Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG080818\
Data File : BG036196.D
Acq On : 8 Aug 2018 13:58
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 09 14:50:43 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	81	-0.02
2		1,4-Dioxane	0.613	0.542	11.6	77	-0.01
3		Pyridine	1.601	1.350	15.7	66	-0.01
4		n-Nitrosodimethylamine	0.687	0.596	13.2	71	-0.01
5	S	2-Fluorophenol	1.205	1.132	6.1	75	-0.01
6		Aniline	2.330	2.038	12.5	71	-0.02
7	S	Phenol-d6	1.797	1.680	6.5	75	-0.02
8		2-Chlorophenol	1.225	1.211	1.1	80	-0.02
9		Benzaldehyde	1.103	1.009	8.5	73	-0.01
10	C	Phenol	1.816	1.698	6.5	75	-0.01
11		bis(2-Chloroethyl)ether	1.422	1.259	11.5	72	-0.02
12		1,3-Dichlorobenzene	1.480	1.441	2.6	80	-0.02
13	C	1,4-Dichlorobenzene	1.503	1.502	0.1	81	-0.02
14		1,2-Dichlorobenzene	1.444	1.403	2.8	80	-0.02
15		Benzyl Alcohol	1.632	1.394	14.6	69	-0.01
16		2,2'-oxybis(1-Chloropropane	1.192	0.958	19.6	65	-0.02
17		2-Methylphenol	1.235	1.155	6.5	74	-0.01
18		Hexachloroethane	0.575	0.573	0.3	83	-0.02
19	P	n-Nitroso-di-n-propylamine	1.513	1.207	20.2	66	-0.02
20		3+4-Methylphenols	1.713	1.571	8.3	71	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	73	-0.02
22		Acetophenone	0.622	0.580	6.8	71	-0.01
23	S	Nitrobenzene-d5	0.479	0.467	2.5	72	-0.02
24		Nitrobenzene	0.481	0.463	3.7	72	-0.02
25		Isophorone	0.889	0.764	14.1	64	-0.02
26	C	2-Nitrophenol	0.171	0.185	-8.2	78	-0.02
27		2,4-Dimethylphenol	0.289	0.275	4.8	70	-0.02
28		bis(2-Chloroethoxy)methane	0.490	0.440	10.2	67	-0.02
29	C	2,4-Dichlorophenol	0.330	0.344	-4.2	75	-0.01
30		1,2,4-Trichlorobenzene	0.405	0.424	-4.7	79	-0.02
31		Naphthalene	1.042	0.994	4.6	73	-0.02
32		Benzoic acid	0.195	0.159	18.5	60	0.00
33		4-Chloroaniline	0.454	0.421	7.3	70	-0.02
34	C	Hexachlorobutadiene	0.316	0.350	-10.8	84	-0.02
35		Caprolactam	0.142	0.132	7.0	72	-0.02
36	C	4-Chloro-3-methylphenol	0.443	0.451	-1.8	74	-0.02
37		2-Methylnaphthalene	0.776	0.750	3.4	72	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	74	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.755	0.757	-0.3	80	-0.02
40	P	Hexachlorocyclopentadiene	0.396	0.339	14.4	66	-0.02
41	S	2,4,6-Tribromophenol	0.264	0.304	-15.2	90	-0.02
42	C	2,4,6-Trichlorophenol	0.441	0.446	-1.1	75	-0.02
43		2,4,5-Trichlorophenol	0.494	0.514	-4.0	83	-0.01
44	S	2-Fluorobiphenyl	1.493	1.417	5.1	76	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.551	1.455	6.2	74	-0.02
46		2-Chloronaphthalene	1.206	1.159	3.9	77	-0.02
47		2-Nitroaniline	0.426	0.428	-0.5	77	-0.01
48		Acenaphthylene	2.020	1.929	4.5	75	-0.02
49		Dimethylphthalate	1.752	1.664	5.0	76	-0.01
50		2,6-Dinitrotoluene	0.357	0.371	-3.9	80	-0.02
51	C	Acenaphthene	1.259	1.188	5.6	75	-0.02
52		3-Nitroaniline	0.365	0.367	-0.5	76	-0.02
53	P	2,4-Dinitrophenol	0.158	0.183	-15.8	90	0.00
54		Dibenzofuran	2.002	1.976	1.3	78	-0.02
55	P	4-Nitrophenol	0.260	0.291	-11.9	85	0.00
56		2,4-Dinitrotoluene	0.512	0.575	-12.3	83	-0.01
57		Fluorene	1.678	1.678	0.0	80	-0.02
58		2,3,4,6-Tetrachlorophenol	0.473	0.502	-6.1	82	-0.02
59		Diethylphthalate	1.802	1.765	2.1	78	-0.01
60		4-Chlorophenyl-phenylether	0.986	1.002	-1.6	81	-0.02
61		4-Nitroaniline	0.382	0.388	-1.6	77	-0.02
62		Azobenzene	1.777	1.640	7.7	73	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	83	-0.01
64		4,6-Dinitro-2-methylphenol	0.111	0.116	-4.5	90	-0.01
65	c	n-Nitrosodiphenylamine	0.569	0.543	4.6	82	-0.01
66		4-Bromophenyl-phenylether	0.232	0.230	0.9	84	-0.02
67		Hexachlorobenzene	0.239	0.238	0.4	86	-0.01
68		Atrazine	0.234	0.198	15.4	70	-0.01
69	C	Pentachlorophenol	0.146	0.123	15.8	70	-0.01
70		Phenanthrene	0.998	0.962	3.6	84	-0.02
71		Anthracene	0.994	0.955	3.9	83	-0.02
72		Carbazole	0.944	0.902	4.4	82	-0.01
73		Di-n-butylphthalate	1.115	1.049	5.9	80	-0.02
74	C	Fluoranthene	1.344	1.324	1.5	85	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	88	-0.01
76		Benzidine	0.526	0.469	10.8	85	-0.01
77		Pyrene	1.265	1.186	6.2	87	-0.02
78	S	Terphenyl-d14	0.907	0.879	3.1	89	-0.02
79		Butylbenzylphthalate	0.452	0.434	4.0	85	-0.01
80		Benzo(a)anthracene	1.246	1.174	5.8	87	-0.02
81		3,3'-Dichlorobenzidine	0.435	0.413	5.1	85	-0.02
82		Chrysene	1.155	1.082	6.3	87	-0.02
83		Bis(2-ethylhexyl)phthalate	0.615	0.611	0.7	88	-0.02
84	c	Di-n-octyl phthalate	1.029	1.019	1.0	87	-0.02
85		Indeno(1,2,3-cd)pyrene	1.279	1.231	3.8	87	-0.04
86	I	Perylene-d12	1.000	1.000	0.0	90	-0.03
87		Benzo(b)fluoranthene	1.216	1.148	5.6	89	-0.02

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88	Benzo(k)fluoranthene	1.188	1.164	2.0	90	-0.02
89 C	Benzo(a)pyrene	1.131	1.075	5.0	87	-0.02
90	Dibenzo(a,h)anthracene	1.110	1.047	5.7	86	-0.05
91	Benzo(g,h,i)perylene	1.086	1.046	3.7	88	-0.04

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