

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082024\
 Data File : BP021570.D
 Acq On : 20 Aug 2024 12:59
 Operator : RC/JU
 Sample : P3650-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RW7-SP100-20240814

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-BP081324.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP021570.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.028	3	7	26	rVB	11837178	15368706	59.67%	12.357%
2	3.258	41	46	54	rBV	5682600	8020585	31.14%	6.449%
3	5.399	404	410	420	rBV	2525934	3771415	14.64%	3.032%
4	6.969	669	677	686	rBV	1429119	2319483	9.01%	1.865%
5	7.805	812	819	826	rBV	1318655	2061125	8.00%	1.657%
6	8.952	1007	1014	1024	rVB	3855909	6501035	25.24%	5.227%
7	10.593	1285	1293	1302	rBV	1705690	2892130	11.23%	2.325%
8	13.063	1705	1713	1725	rBV	9320075	14201941	55.14%	11.419%
9	14.451	1943	1949	1958	rBV	3015305	4062603	15.77%	3.266%
10	15.286	2082	2091	2097	rBV	321446	637518	2.48%	0.513%
11	15.839	2174	2185	2194	rBV	236533	410962	1.60%	0.330%
12	15.969	2200	2207	2216	rBV	8797653	12560256	48.77%	10.099%
13	17.263	2421	2427	2434	rVB	3440227	5115186	19.86%	4.113%
14	18.216	2581	2589	2596	rBV	2562864	3858080	14.98%	3.102%
15	18.280	2596	2600	2613	rVB	913537	1638583	6.36%	1.317%
16	19.410	2788	2792	2801	rBV2	349065	526333	2.04%	0.423%
17	19.545	2810	2815	2827	rBV	813023	1312338	5.10%	1.055%
18	19.992	2886	2891	2902	rVB	19688417	25754660	100.00%	20.708%
19	21.663	3171	3175	3180	rBV	613054	836949	3.25%	0.673%
20	21.727	3180	3186	3198	rVB	3429019	6009791	23.33%	4.832%
21	25.162	3761	3770	3784	rVB	2499847	6513285	25.29%	5.237%

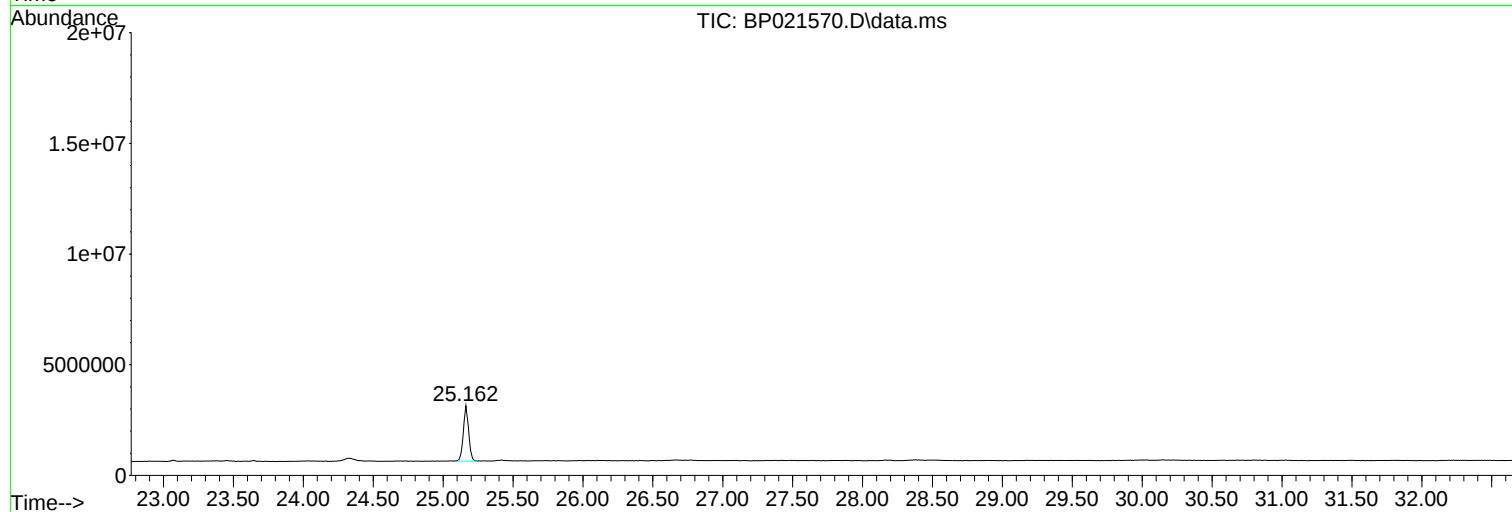
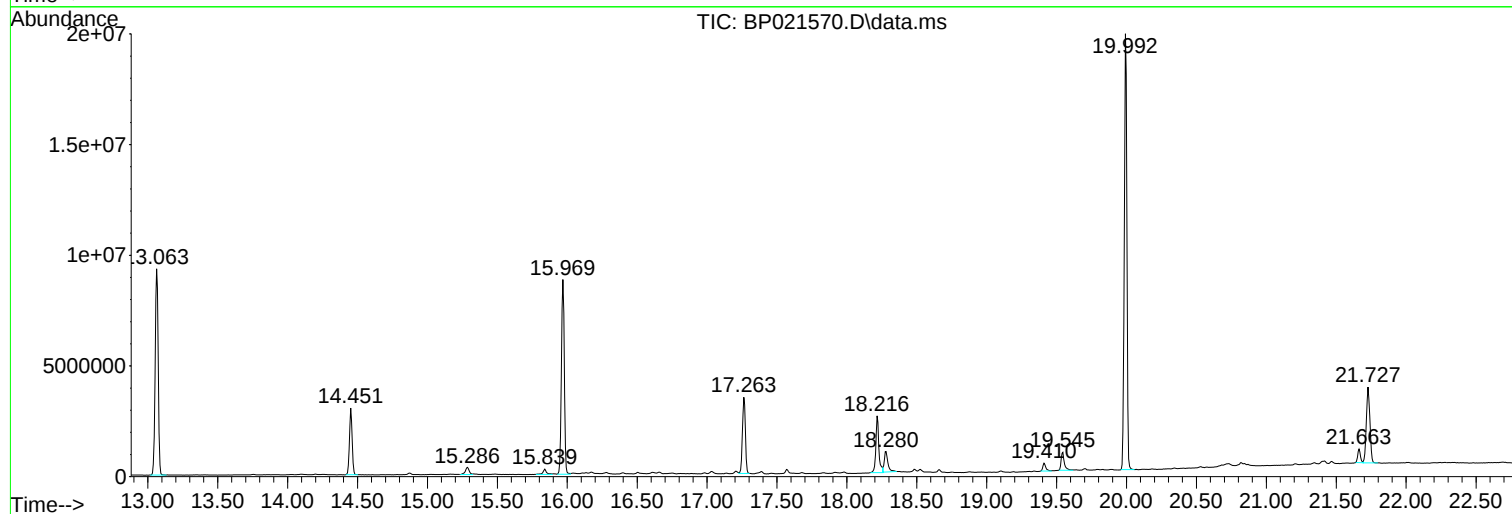
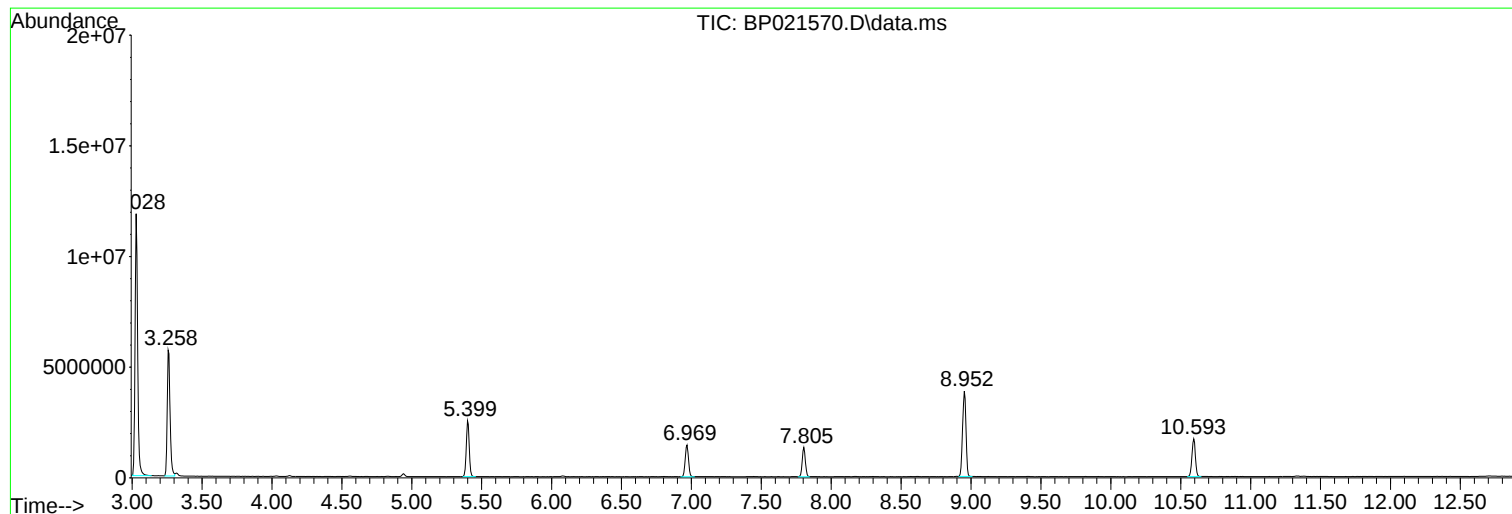
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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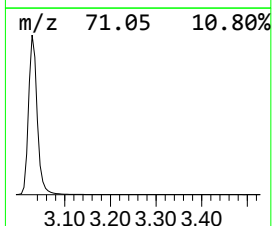
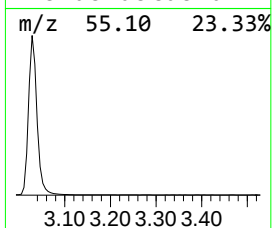
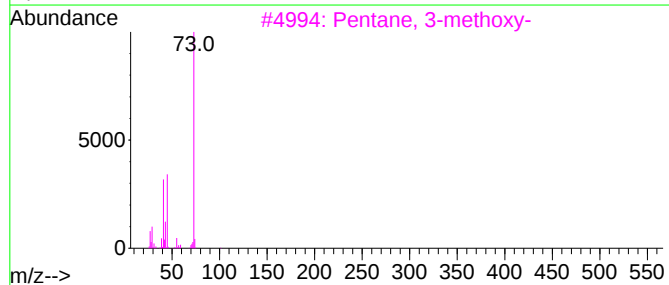
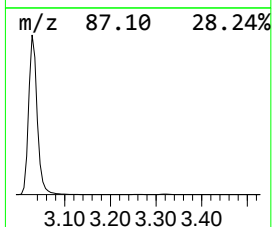
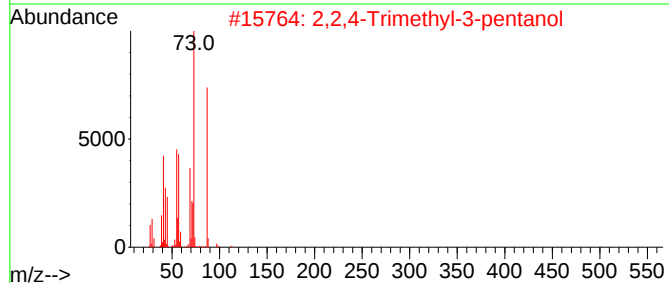
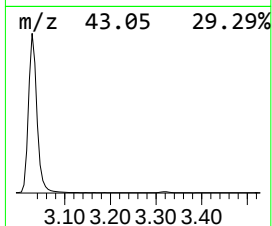
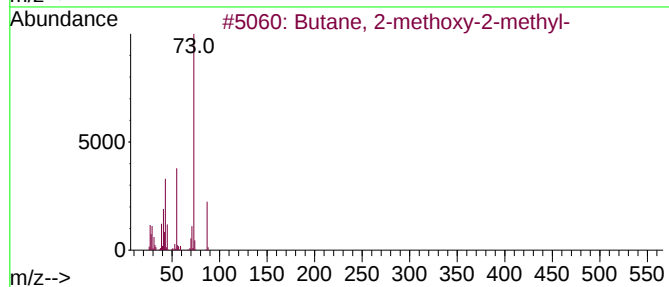
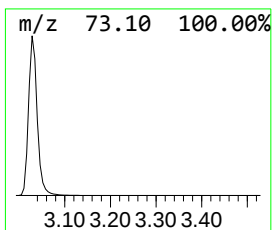
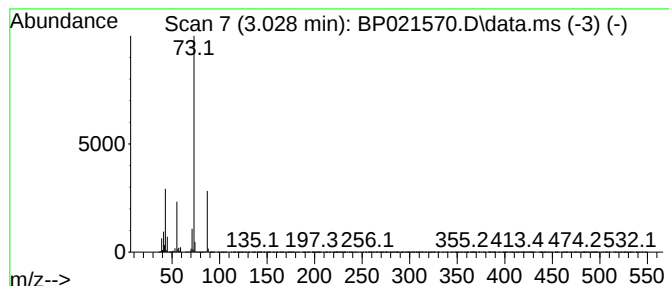
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.028	149.13 ng	15368700	1,4-Dichlorobenzene-d4	7.805

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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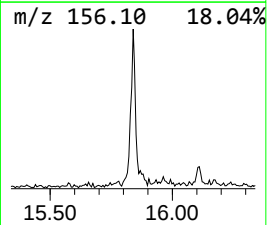
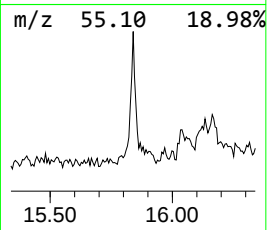
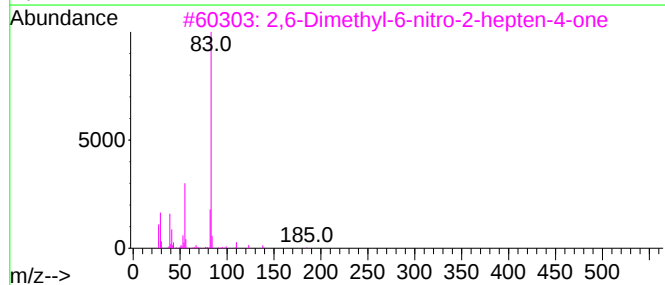
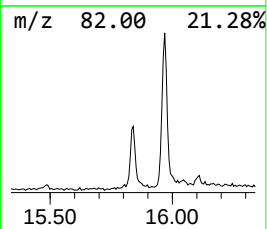
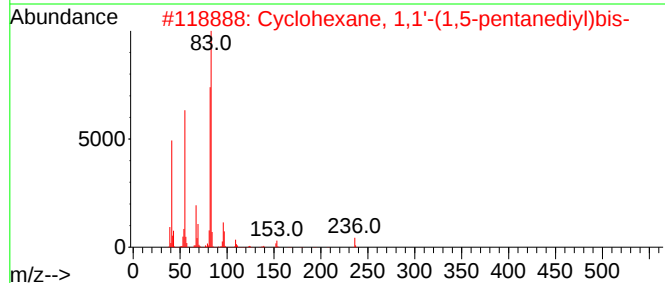
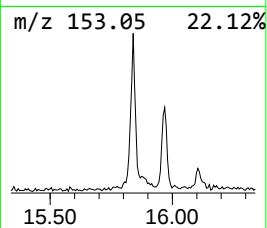
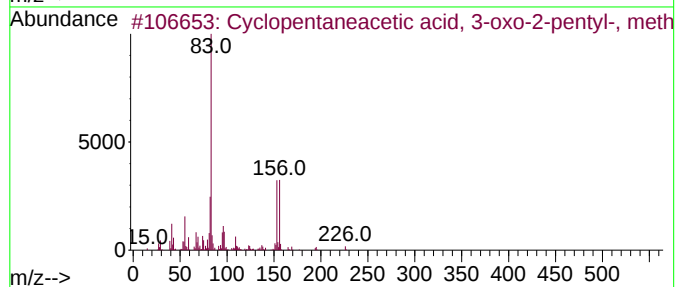
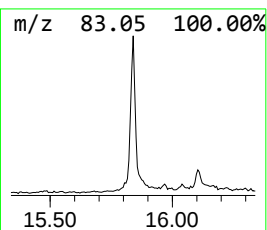
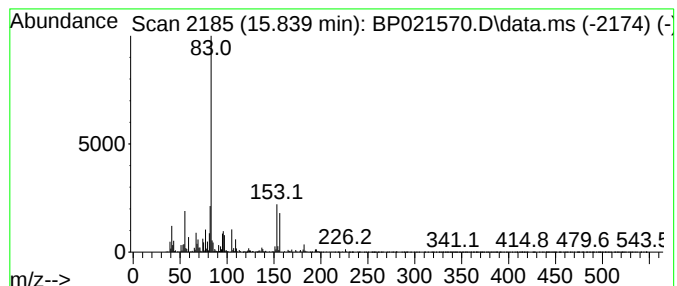
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Cyclopentaneacetic acid, 3-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.839	2.02 ng	410962	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentaneacetic acid, 3-oxo-2...	226	C13H22O3	024851-98-7	91
2			Cyclohexane, 1,1'-(1,5-pentaned...	236	C17H32	054833-31-7	47
3			2,6-Dimethyl-6-nitro-2-hepten-4-one	185	C9H15NO3	073583-56-9	47
4			6-Hydroxy-10a-methyl-4,7-dimethy...	376	C20H24O7	1798297-60-5	38
5			Hexane, 1,6-dicyclohexyl-	250	C18H34	001610-23-7	38



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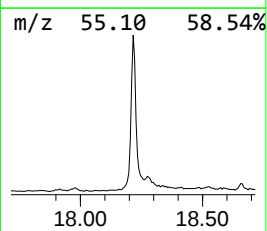
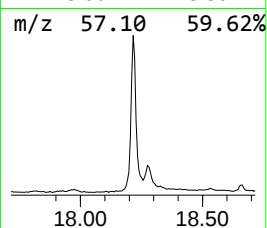
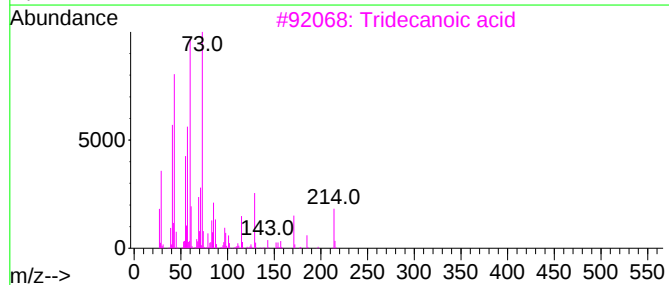
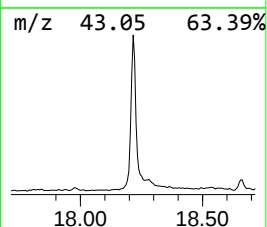
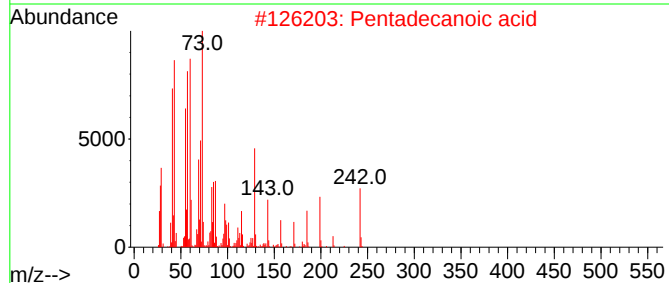
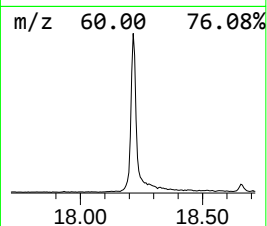
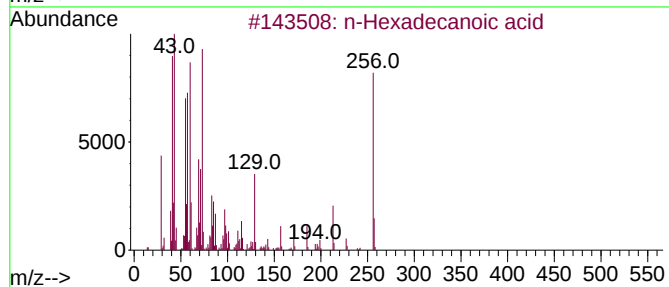
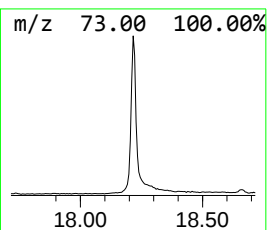
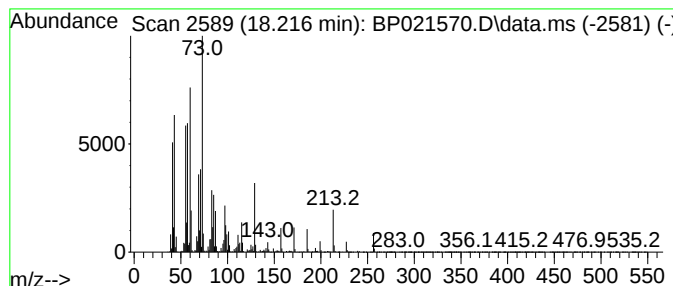
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	15.08 ng	3858080	Phenanthrene-d10	17.263

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	83
3			Tridecanoic acid	214	C13H26O2	000638-53-9	64
4			8-Methylnonanoic acid	172	C10H20O2	005963-14-4	43
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	38



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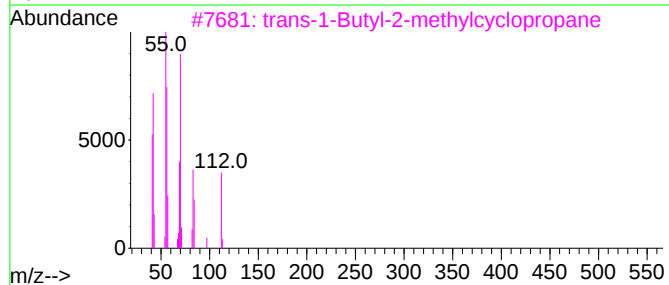
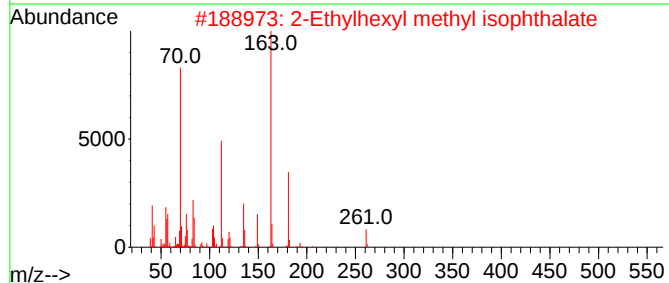
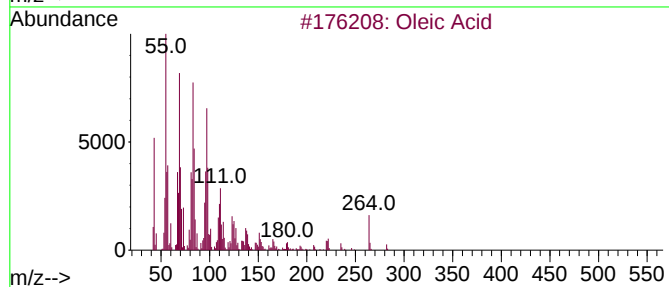
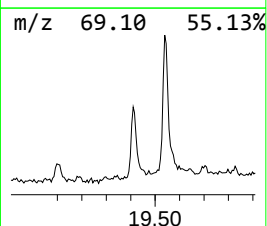
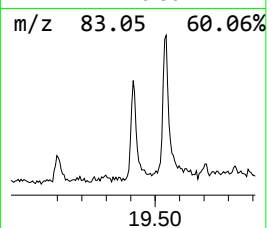
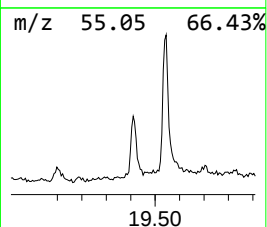
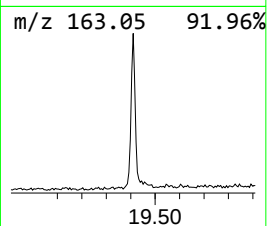
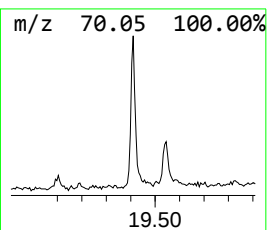
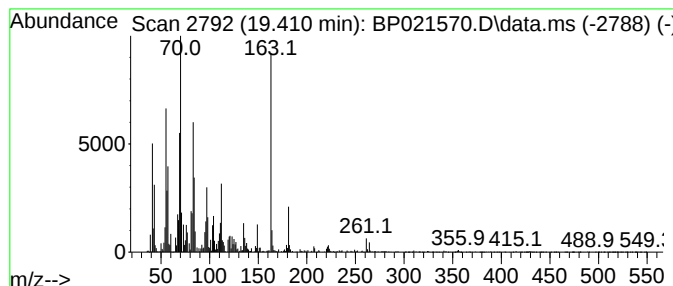
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TIC Library : C:\Database\NIST20.L
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Peak Number 5 Oleic Acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.410	2.06 ng	526333	Phenanthrene-d10	17.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oleic Acid	282	C18H34O2	000112-80-1	59
2			2-Ethylhexyl methyl isophthalate	292	C17H24O4	2135327-80-7	43
3			trans-1-Butyl-2-methylcyclopropane	112	C8H16	038851-70-6	43
4			5-Dodecene, (Z)-	168	C12H24	007206-28-2	38
5			3-Dodecene, (E)-	168	C12H24	007206-14-6	38



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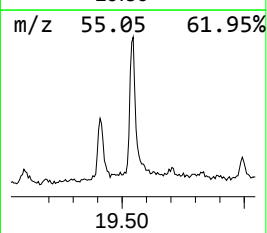
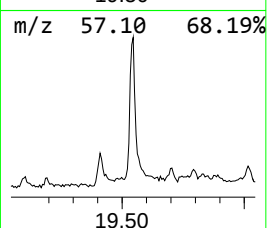
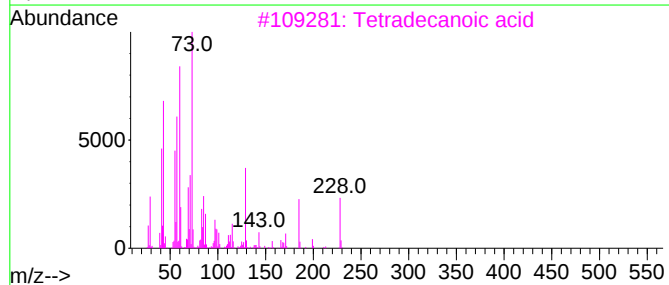
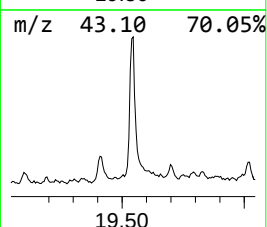
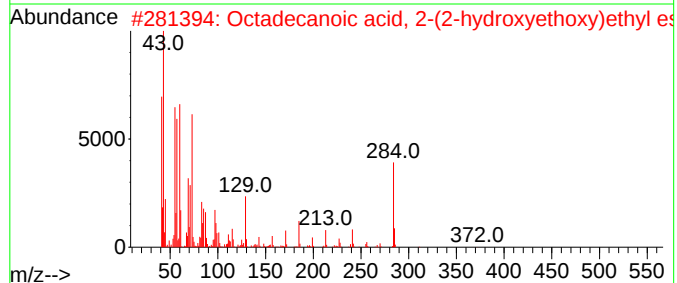
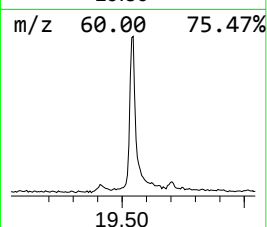
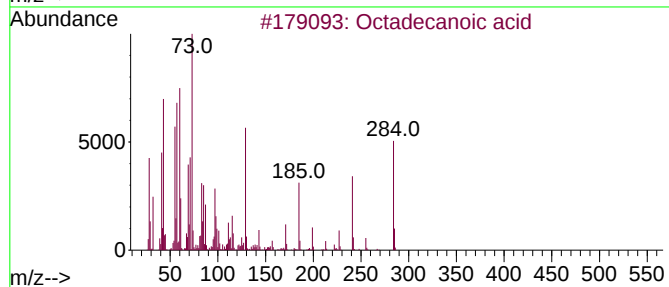
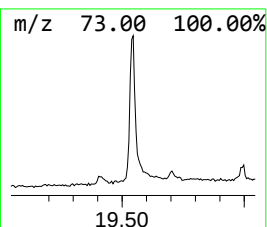
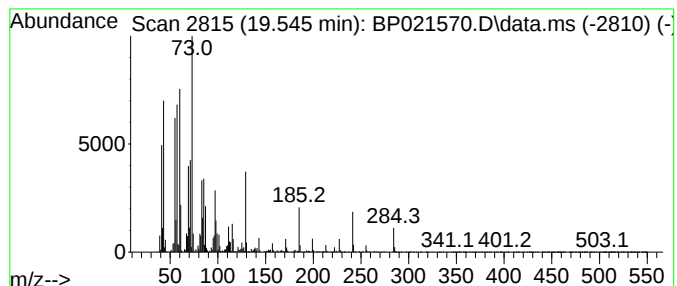
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.545	4.37 ng	1312340	Chrysene-d12	21.727

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	86
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4			n-Decanoic acid	172	C10H20O2	000334-48-5	76
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	72



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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
Butane, 2-metho...	3.028	149.1	ng	15368700	1	7.805	2061130	20.0
Cyclopentaneace...	15.839	2.0	ng	410962	3	14.451	4062600	20.0
n-Hexadecanoic ...	18.216	15.1	ng	3858080	4	17.263	5115190	20.0
Oleic Acid	19.410	2.1	ng	526333	4	17.263	5115190	20.0
Octadecanoic acid	19.545	4.4	ng	1312340	5	21.727	6009790	20.0