

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021522\
 Data File : BP009166.D
 Acq On : 15 Feb 2022 21:50
 Operator : CG/JU
 Sample : N1521-16
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP3-B

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_P\Methods\8270E-BP020722.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP009166.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.911	322	327	334	rBV	82836	137446	1.25%	0.275%
2	5.381	401	407	435	rBV	2904020	4929649	45.01%	9.851%
3	6.975	671	678	722	rBV	2556043	5011281	45.75%	10.014%
4	7.781	808	815	833	rBV	568264	1032133	9.42%	2.062%
5	8.946	1006	1013	1042	rBV	1634508	3259313	29.76%	6.513%
6	10.587	1284	1292	1310	rBV	590148	1270899	11.60%	2.540%
7	13.045	1703	1710	1726	rBV	4765118	7716947	70.46%	15.420%
8	14.428	1938	1945	1964	rBV	984687	1630488	14.89%	3.258%
9	15.928	2193	2200	2222	rBV	3157453	5583980	50.98%	11.158%
10	17.192	2408	2415	2432	rBV	1082478	1897250	17.32%	3.791%
11	18.086	2562	2567	2577	rBV	245710	460786	4.21%	0.921%
12	19.322	2772	2777	2789	rBV	217466	438654	4.00%	0.877%
13	19.751	2844	2850	2859	rBV	9169505	10952757	100.00%	21.886%
14	20.992	3057	3061	3071	rBV	349001	487497	4.45%	0.974%
15	21.286	3106	3111	3122	rBV	1640836	2343716	21.40%	4.683%
16	21.421	3131	3134	3140	rVB2	138100	150110	1.37%	0.300%
17	21.880	3209	3212	3220	rVB2	116300	155815	1.42%	0.311%
18	23.674	3511	3517	3529	rVB2	1171088	2585184	23.60%	5.166%

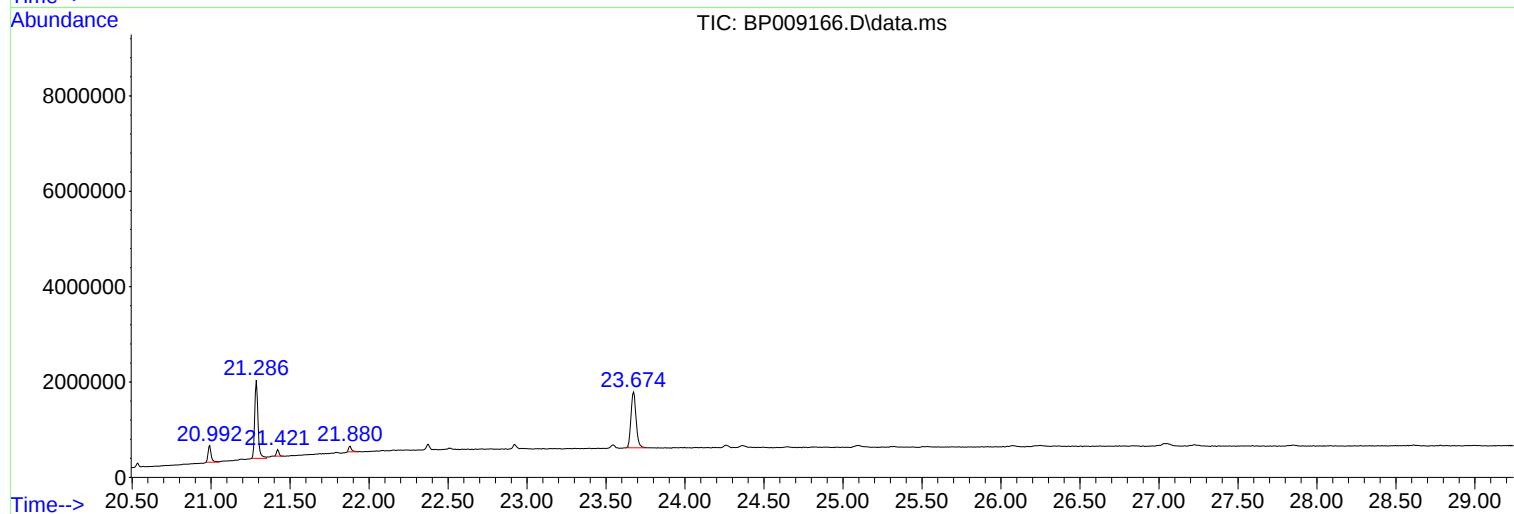
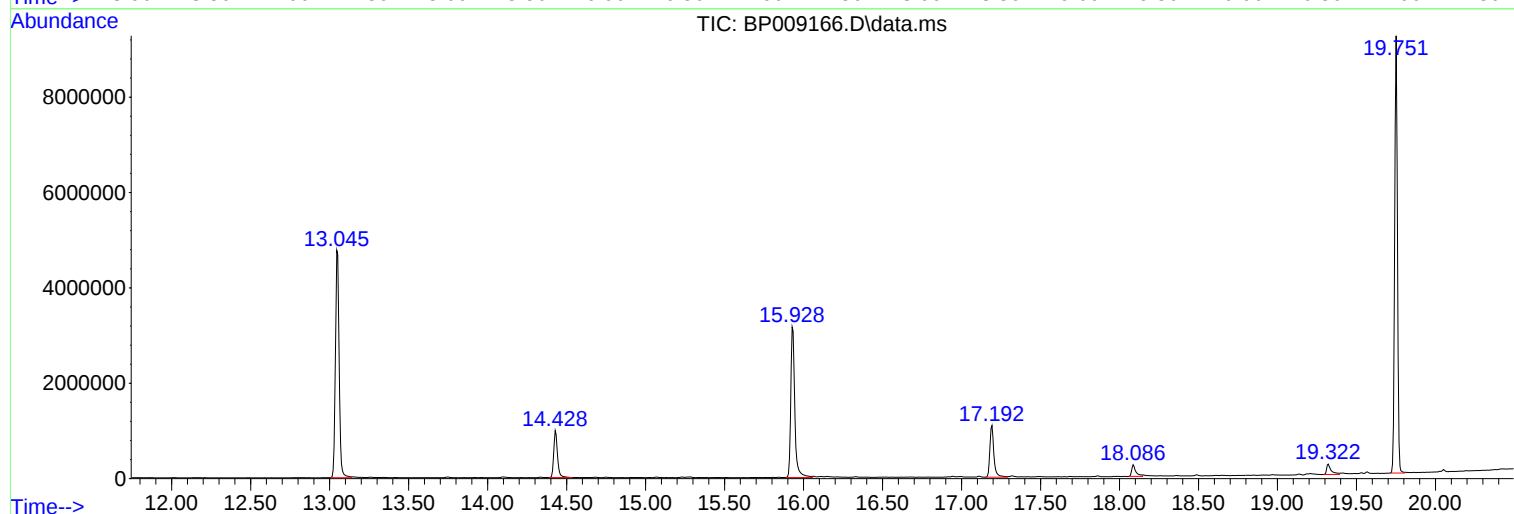
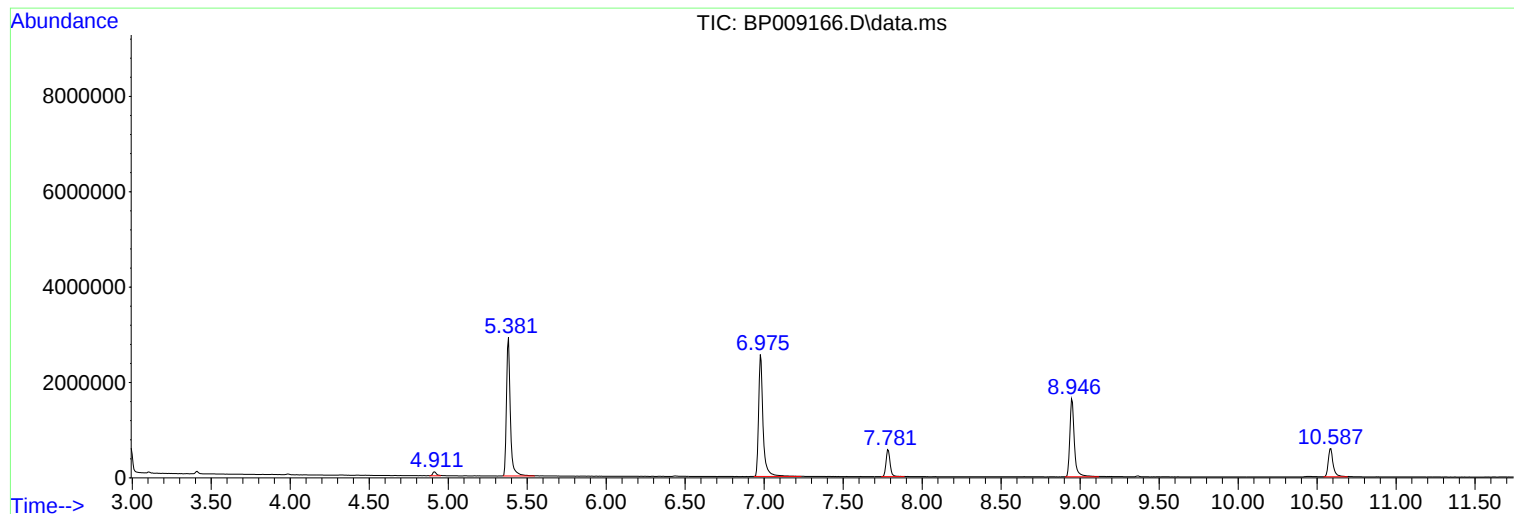
Sum of corrected areas: 50043905

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.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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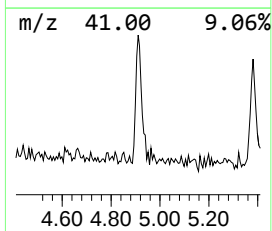
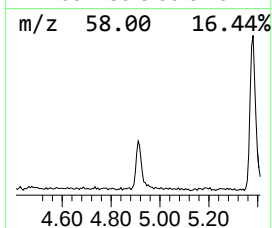
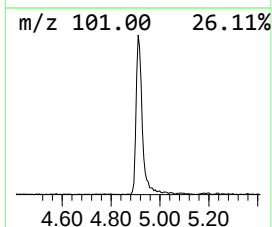
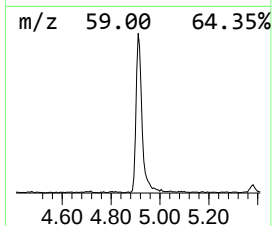
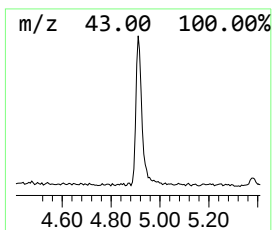
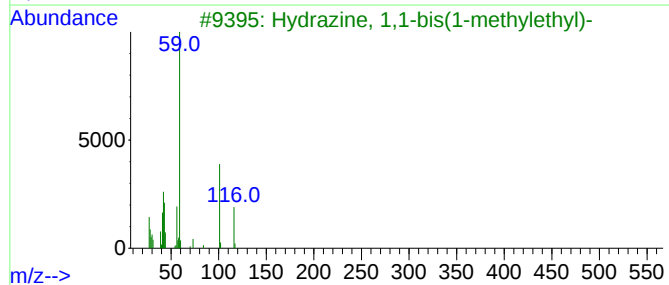
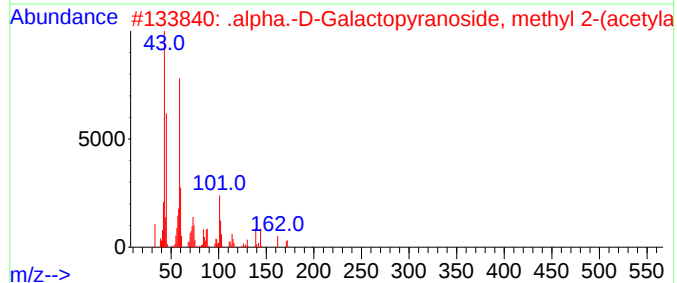
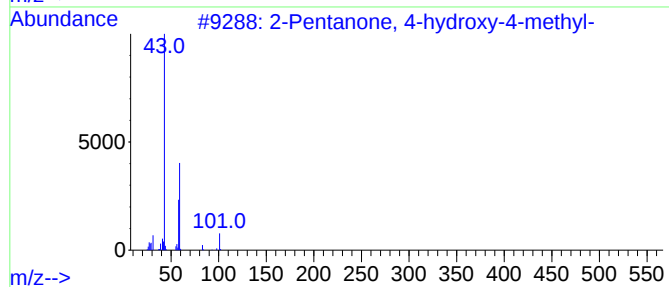
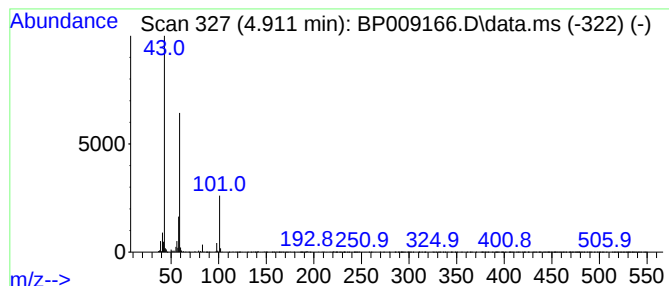
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.911	2.66 ng	137446	1,4-Dichlorobenzene-d4	7.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			.alpha.-D-Galactopyranoside, met...	249	C10H19NO6	017296-07-0	36
3			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	25
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
5			1-Cyclohexylethanol, methyl ether	142	C9H18O	1000365-12-8	9



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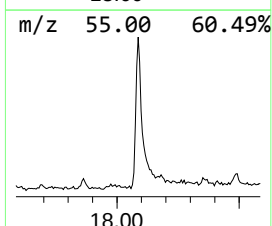
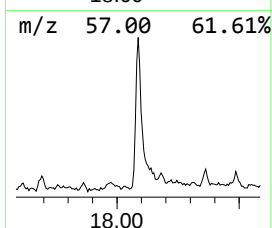
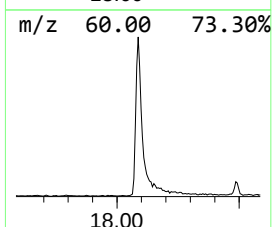
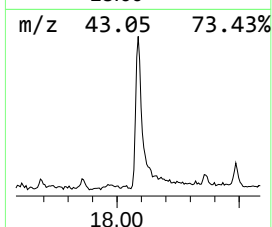
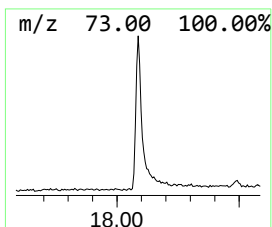
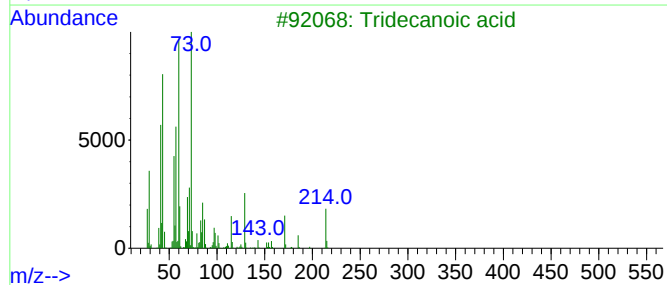
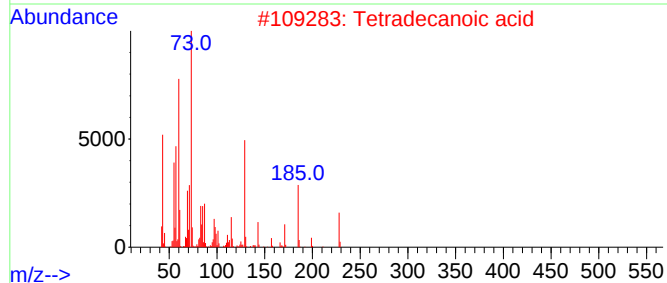
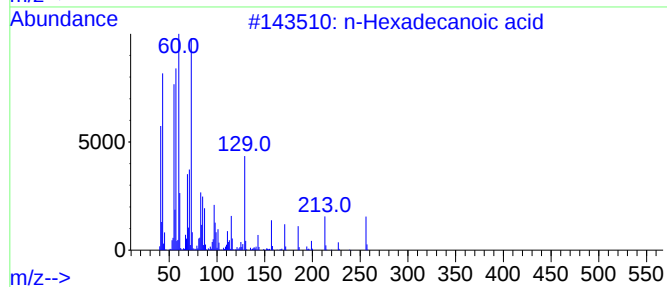
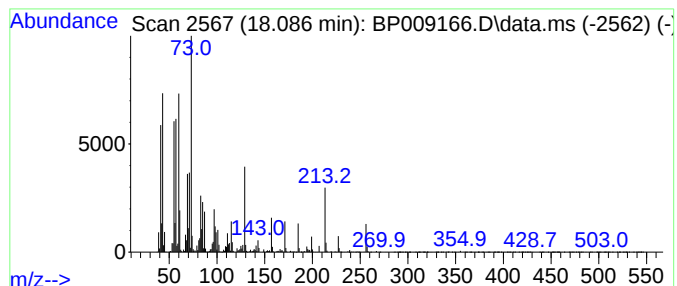
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Peak Number 2 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.086	4.86 ng	460786	Phenanthrene-d10	17.192

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	90
3		Tridecanoic acid	214	C13H26O2	000638-53-9	89
4		Octadecanoic acid	284	C18H36O2	000057-11-4	87
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	87



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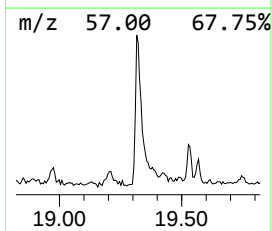
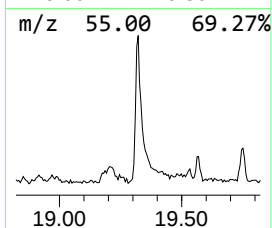
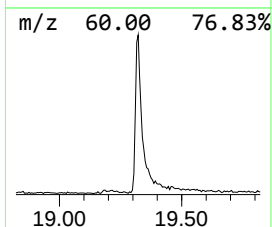
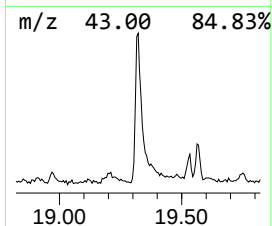
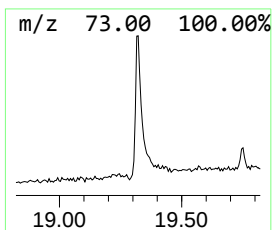
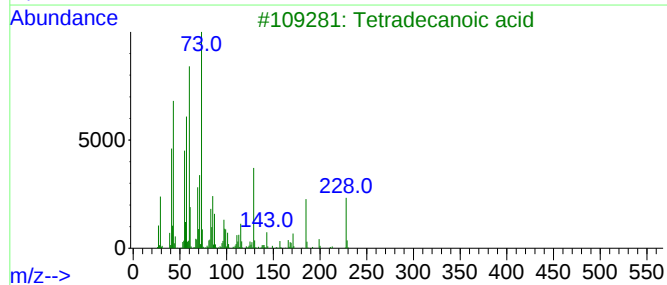
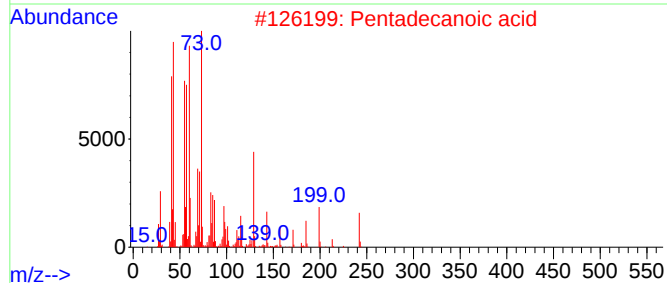
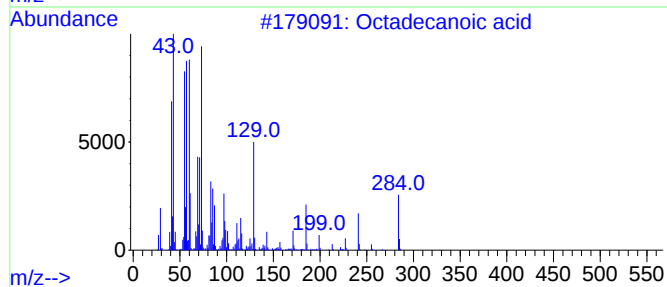
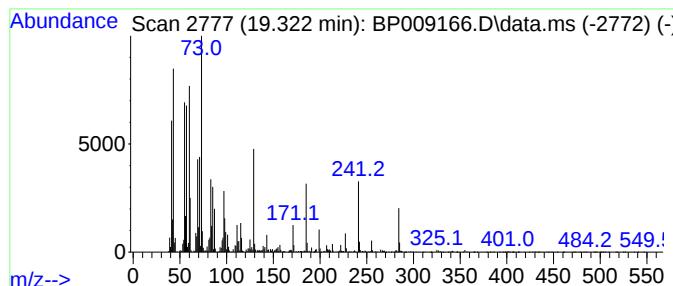
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Peak Number 3 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.322	3.74 ng	438654	Chrysene-d12	21.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
4			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	70
5			n-Decanoic acid	172	C10H20O2	000334-48-5	64



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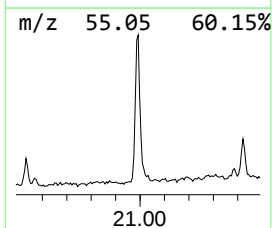
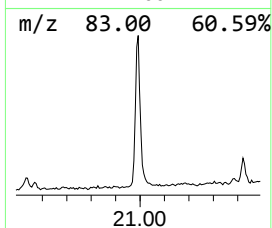
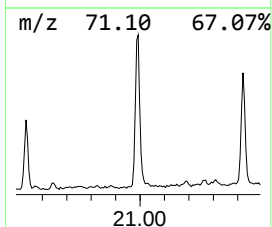
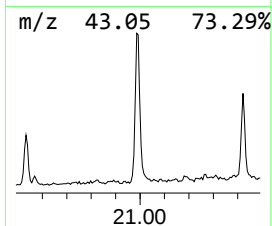
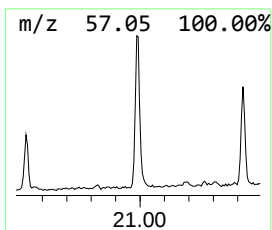
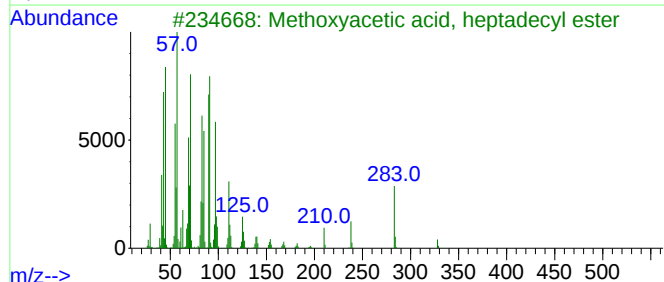
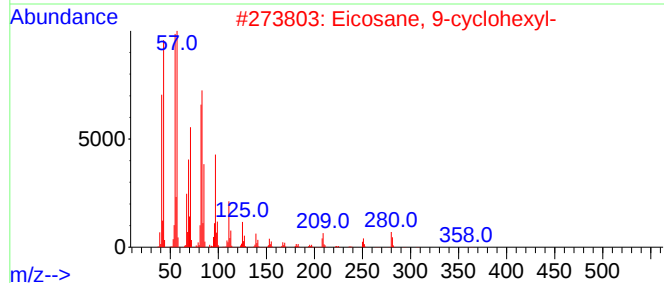
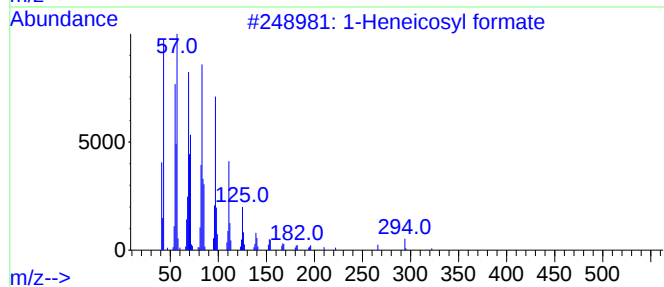
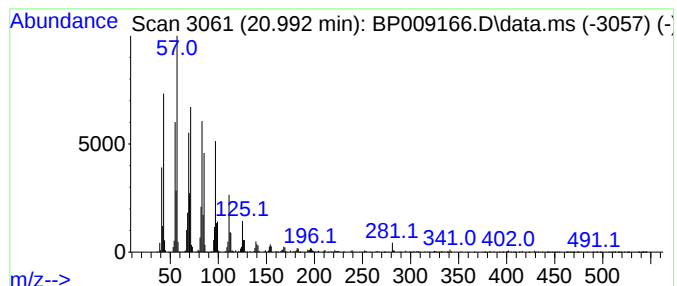
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Peak Number 4 1-Heneicosyl formate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.992	4.16 ng	487497	Chrysene-d12	21.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
2			Eicosane, 9-cyclohexyl-	364	C26H52	004443-61-2	93
3			Methoxyacetic acid, heptadecyl e...	328	C20H40O3	1000282-99-1	91
4			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91
5			Triacetyl acetate	480	C32H64O2	041755-58-2	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.911	2.7	ng	137446	1	7.781	1032130	20.0
n-Hexadecanoic ...	18.086	4.9	ng	460786	4	17.192	1897250	20.0
Octadecanoic acid	19.322	3.7	ng	438654	5	21.286	2343720	20.0
1-Heneicosyl fo...	20.992	4.2	ng	487497	5	21.286	2343720	20.0