Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG061418\
Data File : BG035156.D
Acq On : 15 Jun 2018 3:35
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 15 07:24:01 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG060718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 08 17:16:28 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	194#	-0.01
2		1,4-Dioxane	0.565	0.525	7.1	181#	0.00
3		Pyridine	1.526	1.616	-5.9	192#	0.00
4		n-Nitrosodimethylamine	0.655	0.623	4.9	173#	-0.01
5	S	2-Fluorophenol	1.185	1.219	-2.9	195#	0.00
6		Aniline	2.261	2.384	-5.4	201#	-0.01
7	S	Phenol-d6	1.749	1.926	-10.1	207#	0.00
8		2-Chlorophenol	1.242	1.327	-6.8	201#	0.00
9		Benzaldehyde	1.347	1.292	4.1	176#	0.00
10	C	Phenol	1.810	1.961	-8.3	206#	0.00
11		bis(2-Chloroethyl)ether	1.405	1.481	-5.4	201#	-0.01
12		1,3-Dichlorobenzene	1.482	1.551	-4.7	205#	0.00
13	C	1,4-Dichlorobenzene	1.530	1.563	-2.2	196#	0.00
14		1,2-Dichlorobenzene	1.437	1.507	-4.9	199#	-0.01
15		Benzyl Alcohol	1.657	1.744	-5.3	199#	-0.01
16		2,2'-oxybis(1-Chloropropane	1.398	1.121	19.8	155#	-0.01
17		2-Methylphenol	1.193	1.319	-10.6	204#	0.00
18		Hexachloroethane	0.548	0.571	-4.2	194#	-0.01
19	P	n-Nitroso-di-n-propylamine	1.486	1.499	-0.9	191#	0.00
20		3+4-Methylphenols	1.711	1.935	-13.1	214#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	204#	-0.01
22		Acetophenone	0.620	0.617	0.5	201#	-0.01
23	S	Nitrobenzene-d5	0.421	0.477	-13.3	219#	-0.01
24		Nitrobenzene	0.431	0.475	-10.2	215#	-0.01
25		Isophorone	0.888	0.863	2.8	196#	-0.01
26	C	2-Nitrophenol	0.149	0.198	-32.9#	269#	-0.01
27		2,4-Dimethylphenol	0.312	0.313	-0.3	205#	-0.01
28		bis(2-Chloroethoxy)methane	0.477	0.487	-2.1	205#	-0.01
29	C	2,4-Dichlorophenol	0.340	0.374	-10.0	218#	0.00
30		1,2,4-Trichlorobenzene	0.407	0.436	-7.1	210#	-0.01
31		Naphthalene	1.039	1.062	-2.2	207#	0.00
32		Benzoic acid	0.209	0.242	-15.8	234#	0.02
33		4-Chloroaniline	0.451	0.480	-6.4	212#	0.00
34	C	Hexachlorobutadiene	0.317	0.324	-2.2	207#	-0.01
35		Caprolactam	0.126	0.140	-11.1	227#	0.01
36	C	4-Chloro-3-methylphenol	0.423	0.471	-11.3	222#	0.00
37		2-Methylnaphthalene	0.788	0.828	-5.1	209#	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	222#	0.00
39		1,2,4,5-Tetrachlorobenzene	0.745	0.769	-3.2	222#	0.00
40	P	Hexachlorocyclopentadiene	0.467	0.439	6.0	197#	-0.02
41	S	2,4,6-Tribromophenol	0.256	0.291	-13.7	246#	0.00
42	C	2,4,6-Trichlorophenol	0.446	0.497	-11.4	236#	0.00
43		2,4,5-Trichlorophenol	0.490	0.537	-9.6	240#	0.00
44	S	2-Fluorobiphenyl	1.538	1.472	4.3	210#	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.608	1.572	2.2	212#	0.00
46		2-Chloronaphthalene	1.216	1.213	0.2	221#	-0.01
47		2-Nitroaniline	0.359	0.436	-21.4	265#	0.00
48		Acenaphthylene	2.039	2.014	1.2	216#	0.00
49		Dimethylphthalate	1.744	1.718	1.5	217#	-0.01
50		2,6-Dinitrotoluene	0.318	0.381	-19.8	255#	0.00
51	C	Acenaphthene	1.332	1.352	-1.5	220#	0.00
52		3-Nitroaniline	0.312	0.381	-22.1	252#	0.00
53	P	2,4-Dinitrophenol	0.129	0.139	-7.8	228#	0.00
54		Dibenzofuran	1.995	1.981	0.7	216#	0.00
55	P	4-Nitrophenol	0.263	0.274	-4.2	214#	0.00
56		2,4-Dinitrotoluene	0.423	0.546	-29.1#	275#	0.00
57		Fluorene	1.667	1.644	1.4	214#	0.00
58		2,3,4,6-Tetrachlorophenol	0.476	0.528	-10.9	239#	0.00
59		Diethylphthalate	1.779	1.740	2.2	216#	-0.01
60		4-Chlorophenyl-phenylether	0.951	0.987	-3.8	229#	-0.02
61		4-Nitroaniline	0.335	0.395	-17.9	245#	0.00
62		Azobenzene	1.744	1.635	6.2	206#	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	223#	0.00
64		4,6-Dinitro-2-methylphenol	0.091	0.116	-27.5#	278#	0.00
65	c	n-Nitrosodiphenylamine	0.601	0.607	-1.0	218#	0.00
66		4-Bromophenyl-phenylether	0.241	0.255	-5.8	234#	-0.02
67		Hexachlorobenzene	0.245	0.259	-5.7	234#	0.00
68		Atrazine	0.256	0.266	-3.9	225#	-0.01
69	C	Pentachlorophenol	0.165	0.178	-7.9	231#	-0.01
70		Phenanthrene	1.016	1.020	-0.4	219#	-0.01
71		Anthracene	1.018	1.019	-0.1	216#	0.00
72		Carbazole	0.944	0.945	-0.1	216#	0.00
73		Di-n-butylphthalate	1.125	1.097	2.5	211#	-0.02
74	C	Fluoranthene	1.322	1.317	0.4	215#	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	226#	-0.01
76		Benzidine	0.744	0.614	17.5	182#	-0.01
77		Pyrene	1.318	1.258	4.6	214#	0.00
78	S	Terphenyl-d14	0.955	0.895	6.3	208#	-0.01
79		Butylbenzylphthalate	0.479	0.487	-1.7	219#	-0.02
80		Benzo(a)anthracene	1.278	1.269	0.7	223#	-0.01
81		3,3'-Dichlorobenzidine	0.481	0.472	1.9	213#	-0.01
82		Chrysene	1.170	1.176	-0.5	225#	-0.01
83		Bis(2-ethylhexyl)phthalate	0.676	0.657	2.8	213#	-0.02
84	c	Di-n-octyl phthalate	1.161	1.145	1.4	215#	-0.04
85		Indeno(1,2,3-cd)pyrene	1.370	1.327	3.1	214#	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	220#	-0.02
87		Benzo(b)fluoranthene	1.232	1.252	-1.6	217#	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.205	1.243	-3.2	231#	-0.01
89 C	Benzo(a)pyrene	1.159	1.196	-3.2	224#	-0.02
90	Dibenzo(a,h)anthracene	1.144	1.092	4.5	209#	-0.03
91	Benzo(g,h,i)perylene	1.120	1.082	3.4	211#	-0.03

(#) = Out of Range SPCC's out = 0 CCC's out = 1