Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF122415\
Data File : BF084082.D
Acq On : 24 Dec 2015 16:31
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF122415

Quant Time: Dec 25 03:54:28 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 24 16:48:39 2015
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	102	0.00
2		1,4-Dioxane	40.000	39.742	0.6	98	0.00
3		Pyridine	40.000	45.042	-12.6	117	-0.01
4		n-Nitrosodimethylamine	40.000	40.712	-1.8	106	-0.01
5	S	2-Fluorophenol	80.000	79.917	0.1	98	0.00
6		Aniline	40.000	39.802	0.5	99	0.00
7	S	Phenol-d6	80.000	74.151	7.3	99	0.00
8		2-Chlorophenol	40.000	39.048	2.4	96	0.00
9		Benzaldehyde	40.000	37.429	6.4	97	0.00
10	C	Phenol	40.000	37.004	7.5	102	0.00
11		bis(2-Chloroethyl)ether	40.000	37.588	6.0	98	0.00
12		1,3-Dichlorobenzene	40.000	38.974	2.6	97	0.00
13	C	1,4-Dichlorobenzene	40.000	39.076	2.3	98	0.00
14		1,2-Dichlorobenzene	40.000	37.584	6.0	99	0.00
15		Benzyl Alcohol	40.000	39.009	2.5	100	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.271	6.8	100	0.00
17		2-Methylphenol	40.000	36.368	9.1	98	0.00
18		Hexachloroethane	40.000	37.816	5.5	96	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	36.823	7.9	98	0.00
20		3+4-Methylphenols	40.000	38.217	4.5	98	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	101	0.00
22		Acetophenone	40.000	38.092	4.8	97	0.00
23	S	Nitrobenzene-d5	80.000	74.422	7.0	98	0.00
24		Nitrobenzene	40.000	39.474	1.3	97	0.00
25		Isophorone	40.000	37.453	6.4	96	0.00
26	C	2-Nitrophenol	40.000	39.362	1.6	100	0.00
27		2,4-Dimethylphenol	40.000	40.995	-2.5	98	0.00
28		bis(2-Chloroethoxy)methane	40.000	34.155	14.6	98	0.00
29	C	2,4-Dichlorophenol	40.000	36.989	7.5	98	0.00
30		1,2,4-Trichlorobenzene	40.000	39.019	2.5	98	0.00
31		Naphthalene	40.000	38.461	3.8	97	0.00
32		Benzoic acid	40.000	36.683	8.3	101	0.00
33		4-Chloroaniline	40.000	36.428	8.9	99	0.01
34	C	Hexachlorobutadiene	40.000	36.658	8.4	100	0.00
35		Caprolactam	40.000	39.139	2.2	93	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.546	1.1	98	0.00
37		2-Methylnaphthalene	40.000	35.968	10.1	99	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	102	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.672	3.3	97	0.00
40	P	Hexachlorocyclopentadiene	40.000	38.727	3.2	107	0.00
41	S	2,4,6-Tribromophenol	80.000	70.346	12.1	97	0.00
42	C	2,4,6-Trichlorophenol	40.000	37.664	5.8	98	0.00
43		2,4,5-Trichlorophenol	40.000	38.622	3.4	97	0.00
44	S	2-Fluorobiphenyl	80.000	70.306	12.1	95	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF122415\
Data File : BF084082.D
Acq On : 24 Dec 2015 16:31
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF122415

Quant Time: Dec 25 03:54:28 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 24 16:48:39 2015
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	40.557	-1.4	98	0.00
46		2-Chloronaphthalene	40.000	38.327	4.2	96	0.00
47		2-Nitroaniline	40.000	40.969	-2.4	101	0.00
48		Acenaphthylene	40.000	36.819	8.0	99	0.00
49		Dimethylphthalate	40.000	37.217	7.0	98	-0.01
50		2,6-Dinitrotoluene	40.000	39.979	0.1	94	0.00
51	C	Acenaphthene	40.000	36.881	7.8	98	0.00
52		3-Nitroaniline	40.000	42.076	-5.2	101	0.00
53	P	2,4-Dinitrophenol	40.000	36.775	8.1	103	0.00
54		Dibenzofuran	40.000	35.558	11.1	98	0.00
55	P	4-Nitrophenol	40.000	39.158	2.1	101	0.00
56		2,4-Dinitrotoluene	40.000	36.314	9.2	98	0.00
57		Fluorene	40.000	36.935	7.7	100	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.599	1.0	101	0.00
59		Diethylphthalate	40.000	33.821	15.4	95	0.00
60		4-Chlorophenyl-phenylether	40.000	33.434	16.4	93	0.00
61		4-Nitroaniline	40.000	40.377	-0.9	101	0.00
62		Azobenzene	40.000	37.315	6.7	98	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	102	0.00
64		4,6-Dinitro-2-methylphenol	40.000	37.589	6.0	104	0.00
65	c	n-Nitrosodiphenylamine	40.000	40.098	-0.2	98	0.00
66		4-Bromophenyl-phenylether	40.000	39.377	1.6	100	0.00
67		Hexachlorobenzene	40.000	35.653	10.9	96	0.00
68		Atrazine	40.000	34.977	12.6	98	0.00
69	C	Pentachlorophenol	40.000	36.518	8.7	99	0.00
70		Phenanthrene	40.000	38.118	4.7	96	0.00
71		Anthracene	40.000	36.893	7.8	97	0.00
72		Carbazole	40.000	36.173	9.6	98	0.00
73		Di-n-butylphthalate	40.000	36.777	8.1	99	0.00
74	C	Fluoranthene	40.000	38.356	4.1	97	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	109	0.00
76		Benzidine	40.000	36.624	8.4	97	0.00
77		Pyrene	40.000	37.872	5.3	99	0.00
78	S	Terphenyl-d14	80.000	67.930	15.1	99	0.00
79		Butylbenzylphthalate	40.000	38.177	4.6	103	0.00
80		Benzo(a)anthracene	40.000	37.713	5.7	103	0.00
81		3,3'-Dichlorobenzidine	40.000	39.555	1.1	99	0.00
82		Chrysene	40.000	40.104	-0.3	103	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	38.205	4.5	100	0.00
84	c	Di-n-octyl phthalate	40.000	38.843	2.9	99	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	37.428	6.4	98	0.01
86	I	Perylene-d12	20.000	20.000	0.0	106	0.00
87		Benzo(b)fluoranthene	40.000	37.367	6.6	99	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF122415\
Data File : BF084082.D
Acq On : 24 Dec 2015 16:31
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF122415

Quant Time: Dec 25 03:54:28 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 24 16:48:39 2015
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	40.460	-1.2	113	0.00
89 C	Benzo(a)pyrene	40.000	36.593	8.5	101	0.00
90	Dibenzo(a,h)anthracene	40.000	37.726	5.7	98	0.00
91	Benzo(g,h,i)perylene	40.000	37.433	6.4	98	0.00

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