Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100516\
Data File : BF090380.D
Acq On : 5 Oct 2016 14:41
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF100516

Quant Time: Oct 05 15:20:58 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF100516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 05 14:57:00 2016
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	108	0.00
2		1,4-Dioxane	40.000	40.757	-1.9	110	0.00
3		Pyridine	40.000	42.031	-5.1	110	0.00
4		n-Nitrosodimethylamine	40.000	41.724	-4.3	111	0.00
5	S	2-Fluorophenol	80.000	80.379	-0.5	106	0.00
6		Aniline	40.000	42.341	-5.9	110	0.01
7	S	Phenol-d6	80.000	81.739	-2.2	112	0.00
8		2-Chlorophenol	40.000	39.838	0.4	111	0.00
9		Benzaldehyde	40.000	40.567	-1.4	111	0.00
10	C	Phenol	40.000	40.194	-0.5	114	0.00
11		bis(2-Chloroethyl)ether	40.000	42.137	-5.3	107	0.00
12		1,3-Dichlorobenzene	40.000	39.873	0.3	110	0.00
13	C	1,4-Dichlorobenzene	40.000	43.648	-9.1	111	0.00
14		1,2-Dichlorobenzene	40.000	38.932	2.7	113	0.01
15		Benzyl Alcohol	40.000	41.076	-2.7	116	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	42.033	-5.1	106	0.01
17		2-Methylphenol	40.000	40.133	-0.3	109	0.00
18		Hexachloroethane	40.000	41.914	-4.8	111	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	43.197	-8.0	111	0.00
20		3+4-Methylphenols	40.000	43.406	-8.5	111	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	106	0.00
22		Acetophenone	40.000	39.361	1.6	109	0.00
23	S	Nitrobenzene-d5	80.000	83.455	-4.3	107	0.00
24		Nitrobenzene	40.000	39.880	0.3	111	0.00
25		Isophorone	40.000	41.942	-4.9	114	0.00
26	C	2-Nitrophenol	40.000	38.753	3.1	106	0.00
27		2,4-Dimethylphenol	40.000	37.498	6.3	108	0.00
28		bis(2-Chloroethoxy)methane	40.000	45.070	-12.7	110	0.00
29	C	2,4-Dichlorophenol	40.000	39.305	1.7	113	0.00
30		1,2,4-Trichlorobenzene	40.000	42.744	-6.9	109	0.00
31		Naphthalene	40.000	37.873	5.3	106	0.00
32		Benzoic acid	40.000	44.895	-12.2	123	0.01
33		4-Chloroaniline	40.000	43.169	-7.9	107	0.00
34	C	Hexachlorobutadiene	40.000	42.706	-6.8	113	0.00
35		Caprolactam	40.000	41.425	-3.6	109	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.824	-4.6	106	0.00
37		2-Methylnaphthalene	40.000	44.198	-10.5	109	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	109	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	39.863	0.3	115	0.00
40	P	Hexachlorocyclopentadiene	40.000	43.937	-9.8	114	0.00
41	S	2,4,6-Tribromophenol	80.000	88.100	-10.1	111	0.00
42	C	2,4,6-Trichlorophenol	40.000	45.752	-14.4	108	0.00
43		2,4,5-Trichlorophenol	40.000	38.590	3.5	108	0.00
44	S	2-Fluorobiphenyl	80.000	89.057	-11.3	108	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100516\
Data File : BF090380.D
Acq On : 5 Oct 2016 14:41
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF100516

Quant Time: Oct 05 15:20:58 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF100516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 05 14:57:00 2016
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	42.771	-6.9	112	0.00
46		2-Chloronaphthalene	40.000	45.939	-14.8	110	0.00
47		2-Nitroaniline	40.000	44.805	-12.0	114	0.00
48		Acenaphthylene	40.000	38.737	3.2	108	0.00
49		Dimethylphthalate	40.000	43.866	-9.7	111	0.00
50		2,6-Dinitrotoluene	40.000	44.294	-10.7	108	0.00
51	C	Acenaphthene	40.000	40.503	-1.3	112	0.00
52		3-Nitroaniline	40.000	47.961	-19.9	117	0.00
53	P	2,4-Dinitrophenol	40.000	45.136	-12.8	124	0.01
54		Dibenzofuran	40.000	39.429	1.4	112	0.00
55	P	4-Nitrophenol	40.000	45.746	-14.4	108	0.00
56		2,4-Dinitrotoluene	40.000	39.358	1.6	103	0.00
57		Fluorene	40.000	42.294	-5.7	114	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	38.668	3.3	110	0.00
59		Diethylphthalate	40.000	40.590	-1.5	110	0.00
60		4-Chlorophenyl-phenylether	40.000	46.411	-16.0	113	0.00
61		4-Nitroaniline	40.000	44.627	-11.6	122	0.00
62		Azobenzene	40.000	44.786	-12.0	108	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	109	0.00
64		4,6-Dinitro-2-methylphenol	40.000	38.957	2.6	108	0.00
65	c	n-Nitrosodiphenylamine	40.000	40.084	-0.2	108	0.00
66		4-Bromophenyl-phenylether	40.000	43.718	-9.3	109	0.00
67		Hexachlorobenzene	40.000	44.383	-11.0	112	0.00
68		Atrazine	40.000	40.202	-0.5	108	0.00
69	C	Pentachlorophenol	40.000	39.164	2.1	112	0.00
70		Phenanthrene	40.000	42.119	-5.3	112	0.00
71		Anthracene	40.000	38.910	2.7	111	0.00
72		Carbazole	40.000	44.252	-10.6	110	0.00
73		Di-n-butylphthalate	40.000	44.625	-11.6	111	0.00
74	C	Fluoranthene	40.000	36.955	7.6	111	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	110	0.00
76		Benzidine	40.000	38.243	4.4	110	0.00
77		Pyrene	40.000	37.721	5.7	111	0.00
78	S	Terphenyl-d14	80.000	78.062	2.4	111	0.00
79		Butylbenzylphthalate	40.000	38.849	2.9	111	0.00
80		Benzo(a)anthracene	40.000	41.279	-3.2	115	0.00
81		3,3'-Dichlorobenzidine	40.000	36.871	7.8	115	0.00
82		Chrysene	40.000	39.996	0.0	112	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	42.548	-6.4	115	0.00
84	c	Di-n-octyl phthalate	40.000	41.212	-3.0	115	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	41.771	-4.4	123	0.01
86	I	Perylene-d12	20.000	20.000	0.0	117	0.00
87		Benzo(b)fluoranthene	40.000	43.384	-8.5	116	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100516\
Data File : BF090380.D
Acq On : 5 Oct 2016 14:41
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF100516

Quant Time: Oct 05 15:20:58 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF100516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 05 14:57:00 2016
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	37.150	7.1	116	0.00
89 C	Benzo(a)pyrene	40.000	41.275	-3.2	117	0.00
90	Dibenzo(a,h)anthracene	40.000	41.951	-4.9	122	0.00
91	Benzo(g,h,i)perylene	40.000	41.808	-4.5	122	0.00

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