Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092921\
Data File : BP007243.D
Acq On : 29 Sep 2021 13:28
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Sep 29 14:46:49 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 29 14:45:49 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	65	0.00
2		1,4-Dioxane	40.000	45.233	-13.1	77	0.00
3		Pyridine	40.000	44.851	-12.1	73	0.00
4		n-Nitrosodimethylamine	40.000	47.578	-18.9	78	0.00
5	S	2-Fluorophenol	80.000	75.122	6.1	63	0.00
6		Aniline	40.000	37.210	7.0	60	0.00
7	S	Phenol-d6	80.000	74.426	7.0	61	0.00
8		2-Chlorophenol	40.000	37.685	5.8	63	0.00
9		Benzaldehyde	40.000	35.868	10.3	58	0.00
10	C	Phenol	40.000	36.742	8.1	60	0.00
11		bis(2-Chloroethyl)ether	40.000	36.188	9.5	59	0.00
12		1,3-Dichlorobenzene	40.000	38.232	4.4	64	0.00
13	C	1,4-Dichlorobenzene	40.000	38.610	3.5	65	0.00
14		1,2-Dichlorobenzene	40.000	38.211	4.5	64	0.00
15		Benzyl Alcohol	40.000	39.035	2.4	63	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	36.255	9.4	60	0.00
17		2-Methylphenol	40.000	38.019	5.0	62	0.00
18		Hexachloroethane	40.000	38.584	3.5	64	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	36.806	8.0	60	0.00
20		3+4-Methylphenols	40.000	37.558	6.1	61	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	61	0.00
22		Acetophenone	40.000	38.939	2.7	61	0.00
23	S	Nitrobenzene-d5	80.000	79.861	0.2	62	0.00
24		Nitrobenzene	40.000	39.461	1.3	60	0.00
25		Isophorone	40.000	39.154	2.1	60	0.00
26	C	2-Nitrophenol	40.000	42.395	-6.0	64	0.00
27		2,4-Dimethylphenol	40.000	38.849	2.9	59	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.283	6.8	58	0.00
29	C	2,4-Dichlorophenol	40.000	40.609	-1.5	62	0.00
30		1,2,4-Trichlorobenzene	40.000	40.393	-1.0	64	0.00
31		Naphthalene	40.000	38.660	3.4	60	0.00
32		Benzoic acid	40.000	40.420	-1.1	59	0.00
33		4-Chloroaniline	40.000	39.095	2.3	60	0.00
34	C	Hexachlorobutadiene	40.000	43.205	-8.0	68	0.00
35		Caprolactam	40.000	42.467	-6.2	63	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.370	-0.9	62	0.00
37		2-Methylnaphthalene	40.000	38.782	3.0	61	0.00
38		1-Methylnaphthalene	40.000	40.035	-0.1	63	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	62	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.381	-1.0	64	0.00
41	P	Hexachlorocyclopentadiene	40.000	33.682	15.8	52	0.00
42	S	2,4,6-Tribromophenol	80.000	91.239	-14.0	71	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.818	-2.0	63	0.00
44		2,4,5-Trichlorophenol	40.000	40.527	-1.3	63	0.00
45	S	2-Fluorobiphenyl	80.000	77.653	2.9	62	0.00
46		1,1'-Biphenyl	40.000	38.210	4.5	61	0.00
47		2-Chloronaphthalene	40.000	38.782	3.0	62	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092921\
Data File : BP007243.D
Acq On : 29 Sep 2021 13:28
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Sep 29 14:46:49 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 29 14:45:49 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	41.604	-4.0	63	0.00
49		Acenaphthylene	40.000	39.431	1.4	62	0.00
50		Dimethylphthalate	40.000	39.584	1.0	63	0.00
51		2,6-Dinitrotoluene	40.000	41.359	-3.4	64	0.00
52	C	Acenaphthene	40.000	38.143	4.6	60	0.00
53		3-Nitroaniline	40.000	40.807	-2.0	62	0.00
54	P	2,4-Dinitrophenol	40.000	43.200	-8.0	63	0.00
55		Dibenzofuran	40.000	38.744	3.1	62	0.00
56	P	4-Nitrophenol	40.000	40.952	-2.4	60	0.00
57		2,4-Dinitrotoluene	40.000	43.359	-8.4	66	0.00
58		Fluorene	40.000	39.288	1.8	62	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	41.293	-3.2	65	0.00
60		Diethylphthalate	40.000	39.989	0.0	64	0.00
61		4-Chlorophenyl-phenylether	40.000	40.188	-0.5	64	0.00
62		4-Nitroaniline	40.000	43.098	-7.7	64	0.00
63		Azobenzene	40.000	38.545	3.6	61	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	65	0.00
65		4,6-Dinitro-2-methylphenol	40.000	43.266	-8.2	66	0.00
66	c	n-Nitrosodiphenylamine	40.000	38.390	4.0	63	0.00
67		4-Bromophenyl-phenylether	40.000	40.625	-1.6	67	0.00
68		Hexachlorobenzene	40.000	41.226	-3.1	69	0.00
69		Atrazine	40.000	40.505	-1.3	65	0.00
70	C	Pentachlorophenol	40.000	38.892	2.8	62	0.00
71		Phenanthrene	40.000	38.203	4.5	63	0.00
72		Anthracene	40.000	38.926	2.7	64	0.00
73		Carbazole	40.000	39.262	1.8	64	0.00
74		Di-n-butylphthalate	40.000	39.842	0.4	64	0.00
75	C	Fluoranthene	40.000	40.248	-0.6	66	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	67	0.00
77		Benzidine	40.000	36.617	8.5	62	0.00
78		Pyrene	40.000	38.163	4.6	66	0.00
79	S	Terphenyl-d14	80.000	78.274	2.2	67	0.00
80		Butylbenzylphthalate	40.000	45.515	-13.8	76	0.00
81		Benzo(a)anthracene	40.000	38.786	3.0	67	0.00
82		3,3'-Dichlorobenzidine	40.000	40.960	-2.4	69	0.00
83		Chrysene	40.000	38.566	3.6	66	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	39.229	1.9	66	0.00
85	c	Di-n-octyl phthalate	40.000	40.040	-0.1	67	0.00
86	I	Perylene-d12	20.000	20.000	0.0	71	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	39.466	1.3	72	0.00
88		Benzo(b)fluoranthene	40.000	38.251	4.4	69	0.00
89		Benzo(k)fluoranthene	40.000	37.258	6.9	68	0.00
90	C	Benzo(a)pyrene	40.000	38.502	3.7	70	0.00
91		Dibenzo(a,h)anthracene	40.000	39.438	1.4	72	0.00
92		Benzo(g,h,i)perylene	40.000	39.704	0.7	72	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092921\
Data File : BP007243.D
Acq On : 29 Sep 2021 13:28
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Sep 29 14:46:49 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 29 14:45:49 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