Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111521\
Data File : BP007978.D
Acq On : 15 Nov 2021 10:18
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Nov 15 11:44:07 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 15 05:36:49 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	84	0.00
2		1,4-Dioxane	40.000	37.128	7.2	86	0.00
3		Pyridine	40.000	38.852	2.9	88	0.00
4		n-Nitrosodimethylamine	40.000	41.770	-4.4	89	0.00
5	S	2-Fluorophenol	80.000	77.210	3.5	87	0.00
6		Aniline	40.000	40.724	-1.8	87	0.00
7	S	Phenol-d6	80.000	78.954	1.3	87	0.00
8		2-Chlorophenol	40.000	40.156	-0.4	85	0.00
9		Benzaldehyde	40.000	42.098	-5.2	85	0.00
10	C	Phenol	40.000	40.713	-1.8	86	0.00
11		bis(2-Chloroethyl)ether	40.000	38.091	4.8	85	0.00
12		1,3-Dichlorobenzene	40.000	40.101	-0.3	85	0.00
13	C	1,4-Dichlorobenzene	40.000	40.129	-0.3	86	0.00
14		1,2-Dichlorobenzene	40.000	38.562	3.6	86	0.00
15		Benzyl Alcohol	40.000	42.256	-5.6	88	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	42.540	-6.3	87	0.00
17		2-Methylphenol	40.000	41.040	-2.6	86	0.00
18		Hexachloroethane	40.000	38.578	3.6	81	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.282	4.3	84	0.00
20		3+4-Methylphenols	40.000	40.304	-0.8	84	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	81	0.00
22		Acetophenone	40.000	41.827	-4.6	84	0.00
23	S	Nitrobenzene-d5	80.000	83.974	-5.0	83	0.00
24		Nitrobenzene	40.000	41.498	-3.7	82	0.00
25		Isophorone	40.000	41.178	-2.9	81	0.00
26	C	2-Nitrophenol	40.000	41.206	-3.0	79	0.00
27		2,4-Dimethylphenol	40.000	40.728	-1.8	80	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.172	-0.4	81	0.00
29	C	2,4-Dichlorophenol	40.000	41.804	-4.5	82	0.00
30		1,2,4-Trichlorobenzene	40.000	40.598	-1.5	82	0.00
31		Naphthalene	40.000	40.081	-0.2	81	0.00
32		Benzoic acid	40.000	42.163	-5.4	78	0.00
33		4-Chloroaniline	40.000	40.739	-1.8	81	0.00
34	C	Hexachlorobutadiene	40.000	41.567	-3.9	82	0.00
35		Caprolactam	40.000	40.837	-2.1	79	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.707	-4.3	83	0.00
37		2-Methylnaphthalene	40.000	40.116	-0.3	81	0.00
38		1-Methylnaphthalene	40.000	40.338	-0.8	81	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	81	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.149	-0.4	81	0.00
41	P	Hexachlorocyclopentadiene	40.000	39.577	1.1	80	0.00
42	S	2,4,6-Tribromophenol	80.000	86.058	-7.6	88	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.634	-1.6	81	0.00
44		2,4,5-Trichlorophenol	40.000	40.354	-0.9	81	0.00
45	S	2-Fluorobiphenyl	80.000	78.798	1.5	83	0.00
46		1,1'-Biphenyl	40.000	39.203	2.0	81	0.00
47		2-Chloronaphthalene	40.000	39.606	1.0	82	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111521\
Data File : BP007978.D
Acq On : 15 Nov 2021 10:18
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Nov 15 11:44:07 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 15 05:36:49 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	42.507	-6.3	83	0.00
49		Acenaphthylene	40.000	40.688	-1.7	81	0.00
50		Dimethylphthalate	40.000	39.634	0.9	82	0.00
51		2,6-Dinitrotoluene	40.000	39.961	0.1	80	0.00
52	C	Acenaphthene	40.000	40.576	-1.4	82	0.00
53		3-Nitroaniline	40.000	41.446	-3.6	81	0.00
54	P	2,4-Dinitrophenol	40.000	43.369	-8.4	80	0.00
55		Dibenzofuran	40.000	40.869	-2.2	88	0.00
56	P	4-Nitrophenol	40.000	42.601	-6.5	83	0.00
57		2,4-Dinitrotoluene	40.000	43.091	-7.7	87	0.00
58		Fluorene	40.000	39.382	1.5	88	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	42.570	-6.4	89	0.00
60		Diethylphthalate	40.000	41.175	-2.9	87	0.00
61		4-Chlorophenyl-phenylether	40.000	40.845	-2.1	88	0.00
62		4-Nitroaniline	40.000	43.225	-8.1	88	0.00
63		Azobenzene	40.000	42.216	-5.5	88	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	87	0.00
65		4,6-Dinitro-2-methylphenol	40.000	43.237	-8.1	85	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.448	-1.1	88	0.00
67		4-Bromophenyl-phenylether	40.000	40.753	-1.9	87	0.00
68		Hexachlorobenzene	40.000	41.585	-4.0	89	0.00
69		Atrazine	40.000	40.778	-1.9	86	0.00
70	C	Pentachlorophenol	40.000	43.526	-8.8	87	0.00
71		Phenanthrene	40.000	39.980	0.1	87	0.00
72		Anthracene	40.000	40.611	-1.5	88	0.00
73		Carbazole	40.000	40.693	-1.7	88	0.00
74		Di-n-butylphthalate	40.000	41.877	-4.7	88	0.00
75	C	Fluoranthene	40.000	40.715	-1.8	88	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	92	0.00
77		Benzidine	40.000	48.229	-20.6	98	0.00
78		Pyrene	40.000	39.337	1.7	89	0.00
79	S	Terphenyl-d14	80.000	78.958	1.3	91	0.00
80		Butylbenzylphthalate	40.000	40.737	-1.8	89	0.00
81		Benzo(a)anthracene	40.000	39.880	0.3	91	0.00
82		3,3'-Dichlorobenzidine	40.000	42.513	-6.3	93	0.00
83		Chrysene	40.000	39.926	0.2	91	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	40.864	-2.2	93	0.00
85	c	Di-n-octyl phthalate	40.000	41.657	-4.1	91	0.00
86	I	Perylene-d12	20.000	20.000	0.0	93	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	41.798	-4.5	97	0.00
88		Benzo(b)fluoranthene	40.000	39.456	1.4	94	0.00
89		Benzo(k)fluoranthene	40.000	40.660	-1.6	93	0.00
90	C	Benzo(a)pyrene	40.000	40.371	-0.9	93	0.00
91		Dibenzo(a,h)anthracene	40.000	41.832	-4.6	97	0.00
92		Benzo(g,h,i)perylene	40.000	42.232	-5.6	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111521\
Data File : BP007978.D
Acq On : 15 Nov 2021 10:18
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Nov 15 11:44:07 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 15 05:36:49 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