Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120916\
Data File : BF091532.D
Acq On : 9 Dec 2016 17:08
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF120916

Quant Time: Dec 09 19:04:42 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 09 19:00:21 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	101	0.00
2		1,4-Dioxane	40.000	38.291	4.3	101	-0.01
3		Pyridine	40.000	40.476	-1.2	99	-0.01
4		n-Nitrosodimethylamine	40.000	37.529	6.2	96	-0.01
5	S	2-Fluorophenol	80.000	81.450	-1.8	103	0.00
6		Aniline	40.000	38.376	4.1	101	0.00
7	S	Phenol-d6	80.000	74.860	6.4	103	0.00
8		2-Chlorophenol	40.000	39.919	0.2	100	0.00
9		Benzaldehyde	40.000	34.442	13.9	98	0.00
10	C	Phenol	40.000	37.221	6.9	112	0.00
11		bis(2-Chloroethyl)ether	40.000	41.573	-3.9	104	0.00
12		1,3-Dichlorobenzene	40.000	36.756	8.1	101	0.00
13	C	1,4-Dichlorobenzene	40.000	38.166	4.6	101	0.00
14		1,2-Dichlorobenzene	40.000	40.155	-0.4	100	0.00
15		Benzyl Alcohol	40.000	36.068	9.8	96	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.821	0.4	99	0.00
17		2-Methylphenol	40.000	39.246	1.9	100	0.00
18		Hexachloroethane	40.000	38.325	4.2	99	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.577	3.6	97	-0.01
20		3+4-Methylphenols	40.000	37.350	6.6	101	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	102	0.00
22		Acetophenone	40.000	38.410	4.0	103	0.00
23	S	Nitrobenzene-d5	80.000	74.399	7.0	97	0.00
24		Nitrobenzene	40.000	42.087	-5.2	101	0.00
25		Isophorone	40.000	38.849	2.9	101	0.00
26	C	2-Nitrophenol	40.000	40.276	-0.7	97	0.00
27		2,4-Dimethylphenol	40.000	36.639	8.4	101	0.00
28		bis(2-Chloroethoxy)methane	40.000	41.390	-3.5	103	0.00
29	C	2,4-Dichlorophenol	40.000	40.959	-2.4	99	0.00
30		1,2,4-Trichlorobenzene	40.000	37.942	5.1	101	0.00
31		Naphthalene	40.000	37.702	5.7	101	0.00
32		Benzoic acid	40.000	39.898	0.3	107	0.00
33		4-Chloroaniline	40.000	38.150	4.6	102	0.00
34	C	Hexachlorobutadiene	40.000	40.601	-1.5	102	0.00
35		Caprolactam	40.000	39.924	0.2	109	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.846	-4.6	104	0.00
37		2-Methylnaphthalene	40.000	38.512	3.7	102	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	101	-0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	35.737	10.7	100	0.00
40	P	Hexachlorocyclopentadiene	40.000	40.183	-0.5	100	0.00
41	S	2,4,6-Tribromophenol	80.000	72.488	9.4	95	0.00
42	C	2,4,6-Trichlorophenol	40.000	42.365	-5.9	101	0.00
43		2,4,5-Trichlorophenol	40.000	36.848	7.9	101	0.00
44	S	2-Fluorobiphenyl	80.000	75.741	5.3	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120916\
Data File : BF091532.D
Acq On : 9 Dec 2016 17:08
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF120916

Quant Time: Dec 09 19:04:42 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 09 19:00:21 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	36.634	8.4	100	0.00
46		2-Chloronaphthalene	40.000	36.514	8.7	100	0.00
47		2-Nitroaniline	40.000	39.403	1.5	97	-0.01
48		Acenaphthylene	40.000	37.367	6.6	99	0.00
49		Dimethylphthalate	40.000	41.692	-4.2	103	0.00
50		2,6-Dinitrotoluene	40.000	43.236	-8.1	105	0.00
51	C	Acenaphthene	40.000	37.212	7.0	99	0.00
52		3-Nitroaniline	40.000	40.023	-0.1	95	0.00
53	P	2,4-Dinitrophenol	40.000	39.447	1.4	106	0.00
54		Dibenzofuran	40.000	37.180	7.1	98	0.00
55	P	4-Nitrophenol	40.000	45.731	-14.3	100	0.00
56		2,4-Dinitrotoluene	40.000	46.706	-16.8	108	0.00
57		Fluorene	40.000	37.048	7.4	102	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	42.839	-7.1	103	0.00
59		Diethylphthalate	40.000	42.139	-5.3	103	0.00
60		4-Chlorophenyl-phenylether	40.000	39.148	2.1	102	0.00
61		4-Nitroaniline	40.000	38.988	2.5	107	0.00
62		Azobenzene	40.000	41.342	-3.4	101	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	103	0.00
64		4,6-Dinitro-2-methylphenol	40.000	35.248	11.9	97	-0.01
65	c	n-Nitrosodiphenylamine	40.000	36.396	9.0	103	0.00
66		4-Bromophenyl-phenylether	40.000	37.772	5.6	99	0.00
67		Hexachlorobenzene	40.000	37.462	6.3	104	0.00
68		Atrazine	40.000	35.938	10.2	93	-0.01
69	C	Pentachlorophenol	40.000	40.937	-2.3	102	0.00
70		Phenanthrene	40.000	37.762	5.6	102	0.00
71		Anthracene	40.000	38.167	4.6	101	0.00
72		Carbazole	40.000	35.973	10.1	103	0.00
73		Di-n-butylphthalate	40.000	40.891	-2.2	102	0.00
74	C	Fluoranthene	40.000	38.773	3.1	99	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	102	0.00
76		Benzidine	40.000	36.291	9.3	90	-0.01
77		Pyrene	40.000	39.959	0.1	98	0.00
78	S	Terphenyl-d14	80.000	79.859	0.2	98	0.00
79		Butylbenzylphthalate	40.000	39.631	0.9	99	0.00
80		Benzo(a)anthracene	40.000	38.011	5.0	98	0.00
81		3,3'-Dichlorobenzidine	40.000	36.688	8.3	95	0.00
82		Chrysene	40.000	37.595	6.0	94	-0.01
83		Bis(2-ethylhexyl)phthalate	40.000	40.836	-2.1	101	0.00
84	c	Di-n-octyl phthalate	40.000	40.982	-2.5	96	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	38.686	3.3	97	-0.01
86	I	Perylene-d12	20.000	20.000	0.0	96	0.00
87		Benzo(b)fluoranthene	40.000	41.986	-5.0	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120916\
Data File : BF091532.D
Acq On : 9 Dec 2016 17:08
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF120916

Quant Time: Dec 09 19:04:42 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 09 19:00:21 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	34.358	14.1	94	0.00
89 C	Benzo(a)pyrene	40.000	39.266	1.8	97	0.00
90	Dibenzo(a,h)anthracene	40.000	38.790	3.0	97	0.00
91	Benzo(g,h,i)perylene	40.000	39.209	2.0	99	0.00

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