Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF072018\
Data File : BF107526.D
Acq On : 20 Jul 2018 16:21
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF072018

Quant Time: Jul 20 16:43:17 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 20 16:29:27 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	106	0.00
2		1,4-Dioxane	40.000	36.970	7.6	98	0.00
3		Pyridine	40.000	38.167	4.6	100	0.00
4		n-Nitrosodimethylamine	40.000	36.628	8.4	99	0.00
5	S	2-Fluorophenol	80.000	76.726	4.1	100	0.00
6		Aniline	40.000	36.923	7.7	98	0.00
7	S	Phenol-d6	80.000	73.703	7.9	100	0.00
8		2-Chlorophenol	40.000	37.998	5.0	101	0.00
9		Benzaldehyde	40.000	36.357	9.1	99	0.00
10	C	Phenol	40.000	36.950	7.6	100	0.00
11		bis(2-Chloroethyl)ether	40.000	39.003	2.5	102	0.00
12		1,3-Dichlorobenzene	40.000	37.978	5.1	103	0.00
13	C	1,4-Dichlorobenzene	40.000	37.860	5.4	102	0.00
14		1,2-Dichlorobenzene	40.000	36.767	8.1	100	0.00
15		Benzyl Alcohol	40.000	39.345	1.6	102	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.068	7.3	100	0.00
17		2-Methylphenol	40.000	36.357	9.1	97	0.00
18		Hexachloroethane	40.000	38.950	2.6	101	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	36.250	9.4	101	0.00
20		3+4-Methylphenols	40.000	36.578	8.6	99	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	106	0.00
22		Acetophenone	40.000	35.803	10.5	101	0.00
23	S	Nitrobenzene-d5	80.000	82.211	-2.8	105	0.00
24		Nitrobenzene	40.000	40.640	-1.6	103	0.00
25		Isophorone	40.000	36.434	8.9	99	0.00
26	C	2-Nitrophenol	40.000	45.221	-13.1	113	0.00
27		2,4-Dimethylphenol	40.000	38.076	4.8	102	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.057	7.4	101	0.00
29	C	2,4-Dichlorophenol	40.000	39.275	1.8	101	0.00
30		1,2,4-Trichlorobenzene	40.000	36.950	7.6	100	0.00
31		Naphthalene	40.000	36.199	9.5	101	0.00
32		Benzoic acid	40.000	39.969	0.1	107	0.00
33		4-Chloroaniline	40.000	33.507	16.2	96	0.00
34	C	Hexachlorobutadiene	40.000	38.656	3.4	105	0.00
35		Caprolactam	40.000	37.337	6.7	97	0.00
36	C	4-Chloro-3-methylphenol	40.000	37.783	5.5	100	0.00
37		2-Methylnaphthalene	40.000	36.359	9.1	101	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	106	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	37.797	5.5	103	0.00
40	P	Hexachlorocyclopentadiene	40.000	41.696	-4.2	105	0.00
41	S	2,4,6-Tribromophenol	80.000	80.989	-1.2	103	0.00
42	C	2,4,6-Trichlorophenol	40.000	39.382	1.5	100	0.00
43		2,4,5-Trichlorophenol	40.000	40.247	-0.6	102	0.00
44	S	2-Fluorobiphenyl	80.000	75.532	5.6	101	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF072018\
Data File : BF107526.D
Acq On : 20 Jul 2018 16:21
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF072018

Quant Time: Jul 20 16:43:17 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 20 16:29:27 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	36.346	9.1	101	0.00
46		2-Chloronaphthalene	40.000	36.717	8.2	101	0.00
47		2-Nitroaniline	40.000	43.289	-8.2	108	0.00
48		Acenaphthylene	40.000	36.442	8.9	101	0.00
49		Dimethylphthalate	40.000	36.256	9.4	100	0.00
50		2,6-Dinitrotoluene	40.000	41.455	-3.6	106	0.00
51	C	Acenaphthene	40.000	37.948	5.1	102	0.00
52		3-Nitroaniline	40.000	40.737	-1.8	105	0.00
53	P	2,4-Dinitrophenol	40.000	45.264	-13.2	130	0.00
54		Dibenzofuran	40.000	36.074	9.8	101	0.00
55	P	4-Nitrophenol	40.000	41.219	-3.0	103	0.00
56		2,4-Dinitrotoluene	40.000	42.931	-7.3	107	0.00
57		Fluorene	40.000	37.511	6.2	101	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.908	0.2	102	0.00
59		Diethylphthalate	40.000	36.280	9.3	100	0.00
60		4-Chlorophenyl-phenylether	40.000	36.470	8.8	102	0.00
61		4-Nitroaniline	40.000	39.286	1.8	98	0.00
62		Azobenzene	40.000	37.291	6.8	100	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	104	0.00
64		4,6-Dinitro-2-methylphenol	40.000	43.227	-8.1	120	0.00
65	c	n-Nitrosodiphenylamine	40.000	36.491	8.8	99	0.00
66		4-Bromophenyl-phenylether	40.000	37.598	6.0	101	0.00
67		Hexachlorobenzene	40.000	37.826	5.4	102	0.00
68		Atrazine	40.000	38.555	3.6	101	0.00
69	C	Pentachlorophenol	40.000	41.057	-2.6	104	0.00
70		Phenanthrene	40.000	35.499	11.3	98	0.00
71		Anthracene	40.000	35.674	10.8	99	0.00
72		Carbazole	40.000	35.942	10.1	100	0.00
73		Di-n-butylphthalate	40.000	38.525	3.7	102	0.00
74	C	Fluoranthene	40.000	36.081	9.8	100	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	106	0.00
76		Benzidine	40.000	29.762	25.6#	82	0.00
77		Pyrene	40.000	35.589	11.0	100	0.00
78	S	Terphenyl-d14	80.000	73.465	8.2	103	0.00
79		Butylbenzylphthalate	40.000	41.892	-4.7	105	0.00
80		Benzo(a)anthracene	40.000	37.090	7.3	101	0.00
81		3,3'-Dichlorobenzidine	40.000	38.645	3.4	105	0.00
82		Chrysene	40.000	36.593	8.5	103	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	39.314	1.7	107	0.00
84	c	Di-n-octyl phthalate	40.000	40.800	-2.0	105	0.01
85		Indeno(1,2,3-cd)pyrene	40.000	40.201	-0.5	118	0.02
86	I	Perylene-d12	20.000	20.000	0.0	119	0.02
87		Benzo(b)fluoranthene	40.000	36.223	9.4	108	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF072018\
Data File : BF107526.D
Acq On : 20 Jul 2018 16:21
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF072018

Quant Time: Jul 20 16:43:17 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 20 16:29:27 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	37.176	7.1	114	0.01
89 C	Benzo(a)pyrene	40.000	37.147	7.1	112	0.02
90	Dibenzo(a,h)anthracene	40.000	37.619	6.0	118	0.02
91	Benzo(g,h,i)perylene	40.000	36.598	8.5	115	0.02

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