	73	0.04
86	I	Perylene-d12	1.000	1.000	0.0	72	0.02
87		Benzo(b)fluoranthene	1.195	1.169	2.2	67	0.02

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88	Benzo(k)fluoranthene	1.099	1.136	-3.4	76	0.02
89 C	Benzo(a)pyrene	1.070	1.058	1.1	70	0.03
90	Dibenzo(a,h)anthracene	0.885	0.906	-2.4	73	0.04
91	Benzo(g,h,i)perylene	0.872	0.896	-2.8	75	0.05

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