Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF030315\
Data File : BF077542.D
Acq On : 3 Mar 2015 13:34
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 04 00:06:19 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 03 23:57:38 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	93	0.00
2		1,4-Dioxane	40.000	38.259	4.4	90	0.00
3		Pyridine	40.000	38.100	4.7	83	0.00
4		n-Nitrosodimethylamine	40.000	37.798	5.5	90	0.00
5	S	2-Fluorophenol	80.000	76.056	4.9	87	0.00
6		Aniline	40.000	36.500	8.8	82	0.00
7	S	Phenol-d6	80.000	72.636	9.2	82	0.00
8		2-Chlorophenol	40.000	36.133	9.7	83	0.00
9		Benzaldehyde	40.000	41.359	-3.4	98	0.00
10	C	Phenol	40.000	36.942	7.6	80	0.00
11		bis(2-Chloroethyl)ether	40.000	34.680	13.3	81	0.00
12		1,3-Dichlorobenzene	40.000	38.011	5.0	86	0.00
13	C	1,4-Dichlorobenzene	40.000	37.193	7.0	84	0.00
14		1,2-Dichlorobenzene	40.000	37.576	6.1	84	0.00
15		Benzyl Alcohol	40.000	38.446	3.9	87	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	35.248	11.9	79	0.00
17		2-Methylphenol	40.000	37.298	6.8	80	0.00
18		Hexachloroethane	40.000	37.273	6.8	85	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	36.048	9.9	80	0.00
20		3+4-Methylphenols	40.000	36.698	8.3	81	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	80	0.00
22		Acetophenone	40.000	40.672	-1.7	84	0.00
23	S	Nitrobenzene-d5	80.000	84.332	-5.4	82	0.00
24		Nitrobenzene	40.000	41.149	-2.9	81	0.00
25		Isophorone	40.000	37.406	6.5	70	0.00
26	C	2-Nitrophenol	40.000	46.289	-15.7	83	0.00
27		2,4-Dimethylphenol	40.000	35.691	10.8	70	0.00
28		bis(2-Chloroethoxy)methane	40.000	35.307	11.7	69	0.00
29	C	2,4-Dichlorophenol	40.000	39.040	2.4	76	0.00
30		1,2,4-Trichlorobenzene	40.000	36.795	8.0	72	0.00
31		Naphthalene	40.000	36.875	7.8	74	0.00
32		Benzoic acid	40.000	42.240	-5.6	77	0.00
33		4-Chloroaniline	40.000	37.447	6.4	71	0.00
34	C	Hexachlorobutadiene	40.000	35.999	10.0	71	0.00
35		Caprolactam	40.000	39.901	0.2	76	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.031	-0.1	75	0.00
37		2-Methylnaphthalene	40.000	35.336	11.7	67	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	70	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	40.359	-0.9	68	0.00
40	P	Hexachlorocyclopentadiene	40.000	39.940	0.2	63	0.00
41	S	2,4,6-Tribromophenol	80.000	84.885	-6.1	70	0.00
42	C	2,4,6-Trichlorophenol	40.000	41.012	-2.5	67	0.00
43		2,4,5-Trichlorophenol	40.000	43.231	-8.1	72	0.00
44	S	2-Fluorobiphenyl	80.000	82.659	-3.3	72	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF030315\
Data File : BF077542.D
Acq On : 3 Mar 2015 13:34
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 04 00:06:19 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 03 23:57:38 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	40.135	-0.3	68	0.00
46		2-Chloronaphthalene	40.000	40.939	-2.3	72	0.00
47		2-Nitroaniline	40.000	45.235	-13.1	76	0.00
48		Acenaphthylene	40.000	42.206	-5.5	73	0.00
49		Dimethylphthalate	40.000	40.665	-1.7	70	0.00
50		2,6-Dinitrotoluene	40.000	45.985	-15.0	72	0.00
51	C	Acenaphthene	40.000	41.089	-2.7	70	0.00
52		3-Nitroaniline	40.000	42.690	-6.7	69	0.00
53	P	2,4-Dinitrophenol	40.000	55.351	-38.4#	92	0.00
54		Dibenzofuran	40.000	40.723	-1.8	70	0.00
55	P	4-Nitrophenol	40.000	47.486	-18.7	75	0.00
56		2,4-Dinitrotoluene	40.000	43.592	-9.0	68	0.00
57		Fluorene	40.000	42.397	-6.0	72	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	42.891	-7.2	69	0.00
59		Diethylphthalate	40.000	39.374	1.6	67	0.00
60		4-Chlorophenyl-phenylether	40.000	38.590	3.5	65	0.00
61		4-Nitroaniline	40.000	44.684	-11.7	74	0.00
62		Azobenzene	40.000	39.976	0.1	66	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	72	0.00
64		4,6-Dinitro-2-methylphenol	40.000	46.564	-16.4	84	0.00
65	c	n-Nitrosodiphenylamine	40.000	37.725	5.7	66	0.00
66		4-Bromophenyl-phenylether	40.000	36.519	8.7	66	0.00
67		Hexachlorobenzene	40.000	36.189	9.5	67	0.00
68		Atrazine	40.000	33.808	15.5	65	0.00
69	C	Pentachlorophenol	40.000	37.584	6.0	64	0.00
70		Phenanthrene	40.000	37.823	5.4	69	0.00
71		Anthracene	40.000	38.659	3.4	71	0.00
72		Carbazole	40.000	38.755	3.1	67	0.00
73		Di-n-butylphthalate	40.000	36.145	9.6	62	0.00
74	C	Fluoranthene	40.000	38.000	5.0	67	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	70	0.03
76		Benzidine	40.000	39.292	1.8	74	0.00
77		Pyrene	40.000	39.343	1.6	70	0.00
78	S	Terphenyl-d14	80.000	71.276	10.9	62	0.00
79		Butylbenzylphthalate	40.000	38.650	3.4	64	0.01
80		Benzo(a)anthracene	40.000	38.693	3.3	69	0.03
81		3,3'-Dichlorobenzidine	40.000	39.684	0.8	68	0.03
82		Chrysene	40.000	37.062	7.3	67	0.04
83		Bis(2-ethylhexyl)phthalate	40.000	35.120	12.2	60	0.04
84	c	Di-n-octyl phthalate	40.000	36.273	9.3	60	0.10
85		Indeno(1,2,3-cd)pyrene	40.000	39.270	1.8	68	0.21
86	I	Perylene-d12	20.000	20.000	0.0	74	0.18
87		Benzo(b)fluoranthene	40.000	38.844	2.9	67	0.14

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF030315\
Data File : BF077542.D
Acq On : 3 Mar 2015 13:34
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 04 00:06:19 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 03 23:57:38 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	37.216	7.0	73	0.14
89 C	Benzo(a)pyrene	40.000	38.006	5.0	68	0.18
90	Dibenzo(a,h)anthracene	40.000	38.655	3.4	69	0.21
91	Benzo(g,h,i)perylene	40.000	43.889	-9.7	79	0.22

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