

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112522\
 Data File : BM037746.D
 Acq On : 26 Nov 2022 23:00
 Operator : CG/JU
 Sample : N5694-04
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BHB40

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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_M\Methods\SFAM-EPA-BM112322.M
 Title : SVOA CALIBRATION

Signal : TIC: BM037746.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.434	5	8	11	rBV	42136	50886	0.73%	0.074%
2	3.469	11	14	22	rVB	57674	83190	1.20%	0.121%
3	3.881	81	84	99	rBV	181766	354440	5.11%	0.517%
4	4.116	120	124	132	rVB3	27953	46566	0.67%	0.068%
5	4.210	137	140	143	rVB2	12787	14343	0.21%	0.021%
6	4.816	240	243	246	rBV4	5599	7997	0.12%	0.012%
7	4.951	264	266	272	rVB5	3823	6963	0.10%	0.010%
8	5.169	301	303	307	rVB4	6471	9177	0.13%	0.013%
9	7.257	652	658	668	rVB	279178	437497	6.31%	0.639%
10	7.428	680	687	695	rBV	821703	1272526	18.34%	1.857%
11	7.634	714	722	729	rBV	835297	1290372	18.60%	1.884%
12	7.716	732	736	740	rBV5	9199	15430	0.22%	0.023%

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37	12.327	1511	1520	1529	rVB8	13099	28228	0.41%	0.041%
38	12.427	1529	1537	1542	rBV9	8358	22347	0.32%	0.033%
39	12.504	1542	1550	1554	rVV3	28191	50784	0.73%	0.074%
40	12.545	1554	1557	1561	rVB4	11086	14597	0.21%	0.021%
41	12.716	1582	1586	1589	rVB5	5762	7712	0.11%	0.011%
42	12.986	1630	1632	1635	rVV4	7042	7720	0.11%	0.011%
43	13.027	1636	1639	1643	rVV6	4820	7404	0.11%	0.011%
44	13.098	1647	1651	1654	rBV5	5830	8893	0.13%	0.013%
45	13.210	1668	1670	1675	rBV5	4790	7751	0.11%	0.011%
46	13.551	1719	1728	1733	rBV3	24136	39547	0.57%	0.058%
47	13.627	1736	1741	1747	rVB4	26639	53745	0.77%	0.078%
48	13.704	1749	1754	1756	rVV4	10006	15567	0.22%	0.023%
49	14.139	1820	1828	1834	rBV	2412855	3215716	46.36%	

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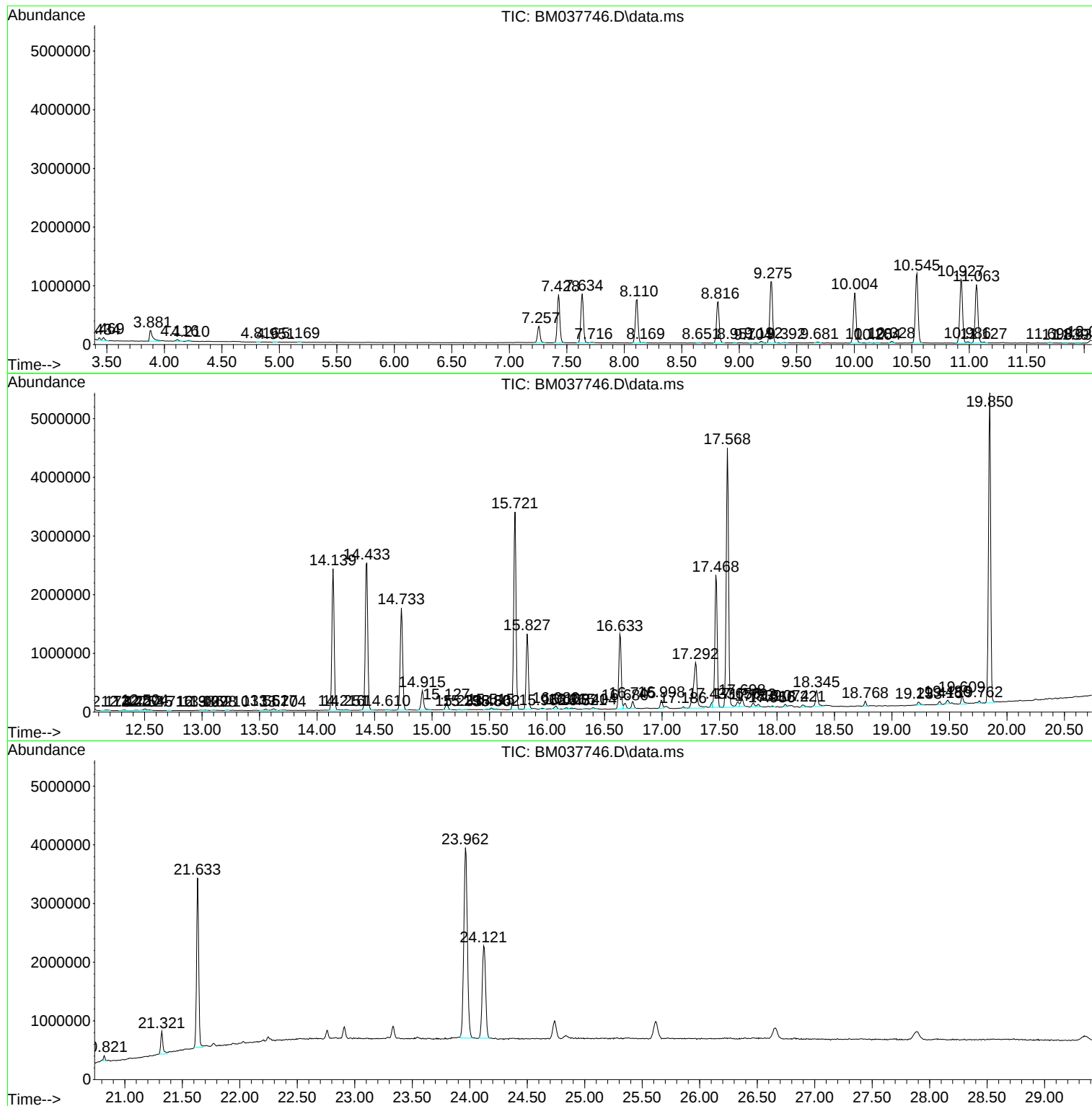
77	17.468	2389	2394	2404	rVB	2257231	3201675	46.15%	4.673%
78	17.568	2404	2411	2420	rBV	4424404	6101706	87.96%	8.907%
79	17.656	2422	2426	2429	rVV2	83096	136656	1.97%	0.199%
80	17.698	2429	2433	2438	rVB	132392	205613	2.96%	0.300%
81	17.792	2445	2449	2454	rBV	67188	109758	1.58%	0.160%
82	17.839	2454	2457	2462	rVB3	43219	60097	0.87%	0.088%
83	17.956	2475	2477	2482	rVB5	18743	22117	0.32%	0.032%
84	18.074	2493	2497	2500	rBV	41214	55389	0.80%	0.081%
85	18.221	2519	2522	2528	rBV3	35855	67534	0.97%	0.099%
86	18.345	2537	2543	2553	rBV3	268913	416920	6.01%	0.609%
87	18.768	2611	2615	2619	rBV2	85866	96821	1.40%	0.141%
88	19.233	2690	2694	2702	rBV	51473	83719	1.21%	0.122%
89	19.415	2721	2725	2729	rBV3	55748	73397	1.0	

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

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



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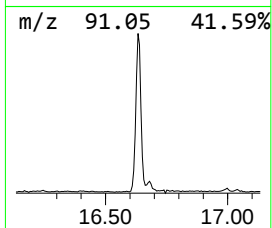
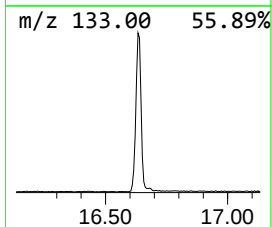
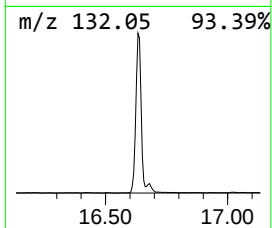
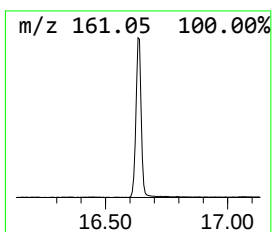
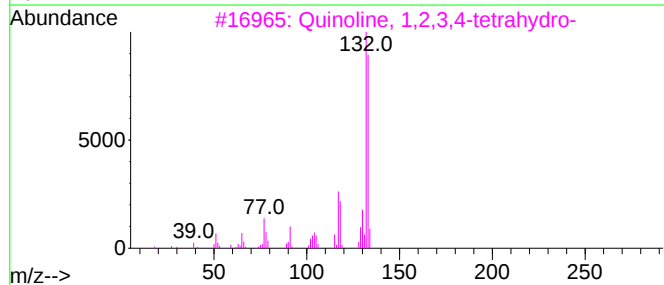
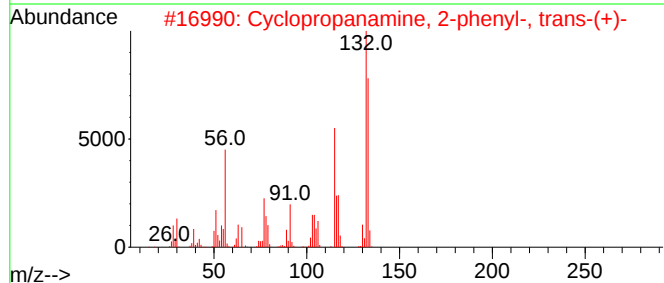
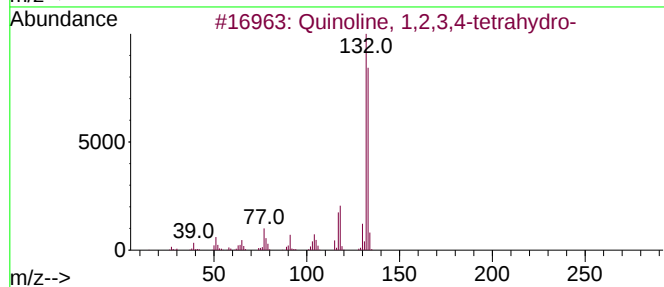
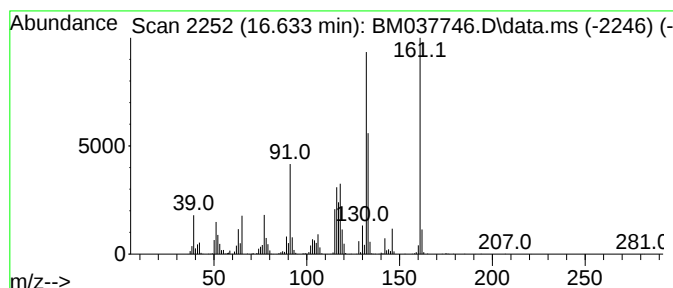
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Quinoline, 1,2,3,4-tetrahydro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.633	11.59 ng/ul	1855180	Phenanthrene-d10	17.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	55
2			Cyclopropanamine, 2-phenyl-, tra...	133	C9H11N	003721-28-6	55
3			Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	50
4			Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	50
5			Tranlylcypromine	133	C9H11N	000155-09-9	49



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Instrument :
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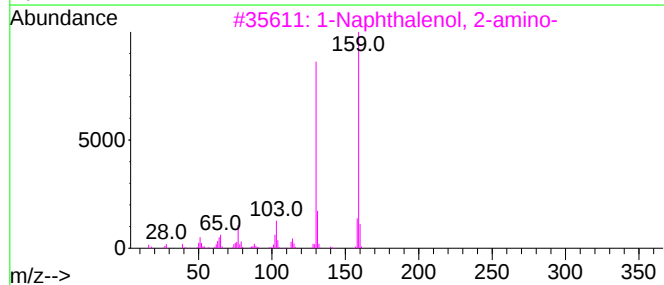
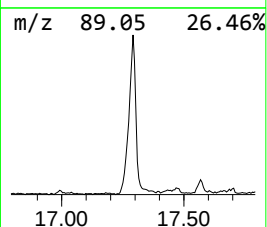
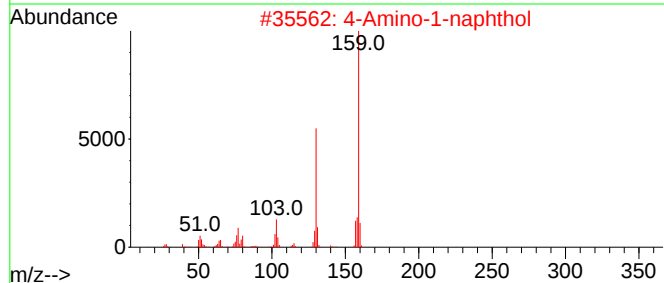
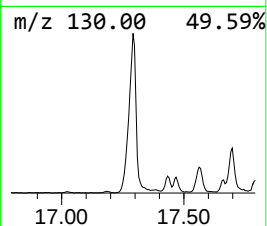
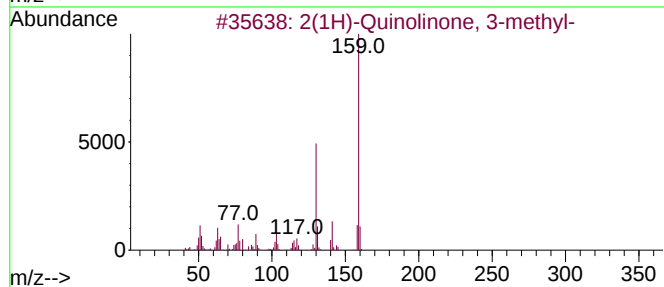
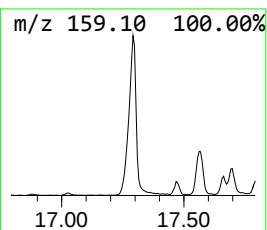
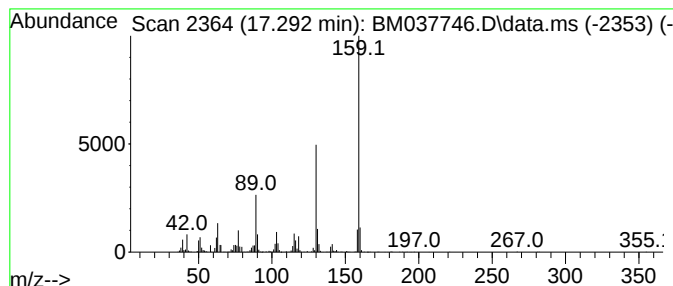
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2(1H)-Quinolinone, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.292	10.65 ng/ul	1704970	Phenanthrene-d10	17.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2(1H)-Quinolinone, 3-methyl-	159	C10H9NO	002721-59-7	83
2			4-Amino-1-naphthol	159	C10H9NO	002834-90-4	81
3			1-Naphthalenol, 2-amino-	159	C10H9NO	000606-41-7	81
4			2-Naphthalenol, 1-amino-	159	C10H9NO	002834-92-6	81
5			1-Naphthalenol, 6-amino-	159	C10H9NO	023894-12-4	76



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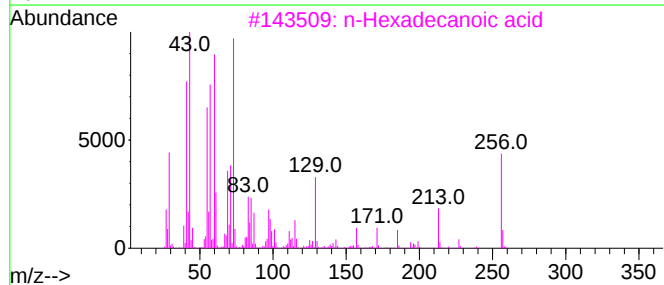
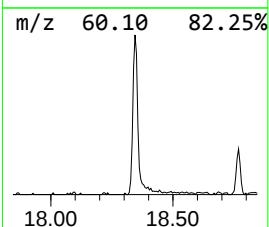
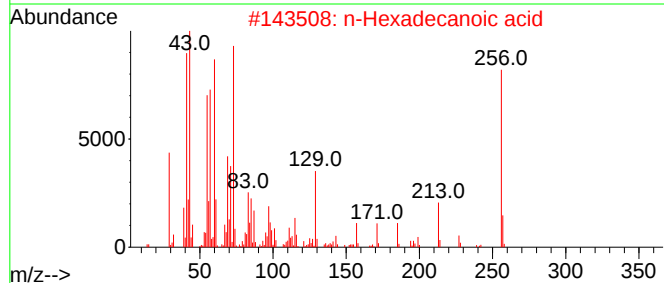
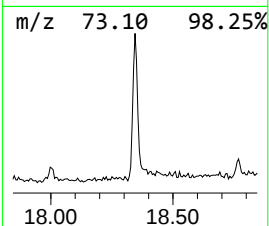
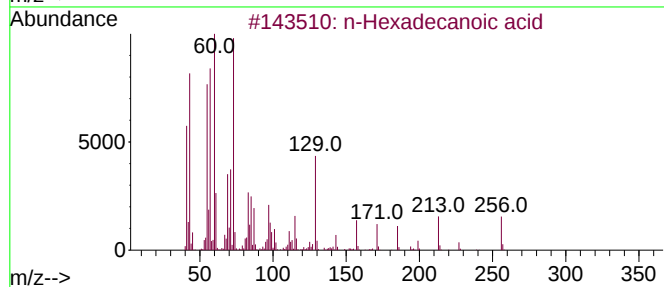
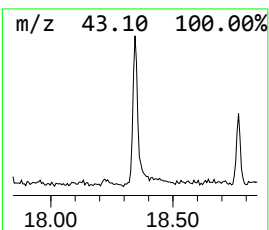
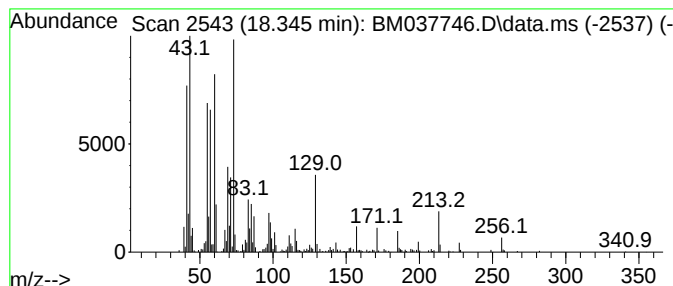
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.345	2.60 ng/ul	416920	Phenanthrene-d10	17.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	92
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	91



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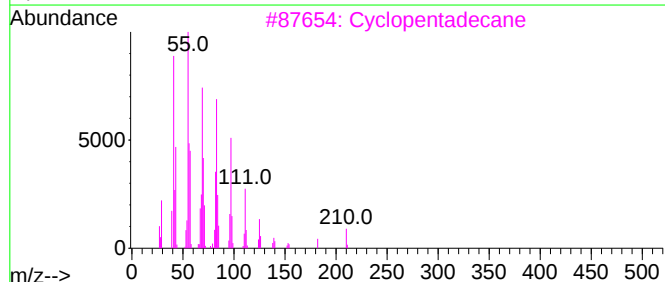
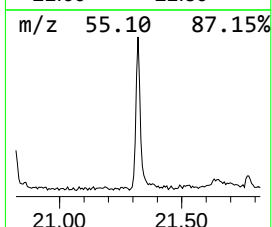
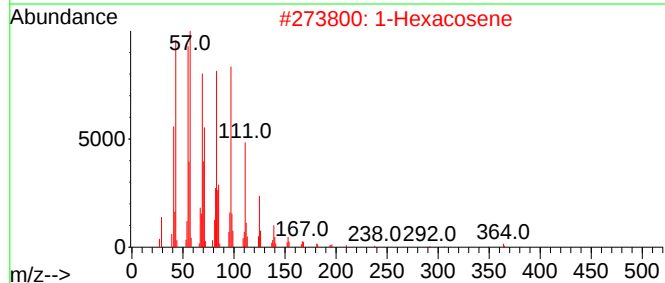
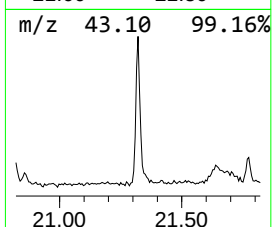
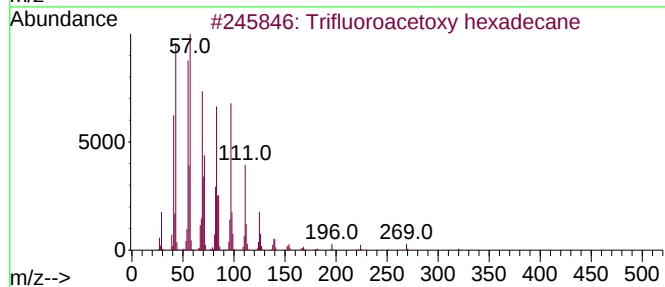
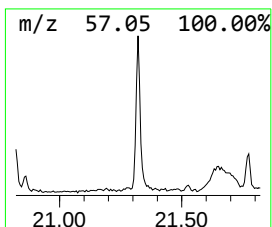
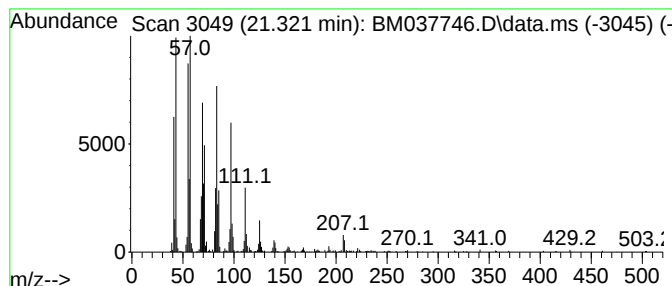
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Quant Title : SVOA CALIBRATION

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TIC Integration Parameters: LSCINT.P

Peak Number 4 Trifluoroacetoxy hexadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.321	2.87 ng/ul	522425	Chrysene-d12	21.633

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
2			1-Hexacosene	364	C26H52	018835-33-1	94
3			Cyclopentadecane	210	C15H30	000295-48-7	93
4			Cyclotetracosane	336	C24H48	000297-03-0	91
5			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	91



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Quinoline, 1,2,...	16.633	11.6	ng/ul	1855180	4	17.468	3201680	20.0
2(1H)-Quinolino...	17.292	10.7	ng/ul	1704970	4	17.468	3201680	20.0
n-Hexadecanoic ...	18.345	2.6	ng/ul	416920	4	17.468	3201680	20.0
Trifluoroacetox...	21.321	2.9	ng/ul	522425	5	21.633	3645330	20.0