

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM031919\
 Data File : BM019253.D
 Acq On : 20 Mar 2019 04:41
 Operator : JU/SJ
 Sample : K1948-02
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PST-025-B126-031319

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_M\METHODS\8270-BM031119.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	5.234	376	382	399	rBV	3363813	5036899	38.31%	7.437%
2	6.822	644	652	671	rBV	3651225	6177876	46.98%	9.122%
3	7.169	704	711	726	rBV	3654738	5864870	44.60%	8.660%
4	7.634	783	790	801	rBV	569201	899846	6.84%	1.329%
5	7.951	837	844	854	rBV	2494102	3936029	29.93%	5.812%
6	8.804	982	989	1012	rBV	1959813	3379041	25.70%	4.989%
7	10.422	1257	1264	1282	rBV	697869	1266795	9.63%	1.870%
8	12.904	1679	1686	1701	rBV	5660808	8310722	63.20%	12.271%
9	13.757	1825	1831	1842	rBV	423428	703241	5.35%	1.038%
10	14.292	1915	1922	1936	rVB2	1255945	1923833	14.63%	2.841%
11	15.792	2171	2177	2192	rBV	5317930	7538600	57.33%	11.131%
12	17.045	2384	2390	2403	rBV2	1589048	2411413	18.34%	3.561%
13	17.939	2537	2542	2553	rBV	322188	538658	4.10%	0.795%
14	19.227	2757	2761	2775	rBV	113610	242148	1.84%	0.358%
15	19.686	2833	2839	2853	rBV	10467911	13149304	100.00%	19.416%
16	20.950	3049	3054	3067	rBV	565448	898908	6.84%	1.327%
17	21.244	3099	3104	3117	rBV	1892328	2724741	20.72%	4.023%
18	22.268	3275	3278	3283	rVB	143039	175140	1.33%	0.259%
19	22.768	3360	3363	3368	rVB	180744	215876	1.64%	0.319%
20	23.327	3455	3458	3463	rVB	144286	194771	1.48%	0.288%
21	23.474	3477	3483	3494	rVB2	1112099	2137050	16.25%	3.155%

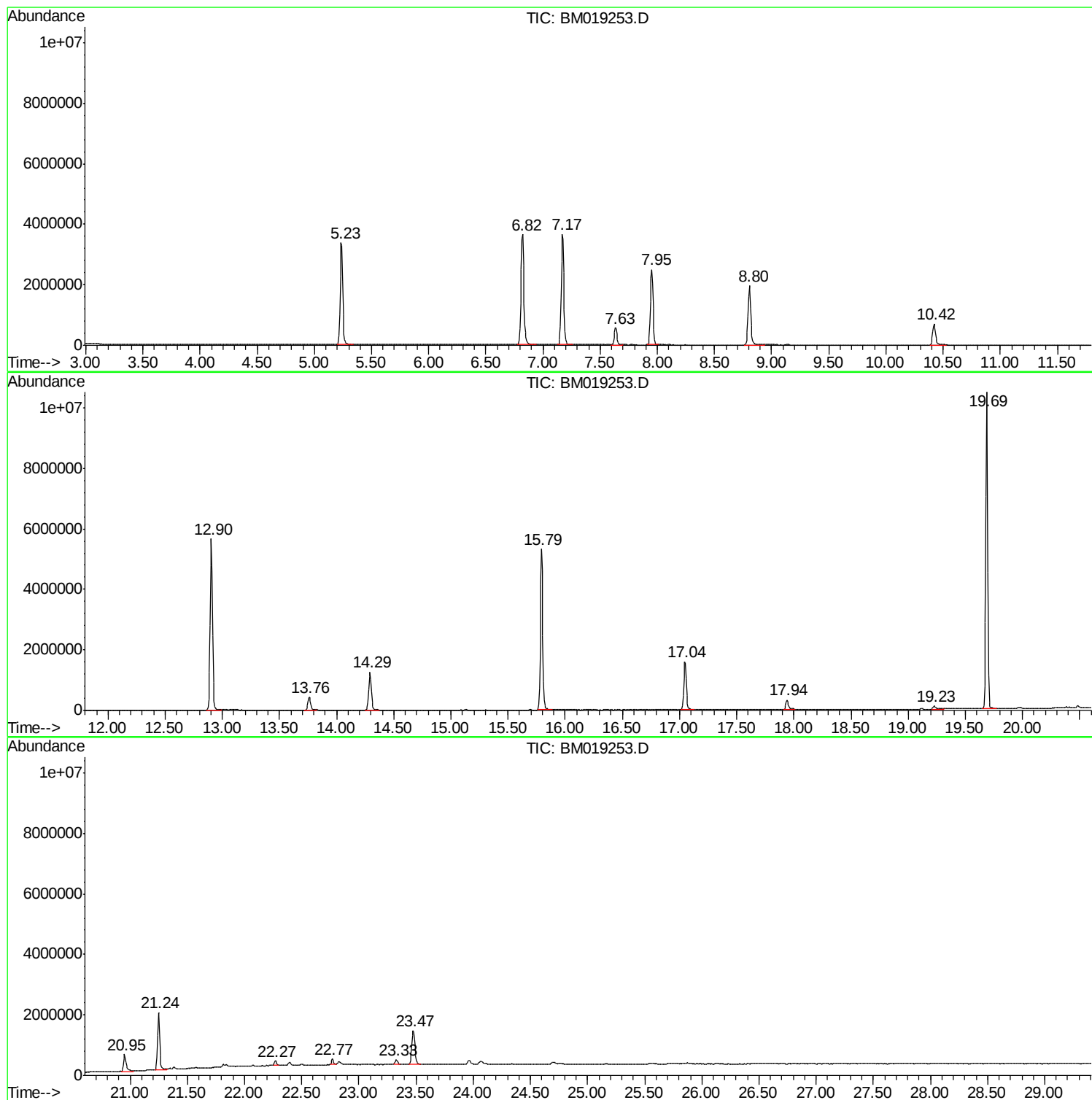
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM031119.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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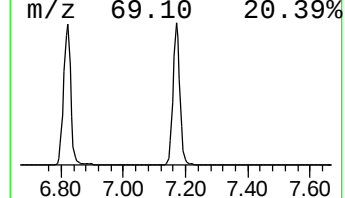
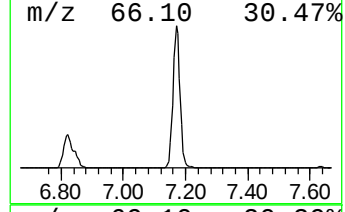
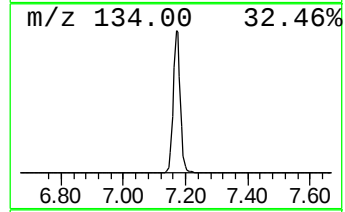
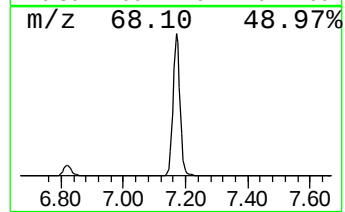
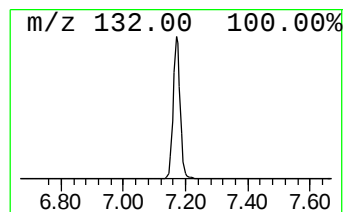
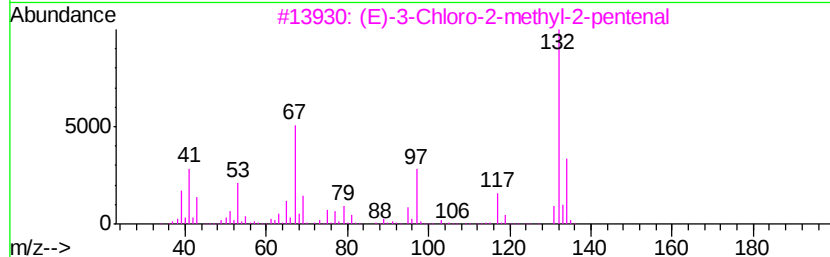
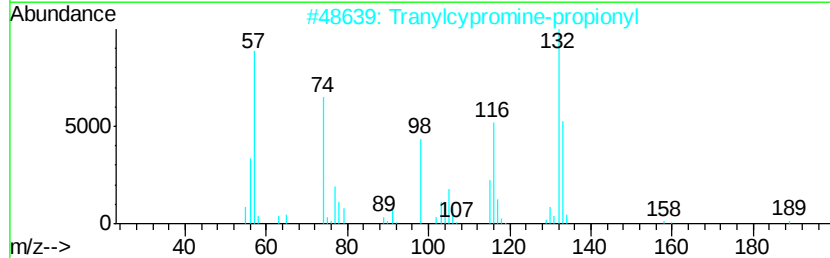
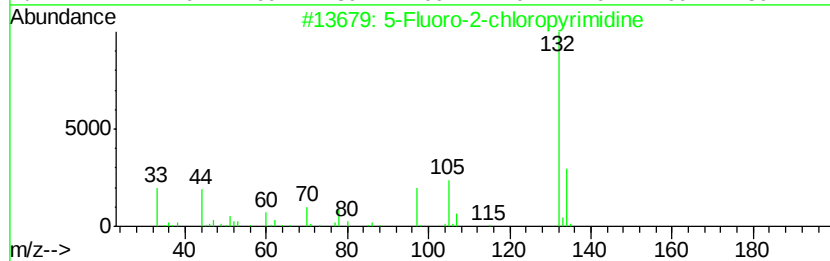
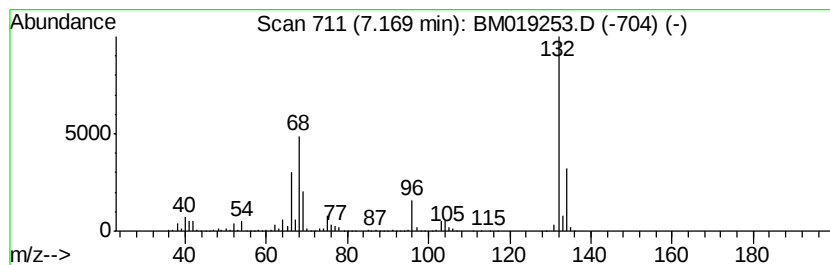
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM031119.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown7.17 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.17	130.35 ng	5864870	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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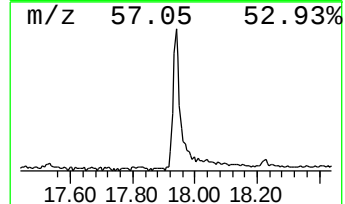
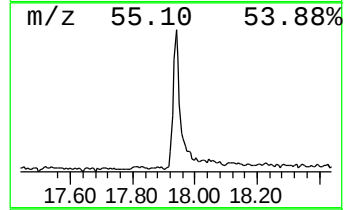
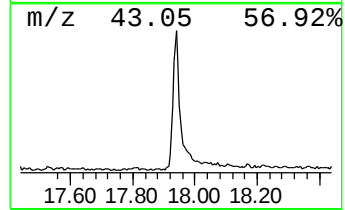
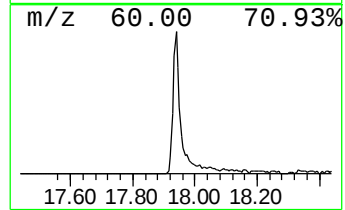
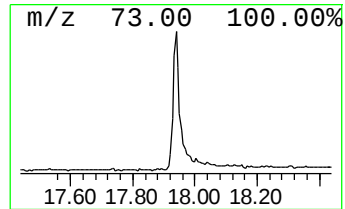
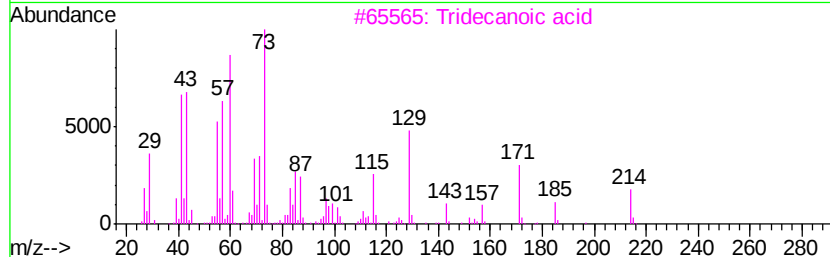
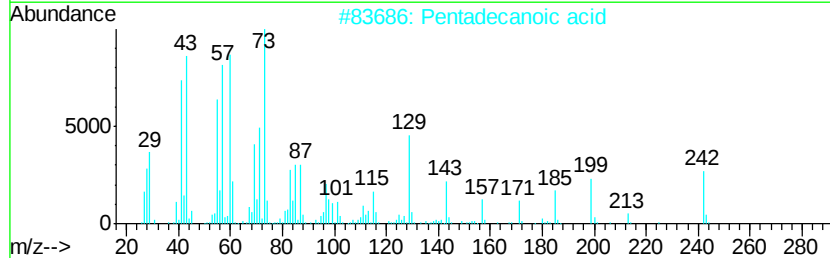
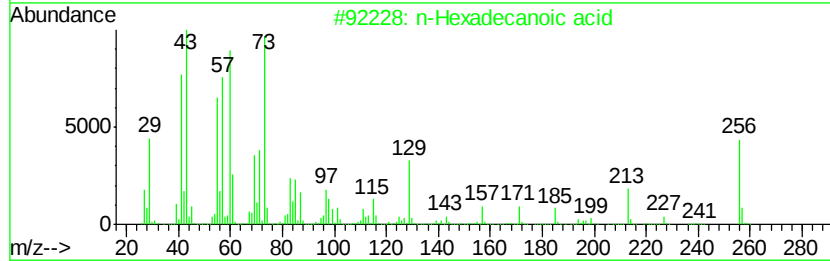
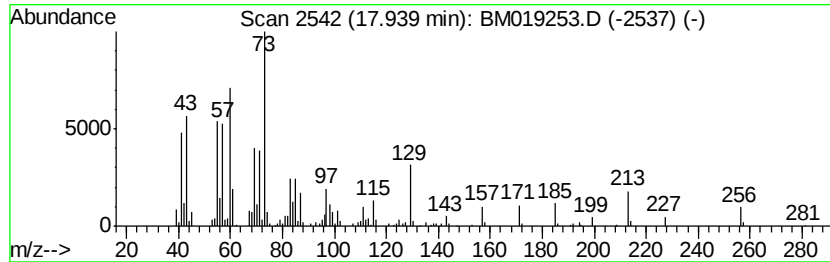
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.94	4.47 ng	538658	Phenanthrene-d10	17.04	
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	72
3	Tridecanoic acid	214	C13H26O2	000638-53-9	55
4	Undecanoic acid	186	C11H22O2	000112-37-8	38
5	Decane, 1-iodo-	268	C10H21I	002050-77-3	18



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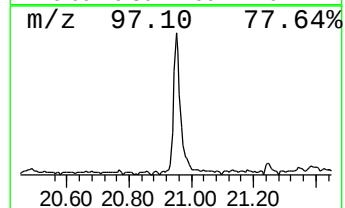
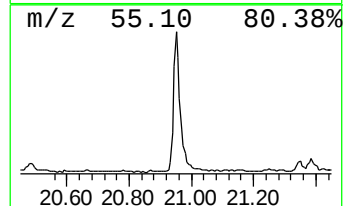
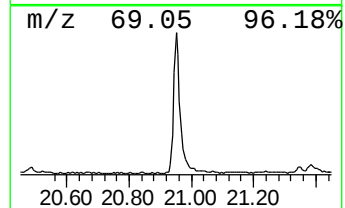
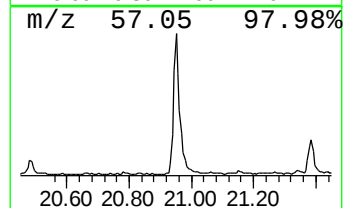
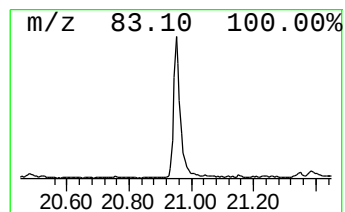
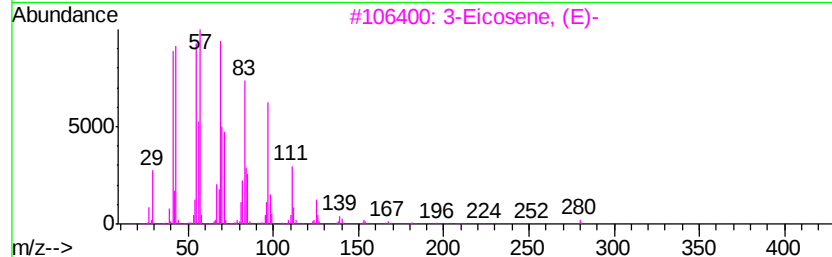
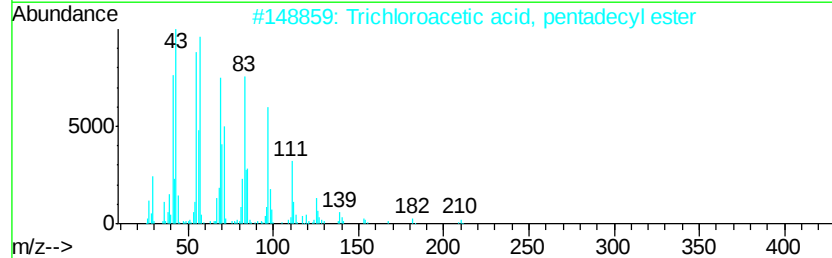
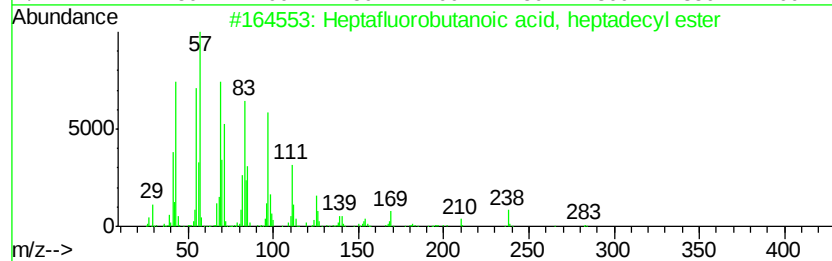
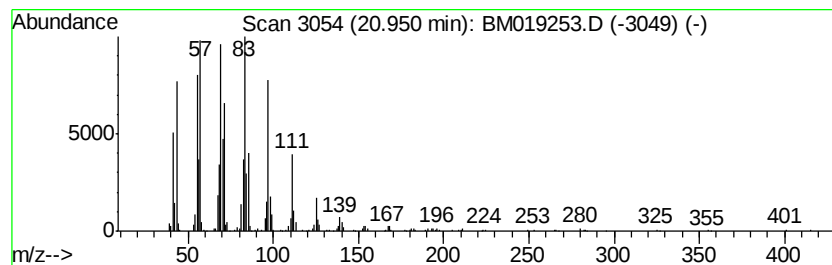
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Peak Number 4 Heptafluorobutanoic acid, h... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.95	6.60 ng	898908	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	91
2		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
3		3-Eicosene, (E)-	280	C20H40	074685-33-9	91
4		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
5		Trifluoroacetic acid, n-octadecy...	366	C20H37F3O2	079392-43-1	91



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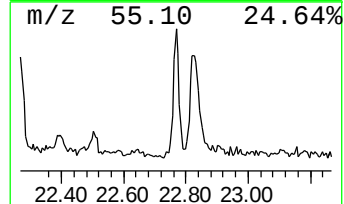
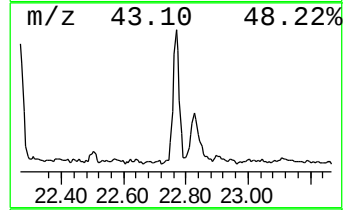
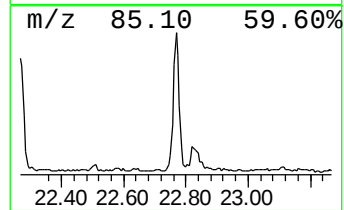
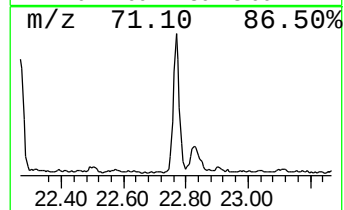
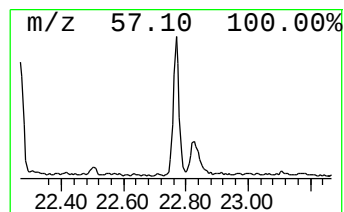
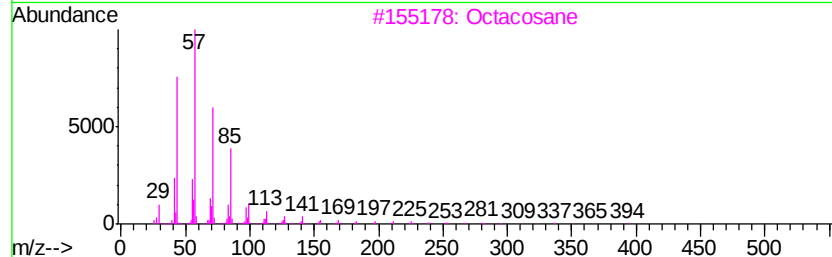
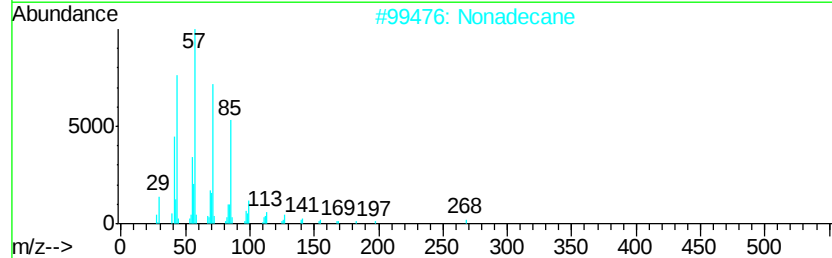
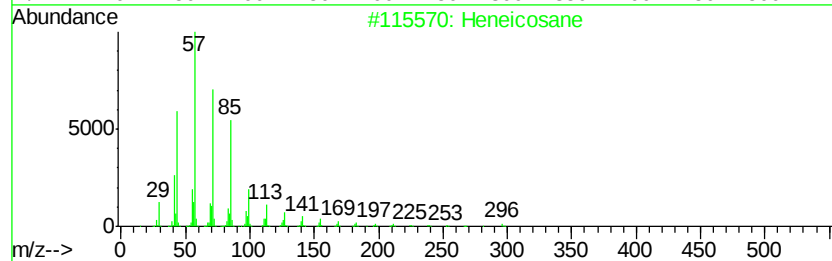
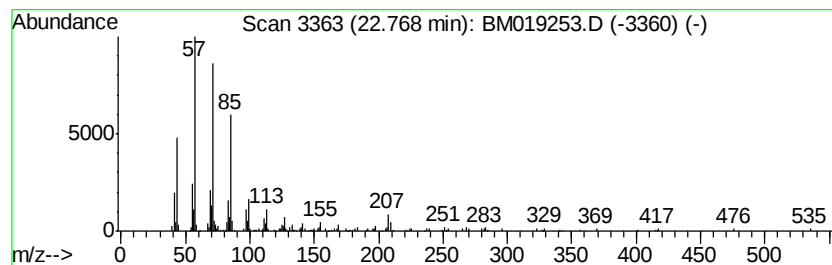
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Peak Number 5 Heneicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.77	2.02 ng	215876	Perylene-d12	23.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	96
2		Nonadecane	268	C19H40	000629-92-5	93
3		Octacosane	394	C28H58	000630-02-4	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	90



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.17	7.17	130.3	ng	5864870	1	7.63	899846	20.0
n-Hexadecanoic acid	17.94	4.5	ng	538658	4	17.04	2411410	20.0
Heptafluorobutano...	20.95	6.6	ng	898908	5	21.24	2724740	20.0
Heneicosane	22.77	2.0	ng	215876	6	23.47	2137050	20.0