

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
 Data File : BM044975.D
 Acq On : 06 Apr 2024 17:51
 Operator : MA/JU
 Sample : P2018-10
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BH5F1

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_M\Methods\SFAM-EPA-BM040524.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM044975.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.557	61	80	83	rBV2	199233	683125	5.24%	0.489%
2	3.963	119	149	152	rBV	389634	2046450	15.70%	1.466%
3	5.210	345	361	368	rBV	145166	450189	3.45%	0.322%
4	6.716	603	617	625	rBV	170035	473801	3.63%	0.339%
5	6.816	627	634	645	rVV	286795	525224	4.03%	0.376%
6	6.987	658	663	680	rVB	221234	369638	2.84%	0.265%
7	7.169	681	694	702	rVV	302686	493416	3.79%	0.353%
8	7.634	766	773	780	rBV	381406	607032	4.66%	0.435%
9	8.345	885	894	900	rVV	333723	580431	4.45%	0.416%
10	8.792	964	970	988	rVB	253511	470875	3.61%	0.337%
11	9.504	1085	1091	1102	rBV	220890	400996	3.08%	0.287%
12	9.751	1115	1133	1143	rBV	247188	898932	6.90%	0.644%
13	10.039	1175	1182	1194	rVB	325367	603423	4.63%	0.432%
14	10.410	1238	1245	1251	rVV	508095	851288	6.53%	0.610%
15	12.551	1603	1609	1616	rVB	124555	211581	1.62%	0.152%
16	12.639	1616	1624	1632	rBV2	98273	209586	1.61%	0.150%
17	13.222	1712	1723	1730	rVV	1528077	2364092	18.14%	1.693%
18	13.457	1757	1763	1769	rVV	122534	276974	2.12%	0.198%
19	13.710	1793	1806	1811	rVV	516269	801673	6.15%	0.574%
20	13.974	1845	1851	1859	rVB	634958	989954	7.59%	0.709%
21	14.163	1878	1883	1888	rVV	207269	296069	2.27%	0.212%
22	14.280	1897	1903	1910	rVV	732940	1130575	8.67%	0.810%
23	14.533	1936	1946	1954	rBV5	310387	651442	5.00%	0.467%
24	14.727	1974	1979	1985	rVV5	115271	237567	1.82%	0.170%
25	14.810	1985	1993	2001	rVV2	401246	884290	6.78%	0.633%
26	14.898	2001	2008	2018	rVB2	182871	396400	3.04%	0.284%
27	15.116	2039	2045	2050	rBV	454437	665859	5.11%	0.477%
28	15.286	2069	2074	2081	rVB	767818	1136104	8.72%	0.814%
29	15.421	2091	2097	2106	rBV	260592	435229	3.34%	0.312%
30	15.680	2129	2141	2147	rBV3	301967	706177	5.42%	0.506%
31	15.780	2147	2158	2162	rBV	8473634	13034999	100.00%	9.337%
32	15.863	2167	2172	2178	rBV2	148982	228433	1.75%	0.164%
33	16.021	2193	2199	2211	rBV2	995031	1530087	11.74%	1.096%
34	16.362	2252	2257	2267	rVB	506057	730640	5.61%	0.523%
35	16.457	2267	2273	2284	rBV2	1089963	1780878	13.66%	1.276%
36	16.933	2345	2354	2364	rBV	790232	1552242	11.91%	1.112%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
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37	17.045	2366	2373	2378	rBV2	1350741	2112641	16.21%	1.513%
38	17.145	2384	2390	2400	rVB	930969	1516938	11.64%	1.087%
39	17.286	2410	2414	2422	rVB2	210848	304803	2.34%	0.218%
40	17.698	2479	2484	2486	rBV	180674	252525	1.94%	0.181%
41	17.733	2486	2490	2495	rVB	210504	285720	2.19%	0.205%
42	17.839	2503	2508	2514	rBV3	150155	292785	2.25%	0.210%
43	17.898	2514	2518	2524	rVB	207412	272858	2.09%	0.195%
44	17.974	2524	2531	2542	rBV	5118813	8086652	62.04%	5.793%
45	18.215	2568	2572	2573	rBV	235679	253295	1.94%	0.181%
46	18.251	2573	2578	2582	rVV	1928217	2725778	20.91%	1.953%
47	18.292	2582	2585	2592	rVB2	232238	358749	2.75%	0.257%
48	18.392	2598	2602	2608	rBV4	166452	245584	1.88%	0.176%
49	18.456	2608	2613	2619	rVB3	293013	505119	3.88%	0.362%
50	18.568	2628	2632	2638	rVB3	187979	268510	2.06%	0.192%
51	18.633	2638	2643	2651	rBV5	284447	582201	4.47%	0.417%
52	18.880	2680	2685	2688	rBV3	422574	613633	4.71%	0.440%
53	18.915	2688	2691	2695	rVB3	438327	543303	4.17%	0.389%
54	19.039	2707	2712	2719	rVV3	379788	756781	5.81%	0.542%
55	19.121	2719	2726	2729	rVV	5203354	10110232	77.56%	7.242%
56	19.151	2729	2731	2745	rVV2	7058095	10298728	79.01%	7.377%
57	19.256	2745	2749	2762	rVB	1426207	1945865	14.93%	1.394%
58	19.421	2773	2777	2779	rBV	472743	645517	4.95%	0.462%
59	19.451	2779	2782	2790	rVV	1136366	1471921	11.29%	1.054%
60	19.780	2830	2838	2842	rBV5	543925	982800	7.54%	0.704%
61	19.856	2848	2851	2856	rVB	241054	296213	2.27%	0.212%
62	19.992	2871	2874	2880	rVB4	180136	292561	2.24%	0.210%
63	20.156	2895	2902	2910	rBV6	394739	1043073	8.00%	0.747%
64	20.250	2910	2918	2921	rBV2	845855	1416647	10.87%	1.015%
65	20.292	2921	2925	2928	rVV4	922213	1428571	10.96%	1.023%
66	20.345	2930	2934	2939	rVV3	1062732	1751650	13.44%	1.255%
67	20.403	2941	2944	2946	rVV	319738	469969	3.61%	0.337%
68	20.445	2946	2951	2953	rVV	2000638	2913689	22.35%	2.087%
69	20.468	2953	2955	2958	rVV	1654001	1898296	14.56%	1.360%
70	20.498	2958	2960	2965	rVV2	485224	795458	6.10%	0.570%
71	20.545	2965	2968	2972	rVV3	268092	451163	3.46%	0.323%
72	20.592	2973	2976	2983	rVB2	290232	487494	3.74%	0.349%
73	20.792	3007	3010	3012	rBV2	295300	381671	2.93%	0.273%
74	20.827	3012	3016	3024	rVB	1712766	2308893	17.71%	1.654%
75	20.956	3033	3038	3047	rBV2	2679902	4410477	33.84%	3.159%
76	21.121	3063	3066	3070	rVB2	155879	205757	1.58%	0.147%

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77	21.203	3075	3080	3084	rVV6	530112	947577	7.27%	0.679%
78	21.250	3084	3088	3091	rVV	1336096	1896697	14.55%	1.359%
79	21.292	3091	3095	3103	rVV2	1250031	1792583	13.75%	1.284%
80	21.362	3103	3107	3112	rVB2	587458	770579	5.91%	0.552%
81	21.580	3141	3144	3148	rVB	354524	364431	2.80%	0.261%
82	21.721	3165	3168	3172	rBV3	255228	363002	2.78%	0.260%
83	21.874	3191	3194	3204	rVB3	893548	1443324	11.07%	1.034%
84	21.997	3212	3215	3220	rVB	235669	272669	2.09%	0.195%
85	22.097	3229	3232	3237	rVB	492900	643461	4.94%	0.461%
86	22.197	3244	3249	3260	rBV2	840851	1454787	11.16%	1.042%
87	22.321	3267	3270	3274	rVB	463831	567884	4.36%	0.407%
88	22.444	3287	3291	3296	rVB	680718	942058	7.23%	0.675%
89	22.827	3352	3356	3360	rVB	737820	1007406	7.73%	0.722%
90	22.874	3360	3364	3375	rVB2	394367	644194	4.94%	0.461%
91	23.203	3415	3420	3425	rBV	1007165	1592883	12.22%	1.141%
92	23.338	3437	3443	3449	rVB	856934	1519169	11.65%	1.088%
93	23.474	3460	3466	3474	rVB	944745	1913625	14.68%	1.371%
94	24.044	3558	3563	3570	rVB	464545	879541	6.75%	0.630%
95	24.127	3572	3577	3587	rVB2	382928	825804	6.34%	0.592%
96	24.533	3640	3646	3656	rVB2	1228677	2825565	21.68%	2.024%
97	24.991	3719	3724	3736	rVB	439778	1033914	7.93%	0.741%
98	26.115	3907	3915	3942	rVB2	853918	2873525	22.04%	2.058%
99	26.527	3975	3985	4000	rVB2	1784547	5767041	44.24%	4.131%
100	27.785	4193	4199	4208	rBV3	542712	1569112	12.04%	1.124%

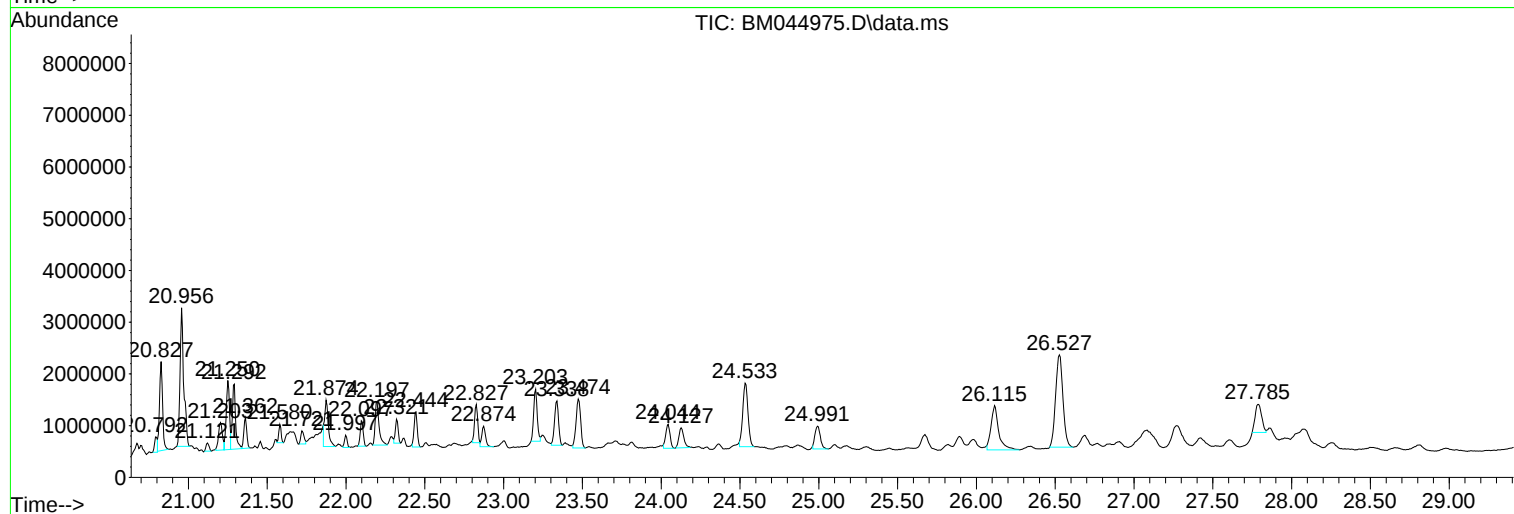
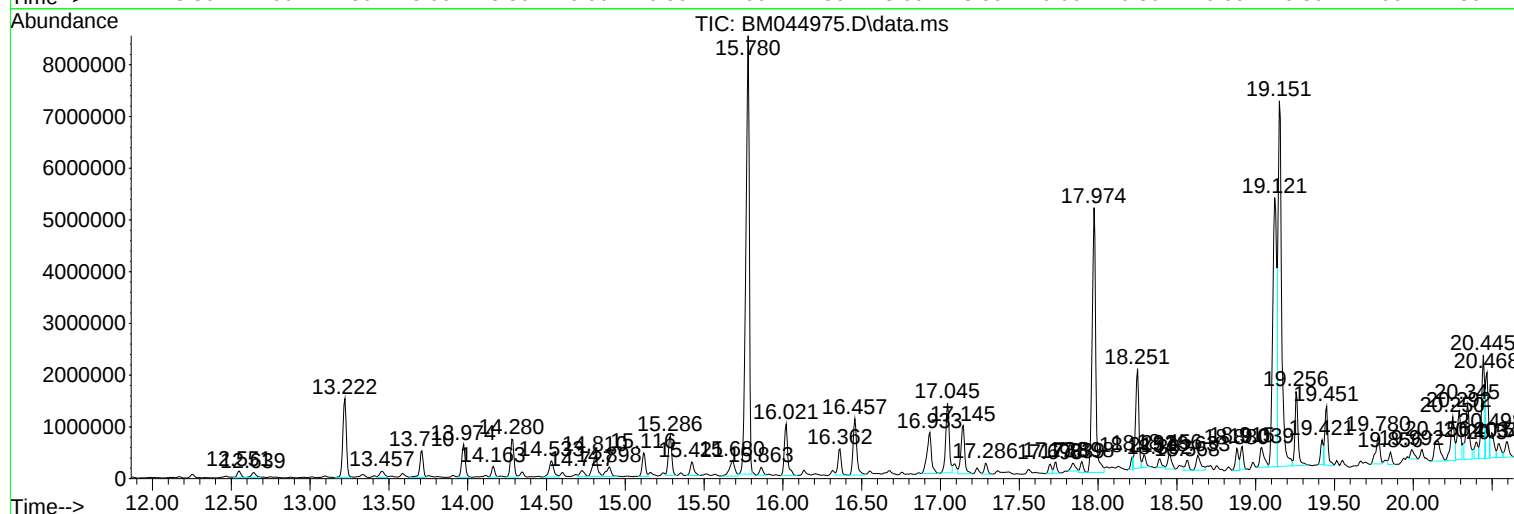
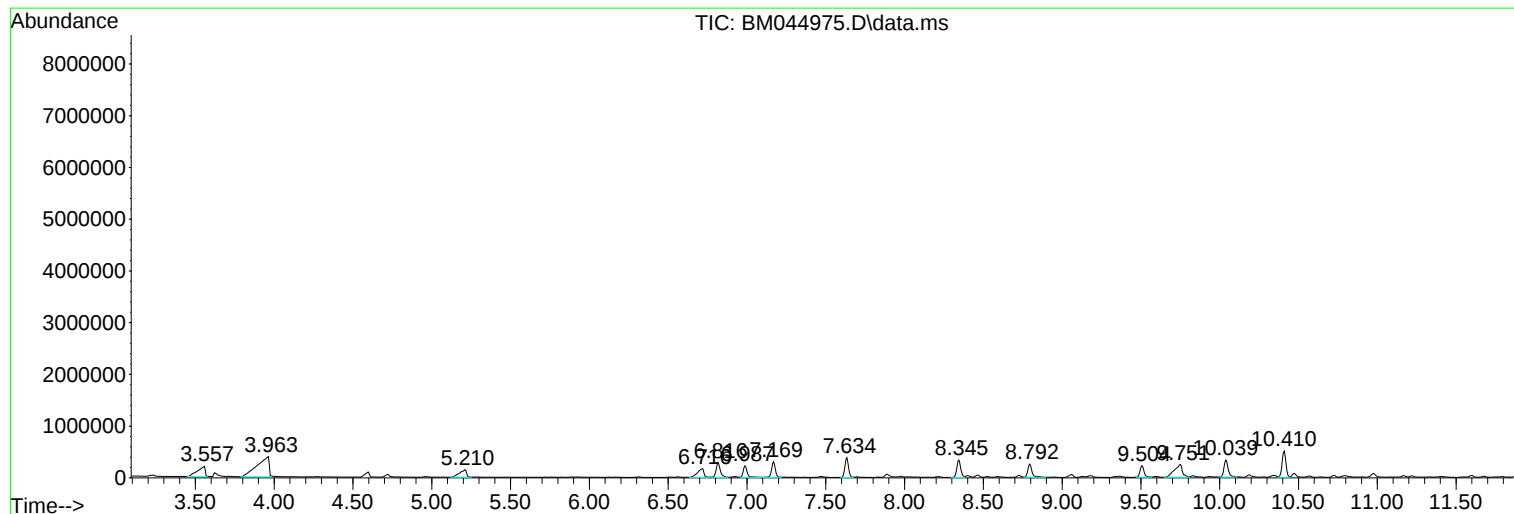
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION

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



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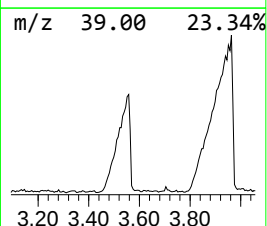
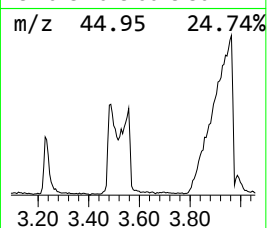
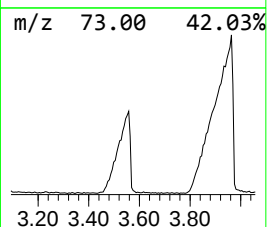
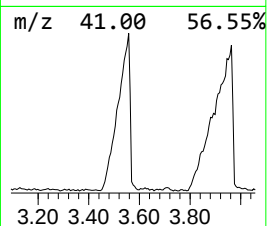
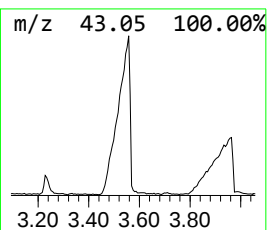
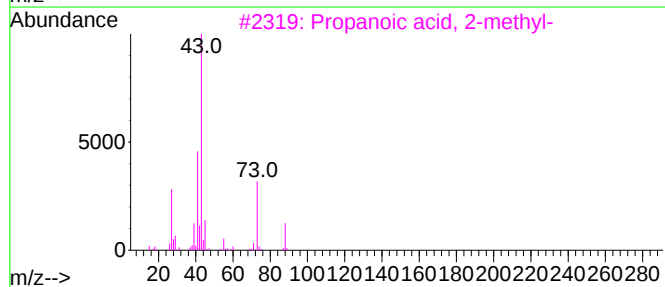
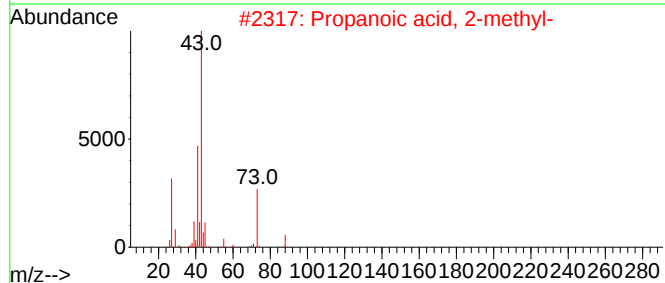
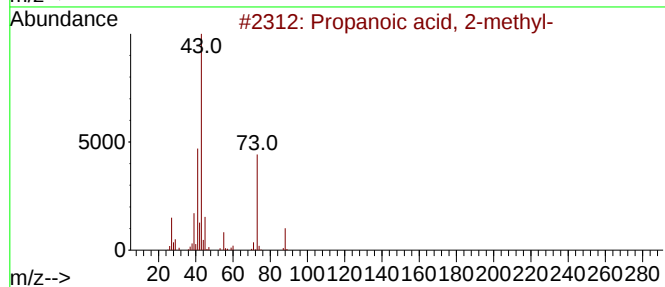
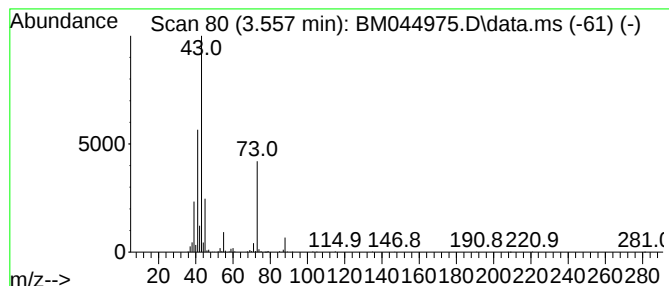
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propanoic acid, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.557	22.51 ng/ul	683125	1,4-Dichlorobenzene-d4	7.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	80
2			Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	72
3			Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	64
4			Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	59
5			Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	45



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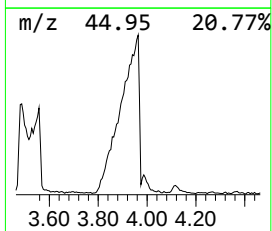
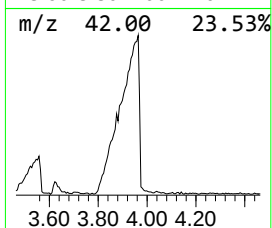
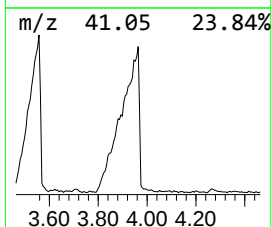
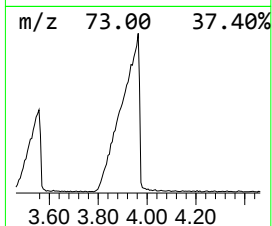
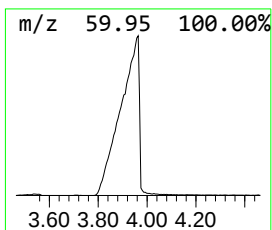
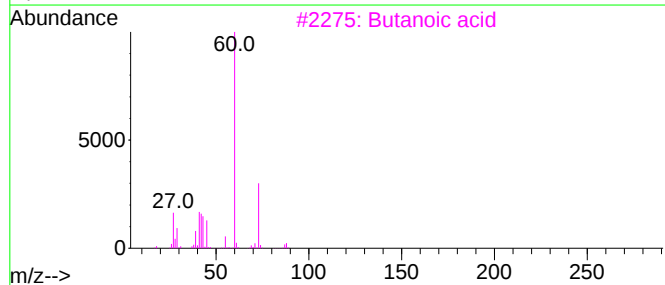
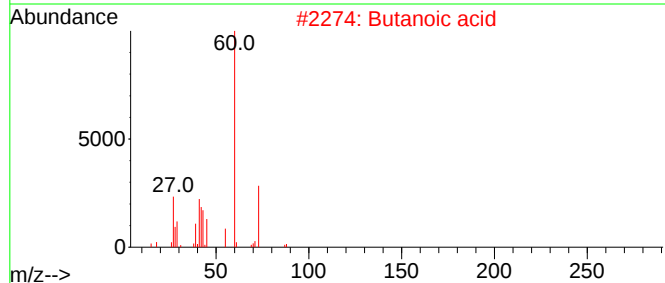
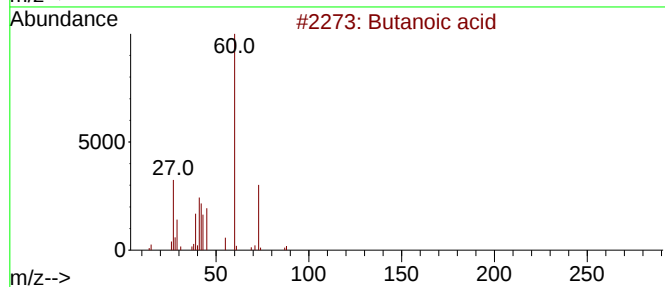
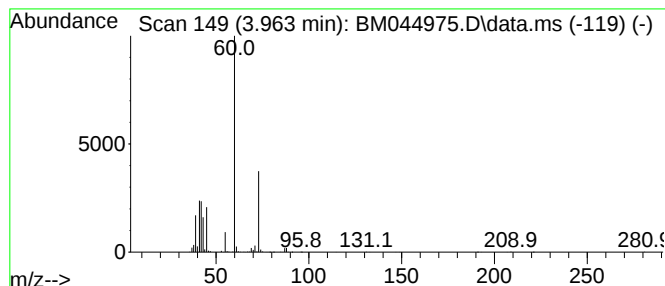
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.963	67.42 ng/ul	2046450	1,4-Dichlorobenzene-d4	7.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid	88	C4H8O2	000107-92-6	91
2			Butanoic acid	88	C4H8O2	000107-92-6	87
3			Butanoic acid	88	C4H8O2	000107-92-6	80
4			Butanoic acid	88	C4H8O2	000107-92-6	72
5			Silver butanoate	194	C4H7AgO2	005076-24-4	40



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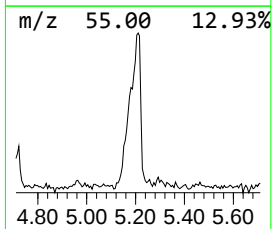
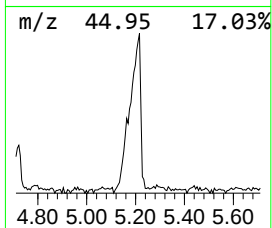
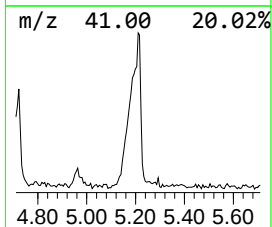
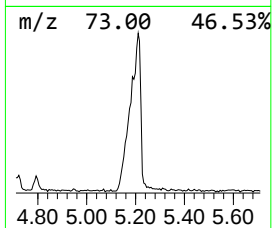
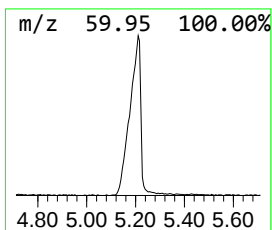
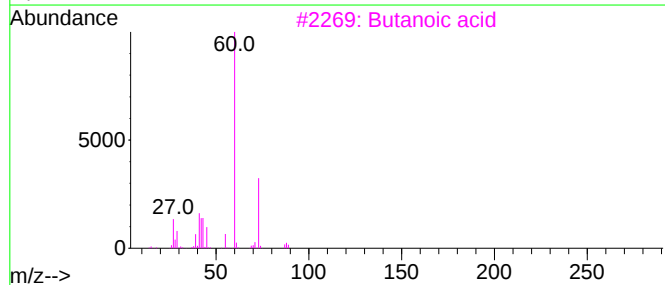
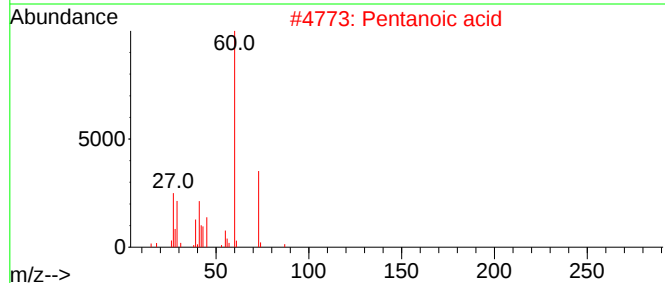
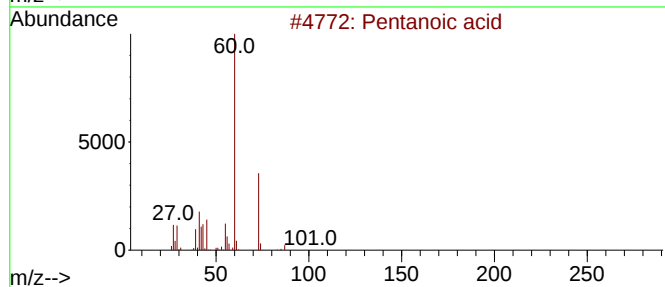
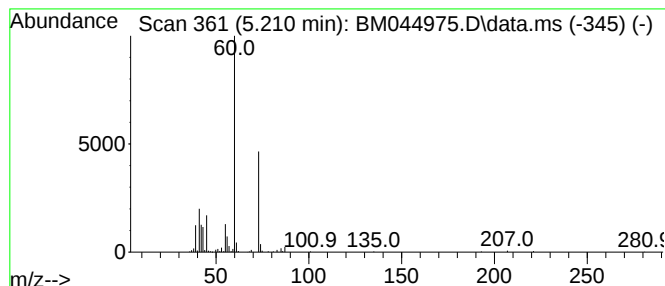
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Pentanoic acid Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.210	14.83 ng/ul	450189	1,4-Dichlorobenzene-d4	7.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanoic acid	102	C5H10O2	000109-52-4	90
2			Pentanoic acid	102	C5H10O2	000109-52-4	83
3			Butanoic acid	88	C4H8O2	000107-92-6	83
4			Pentanoic acid	102	C5H10O2	000109-52-4	78
5			Butanoic acid	88	C4H8O2	000107-92-6	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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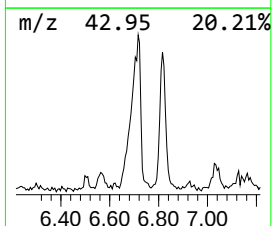
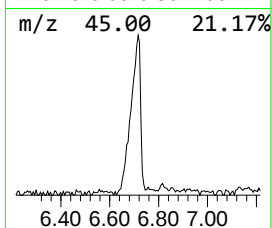
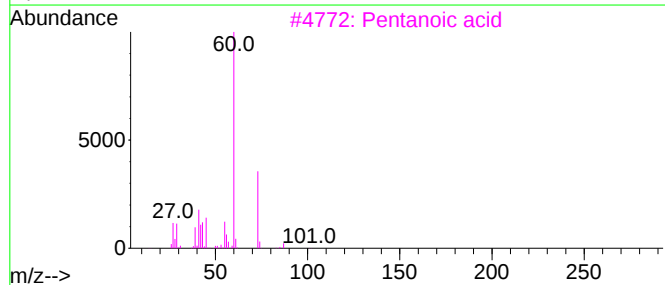
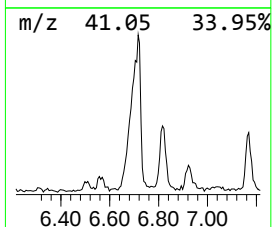
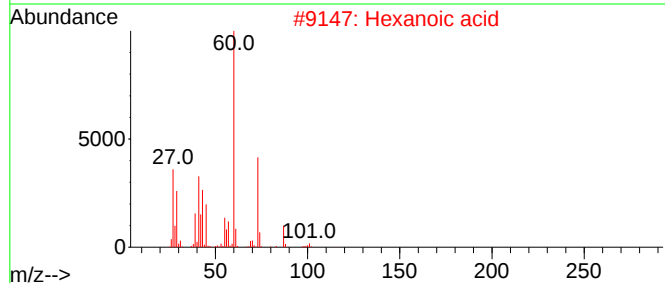
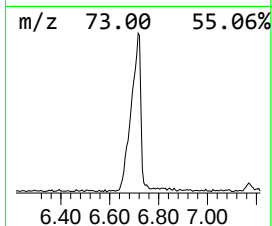
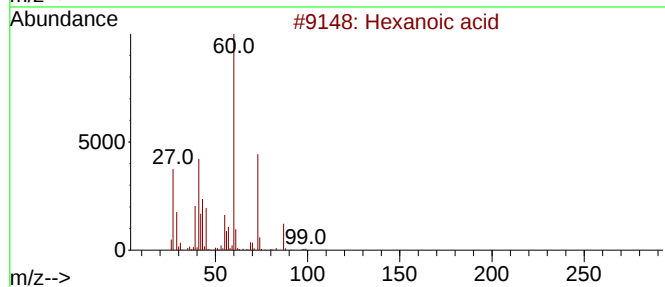
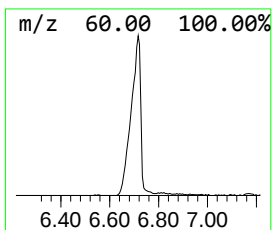
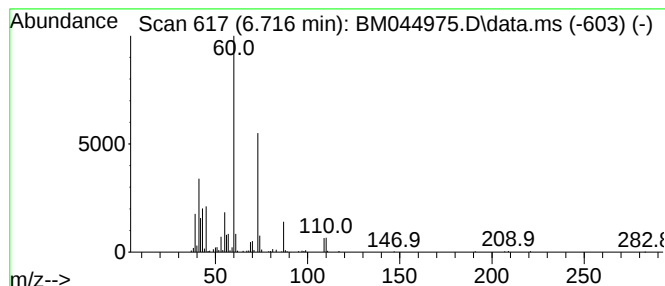
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Hexanoic acid Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.716	15.61 ng/ul	473801	1,4-Dichlorobenzene-d4	7.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid	116	C6H12O2	000142-62-1	90
2			Hexanoic acid	116	C6H12O2	000142-62-1	78
3			Pentanoic acid	102	C5H10O2	000109-52-4	72
4			Hexanoic acid	116	C6H12O2	000142-62-1	72
5			Hexanoic acid	116	C6H12O2	000142-62-1	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
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Sample : P2018-10
Misc :
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ClientSampleId :
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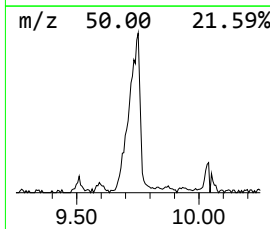
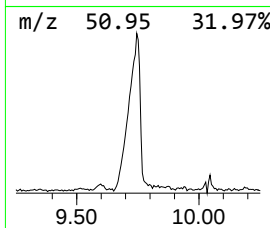
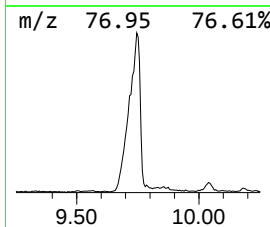
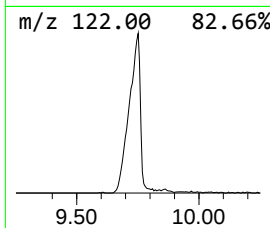
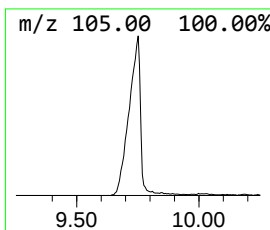
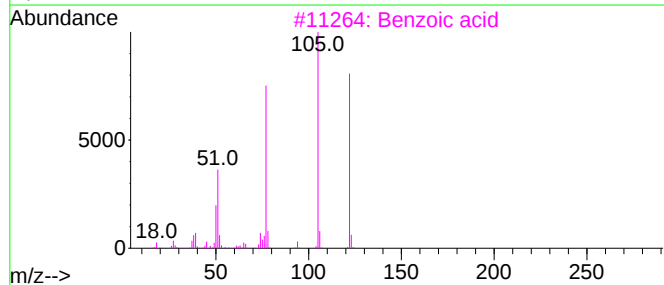
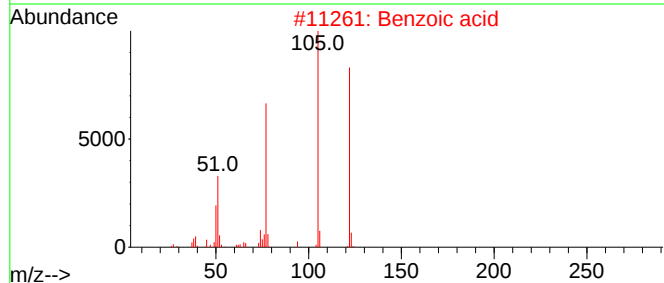
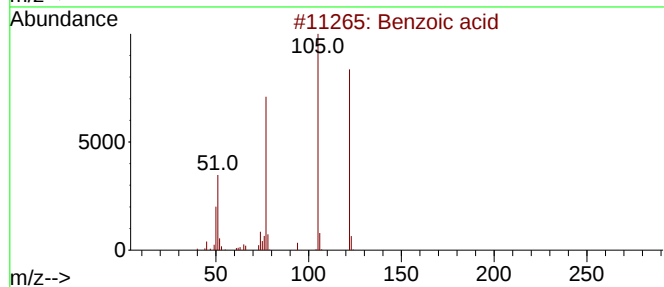
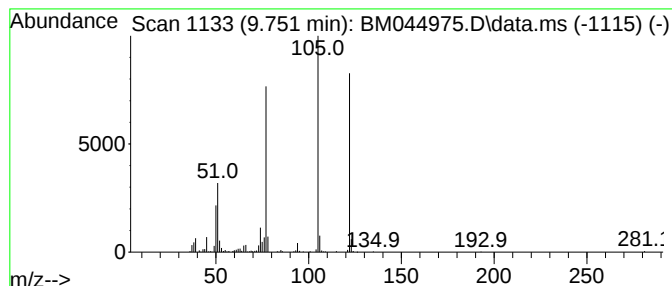
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.751	21.12 ng/ul	898932	Naphthalene-d8	10.410

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, silver(1+) salt	228	C7H5AgO2	000532-31-0	90
5			Heptanediamide, N,N'-di-benzoyloxy-	398	C21H22N2O6	1000253-26-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
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ClientSampleId :
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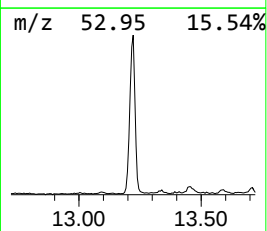
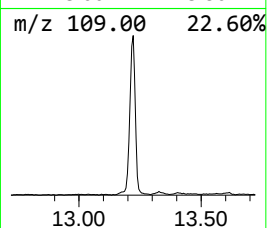
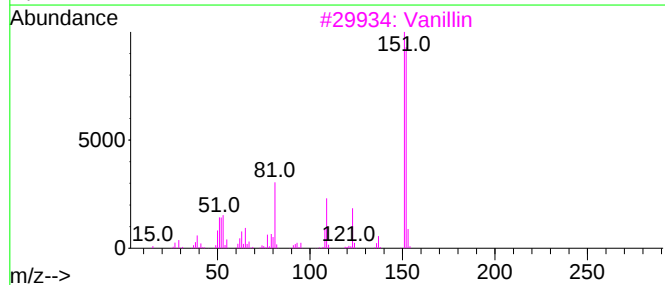
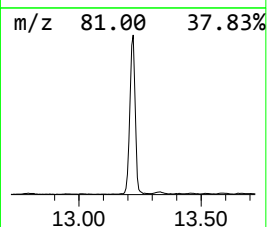
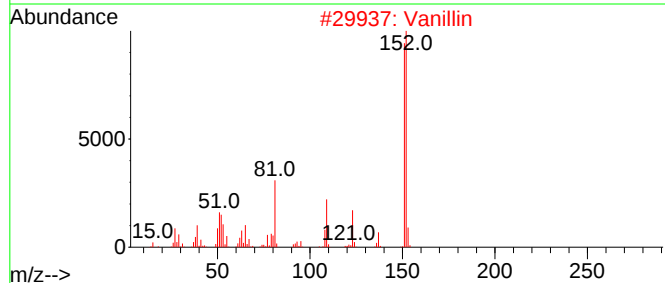
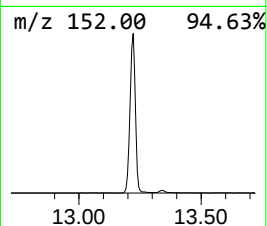
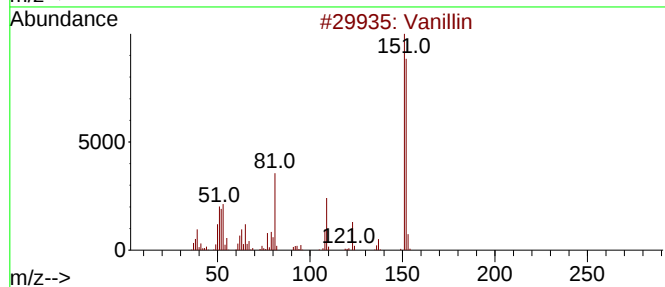
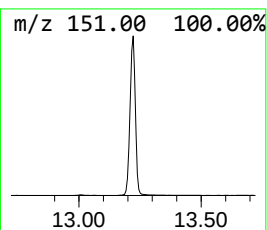
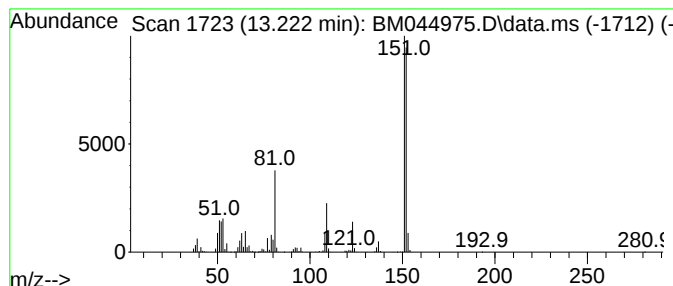
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Vanillin Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.222	41.82 ng/ul	2364090	Acenaphthene-d10	14.280

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Vanillin	152	C8H8O3	000121-33-5	98
2			Vanillin	152	C8H8O3	000121-33-5	97
3			Vanillin	152	C8H8O3	000121-33-5	97
4			Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	97
5			Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
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Sample : P2018-10
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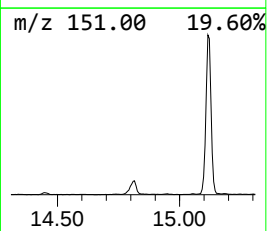
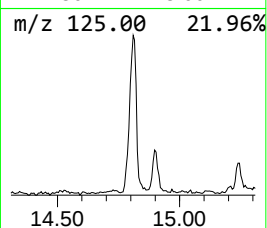
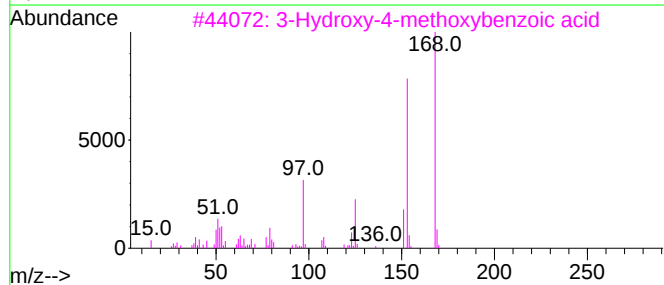
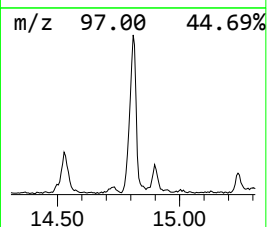
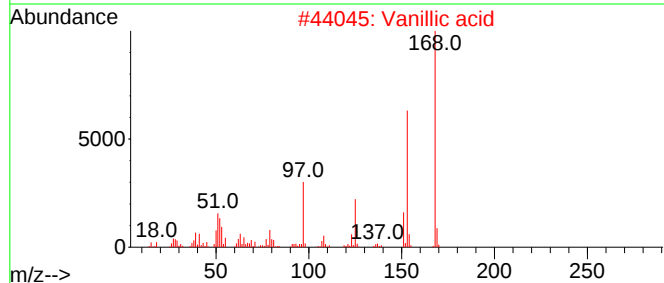
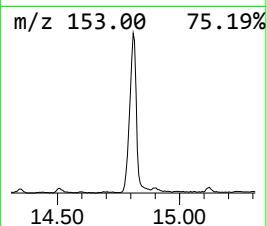
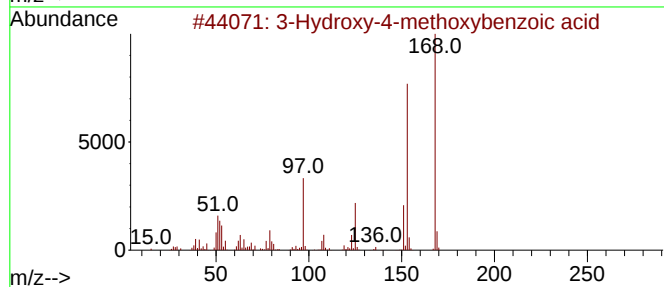
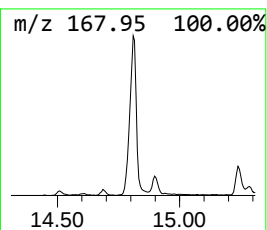
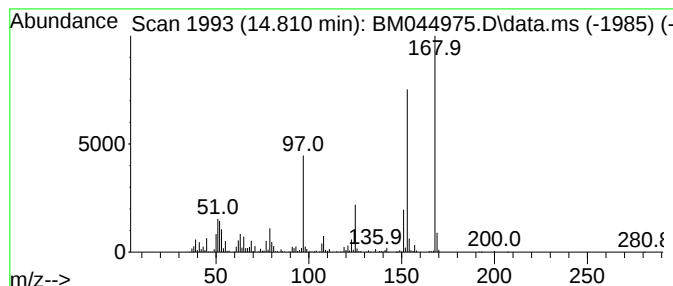
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 3-Hydroxy-4-methoxybenzoic ... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.810	15.64 ng/ul	884290	Acenaphthene-d10		14.280	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	3-Hydroxy-4-methoxybenzoic acid		168	C8H8O4	000645-08-9	96
2	Vanillic acid		168	C8H8O4	000121-34-6	96
3	3-Hydroxy-4-methoxybenzoic acid		168	C8H8O4	000645-08-9	95
4	3-Hydroxy-4-methoxybenzoic acid		168	C8H8O4	000645-08-9	93
5	1,2,4-Trimethoxybenzene		168	C9H12O3	000135-77-3	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
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Sample : P2018-10
Misc :
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Instrument :
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ClientSampleId :
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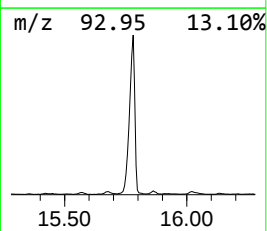
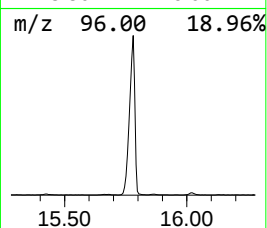
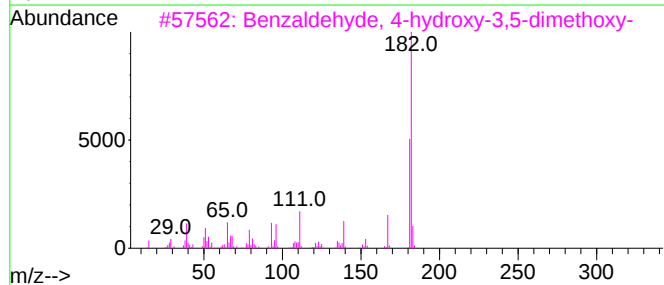
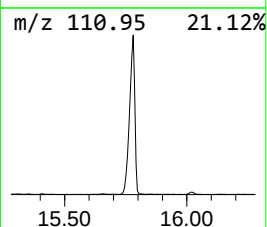
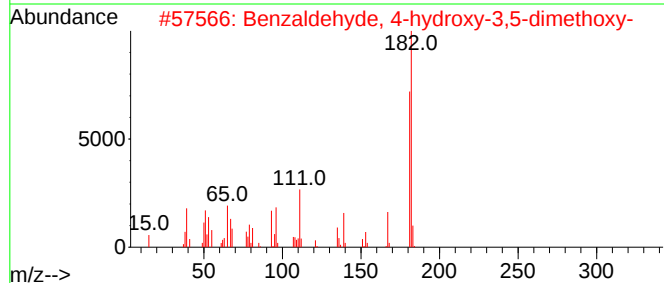
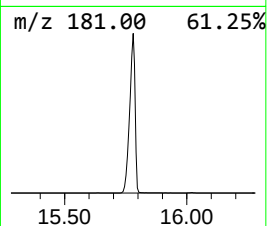
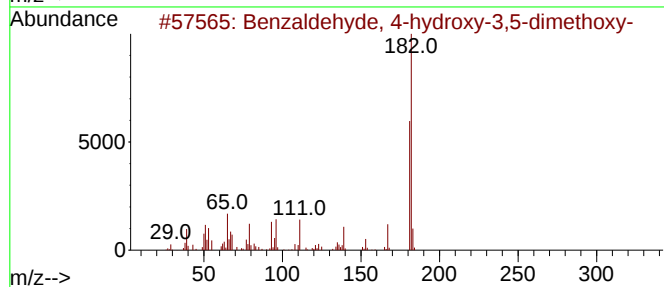
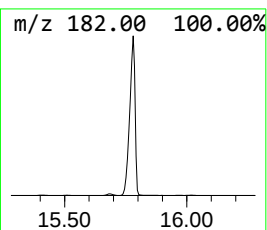
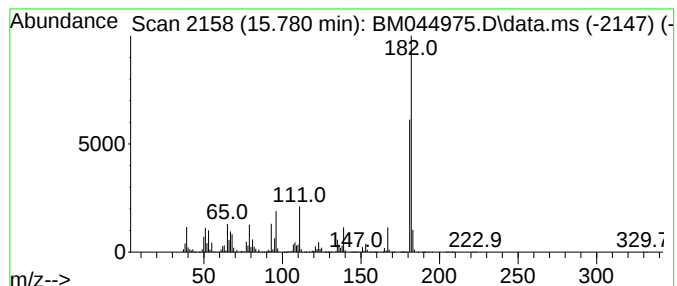
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Benzaldehyde, 4-hydroxy-3,5... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.780	123.40 ng/ul	13035000	Phenanthrene-d10	17.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	96
2			Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	94
3			Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	93
4			Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	93
5			Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
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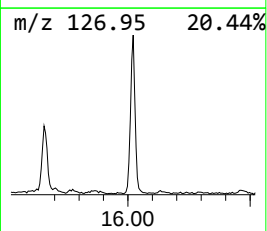
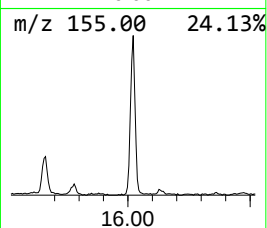
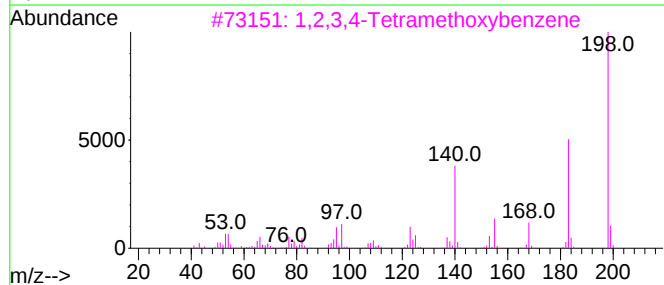
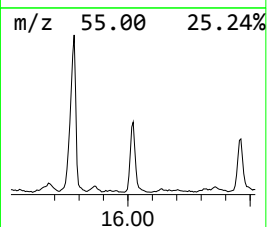
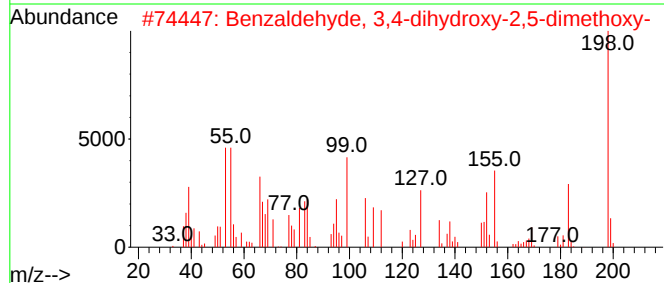
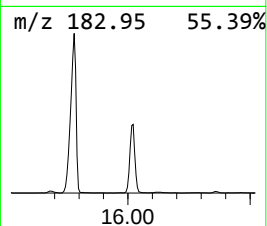
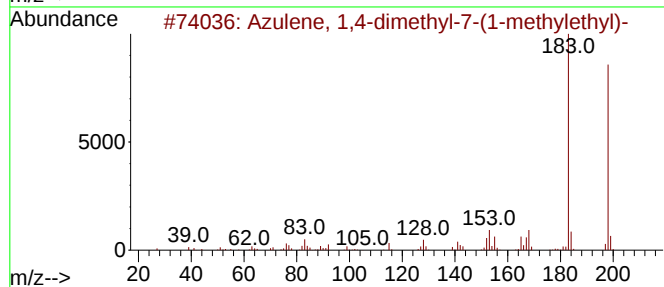
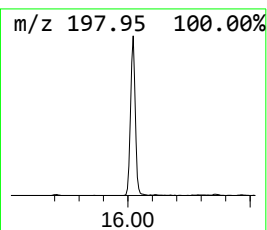
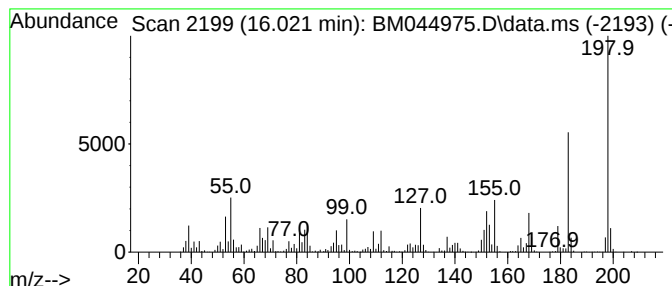
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Azulene, 1,4-dimethyl-7-(1-... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.021	14.49 ng/ul	1530090	Phenanthrene-d10	17.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Azulene, 1,4-dimethyl-7-(1-methy...	198	C15H18	000489-84-9	76
2			Benzaldehyde, 3,4-dihydroxy-2,5-...	198	C9H10O5	1000362-67-7	70
3			1,2,3,4-Tetramethoxybenzene	198	C10H14O4	021450-56-6	68
4			ethanone, 1-(5,8-dimethyl-1-naph...	198	C14H14O	1000396-02-9	68
5			.beta.-Carboline, 6-methoxy-	198	C12H10N2O	030684-42-5	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
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Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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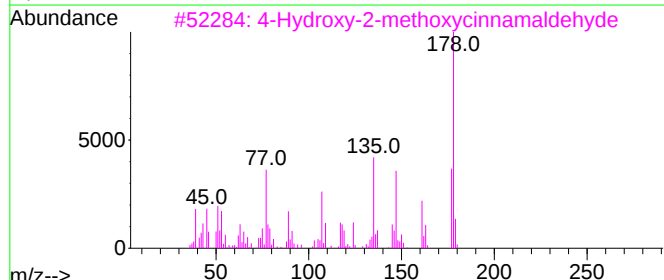
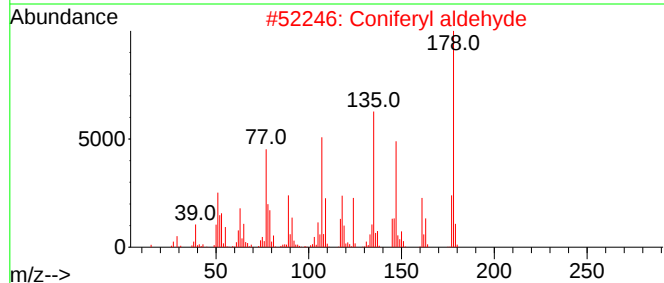
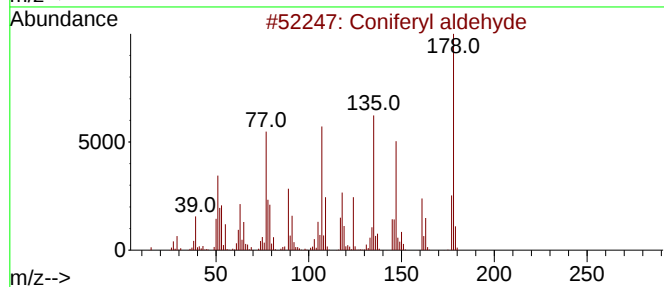
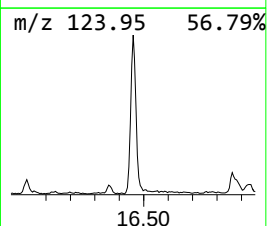
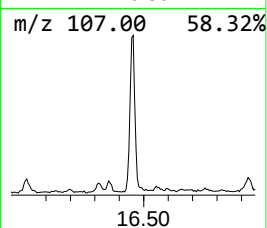
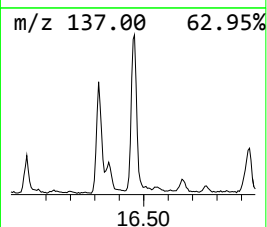
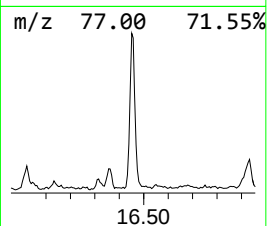
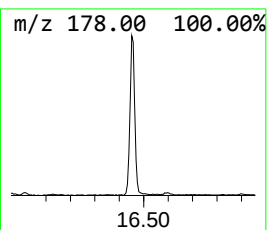
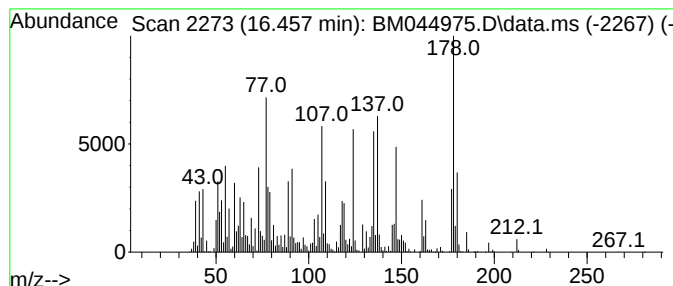
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Coniferyl aldehyde Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.457	16.86 ng/ul	1780880	Phenanthrene-d10	17.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Coniferyl aldehyde	178	C10H10O3	000458-36-6	99
2			Coniferyl aldehyde	178	C10H10O3	000458-36-6	98
3			4-Hydroxy-2-methoxycinnamaldehyde	178	C10H10O3	127321-19-1	43
4			1-(2-Ethoxyphenyl)acetone	178	C11H14O2	086432-38-4	42
5			1-(2-Ethoxyphenyl)acetone	178	C11H14O2	086432-38-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

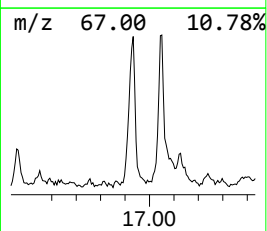
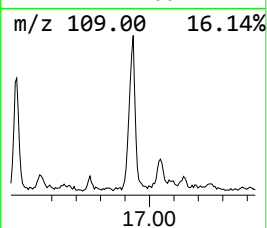
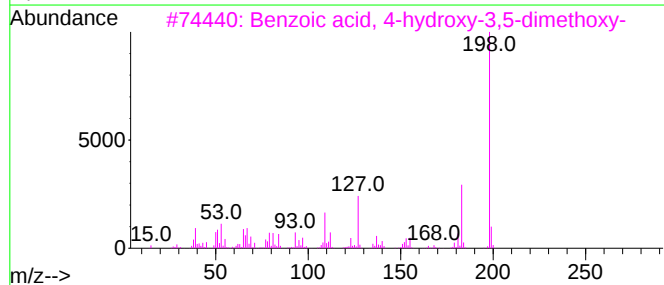
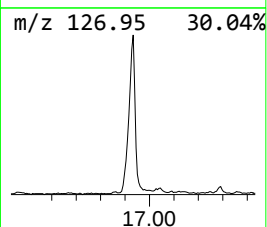
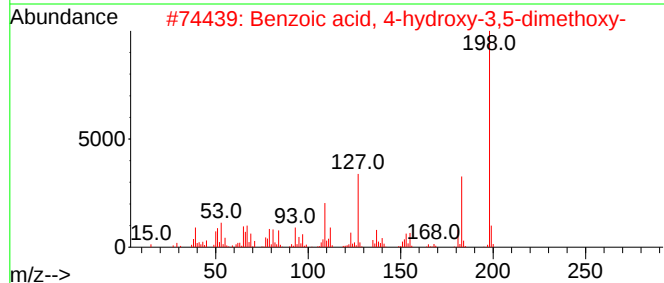
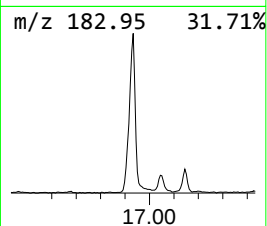
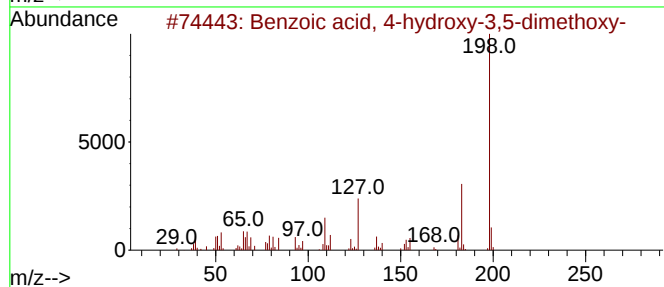
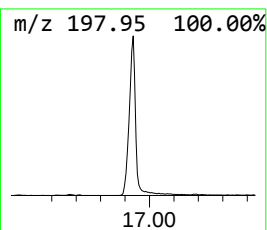
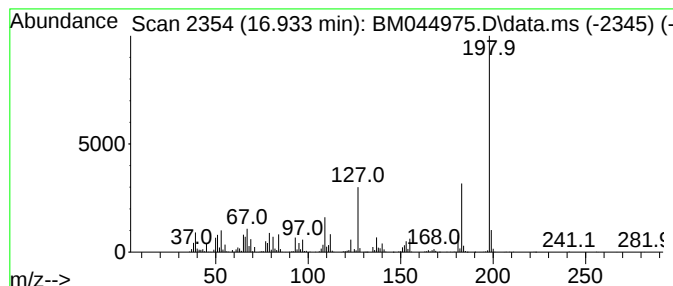
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Benzoic acid, 4-hydroxy-3,5... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.933	14.69 ng/ul	1552240	Phenanthrene-d10	17.045	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 4-hydroxy-3,5-dime...	198	C9H10O5	000530-57-4	98
2	Benzoic acid, 4-hydroxy-3,5-dime...	198	C9H10O5	000530-57-4	98
3	Benzoic acid, 4-hydroxy-3,5-dime...	198	C9H10O5	000530-57-4	98
4	Benzoic acid, 4-hydroxy-3,5-dime...	198	C9H10O5	000530-57-4	93
5	Benzoic acid, 4-hydroxy-3,5-dime...	198	C9H10O5	000530-57-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
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Instrument :
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ClientSampleId :
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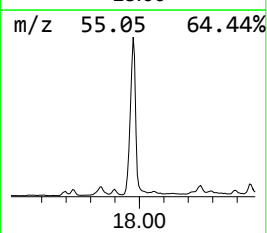
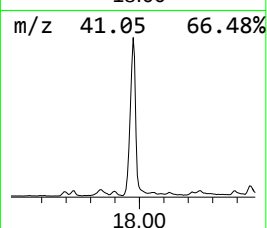
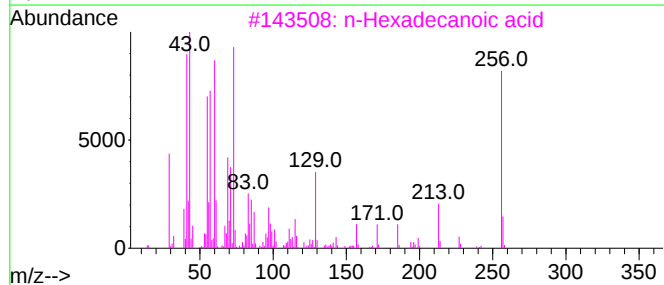
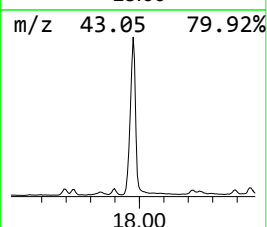
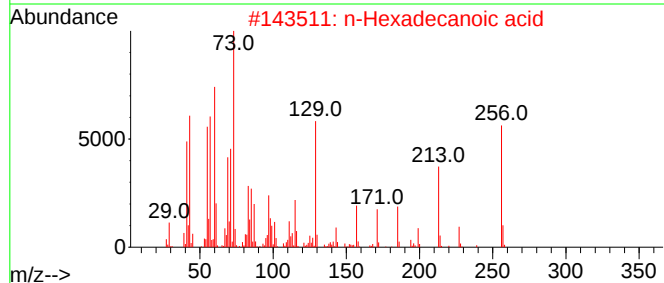
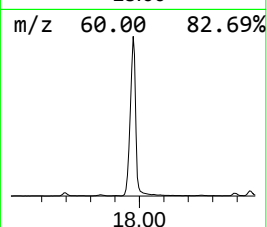
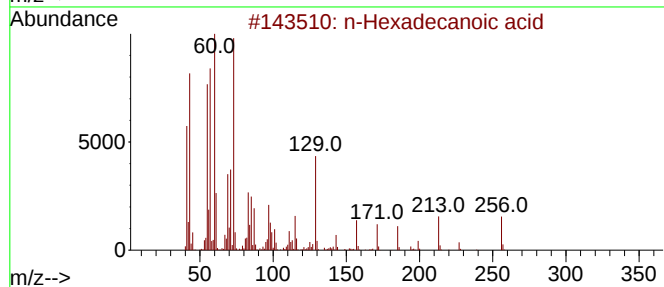
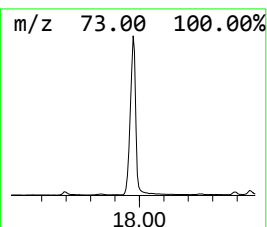
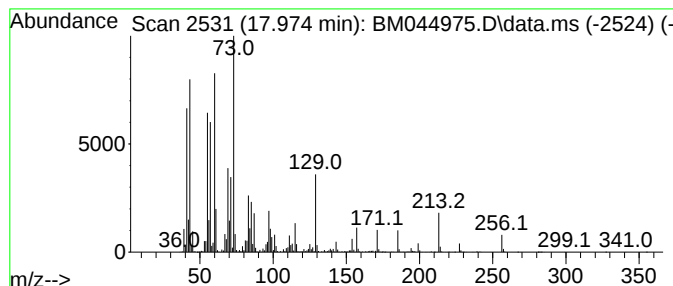
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.974	76.55 ng/ul	8086650	Phenanthrene-d10	17.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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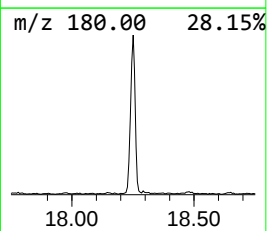
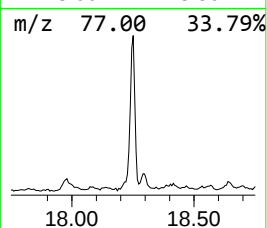
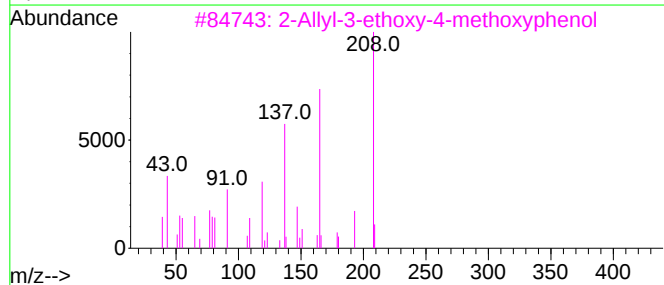
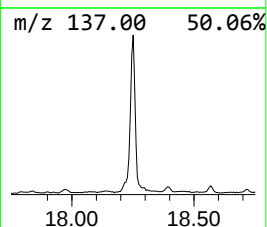
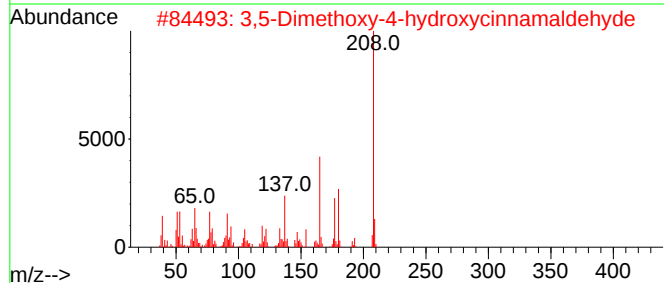
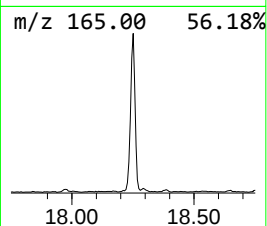
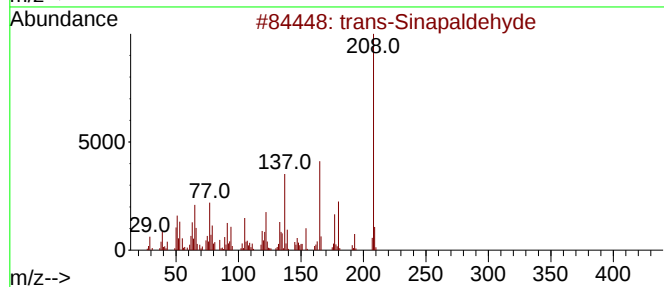
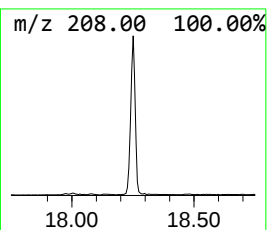
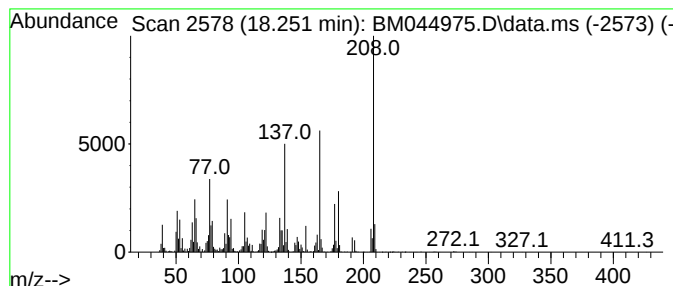
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 trans-Sinapaldehyde Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.251	25.80 ng/ul	2725780	Phenanthrene-d10	17.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-Sinapaldehyde	208	C11H12O4	004206-58-0	93
2			3,5-Dimethoxy-4-hydroxycinnamal...	208	C11H12O4	087345-53-7	86
3			2-Allyl-3-ethoxy-4-methoxyphenol	208	C12H16O3	1000122-70-6	58
4			Benzene, 1,2,4-trimethoxy-5-(1-p...	208	C12H16O3	000494-40-6	53
5			.beta.-Asarone	208	C12H16O3	005273-86-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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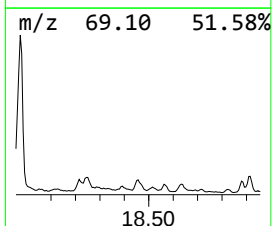
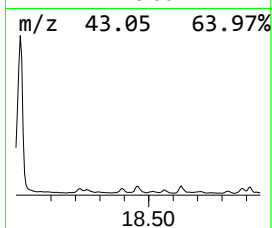
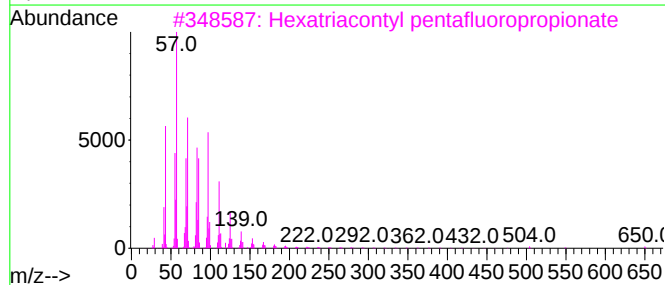
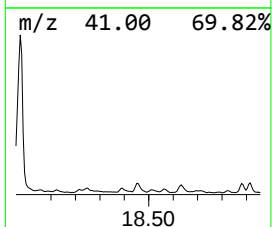
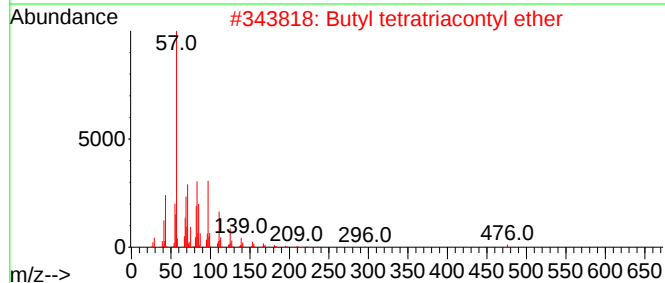
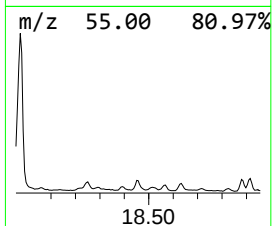
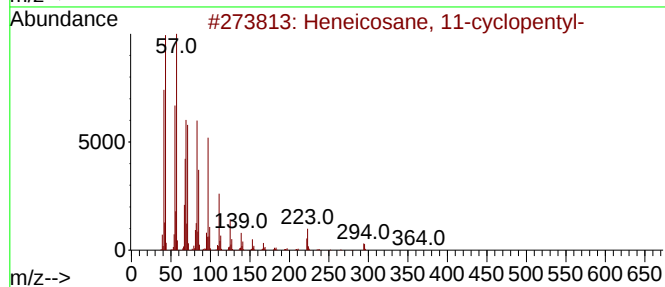
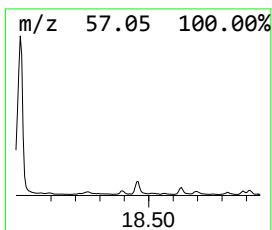
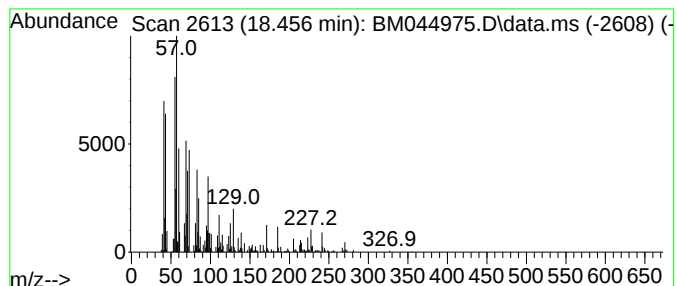
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Quant Title : SVOA CALIBRATION

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

Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.456	4.78 ng/ul	505119	Phenanthrene-d10	17.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane, 11-cyclopentyl-	364	C26H52	006703-81-7	42
2			Butyl tetratriacontyl ether	551	C38H78O	1000406-41-4	38
3			Hexatriacontyl pentafluoropropio...	669	C39H73F5O2	1000351-89-0	38
4			Butyl docosyl ether	382	C26H54O	1000406-40-8	30
5			Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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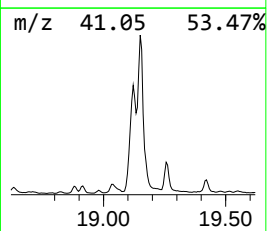
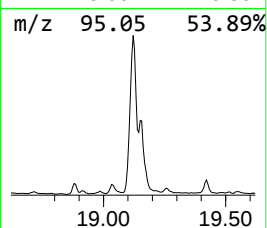
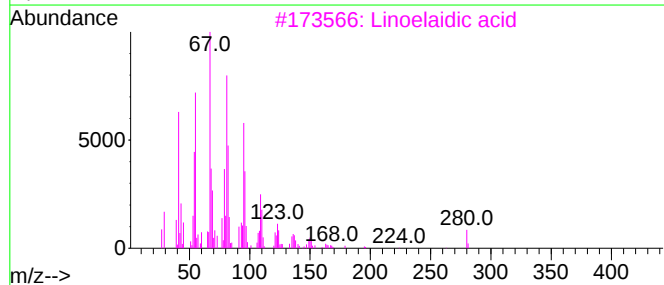
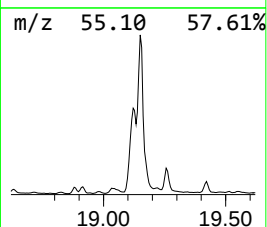
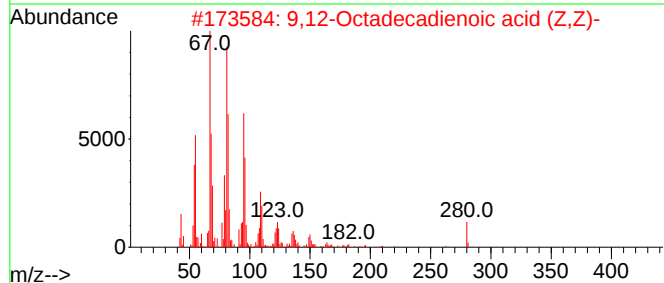
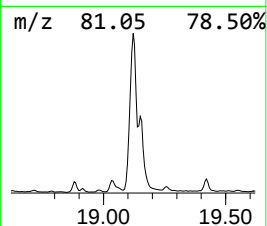
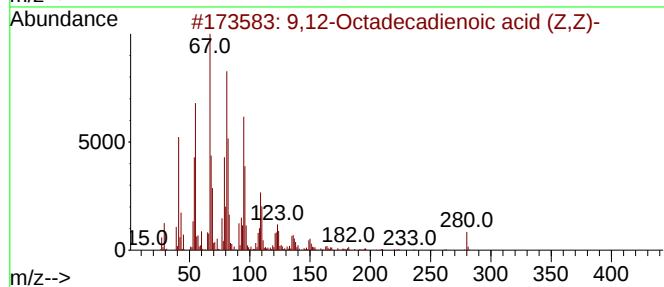
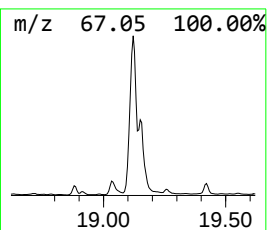
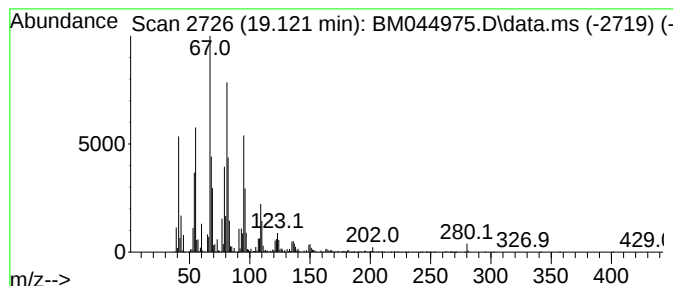
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 9,12-Octadecadienoic acid (... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.121	95.71 ng/ul	10110200	Phenanthrene-d10	17.045

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	99
2		9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	99
3		Linoleic acid	280	C18H32O2	000506-21-8	98
4		10E,12Z-Octadecadienoic acid	280	C18H32O2	002420-56-6	98
5		(9E,11E)-Octadecadienoic acid	280	C18H32O2	000544-71-8	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

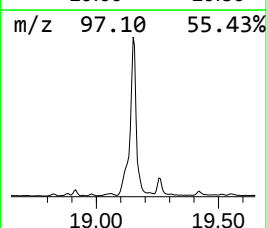
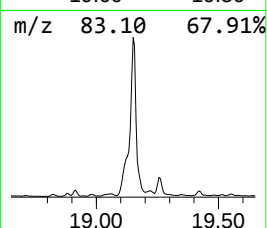
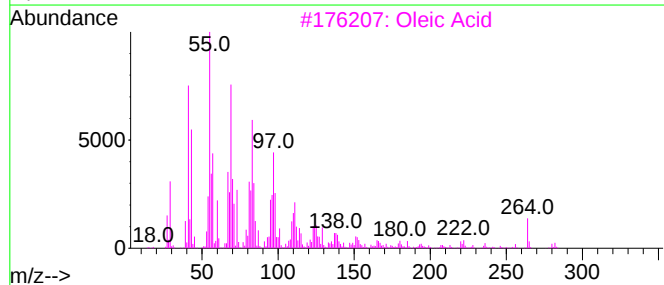
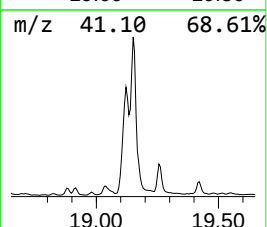
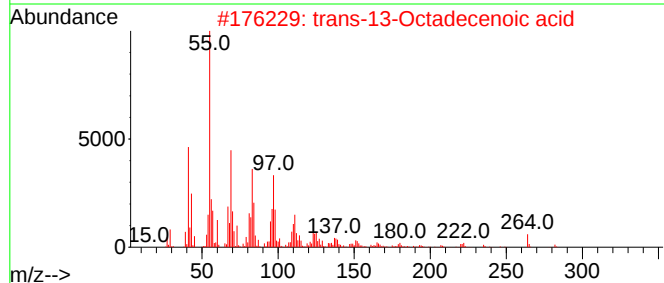
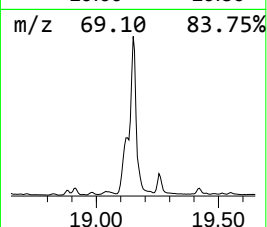
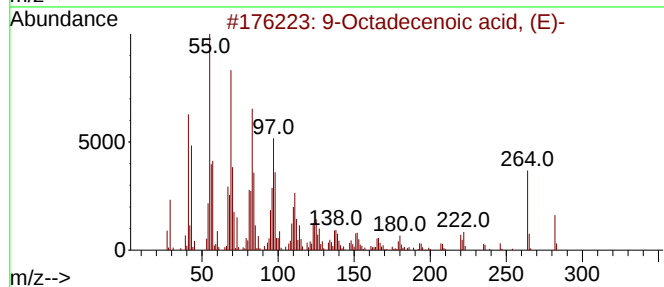
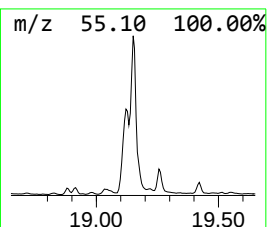
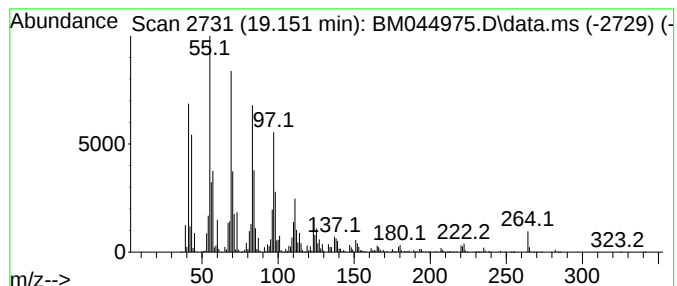
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 9-Octadecenoic acid, (E)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.151	108.60 ng/ul	10298700	Chrysene-d12	21.250	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	99
2	trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	87
3	Oleic Acid	282	C18H34O2	000112-80-1	87
4	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	78
5	cis-Vaccenic acid	282	C18H34O2	000506-17-2	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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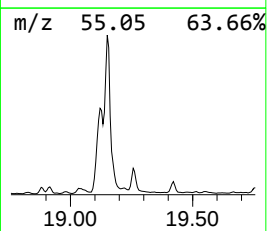
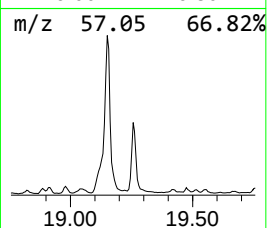
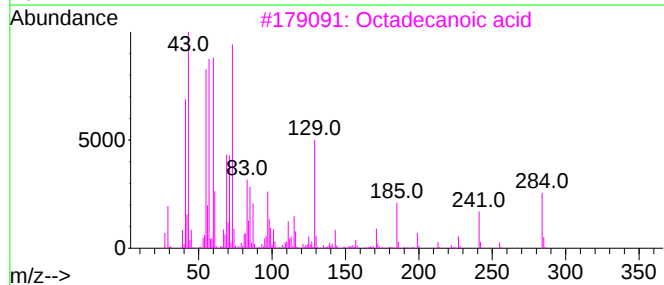
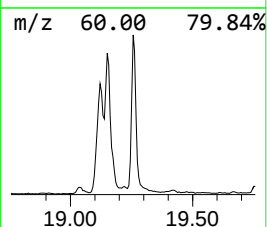
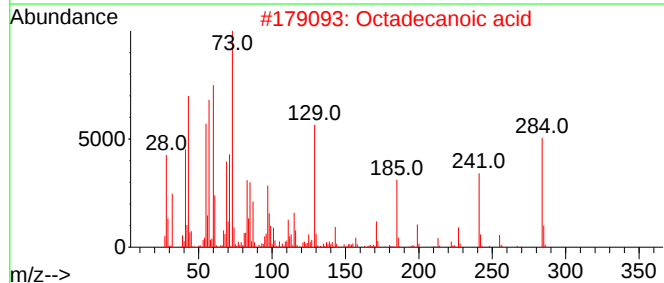
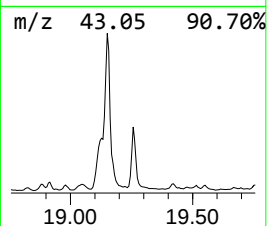
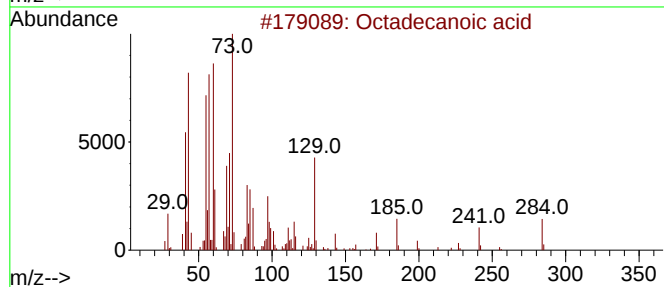
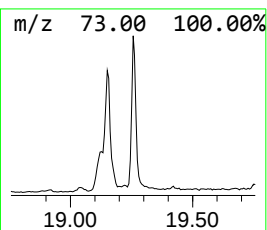
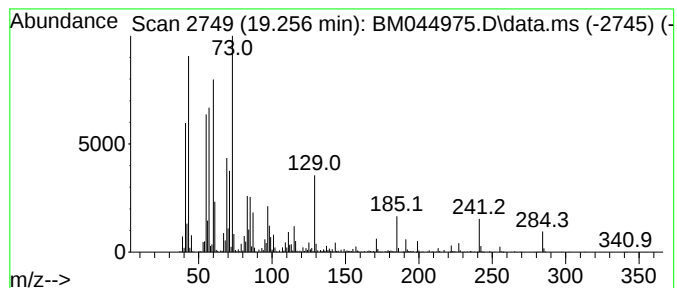
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Quant Title : SVOA CALIBRATION

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

Peak Number 40 Octadecanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.256	20.52 ng/ul	1945870	Chrysene-d12	21.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	96
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

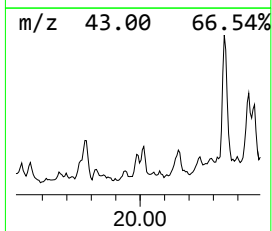
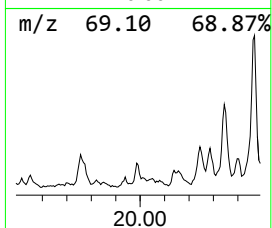
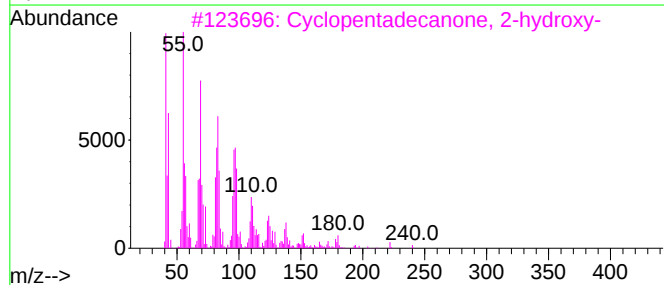
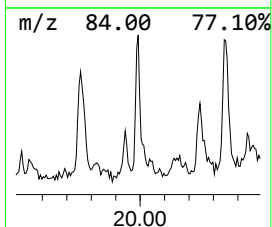
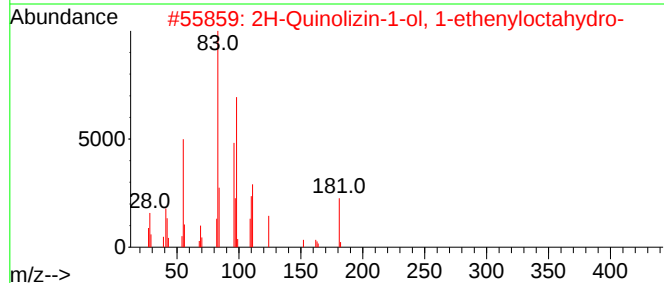
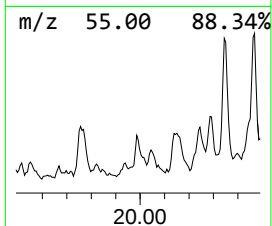
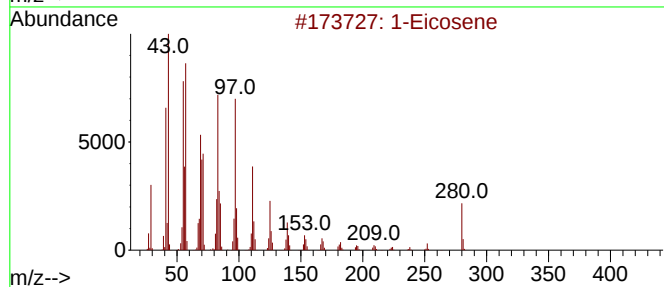
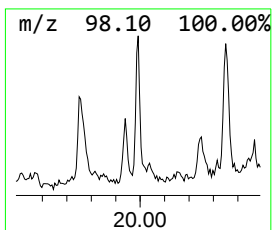
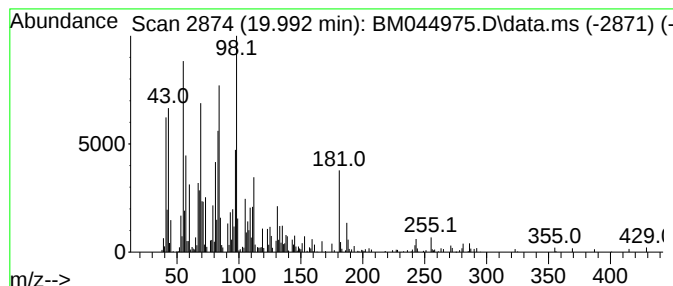
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 65

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.992	3.08 ng/ul	292561	Chrysene-d12	21.250	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Eicosene	280	C20H40	003452-07-1	58
2	2H-Quinolizin-1-ol, 1-ethenyloct...	181	C11H19NO	069597-57-5	46
3	Cyclopentadecanone, 2-hydroxy-	240	C15H28O2	004727-18-8	41
4	5-Eicosene, (E)-	280	C20H40	074685-30-6	38
5	Cyclohexanone, 2-ethyl-	126	C8H14O	004423-94-3	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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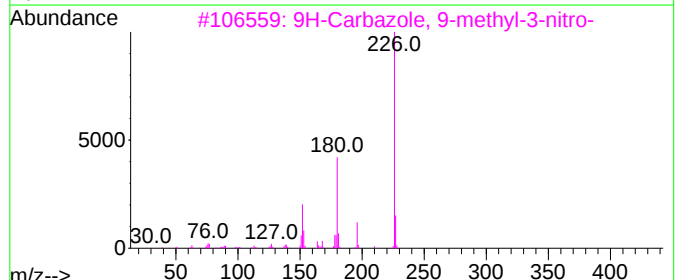
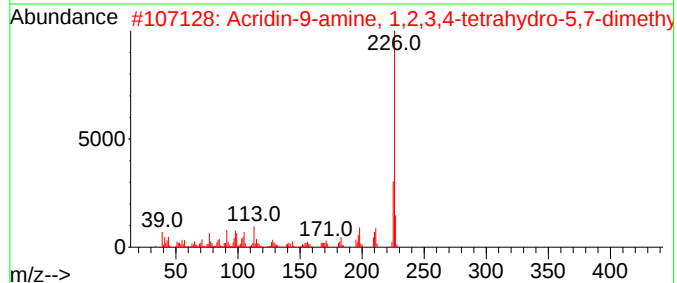
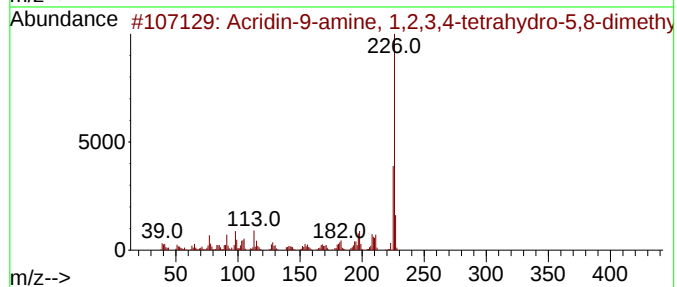
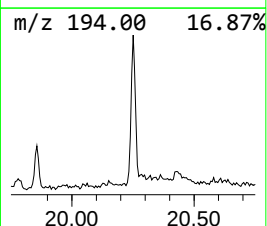
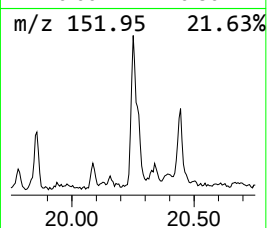
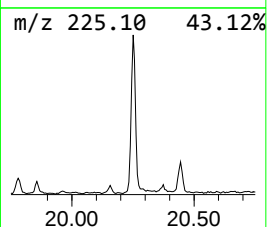
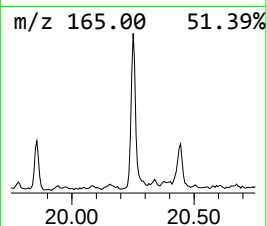
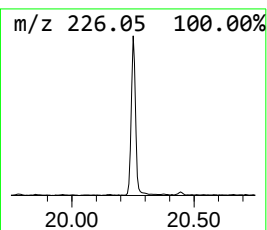
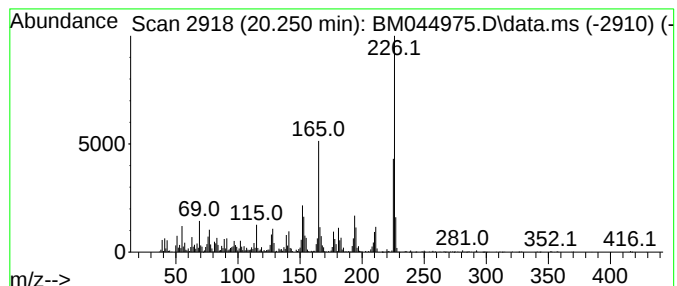
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Quant Title : SVOA CALIBRATION

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

Peak Number 45 Acridin-9-amine, 1,2,3,4-te... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.250	14.94 ng/ul	1416650	Chrysene-d12	21.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acridin-9-amine, 1,2,3,4-tetrahy...	226	C15H18N2	297758-19-1	64
2			Acridin-9-amine, 1,2,3,4-tetrahy...	226	C15H18N2	1000300-57-6	60
3			9H-Carbazole, 9-methyl-3-nitro-	226	C13H10N2O2	061166-05-0	50
4			Anthralin	226	C14H10O3	001143-38-0	50
5			9H-Pyrido[3,4-b]indole, 7-methox...	226	C14H14N2O	143502-37-8	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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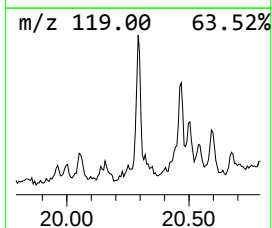
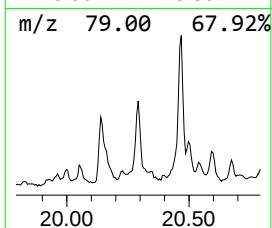
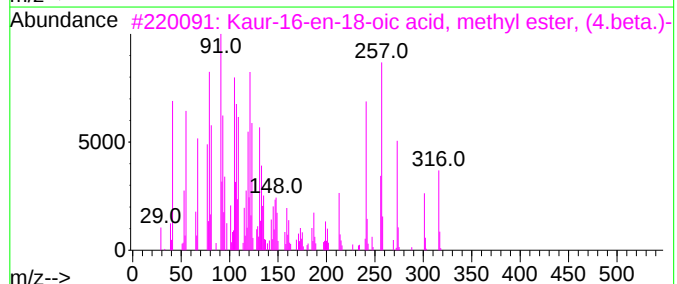
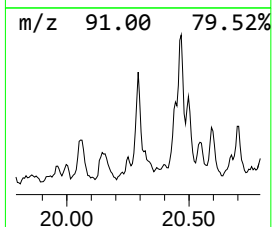
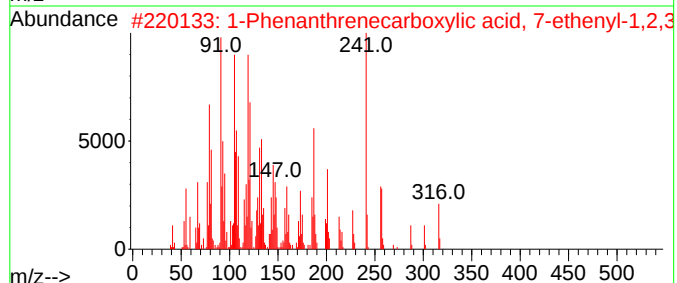
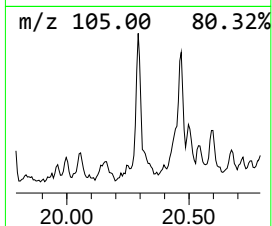
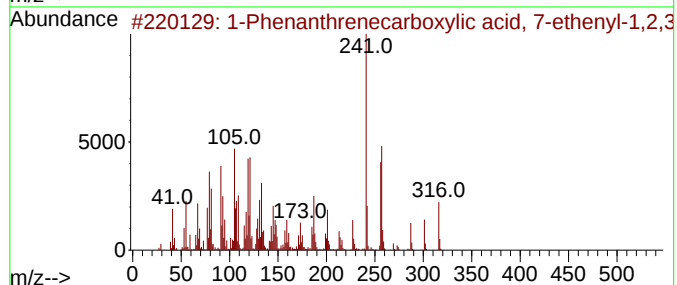
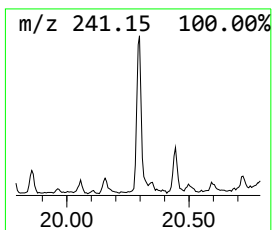
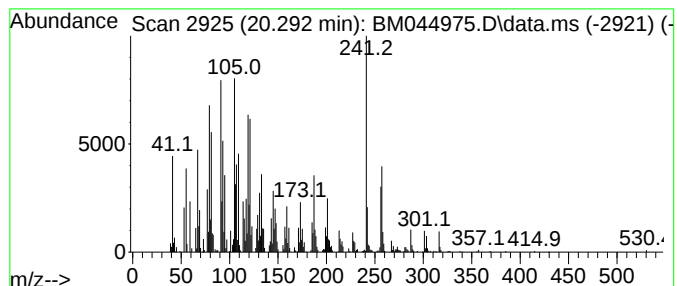
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Quant Title : SVOA CALIBRATION

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

Peak Number 46 1-Phenanthrenecarboxylic ac... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.292	15.06 ng/ul	1428570	Chrysene-d12	21.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenanthrenecarboxylic acid, 7...	316	C21H32O2	001686-62-0	91
2			1-Phenanthrenecarboxylic acid, 7...	316	C21H32O2	001686-62-0	89
3			Kaur-16-en-18-oic acid, methyl e...	316	C21H32O2	005524-25-4	70
4			1-Phenanthrenecarboxylic acid, 7...	316	C21H32O2	001686-62-0	53
5			7-Ethenyl-1,4a,7-trimethyl-3,4,6...	316	C21H32O2	1000504-06-6	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
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Sample : P2018-10
Misc :
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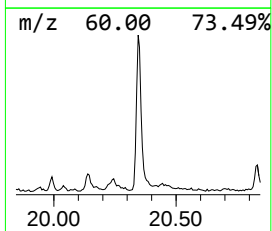
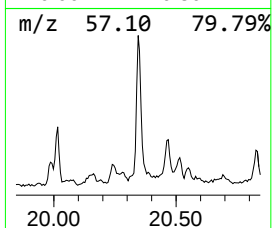
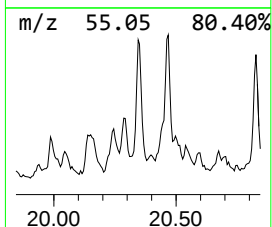
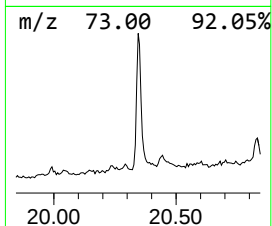
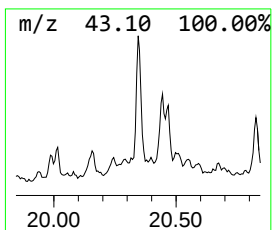
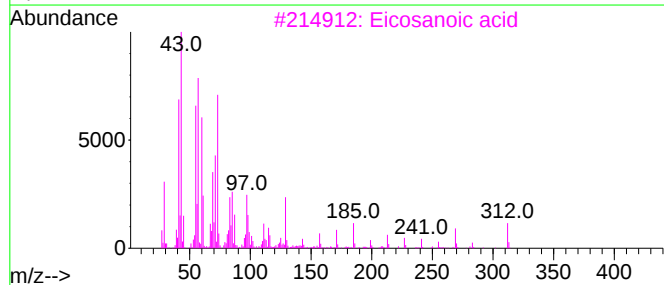
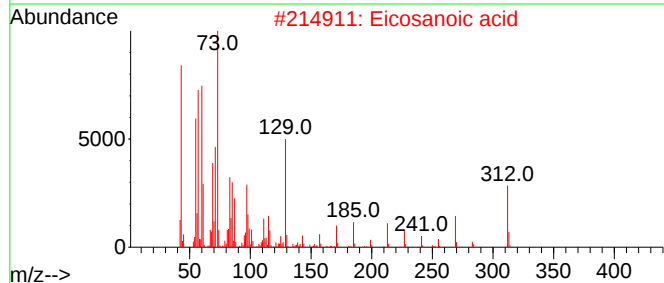
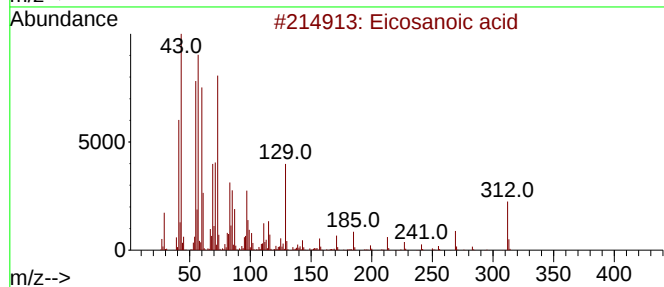
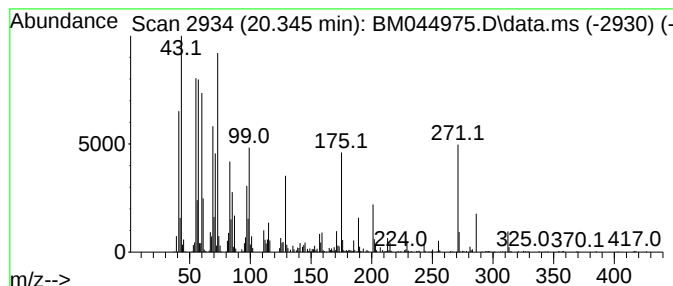
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Quant Title : SVOA CALIBRATION

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

Peak Number 47 Eicosanoic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.345	18.47 ng/ul	1751650	Chrysene-d12	21.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosanoic acid	312	C20H40O2	000506-30-9	95
2			Eicosanoic acid	312	C20H40O2	000506-30-9	84
3			Eicosanoic acid	312	C20H40O2	000506-30-9	83
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	78
5			n-Decanoic acid	172	C10H20O2	000334-48-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
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Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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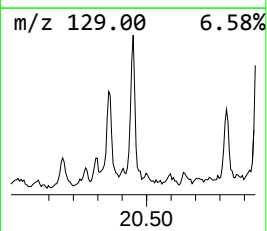
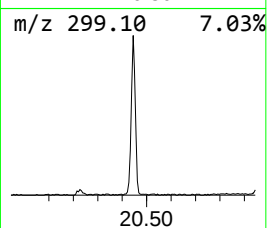
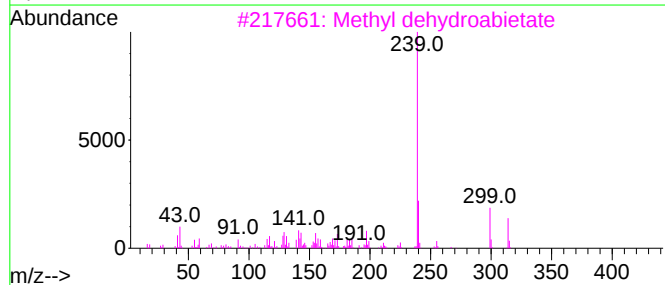
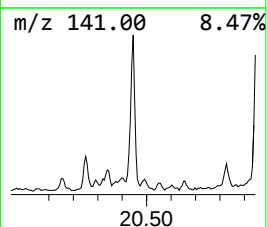
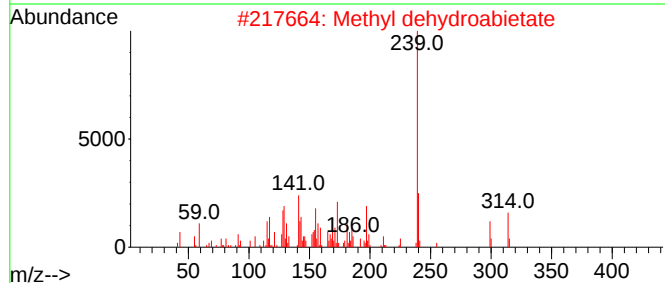
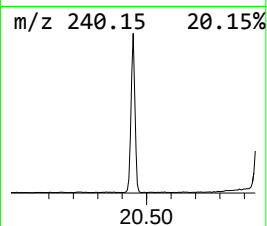
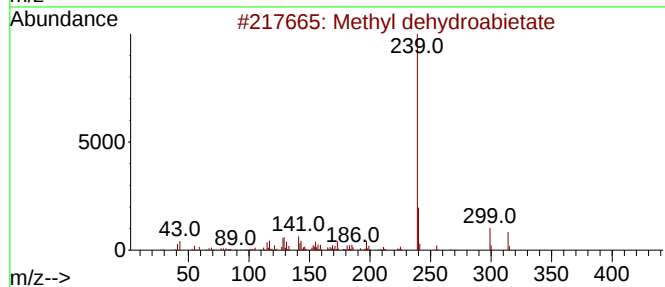
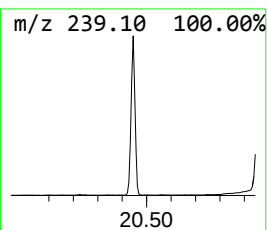
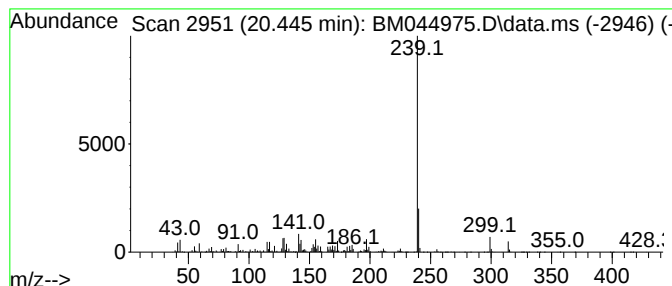
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Quant Title : SVOA CALIBRATION

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

Peak Number 49 Methyl dehydroabietate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.445	30.72 ng/ul	2913690	Chrysene-d12	21.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl dehydroabietate	314	C21H30O2	001235-74-1	95
2			Methyl dehydroabietate	314	C21H30O2	001235-74-1	92
3			Methyl dehydroabietate	314	C21H30O2	001235-74-1	91
4			Methyl dehydroabietate	314	C21H30O2	001235-74-1	90
5			9,10-Anthracenedione, 1-amino-4-...	239	C14H9NO3	000116-85-8	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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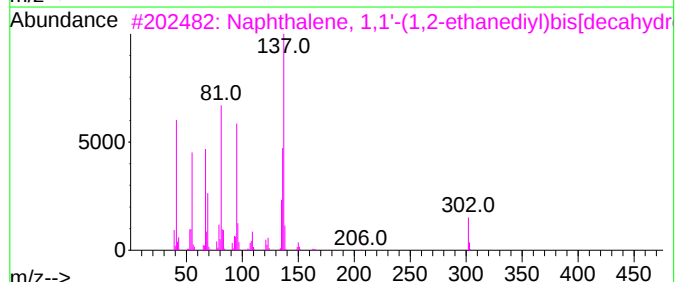
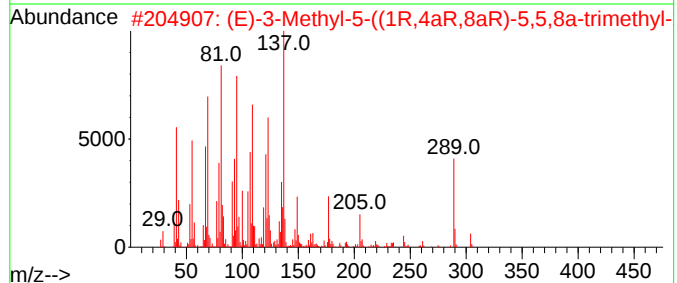
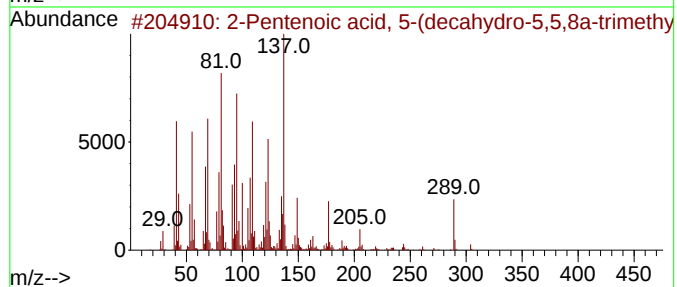
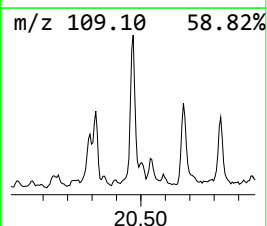
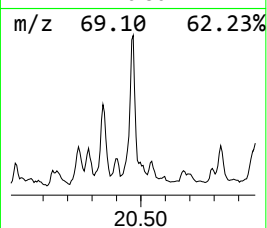
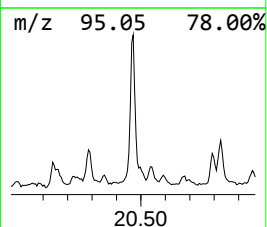
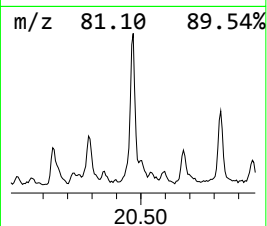
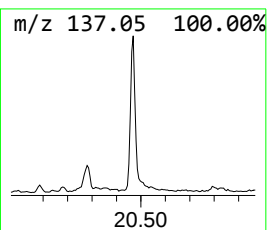
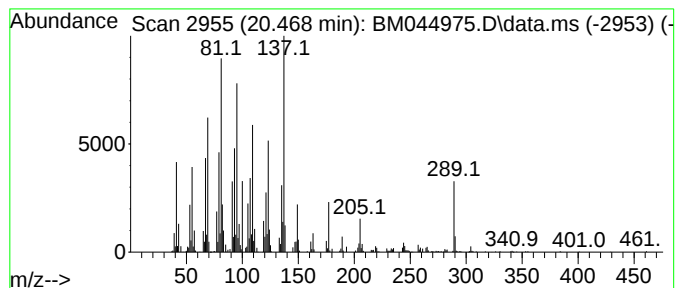
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Quant Title : SVOA CALIBRATION

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

Peak Number 50 2-Pentenoic acid, 5-(decahy... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.468	20.02 ng/ul	1898300	Chrysene-d12	21.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentenoic acid, 5-(decahydro-5...	304	C20H32O2	024470-48-2	98
2			(E)-3-Methyl-5-((1R,4aR,8aR)-5,5...	304	C20H32O2	020257-75-4	94
3			Naphthalene, 1,1'-(1,2-ethanediyl...	302	C22H38	054934-69-9	64
4			3-Iodomethyl-3,6,6-trimethyl-cyc...	264	C10H17I	1000193-08-2	50
5			Bicyclo[6.1.0]nonane, 9-bromo-9...	216	C10H17Br	059474-03-2	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
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Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

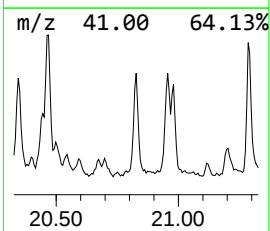
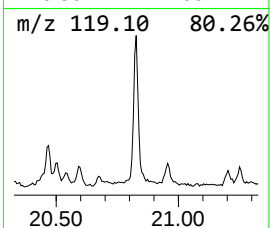
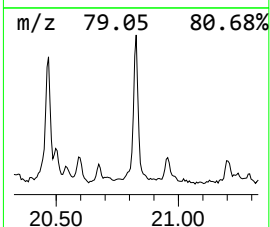
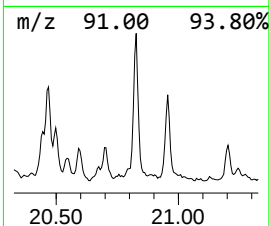
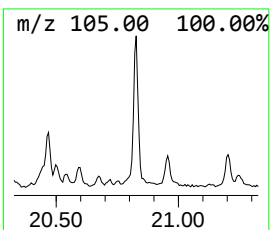
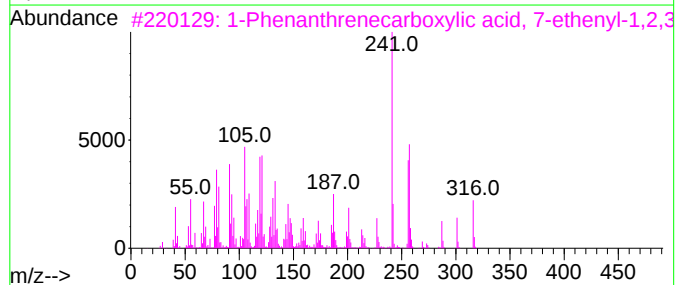
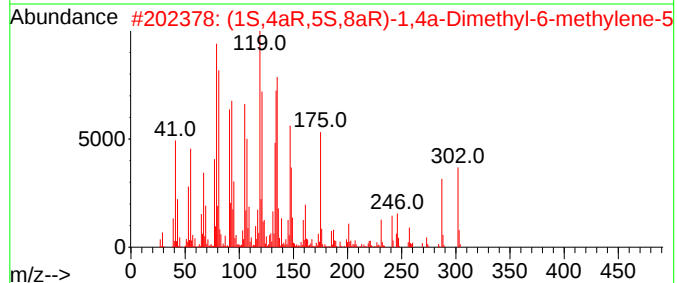
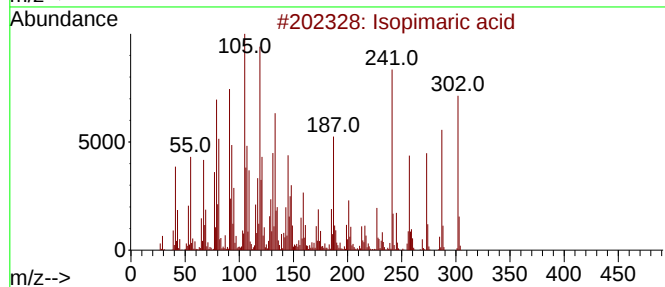
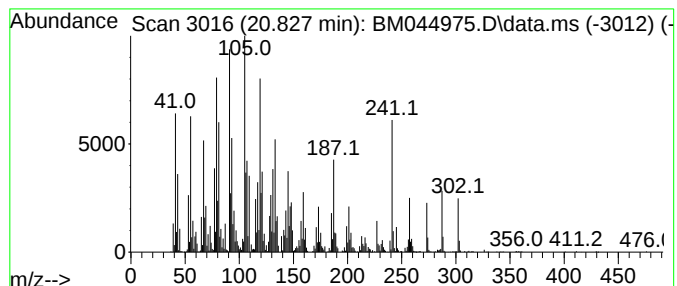
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 55 Isopimaric acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.827	24.35 ng/ul	2308890	Chrysene-d12		21.250	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Isopimaric acid		302	C20H30O2	005835-26-7	99
2	(1S,4aR,5S,8aR)-1,4a-Dimethyl-6-...		302	C20H30O2	002761-77-5	64
3	1-Phenanthrenecarboxylic acid, 7...		316	C21H32O2	001686-62-0	53
4	1-Phenanthrenecarboxylic acid, 7...		316	C21H32O2	001686-62-0	47
5	9,12-Octadecadiynoic acid, methy...		290	C19H30O2	1000333-56-4	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
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Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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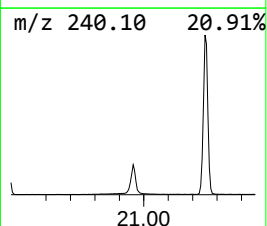
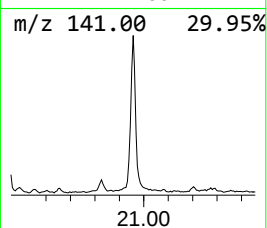
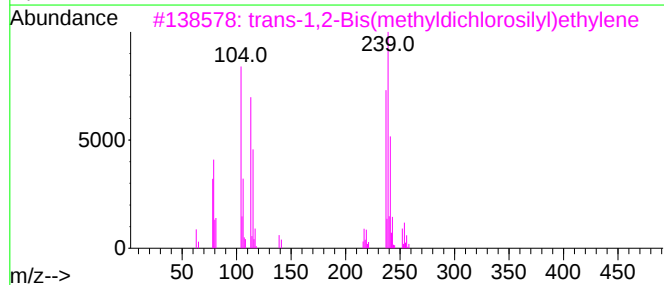
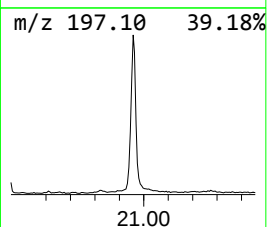
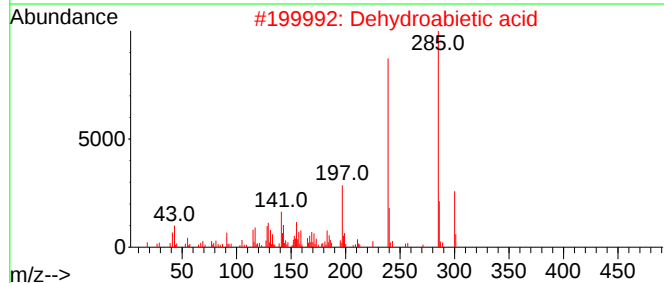
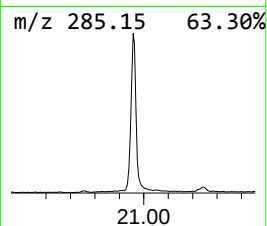
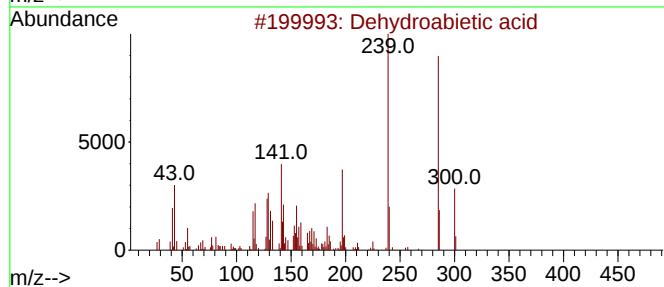
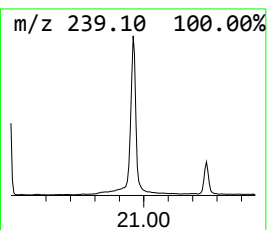
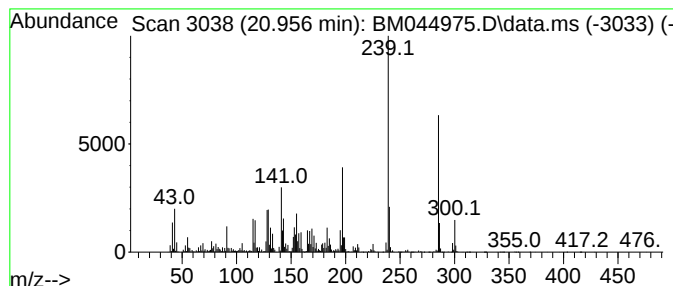
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Quant Title : SVOA CALIBRATION

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

Peak Number 56 Dehydroabiatic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.956	46.51 ng/ul	4410480	Chrysene-d12	21.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dehydroabiatic acid	300	C20H28O2	001740-19-8	99
2			Dehydroabiatic acid	300	C20H28O2	001740-19-8	98
3			trans-1,2-Bis(methyldichlorosilyl)...	252	C4H8Cl4Si2	065899-10-7	92
4			Dehydroabiatic acid	300	C20H28O2	001740-19-8	87
5			1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

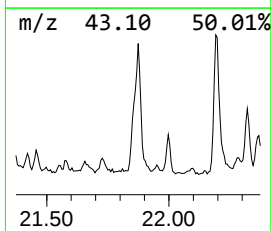
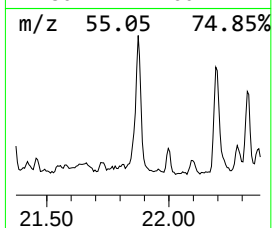
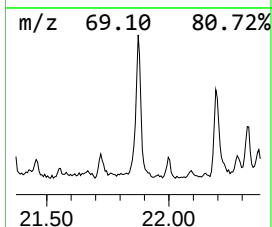
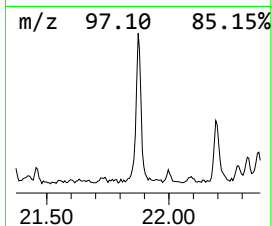
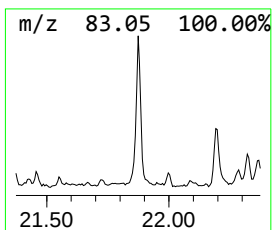
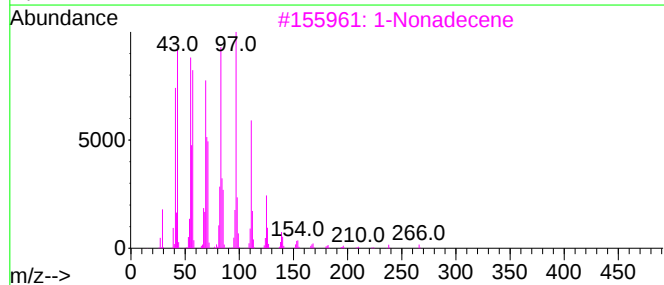
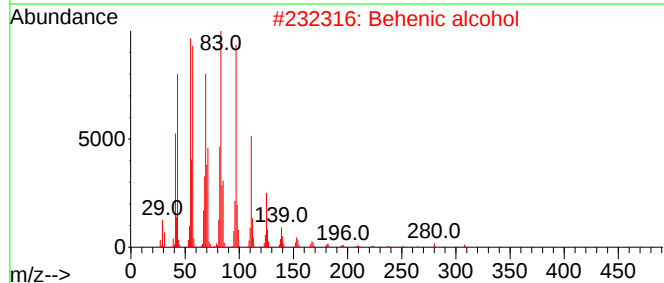
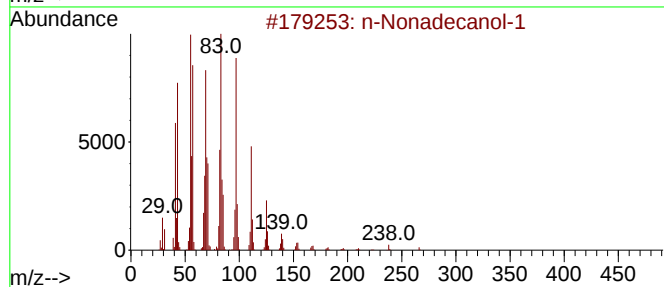
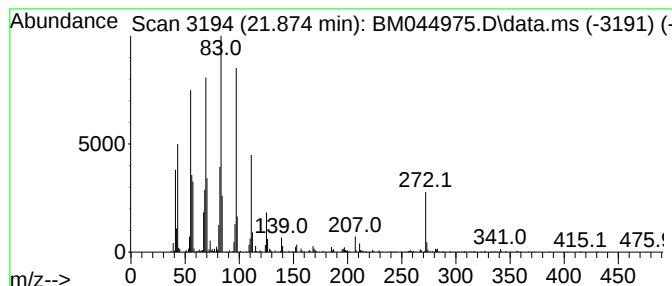
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 61 n-Nonadecanol-1 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.874	15.22 ng/ul	1443320	Chrysene-d12	21.250	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Nonadecanol-1	284	C19H40O	001454-84-8	58
2	Behenic alcohol	326	C22H46O	000661-19-8	49
3	1-Nonadecene	266	C19H38	018435-45-5	46
4	1-Octadecanol	270	C18H38O	000112-92-5	45
5	1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
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Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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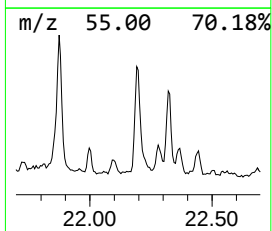
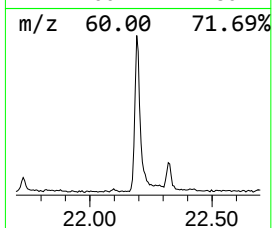
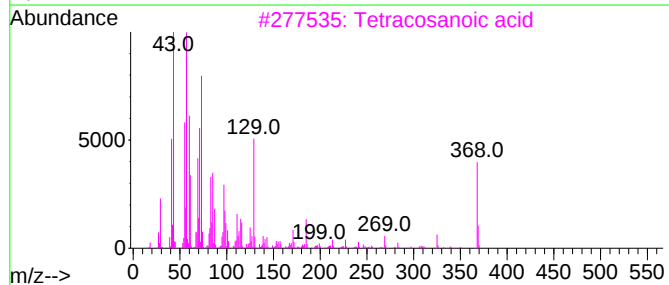
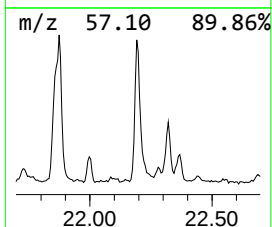
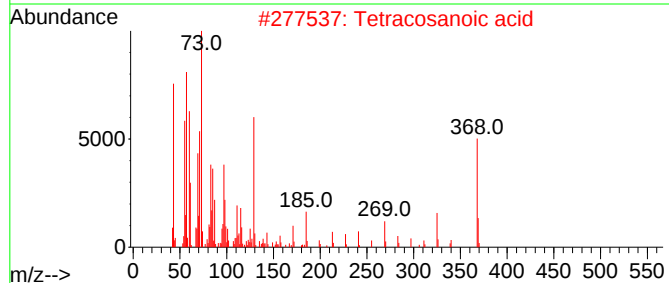
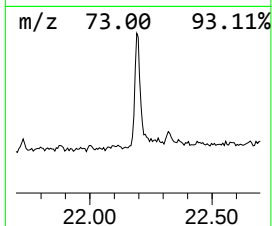
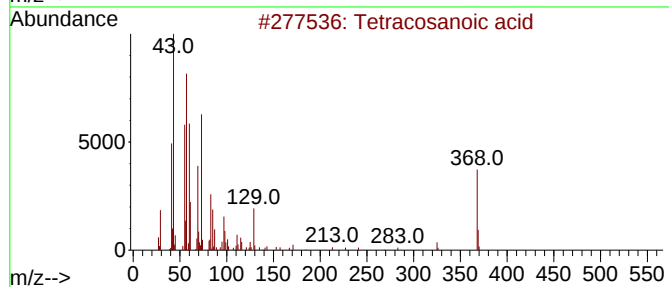
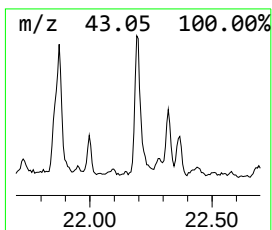
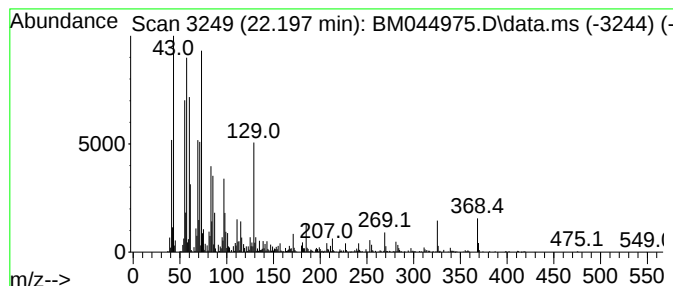
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Quant Title : SVOA CALIBRATION

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

Peak Number 64 Tetracosanoic acid Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.197	15.34 ng/ul	1454790	Chrysene-d12	21.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosanoic acid	368	C24H48O2	000557-59-5	99
2			Tetracosanoic acid	368	C24H48O2	000557-59-5	96
3			Tetracosanoic acid	368	C24H48O2	000557-59-5	94
4			Octadecanoic acid	284	C18H36O2	000057-11-4	53
5			n-Decanoic acid	172	C10H20O2	000334-48-5	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
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Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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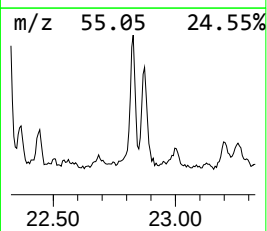
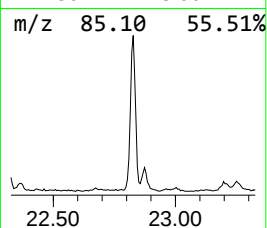
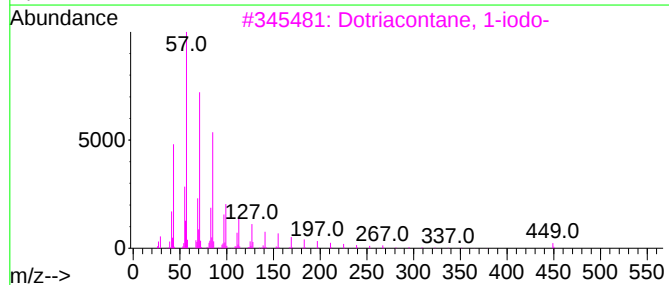
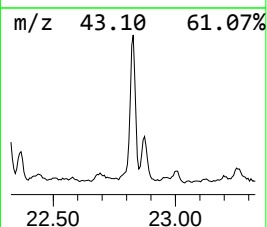
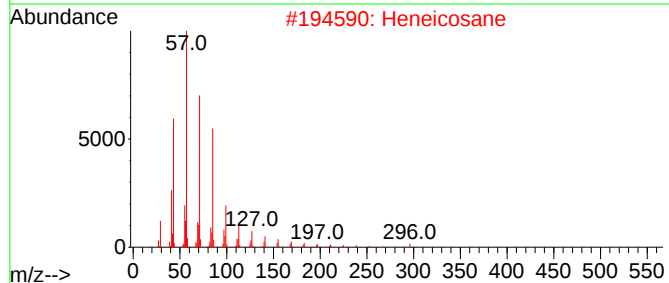
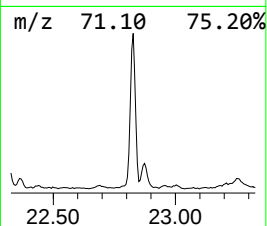
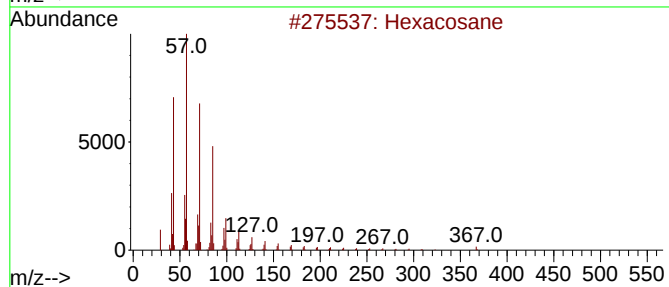
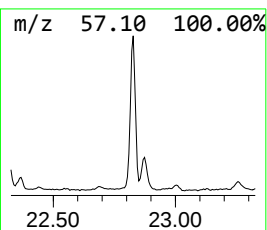
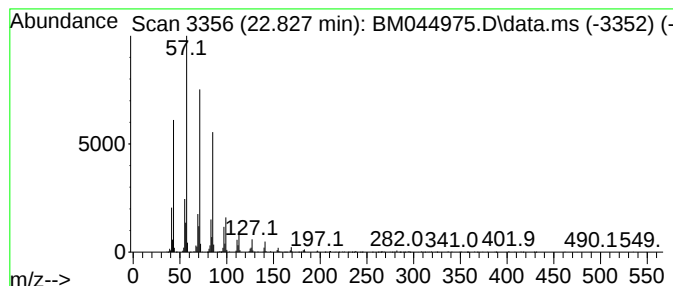
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Quant Title : SVOA CALIBRATION

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

Peak Number 67 (DEL) Alkane: Straight-Chai... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.827	10.53 ng/ul	1007410	Perylene-d12	23.480

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexacosane	366	C26H54	000630-01-3	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	90
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
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ClientSampleId :
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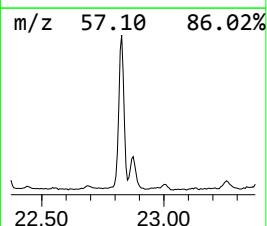
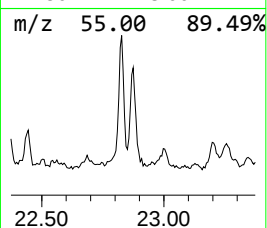
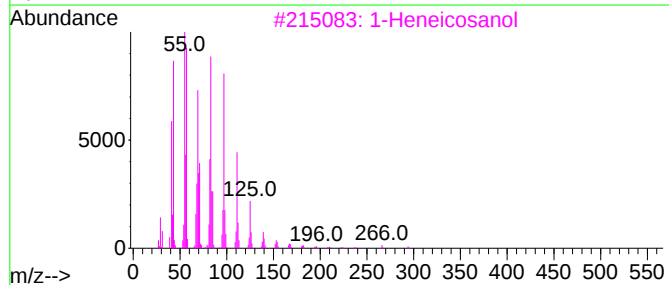
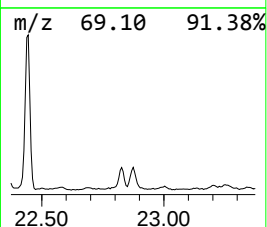
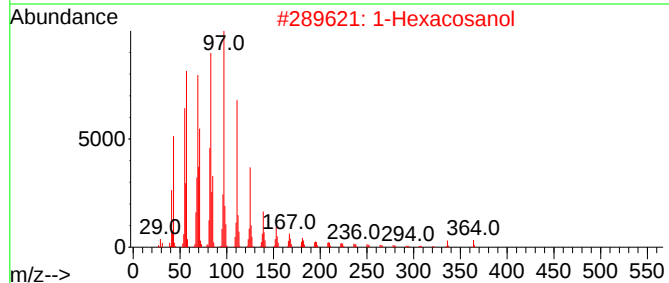
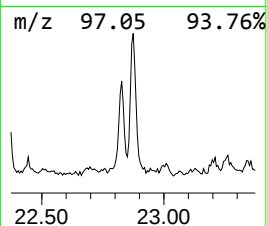
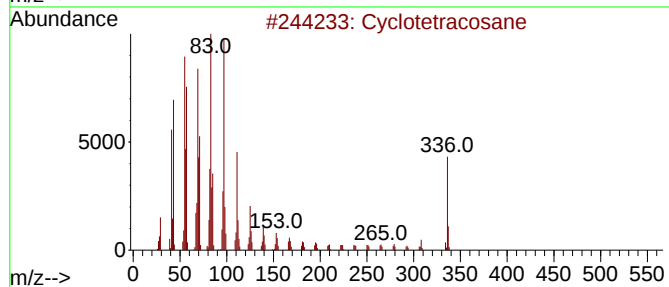
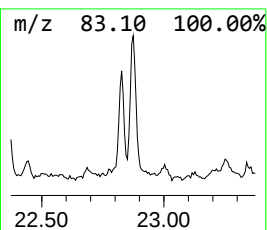
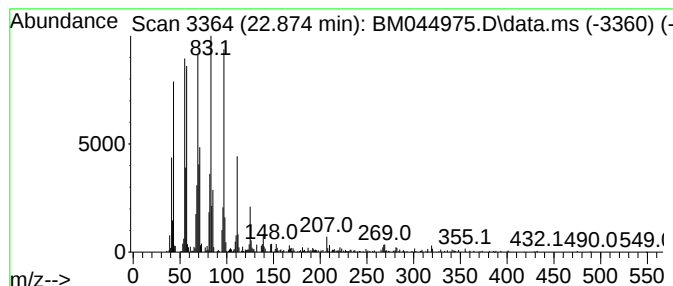
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Quant Title : SVOA CALIBRATION

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

Peak Number 68 (DEL) Alkane: Cyclic22.874 Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.874	6.73 ng/ul	644194	Perylene-d12	23.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	95
2			1-Hexacosanol	382	C26H54O	000506-52-5	91
3			1-Heneicosanol	312	C21H44O	015594-90-8	91
4			Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	91
5			1-Tetracosene	336	C24H48	010192-32-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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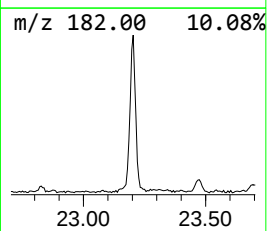
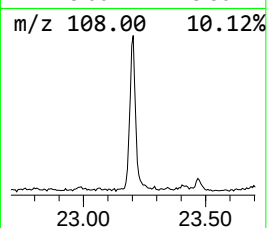
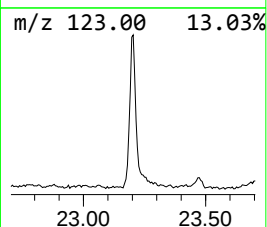
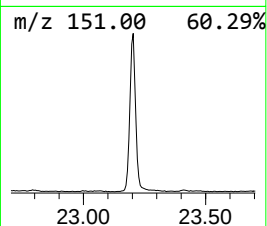
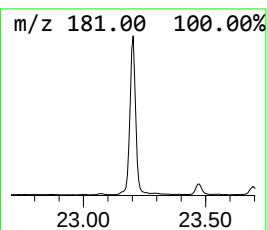
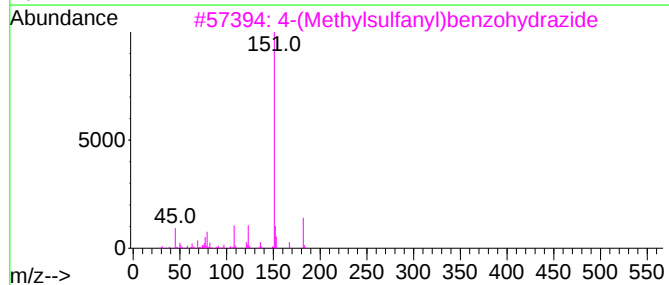
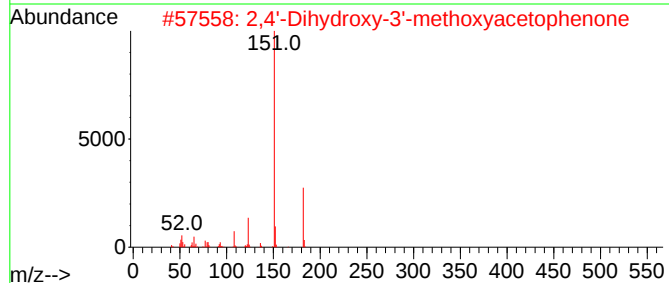
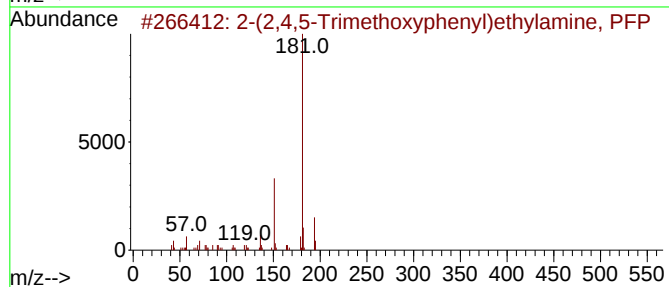
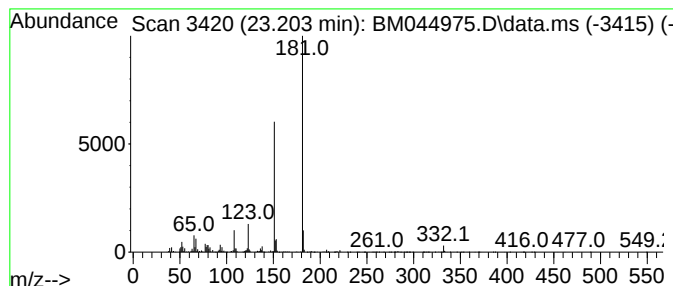
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Quant Title : SVOA CALIBRATION

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

Peak Number 69 2-(2,4,5-Trimethoxyphenyl)e... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.203	16.65 ng/ul	1592880	Perylene-d12	23.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(2,4,5-Trimethoxyphenyl)ethyla...	357	C14H16F5N04	1000071-83-4	64		
2	2,4'-Dihydroxy-3'-methoxyacetoph...	182	C9H10O4	018256-48-9	46		
3	4-(Methylsulfanyl)benzohydrazide	182	C8H10N2OS	1000484-57-9	41		
4	Acetonitrile, (3-chloro-5,5-dime...	181	C10H12ClN	076399-17-2	38		
5	Benzoic acid, 4-hydroxy-3-methox...	182	C9H10O4	003943-74-6	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

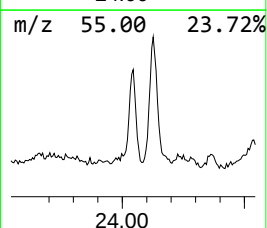
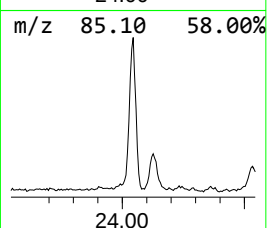
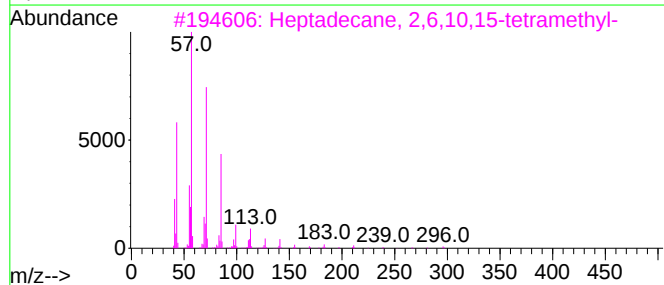
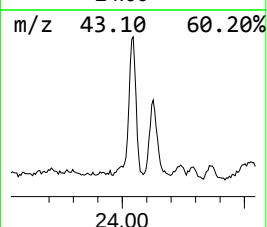
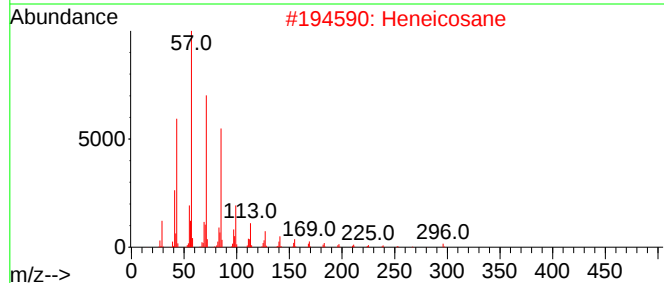
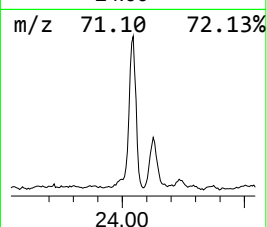
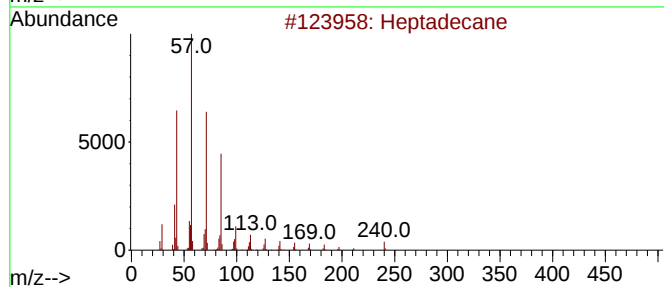
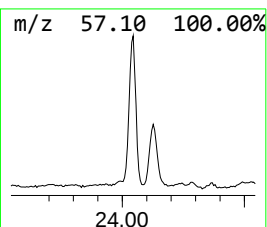
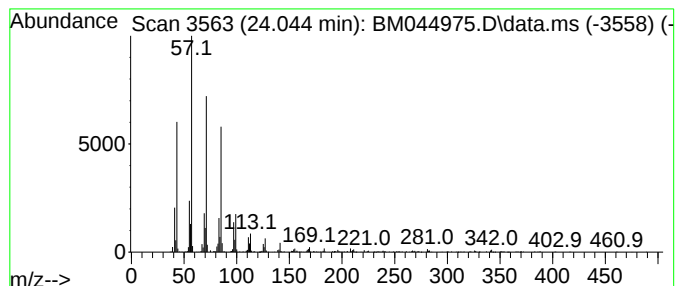
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 70 (DEL) Alkane: Straight-Chai... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD		R.T.	
24.044	9.19 ng/ul	879541	Perylene-d12		23.480	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	86
2	Heneicosane		296	C21H44	000629-94-7	83
3	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	80
4	Nonadecane		268	C19H40	000629-92-5	80
5	Borane, diethyl(decyloxy)-		226	C14H31BO	1000152-34-3	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

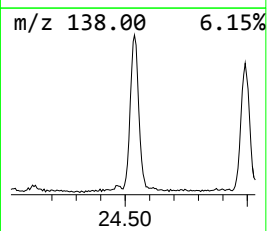
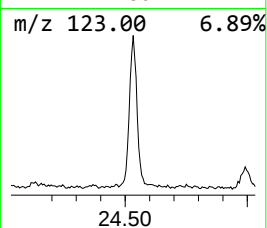
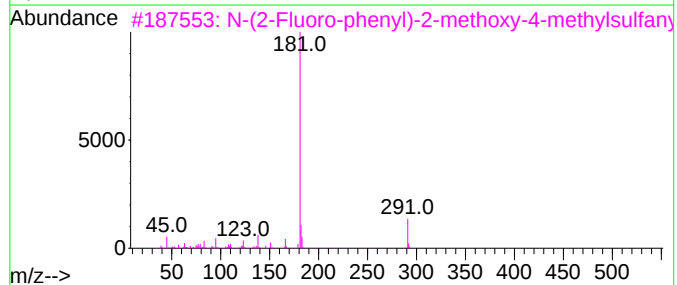
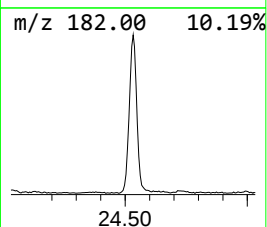
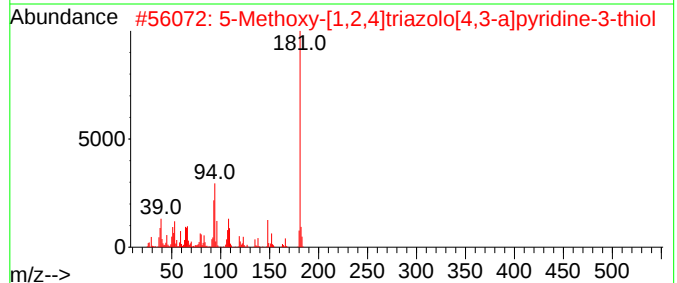
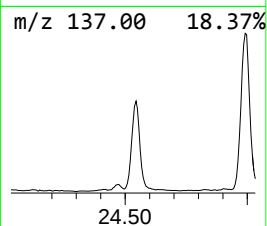
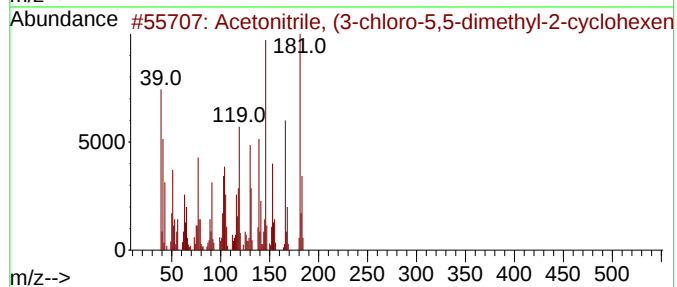
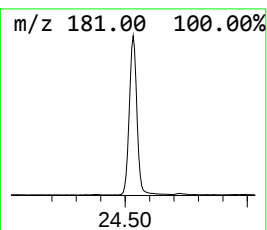
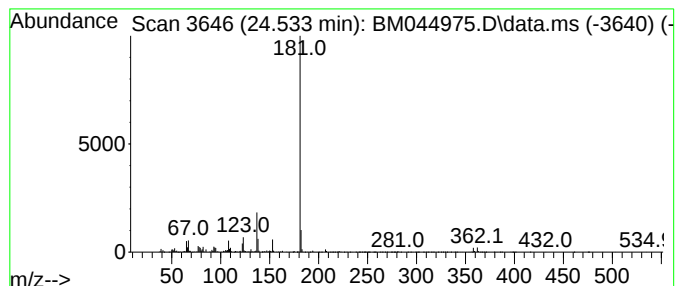
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 72 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD			R.T.
24.532	29.53 ng/ul	2825570	Perylene-d12			23.480
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Acetonitrile, (3-chloro-5,5-dime...	181	C10H12ClN	076399-17-2	47
2		5-Methoxy-[1,2,4]triazolo[4,3-a]...	181	C7H7N3OS	1000186-81-4	42
3		N-(2-Fluoro-phenyl)-2-methoxy-4-...	291	C15H14FN02S	1010294-30-3	39
4		8-Methoxy-[1,2,4]triazolo[4,3-a]...	181	C7H7N3OS	1000186-80-8	38
5		Danielone	212	C10H12O5	090426-22-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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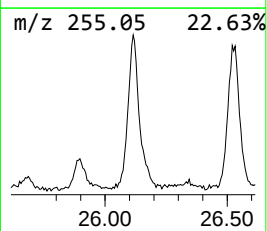
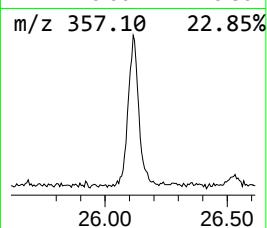
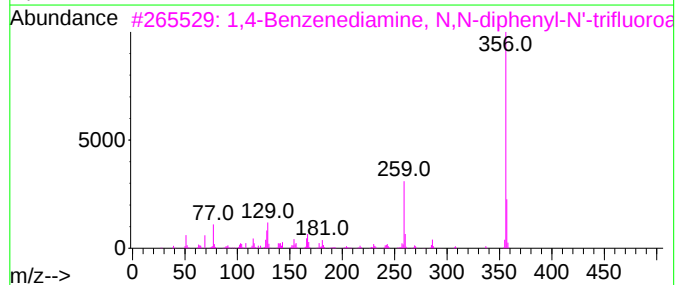
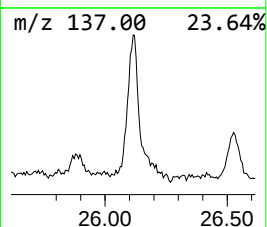
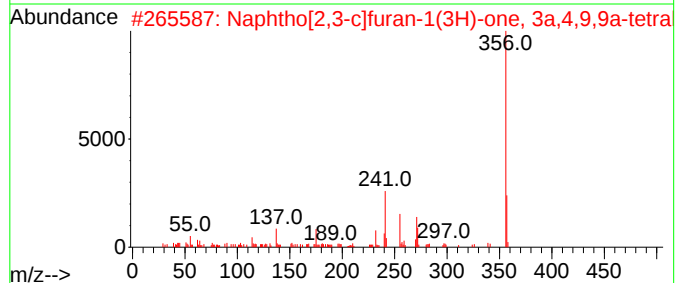
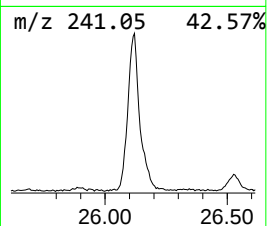
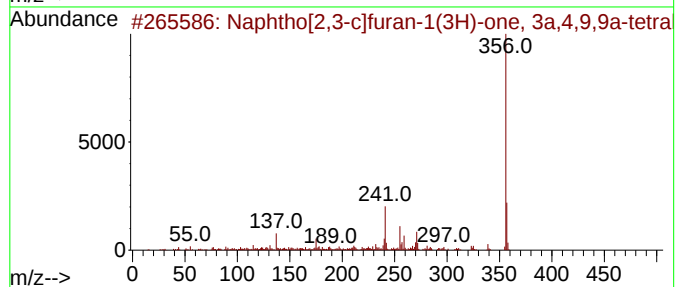
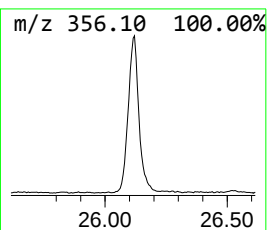
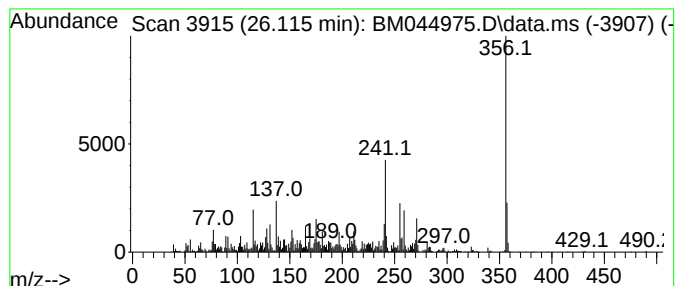
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Quant Title : SVOA CALIBRATION

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

Peak Number 74 Naphtho[2,3-c]furan-1(3H)-o... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.115	30.03 ng/ul	2873530	Perylene-d12	23.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphtho[2,3-c]furan-1(3H)-one, 3...	356	C20H20O6	000518-55-8	64
2			Naphtho[2,3-c]furan-1(3H)-one, 3...	356	C20H20O6	000518-55-8	64
3			1,4-Benzenediamine, N,N-diphenyl...	356	C20H15F3N2O	1000495-45-7	46
4			Cobalt, hexamethylbenzene-pentam...	356	C22H33Co	1000161-46-2	41
5			N-(3-Cyano-4,5-diphenylfuran-2-y...	356	C19H11F3N2O2	1000311-96-1	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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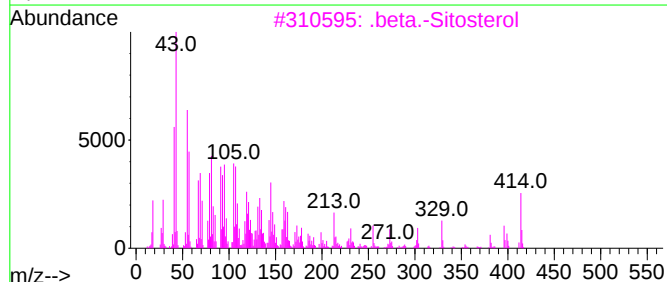
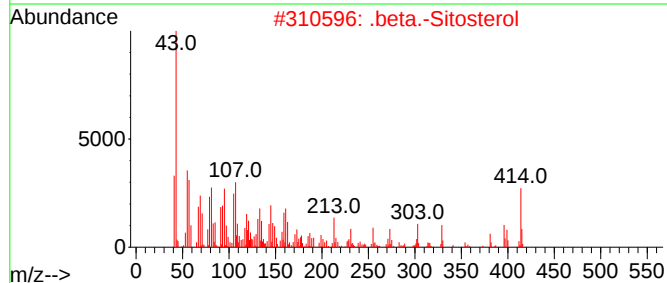
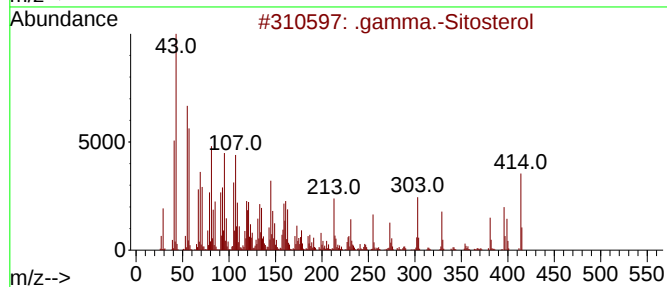
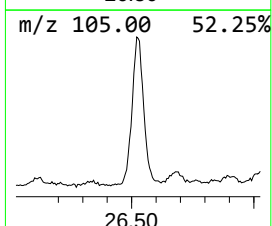
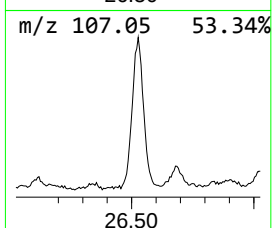
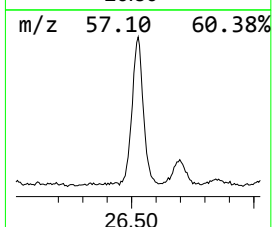
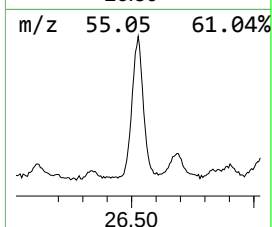
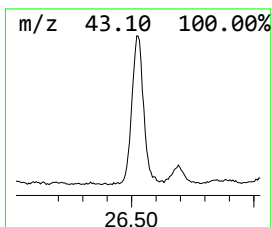
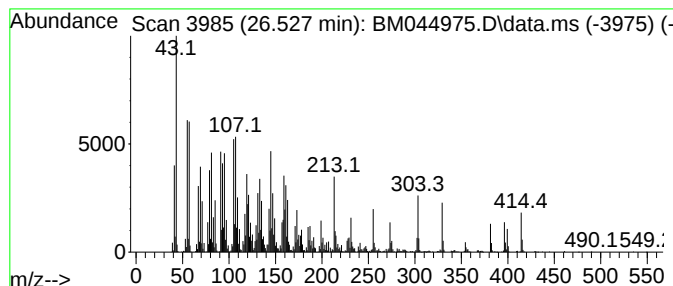
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Quant Title : SVOA CALIBRATION

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

Peak Number 75 .gamma.-Sitosterol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.526	60.27 ng/ul	5767040	Perylene-d12	23.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	gamma.-Sitosterol	414	C29H50O	000083-47-6	99	
2	.	beta.-Sitosterol	414	C29H50O	000083-46-5	99	
3	.	beta.-Sitosterol	414	C29H50O	000083-46-5	91	
4	.	gamma.-Sitosterol	414	C29H50O	000083-47-6	83	
5	(3S,8S,9S,10R,13R,14S,17R)-17-((...	414	C29H50O	029365-29-5	56		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
Data File : BM044975.D
Acq On : 06 Apr 2024 17:51
Operator : MA/JU
Sample : P2018-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BH5F1

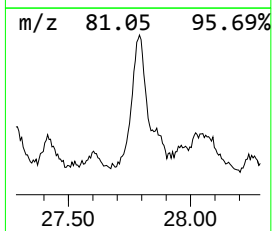
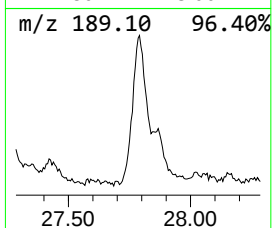
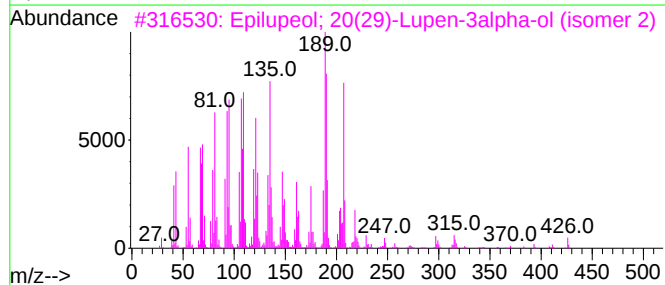
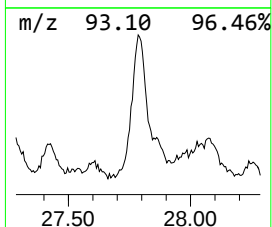
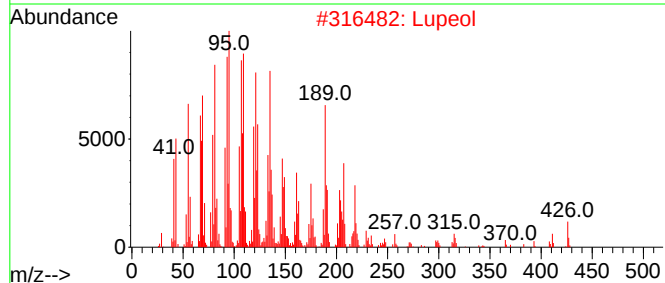
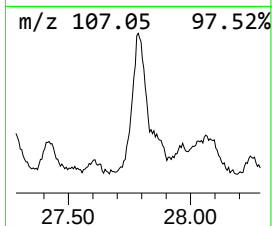
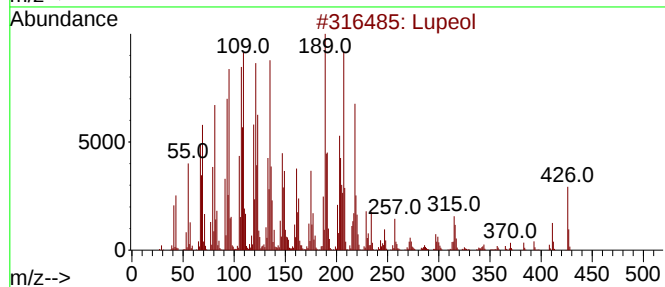
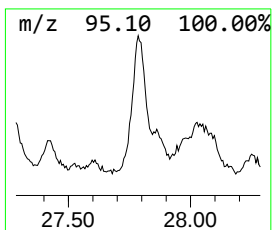
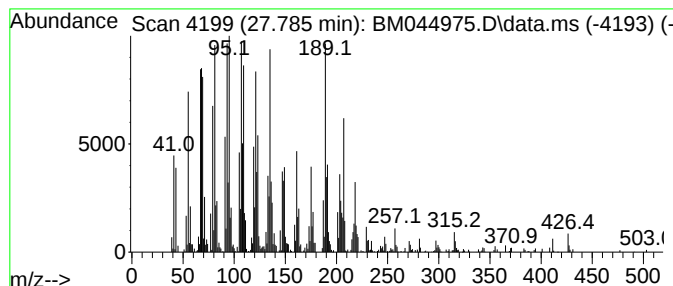
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Quant Title : SVOA CALIBRATION

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

Peak Number 76 Lupeol Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.785	16.40 ng/ul	1569110	Perylene-d12	23.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Lupeol	426	C30H50O	000545-47-1	97
2			Lupeol	426	C30H50O	000545-47-1	97
3			Epilupeol; 20(29)-Lupen-3alpha-o...	426	C30H50O	1000513-01-4	91
4			Epilupeol; 20(29)-Lupen-3alpha-o...	468	C32H52O2	1010513-01-8	64
5			Guaia-9,11-diene	204	C15H24	1000374-19-8	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
 Data File : BM044975.D
 Acq On : 06 Apr 2024 17:51
 Operator : MA/JU
 Sample : P2018-10
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propanoic acid,...	3.557	22.5	ng/ul	683125	1	7.634	607032	20.0
Butanoic acid	3.963	67.4	ng/ul	2046450	1	7.634	607032	20.0
Pentanoic acid	5.210	14.8	ng/ul	450189	1	7.634	607032	20.0
Hexanoic acid	6.716	15.6	ng/ul	473801	1	7.634	607032	20.0
Benzoic acid	9.751	21.1	ng/ul	898932	2	10.410	851288	20.0
Vanillin	13.222	41.8	ng/ul	2364090	3	14.280	1130580	20.0
3-Hydroxy-4-met...	14.810	15.6	ng/ul	884290	3	14.280	1130580	20.0
Benzaldehyde, 4...	15.780	123.4	ng/ul	13035000	4	17.045	2112640	20.0
Azulene, 1,4-di...	16.021	14.5	ng/ul	1530090	4	17.045	2112640	20.0
Coniferyl aldehyde	16.457	16.9	ng/ul	1780880	4	17.045	2112640	20.0
Benzoic acid, 4...	16.933	14.7	ng/ul	1552240	4	17.045	2112640	20.0
n-Hexadecanoic ...	17.974	76.5	ng/ul	8086650	4	17.045	2112640	20.0
trans-Sinapalde...	18.251	25.8	ng/ul	2725780	4	17.045	2112640	20.0
(DEL) Alkane: S...	18.456	4.8	ng/ul	505119	4	17.045	2112640	20.0
9,12-Octadecadi...	19.121	95.7	ng/ul	10110200	4	17.045	2112640	20.0
9-Octadecenoic ...	19.151	108.6	ng/ul	10298700	5	21.250	1896700	20.0
Octadecanoic acid	19.256	20.5	ng/ul	1945870	5	21.250	1896700	20.0
(DEL) Alkane: S...	19.992	3.1	ng/ul	292561	5	21.250	1896700	20.0
Acridin-9-amine...	20.250	14.9	ng/ul	1416650	5	21.250	1896700	20.0
1-Phenanthrenec...	20.292	15.1	ng/ul	1428570	5	21.250	1896700	20.0
Eicosanoic acid	20.345	18.5	ng/ul	1751650	5	21.250	1896700	20.0
Methyl dehydroa...	20.445	30.7	ng/ul	2913690	5	21.250	1896700	20.0
2-Pentenoic aci...	20.468	20.0	ng/ul	1898300	5	21.250	1896700	20.0
Isopimaric acid	20.827	24.4	ng/ul	2308890	5	21.250	1896700	20.0
Dehydroabietic ...	20.956	46.5	ng/ul	4410480	5	21.250	1896700	20.0
n-Nonadecanol-1	21.874	15.2	ng/ul	1443320	5	21.250	1896700	20.0
Tetracosanoic acid	22.197	15.3	ng/ul	1454790	5	21.250	1896700	20.0
(DEL) Alkane: S...	22.827	10.5	ng/ul	1007410	6	23.480	1913630	20.0
(DEL) Alkane: C...	22.874	6.7	ng/ul	644194	6	23.480	1913630	20.0
2-(2,4,5-Trimet...	23.203	16.6	ng/ul	1592880	6	23.480	1913630	20.0
(DEL) Alkane: S...	24.044	9.2	ng/ul	879541	6	23.480	1913630	20.0
unknown-01	24.532	29.5	ng/ul	2825570	6	23.480	1913630	20.0
Naphtho[2,3-c]f...	26.115	30.0	ng/ul	2873530	6	23.480	1913630	20.0
.gamma.-Sitosterol	26.526	60.3	ng/ul	5767040	6	23.480	1913630	20.0
Lupeol	27.785	16.4	ng/ul	1569110	6	23.480	1913630	20.0