Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG082318\
Data File : BG036360.D
Acq On : 23 Aug 2018 2:26
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG082318

Quant Time: Aug 23 12:28:28 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 23 12:07:02 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	115	0.00
2		1,4-Dioxane	40.000	35.413	11.5	106	0.00
3		Pyridine	40.000	38.007	5.0	107	0.00
4		n-Nitrosodimethylamine	40.000	37.886	5.3	109	0.00
5	S	2-Fluorophenol	80.000	76.113	4.9	109	0.00
6		Aniline	40.000	38.588	3.5	111	0.00
7	S	Phenol-d6	80.000	77.213	3.5	111	0.00
8		2-Chlorophenol	40.000	37.576	6.1	108	0.00
9		Benzaldehyde	40.000	36.100	9.7	111	0.00
10	C	Phenol	40.000	38.641	3.4	111	0.00
11		bis(2-Chloroethyl)ether	40.000	38.039	4.9	111	0.00
12		1,3-Dichlorobenzene	40.000	37.995	5.0	112	0.00
13	C	1,4-Dichlorobenzene	40.000	38.159	4.6	111	0.00
14		1,2-Dichlorobenzene	40.000	37.371	6.6	110	0.00
15		Benzyl Alcohol	40.000	39.789	0.5	113	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.495	6.3	110	0.00
17		2-Methylphenol	40.000	38.441	3.9	112	0.00
18		Hexachloroethane	40.000	38.307	4.2	110	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.888	2.8	113	0.00
20		3+4-Methylphenols	40.000	38.353	4.1	111	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	117	0.00
22		Acetophenone	40.000	37.248	6.9	109	0.00
23	S	Nitrobenzene-d5	80.000	79.948	0.1	114	0.00
24		Nitrobenzene	40.000	39.272	1.8	112	0.00
25		Isophorone	40.000	38.550	3.6	112	0.00
26	C	2-Nitrophenol	40.000	40.196	-0.5	117	0.00
27		2,4-Dimethylphenol	40.000	37.554	6.1	110	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.773	5.6	110	0.00
29	C	2,4-Dichlorophenol	40.000	39.458	1.4	113	0.00
30		1,2,4-Trichlorobenzene	40.000	38.094	4.8	113	0.00
31		Naphthalene	40.000	38.007	5.0	111	0.00
32		Benzoic acid	40.000	38.760	3.1	116	0.00
33		4-Chloroaniline	40.000	38.208	4.5	111	0.00
34	C	Hexachlorobutadiene	40.000	38.656	3.4	112	0.00
35		Caprolactam	40.000	38.816	3.0	109	0.00
36	C	4-Chloro-3-methylphenol	40.000	38.772	3.1	111	0.00
37		2-Methylnaphthalene	40.000	38.518	3.7	112	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	114	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.966	2.6	111	0.00
40	P	Hexachlorocyclopentadiene	40.000	39.853	0.4	112	0.00
41	S	2,4,6-Tribromophenol	80.000	78.711	1.6	106	0.00
42	C	2,4,6-Trichlorophenol	40.000	39.678	0.8	112	0.00
43		2,4,5-Trichlorophenol	40.000	40.046	-0.1	111	0.00
44	S	2-Fluorobiphenyl	80.000	76.708	4.1	111	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG082318\
Data File : BG036360.D
Acq On : 23 Aug 2018 2:26
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG082318

Quant Time: Aug 23 12:28:28 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 23 12:07:02 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	38.065	4.8	111	0.00
46		2-Chloronaphthalene	40.000	38.470	3.8	111	0.00
47		2-Nitroaniline	40.000	41.293	-3.2	112	0.00
48		Acenaphthylene	40.000	38.238	4.4	110	0.00
49		Dimethylphthalate	40.000	38.614	3.5	110	0.00
50		2,6-Dinitrotoluene	40.000	38.782	3.0	109	0.00
51	C	Acenaphthene	40.000	37.768	5.6	108	0.00
52		3-Nitroaniline	40.000	39.808	0.5	105	0.00
53	P	2,4-Dinitrophenol	40.000	39.696	0.8	112	0.00
54		Dibenzofuran	40.000	38.115	4.7	110	0.00
55	P	4-Nitrophenol	40.000	37.322	6.7	101	0.00
56		2,4-Dinitrotoluene	40.000	40.148	-0.4	112	0.00
57		Fluorene	40.000	37.352	6.6	106	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.620	1.0	110	0.00
59		Diethylphthalate	40.000	37.813	5.5	107	0.00
60		4-Chlorophenyl-phenylether	40.000	37.322	6.7	108	0.00
61		4-Nitroaniline	40.000	39.116	2.2	104	0.00
62		Azobenzene	40.000	37.389	6.5	107	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	108	0.00
64		4,6-Dinitro-2-methylphenol	40.000	40.297	-0.7	109	0.00
65	c	n-Nitrosodiphenylamine	40.000	38.360	4.1	106	0.00
66		4-Bromophenyl-phenylether	40.000	39.384	1.5	107	0.00
67		Hexachlorobenzene	40.000	38.965	2.6	106	0.00
68		Atrazine	40.000	36.084	9.8	100	0.00
69	C	Pentachlorophenol	40.000	41.529	-3.8	105	0.00
70		Phenanthrene	40.000	38.402	4.0	105	0.00
71		Anthracene	40.000	37.963	5.1	103	0.00
72		Carbazole	40.000	37.368	6.6	100	0.00
73		Di-n-butylphthalate	40.000	38.059	4.9	101	0.00
74	C	Fluoranthene	40.000	37.611	6.0	101	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	106	0.00
76		Benzidine	40.000	34.903	12.7	99	0.00
77		Pyrene	40.000	37.881	5.3	100	0.00
78	S	Terphenyl-d14	80.000	75.911	5.1	102	0.00
79		Butylbenzylphthalate	40.000	39.694	0.8	102	0.00
80		Benzo(a)anthracene	40.000	37.717	5.7	102	0.00
81		3,3'-Dichlorobenzidine	40.000	38.302	4.2	99	0.00
82		Chrysene	40.000	37.604	6.0	100	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	39.180	2.1	102	0.00
84	c	Di-n-octyl phthalate	40.000	39.809	0.5	102	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	38.598	3.5	102	0.00
86	I	Perylene-d12	20.000	20.000	0.0	107	0.00
87		Benzo(b)fluoranthene	40.000	38.386	4.0	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG082318\
Data File : BG036360.D
Acq On : 23 Aug 2018 2:26
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG082318

Quant Time: Aug 23 12:28:28 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 23 12:07:02 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	37.915	5.2	102	0.00
89 C	Benzo(a)pyrene	40.000	38.026	4.9	101	0.00
90	Dibenzo(a,h)anthracene	40.000	38.000	5.0	101	0.00
91	Benzo(g,h,i)perylene	40.000	38.306	4.2	102	0.00

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