Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082521\
Data File : BM031898.D
Acq On : 25 Aug 2021 20:01
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 26 02:54:18 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM082521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 25 14:16:52 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	92	0.00
2		1,4-Dioxane	0.493	0.422	14.4	91	0.00
3		Pyridine	1.231	1.135	7.8	95	0.00
4		n-Nitrosodimethylamine	0.836	0.848	-1.4	100	0.00
5	S	2-Fluorophenol	1.210	1.098	9.3	91	0.00
6		Aniline	1.818	1.643	9.6	89	0.00
7	S	Phenol-d6	1.712	1.554	9.2	90	0.00
8		2-Chlorophenol	1.408	1.301	7.6	91	0.00
9		Benzaldehyde	1.135	0.983	13.4	94	0.00
10	C	Phenol	1.693	1.538	9.2	90	0.00
11		bis(2-Chloroethyl)ether	1.235	1.117	9.6	89	0.00
12		1,3-Dichlorobenzene	1.556	1.440	7.5	93	0.00
13	C	1,4-Dichlorobenzene	1.594	1.478	7.3	94	0.00
14		1,2-Dichlorobenzene	1.552	1.434	7.6	93	0.00
15		Benzyl Alcohol	1.460	1.413	3.2	93	0.00
16		2,2'-oxybis(1-Chloropropane	1.657	1.531	7.6	93	0.00
17		2-Methylphenol	1.208	1.124	7.0	92	0.00
18		Hexachloroethane	0.660	0.635	3.8	95	0.00
19	P	n-Nitroso-di-n-propylamine	1.319	1.260	4.5	95	0.00
20		3+4-Methylphenols	1.661	1.604	3.4	93	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
22		Acetophenone	0.544	0.494	9.2	94	0.00
23	S	Nitrobenzene-d5	0.462	0.423	8.4	94	0.00
24		Nitrobenzene	0.468	0.418	10.7	95	0.00
25		Isophorone	0.797	0.716	10.2	93	0.00
26	C	2-Nitrophenol	0.187	0.174	7.0	94	0.00
27		2,4-Dimethylphenol	0.282	0.258	8.5	93	0.00
28		bis(2-Chloroethoxy)methane	0.389	0.351	9.8	93	0.00
29	C	2,4-Dichlorophenol	0.331	0.311	6.0	96	0.00
30		1,2,4-Trichlorobenzene	0.363	0.341	6.1	98	0.00
31		Naphthalene	1.118	1.030	7.9	97	0.00
32		Benzoic acid	0.240	0.225	6.2	89	0.01
33		4-Chloroaniline	0.447	0.417	6.7	95	0.00
34	C	Hexachlorobutadiene	0.240	0.226	5.8	99	0.00
35		Caprolactam	0.101	0.097	4.0	94	0.01
36	C	4-Chloro-3-methylphenol	0.392	0.374	4.6	97	0.00
37		2-Methylnaphthalene	0.832	0.794	4.6	99	0.00
38		1-Methylnaphthalene	0.787	0.757	3.8	99	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
40		1,2,4,5-Tetrachlorobenzene	0.586	0.520	11.3	99	0.00
41	P	Hexachlorocyclopentadiene	0.286	0.273	4.5	104	0.00
42	S	2,4,6-Tribromophenol	0.236	0.227	3.8	103	0.00
43	C	2,4,6-Trichlorophenol	0.421	0.378	10.2	97	0.00
44		2,4,5-Trichlorophenol	0.455	0.411	9.7	97	0.00
45	S	2-Fluorobiphenyl	1.516	1.377	9.2	101	0.00
46		1,1'-Biphenyl	1.548	1.400	9.6	100	0.00
47		2-Chloronaphthalene	1.219	1.096	10.1	100	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082521\
Data File : BM031898.D
Acq On : 25 Aug 2021 20:01
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 26 02:54:18 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM082521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 25 14:16:52 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.424	0.402	5.2	100	0.00
49		Acenaphthylene	1.960	1.806	7.9	101	0.00
50		Dimethylphthalate	1.585	1.440	9.1	99	0.00
51		2,6-Dinitrotoluene	0.349	0.322	7.7	98	0.00
52	C	Acenaphthene	1.194	1.099	8.0	101	0.00
53		3-Nitroaniline	0.340	0.320	5.9	97	0.00
54	P	2,4-Dinitrophenol	0.198	0.197	0.5	100	0.00
55		Dibenzofuran	2.010	1.865	7.2	101	0.00
56	P	4-Nitrophenol	0.268	0.269	-0.4	100	0.00
57		2,4-Dinitrotoluene	0.497	0.476	4.2	100	0.00
58		Fluorene	1.648	1.568	4.9	103	0.00
59		2,3,4,6-Tetrachlorophenol	0.397	0.386	2.8	104	0.00
60		Diethylphthalate	1.730	1.634	5.5	102	0.00
61		4-Chlorophenyl-phenylether	0.818	0.790	3.4	104	0.00
62		4-Nitroaniline	0.328	0.324	1.2	99	0.00
63		Azobenzene	1.908	1.854	2.8	104	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
65		4,6-Dinitro-2-methylphenol	0.138	0.130	5.8	100	0.00
66	c	n-Nitrosodiphenylamine	0.667	0.601	9.9	102	0.00
67		4-Bromophenyl-phenylether	0.223	0.203	9.0	102	0.00
68		Hexachlorobenzene	0.236	0.215	8.9	104	0.00
69		Atrazine	0.225	0.210	6.7	101	0.00
70	C	Pentachlorophenol	0.148	0.143	3.4	101	0.00
71		Phenanthrene	1.178	1.083	8.1	102	0.00
72		Anthracene	1.152	1.073	6.9	103	0.00
73		Carbazole	1.088	1.012	7.0	101	0.00
74		Di-n-butylphthalate	1.365	1.302	4.6	100	0.00
75	C	Fluoranthene	1.286	1.187	7.7	98	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	92	0.00
77		Benzidine	0.523	0.565	-8.0	95	0.00
78		Pyrene	1.673	1.622	3.0	98	0.00
79	S	Terphenyl-d14	1.386	1.373	0.9	98	0.00
80		Butylbenzylphthalate	0.719	0.698	2.9	92	0.00
81		Benzo(a)anthracene	1.439	1.335	7.2	92	0.00
82		3,3'-Dichlorobenzidine	0.414	0.395	4.6	93	0.00
83		Chrysene	1.387	1.276	8.0	92	0.00
84		Bis(2-ethylhexyl)phthalate	1.035	0.986	4.7	88	0.00
85	c	Di-n-octyl phthalate	1.560	1.400	10.3	87	0.00
86	I	Perylene-d12	1.000	1.000	0.0	94	0.00
87		Indeno(1,2,3-cd)pyrene	1.384	1.276	7.8	102	0.00
88		Benzo(b)fluoranthene	1.514	1.448	4.4	95	0.00
89		Benzo(k)fluoranthene	1.435	1.303	9.2	91	0.00
90	C	Benzo(a)pyrene	1.312	1.214	7.5	95	0.00
91		Dibenzo(a,h)anthracene	1.199	1.112	7.3	102	0.00
92		Benzo(g,h,i)perylene	1.133	1.049	7.4	103	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082521\
Data File : BM031898.D
Acq On : 25 Aug 2021 20:01
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Aug 26 02:54:18 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM082521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 25 14:16:52 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	AvgRF	CCRF	%Dev Area%	Dev(min)

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