Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060419\
Data File : BF114795.D
Acq On : 4 Jun 2019 16:52
Operator : HP/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF060419

Quant Time: Jun 04 17:24:06 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF060419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 04 16:57:15 2019
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	90	0.00
2		1,4-Dioxane	40.000	39.094	2.3	88	0.00
3		Pyridine	40.000	40.043	-0.1	89	-0.01
4		n-Nitrosodimethylamine	40.000	38.839	2.9	87	-0.01
5	S	2-Fluorophenol	80.000	79.199	1.0	89	0.00
6		Aniline	40.000	39.993	0.0	88	0.00
7	S	Phenol-d6	80.000	80.275	-0.3	89	0.00
8		2-Chlorophenol	40.000	40.016	-0.0	87	0.00
9		Benzaldehyde	40.000	39.187	2.0	86	0.00
10	C	Phenol	40.000	40.684	-1.7	91	0.00
11		bis(2-Chloroethyl)ether	40.000	40.308	-0.8	87	0.00
12		1,3-Dichlorobenzene	40.000	40.221	-0.6	88	0.00
13	C	1,4-Dichlorobenzene	40.000	40.450	-1.1	88	0.00
14		1,2-Dichlorobenzene	40.000	41.188	-3.0	88	0.00
15		Benzyl Alcohol	40.000	41.119	-2.8	88	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.198	-0.5	86	0.00
17		2-Methylphenol	40.000	39.796	0.5	87	0.00
18		Hexachloroethane	40.000	39.598	1.0	86	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.286	1.8	86	0.00
20		3+4-Methylphenols	40.000	40.714	-1.8	90	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	89	0.00
22		Acetophenone	40.000	39.951	0.1	88	0.00
23	S	Nitrobenzene-d5	80.000	80.205	-0.3	87	0.00
24		Nitrobenzene	40.000	40.646	-1.6	88	0.00
25		Isophorone	40.000	38.982	2.5	86	0.00
26	C	2-Nitrophenol	40.000	40.707	-1.8	87	0.00
27		2,4-Dimethylphenol	40.000	39.971	0.1	87	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.949	0.1	86	0.00
29	C	2,4-Dichlorophenol	40.000	39.736	0.7	85	0.00
30		1,2,4-Trichlorobenzene	40.000	40.555	-1.4	87	0.00
31		Naphthalene	40.000	41.069	-2.7	90	0.00
32		Benzoic acid	40.000	43.982	-10.0	87	0.00
33		4-Chloroaniline	40.000	40.126	-0.3	87	0.00
34	C	Hexachlorobutadiene	40.000	40.510	-1.3	86	0.00
35		Caprolactam	40.000	38.904	2.7	87	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.685	0.8	87	0.00
37		2-Methylnaphthalene	40.000	40.727	-1.8	88	0.00
38		1-Methylnaphthalene	40.000	40.696	-1.7	88	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	89	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.695	-1.7	88	0.00
41	P	Hexachlorocyclopentadiene	40.000	40.263	-0.7	86	0.00
42	S	2,4,6-Tribromophenol	80.000	79.681	0.4	87	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.366	1.6	86	0.00
44		2,4,5-Trichlorophenol	40.000	39.739	0.7	86	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060419\
Data File : BF114795.D
Acq On : 4 Jun 2019 16:52
Operator : HP/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF060419

Quant Time: Jun 04 17:24:06 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF060419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 04 16:57:15 2019
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	S	2-Fluorobiphenyl	80.000	85.782	-7.2	88	0.00
46		1,1'-Biphenyl	40.000	40.979	-2.4	89	0.00
47		2-Chloronaphthalene	40.000	41.010	-2.5	88	0.00
48		2-Nitroaniline	40.000	39.761	0.6	88	0.00
49		Acenaphthylene	40.000	41.152	-2.9	90	0.00
50		Dimethylphthalate	40.000	40.430	-1.1	89	0.00
51		2,6-Dinitrotoluene	40.000	39.985	0.0	87	0.00
52	C	Acenaphthene	40.000	40.408	-1.0	88	0.00
53		3-Nitroaniline	40.000	40.086	-0.2	89	0.00
54	P	2,4-Dinitrophenol	40.000	41.413	-3.5	89	0.00
55		Dibenzofuran	40.000	39.251	1.9	89	0.00
56	P	4-Nitrophenol	40.000	39.448	1.4	87	0.00
57		2,4-Dinitrotoluene	40.000	41.392	-3.5	88	0.00
58		Fluorene	40.000	42.973	-7.4	90	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	39.800	0.5	86	0.00
60		Diethylphthalate	40.000	40.669	-1.7	89	0.00
61		4-Chlorophenyl-phenylether	40.000	37.466	6.3	89	0.00
62		4-Nitroaniline	40.000	39.156	2.1	86	0.00
63		Azobenzene	40.000	40.626	-1.6	88	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	90	0.00
65		4,6-Dinitro-2-methylphenol	40.000	41.443	-3.6	90	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.879	0.3	87	0.00
67		4-Bromophenyl-phenylether	40.000	39.709	0.7	86	0.00
68		Hexachlorobenzene	40.000	40.261	-0.7	88	0.00
69		Atrazine	40.000	39.441	1.4	86	0.00
70	C	Pentachlorophenol	40.000	40.376	-0.9	87	0.00
71		Phenanthrene	40.000	40.539	-1.3	91	0.00
72		Anthracene	40.000	38.924	2.7	89	0.00
73		Carbazole	40.000	39.714	0.7	89	0.00
74		Di-n-butylphthalate	40.000	39.534	1.2	90	0.00
75	C	Fluoranthene	40.000	38.539	3.7	91	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	96	0.00
77		Benzidine	40.000	34.684	13.3	81	0.00
78		Pyrene	40.000	37.817	5.5	88	0.00
79	S	Terphenyl-d14	80.000	75.709	5.4	89	0.00
80		Butylbenzylphthalate	40.000	37.775	5.6	89	0.00
81		Benzo(a)anthracene	40.000	38.663	3.3	91	0.00
82		3,3'-Dichlorobenzidine	40.000	37.063	7.3	87	0.00
83		Chrysene	40.000	37.723	5.7	91	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	39.058	2.4	89	0.00
85	c	Di-n-octyl phthalate	40.000	38.876	2.8	91	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	36.492	8.8	91	0.00
87	I	Perylene-d12	20.000	20.000	0.0	94	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060419\
Data File : BF114795.D
Acq On : 4 Jun 2019 16:52
Operator : HP/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF060419

Quant Time: Jun 04 17:24:06 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF060419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 04 16:57:15 2019
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(b)fluoranthene	40.000	40.558	-1.4	99	0.00
89	Benzo(k)fluoranthene	40.000	38.218	4.5	82	0.00
90 C	Benzo(a)pyrene	40.000	39.914	0.2	90	0.00
91	Dibenzo(a,h)anthracene	40.000	40.018	-0.0	91	0.01
92	Benzo(g,h,i)perylene	40.000	39.179	2.1	91	0.00

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