

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040820\
 Data File : BF119853.D
 Acq On : 9 Apr 2020 4:56
 Operator : MA/SJ
 Sample : L2151-03
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CBH-139

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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-BF040820.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	4.963	482	489	497	rBV	100107	133827	1.39%	0.250%
2	5.375	554	559	562	rBV	3387921	3079175	32.09%	5.753%
3	6.393	727	732	739	rBV	2055540	1910568	19.91%	3.569%
4	6.763	791	795	799	rBV	1422012	1199252	12.50%	2.241%
5	6.922	817	822	826	rBV	5297309	5424211	56.53%	10.134%
6	7.334	885	892	895	rBV	4250635	4920197	51.28%	9.192%
7	8.051	1009	1014	1017	rBV	1780890	1631872	17.01%	3.049%
8	9.134	1192	1198	1201	rBV	8017014	8872282	92.46%	16.576%
9	9.522	1259	1264	1267	rBV	1028421	854180	8.90%	1.596%
10	9.804	1305	1312	1316	rBV	2031594	2030180	21.16%	3.793%
11	10.604	1440	1448	1451	rBV	5702948	6220442	64.83%	11.622%
12	11.292	1561	1565	1569	rBV	2559519	2147090	22.38%	4.011%
13	11.827	1652	1656	1666	rBV	422028	463753	4.83%	0.866%
14	12.610	1786	1789	1798	rBV	170695	202886	2.11%	0.379%
15	12.892	1830	1837	1840	rBV	8185932	9595411	100.00%	17.927%
16	13.363	1914	1917	1921	rVB	282710	249957	2.60%	0.467%
17	13.816	1991	1994	2006	rVB	191644	417362	4.35%	0.780%
18	13.933	2010	2014	2018	rBV	2087088	1810305	18.87%	3.382%
19	14.410	2092	2095	2099	rBV	111406	106131	1.11%	0.198%
20	14.521	2111	2114	2118	rVB	201308	167549	1.75%	0.313%
21	14.710	2143	2146	2150	rVB	126434	124811	1.30%	0.233%
22	15.039	2198	2202	2208	rBV	128727	148046	1.54%	0.277%
23	15.374	2249	2259	2262	rBV	1185825	1537509	16.02%	2.873%
24	15.410	2262	2265	2275	rVB	119828	157221	1.64%	0.294%
25	15.833	2332	2337	2342	rVB	82666	120763	1.26%	0.226%

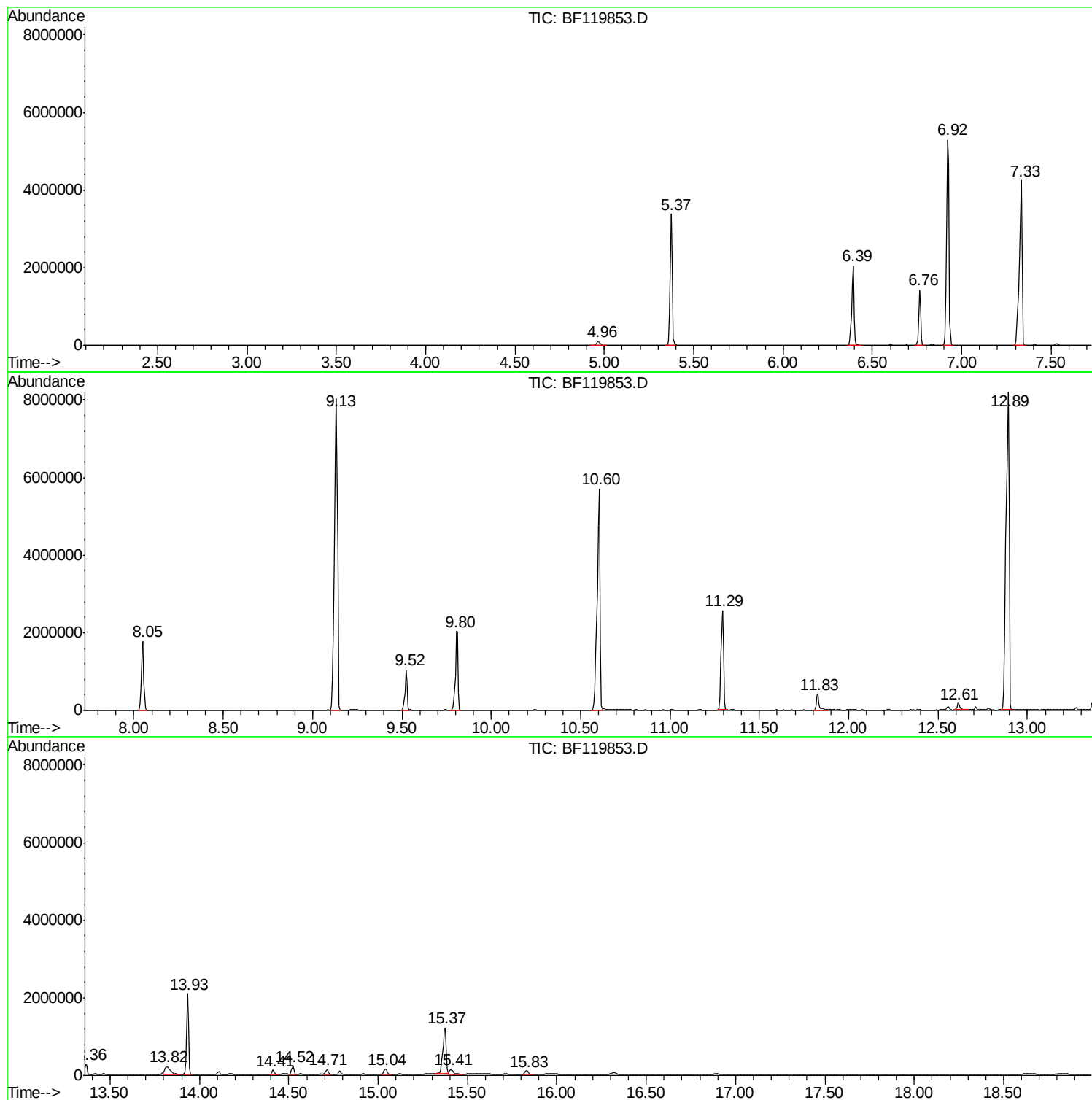
Sum of corrected areas: 53524980

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



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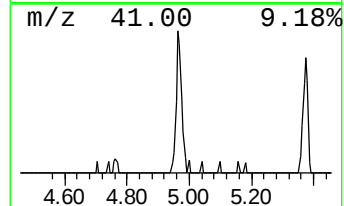
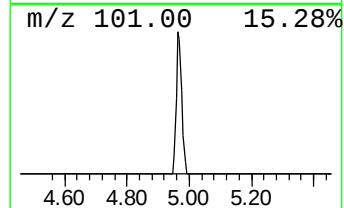
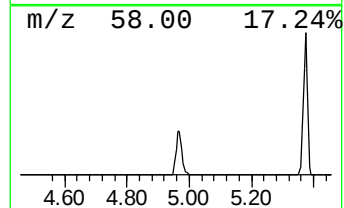
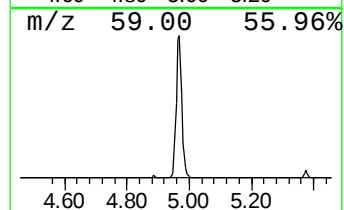
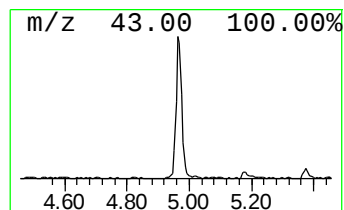
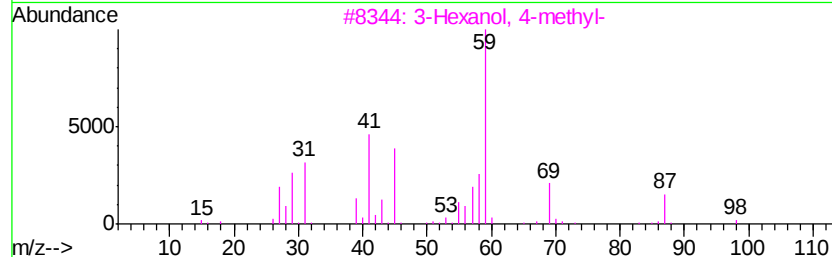
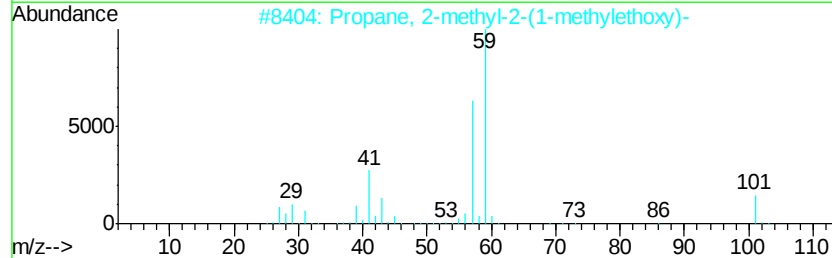
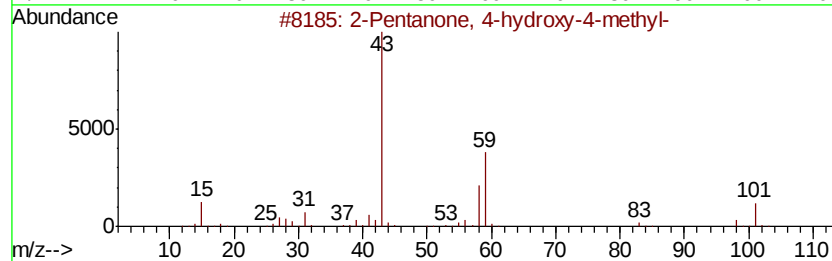
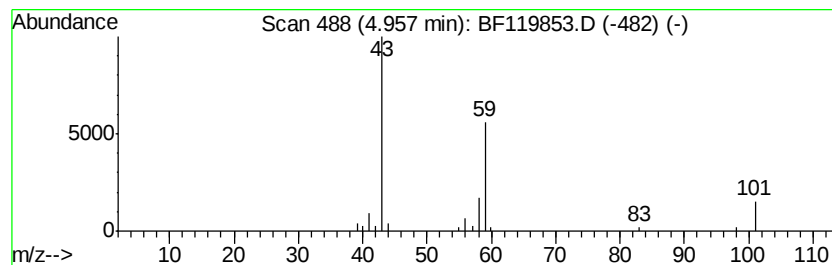
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	2.23 ng	133827	1,4-Dichlorobenzene-d4	6.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	36
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
5			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	23



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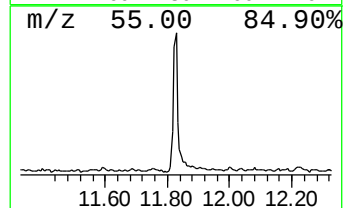
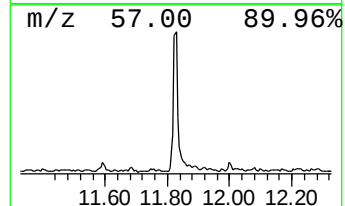
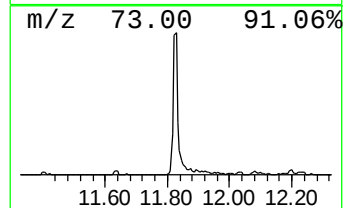
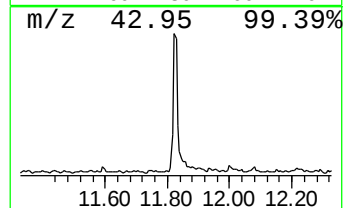
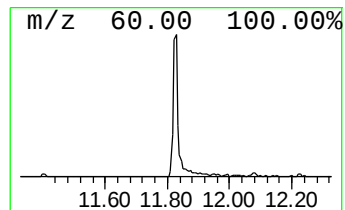
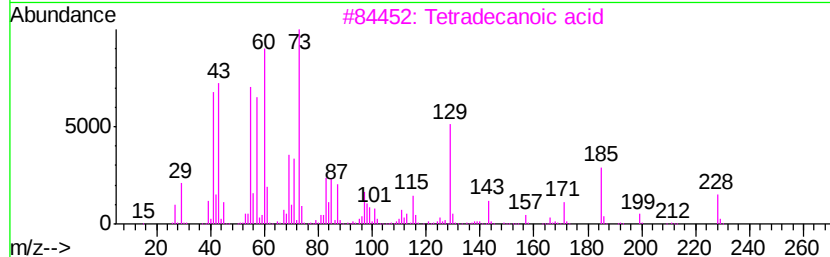
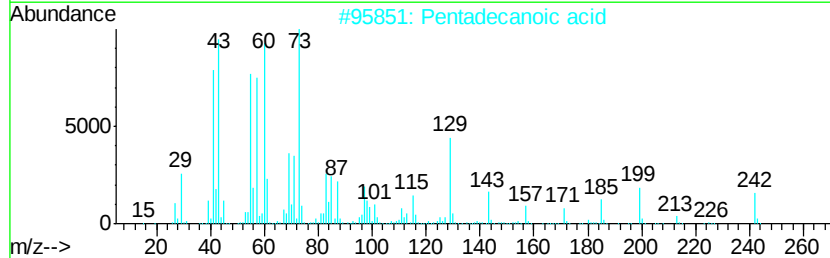
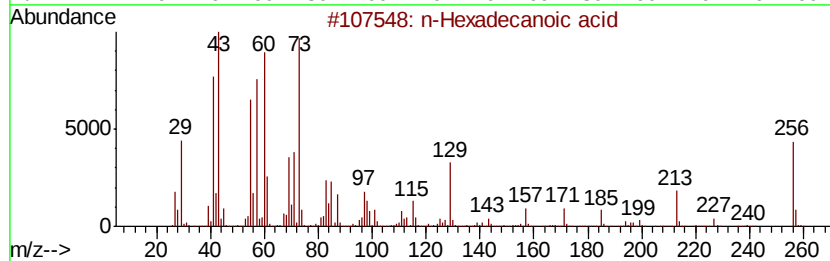
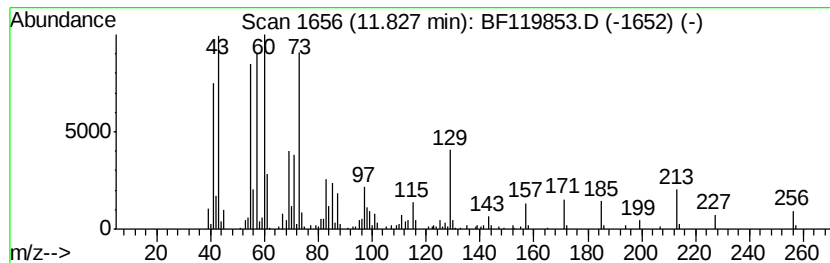
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.83	4.32 ng	463753	Phenanthrene-d10	11.29	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
4	Tridecanoic acid	214	C13H26O2	000638-53-9	64
5	n-Decanoic acid	172	C10H20O2	000334-48-5	53



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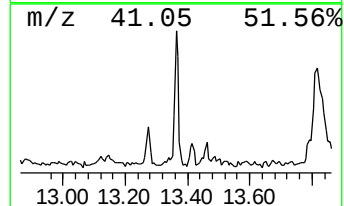
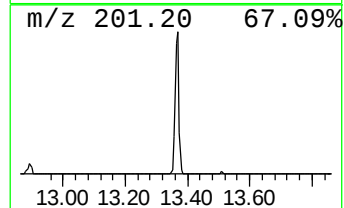
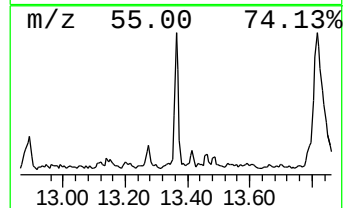
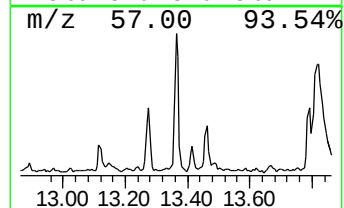
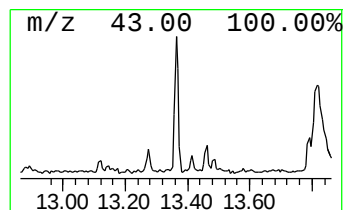
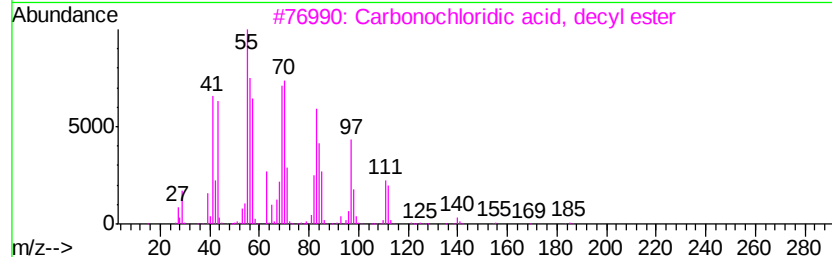
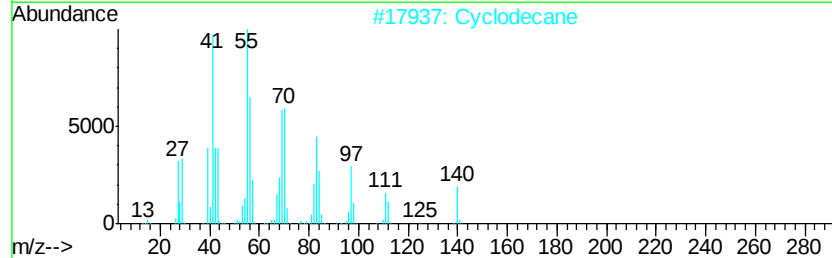
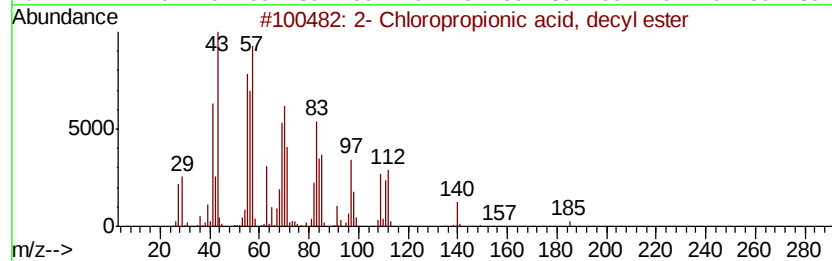
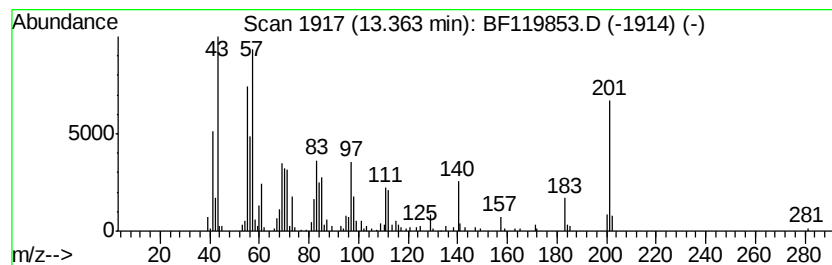
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Peak Number 4 unknown13.36 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.36	2.76 ng	249957	Chrysene-d12	13.93

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-	Chloropropionic acid, decyl e...	248	C13H25ClO2	086711-78-6	46
2		Cyclodecane	140	C10H20	000293-96-9	46
3		Carbonochloridic acid, decyl ester	220	C11H21ClO2	055488-51-2	46
4		Decane, 1-fluoro-	160	C10H21F	000334-56-5	43
5		5-Bromovaleric acid, decyl ester	320	C15H29BrO2	175083-30-4	38



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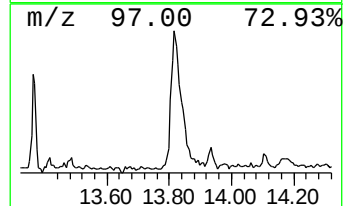
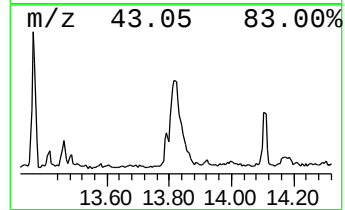
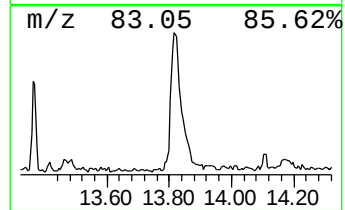
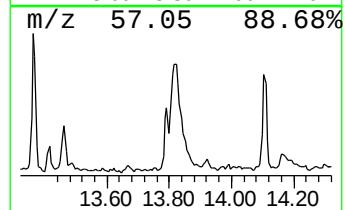
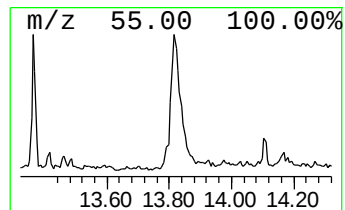
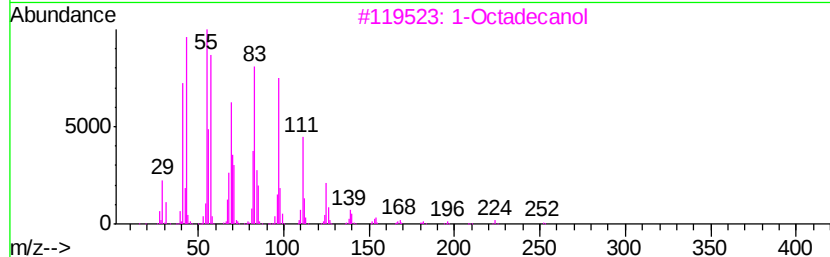
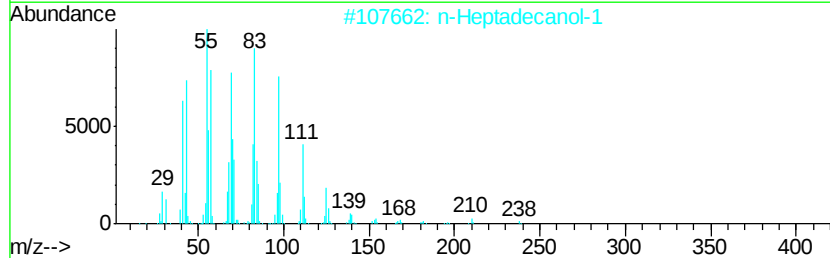
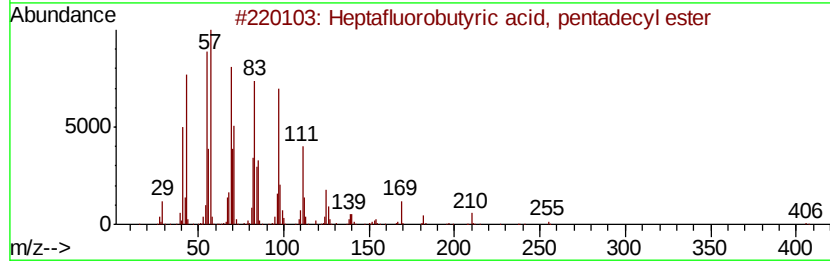
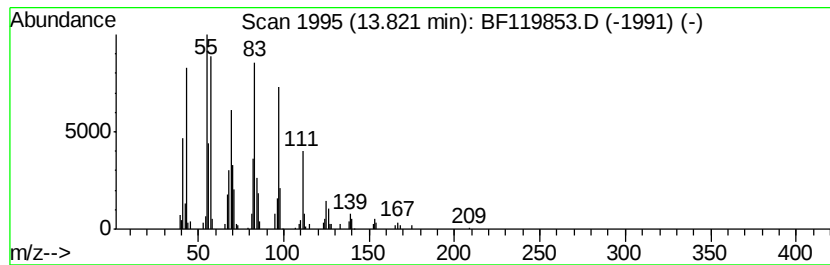
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Peak Number 5 Heptafluorobutyric acid, pe... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	4.61 ng	417362	Chrysene-d12	13.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
2		n-Heptadecanol-1	256	C17H36O	001454-85-9	91
3		1-Octadecanol	270	C18H38O	000112-92-5	91
4		2-Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	91
5		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.96	2.2	ng	133827	1	6.76	1199250	20.0
n-Hexadecanoic acid	11.83	4.3	ng	463753	4	11.29	2147090	20.0
unknown13.36	13.36	2.8	ng	249957	5	13.93	1810310	20.0
Heptafluorobutyri...	13.82	4.6	ng	417362	5	13.93	1810310	20.0