Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121821\
Data File : BP008488.D
Acq On : 18 Dec 2021 14:24
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Dec 20 04:19:22 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP121421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Dec 18 06:49:58 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	86	0.00
2		1,4-Dioxane	40.000	46.514	-16.3	90	0.00
3		Pyridine	40.000	42.250	-5.6	89	0.00
4		n-Nitrosodimethylamine	40.000	45.224	-13.1	96	0.00
5	S	2-Fluorophenol	80.000	79.004	1.2	84	0.00
6		Aniline	40.000	42.310	-5.8	86	0.00
7	S	Phenol-d6	80.000	84.338	-5.4	86	0.00
8		2-Chlorophenol	40.000	41.451	-3.6	85	0.00
9		Benzaldehyde	40.000	41.525	-3.8	85	0.00
10	C	Phenol	40.000	41.965	-4.9	86	0.00
11		bis(2-Chloroethyl)ether	40.000	40.432	-1.1	88	0.00
12		1,3-Dichlorobenzene	40.000	39.462	1.3	83	0.00
13	C	1,4-Dichlorobenzene	40.000	40.040	-0.1	84	0.00
14		1,2-Dichlorobenzene	40.000	40.125	-0.3	84	-0.01
15		Benzyl Alcohol	40.000	43.595	-9.0	85	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	42.731	-6.8	92	-0.02
17		2-Methylphenol	40.000	43.123	-7.8	87	0.00
18		Hexachloroethane	40.000	41.048	-2.6	87	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	42.483	-6.2	90	0.00
20		3+4-Methylphenols	40.000	43.308	-8.3	87	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	92	0.00
22		Acetophenone	40.000	38.996	2.5	88	0.00
23	S	Nitrobenzene-d5	80.000	78.472	1.9	88	-0.01
24		Nitrobenzene	40.000	39.636	0.9	89	0.00
25		Isophorone	40.000	41.618	-4.0	93	0.00
26	C	2-Nitrophenol	40.000	40.108	-0.3	89	0.00
27		2,4-Dimethylphenol	40.000	40.029	-0.1	89	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.770	-1.9	90	0.00
29	C	2,4-Dichlorophenol	40.000	39.420	1.4	87	0.00
30		1,2,4-Trichlorobenzene	40.000	38.507	3.7	88	0.00
31		Naphthalene	40.000	38.988	2.5	89	0.00
32		Benzoic acid	40.000	42.769	-6.9	86	0.00
33		4-Chloroaniline	40.000	41.627	-4.1	92	0.00
34	C	Hexachlorobutadiene	40.000	37.306	6.7	85	-0.01
35		Caprolactam	40.000	49.292	-23.2	108	-0.01
36	C	4-Chloro-3-methylphenol	40.000	43.285	-8.2	95	0.00
37		2-Methylnaphthalene	40.000	40.926	-2.3	91	0.00
38		1-Methylnaphthalene	40.000	41.576	-3.9	92	-0.01
39	I	Acenaphthene-d10	20.000	20.000	0.0	107	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	34.667	13.3	91	0.00
41	P	Hexachlorocyclopentadiene	40.000	41.102	-2.8	113	0.00
42	S	2,4,6-Tribromophenol	80.000	84.831	-6.0	111	0.00
43	C	2,4,6-Trichlorophenol	40.000	36.812	8.0	93	0.00
44		2,4,5-Trichlorophenol	40.000	36.858	7.9	94	0.00
45	S	2-Fluorobiphenyl	80.000	72.439	9.5	94	0.00
46		1,1'-Biphenyl	40.000	36.945	7.6	96	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121821\
Data File : BP008488.D
Acq On : 18 Dec 2021 14:24
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Dec 20 04:19:22 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP121421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Dec 18 06:49:58 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)
47		2-Chloronaphthalene	40.000	36.666	8.3	96	0.00
48		2-Nitroaniline	40.000	40.804	-2.0	104	0.00
49		Acenaphthylene	40.000	38.452	3.9	100	0.00
50		Dimethylphthalate	40.000	41.317	-3.3	108	-0.01
51		2,6-Dinitrotoluene	40.000	41.326	-3.3	108	0.00
52	C	Acenaphthene	40.000	38.775	3.1	100	-0.01
53		3-Nitroaniline	40.000	43.429	-8.6	112	0.00
54	P	2,4-Dinitrophenol	40.000	46.013	-15.0	115	0.00
55		Dibenzofuran	40.000	39.370	1.6	104	0.00
56	P	4-Nitrophenol	40.000	42.521	-6.3	105	0.00
57		2,4-Dinitrotoluene	40.000	44.044	-10.1	116	0.00
58		Fluorene	40.000	40.394	-1.0	107	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	41.119	-2.8	108	0.00
60		Diethylphthalate	40.000	43.892	-9.7	117	0.00
61		4-Chlorophenyl-phenylether	40.000	39.850	0.4	105	0.00
62		4-Nitroaniline	40.000	46.744	-16.9	122	0.00
63		Azobenzene	40.000	43.003	-7.5	112	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	125	0.00
65		4,6-Dinitro-2-methylphenol	40.000	43.953	-9.9	125	0.00
66	c	n-Nitrosodiphenylamine	40.000	37.191	7.0	112	0.00
67		4-Bromophenyl-phenylether	40.000	36.493	8.8	110	0.00
68		Hexachlorobenzene	40.000	36.500	8.8	111	0.00
69		Atrazine	40.000	42.708	-6.8	125	0.00
70	C	Pentachlorophenol	40.000	40.749	-1.9	117	0.00
71		Phenanthrene	40.000	38.684	3.3	120	0.00
72		Anthracene	40.000	39.212	2.0	121	0.00
73		Carbazole	40.000	40.847	-2.1	128	0.00
74		Di-n-butylphthalate	40.000	44.994	-12.5	141	0.00
75	C	Fluoranthene	40.000	43.281	-8.2	140	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	168	0.00
77		Benzidine	40.000	48.647	-21.6	184	0.00
78		Pyrene	40.000	33.977	15.1	143	0.00
79	S	Terphenyl-d14	80.000	74.634	6.7	150	0.00
80		Butylbenzylphthalate	40.000	41.130	-2.8	171	0.00
81		Benzo(a)anthracene	40.000	38.251	4.4	158	0.00
82		3,3'-Dichlorobenzidine	40.000	39.684	0.8	165	0.00
83		Chrysene	40.000	38.181	4.5	160	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	43.610	-9.0	180	0.00
85	c	Di-n-octyl phthalate	40.000	43.933	-9.8	182	0.00
86	I	Perylene-d12	20.000	20.000	0.0	156	-0.01
87		Indeno(1,2,3-cd)pyrene	40.000	33.596	16.0	127	-0.01
88		Benzo(b)fluoranthene	40.000	39.967	0.1	154	-0.01
89		Benzo(k)fluoranthene	40.000	41.493	-3.7	159	0.00
90	C	Benzo(a)pyrene	40.000	39.494	1.3	152	-0.01
91		Dibenzo(a,h)anthracene	40.000	33.474	16.3	126	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121821\
Data File : BP008488.D
Acq On : 18 Dec 2021 14:24
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Dec 20 04:19:22 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP121421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Dec 18 06:49:58 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)
92	Benzo(g,h,i)perylene	40.000	31.652	20.9	119	-0.02

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