Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060321\
Data File : BF124126.D
Acq On : 3 Jun 2021 17:47
Operator : JU/CG
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF060321

Quant Time: Jun 04 09:59:43 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 04 09:56:24 2021
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	112	0.00
2		1,4-Dioxane	40.000	36.321	9.2	104	0.00
3		Pyridine	40.000	36.311	9.2	104	0.00
4		n-Nitrosodimethylamine	40.000	35.639	10.9	102	0.00
5	S	2-Fluorophenol	80.000	72.115	9.9	104	0.00
6		Aniline	40.000	36.077	9.8	107	0.00
7	S	Phenol-d6	80.000	71.310	10.9	104	0.00
8		2-Chlorophenol	40.000	37.378	6.6	106	0.00
9		Benzaldehyde	40.000	34.961	12.6	106	0.00
10	C	Phenol	40.000	35.044	12.4	103	0.00
11		bis(2-Chloroethyl)ether	40.000	34.695	13.3	99	0.00
12		1,3-Dichlorobenzene	40.000	35.678	10.8	103	0.00
13	C	1,4-Dichlorobenzene	40.000	36.718	8.2	106	0.00
14		1,2-Dichlorobenzene	40.000	35.717	10.7	103	0.00
15		Benzyl Alcohol	40.000	36.983	7.5	109	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	35.386	11.5	105	0.00
17		2-Methylphenol	40.000	33.610	16.0	98	0.00
18		Hexachloroethane	40.000	36.201	9.5	105	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	36.238	9.4	110	0.00
20		3+4-Methylphenols	40.000	35.443	11.4	103	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	109	0.00
22		Acetophenone	40.000	37.282	6.8	104	0.00
23	S	Nitrobenzene-d5	80.000	73.622	8.0	105	0.00
24		Nitrobenzene	40.000	36.585	8.5	101	0.00
25		Isophorone	40.000	36.418	9.0	104	0.00
26	C	2-Nitrophenol	40.000	36.548	8.6	103	0.00
27		2,4-Dimethylphenol	40.000	36.416	9.0	103	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.569	6.1	111	0.00
29	C	2,4-Dichlorophenol	40.000	37.440	6.4	106	0.00
30		1,2,4-Trichlorobenzene	40.000	37.446	6.4	108	0.00
31		Naphthalene	40.000	36.701	8.2	104	0.00
32		Benzoic acid	40.000	36.060	9.8	108	0.00
33		4-Chloroaniline	40.000	37.303	6.7	106	0.00
34	C	Hexachlorobutadiene	40.000	36.939	7.7	104	0.00
35		Caprolactam	40.000	36.410	9.0	105	0.00
36	C	4-Chloro-3-methylphenol	40.000	36.477	8.8	109	0.00
37		2-Methylnaphthalene	40.000	37.021	7.4	104	0.00
38		1-Methylnaphthalene	40.000	36.983	7.5	104	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	103	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	38.066	4.8	107	0.00
41	P	Hexachlorocyclopentadiene	40.000	36.412	9.0	103	0.00
42	S	2,4,6-Tribromophenol	80.000	76.460	4.4	105	0.00
43	C	2,4,6-Trichlorophenol	40.000	38.387	4.0	103	0.00
44		2,4,5-Trichlorophenol	40.000	38.314	4.2	105	0.00
45	S	2-Fluorobiphenyl	80.000	75.798	5.3	104	0.00
46		1,1'-Biphenyl	40.000	37.613	6.0	105	0.00
47		2-Chloronaphthalene	40.000	36.818	8.0	104	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060321\
Data File : BF124126.D
Acq On : 3 Jun 2021 17:47
Operator : JU/CG
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF060321

Quant Time: Jun 04 09:59:43 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 04 09:56:24 2021
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)
48		2-Nitroaniline	40.000	38.555	3.6	106	0.00
49		Acenaphthylene	40.000	38.310	4.2	106	0.00
50		Dimethylphthalate	40.000	38.016	5.0	109	0.00
51		2,6-Dinitrotoluene	40.000	37.794	5.5	109	0.00
52	C	Acenaphthene	40.000	37.786	5.5	105	0.00
53		3-Nitroaniline	40.000	36.532	8.7	103	0.00
54	P	2,4-Dinitrophenol	40.000	40.341	-0.9	116	0.00
55		Dibenzofuran	40.000	38.011	5.0	106	0.00
56	P	4-Nitrophenol	40.000	39.997	0.0	106	0.00
57		2,4-Dinitrotoluene	40.000	38.078	4.8	101	0.00
58		Fluorene	40.000	38.199	4.5	104	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	39.380	1.5	112	0.00
60		Diethylphthalate	40.000	37.961	5.1	105	0.00
61		4-Chlorophenyl-phenylether	40.000	38.724	3.2	109	0.00
62		4-Nitroaniline	40.000	37.885	5.3	103	0.00
63		Azobenzene	40.000	37.497	6.3	107	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	104	0.00
65		4,6-Dinitro-2-methylphenol	40.000	39.060	2.3	109	0.00
66	c	n-Nitrosodiphenylamine	40.000	37.134	7.2	104	0.00
67		4-Bromophenyl-phenylether	40.000	37.016	7.5	108	0.00
68		Hexachlorobenzene	40.000	36.993	7.5	105	0.00
69		Atrazine	40.000	39.997	0.0	102	0.00
70	C	Pentachlorophenol	40.000	37.973	5.1	108	0.00
71		Phenanthrene	40.000	37.198	7.0	104	0.00
72		Anthracene	40.000	37.523	6.2	108	0.00
73		Carbazole	40.000	37.181	7.0	105	0.00
74		Di-n-butylphthalate	40.000	37.283	6.8	106	0.00
75	C	Fluoranthene	40.000	37.160	7.1	104	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	106	0.00
77		Benzidine	40.000	39.338	1.7	97	0.00
78		Pyrene	40.000	37.458	6.4	105	0.00
79	S	Terphenyl-d14	80.000	76.280	4.6	106	0.00
80		Butylbenzylphthalate	40.000	38.601	3.5	108	0.00
81		Benzo(a)anthracene	40.000	37.592	6.0	107	0.00
82		3,3'-Dichlorobenzidine	40.000	37.243	6.9	106	0.00
83		Chrysene	40.000	37.067	7.3	105	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.045	-2.6	114	0.00
85	c	Di-n-octyl phthalate	40.000	42.713	-6.8	122	0.00
86	I	Perylene-d12	20.000	20.000	0.0	111	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	34.324	14.2	104	0.00
88		Benzo(b)fluoranthene	40.000	36.002	10.0	112	0.00
89		Benzo(k)fluoranthene	40.000	37.670	5.8	114	0.00
90	C	Benzo(a)pyrene	40.000	36.342	9.1	110	0.00
91		Dibenzo(a,h)anthracene	40.000	34.141	14.6	105	0.00
92		Benzo(g,h,i)perylene	40.000	34.049	14.9	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060321\
Data File : BF124126.D
Acq On : 3 Jun 2021 17:47
Operator : JU/CG
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF060321

Quant Time: Jun 04 09:59:43 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 04 09:56:24 2021
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)

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