

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100225\
 Data File : BF143871.D
 Acq On : 02 Oct 2025 23:01
 Operator : RC/JU
 Sample : Q3262-08 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 279887

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-BF092325.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF143871.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.487	556	560	576	rBV	119475	176005	3.66%	0.873%
2	6.498	726	732	748	rBV	110620	162546	3.38%	0.807%
3	6.881	792	797	813	rBV	542628	708259	14.72%	3.514%
4	7.440	882	892	903	rBV	68382	101579	2.11%	0.504%
5	8.163	1010	1015	1021	rBV	647209	850827	17.68%	4.222%
6	9.045	1162	1165	1168	rBV4	37792	57834	1.20%	0.287%
7	9.163	1182	1185	1186	rBV	58755	63847	1.33%	0.317%
8	9.187	1186	1189	1193	rVB	139464	168718	3.51%	0.837%
9	9.239	1194	1198	1202	rBV	174657	226546	4.71%	1.124%
10	9.322	1208	1212	1216	rBV4	139306	227857	4.73%	1.131%
11	9.639	1261	1266	1270	rBV2	600400	1011657	21.02%	5.020%
12	9.739	1280	1283	1287	rVB3	146004	182485	3.79%	0.905%
13	9.792	1288	1292	1295	rBV3	260639	411214	8.54%	2.040%
14	9.834	1296	1299	1303	rVB	454695	592855	12.32%	2.942%
15	9.922	1308	1314	1318	rVV3	692915	1177834	24.47%	5.844%
16	10.181	1355	1358	1365	rBV4	322672	550024	11.43%	2.729%
17	10.333	1381	1384	1388	rBV2	489708	596843	12.40%	2.961%
18	10.581	1422	1426	1430	rBV	1088221	1477376	30.70%	7.330%
19	10.851	1467	1472	1479	rBV	3122650	4812439	100.00%	23.878%
20	11.322	1548	1552	1558	rVB	2130663	2943923	61.17%	14.607%
21	14.057	2014	2017	2041	rVB	872291	1928870	40.08%	9.571%
22	15.533	2262	2268	2284	rVB	777040	1353343	28.12%	6.715%
23	16.010	2344	2349	2355	rVB	63081	110699	2.30%	0.549%
24	16.527	2432	2437	2448	rVB	73115	158821	3.30%	0.788%
25	17.133	2536	2540	2548	rVB	47647	101490	2.11%	0.504%

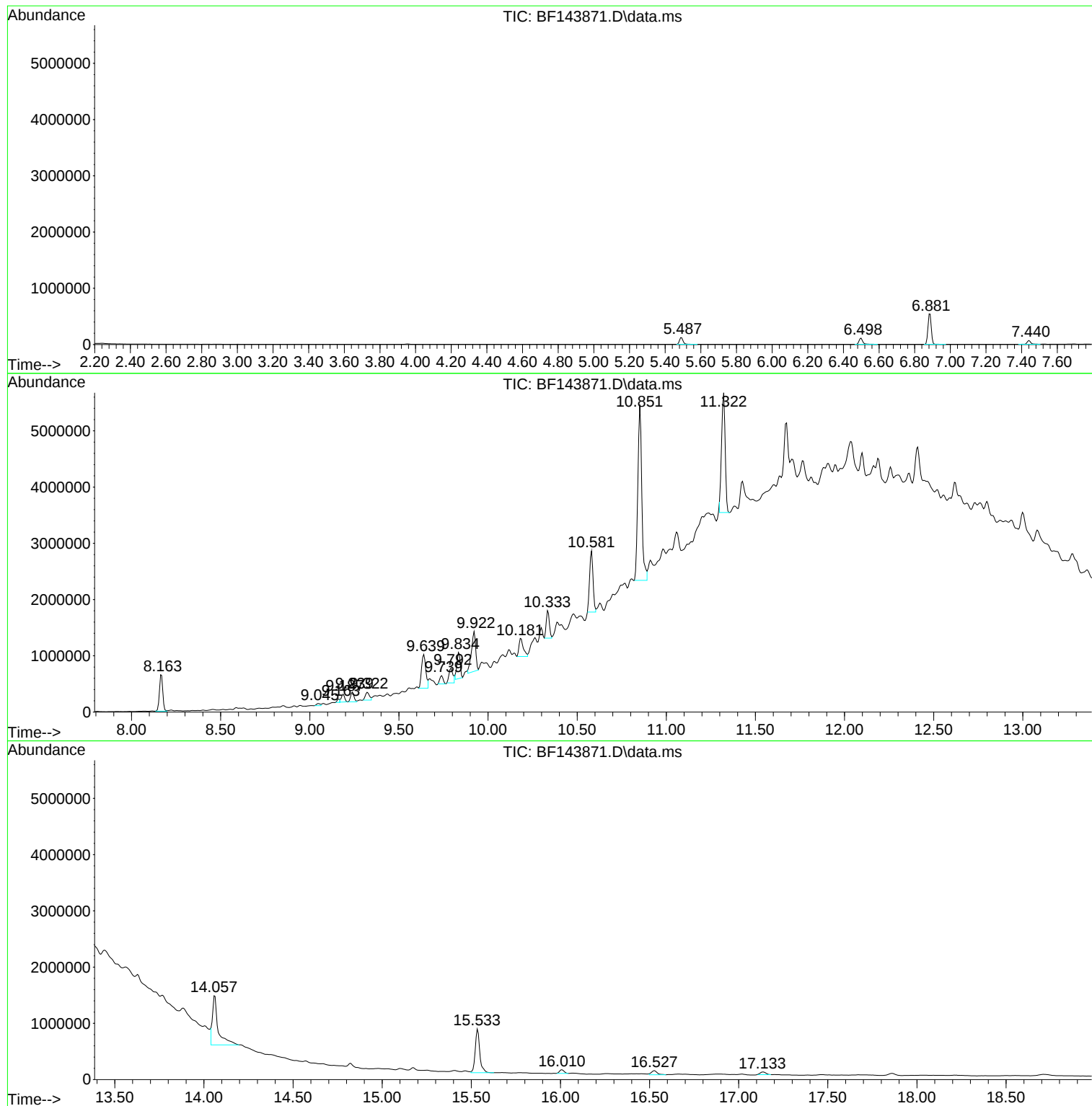
Sum of corrected areas: 20153891

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF092325.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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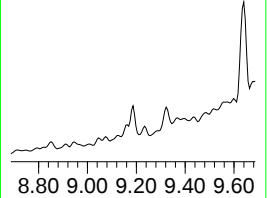
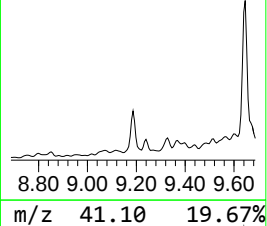
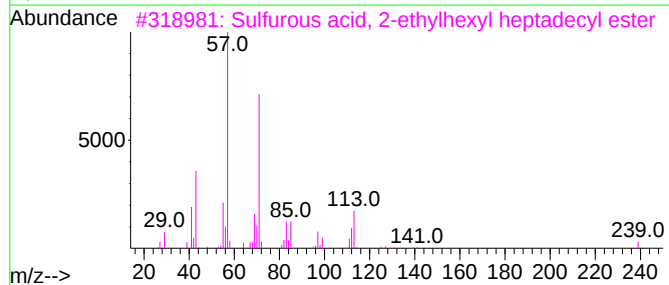
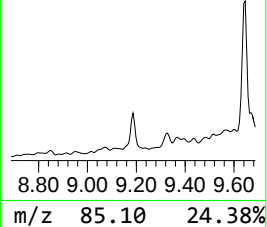
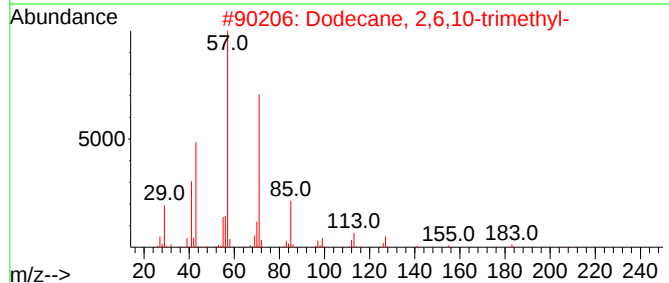
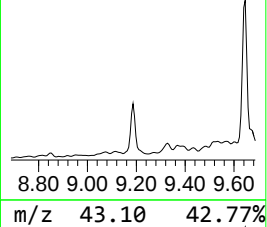
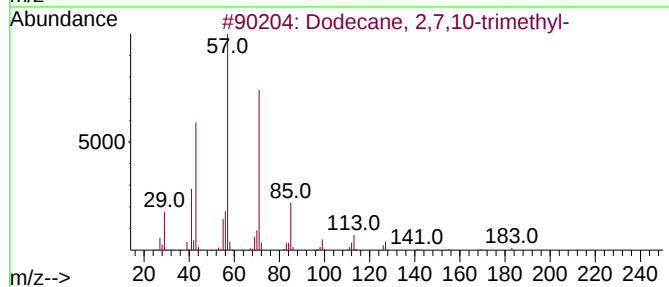
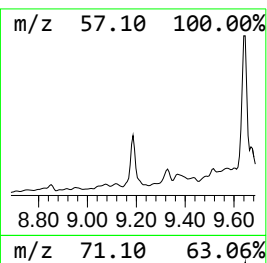
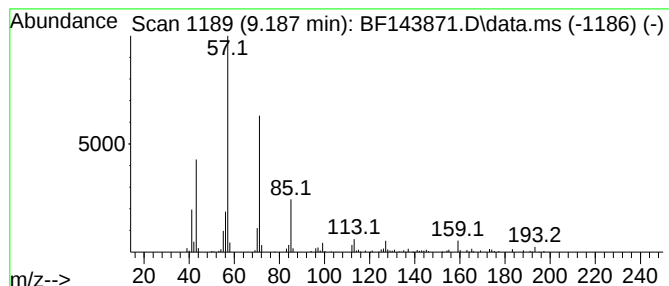
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF092325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Dodecane, 2,7,10-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
9.187	2.86 ng	168718	Acenaphthene-d10		9.922	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Dodecane, 2,7,10-trimethyl-		212	C15H32	074645-98-0	86
2	Dodecane, 2,6,10-trimethyl-		212	C15H32	003891-98-3	86
3	Sulfurous acid, 2-ethylhexyl hep...		432	C25H52O3S	1000309-20-0	59
4	Undecane, 6-ethyl-		184	C13H28	017312-60-6	59
5	Nonane, 3,7-dimethyl-		156	C11H24	017302-32-8	59



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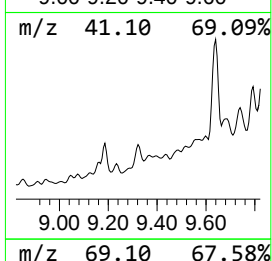
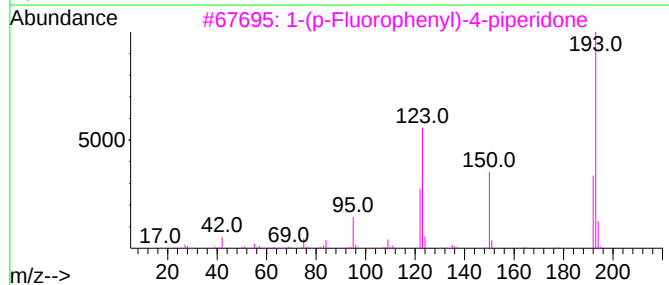
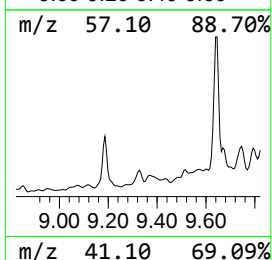
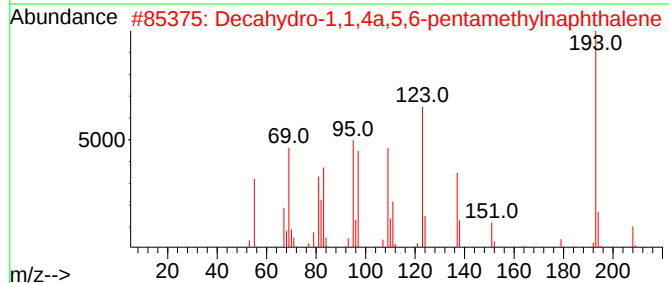
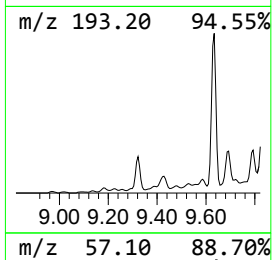
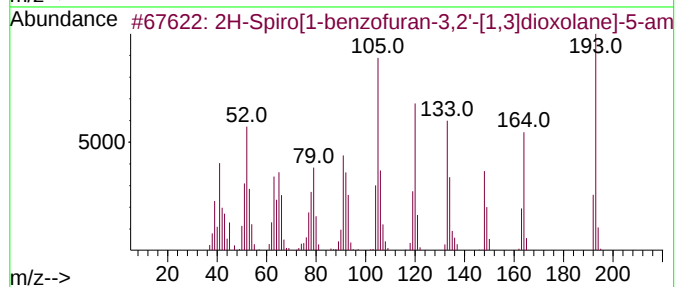
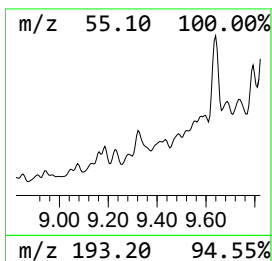
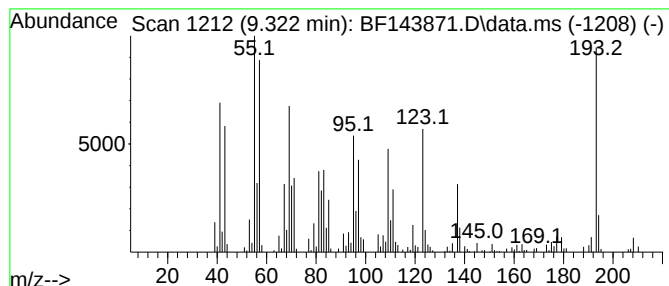
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 2H-Spiro[1-benzofuran-3,2'-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD			R.T.
9.322	3.87 ng	227857	Acenaphthene-d10			9.922
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2H-Spiro[1-benzofuran-3,2'-[1,3]...		193	C10H11NO3	1000387-33-4	90
2	Decahydro-1,1,4a,5,6-pentamethyl...		208	C15H28	080655-44-3	68
3	1-(p-Fluorophenyl)-4-piperidone		193	C11H12FNO	1000238-56-7	38
4	[1,2,4]Triazolo[1,5-a]pyrimidin-...		193	C8H11N5O	1000351-50-1	38
5	2-Anthracenamine		193	C14H11N	000613-13-8	27



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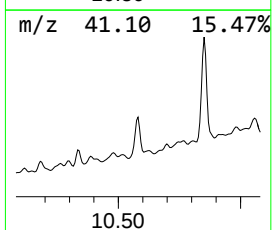
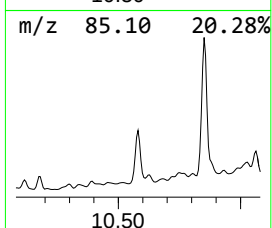
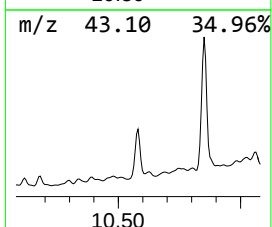
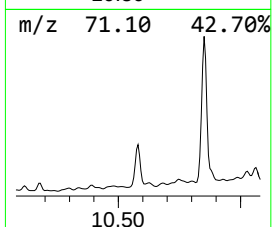
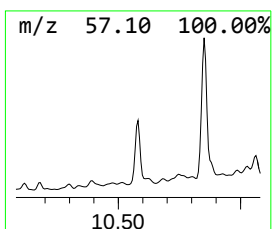
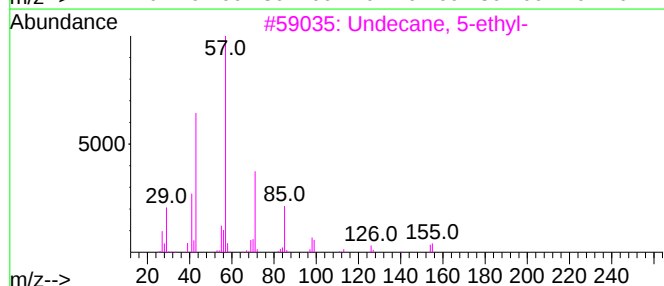
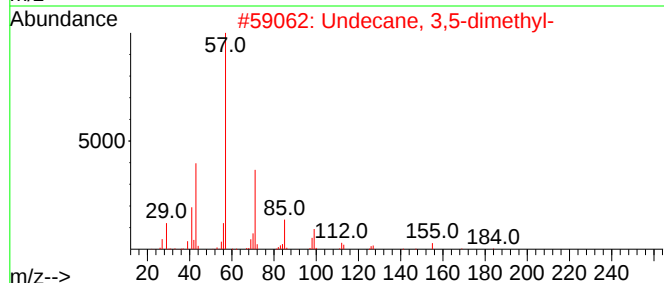
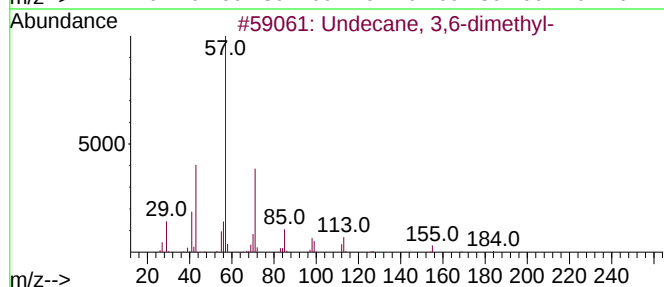
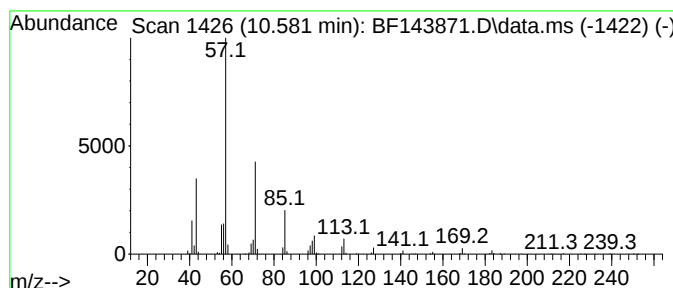
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BNA_F
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Undecane, 3,6-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.581	25.09 ng	1477380	Acenaphthene-d10	9.922	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	93
2	Undecane, 3,5-dimethyl-	184	C13H28	017312-81-1	87
3	Undecane, 5-ethyl-	184	C13H28	017453-94-0	78
4	Decane, 2,6,8-trimethyl-	184	C13H28	062108-26-3	74
5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	70



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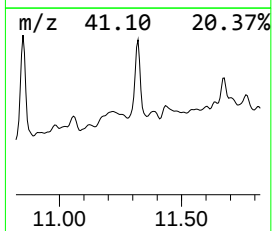
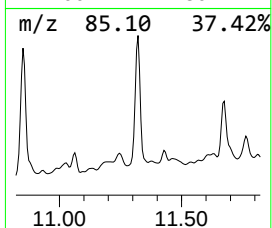
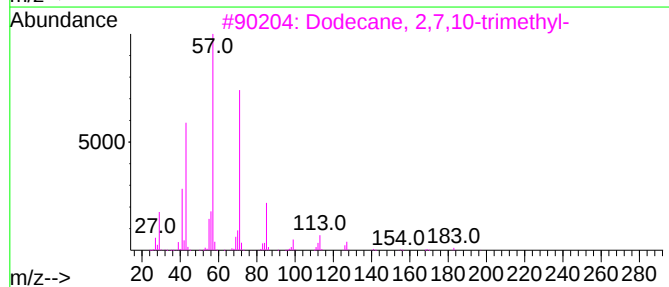
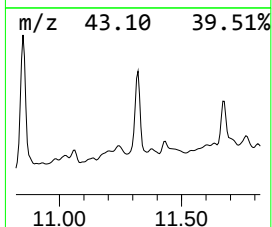
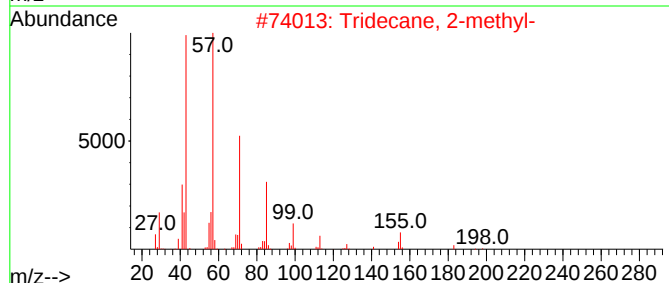
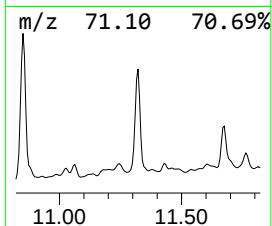
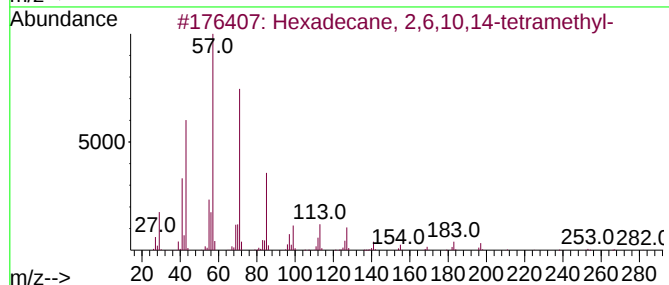
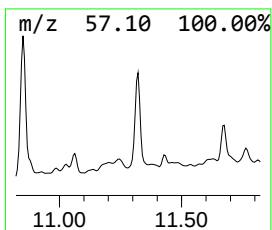
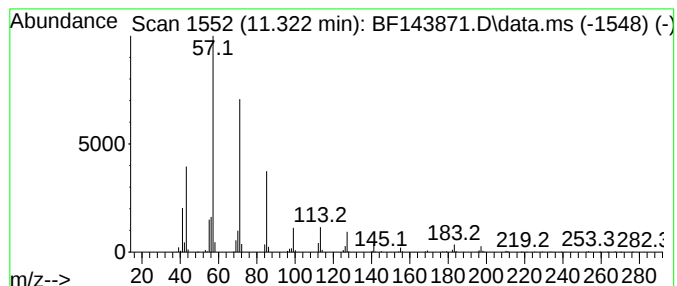
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Hexadecane, 2,6,10,14-tetra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.322	20.00 ng	2943920	Phenanthrene-d10	11.422	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	90
2	Tridecane, 2-methyl-	198	C14H30	001560-96-9	87
3	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	80
4	Hexadecane	226	C16H34	000544-76-3	78
5	Decane, 5-ethyl-5-methyl-	184	C13H28	017312-74-2	72



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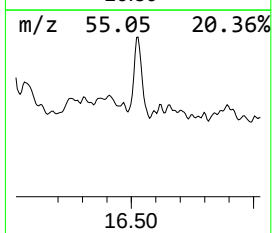
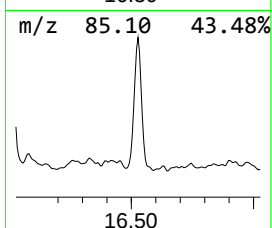
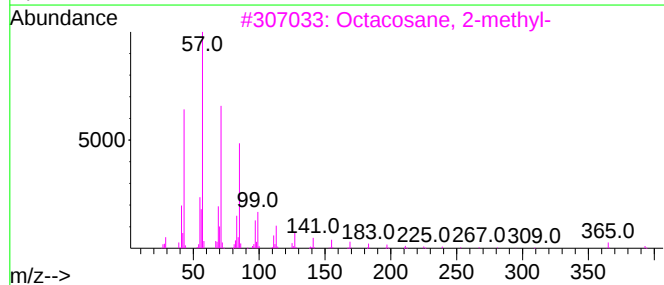
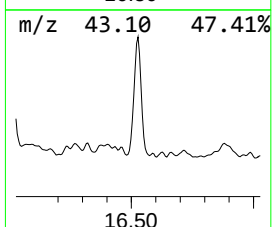
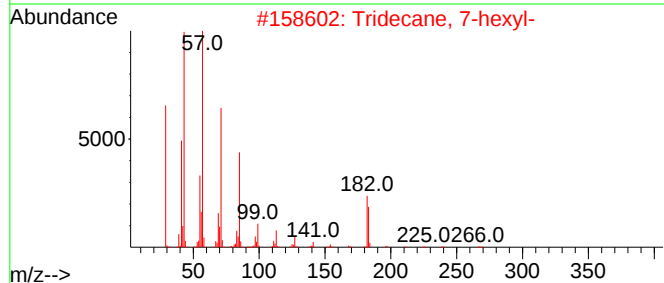
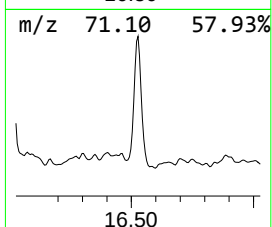
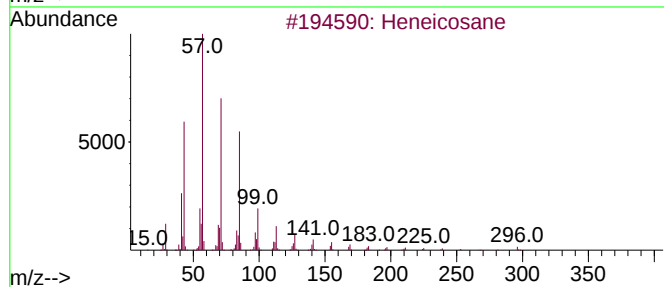
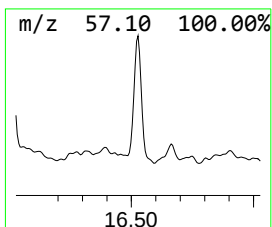
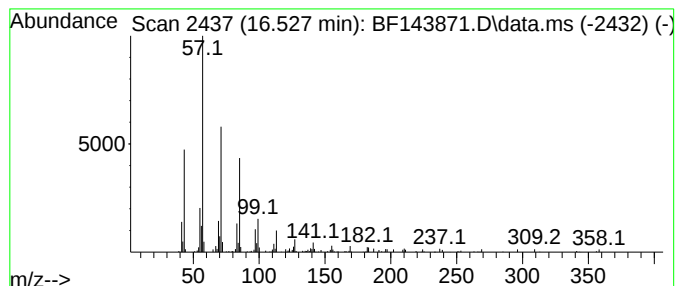
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Heneicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.527	2.35 ng	158821	Perylene-d12	15.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	94
2			Tridecane, 7-hexyl-	268	C19H40	007225-66-3	90
3			Octacosane, 2-methyl-	408	C29H60	001560-98-1	90
4			Pentacosane	352	C25H52	000629-99-2	87
5			Octacosane	394	C28H58	000630-02-4	87



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Dodecane, 2,7,1...	9.187	2.9	ng	168718	3	9.922	1177830	20.0
2H-Spiro[1-benz...	9.322	3.9	ng	227857	3	9.922	1177830	20.0
Undecane, 3,6-d...	10.581	25.1	ng	1477380	3	9.922	1177830	20.0
Hexadecane, 2,6...	11.322	20.0	ng	2943920	4	11.422	2943920	20.0
Heneicosane	16.527	2.4	ng	158821	6	15.533	1353340	20.0