

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
 Data File : BP021696.D
 Acq On : 28 Aug 2024 09:35
 Operator : RC/JU
 Sample : P3675-17
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCXZ1

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-BP082324.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021696.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.017	3	5	24	rVB	4031011	4785925	39.03%	2.594%
2	3.275	45	49	58	rVB	80570	116421	0.95%	0.063%
3	3.540	90	94	104	rBV	374475	513823	4.19%	0.279%
4	3.687	115	119	133	rBV	942450	1323602	10.79%	0.718%
5	5.134	360	365	374	rBV	163623	271249	2.21%	0.147%
6	5.893	482	494	503	rBV2	60020	111103	0.91%	0.060%
7	6.499	590	597	606	rVB	656837	1010427	8.24%	0.548%
8	6.799	644	648	652	rVV2	101802	169835	1.38%	0.092%
9	6.963	667	676	690	rVB	1267106	2067074	16.86%	1.121%
10	7.122	695	703	709	rBV	1123780	1801392	14.69%	0.977%
11	7.246	717	724	731	rBV	332529	563451	4.59%	0.305%
12	7.328	731	738	744	rVV	1348560	2080519	16.96%	1.128%
13	7.393	744	749	756	rVB	245489	385662	3.14%	0.209%
14	7.793	810	817	824	rVV	1841771	2861565	23.33%	1.551%
15	8.487	927	935	943	rBV	1259071	2067508	16.86%	1.121%
16	8.934	1001	1011	1018	rBV	1311116	2132015	17.38%	1.156%
17	9.363	1078	1084	1099	rVB5	65186	162582	1.33%	0.088%
18	9.499	1099	1107	1114	rBV	136452	242838	1.98%	0.132%
19	9.657	1127	1134	1140	rBV	987001	1688864	13.77%	0.916%
20	9.828	1155	1163	1168	rBV	236723	515561	4.20%	0.279%
21	9.887	1168	1173	1183	rVB2	288635	576540	4.70%	0.313%
22	10.199	1216	1226	1247	rVV	1509799	2954254	24.09%	1.602%
23	10.575	1282	1290	1303	rVB	2428791	4230823	34.50%	2.294%
24	10.704	1303	1312	1329	rBV	710215	1434670	11.70%	0.778%
25	10.881	1337	1342	1349	rBV2	70827	120528	0.98%	0.065%
26	11.116	1376	1382	1389	rBV	120283	244017	1.99%	0.132%
27	11.316	1406	1416	1433	rBV	319650	764753	6.24%	0.415%
28	12.798	1663	1668	1676	rVB	104882	183873	1.50%	0.100%
29	13.169	1724	1731	1745	rBV2	339954	623207	5.08%	0.338%
30	13.316	1750	1756	1765	rVV3	651400	1146795	9.35%	0.622%
31	13.716	1818	1824	1828	rVV4	114143	197642	1.61%	0.107%
32	13.769	1828	1833	1838	rVV	545712	789105	6.43%	0.428%
33	13.834	1838	1844	1852	rVV	2906758	4342971	35.41%	2.354%
34	14.128	1884	1894	1904	rVV	3426978	5040379	41.10%	2.732%
35	14.287	1915	1921	1928	rVB5	93436	169270	1.38%	0.092%
36	14.440	1934	1947	1954	rBV	4209334	6312984	51.48%	3.422%

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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-BP082324.MA.M
 Title : SVOA CALIBRATION

37	14.516	1955	1960	1969	rBV2	293167	606206	4.94%	0.329%
38	14.663	1976	1985	1986	rBV2	7926937	12263624	100.00%	6.648%
39	14.681	1986	1988	1995	rVV	8930532	11223134	91.52%	6.084%
40	14.734	1995	1997	2001	rVB	228331	241260	1.97%	0.131%
41	14.869	2016	2020	2025	rBV3	89514	152674	1.24%	0.083%
42	15.345	2094	2101	2102	rBV6	89814	136832	1.12%	0.074%
43	15.375	2103	2106	2111	rVV2	209506	363743	2.97%	0.197%
44	15.451	2112	2119	2131	rVB	4324129	6584335	53.69%	3.569%
45	15.563	2132	2138	2147	rBV	1149200	1749312	14.26%	0.948%
46	15.645	2148	2152	2158	rVV3	79227	139321	1.14%	0.076%
47	15.951	2199	2204	2209	rVB4	87698	182579	1.49%	0.099%
48	16.034	2214	2218	2228	rVB5	79569	194911	1.59%	0.106%
49	16.257	2251	2256	2261	rVB2	326510	470201	3.83%	0.255%
50	16.645	2319	2322	2332	rVB	164880	271096	2.21%	0.147%
51	16.863	2353	2359	2366	rBV8	43297	106565	0.87%	0.058%
52	17.057	2387	2392	2400	rBV	806834	1098380	8.96%	0.595%
53	17.163	2406	2410	2412	rBV3	105707	151229	1.23%	0.082%
54	17.192	2412	2415	2418	rVV3	165285	307342	2.51%	0.167%
55	17.245	2418	2424	2431	rVV	4512158	7506932	61.21%	4.070%
56	17.351	2432	2442	2452	rVB	4290655	6951236	56.68%	3.768%
57	17.445	2453	2458	2464	rBV3	143121	258781	2.11%	0.140%
58	17.563	2474	2478	2481	rVV4	110125	171652	1.40%	0.093%
59	17.604	2483	2485	2488	rVV2	93958	115213	0.94%	0.062%
60	17.651	2489	2493	2502	rVB3	188319	342891	2.80%	0.186%
61	17.922	2534	2539	2543	rBV	766574	1144954	9.34%	0.621%
62	18.133	2571	2575	2580	rBV3	109855	186041	1.52%	0.101%
63	18.204	2581	2587	2597	rBV	3094299	5052067	41.20%	2.739%
64	18.286	2598	2601	2604	rVV2	148311	178056	1.45%	0.097%
65	18.445	2624	2628	2630	rBV	147056	212421	1.73%	0.115%
66	18.639	2658	2661	2667	rBV4	110301	265034	2.16%	0.144%
67	19.080	2731	2736	2742	rBV4	361278	618734	5.05%	0.335%
68	19.139	2742	2746	2749	rVV4	130803	187699	1.53%	0.102%
69	19.180	2749	2753	2757	rVB2	371102	465561	3.80%	0.252%
70	19.275	2766	2769	2771	rBV3	128796	142174	1.16%	0.077%
71	19.322	2773	2777	2779	rBV2	149946	194310	1.58%	0.105%
72	19.363	2781	2784	2786	rBV2	319223	359452	2.93%	0.195%
73	19.398	2786	2790	2791	rBV2	247102	311486	2.54%	0.169%
74	19.522	2806	2811	2821	rBV	2874443	4038288	32.93%	2.189%
75	19.728	2841	2846	2855	rBV	5809383	8035637	65.52%	4.356%
76	19.992	2888	2891	2897	rVB5	264968	368886	3.01%	0.200%

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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-BP082324.MA.M
 Title : SVOA CALIBRATION

77	20.151	2915	2918	2923	rBV	289972	380066	3.10%	0.206%
78	20.198	2924	2926	2935	rVV4	377480	581672	4.74%	0.315%
79	20.275	2936	2939	2951	rVB2	599908	1076092	8.77%	0.583%
80	20.451	2964	2969	2974	rBV8	173510	446005	3.64%	0.242%
81	20.539	2981	2984	2988	rBV5	186435	246341	2.01%	0.134%
82	20.610	2993	2996	2999	rVV2	256389	338378	2.76%	0.183%
83	20.651	2999	3003	3008	rVV	805947	1058854	8.63%	0.574%
84	20.704	3009	3012	3016	rVV	301534	399697	3.26%	0.217%
85	20.757	3018	3021	3025	rVV	278906	397810	3.24%	0.216%
86	20.816	3028	3031	3038	rVB4	484142	831788	6.78%	0.451%
87	21.151	3084	3088	3096	rBV	2292092	3177422	25.91%	1.722%
88	21.322	3112	3117	3121	rBV	2029758	3105850	25.33%	1.684%
89	21.369	3121	3125	3129	rVB	1076139	1732741	14.13%	0.939%
90	21.563	3155	3158	3161	rBV	402230	547652	4.47%	0.297%
91	21.598	3161	3164	3169	rVB	551577	745284	6.08%	0.404%
92	21.698	3175	3181	3189	rVV2	4488270	7788585	63.51%	4.222%
93	21.769	3191	3193	3201	rVB3	286752	439816	3.59%	0.238%
94	22.645	3336	3342	3358	rVB	1423711	3449785	28.13%	1.870%
95	24.280	3614	3620	3624	rBV2	582223	1298868	10.59%	0.704%
96	24.880	3712	3722	3732	rBV	2660612	7412066	60.44%	4.018%
97	25.110	3753	3761	3774	rVB	2948232	8211820	66.96%	4.452%
98	26.557	4000	4007	4016	rVB3	458604	1561282	12.73%	0.846%
99	26.680	4019	4028	4045	rVB	1916087	6085739	49.62%	3.299%
100	30.880	4734	4742	4763	rVB3	1318067	5798868	47.29%	3.144%

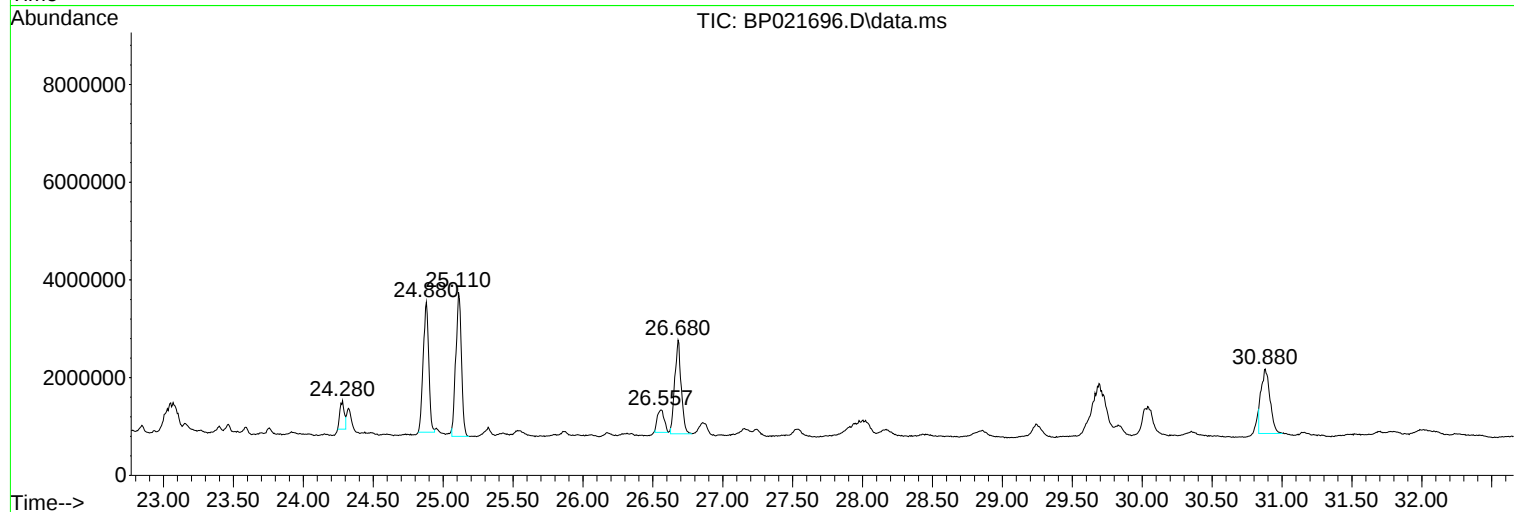
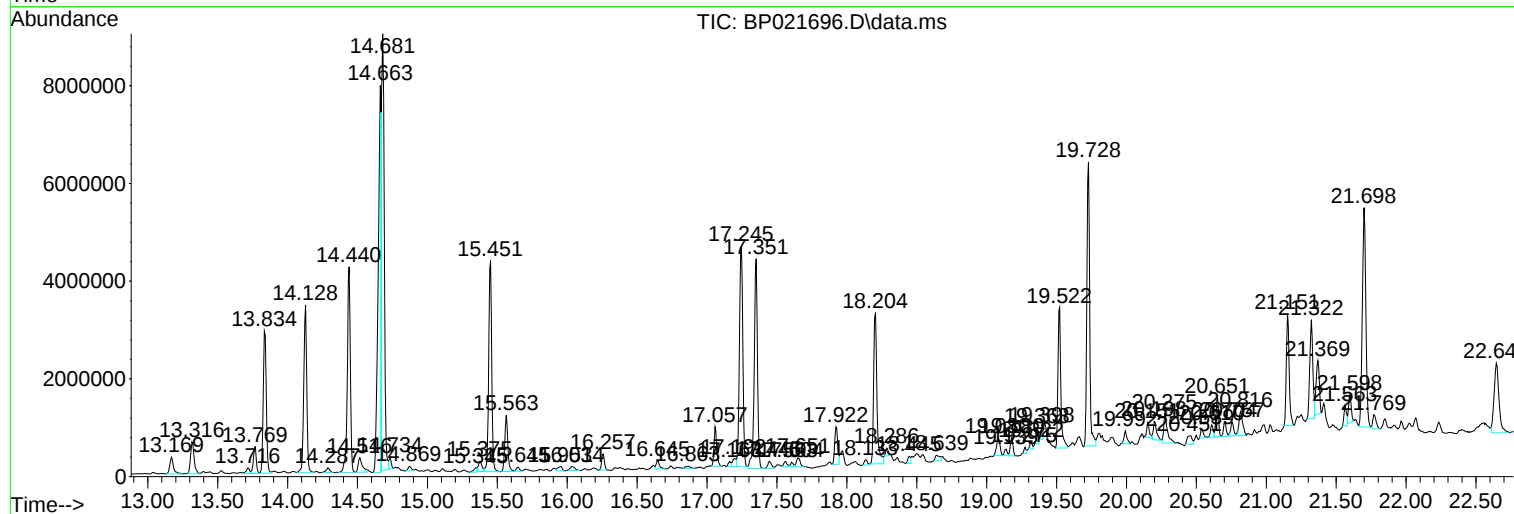
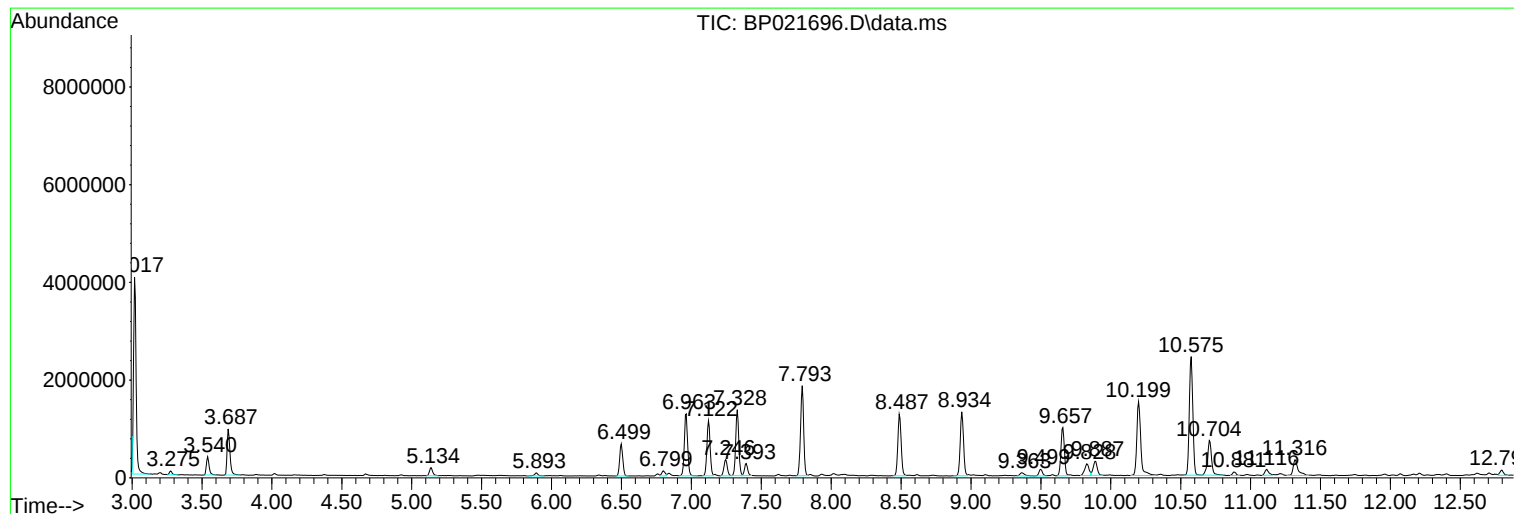
Sum of corrected areas: 184465957

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Instrument :
BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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



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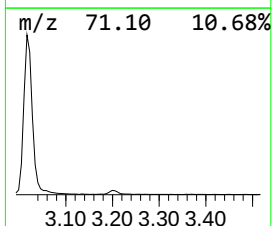
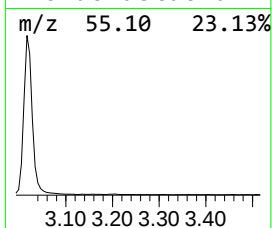
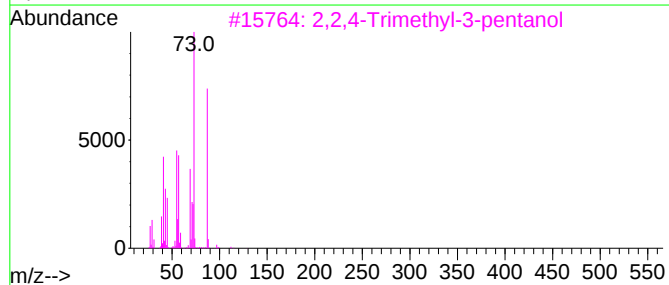
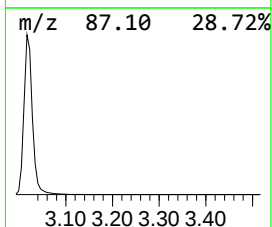
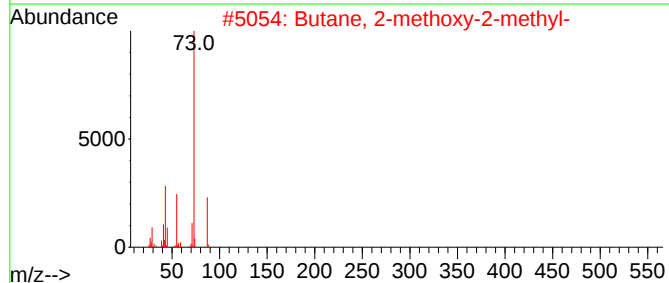
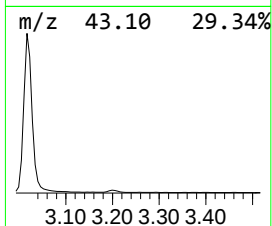
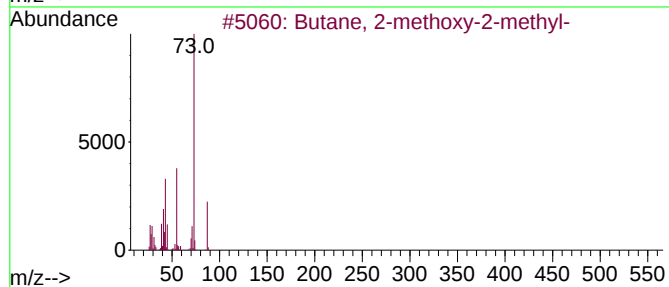
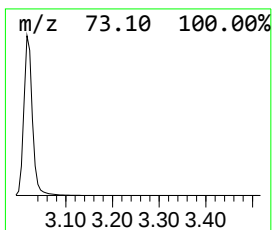
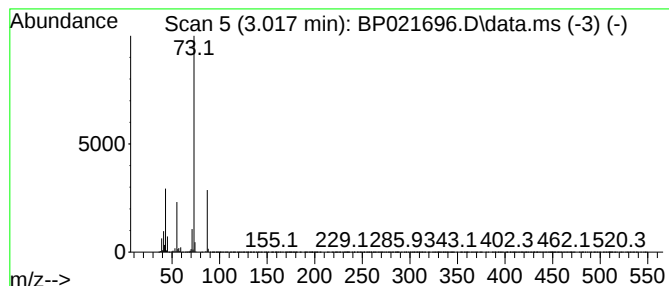
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.017	33.45 ng/ul	4785930	1,4-Dichlorobenzene-d4	7.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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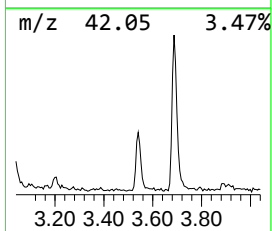
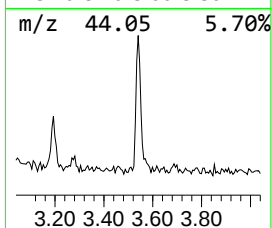
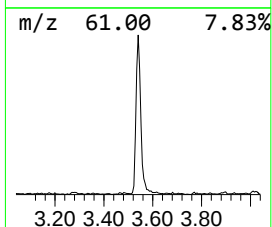
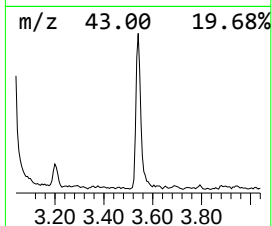
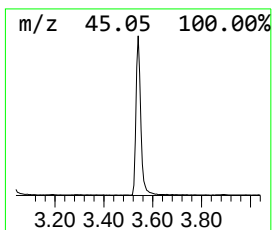
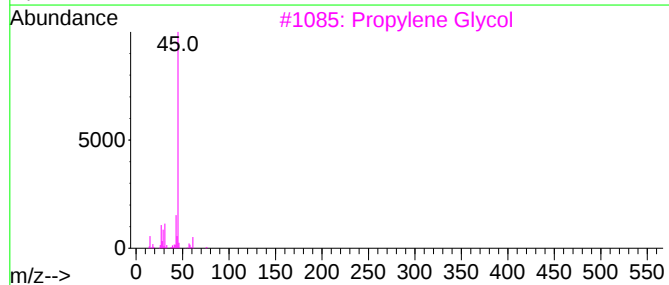
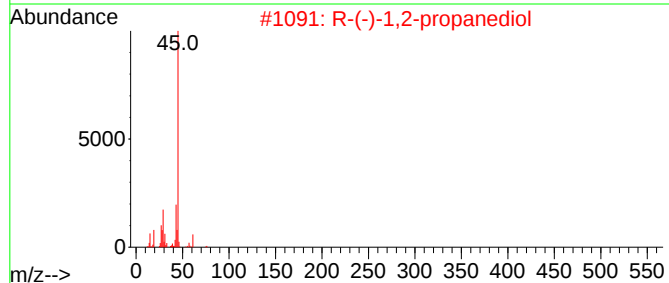
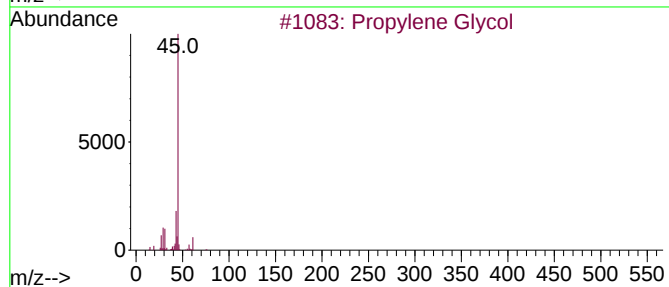
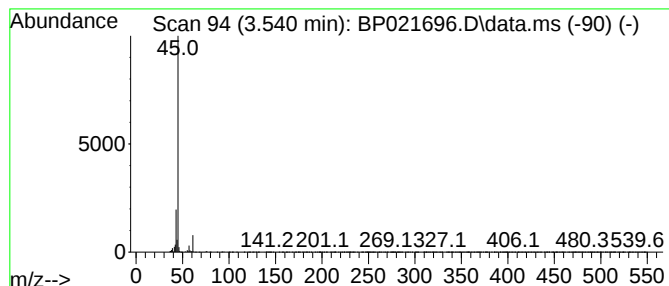
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Propylene Glycol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.540	3.59 ng/ul	513823	1,4-Dichlorobenzene-d4	7.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	90
2			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	78
3			Propylene Glycol	76	C3H8O2	000057-55-6	64
4			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
5			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40



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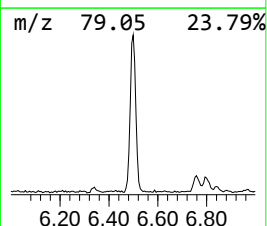
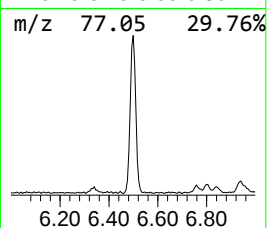
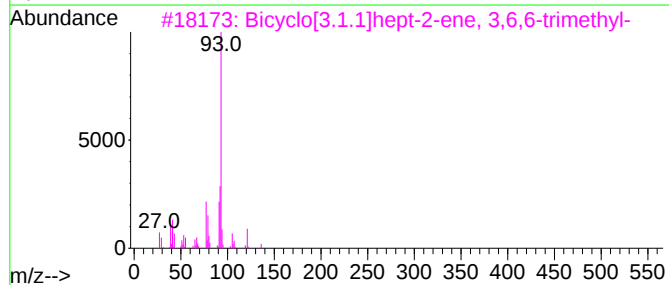
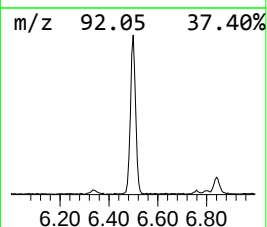
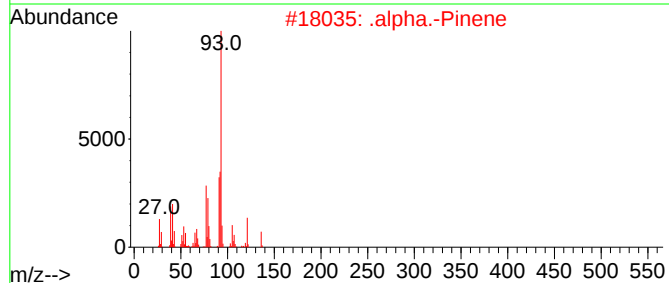
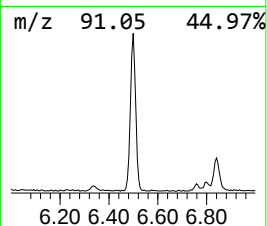
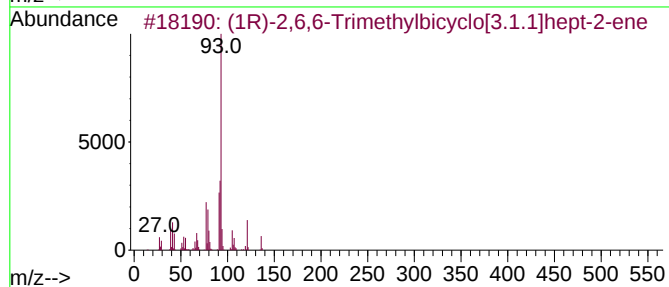
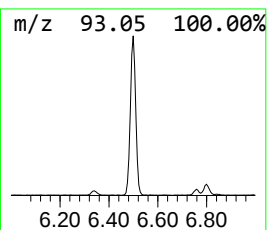
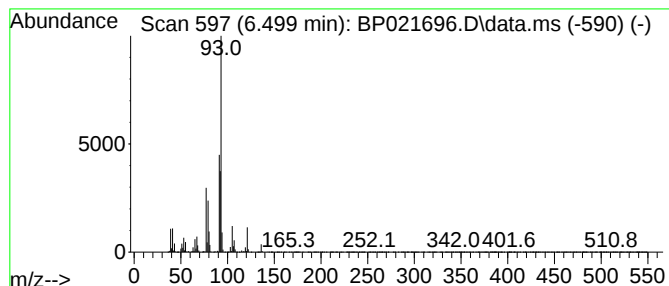
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.499	7.06 ng/ul	1010430	1,4-Dichlorobenzene-d4	7.793

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	94		
2	.alpha.-Pinene	136	C10H16	000080-56-8	94		
3	Bicyclo[3.1.1]hept-2-ene, 3,6,6-trimethyl-	136	C10H16	004889-83-2	91		
4	.alpha.-Pinene	136	C10H16	000080-56-8	91		
5	(1S)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-26-4	90		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
Operator : RC/JU
Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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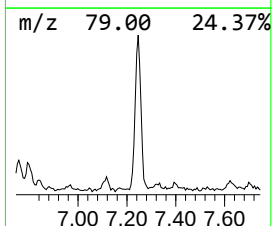
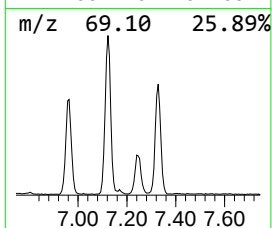
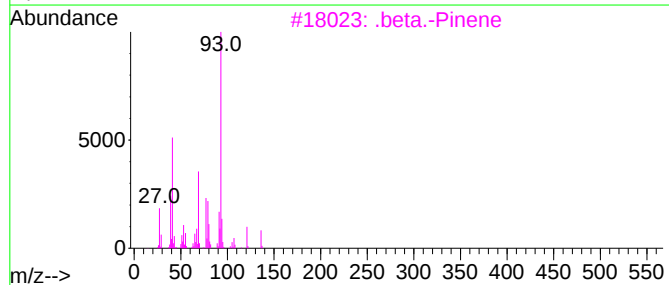
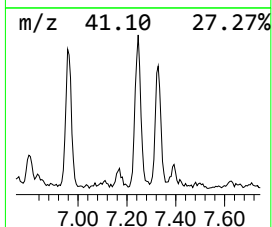
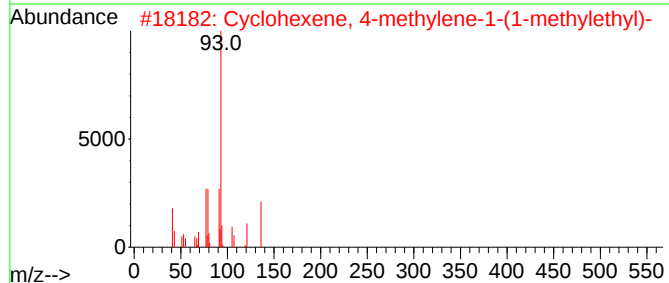
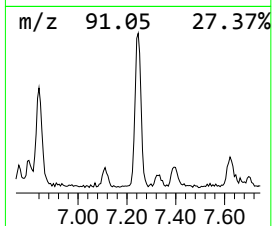
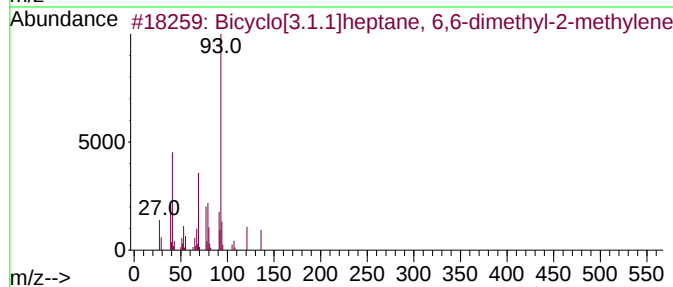
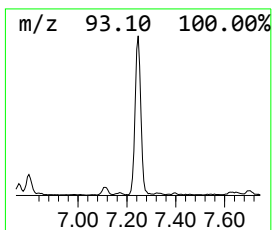
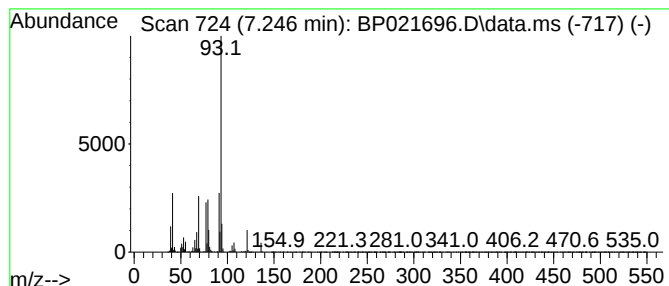
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Bicyclo[3.1.1]heptane, 6,6-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.246	3.94 ng/ul	563451	1,4-Dichlorobenzene-d4	7.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	91
2			Cyclohexene, 4-methylene-1-(1-me...	136	C10H16	000099-84-3	91
3			.beta.-Pinene	136	C10H16	000127-91-3	91
4			Cyclohexane, 1-methylene-4-(1-me...	136	C10H16	000499-97-8	87
5			.alpha.-Pinene	136	C10H16	000080-56-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
Operator : RC/JU
Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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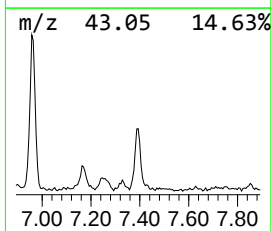
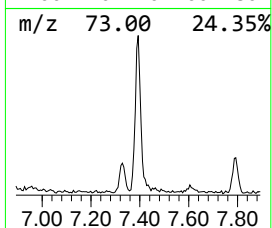
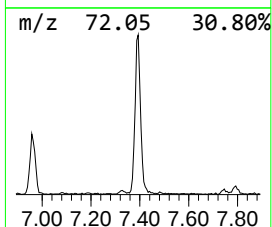
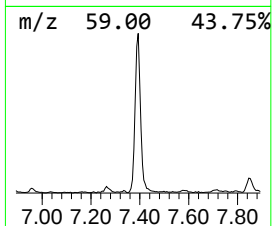
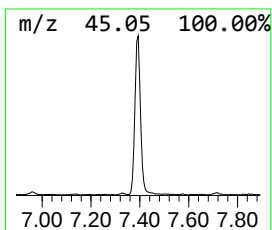
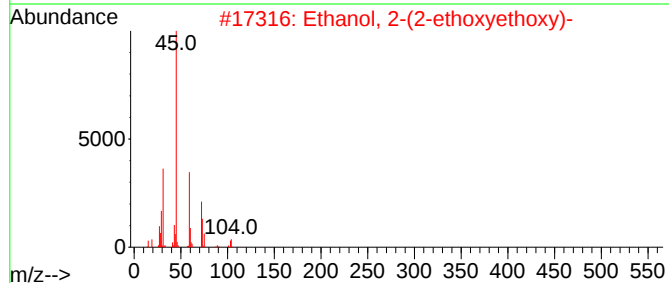
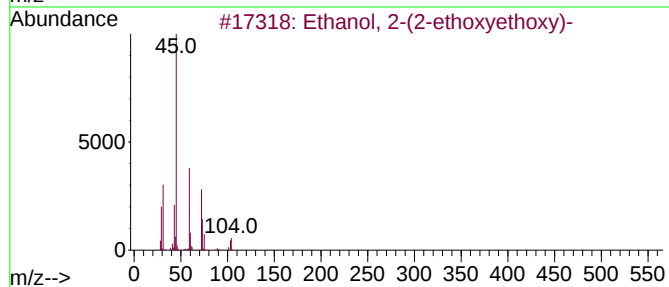
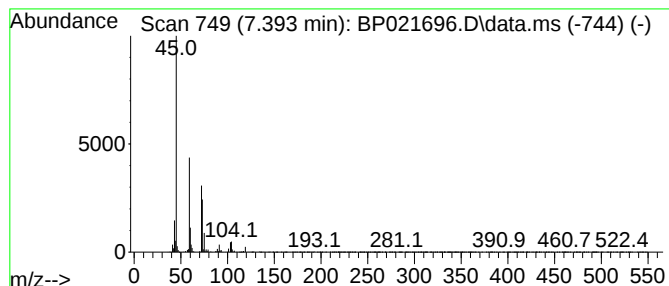
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.393	2.70 ng/ul	385662	1,4-Dichlorobenzene-d4	7.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	86
2			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	83
3			Diethyl carbitol	162	C8H18O3	000112-36-7	72
4			Ethanol, 2-[2-(2-ethoxyethoxy)et...	178	C8H18O4	000112-50-5	72
5			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
Operator : RC/JU
Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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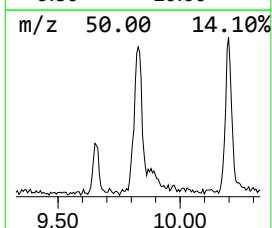
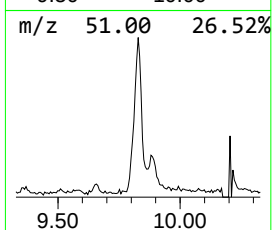
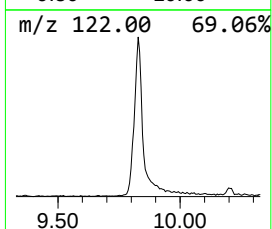
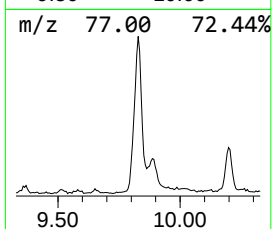
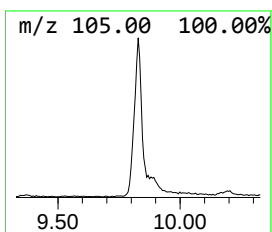
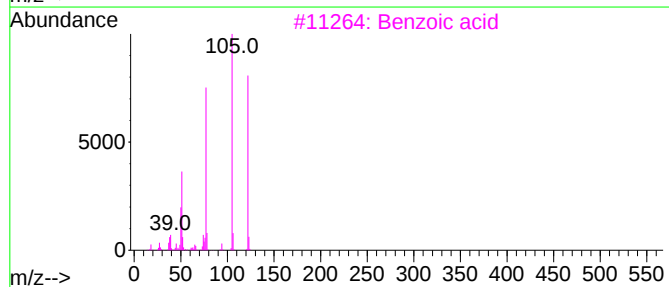
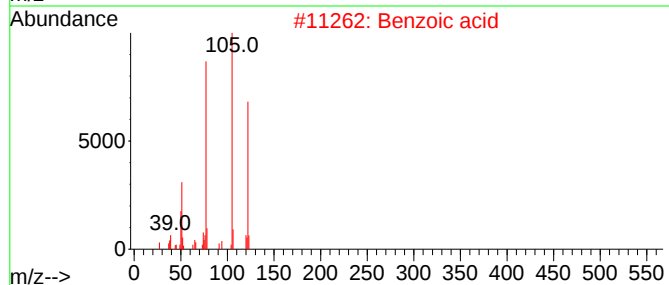
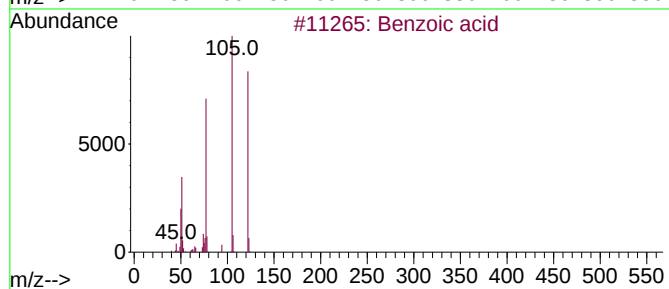
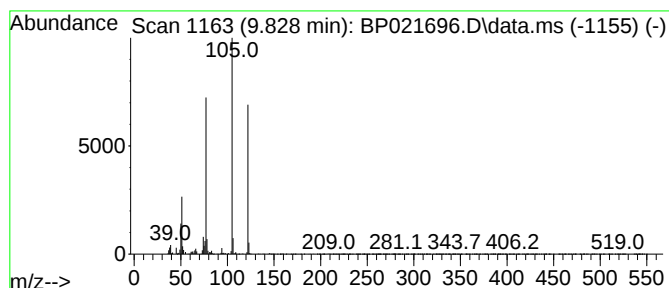
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzoic acid Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.828	2.44 ng/ul	515561	Naphthalene-d8	10.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid			122	C7H6O2	000065-85-0	96
2	Benzoic acid			122	C7H6O2	000065-85-0	94
3	Benzoic acid			122	C7H6O2	000065-85-0	94
4	Benzoic acid			122	C7H6O2	000065-85-0	93
5	Benzoic acid, silver(1+) salt			228	C7H5AgO2	000532-31-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
Operator : RC/JU
Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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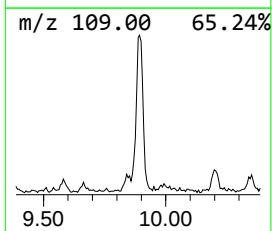
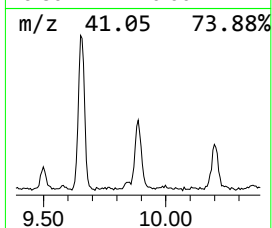
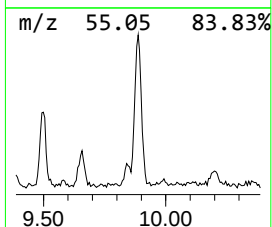
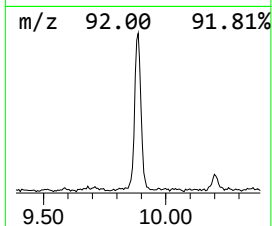
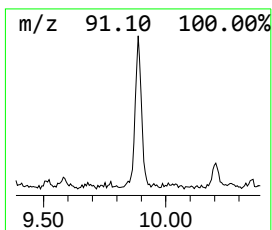
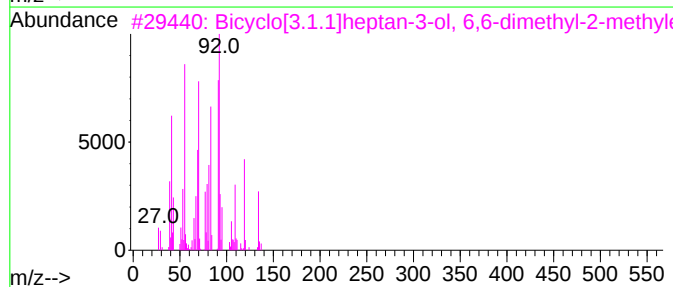
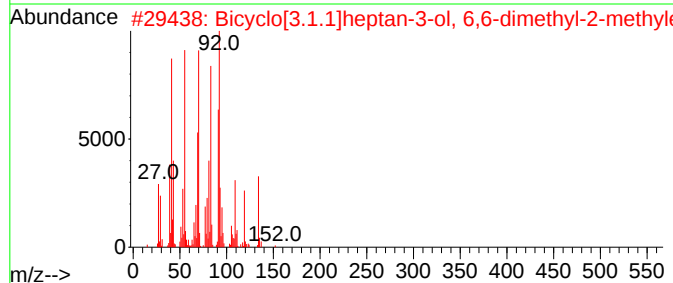
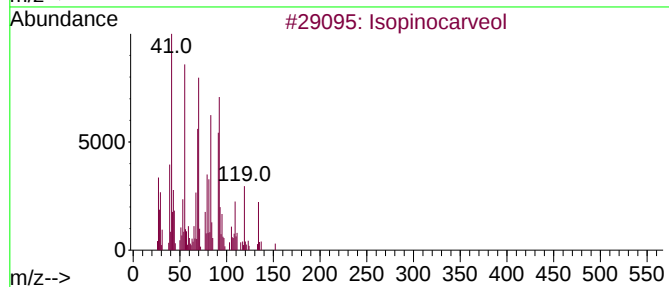
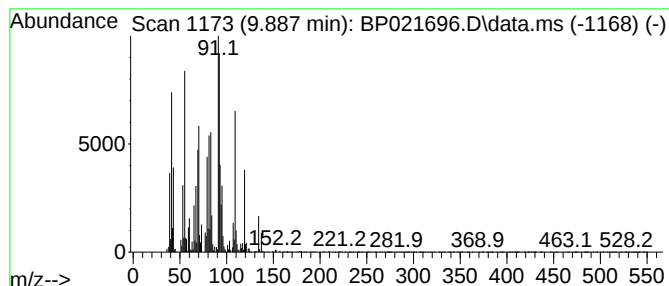
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-01 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.887	2.73 ng/ul	576540	Naphthalene-d8	10.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopinocarveol	152	C10H16O	006712-79-4	43
2			Bicyclo[3.1.1]heptan-3-ol, 6,6-d...	152	C10H16O	000547-61-5	41
3			Bicyclo[3.1.1]heptan-3-ol, 6,6-d...	152	C10H16O	000547-61-5	40
4			1-Adamantyl fluoroformate	198	C11H15FO2	062087-82-5	32
5			Bicyclo[3.1.1]heptan-3-ol, 6,6-d...	152	C10H16O	005947-36-4	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
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Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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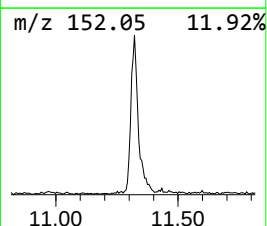
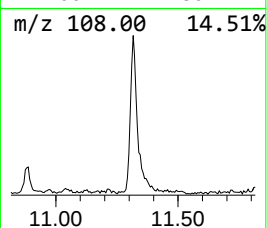
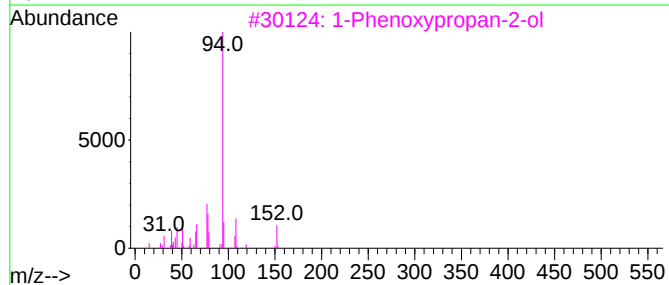
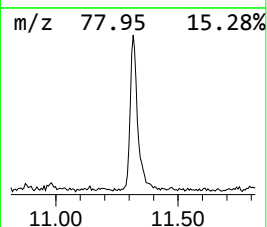
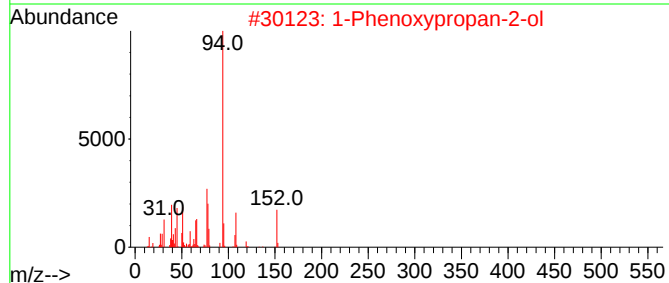
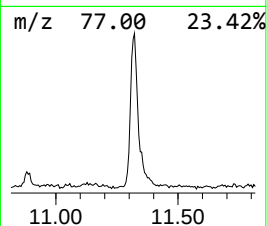
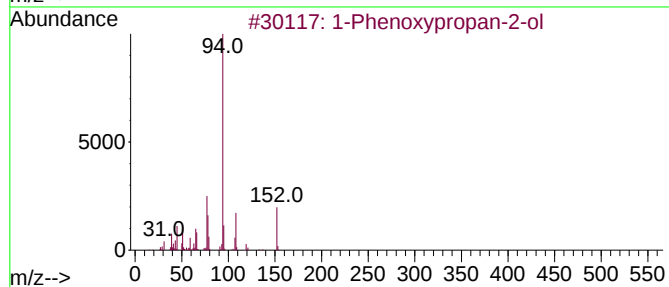
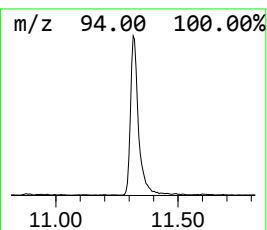
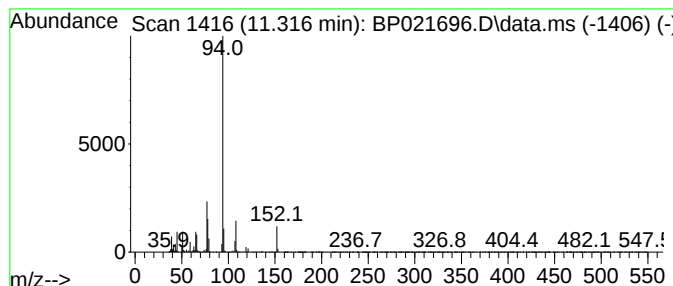
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1-Phenoxypropan-2-ol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.316	3.62 ng/ul	764753	Naphthalene-d8	10.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	90
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
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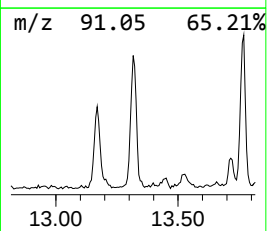
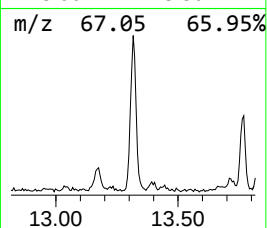
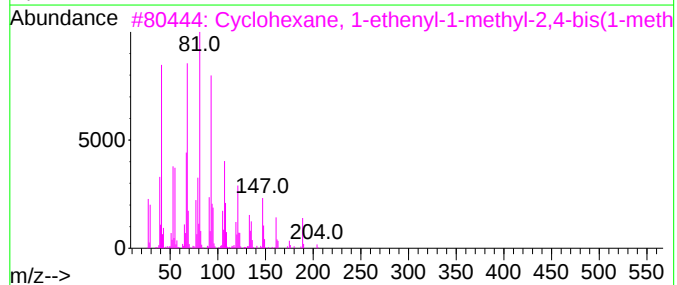
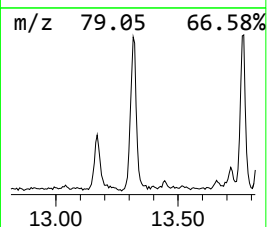
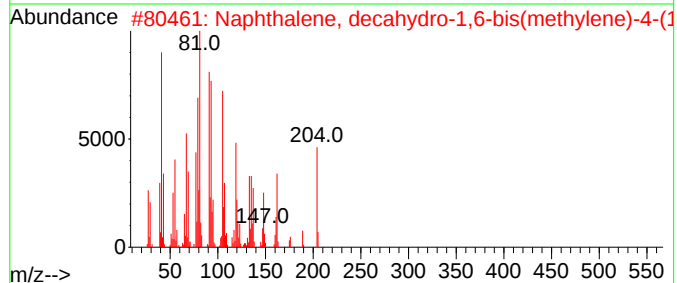
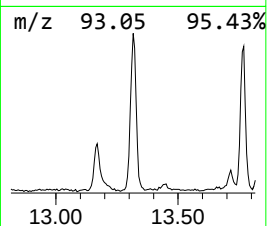
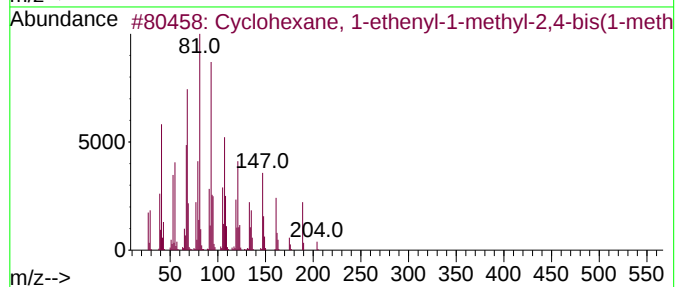
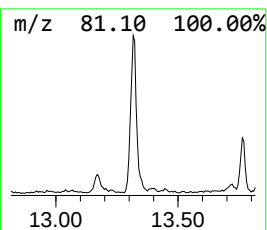
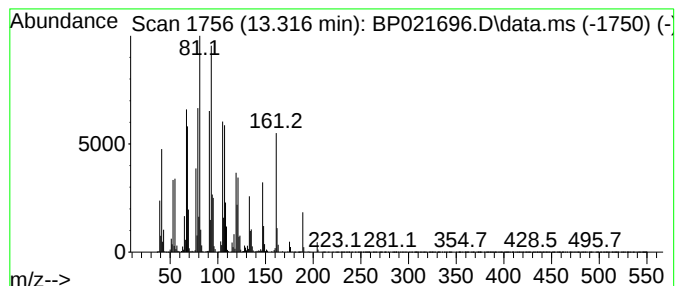
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Cyclohexane, 1-ethenyl-1-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.316	3.63 ng/ul	1146800	Acenaphthene-d10	14.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-ethenyl-1-methyl-...	204	C15H24	000515-13-9	91
2			Naphthalene, decahydro-1,6-bis(m...	204	C15H24	030021-46-6	86
3			Cyclohexane, 1-ethenyl-1-methyl-...	204	C15H24	033880-83-0	74
4			Cyclohexane, 1-ethenyl-1-methyl-...	204	C15H24	000515-13-9	74
5			2,6,10,10-Tetramethylbicyclo[7.2...	204	C15H24	136296-37-2	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
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Sample : P3675-17
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ALS Vial : 28 Sample Multiplier: 1

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ClientSampleId :
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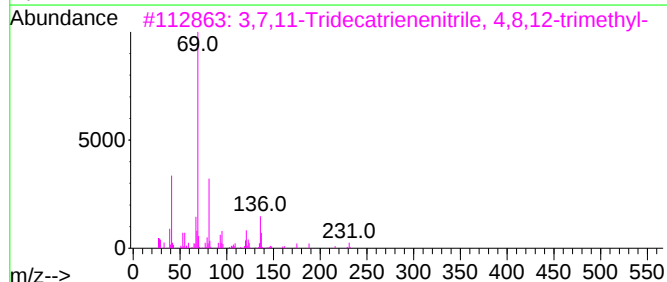
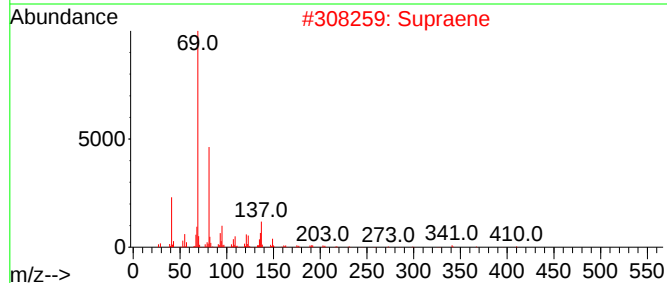
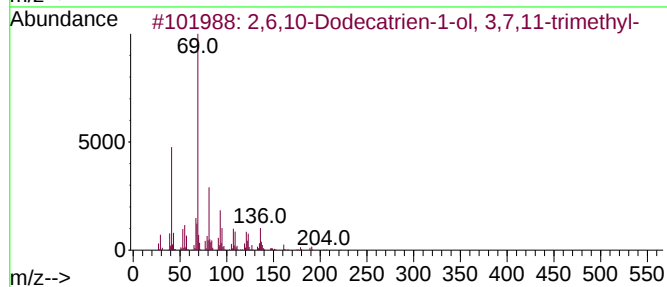
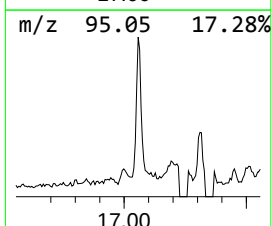
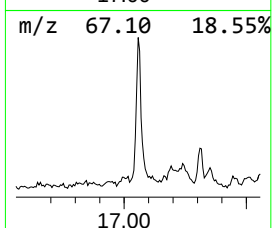
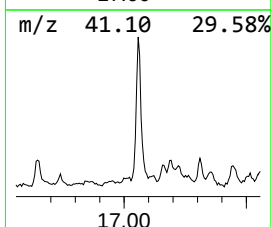
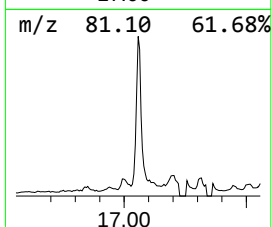
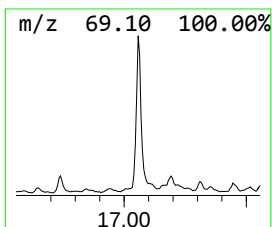
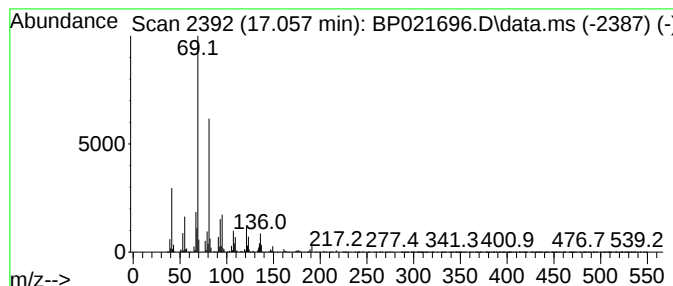
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 2,6,10-Dodecatrien-1-ol, 3,... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.057	2.93 ng/ul	1098380	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	86		
2	Supraene	410	C30H50	007683-64-9	78		
3	3,7,11-Tridecatrienitrile, 4,8...	231	C16H25N	006006-01-5	64		
4	Cyclopropanemethanol, 2-methyl-2...	168	C11H20O	098678-70-7	59		
5	5,9,13-Pentadecatrien-2-one, 6,1...	262	C18H30O	001117-52-8	56		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
Operator : RC/JU
Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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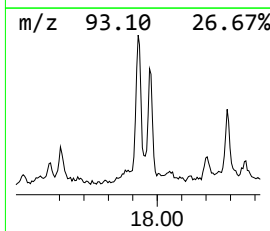
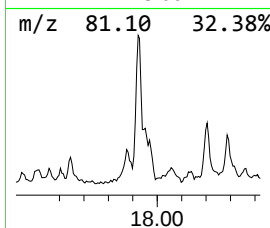
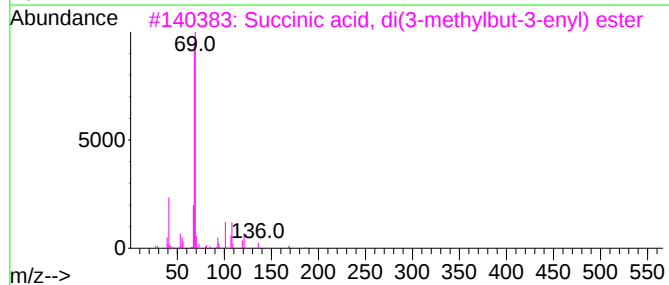
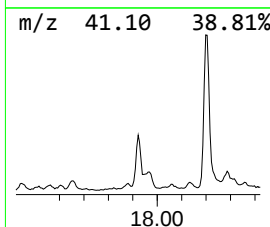
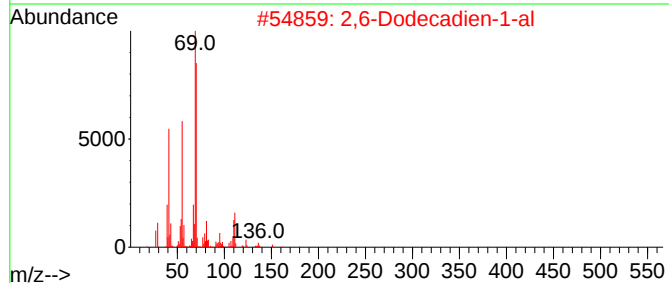
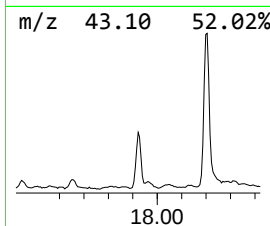
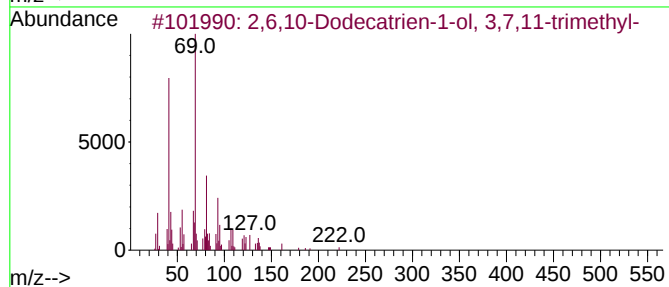
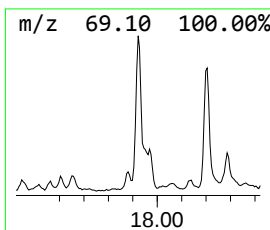
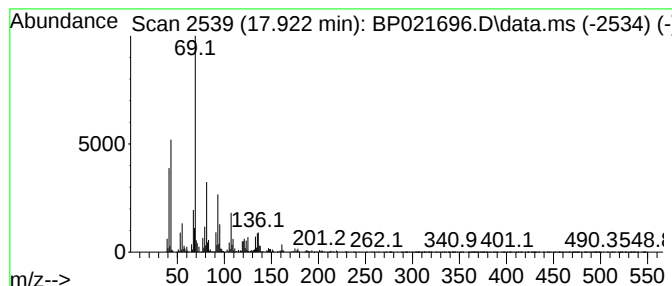
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 2,6-Dodecadien-1-al Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.922	3.05 ng/ul	1144950	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	80		
2	2,6-Dodecadien-1-al	180	C12H20O	021662-13-5	52		
3	Succinic acid, di(3-methylbut-3-...	254	C14H22O4	1010353-45-6	50		
4	1,3,5-Tris(cyanomethyl)-S-hexahy...	204	C9H12N6	004560-87-6	47		
5	Cyclopentane, bromo-	148	C5H9Br	000137-43-9	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
Operator : RC/JU
Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ1

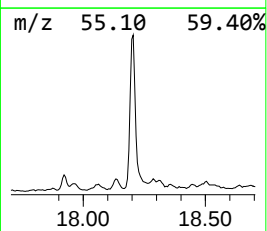
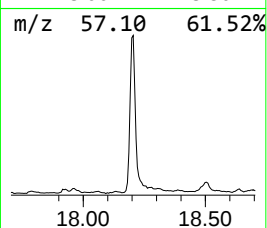
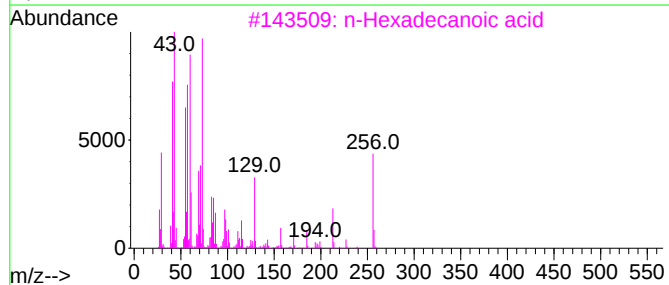
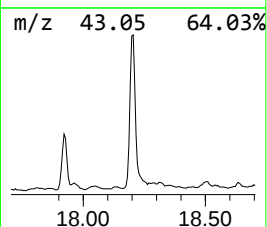
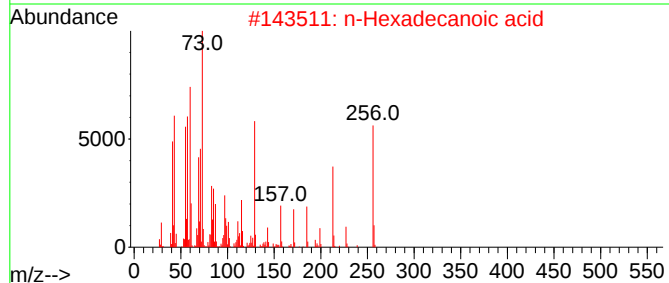
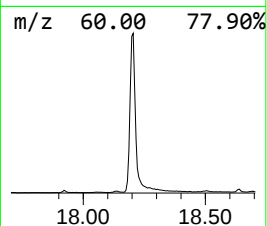
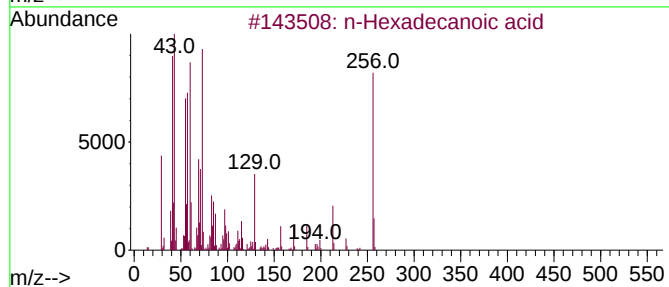
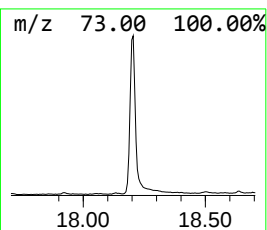
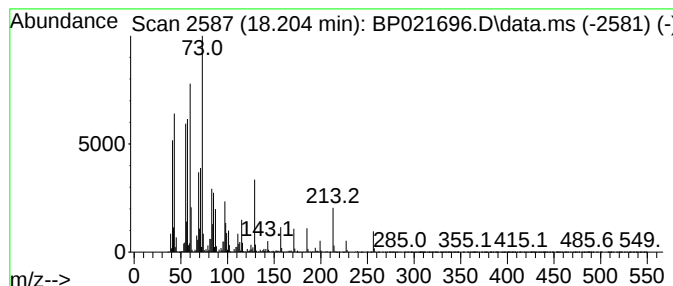
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.204	13.46 ng/ul	5052070	Phenanthrene-d10	17.245

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	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
Operator : RC/JU
Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ1

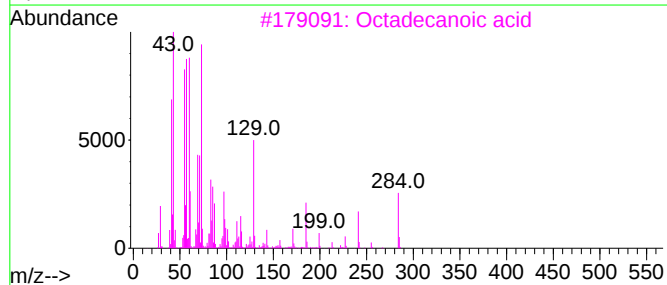
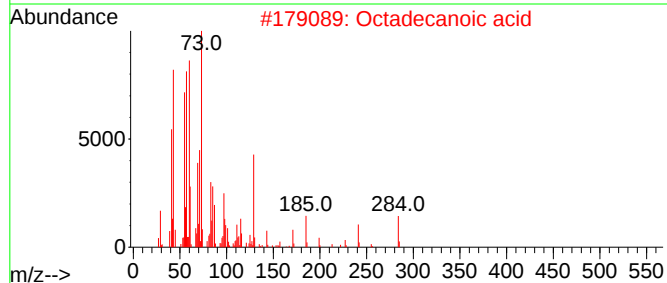
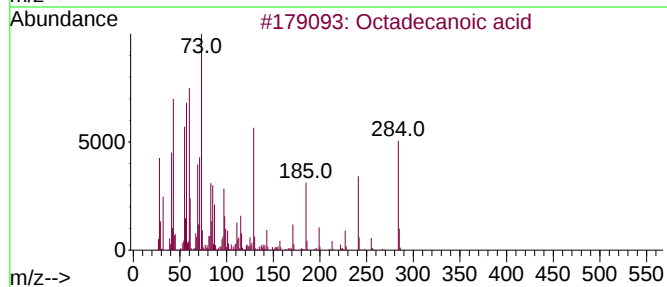
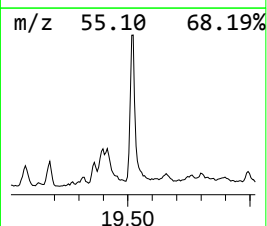
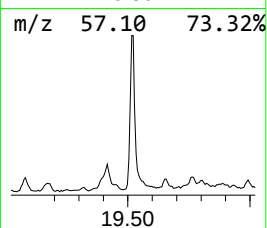
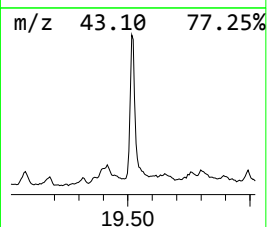
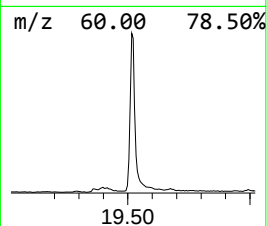
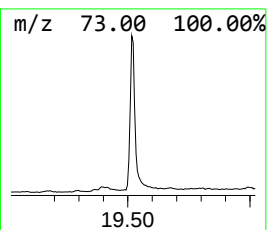
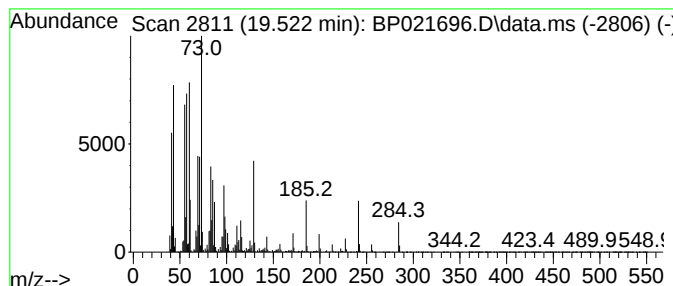
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.522	10.37 ng/ul	4038290	Chrysene-d12	21.698

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	98
4		Octadecanoic acid	284	C18H36O2	000057-11-4	95
5		Octadecanoic acid	284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
Operator : RC/JU
Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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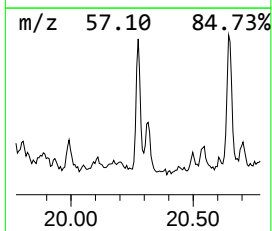
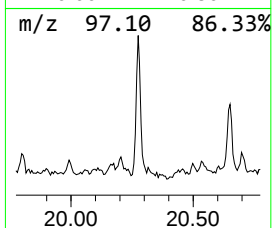
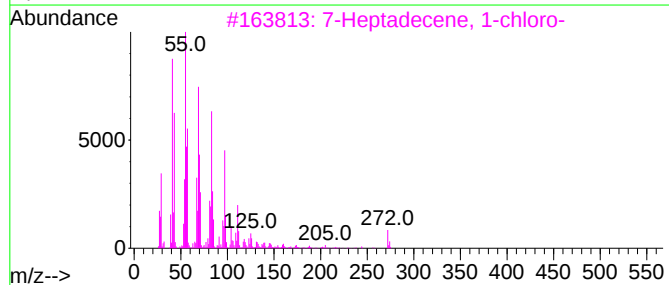
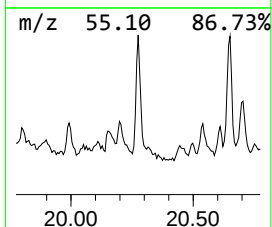
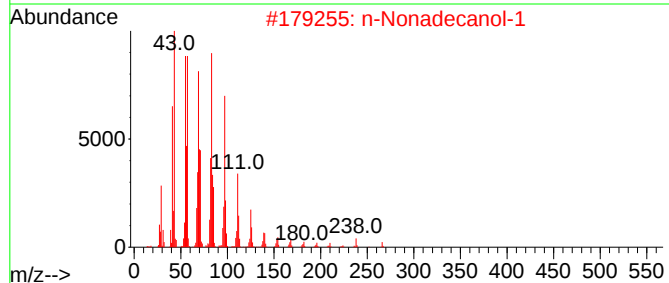
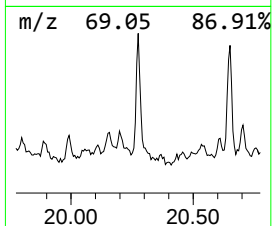
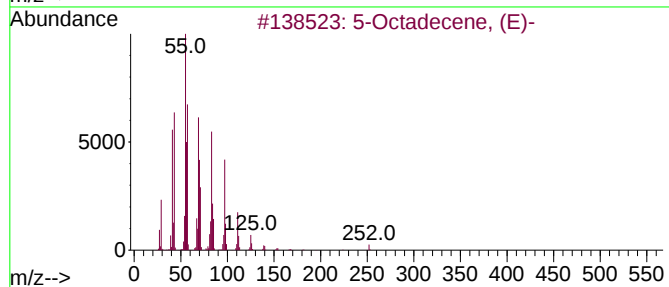
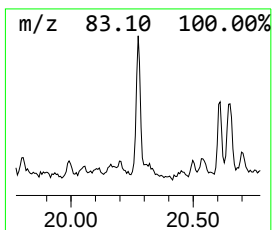
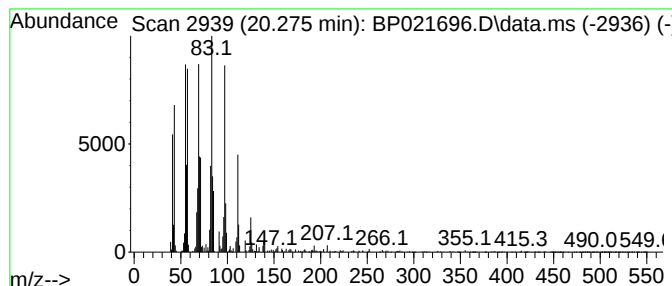
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.275	2.76 ng/ul	1076090	Chrysene-d12	21.698

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Octadecene, (E)-		252	C18H36	007206-21-5	98
2	n-Nonadecanol-1		284	C19H40O	001454-84-8	94
3	7-Heptadecene, 1-chloro-		272	C17H33Cl	056554-78-0	93
4	n-Heptadecanol-1		256	C17H36O	001454-85-9	91
5	1-Heneicosanol		312	C21H44O	015594-90-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
Operator : RC/JU
Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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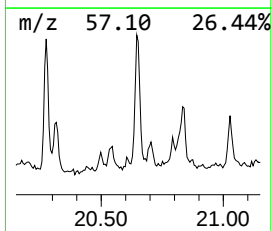
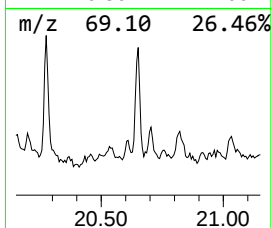
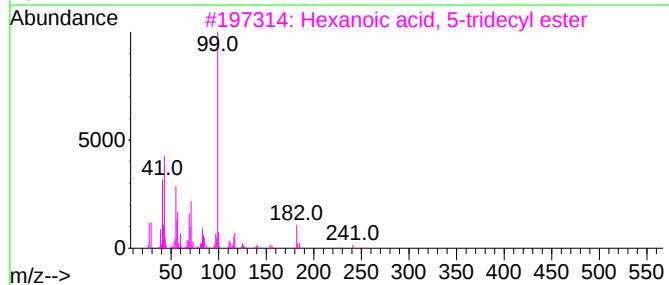
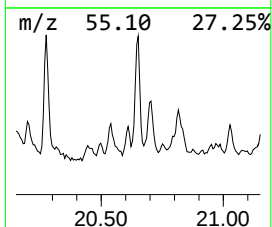
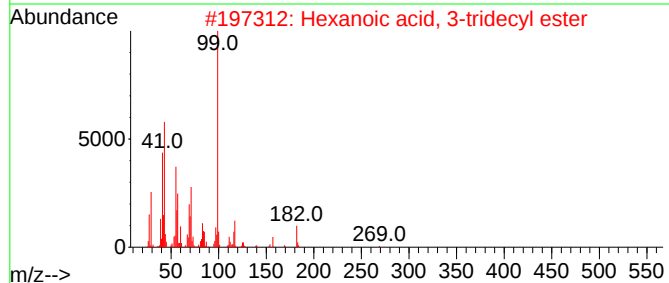
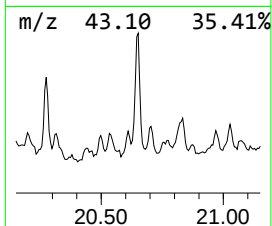
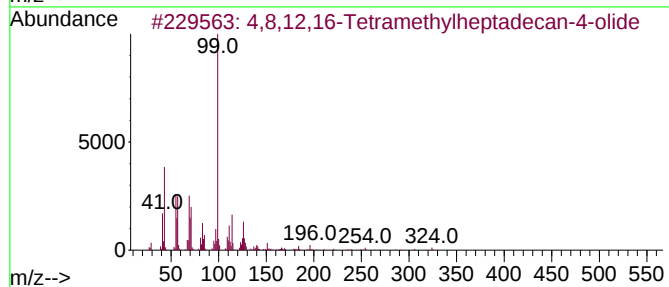
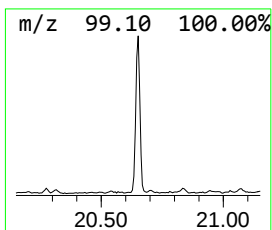
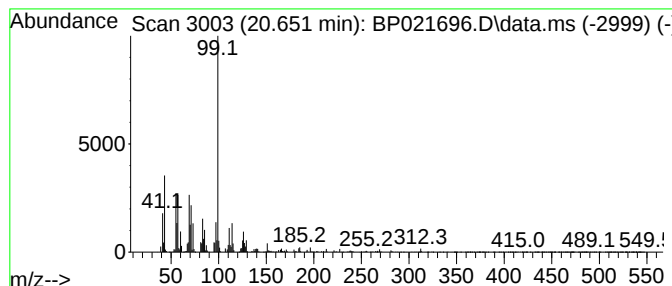
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 4,8,12,16-Tetramethylheptad... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.651	2.72 ng/ul	1058850	Chrysene-d12	21.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,8,12,16-Tetramethylheptadecan-...	324	C21H40O2	096168-15-9	74		
2	Hexanoic acid, 3-tridecyl ester	298	C19H38O2	1000279-29-2	64		
3	Hexanoic acid, 5-tridecyl ester	298	C19H38O2	1000279-29-4	64		
4	4,8,12,16-Tetramethylheptadecan-...	324	C21H40O2	096168-15-9	58		
5	Hexanoic acid, 6-tridecyl ester	298	C19H38O2	1000279-29-5	53		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
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Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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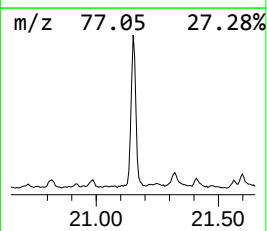
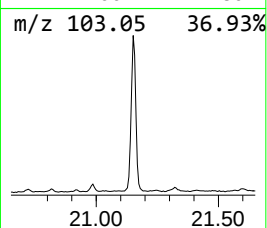
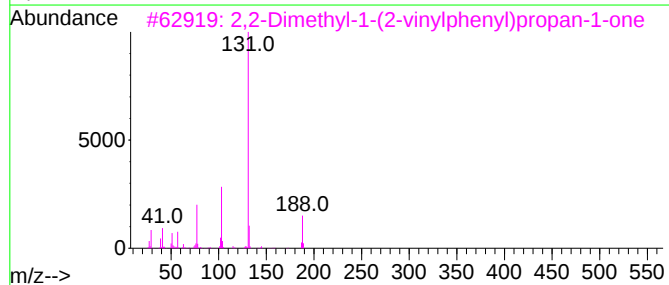
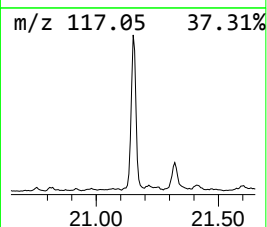
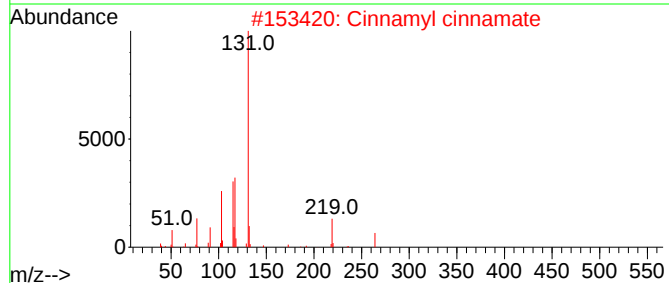
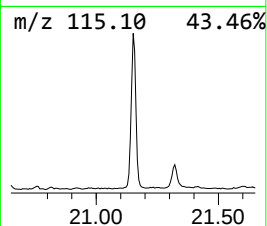
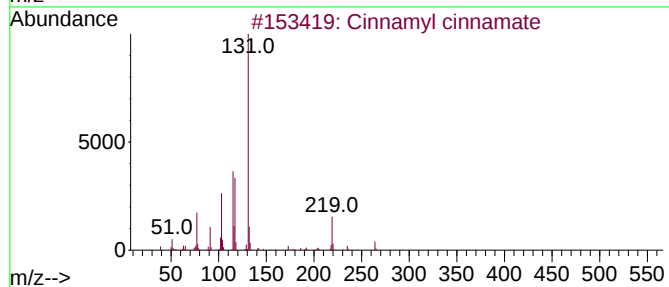
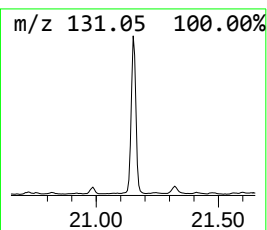
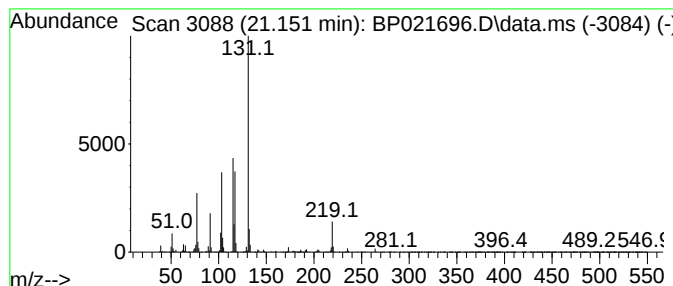
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Cinnamyl cinnamate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.151	8.16 ng/ul	3177420	Chrysene-d12	21.698

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		2,2-Dimethyl-1-(2-vinylphenyl)pr...	188	C13H16O	1000210-99-9	60
4		(E)-4-Bromo-4-methyl-1-phenylpen...	252	C12H13BrO	1000443-02-8	50
5		Cinnamoylglycine, methyl ester	219	C12H13NO3	040778-04-9	50



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Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
Operator : RC/JU
Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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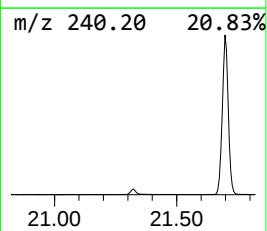
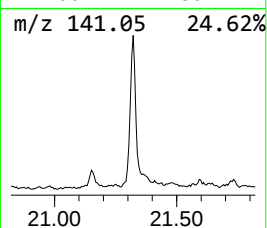
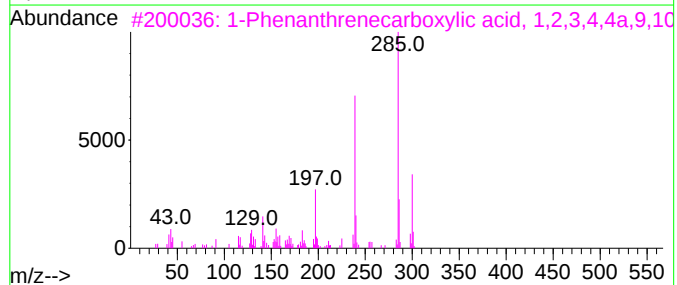
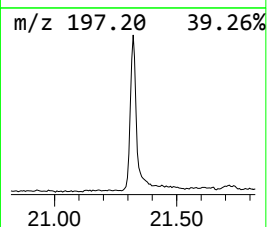
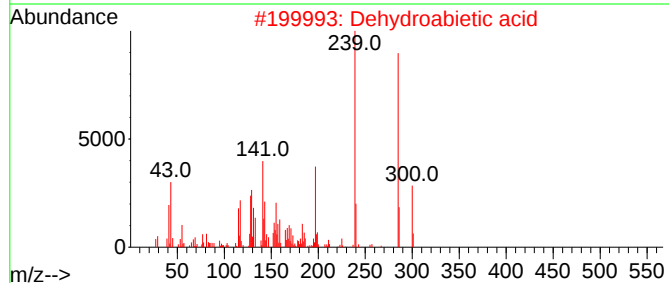
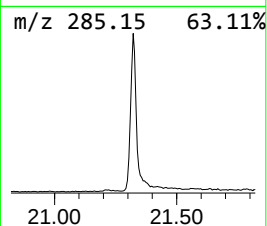
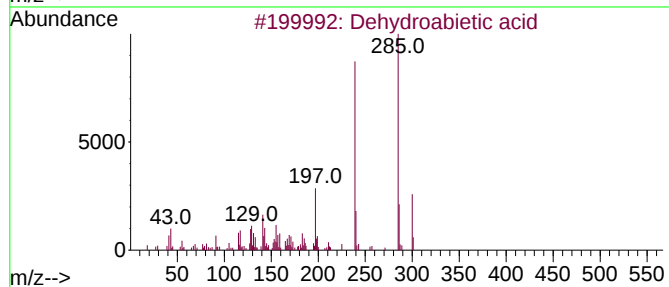
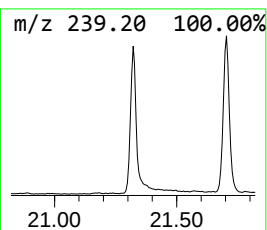
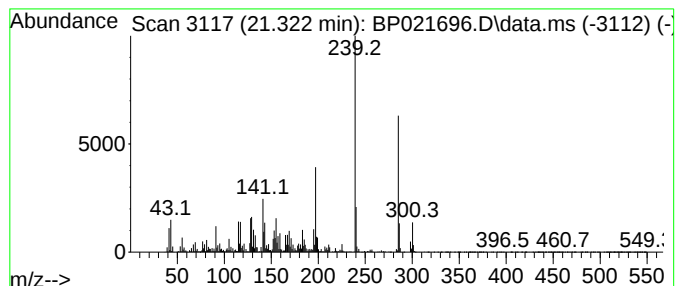
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Dehydroabietic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.322	7.98 ng/ul	3105850	Chrysene-d12	21.698

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	99
3			1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	94
4			Dehydroabietic acid	300	C20H28O2	001740-19-8	94
5			1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
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Sample : P3675-17
Misc :
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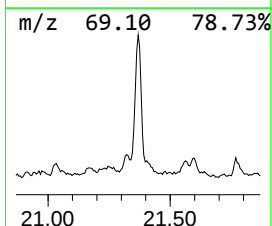
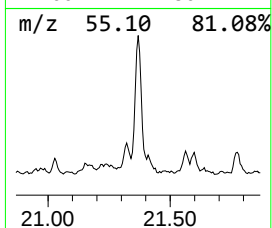
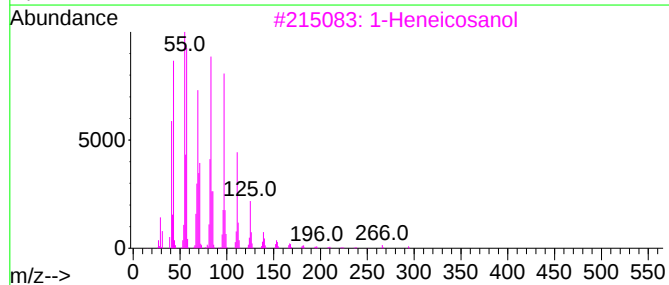
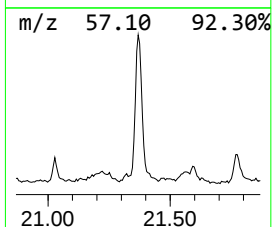
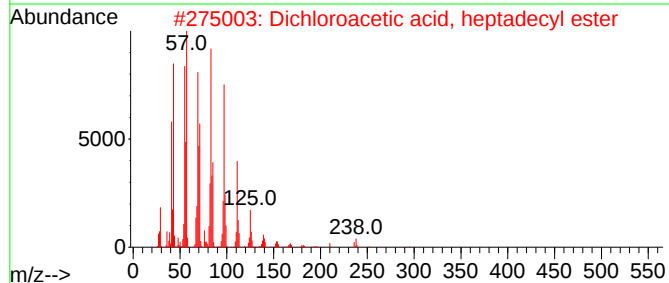
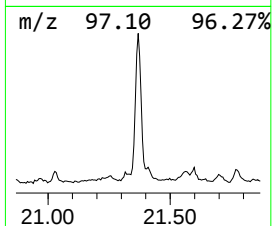
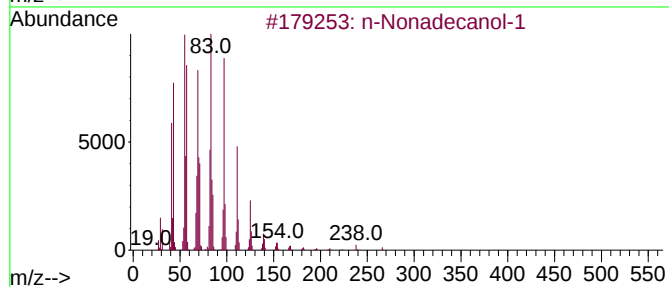
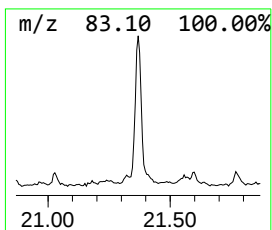
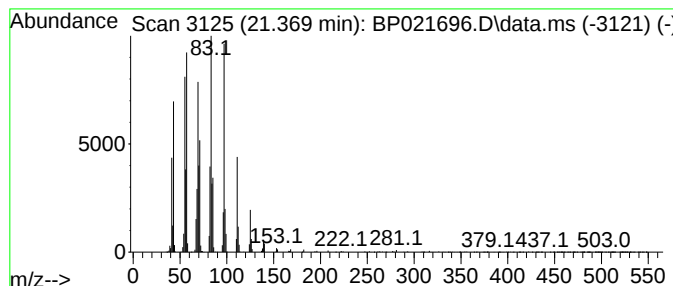
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 n-Nonadecanol-1 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.369	4.45 ng/ul	1732740	Chrysene-d12	21.698

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Nonadecanol-1	284	C19H40O	001454-84-8	94
2		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	93
3		1-Heneicosanol	312	C21H44O	015594-90-8	91
4		Behenic alcohol	326	C22H46O	000661-19-8	91
5		1-Nonadecene	266	C19H38	018435-45-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
Operator : RC/JU
Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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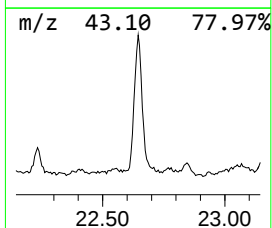
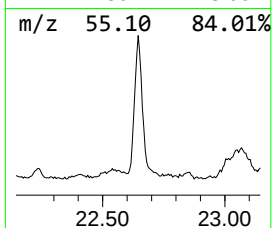
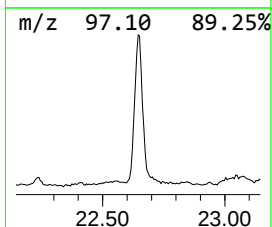
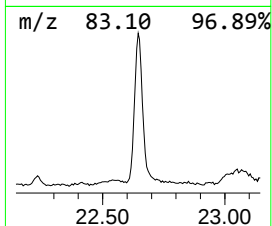
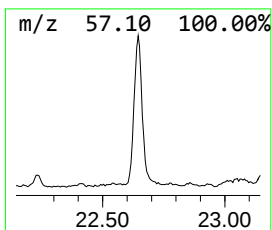
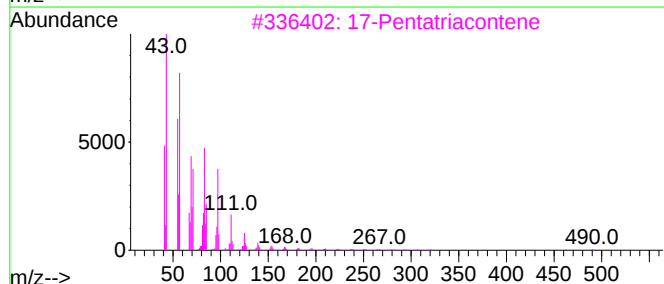
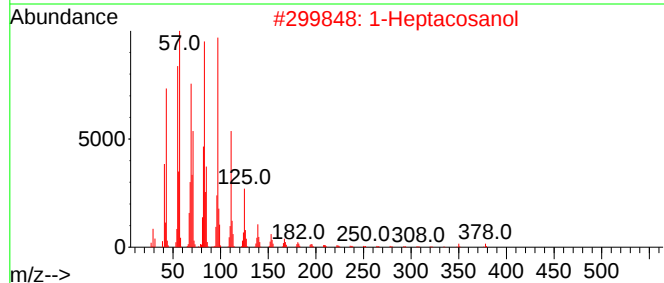
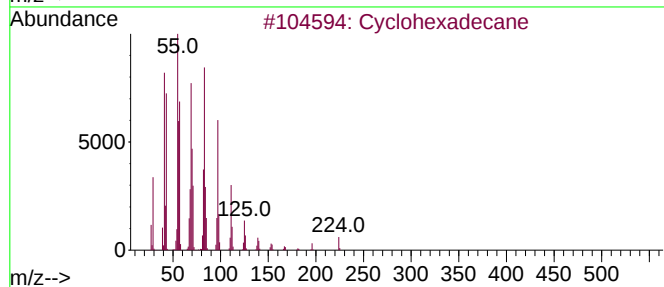
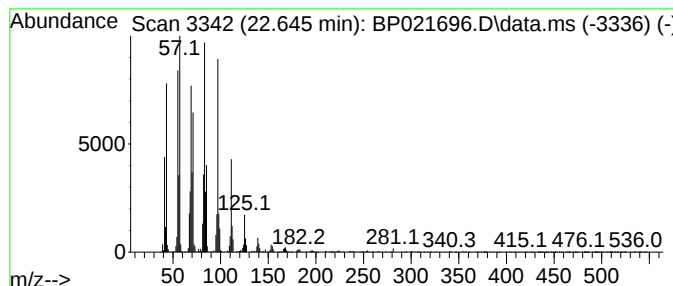
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Cyclic22.645 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.645	8.86 ng/ul	3449790	Chrysene-d12	21.698

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexadecane	224	C16H32	000295-65-8	94
2		1-Heptacosanol	396	C27H56O	002004-39-9	94
3		17-Pentatriacontene	491	C35H70	006971-40-0	91
4		1-Heneicosanol	312	C21H44O	015594-90-8	91
5		Behenic alcohol	326	C22H46O	000661-19-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
Operator : RC/JU
Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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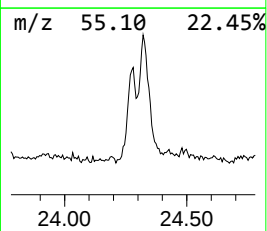
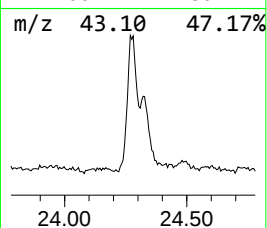
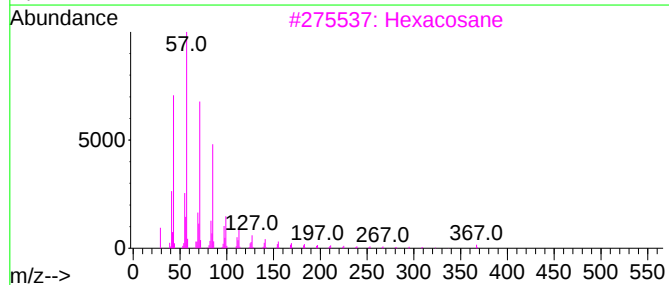
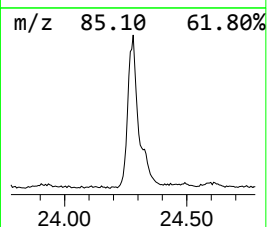
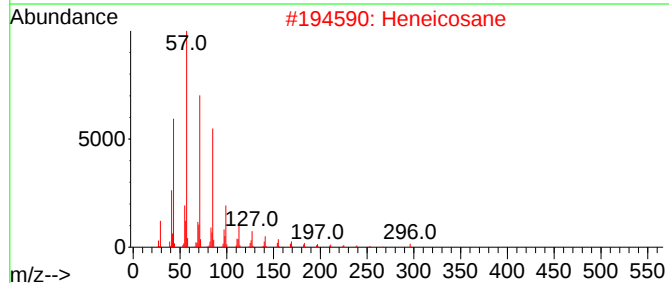
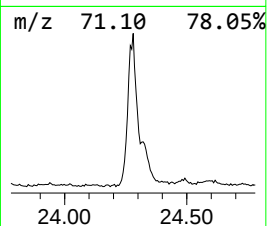
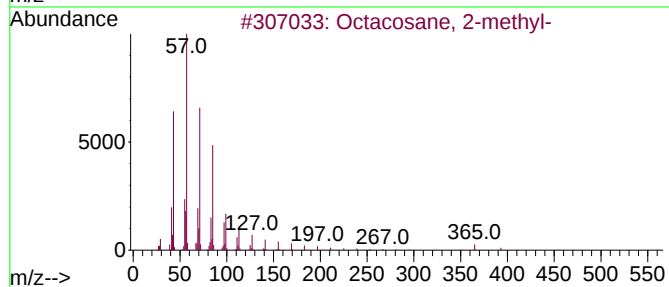
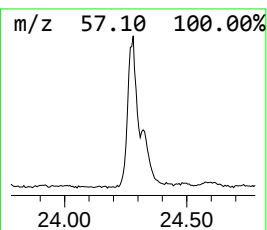
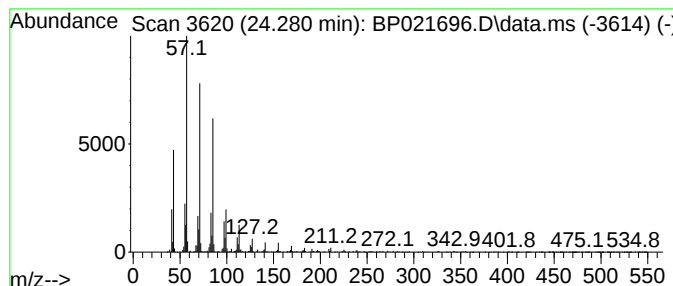
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.280	3.16 ng/ul	1298870	Perylene-d12	25.110

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Hexacosane	366	C26H54	000630-01-3	90
4		2-Methyltetracosane	352	C25H52	001560-78-7	90
5		Tetradecane	198	C14H30	000629-59-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
Operator : RC/JU
Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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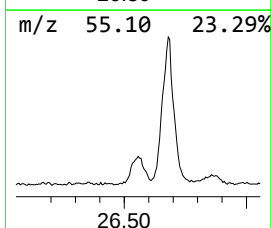
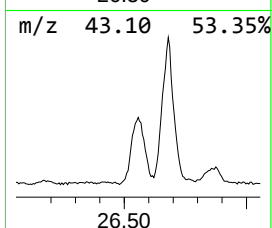
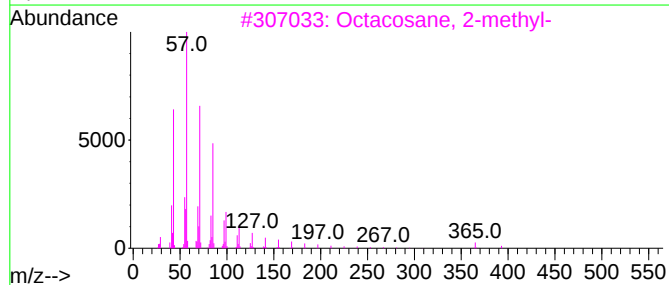
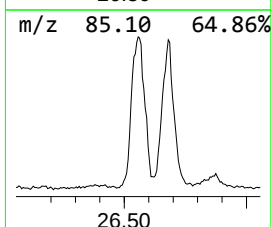
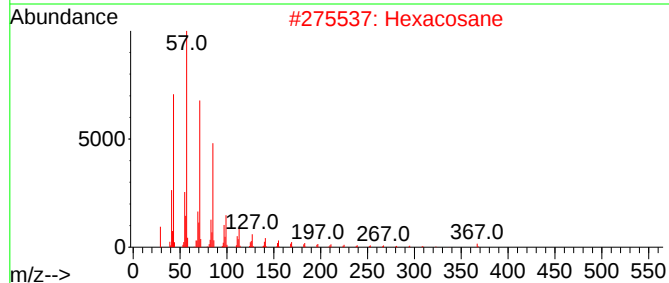
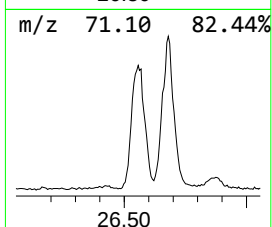
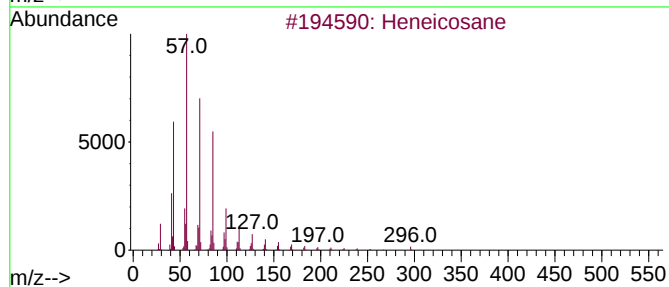
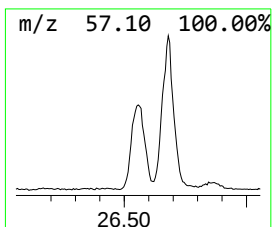
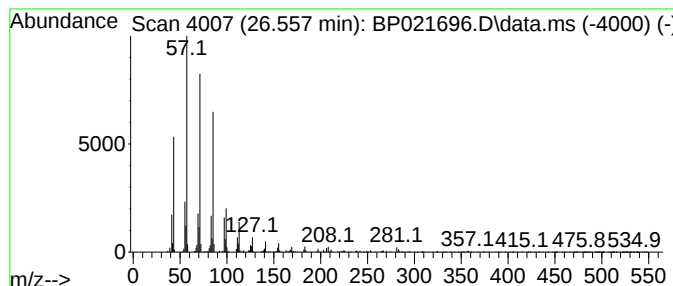
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.557	3.80 ng/ul	1561280	Perylene-d12	25.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Hexacosane	366	C26H54	000630-01-3	91
3			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4			2-Methyltetracosane	352	C25H52	001560-78-7	90
5			Pentacosane	352	C25H52	000629-99-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
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Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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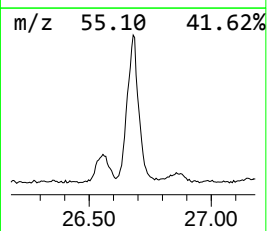
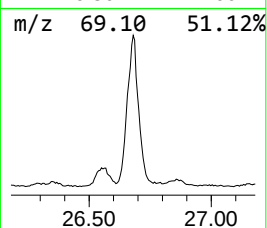
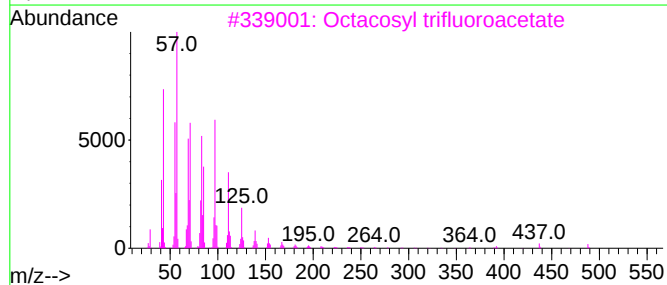
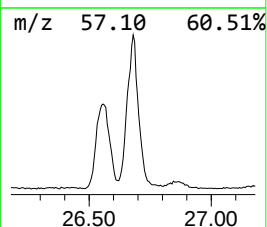
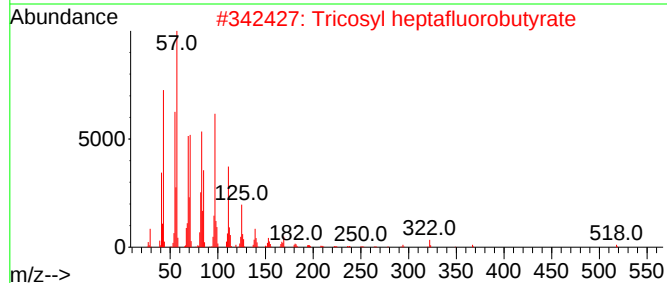
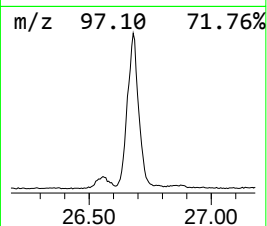
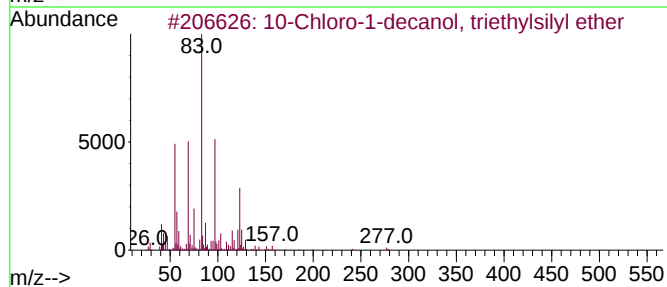
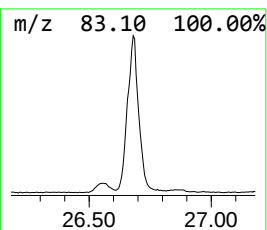
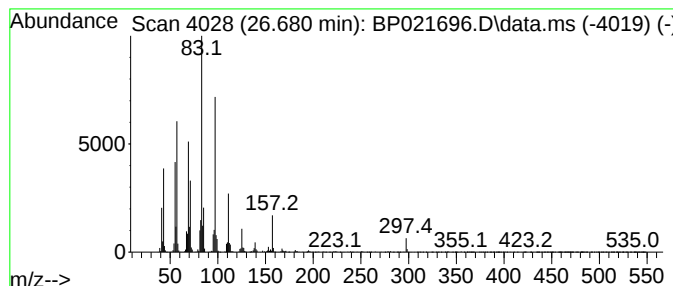
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 unknown-02 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.680	14.82 ng/ul	6085740	Perylene-d12	25.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10-Chloro-1-decanol, triethylsil...	306	C16H35ClOSi	1000447-39-3	47		
2	Tricosyl heptafluorobutyrate	536	C27H47F7O2	1000351-83-4	46		
3	Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	45		
4	Tetratriacontyl heptafluorobutyrate	691	C38H69F7O2	1000351-84-1	43		
5	Tetratriacontyl pentafluoropropi...	641	C37H69F5O2	1000351-81-5	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021696.D
Acq On : 28 Aug 2024 09:35
Operator : RC/JU
Sample : P3675-17
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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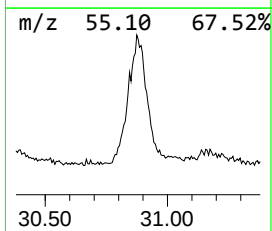
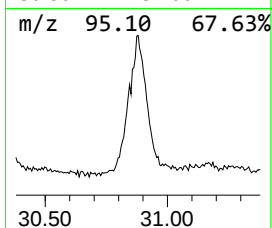
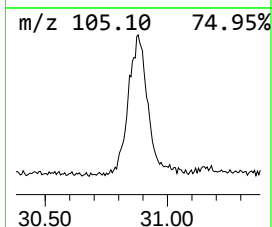
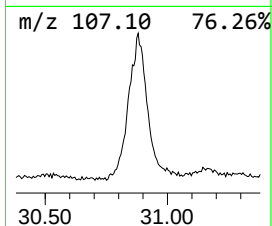
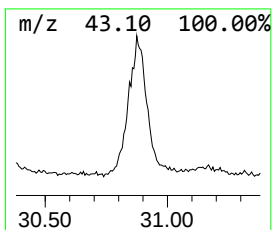
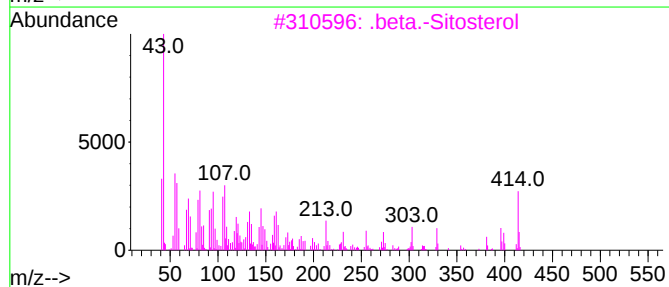
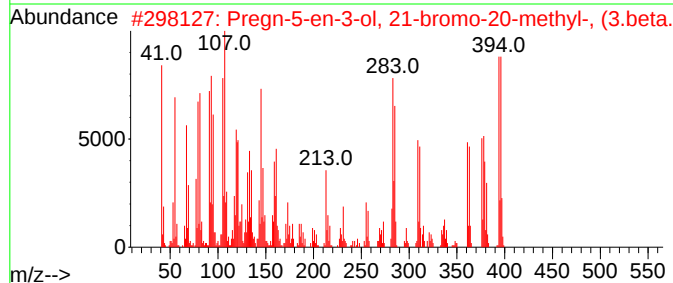
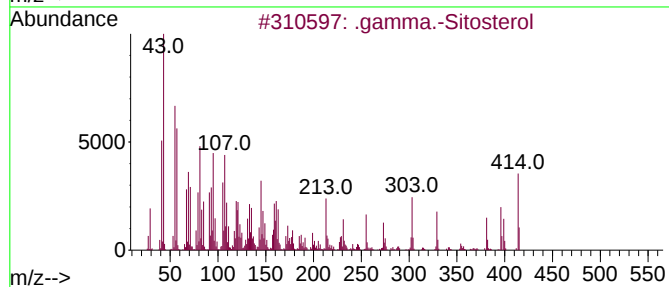
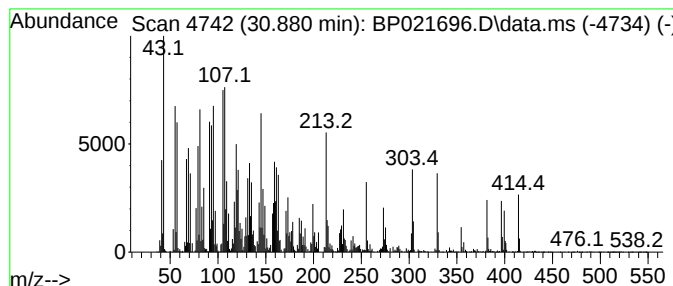
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 25 .gamma.-Sitosterol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.880	14.12 ng/ul	5798870	Perylene-d12	25.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	gamma.-Sitosterol	414	C29H50O	000083-47-6	99	
2	Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	95		
3	.beta.-Sitosterol	414	C29H50O	000083-46-5	78		
4	(3S,8S,9S,10R,13R,14S,17R)-17-((...	414	C29H50O	029365-29-5	55		
5	.	gamma.-Sitosterol	414	C29H50O	000083-47-6	49	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
 Data File : BP021696.D
 Acq On : 28 Aug 2024 09:35
 Operator : RC/JU
 Sample : P3675-17
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCXZ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.017	33.5	ng/ul	4785930	1	7.793	2861570	20.0
Propylene Glycol	3.540	3.6	ng/ul	513823	1	7.793	2861570	20.0
(1R)-2,6,6-Trim...	6.499	7.1	ng/ul	1010430	1	7.793	2861570	20.0
Bicyclo[3.1.1]h...	7.246	3.9	ng/ul	563451	1	7.793	2861570	20.0
Ethanol, 2-(2-e...	7.393	2.7	ng/ul	385662	1	7.793	2861570	20.0
Benzoic acid	9.828	2.4	ng/ul	515561	2	10.575	4230820	20.0
unknown-01	9.887	2.7	ng/ul	576540	2	10.575	4230820	20.0
1-Phenoxypropan...	11.316	3.6	ng/ul	764753	2	10.575	4230820	20.0
Cyclohexane, 1-...	13.316	3.6	ng/ul	1146800	3	14.440	6312980	20.0
Caryophyllene	13.769	2.5	ng/ul	789105	3	14.440	6312980	20.0
2,6,10-Dodecatr...	17.057	2.9	ng/ul	1098380	4	17.245	7506930	20.0
2,6-Dodecadien-...	17.922	3.0	ng/ul	1144950	4	17.245	7506930	20.0
n-Hexadecanoic ...	18.204	13.5	ng/ul	5052070	4	17.245	7506930	20.0
Octadecanoic acid	19.522	10.4	ng/ul	4038290	5	21.698	7788590	20.0
(DEL) Alkane: S...	20.275	2.8	ng/ul	1076090	5	21.698	7788590	20.0
4,8,12,16-Tetra...	20.651	2.7	ng/ul	1058850	5	21.698	7788590	20.0
Cinnamyl cinnamate	21.151	8.2	ng/ul	3177420	5	21.698	7788590	20.0
Dehydroabietic ...	21.322	8.0	ng/ul	3105850	5	21.698	7788590	20.0
n-Nonadecanol-1	21.369	4.5	ng/ul	1732740	5	21.698	7788590	20.0
(DEL) Alkane: C...	22.645	8.9	ng/ul	3449790	5	21.698	7788590	20.0
(DEL) Alkane: S...	24.280	3.2	ng/ul	1298870	6	25.110	8211820	20.0
(DEL) Alkane: S...	26.557	3.8	ng/ul	1561280	6	25.110	8211820	20.0
unknown-02	26.680	14.8	ng/ul	6085740	6	25.110	8211820	20.0
.gamma.-Sitosterol	30.880	14.1	ng/ul	5798870	6	25.110	8211820	20.0