

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010722\
 Data File : BF126544.D
 Acq On : 7 Jan 2022 16:21
 Operator : JU/CG
 Sample : N1065-34
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AMI-9

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-BF122921.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF126544.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.998	493	495	503	rVB	40971	50247	2.11%	0.279%
2	5.398	558	563	566	rBV	2336039	2291938	96.18%	12.721%
3	6.410	730	735	738	rBV	2087956	2382867	100.00%	13.226%
4	6.775	794	797	801	rBV	708804	640505	26.88%	3.555%
5	7.339	888	893	896	rBV	1474924	1595760	66.97%	8.857%
6	7.969	996	1000	1010	rBV	134106	163844	6.88%	0.909%
7	8.063	1010	1016	1019	rVV	830515	884375	37.11%	4.909%
8	9.139	1193	1199	1202	rBV	2305929	2382313	99.98%	13.223%
9	9.816	1309	1314	1318	rBV	984149	936004	39.28%	5.195%
10	10.610	1442	1449	1452	rBV	1494453	1495150	62.75%	8.299%
11	11.304	1562	1567	1571	rBV	1058129	921141	38.66%	5.113%
12	11.839	1655	1658	1662	rBV2	33217	35058	1.47%	0.195%
13	12.515	1769	1773	1777	rBV	23985	24363	1.02%	0.135%
14	12.898	1833	1838	1841	rBV	2232425	2137115	89.69%	11.862%
15	13.821	1986	1995	2012	rBV6	61999	210990	8.85%	1.171%
16	13.951	2012	2017	2020	rBV	852444	776012	32.57%	4.307%
17	14.127	2042	2047	2051	rBV	33223	31012	1.30%	0.172%
18	14.439	2095	2100	2104	rBV	35930	34936	1.47%	0.194%
19	14.745	2149	2152	2156	rVB	42729	41573	1.74%	0.231%
20	15.080	2204	2209	2215	rBV	57204	65752	2.76%	0.365%
21	15.409	2259	2265	2269	rBV	507323	636325	26.70%	3.532%
22	15.451	2269	2272	2277	rVB	57176	70955	2.98%	0.394%
23	15.880	2339	2345	2350	rVB	46438	66884	2.81%	0.371%
24	16.368	2423	2428	2436	rBV3	34920	61656	2.59%	0.342%
25	16.950	2520	2527	2534	rVB3	20320	42565	1.79%	0.236%
26	17.633	2635	2643	2648	rBV3	14637	37068	1.56%	0.206%

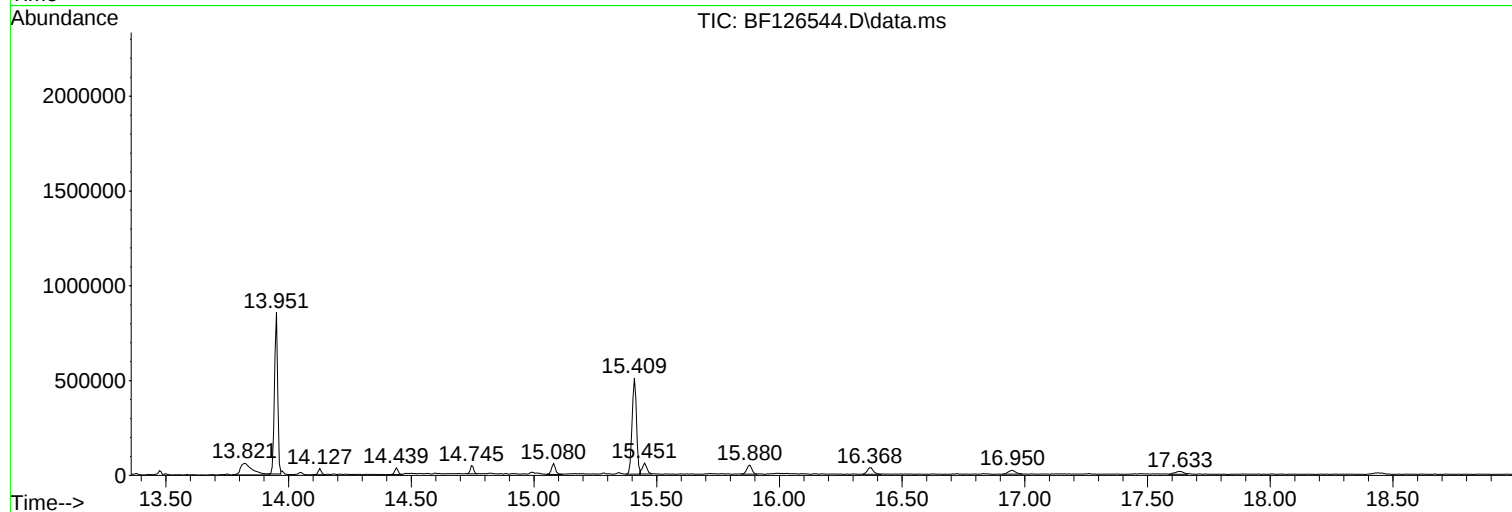
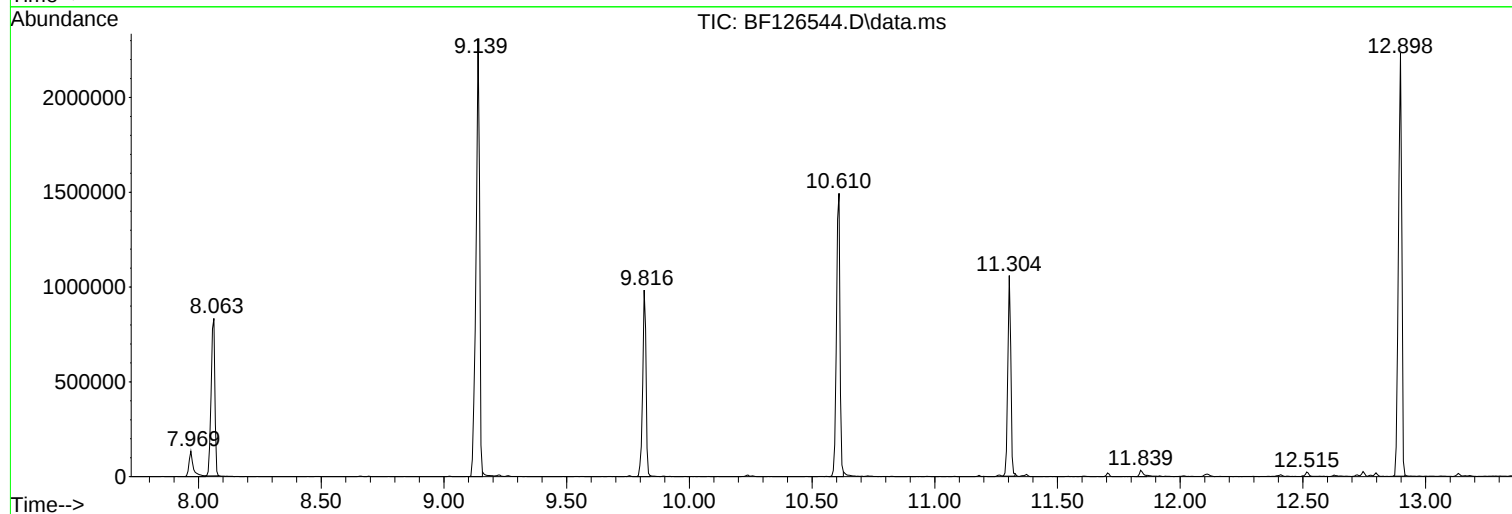
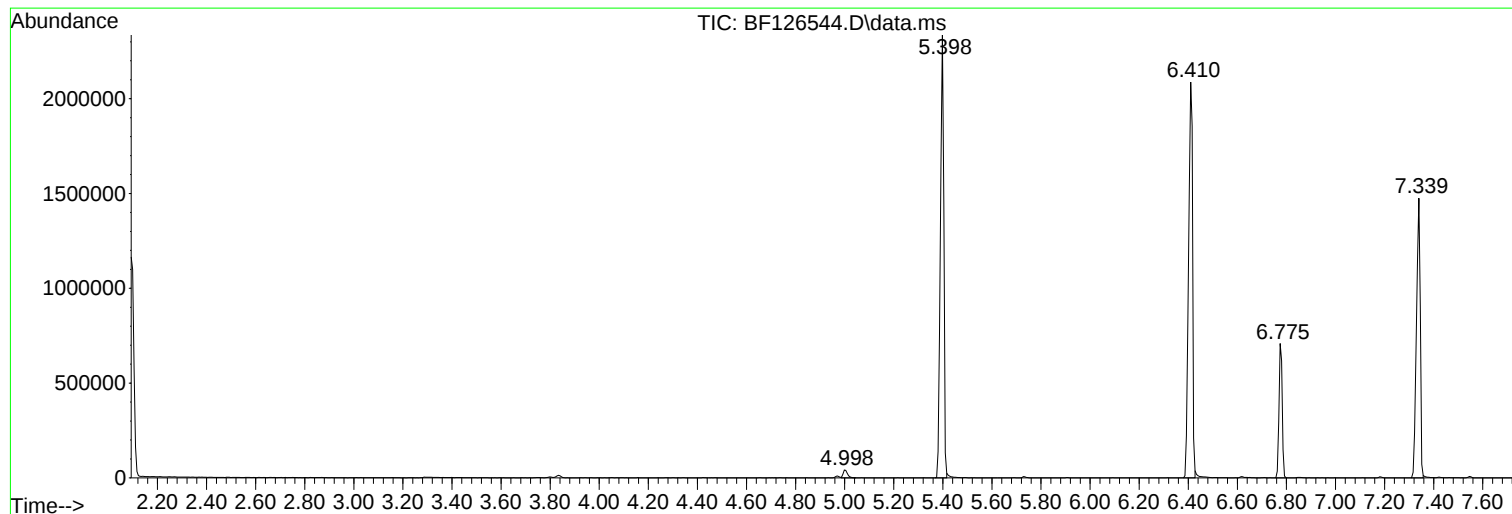
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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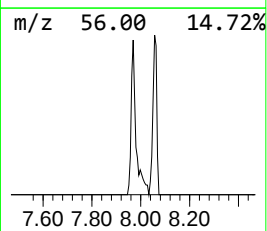
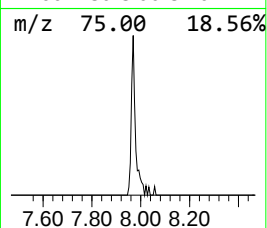
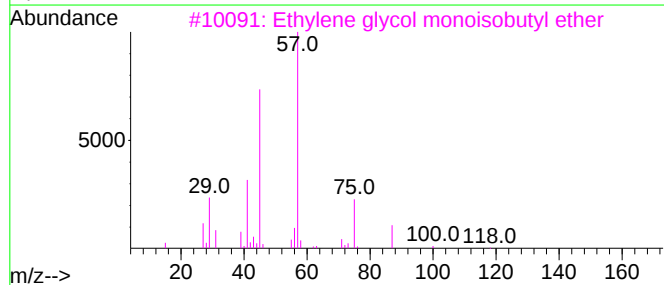
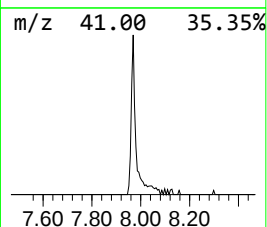
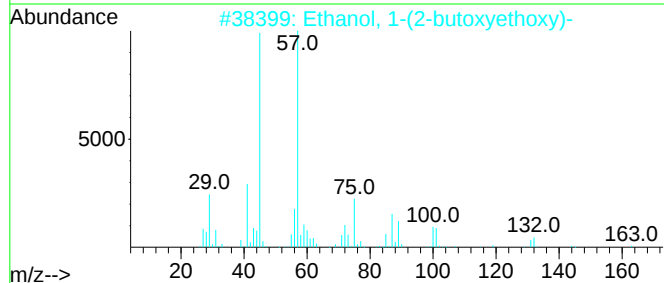
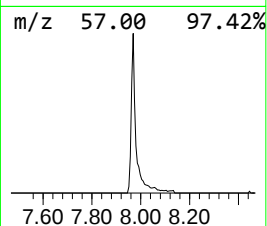
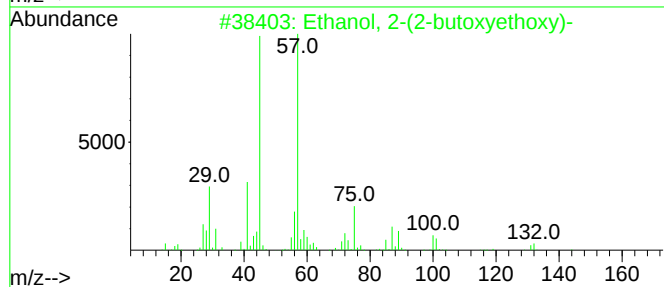
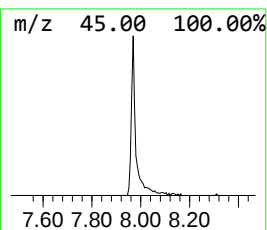
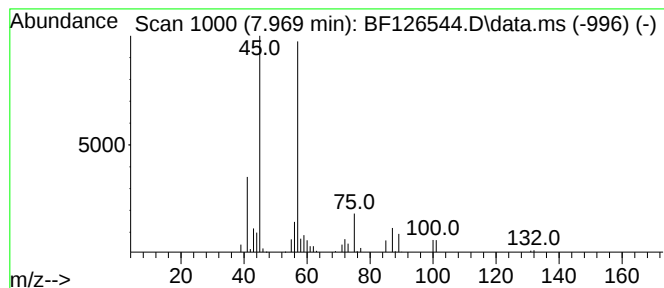
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.969	3.71 ng	163844	Naphthalene-d8	8.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
3			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
4			Di-sec-Butyl ether	130	C8H18O	006863-58-7	53
5			Pentaethylene glycol	238	C10H22O6	004792-15-8	50



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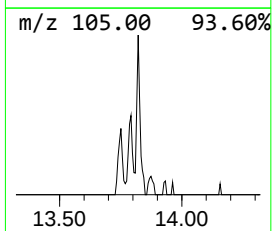
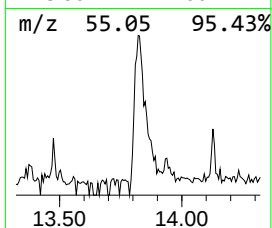
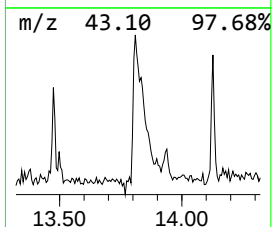
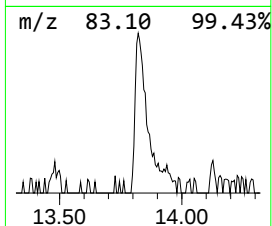
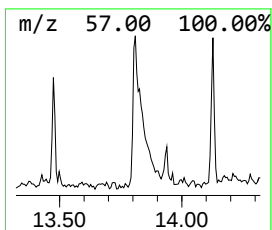
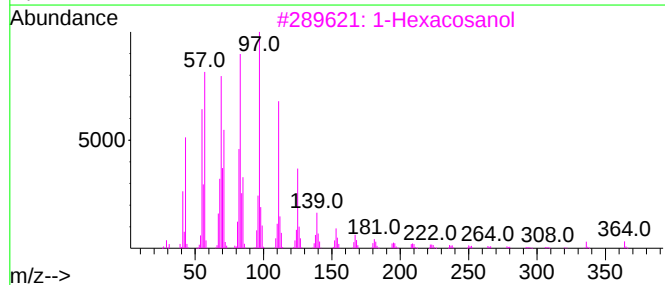
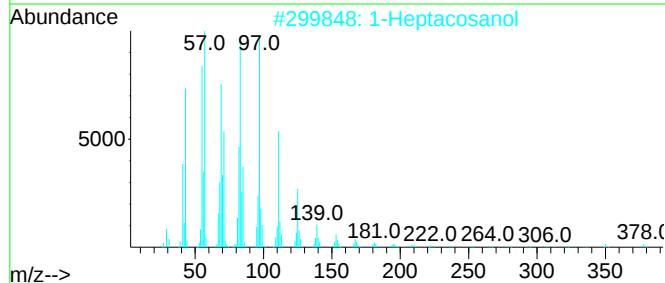
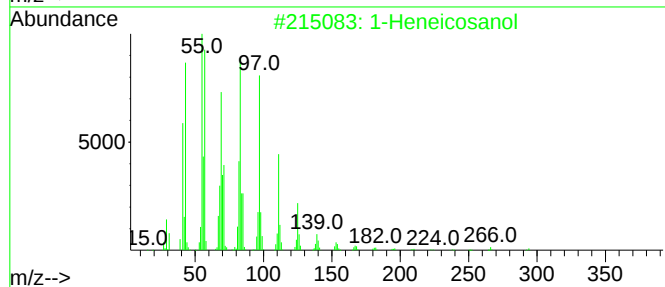
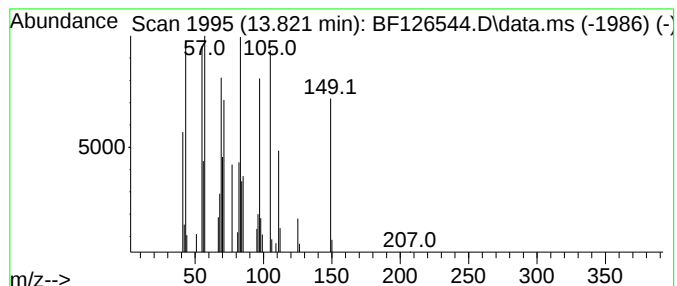
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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TIC Integration Parameters: LSCINT.P

Peak Number 2 1-Heneicosanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.821	5.44 ng	210990	Chrysene-d12	13.951	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	81
2	1-Heptacosanol	396	C27H56O	002004-39-9	81
3	1-Hexacosanol	382	C26H54O	000506-52-5	81
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	81
5	n-Nonadecanol-1	284	C19H40O	001454-84-8	81



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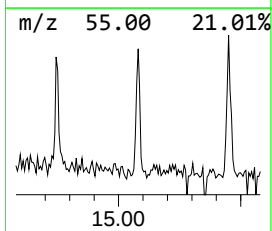
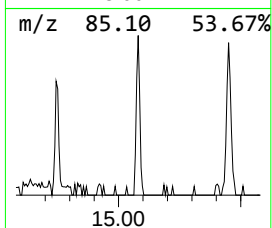
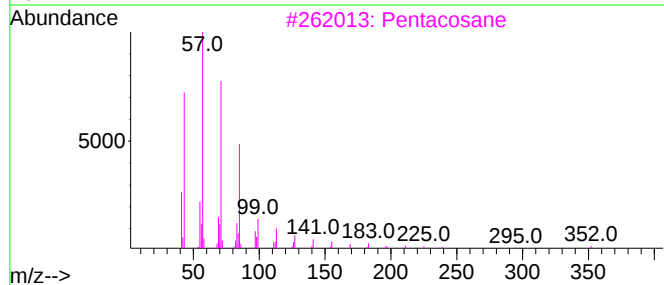
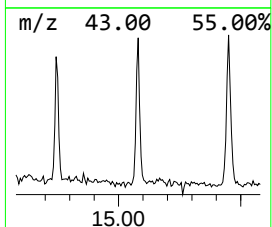
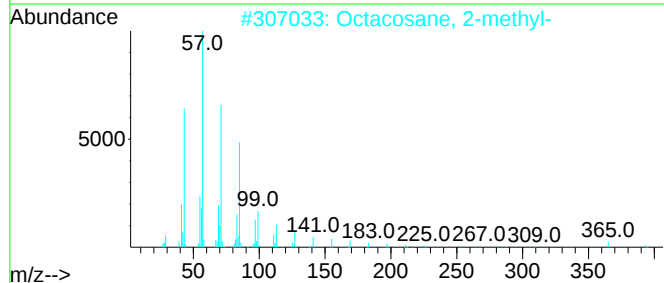
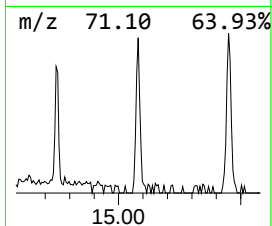
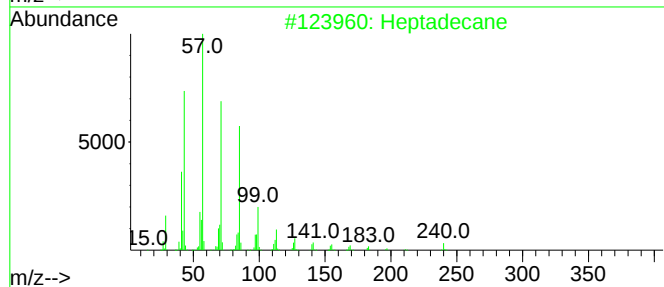
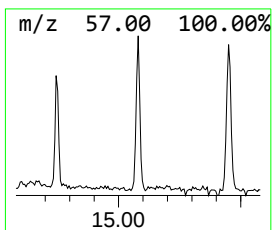
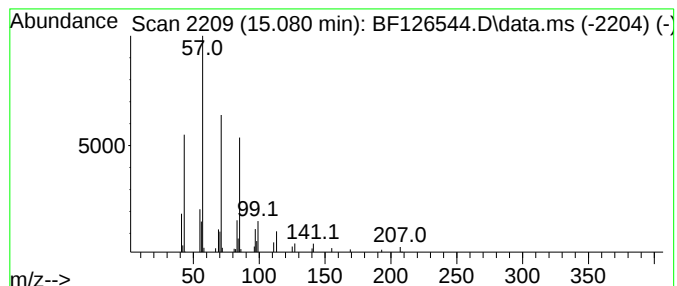
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Peak Number 3 Heptadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.080	2.07 ng	65752	Perylene-d12	15.409

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	91
2			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
3			Pentacosane	352	C25H52	000629-99-2	91
4			2-Methyltetracosane	352	C25H52	001560-78-7	91
5			Hexacosane	366	C26H54	000630-01-3	91



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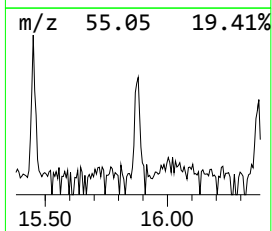
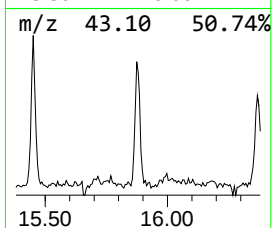
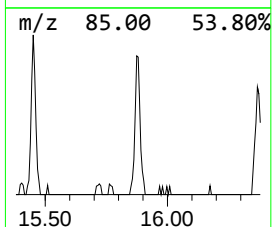
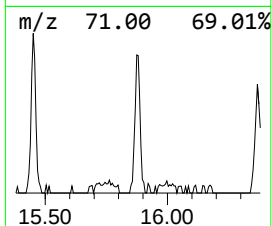
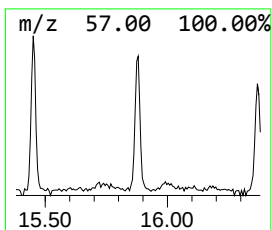
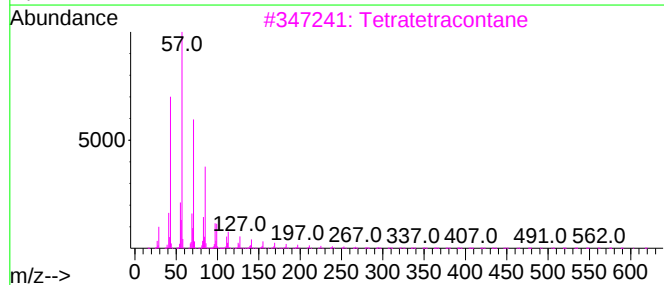
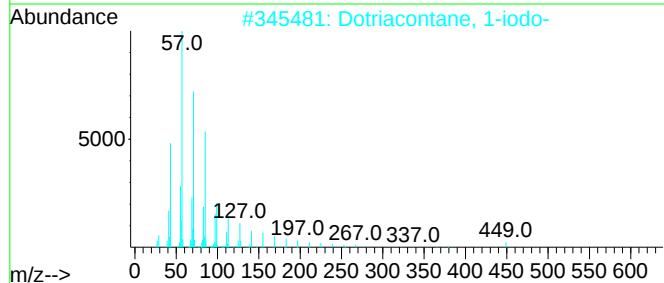
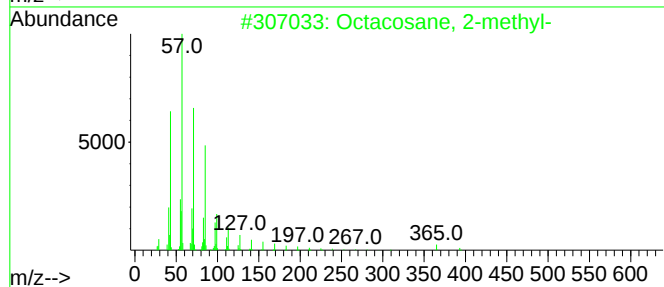
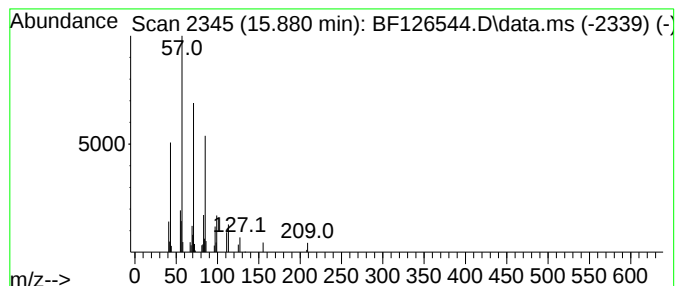
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Peak Number 4 Octacosane, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.880	2.10 ng	66884	Perylene-d12	15.409

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
2			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	90
3			Tetratetracontane	619	C44H90	007098-22-8	90
4			Hentriacontane	437	C31H64	000630-04-6	86
5			Hexacosane	366	C26H54	000630-01-3	86



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(2-b...	7.969	3.7	ng	163844	2	8.063	884375	20.0
1-Heneicosanol	13.821	5.4	ng	210990	5	13.951	776012	20.0
Heptadecane	15.080	2.1	ng	65752	6	15.409	636325	20.0
Octacosane, 2-m...	15.880	2.1	ng	66884	6	15.409	636325	20.0