

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
 Data File : BP015894.D
 Acq On : 01 Jul 2023 21:03
 Operator : MA/JU
 Sample : 03242-10
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGB1

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-BP062623.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015894.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	29	32	40	rVB	66866	89374	0.16%	0.041%
2	3.828	106	109	122	rBV	360999	614249	1.11%	0.279%
3	4.040	141	145	150	rVB	52070	77591	0.14%	0.035%
4	5.122	320	329	336	rVB2	36555	76018	0.14%	0.034%
5	6.528	563	568	574	rBV	71716	117195	0.21%	0.053%
6	6.693	585	596	605	rBV	2016095	3095703	5.60%	1.405%
7	6.999	644	648	653	rVV	149733	229007	0.41%	0.104%
8	7.046	653	656	662	rVB	56296	91075	0.16%	0.041%
9	7.234	679	688	701	rBV	649601	1558491	2.82%	0.707%
10	7.375	706	712	721	rVV	737815	1291104	2.34%	0.586%
11	7.458	721	726	739	rVB	351556	570685	1.03%	0.259%
12	7.581	740	747	763	rBV	907731	1676022	3.03%	0.761%
13	8.046	820	826	841	rBV	1075161	1939678	3.51%	0.880%
14	8.793	946	953	971	rBV	608062	1528376	2.77%	0.694%
15	9.240	1018	1029	1048	rVV	828852	1760909	3.19%	0.799%
16	9.634	1090	1096	1101	rBV2	99003	160273	0.29%	0.073%
17	9.857	1128	1134	1142	rBV	94654	152295	0.28%	0.069%
18	9.969	1146	1153	1170	rBV	566211	1366478	2.47%	0.620%
19	10.175	1174	1188	1198	rBV4	408577	1109689	2.01%	0.504%
20	10.275	1199	1205	1211	rVB2	87616	156357	0.28%	0.071%
21	10.487	1232	1241	1244	rBV4	73055	140986	0.26%	0.064%
22	10.540	1244	1250	1267	rBV	548137	1617481	2.93%	0.734%
23	10.875	1300	1307	1320	rVV	1440812	2814430	5.09%	1.277%
24	11.004	1320	1329	1336	rVV2	127316	341863	0.62%	0.155%
25	11.081	1336	1342	1354	rVV	50780	171106	0.31%	0.078%
26	11.187	1354	1360	1372	rVV	213690	403205	0.73%	0.183%
27	11.293	1373	1378	1392	rVB10	27419	78384	0.14%	0.036%
28	12.057	1503	1508	1515	rVB5	41463	77013	0.14%	0.035%
29	13.040	1668	1675	1694	rVB10	42341	153301	0.28%	0.070%
30	13.487	1746	1751	1757	rVV4	74707	160239	0.29%	0.073%
31	13.557	1757	1763	1773	rVV5	117910	255131	0.46%	0.116%
32	13.681	1776	1784	1793	rVV6	44639	107346	0.19%	0.049%
33	13.998	1833	1838	1847	rVV2	357433	549376	0.99%	0.249%
34	14.110	1849	1857	1868	rVV	1815544	2926886	5.30%	1.328%
35	14.204	1868	1873	1881	rVB3	104509	165171	0.30%	0.075%
36	14.392	1899	1905	1919	rVB2	2158728	3490195	6.32%	1.584%

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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 >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Title : SVOA CALIBRATION

37	14.516	1920	1926	1931	rBV	134541	190554	0.34%	0.086%
38	14.640	1942	1947	1950	rBV3	50923	78160	0.14%	0.035%
39	14.698	1951	1957	1962	rVV	2002766	3080254	5.57%	1.398%
40	14.745	1962	1965	1971	rVV2	285135	483093	0.87%	0.219%
41	14.922	1990	1995	1999	rBV2	399879	546074	0.99%	0.248%
42	14.992	2003	2007	2021	rVV2	209781	594548	1.08%	0.270%
43	15.116	2023	2028	2040	rVB6	142343	437240	0.79%	0.198%
44	15.516	2092	2096	2102	rBV2	42311	78384	0.14%	0.036%
45	15.592	2102	2109	2116	rBV	808947	1180976	2.14%	0.536%
46	15.698	2117	2127	2135	rBV	2442281	4894371	8.86%	2.221%
47	15.869	2151	2156	2162	rBV	238141	472585	0.86%	0.214%
48	15.928	2162	2166	2171	rVB4	187876	282364	0.51%	0.128%
49	16.063	2184	2189	2195	rBV	459844	715600	1.29%	0.325%
50	16.151	2195	2204	2212	rVV3	307506	798771	1.45%	0.362%
51	16.251	2217	2221	2229	rVV	476452	684705	1.24%	0.311%
52	16.334	2232	2235	2239	rVB	63601	88933	0.16%	0.040%
53	16.404	2239	2247	2257	rBV5	148323	416424	0.75%	0.189%
54	16.545	2267	2271	2276	rBV2	75462	98081	0.18%	0.045%
55	16.616	2277	2283	2288	rBV5	94693	197926	0.36%	0.090%
56	16.839	2317	2321	2324	rBV	92394	135386	0.24%	0.061%
57	16.881	2324	2328	2333	rVV2	78763	134928	0.24%	0.061%
58	16.945	2335	2339	2344	rVB	216145	292833	0.53%	0.133%
59	17.175	2375	2378	2382	rBV	46928	70639	0.13%	0.032%
60	17.286	2390	2397	2402	rBV2	42035	106935	0.19%	0.049%
61	17.463	2421	2427	2432	rBV	1827387	2767879	5.01%	1.256%
62	17.510	2432	2435	2439	rVV2	255599	401811	0.73%	0.182%
63	17.563	2439	2444	2456	rVB	1945659	3049616	5.52%	1.384%
64	18.204	2549	2553	2558	rBV4	73567	130111	0.24%	0.059%
65	18.257	2558	2562	2567	rVV	504541	673637	1.22%	0.306%
66	18.322	2567	2573	2594	rVV	3743739	5589506	10.11%	2.536%
67	18.586	2615	2618	2624	rBV2	64522	96286	0.17%	0.044%
68	18.957	2678	2681	2682	rBV3	64665	71243	0.13%	0.032%
69	19.204	2718	2723	2726	rBV2	161554	271154	0.49%	0.123%
70	19.298	2736	2739	2744	rVB	204048	255179	0.46%	0.116%
71	19.404	2753	2757	2759	rBV	833664	1102341	1.99%	0.500%
72	19.428	2759	2761	2766	rVV3	575706	874698	1.58%	0.397%
73	19.539	2776	2780	2792	rBV	1793076	2534425	4.59%	1.150%
74	19.680	2800	2804	2812	rBV3	251125	470500	0.85%	0.213%
75	19.792	2818	2823	2835	rVB2	2641560	4434761	8.03%	2.012%
76	20.151	2881	2884	2889	rVB2	186916	236276	0.43%	0.107%

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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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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
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77	20.263	2900	2903	2907	rVB	331867	392666	0.71%	0.178%
78	20.586	2955	2958	2964	rVB4	584204	744546	1.35%	0.338%
79	20.710	2976	2979	2983	rVB	426607	472898	0.86%	0.215%
80	20.774	2987	2990	2995	rVB2	583374	725519	1.31%	0.329%
81	21.216	3060	3065	3074	rBV2	3859091	5787269	10.47%	2.626%
82	21.404	3094	3097	3102	rVB2	1013359	1315679	2.38%	0.597%
83	21.516	3113	3116	3118	rBV	735559	903968	1.64%	0.410%
84	21.551	3118	3122	3127	rVB	1945062	2771558	5.02%	1.258%
85	21.851	3170	3173	3178	rVB2	545568	680279	1.23%	0.309%
86	22.127	3216	3220	3223	rBV	823298	1046560	1.89%	0.475%
87	22.174	3223	3228	3237	rVB	2729462	4204402	7.61%	1.908%
88	22.704	3315	3318	3324	rVB	582581	830868	1.50%	0.377%
89	22.916	3350	3354	3359	rVB	1000881	1466240	2.65%	0.665%
90	23.233	3402	3408	3415	rVB	3658644	6293335	11.39%	2.856%
91	23.316	3417	3422	3439	rVB	2085968	4738846	8.58%	2.150%
92	23.927	3522	3526	3532	rVB2	1453176	2602429	4.71%	1.181%
93	24.098	3549	3555	3570	rVB	1403894	3325730	6.02%	1.509%
94	24.268	3581	3584	3595	rVB	897159	1691829	3.06%	0.768%
95	24.668	3645	3652	3663	rVB	3645426	8205662	14.85%	3.723%
96	24.786	3664	3672	3687	rVV	7435304	21721641	39.31%	9.857%
97	26.121	3892	3899	3909	rVB	1462920	4067099	7.36%	1.846%
98	26.633	3977	3986	3994	rBV	2557964	7633839	13.81%	3.464%
99	27.733	4163	4173	4190	rVB3	3345173	13359954	24.18%	6.062%
100	28.809	4342	4356	4374	rVB	11835325	55261115	100.00%	25.076%

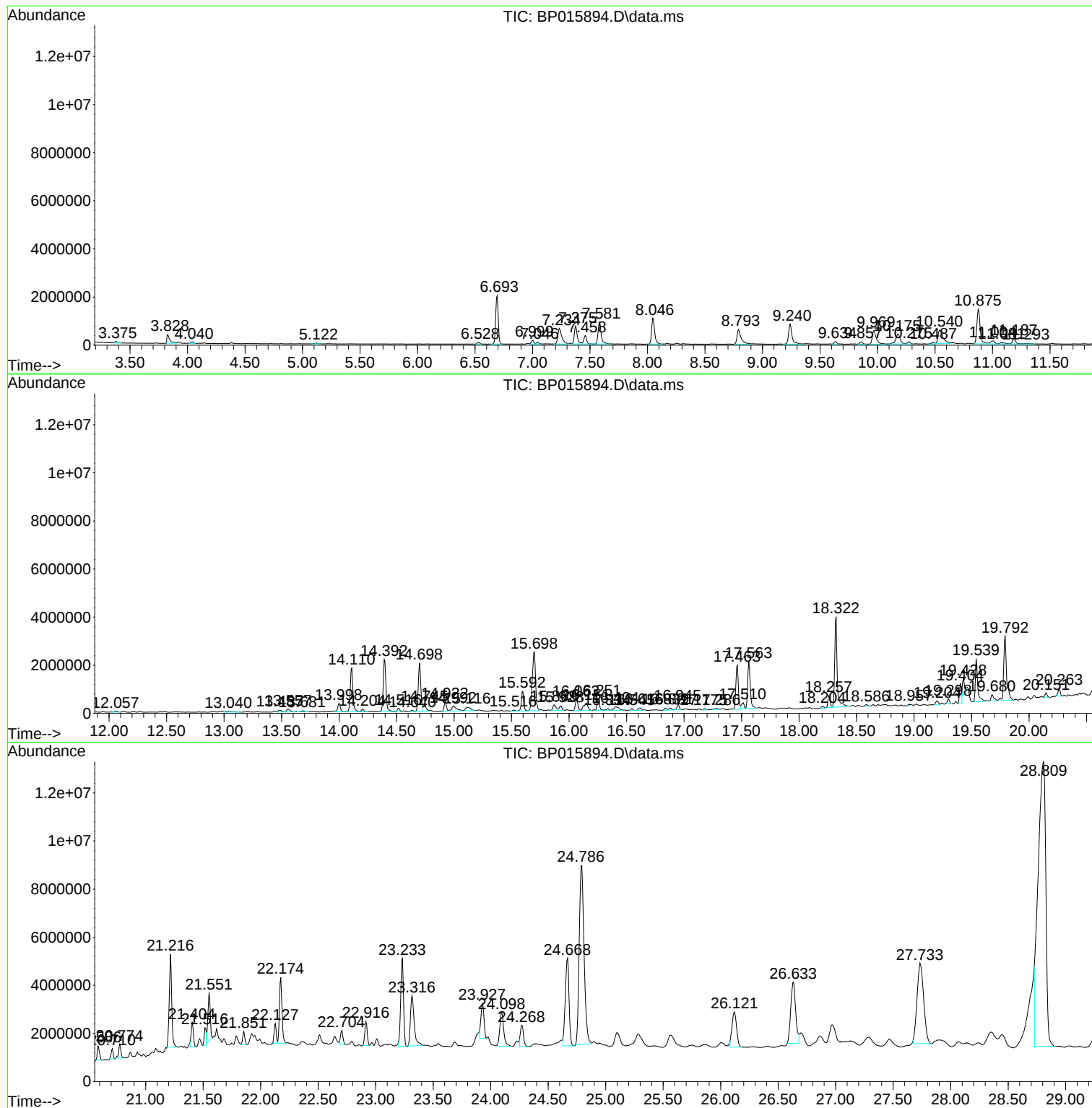
Sum of corrected areas: 220375800

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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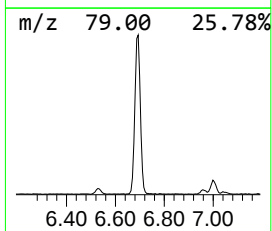
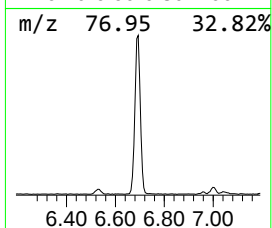
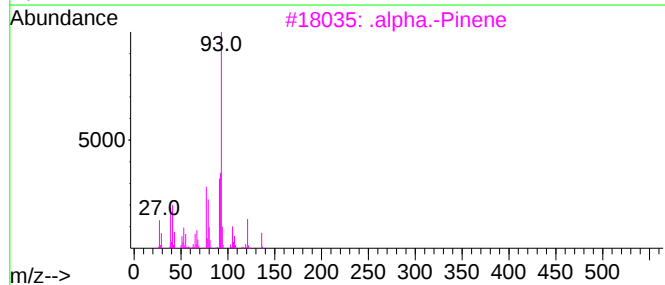
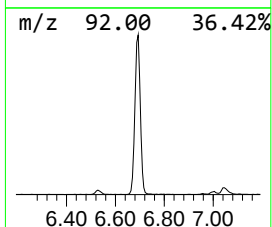
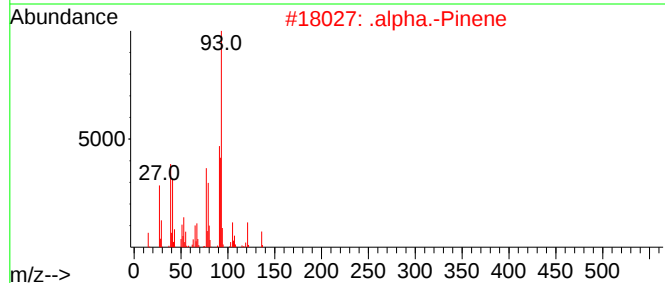
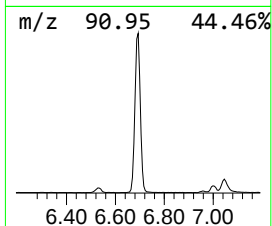
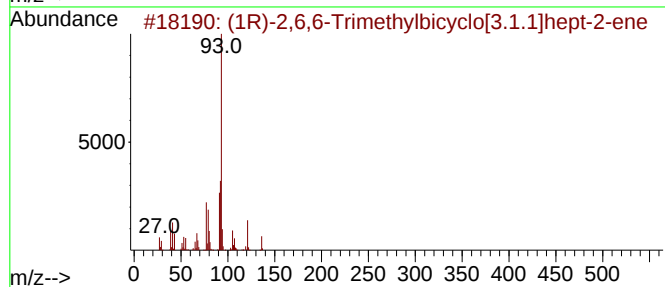
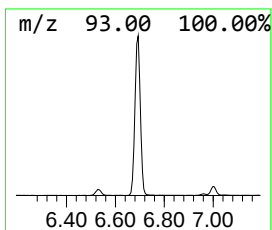
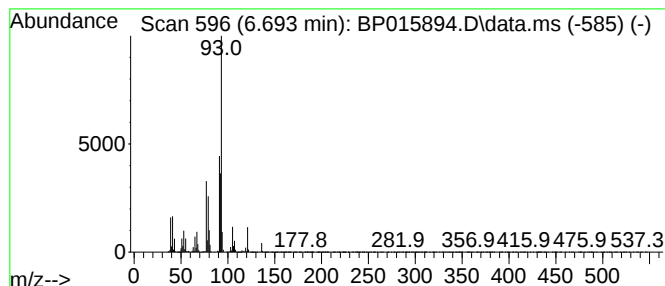
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.693	31.92 ng/ul	3095700	1,4-Dichlorobenzene-d4	8.046

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1R)-2,6,6-Trimethylbicyclo[3.1...	136	C10H16	007785-70-8	95	
2	.alpha.-Pinene	136	C10H16	000080-56-8	95	
3	.alpha.-Pinene	136	C10H16	000080-56-8	94	
4	Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	91	
5	.alpha.-Pinene	136	C10H16	000080-56-8	91	



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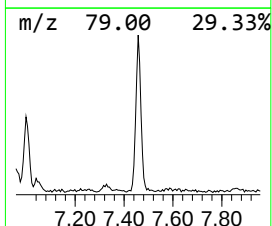
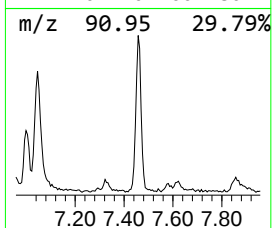
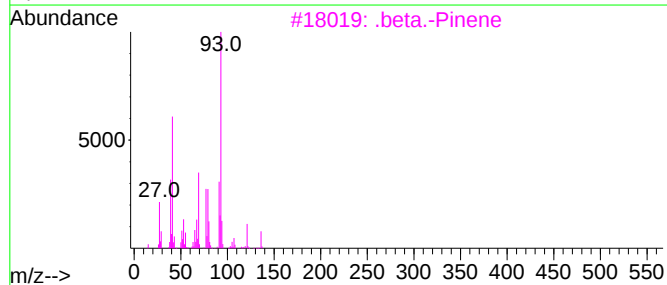
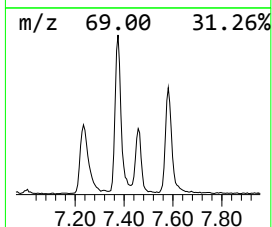
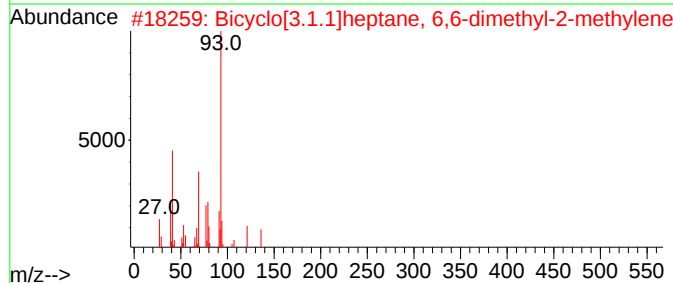
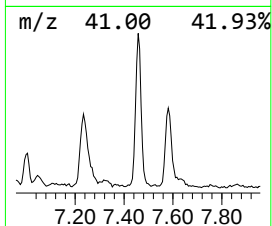
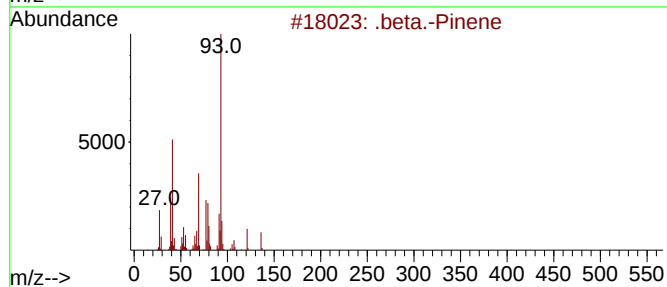
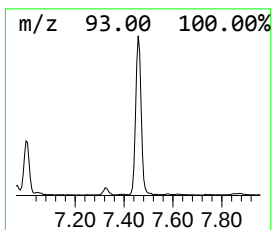
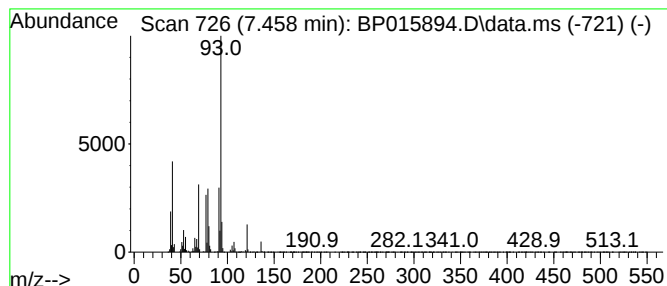
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 .beta.-Pinene Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.458	5.88 ng/ul	570685	1,4-Dichlorobenzene-d4	8.046

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		.beta.-Pinene	136	C10H16	000127-91-3	91
2		Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	91
3		.beta.-Pinene	136	C10H16	000127-91-3	91
4		Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	91
5		Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	91



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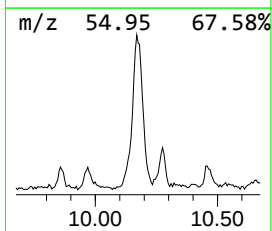
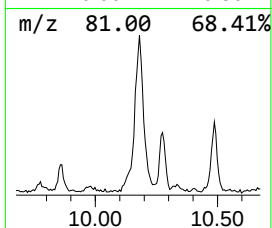
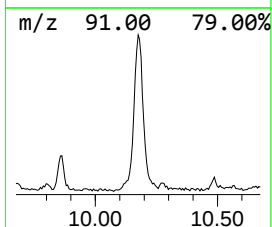
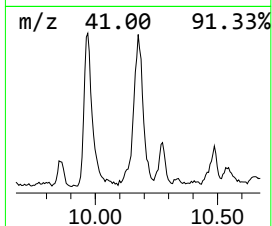
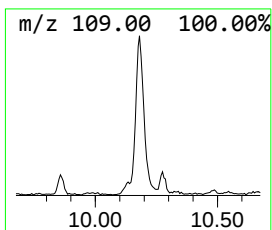
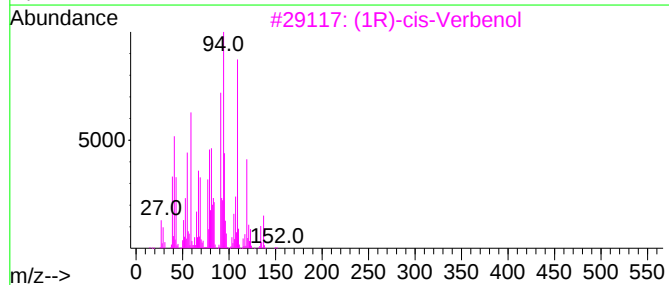
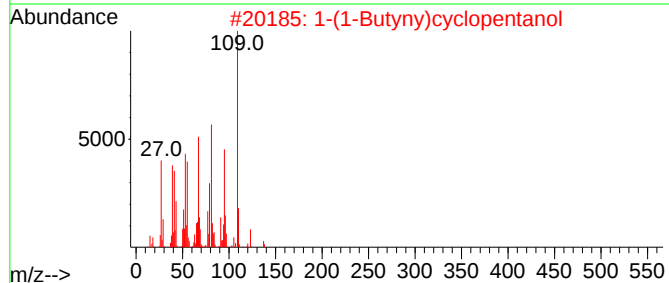
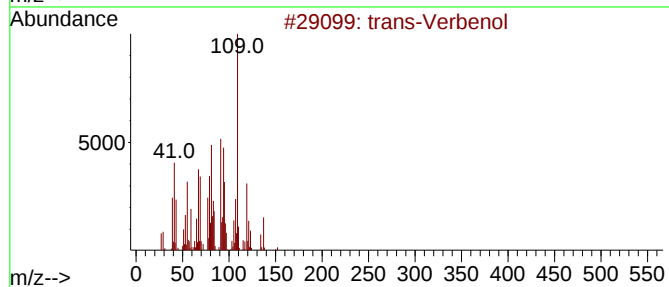
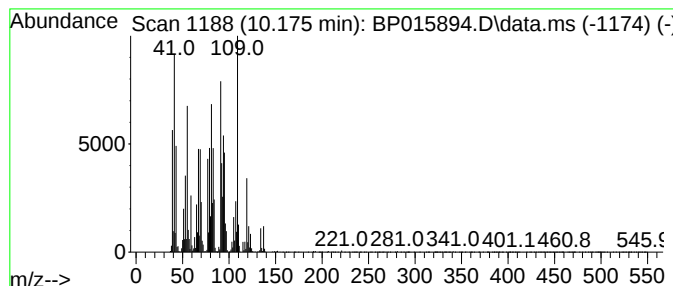
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 trans-Verbenol Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.175	7.89 ng/ul	1109690	Naphthalene-d8	10.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-Verbenol	152	C10H16O	001820-09-3	96
2			1-(1-Butyny)cyclopentanol	138	C9H14O	1000342-86-5	80
3			(1R)-cis-Verbenol	152	C10H16O	013040-03-4	70
4			cis-Verbenol	152	C10H16O	001845-30-3	43
5			Cycloheptane, 1,3,5-tris(methyle...	134	C10H14	068284-24-2	38



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Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
ALS Vial : 62 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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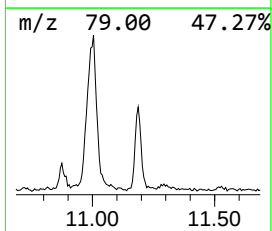
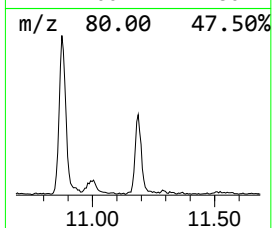
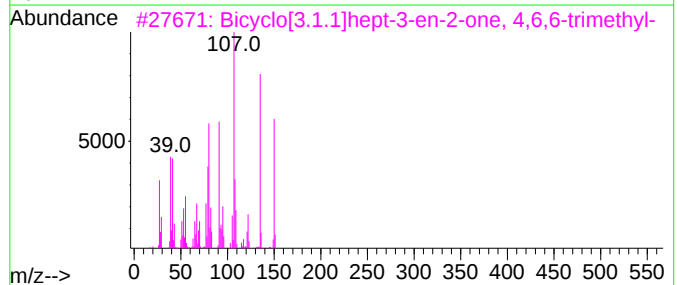
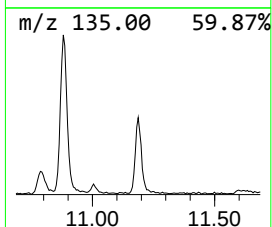
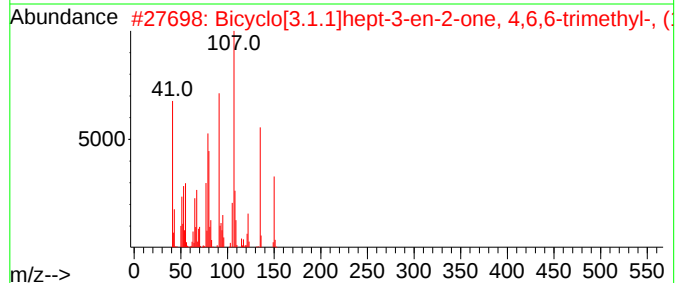
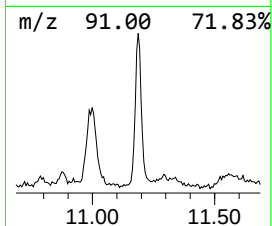
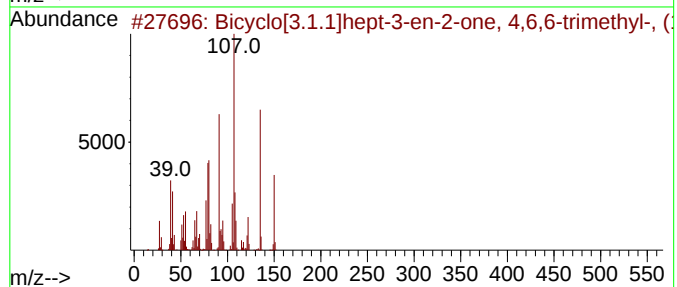
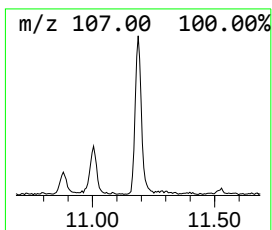
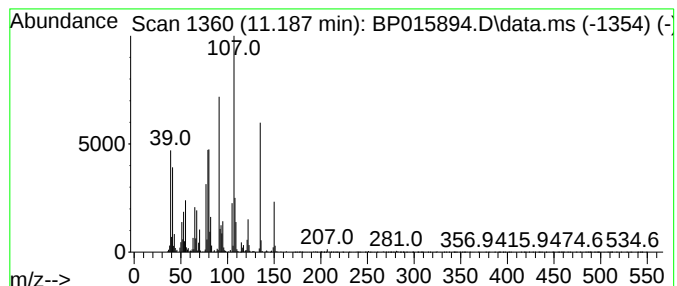
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Bicyclo[3.1.1]hept-3-en-2-o... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.187	2.87 ng/ul	403205	Naphthalene-d8	10.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[3.1.1]hept-3-en-2-one, 4...	150	C10H14O	001196-01-6	98
2			Bicyclo[3.1.1]hept-3-en-2-one, 4...	150	C10H14O	001196-01-6	97
3			Bicyclo[3.1.1]hept-3-en-2-one, 4...	150	C10H14O	000080-57-9	93
4			Bicyclo[3.1.1]hept-3-en-2-one, 4...	150	C10H14O	000080-57-9	93
5			Bicyclo[3.1.1]hept-3-en-2-one, 4...	150	C10H14O	000080-57-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
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Misc :
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ClientSampleId :
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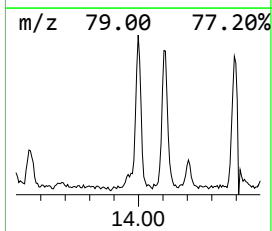
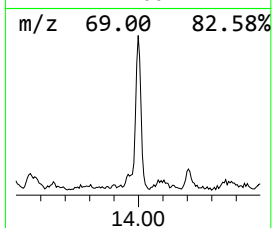
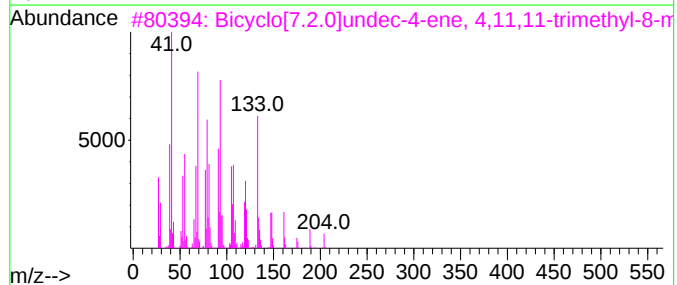
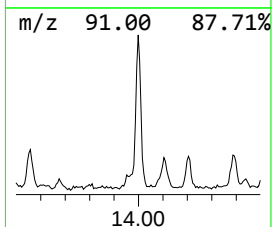
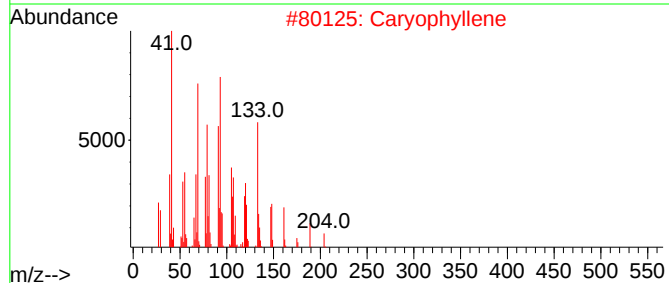
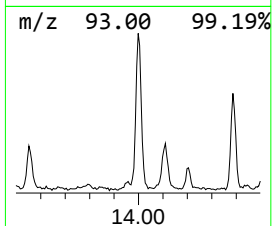
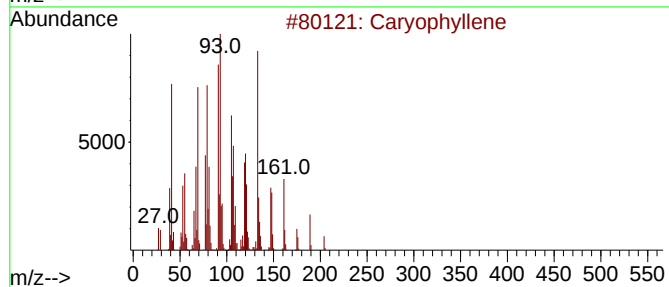
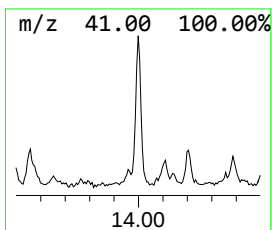
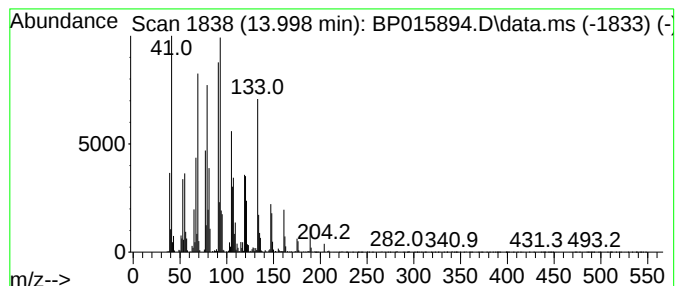
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Caryophyllene Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.998	3.57 ng/ul	549376	Acenaphthene-d10	14.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Caryophyllene	204	C15H24	000087-44-5	99
2			Caryophyllene	204	C15H24	000087-44-5	97
3			Bicyclo[7.2.0]undec-4-ene, 4,11,...	204	C15H24	000118-65-0	95
4			Caryophyllene	204	C15H24	000087-44-5	93
5			Caryophyllene	204	C15H24	000087-44-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
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Sample : 03242-10
Misc :
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ClientSampleId :
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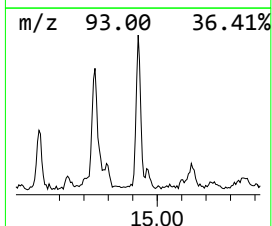
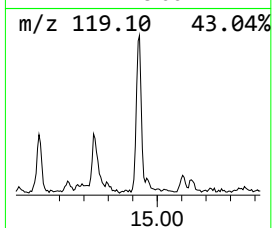
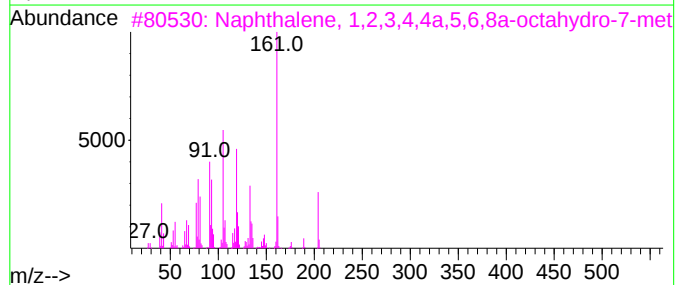
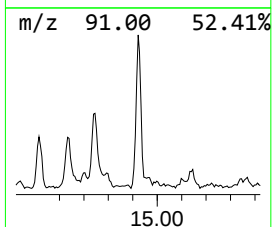
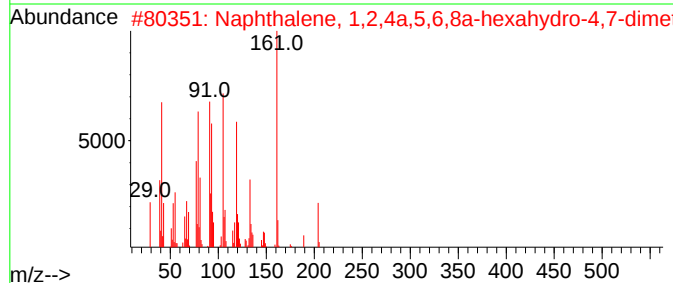
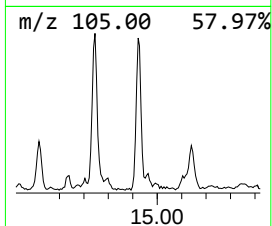
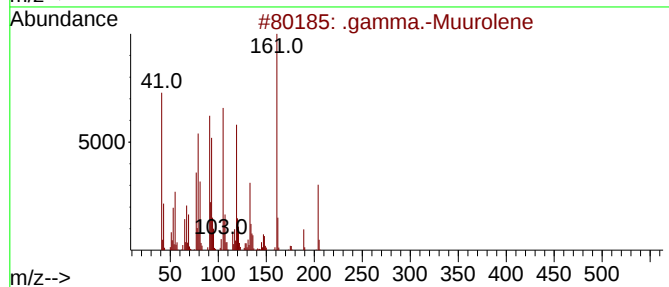
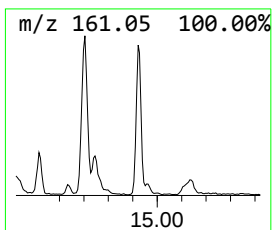
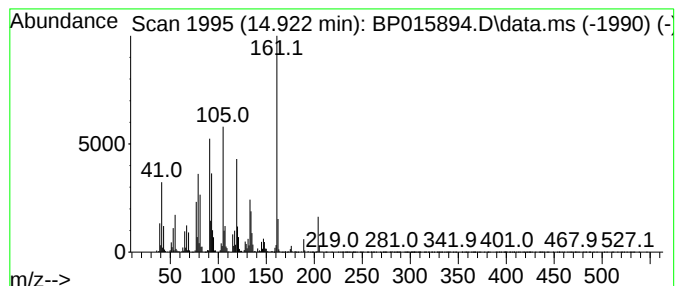
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 .gamma.-Muurolene Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.922	3.55 ng/ul	546074	Acenaphthene-d10	14.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.gamma.-Muurolene	204	C15H24	030021-74-0	99
2			Naphthalene, 1,2,4a,5,6,8a-hexah...	204	C15H24	000483-75-0	96
3			Naphthalene, 1,2,3,4,4a,5,6,8a-o...	204	C15H24	039029-41-9	95
4			(1S,4S,4aS)-1-Isopropyl-4,7-dime...	204	C15H24	267665-20-3	90
5			1H-Cyclopenta[1,3]cyclopropa[1,2...	204	C15H24	013744-15-5	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
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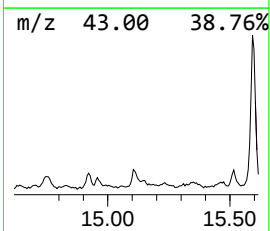
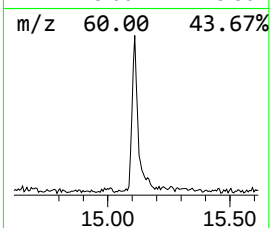
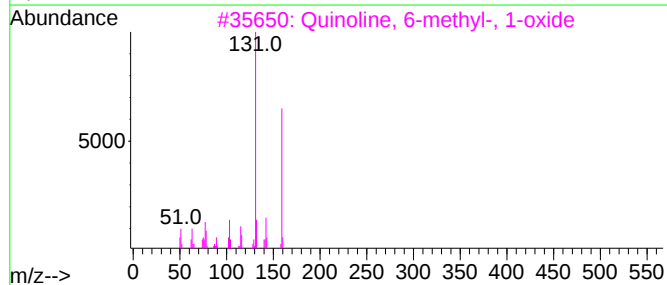
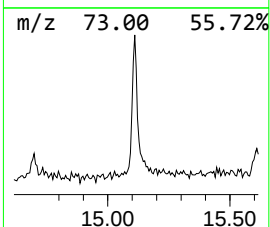
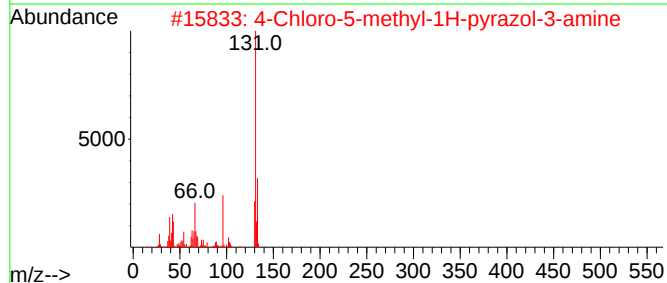
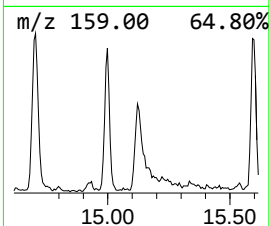
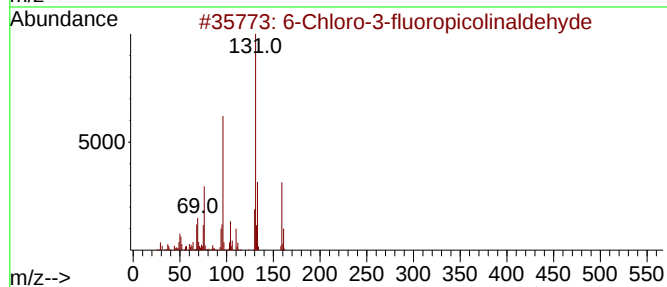
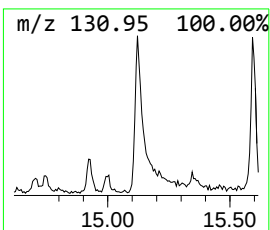
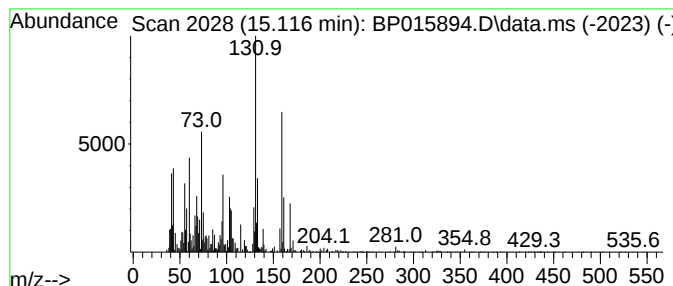
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.116	2.84 ng/ul	437240	Acenaphthene-d10	14.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Chloro-3-fluoropicolinaldehyde	159	C6H3ClFNO	884494-77-3	46
2			4-Chloro-5-methyl-1H-pyrazol-3-a...	131	C4H6ClN3	110580-44-4	38
3			Quinoline, 6-methyl-, 1-oxide	159	C10H9NO	004053-42-3	35
4			2,2-Dimethyl-3-pentanol, TMS der...	188	C10H24OSi	1000364-14-3	27
5			1,2,4-Triazine-3,5(2H,4H)-dione,...	131	C3H2FN3O2	082736-95-6	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
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ClientSampleId :
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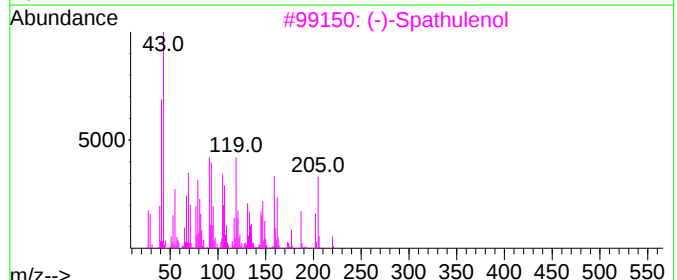
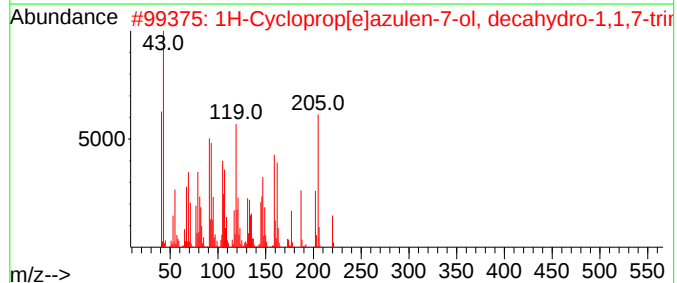
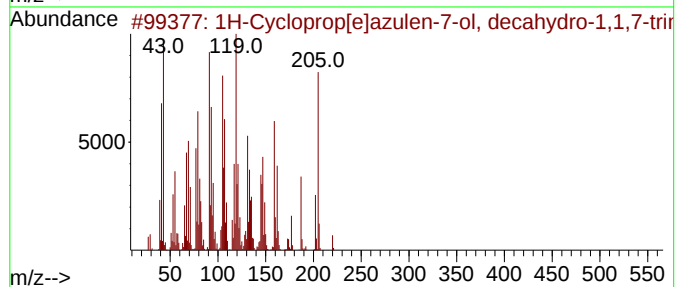
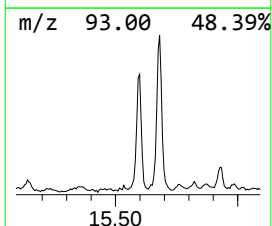
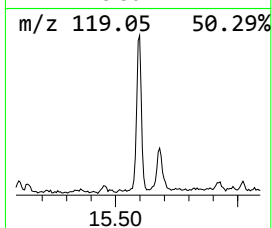
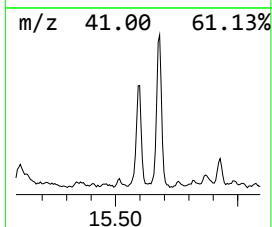
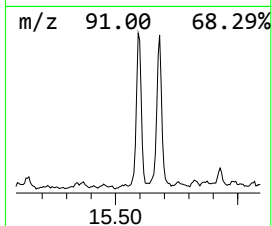
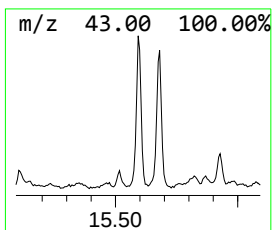
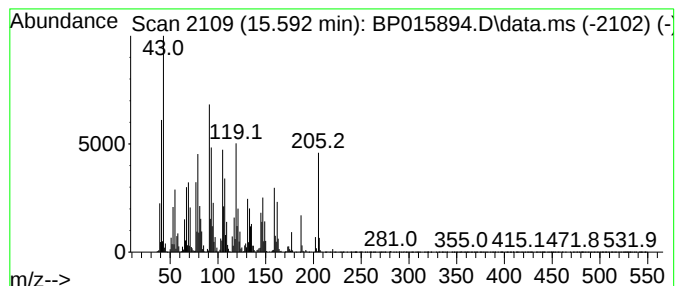
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 1H-Cycloprop[e]azulen-7-ol,... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.592	7.67 ng/ul	1180980	Acenaphthene-d10	14.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cycloprop[e]azulen-7-ol, deca...	220	C15H24O	006750-60-3	93
2			1H-Cycloprop[e]azulen-7-ol, deca...	220	C15H24O	006750-60-3	87
3			(-)-Spathulenol	220	C15H24O	077171-55-2	81
4			Isospathulenol	220	C15H24O	088395-46-4	43
5			Icosa-9,11-diyne	274	C20H34	028393-07-9	22



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Misc :
ALS Vial : 62 Sample Multiplier: 1

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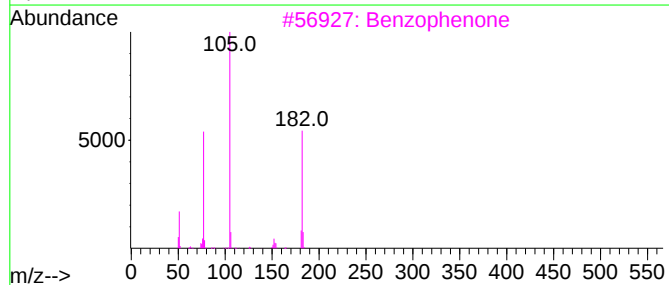
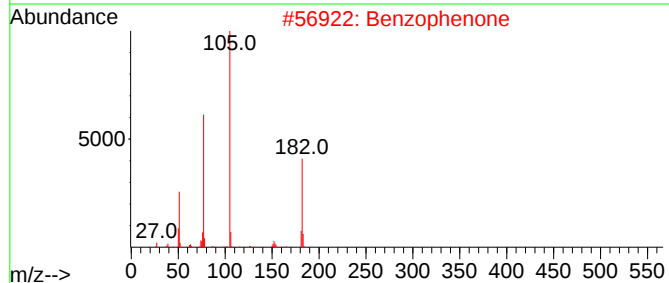
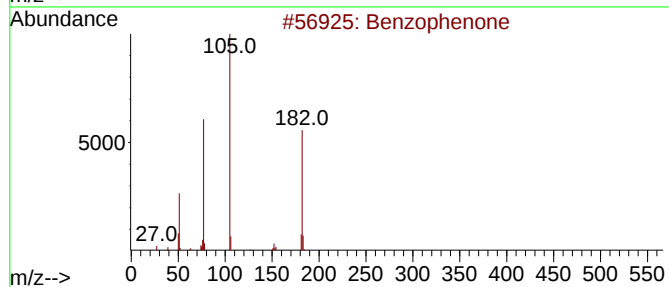
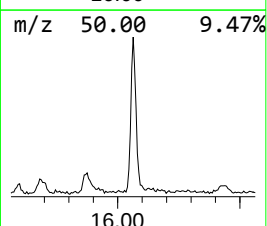
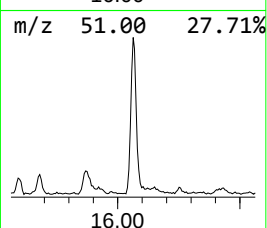
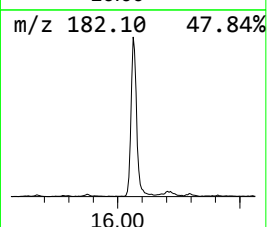
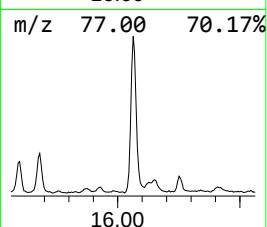
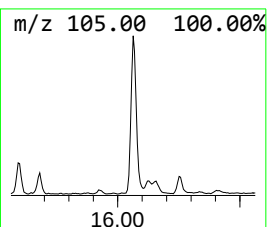
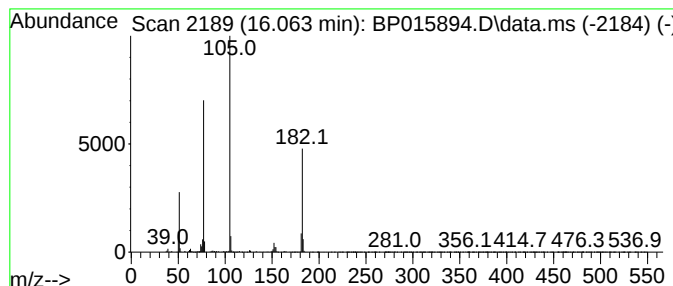
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzophenone Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.063	4.65 ng/ul	715600	Acenaphthene-d10	14.698

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
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Misc :
ALS Vial : 62 Sample Multiplier: 1

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ClientSampleId :
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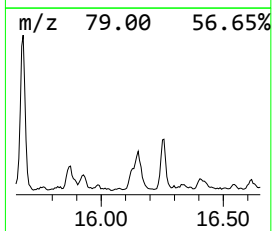
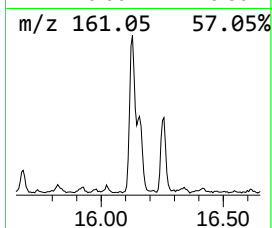
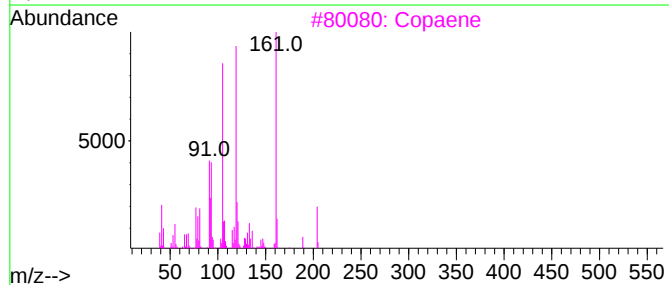
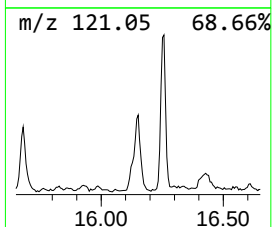
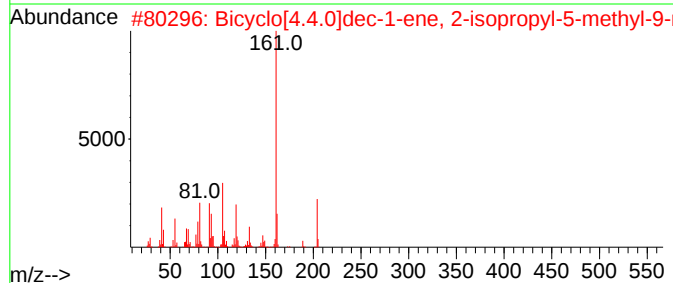
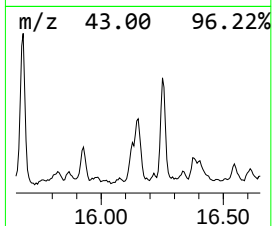
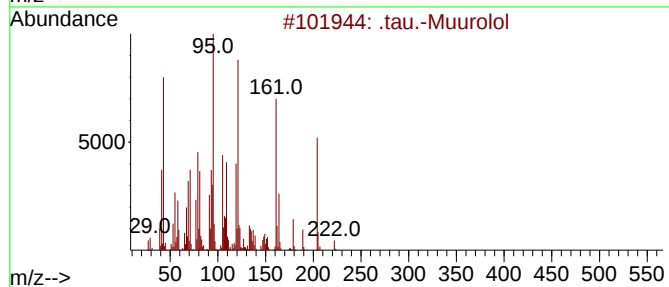
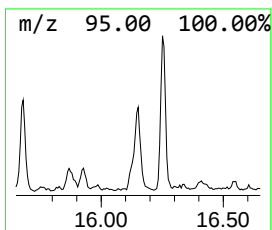
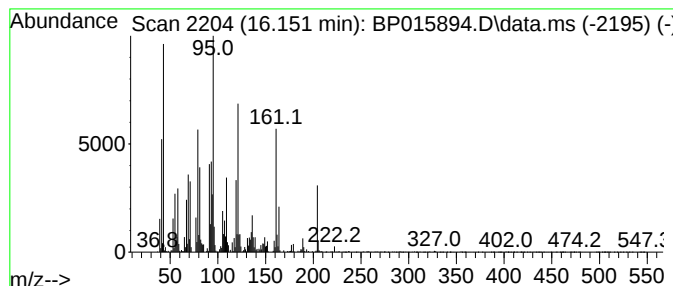
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 .tau.-Muurolo1 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.151	5.77 ng/ul	798771	Phenanthrene-d10	17.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	tau.-Muurolo1	222	C15H26O	019912-62-0	91	
2	Bicyclo[4.4.0]dec-1-ene, 2-isopr...	204	C15H24	150320-52-8	50		
3	Copaene	204	C15H24	003856-25-5	50		
4	.alpha.-Cadinol	222	C15H26O	000481-34-5	46		
5	.gamma.-Muurolene	204	C15H24	030021-74-0	45		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
ALS Vial : 62 Sample Multiplier: 1

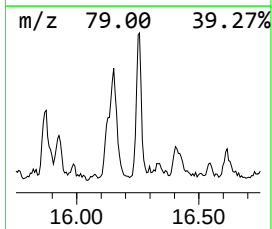
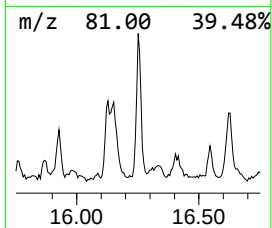
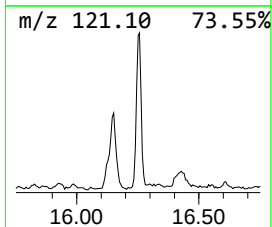
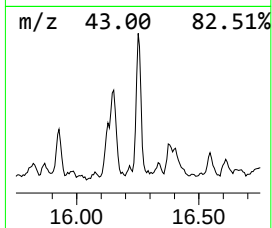
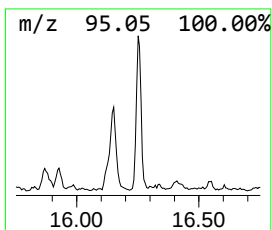
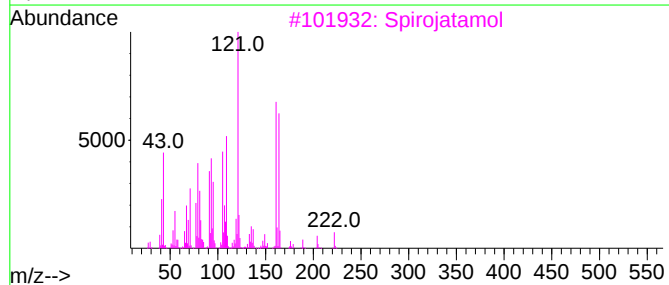
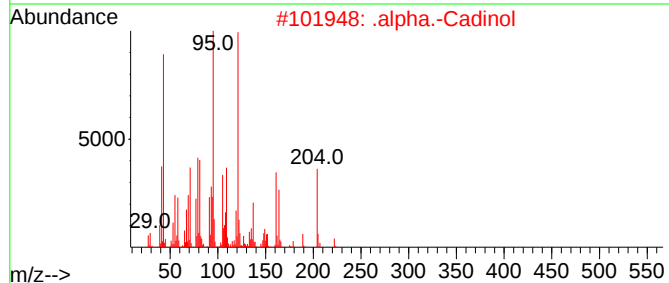
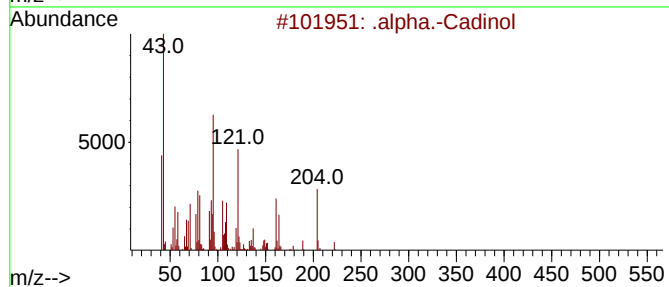
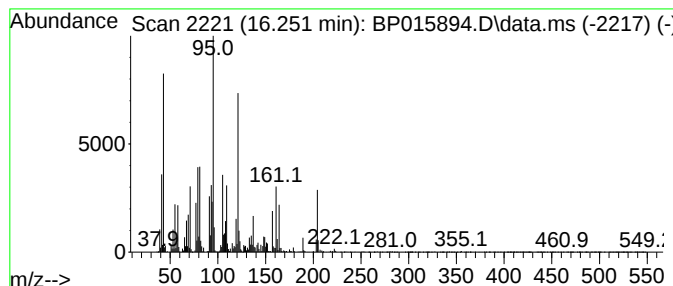
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 .alpha.-Cadinol Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.251	4.95 ng/ul	684705	Phenanthrene-d10	17.463	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Cadinol	222	C15H26O	000481-34-5	96
2	.alpha.-Cadinol	222	C15H26O	000481-34-5	94
3	Spirojatamol	222	C15H26O	128487-46-7	49
4	Cyclohexane, 1-ethenyl-1-methyl-...	204	C15H24	003242-08-8	49
5	Tricyclo[3.1.0.0(2,4)]hexane, 3,...	136	C10H16	058987-01-2	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
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Sample : 03242-10
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ALS Vial : 62 Sample Multiplier: 1

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ClientSampleId :
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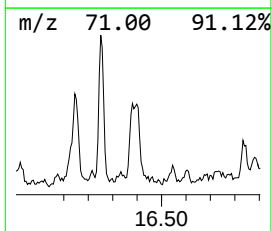
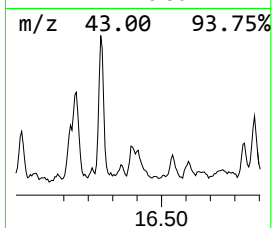
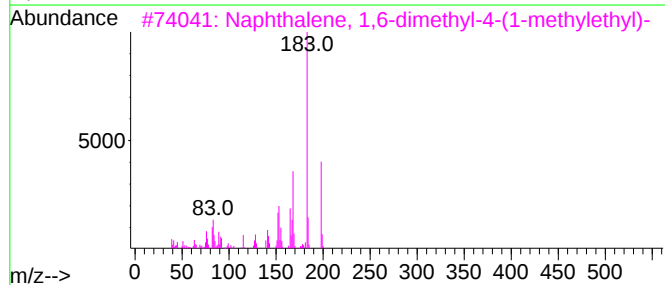
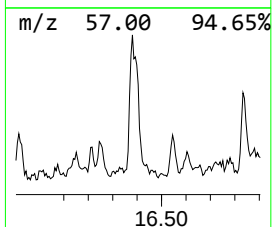
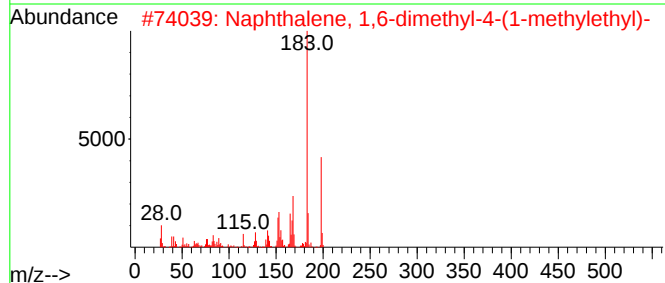
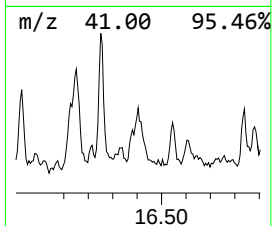
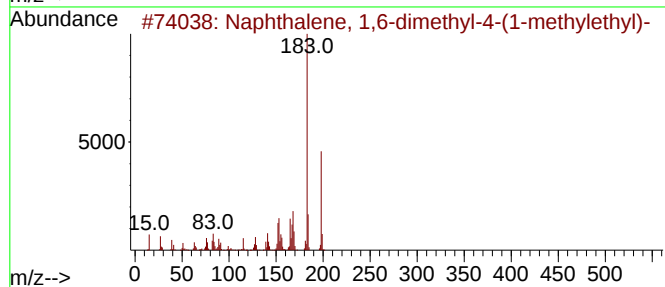
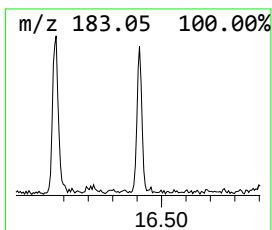
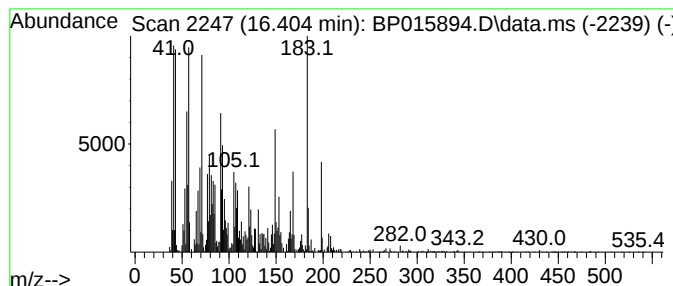
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Naphthalene, 1,6-dimethyl-4... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.404	3.01 ng/ul	416424	Phenanthrene-d10	17.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-4-(1-m...	198	C15H18	000483-78-3	89
2			Naphthalene, 1,6-dimethyl-4-(1-m...	198	C15H18	000483-78-3	64
3			Naphthalene, 1,6-dimethyl-4-(1-m...	198	C15H18	000483-78-3	64
4			6-Isopropyl-1,4-dimethylnaphthalene	198	C15H18	000489-77-0	60
5			Naphthalene, 1,6-dimethyl-4-(1-m...	198	C15H18	000483-78-3	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
ALS Vial : 62 Sample Multiplier: 1

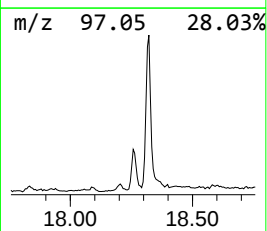
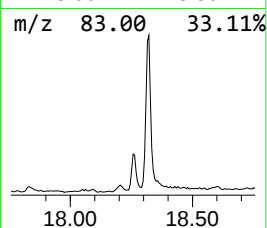
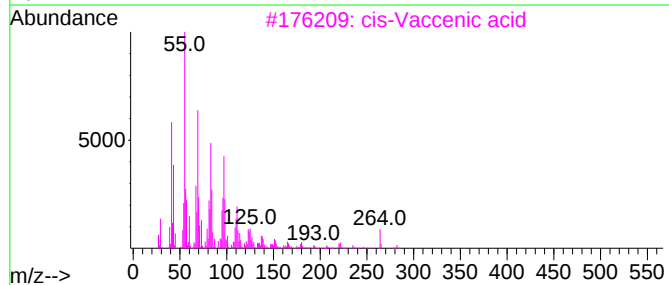
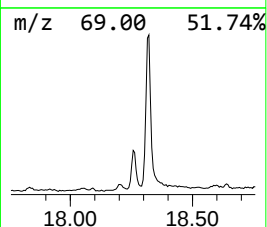
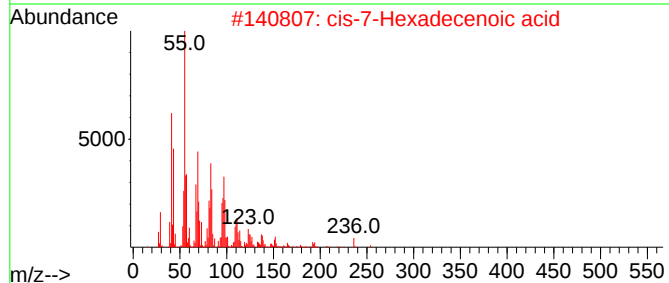
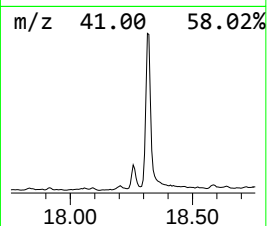
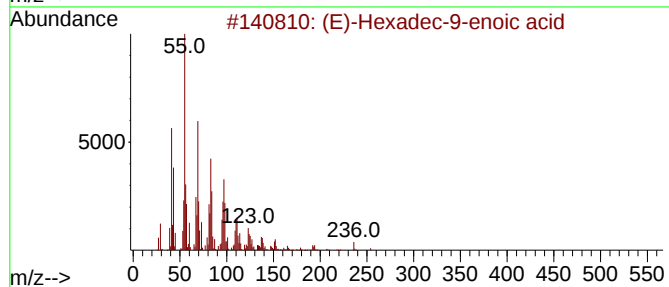
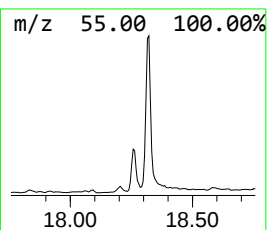
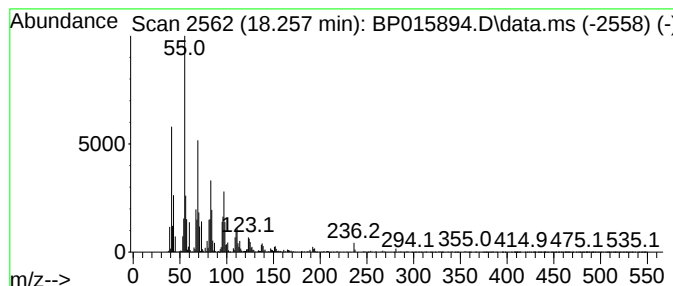
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (E)-Hexadec-9-enoic acid Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.257	4.87 ng/ul	673637	Phenanthrene-d10	17.463	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	99
2	cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	99
3	cis-Vaccenic acid	282	C18H34O2	000506-17-2	98
4	Erucic acid	338	C22H42O2	000112-86-7	95
5	Palmitoleic acid	254	C16H30O2	000373-49-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
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Sample : 03242-10
Misc :
ALS Vial : 62 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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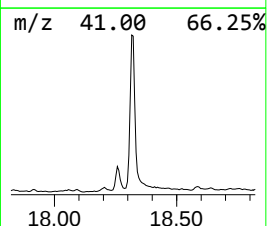
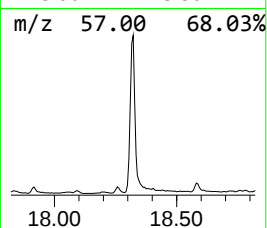
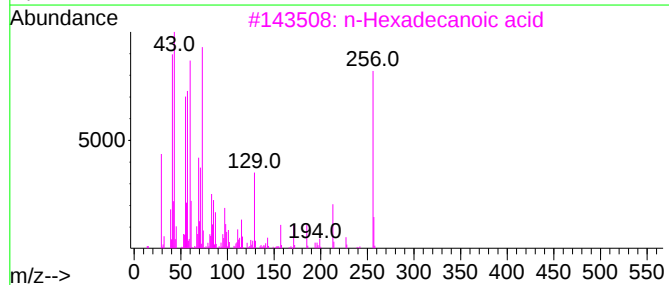
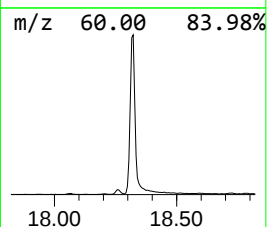
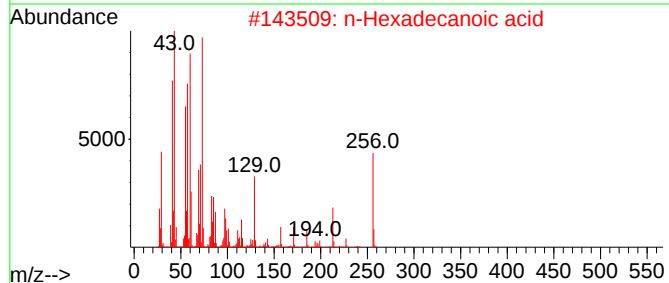
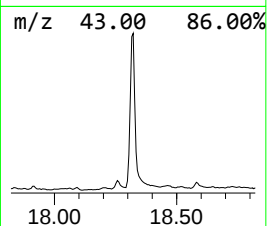
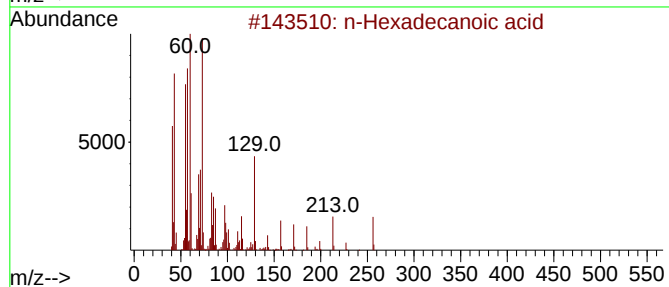
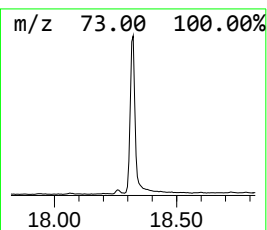
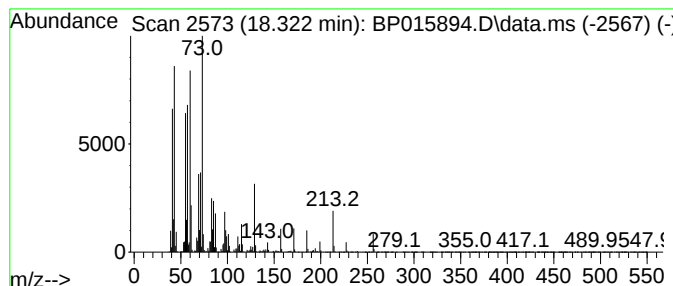
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.322	40.39 ng/ul	5589510	Phenanthrene-d10	17.463

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			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
ALS Vial : 62 Sample Multiplier: 1

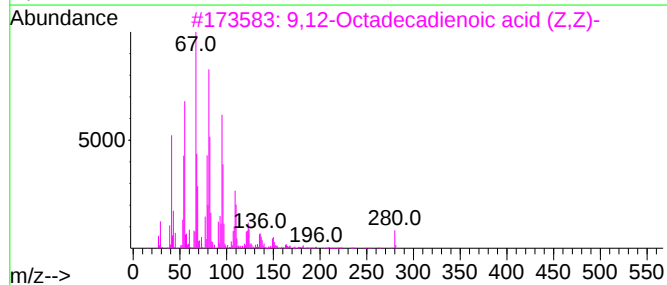
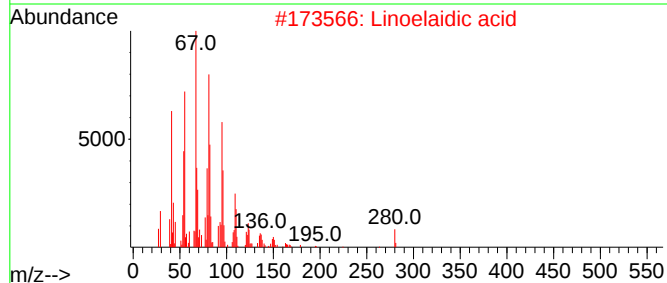
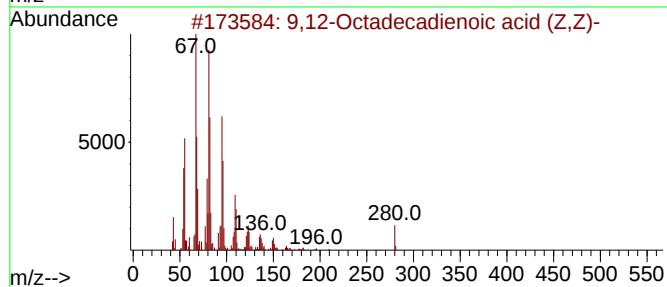
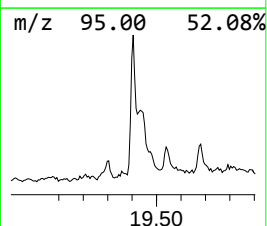
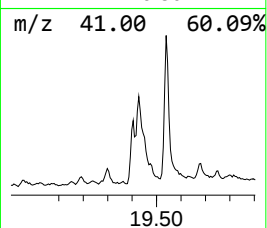
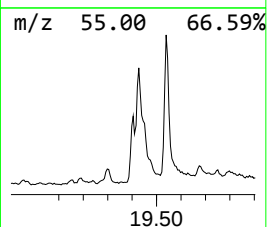
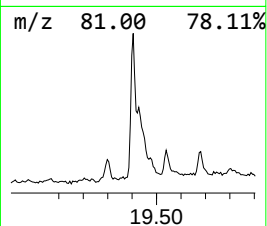
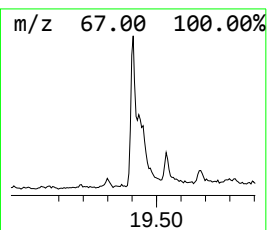
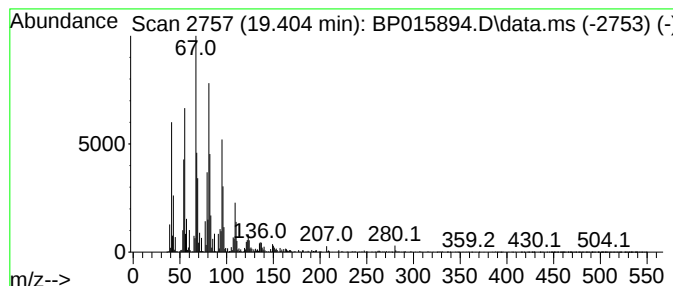
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 9,12-Octadecadienoic acid (... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.404	7.97 ng/ul	1102340	Phenanthrene-d10	17.463	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	99
2	Linoelaidic acid	280	C18H32O2	000506-21-8	99
3	9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	99
4	10E,12Z-Octadecadienoic acid	280	C18H32O2	002420-56-6	98
5	(9E,11E)-Octadecadienoic acid	280	C18H32O2	000544-71-8	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
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ClientSampleId :
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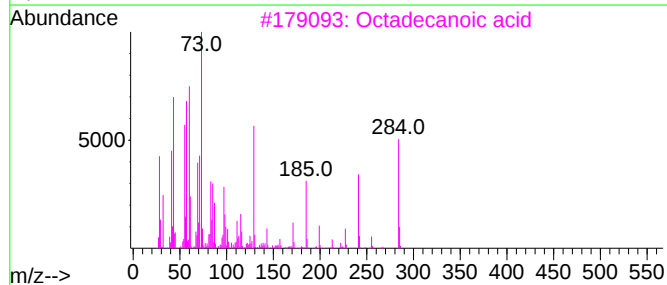
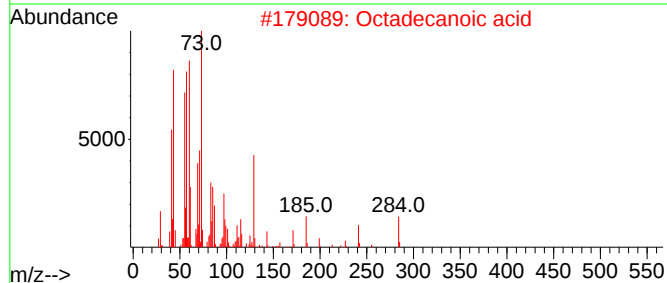
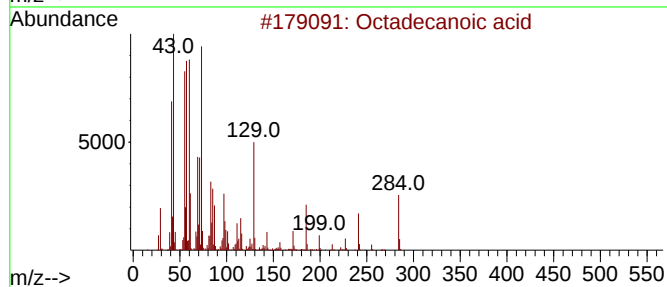
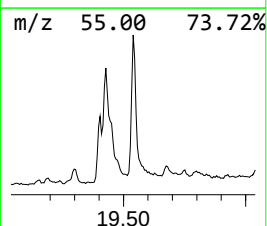
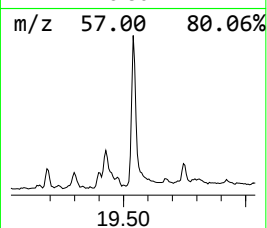
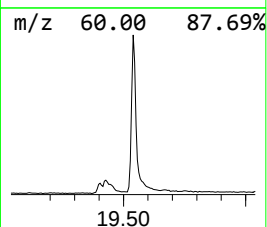
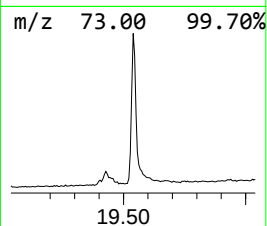
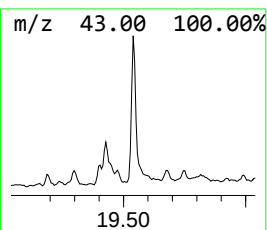
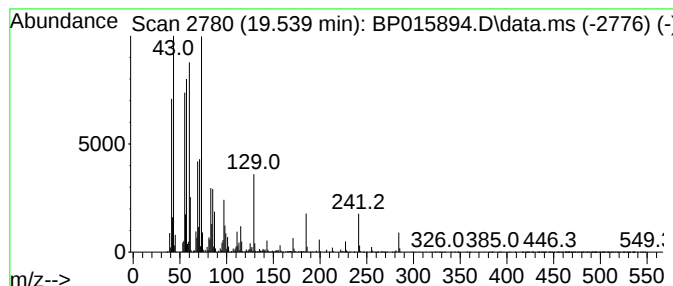
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Octadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.539	18.29 ng/ul	2534430	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
ALS Vial : 62 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGB1

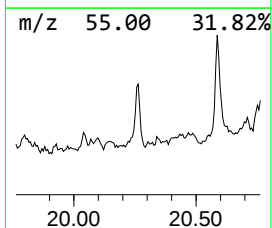
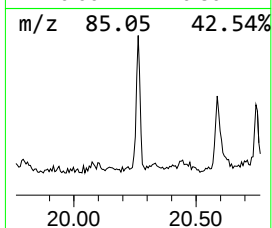
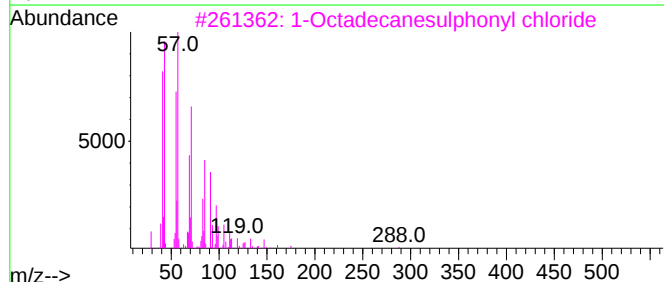
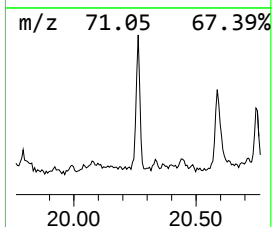
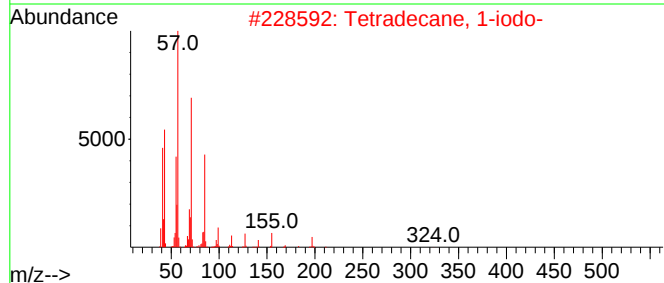
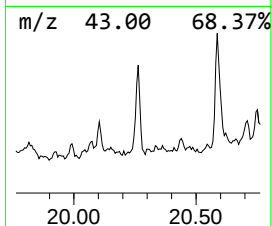
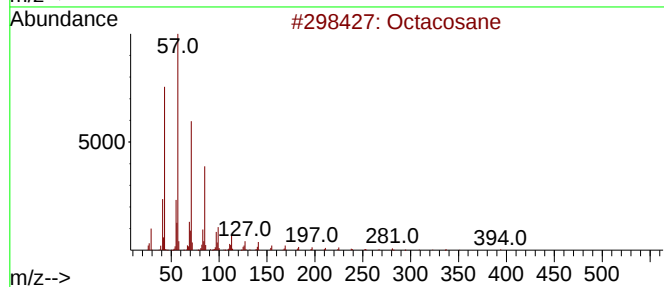
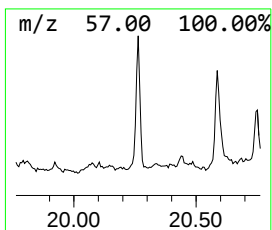
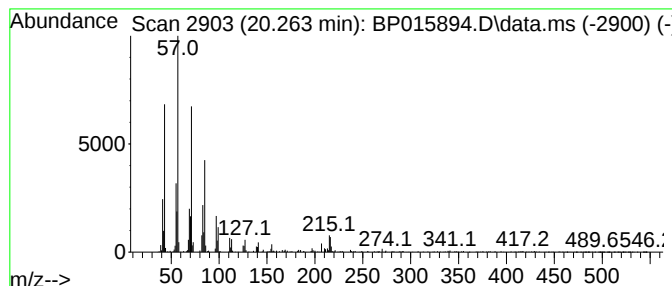
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.263	2.83 ng/ul	392666	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	93
2			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	90
3			1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	87
4			Tetratriacontane, 17-hexadecyl-	703	C50H102	055256-07-0	87
5			Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
ALS Vial : 62 Sample Multiplier: 1

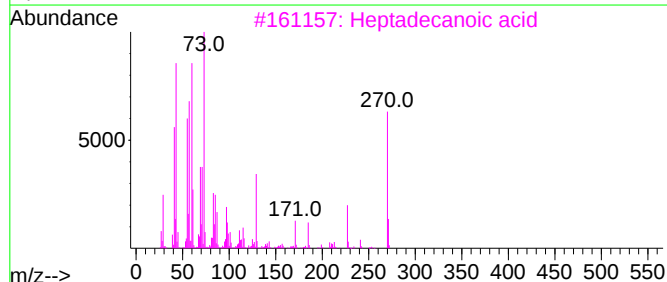
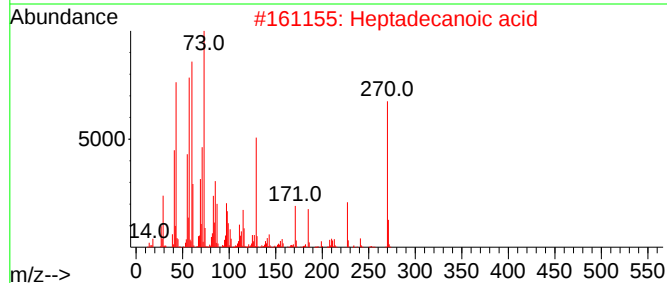
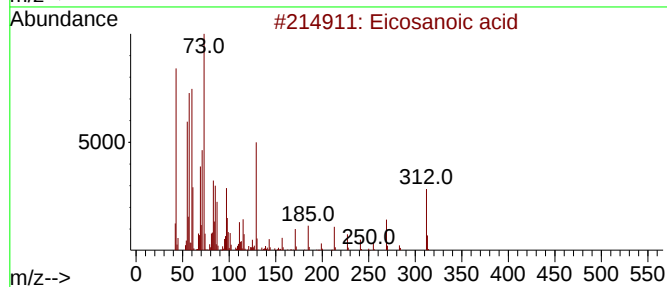
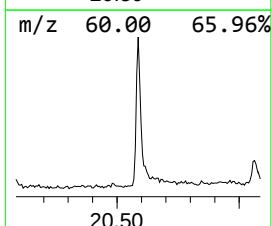
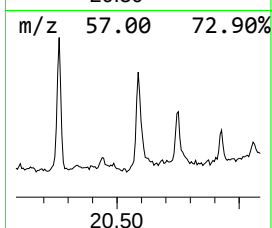
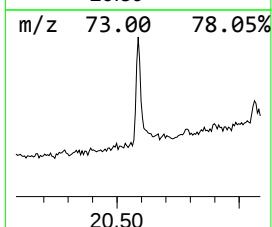
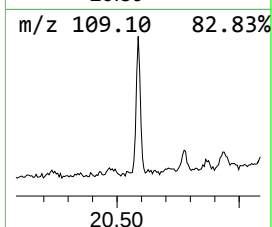
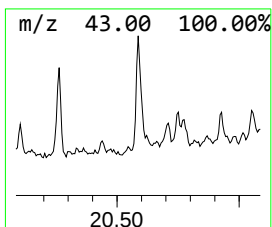
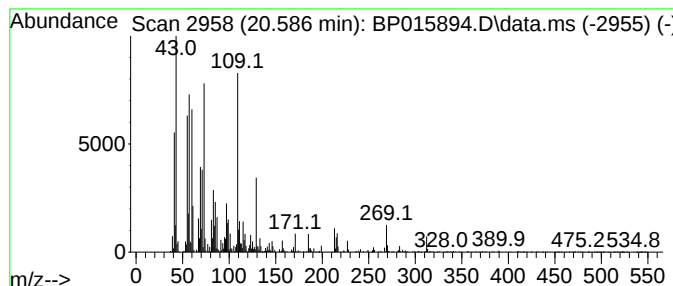
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Eicosanoic acid Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.586	5.37 ng/ul	744546	Chrysene-d12	21.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosanoic acid	312	C20H40O2	000506-30-9	94
2	Heptadecanoic acid	270	C17H34O2	000506-12-7	91
3	Heptadecanoic acid	270	C17H34O2	000506-12-7	91
4	Eicosanoic acid	312	C20H40O2	000506-30-9	89
5	Eicosanoic acid	312	C20H40O2	000506-30-9	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
ALS Vial : 62 Sample Multiplier: 1

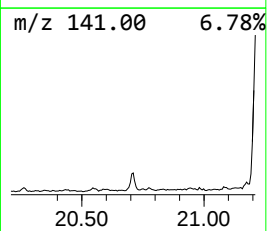
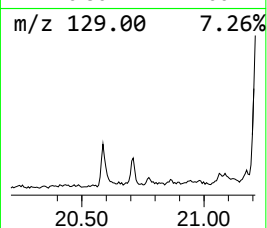
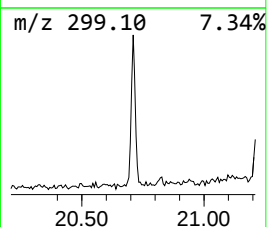
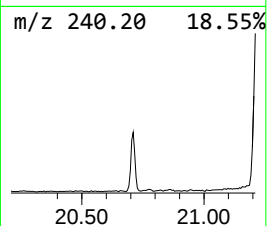
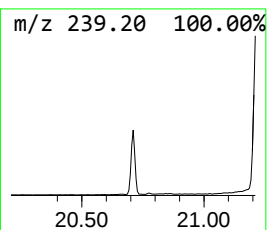
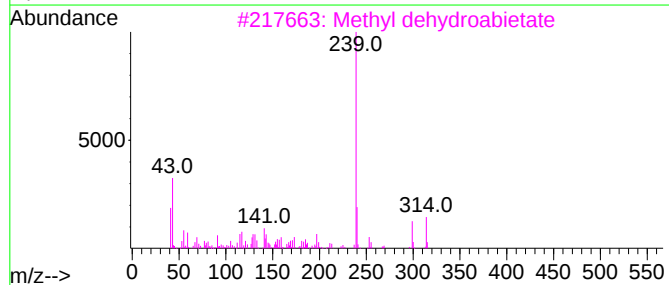
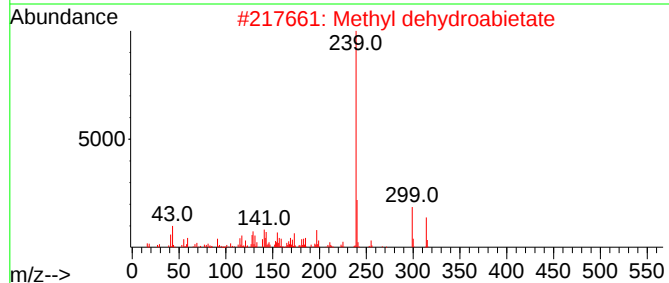
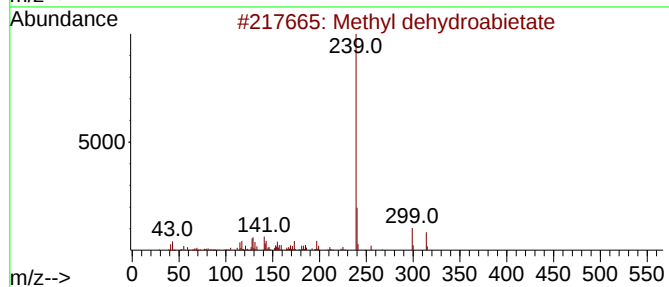
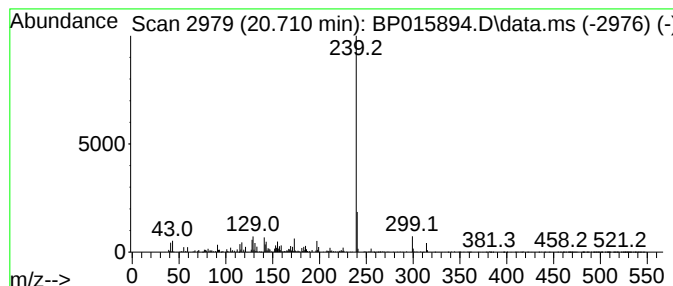
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Methyl dehydroabietate Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.710	3.41 ng/ul	472898	Chrysene-d12	21.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl dehydroabietate	314	C21H30O2	001235-74-1	98
2	Methyl dehydroabietate	314	C21H30O2	001235-74-1	78
3	Methyl dehydroabietate	314	C21H30O2	001235-74-1	78
4	Imidazolo[2,1-b]thiazole, 5-(3-i...	239	C13H9N3S	327097-37-0	59
5	Pyrano[3,2-c]chromene-3-carbonit...	361	C19H11N3O5	177028-91-0	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
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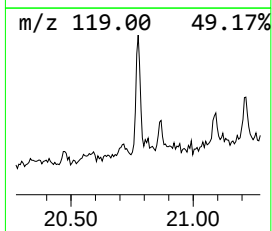
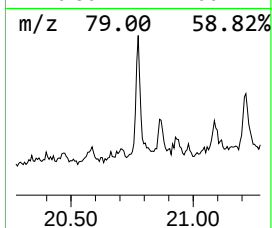
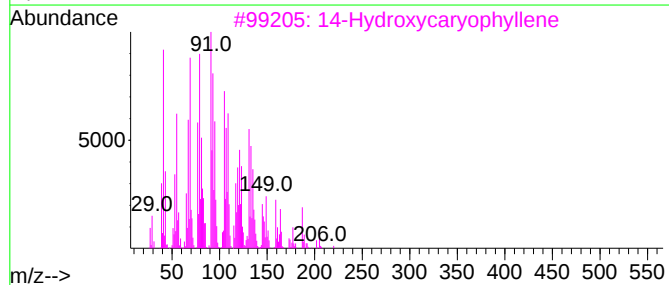
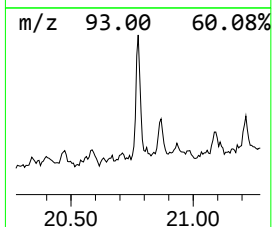
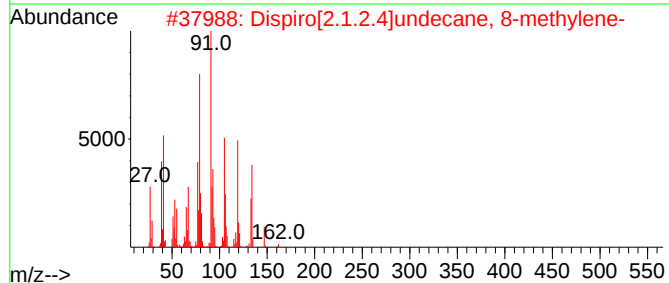
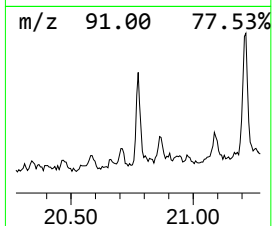
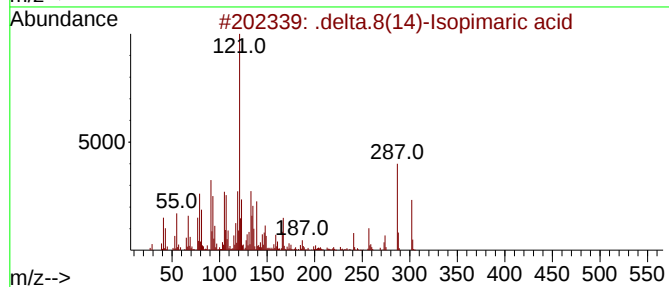
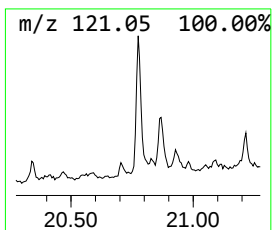
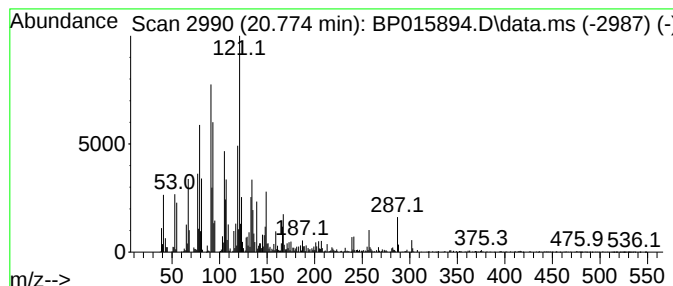
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 .delta.8(14)-Isopimaric acid Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.775	5.24 ng/ul	725519	Chrysene-d12	21.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.delta.8(14)-Isopimaric acid	302	C20H30O2	000471-74-9	93
2	Dispiro[2.1.2.4]undecane, 8-meth...	162	C12H18	051567-08-9	53
3	14-Hydroxycaryophyllene	220	C15H24O	050277-33-3	35
4	(+)-3-Bromocamphor-8-sulfonic acid	310	C10H15BrO4S	005344-58-1	35
5	Santolina triene	136	C10H16	002153-66-4	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
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Instrument :
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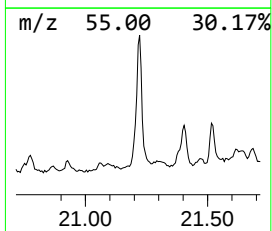
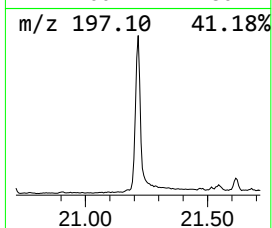
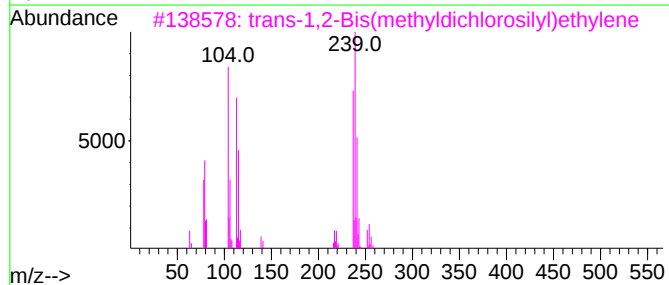
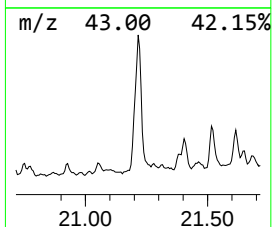
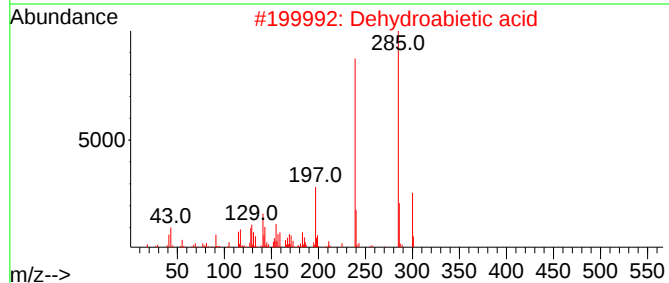
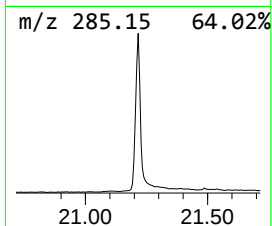
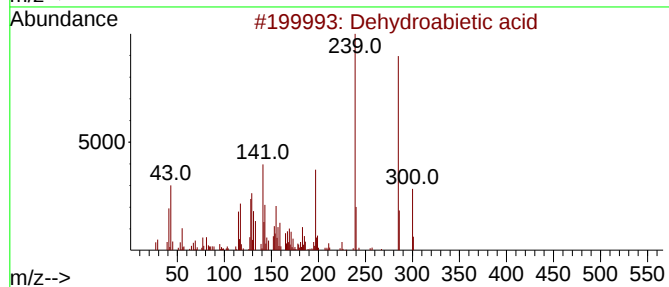
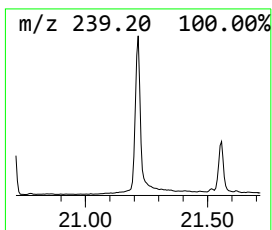
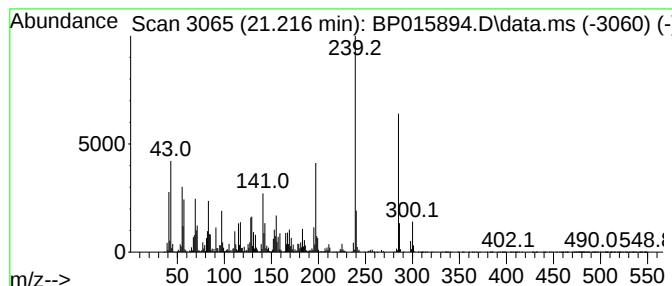
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 Dehydroabietic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.216	41.76 ng/ul	5787270	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dehydroabietic acid	300	C20H28O2	001740-19-8	99
2			Dehydroabietic acid	300	C20H28O2	001740-19-8	96
3			trans-1,2-Bis(methyldichlorosilyl)...	252	C4H8Cl4Si2	065899-10-7	93
4			Dehydroabietic acid	300	C20H28O2	001740-19-8	91
5			1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
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Misc :
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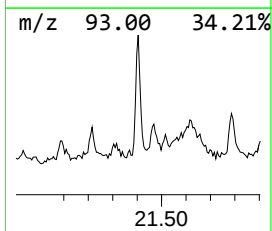
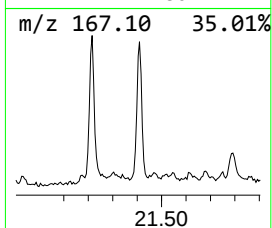
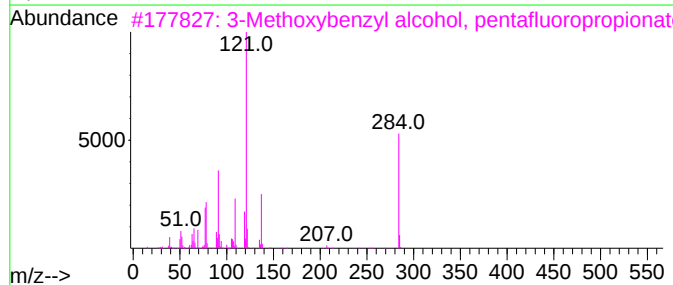
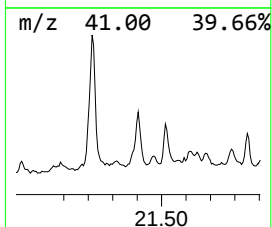
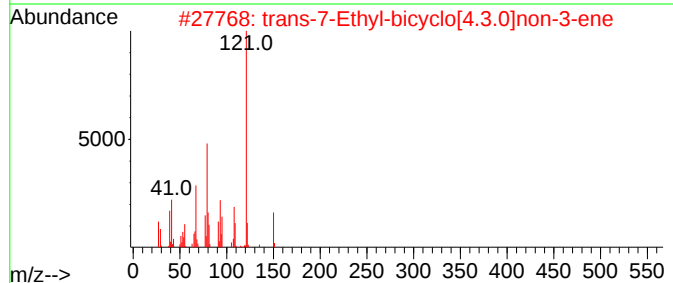
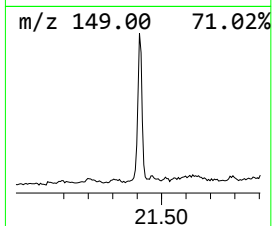
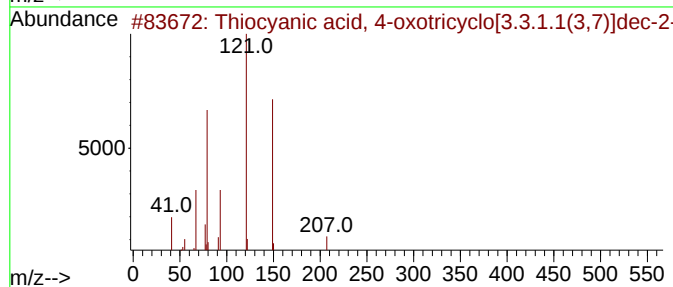
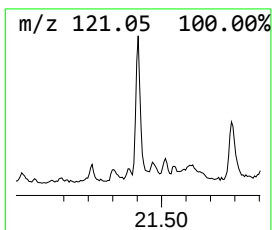
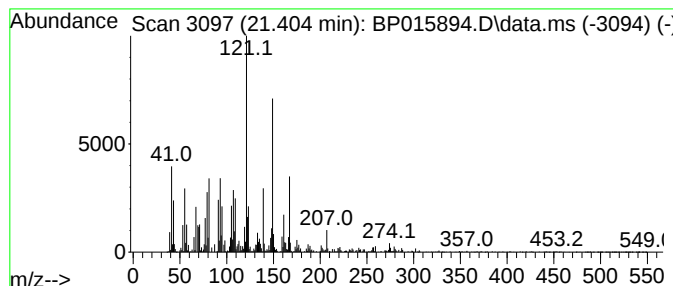
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 unknown-02 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.404	9.49 ng/ul	1315680	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiocyanic acid, 4-oxotricyclo[3... 207	C11H13NOS	056781-89-6	43	
2			trans-7-Ethyl-bicyclo[4.3.0]non-... 150	C11H18	1000145-84-7	35	
3			3-Methoxybenzyl alcohol, pentafl... 284	C11H9F5O3	1000365-49-8	30	
4			1,2,4-Oxadiazol-5-amine, 3-[4-(c... 328	C15H16N6O3	1000351-08-7	22	
5			9-t-Butyltricyclo[4.2.1.1(2,5)]d... 224	C14H24O2	1000186-30-2	22	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
ALS Vial : 62 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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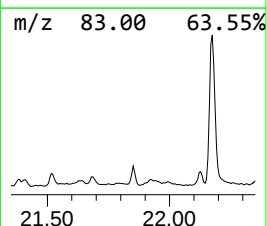
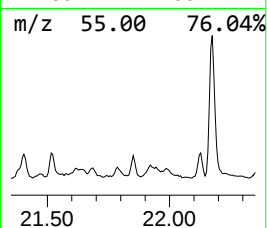
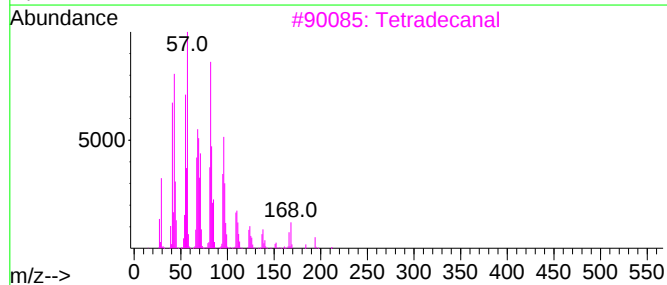
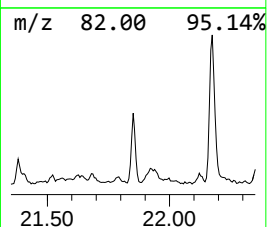
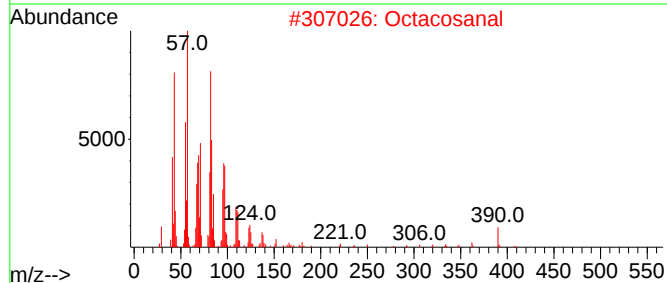
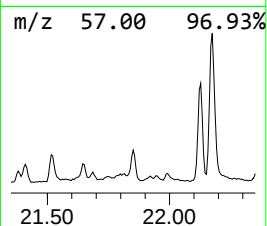
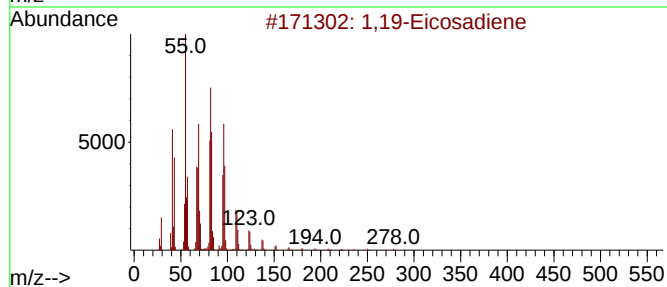
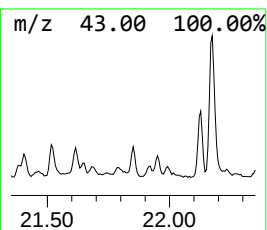
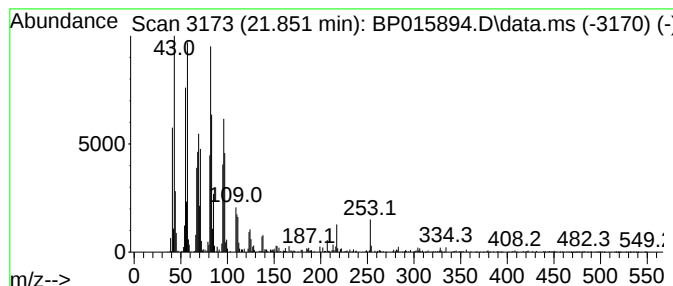
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 1,19-Eicosadiene Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.851	4.91 ng/ul	680279	Chrysene-d12	21.551

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,19-Eicosadiene		278	C20H38	014811-95-1	95
2	Octacosanal		408	C28H56O	022725-64-0	95
3	Tetradecanal		212	C14H28O	000124-25-4	95
4	Eicosanal-		296	C20H40O	002400-66-0	94
5	Hexacosanal		380	C26H52O	026627-85-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
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Instrument :
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ClientSampleId :
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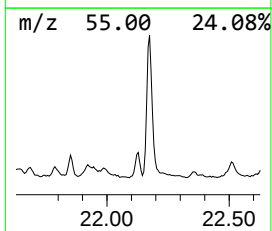
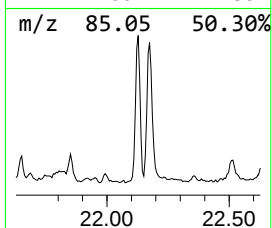
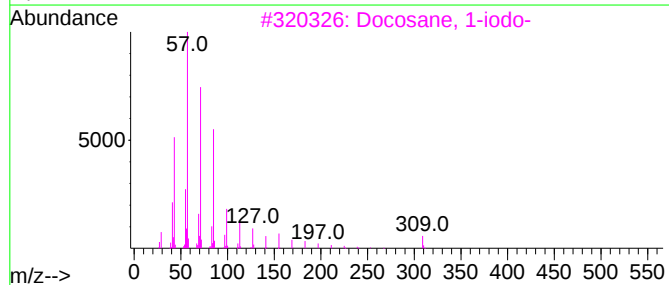
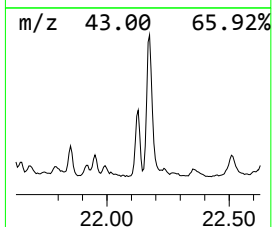
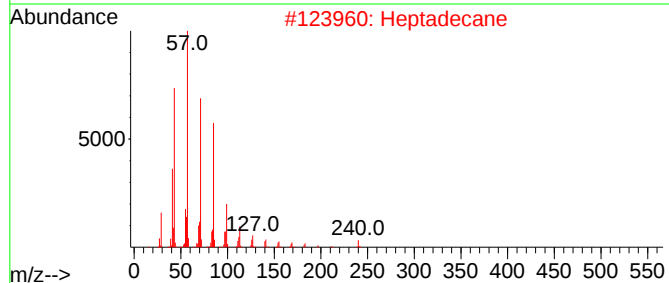
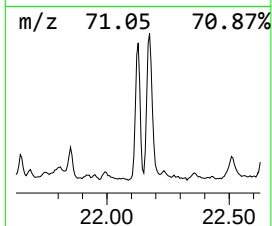
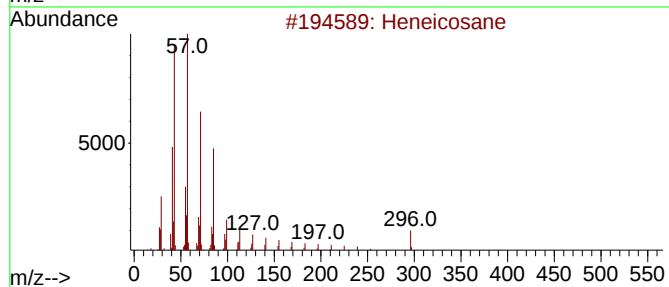
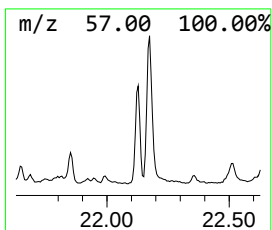
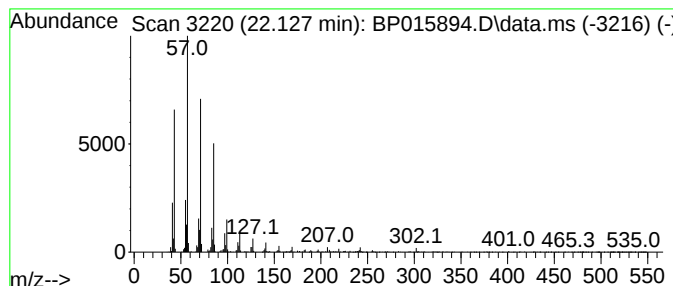
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.127	7.55 ng/ul	1046560	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	94
2			Heptadecane	240	C17H36	000629-78-7	93
3			Docosane, 1-iodo-	436	C22H45I	1000406-31-9	91
4			Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
5			Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
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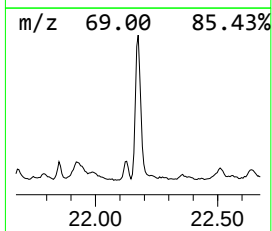
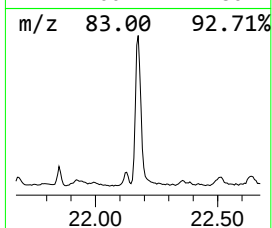
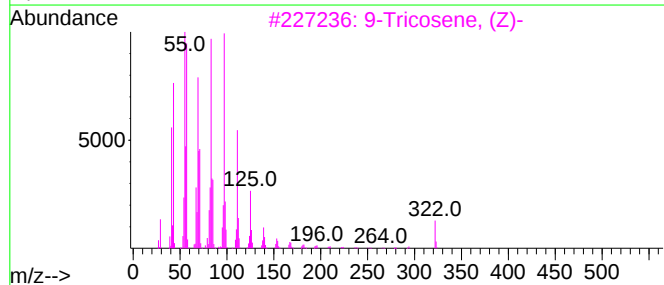
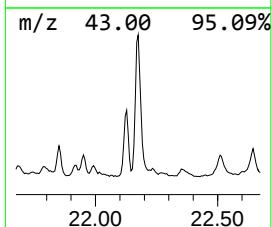
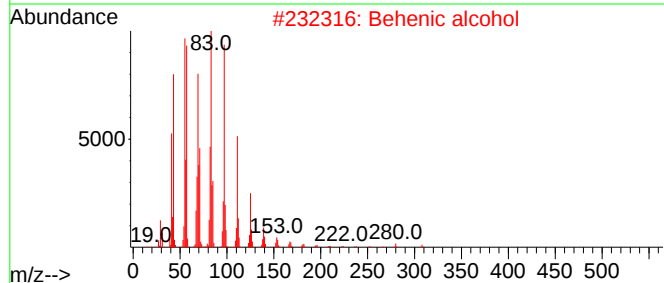
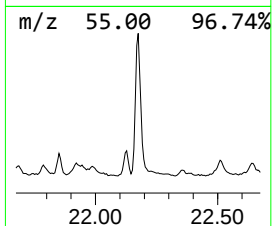
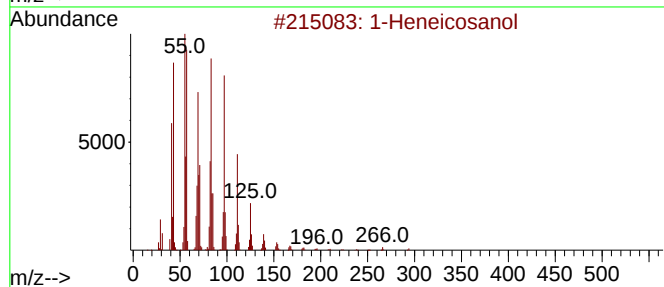
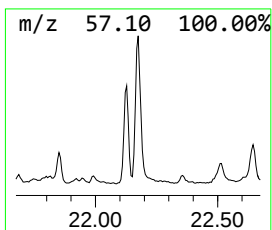
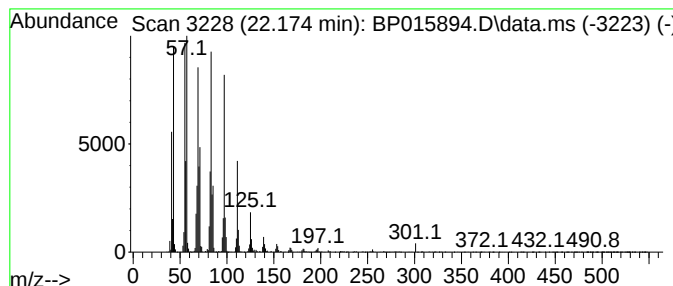
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 1-Heneicosanol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.174	30.34 ng/ul	4204400	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	95
2			Behenic alcohol	326	C22H46O	000661-19-8	94
3			9-Tricosene, (Z)-	322	C23H46	027519-02-4	94
4			n-Nonadecanol-1	284	C19H40O	001454-84-8	94
5			3-Eicosene, (E)-	280	C20H40	074685-33-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
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ClientSampleId :
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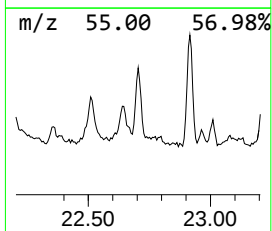
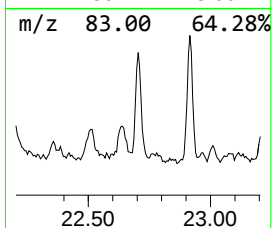
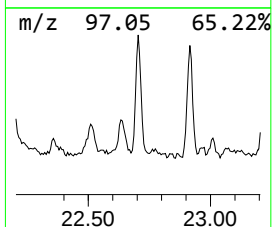
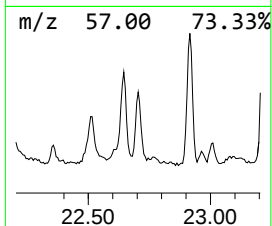
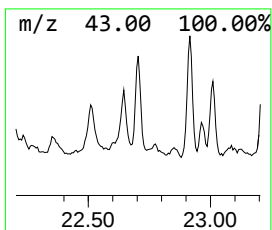
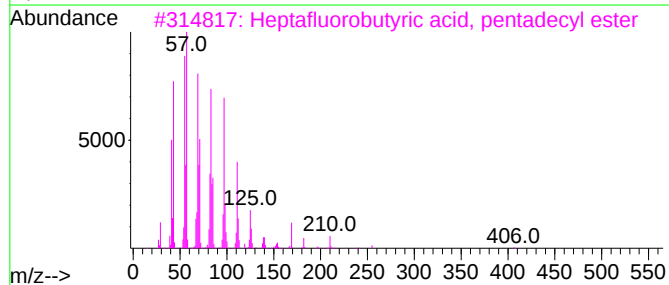
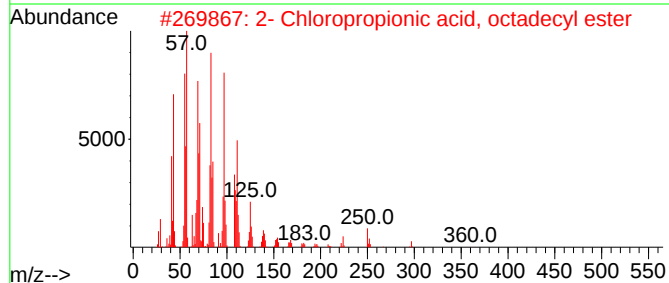
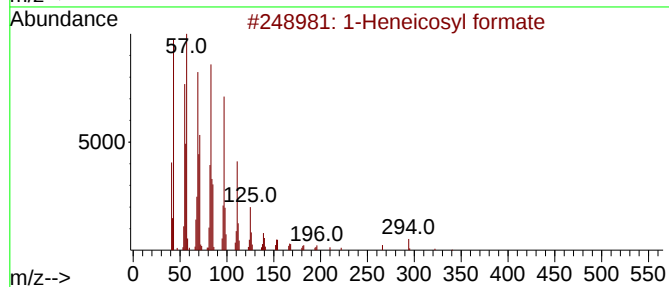
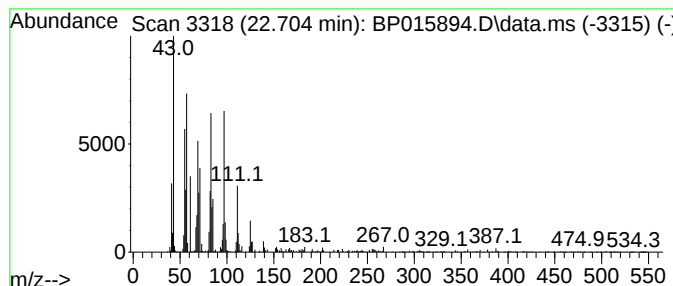
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 1-Heneicosyl formate Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.704	6.00 ng/ul	830868	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosyl formate	340	C22H44O2	077899-03-7	95
2			2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	91
3			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	91
4			Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91
5			Heptacos-1-ene	378	C27H54	015306-27-1	91



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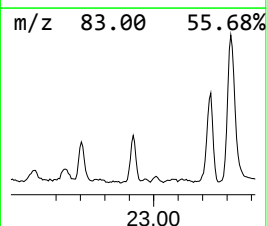
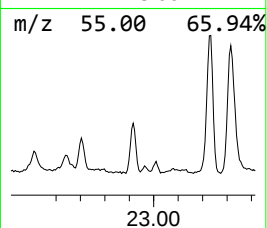
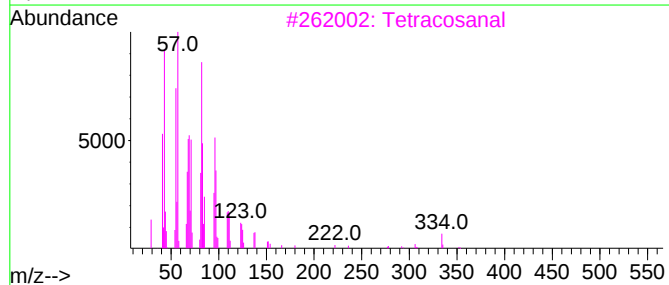
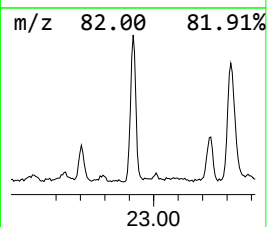
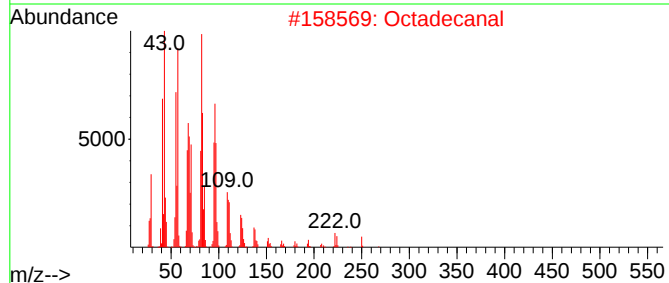
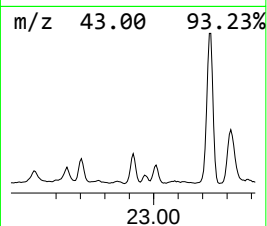
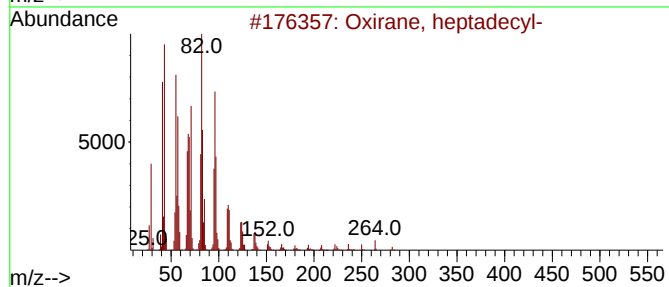
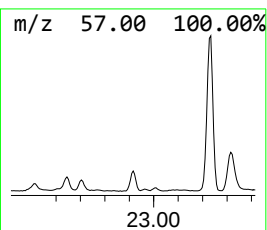
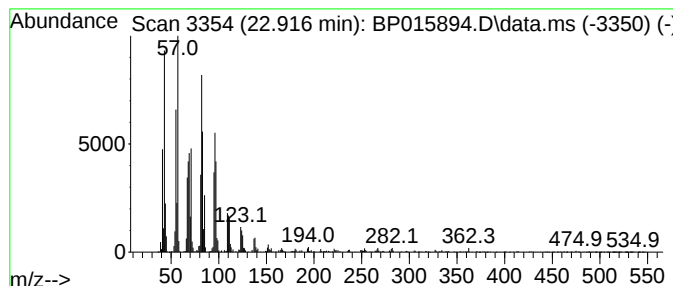
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 Oxirane, heptadecyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.916	8.82 ng/ul	1466240	Perylene-d12	24.098

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	97
2		Octadecanal	268	C18H36O	000638-66-4	94
3		Tetracosanal	352	C24H48O	057866-08-7	94
4		Octadecanal	268	C18H36O	000638-66-4	93
5		Octadecanal	268	C18H36O	000638-66-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
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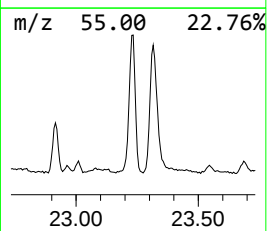
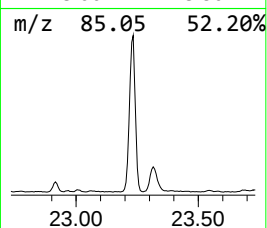
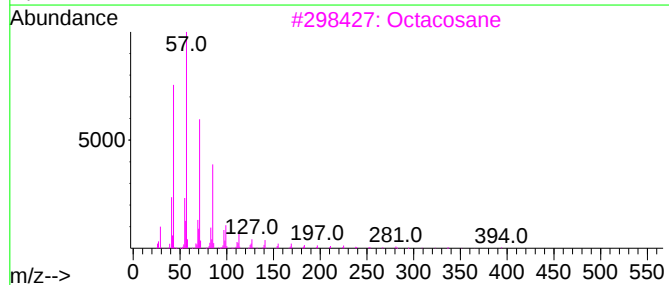
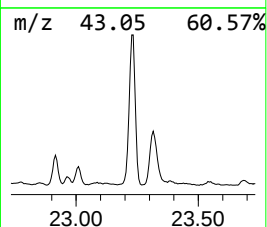
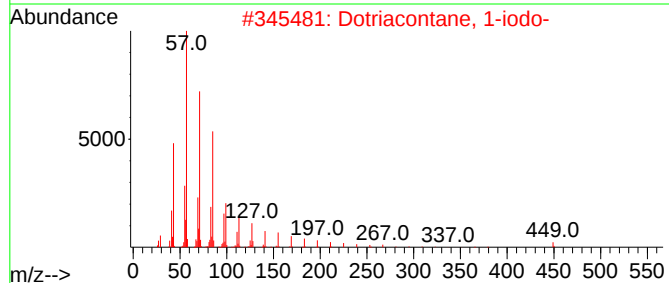
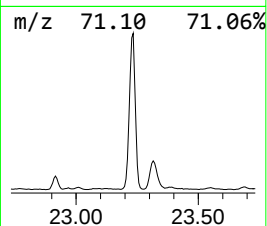
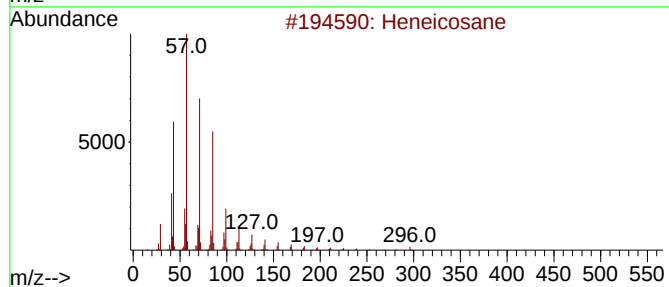
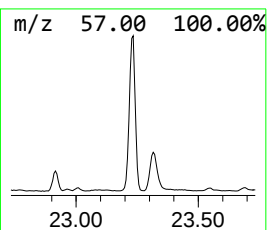
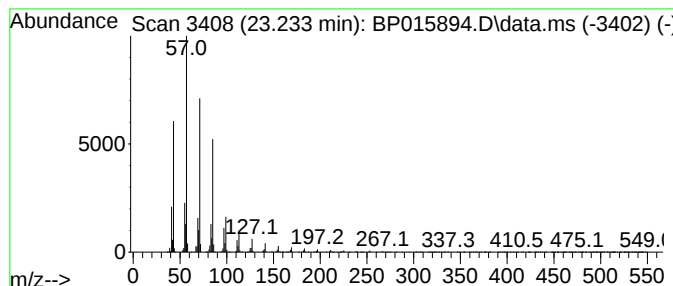
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Quant Title : SVOA CALIBRATION

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

Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.233	37.85 ng/ul	6293340	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
3			Octacosane	394	C28H58	000630-02-4	91
4			Hexatriacontane	507	C36H74	000630-06-8	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
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Acq On : 01 Jul 2023 21:03
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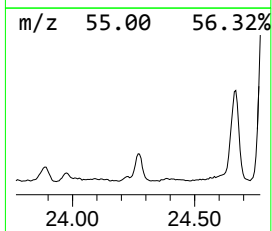
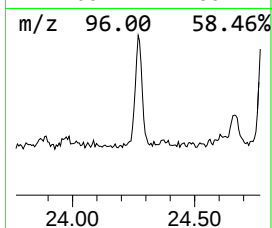
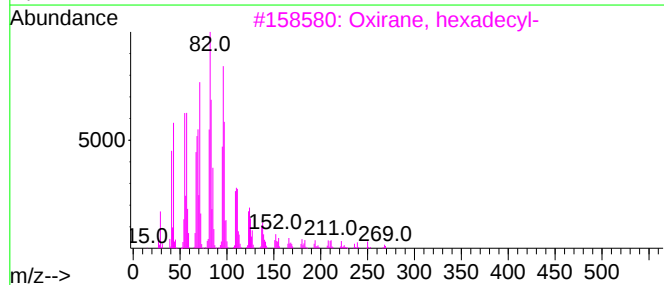
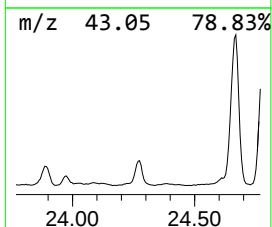
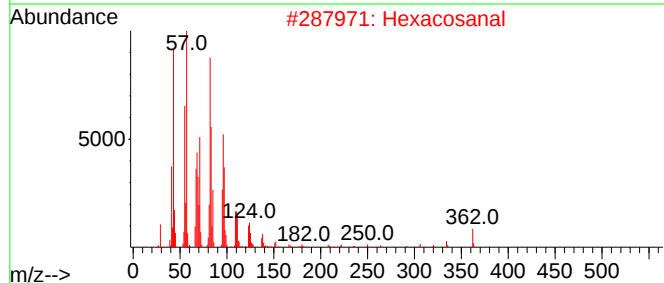
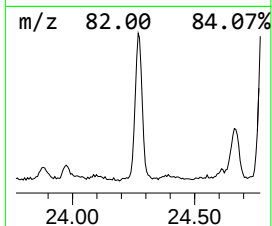
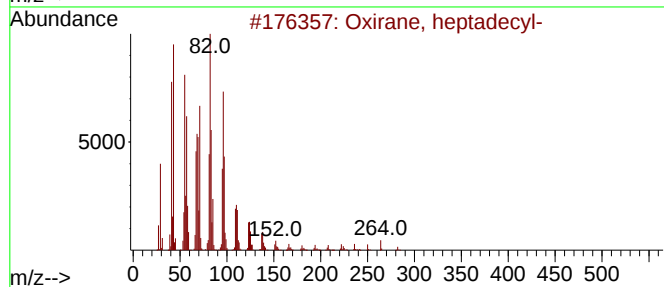
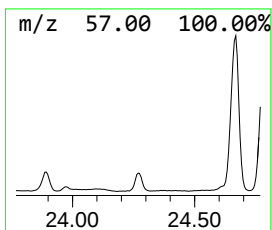
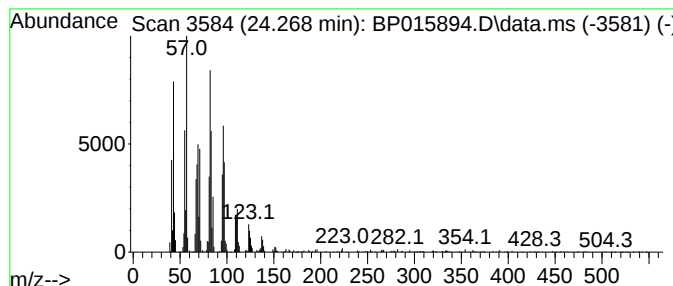
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Quant Title : SVOA CALIBRATION

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

Peak Number 33 Hexacosanal Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.268	10.17 ng/ul	1691830	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, heptadecyl-			282	C19H38O	067860-04-2	95
2	Hexacosanal			380	C26H52O	026627-85-0	94
3	Oxirane, hexadecyl-			268	C18H36O	007390-81-0	91
4	Octadecanal			268	C18H36O	000638-66-4	91
5	Heptadecanal			254	C17H34O	1000376-70-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
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ALS Vial : 62 Sample Multiplier: 1

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ClientSampleId :
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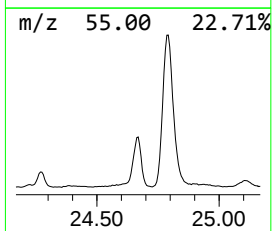
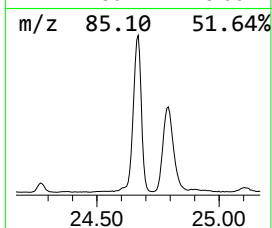
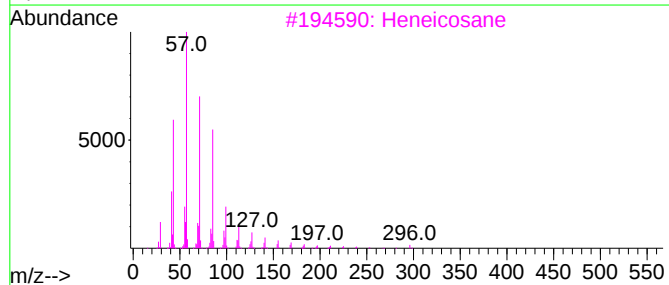
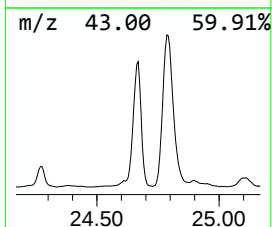
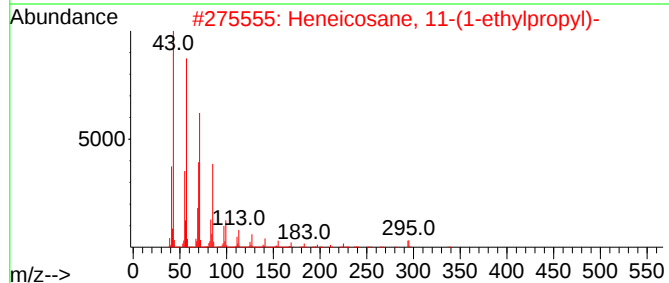
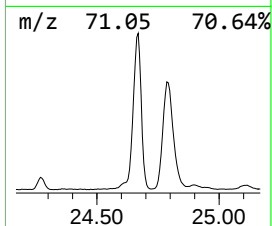
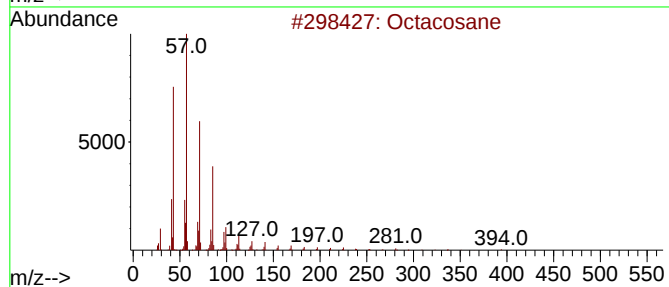
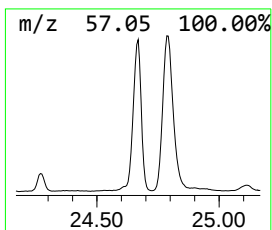
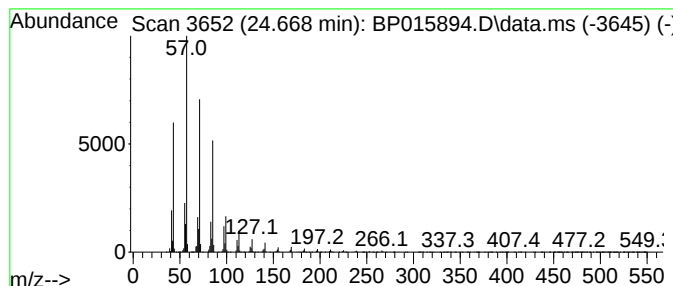
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Quant Title : SVOA CALIBRATION

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

Peak Number 34 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.668	49.35 ng/ul	8205660	Perylene-d12	24.098

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosane	394	C28H58	000630-02-4	91
2		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Hexatriacontane	507	C36H74	000630-06-8	91
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
ALS Vial : 62 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGB1

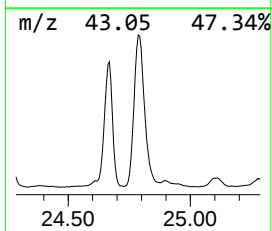
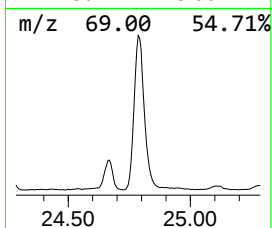
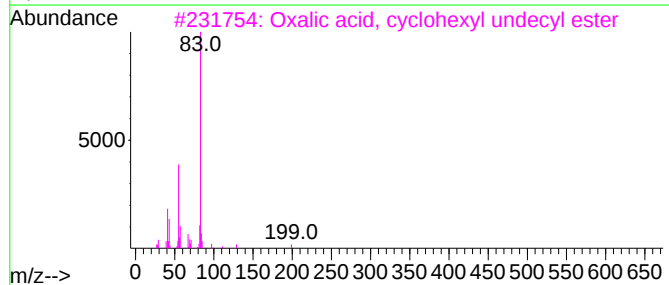
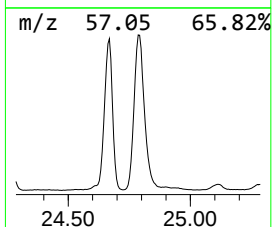
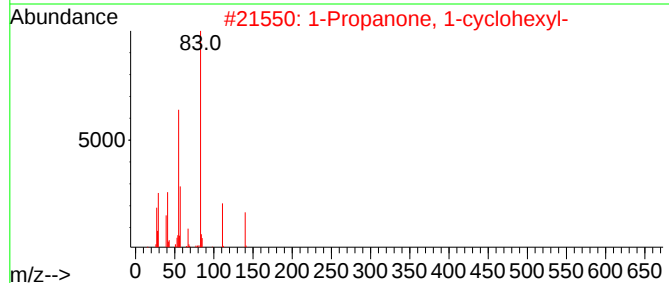
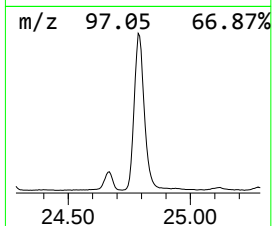
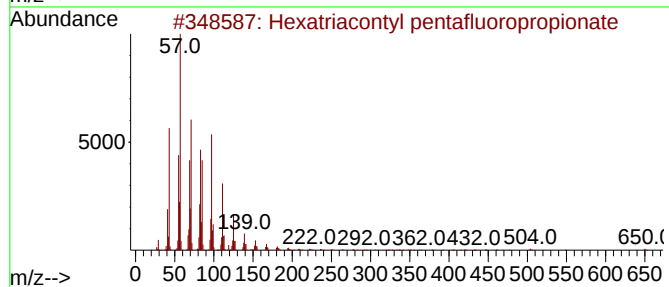
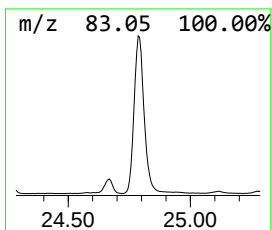
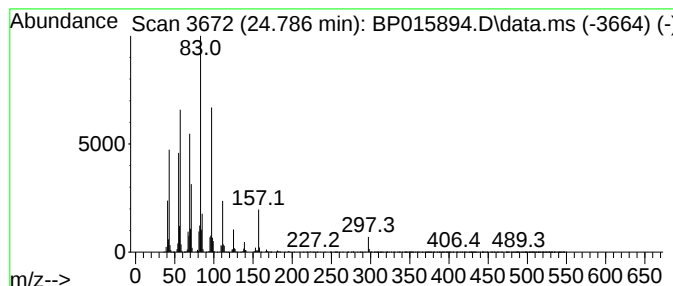
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Quant Title : SVOA CALIBRATION

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

Peak Number 35 unknown-03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.786	130.63 ng/ul	21721600	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexatriacontyl pentafluoropropio...	669	C39H73F5O2	1000351-89-0	38
2			1-Propanone, 1-cyclohexyl-	140	C9H16O	001123-86-0	38
3			Oxalic acid, cyclohexyl undecyl ...	326	C19H34O4	1000309-31-3	35
4			Octacosanol	410	C28H58O	000557-61-9	30
5			1-Heptacosanol	396	C27H56O	002004-39-9	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
ALS Vial : 62 Sample Multiplier: 1

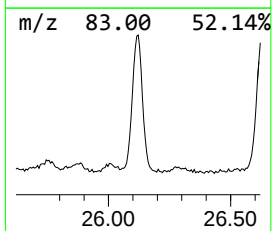
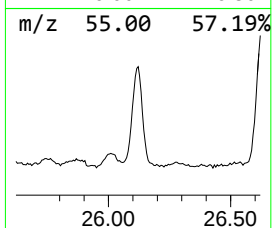
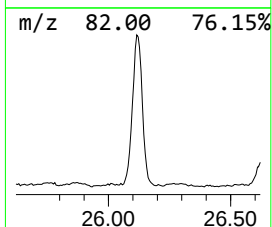
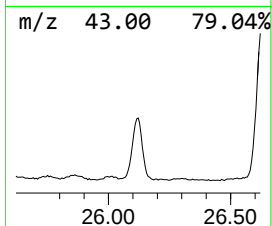
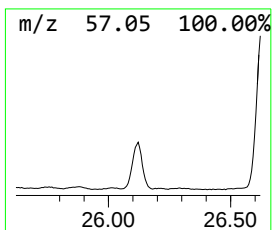
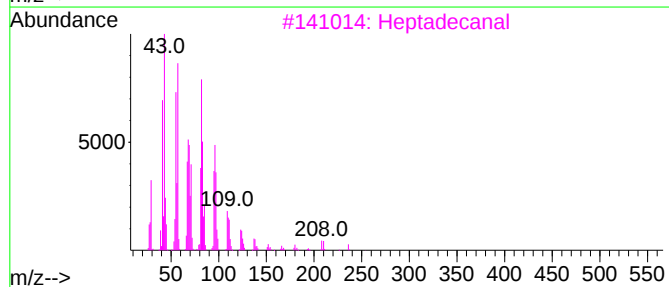
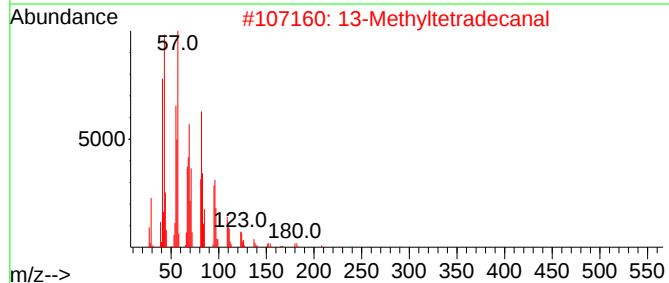
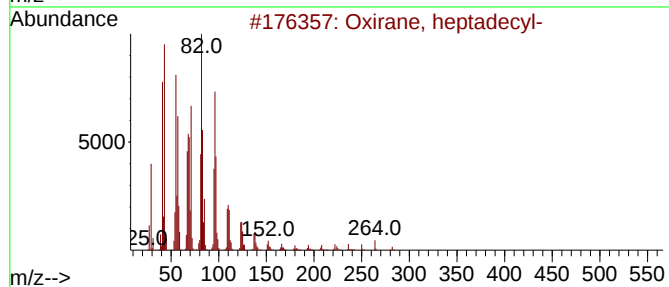
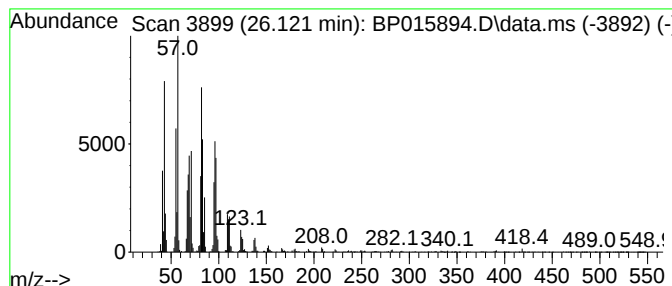
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 36 13-Methyltetradecanal Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
26.121	24.46 ng/ul	4067100	Perylene-d12	24.098	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	93
2	13-Methyltetradecanal	226	C15H30O	075853-51-9	91
3	Heptadecanal	254	C17H34O	1000376-70-0	90
4	Eicosanal-	296	C20H40O	002400-66-0	86
5	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
ALS Vial : 62 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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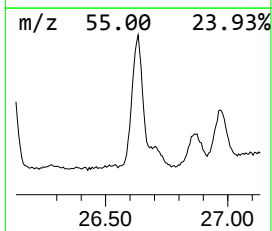
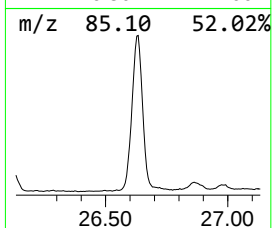
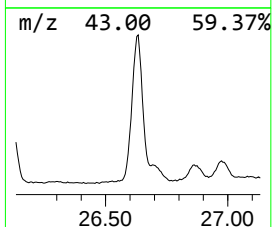
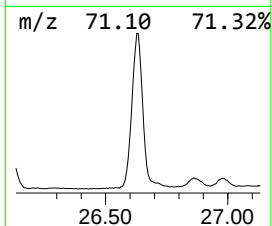
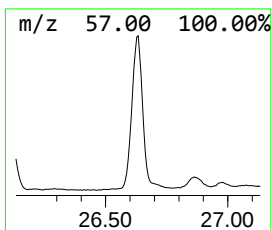
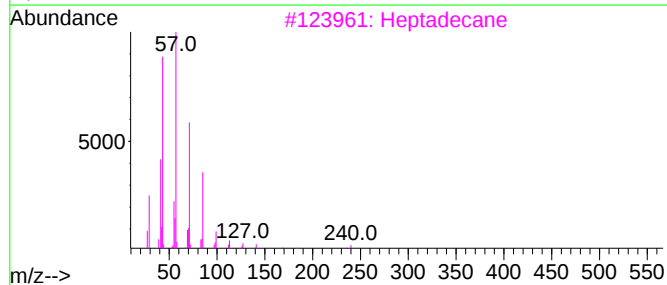
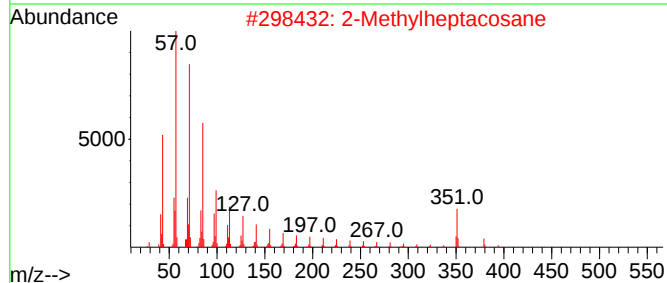
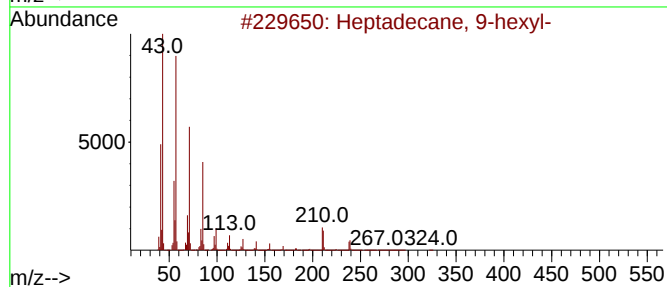
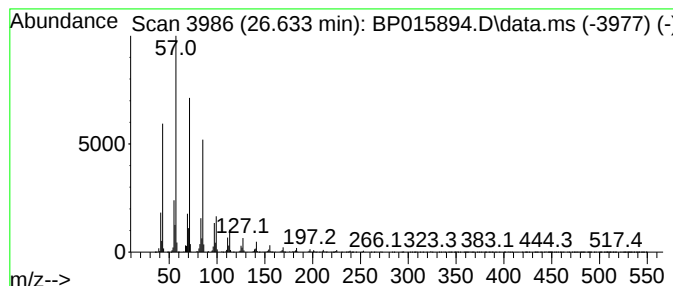
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Quant Title : SVOA CALIBRATION

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

Peak Number 37 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.633	45.91 ng/ul	7633840	Perylene-d12	24.098

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
2		2-Methylheptacosane	394	C28H58	001561-00-8	94
3		Heptadecane	240	C17H36	000629-78-7	93
4		Heneicosane	296	C21H44	000629-94-7	91
5		Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
ALS Vial : 62 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGB1

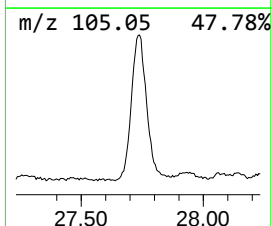
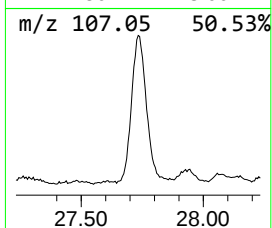
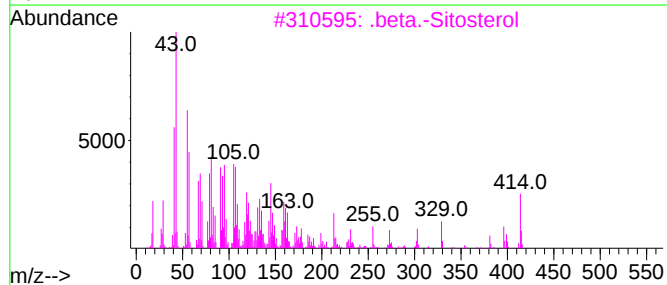
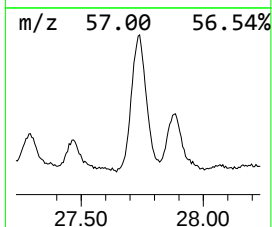
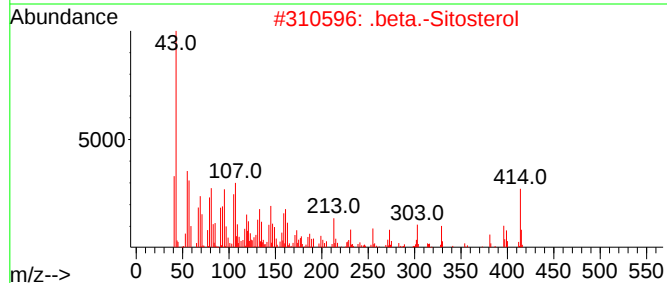
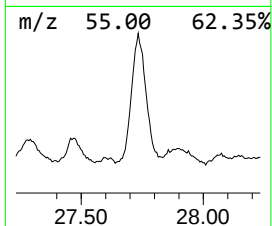
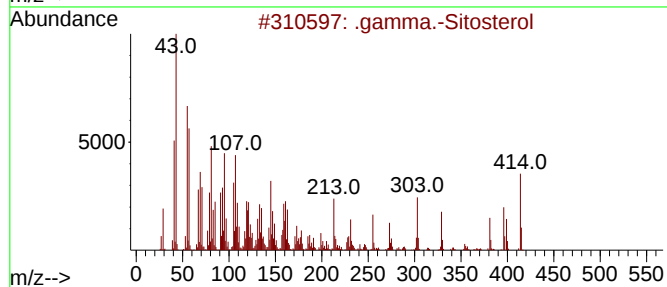
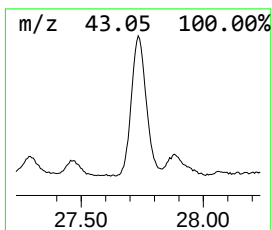
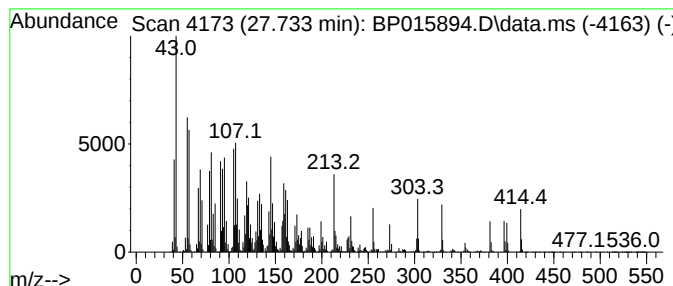
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 .gamma.-Sitosterol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.733	80.34 ng/ul	13360000	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	gamma.-Sitosterol	414	C29H50O	000083-47-6	99	
2	.	beta.-Sitosterol	414	C29H50O	000083-46-5	95	
3	.	beta.-Sitosterol	414	C29H50O	000083-46-5	90	
4	Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	83		
5	.	gamma.-Sitosterol	414	C29H50O	000083-47-6	83	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015894.D
Acq On : 01 Jul 2023 21:03
Operator : MA/JU
Sample : 03242-10
Misc :
ALS Vial : 62 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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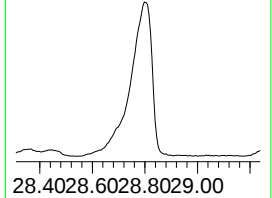
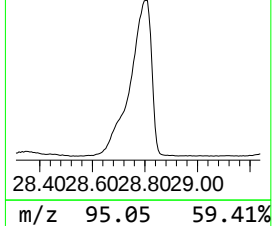
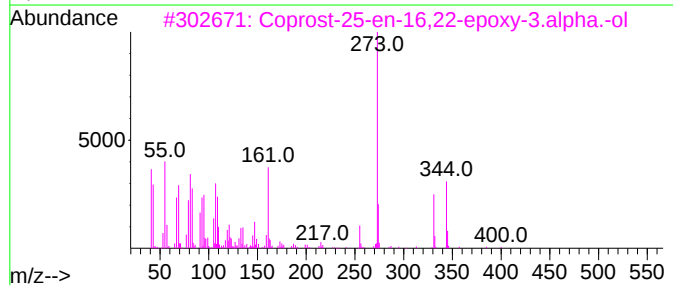
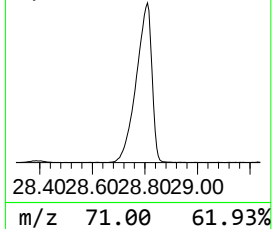
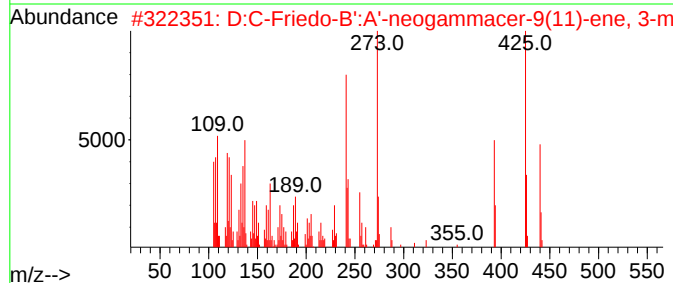
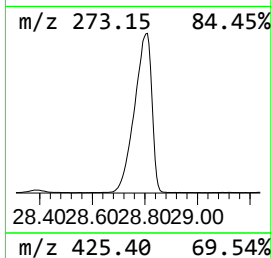
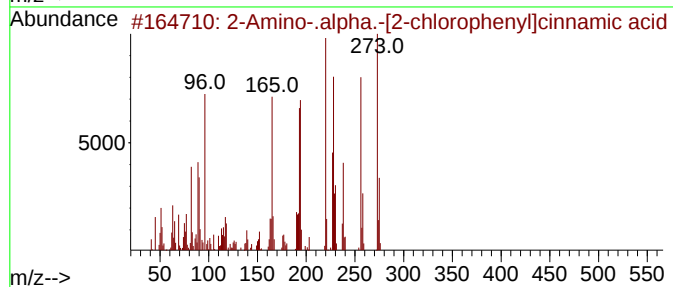
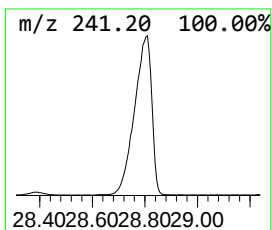
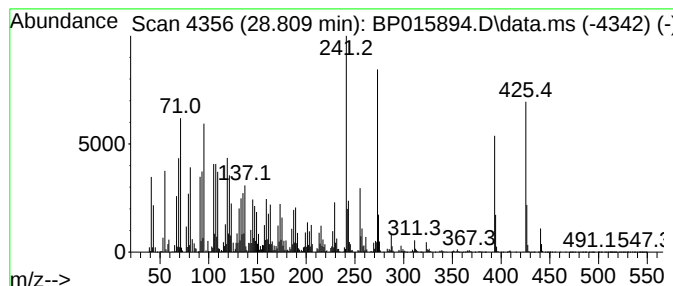
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 2-Amino-.alpha.-[2-chloroph... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.809	332.32 ng/ul	55261100	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Amino-.alpha.-[2-chlorophenyl]...	273	C15H12ClNO2	1000254-08-5	59
2			D:C-Friedo-B':A'-neogammacer-9(1...	440	C31H52O	004555-56-0	53
3			Coprost-25-en-16,22-epoxy-3.alpha...	400	C27H44O2	122337-93-3	30
4			Phenol, 2,6-dimethyl-, phosphite...	394	C24H27O3P	052830-49-6	25
5			Pyrazolo[1,5-a]pyrimidine, 6,7-d...	273	C18H15N3	186956-30-9	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
 Data File : BP015894.D
 Acq On : 01 Jul 2023 21:03
 Operator : MA/JU
 Sample : 03242-10
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGB1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.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
(1R)-2,6,6-Trim...	6.693	31.9	ng/ul	3095700	1	8.046	1939680	20.0
.beta.-Pinene	7.458	5.9	ng/ul	570685	1	8.046	1939680	20.0
trans-Verbenol	10.175	7.9	ng/ul	1109690	2	10.875	2814430	20.0
Bicyclo[3.1.1]h...	11.187	2.9	ng/ul	403205	2	10.875	2814430	20.0
Caryophyllene	13.998	3.6	ng/ul	549376	3	14.698	3080250	20.0
.gamma.-Muurolene	14.922	3.5	ng/ul	546074	3	14.698	3080250	20.0
unknown-01	15.116	2.8	ng/ul	437240	3	14.698	3080250	20.0
1H-Cycloprop[e]...	15.592	7.7	ng/ul	1180980	3	14.698	3080250	20.0
Benzophenone	16.063	4.7	ng/ul	715600	3	14.698	3080250	20.0
.tau.-Muurolol	16.151	5.8	ng/ul	798771	4	17.463	2767880	20.0
.alpha.-Cadinol	16.251	5.0	ng/ul	684705	4	17.463	2767880	20.0
Naphthalene, 1,...	16.404	3.0	ng/ul	416424	4	17.463	2767880	20.0
(E)-Hexadec-9-e...	18.257	4.9	ng/ul	673637	4	17.463	2767880	20.0
n-Hexadecanoic ...	18.322	40.4	ng/ul	5589510	4	17.463	2767880	20.0
9,12-Octadecadi...	19.404	8.0	ng/ul	1102340	4	17.463	2767880	20.0
Octadecanoic acid	19.539	18.3	ng/ul	2534430	5	21.551	2771560	20.0
(DEL) Alkane: S...	20.263	2.8	ng/ul	392666	5	21.551	2771560	20.0
Eicosanoic acid	20.586	5.4	ng/ul	744546	5	21.551	2771560	20.0
Methyl dehydroa...	20.710	3.4	ng/ul	472898	5	21.551	2771560	20.0
.delta.8(14)-Is...	20.775	5.2	ng/ul	725519	5	21.551	2771560	20.0
Dehydroabietic ...	21.216	41.8	ng/ul	5787270	5	21.551	2771560	20.0
unknown-02	21.404	9.5	ng/ul	1315680	5	21.551	2771560	20.0
1,19-Eicosadiene	21.851	4.9	ng/ul	680279	5	21.551	2771560	20.0
(DEL) Alkane: S...	22.127	7.5	ng/ul	1046560	5	21.551	2771560	20.0
1-Heneicosanol	22.174	30.3	ng/ul	4204400	5	21.551	2771560	20.0
1-Heneicosyl fo...	22.704	6.0	ng/ul	830868	5	21.551	2771560	20.0
Oxirane, heptad...	22.916	8.8	ng/ul	1466240	6	24.098	3325730	20.0
(DEL) Alkane: S...	23.233	37.9	ng/ul	6293340	6	24.098	3325730	20.0
Hexacosanal	24.268	10.2	ng/ul	1691830	6	24.098	3325730	20.0
(DEL) Alkane: S...	24.668	49.4	ng/ul	8205660	6	24.098	3325730	20.0
unknown-03	24.786	130.6	ng/ul	21721600	6	24.098	3325730	20.0
13-Methyltetrad...	26.121	24.5	ng/ul	4067100	6	24.098	3325730	20.0
(DEL) Alkane: S...	26.633	45.9	ng/ul	7633840	6	24.098	3325730	20.0
.gamma.-Sitosterol	27.733	80.3	ng/ul	13360000	6	24.098	3325730	20.0
2-Amino-.alpha....	28.809	332.3	ng/ul	55261100	6	24.098	3325730	20.0