

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
 Data File : BP013299.D
 Acq On : 24 Dec 2022 14:41
 Operator : CG/JU
 Sample : N6016-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBYC2

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-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP013299.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.328	20	24	34	rVB	172453	234180	1.05%	0.118%
2	3.764	95	98	111	rBV	632789	918605	4.11%	0.465%
3	3.864	111	115	119	rVB	132962	166617	0.75%	0.084%
4	3.987	130	136	144	rVB	105478	162367	0.73%	0.082%
5	4.087	149	153	157	rVB	51812	67463	0.30%	0.034%
6	4.652	245	249	253	rVB	38697	47150	0.21%	0.024%
7	4.699	253	257	264	rVB	130685	171850	0.77%	0.087%
8	5.022	305	312	321	rBV3	379069	738804	3.31%	0.374%
9	5.117	324	328	336	rVB	40853	60231	0.27%	0.030%
10	6.540	565	570	576	rVB3	40901	67811	0.30%	0.034%
11	6.611	576	582	588	rBV	414530	608933	2.73%	0.308%
12	7.105	654	666	679	rBV	2439470	4110355	18.40%	2.079%
13	7.269	683	694	706	rVV	2022708	3242838	14.52%	1.640%
14	7.369	707	711	721	rVB2	32431	58229	0.26%	0.029%
15	7.469	721	728	738	rBV	2592798	4029072	18.03%	2.037%
16	7.940	800	808	815	rBV	2478270	3860003	17.28%	1.952%
17	8.652	922	929	939	rBV	2597152	4122443	18.45%	2.085%
18	8.769	943	949	957	rVB	231374	371954	1.66%	0.188%
19	9.110	999	1007	1015	rBV	2342540	3852137	17.24%	1.948%
20	9.505	1069	1074	1084	rVB2	76200	123038	0.55%	0.062%
21	9.834	1120	1130	1145	rBV	1979385	3290576	14.73%	1.664%
22	9.987	1149	1156	1162	rVV	70303	134408	0.60%	0.068%
23	10.052	1162	1167	1173	rVB5	34942	64836	0.29%	0.033%
24	10.193	1184	1191	1197	rVB2	41229	71716	0.32%	0.036%
25	10.375	1214	1222	1242	rBV	2875259	4924861	22.04%	2.490%
26	10.752	1278	1286	1294	rBV	3081032	5238475	23.45%	2.649%
27	10.893	1302	1310	1331	rBV	1770036	3159115	14.14%	1.598%
28	11.287	1371	1377	1385	rBV8	32297	73137	0.33%	0.037%
29	11.546	1415	1421	1426	rBV3	22833	48151	0.22%	0.024%
30	11.675	1436	1443	1449	rBV4	20901	49528	0.22%	0.025%
31	12.340	1550	1556	1562	rBV2	49838	83303	0.37%	0.042%
32	12.946	1654	1659	1667	rVB9	32153	65482	0.29%	0.033%
33	13.340	1717	1726	1731	rBV4	50982	114477	0.51%	0.058%
34	13.398	1731	1736	1741	rVB	73116	104927	0.47%	0.053%
35	13.457	1741	1746	1752	rBV3	111853	184601	0.83%	0.093%
36	13.904	1817	1822	1829	rVB2	167995	247818	1.11%	0.125%

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37	13.993	1829	1837	1846	rBV	4747727	6435650	28.81%	3.254%
38	14.281	1879	1886	1893	rVV	5285246	8019936	35.90%	4.056%
39	14.340	1893	1896	1902	rVV3	109377	171141	0.77%	0.087%
40	14.422	1905	1910	1914	rVV	85659	139335	0.62%	0.070%
41	14.469	1914	1918	1923	rVV2	68910	103162	0.46%	0.052%
42	14.540	1925	1930	1933	rVV	591526	825606	3.70%	0.417%
43	14.587	1933	1938	1944	rVV	4023009	5946382	26.62%	3.007%
44	14.645	1944	1948	1960	rVB5	113070	276285	1.24%	0.140%
45	14.787	1966	1972	1977	rBV	1528554	2350046	10.52%	1.188%
46	14.828	1977	1979	1983	rVV	242174	380190	1.70%	0.192%
47	14.893	1987	1990	1995	rVV	121752	180231	0.81%	0.091%
48	14.934	1995	1997	2002	rVB2	60089	80215	0.36%	0.041%
49	14.998	2002	2008	2012	rBV6	60352	114462	0.51%	0.058%
50	15.581	2100	2107	2113	rBV	6410415	9219369	41.27%	4.662%
51	15.698	2122	2127	2137	rVB	1473477	1953392	8.74%	0.988%
52	15.822	2143	2148	2153	rVB3	76476	121505	0.54%	0.061%
53	15.910	2158	2163	2166	rVV5	64089	116644	0.52%	0.059%
54	15.945	2166	2169	2173	rVV	85429	117041	0.52%	0.059%
55	15.998	2174	2178	2179	rVV4	46091	53099	0.24%	0.027%
56	16.028	2179	2183	2184	rVV3	56701	84060	0.38%	0.043%
57	16.057	2184	2188	2195	rVB2	155744	258657	1.16%	0.131%
58	16.151	2199	2204	2210	rBV3	96554	142620	0.64%	0.072%
59	16.304	2224	2230	2234	rVB2	46553	84178	0.38%	0.043%
60	16.434	2247	2252	2256	rBV4	62707	105001	0.47%	0.053%
61	16.475	2256	2259	2269	rVB5	38136	80318	0.36%	0.041%
62	16.728	2298	2302	2306	rBV3	39310	46374	0.21%	0.023%
63	17.222	2382	2386	2391	rBV5	47748	72693	0.33%	0.037%
64	17.292	2394	2398	2401	rBV5	32836	52472	0.23%	0.027%
65	17.345	2401	2407	2413	rVV	4605244	6456857	28.90%	3.265%
66	17.445	2417	2424	2431	rBV	6337639	9082329	40.65%	4.593%
67	17.534	2437	2439	2446	rVB7	31568	55570	0.25%	0.028%
68	17.722	2468	2471	2480	rVB6	55143	102911	0.46%	0.052%
69	17.998	2515	2518	2521	rBV4	46291	51826	0.23%	0.026%
70	18.098	2530	2535	2540	rBV6	65775	112480	0.50%	0.057%
71	18.157	2541	2545	2549	rBV2	103108	135062	0.60%	0.068%
72	18.216	2550	2555	2565	rBV	1471133	2162394	9.68%	1.093%
73	18.428	2587	2591	2595	rVB2	114295	138885	0.62%	0.070%
74	18.539	2605	2610	2620	rVV3	626563	1013449	4.54%	0.512%
75	18.622	2620	2624	2631	rVB	678275	816629	3.66%	0.413%
76	18.992	2685	2687	2692	rVB	44408	49448	0.22%	0.025%

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77	19.045	2693	2696	2701	rBV4	75542	116560	0.52%	0.059%
78	19.104	2703	2706	2709	rBV4	38097	46510	0.21%	0.024%
79	19.210	2720	2724	2727	rBV4	95001	148396	0.66%	0.075%
80	19.251	2728	2731	2737	rVB4	138655	202804	0.91%	0.103%
81	19.333	2737	2745	2758	rBV3	448239	1268519	5.68%	0.641%
82	19.445	2760	2764	2778	rVV	435487	691558	3.10%	0.350%
83	19.675	2798	2803	2817	rVB	8970756	12575270	56.29%	6.359%
84	19.828	2825	2829	2833	rBV3	208272	268949	1.20%	0.136%
85	20.051	2864	2867	2871	rVB	314194	354008	1.58%	0.179%
86	20.192	2887	2891	2899	rVB5	320438	543446	2.43%	0.275%
87	20.669	2969	2972	2976	rBV2	235834	267623	1.20%	0.135%
88	20.975	3020	3024	3031	rBV	4521594	4982345	22.30%	2.519%
89	21.127	3046	3050	3059	rBV	2038190	2989577	13.38%	1.512%
90	21.433	3097	3102	3115	rVB	6263201	8297083	37.14%	4.196%
91	22.039	3202	3205	3207	rBV	464765	556160	2.49%	0.281%
92	22.069	3208	3210	3221	rVB2	664242	1098158	4.92%	0.555%
93	22.551	3290	3292	3298	rVB	471248	600024	2.69%	0.303%
94	22.698	3315	3317	3322	rVB	409820	559033	2.50%	0.283%
95	23.127	3385	3390	3396	rBV	3301029	5231532	23.42%	2.645%
96	23.757	3490	3497	3509	rVB2	6928035	14463508	64.74%	7.314%
97	23.916	3517	3524	3539	rVB2	4299053	9038136	40.46%	4.570%
98	24.533	3623	3629	3640	rVB	1792715	3906610	17.49%	1.976%
99	26.439	3948	3953	3974	rVB	1366947	4628647	20.72%	2.341%
100	28.504	4293	4304	4329	rVB2	5406016	22340826	100.00%	11.297%

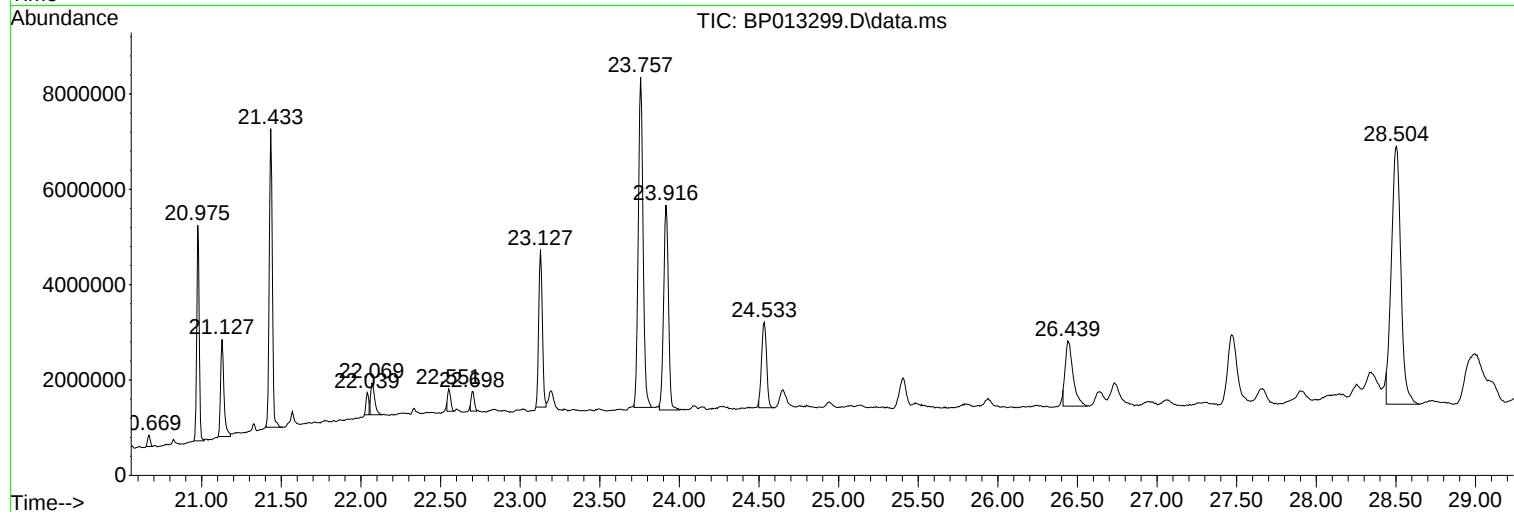
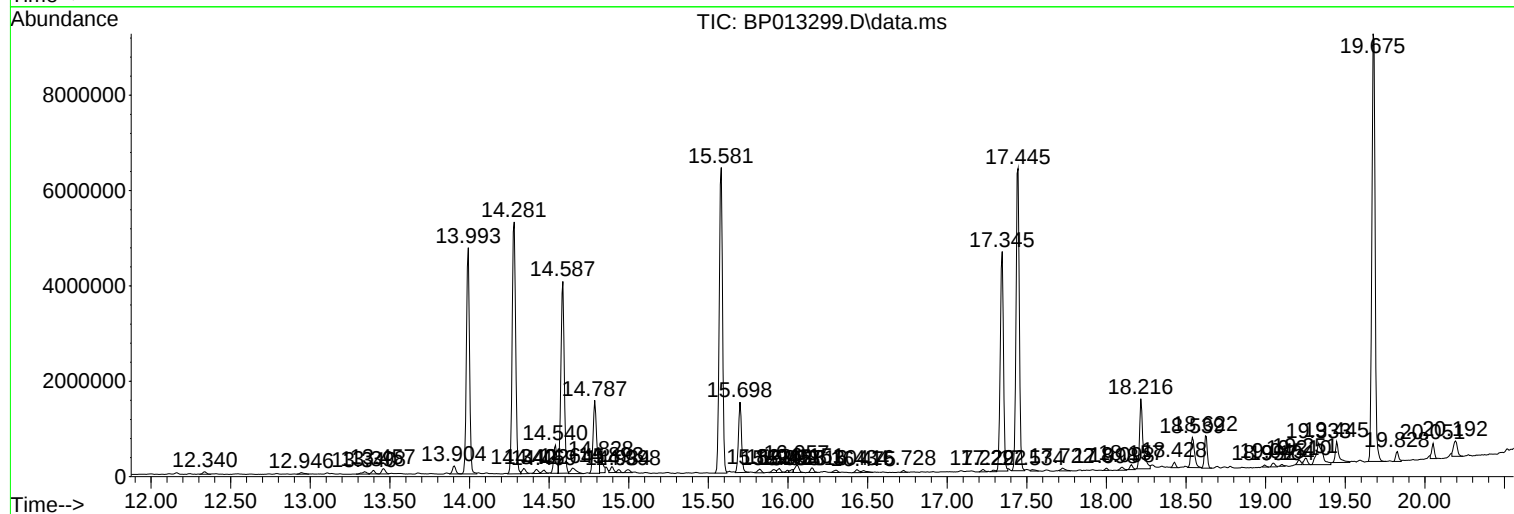
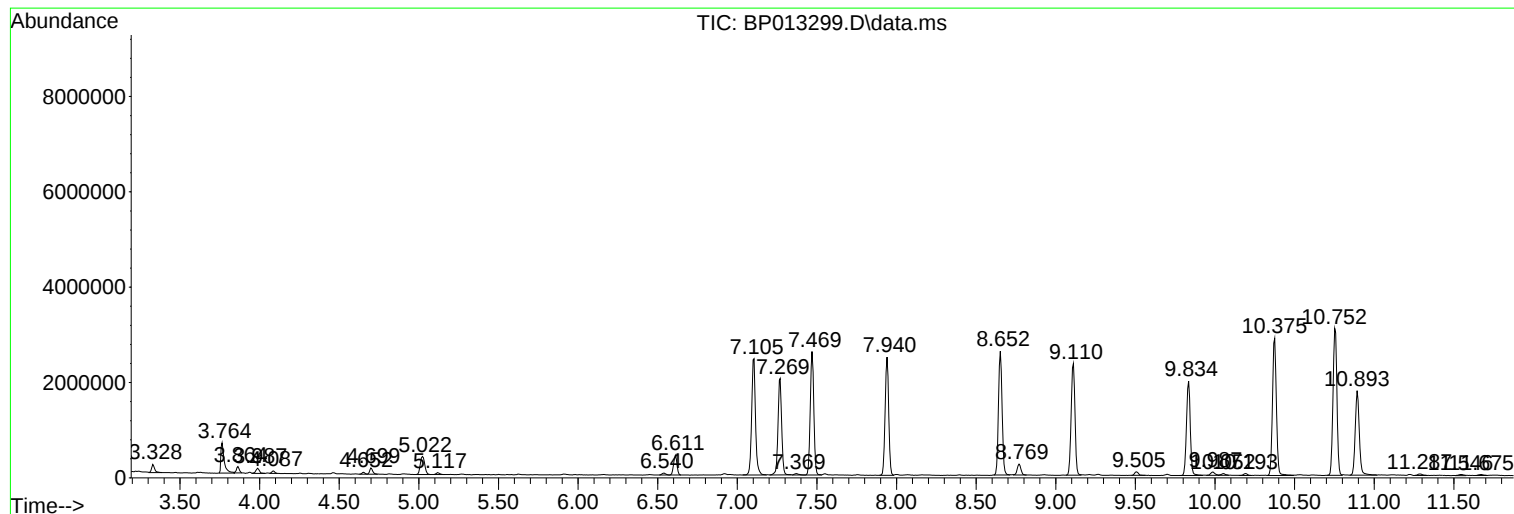
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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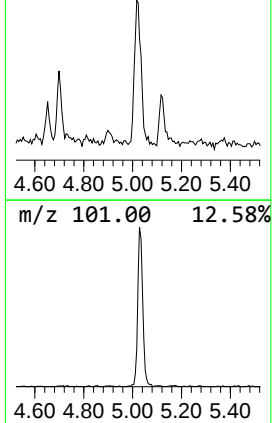
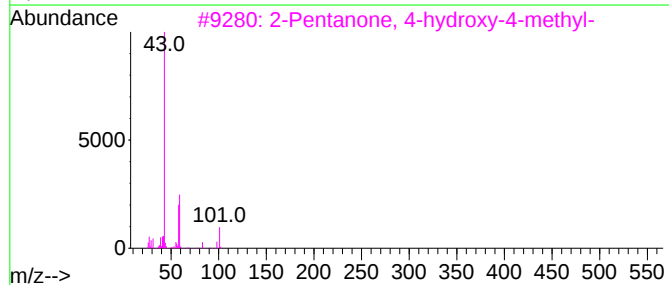
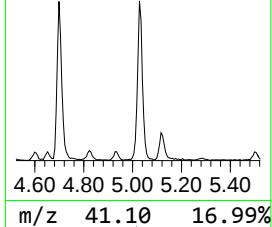
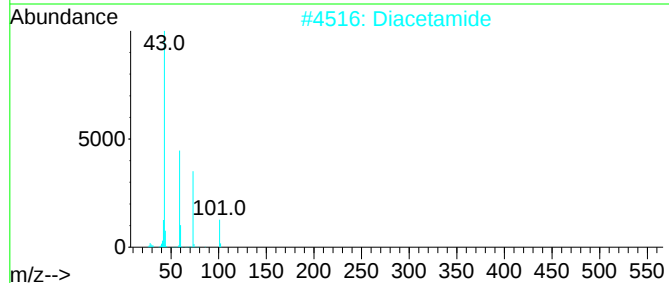
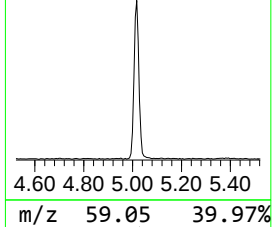
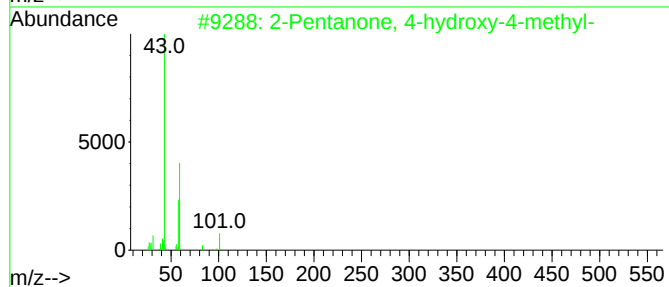
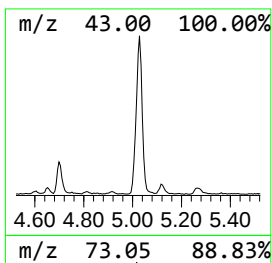
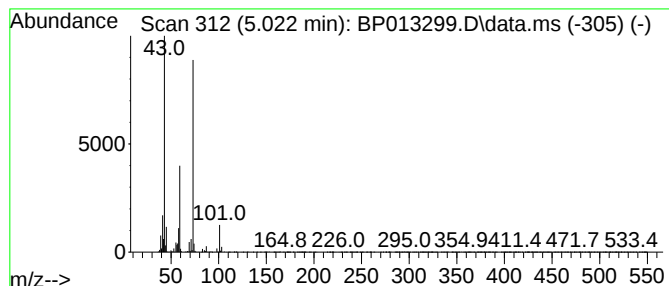
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.022	3.83 ng/ul	738804	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		
2	Diacetamide	101	C4H7NO2	000625-77-4	38		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	37		
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28		
5	Propylamine	59	C3H9N	000107-10-8	25		



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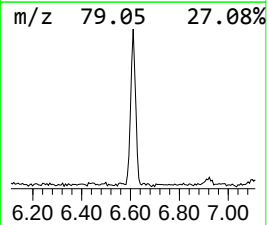
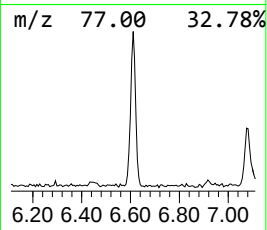
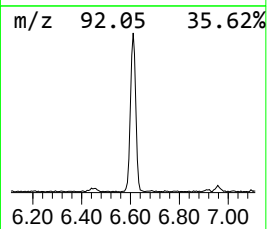
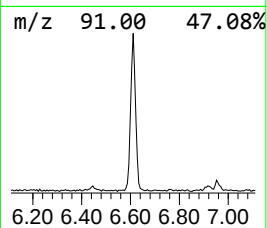
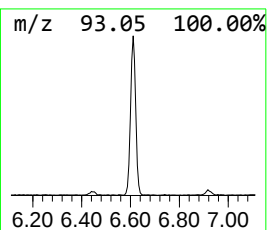
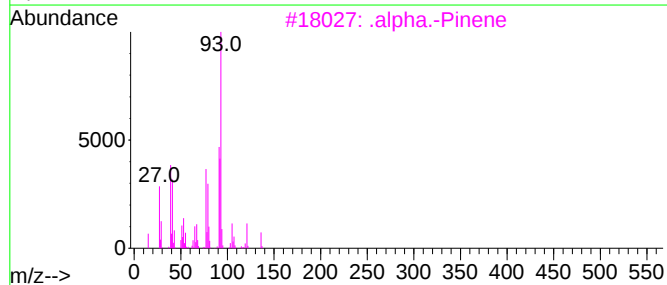
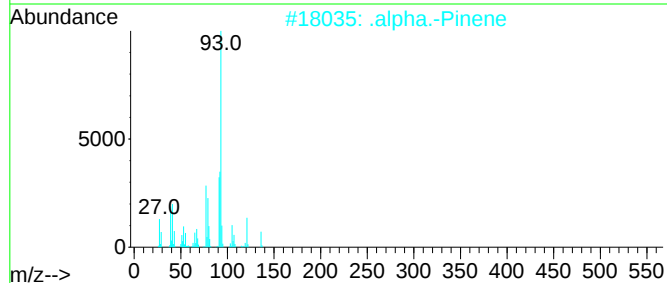
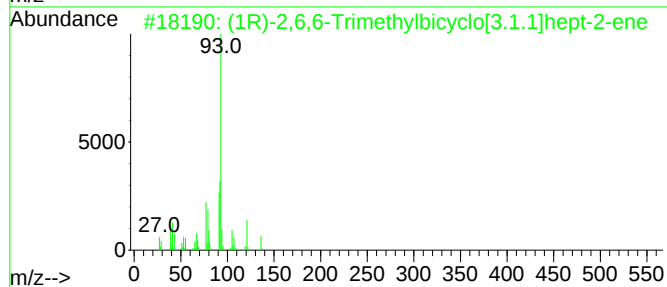
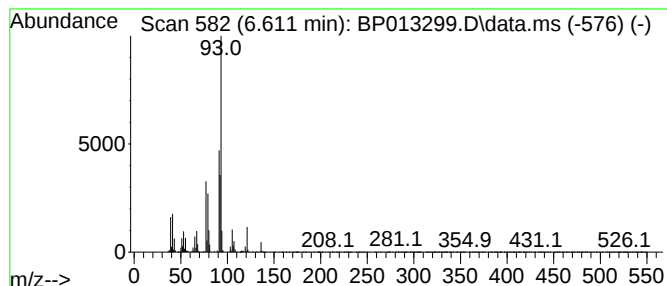
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.611	3.16 ng/ul	608933	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-70-8	96		
2	.alpha.-Pinene	136	C10H16	000080-56-8	95		
3	.alpha.-Pinene	136	C10H16	000080-56-8	95		
4	3-Carene	136	C10H16	013466-78-9	91		
5	(1S)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-26-4	91		



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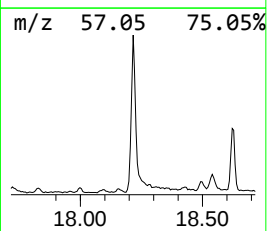
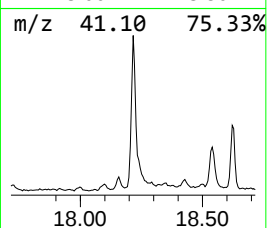
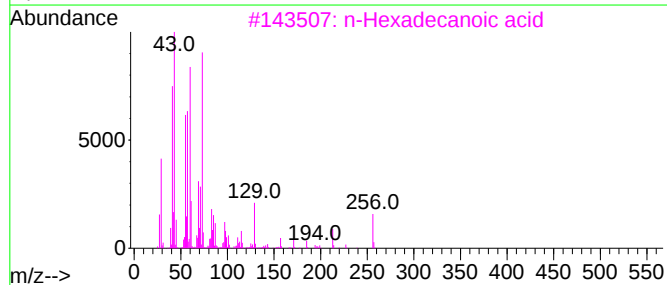
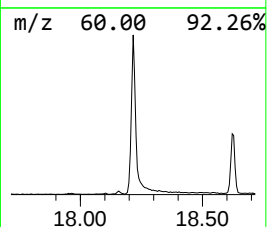
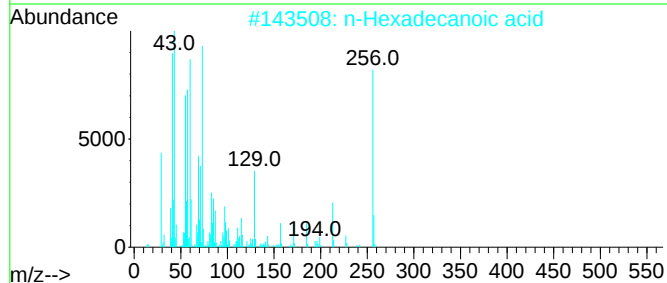
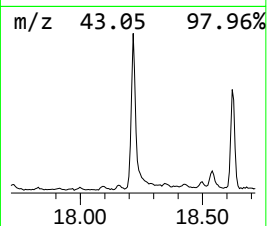
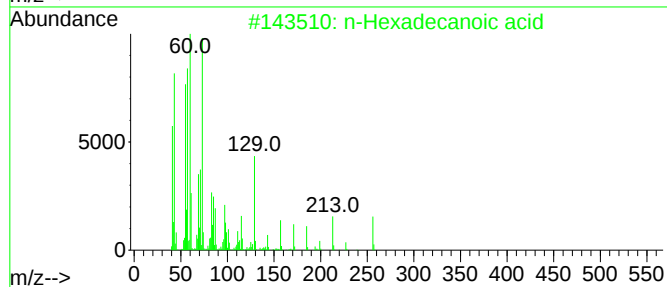
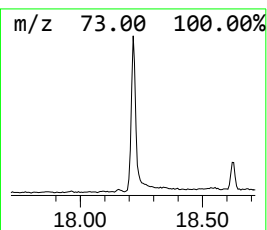
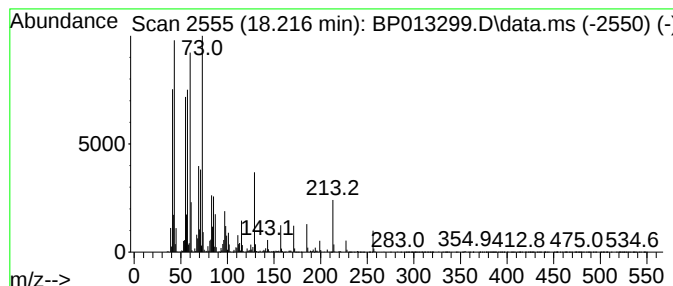
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	6.70 ng/ul	2162390	Phenanthrene-d10	17.345

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	93
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
5			Tridecanoic acid	214	C13H26O2	000638-53-9	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013299.D
Acq On : 24 Dec 2022 14:41
Operator : CG/JU
Sample : N6016-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DBYC2

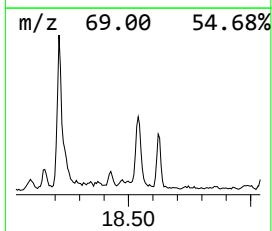
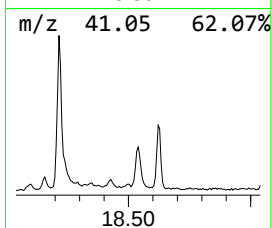
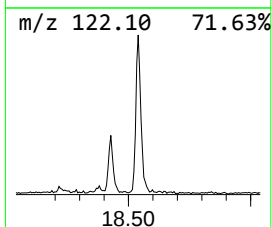
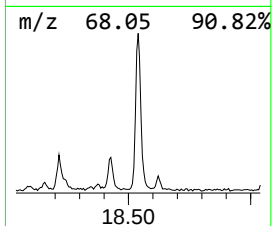
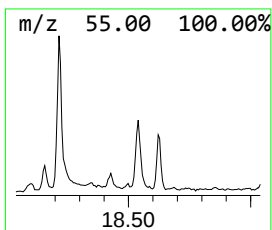
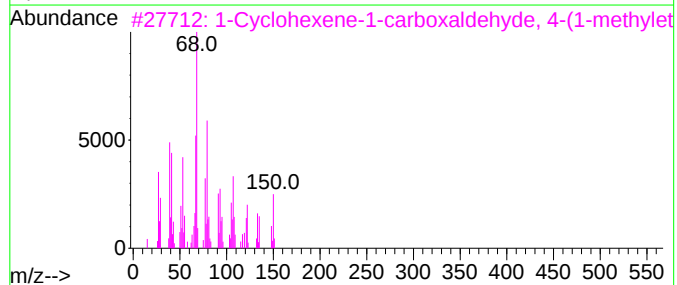
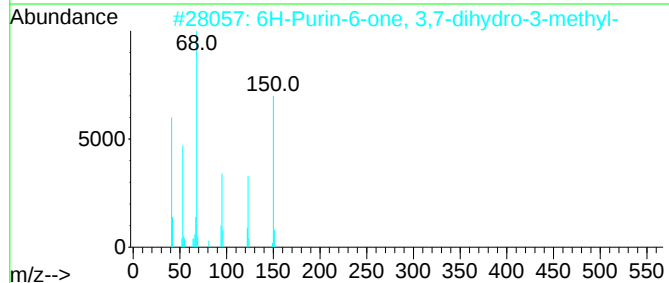
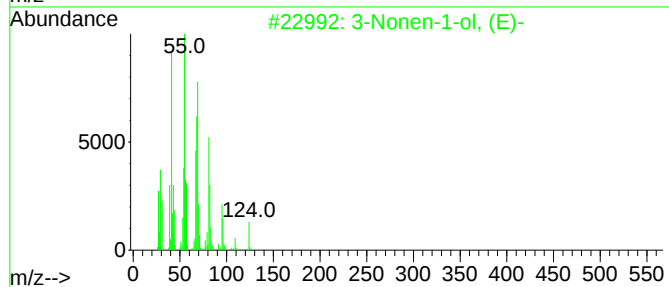
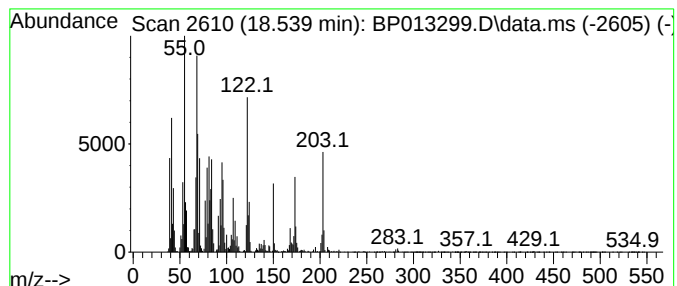
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.539	3.14 ng/ul	1013450	Phenanthrene-d10	17.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Nonen-1-ol, (E)-	142	C9H18O	010339-61-4	35
2			6H-Purin-6-one, 3,7-dihydro-3-me...	150	C6H6N4O	001006-11-7	35
3			1-Cyclohexene-1-carboxaldehyde, ...	150	C10H14O	018031-40-8	27
4			4-Heptenal, (Z)-	112	C7H12O	006728-31-0	25
5			4-Heptenal, (Z)-	112	C7H12O	006728-31-0	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013299.D
Acq On : 24 Dec 2022 14:41
Operator : CG/JU
Sample : N6016-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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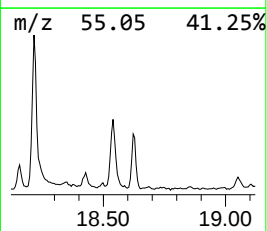
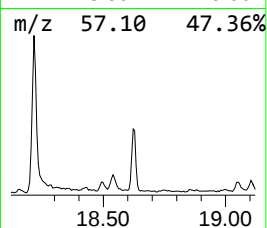
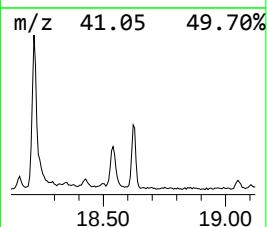
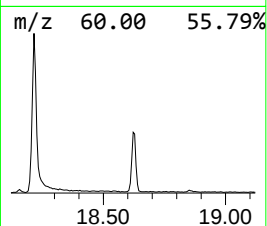
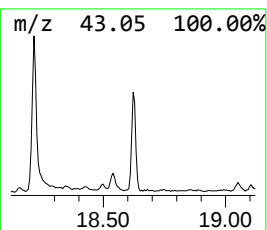
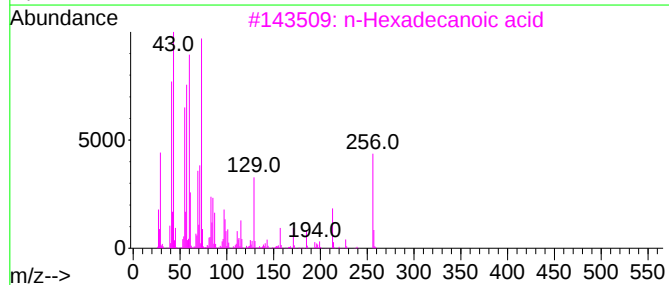
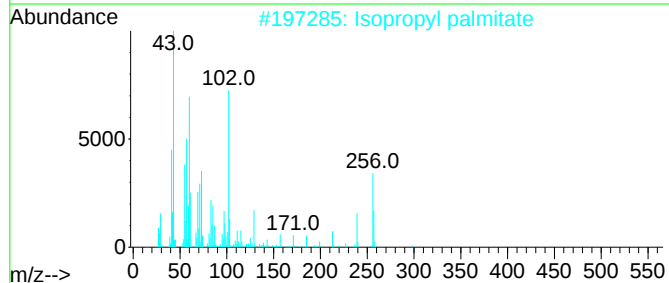
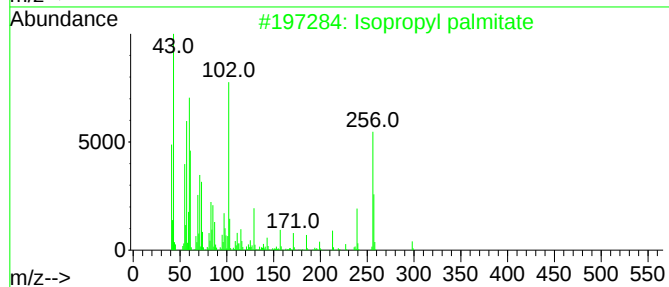
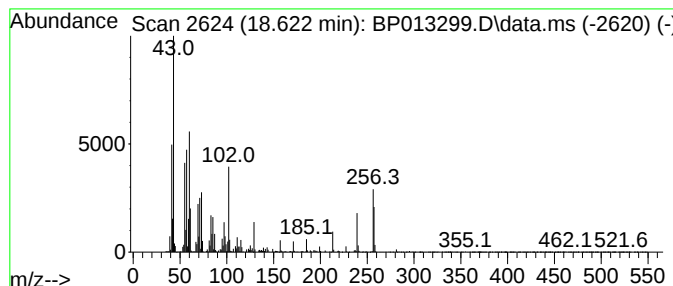
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Isopropyl palmitate Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.622	2.53 ng/ul	816629	Phenanthrene-d10	17.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl palmitate	298	C19H38O2	000142-91-6	72
2			Isopropyl palmitate	298	C19H38O2	000142-91-6	62
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	53
4			Hexadecanoic acid, propyl ester	298	C19H38O2	002239-78-3	50
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013299.D
Acq On : 24 Dec 2022 14:41
Operator : CG/JU
Sample : N6016-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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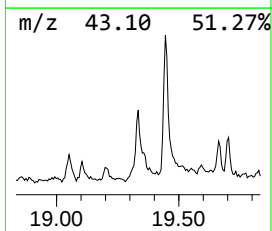
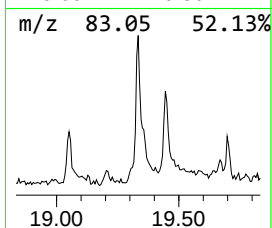
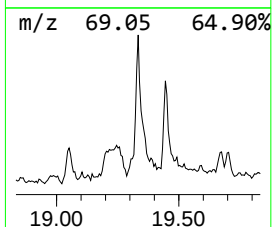
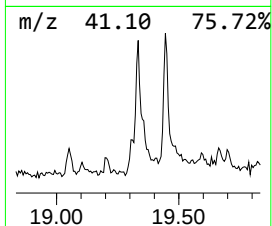
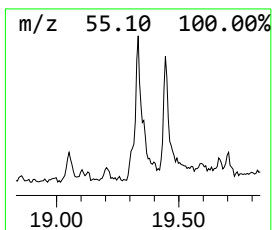
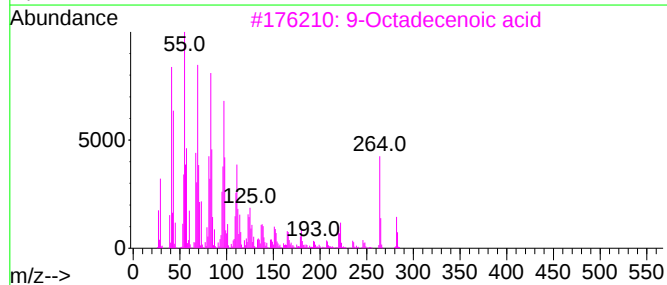
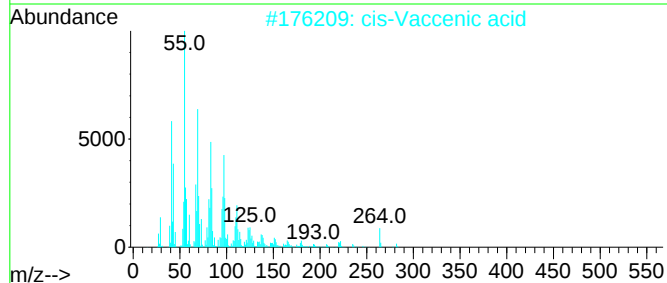
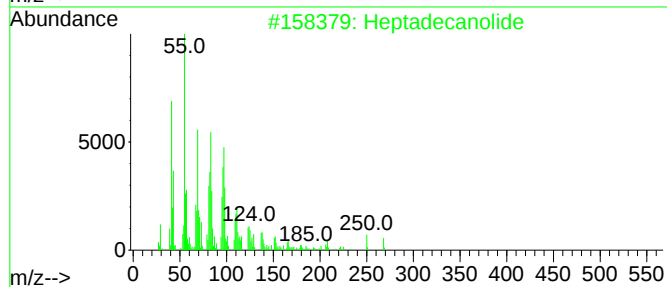
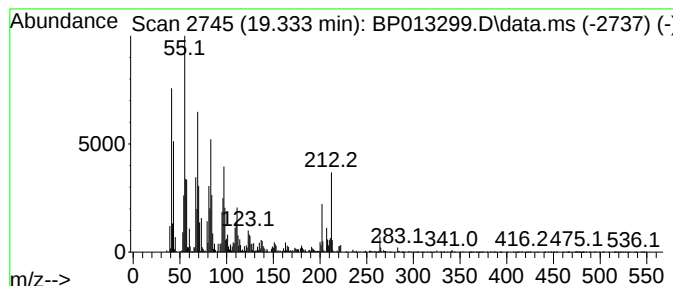
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Heptadecanolide Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.333	3.93 ng/ul	1268520	Phenanthrene-d10	17.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecanolide	268	C17H32O2	005637-97-8	86
2			cis-Vaccenic acid	282	C18H34O2	000506-17-2	76
3			9-Octadecenoic acid	282	C18H34O2	002027-47-6	76
4			Palmitoleic acid	254	C16H30O2	000373-49-9	76
5			trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013299.D
Acq On : 24 Dec 2022 14:41
Operator : CG/JU
Sample : N6016-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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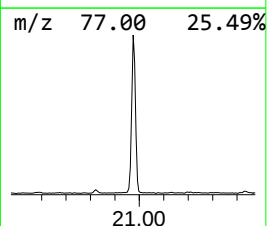
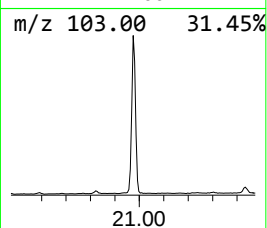
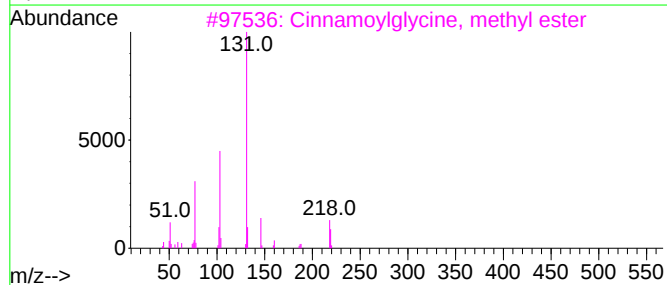
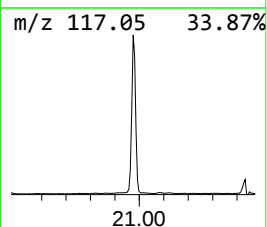
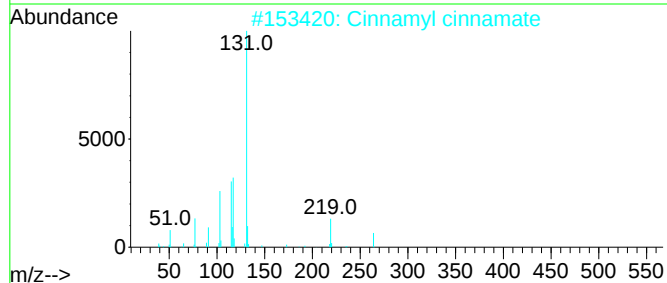
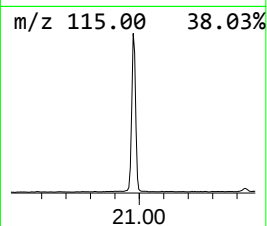
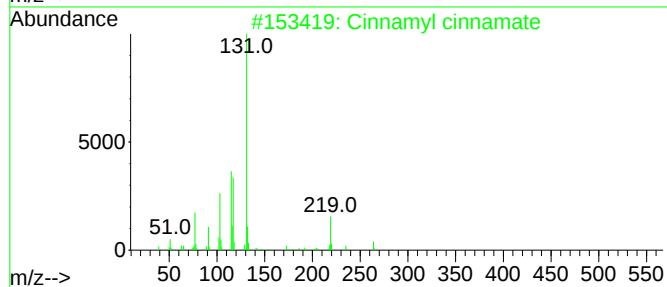
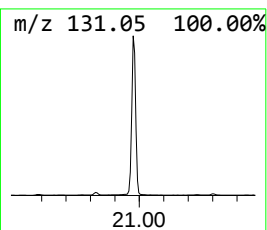
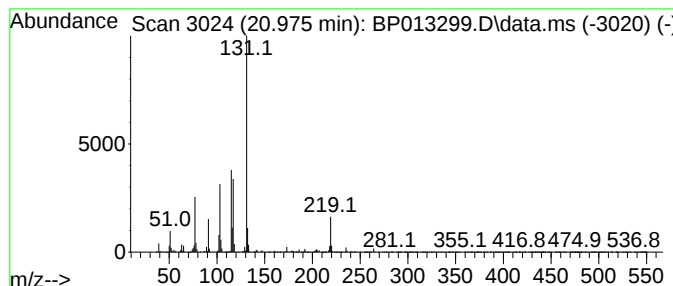
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Cinnamyl cinnamate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.974	12.01 ng/ul	4982350	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cinnamyl cinnamate	264	C18H16O2	000122-69-0	91
2			Cinnamyl cinnamate	264	C18H16O2	000122-69-0	87
3			Cinnamoylglycine, methyl ester	219	C12H13NO3	040778-04-9	53
4			2,2-Dimethyl-1-(2-vinylphenyl)pr...	188	C13H16O	1000210-99-9	52
5			(E)-4-Bromo-4-methyl-1-phenylpen...	252	C12H13BrO	1000443-02-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013299.D
Acq On : 24 Dec 2022 14:41
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Sample : N6016-06
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ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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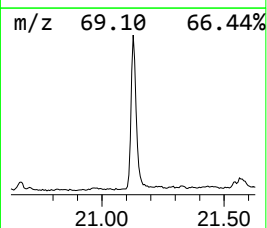
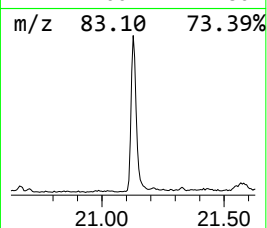
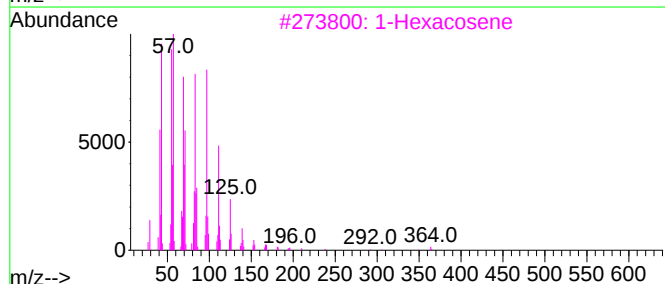
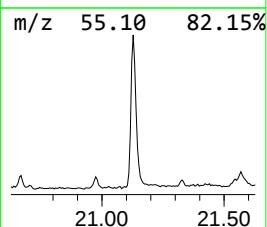
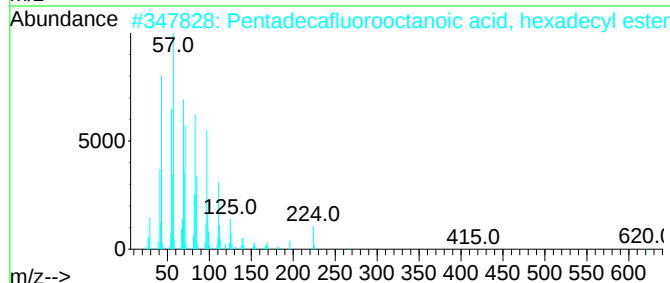
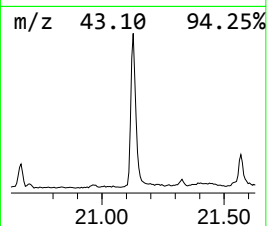
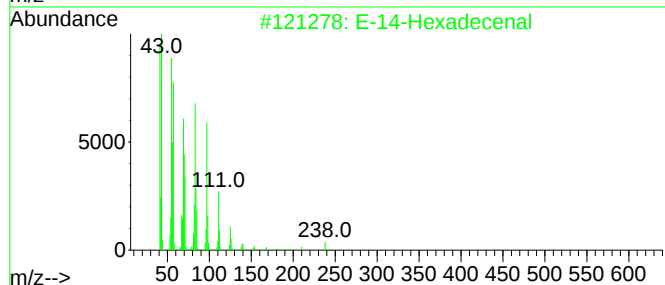
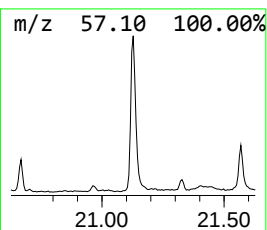
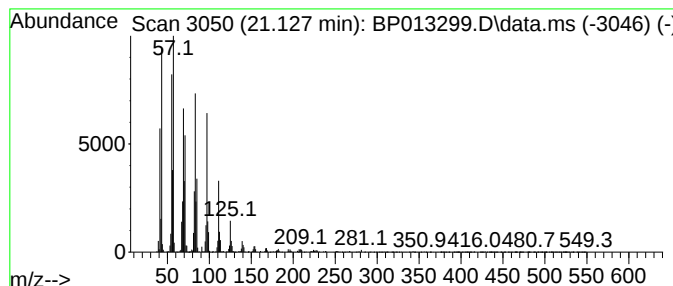
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 E-14-Hexadecenal Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.128	7.21 ng/ul	2989580	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			E-14-Hexadecenal	238	C16H30O	330207-53-9	94
2			Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	94
3			1-Hexacosene	364	C26H52	018835-33-1	94
4			Cyclotetracosane	336	C24H48	000297-03-0	94
5			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013299.D
Acq On : 24 Dec 2022 14:41
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Misc :
ALS Vial : 10 Sample Multiplier: 1

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ClientSampleId :
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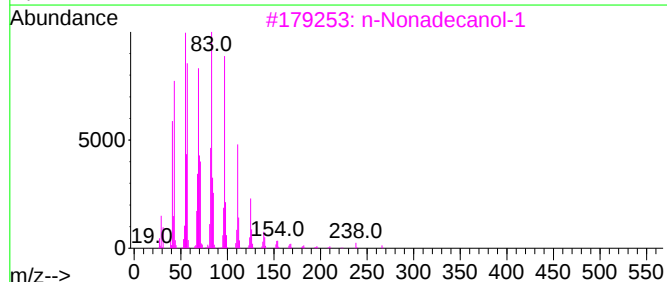
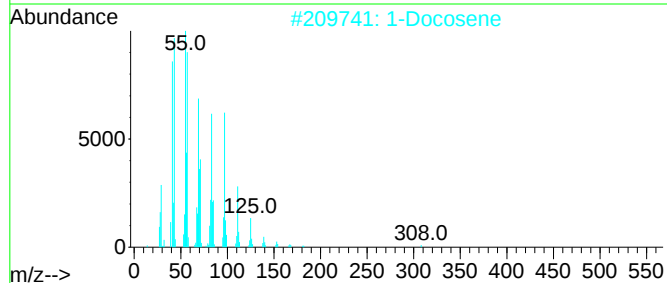
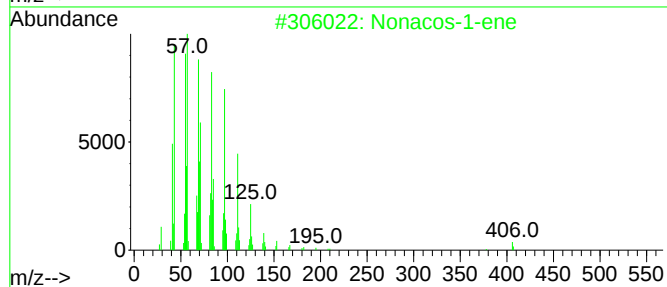
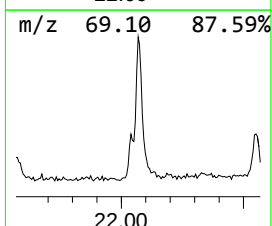
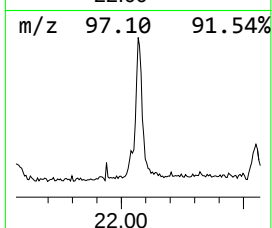
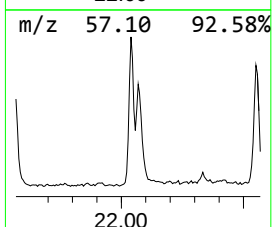
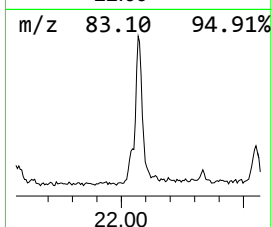
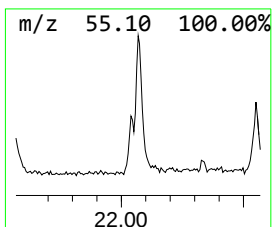
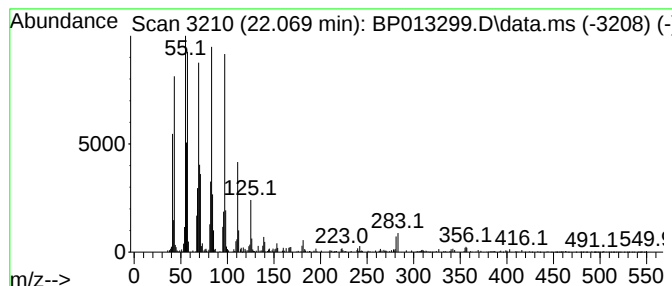
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.069	2.65 ng/ul	1098160	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonacos-1-ene	406	C29H58	018835-35-3	94
2			1-Docosene	308	C22H44	001599-67-3	91
3			n-Nonadecanol-1	284	C19H40O	001454-84-8	91
4			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91
5			1-Heneicosanol	312	C21H44O	015594-90-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013299.D
Acq On : 24 Dec 2022 14:41
Operator : CG/JU
Sample : N6016-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DBYC2

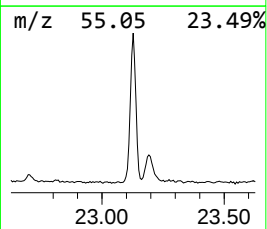
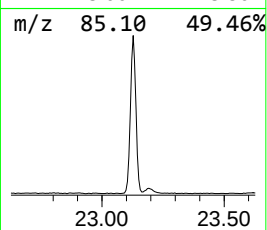
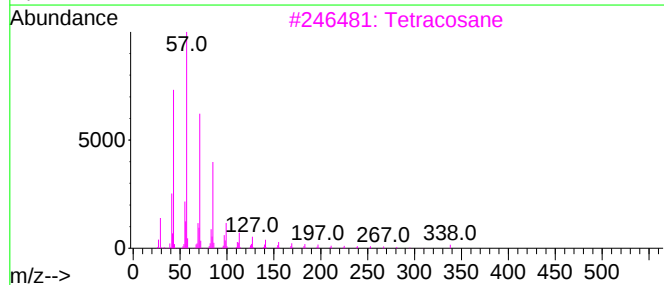
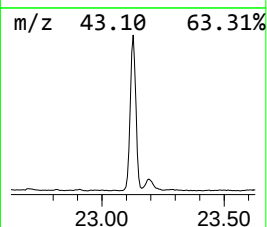
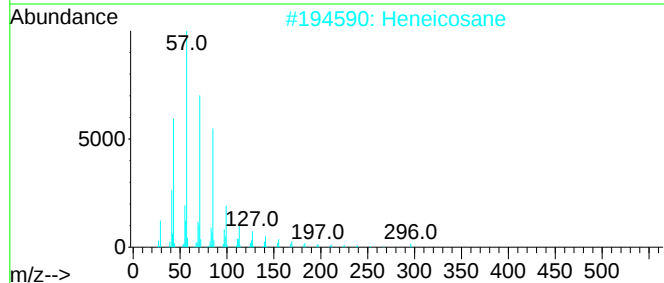
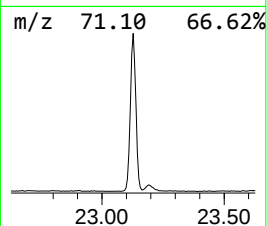
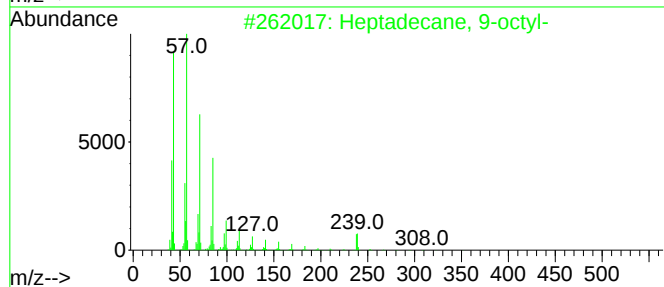
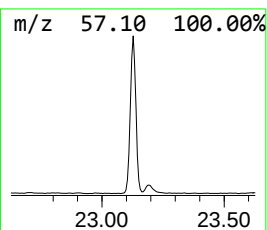
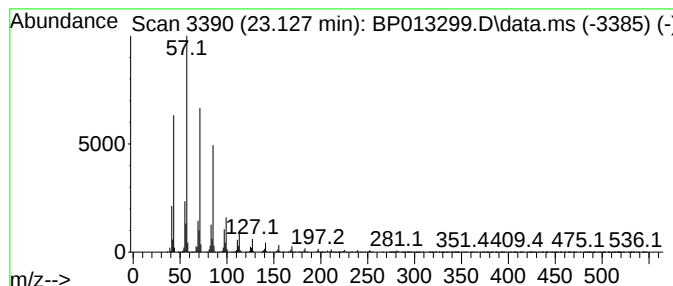
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.127	11.58 ng/ul	5231530	Perylene-d12	23.916

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
2		Heneicosane	296	C21H44	000629-94-7	91
3		Tetracosane	338	C24H50	000646-31-1	91
4		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5		Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013299.D
Acq On : 24 Dec 2022 14:41
Operator : CG/JU
Sample : N6016-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYC2

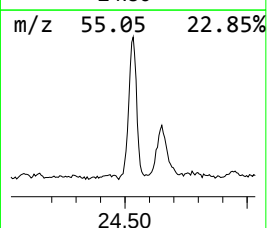
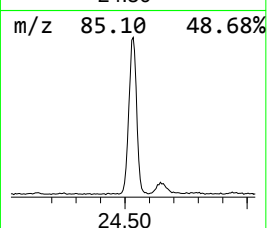
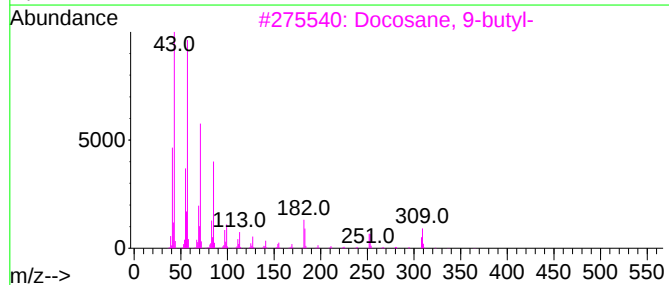
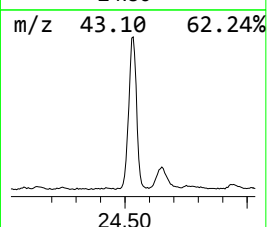
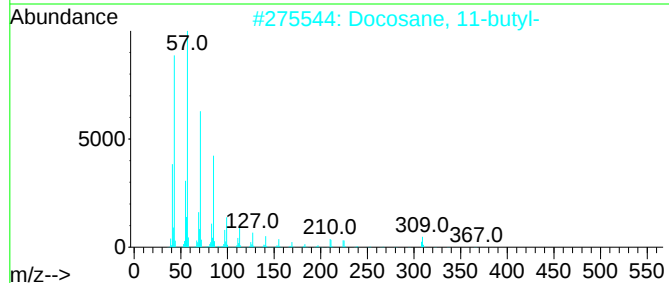
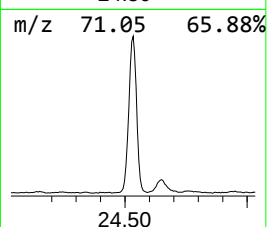
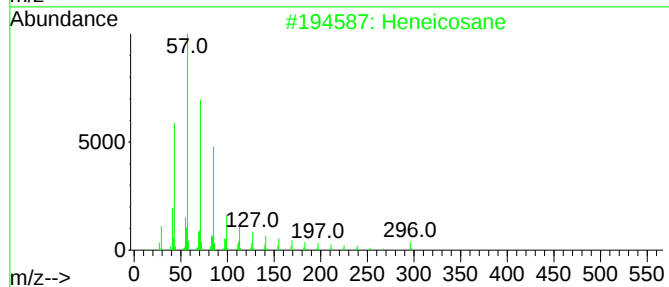
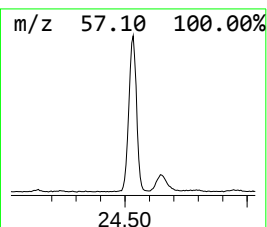
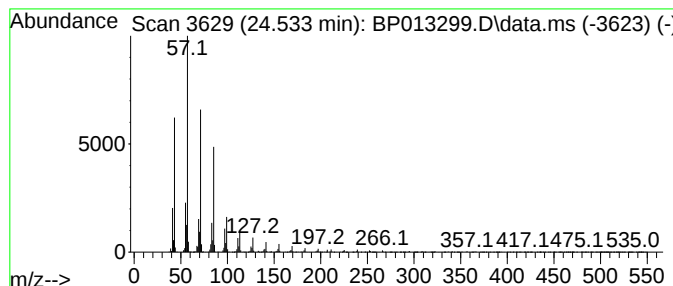
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.533	8.64 ng/ul	3906610	Perylene-d12	23.916

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	97
2		Docosane, 11-butyl-	366	C26H54	013475-76-8	94
3		Docosane, 9-butyl-	366	C26H54	055282-14-9	94
4		2-Bromo dodecane	248	C12H25Br	013187-99-0	91
5		triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013299.D
Acq On : 24 Dec 2022 14:41
Operator : CG/JU
Sample : N6016-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYC2

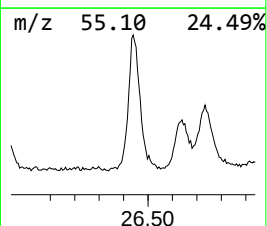
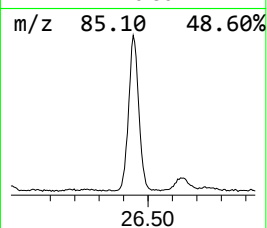
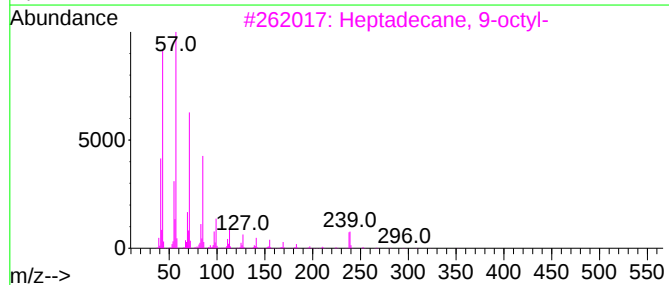
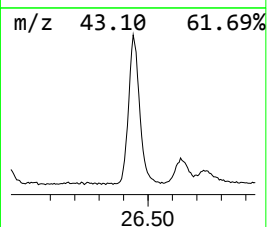
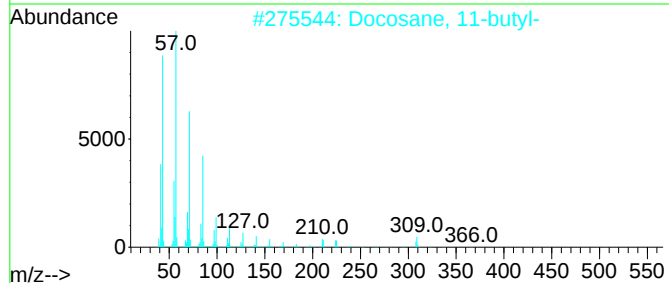
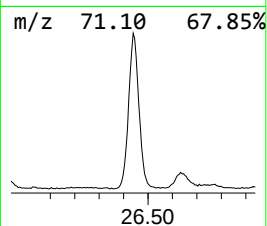
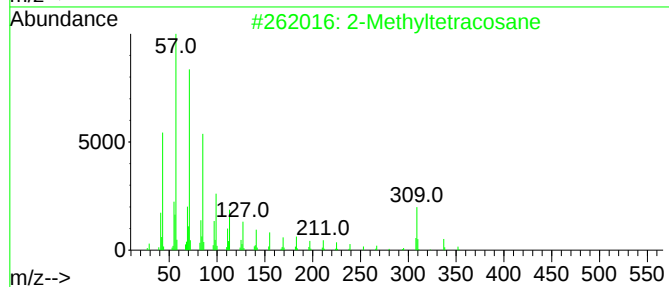
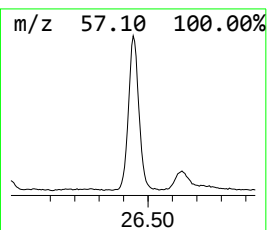
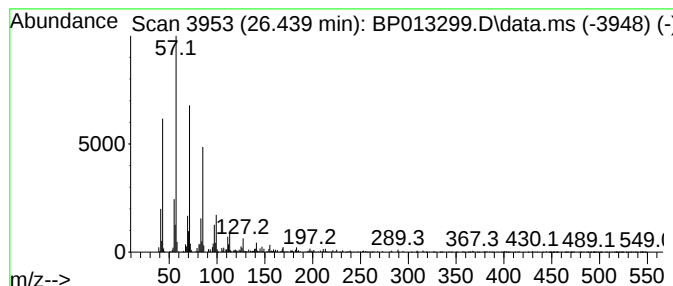
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.439	10.24 ng/ul	4628650	Perylene-d12	23.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyltetracosane			352	C25H52	001560-78-7	95
2	Docosane, 11-butyl-			366	C26H54	013475-76-8	93
3	Heptadecane, 9-octyl-			352	C25H52	007225-64-1	93
4	Pentadecane, 8-hexyl-			296	C21H44	013475-75-7	93
5	2-Bromo dodecane			248	C12H25Br	013187-99-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013299.D
Acq On : 24 Dec 2022 14:41
Operator : CG/JU
Sample : N6016-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYC2

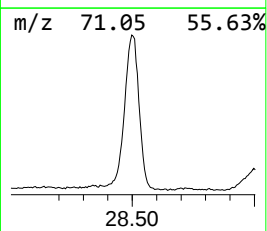
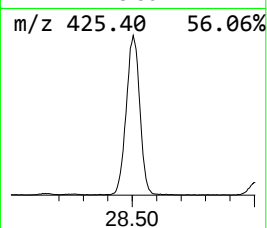
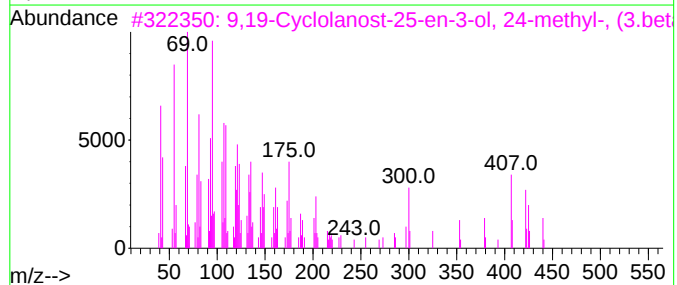
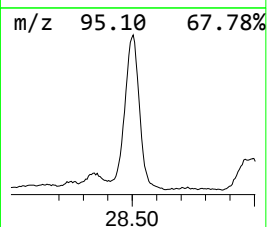
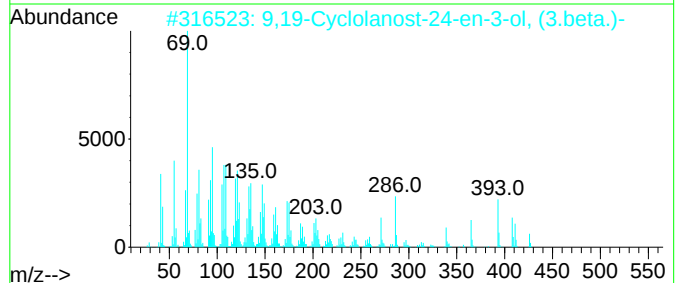
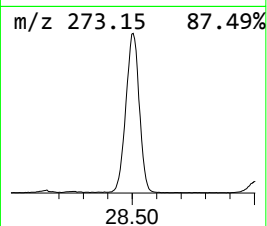
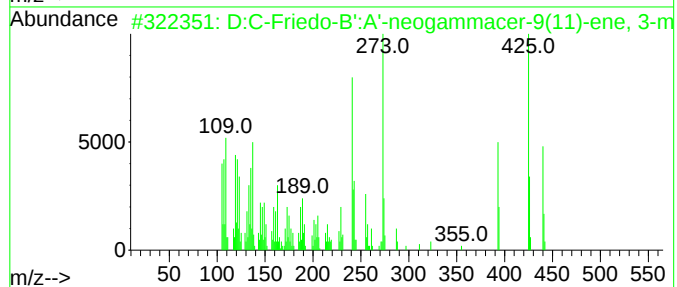
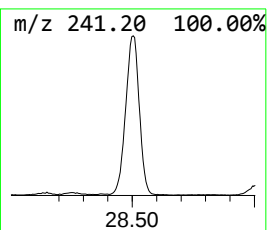
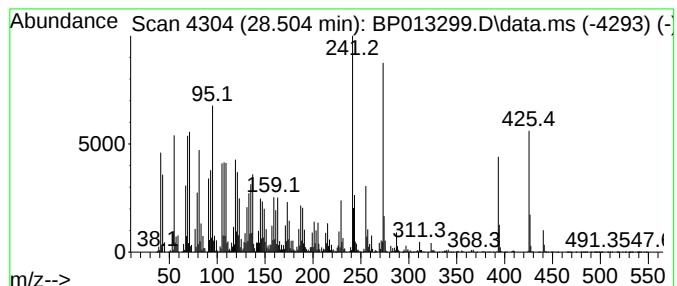
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 D:C-Friedo-B':A'-neogammace... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.504	49.44 ng/ul	22340800	Perylene-d12	23.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			D:C-Friedo-B':A'-neogammacer-9(1...	440	C31H52O	004555-56-0	53
2			9,19-Cyclolanost-24-en-3-ol, (3...	426	C30H50O	000469-38-5	38
3			9,19-Cyclolanost-25-en-3-ol, 24-...	440	C31H52O	000511-61-5	35
4			Lupeol	426	C30H50O	000545-47-1	30
5			Imidazolo[1,2-a]pyridine, 6-chlo...	273	C13H8ClN3O2	118000-62-7	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
 Data File : BP013299.D
 Acq On : 24 Dec 2022 14:41
 Operator : CG/JU
 Sample : N6016-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBYC2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.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
2-Pentanone, 4-...	5.022	3.8	ng/ul	738804	1	7.940	3860000	20.0
(1R)-2,6,6-Trim...	6.611	3.2	ng/ul	608933	1	7.940	3860000	20.0
n-Hexadecanoic ...	18.216	6.7	ng/ul	2162390	4	17.345	6456860	20.0
unknown-01	18.539	3.1	ng/ul	1013450	4	17.345	6456860	20.0
Isopropyl palmi...	18.622	2.5	ng/ul	816629	4	17.345	6456860	20.0
Heptadecanolide	19.333	3.9	ng/ul	1268520	4	17.345	6456860	20.0
Cinnamyl cinnamate	20.974	12.0	ng/ul	4982350	5	21.433	8297080	20.0
E-14-Hexadecenal	21.128	7.2	ng/ul	2989580	5	21.433	8297080	20.0
(DEL) Alkane: S...	22.069	2.6	ng/ul	1098160	5	21.433	8297080	20.0
(DEL) Alkane: S...	23.127	11.6	ng/ul	5231530	6	23.916	9038140	20.0
(DEL) Alkane: S...	24.533	8.6	ng/ul	3906610	6	23.916	9038140	20.0
(DEL) Alkane: S...	26.439	10.2	ng/ul	4628650	6	23.916	9038140	20.0
D:C-Friedo-B':A...	28.504	49.4	ng/ul	22340800	6	23.916	9038140	20.0