

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM013020\
 Data File : BM024666.D
 Acq On : 30 Jan 2020 19:44
 Operator : CG/JU
 Sample : L1318-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RAINEY-PARK-BH-03-B

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM012320.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.069	365	371	379	rBV	2146668	2970125	34.21%	6.267%
2	6.628	627	636	647	rBV	2065612	3118077	35.92%	6.580%
3	7.428	765	772	782	rBV	605821	905052	10.42%	1.910%
4	7.745	819	826	834	rBV	2010469	3057421	35.22%	6.452%
5	8.569	958	966	976	rBV	1292412	2109068	24.29%	4.450%
6	10.186	1234	1241	1251	rBV	719941	1166377	13.44%	2.461%
7	12.692	1659	1667	1678	rBV	3553574	5146035	59.28%	10.859%
8	13.551	1808	1813	1819	rBV	99315	136052	1.57%	0.287%
9	14.080	1897	1903	1910	rBV	1169631	1655802	19.07%	3.494%
10	15.574	2151	2157	2170	rBV	3072395	4474576	51.54%	9.442%
11	16.833	2364	2371	2375	rBV	1469565	2053701	23.66%	4.334%
12	16.874	2375	2378	2384	rVB	99577	126951	1.46%	0.268%
13	17.774	2524	2531	2538	rBV2	285353	412509	4.75%	0.870%
14	18.897	2717	2722	2727	rBV	210891	293804	3.38%	0.620%
15	19.262	2776	2784	2789	rBV	204603	303386	3.49%	0.640%
16	19.492	2817	2823	2835	rVB	7038694	8681569	100.00%	18.319%
17	20.209	2939	2945	2950	rBV	322813	398239	4.59%	0.840%
18	20.797	3041	3045	3056	rVB2	593837	794308	9.15%	1.676%
19	21.038	3081	3086	3091	rBV	2359933	3183405	36.67%	6.717%
20	21.668	3190	3193	3200	rBV2	154985	206549	2.38%	0.436%
21	22.103	3265	3267	3271	rVB2	153145	167285	1.93%	0.353%
22	22.221	3284	3287	3293	rVB	364990	436586	5.03%	0.921%
23	23.150	3440	3445	3457	rVB	1834410	3229806	37.20%	6.815%
24	24.362	3645	3651	3663	rVB4	1015800	2363392	27.22%	4.987%

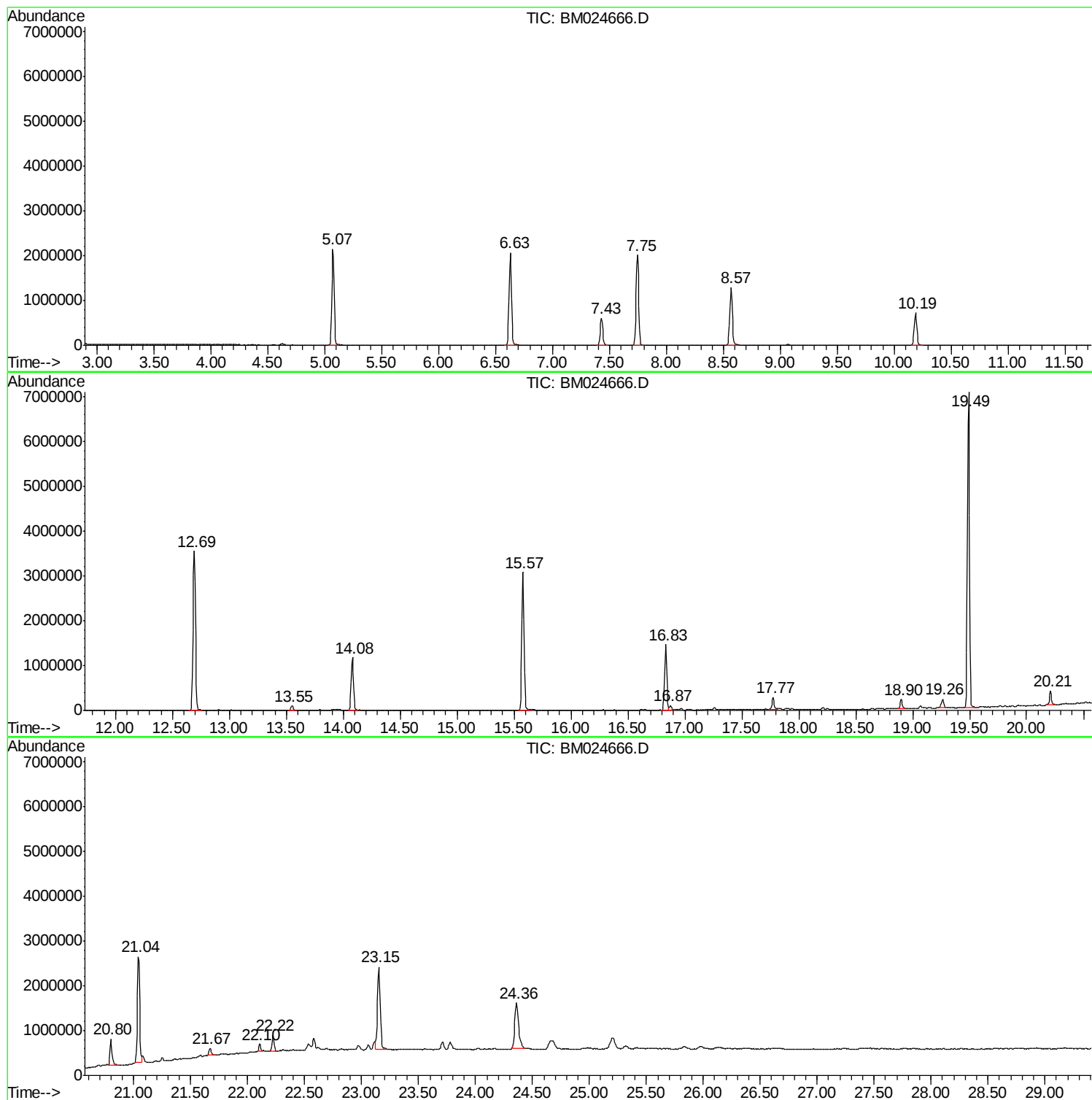
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM012320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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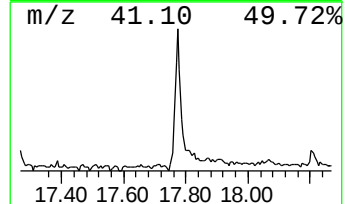
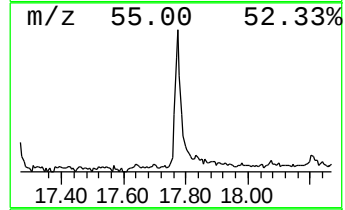
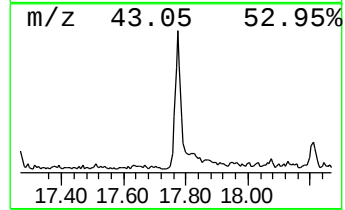
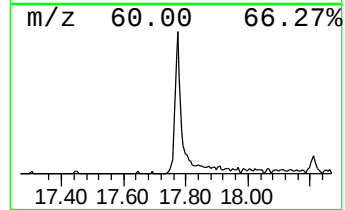
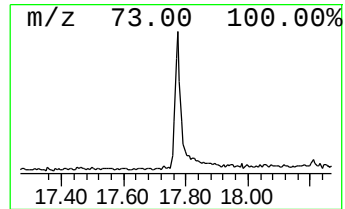
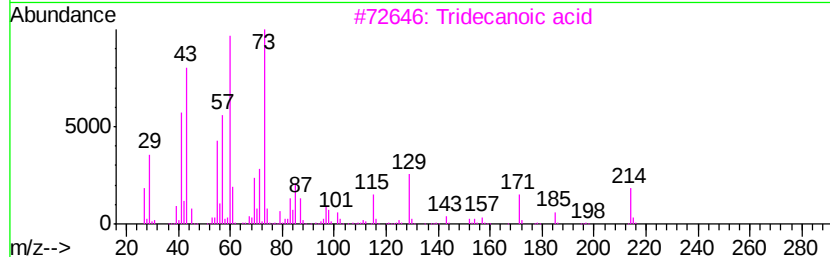
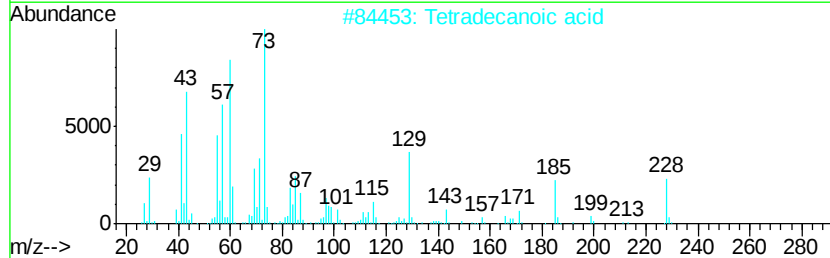
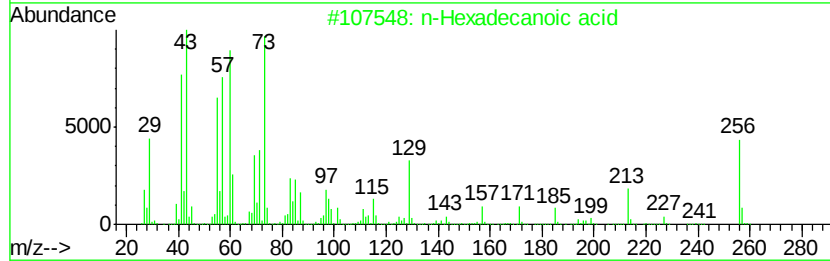
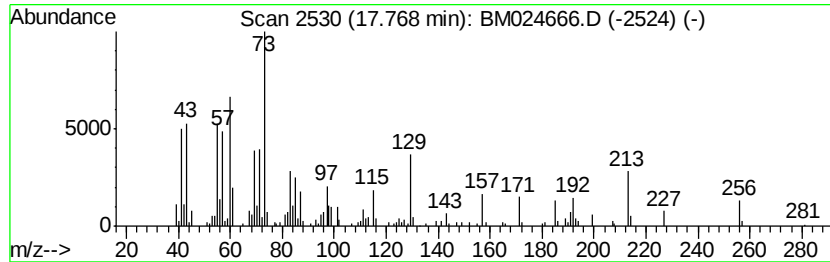
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM012320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.77	4.02 ng	412509	Phenanthrene-d10	16.83

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
3	Tridecanoic acid	214	C13H26O2	000638-53-9	64
4	n-Decanoic acid	172	C10H20O2	000334-48-5	58
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	49



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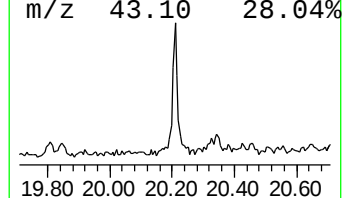
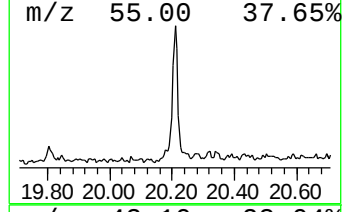
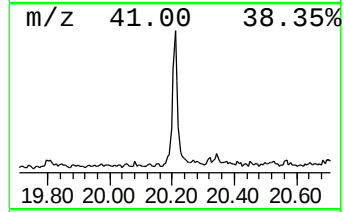
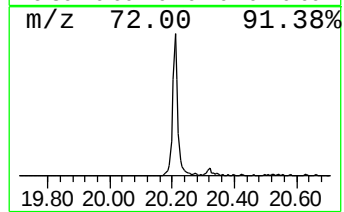
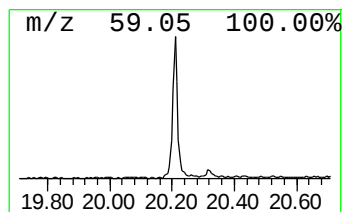
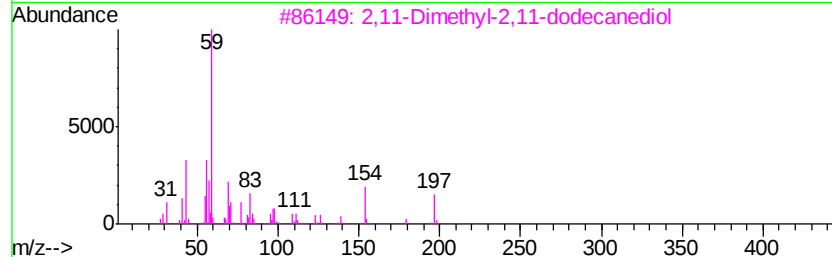
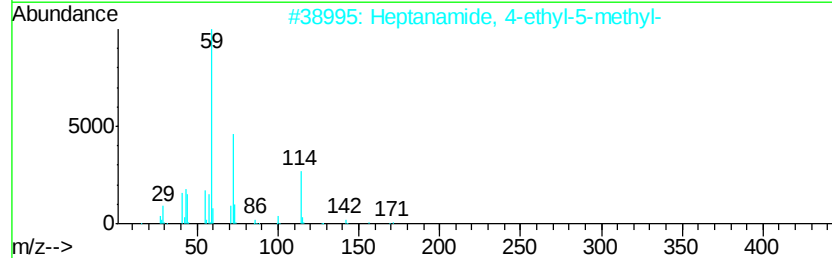
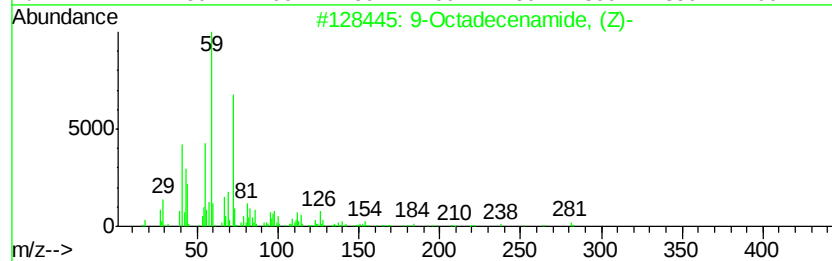
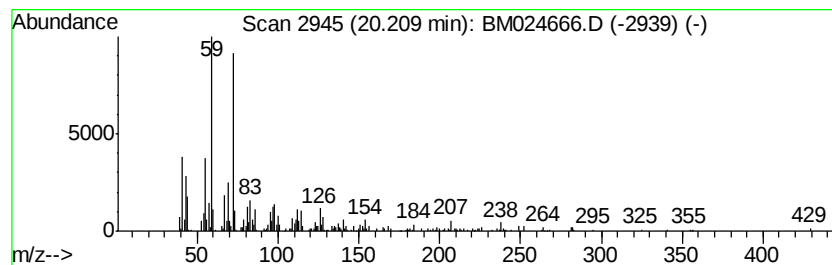
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TIC Integration Parameters: LSCINT.P

Peak Number 3 9-Octadecenamide, (Z)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.21	2.50 ng	398239	Chrysene-d12	21.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	98
2		Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1	43
3		2,11-Dimethyl-2,11-dodecanediol	230	C14H30O2	022092-59-7	35
4		2-Hydroxy-2-methylundec-5-en-3-one	198	C12H22O2	1000191-46-2	27
5		Bicyclo[3.2.1]heptane, 1-methoxy-	126	C8H14O	1000156-43-7	27



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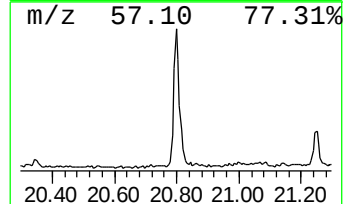
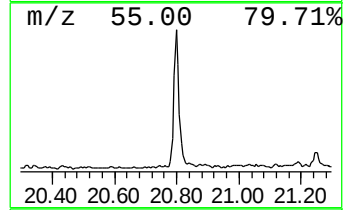
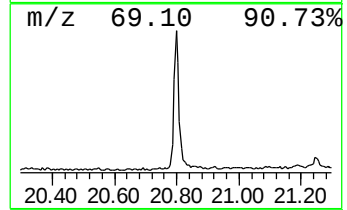
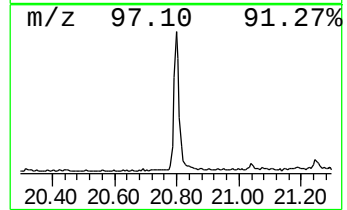
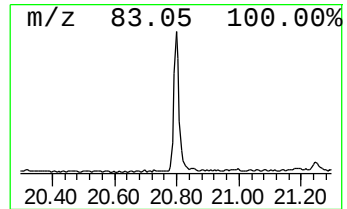
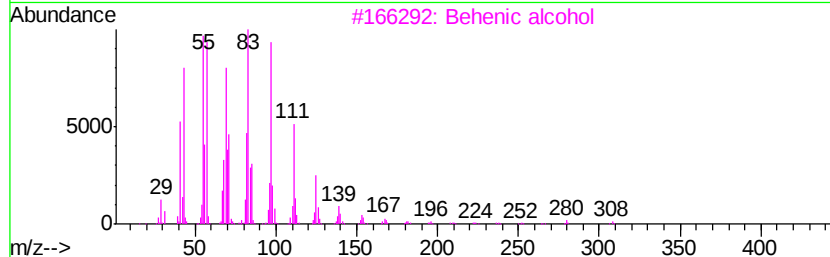
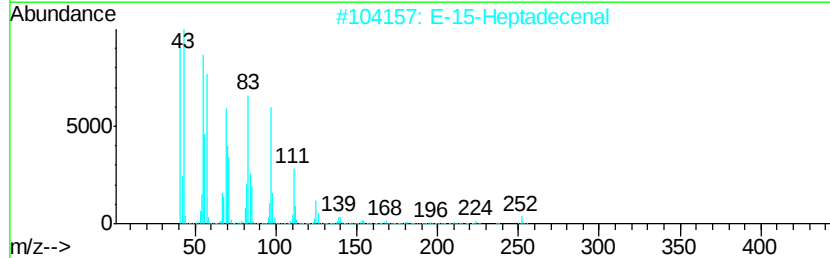
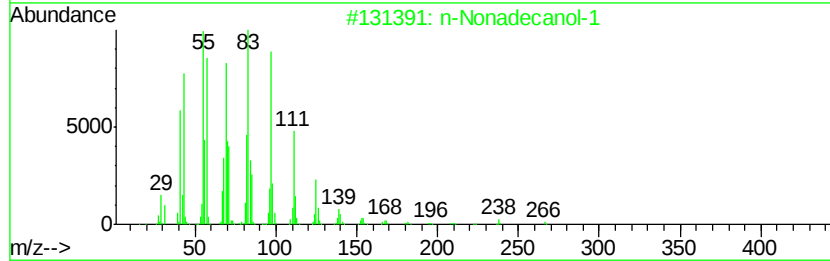
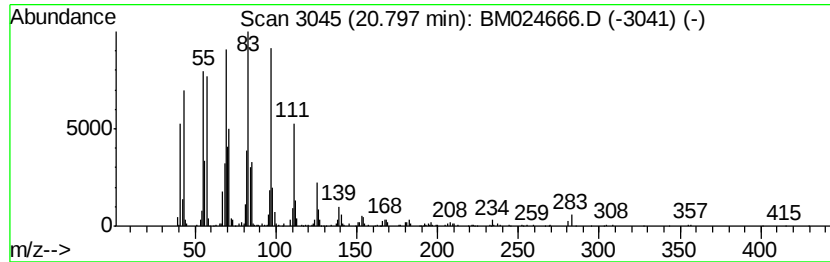
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 TIC Integration Parameters: LSCINT.P

 Peak Number 4 n-Nonadecanol-1 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.80	4.99 ng	794308	Chrysene-d12		21.04	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Nonadecanol-1		284	C19H40O	001454-84-8	94
2	E-15-Heptadecenal		252	C17H32O	1000130-97-9	92
3	Behenic alcohol		326	C22H46O	000661-19-8	91
4	1-Heneicosyl formate		340	C22H44O2	077899-03-7	91
5	Trifluoroacetoxy hexadecane		338	C18H33F3O2	006222-03-3	91



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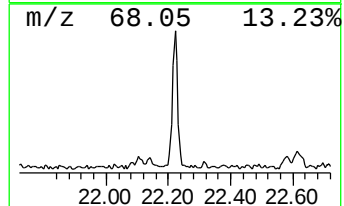
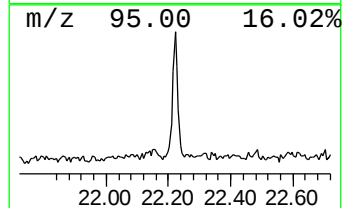
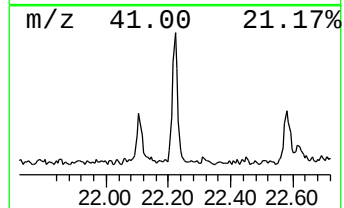
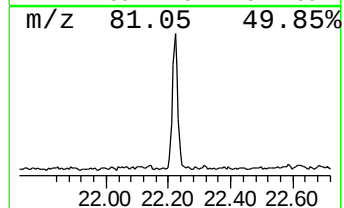
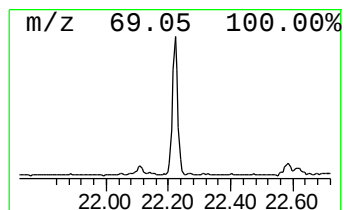
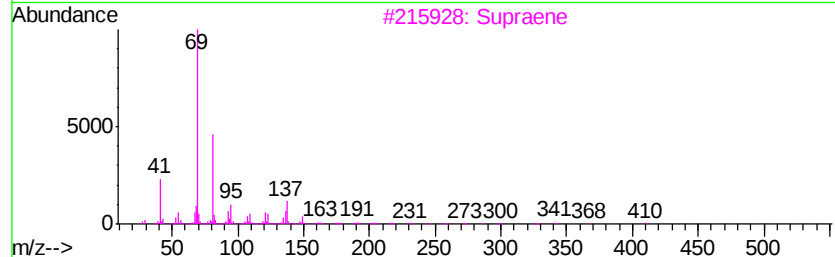
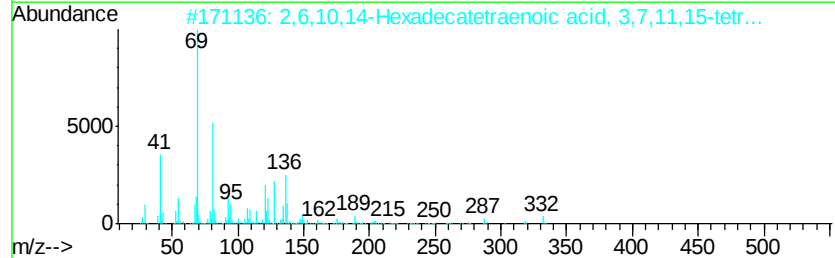
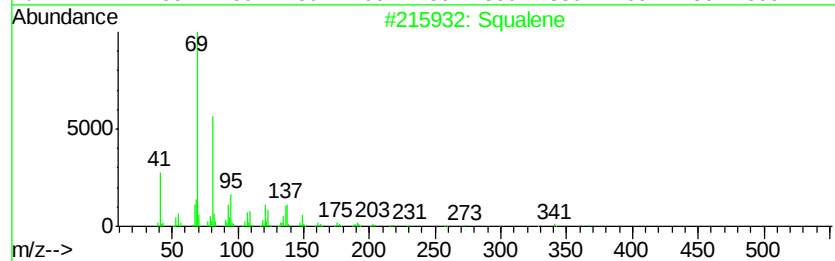
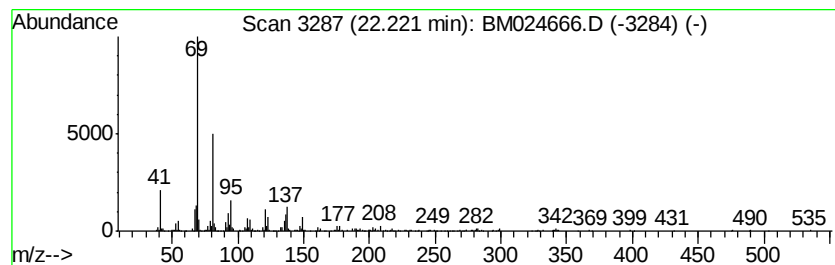
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Peak Number 5 Squalene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.22	2.70 ng	436586	Perylene-d12	23.15
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Squalene		410 C30H50	000111-02-4 87
2	2,6,10,14-Hexadecatetraenoic aci...		332 C22H36O2	024035-35-6 76
3	Supraene		410 C30H50	007683-64-9 74
4	1,5,9-Undecatriene, 2,6,10-trime...		192 C14H24	062951-96-6 72
5	7,11-Dimethyldodeca-2,6,10-trien...		208 C14H24O	125529-10-4 64



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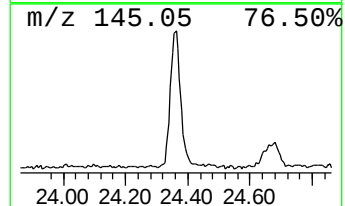
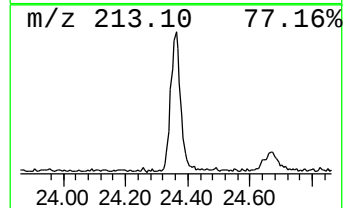
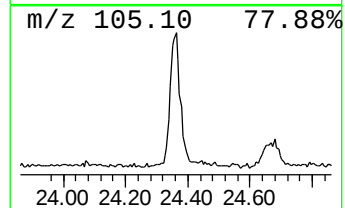
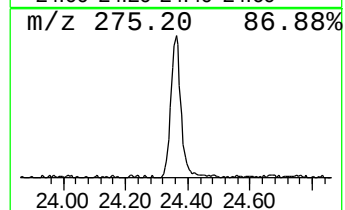
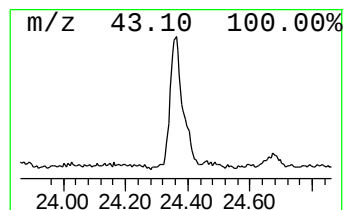
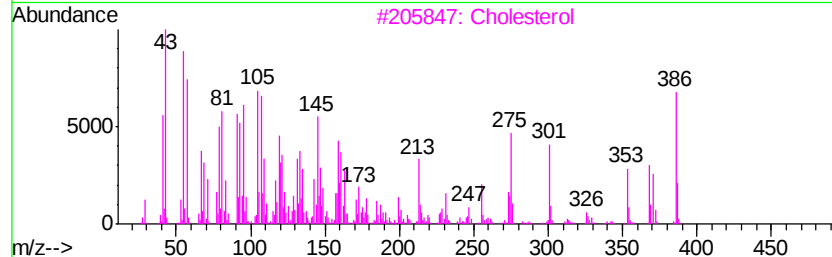
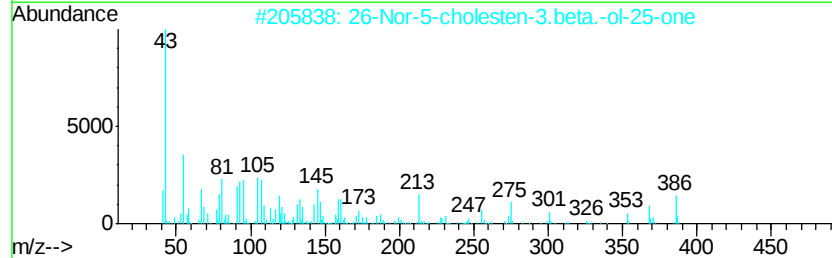
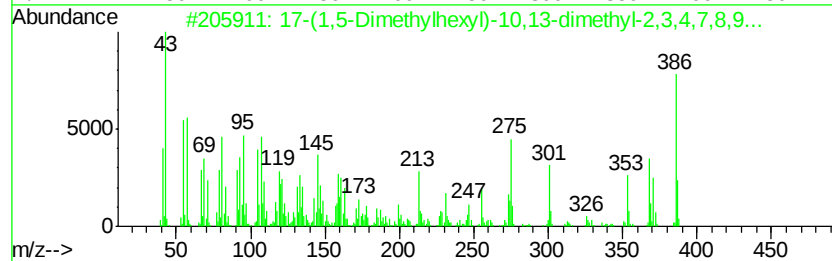
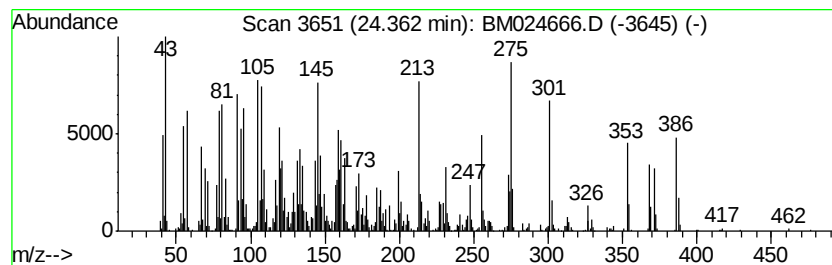
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Peak Number 6 17-(1,5-Dimethylhexyl)-10,1... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
24.36	14.63 ng	2363390	Perylene-d12	23.15		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		17-(1,5-Dimethylhexyl)-10,13-dim...	386	C27H46O	1000210-38-4	99
2		26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	99
3		Cholesterol	386	C27H46O	000057-88-5	98
4		Cholest-5-en-3-ol, (3.alpha.)-	386	C27H46O	000474-77-1	68
5		Cholestane-3,5-diol, 5-acetate, ...	446	C29H50O3	041721-93-1	43



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
n-Hexadecanoic acid	17.77	4.0	ng	412509	4	16.83	2053700	20.0
9-Octadecenamide,...	20.21	2.5	ng	398239	5	21.04	3183410	20.0
n-Nonadecanol-1	20.80	5.0	ng	794308	5	21.04	3183410	20.0
Squalene	22.22	2.7	ng	436586	6	23.15	3229810	20.0
17-(1,5-Dimethylh...	24.36	14.6	ng	2363390	6	23.15	3229810	20.0