

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP061920\
 Data File : BP002471.D
 Acq On : 19 Jun 2020 22:16
 Operator : CG/JU
 Sample : L3037-06
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SL-WC-03C

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_P\METHODS\8270-BP060520.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	5.211	388	395	404	rBV	4418765	6773647	77.58%	9.647%
2	6.781	650	662	677	rBV	4097026	7040335	80.63%	10.027%
3	7.587	792	799	806	rBV	1046576	1585137	18.15%	2.258%
4	7.905	845	853	865	rBV	3735907	6172456	70.69%	8.791%
5	8.734	980	994	1008	rBV	2903190	4930254	56.47%	7.022%
6	10.357	1263	1270	1282	rBV	1206748	2028390	23.23%	2.889%
7	12.851	1687	1694	1711	rVB	5532109	8731513	100.00%	12.436%
8	13.693	1829	1837	1844	rBV	333767	538379	6.17%	0.767%
9	13.957	1876	1882	1887	rBV2	119648	196225	2.25%	0.279%
10	14.075	1896	1902	1915	rVB2	148745	334200	3.83%	0.476%
11	14.234	1922	1929	1936	rBV	1482326	2183984	25.01%	3.110%
12	15.734	2177	2184	2192	rBV	3438669	5152087	59.01%	7.338%
13	16.028	2231	2234	2243	rVB	78384	109656	1.26%	0.156%
14	16.992	2392	2398	2402	rBV	1451894	2134114	24.44%	3.039%
15	17.034	2402	2405	2410	rVB	465268	588360	6.74%	0.838%
16	17.122	2416	2420	2426	rVB	199643	271637	3.11%	0.387%
17	17.869	2543	2547	2551	rBV	84302	114085	1.31%	0.162%
18	17.922	2551	2556	2563	rBV2	1062642	1460272	16.72%	2.080%
19	18.051	2575	2578	2583	rBV2	100421	155464	1.78%	0.221%
20	18.763	2695	2699	2704	rBV2	104400	163255	1.87%	0.233%
21	19.016	2737	2742	2746	rBV	1008273	1297960	14.87%	1.849%
22	19.163	2764	2767	2774	rBV2	375695	533545	6.11%	0.760%
23	19.369	2794	2802	2811	rVB	942233	1419792	16.26%	2.022%
24	19.580	2833	2838	2851	rVB	6387889	8490717	97.24%	12.093%
25	19.957	2899	2902	2906	rVB	108860	134376	1.54%	0.191%
26	20.839	3048	3052	3058	rBV4	797272	1257666	14.40%	1.791%
27	21.098	3089	3096	3099	rBV2	1862801	3066184	35.12%	4.367%
28	21.133	3099	3102	3110	rVB	562972	736828	8.44%	1.049%
29	22.639	3354	3358	3361	rBV	415584	744326	8.52%	1.060%
30	23.298	3465	3470	3476	rVB	1056295	1868624	21.40%	2.661%

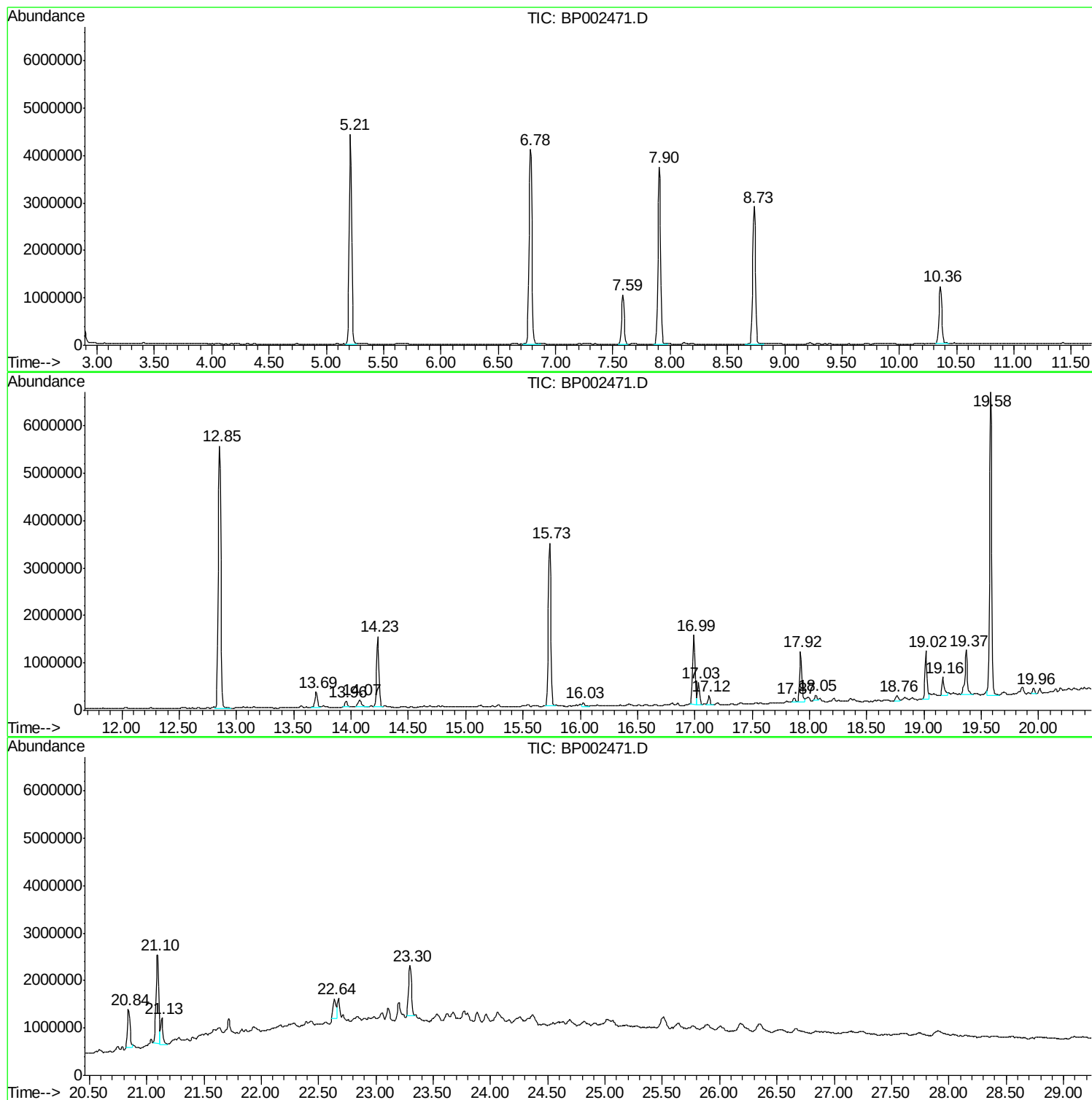
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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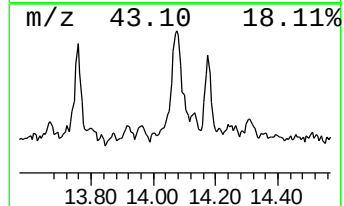
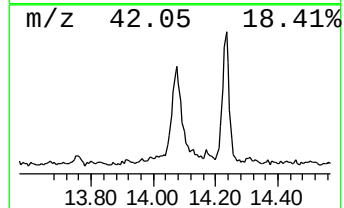
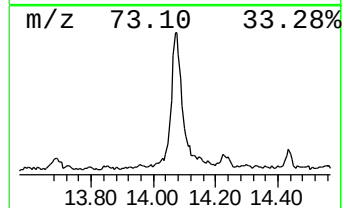
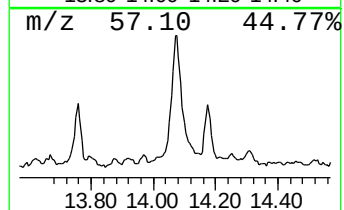
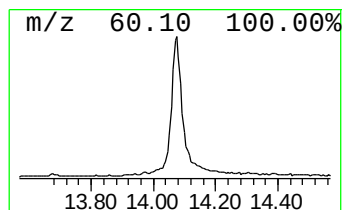
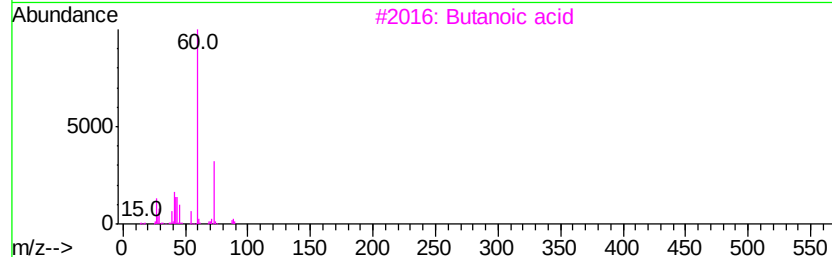
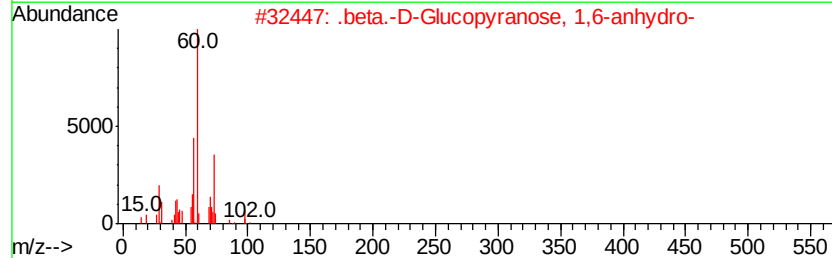
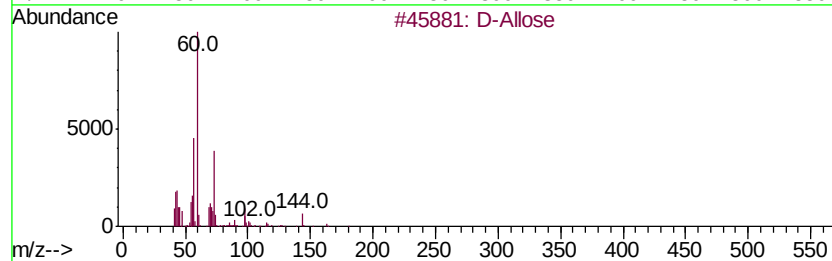
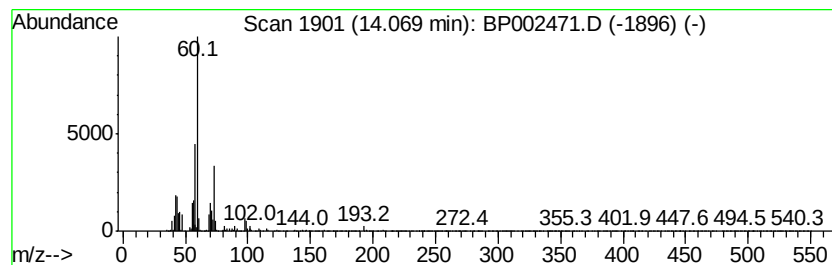
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 D-Allose Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.07	3.06 ng	334200	Acenaphthene-d10	14.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	D-Allose	180	C6H12O6	002595-97-3	72
2		.beta.-D-Glucopyranose, 1,6-anhy...	162	C6H10O5	000498-07-7	64
3		Butanoic acid	88	C4H8O2	000107-92-6	59
4		3,4-Altrosan	162	C6H10O5	1000129-76-4	53
5		Hexanoic acid	116	C6H12O2	000142-62-1	50



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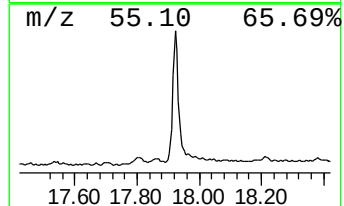
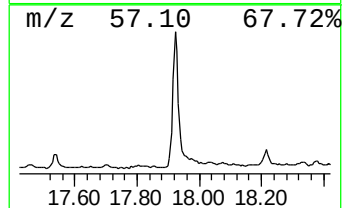
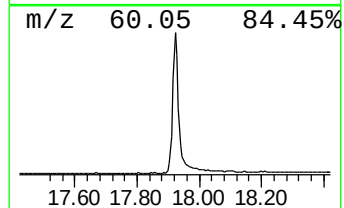
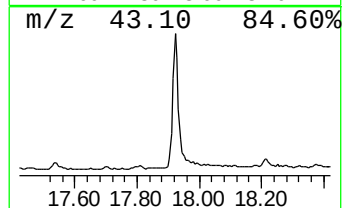
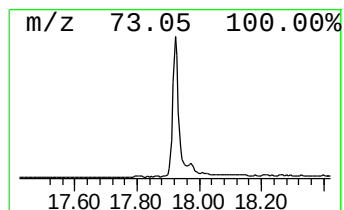
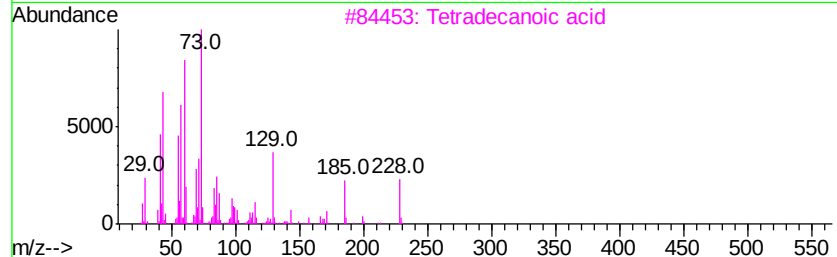
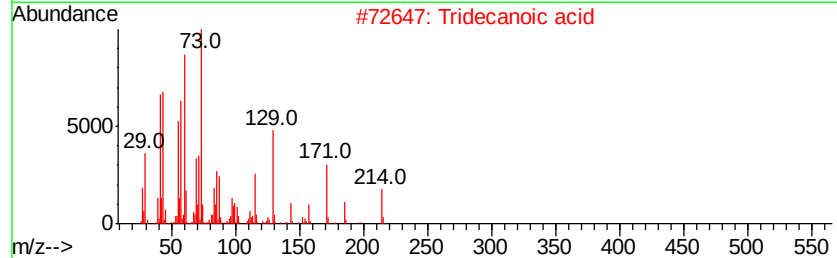
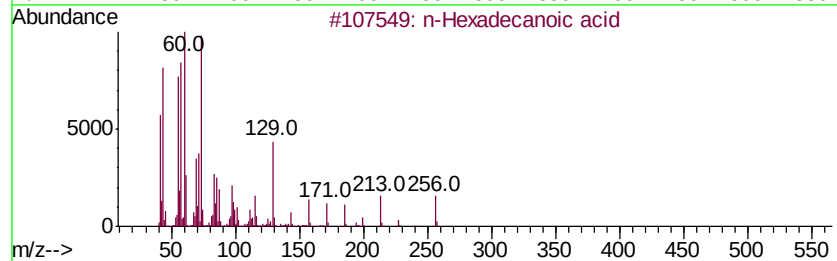
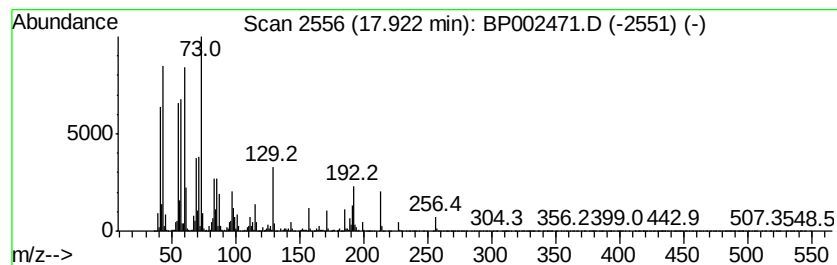
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.92	13.69 ng	1460270	Phenanthrene-d10	16.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	91
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	90
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5		Methyl-.alpha.-d-ribofuranoside	164	C6H12O5	1000129-83-1	38



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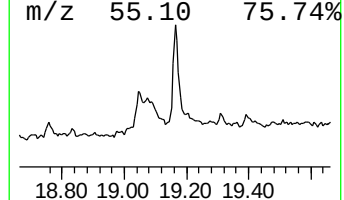
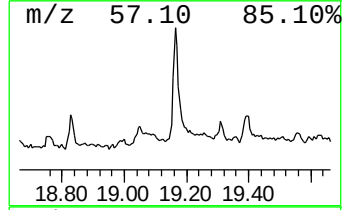
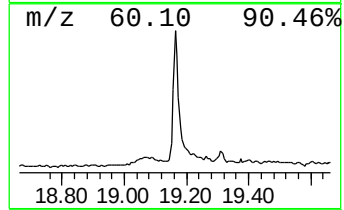
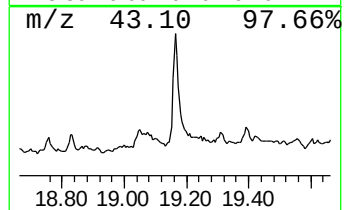
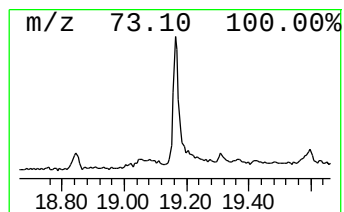
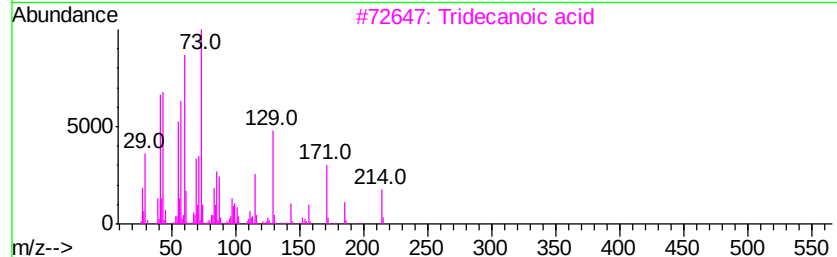
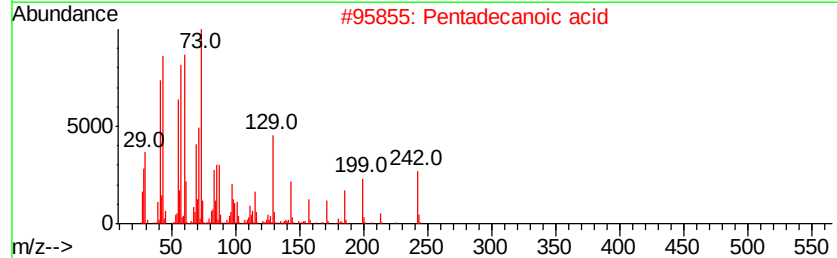
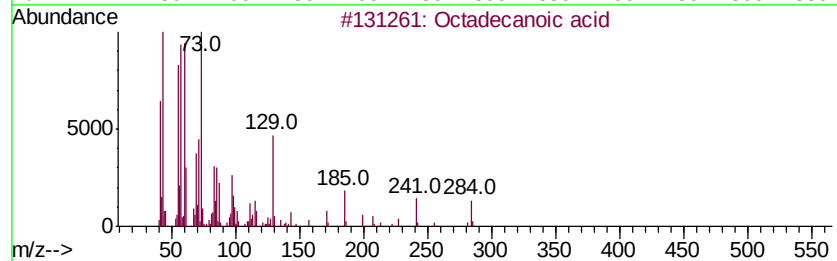
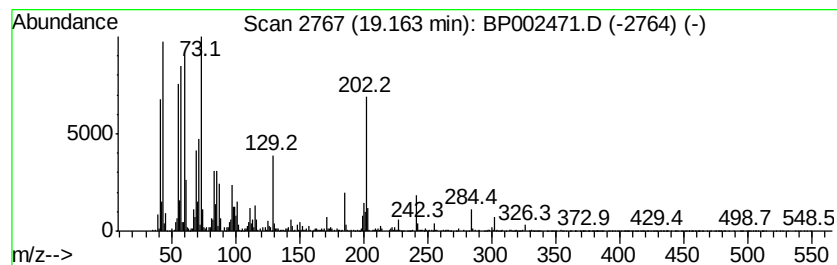
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.16	3.48 ng	533545	Chrysene-d12	21.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3		Tridecanoic acid	214	C13H26O2	000638-53-9	89
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	81
5		Octadecanoic acid, 2-(2-hydroxy...	372	C22H44O4	000106-11-6	74



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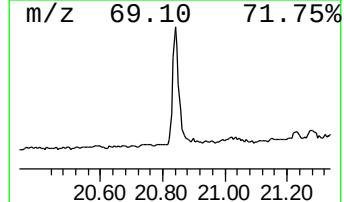
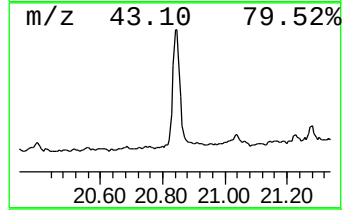
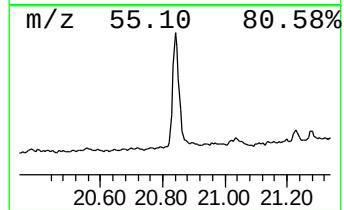
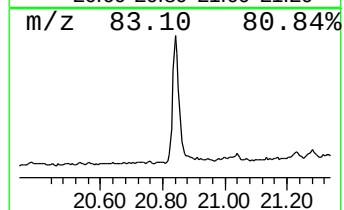
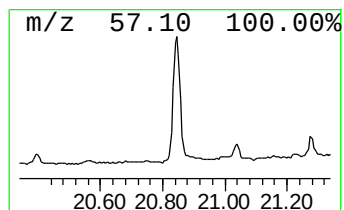
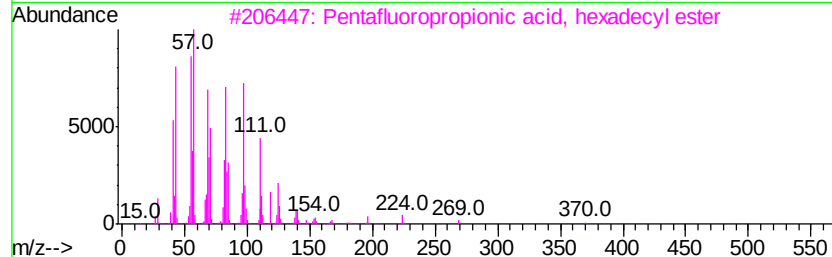
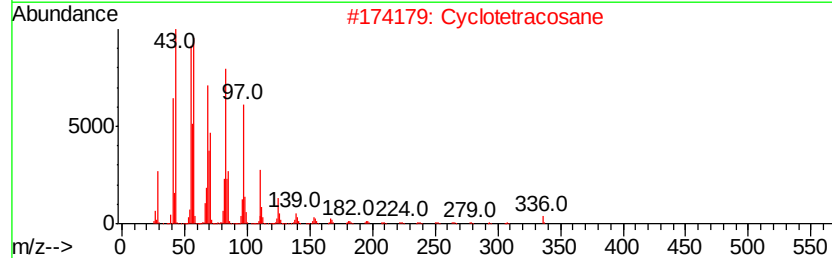
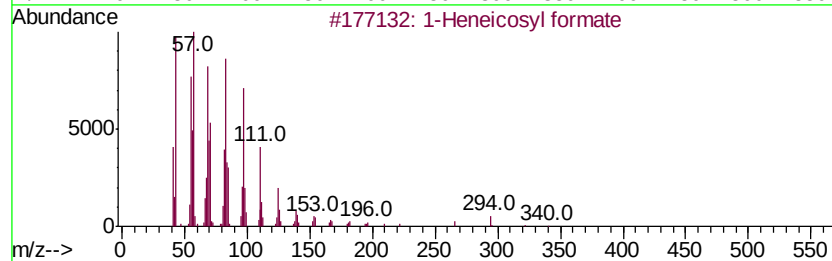
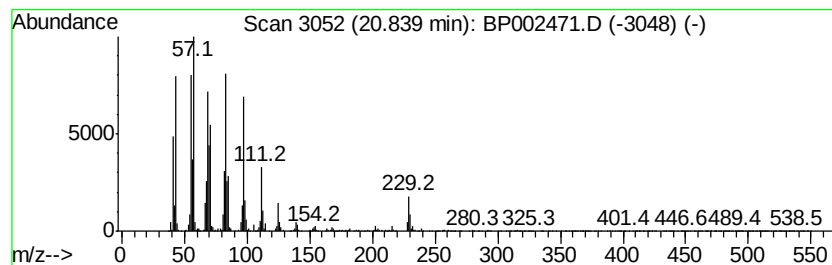
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1-Heneicosyl formate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.84	8.20 ng	1257670	Chrysene-d12		21.10	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heneicosyl formate	340	C22H44O2	077899-03-7	97
2		Cyclotetracosane	336	C24H48	000297-03-0	93
3		Pentafluoropropionic acid, hexadecyl ester	388	C19H33F5O2	006222-07-7	91
4		Trichloroacetic acid, hexadecyl ester	386	C18H33Cl3O2	074339-54-1	91
5		Dichloroacetic acid, heptadecyl ester	366	C19H36Cl2O2	1000282-98-2	91



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TIC Library : C:\DATABASE\NIST11.L
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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
D-Allose	14.07	3.1 ng		334200	3	14.23	2183980	20.0
n-Hexadecanoic acid	17.92	13.7 ng		1460270	4	16.99	2134110	20.0
Octadecanoic acid	19.16	3.5 ng		533545	5	21.10	3066180	20.0
1-Heneicosyl formate	20.84	8.2 ng		1257670	5	21.10	3066180	20.0