Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083018\
Data File : BF108639.D
Acq On : 30 Aug 2018 14:37
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF083018

Quant Time: Aug 30 15:12:42 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 30 15:06:55 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	100	0.00
2		1,4-Dioxane	40.000	35.075	12.3	90	0.00
3		Pyridine	40.000	34.050	14.9	86	0.00
4		n-Nitrosodimethylamine	40.000	34.565	13.6	85	0.00
5	S	2-Fluorophenol	80.000	68.355	14.6	86	0.00
6		Aniline	40.000	39.052	2.4	94	0.00
7	S	Phenol-d6	80.000	75.100	6.1	94	0.00
8		2-Chlorophenol	40.000	38.715	3.2	94	0.00
9		Benzaldehyde	40.000	38.476	3.8	96	0.00
10	C	Phenol	40.000	38.395	4.0	95	0.00
11		bis(2-Chloroethyl)ether	40.000	37.627	5.9	93	0.00
12		1,3-Dichlorobenzene	40.000	38.256	4.4	94	0.00
13	C	1,4-Dichlorobenzene	40.000	38.749	3.1	96	0.00
14		1,2-Dichlorobenzene	40.000	37.337	6.7	95	0.00
15		Benzyl Alcohol	40.000	38.450	3.9	93	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.026	4.9	94	0.00
17		2-Methylphenol	40.000	40.186	-0.5	95	0.00
18		Hexachloroethane	40.000	38.193	4.5	96	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.122	2.2	94	0.00
20		3+4-Methylphenols	40.000	42.013	-5.0	96	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	100	0.00
22		Acetophenone	40.000	38.672	3.3	97	0.00
23	S	Nitrobenzene-d5	80.000	76.302	4.6	95	0.00
24		Nitrobenzene	40.000	39.345	1.6	96	0.00
25		Isophorone	40.000	38.118	4.7	95	0.00
26	C	2-Nitrophenol	40.000	39.473	1.3	96	0.00
27		2,4-Dimethylphenol	40.000	37.387	6.5	95	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.004	5.0	95	0.00
29	C	2,4-Dichlorophenol	40.000	38.601	3.5	96	0.00
30		1,2,4-Trichlorobenzene	40.000	38.674	3.3	97	0.00
31		Naphthalene	40.000	37.638	5.9	94	0.00
32		Benzoic acid	40.000	33.432	16.4	83	0.00
33		4-Chloroaniline	40.000	38.685	3.3	97	0.00
34	C	Hexachlorobutadiene	40.000	38.168	4.6	96	0.00
35		Caprolactam	40.000	39.383	1.5	96	0.00
36	C	4-Chloro-3-methylphenol	40.000	38.501	3.7	95	0.00
37		2-Methylnaphthalene	40.000	38.562	3.6	96	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	102	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.012	5.0	96	0.00
40	P	Hexachlorocyclopentadiene	40.000	37.380	6.5	96	0.00
41	S	2,4,6-Tribromophenol	80.000	76.707	4.1	98	0.00
42	C	2,4,6-Trichlorophenol	40.000	38.954	2.6	95	0.00
43		2,4,5-Trichlorophenol	40.000	39.945	0.1	103	0.00
44	S	2-Fluorobiphenyl	80.000	75.584	5.5	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083018\
Data File : BF108639.D
Acq On : 30 Aug 2018 14:37
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF083018

Quant Time: Aug 30 15:12:42 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 30 15:06:55 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	37.743	5.6	97	0.00
46		2-Chloronaphthalene	40.000	37.794	5.5	97	0.00
47		2-Nitroaniline	40.000	38.021	4.9	95	0.00
48		Acenaphthylene	40.000	37.557	6.1	96	0.00
49		Dimethylphthalate	40.000	37.766	5.6	97	0.00
50		2,6-Dinitrotoluene	40.000	38.630	3.4	97	0.00
51	C	Acenaphthene	40.000	37.395	6.5	98	0.00
52		3-Nitroaniline	40.000	38.138	4.7	96	0.00
53	P	2,4-Dinitrophenol	40.000	39.328	1.7	97	0.00
54		Dibenzofuran	40.000	37.600	6.0	97	0.00
55	P	4-Nitrophenol	40.000	40.007	-0.0	97	0.04
56		2,4-Dinitrotoluene	40.000	38.591	3.5	98	0.00
57		Fluorene	40.000	37.339	6.7	96	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	38.683	3.3	97	0.00
59		Diethylphthalate	40.000	37.948	5.1	98	0.00
60		4-Chlorophenyl-phenylether	40.000	37.716	5.7	97	0.00
61		4-Nitroaniline	40.000	38.241	4.4	98	0.00
62		Azobenzene	40.000	37.181	7.0	97	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	102	0.00
64		4,6-Dinitro-2-methylphenol	40.000	38.774	3.1	98	0.00
65	c	n-Nitrosodiphenylamine	40.000	37.083	7.3	96	0.00
66		4-Bromophenyl-phenylether	40.000	38.234	4.4	98	0.00
67		Hexachlorobenzene	40.000	37.579	6.1	98	0.00
68		Atrazine	40.000	36.866	7.8	97	0.00
69	C	Pentachlorophenol	40.000	39.886	0.3	95	0.00
70		Phenanthrene	40.000	37.399	6.5	97	0.00
71		Anthracene	40.000	37.529	6.2	97	0.00
72		Carbazole	40.000	37.733	5.7	98	0.00
73		Di-n-butylphthalate	40.000	37.253	6.9	97	0.00
74	C	Fluoranthene	40.000	37.246	6.9	97	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	104	0.00
76		Benzidine	40.000	42.629	-6.6	108	0.00
77		Pyrene	40.000	37.618	6.0	98	0.00
78	S	Terphenyl-d14	80.000	75.439	5.7	99	0.00
79		Butylbenzylphthalate	40.000	38.213	4.5	100	0.00
80		Benzo(a)anthracene	40.000	37.259	6.9	99	0.00
81		3,3'-Dichlorobenzidine	40.000	37.138	7.2	96	0.00
82		Chrysene	40.000	37.093	7.3	97	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	37.972	5.1	100	0.00
84	c	Di-n-octyl phthalate	40.000	37.860	5.4	100	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	34.616	13.5	94	0.00
86	I	Perylene-d12	20.000	20.000	0.0	103	0.00
87		Benzo(b)fluoranthene	40.000	36.511	8.7	91	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083018\
Data File : BF108639.D
Acq On : 30 Aug 2018 14:37
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF083018

Quant Time: Aug 30 15:12:42 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 30 15:06:55 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	40.139	-0.3	107	0.00
89 C	Benzo(a)pyrene	40.000	38.147	4.6	99	0.00
90	Dibenzo(a,h)anthracene	40.000	37.608	6.0	95	0.00
91	Benzo(g,h,i)perylene	40.000	37.515	6.2	93	0.00

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