Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG072518\
Data File : BG035875.D
Acq On : 25 Jul 2018 15:52
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 25 17:20:53 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 25 11:04:14 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	103	0.00
2		1,4-Dioxane	40.000	32.833	17.9	91	0.00
3		Pyridine	40.000	35.449	11.4	88	0.00
4		n-Nitrosodimethylamine	40.000	32.699	18.3	85	0.00
5	S	2-Fluorophenol	80.000	71.621	10.5	91	0.00
6		Aniline	40.000	37.403	6.5	96	0.00
7	S	Phenol-d6	80.000	75.521	5.6	96	0.00
8		2-Chlorophenol	40.000	39.140	2.1	100	0.00
9		Benzaldehyde	40.000	33.463	16.3	84	0.00
10	C	Phenol	40.000	39.366	1.6	99	0.00
11		bis(2-Chloroethyl)ether	40.000	38.879	2.8	100	0.00
12		1,3-Dichlorobenzene	40.000	38.213	4.5	99	0.00
13	C	1,4-Dichlorobenzene	40.000	38.805	3.0	100	0.00
14		1,2-Dichlorobenzene	40.000	38.844	2.9	101	0.00
15		Benzyl Alcohol	40.000	38.853	2.9	99	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.029	7.4	95	0.00
17		2-Methylphenol	40.000	38.543	3.6	97	0.00
18		Hexachloroethane	40.000	38.313	4.2	101	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.920	2.7	102	0.00
20		3+4-Methylphenols	40.000	41.403	-3.5	101	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	106	0.00
22		Acetophenone	40.000	37.774	5.6	103	0.00
23	S	Nitrobenzene-d5	80.000	73.809	7.7	99	0.00
24		Nitrobenzene	40.000	35.932	10.2	98	0.00
25		Isophorone	40.000	38.007	5.0	103	0.00
26	C	2-Nitrophenol	40.000	43.694	-9.2	114	0.00
27		2,4-Dimethylphenol	40.000	39.747	0.6	106	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.510	3.7	104	0.00
29	C	2,4-Dichlorophenol	40.000	42.173	-5.4	109	0.00
30		1,2,4-Trichlorobenzene	40.000	40.134	-0.3	109	0.00
31		Naphthalene	40.000	39.049	2.4	109	0.00
32		Benzoic acid	40.000	33.259	16.9	89	0.00
33		4-Chloroaniline	40.000	39.044	2.4	107	0.00
34	C	Hexachlorobutadiene	40.000	38.940	2.7	106	0.00
35		Caprolactam	40.000	38.133	4.7	107	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.661	-1.7	108	0.00
37		2-Methylnaphthalene	40.000	41.110	-2.8	111	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	111	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.288	4.3	114	0.00
40	P	Hexachlorocyclopentadiene	40.000	37.892	5.3	109	0.00
41	S	2,4,6-Tribromophenol	80.000	82.154	-2.7	119	0.00
42	C	2,4,6-Trichlorophenol	40.000	41.209	-3.0	115	0.00
43		2,4,5-Trichlorophenol	40.000	40.381	-1.0	121	0.00
44	S	2-Fluorobiphenyl	80.000	77.201	3.5	115	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG072518\
Data File : BG035875.D
Acq On : 25 Jul 2018 15:52
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 25 17:20:53 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 25 11:04:14 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	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	38.771	3.1	114	0.00
46		2-Chloronaphthalene	40.000	38.379	4.1	114	0.00
47		2-Nitroaniline	40.000	38.346	4.1	110	0.00
48		Acenaphthylene	40.000	38.878	2.8	114	0.00
49		Dimethylphthalate	40.000	38.452	3.9	116	0.00
50		2,6-Dinitrotoluene	40.000	42.253	-5.6	121	0.00
51	C	Acenaphthene	40.000	38.753	3.1	116	0.00
52		3-Nitroaniline	40.000	38.623	3.4	110	0.00
53	P	2,4-Dinitrophenol	40.000	34.381	14.0	102	0.00
54		Dibenzofuran	40.000	38.906	2.7	115	0.00
55	P	4-Nitrophenol	40.000	37.428	6.4	107	0.00
56		2,4-Dinitrotoluene	40.000	42.168	-5.4	117	0.00
57		Fluorene	40.000	38.847	2.9	117	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	40.702	-1.8	118	0.00
59		Diethylphthalate	40.000	37.546	6.1	111	0.00
60		4-Chlorophenyl-phenylether	40.000	39.239	1.9	118	0.00
61		4-Nitroaniline	40.000	37.531	6.2	106	0.00
62		Azobenzene	40.000	36.376	9.1	108	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	111	0.00
64		4,6-Dinitro-2-methylphenol	40.000	41.124	-2.8	119	0.00
65	c	n-Nitrosodiphenylamine	40.000	39.912	0.2	115	0.00
66		4-Bromophenyl-phenylether	40.000	41.085	-2.7	117	0.00
67		Hexachlorobenzene	40.000	40.452	-1.1	117	0.00
68		Atrazine	40.000	36.925	7.7	103	0.00
69	C	Pentachlorophenol	40.000	41.866	-4.7	116	0.00
70		Phenanthrene	40.000	38.527	3.7	112	0.00
71		Anthracene	40.000	38.708	3.2	112	0.00
72		Carbazole	40.000	36.815	8.0	106	0.00
73		Di-n-butylphthalate	40.000	36.646	8.4	105	0.00
74	C	Fluoranthene	40.000	36.264	9.3	105	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	106	0.00
76		Benzidine	40.000	38.592	3.5	111	0.00
77		Pyrene	40.000	37.900	5.3	106	0.00
78	S	Terphenyl-d14	80.000	75.607	5.5	105	0.00
79		Butylbenzylphthalate	40.000	39.262	1.8	106	0.00
80		Benzo(a)anthracene	40.000	37.597	6.0	105	0.00
81		3,3'-Dichlorobenzidine	40.000	39.159	2.1	106	0.00
82		Chrysene	40.000	38.136	4.7	107	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	40.151	-0.4	108	0.00
84	c	Di-n-octyl phthalate	40.000	41.512	-3.8	111	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	39.297	1.8	108	0.00
86	I	Perylene-d12	20.000	20.000	0.0	107	0.00
87		Benzo(b)fluoranthene	40.000	38.649	3.4	109	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG072518\
Data File : BG035875.D
Acq On : 25 Jul 2018 15:52
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 25 17:20:53 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 25 11:04:14 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	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	38.917	2.7	106	0.00
89 C	Benzo(a)pyrene	40.000	38.982	2.5	106	0.00
90	Dibenzo(a,h)anthracene	40.000	39.564	1.1	108	0.00
91	Benzo(g,h,i)perylene	40.000	39.260	1.9	107	0.00

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