Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090921\
Data File : BM032105.D
Acq On : 09 Sep 2021 10:06
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Sep 09 12:29:55 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM090821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 08 19:23:17 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	76	0.00
2		1,4-Dioxane	0.507	0.502	1.0	82	0.00
3		Pyridine	1.323	1.376	-4.0	82	0.00
4		n-Nitrosodimethylamine	0.690	0.790	-14.5	90	0.00
5	S	2-Fluorophenol	1.121	1.080	3.7	77	0.00
6		Aniline	1.886	1.899	-0.7	79	0.00
7	S	Phenol-d6	1.705	1.689	0.9	79	0.00
8		2-Chlorophenol	1.252	1.223	2.3	76	0.00
9		Benzaldehyde	1.109	1.010	8.9	78	0.00
10	C	Phenol	1.729	1.715	0.8	78	0.00
11		bis(2-Chloroethyl)ether	1.305	1.320	-1.1	80	0.00
12		1,3-Dichlorobenzene	1.548	1.486	4.0	78	0.00
13	C	1,4-Dichlorobenzene	1.568	1.521	3.0	78	0.00
14		1,2-Dichlorobenzene	1.533	1.486	3.1	78	0.00
15		Benzyl Alcohol	1.342	1.453	-8.3	84	0.00
16		2,2'-oxybis(1-Chloropropane	1.879	1.984	-5.6	85	0.00
17		2-Methylphenol	1.158	1.168	-0.9	79	-0.01
18		Hexachloroethane	0.590	0.609	-3.2	82	0.00
19	P	n-Nitroso-di-n-propylamine	1.247	1.362	-9.2	86	0.00
20		3+4-Methylphenols	1.534	1.603	-4.5	81	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	83	0.00
22		Acetophenone	0.576	0.540	6.2	83	0.00
23	S	Nitrobenzene-d5	0.449	0.448	0.2	86	0.00
24		Nitrobenzene	0.478	0.472	1.3	86	0.00
25		Isophorone	0.833	0.810	2.8	86	0.00
26	C	2-Nitrophenol	0.092	0.101	-9.8	88	0.00
27		2,4-Dimethylphenol	0.278	0.262	5.8	84	0.00
28		bis(2-Chloroethoxy)methane	0.417	0.402	3.6	85	0.00
29	C	2,4-Dichlorophenol	0.302	0.295	2.3	82	0.00
30		1,2,4-Trichlorobenzene	0.410	0.385	6.1	84	0.00
31		Naphthalene	1.119	1.043	6.8	83	0.00
32		Benzoic acid	0.096	0.093	3.1	81	0.00
33		4-Chloroaniline	0.430	0.413	4.0	84	0.00
34	C	Hexachlorobutadiene	0.275	0.262	4.7	85	0.00
35		Caprolactam	0.083	0.086	-3.6	85	0.00
36	C	4-Chloro-3-methylphenol	0.354	0.360	-1.7	88	0.00
37		2-Methylnaphthalene	0.810	0.788	2.7	88	0.00
38		1-Methylnaphthalene	0.766	0.740	3.4	86	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	94	0.00
40		1,2,4,5-Tetrachlorobenzene	0.669	0.586	12.4	87	0.00
41	P	Hexachlorocyclopentadiene	0.206	0.191	7.3	90	0.00
42	S	2,4,6-Tribromophenol	0.159	0.166	-4.4	106	0.00
43	C	2,4,6-Trichlorophenol	0.272	0.258	5.1	93	0.00
44		2,4,5-Trichlorophenol	0.331	0.334	-0.9	93	0.00
45	S	2-Fluorobiphenyl	1.525	1.371	10.1	90	0.00
46		1,1'-Biphenyl	1.526	1.370	10.2	89	0.00
47		2-Chloronaphthalene	1.237	1.112	10.1	90	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090921\
Data File : BM032105.D
Acq On : 09 Sep 2021 10:06
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Sep 09 12:29:55 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM090821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 08 19:23:17 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	0.314	0.345	-9.9	102	0.00
49		Acenaphthylene	1.855	1.729	6.8	92	0.00
50		Dimethylphthalate	1.452	1.361	6.3	94	0.00
51		2,6-Dinitrotoluene	0.295	0.293	0.7	93	0.00
52	C	Acenaphthene	1.153	1.069	7.3	93	0.00
53		3-Nitroaniline	0.258	0.260	-0.8	94	0.00
54	P	2,4-Dinitrophenol	0.084	0.076	9.5	90	0.00
55		Dibenzofuran	1.922	1.812	5.7	95	0.00
56	P	4-Nitrophenol	0.144	0.137	4.9	97	0.00
57		2,4-Dinitrotoluene	0.362	0.400	-10.5	101	0.00
58		Fluorene	1.510	1.445	4.3	96	0.00
59		2,3,4,6-Tetrachlorophenol	0.242	0.250	-3.3	109	0.00
60		Diethylphthalate	1.410	1.390	1.4	98	0.00
61		4-Chlorophenyl-phenylether	0.811	0.778	4.1	97	0.00
62		4-Nitroaniline	0.227	0.248	-9.3	104	0.00
63		Azobenzene	1.649	1.707	-3.5	102	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
65		4,6-Dinitro-2-methylphenol	0.067	0.062	7.5	97	0.00
66	c	n-Nitrosodiphenylamine	0.655	0.598	8.7	96	0.00
67		4-Bromophenyl-phenylether	0.235	0.217	7.7	97	0.00
68		Hexachlorobenzene	0.266	0.238	10.5	96	0.00
69		Atrazine	0.201	0.202	-0.5	100	0.00
70	C	Pentachlorophenol	0.087	0.079	9.2	90	0.00
71		Phenanthrene	1.144	1.061	7.3	99	0.00
72		Anthracene	1.118	1.047	6.4	99	0.00
73		Carbazole	1.027	0.966	5.9	99	0.00
74		Di-n-butylphthalate	1.068	1.083	-1.4	104	0.00
75	C	Fluoranthene	1.275	1.205	5.5	100	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
77		Benzidine	0.368	0.368	0.0	114	0.00
78		Pyrene	1.524	1.421	6.8	102	0.00
79	S	Terphenyl-d14	1.226	1.164	5.1	103	0.00
80		Butylbenzylphthalate	0.447	0.475	-6.3	108	0.00
81		Benzo(a)anthracene	1.340	1.247	6.9	101	0.00
82		3,3'-Dichlorobenzidine	0.369	0.357	3.3	101	0.00
83		Chrysene	1.422	1.327	6.7	104	0.00
84		Bis(2-ethylhexyl)phthalate	0.681	0.723	-6.2	107	0.00
85	c	Di-n-octyl phthalate	1.150	1.146	0.3	107	0.00
86	I	Perylene-d12	1.000	1.000	0.0	103	0.00
87		Indeno(1,2,3-cd)pyrene	1.356	1.250	7.8	99	0.00
88		Benzo(b)fluoranthene	1.338	1.213	9.3	98	0.00
89		Benzo(k)fluoranthene	1.447	1.397	3.5	104	0.00
90	C	Benzo(a)pyrene	1.229	1.185	3.6	108	0.00
91		Dibenzo(a,h)anthracene	1.169	1.108	5.2	102	0.00
92		Benzo(g,h,i)perylene	1.158	1.075	7.2	101	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090921\
Data File : BM032105.D
Acq On : 09 Sep 2021 10:06
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Sep 09 12:29:55 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM090821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 08 19:23:17 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