

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091324\
 Data File : BP021914.D
 Acq On : 13 Sep 2024 18:32
 Operator : RC/JU
 Sample : P3952-06
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B65

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_P\Methods\SFAM-EPA-BP091124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021914.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.270	44	48	52	rVB	24090	28922	0.88%	0.096%
2	3.540	90	94	102	rVB	32199	49812	1.51%	0.166%
3	3.681	114	118	137	rVB	136983	219441	6.65%	0.730%
4	3.999	169	172	177	rVB3	5121	8539	0.26%	0.028%
5	4.446	244	248	254	rVB8	2627	4097	0.12%	0.014%
6	4.793	302	307	311	rBV7	3325	6287	0.19%	0.021%
7	4.917	322	328	334	rBV2	14573	22185	0.67%	0.074%
8	6.046	515	520	524	rVB8	4634	6778	0.21%	0.023%
9	6.934	662	671	681	rBV	84996	148092	4.49%	0.492%
10	7.093	690	698	709	rBV	206675	319455	9.68%	1.062%
11	7.299	723	733	742	rBV	278022	442688	13.41%	1.472%
12	7.375	742	746	755	rVB7	4293	10153	0.31%	0.034%
13	7.581	776	781	787	rBV9	1729	4008	0.12%	0.013%
14	7.758	803	811	818	rBV	412472	689757	20.89%	2.294%
15	7.822	820	822	827	rVB5	2959	4150	0.13%	0.014%
16	8.452	922	929	939	rBV	340729	569847	17.26%	1.895%
17	8.899	997	1005	1017	rBV	355130	600344	18.19%	1.996%
18	9.328	1073	1078	1084	rBV2	8811	13810	0.42%	0.046%
19	9.457	1095	1100	1106	rVB2	12539	19038	0.58%	0.063%
20	9.610	1118	1126	1140	rBV	266392	472130	14.30%	1.570%
21	10.152	1210	1218	1236	rBV	518376	864520	26.19%	2.875%
22	10.522	1273	1281	1293	rBV	570649	978658	29.64%	3.255%
23	10.652	1295	1303	1317	rBV	466771	831685	25.19%	2.766%
24	11.057	1368	1372	1378	rBV9	4125	7611	0.23%	0.025%
25	11.263	1399	1407	1423	rBV2	39596	103677	3.14%	0.345%
26	11.810	1494	1500	1507	rBV3	6955	13143	0.40%	0.044%
27	12.104	1543	1550	1558	rVB4	7742	15664	0.47%	0.052%
28	12.975	1694	1698	1702	rBV7	3314	5101	0.15%	0.017%
29	13.263	1743	1747	1754	rVB9	3615	7147	0.22%	0.024%
30	13.769	1826	1833	1846	rBV	877974	1379428	41.78%	4.587%
31	14.057	1871	1882	1893	rBV2	899515	1513042	45.83%	5.032%
32	14.369	1926	1935	1944	rBV	944368	1460293	44.23%	4.856%
33	14.575	1964	1970	1985	rBV2	116296	273981	8.30%	0.911%
34	14.792	2000	2007	2016	rVB10	10388	19969	0.60%	0.066%
35	15.187	2070	2074	2080	rVB6	4742	9851	0.30%	0.033%

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36	15.375	2099	2106	2119	rBV	1294269	2040953	61.82%	6.787%
37	15.492	2120	2126	2132	rBV	337069	485445	14.70%	1.614%
38	15.616	2142	2147	2157	rVB7	9326	19327	0.59%	0.064%
39	16.116	2228	2232	2234	rBV4	4160	4556	0.14%	0.015%
40	16.486	2292	2295	2300	rVB7	7042	9587	0.29%	0.032%
41	16.922	2365	2369	2375	rBV9	10333	19518	0.59%	0.065%
42	17.151	2401	2408	2414	rBV	1219472	1877271	56.87%	6.243%
43	17.251	2419	2425	2437	rVV	1686106	2399903	72.70%	7.981%
44	17.475	2461	2463	2468	rVB6	6361	9112	0.28%	0.030%
45	17.851	2523	2527	2532	rVB7	10367	14221	0.43%	0.047%
46	18.081	2560	2566	2575	rBV	123176	188357	5.71%	0.626%
47	19.169	2747	2751	2756	rVB2	26575	29811	0.90%	0.099%
48	19.239	2759	2763	2766	rBV	20684	26917	0.82%	0.090%
49	19.398	2786	2790	2799	rBV4	51138	79377	2.40%	0.264%
50	19.610	2821	2826	2834	rBV	2457494	3116699	94.41%	10.365%
51	21.251	3101	3105	3115	rBV3	138185	303351	9.19%	1.009%
52	21.574	3154	3160	3172	rVB2	1404859	2392874	72.48%	7.958%
53	24.674	3678	3687	3701	rVB	1319213	3301266	100.00%	10.979%
54	24.904	3717	3726	3743	rVB	963940	2628130	79.61%	8.740%

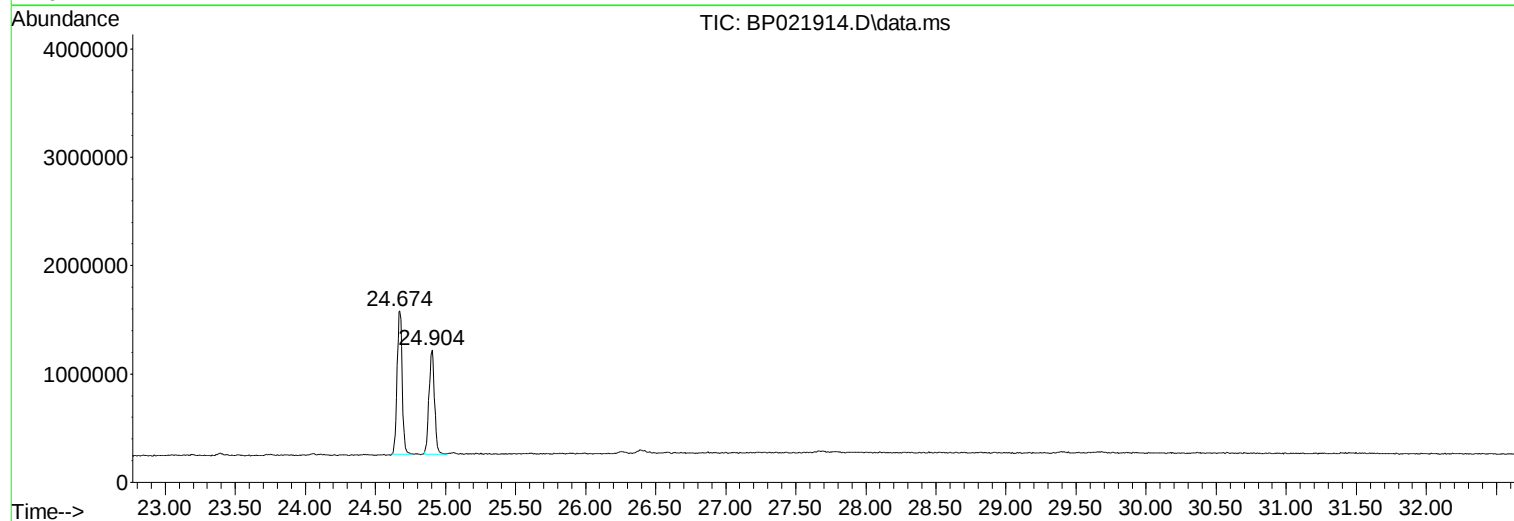
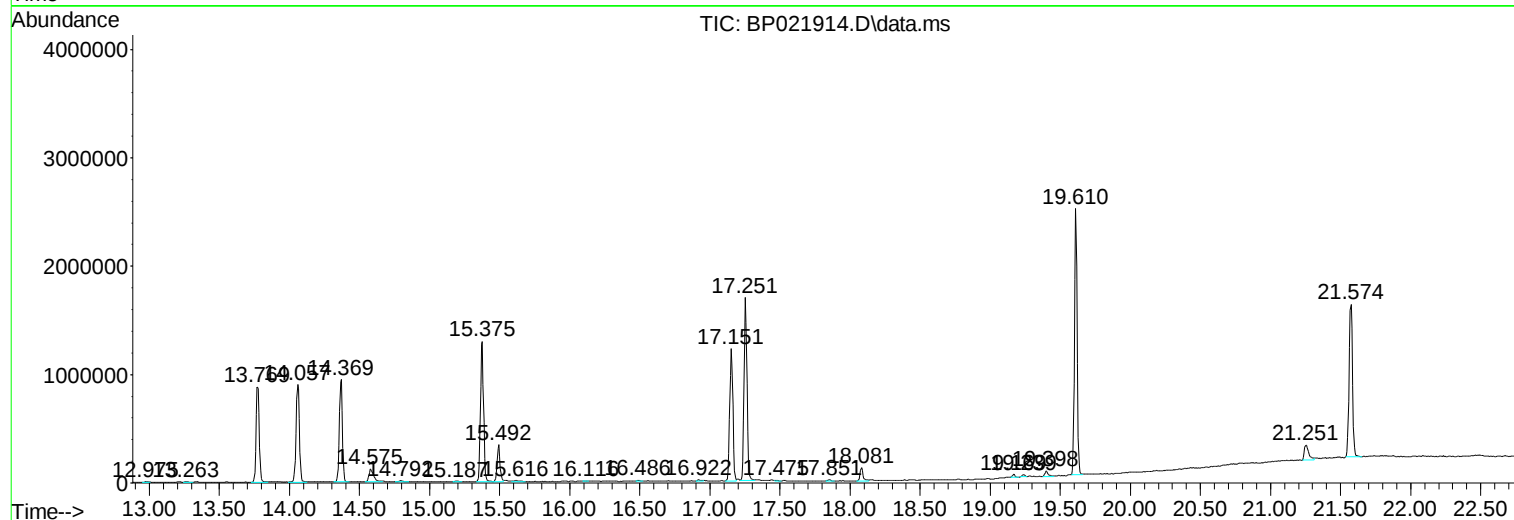
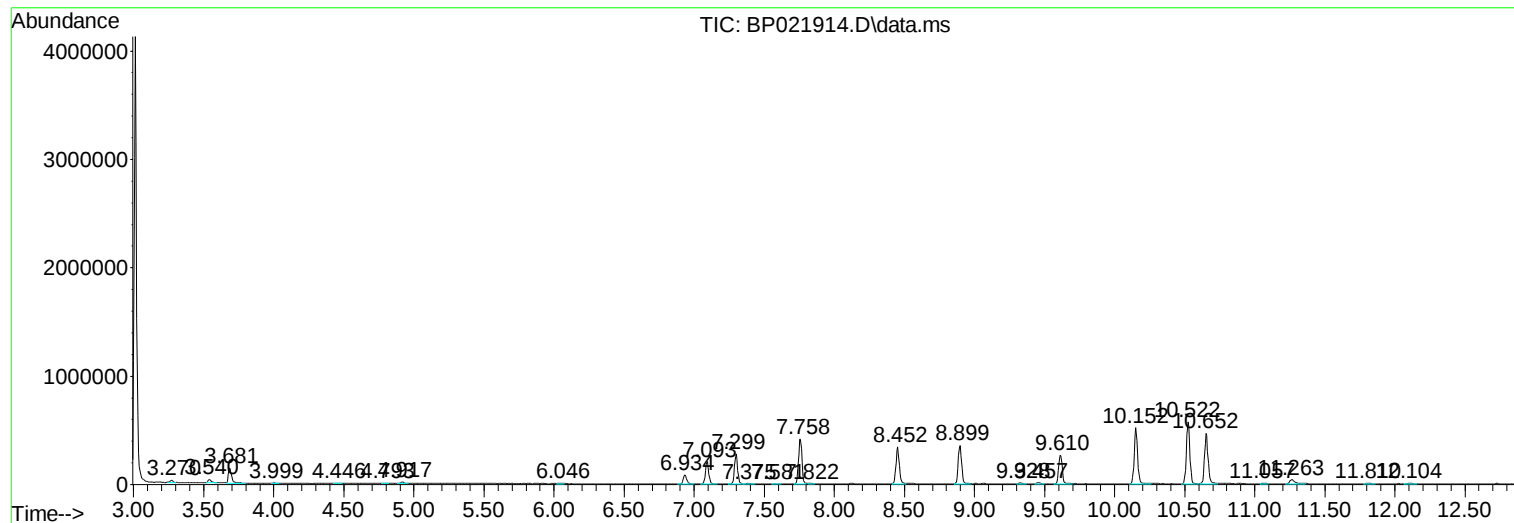
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Quant Title : SVOA CALIBRATION

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



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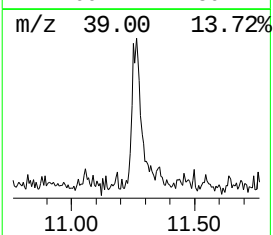
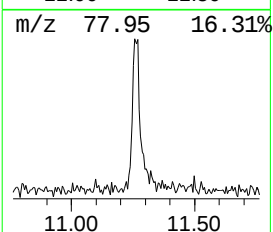
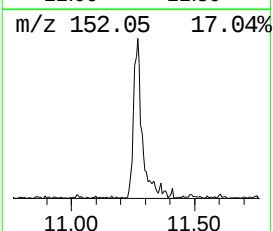
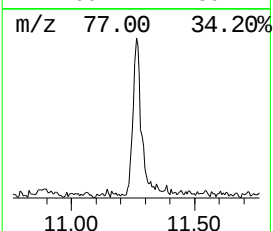
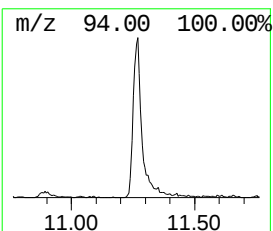
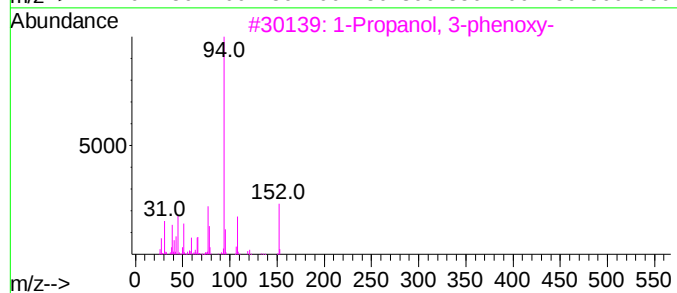
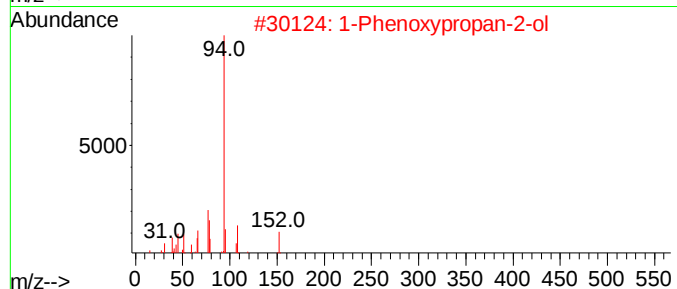
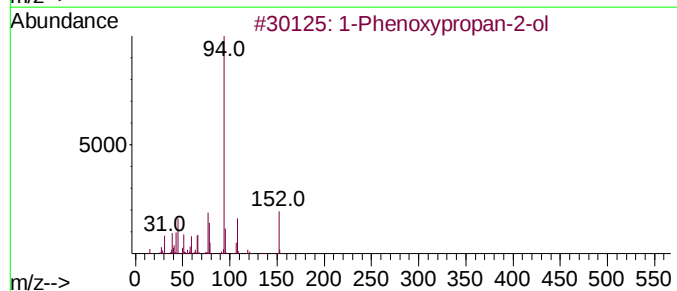
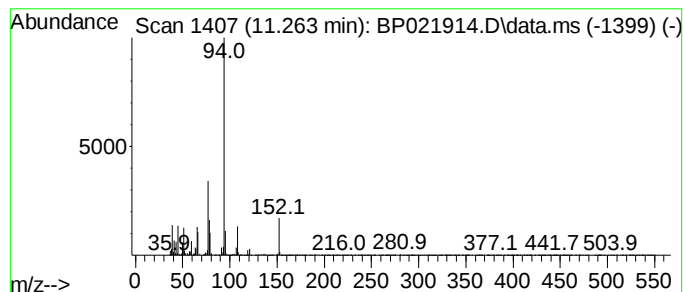
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Phenoxypropan-2-ol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.263	2.12 ng/ul	103677	Naphthalene-d8	10.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	81
3			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	81
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	64
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	64



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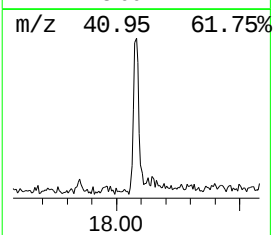
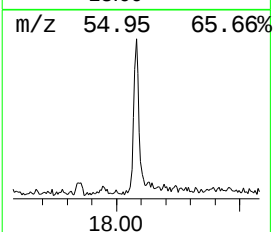
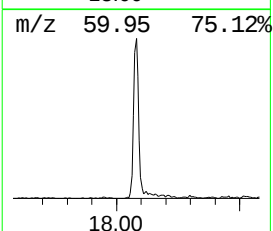
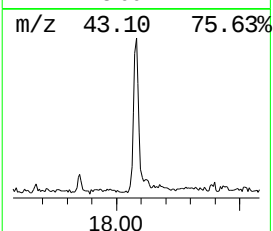
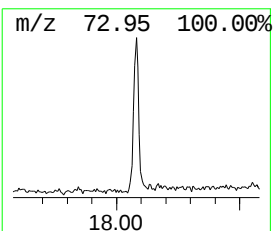
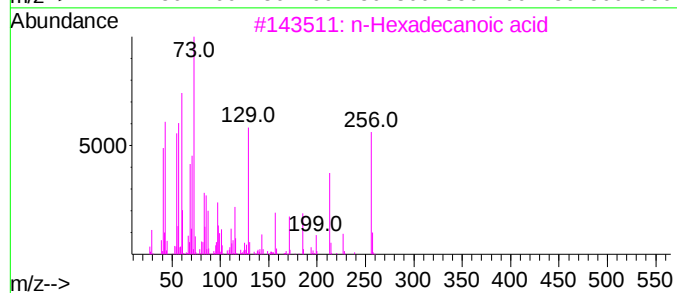
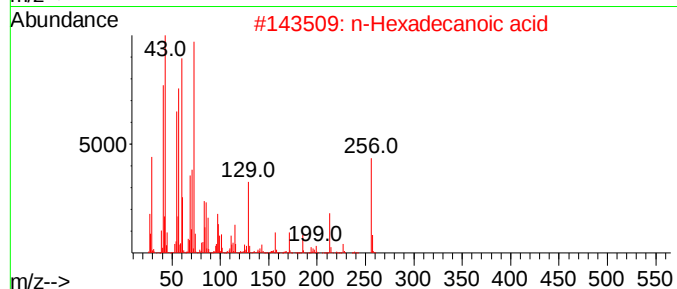
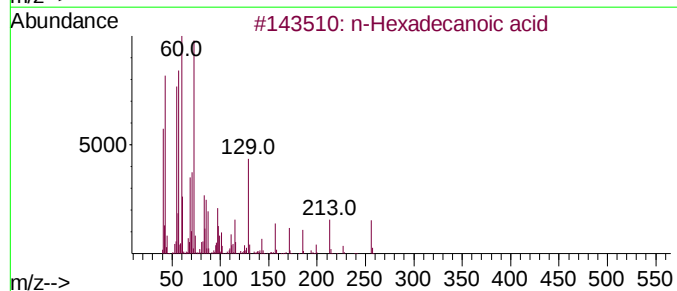
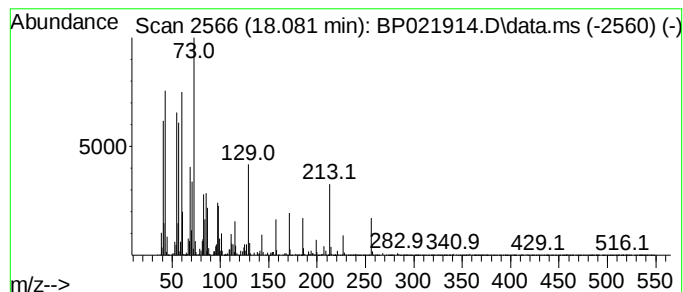
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.081	2.01 ng/ul	188357	Phenanthrene-d10	17.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
5			Tridecanoic acid	214	C13H26O2	000638-53-9	91



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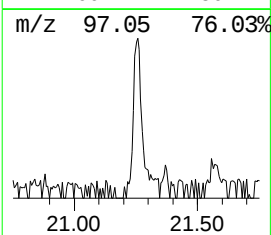
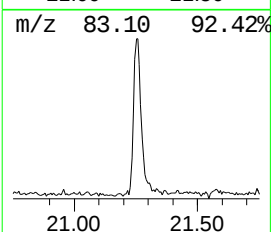
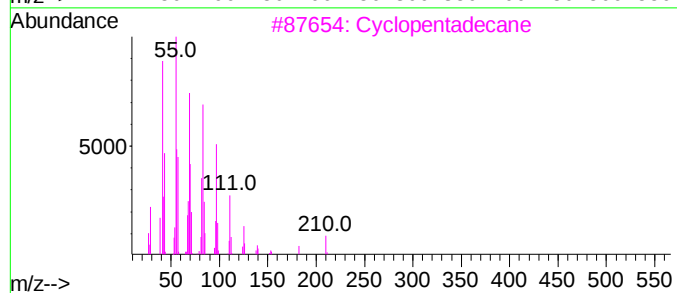
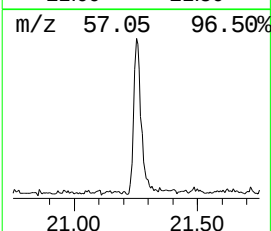
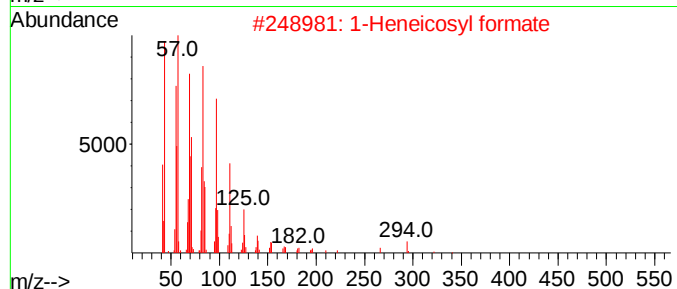
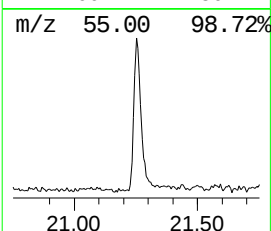
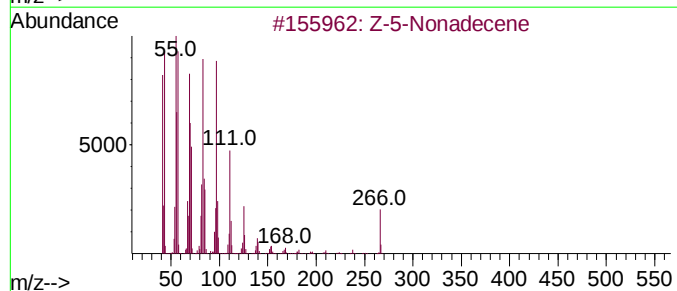
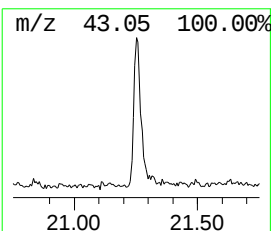
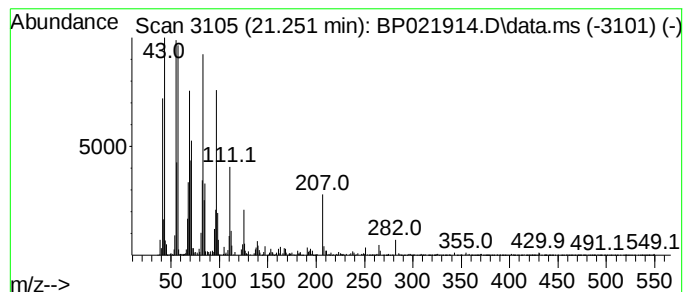
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.251	2.54 ng/ul	303351	Chrysene-d12	21.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Z-5-Nonadecene	266	C19H38	1000131-11-8	96
2			1-Heneicosyl formate	340	C22H44O2	077899-03-7	95
3			Cyclopentadecane	210	C15H30	000295-48-7	93
4			1-Heneicosanol	312	C21H44O	015594-90-8	93
5			1-Heptacosanol	396	C27H56O	002004-39-9	93



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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
1-Phenoxypropan...	11.263	2.1	ng/ul	103677	2	10.522	978658	20.0
n-Hexadecanoic ...	18.081	2.0	ng/ul	188357	4	17.151	1877270	20.0
(DEL) Alkane: S...	21.251	2.5	ng/ul	303351	5	21.575	2392870	20.0