

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
 Data File : BN023301.D
 Acq On : 20 Dec 2022 16:42
 Operator : CG/JU
 Sample : N6003-23
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBXV2

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_N\Methods\SFAM-EPA-BN121522.M
 Title : SVOA CALIBRATION

Signal : TIC: BN023301.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.340	40	43	51	rVB	43020	61347	0.94%	0.071%
2	3.775	114	117	131	rBV	183706	324143	4.94%	0.375%
3	3.887	132	136	143	rVB	43938	62073	0.95%	0.072%
4	4.011	153	157	163	rVB	22417	36509	0.56%	0.042%
5	4.111	170	174	179	rVB	15660	24550	0.37%	0.028%
6	4.499	235	240	244	rBV6	6413	10161	0.15%	0.012%
7	5.063	329	336	344	rBV3	29899	54426	0.83%	0.063%
8	5.152	349	351	360	rVB6	7861	12474	0.19%	0.014%
9	6.975	655	661	666	rBV8	6823	14188	0.22%	0.016%
10	7.152	682	691	701	rBV	897609	1418404	21.63%	1.642%
11	7.322	713	720	728	rBV	750605	1139501	17.37%	1.319%
12	7.522	744	754	764	rBV	947555	1475751	22.50%	1.708%
13	7.610	766	769	776	rVB7	8684	13946	0.21%	0.016%
14	7.999	828	835	841	rBV	881013	1368282	20.86%	1.584%
15	8.057	843	845	850	rVB2	8639	12572	0.19%	0.015%
16	8.710	948	956	967	rBV	990351	1559445	23.78%	1.805%
17	8.828	972	976	983	rVB2	22608	38760	0.59%	0.045%
18	9.169	1025	1034	1043	rBV	928379	1501445	22.89%	1.738%
19	9.328	1056	1061	1071	rVB9	7611	16997	0.26%	0.020%
20	9.587	1097	1105	1110	rBV	22546	36891	0.56%	0.043%
21	9.893	1146	1157	1173	rBV	795869	1321632	20.15%	1.530%
22	10.122	1191	1196	1205	rVB8	5518	12356	0.19%	0.014%
23	10.434	1241	1249	1262	rBV	1264824	2035763	31.04%	2.356%
24	10.598	1271	1277	1282	rVB8	6141	10654	0.16%	0.012%
25	10.816	1306	1314	1327	rBV	1362532	2226209	33.94%	2.577%
26	10.957	1327	1338	1352	rBV	1009460	1744378	26.60%	2.019%
27	11.351	1399	1405	1409	rVB8	7853	13389	0.20%	0.015%
28	11.598	1441	1447	1455	rBV8	6345	13902	0.21%	0.016%
29	12.234	1547	1555	1560	rBV	21703	32689	0.50%	0.038%
30	12.404	1577	1584	1589	rVV2	22416	38880	0.59%	0.045%
31	12.969	1675	1680	1684	rBV6	7087	10023	0.15%	0.012%
32	13.181	1711	1716	1721	rVB3	9014	11472	0.17%	0.013%
33	13.469	1761	1765	1773	rVV	42120	59536	0.91%	0.069%
34	13.563	1773	1781	1786	rVV6	13133	29779	0.45%	0.034%
35	13.610	1786	1789	1796	rVB6	7570	17905	0.27%	0.021%
36	13.822	1821	1825	1830	rVB5	9309	13062	0.20%	0.015%

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Integrator: RTE

Smoothing : OFF

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Sampling : 1

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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_N\Methods\SFAM-EPA-BN121522.M
 Title : SVOA CALIBRATION

37	13.969	1846	1850	1854	rBV5	12804	18106	0.28%	0.021%
38	14.057	1858	1865	1874	rVV	2243922	3076039	46.90%	3.560%
39	14.122	1874	1876	1880	rVB4	9572	12558	0.19%	0.015%
40	14.169	1880	1884	1893	rVB8	12148	21147	0.32%	0.024%

41	14.339	1907	1913	1924	rVB	2627316	3801169	57.96%	4.400%
42	14.539	1943	1947	1952	rVB2	22847	29449	0.45%	0.034%
43	14.645	1955	1965	1973	rBV	2117785	3094185	47.18%	3.581%
44	14.839	1991	1998	2013	rBV	1091157	1535718	23.42%	1.778%
45	14.992	2020	2024	2029	rVV4	18652	29283	0.45%	0.034%

46	15.057	2029	2035	2041	rVB3	51404	81463	1.24%	0.094%
47	15.210	2056	2061	2062	rBV4	7221	10345	0.16%	0.012%
48	15.486	2105	2108	2112	rVB2	10607	12552	0.19%	0.015%
49	15.639	2127	2134	2140	rBV	3374669	4782190	72.92%	5.535%
50	15.692	2140	2143	2147	rVB4	11416	15024	0.23%	0.017%

51	15.751	2147	2153	2162	rBV	1019980	1369979	20.89%	1.586%
52	15.875	2170	2174	2182	rVB3	16998	29890	0.46%	0.035%
53	15.963	2185	2189	2190	rBV3	8480	10722	0.16%	0.012%
54	16.004	2190	2196	2205	rVB	1107920	1456491	22.21%	1.686%
55	16.351	2247	2255	2262	rVV	12355	36881	0.56%	0.043%

56	16.510	2276	2282	2285	rBV6	10181	18201	0.28%	0.021%
57	16.569	2288	2292	2296	rVB2	11949	16766	0.26%	0.019%
58	16.681	2306	2311	2315	rBV8	4838	9670	0.15%	0.011%
59	16.786	2326	2329	2337	rVB9	10050	17784	0.27%	0.021%
60	16.928	2349	2353	2357	rBV4	14317	19678	0.30%	0.023%

61	17.045	2369	2373	2375	rBV5	7758	10325	0.16%	0.012%
62	17.280	2409	2413	2421	rBV4	35174	63453	0.97%	0.073%
63	17.351	2421	2425	2426	rBV3	16596	18791	0.29%	0.022%
64	17.392	2426	2432	2437	rVV	2846489	3858455	58.83%	4.466%
65	17.433	2437	2439	2443	rVV	75811	93498	1.43%	0.108%

66	17.492	2443	2449	2458	rVV	3666479	4994422	76.15%	5.781%
67	17.580	2459	2464	2473	rVB5	77451	125182	1.91%	0.145%
68	17.769	2493	2496	2499	rBV4	18438	27797	0.42%	0.032%
69	18.057	2543	2545	2550	rVB5	16847	19360	0.30%	0.022%
70	18.110	2552	2554	2558	rVB5	9890	9682	0.15%	0.011%

71	18.169	2558	2564	2569	rBV3	43239	79806	1.22%	0.092%
72	18.227	2570	2574	2579	rBV	102059	155596	2.37%	0.180%
73	18.286	2579	2584	2594	rBV	898972	1335516	20.36%	1.546%
74	18.586	2631	2635	2637	rBV4	27111	40312	0.61%	0.047%
75	18.616	2638	2640	2648	rVB3	35209	51144	0.78%	0.059%

76	18.710	2652	2656	2661	rBV	140914	163055	2.49%	0.189%
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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 >
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 Title : SVOA CALIBRATION

77	18.769	2663	2666	2671	rVV3	27938	38405	0.59%	0.044%
78	19.139	2726	2729	2733	rBV5	17594	25520	0.39%	0.030%
79	19.186	2733	2737	2739	rBV3	41380	62240	0.95%	0.072%
80	19.351	2761	2765	2771	rVB	52234	76359	1.16%	0.088%
81	19.416	2772	2776	2778	rVV2	124170	159906	2.44%	0.185%
82	19.445	2778	2781	2794	rVB	268934	456142	6.96%	0.528%
83	19.563	2796	2801	2812	rBV	224476	364045	5.55%	0.421%
84	19.786	2833	2839	2851	rBV	4782905	6558351	100.00%	7.591%
85	20.292	2922	2925	2926	rBV2	37885	41927	0.64%	0.049%
86	20.816	3011	3014	3018	rVB2	85526	96094	1.47%	0.111%
87	21.274	3088	3092	3102	rVB	1412884	1846718	28.16%	2.138%
88	21.480	3124	3127	3132	rBV	678039	752302	11.47%	0.871%
89	21.574	3138	3143	3148	rBV	3112236	4364777	66.55%	5.052%
90	22.198	3245	3249	3250	rBV3	168525	218505	3.33%	0.253%
91	23.268	3426	3431	3436	rBV2	625020	1031033	15.72%	1.193%
92	23.868	3526	3533	3546	rVB	2608633	5720283	87.22%	6.621%
93	24.027	3553	3560	3568	rBV	1929137	3737943	57.00%	4.327%
94	24.268	3598	3601	3613	rVB6	227800	454432	6.93%	0.526%
95	24.545	3640	3648	3657	rVB2	2956025	6425070	97.97%	7.437%
96	24.651	3661	3666	3675	rVB	709612	1456843	22.21%	1.686%
97	24.751	3678	3683	3690	rVV2	357368	878506	13.40%	1.017%
98	24.827	3692	3696	3709	rVB8	410313	923916	14.09%	1.069%
99	26.039	3896	3902	3913	rVB3	639310	1657756	25.28%	1.919%
100	26.545	3981	3988	4003	rVB3	550045	1722928	26.27%	1.994%

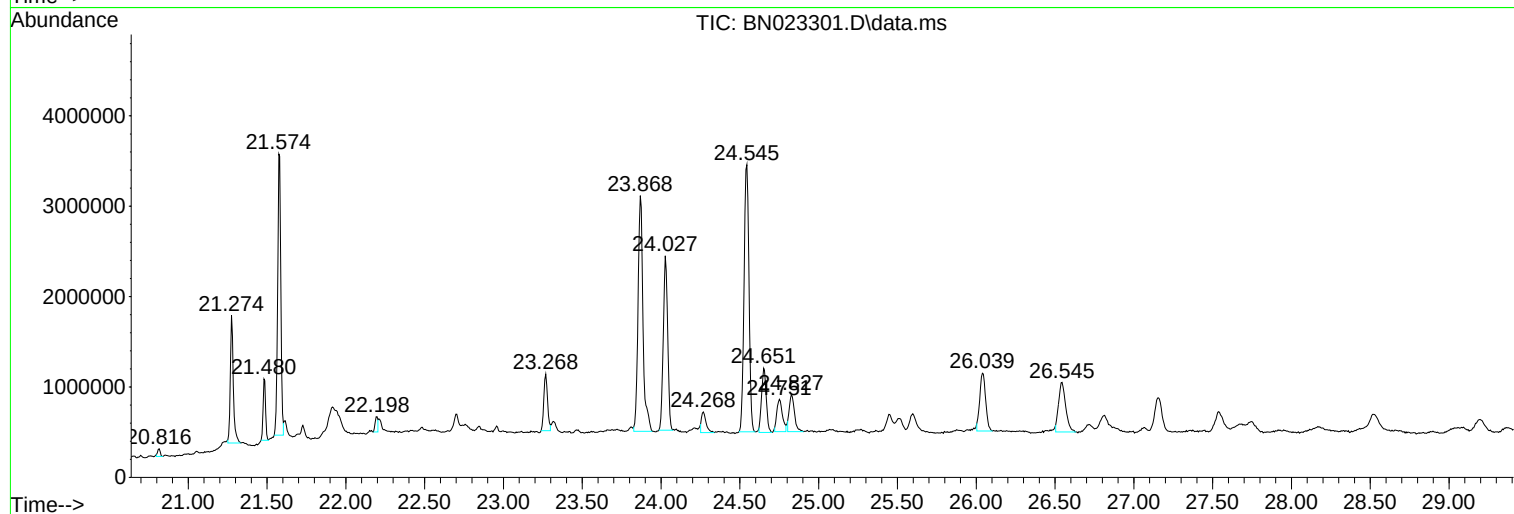
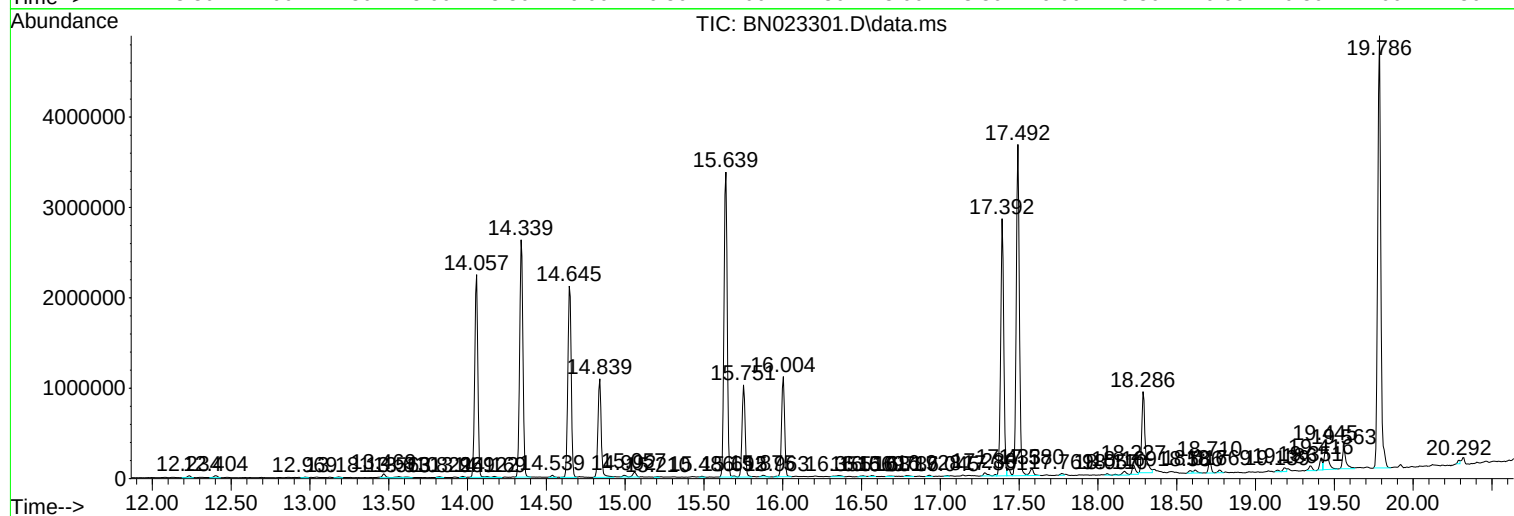
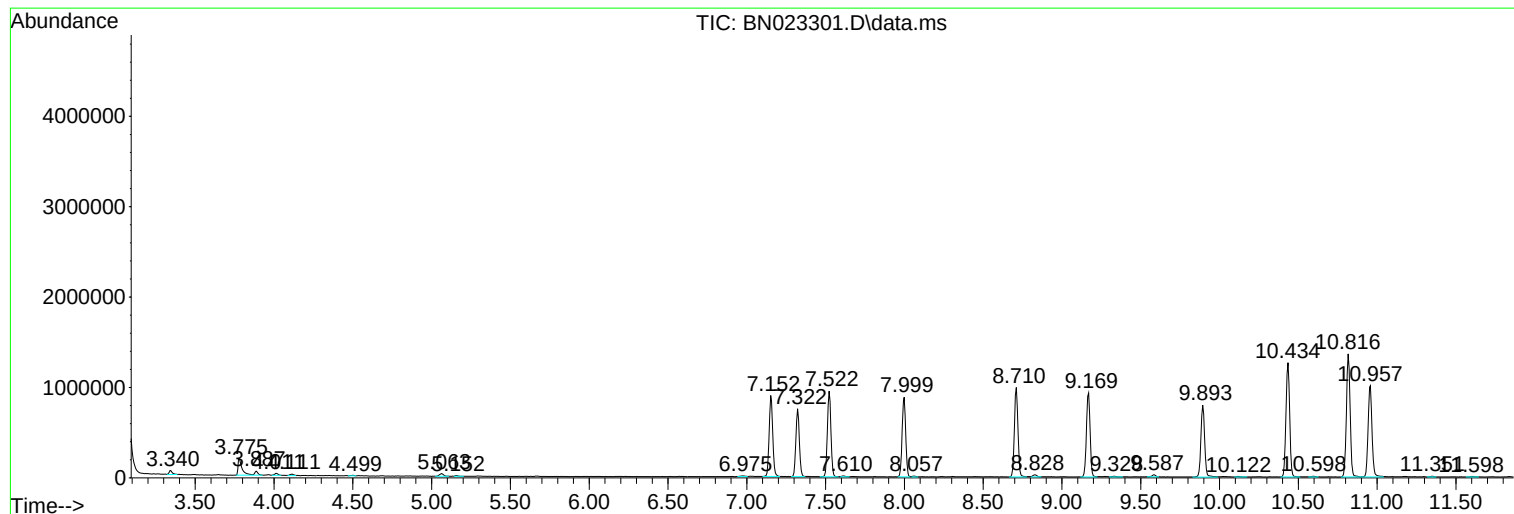
Sum of corrected areas: 86395179

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
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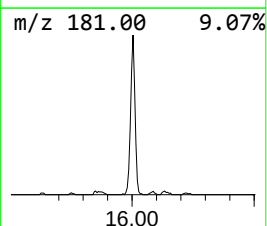
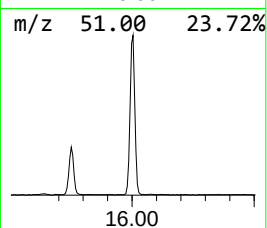
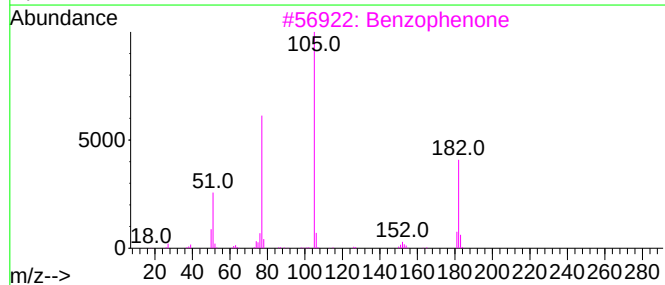
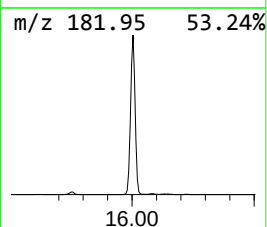
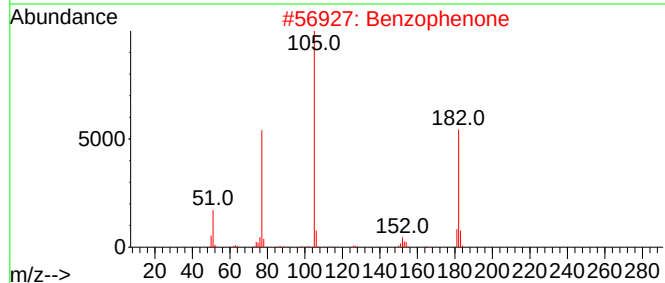
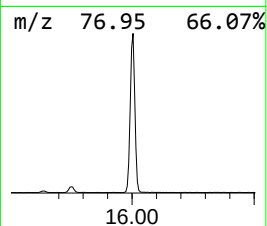
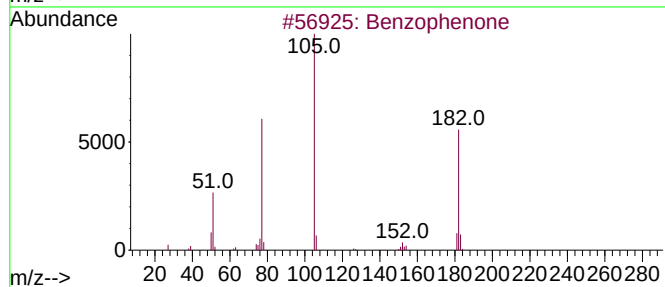
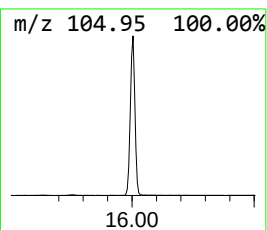
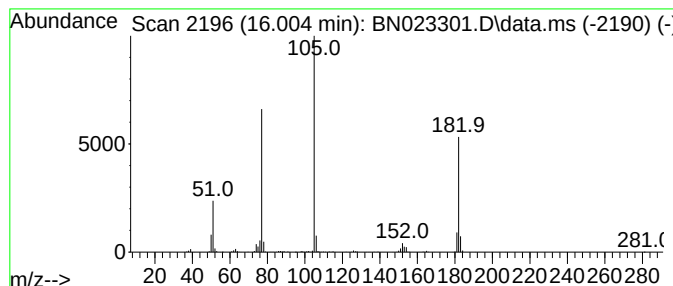
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.004	9.41 ng/ul	1456490	Acenaphthene-d10	14.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	93
5			Benzophenone	182	C13H10O	000119-61-9	91



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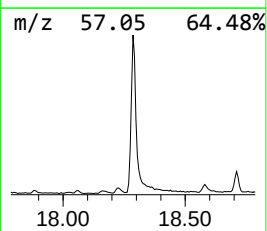
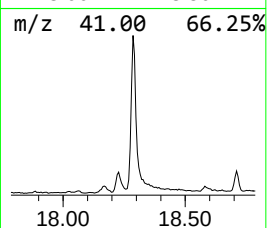
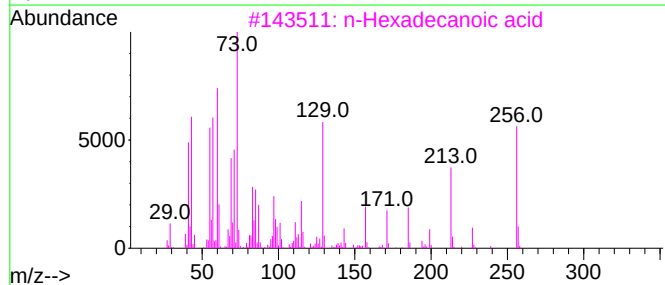
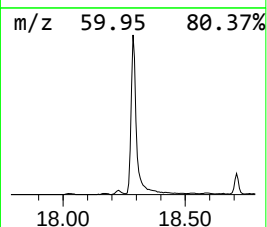
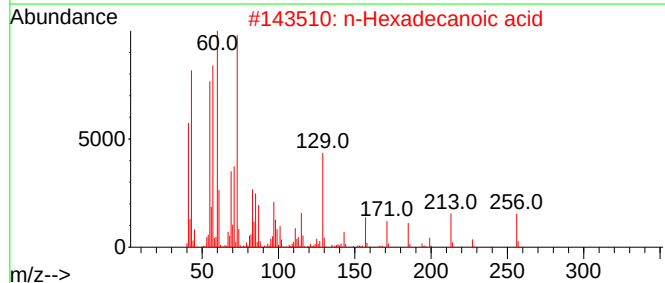
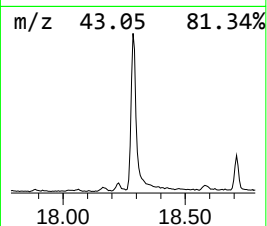
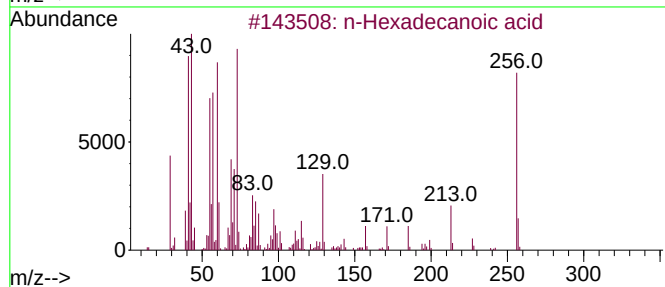
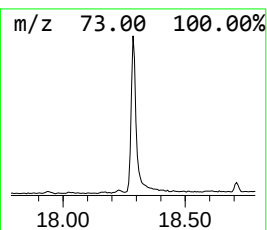
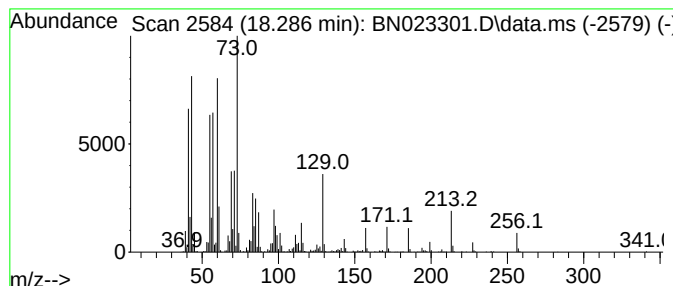
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.286	6.92 ng/ul	1335520	Phenanthrene-d10	17.392

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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	91



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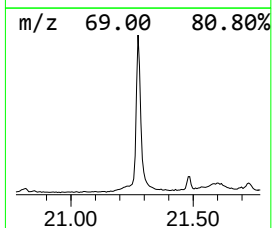
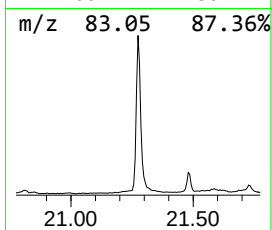
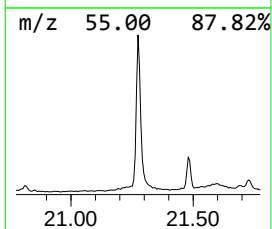
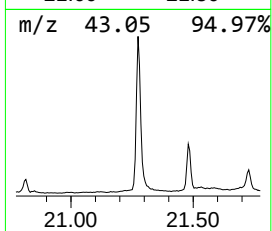
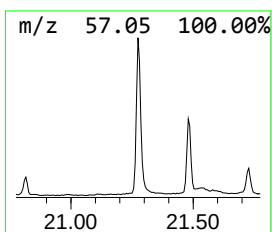
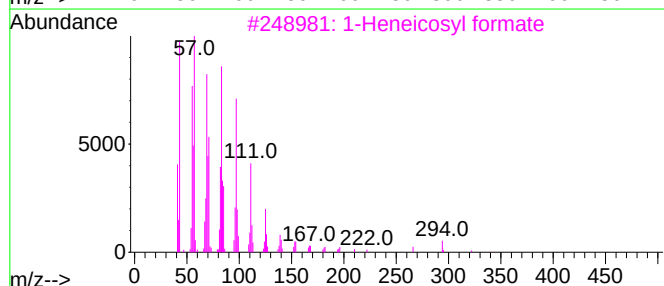
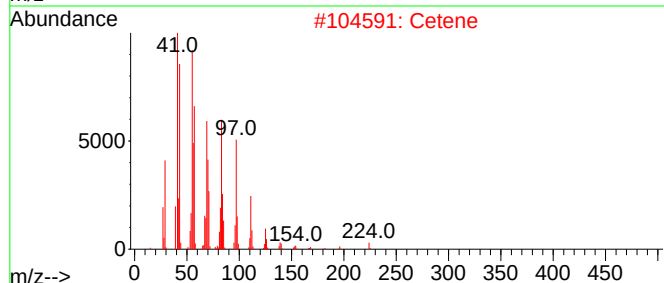
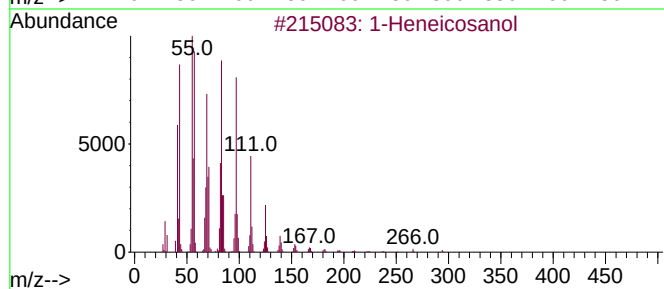
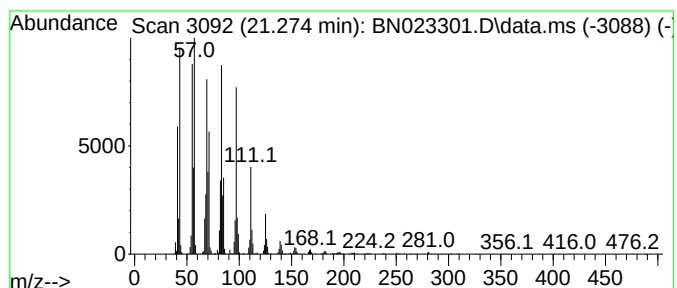
Instrument :
BNA_N
ClientSampleId :
DBXV2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Heneicosanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.274	8.46 ng/ul	1846720	Chrysene-d12	21.574	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	Cetene	224	C16H32	000629-73-2	94
3	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94



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Acq On : 20 Dec 2022 16:42
Operator : CG/JU
Sample : N6003-23
Misc :
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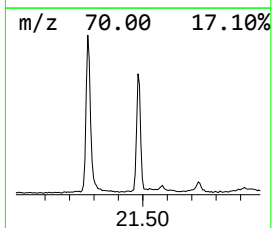
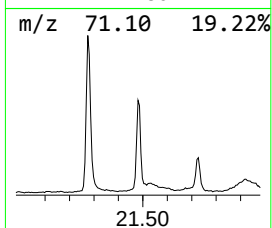
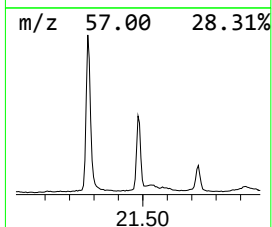
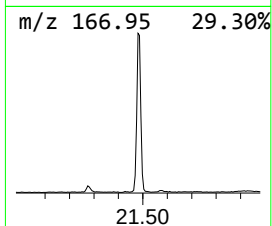
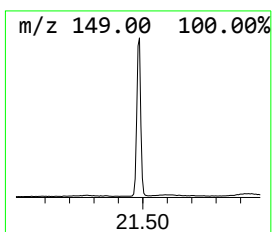
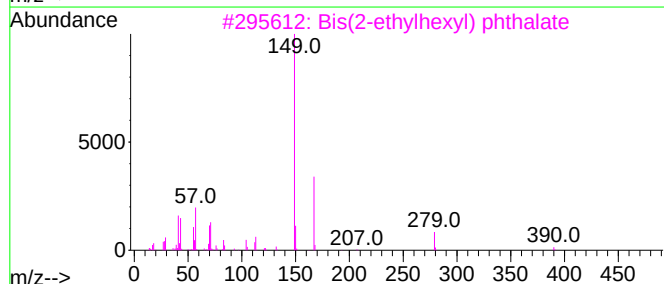
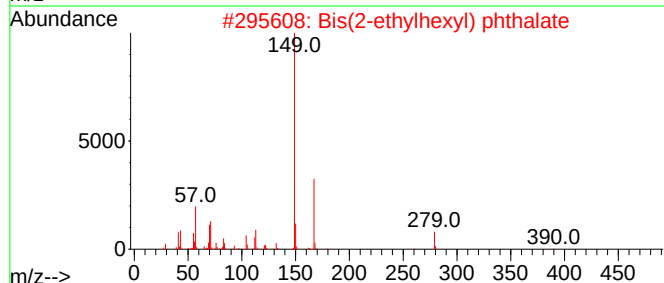
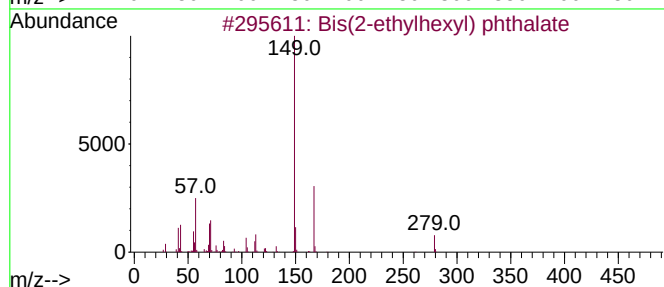
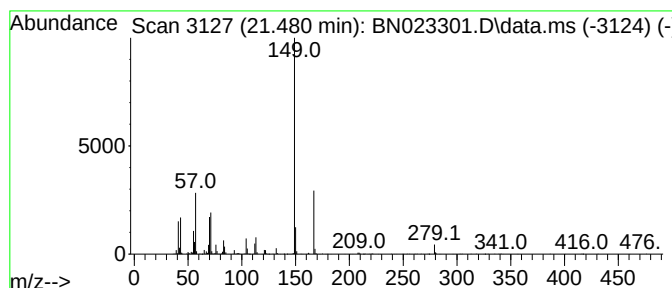
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Bis(2-ethylhexyl) phthalate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.480	3.45 ng/ul	752302	Chrysene-d12	21.574

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	91
2			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	91
3			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	91
4			Phthalic acid, 2-ethylhexyl isoh...	362	C22H34O4	1000308-98-5	80
5			Phthalic acid, cyclohexyl neopen...	318	C19H26O4	1000315-54-2	72



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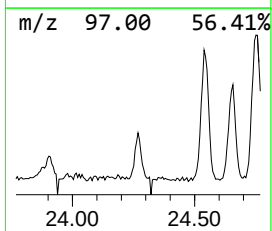
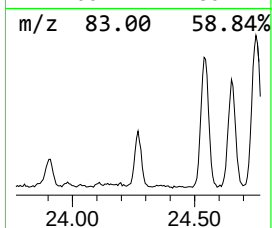
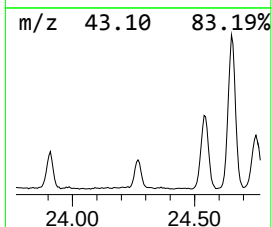
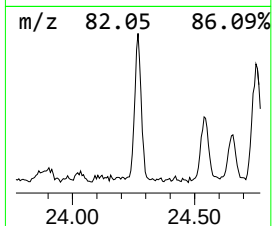
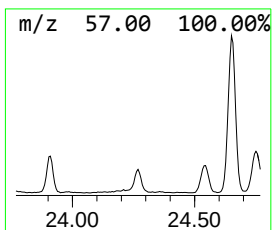
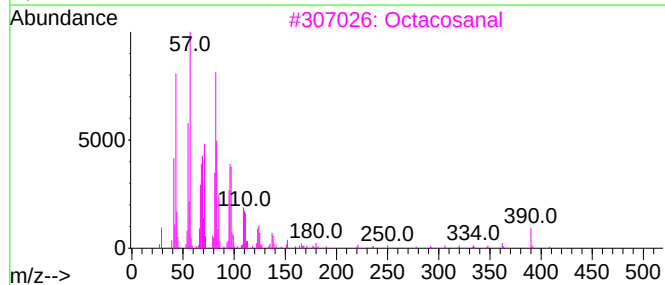
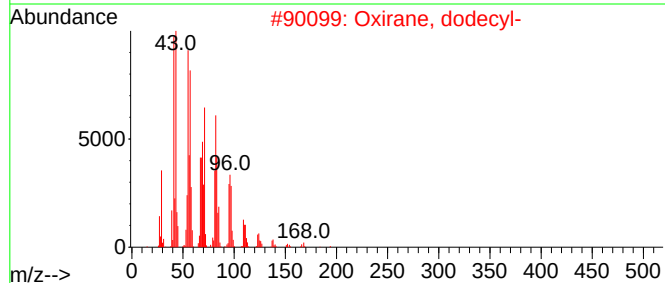
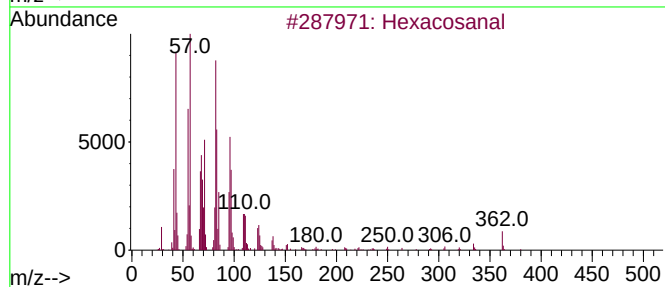
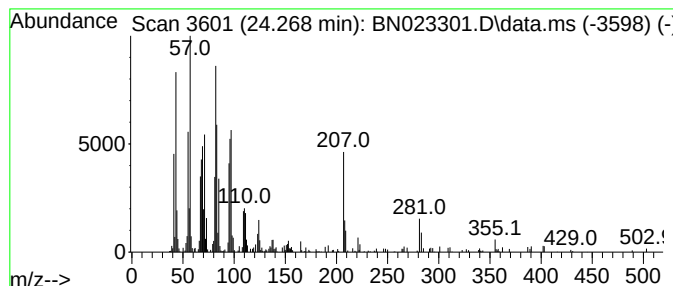
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Hexacosanal Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.268	2.43 ng/ul	454432	Perylene-d12	24.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosanal	380	C26H52O	026627-85-0	64
2			Oxirane, dodecyl-	212	C14H28O	003234-28-4	64
3			Octacosanal	408	C28H56O	022725-64-0	64
4			Octadecanal	268	C18H36O	000638-66-4	62
5			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	58



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Misc :
ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
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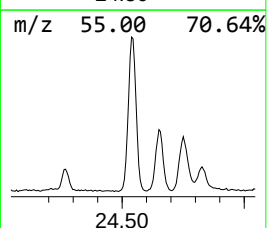
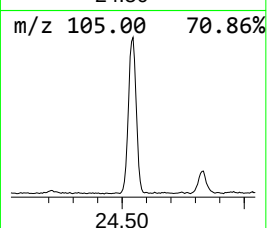
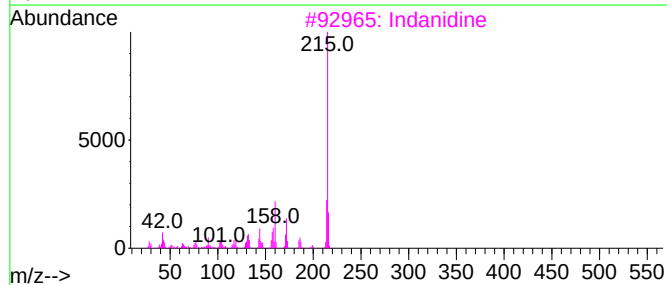
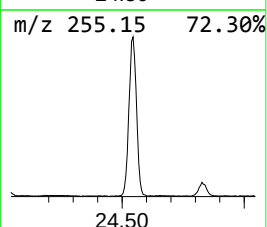
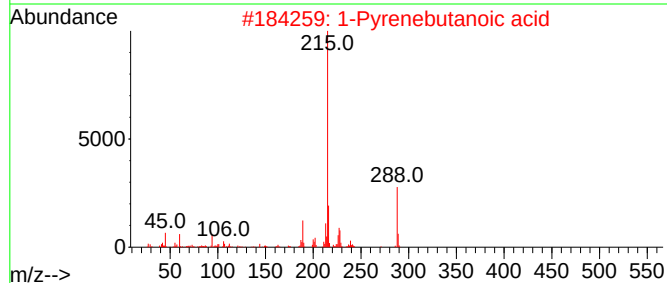
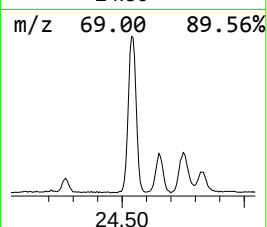
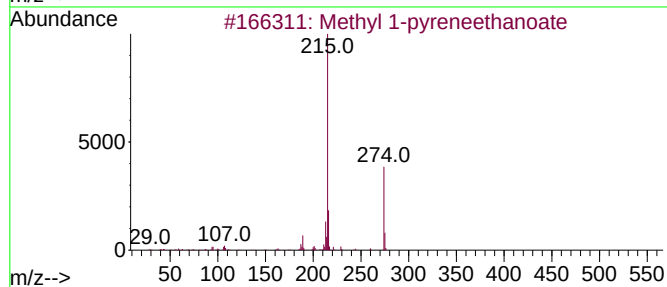
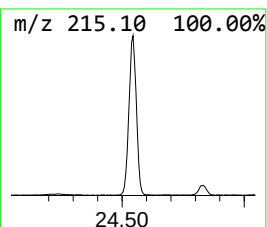
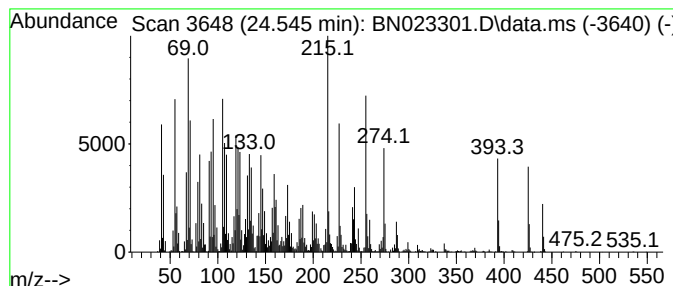
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.545	34.38 ng/ul	6425070	Perylene-d12	24.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	48
2			1-Pyrenebutanoic acid	288	C20H16O2	003443-45-6	25
3			Indanidine	215	C11H13N5	085392-79-6	15
4			4-(1H-Indol-3-yl)-1,3-thiazol-2-...	215	C11H9N3S	022258-56-6	15
5			Chlorflurecol methyl ester	274	C15H11ClO3	002536-31-4	14



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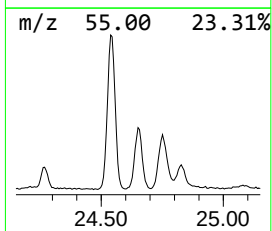
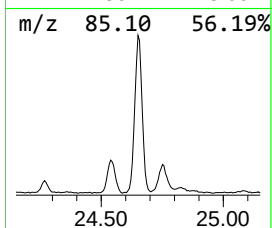
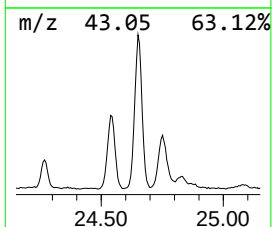
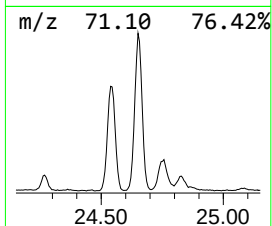
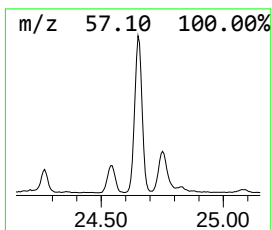
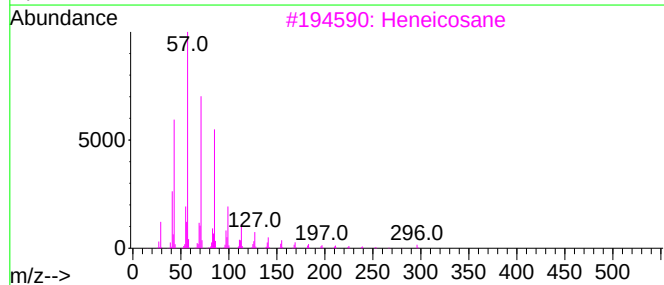
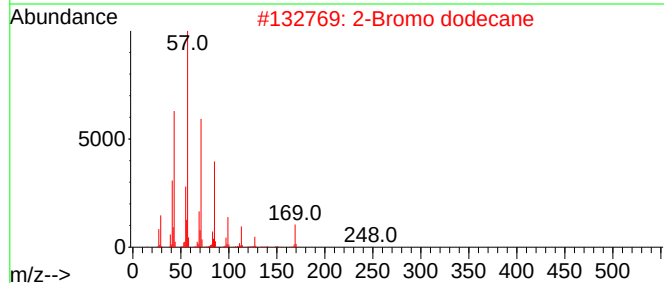
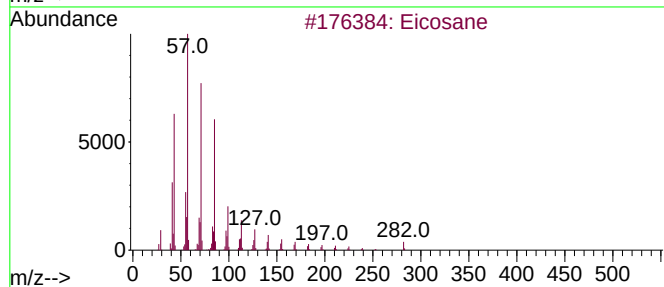
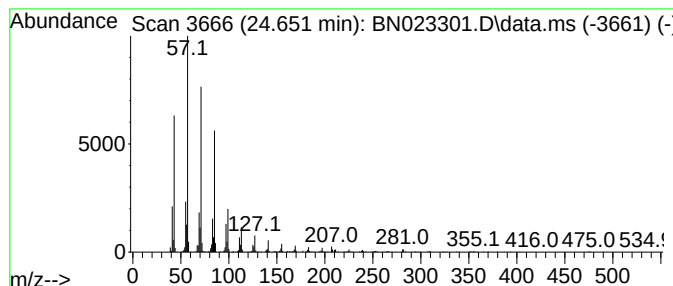
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.651	7.79 ng/ul	1456840	Perylene-d12	24.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	95
2			2-Bromo dodecane	248	C12H25Br	013187-99-0	93
3			Heneicosane	296	C21H44	000629-94-7	91
4			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91



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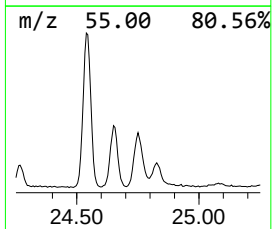
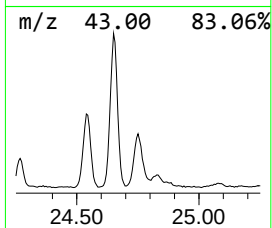
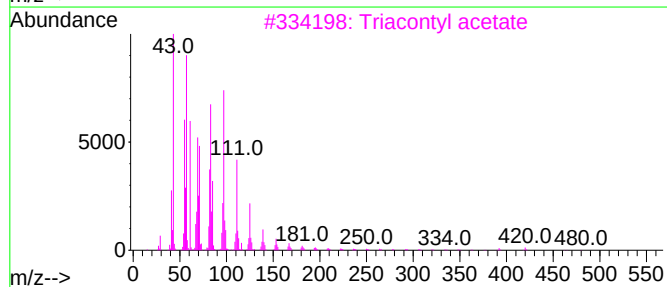
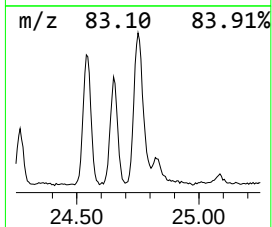
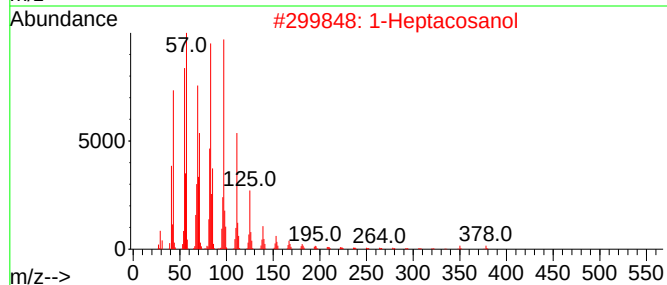
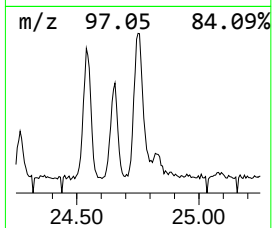
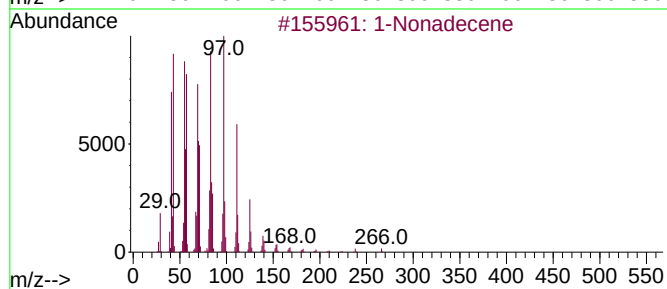
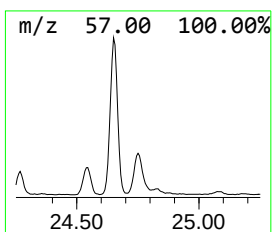
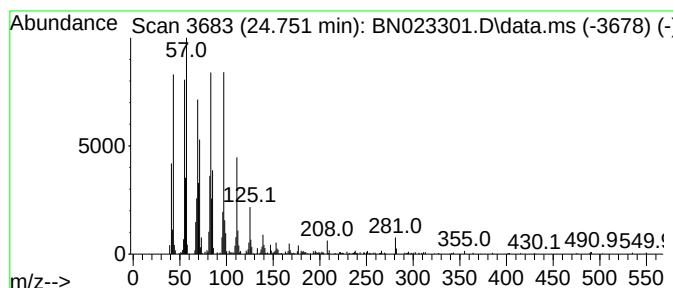
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.751	4.70 ng/ul	878506	Perylene-d12	24.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene			266	C19H38	018435-45-5	99
2	1-Heptacosanol			396	C27H56O	002004-39-9	95
3	Triacontyl acetate			480	C32H64O2	041755-58-2	95
4	Heneicosyl heptafluorobutyrate			508	C25H43F7O2	1000351-83-8	94
5	1-Triacontanol			438	C30H62O	000593-50-0	94



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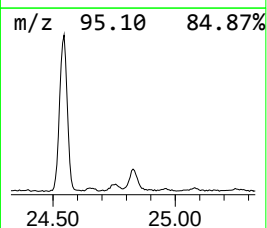
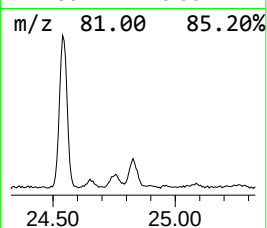
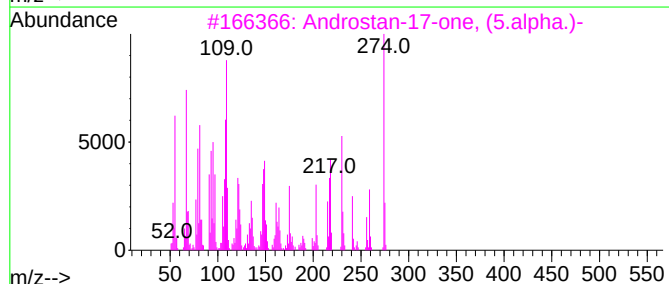
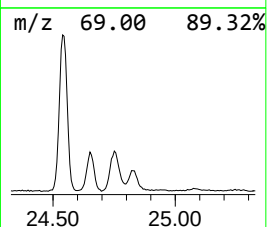
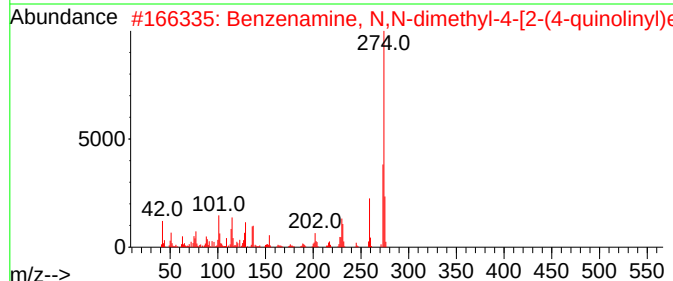
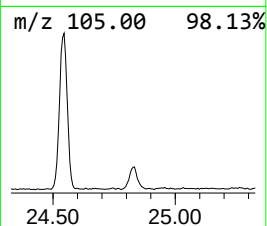
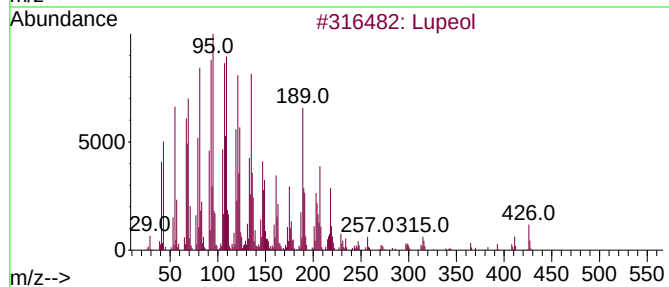
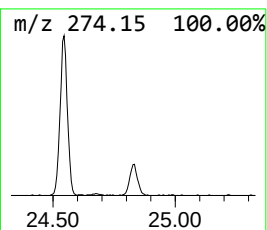
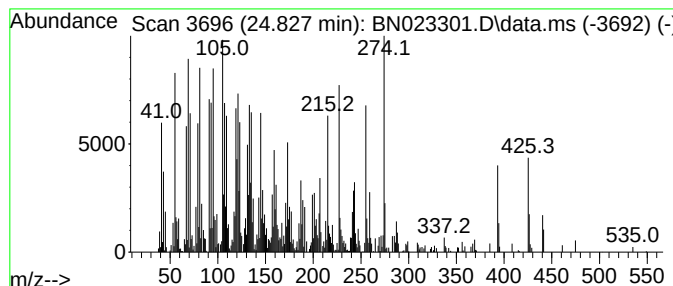
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Lupeol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.827	4.94 ng/ul	923916	Perylene-d12	24.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Lupeol	426	C30H50O	000545-47-1	53
2			Benzenamine, N,N-dimethyl-4-[2-(...	274	C19H18N2	000897-55-2	44
3			Androstan-17-one, (5.alpha.)-	274	C19H30O	000963-74-6	42
4			Androstan-17-one, (5.alpha.)-	274	C19H30O	000963-74-6	42
5			Lupeol	426	C30H50O	000545-47-1	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
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ALS Vial : 13 Sample Multiplier: 1

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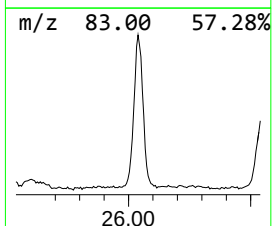
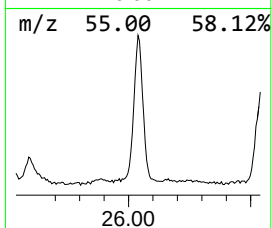
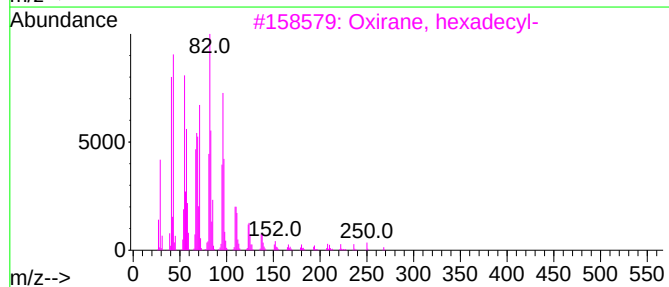
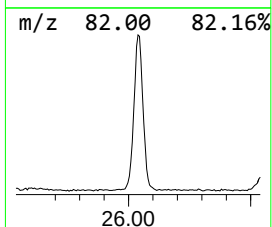
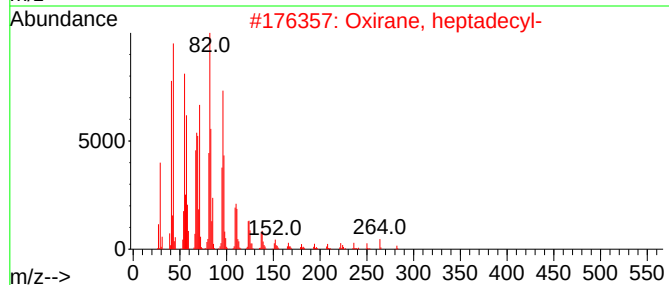
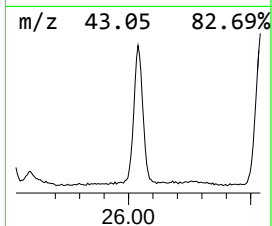
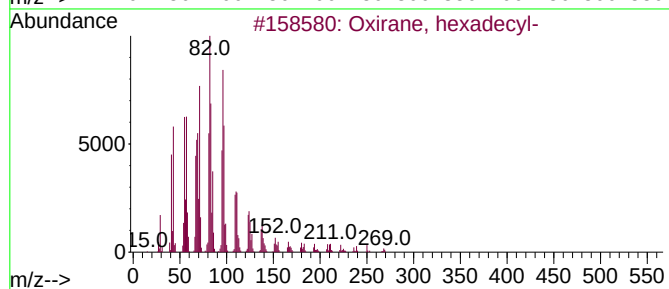
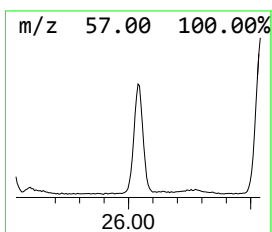
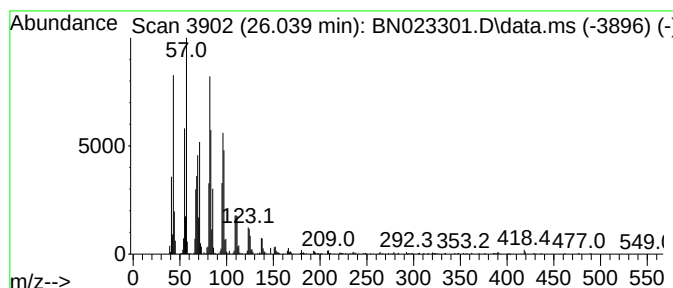
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Oxirane, hexadecyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.039	8.87 ng/ul	1657760	Perylene-d12	24.027

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	94
2		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	91
3		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
4		Hexacosanal	380	C26H52O	026627-85-0	90
5		Octadecanal	268	C18H36O	000638-66-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023301.D
Acq On : 20 Dec 2022 16:42
Operator : CG/JU
Sample : N6003-23
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBXV2

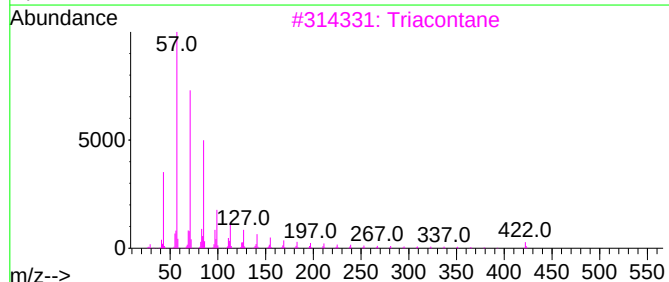
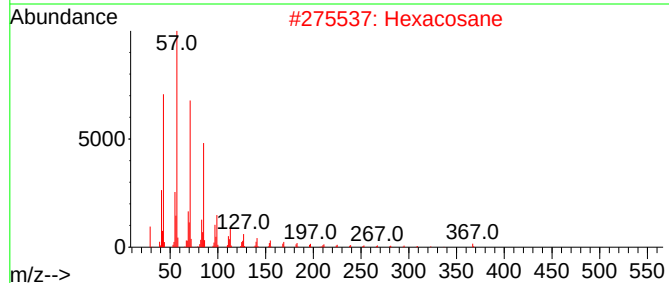
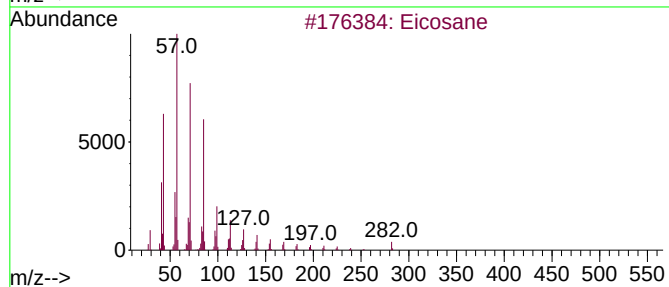
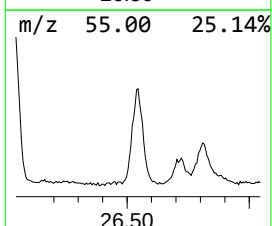
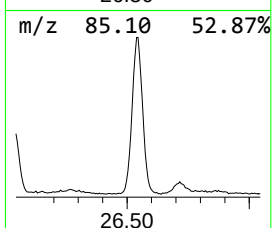
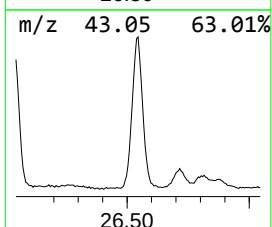
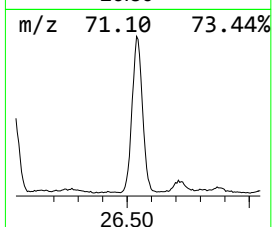
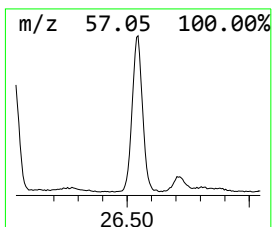
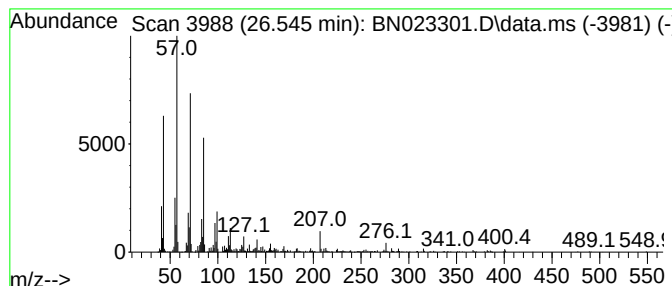
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.545	9.22 ng/ul	1722930	Perylene-d12	24.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	95
2			Hexacosane	366	C26H54	000630-01-3	90
3			Triacontane	422	C30H62	000638-68-6	90
4			Heneicosane	296	C21H44	000629-94-7	90
5			Octacosane, 2-methyl-	408	C29H60	001560-98-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
 Data File : BN023301.D
 Acq On : 20 Dec 2022 16:42
 Operator : CG/JU
 Sample : N6003-23
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBXV2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	16.004	9.4	ng/ul	1456490	3	14.645	3094190	20.0
n-Hexadecanoic ...	18.286	6.9	ng/ul	1335520	4	17.392	3858460	20.0
1-Heneicosanol	21.274	8.5	ng/ul	1846720	5	21.574	4364780	20.0
Bis(2-ethylhexy...	21.480	3.5	ng/ul	752302	5	21.574	4364780	20.0
Hexacosanal	24.268	2.4	ng/ul	454432	6	24.027	3737940	20.0
unknown-01	24.545	34.4	ng/ul	6425070	6	24.027	3737940	20.0
(DEL) Alkane: S...	24.651	7.8	ng/ul	1456840	6	24.027	3737940	20.0
(DEL) Alkane: S...	24.751	4.7	ng/ul	878506	6	24.027	3737940	20.0
Lupeol	24.827	4.9	ng/ul	923916	6	24.027	3737940	20.0
Oxirane, hexade...	26.039	8.9	ng/ul	1657760	6	24.027	3737940	20.0
(DEL) Alkane: S...	26.545	9.2	ng/ul	1722930	6	24.027	3737940	20.0