

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102221\
 Data File : BG050704.D
 Acq On : 23 Oct 2021 1:07
 Operator : CG/JU
 Sample : M4298-01 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NB-08-102021

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG050704.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.800	385	393	410	rBV	411541	755940	53.34%	5.477%
2	7.398	656	665	678	rBV	450329	835054	58.93%	6.050%
3	8.255	804	811	822	rVB2	228250	405864	28.64%	2.940%
4	9.425	1002	1010	1023	rVB	291169	553910	39.09%	4.013%
5	10.465	1182	1187	1192	rBV6	8028	14238	1.00%	0.103%
6	10.823	1241	1248	1259	rBV	50077	100075	7.06%	0.725%
7	11.017	1276	1281	1285	rBV	23913	40629	2.87%	0.294%
8	11.082	1285	1292	1298	rVV	302277	554212	39.11%	4.015%
9	11.205	1306	1313	1320	rVB5	10488	24185	1.71%	0.175%
10	11.757	1399	1407	1412	rBV10	9172	21150	1.49%	0.153%
11	11.892	1423	1430	1435	rBV9	7279	17412	1.23%	0.126%
12	12.057	1453	1458	1463	rVB3	12450	22872	1.61%	0.166%
13	12.251	1484	1491	1497	rVB5	16240	30965	2.19%	0.224%
14	12.445	1518	1524	1531	rBV	49764	78406	5.53%	0.568%
15	12.721	1566	1571	1577	rVB	16484	26749	1.89%	0.194%
16	12.938	1602	1608	1609	rBV	17999	29193	2.06%	0.211%
17	13.244	1658	1660	1668	rVB7	10908	20097	1.42%	0.146%
18	13.385	1679	1684	1689	rBV5	17499	29315	2.07%	0.212%
19	13.496	1696	1703	1713	rVB	659606	1046333	73.84%	7.580%
20	13.673	1724	1733	1738	rBV	71191	119314	8.42%	0.864%
21	13.814	1753	1757	1762	rVB5	12532	20216	1.43%	0.146%
22	14.049	1787	1797	1803	rBV5	16766	50074	3.53%	0.363%
23	14.190	1816	1821	1826	rVB3	26749	42839	3.02%	0.310%
24	14.243	1826	1830	1834	rVB5	19257	29105	2.05%	0.211%
25	14.290	1834	1838	1841	rBV3	30364	47035	3.32%	0.341%
26	14.319	1841	1843	1847	rVB3	27967	34540	2.44%	0.250%
27	14.372	1847	1852	1856	rBV3	12284	22017	1.55%	0.160%
28	14.431	1858	1862	1870	rVB4	18379	35419	2.50%	0.257%
29	14.607	1886	1892	1896	rBV7	17702	40758	2.88%	0.295%
30	14.695	1903	1907	1910	rBV5	19077	26409	1.86%	0.191%
31	14.736	1910	1914	1920	rVB	71733	92559	6.53%	0.671%
32	14.871	1931	1937	1945	rVB	419067	695459	49.08%	5.038%
33	14.942	1945	1949	1955	rBV9	13084	27128	1.91%	0.197%
34	15.077	1966	1972	1976	rBV8	12311	26988	1.90%	0.196%
35	15.183	1984	1990	1993	rBV8	8872	18810	1.33%	0.136%
36	15.347	2012	2018	2021	rBV4	19800	34776	2.45%	0.252%

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 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
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37	15.412	2027	2029	2033	rVB4	11861	14261	1.01%	0.103%
38	15.482	2036	2041	2045	rBV6	16609	36041	2.54%	0.261%
39	15.576	2054	2057	2062	rVB6	11327	19336	1.36%	0.140%
40	15.653	2066	2070	2071	rVV3	18777	26085	1.84%	0.189%
41	15.682	2071	2075	2080	rVB	90686	118039	8.33%	0.855%
42	15.911	2111	2114	2120	rVB5	11685	17358	1.22%	0.126%
43	16.082	2138	2143	2148	rVB2	26774	45172	3.19%	0.327%
44	16.305	2177	2181	2184	rBV6	15008	19034	1.34%	0.138%
45	16.358	2184	2190	2198	rBV	366504	612772	43.24%	4.439%
46	16.540	2217	2221	2232	rBV	87379	190684	13.46%	1.381%
47	17.051	2304	2308	2314	rBV8	12960	24716	1.74%	0.179%
48	17.333	2352	2356	2361	rBV	81056	106459	7.51%	0.771%
49	17.386	2361	2365	2370	rVV2	28568	43913	3.10%	0.318%
50	17.615	2398	2404	2409	rBV	440507	721596	50.92%	5.228%
51	17.662	2409	2412	2417	rVB	68152	96946	6.84%	0.702%
52	18.073	2478	2482	2487	rBV	69915	91866	6.48%	0.666%
53	18.250	2507	2512	2519	rVB	520547	656993	46.36%	4.760%
54	18.761	2595	2599	2602	rVB	61054	68499	4.83%	0.496%
55	19.378	2699	2704	2707	rBV2	874288	1204675	85.01%	8.728%
56	19.413	2707	2710	2719	rVB	1052288	1417093	100.00%	10.266%
57	19.536	2727	2731	2736	rBV	242766	305093	21.53%	2.210%
58	19.660	2747	2752	2758	rBV	100518	175407	12.38%	1.271%
59	20.018	2810	2813	2818	rVB	98766	139219	9.82%	1.009%
60	20.206	2840	2845	2850	rVB	838301	1082561	76.39%	7.843%
61	21.922	3131	3137	3143	rBV	320902	599278	42.29%	4.342%

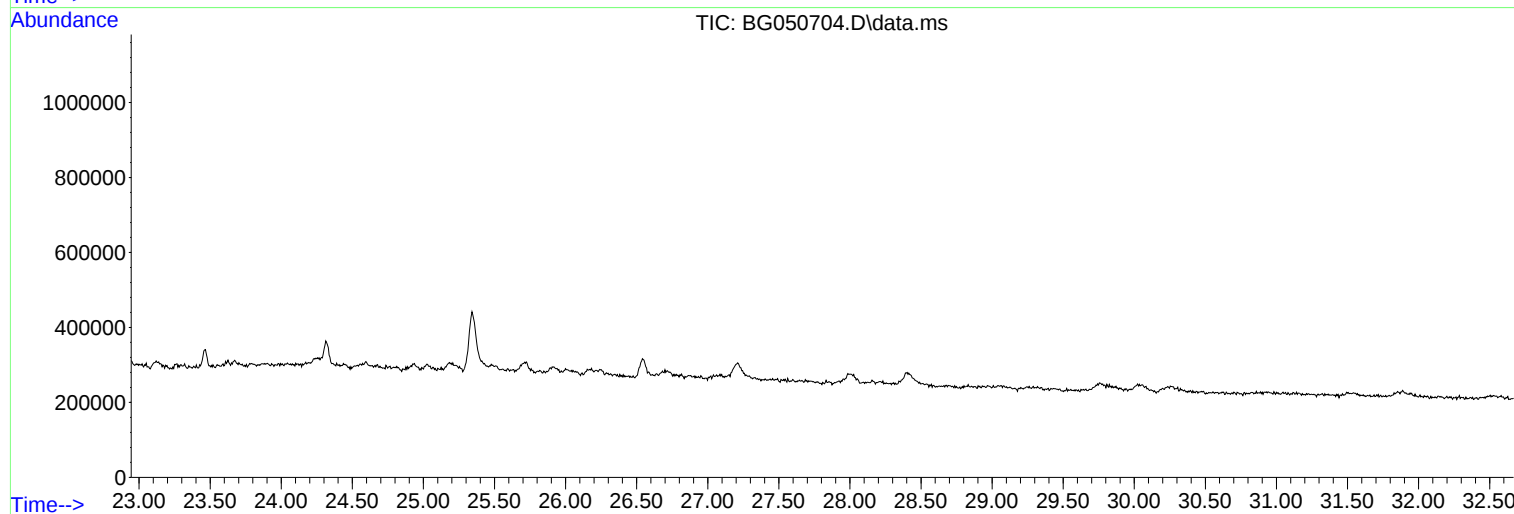
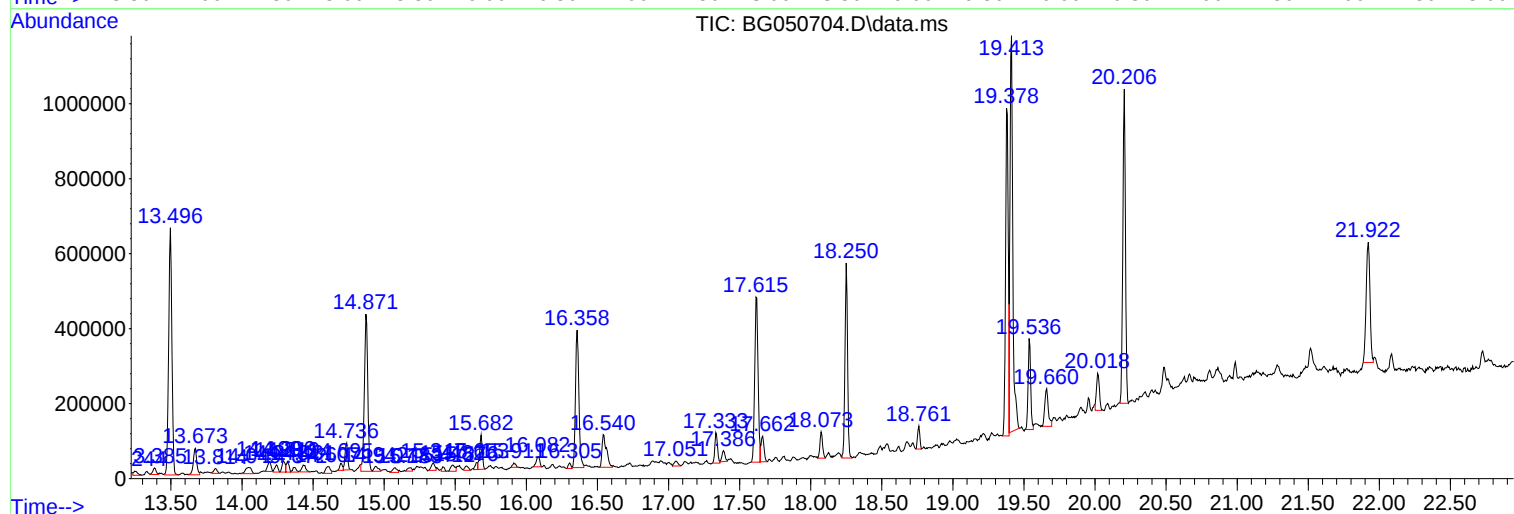
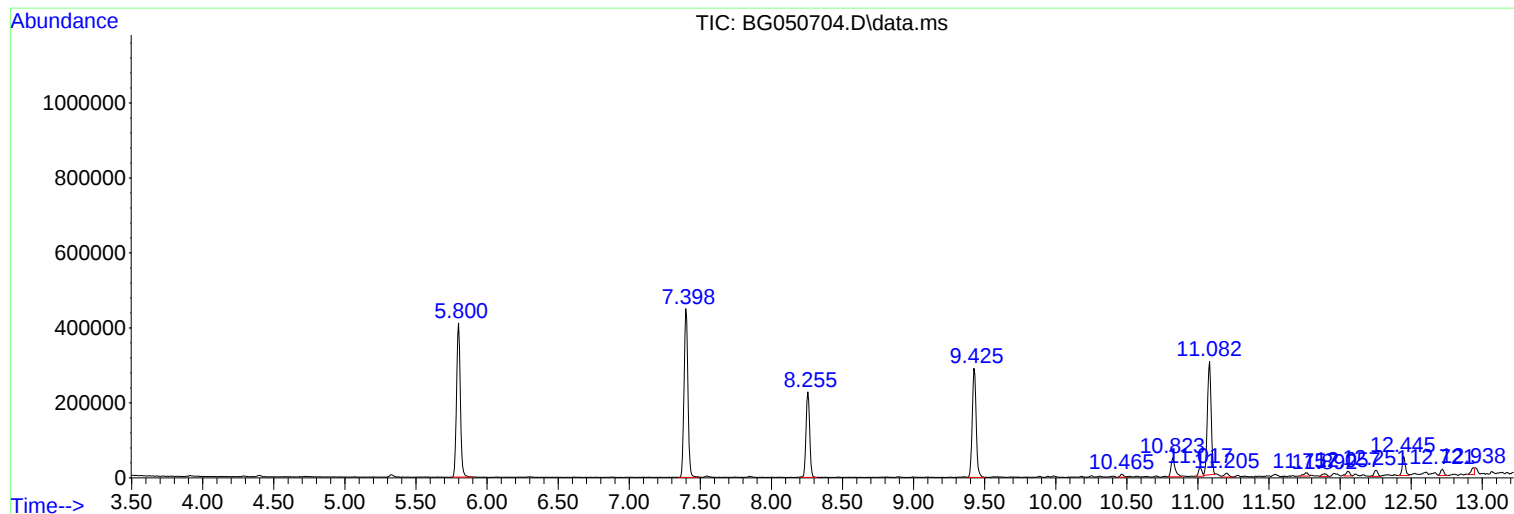
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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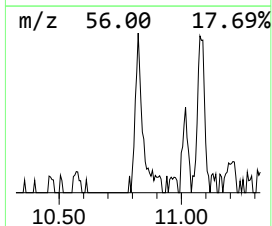
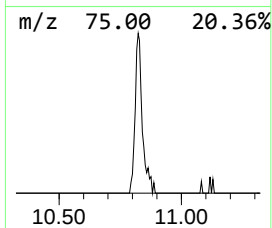
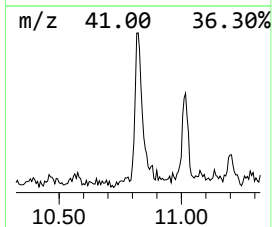
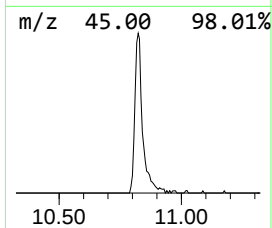
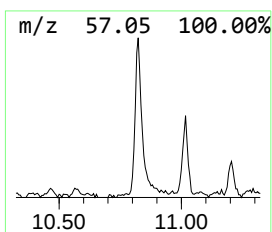
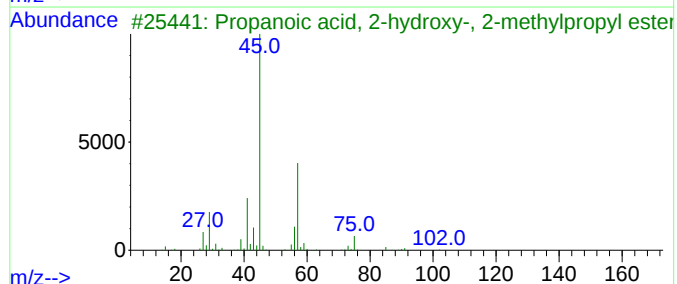
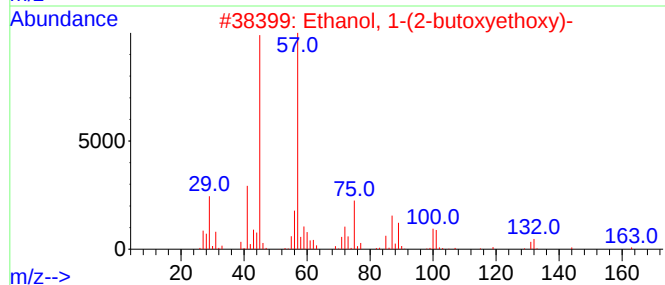
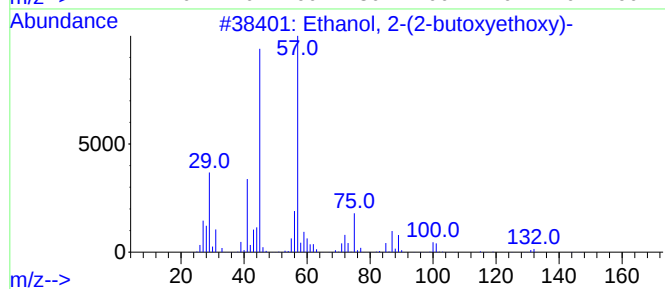
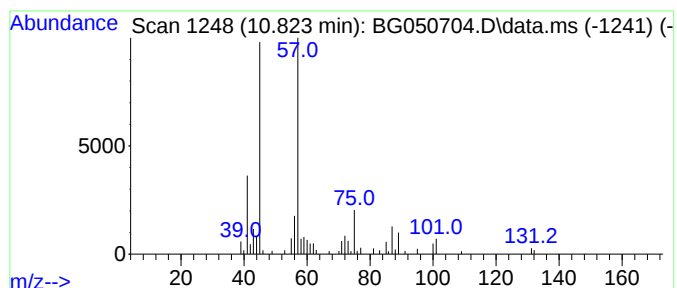
Instrument :
BNA_G
ClientSampleId :
NB-08-102021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.823	3.61 ng	100075	Naphthalene-d8	11.082	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3	Propanoic acid, 2-hydroxy-, 2-me...	146	C7H14O3	000585-24-0	59
4	Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	50
5	Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	50



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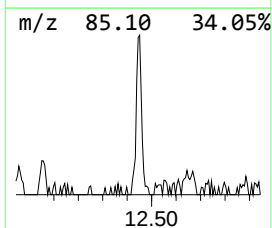
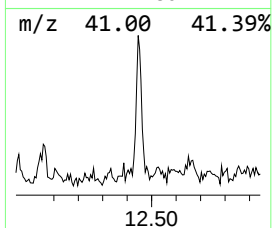
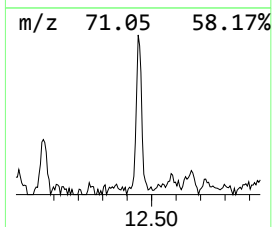
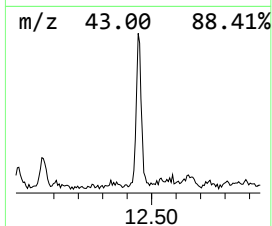
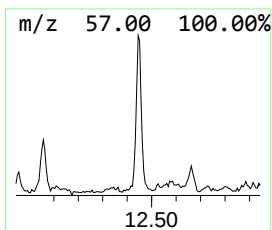
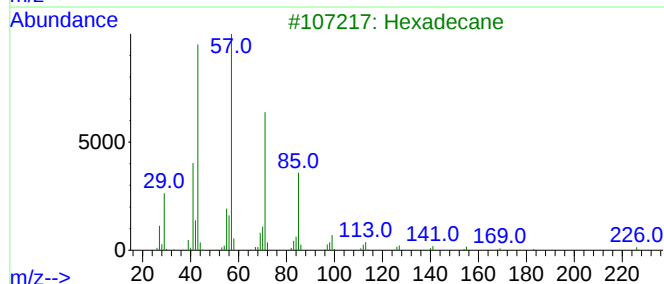
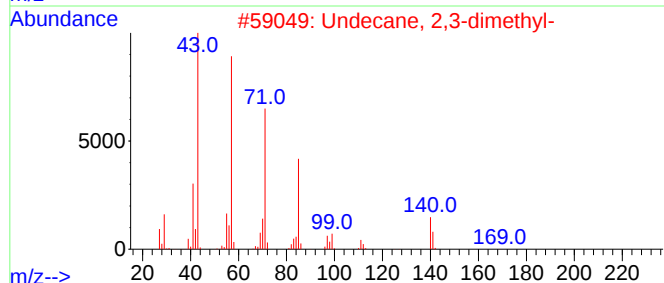
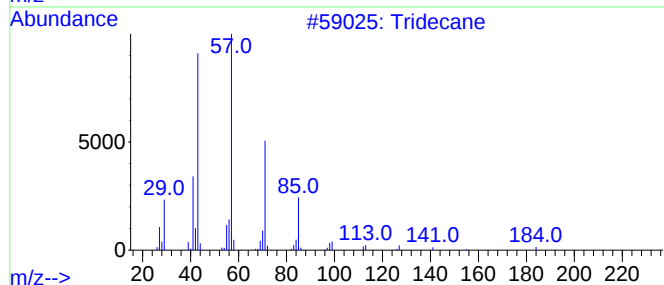
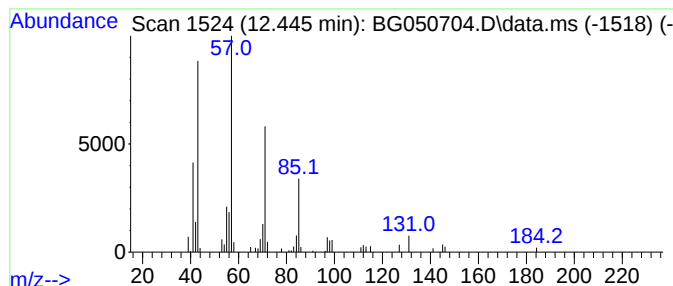
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Tridecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.445	2.83 ng	78406	Naphthalene-d8	11.082

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	95
2			Undecane, 2,3-dimethyl-	184	C13H28	017312-77-5	80
3			Hexadecane	226	C16H34	000544-76-3	80
4			Carbonic acid, tetradecyl vinyl ...	284	C17H32O3	1000382-54-5	72
5			Undecane	156	C11H24	001120-21-4	72



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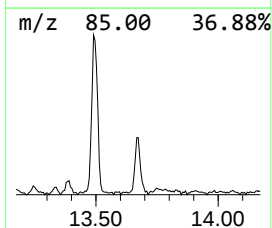
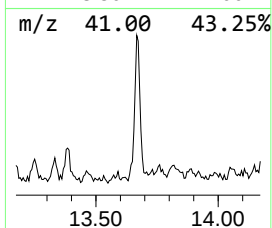
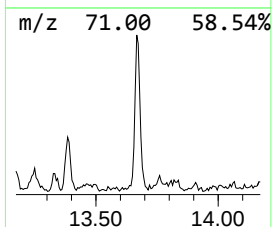
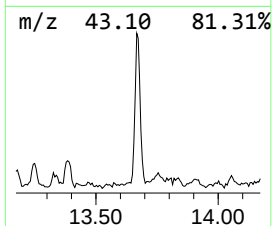
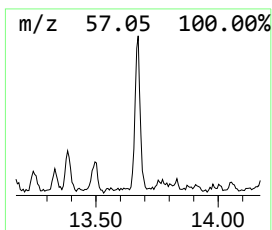
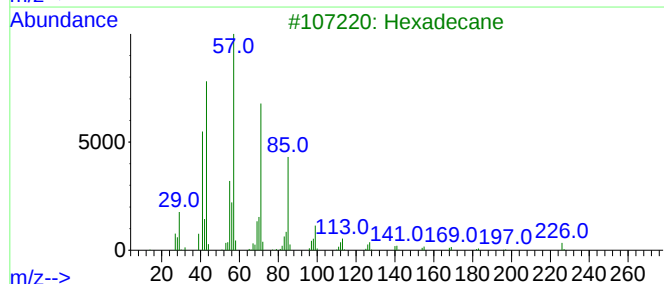
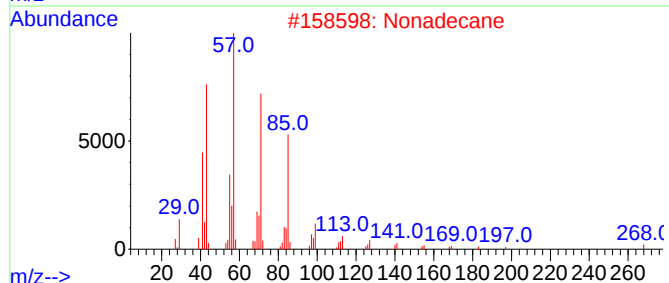
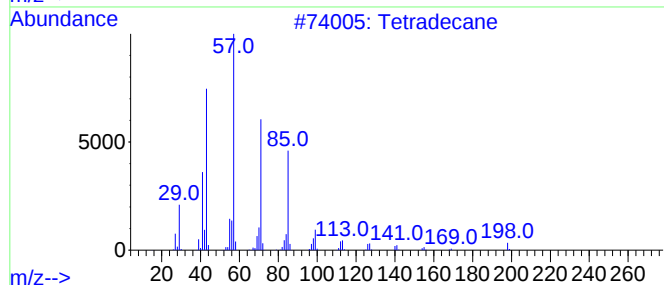
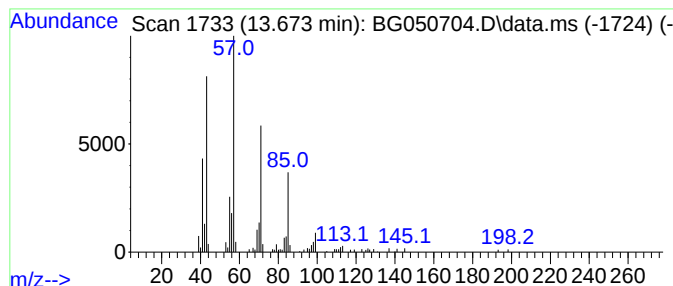
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Tetradecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.673	3.43 ng	119314	Acenaphthene-d10	14.871

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	91
2		Nonadecane	268	C19H40	000629-92-5	90
3		Hexadecane	226	C16H34	000544-76-3	90
4		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	90
5		Tridecane, 7-methyl-	198	C14H30	026730-14-3	81



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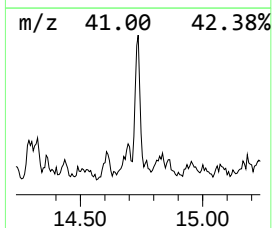
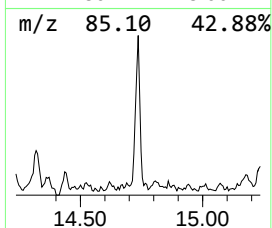
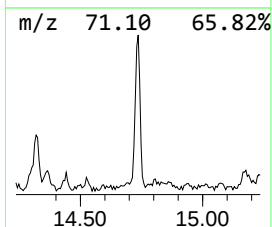
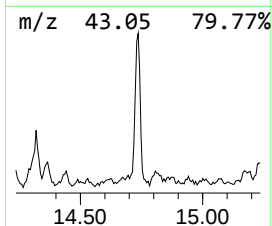
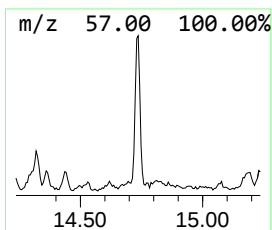
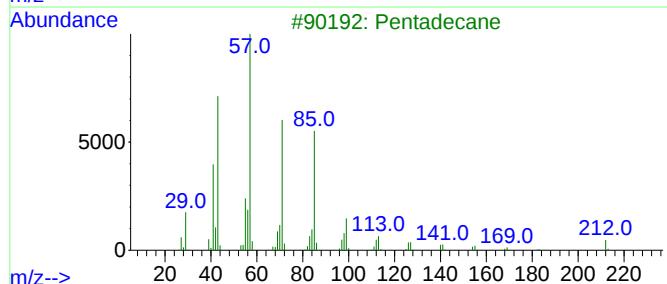
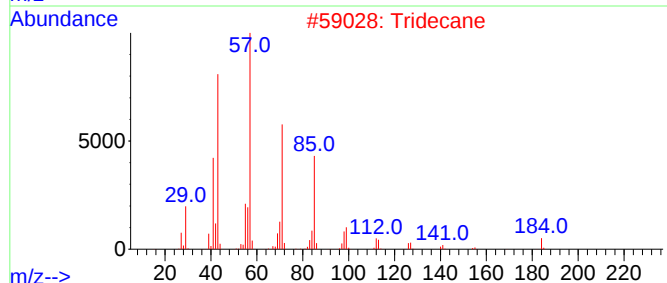
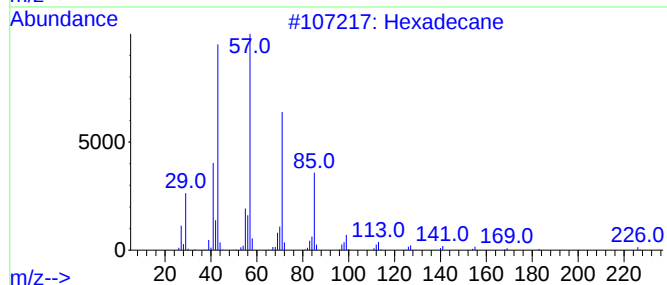
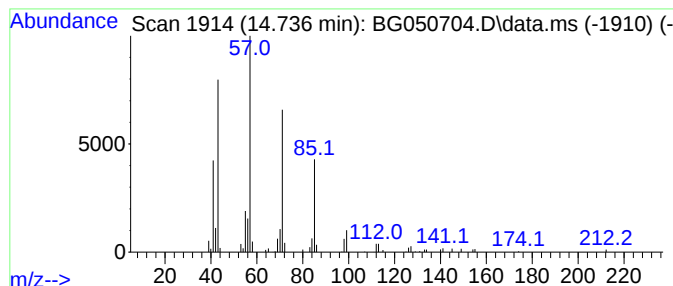
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.736	2.66 ng	92559	Acenaphthene-d10	14.871

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	91
2			Tridecane	184	C13H28	000629-50-5	91
3			Pentadecane	212	C15H32	000629-62-9	90
4			Tetradecane	198	C14H30	000629-59-4	90
5			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	86



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ClientSampleId :
NB-08-102021

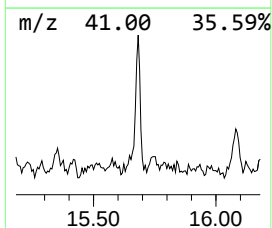
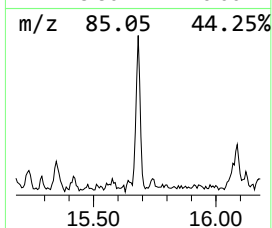
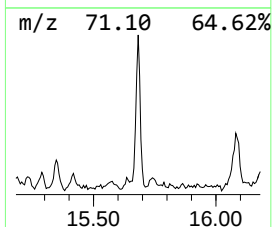
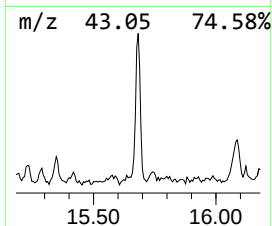
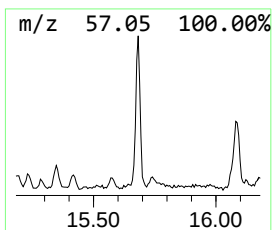
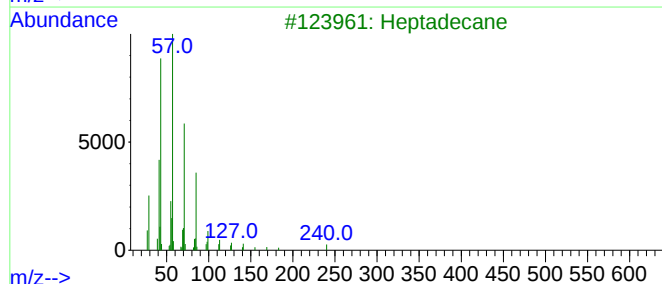
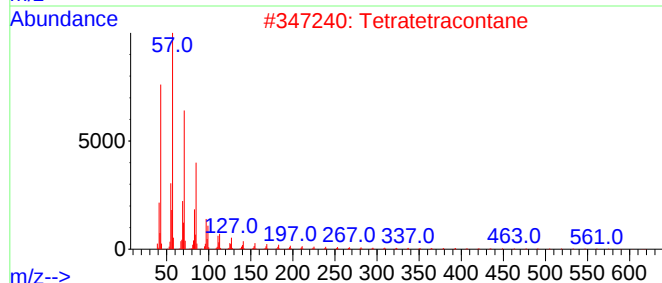
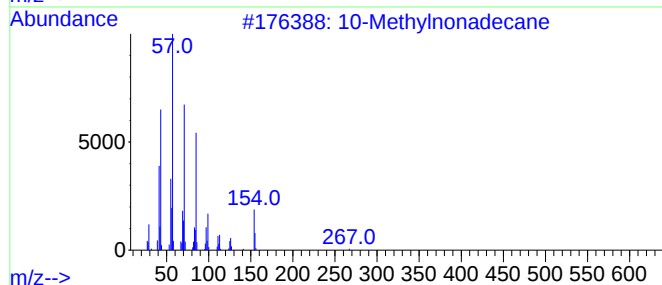
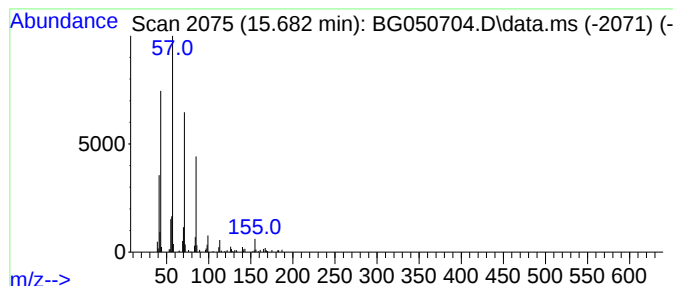
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 10-Methylnonadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.682	3.39 ng	118039	Acenaphthene-d10	14.871

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			10-Methylnonadecane	282	C20H42	056862-62-5	90
2			Tetratetracontane	619	C44H90	007098-22-8	83
3			Heptadecane	240	C17H36	000629-78-7	78
4			Hexadecane	226	C16H34	000544-76-3	72
5			Decane, 2,3,8-trimethyl-	184	C13H28	062238-14-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102221\
Data File : BG050704.D
Acq On : 23 Oct 2021 1:07
Operator : CG/JU
Sample : M4298-01 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
NB-08-102021

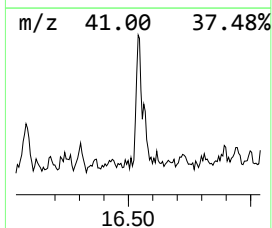
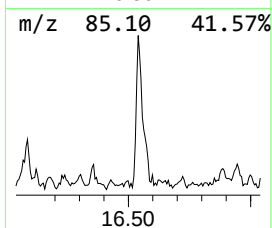
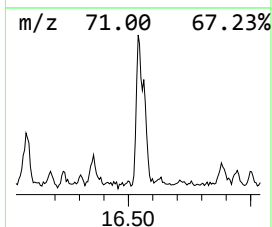
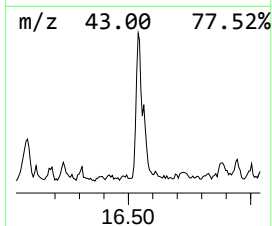
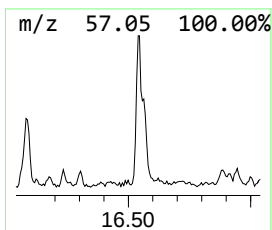
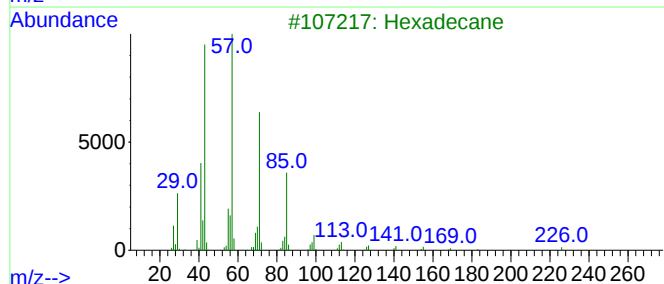
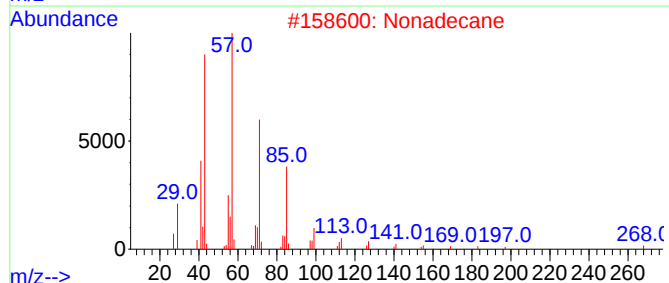
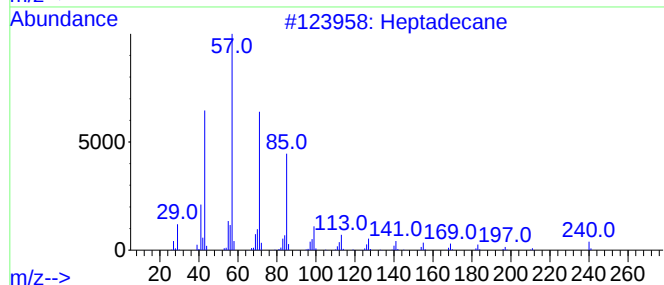
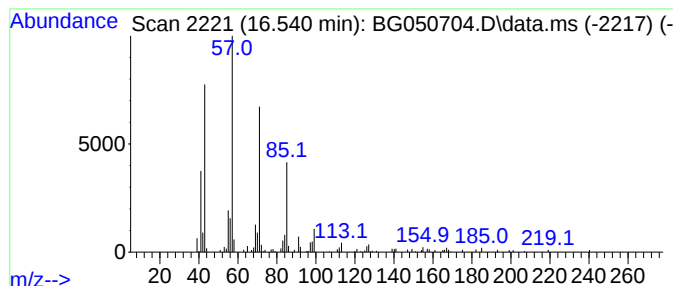
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Heptadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.540	5.29 ng	190684	Phenanthrene-d10	17.621

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	94
2		Nonadecane	268	C19H40	000629-92-5	91
3		Hexadecane	226	C16H34	000544-76-3	91
4		Pentadecane	212	C15H32	000629-62-9	90
5		Eicosane	282	C20H42	000112-95-8	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102221\
Data File : BG050704.D
Acq On : 23 Oct 2021 1:07
Operator : CG/JU
Sample : M4298-01 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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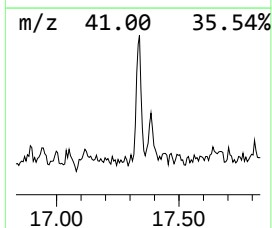
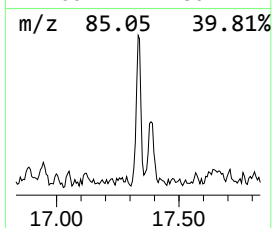
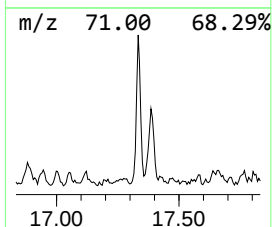
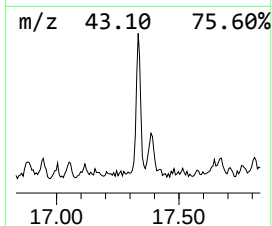
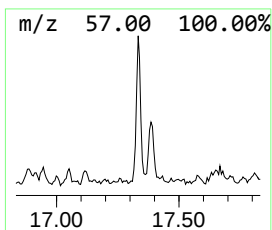
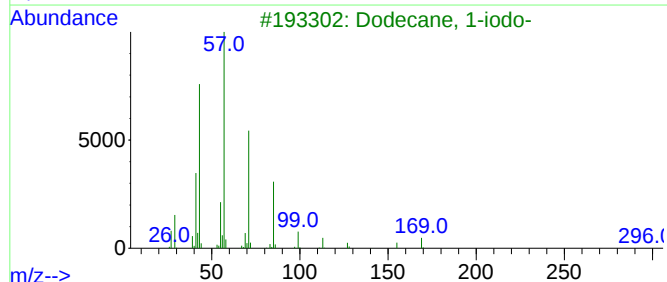
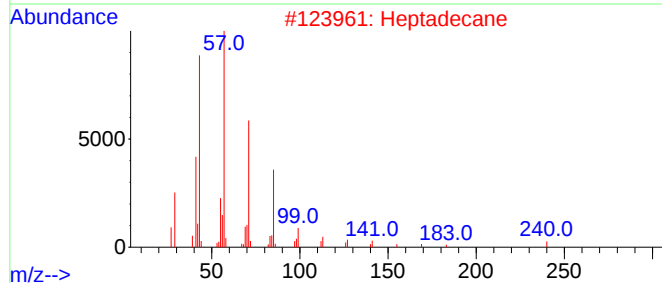
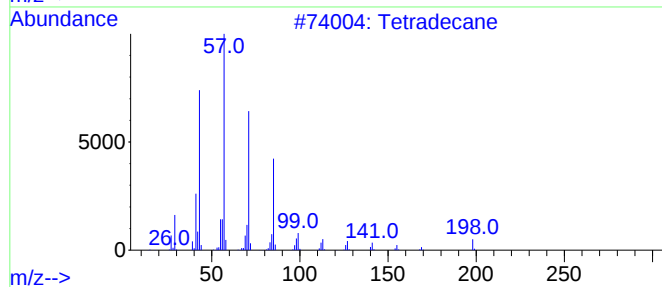
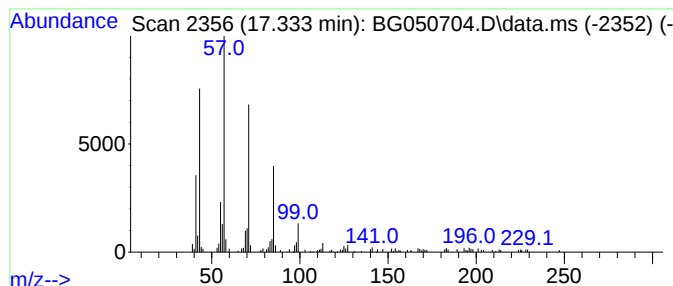
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Dodecane, 1-iodo- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.333	2.95 ng	106459	Phenanthrene-d10	17.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	93
2			Heptadecane	240	C17H36	000629-78-7	86
3			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86
4			Undecane, 2-methyl-	170	C12H26	007045-71-8	86
5			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102221\
Data File : BG050704.D
Acq On : 23 Oct 2021 1:07
Operator : CG/JU
Sample : M4298-01 2X
Misc :
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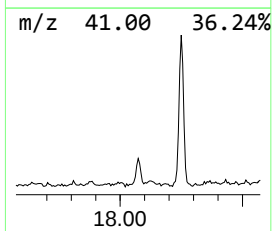
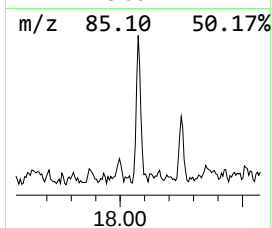
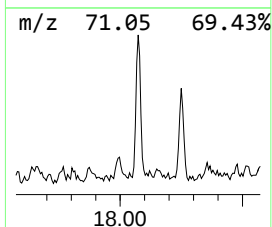
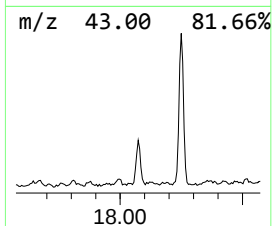
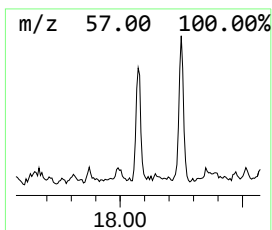
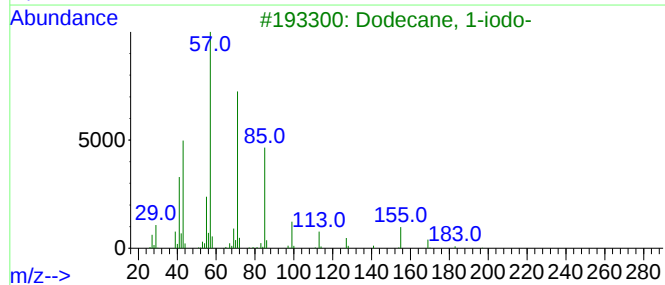
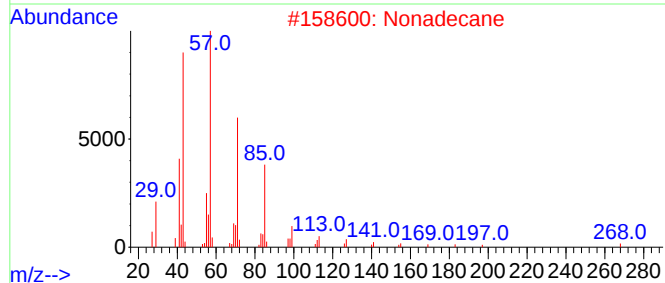
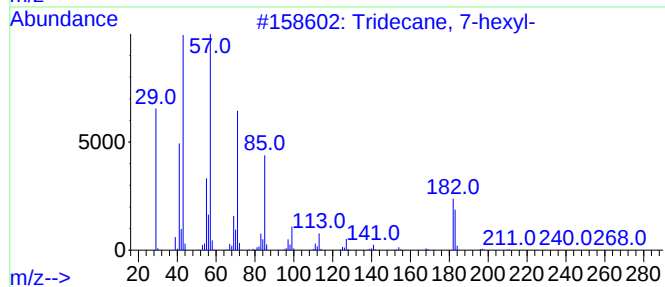
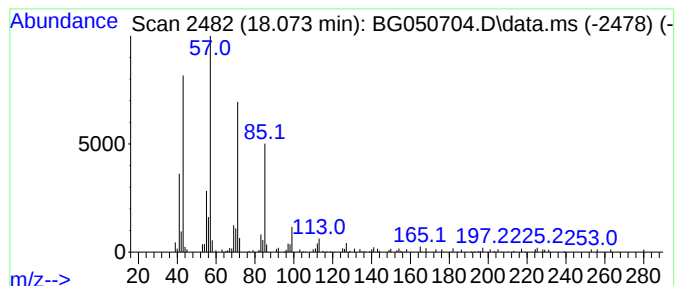
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Tridecane, 7-hexyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.073	2.55 ng	91866	Phenanthrene-d10		17.621	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tridecane, 7-hexyl-		268	C19H40	007225-66-3	87
2	Nonadecane		268	C19H40	000629-92-5	87
3	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	86
4	Pentadecane		212	C15H32	000629-62-9	86
5	Tetradecane, 1-iodo-		324	C14H29I	019218-94-1	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102221\
Data File : BG050704.D
Acq On : 23 Oct 2021 1:07
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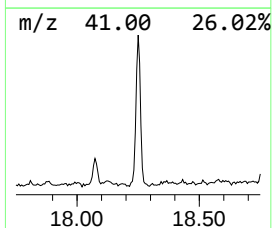
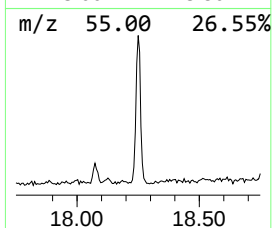
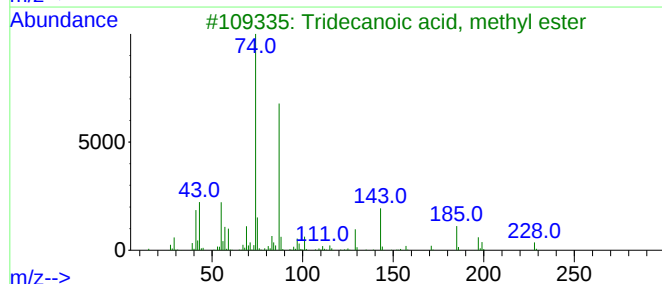
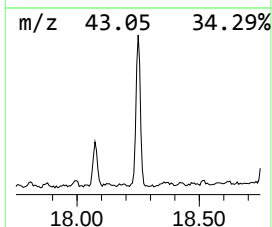
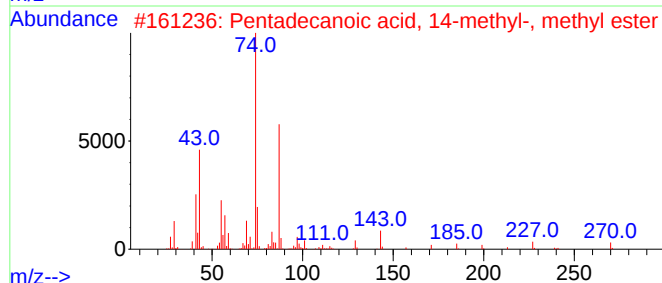
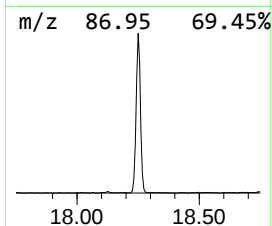
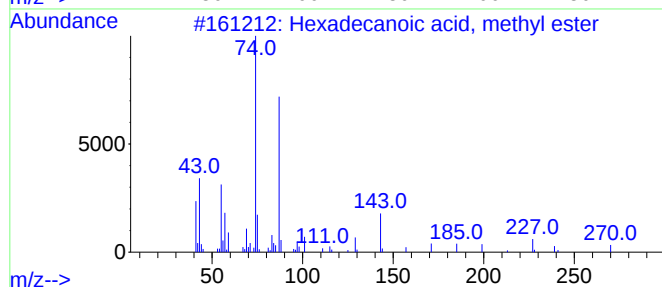
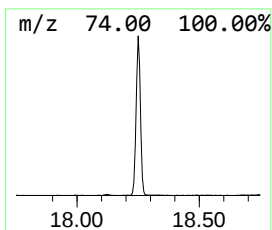
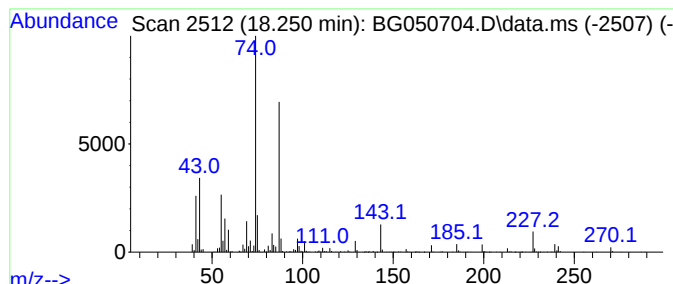
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.250	18.21 ng	656993	Phenanthrene-d10	17.621

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	95
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	94
3		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	93
4		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102221\
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Sample : M4298-01 2X
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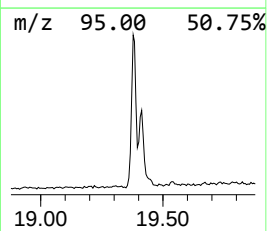
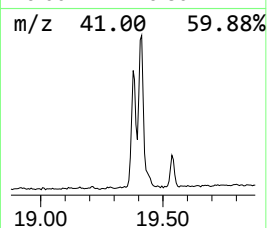
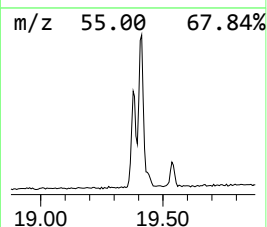
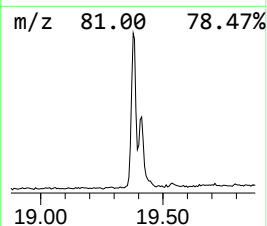
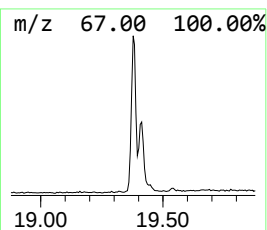
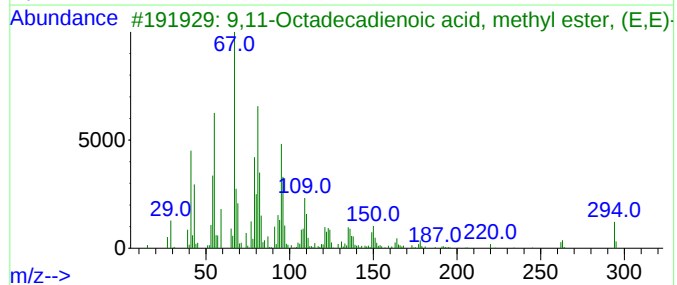
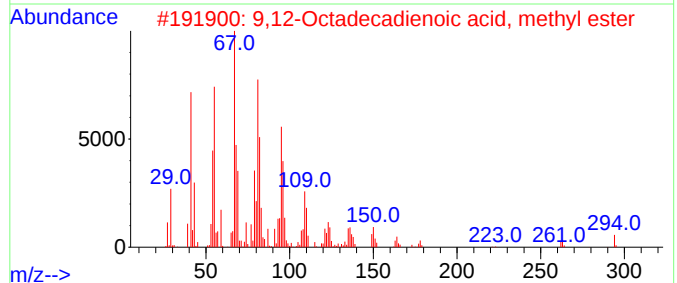
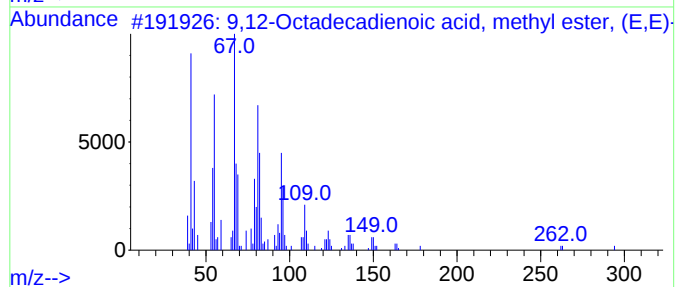
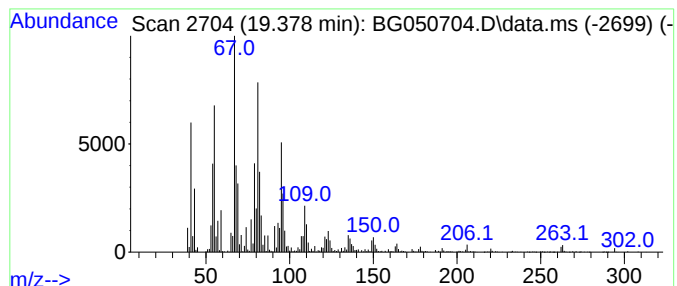
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 9,12-Octadecadienoic acid, ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.378	33.39 ng	1204680	Phenanthrene-d10	17.621	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
3	9,11-Octadecadienoic acid, methy...	294	C19H34O2	013038-47-6	98
4	Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	96
5	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102221\
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Sample : M4298-01 2X
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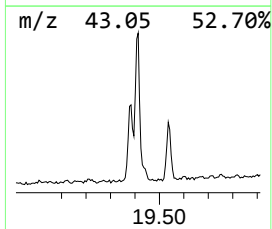
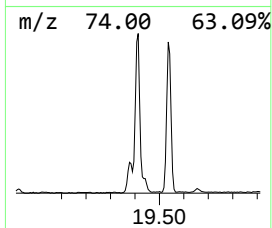
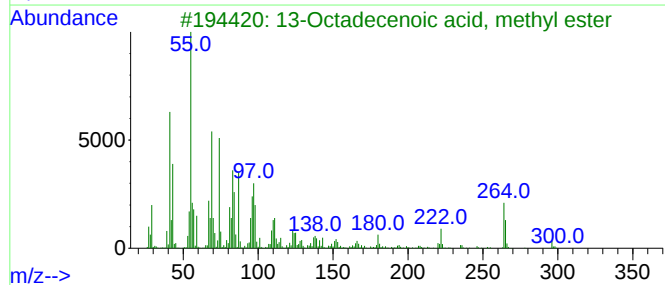
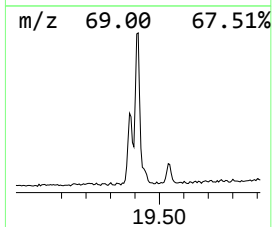
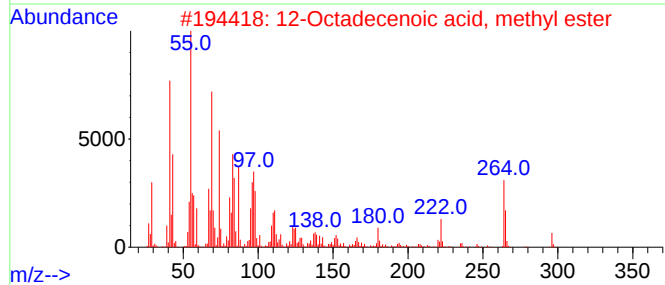
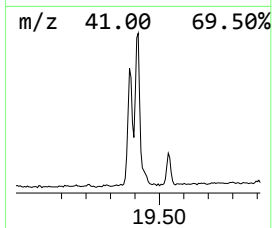
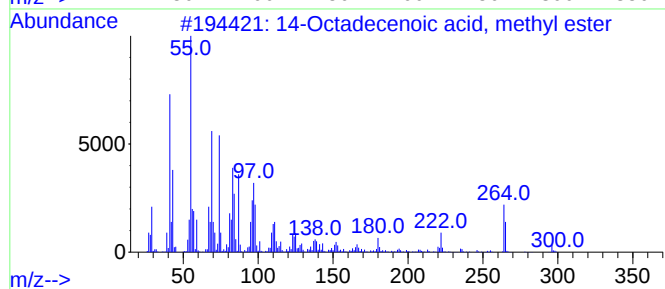
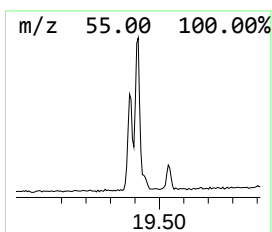
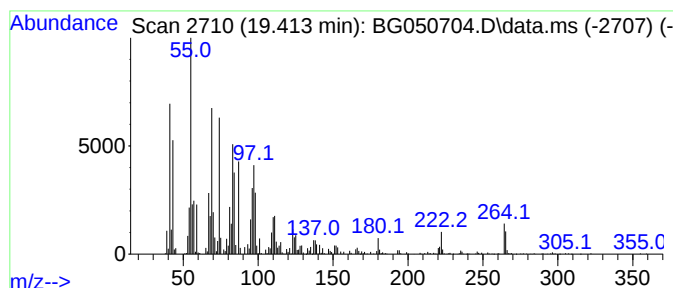
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TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 14-Octadecenoic acid, methyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.413	39.28 ng	1417090	Phenanthrene-d10	17.621	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	99
2	12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	99
3	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	99
4	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	98
5	9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102221\
Data File : BG050704.D
Acq On : 23 Oct 2021 1:07
Operator : CG/JU
Sample : M4298-01 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
NB-08-102021

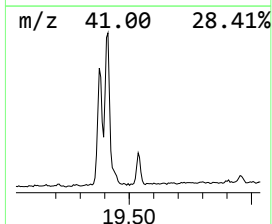
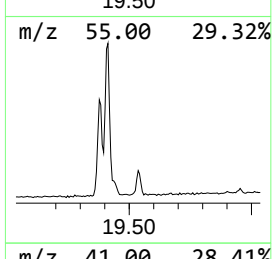
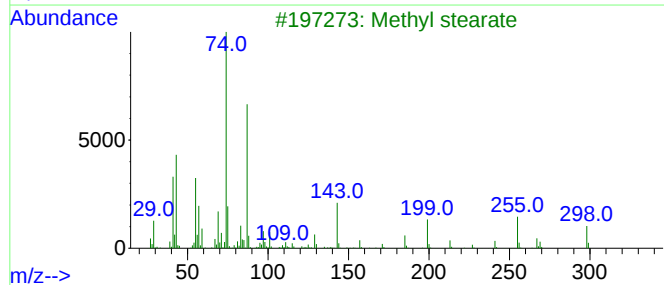
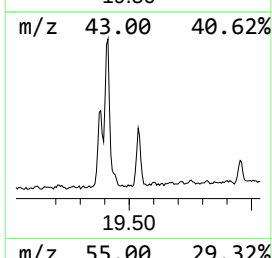
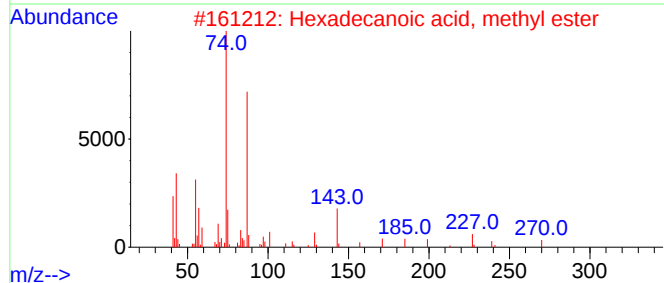
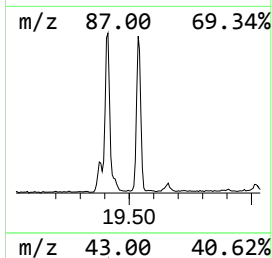
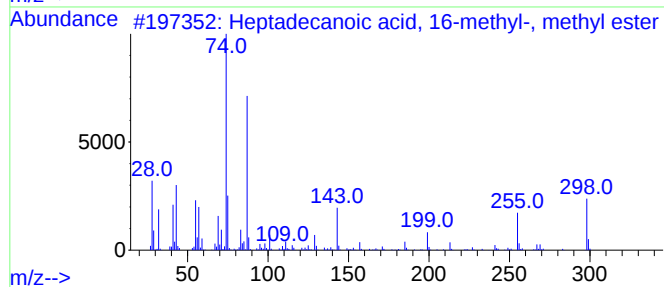
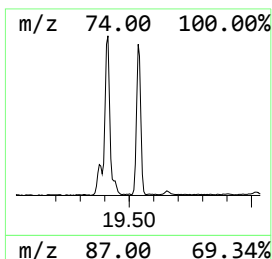
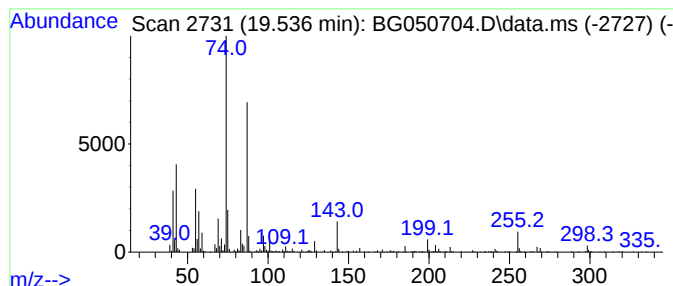
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Heptadecanoic acid, 16-meth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.536	8.46 ng	305093	Phenanthrene-d10	17.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	95
2			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	94
3			Methyl stearate	298	C19H38O2	000112-61-8	93
4			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	93
5			Methyl tetradecanoate	242	C15H30O2	000124-10-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102221\
 Data File : BG050704.D
 Acq On : 23 Oct 2021 1:07
 Operator : CG/JU
 Sample : M4298-01 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NB-08-102021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(2-b...	10.823	3.6	ng	100075	2	11.082	554212	20.0
Tridecane	12.445	2.8	ng	78406	2	11.082	554212	20.0
Tetradecane	13.673	3.4	ng	119314	3	14.871	695459	20.0
Hexadecane	14.736	2.7	ng	92559	3	14.871	695459	20.0
10-Methylnonade...	15.682	3.4	ng	118039	3	14.871	695459	20.0
Heptadecane	16.540	5.3	ng	190684	4	17.621	721596	20.0
Dodecane, 1-iodo-	17.333	3.0	ng	106459	4	17.621	721596	20.0
Tridecane, 7-he...	18.073	2.5	ng	91866	4	17.621	721596	20.0
Hexadecanoic ac...	18.250	18.2	ng	656993	4	17.621	721596	20.0
9,12-Octadecadi...	19.378	33.4	ng	1204680	4	17.621	721596	20.0
14-Octadecenoic...	19.413	39.3	ng	1417090	4	17.621	721596	20.0
Heptadecanoic a...	19.536	8.5	ng	305093	4	17.621	721596	20.0