Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040621\
Data File : BP005191.D
Acq On : 06 Apr 2021 11:31
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Apr 06 12:35:21 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 06 12:33:08 2021
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	57	0.00
2		1,4-Dioxane	40.000	47.049	-17.6	73	0.00
3		Pyridine	40.000	42.520	-6.3	58	0.00
4		n-Nitrosodimethylamine	40.000	43.437	-8.6	62	0.00
5	S	2-Fluorophenol	80.000	82.179	-2.7	60	0.00
6		Aniline	40.000	41.378	-3.4	59	0.00
7	S	Phenol-d6	80.000	81.863	-2.3	58	0.00
8		2-Chlorophenol	40.000	41.843	-4.6	60	0.00
9		Benzaldehyde	40.000	35.463	11.3	54	0.00
10	C	Phenol	40.000	40.743	-1.9	58	0.00
11		bis(2-Chloroethyl)ether	40.000	40.783	-2.0	59	0.00
12		1,3-Dichlorobenzene	40.000	39.986	0.0	57	0.00
13	C	1,4-Dichlorobenzene	40.000	39.255	1.9	57	0.00
14		1,2-Dichlorobenzene	40.000	39.518	1.2	58	0.00
15		Benzyl Alcohol	40.000	39.994	0.0	56	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	45.323	-13.3	64	0.00
17		2-Methylphenol	40.000	41.766	-4.4	59	0.00
18		Hexachloroethane	40.000	38.979	2.6	57	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.959	0.1	56	0.00
20		3+4-Methylphenols	40.000	41.849	-4.6	58	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	55	0.00
22		Acetophenone	40.000	41.527	-3.8	57	0.00
23	S	Nitrobenzene-d5	80.000	80.468	-0.6	55	0.00
24		Nitrobenzene	40.000	39.081	2.3	54	0.00
25		Isophorone	40.000	42.086	-5.2	57	0.00
26	C	2-Nitrophenol	40.000	45.998	-15.0	62	0.00
27		2,4-Dimethylphenol	40.000	41.464	-3.7	57	0.00
28		bis(2-Chloroethoxy)methane	40.000	42.157	-5.4	57	0.00
29	C	2,4-Dichlorophenol	40.000	40.492	-1.2	55	0.00
30		1,2,4-Trichlorobenzene	40.000	39.842	0.4	55	0.00
31		Naphthalene	40.000	40.951	-2.4	57	0.00
32		Benzoic acid	40.000	44.041	-10.1	57	0.00
33		4-Chloroaniline	40.000	41.936	-4.8	58	0.00
34	C	Hexachlorobutadiene	40.000	36.962	7.6	50	0.00
35		Caprolactam	40.000	48.720	-21.8	66	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.014	-5.0	57	0.00
37		2-Methylnaphthalene	40.000	40.337	-0.8	55	0.00
38		1-Methylnaphthalene	40.000	40.463	-1.2	55	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	56	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	37.087	7.3	52	0.00
41	P	Hexachlorocyclopentadiene	40.000	34.044	14.9	46	0.00
42	S	2,4,6-Tribromophenol	80.000	77.809	2.7	55	0.00
43	C	2,4,6-Trichlorophenol	40.000	38.983	2.5	54	0.00
44		2,4,5-Trichlorophenol	40.000	39.270	1.8	55	0.00
45	S	2-Fluorobiphenyl	80.000	74.904	6.4	53	0.00
46		1,1'-Biphenyl	40.000	38.313	4.2	54	0.00
47		2-Chloronaphthalene	40.000	38.788	3.0	56	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040621\
Data File : BP005191.D
Acq On : 06 Apr 2021 11:31
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Apr 06 12:35:21 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 06 12:33:08 2021
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)
48		2-Nitroaniline	40.000	41.092	-2.7	56	0.00
49		Acenaphthylene	40.000	39.434	1.4	56	0.00
50		Dimethylphthalate	40.000	38.242	4.4	55	0.00
51		2,6-Dinitrotoluene	40.000	39.331	1.7	55	0.00
52	C	Acenaphthene	40.000	38.485	3.8	55	0.00
53		3-Nitroaniline	40.000	44.718	-11.8	60	0.00
54	P	2,4-Dinitrophenol	40.000	40.948	-2.4	56	0.00
55		Dibenzofuran	40.000	38.484	3.8	55	0.00
56	P	4-Nitrophenol	40.000	43.763	-9.4	58	0.00
57		2,4-Dinitrotoluene	40.000	41.862	-4.7	57	0.00
58		Fluorene	40.000	39.018	2.5	55	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	38.254	4.4	53	0.00
60		Diethylphthalate	40.000	39.394	1.5	55	0.00
61		4-Chlorophenyl-phenylether	40.000	36.979	7.6	54	0.00
62		4-Nitroaniline	40.000	46.457	-16.1	62	0.00
63		Azobenzene	40.000	37.290	6.8	53	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	57	0.00
65		4,6-Dinitro-2-methylphenol	40.000	43.330	-8.3	57	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.810	0.5	56	0.00
67		4-Bromophenyl-phenylether	40.000	38.113	4.7	53	0.00
68		Hexachlorobenzene	40.000	37.107	7.2	52	0.00
69		Atrazine	40.000	40.613	-1.5	54	0.00
70	C	Pentachlorophenol	40.000	38.402	4.0	54	0.00
71		Phenanthrene	40.000	38.867	2.8	54	0.00
72		Anthracene	40.000	40.069	-0.2	56	0.00
73		Carbazole	40.000	40.978	-2.4	58	0.00
74		Di-n-butylphthalate	40.000	42.714	-6.8	58	0.00
75	C	Fluoranthene	40.000	40.437	-1.1	57	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	61	0.00
77		Benzidine	40.000	47.485	-18.7	62	0.00
78		Pyrene	40.000	38.789	3.0	57	0.00
79	S	Terphenyl-d14	80.000	76.015	5.0	56	0.00
80		Butylbenzylphthalate	40.000	44.732	-11.8	63	0.00
81		Benzo(a)anthracene	40.000	40.253	-0.6	61	0.00
82		3,3'-Dichlorobenzidine	40.000	42.769	-6.9	62	0.00
83		Chrysene	40.000	40.386	-1.0	61	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	44.060	-10.2	62	0.00
85	c	Di-n-octyl phthalate	40.000	46.152	-15.4	66	0.00
86	I	Perylene-d12	20.000	20.000	0.0	65	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	39.170	2.1	63	0.00
88		Benzo(b)fluoranthene	40.000	37.086	7.3	60	0.00
89		Benzo(k)fluoranthene	40.000	38.728	3.2	61	0.00
90	C	Benzo(a)pyrene	40.000	39.085	2.3	62	0.00
91		Dibenzo(a,h)anthracene	40.000	39.544	1.1	63	0.00
92		Benzo(g,h,i)perylene	40.000	39.516	1.2	63	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040621\
Data File : BP005191.D
Acq On : 06 Apr 2021 11:31
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Apr 06 12:35:21 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 06 12:33:08 2021
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)

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