Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080222\
Data File : BM036043.D
Acq On : 02 Aug 2022 09:34
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Aug 03 00:50:54 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 01 17:50:14 2022
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	82	0.00
2		1,4-Dioxane	0.497	0.460	7.4	85	0.00
3		Pyridine	1.333	1.269	4.8	90	0.00
4		n-Nitrosodimethylamine	0.548	0.522	4.7	88	0.00
5	S	2-Fluorophenol	1.234	1.147	7.1	84	0.00
6		Aniline	1.738	1.655	4.8	87	0.00
7	S	Phenol-d6	1.570	1.478	5.9	86	0.00
8		2-Chlorophenol	1.315	1.218	7.4	83	0.00
9		Benzaldehyde	0.831	0.696	16.2	86	0.00
10	C	Phenol	1.580	1.493	5.5	87	0.00
11		bis(2-Chloroethyl)ether	1.302	1.216	6.6	83	0.00
12		1,3-Dichlorobenzene	1.458	1.335	8.4	82	0.00
13	C	1,4-Dichlorobenzene	1.486	1.370	7.8	82	0.00
14		1,2-Dichlorobenzene	1.434	1.332	7.1	82	0.00
15		Benzyl Alcohol	1.053	0.979	7.0	84	0.00
16		2,2'-oxybis(1-Chloropropane	1.731	1.623	6.2	82	0.00
17		2-Methylphenol	1.046	0.971	7.2	85	0.00
18		Hexachloroethane	0.535	0.495	7.5	80	0.00
19	P	n-Nitroso-di-n-propylamine	0.953	0.918	3.7	86	0.00
20		3+4-Methylphenols	1.421	1.339	5.8	86	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	84	0.00
22		Acetophenone	0.473	0.421	11.0	84	0.00
23	S	Nitrobenzene-d5	0.357	0.322	9.8	86	0.00
24		Nitrobenzene	0.336	0.302	10.1	85	0.00
25		Isophorone	0.581	0.517	11.0	84	0.00
26	C	2-Nitrophenol	0.158	0.138	12.7	81	0.00
27		2,4-Dimethylphenol	0.245	0.220	10.2	85	0.00
28		bis(2-Chloroethoxy)methane	0.414	0.369	10.9	83	0.00
29	C	2,4-Dichlorophenol	0.276	0.247	10.5	84	0.00
30		1,2,4-Trichlorobenzene	0.313	0.275	12.1	82	0.00
31		Naphthalene	0.994	0.884	11.1	85	0.00
32		Benzoic acid	0.195	0.167	14.4	81	0.00
33		4-Chloroaniline	0.401	0.359	10.5	86	0.00
34	C	Hexachlorobutadiene	0.184	0.159	13.6	79	0.00
35		Caprolactam	0.085	0.076	10.6	89	-0.01
36	C	4-Chloro-3-methylphenol	0.290	0.259	10.7	86	0.00
37		2-Methylnaphthalene	0.662	0.589	11.0	84	0.00
38		1-Methylnaphthalene	0.648	0.577	11.0	85	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	87	0.00
40		1,2,4,5-Tetrachlorobenzene	0.542	0.467	13.8	82	0.00
41	P	Hexachlorocyclopentadiene	0.287	0.246	14.3	78	0.00
42	S	2,4,6-Tribromophenol	0.235	0.211	10.2	89	0.00
43	C	2,4,6-Trichlorophenol	0.369	0.323	12.5	84	0.00
44		2,4,5-Trichlorophenol	0.389	0.346	11.1	86	0.00
45	S	2-Fluorobiphenyl	1.355	1.189	12.3	85	0.00
46		1,1'-Biphenyl	1.458	1.268	13.0	85	0.00
47		2-Chloronaphthalene	1.116	0.966	13.4	85	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.304	0.277	8.9	92	0.00
49		Acenaphthylene	1.724	1.516	12.1	87	0.00
50		Dimethylphthalate	1.360	1.181	13.2	87	0.00
51		2,6-Dinitrotoluene	0.273	0.248	9.2	89	0.00
52	C	Acenaphthene	1.128	0.991	12.1	88	0.00
53		3-Nitroaniline	0.318	0.295	7.2	94	0.00
54	P	2,4-Dinitrophenol	0.144	0.130	9.7	90	0.00
55		Dibenzofuran	1.689	1.491	11.7	88	0.00
56	P	4-Nitrophenol	0.255	0.241	5.5	97	0.00
57		2,4-Dinitrotoluene	0.358	0.333	7.0	92	0.00
58		Fluorene	1.332	1.175	11.8	89	0.00
59		2,3,4,6-Tetrachlorophenol	0.330	0.295	10.6	89	0.00
60		Diethylphthalate	1.398	1.230	12.0	87	0.00
61		4-Chlorophenyl-phenylether	0.664	0.581	12.5	86	0.00
62		4-Nitroaniline	0.328	0.314	4.3	97	0.00
63		Azobenzene	1.329	1.200	9.7	87	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	93	0.00
65		4,6-Dinitro-2-methylphenol	0.102	0.092	9.8	91	0.00
66	c	n-Nitrosodiphenylamine	0.565	0.485	14.2	89	0.00
67		4-Bromophenyl-phenylether	0.200	0.168	16.0	87	0.00
68		Hexachlorobenzene	0.230	0.196	14.8	89	0.00
69		Atrazine	0.158	0.147	7.0	91	0.00
70	C	Pentachlorophenol	0.134	0.120	10.4	92	0.00
71		Phenanthrene	1.019	0.890	12.7	93	0.00
72		Anthracene	0.986	0.871	11.7	93	0.00
73		Carbazole	1.000	0.906	9.4	98	0.00
74		Di-n-butylphthalate	1.193	1.056	11.5	90	0.00
75	C	Fluoranthene	1.141	1.037	9.1	98	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
77		Benzidine	0.300	0.347	-15.7	123	0.00
78		Pyrene	1.240	1.044	15.8	99	0.00
79	S	Terphenyl-d14	1.014	0.863	14.9	96	0.00
80		Butylbenzylphthalate	0.535	0.454	15.1	94	0.00
81		Benzo(a)anthracene	1.232	1.088	11.7	103	0.00
82		3,3'-Dichlorobenzidine	0.299	0.272	9.0	104	0.00
83		Chrysene	1.171	1.031	12.0	103	0.00
84		Bis(2-ethylhexyl)phthalate	0.808	0.688	14.9	93	0.00
85	c	Di-n-octyl phthalate	1.318	1.132	14.1	93	0.00
86	I	Perylene-d12	1.000	1.000	0.0	102	0.00
87		Indeno(1,2,3-cd)pyrene	1.405	1.225	12.8	102	-0.01
88		Benzo(b)fluoranthene	1.172	1.017	13.2	101	0.00
89		Benzo(k)fluoranthene	1.184	1.056	10.8	104	0.00
90	C	Benzo(a)pyrene	0.983	0.864	12.1	103	0.00
91		Dibenzo(a,h)anthracene	1.176	1.024	12.9	102	0.00
92		Benzo(g,h,i)perylene	1.139	0.999	12.3	103	0.00

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