

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052325\
 Data File : BF142537.D
 Acq On : 23 May 2025 16:05
 Operator : RC/JU
 Sample : Q2100-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-16

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_F\Methods\8270-BF052025.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF142537.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.116	488	497	507	rBV	143192	213128	8.69%	1.228%
2	5.510	558	564	569	rBV	1446528	1915190	78.09%	11.036%
3	6.516	728	735	740	rBV	1460947	1978380	80.66%	11.400%
4	6.892	794	799	805	rBV	505448	643300	26.23%	3.707%
5	7.457	888	895	900	rBV	957563	1270657	51.81%	7.322%
6	7.710	933	938	946	rVB	25095	31754	1.29%	0.183%
7	8.181	1012	1018	1035	rVB	654022	862248	35.16%	4.969%
8	9.251	1194	1200	1206	rBV	1882534	2452664	100.00%	14.133%
9	9.933	1311	1316	1321	rBV	789156	967697	39.45%	5.576%
10	10.645	1432	1437	1445	rVB	348678	432081	17.62%	2.490%
11	10.722	1445	1450	1462	rBV	1192125	1514393	61.74%	8.726%
12	11.080	1507	1511	1522	rBV	31571	47445	1.93%	0.273%
13	11.422	1564	1569	1574	rBV	752976	927984	37.84%	5.347%
14	11.869	1640	1645	1652	rBV	43041	66957	2.73%	0.386%
15	11.945	1653	1658	1681	rVB	338176	533569	21.75%	3.075%
16	12.198	1697	1701	1709	rVB2	18033	24639	1.00%	0.142%
17	12.651	1773	1778	1786	rBV4	14925	30400	1.24%	0.175%
18	12.733	1787	1792	1803	rVB	35899	62720	2.56%	0.361%
19	13.010	1833	1839	1844	rBV	1522341	1914350	78.05%	11.031%
20	13.904	1986	1991	2000	rBV	61466	109949	4.48%	0.634%
21	14.063	2013	2018	2024	rBV	436408	549583	22.41%	3.167%
22	14.921	2159	2164	2172	rBV	137872	189139	7.71%	1.090%
23	15.557	2265	2272	2287	rBV	377783	615981	25.11%	3.549%

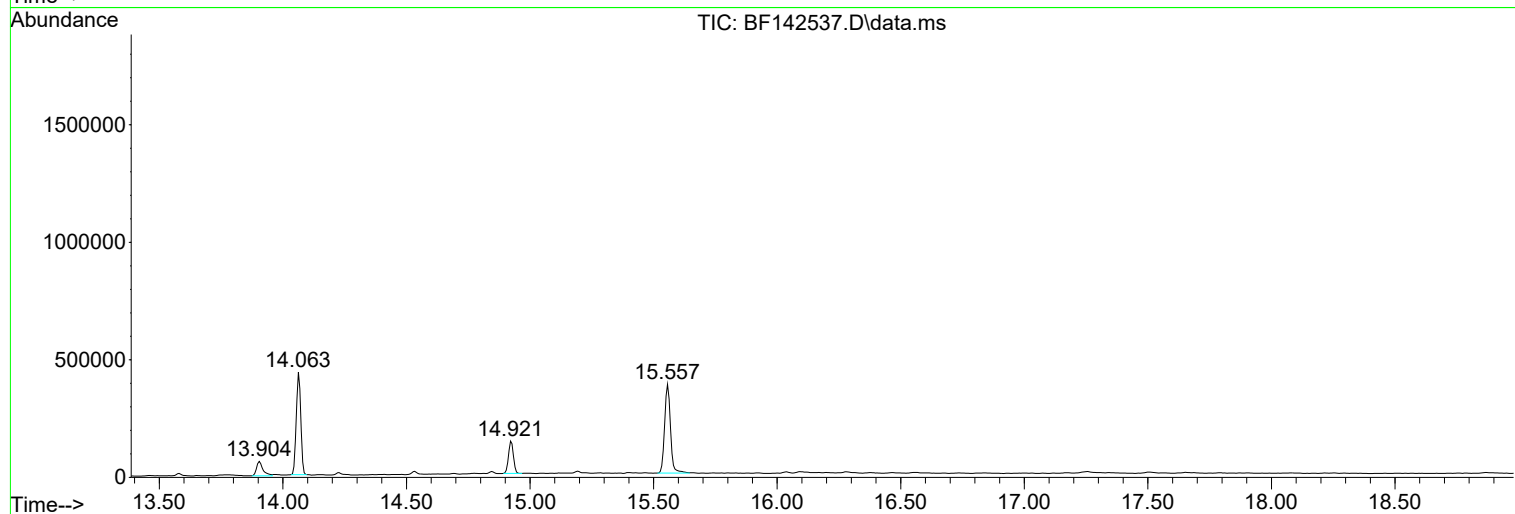
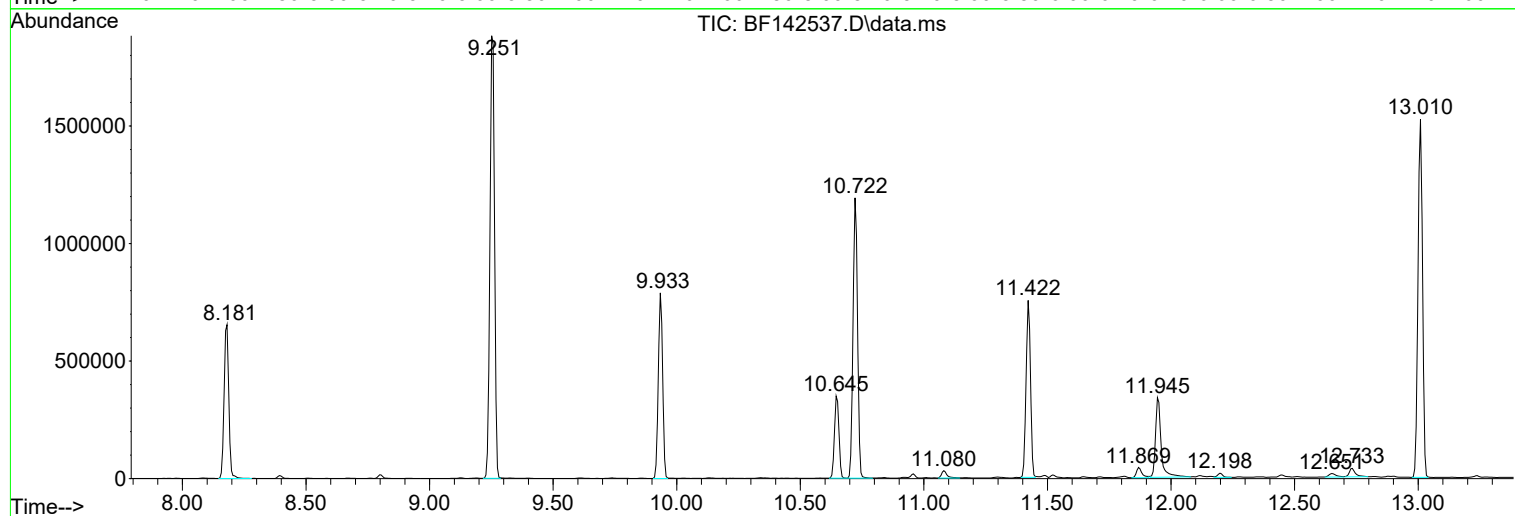
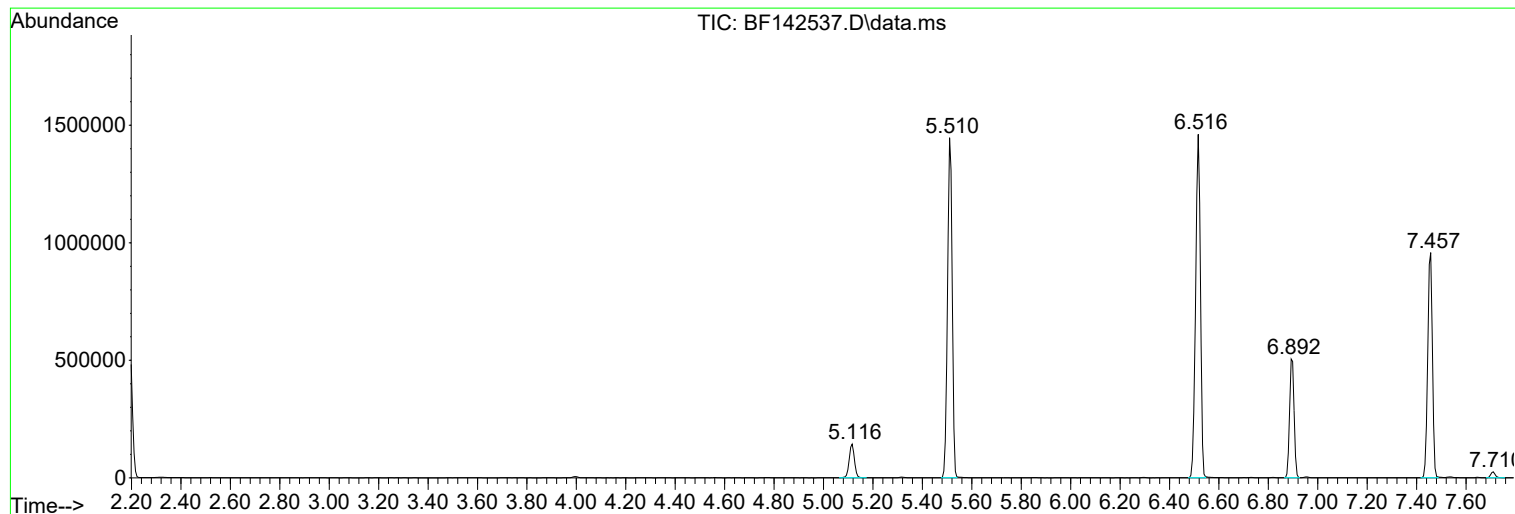
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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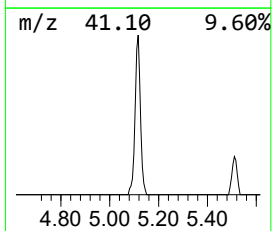
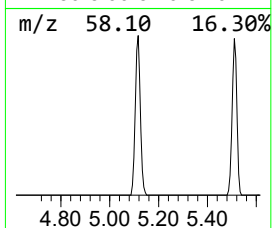
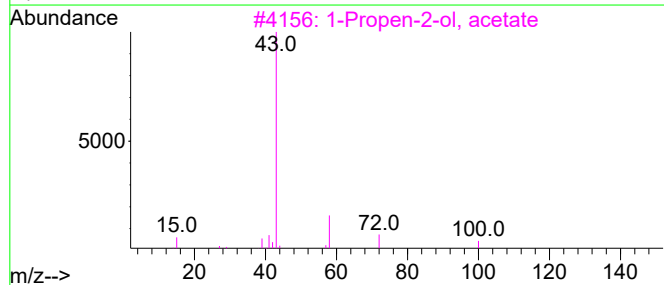
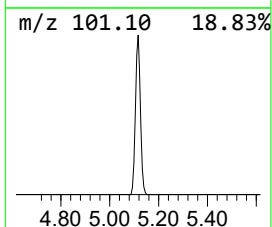
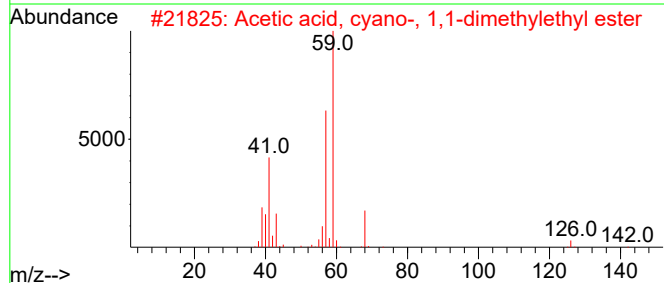
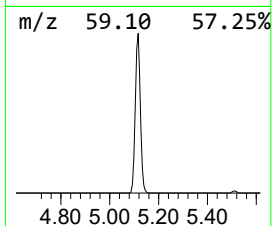
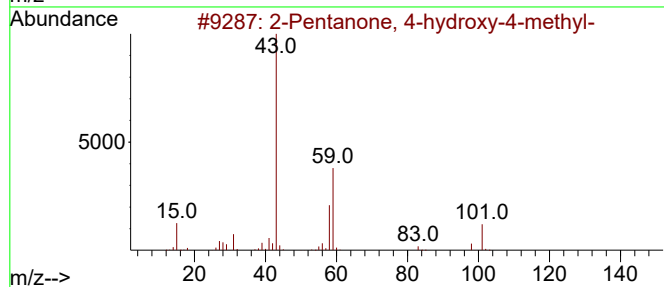
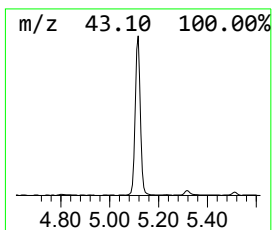
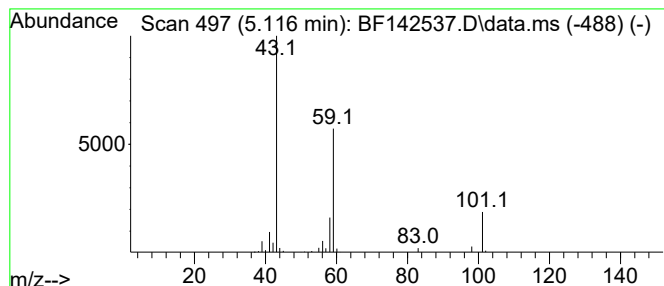
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.116	6.63 ng	213128	1,4-Dichlorobenzene-d4	6.898

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9
5			Propanoic acid, 2-nitro-, methyl...	133	C4H7NO4	006118-50-9	9



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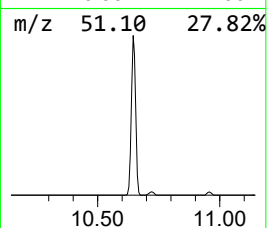
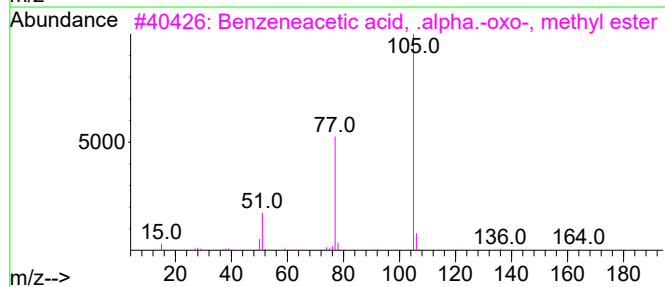
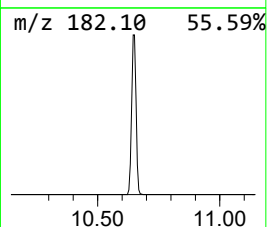
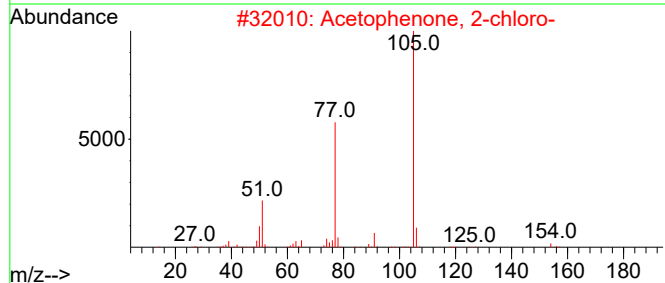
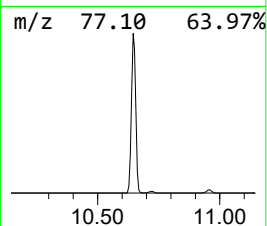
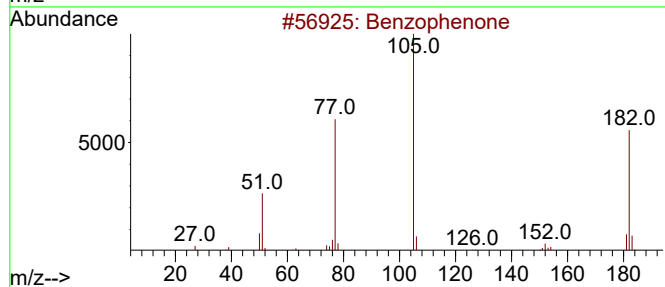
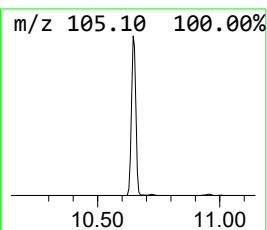
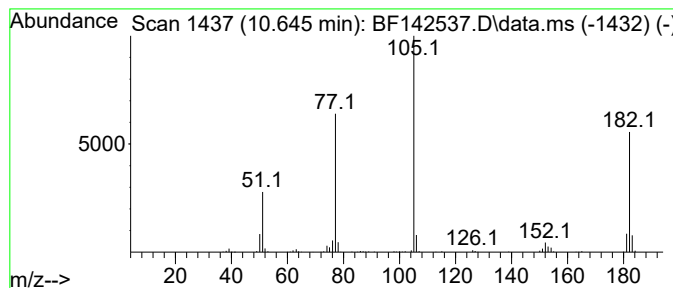
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.645	8.93 ng	432081	Acenaphthene-d10	9.933

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	49
3			Benzeneacetic acid, .alpha.-oxo-...	164	C9H8O3	015206-55-0	49
4			1-Propanone, 2-bromo-1-phenyl-	212	C9H9BrO	002114-00-3	47
5			Benzenebutanoic acid, .gamma.-oxo-	178	C10H10O3	002051-95-8	47



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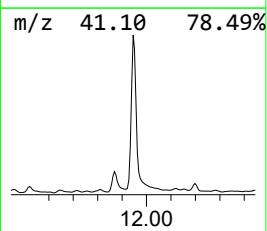
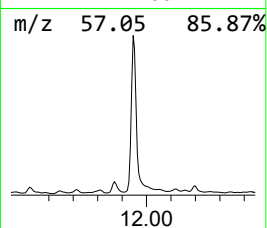
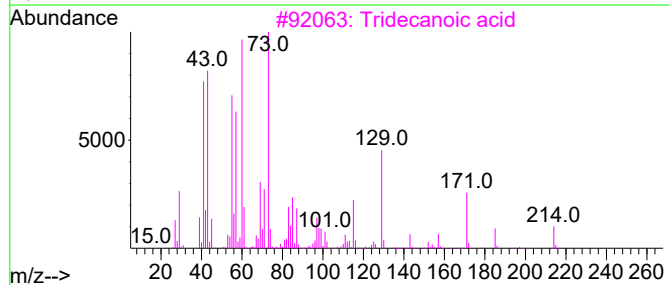
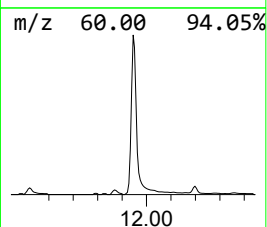
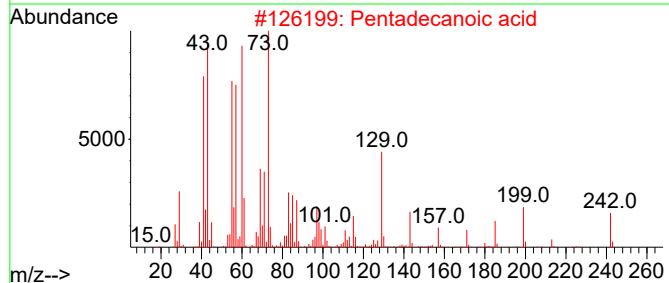
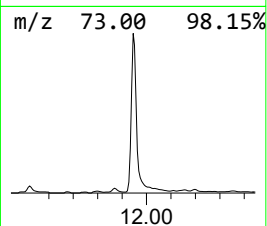
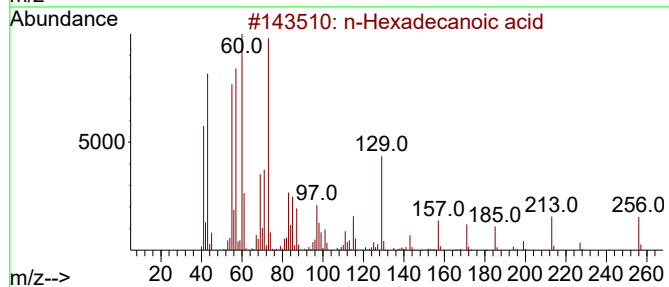
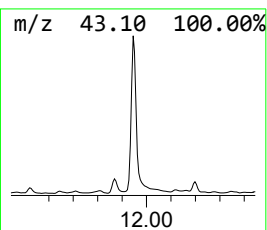
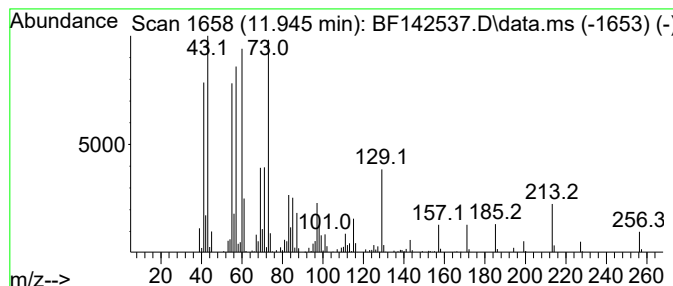
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.945	11.50 ng	533569	Phenanthrene-d10	11.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3			Tridecanoic acid	214	C13H26O2	000638-53-9	89
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5			n-Decanoic acid	172	C10H20O2	000334-48-5	62



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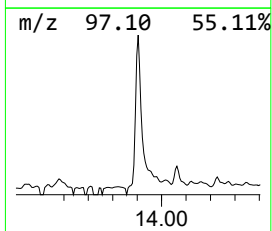
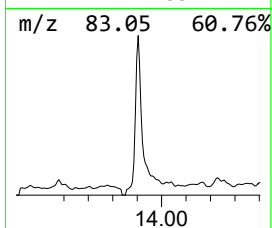
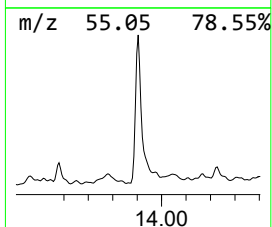
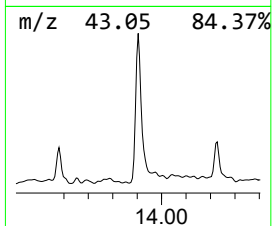
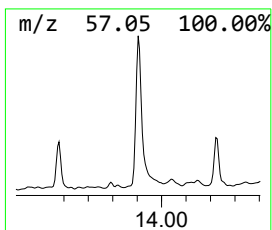
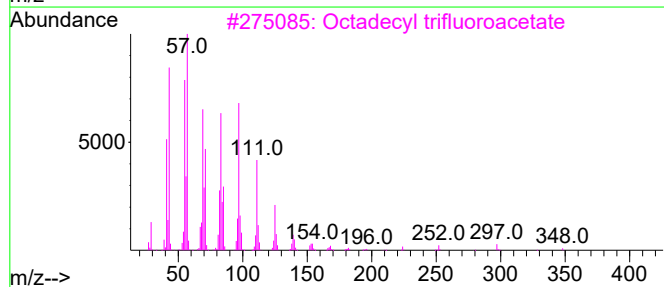
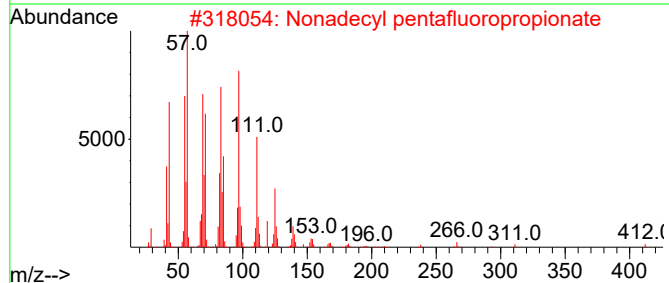
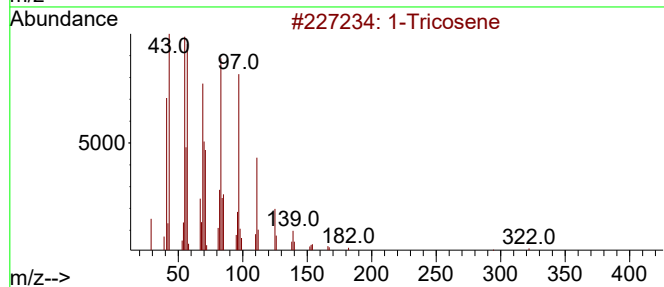
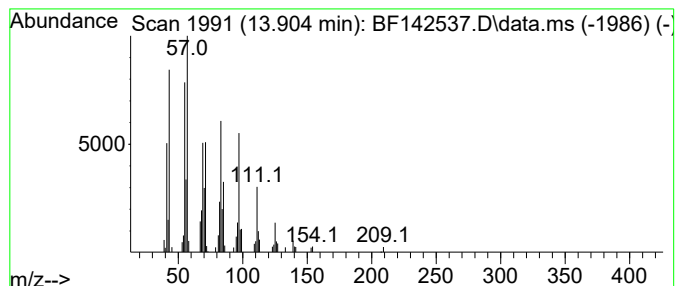
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Peak Number 4 1-Tricosene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.904	4.00 ng	109949	Chrysene-d12	14.063	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tricosene	322	C23H46	018835-32-0	93
2	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
3	Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
4	Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	91
5	Heneicosyl heptafluorobutyrate	508	C25H43F7O2	1000351-83-8	91



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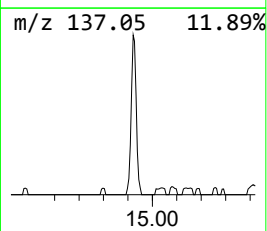
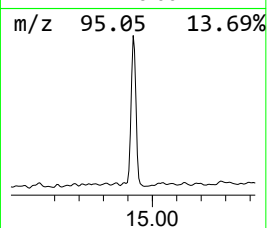
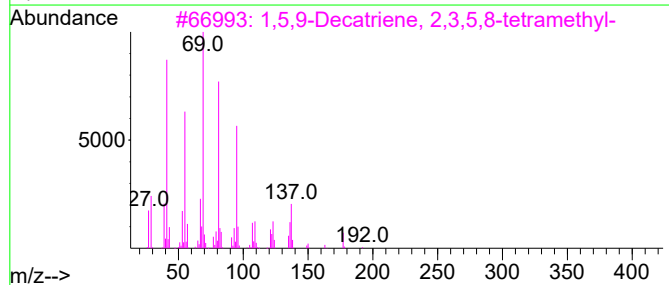
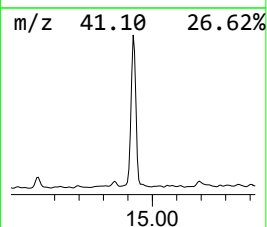
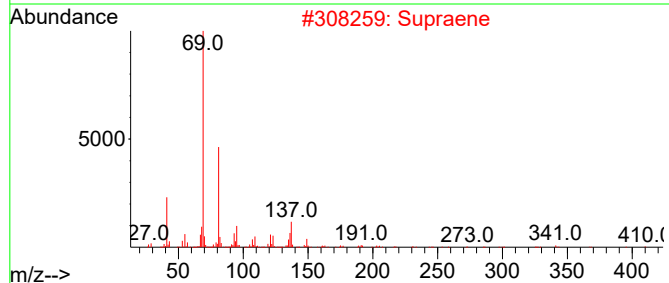
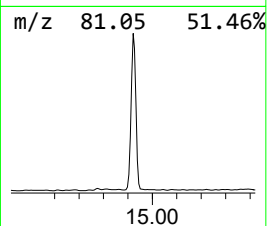
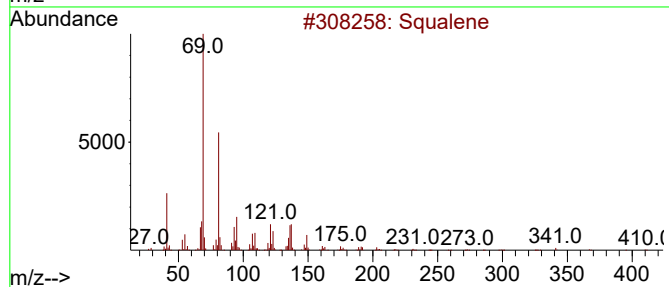
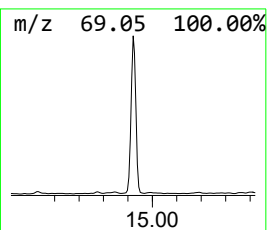
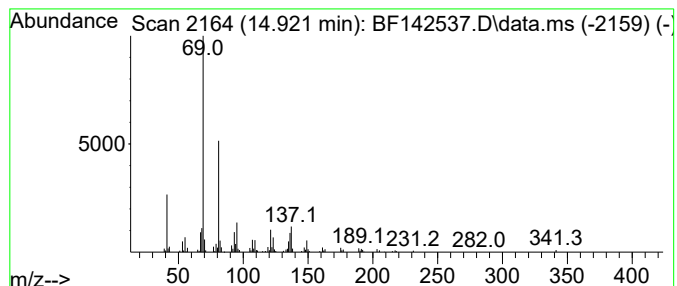
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Squalene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.921	6.14 ng	189139	Perylene-d12	15.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Squalene	410	C30H50	000111-02-4	91
2			Supraene	410	C30H50	007683-64-9	91
3			1,5,9-Decatriene, 2,3,5,8-tetram...	192	C14H24	230646-72-7	72
4			Farnesol isomer a	222	C15H26O	1000108-92-4	64
5			1,5-Heptadiene, 3,3,6-trimethyl-	138	C10H18	035387-63-4	52



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.116	6.6	ng	213128	1	6.898	643300	20.0
Benzophenone	10.645	8.9	ng	432081	3	9.933	967697	20.0
n-Hexadecanoic ...	11.945	11.5	ng	533569	4	11.422	927984	20.0
1-Tricosene	13.904	4.0	ng	109949	5	14.063	549583	20.0
Squalene	14.921	6.1	ng	189139	6	15.557	615981	20.0