

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN011922\
 Data File : BN018418.D
 Acq On : 19 Jan 2022 15:40
 Operator : CG/JU
 Sample : N1129-08
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBM01

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-BN010622.M
 Title : SVOA CALIBRATION

Signal : TIC: BN018418.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.058	9	12	18	rBV	53907	72831	1.01%	0.066%
2	3.470	80	82	93	rBV	319407	530052	7.37%	0.483%
3	3.711	118	123	131	rVB	71696	119854	1.67%	0.109%
4	3.805	135	139	145	rVB	26917	39322	0.55%	0.036%
5	3.893	151	154	161	rVB6	21572	40580	0.56%	0.037%
6	4.034	173	178	185	rBV	57112	82639	1.15%	0.075%
7	4.446	244	248	252	rBV2	34174	46726	0.65%	0.043%
8	4.711	287	293	301	rBV	111778	185175	2.58%	0.169%
9	5.593	438	443	449	rBV	23330	39501	0.55%	0.036%
10	6.287	554	561	571	rVB	610677	913136	12.70%	0.832%
11	6.587	607	612	618	rVB2	55067	89394	1.24%	0.081%
12	6.775	637	644	653	rBV	1108572	1845131	25.66%	1.682%
13	6.928	663	670	681	rBV	951706	1484878	20.65%	1.354%
14	7.034	683	688	692	rVV3	42483	69119	0.96%	0.063%
15	7.122	697	703	717	rVV	1197912	1864840	25.94%	1.700%
16	7.587	776	782	792	rBV	1621360	2503498	34.82%	2.282%
17	8.305	898	904	917	rBV	1157325	1905673	26.50%	1.737%
18	8.411	917	922	928	rVB3	27545	52088	0.72%	0.047%
19	8.734	970	977	988	rBV	1122017	1841158	25.61%	1.678%
20	8.922	1005	1009	1017	rBV7	16534	33387	0.46%	0.030%
21	9.211	1052	1058	1065	rVB2	38601	64383	0.90%	0.059%
22	9.458	1089	1100	1108	rBV	965497	1589977	22.11%	1.449%
23	9.993	1185	1191	1206	rBV	1343645	2380016	33.10%	2.170%
24	10.363	1246	1254	1264	rBV	2276997	3786818	52.67%	3.452%
25	10.505	1271	1278	1295	rBV	1052257	1969763	27.40%	1.796%
26	10.863	1334	1339	1344	rBV2	38765	64747	0.90%	0.059%
27	11.969	1522	1527	1537	rBV3	25536	45987	0.64%	0.042%
28	12.604	1627	1635	1640	rBV7	37385	80846	1.12%	0.074%
29	12.981	1690	1699	1703	rBV8	27779	56754	0.79%	0.052%
30	13.081	1711	1716	1719	rBV	61383	79700	1.11%	0.073%
31	13.122	1719	1723	1729	rVB	206193	279820	3.89%	0.255%
32	13.563	1792	1798	1804	rVB	2056834	2795348	38.88%	2.548%
33	13.651	1807	1813	1821	rBV	2867845	4028875	56.03%	3.673%
34	13.922	1851	1859	1870	rVV	3175576	4830843	67.19%	4.404%
35	14.093	1882	1888	1892	rVB2	44742	60595	0.84%	0.055%
36	14.163	1896	1900	1902	rVV2	41211	53962	0.75%	0.049%

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN010622.M
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37	14.234	1902	1912	1919	rVV	3514542	6071709	84.45%	5.535%
38	14.322	1921	1927	1938	rVB4	148025	293842	4.09%	0.268%
39	14.451	1942	1949	1955	rBV	544800	1009036	14.03%	0.920%
40	14.522	1955	1961	1966	rVV3	222763	523529	7.28%	0.477%
41	14.581	1966	1971	1975	rVB5	38163	78023	1.09%	0.071%
42	14.669	1981	1986	1992	rBV7	40560	62697	0.87%	0.057%
43	14.875	2013	2021	2027	rVB5	28737	58825	0.82%	0.054%
44	15.063	2048	2053	2058	rVV2	30787	50084	0.70%	0.046%
45	15.116	2058	2062	2066	rVV4	32011	57594	0.80%	0.053%
46	15.228	2074	2081	2095	rBV	4010213	5621607	78.19%	5.125%
47	15.351	2095	2102	2114	rVV	1065681	1516287	21.09%	1.382%
48	15.575	2135	2140	2145	rBV4	71795	123881	1.72%	0.113%
49	15.669	2151	2156	2159	rBV6	19493	33853	0.47%	0.031%
50	15.722	2159	2165	2173	rVB2	251993	375312	5.22%	0.342%
51	15.810	2176	2180	2186	rBV8	20913	37946	0.53%	0.035%
52	15.928	2195	2200	2205	rBV8	27153	60453	0.84%	0.055%
53	16.145	2232	2237	2241	rVV6	23560	39581	0.55%	0.036%
54	16.728	2330	2336	2339	rBV6	27792	46636	0.65%	0.043%
55	16.769	2339	2343	2348	rVB3	34528	47919	0.67%	0.044%
56	16.922	2363	2369	2372	rBV5	44884	89302	1.24%	0.081%
57	16.975	2372	2378	2389	rVV	4213198	6130481	85.26%	5.589%
58	17.075	2389	2395	2408	rVV2	4147677	5971649	83.05%	5.444%
59	17.187	2409	2414	2420	rVB7	48654	83439	1.16%	0.076%
60	17.281	2426	2430	2436	rBV7	36660	63277	0.88%	0.058%
61	17.675	2493	2497	2502	rVB2	59672	78418	1.09%	0.071%
62	17.787	2510	2516	2521	rBV7	37248	72086	1.00%	0.066%
63	17.839	2521	2525	2531	rBV6	52653	118943	1.65%	0.108%
64	17.898	2531	2535	2543	rBV2	303400	516542	7.18%	0.471%
65	17.998	2549	2552	2558	rVB4	57667	81857	1.14%	0.075%
66	18.198	2584	2586	2591	rVV5	21971	38071	0.53%	0.035%
67	18.663	2662	2665	2671	rVV4	30797	42118	0.59%	0.038%
68	18.939	2707	2712	2717	rBV2	65226	89350	1.24%	0.081%
69	19.010	2717	2724	2726	rBV4	61535	109956	1.53%	0.100%
70	19.039	2726	2729	2733	rBV2	63251	76834	1.07%	0.070%
71	19.186	2750	2754	2761	rBV2	111114	196812	2.74%	0.179%
72	19.263	2765	2767	2770	rVV4	28961	36276	0.50%	0.033%
73	19.304	2772	2774	2777	rVV4	25189	34860	0.48%	0.032%
74	19.375	2780	2786	2796	rVV	5269259	7190104	100.00%	6.555%
75	19.451	2796	2799	2804	rVB3	44058	56674	0.79%	0.052%
76	19.786	2850	2856	2861	rBV	453780	587196	8.17%	0.535%

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77	19.833	2861	2864	2869	rBV4	40033	56878	0.79%	0.052%
78	19.933	2878	2881	2883	rVV	161720	182342	2.54%	0.166%
79	19.957	2883	2885	2890	rVB2	78369	89877	1.25%	0.082%
80	20.345	2945	2951	2953	rBV5	69587	136379	1.90%	0.124%
81	20.375	2953	2956	2960	rVV	189125	230140	3.20%	0.210%
82	20.451	2964	2969	2972	rVB2	85939	113318	1.58%	0.103%
83	20.528	2979	2982	2986	rVB5	40862	50432	0.70%	0.046%
84	20.581	2987	2991	2997	rBV	296278	324379	4.51%	0.296%
85	20.733	3012	3017	3030	rBV	6074524	6999591	97.35%	6.381%
86	20.916	3043	3048	3055	rBV2	306289	536259	7.46%	0.489%
87	21.175	3086	3092	3102	rBV	5082890	6587371	91.62%	6.005%
88	21.351	3119	3122	3126	rBV	167670	183774	2.56%	0.168%
89	21.498	3144	3147	3151	rBV2	112846	119970	1.67%	0.109%
90	21.786	3191	3196	3204	rBV	287314	493747	6.87%	0.450%
91	22.245	3269	3274	3280	rBV	335976	511240	7.11%	0.466%
92	22.369	3292	3295	3301	rVB	104286	133514	1.86%	0.122%
93	22.751	3354	3360	3378	rVB	1012318	1689207	23.49%	1.540%
94	23.251	3438	3445	3452	rBV2	3647280	6632502	92.24%	6.047%
95	23.322	3452	3457	3462	rVV	422118	773000	10.75%	0.705%
96	23.392	3462	3469	3487	rVB2	3312516	6534985	90.89%	5.958%
97	23.969	3561	3567	3577	rBV	952706	1943391	27.03%	1.772%
98	24.721	3689	3695	3706	rVB	281953	659048	9.17%	0.601%
99	25.610	3839	3846	3855	rBV2	295692	774137	10.77%	0.706%
100	26.480	3989	3994	4002	rVB10	93884	235504	3.28%	0.215%

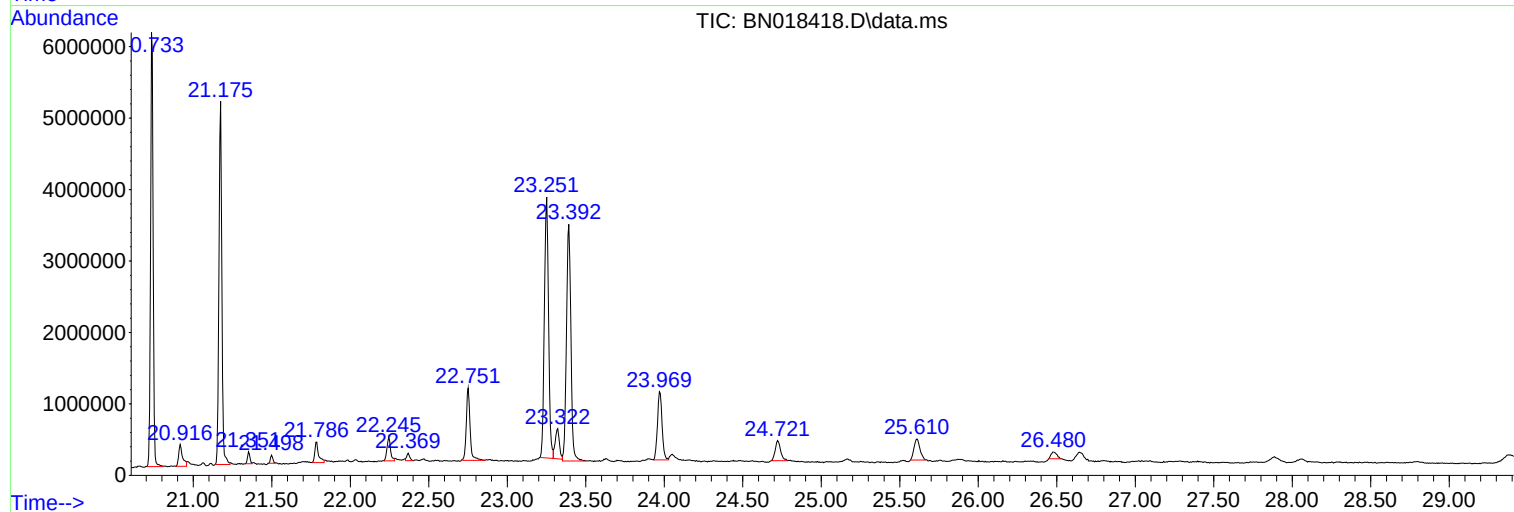
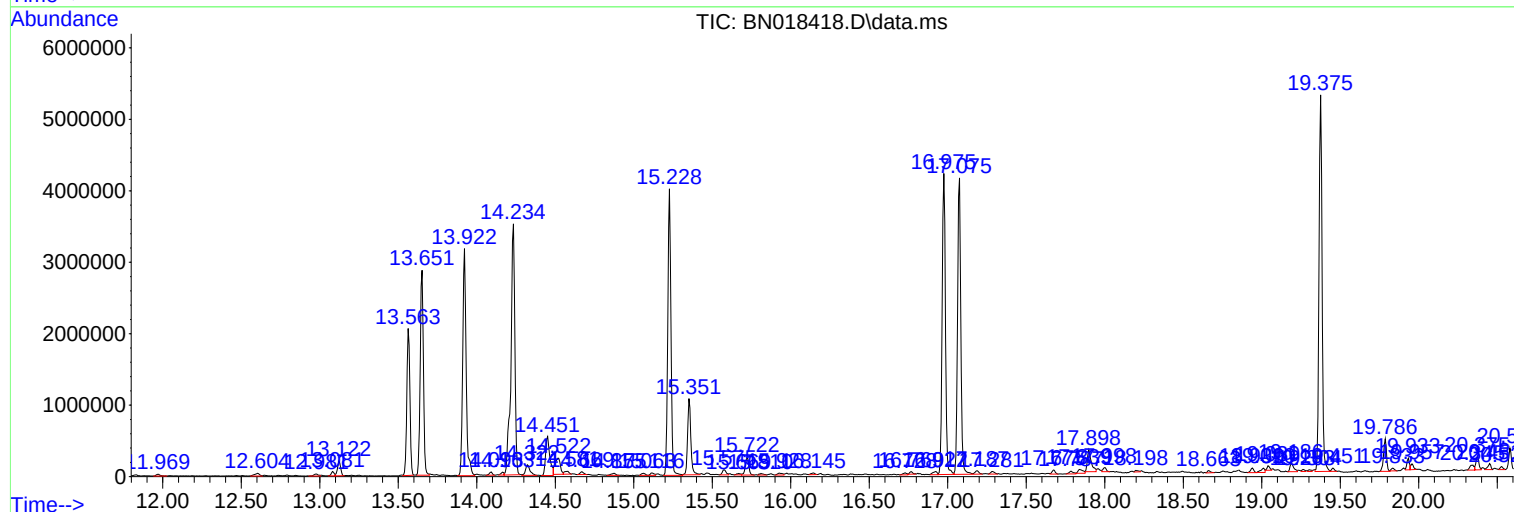
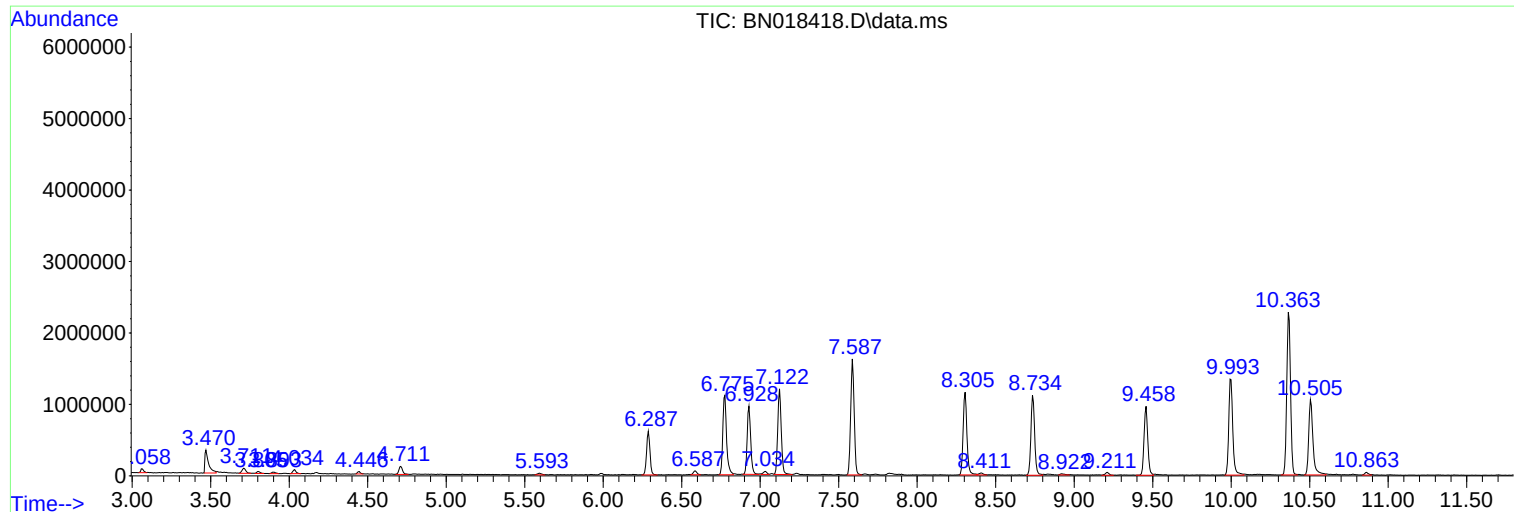
Sum of corrected areas: 109691460

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN010622.M
Quant Title : SVOA CALIBRATION

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



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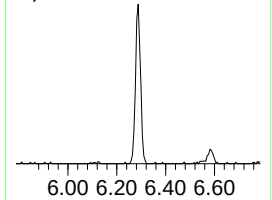
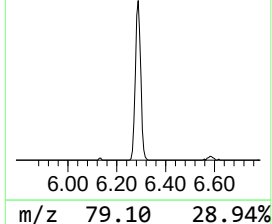
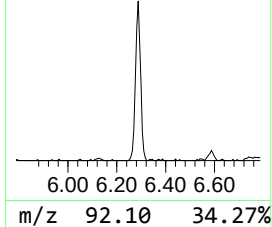
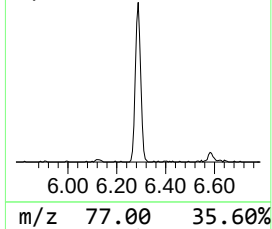
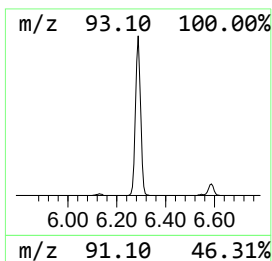
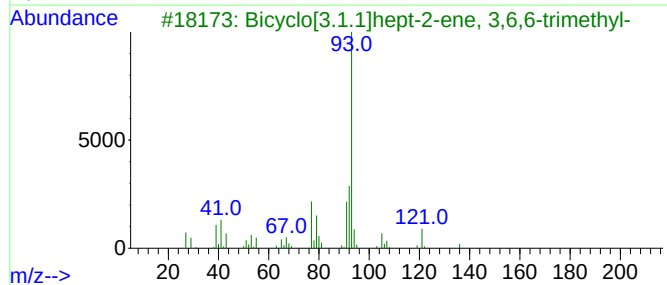
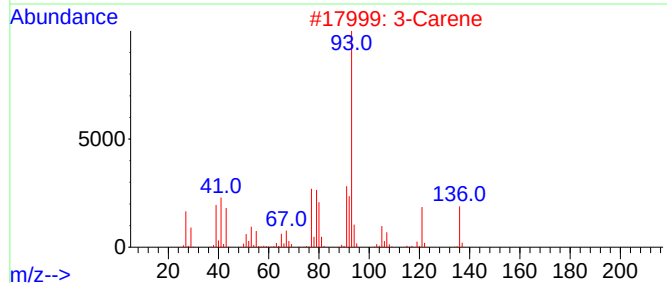
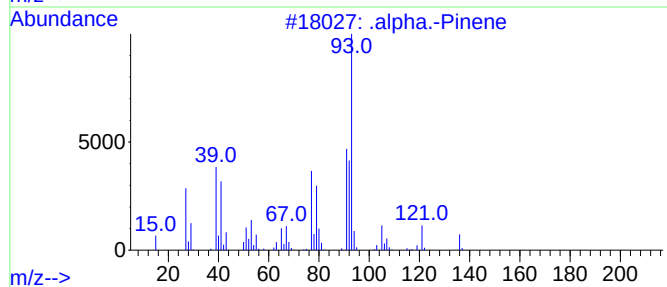
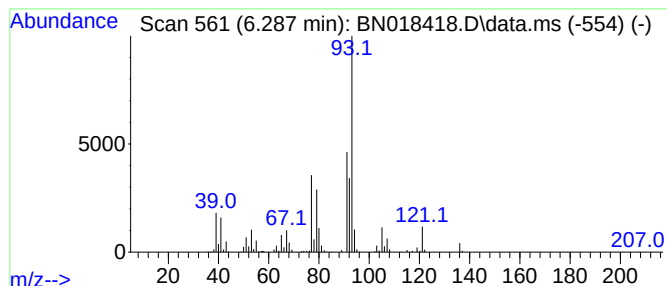
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN010622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 .alpha.-Pinene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.287	7.29 ng/ul	913136	1,4-Dichlorobenzene-d4	7.587

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.alpha.-Pinene	136	C10H16	000080-56-8	95
2			3-Carene	136	C10H16	013466-78-9	91
3			Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	91
4			Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	91
5			3-Carene	136	C10H16	013466-78-9	91



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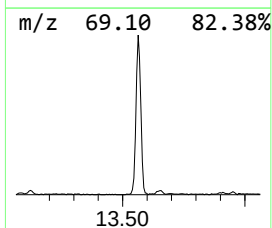
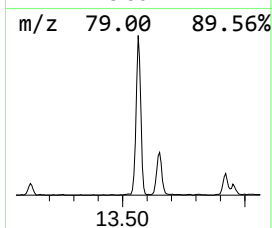
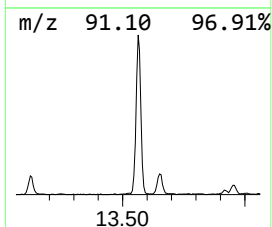
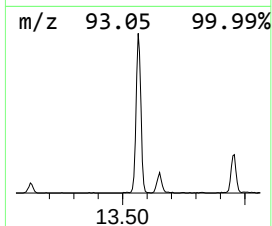
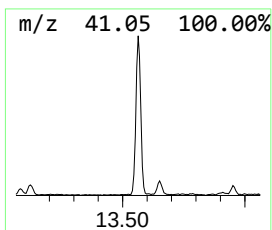
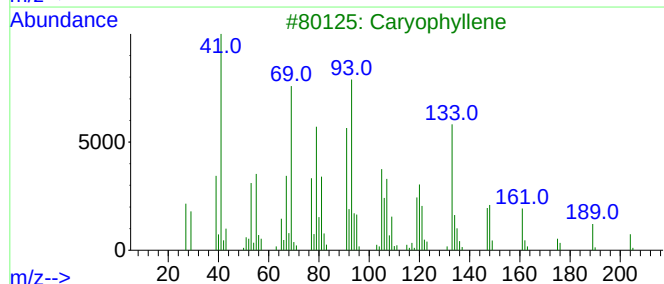
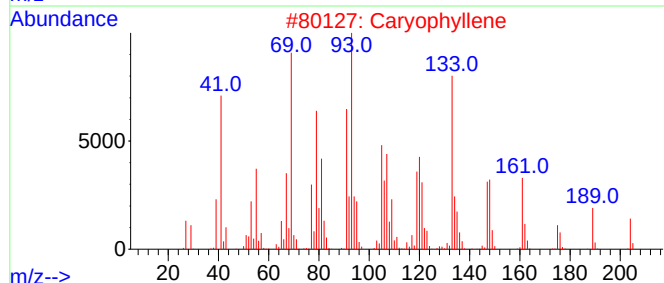
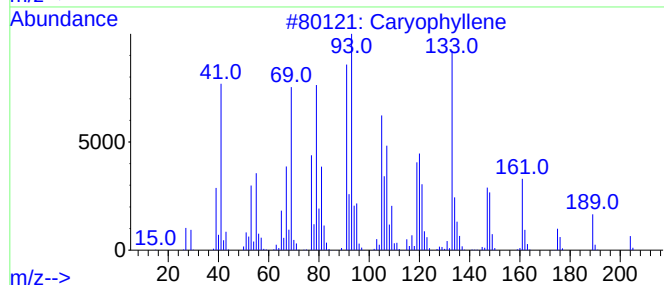
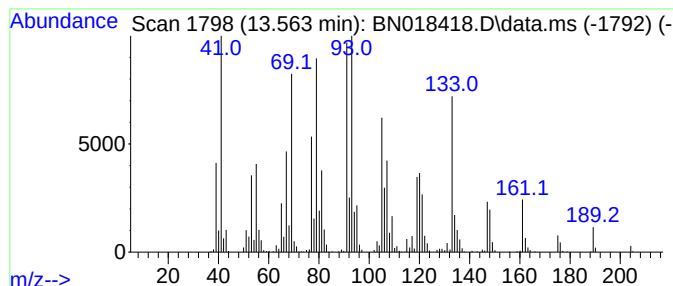
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN010622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Caryophyllene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.563	9.21 ng/ul	2795350	Acenaphthene-d10	14.234

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Caryophyllene	204	C15H24	000087-44-5	99
2			Caryophyllene	204	C15H24	000087-44-5	99
3			Caryophyllene	204	C15H24	000087-44-5	98
4			Bicyclo[7.2.0]undec-4-ene, 4,11,...	204	C15H24	000118-65-0	91
5			Caryophyllene	204	C15H24	000087-44-5	90



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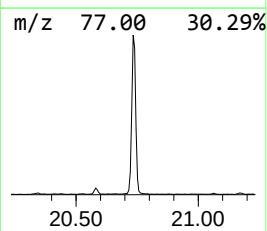
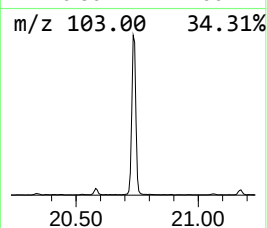
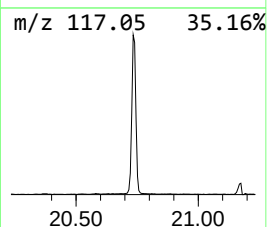
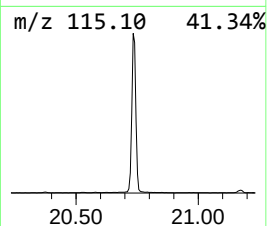
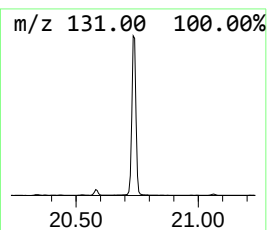
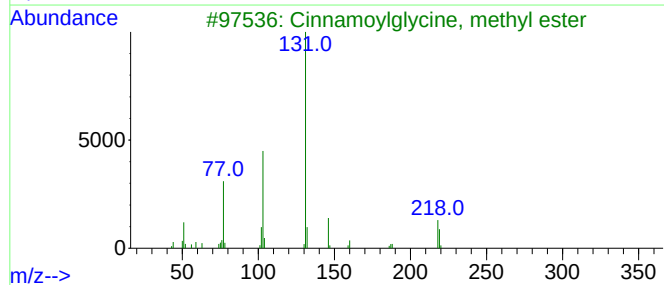
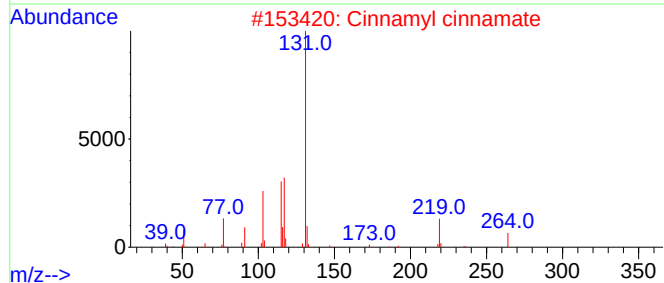
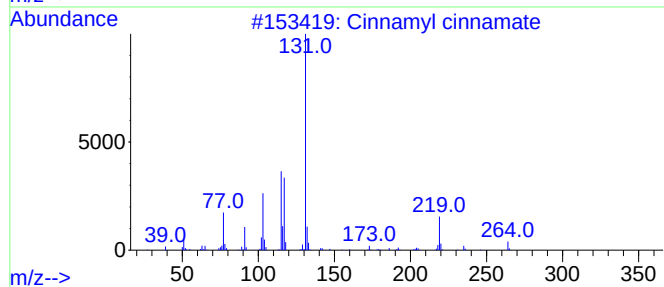
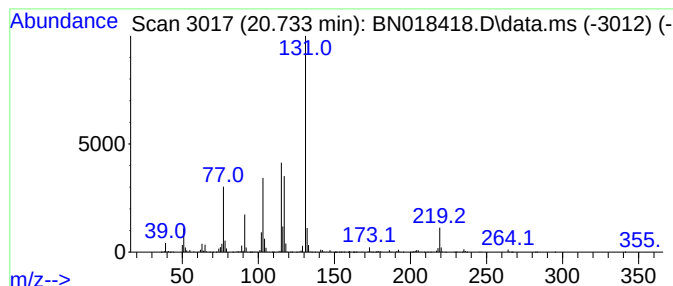
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN010622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Cinnamyl cinnamate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.733	21.25 ng/ul	6999590	Chrysene-d12	21.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cinnamyl cinnamate	264	C18H16O2	000122-69-0	91
2			Cinnamyl cinnamate	264	C18H16O2	000122-69-0	83
3			Cinnamoylglycine, methyl ester	219	C12H13NO3	040778-04-9	64
4			N-(p-Vinylbenzoyl)-l-alanine	219	C12H13NO3	139184-60-4	50
5			(E)-4-Bromo-4-methyl-1-phenylpen...	252	C12H13BrO	1000443-02-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN011922\
Data File : BN018418.D
Acq On : 19 Jan 2022 15:40
Operator : CG/JU
Sample : N1129-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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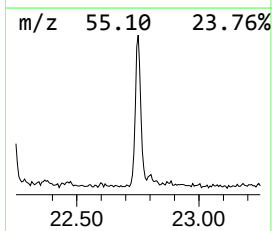
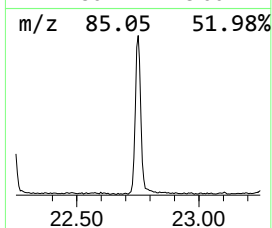
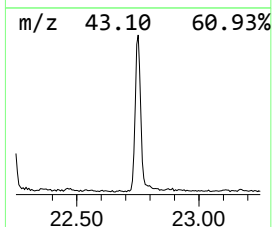
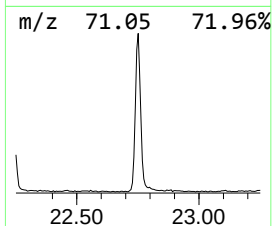
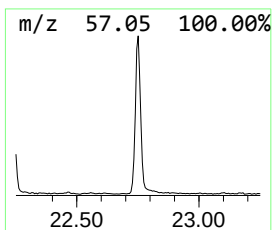
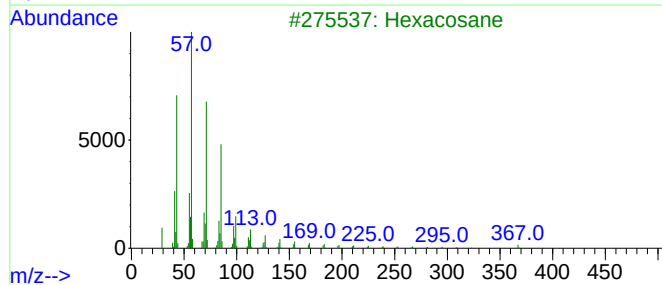
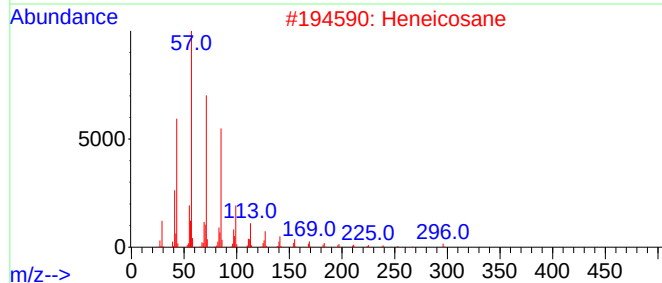
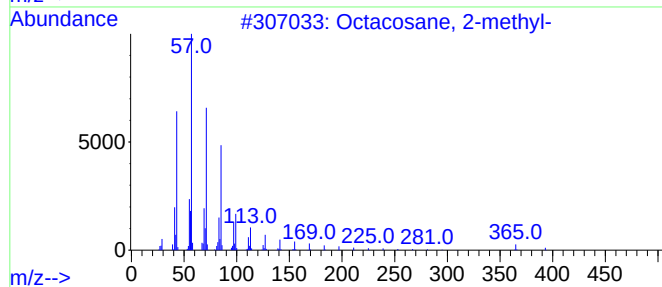
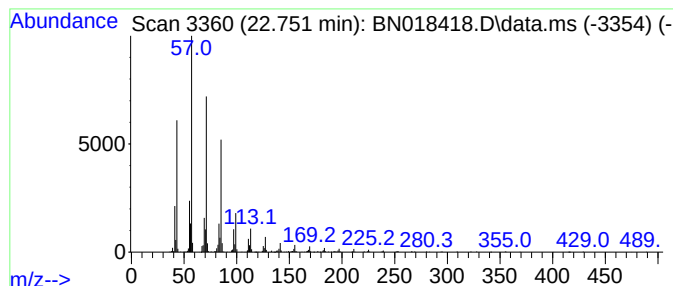
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.751	5.17 ng/ul	1689210	Perylene-d12	23.392

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
5		Tetracosane, 1-iodo-	464	C24H49I	1000406-32-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN011922\
Data File : BN018418.D
Acq On : 19 Jan 2022 15:40
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Sample : N1129-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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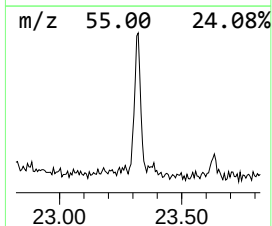
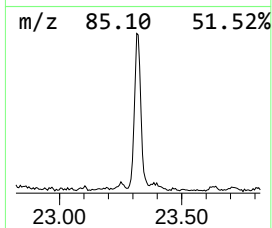
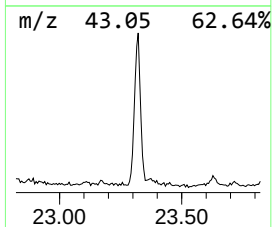
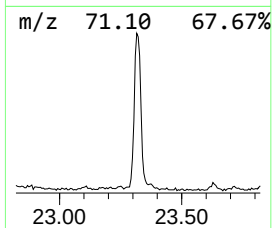
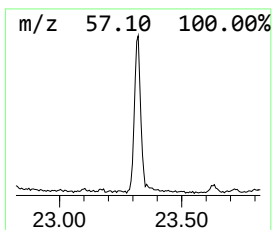
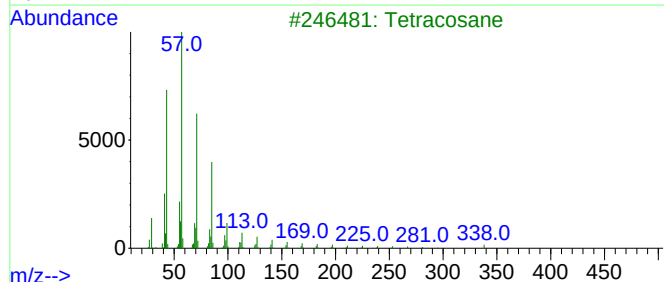
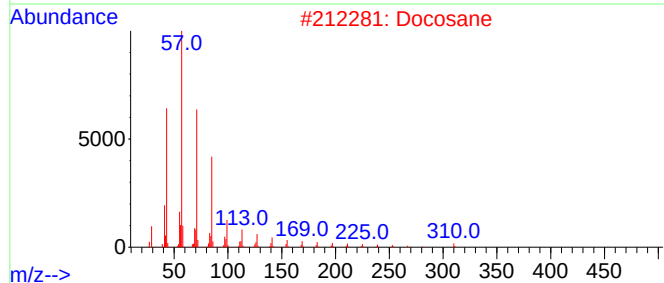
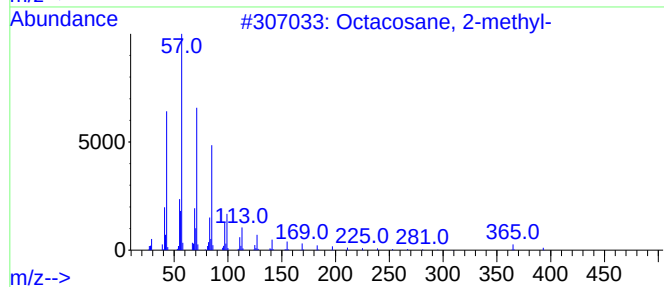
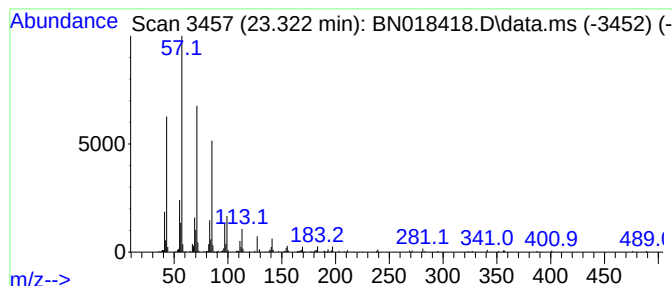
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.322	2.37 ng/ul	773000	Perylene-d12	23.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane, 2-methyl-	408	C29H60	001560-98-1	90
2			Docosane	310	C22H46	000629-97-0	87
3			Tetracosane	338	C24H50	000646-31-1	87
4			Hexatriacontane	507	C36H74	000630-06-8	87
5			Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN011922\
Data File : BN018418.D
Acq On : 19 Jan 2022 15:40
Operator : CG/JU
Sample : N1129-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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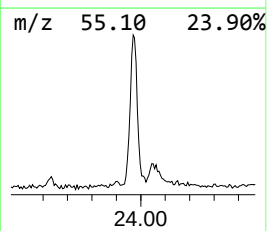
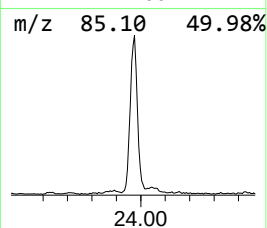
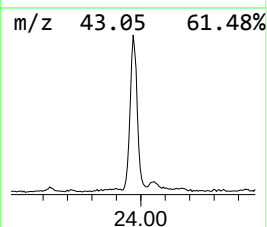
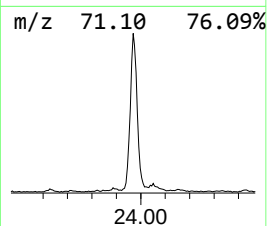
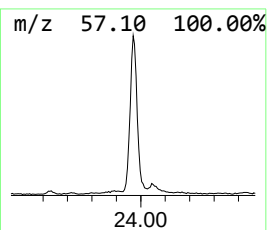
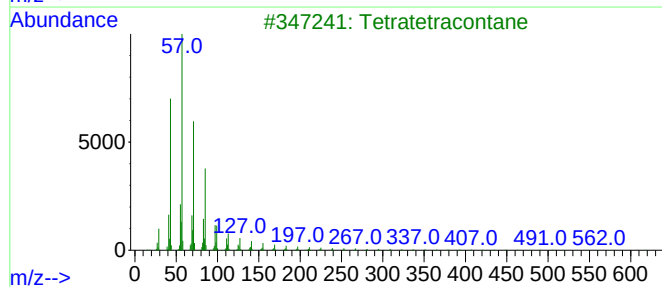
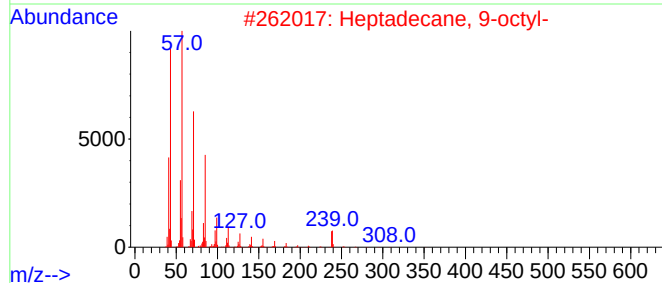
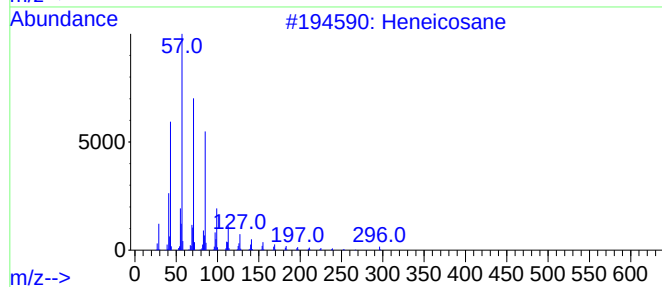
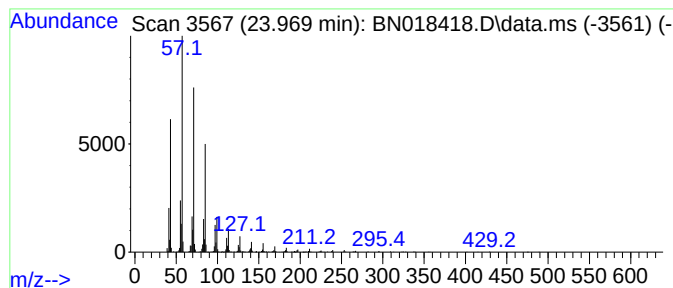
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.969	5.95 ng/ul	1943390	Perylene-d12	23.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3			Tetratetracontane	619	C44H90	007098-22-8	91
4			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5			Hexatriacontane	507	C36H74	000630-06-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN011922\
Data File : BN018418.D
Acq On : 19 Jan 2022 15:40
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Sample : N1129-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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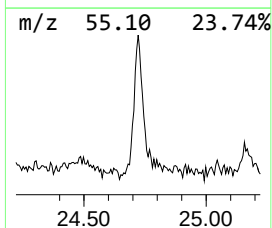
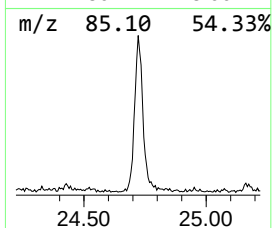
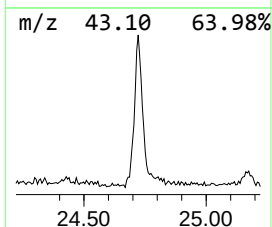
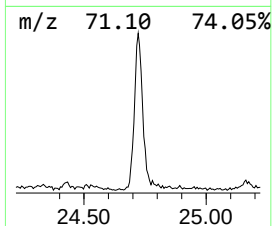
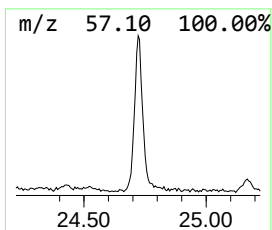
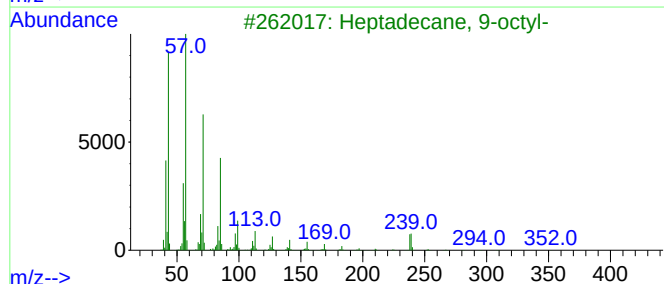
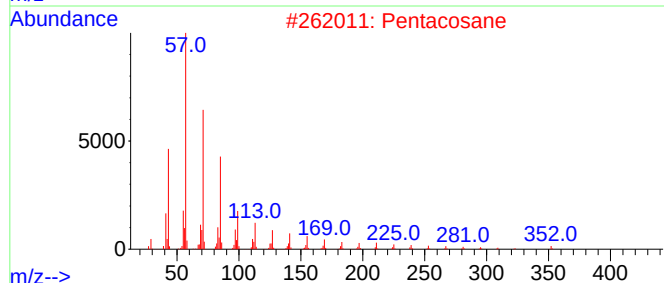
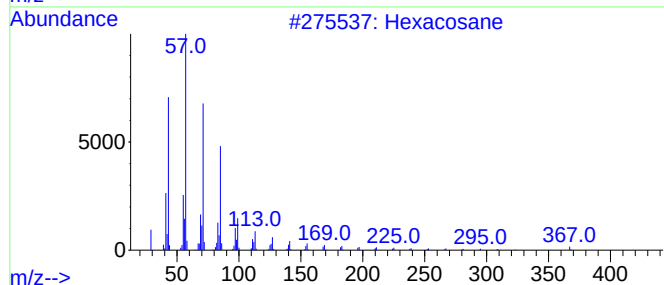
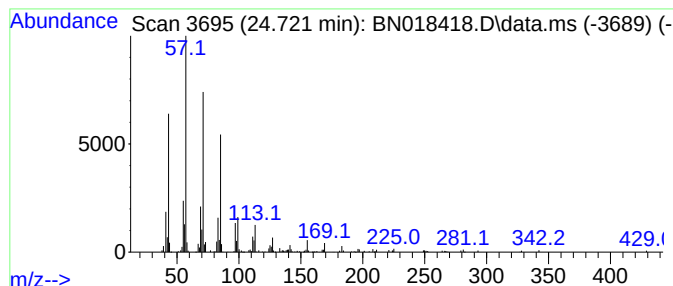
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.721	2.02 ng/ul	659048	Perylene-d12	23.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	91
2			Pentacosane	352	C25H52	000629-99-2	91
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4			Heneicosane	296	C21H44	000629-94-7	91
5			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN011922\
Data File : BN018418.D
Acq On : 19 Jan 2022 15:40
Operator : CG/JU
Sample : N1129-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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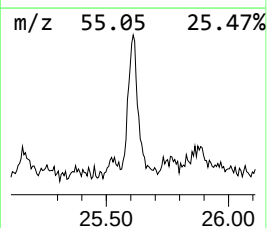
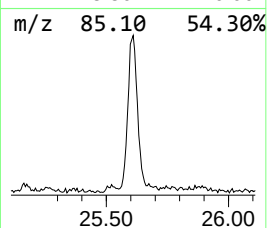
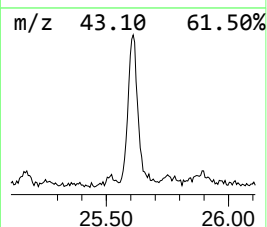
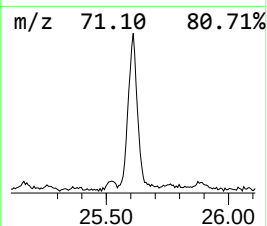
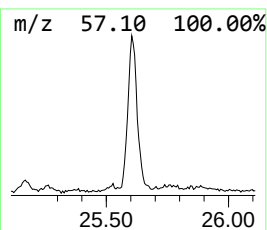
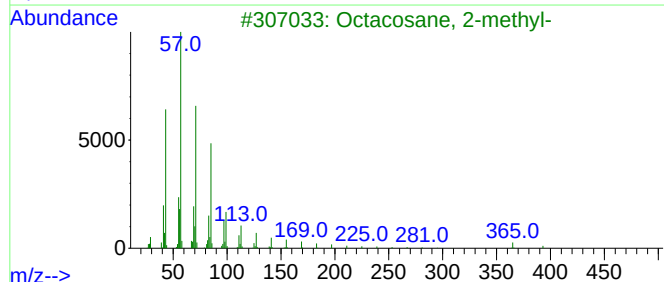
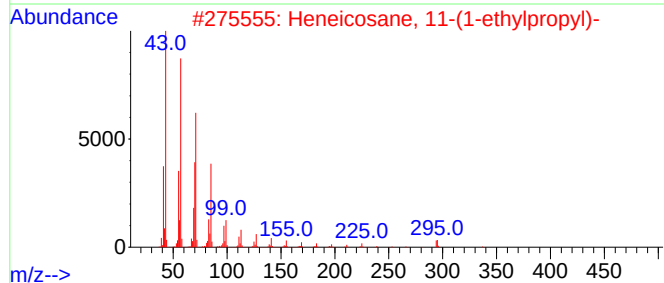
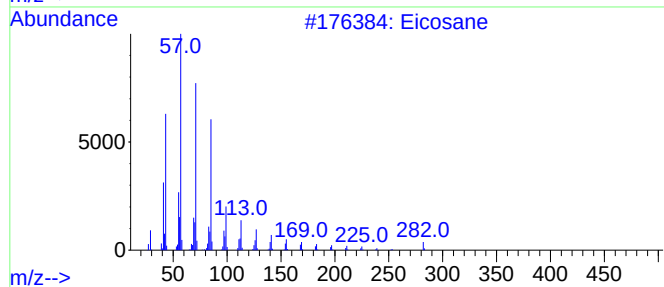
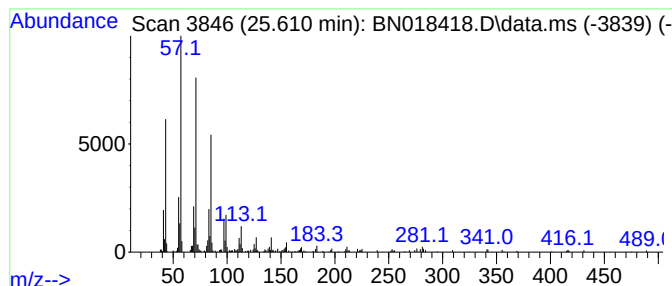
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.610	2.37 ng/ul	774137	Perylene-d12	23.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	93
2			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
3			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4			Hexacosane	366	C26H54	000630-01-3	90
5			Sulfurous acid, 2-propyl undecyl...	278	C14H30O3S	1000309-12-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN011922\
Data File : BN018418.D
Acq On : 19 Jan 2022 15:40
Operator : CG/JU
Sample : N1129-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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					#	RT	Resp	Conc
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Caryophyllene	13.563	9.2	ng/u1	2795350	3	14.234	6071710	20.0
Cinnamyl cinnamate	20.733	21.3	ng/u1	6999590	5	21.175	6587370	20.0
(DEL) Alkane: S...	22.751	5.2	ng/u1	1689210	6	23.392	6534990	20.0
(DEL) Alkane: S...	23.322	2.4	ng/u1	773000	6	23.392	6534990	20.0
(DEL) Alkane: S...	23.969	6.0	ng/u1	1943390	6	23.392	6534990	20.0
(DEL) Alkane: S...	24.721	2.0	ng/u1	659048	6	23.392	6534990	20.0
(DEL) Alkane: S...	25.610	2.4	ng/u1	774137	6	23.392	6534990	20.0