Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG080618\
Data File : BG036146.D
Acq On : 7 Aug 2018 3:57
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 07 04:39:26 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	88	-0.01
2		1,4-Dioxane	40.000	35.209	12.0	84	-0.01
3		Pyridine	40.000	35.568	11.1	76	-0.01
4		n-Nitrosodimethylamine	40.000	34.323	14.2	77	-0.01
5	S	2-Fluorophenol	80.000	79.090	1.1	87	-0.01
6		Aniline	40.000	36.609	8.5	81	-0.02
7	S	Phenol-d6	80.000	78.106	2.4	86	-0.01
8		2-Chlorophenol	40.000	39.703	0.7	87	-0.02
9		Benzaldehyde	40.000	37.243	6.9	81	-0.01
10	C	Phenol	40.000	40.456	-1.1	88	-0.01
11		bis(2-Chloroethyl)ether	40.000	38.212	4.5	84	-0.02
12		1,3-Dichlorobenzene	40.000	40.799	-2.0	91	-0.02
13	C	1,4-Dichlorobenzene	40.000	41.017	-2.5	91	-0.02
14		1,2-Dichlorobenzene	40.000	41.033	-2.6	92	-0.02
15		Benzyl Alcohol	40.000	37.233	6.9	82	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	33.064	17.3	73	-0.02
17		2-Methylphenol	40.000	38.033	4.9	82	-0.01
18		Hexachloroethane	40.000	39.034	2.4	88	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	35.845	10.4	81	-0.02
20		3+4-Methylphenols	40.000	39.195	2.0	83	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	86	-0.02
22		Acetophenone	40.000	37.286	6.8	83	-0.01
23	S	Nitrobenzene-d5	80.000	75.535	5.6	82	-0.01
24		Nitrobenzene	40.000	37.169	7.1	82	-0.01
25		Isophorone	40.000	35.432	11.4	78	-0.02
26	C	2-Nitrophenol	40.000	43.527	-8.8	92	-0.02
27		2,4-Dimethylphenol	40.000	37.947	5.1	82	-0.01
28		bis(2-Chloroethoxy)methane	40.000	36.245	9.4	79	-0.02
29	C	2,4-Dichlorophenol	40.000	42.907	-7.3	90	-0.01
30		1,2,4-Trichlorobenzene	40.000	41.778	-4.4	92	-0.02
31		Naphthalene	40.000	38.372	4.1	87	-0.02
32		Benzoic acid	40.000	33.309	16.7	72	0.00
33		4-Chloroaniline	40.000	38.232	4.4	85	-0.02
34	C	Hexachlorobutadiene	40.000	43.487	-8.7	96	-0.02
35		Caprolactam	40.000	34.695	13.3	79	-0.01
36	C	4-Chloro-3-methylphenol	40.000	39.576	1.1	85	-0.01
37		2-Methylnaphthalene	40.000	39.470	1.3	87	-0.02
38	I	Acenaphthene-d10	20.000	20.000	0.0	87	-0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	39.984	0.0	93	-0.02
40	P	Hexachlorocyclopentadiene	40.000	36.087	9.8	81	-0.01
41	S	2,4,6-Tribromophenol	80.000	90.009	-12.5	102	-0.02
42	C	2,4,6-Trichlorophenol	40.000	41.994	-5.0	91	-0.01
43		2,4,5-Trichlorophenol	40.000	41.279	-3.2	96	-0.01
44	S	2-Fluorobiphenyl	80.000	77.574	3.0	90	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG080618\
Data File : BG036146.D
Acq On : 7 Aug 2018 3:57
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 07 04:39:26 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	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	38.028	4.9	87	-0.02
46		2-Chloronaphthalene	40.000	38.032	4.9	88	-0.01
47		2-Nitroaniline	40.000	37.460	6.3	84	-0.01
48		Acenaphthylene	40.000	38.418	4.0	88	-0.02
49		Dimethylphthalate	40.000	38.712	3.2	91	-0.01
50		2,6-Dinitrotoluene	40.000	42.858	-7.1	96	-0.01
51	C	Acenaphthene	40.000	38.080	4.8	89	-0.01
52		3-Nitroaniline	40.000	39.309	1.7	87	-0.02
53	P	2,4-Dinitrophenol	40.000	40.545	-1.4	97	0.00
54		Dibenzofuran	40.000	38.462	3.8	89	-0.02
55	P	4-Nitrophenol	40.000	39.197	2.0	87	0.00
56		2,4-Dinitrotoluene	40.000	41.847	-4.6	91	-0.01
57		Fluorene	40.000	38.887	2.8	91	-0.01
58		2,3,4,6-Tetrachlorophenol	40.000	42.555	-6.4	96	-0.02
59		Diethylphthalate	40.000	38.687	3.3	89	-0.01
60		4-Chlorophenyl-phenylether	40.000	40.235	-0.6	94	-0.02
61		4-Nitroaniline	40.000	36.545	8.6	80	-0.01
62		Azobenzene	40.000	35.269	11.8	82	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	91	-0.01
64		4,6-Dinitro-2-methylphenol	40.000	43.130	-7.8	102	-0.01
65	c	n-Nitrosodiphenylamine	40.000	38.136	4.7	89	-0.01
66		4-Bromophenyl-phenylether	40.000	41.105	-2.8	95	-0.02
67		Hexachlorobenzene	40.000	40.537	-1.3	96	-0.01
68		Atrazine	40.000	34.151	14.6	77	-0.01
69	C	Pentachlorophenol	40.000	33.329	16.7	75	-0.01
70		Phenanthrene	40.000	38.354	4.1	91	-0.02
71		Anthracene	40.000	38.238	4.4	90	-0.01
72		Carbazole	40.000	36.636	8.4	86	-0.01
73		Di-n-butylphthalate	40.000	37.410	6.5	87	-0.02
74	C	Fluoranthene	40.000	36.833	7.9	87	-0.01
75	I	Chrysene-d12	20.000	20.000	0.0	86	-0.02
76		Benzidine	40.000	28.209	29.5#	66	-0.01
77		Pyrene	40.000	38.801	3.0	88	-0.02
78	S	Terphenyl-d14	80.000	79.803	0.2	90	-0.02
79		Butylbenzylphthalate	40.000	39.193	2.0	86	-0.01
80		Benzo(a)anthracene	40.000	37.943	5.1	86	-0.02
81		3,3'-Dichlorobenzidine	40.000	38.896	2.8	85	-0.02
82		Chrysene	40.000	38.061	4.8	86	-0.02
83		Bis(2-ethylhexyl)phthalate	40.000	39.814	0.5	87	-0.02
84	c	Di-n-octyl phthalate	40.000	40.396	-1.0	87	-0.02
85		Indeno(1,2,3-cd)pyrene	40.000	38.905	2.7	86	-0.05
86	I	Perylene-d12	20.000	20.000	0.0	88	-0.04
87		Benzo(b)fluoranthene	40.000	37.497	6.3	87	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG080618\
Data File : BG036146.D
Acq On : 7 Aug 2018 3:57
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 07 04:39:26 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	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	39.038	2.4	88	-0.02
89 C	Benzo(a)pyrene	40.000	37.901	5.2	85	-0.03
90	Dibenzo(a,h)anthracene	40.000	38.200	4.5	86	-0.06
91	Benzo(g,h,i)perylene	40.000	38.606	3.5	86	-0.05

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