Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG060319\

Evaluate Continuing Calibration Report

Data File : BG041036.D
Acq On : 3 Jun 2019 16:15
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 03 17:46:58 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 29 18:39:57 2019
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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	100	-0.01
2		1,4-Dioxane	40.000	40.443	-1.1	101	-0.01
3		Pyridine	40.000	41.670	-4.2	103	0.00
4		n-Nitrosodimethylamine	40.000	44.361	-10.9	109	-0.01
5	S	2-Fluorophenol	80.000	82.015	-2.5	101	0.00
6		Aniline	40.000	39.832	0.4	100	-0.01
7	S	Phenol-d6	80.000	78.354	2.1	98	0.00
8		2-Chlorophenol	40.000	40.926	-2.3	103	-0.01
9		Benzaldehyde	40.000	41.486	-3.7	103	-0.01
10	C	Phenol	40.000	40.094	-0.2	100	-0.01
11		bis(2-Chloroethyl)ether	40.000	40.913	-2.3	101	0.00
12		1,3-Dichlorobenzene	40.000	41.994	-5.0	104	0.00
13	C	1,4-Dichlorobenzene	40.000	41.971	-4.9	103	-0.01
14		1,2-Dichlorobenzene	40.000	40.968	-2.4	101	-0.01
15		Benzyl Alcohol	40.000	38.670	3.3	97	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	40.361	-0.9	101	-0.02
17		2-Methylphenol	40.000	37.405	6.5	93	-0.01
18		Hexachloroethane	40.000	42.642	-6.6	107	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	39.120	2.2	98	-0.01
20		3+4-Methylphenols	40.000	37.785	5.5	96	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	100	-0.01
22		Acetophenone	40.000	39.401	1.5	97	-0.01
23	S	Nitrobenzene-d5	80.000	80.519	-0.6	100	0.00
24		Nitrobenzene	40.000	39.704	0.7	97	0.00
25		Isophorone	40.000	38.357	4.1	94	-0.01
26	C	2-Nitrophenol	40.000	42.032	-5.1	102	-0.01
27		2,4-Dimethylphenol	40.000	38.133	4.7	94	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.747	3.1	95	-0.01
29	C	2,4-Dichlorophenol	40.000	37.622	5.9	92	0.00
30		1,2,4-Trichlorobenzene	40.000	40.064	-0.2	98	-0.01
31		Naphthalene	40.000	39.204	2.0	96	-0.01
32		Benzoic acid	40.000	38.214	4.5	96	0.00
33		4-Chloroaniline	40.000	37.769	5.6	93	-0.01
34	C	Hexachlorobutadiene	40.000	40.803	-2.0	103	0.00
35		Caprolactam	40.000	36.391	9.0	89	-0.01
36	C	4-Chloro-3-methylphenol	40.000	36.381	9.0	90	-0.01
37		2-Methylnaphthalene	40.000	38.098	4.8	94	-0.01
38		1-Methylnaphthalene	40.000	37.715	5.7	92	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	91	-0.01
40		1,2,4,5-Tetrachlorobenzene	40.000	42.293	-5.7	94	-0.01
41	P	Hexachlorocyclopentadiene	40.000	39.867	0.3	95	-0.01
42	S	2,4,6-Tribromophenol	80.000	80.572	-0.7	90	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.674	-1.7	89	-0.01
44		2,4,5-Trichlorophenol	40.000	40.427	-1.1	90	-0.01

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Quant Time: Jun 03 17:46:58 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	80.000	84.684	-5.9	93	-0.01
46		1,1'-Biphenyl	40.000	41.721	-4.3	91	-0.01
47		2-Chloronaphthalene	40.000	42.044	-5.1	93	0.00
48		2-Nitroaniline	40.000	42.420	-6.1	94	0.00
49		Acenaphthylene	40.000	40.960	-2.4	89	-0.01
50		Dimethylphthalate	40.000	40.399	-1.0	88	0.00
51		2,6-Dinitrotoluene	40.000	40.524	-1.3	90	0.00
52	C	Acenaphthene	40.000	40.283	-0.7	89	0.00
53		3-Nitroaniline	40.000	41.755	-4.4	90	-0.01
54	P	2,4-Dinitrophenol	40.000	40.409	-1.0	94	-0.01
55		Dibenzofuran	40.000	40.750	-1.9	88	0.00
56	P	4-Nitrophenol	40.000	42.078	-5.2	91	0.00
57		2,4-Dinitrotoluene	40.000	41.303	-3.3	90	0.00
58		Fluorene	40.000	39.767	0.6	88	-0.01
59		2,3,4,6-Tetrachlorophenol	40.000	40.559	-1.4	89	-0.01
60		Diethylphthalate	40.000	40.156	-0.4	88	-0.01
61		4-Chlorophenyl-phenylether	40.000	39.932	0.2	89	0.00
62		4-Nitroaniline	40.000	41.644	-4.1	92	-0.01
63		Azobenzene	40.000	40.200	-0.5	87	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	90	0.00
65		4,6-Dinitro-2-methylphenol	40.000	39.059	2.4	92	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.976	0.1	88	0.00
67		4-Bromophenyl-phenylether	40.000	39.439	1.4	86	-0.01
68		Hexachlorobenzene	40.000	39.091	2.3	87	-0.01
69		Atrazine	40.000	42.898	-7.2	87	-0.01
70	C	Pentachlorophenol	40.000	37.988	5.0	83	-0.01
71		Phenanthrene	40.000	41.153	-2.9	90	-0.01
72		Anthracene	40.000	41.730	-4.3	91	0.00
73		Carbazole	40.000	42.119	-5.3	92	-0.01
74		Di-n-butylphthalate	40.000	41.351	-3.4	88	-0.01
75	C	Fluoranthene	40.000	42.433	-6.1	92	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	93	-0.01
77		Benzidine	40.000	40.437	-1.1	89	-0.01
78		Pyrene	40.000	41.856	-4.6	93	-0.01
79	S	Terphenyl-d14	80.000	83.935	-4.9	92	-0.01
80		Butylbenzylphthalate	40.000	41.249	-3.1	92	0.00
81		Benzo(a)anthracene	40.000	40.864	-2.2	93	0.00
82		3,3'-Dichlorobenzidine	40.000	40.445	-1.1	93	-0.01
83		Chrysene	40.000	41.110	-2.8	92	-0.01
84		Bis(2-ethylhexyl)phthalate	40.000	39.958	0.1	90	-0.01
85	c	Di-n-octyl phthalate	40.000	40.120	-0.3	90	-0.01
86		Indeno(1,2,3-cd)pyrene	40.000	38.282	4.3	89	-0.03
87	I	Perylene-d12	20.000	20.000	0.0	94	-0.02

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	40.000	40.171	-0.4	92	-0.01
89	Benzo(k)fluoranthene	40.000	39.927	0.2	91	-0.02
90 C	Benzo(a)pyrene	40.000	39.971	0.1	91	-0.01
91	Dibenzo(a,h)anthracene	40.000	38.106	4.7	86	-0.02
92	Benzo(g,h,i)perylene	40.000	38.467	3.8	89	-0.02

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