

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082421\
 Data File : BM031882.D
 Acq On : 24 Aug 2021 17:03
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 24 17:53:08 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM082321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 24 16:34:26 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	100	0.00
2	1,4-Dioxane	0.442	0.428	3.2	109	0.00
3	Pyridine	1.331	1.143	14.1	95	0.00
4	n-Nitrosodimethylamine	0.581	0.782	-34.6#	143	0.00
5 S	2-Fluorophenol	1.360	1.154	15.1	90	0.00
6	Aniline	1.943	1.763	9.3	95	0.00
7 S	Phenol-d6	1.826	1.665	8.8	96	0.00
8	2-Chlorophenol	1.458	1.372	5.9	99	-0.01
9	Benzaldehyde	1.207	1.021	15.4	93	-0.01
10 C	Phenol	1.824	1.644	9.9	93	-0.01
11	bis(2-Chloroethyl)ether	1.380	1.205	12.7	91	0.00
12	1,3-Dichlorobenzene	1.577	1.480	6.2	101	0.00
13 C	1,4-Dichlorobenzene	1.602	1.504	6.1	101	0.00
14	1,2-Dichlorobenzene	1.560	1.471	5.7	100	0.00
15	Benzyl Alcohol	1.501	1.456	3.0	99	0.00
16	2,2'-oxybis(1-Chloropropane	2.062	1.553	24.7	79	0.00
17	2-Methylphenol	1.228	1.159	5.6	97	0.00
18	Hexachloroethane	0.682	0.624	8.5	96	-0.01
19 P	n-Nitroso-di-n-propylamine	1.375	1.274	7.3	93	0.00
20	3+4-Methylphenols	1.713	1.622	5.3	96	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
22	Acetophenone	0.583	0.518	11.1	97	0.00
23 S	Nitrobenzene-d5	0.491	0.435	11.4	95	-0.01
24	Nitrobenzene	0.503	0.436	13.3	94	-0.01
25	Isophorone	0.881	0.769	12.7	94	0.00
26 C	2-Nitrophenol	0.190	0.184	3.2	102	0.00
27	2,4-Dimethylphenol	0.292	0.275	5.8	100	0.00
28	bis(2-Chloroethoxy)methane	0.428	0.370	13.6	93	0.00
29 C	2,4-Dichlorophenol	0.331	0.322	2.7	105	0.00
30	1,2,4-Trichlorobenzene	0.352	0.348	1.1	110	0.00
31	Naphthalene	1.147	1.048	8.6	100	0.00
32	Benzoic acid	0.248	0.250	-0.8	105	0.00
33	4-Chloroaniline	0.455	0.437	4.0	103	-0.01
34 C	Hexachlorobutadiene	0.220	0.226	-2.7	112	0.00
35	Caprolactam	0.105	0.106	-1.0	103	0.00
36 C	4-Chloro-3-methylphenol	0.393	0.384	2.3	103	-0.01
37	2-Methylnaphthalene	0.827	0.795	3.9	103	0.00
38	1-Methylnaphthalene	0.778	0.756	2.8	105	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	109	0.00
40	1,2,4,5-Tetrachlorobenzene	0.575	0.566	1.6	113	0.00
41 P	Hexachlorocyclopentadiene	0.279	0.277	0.7	113	0.00
42 S	2,4,6-Tribromophenol	0.220	0.242	-10.0	122	0.00
43 C	2,4,6-Trichlorophenol	0.424	0.418	1.4	113	0.00
44	2,4,5-Trichlorophenol	0.453	0.448	1.1	113	0.00
45 S	2-Fluorobiphenyl	1.554	1.442	7.2	107	0.00
46	1,1'-Biphenyl	1.613	1.490	7.6	108	0.00
47	2-Chloronaphthalene	1.253	1.162	7.3	107	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082421\
 Data File : BM031882.D
 Acq On : 24 Aug 2021 17:03
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 24 17:53:08 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM082321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 24 16:34:26 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	0.436	0.411	5.7	102	0.00
49	Acenaphthylene	2.032	1.891	6.9	107	0.00
50	Dimethylphthalate	1.625	1.550	4.6	109	0.00
51	2,6-Dinitrotoluene	0.354	0.344	2.8	108	0.00
52 C	Acenaphthene	1.223	1.137	7.0	107	0.00
53	3-Nitroaniline	0.347	0.348	-0.3	110	0.00
54 P	2,4-Dinitrophenol	0.198	0.215	-8.6	114	0.00
55	Dibenzofuran	2.024	1.931	4.6	110	0.00
56 P	4-Nitrophenol	0.286	0.292	-2.1	110	0.00
57	2,4-Dinitrotoluene	0.493	0.504	-2.2	110	0.00
58	Fluorene	1.658	1.601	3.4	109	0.00
59	2,3,4,6-Tetrachlorophenol	0.388	0.399	-2.8	117	0.00
60	Diethylphthalate	1.769	1.704	3.7	108	0.00
61	4-Chlorophenyl-phenylether	0.795	0.796	-0.1	115	0.00
62	4-Nitroaniline	0.330	0.348	-5.5	112	0.00
63	Azobenzene	2.040	1.816	11.0	100	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	115	0.00
65	4,6-Dinitro-2-methylphenol	0.139	0.143	-2.9	116	0.00
66 c	n-Nitrosodiphenylamine	0.675	0.628	7.0	113	0.00
67	4-Bromophenyl-phenylether	0.217	0.211	2.8	117	-0.01
68	Hexachlorobenzene	0.230	0.227	1.3	119	0.00
69	Atrazine	0.226	0.226	0.0	116	0.00
70 C	Pentachlorophenol	0.150	0.155	-3.3	121	0.00
71	Phenanthrene	1.205	1.137	5.6	114	0.00
72	Anthracene	1.185	1.126	5.0	113	0.00
73	Carbazole	1.140	1.086	4.7	114	0.00
74	Di-n-butylphthalate	1.518	1.402	7.6	108	0.00
75 C	Fluoranthene	1.339	1.329	0.7	119	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	128	0.00
77	Benzydine	0.536	0.530	1.1	117	0.00
78	Pyrene	1.603	1.498	6.6	118	0.00
79 S	Terphenyl-d14	1.329	1.287	3.2	121	0.00
80	Butylbenzylphthalate	0.790	0.717	9.2	115	0.00
81	Benzo(a)anthracene	1.466	1.397	4.7	127	0.00
82	3,3'-Dichlorobenzidine	0.434	0.436	-0.5	134	0.00
83	Chrysene	1.411	1.343	4.8	128	0.00
84	Bis(2-ethylhexyl)phthalate	1.229	1.100	10.5	114	0.00
85 c	Di-n-octyl phthalate	1.967	1.701	13.5	120	0.00
86 I	Perylene-d12	1.000	1.000	0.0	131	0.00
87	Indeno(1,2,3-cd)pyrene	1.553	1.258	19.0	118	0.00
88	Benzo(b)fluoranthene	1.473	1.524	-3.5	141	0.00
89	Benzo(k)fluoranthene	1.418	1.390	2.0	128	0.00
90 C	Benzo(a)pyrene	1.323	1.272	3.9	132	0.00
91	Dibenzo(a,h)anthracene	1.356	1.104	18.6	118	0.00
92	Benzo(g,h,i)perylene	1.262	0.998	20.9	116	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082421\
Data File : BM031882.D
Acq On : 24 Aug 2021 17:03
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Aug 24 17:53:08 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM082321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 24 16:34:26 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