

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060923\
 Data File : BF133679.D
 Acq On : 09 Jun 2023 16:51
 Operator : CG\JU
 Sample : 03071-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-1

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-BF060923.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF133679.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.122	3	6	12	rVB	570215	670552	35.81%	5.014%
2	5.057	495	505	512	rBV	23388	32938	1.76%	0.246%
3	5.451	567	572	575	rBV	1718001	1712715	91.47%	12.806%
4	6.463	739	744	747	rBV	1621256	1761482	94.08%	13.171%
5	6.839	804	808	811	rBV	661349	589868	31.50%	4.411%
6	7.398	899	903	906	rBV	1157901	1121444	59.90%	8.385%
7	8.122	1022	1026	1029	rBV	895521	724756	38.71%	5.419%
8	9.198	1205	1209	1212	rBV	2414846	1872345	100.00%	14.000%
9	9.875	1321	1324	1328	rBV	826324	684686	36.57%	5.119%
10	10.592	1442	1446	1449	rBV	136446	118722	6.34%	0.888%
11	10.663	1454	1458	1462	rBV	1052511	1014686	54.19%	7.587%
12	11.363	1574	1577	1580	rBV	639795	590283	31.53%	4.414%
13	11.386	1580	1581	1585	rVB	78308	55940	2.99%	0.418%
14	11.892	1664	1667	1674	rBV2	116584	106514	5.69%	0.796%
15	12.580	1781	1784	1787	rBV	116744	90677	4.84%	0.678%
16	12.674	1797	1800	1804	rBV5	21925	27829	1.49%	0.208%
17	12.810	1818	1823	1826	rBV	97406	83983	4.49%	0.628%
18	12.957	1844	1848	1851	rBV	1458648	1220082	65.16%	9.123%
19	14.010	2023	2027	2030	rBV	473696	449201	23.99%	3.359%
20	14.480	2105	2107	2112	rBV6	27532	40295	2.15%	0.301%
21	14.768	2153	2156	2162	rVB	113870	129245	6.90%	0.966%
22	15.474	2272	2276	2280	rBV	215377	275887	14.73%	2.063%

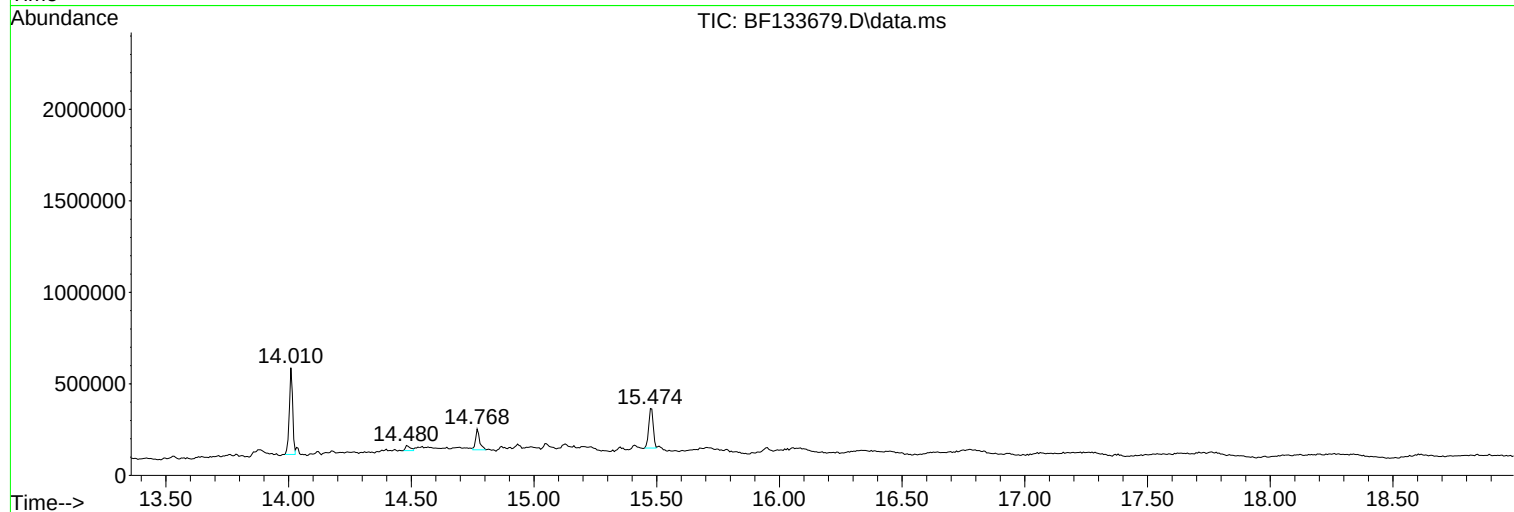
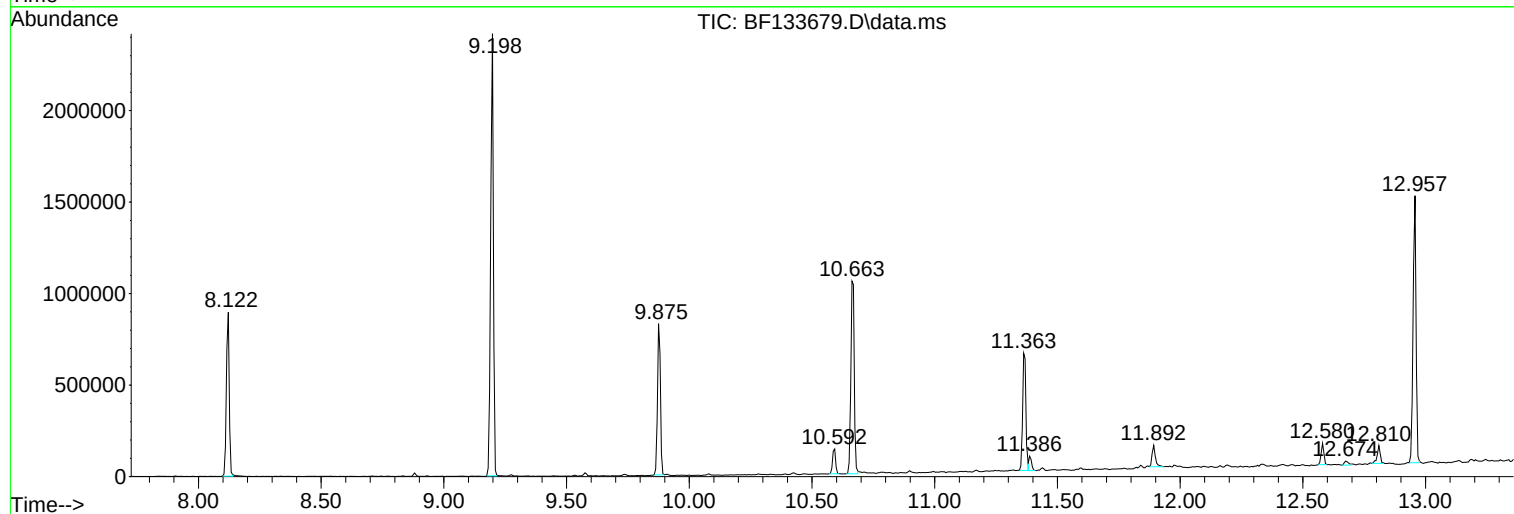
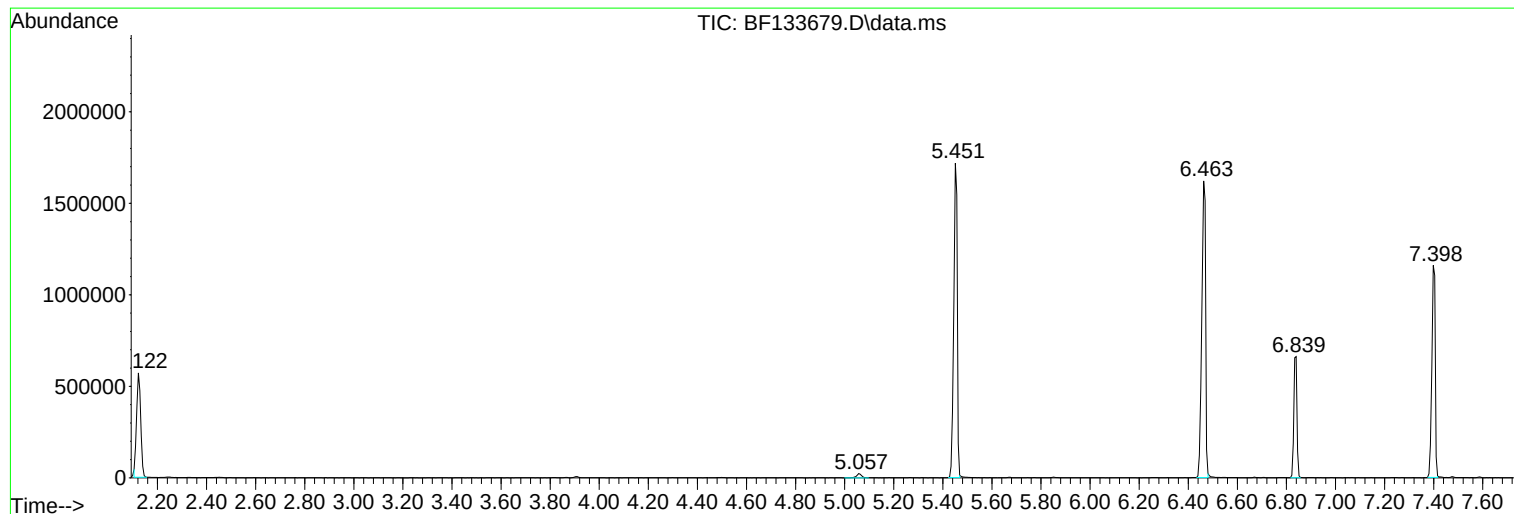
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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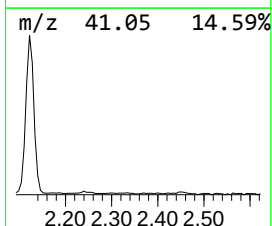
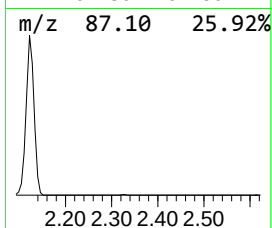
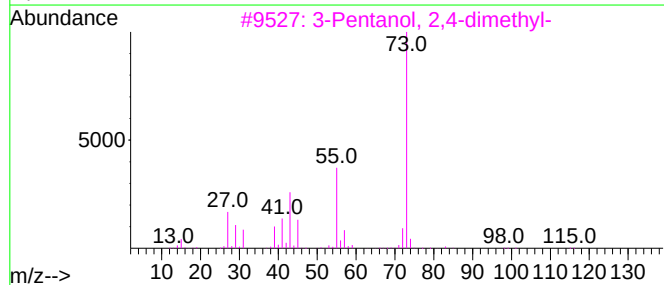
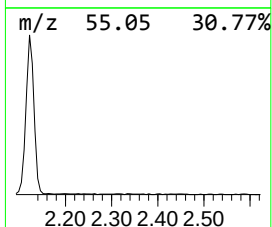
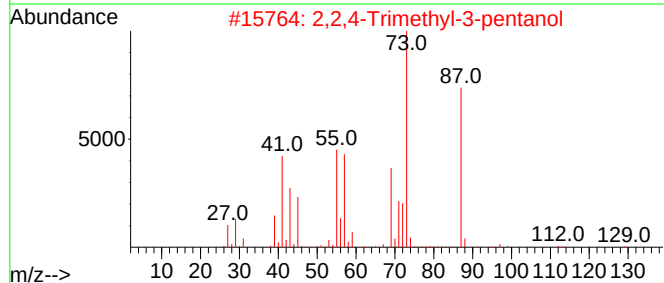
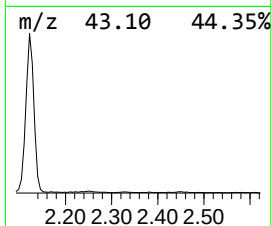
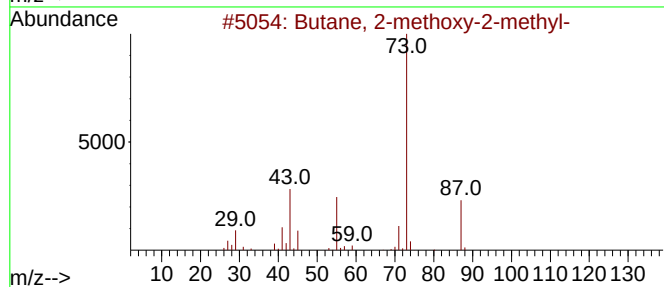
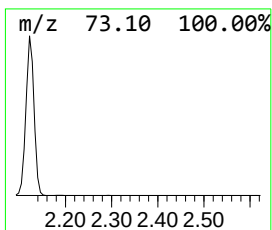
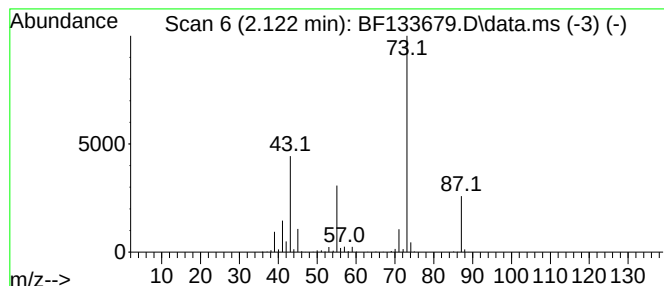
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.122	22.74 ng	670552	1,4-Dichlorobenzene-d4	6.839

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	32
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23



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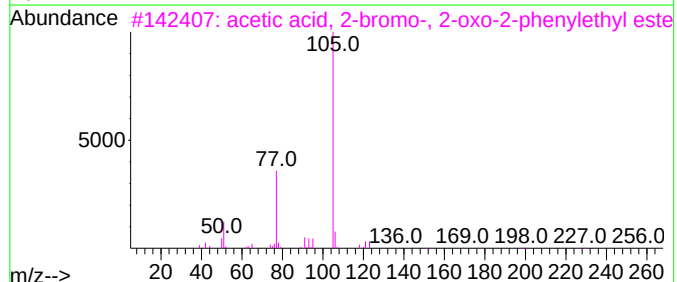
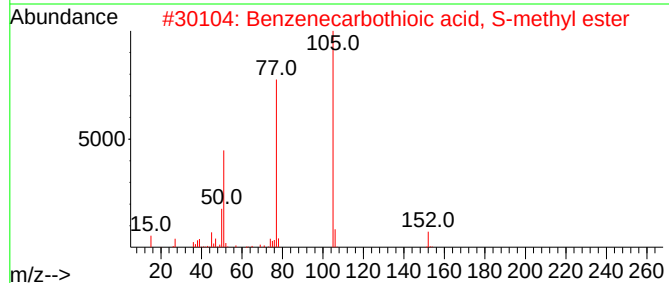
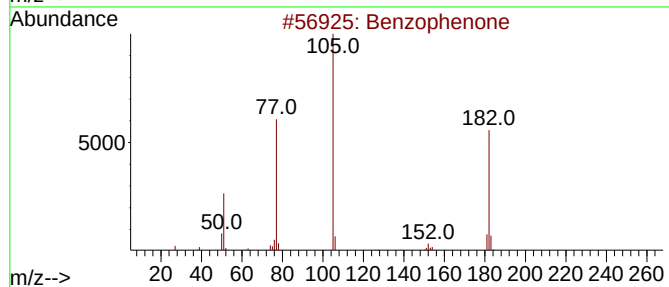
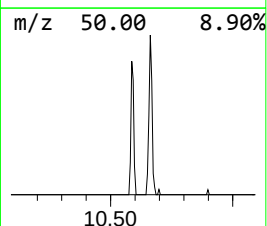
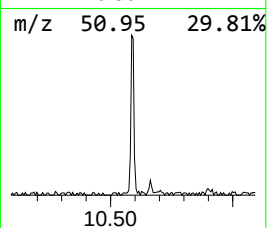
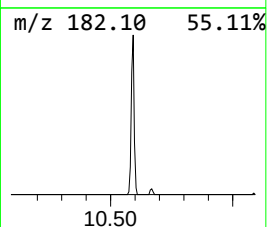
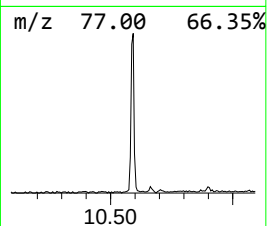
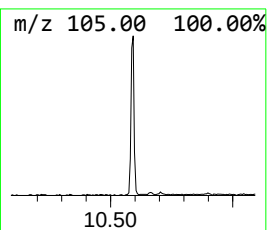
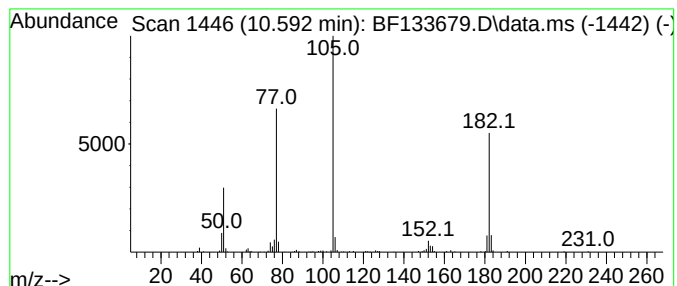
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.592	3.47 ng	118722	Acenaphthene-d10	9.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	47
3			acetic acid, 2-bromo-, 2-oxo-2-p...	256	C10H9BrO3	1000401-50-0	47
4			Benzilic acid	228	C14H12O3	000076-93-7	47
5			N-[2,2-Dichloro-1-(toluene-4-sul...	371	C16H15Cl2NO3S	136800-77-6	47



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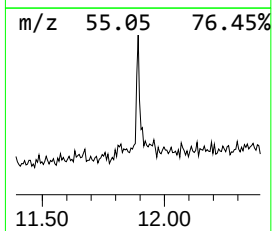
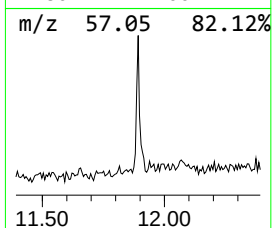
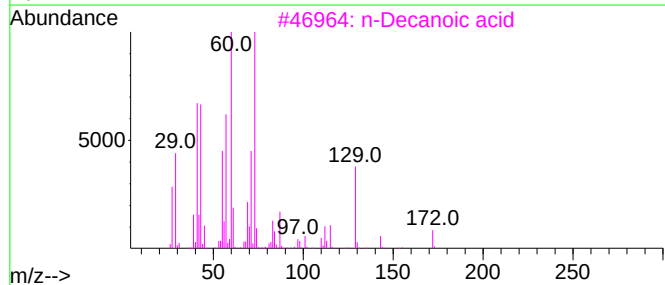
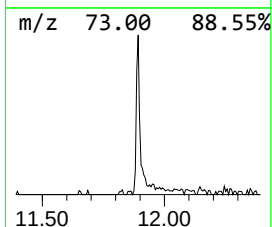
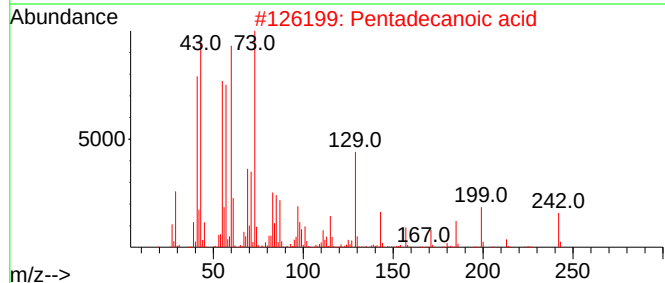
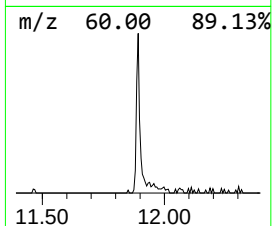
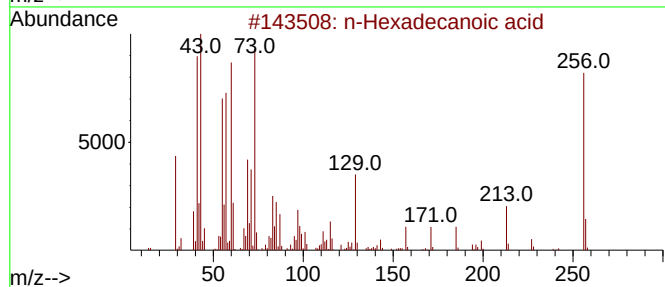
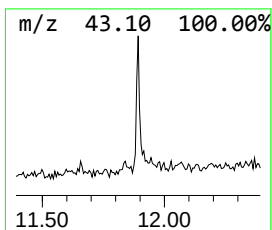
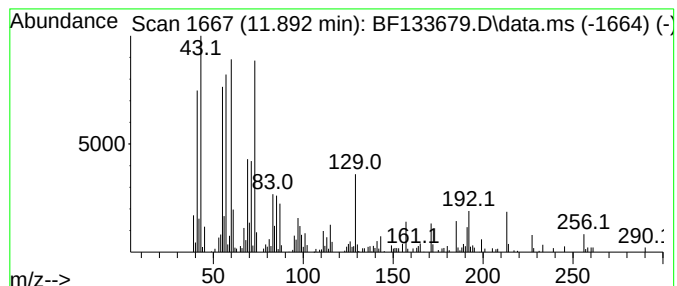
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.892	3.61 ng	106514	Phenanthrene-d10	11.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98		
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	91		
3	n-Decanoic acid	172	C10H20O2	000334-48-5	81		
4	Undecanoic acid	186	C11H22O2	000112-37-8	74		
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	74		



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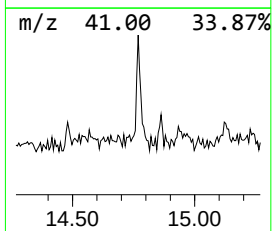
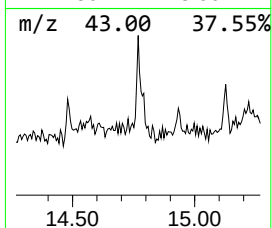
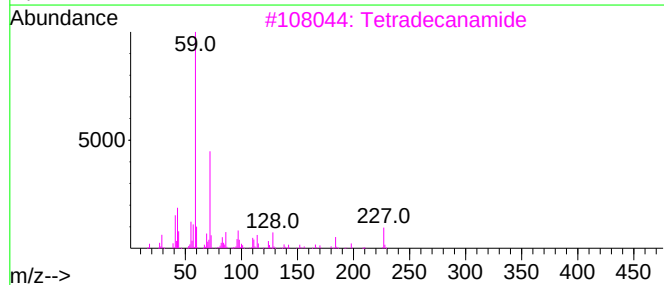
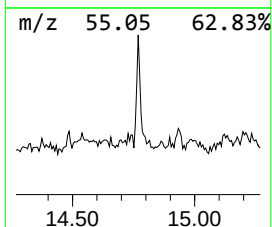
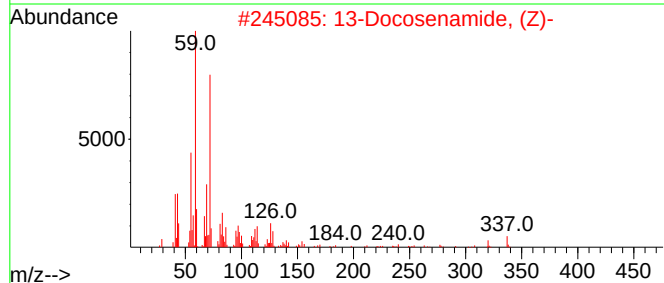
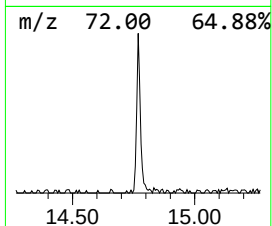
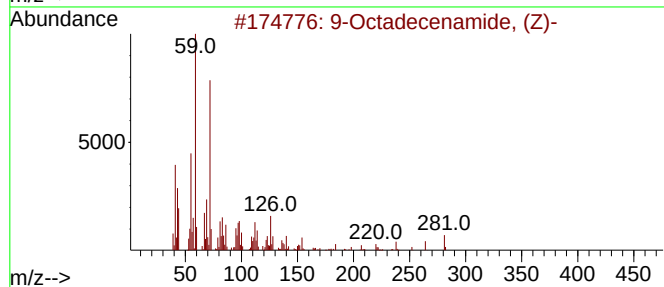
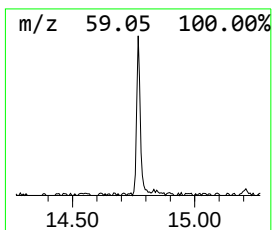
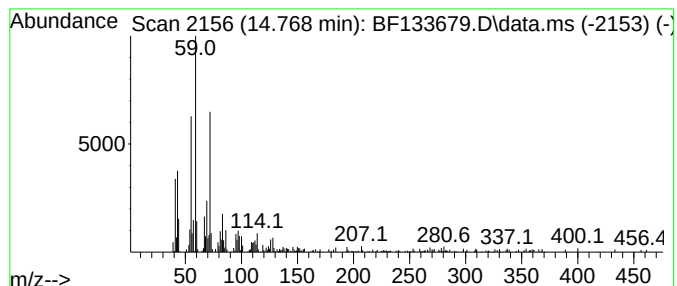
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Peak Number 4 9-Octadecenamide, (Z)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.768	9.37 ng	129245	Perylene-d12	15.474

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	95
2		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	90
3		Tetradecanamide	227	C14H29NO	000638-58-4	78
4		Palmitoleamide	253	C16H31NO	106010-22-4	74
5		Octadecanamide	283	C18H37NO	000124-26-5	72



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.122	22.7	ng	670552	1	6.839	589868	20.0
Benzophenone	10.592	3.5	ng	118722	3	9.875	684686	20.0
n-Hexadecanoic ...	11.892	3.6	ng	106514	4	11.363	590283	20.0
9-Octadecenamid...	14.768	9.4	ng	129245	6	15.474	275887	20.0