

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080321\
 Data File : BG049631.D
 Acq On : 3 Aug 2021 19:53
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 04 02:23:36 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 03 15:15:33 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	102	0.00
2	1,4-Dioxane	0.488	0.498	-2.0	101	0.00
3	Pyridine	1.336	1.373	-2.8	97	0.00
4	n-Nitrosodimethylamine	0.719	0.780	-8.5	105	0.00
5 S	2-Fluorophenol	1.116	1.134	-1.6	96	0.00
6	Aniline	1.790	1.840	-2.8	98	0.00
7 S	Phenol-d6	1.689	1.735	-2.7	97	0.00
8	2-Chlorophenol	1.216	1.239	-1.9	100	0.00
9	Benzaldehyde	1.127	1.090	3.3	95	0.00
10 C	Phenol	1.636	1.656	-1.2	95	0.00
11	bis(2-Chloroethyl)ether	1.105	1.137	-2.9	99	0.00
12	1,3-Dichlorobenzene	1.393	1.408	-1.1	100	0.00
13 C	1,4-Dichlorobenzene	1.410	1.444	-2.4	100	0.00
14	1,2-Dichlorobenzene	1.336	1.329	0.5	95	0.00
15	Benzyl Alcohol	1.418	1.467	-3.5	97	0.00
16	2,2'-oxybis(1-Chloropropane	1.255	1.249	0.5	95	0.00
17	2-Methylphenol	1.078	1.137	-5.5	99	0.00
18	Hexachloroethane	0.542	0.542	0.0	94	0.00
19 P	n-Nitroso-di-n-propylamine	1.130	1.141	-1.0	95	0.00
20	3+4-Methylphenols	1.513	1.555	-2.8	98	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
22	Acetophenone	0.543	0.556	-2.4	96	0.00
23 S	Nitrobenzene-d5	0.452	0.477	-5.5	98	0.00
24	Nitrobenzene	0.477	0.505	-5.9	98	0.00
25	Isophorone	0.807	0.857	-6.2	99	0.00
26 C	2-Nitrophenol	0.182	0.197	-8.2	99	0.00
27	2,4-Dimethylphenol	0.269	0.290	-7.8	100	0.00
28	bis(2-Chloroethoxy)methane	0.353	0.371	-5.1	98	0.00
29 C	2,4-Dichlorophenol	0.330	0.350	-6.1	99	0.00
30	1,2,4-Trichlorobenzene	0.363	0.375	-3.3	98	0.00
31	Naphthalene	1.047	1.084	-3.5	98	0.00
32	Benzoic acid	0.174	0.188	-8.0	95	0.00
33	4-Chloroaniline	0.414	0.441	-6.5	98	0.00
34 C	Hexachlorobutadiene	0.264	0.272	-3.0	98	0.00
35	Caprolactam	0.102	0.113	-10.8	105	0.00
36 C	4-Chloro-3-methylphenol	0.381	0.417	-9.4	101	0.00
37	2-Methylnaphthalene	0.789	0.833	-5.6	97	0.00
38	1-Methylnaphthalene	0.742	0.778	-4.9	99	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
40	1,2,4,5-Tetrachlorobenzene	0.589	0.592	-0.5	96	0.00
41 P	Hexachlorocyclopentadiene	0.243	0.305	-25.5#	98	0.00
42 S	2,4,6-Tribromophenol	0.294	0.313	-6.5	103	0.00
43 C	2,4,6-Trichlorophenol	0.397	0.424	-6.8	100	0.00
44	2,4,5-Trichlorophenol	0.431	0.451	-4.6	98	0.00
45 S	2-Fluorobiphenyl	1.374	1.407	-2.4	98	0.00
46	1,1'-Biphenyl	1.424	1.479	-3.9	100	0.00
47	2-Chloronaphthalene	1.138	1.178	-3.5	98	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080321\
 Data File : BG049631.D
 Acq On : 3 Aug 2021 19:53
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 04 02:23:36 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 03 15:15:33 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	0.418	0.443	-6.0	100	0.00
49	Acenaphthylene	1.803	1.888	-4.7	101	0.00
50	Dimethylphthalate	1.509	1.547	-2.5	100	0.00
51	2,6-Dinitrotoluene	0.332	0.335	-0.9	94	0.00
52 C	Acenaphthene	1.086	1.129	-4.0	100	0.00
53	3-Nitroaniline	0.324	0.343	-5.9	101	0.00
54 P	2,4-Dinitrophenol	0.189	0.202	-6.9	102	0.00
55	Dibenzofuran	1.762	1.836	-4.2	99	0.00
56 P	4-Nitrophenol	0.237	0.257	-8.4	98	0.00
57	2,4-Dinitrotoluene	0.467	0.487	-4.3	102	0.00
58	Fluorene	1.432	1.482	-3.5	100	0.00
59	2,3,4,6-Tetrachlorophenol	0.400	0.432	-8.0	105	0.00
60	Diethylphthalate	1.518	1.585	-4.4	102	0.00
61	4-Chlorophenyl-phenylether	0.764	0.785	-2.7	99	0.00
62	4-Nitroaniline	0.333	0.356	-6.9	100	0.00
63	Azobenzene	1.606	1.655	-3.1	100	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	104	0.00
65	4,6-Dinitro-2-methylphenol	0.129	0.134	-3.9	101	0.00
66 c	n-Nitrosodiphenylamine	0.565	0.573	-1.4	100	0.00
67	4-Bromophenyl-phenylether	0.236	0.239	-1.3	101	0.00
68	Hexachlorobenzene	0.262	0.265	-1.1	101	0.00
69	Atrazine	0.226	0.235	-4.0	103	0.00
70 C	Pentachlorophenol	0.140	0.146	-4.3	102	0.00
71	Phenanthrene	1.052	1.065	-1.2	101	0.00
72	Anthracene	1.033	1.047	-1.4	102	0.00
73	Carbazole	0.979	1.004	-2.6	102	0.00
74	Di-n-butylphthalate	1.141	1.170	-2.5	103	0.00
75 C	Fluoranthene	1.295	1.340	-3.5	104	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	104	0.00
77	Benzidine	0.512	0.547	-6.8	94	0.00
78	Pyrene	1.363	1.368	-0.4	101	0.00
79 S	Terphenyl-d14	1.101	1.128	-2.5	103	0.00
80	Butylbenzylphthalate	0.519	0.535	-3.1	103	0.00
81	Benzo(a)anthracene	1.302	1.341	-3.0	103	0.00
82	3,3'-Dichlorobenzidine	0.380	0.393	-3.4	101	0.00
83	Chrysene	1.245	1.276	-2.5	105	0.00
84	Bis(2-ethylhexyl)phthalate	0.735	0.746	-1.5	101	0.00
85 c	Di-n-octyl phthalate	1.242	1.301	-4.8	105	0.00
86 I	Perylene-d12	1.000	1.000	0.0	103	0.00
87	Indeno(1,2,3-cd)pyrene	1.442	1.496	-3.7	105	0.00
88	Benzo(b)fluoranthene	1.312	1.350	-2.9	105	0.00
89	Benzo(k)fluoranthene	1.224	1.263	-3.2	103	0.00
90 C	Benzo(a)pyrene	1.187	1.222	-2.9	104	0.00
91	Dibenzo(a,h)anthracene	1.215	1.267	-4.3	104	0.00
92	Benzo(g,h,i)perylene	1.191	1.224	-2.8	104	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080321\
Data File : BG049631.D
Acq On : 3 Aug 2021 19:53
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Aug 04 02:23:36 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 03 15:15:33 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