Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032521\
Data File : BM029214.D
Acq On : 24 Mar 2021 21:27
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM032521

Quant Time: Mar 25 03:12:22 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 25 02:53:41 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	111	0.00
2		1,4-Dioxane	40.000	37.653	5.9	108	0.00
3		Pyridine	40.000	40.226	-0.6	109	0.00
4		n-Nitrosodimethylamine	40.000	39.427	1.4	104	0.00
5	S	2-Fluorophenol	80.000	80.508	-0.6	111	0.00
6		Aniline	40.000	40.362	-0.9	110	0.00
7	S	Phenol-d6	80.000	81.232	-1.5	110	0.00
8		2-Chlorophenol	40.000	40.278	-0.7	111	0.00
9		Benzaldehyde	40.000	38.768	3.1	109	0.00
10	C	Phenol	40.000	40.212	-0.5	110	0.00
11		bis(2-Chloroethyl)ether	40.000	39.877	0.3	111	0.00
12		1,3-Dichlorobenzene	40.000	39.273	1.8	111	0.00
13	C	1,4-Dichlorobenzene	40.000	39.598	1.0	111	0.00
14		1,2-Dichlorobenzene	40.000	39.149	2.1	111	0.00
15		Benzyl Alcohol	40.000	40.804	-2.0	111	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.316	4.2	105	0.00
17		2-Methylphenol	40.000	41.782	-4.5	113	0.00
18		Hexachloroethane	40.000	39.922	0.2	110	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	40.638	-1.6	110	0.00
20		3+4-Methylphenols	40.000	41.302	-3.3	112	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	112	0.00
22		Acetophenone	40.000	39.436	1.4	109	0.00
23	S	Nitrobenzene-d5	80.000	80.880	-1.1	110	0.00
24		Nitrobenzene	40.000	40.164	-0.4	111	0.00
25		Isophorone	40.000	41.548	-3.9	112	0.00
26	C	2-Nitrophenol	40.000	42.296	-5.7	118	0.00
27		2,4-Dimethylphenol	40.000	40.905	-2.3	112	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.554	1.1	111	0.00
29	C	2,4-Dichlorophenol	40.000	42.018	-5.0	114	0.00
30		1,2,4-Trichlorobenzene	40.000	39.861	0.3	111	0.00
31		Naphthalene	40.000	39.666	0.8	111	0.00
32		Benzoic acid	40.000	42.369	-5.9	124	0.00
33		4-Chloroaniline	40.000	40.817	-2.0	112	0.00
34	C	Hexachlorobutadiene	40.000	39.886	0.3	115	0.00
35		Caprolactam	40.000	42.992	-7.5	121	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.397	-6.0	116	0.00
37		2-Methylnaphthalene	40.000	40.323	-0.8	113	0.00
38		1-Methylnaphthalene	40.000	40.092	-0.2	112	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	114	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.137	-0.3	114	0.00
41	P	Hexachlorocyclopentadiene	40.000	41.777	-4.4	117	0.00
42	S	2,4,6-Tribromophenol	80.000	83.804	-4.8	116	0.00
43	C	2,4,6-Trichlorophenol	40.000	42.351	-5.9	116	0.00
44		2,4,5-Trichlorophenol	40.000	42.729	-6.8	118	0.00
45	S	2-Fluorobiphenyl	80.000	78.935	1.3	113	0.00
46		1,1'-Biphenyl	40.000	39.021	2.4	111	0.00
47		2-Chloronaphthalene	40.000	39.309	1.7	111	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032521\
Data File : BM029214.D
Acq On : 24 Mar 2021 21:27
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM032521

Quant Time: Mar 25 03:12:22 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 25 02:53:41 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	39.650	0.9	116	0.00
49		Acenaphthylene	40.000	40.743	-1.9	114	0.00
50		Dimethylphthalate	40.000	40.386	-1.0	115	0.00
51		2,6-Dinitrotoluene	40.000	42.685	-6.7	116	0.00
52	C	Acenaphthene	40.000	39.640	0.9	113	0.00
53		3-Nitroaniline	40.000	44.930	-12.3	118	0.00
54	P	2,4-Dinitrophenol	40.000	45.127	-12.8	125	0.00
55		Dibenzofuran	40.000	39.396	1.5	113	0.00
56	P	4-Nitrophenol	40.000	42.078	-5.2	119	0.00
57		2,4-Dinitrotoluene	40.000	41.417	-3.5	118	0.00
58		Fluorene	40.000	40.335	-0.8	115	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	42.509	-6.3	119	0.00
60		Diethylphthalate	40.000	40.852	-2.1	115	0.00
61		4-Chlorophenyl-phenylether	40.000	40.076	-0.2	115	0.00
62		4-Nitroaniline	40.000	41.220	-3.0	119	0.00
63		Azobenzene	40.000	40.325	-0.8	111	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	116	0.00
65		4,6-Dinitro-2-methylphenol	40.000	44.063	-10.2	124	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.464	-1.2	115	0.00
67		4-Bromophenyl-phenylether	40.000	40.705	-1.8	117	0.00
68		Hexachlorobenzene	40.000	39.769	0.6	114	0.00
69		Atrazine	40.000	41.947	-4.9	117	0.00
70	C	Pentachlorophenol	40.000	41.709	-4.3	122	0.00
71		Phenanthrene	40.000	39.644	0.9	115	0.00
72		Anthracene	40.000	40.695	-1.7	115	0.00
73		Carbazole	40.000	40.946	-2.4	115	0.00
74		Di-n-butylphthalate	40.000	42.484	-6.2	117	0.00
75	C	Fluoranthene	40.000	40.635	-1.6	117	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	117	0.00
77		Benzidine	40.000	45.930	-14.8	129	0.00
78		Pyrene	40.000	40.997	-2.5	117	0.00
79	S	Terphenyl-d14	80.000	80.712	-0.9	116	0.00
80		Butylbenzylphthalate	40.000	41.051	-2.6	121	0.00
81		Benzo(a)anthracene	40.000	40.681	-1.7	117	0.00
82		3,3'-Dichlorobenzidine	40.000	42.885	-7.2	119	0.00
83		Chrysene	40.000	40.126	-0.3	118	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.022	-2.6	119	0.00
85	c	Di-n-octyl phthalate	40.000	44.184	-10.5	121	0.00
86	I	Perylene-d12	20.000	20.000	0.0	118	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	40.654	-1.6	117	0.00
88		Benzo(b)fluoranthene	40.000	41.202	-3.0	120	0.00
89		Benzo(k)fluoranthene	40.000	39.844	0.4	115	0.00
90	C	Benzo(a)pyrene	40.000	41.170	-2.9	118	0.00
91		Dibenzo(a,h)anthracene	40.000	40.320	-0.8	117	0.00
92		Benzo(g,h,i)perylene	40.000	40.057	-0.1	117	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032521\
Data File : BM029214.D
Acq On : 24 Mar 2021 21:27
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM032521

Quant Time: Mar 25 03:12:22 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 25 02:53:41 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