Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080718\
Data File : BG036183.D
Acq On : 8 Aug 2018 5:32
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 09 09:08:41 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 31 12:30:06 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	85	-0.02
2		1,4-Dioxane	0.613	0.523	14.7	78	0.00
3		Pyridine	1.601	1.393	13.0	71	0.00
4		n-Nitrosodimethylamine	0.687	0.573	16.6	71	0.00
5	S	2-Fluorophenol	1.205	1.109	8.0	77	-0.01
6		Aniline	2.330	2.049	12.1	74	-0.02
7	S	Phenol-d6	1.797	1.686	6.2	79	-0.01
8		2-Chlorophenol	1.225	1.155	5.7	79	-0.02
9		Benzaldehyde	1.103	1.002	9.2	75	-0.01
10	C	Phenol	1.816	1.736	4.4	80	0.00
11		bis(2-Chloroethyl)ether	1.422	1.240	12.8	74	-0.02
12		1,3-Dichlorobenzene	1.480	1.450	2.0	84	-0.02
13	C	1,4-Dichlorobenzene	1.503	1.468	2.3	83	-0.02
14		1,2-Dichlorobenzene	1.444	1.389	3.8	82	-0.02
15		Benzyl Alcohol	1.632	1.411	13.5	73	-0.01
16		2,2'-oxybis(1-Chloropropane	1.192	1.128	5.4	80	-0.02
17		2-Methylphenol	1.235	1.112	10.0	75	-0.01
18		Hexachloroethane	0.575	0.552	4.0	83	-0.02
19	P	n-Nitroso-di-n-propylamine	1.513	1.239	18.1	71	-0.02
20		3+4-Methylphenols	1.713	1.578	7.9	74	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	79	-0.02
22		Acetophenone	0.622	0.565	9.2	74	-0.02
23	S	Nitrobenzene-d5	0.479	0.448	6.5	75	-0.02
24		Nitrobenzene	0.481	0.440	8.5	74	-0.02
25		Isophorone	0.889	0.761	14.4	69	-0.02
26	C	2-Nitrophenol	0.171	0.179	-4.7	82	-0.02
27		2,4-Dimethylphenol	0.289	0.269	6.9	74	-0.01
28		bis(2-Chloroethoxy)methane	0.490	0.431	12.0	71	-0.02
29	C	2,4-Dichlorophenol	0.330	0.339	-2.7	79	-0.02
30		1,2,4-Trichlorobenzene	0.405	0.417	-3.0	84	-0.02
31		Naphthalene	1.042	0.960	7.9	76	-0.02
32		Benzoic acid	0.195	0.161	17.4	66	0.00
33		4-Chloroaniline	0.454	0.407	10.4	73	-0.02
34	C	Hexachlorobutadiene	0.316	0.342	-8.2	88	-0.02
35		Caprolactam	0.142	0.119	16.2	70	-0.02
36	C	4-Chloro-3-methylphenol	0.443	0.423	4.5	75	-0.01
37		2-Methylnaphthalene	0.776	0.739	4.8	77	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	76	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.755	0.766	-1.5	83	-0.02
40	P	Hexachlorocyclopentadiene	0.396	0.360	9.1	71	-0.02
41	S	2,4,6-Tribromophenol	0.264	0.296	-12.1	89	-0.02
42	C	2,4,6-Trichlorophenol	0.441	0.459	-4.1	79	-0.02
43		2,4,5-Trichlorophenol	0.494	0.514	-4.0	85	-0.01
44	S	2-Fluorobiphenyl	1.493	1.481	0.8	81	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.551	1.495	3.6	77	-0.02
46		2-Chloronaphthalene	1.206	1.185	1.7	80	-0.02
47		2-Nitroaniline	0.426	0.404	5.2	75	-0.01
48		Acenaphthylene	2.020	1.940	4.0	77	-0.02
49		Dimethylphthalate	1.752	1.668	4.8	78	-0.01
50		2,6-Dinitrotoluene	0.357	0.388	-8.7	85	-0.02
51	C	Acenaphthene	1.259	1.198	4.8	78	-0.02
52		3-Nitroaniline	0.365	0.365	0.0	78	-0.01
53	P	2,4-Dinitrophenol	0.158	0.190	-20.3	96	0.00
54		Dibenzofuran	2.002	1.949	2.6	79	-0.02
55	P	4-Nitrophenol	0.260	0.231	11.2	69	0.00
56		2,4-Dinitrotoluene	0.512	0.557	-8.8	83	-0.01
57		Fluorene	1.678	1.675	0.2	82	-0.02
58		2,3,4,6-Tetrachlorophenol	0.473	0.491	-3.8	82	-0.02
59		Diethylphthalate	1.802	1.701	5.6	77	-0.01
60		4-Chlorophenyl-phenylether	0.986	0.977	0.9	81	-0.02
61		4-Nitroaniline	0.382	0.358	6.3	73	-0.02
62		Azobenzene	1.777	1.604	9.7	73	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	80	-0.02
64		4,6-Dinitro-2-methylphenol	0.111	0.124	-11.7	93	-0.01
65	c	n-Nitrosodiphenylamine	0.569	0.549	3.5	80	-0.01
66		4-Bromophenyl-phenylether	0.232	0.241	-3.9	85	-0.02
67		Hexachlorobenzene	0.239	0.243	-1.7	85	-0.01
68		Atrazine	0.234	0.197	15.8	67	-0.01
69	C	Pentachlorophenol	0.146	0.123	15.8	68	-0.01
70		Phenanthrene	0.998	0.965	3.3	82	-0.02
71		Anthracene	0.994	0.971	2.3	82	-0.02
72		Carbazole	0.944	0.894	5.3	79	-0.01
73		Di-n-butylphthalate	1.115	1.061	4.8	79	-0.02
74	C	Fluoranthene	1.344	1.301	3.2	81	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	83	-0.02
76		Benzidine	0.526	0.470	10.6	81	-0.02
77		Pyrene	1.265	1.188	6.1	82	-0.02
78	S	Terphenyl-d14	0.907	0.885	2.4	85	-0.02
79		Butylbenzylphthalate	0.452	0.451	0.2	84	-0.01
80		Benzo(a)anthracene	1.246	1.181	5.2	83	-0.02
81		3,3'-Dichlorobenzidine	0.435	0.422	3.0	82	-0.02
82		Chrysene	1.155	1.109	4.0	84	-0.02
83		Bis(2-ethylhexyl)phthalate	0.615	0.612	0.5	84	-0.02
84	c	Di-n-octyl phthalate	1.029	1.034	-0.5	84	-0.03
85		Indeno(1,2,3-cd)pyrene	1.279	1.237	3.3	83	-0.04
86	I	Perylene-d12	1.000	1.000	0.0	85	-0.04
87		Benzo(b)fluoranthene	1.216	1.156	4.9	85	-0.03

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88	Benzo(k)fluoranthene	1.188	1.132	4.7	82	-0.02
89 C	Benzo(a)pyrene	1.131	1.086	4.0	83	-0.03
90	Dibenzo(a,h)anthracene	1.110	1.076	3.1	84	-0.05
91	Benzo(g,h,i)perylene	1.086	1.051	3.2	83	-0.05

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