

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF080219\
 Data File : BF115950.D
 Acq On : 2 Aug 2019 16:55
 Operator : HP/JU
 Sample : PB121932BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB121932BL

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-BF073019.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	2.199	12	19	28	rVB	120888	189069	3.83%	0.491%
2	5.487	573	578	581	rBV	3575348	3641750	73.78%	9.453%
3	6.493	743	749	751	rBV	3213022	3867472	78.35%	10.039%
4	6.628	767	772	775	rBV	3879662	3948095	79.98%	10.248%
5	6.851	806	810	820	rVB	1189282	992551	20.11%	2.576%
6	7.010	832	837	840	rBV	3413027	3226280	65.36%	8.374%
7	7.416	900	906	909	rBV	2359445	2595131	52.57%	6.736%
8	8.134	1023	1028	1046	rBV	1246216	1256343	25.45%	3.261%
9	9.216	1205	1212	1215	rBV	4457846	4660096	94.41%	12.096%
10	9.892	1322	1327	1344	rBV	1501982	1489672	30.18%	3.867%
11	10.686	1457	1462	1494	rBV	2700241	2950322	59.77%	7.658%
12	11.380	1576	1580	1599	rBV	1544751	1504044	30.47%	3.904%
13	12.969	1845	1850	1853	rBV	4649965	4936235	100.00%	12.813%
14	14.021	2025	2029	2043	rVB	1255762	1270747	25.74%	3.298%
15	14.786	2154	2159	2163	rBV	103095	110267	2.23%	0.286%
16	15.127	2211	2217	2222	rBV	140074	178844	3.62%	0.464%
17	15.498	2274	2280	2297	rBV2	776291	1293084	26.20%	3.356%
18	15.939	2348	2355	2361	rBV	139401	223406	4.53%	0.580%
19	16.439	2435	2440	2448	rVB	78657	140017	2.84%	0.363%
20	17.033	2536	2541	2549	rBV3	25430	51807	1.05%	0.134%

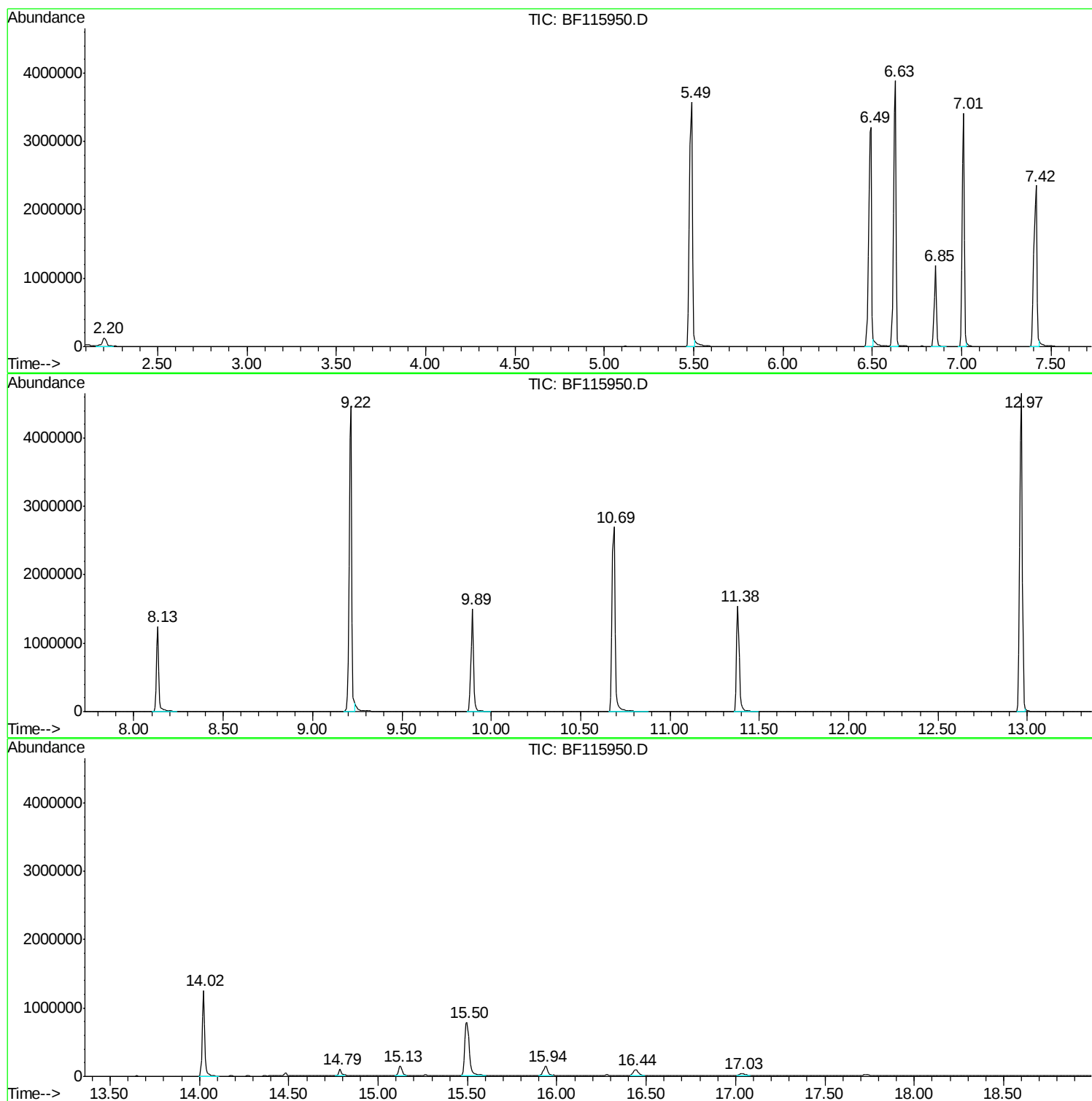
Sum of corrected areas: 38525232

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.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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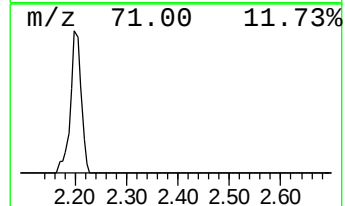
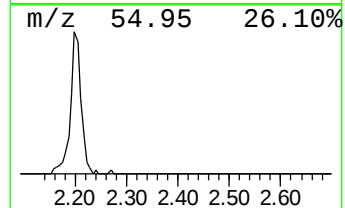
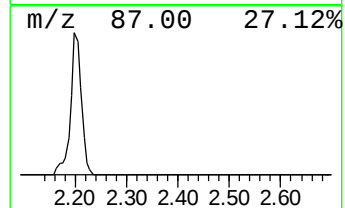
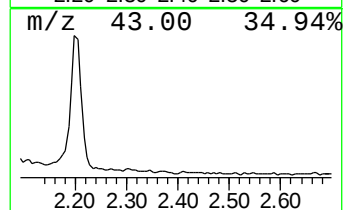
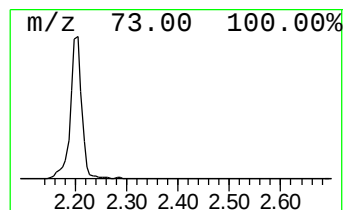
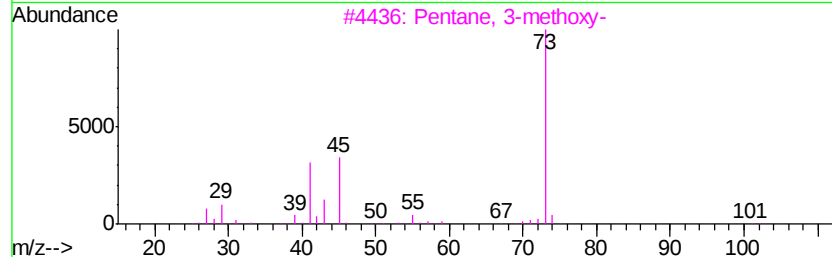
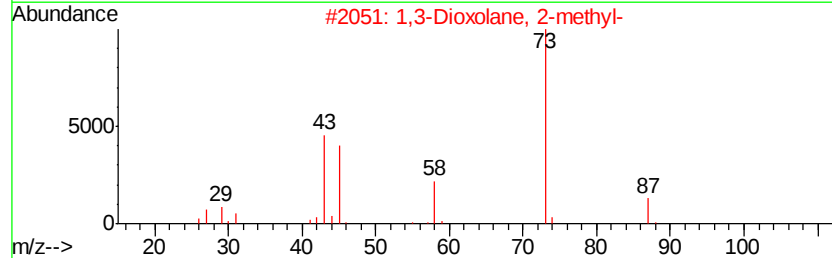
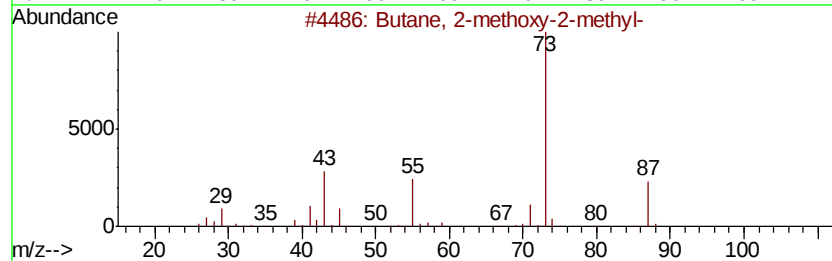
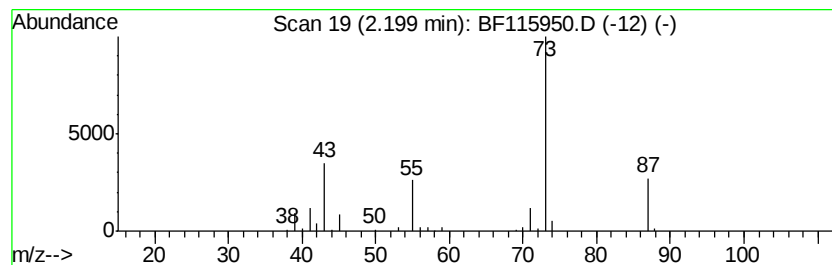
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.20	3.81 ng	189069	1,4-Dichlorobenzene-d4	6.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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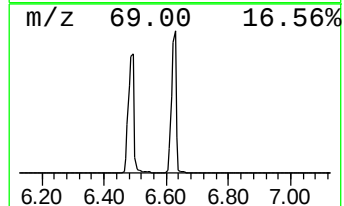
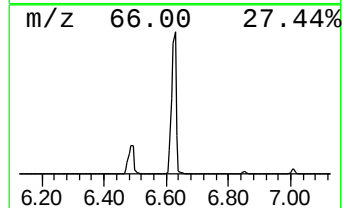
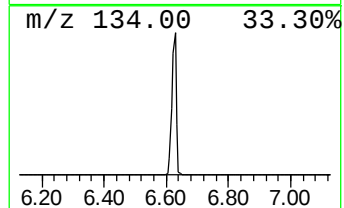
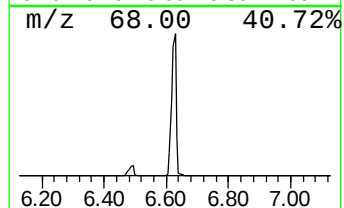
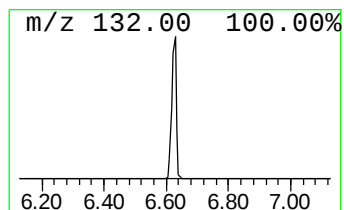
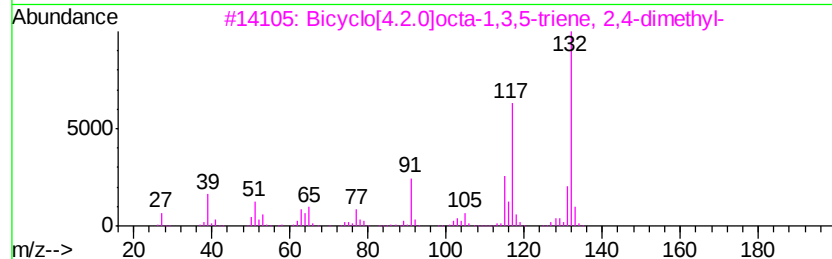
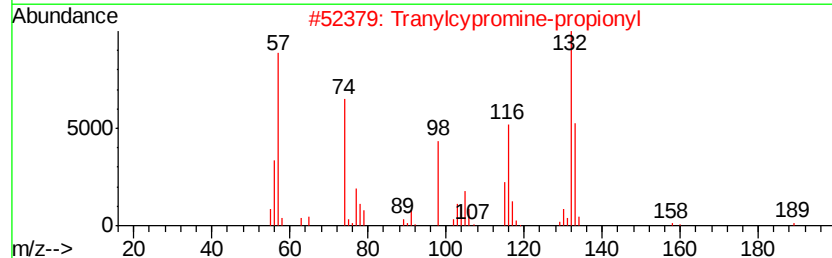
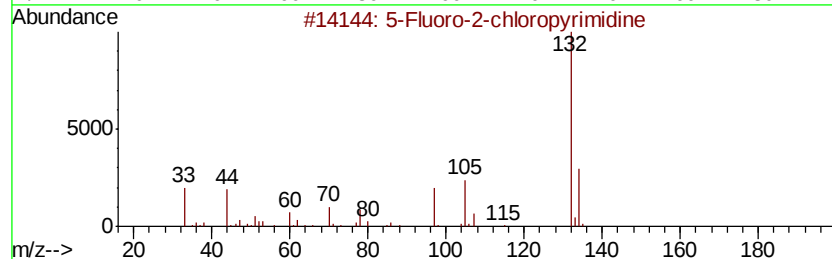
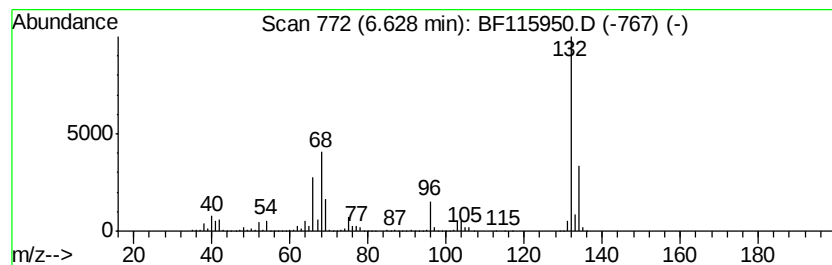
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	79.55 ng	3948100	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Xenon	132	Xe	007440-63-3	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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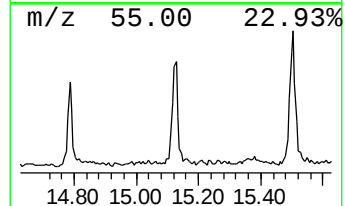
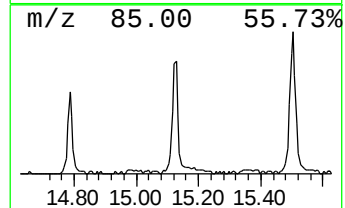
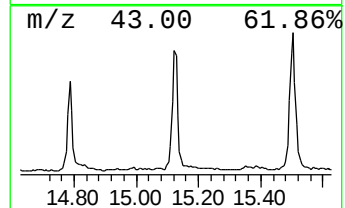
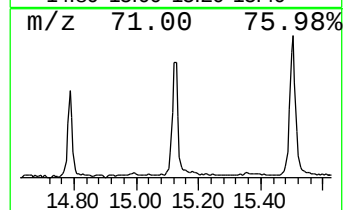
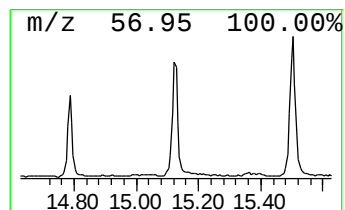
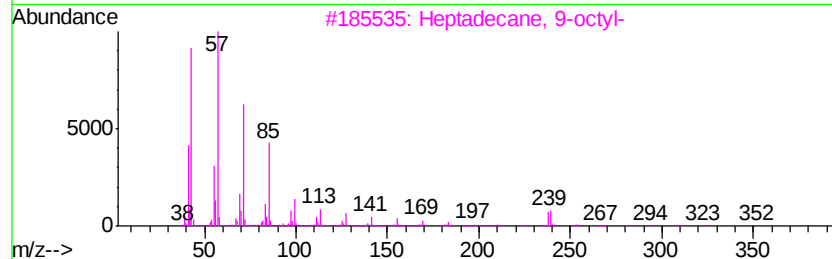
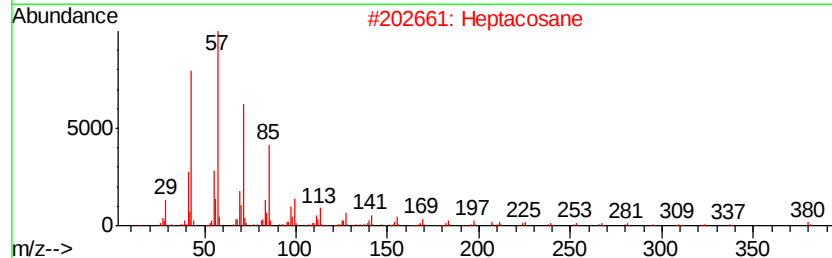
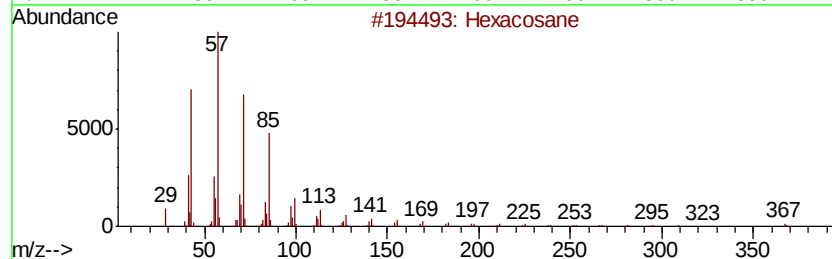
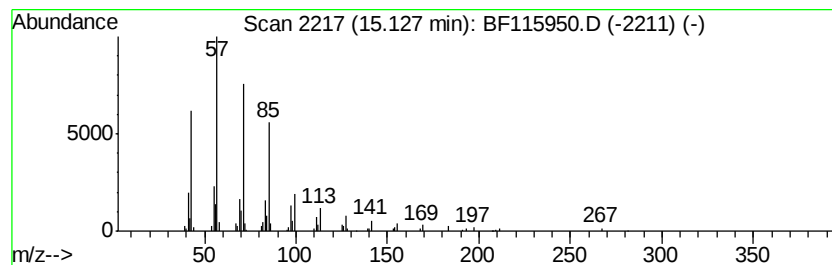
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 Peak Number 4 Hexacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.13	2.77 ng	178844	Perylene-d12	15.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane		366	C26H54	000630-01-3	91
2	Heptacosane		380	C27H56	000593-49-7	91
3	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
4	Heneicosane		296	C21H44	000629-94-7	91
5	2-methyloctacosane		408	C29H60	1000376-72-8	91



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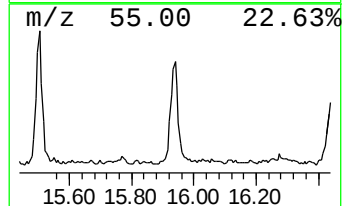
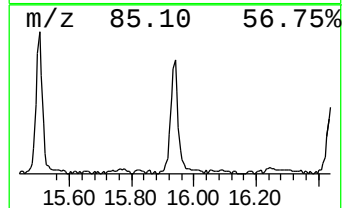
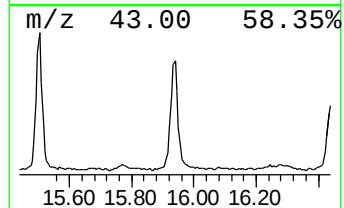
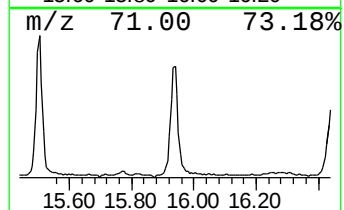
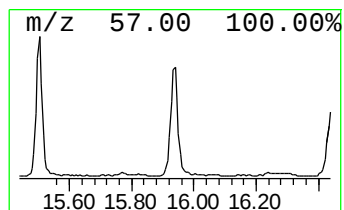
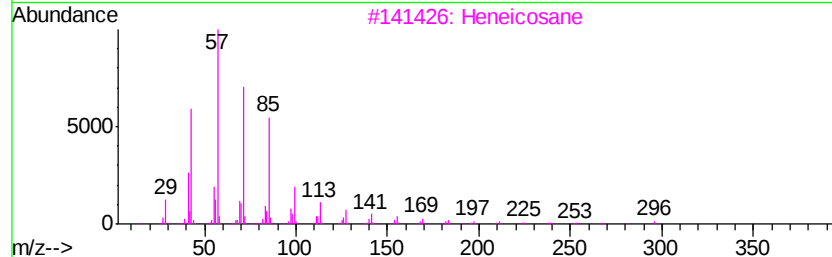
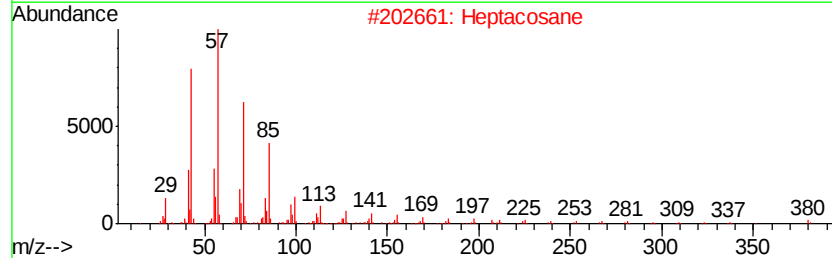
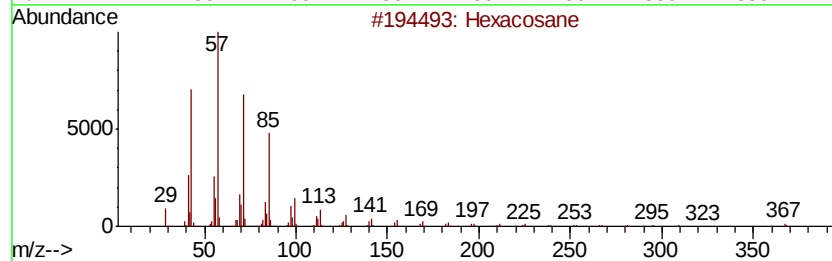
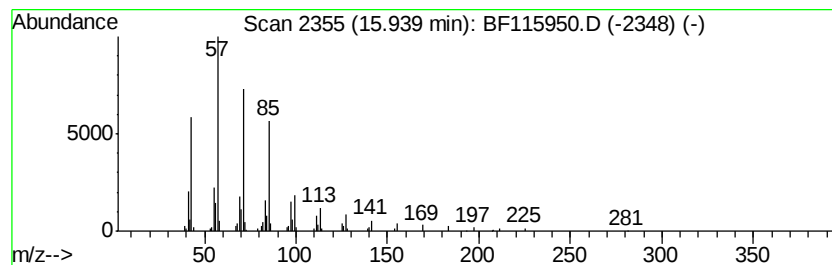
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Heptacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.94	3.46 ng	223406	Perylene-d12	15.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane		366	C26H54	000630-01-3	91
2	Heptacosane		380	C27H56	000593-49-7	91
3	Heneicosane		296	C21H44	000629-94-7	91
4	2-methyloctacosane		408	C29H60	1000376-72-8	91
5	Octacosane		394	C28H58	000630-02-4	91



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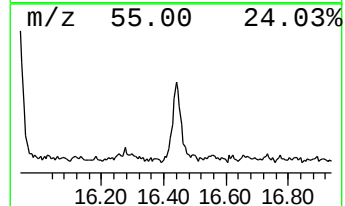
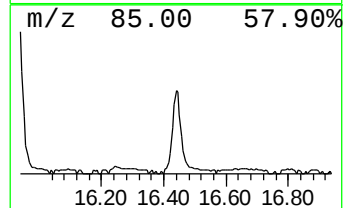
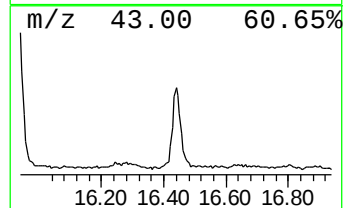
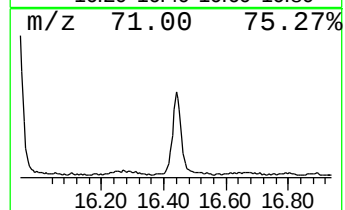
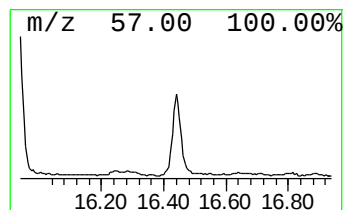
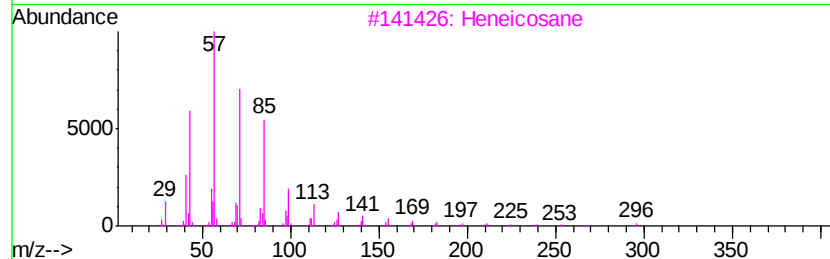
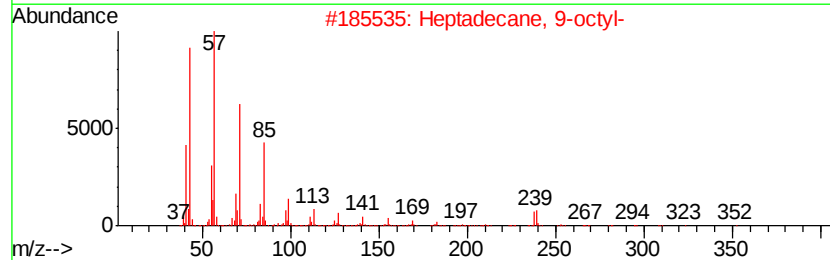
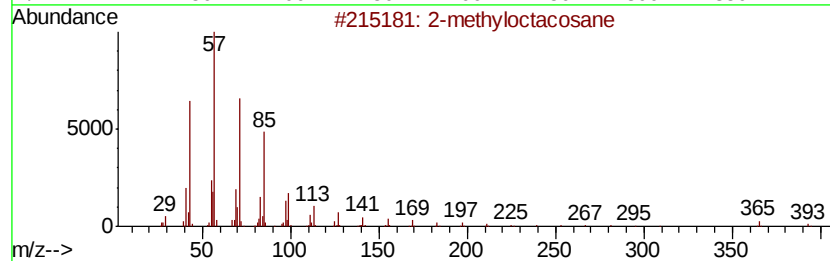
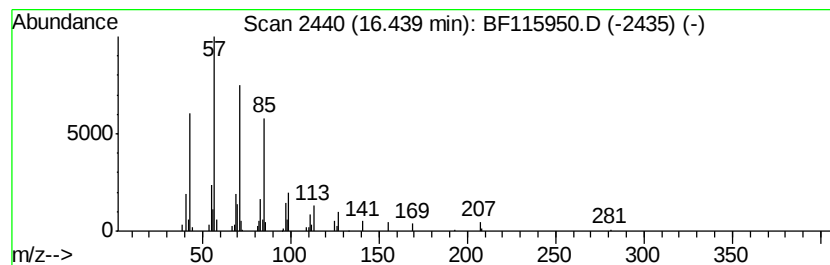
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2-methyloctacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.44	2.17 ng	140017	Perylene-d12	15.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3		Heneicosane	296	C21H44	000629-94-7	90
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
5		Triacontane	422	C30H62	000638-68-6	90



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.20	3.8	ng	189069	1	6.85	992551	20.0
unknown6.63	6.63	79.5	ng	3948100	1	6.85	992551	20.0
Hexacosane	15.13	2.8	ng	178844	6	15.49	1293080	20.0
Heptacosane	15.94	3.5	ng	223406	6	15.49	1293080	20.0
2-methyloctacosane	16.44	2.2	ng	140017	6	15.49	1293080	20.0