

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
 Data File : BN033408.D
 Acq On : 15 Aug 2024 20:52
 Operator : MA/JU
 Sample : P3502-16
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DCXU9

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN033408.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.158	25	29	40	rVB	72496	104585	0.18%	0.043%
2	3.411	68	72	83	rBV	279689	403636	0.68%	0.167%
3	3.546	91	95	110	rBV	419507	616764	1.04%	0.255%
4	5.693	448	460	468	rVB	85304	132070	0.22%	0.055%
5	6.287	555	561	569	rVB	348029	522992	0.88%	0.216%
6	6.752	628	640	654	rVB	1569293	2492944	4.21%	1.029%
7	6.916	660	668	674	rBV	1219492	1818953	3.07%	0.751%
8	7.105	694	700	709	rVV	1584307	2385782	4.03%	0.985%
9	7.181	709	713	726	rVB	75586	121945	0.21%	0.050%
10	7.569	771	779	787	rBV	2031678	3143880	5.31%	1.298%
11	7.646	787	792	799	rVB	61243	88159	0.15%	0.036%
12	8.269	891	898	911	rBV	1661996	2560726	4.32%	1.057%
13	8.705	965	972	979	rBV	1427621	2302696	3.89%	0.951%
14	9.287	1064	1071	1077	rBV	260129	379896	0.64%	0.157%
15	9.422	1087	1094	1101	rBV	1148379	1834370	3.10%	0.757%
16	9.663	1129	1135	1142	rVB6	47694	98020	0.17%	0.040%
17	9.957	1178	1185	1201	rVB	1921161	3095978	5.23%	1.278%
18	10.334	1241	1249	1257	rVB	2903778	4686386	7.91%	1.935%
19	10.469	1263	1272	1282	rBV	562956	957613	1.62%	0.395%
20	10.893	1337	1344	1351	rVV2	114617	202607	0.34%	0.084%
21	11.081	1370	1376	1383	rBV	398513	674063	1.14%	0.278%
22	12.581	1623	1631	1636	rVB2	53296	89817	0.15%	0.037%
23	13.098	1715	1719	1727	rVB4	113316	218473	0.37%	0.090%
24	13.540	1789	1794	1802	rVB	189864	280922	0.47%	0.116%
25	13.628	1803	1809	1815	rVV	3197474	4286184	7.24%	1.770%
26	13.693	1815	1820	1825	rVV	48373	88923	0.15%	0.037%
27	13.892	1847	1854	1865	rVB2	3535677	5381971	9.09%	2.222%
28	14.069	1876	1884	1889	rBV4	102519	188004	0.32%	0.078%
29	14.204	1900	1907	1915	rVB	4036527	6151132	10.39%	2.540%
30	14.292	1917	1922	1929	rBV4	108709	229707	0.39%	0.095%
31	14.410	1936	1942	1946	rBV	1391327	2063489	3.48%	0.852%
32	14.463	1949	1951	1961	rVV6	177155	366793	0.62%	0.151%
33	14.557	1961	1967	1972	rVB2	149618	236885	0.40%	0.098%
34	14.657	1979	1984	1988	rBV4	81689	121695	0.21%	0.050%
35	14.969	2032	2037	2043	rVB2	82066	126416	0.21%	0.052%
36	15.139	2062	2066	2071	rVV4	92549	143345	0.24%	0.059%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
 Title : SVOA CALIBRATION

37	15.204	2071	2077	2089	rVB	4479565	6478263	10.94%	2.675%
38	15.322	2091	2097	2106	rBV	1177567	1687522	2.85%	0.697%
39	15.563	2128	2138	2143	rBV10	54607	124594	0.21%	0.051%
40	15.710	2149	2163	2168	rBV	175595	376944	0.64%	0.156%
41	15.798	2173	2178	2183	rVB3	100389	143998	0.24%	0.059%
42	15.951	2199	2204	2210	rBV2	60422	122432	0.21%	0.051%
43	16.110	2227	2231	2241	rVB2	122464	208934	0.35%	0.086%
44	16.198	2241	2246	2250	rBV	103385	130455	0.22%	0.054%
45	16.357	2265	2273	2276	rBV3	67206	112293	0.19%	0.046%
46	16.392	2276	2279	2288	rVB	113286	158366	0.27%	0.065%
47	16.610	2310	2316	2327	rVB	404119	637786	1.08%	0.263%
48	16.710	2329	2333	2338	rBV2	72614	107574	0.18%	0.044%
49	16.892	2359	2364	2366	rBV	105595	135001	0.23%	0.056%
50	16.957	2369	2375	2385	rVB	4778620	6908777	11.67%	2.853%
51	17.051	2385	2391	2396	rBV2	4365154	6540360	11.05%	2.701%
52	17.345	2434	2441	2465	rBV	2874319	6552090	11.07%	2.705%
53	17.769	2508	2513	2517	rBV4	90347	149517	0.25%	0.062%
54	17.828	2519	2523	2529	rBV	105062	139431	0.24%	0.058%
55	17.892	2529	2534	2548	rBV	2255371	3386858	5.72%	1.398%
56	18.192	2580	2585	2594	rBV4	171223	328741	0.56%	0.136%
57	18.328	2604	2608	2613	rBV4	86811	128655	0.22%	0.053%
58	18.563	2645	2648	2652	rBV4	68010	91453	0.15%	0.038%
59	18.757	2678	2681	2685	rBV3	78189	115130	0.19%	0.048%
60	18.839	2691	2695	2700	rVB	151085	204461	0.35%	0.084%
61	18.916	2704	2708	2714	rVB5	134080	200849	0.34%	0.083%
62	18.986	2714	2720	2723	rBV2	89954	154925	0.26%	0.064%
63	19.033	2723	2728	2730	rBV2	238612	350608	0.59%	0.145%
64	19.063	2730	2733	2735	rVV	259555	323951	0.55%	0.134%
65	19.086	2735	2737	2741	rVB	270686	314708	0.53%	0.130%
66	19.192	2750	2755	2766	rBV	1703769	2490299	4.21%	1.028%
67	19.357	2777	2783	2788	rBV	4996742	6680122	11.28%	2.758%
68	19.439	2788	2797	2822	rBV	28154765	59212813	100.00%	24.450%
69	19.922	2875	2879	2883	rBV2	501904	704465	1.19%	0.291%
70	19.957	2883	2885	2890	rVB	377141	404967	0.68%	0.167%
71	20.275	2935	2939	2945	rBV2	320421	430740	0.73%	0.178%
72	20.369	2952	2955	2960	rVB	240836	264806	0.45%	0.109%
73	20.427	2960	2965	2968	rBV4	276136	512148	0.86%	0.211%
74	20.580	2988	2991	2995	rVB2	169286	203410	0.34%	0.084%
75	20.627	2996	2999	3006	rVV4	252646	325941	0.55%	0.135%
76	20.733	3012	3017	3021	rBV	3125906	3548292	5.99%	1.465%

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Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
 Title : SVOA CALIBRATION

77	20.869	3035	3040	3044	rBV	1870991	2481292	4.19%	1.025%
78	20.916	3044	3048	3061	rVB	5187488	7081230	11.96%	2.924%
79	21.080	3072	3076	3080	rBV	319258	454193	0.77%	0.188%
80	21.139	3084	3086	3089	rVV3	235546	295650	0.50%	0.122%
81	21.186	3089	3094	3102	rVV2	3664913	5324062	8.99%	2.198%
82	21.257	3102	3106	3113	rVB	730350	960090	1.62%	0.396%
83	21.427	3131	3135	3140	rBV	598938	828128	1.40%	0.342%
84	21.645	3168	3172	3181	rVB2	265761	386009	0.65%	0.159%
85	21.798	3195	3198	3207	rVB3	187387	309050	0.52%	0.128%
86	21.992	3224	3231	3245	rBV	7106139	11749649	19.84%	4.852%
87	22.404	3296	3301	3312	rBV2	676992	1221344	2.06%	0.504%
88	22.598	3331	3334	3353	rVB3	455591	1414182	2.39%	0.584%
89	22.898	3381	3385	3392	rVB2	219717	372618	0.63%	0.154%
90	23.310	3447	3455	3459	rBV	4642883	9495447	16.04%	3.921%
91	23.351	3459	3462	3474	rVB	2070142	3827704	6.46%	1.580%
92	23.827	3534	3543	3551	rBV	2008082	4973340	8.40%	2.054%
93	24.015	3568	3575	3585	rVB	2176671	5014037	8.47%	2.070%
94	24.133	3588	3595	3602	rVB3	408976	1068144	1.80%	0.441%
95	24.580	3667	3671	3683	rVB3	298692	693331	1.17%	0.286%
96	25.127	3754	3764	3773	rBV	3738956	10278910	17.36%	4.244%
97	25.221	3773	3780	3795	rVB2	1693265	4952373	8.36%	2.045%
98	26.227	3945	3951	3958	rBV4	264400	623344	1.05%	0.257%
99	26.974	4070	4078	4096	rVB2	604692	2246894	3.79%	0.928%
100	28.656	4353	4364	4378	rVB3	1556368	6441603	10.88%	2.660%

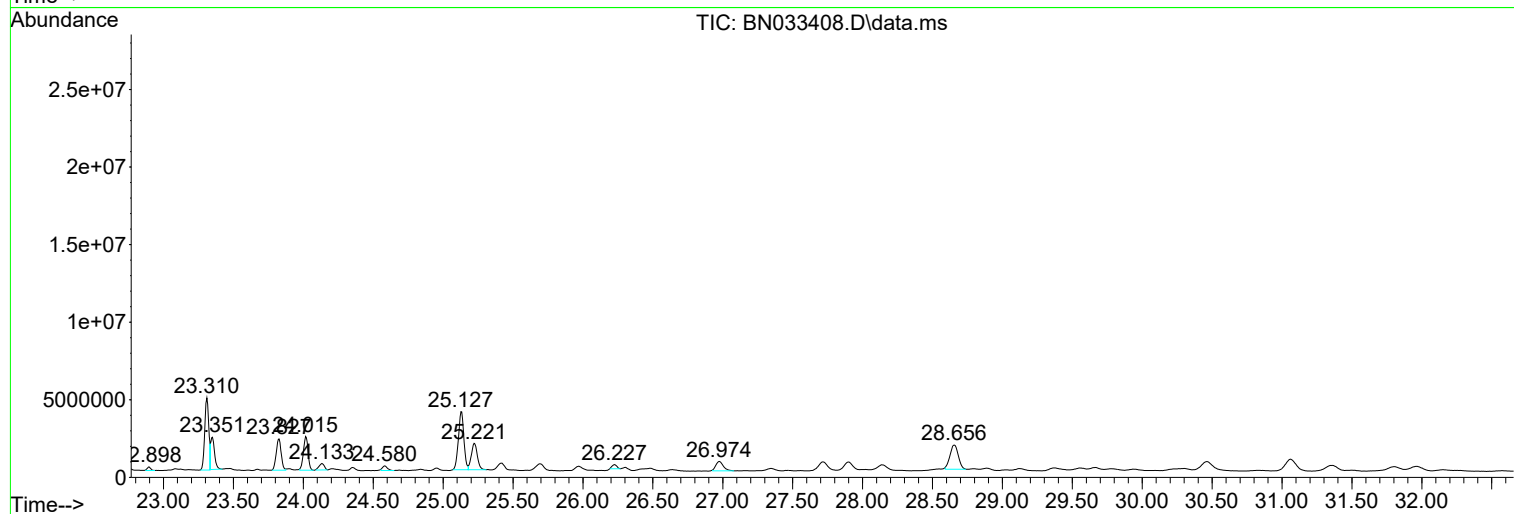
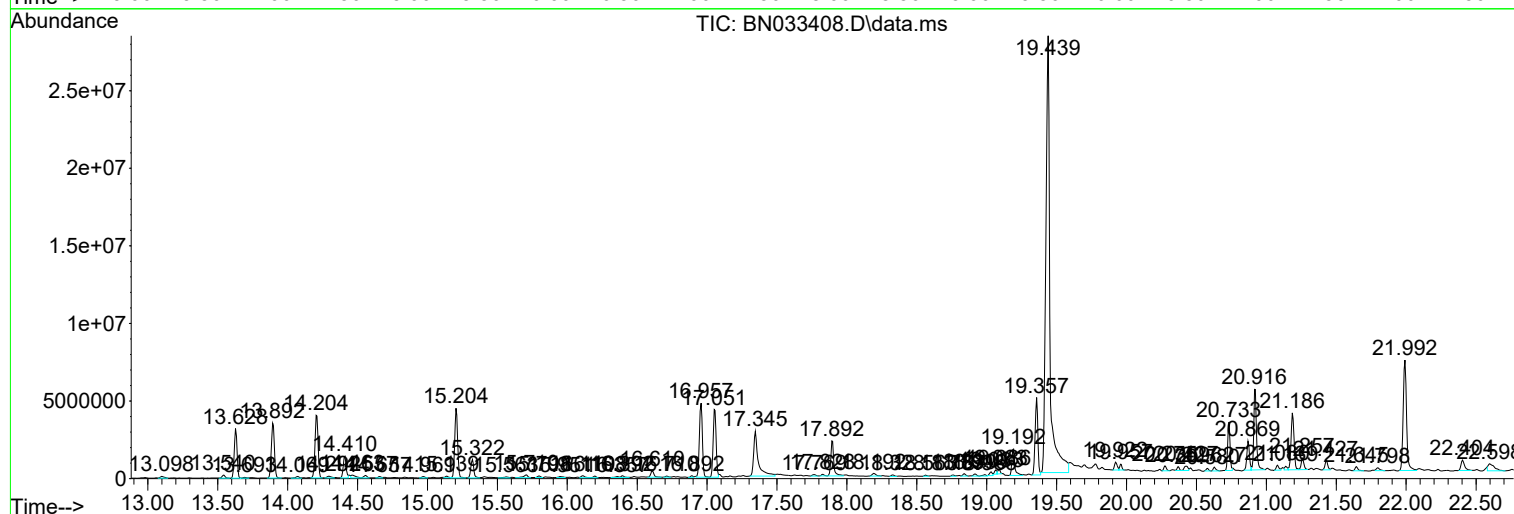
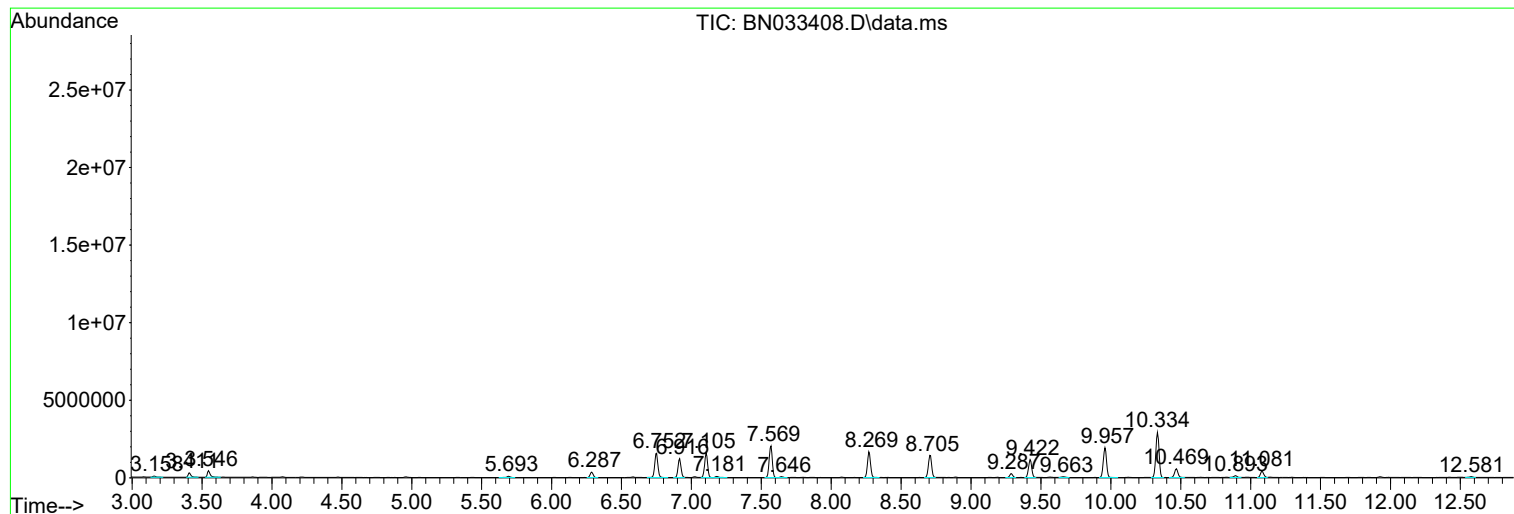
Sum of corrected areas: 242183124

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ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
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Instrument :
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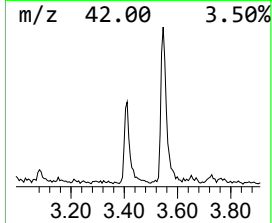
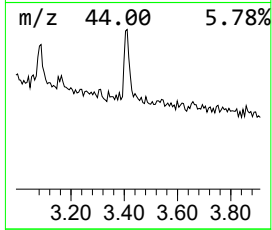
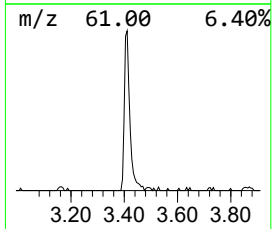
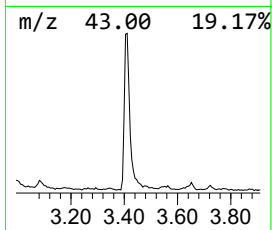
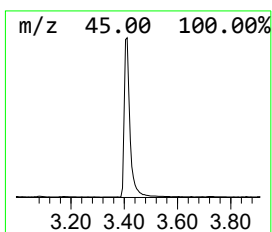
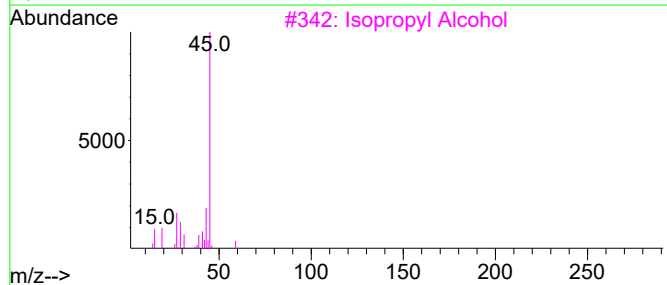
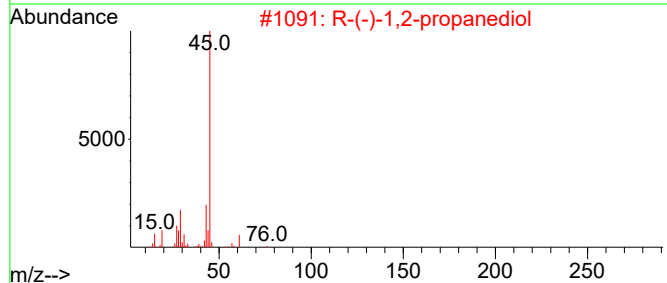
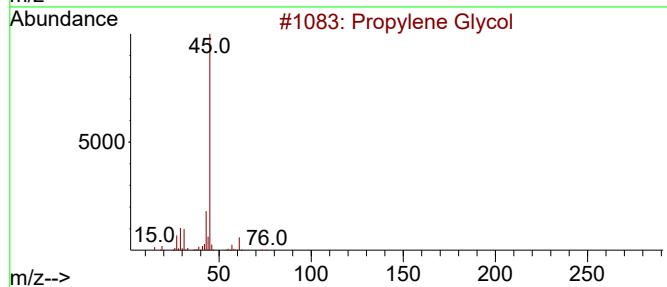
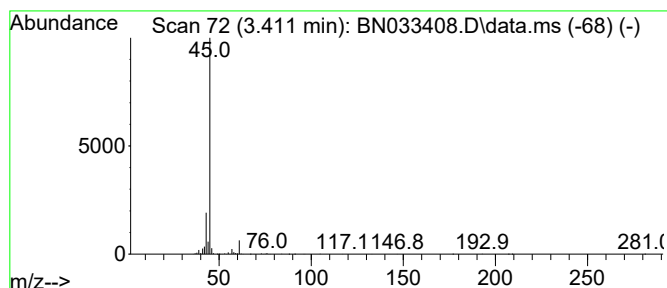
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.411	2.57 ng/ul	403636	1,4-Dichlorobenzene-d4	7.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	90
2			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	78
3			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
4			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
5			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	40



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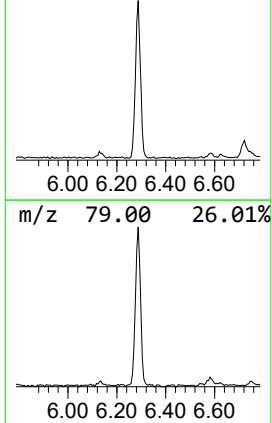
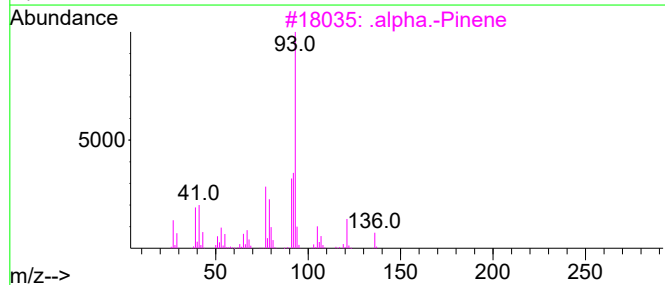
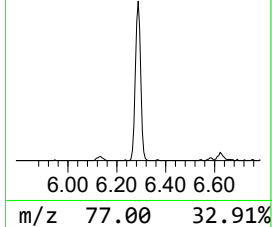
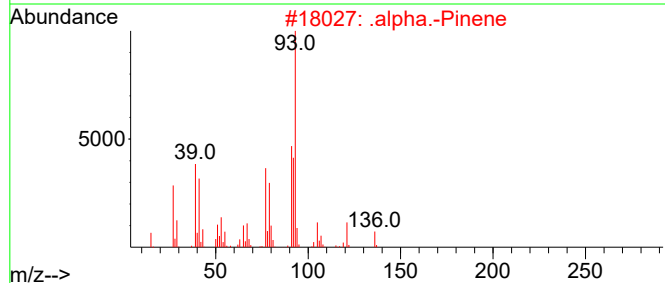
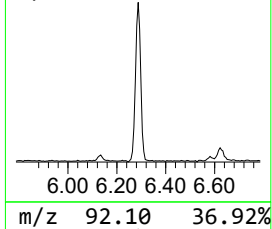
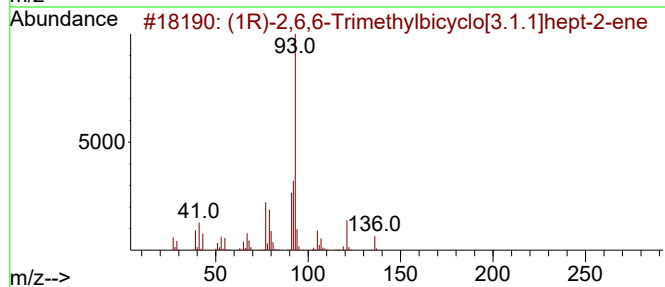
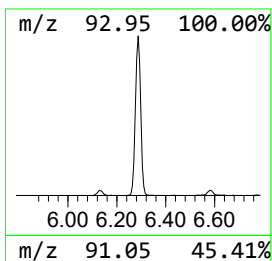
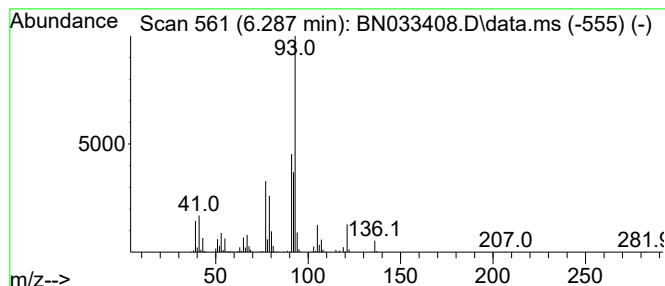
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.287	3.33 ng/ul	522992	1,4-Dichlorobenzene-d4	7.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-70-8	95		
2	.alpha.-Pinene	136	C10H16	000080-56-8	95		
3	.alpha.-Pinene	136	C10H16	000080-56-8	94		
4	Bicyclo[3.1.1]hept-2-ene, 3,6,6-trimethyl-	136	C10H16	004889-83-2	94		
5	(1S)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-26-4	91		



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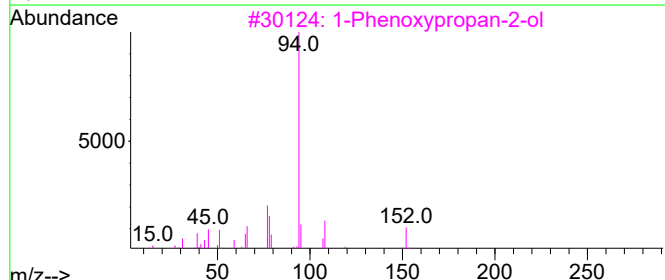
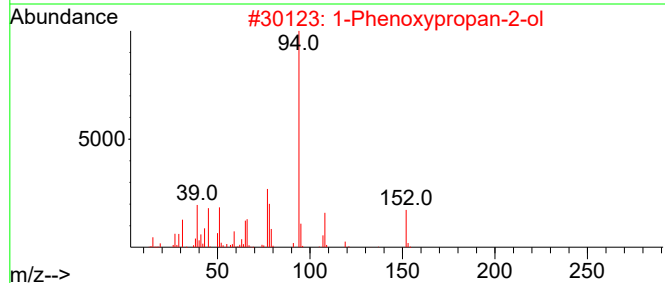
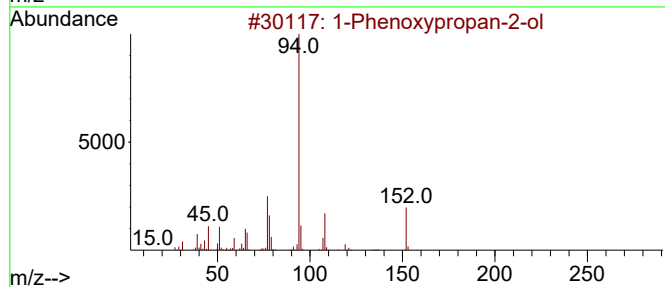
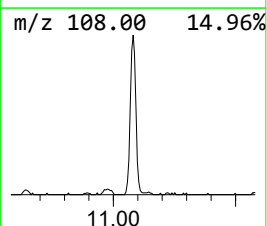
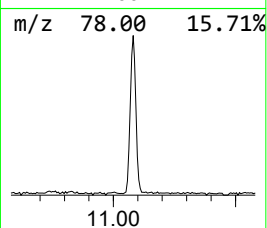
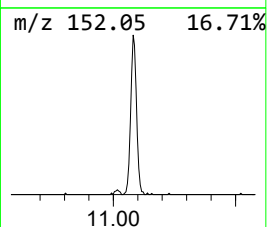
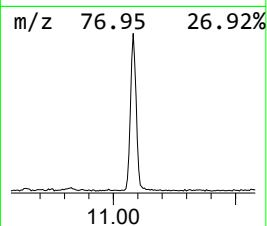
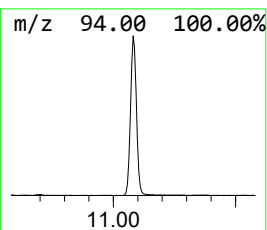
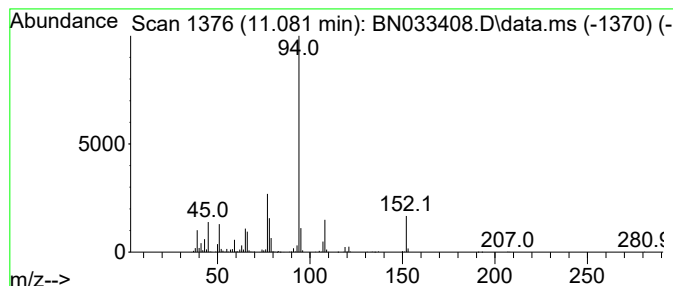
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.081	2.88 ng/ul	674063	Naphthalene-d8	10.334

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
Operator : MA/JU
Sample : P3502-16
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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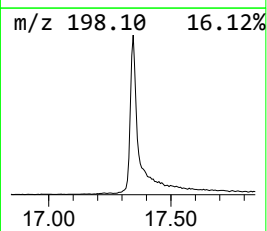
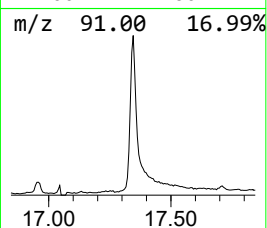
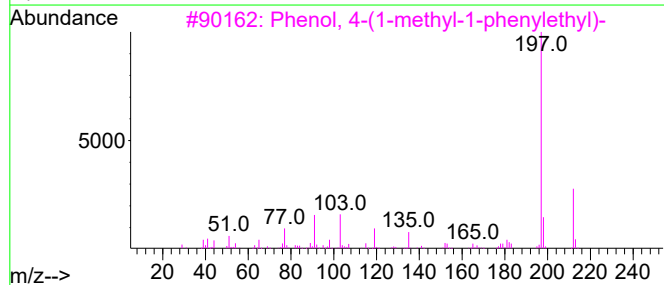
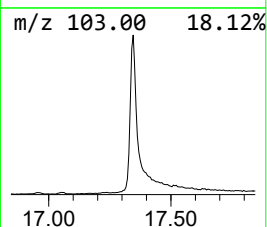
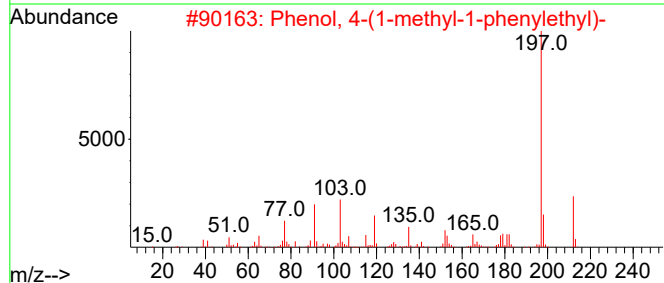
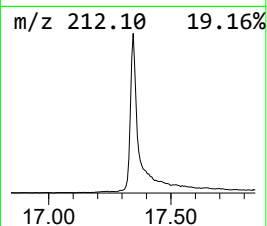
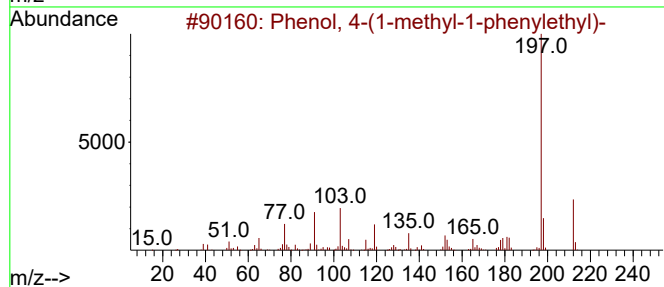
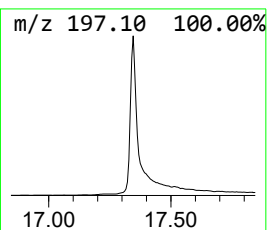
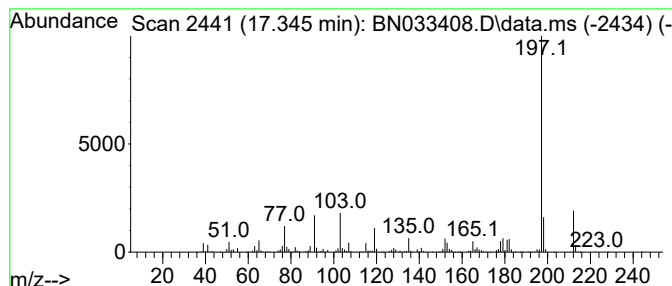
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenol, 4-(1-methyl-1-pheny... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.345	18.97 ng/ul	6552090	Phenanthrene-d10	16.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1-methyl-1-phenylethyl)-	212	C15H16O	000599-64-4	99
2			Phenol, 4-(1-methyl-1-phenylethyl)-	212	C15H16O	000599-64-4	98
3			Phenol, 4-(1-methyl-1-phenylethyl)-	212	C15H16O	000599-64-4	91
4			Dimethyl diphenylsilane	212	C14H16Si	000778-24-5	68
5			2,3-Dimethyl-4-biphenylamine	197	C14H15N	022750-87-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
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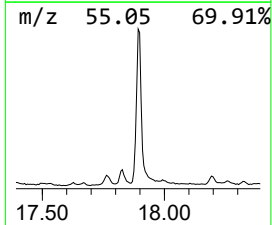
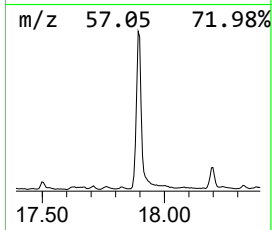
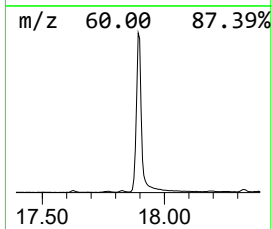
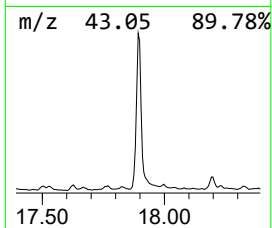
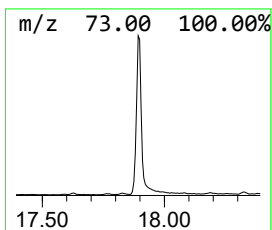
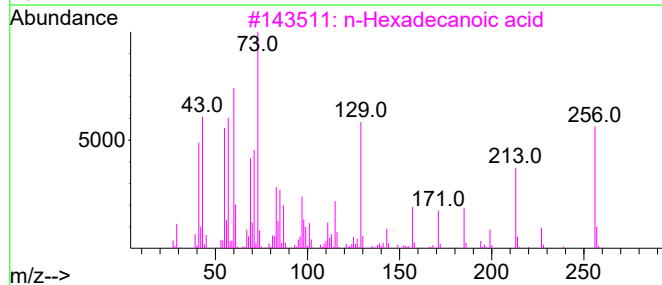
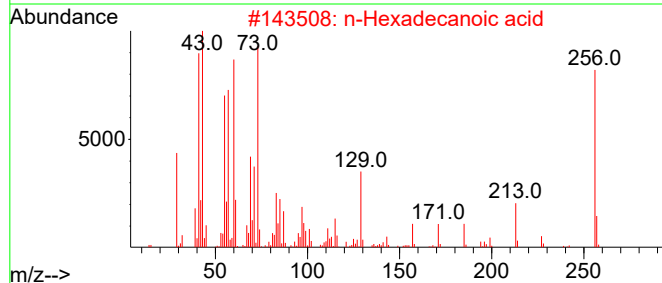
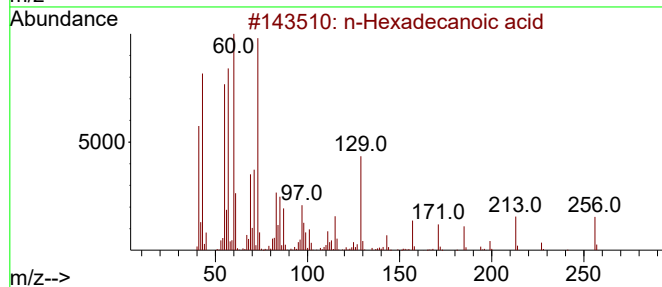
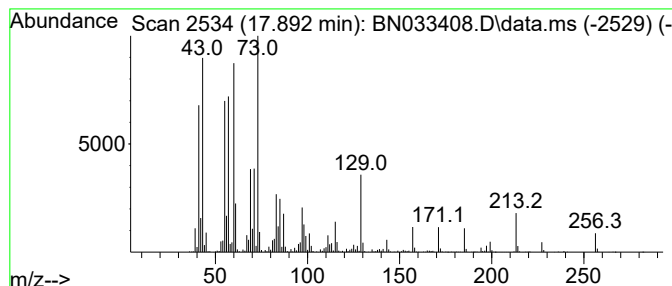
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.892	9.80 ng/ul	3386860	Phenanthrene-d10	16.957

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			Tridecanoic acid	214	C13H26O2	000638-53-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
Operator : MA/JU
Sample : P3502-16
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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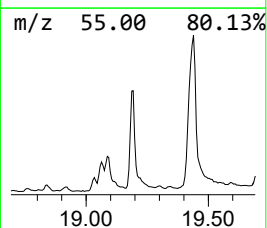
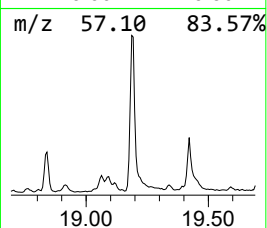
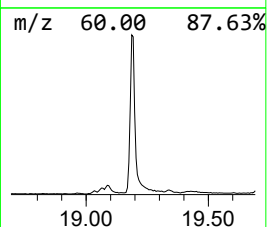
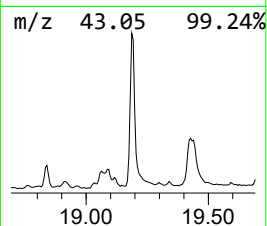
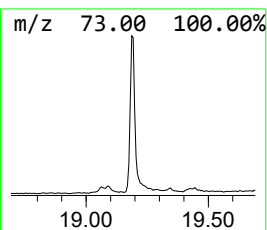
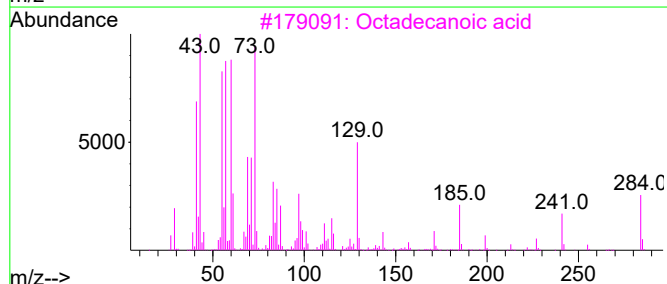
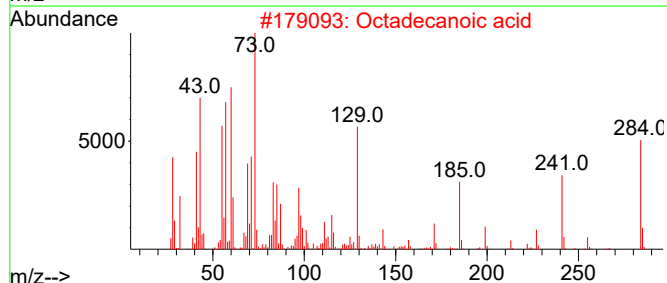
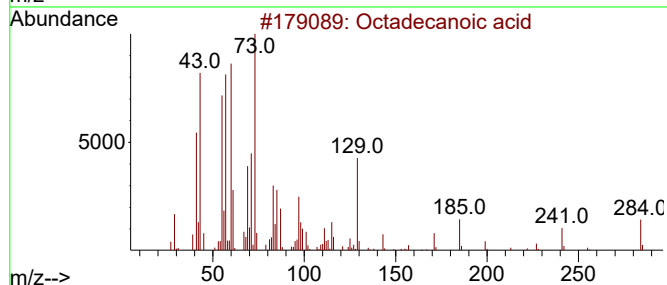
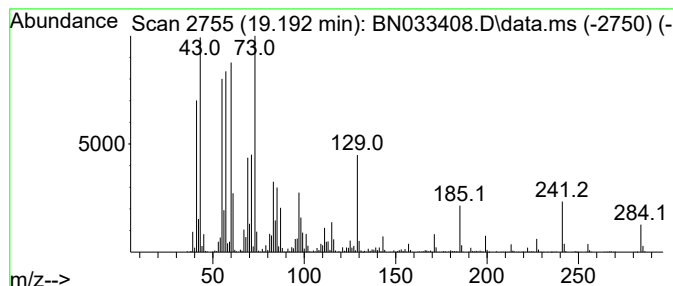
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.192	9.35 ng/ul	2490300	Chrysene-d12	21.186

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
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Sample : P3502-16
Misc :
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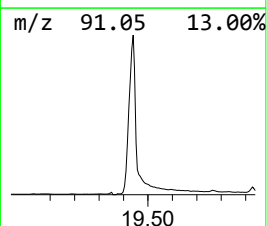
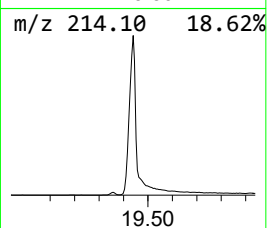
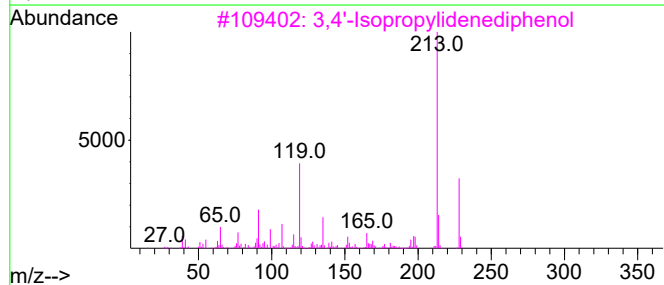
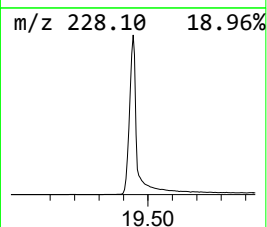
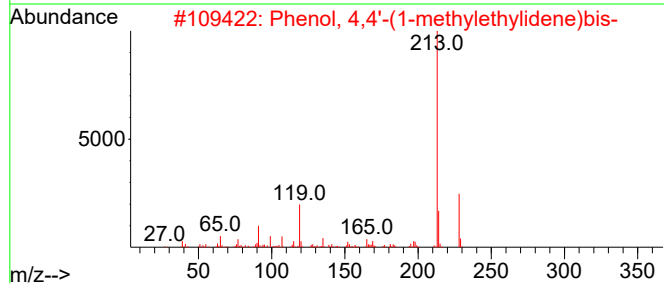
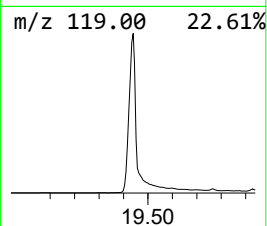
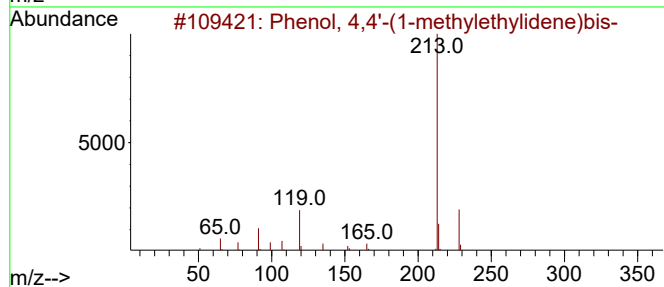
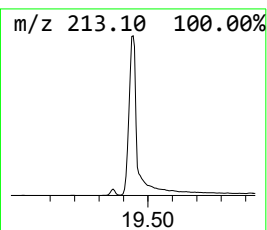
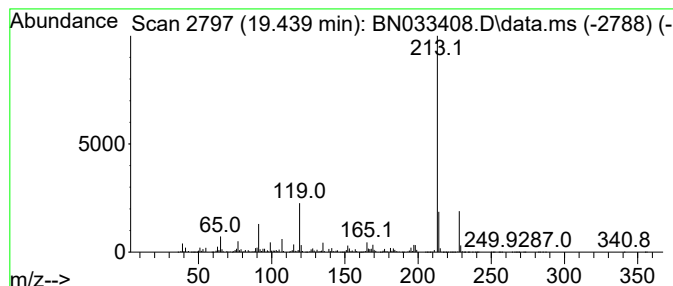
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Phenol, 4,4'-(1-methylethyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.439	222.44 ng/ul	59212800	Chrysene-d12	21.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	98
2			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	96
3			3,4'-Isopropylidenediphenol	228	C15H16O2	046765-25-7	91
4			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	90
5			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
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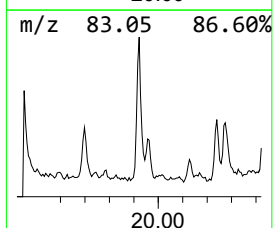
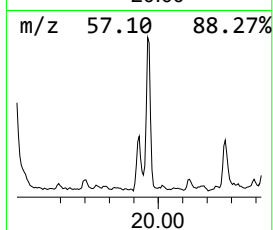
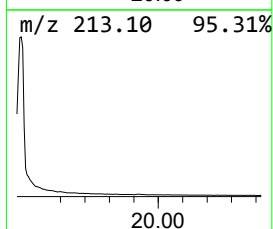
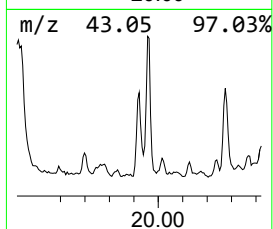
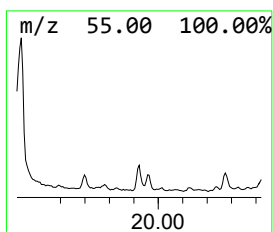
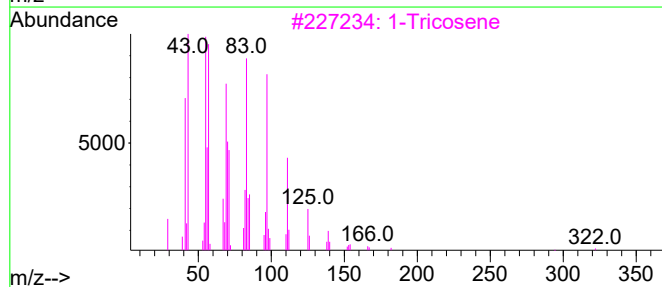
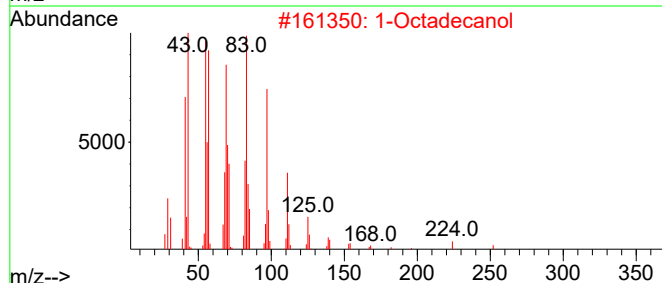
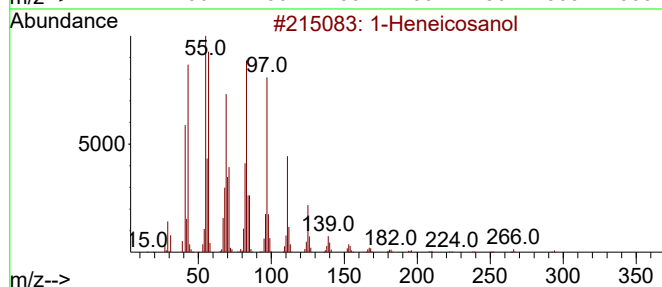
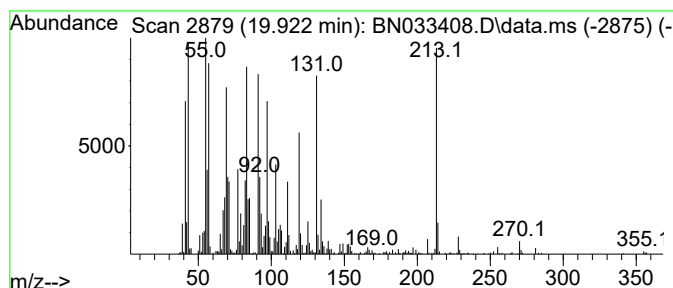
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 1-Heneicosanol Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.922	2.65 ng/ul	704465	Chrysene-d12	21.186	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	62
2	1-Octadecanol	270	C18H38O	000112-92-5	60
3	1-Tricosene	322	C23H46	018835-32-0	52
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	52
5	Behenic alcohol	326	C22H46O	000661-19-8	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
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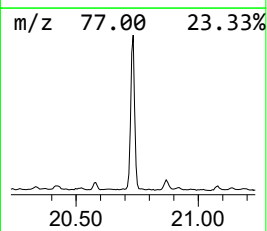
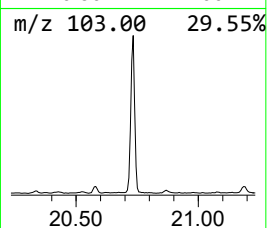
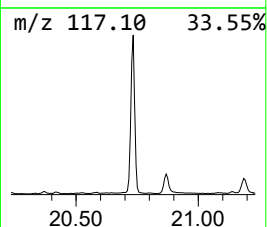
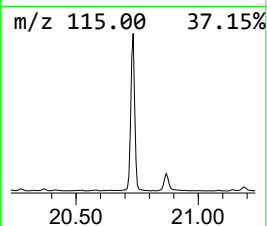
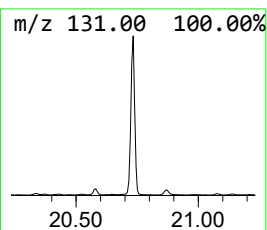
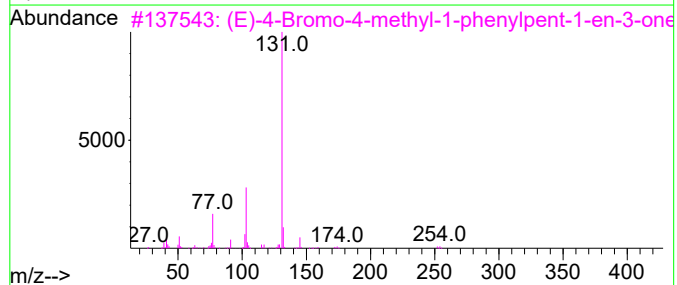
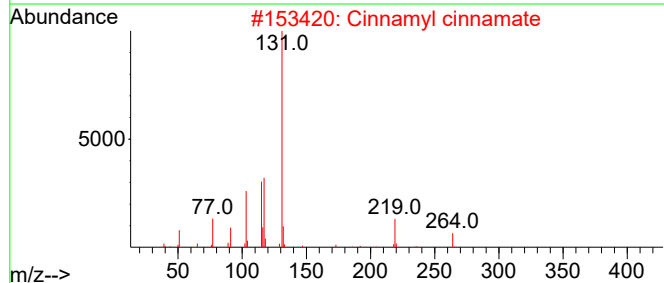
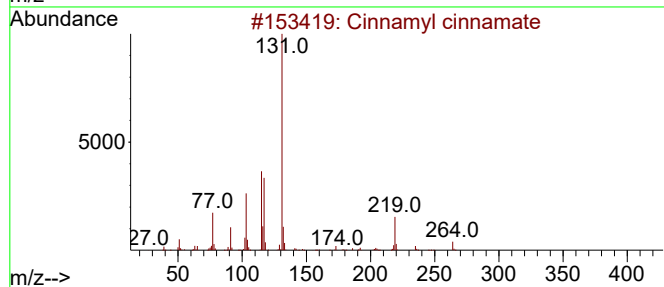
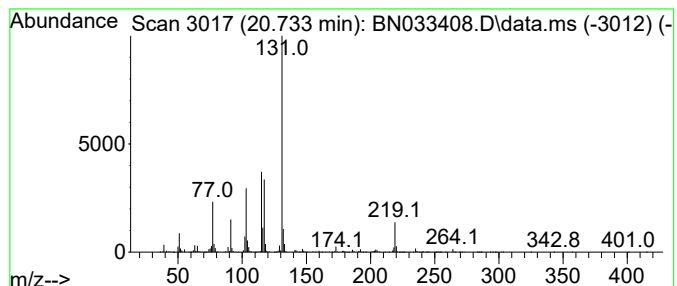
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Cinnamyl cinnamate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.733	13.33 ng/ul	3548290	Chrysene-d12	21.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cinnamyl cinnamate	264	C18H16O2	000122-69-0	90
2			Cinnamyl cinnamate	264	C18H16O2	000122-69-0	87
3			(E)-4-Bromo-4-methyl-1-phenylpen...	252	C12H13BrO	1000443-02-8	53
4			Cinnamoylglycine, methyl ester	219	C12H13NO3	040778-04-9	53
5			2,2-Dimethyl-1-(2-vinylphenyl)pr...	188	C13H16O	1000210-99-9	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
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ClientSampleId :
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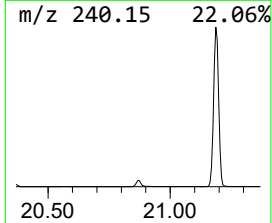
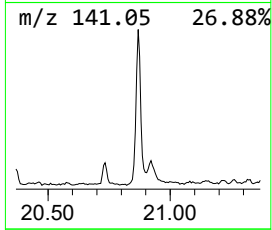
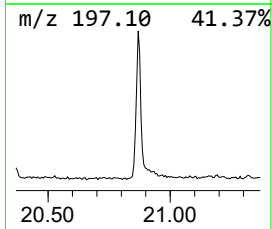
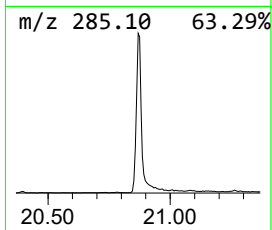
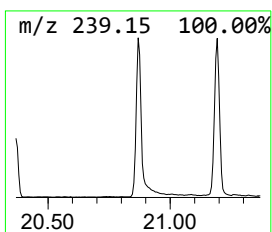
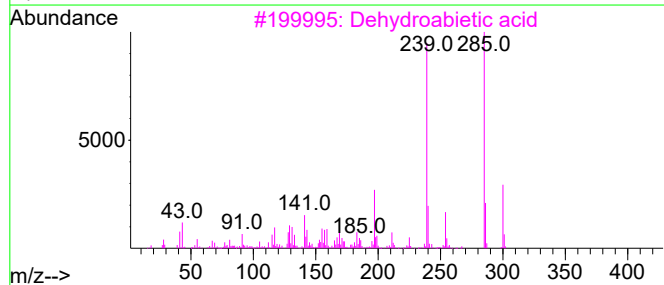
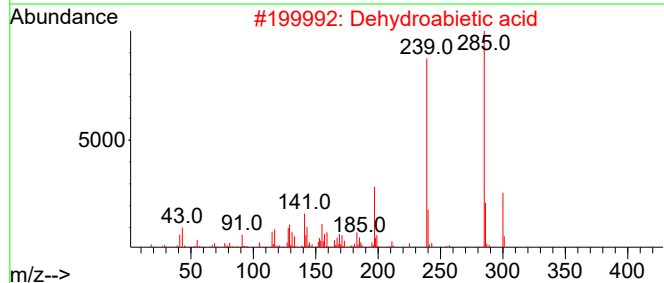
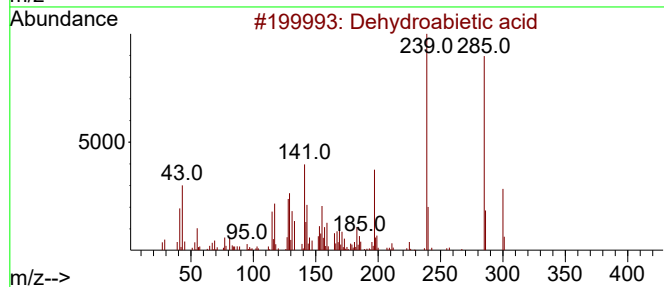
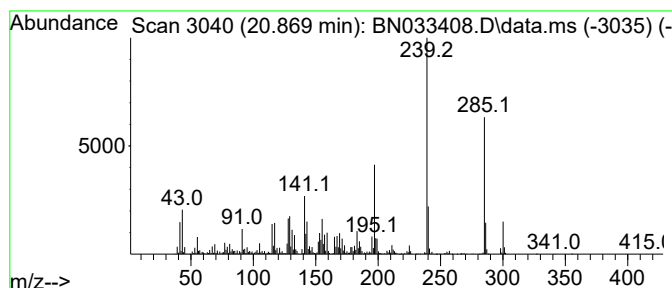
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Dehydroabietic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.869	9.32 ng/ul	2481290	Chrysene-d12	21.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dehydroabietic acid	300	C20H28O2	001740-19-8	99
2			Dehydroabietic acid	300	C20H28O2	001740-19-8	98
3			Dehydroabietic acid	300	C20H28O2	001740-19-8	93
4			1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	76
5			Kaura-9(11),16-dien-18-oic acid,...	300	C20H28O2	022338-67-6	59



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Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
Operator : MA/JU
Sample : P3502-16
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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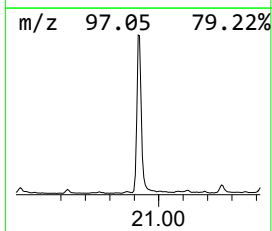
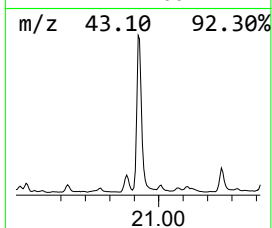
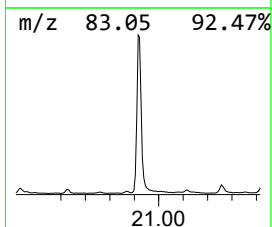
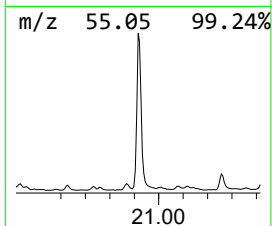
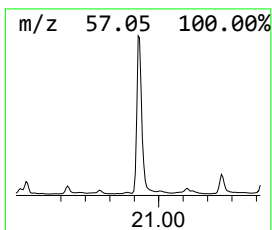
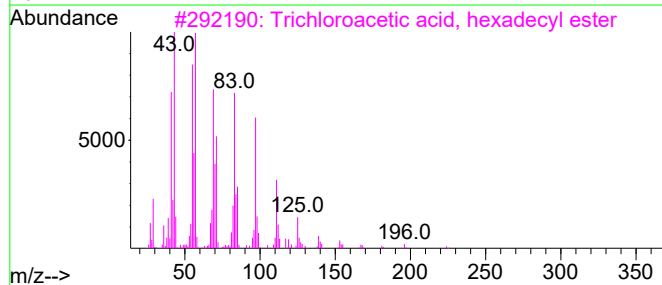
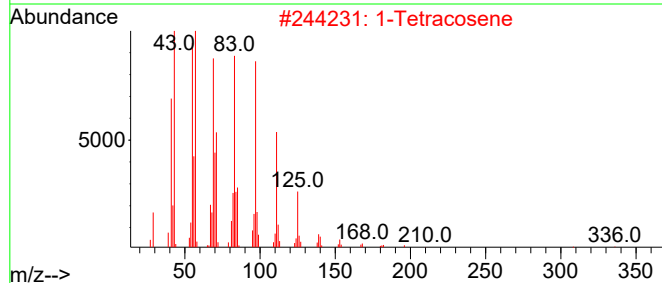
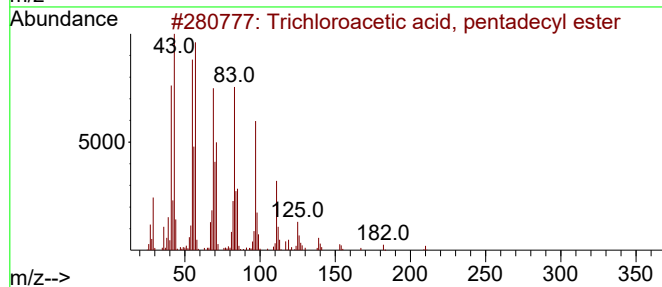
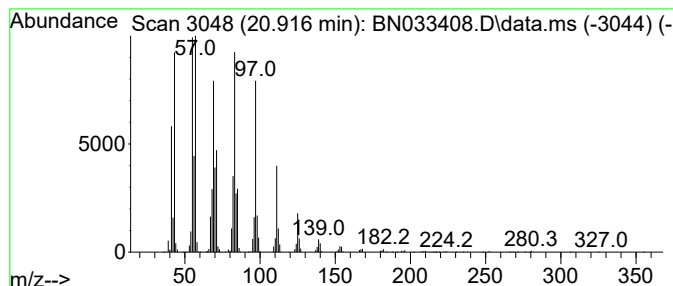
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Trichloroacetic acid, penta... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.916	26.60 ng/ul	7081230	Chrysene-d12	21.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	94
2			1-Tetracosene	336	C24H48	010192-32-2	94
3			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
4			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5			1-Heneicosanol	312	C21H44O	015594-90-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
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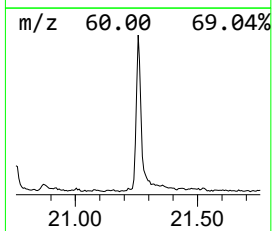
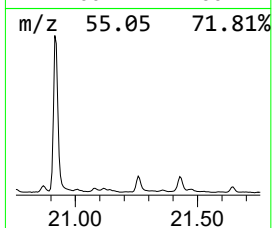
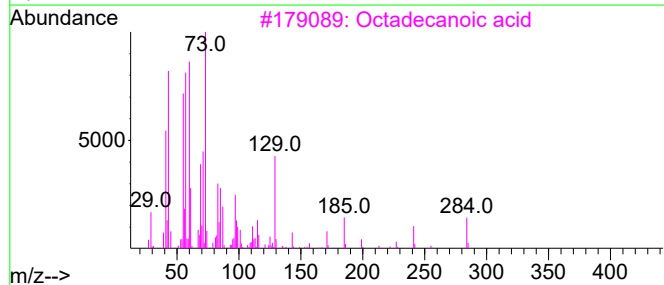
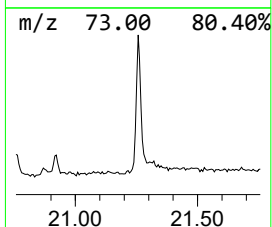
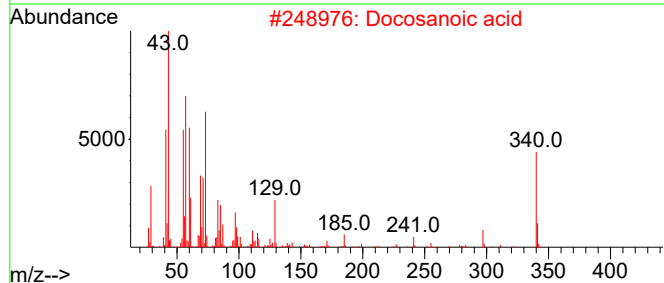
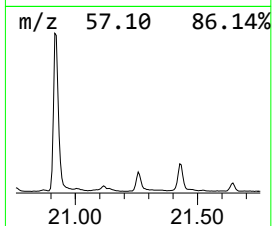
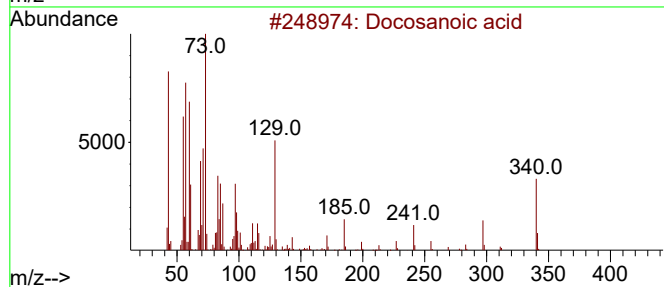
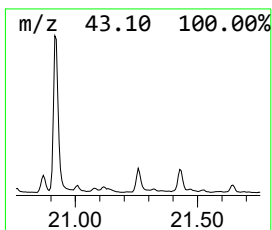
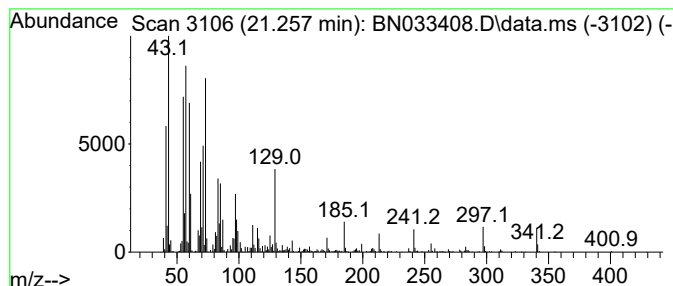
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Docosanoic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.257	3.61 ng/ul	960090	Chrysene-d12	21.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosanoic acid	340	C22H44O2	000112-85-6	99
2			Docosanoic acid	340	C22H44O2	000112-85-6	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	93
4			Nonadecanoic acid	298	C19H38O2	000646-30-0	93
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
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Sample : P3502-16
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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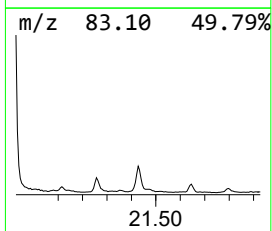
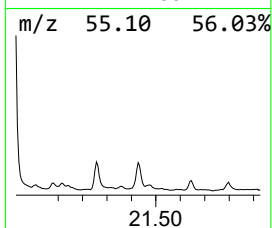
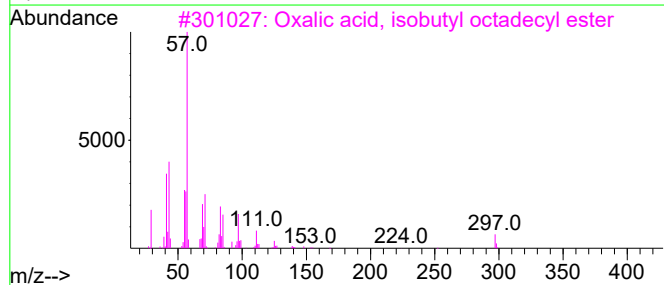
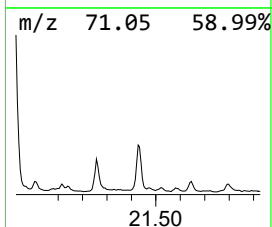
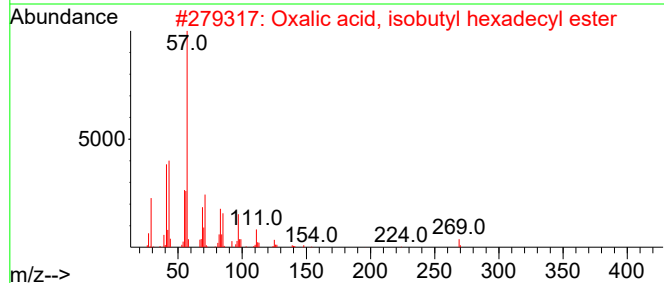
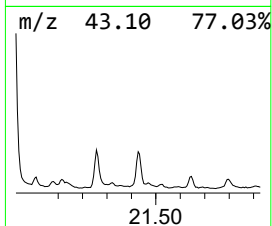
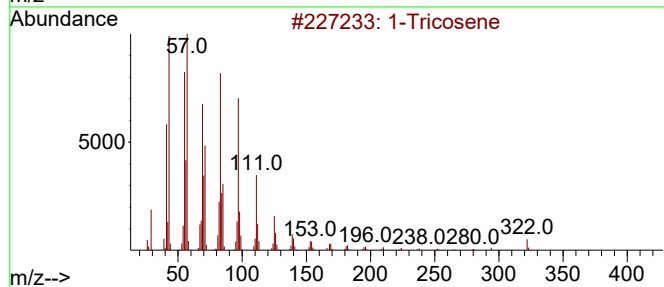
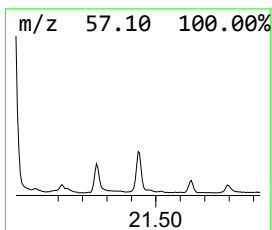
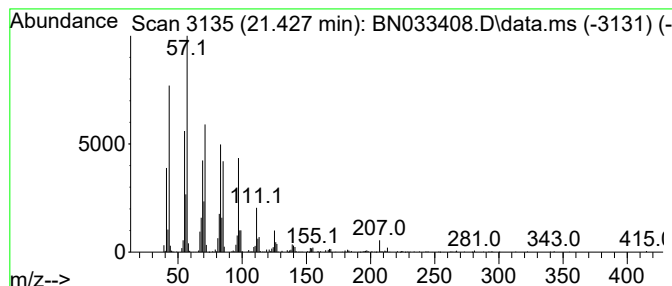
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.427	3.11 ng/ul	828128	Chrysene-d12	21.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Tricosene	322	C23H46	018835-32-0	97
2			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
3			Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	91
4			9-Tricosene, (Z)-	322	C23H46	027519-02-4	89
5			Tetracosyl heptafluorobutyrate	550	C28H49F7O2	1000351-83-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
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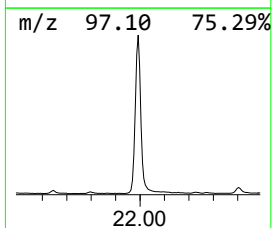
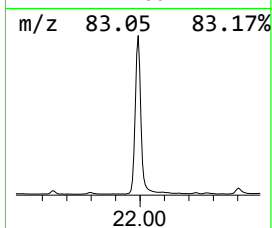
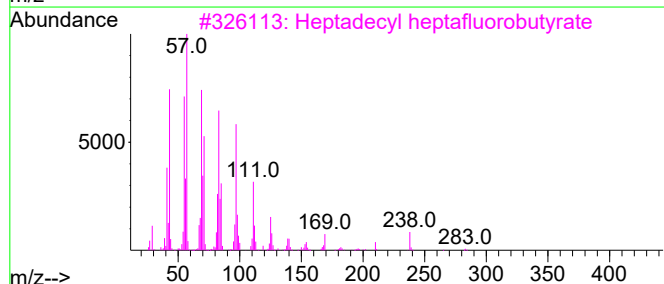
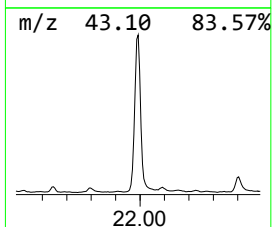
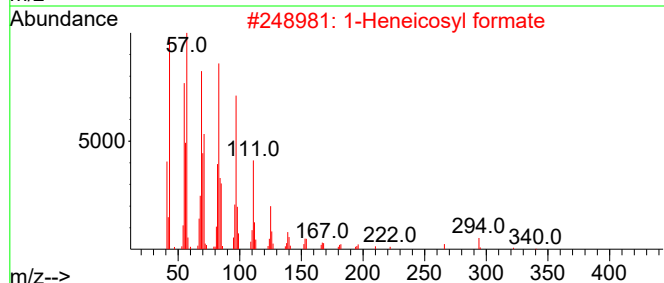
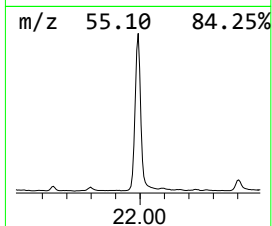
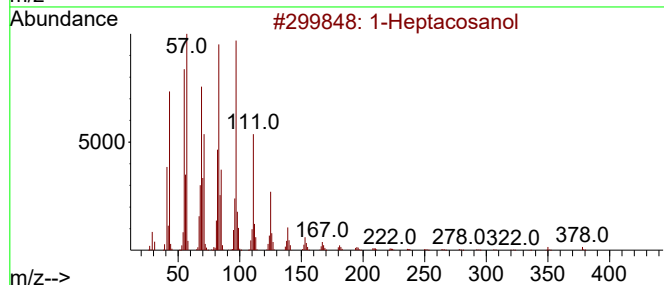
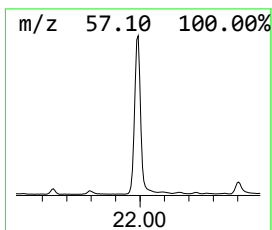
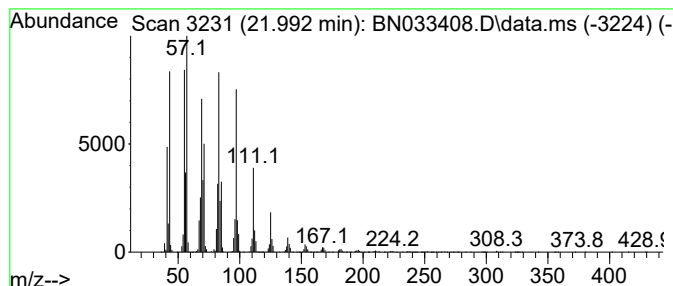
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 1-Heptacosanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.992	44.14 ng/ul	11749600	Chrysene-d12	21.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heptacosanol	396	C27H56O	002004-39-9	95
2			1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
3			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
4			1-Hexacosene	364	C26H52	018835-33-1	93
5			Behenic alcohol	326	C22H46O	000661-19-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
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Sample : P3502-16
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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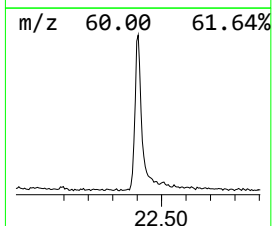
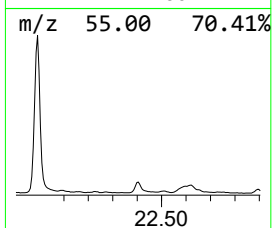
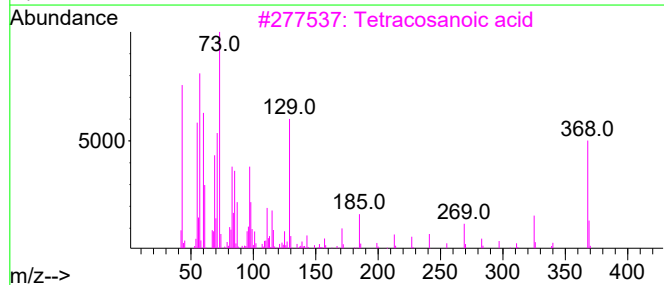
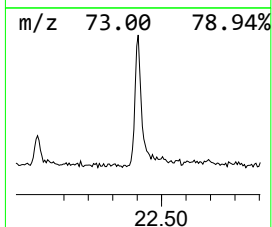
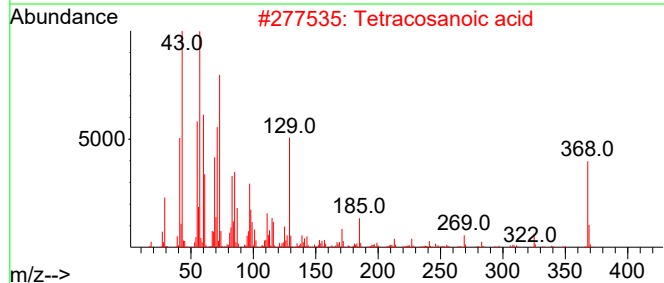
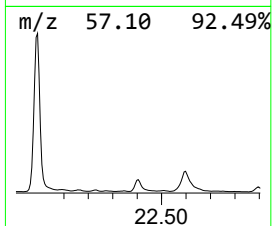
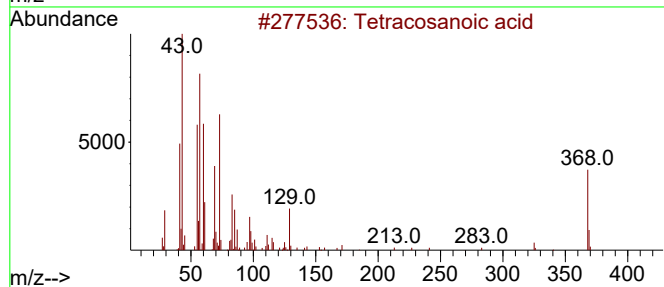
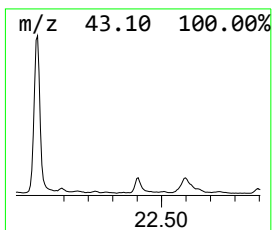
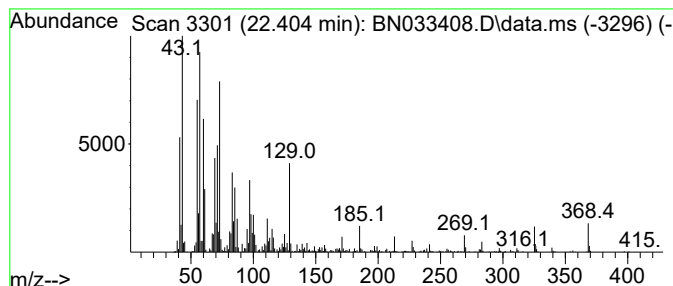
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Tetracosanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.404	4.59 ng/ul	1221340	Chrysene-d12	21.186

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	96
4			Heneicosanoic acid	326	C21H42O2	002363-71-5	89
5			Octadecanoic acid	284	C18H36O2	000057-11-4	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
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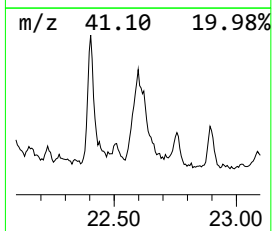
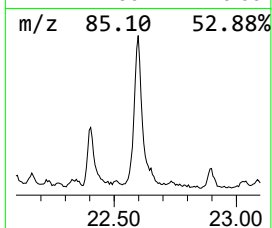
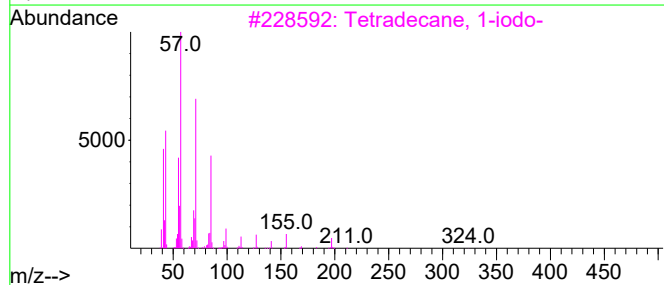
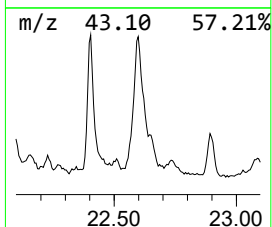
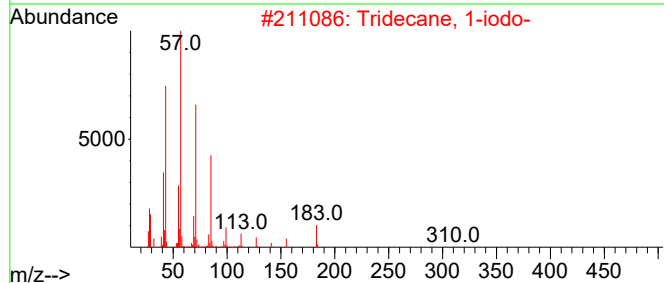
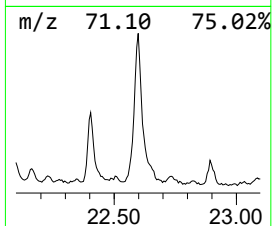
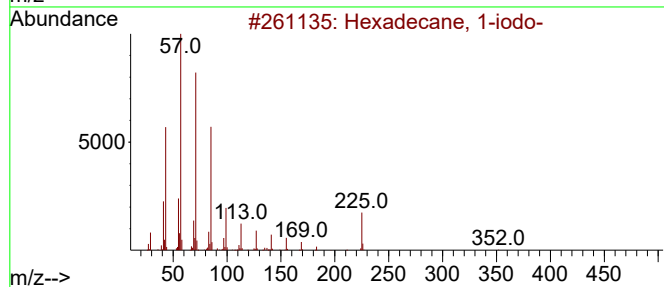
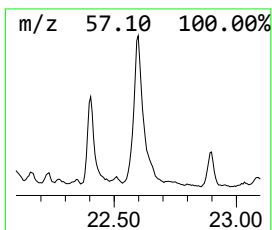
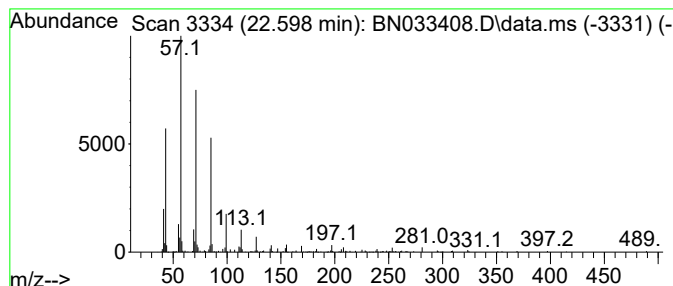
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Hexadecane, 1-iodo- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.598	5.31 ng/ul	1414180	Chrysene-d12	21.186	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	90
2	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
3	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	86
4	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	86
5	Tridecane	184	C13H28	000629-50-5	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
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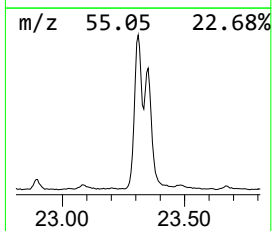
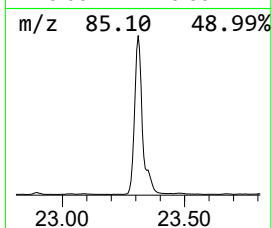
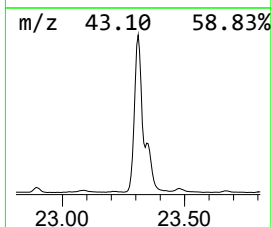
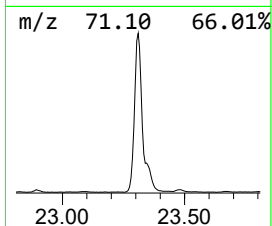
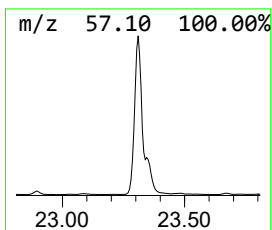
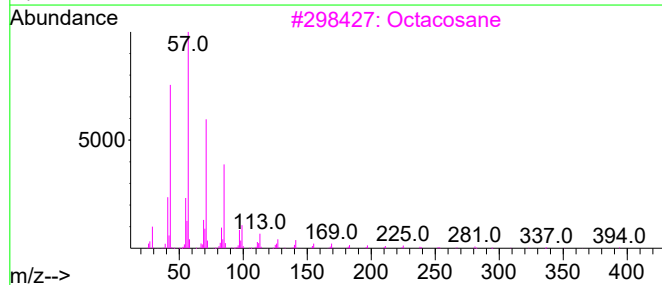
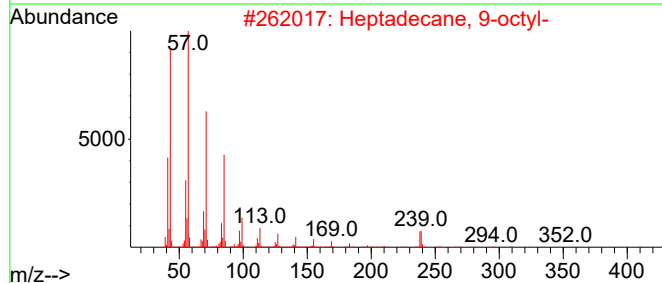
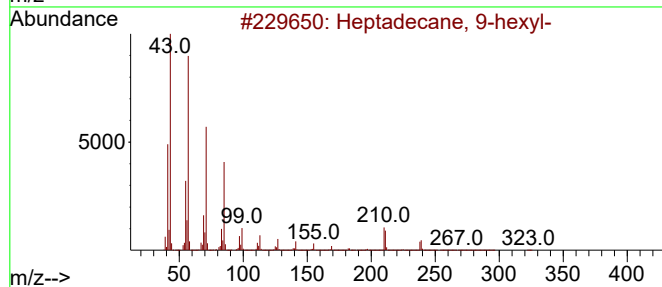
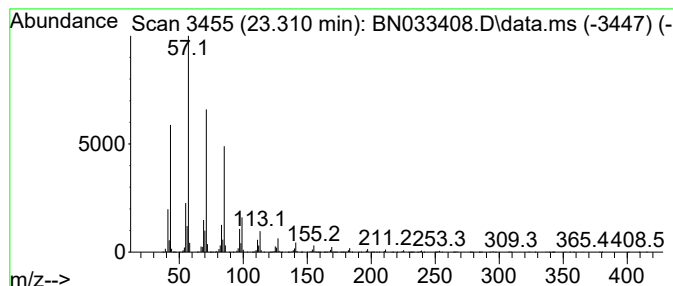
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.310	37.88 ng/ul	9495450	Perylene-d12	24.015

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
Operator : MA/JU
Sample : P3502-16
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DCXU9

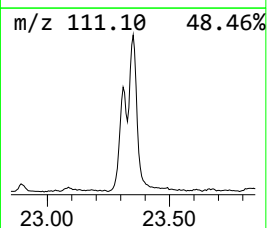
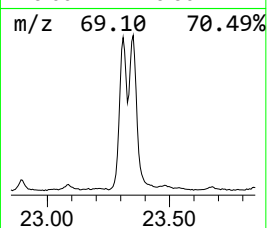
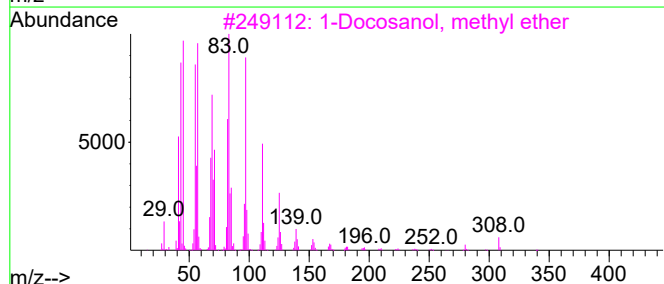
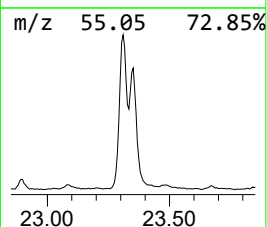
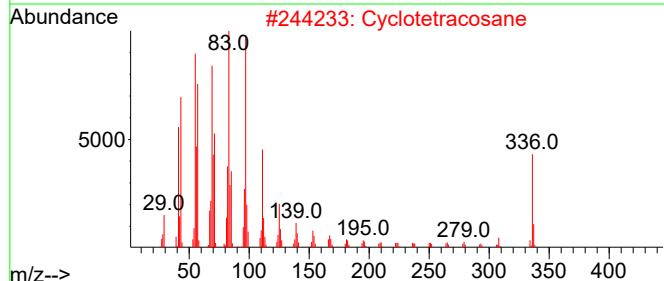
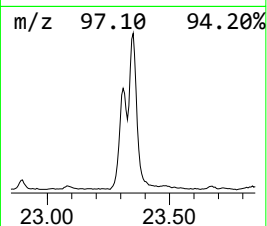
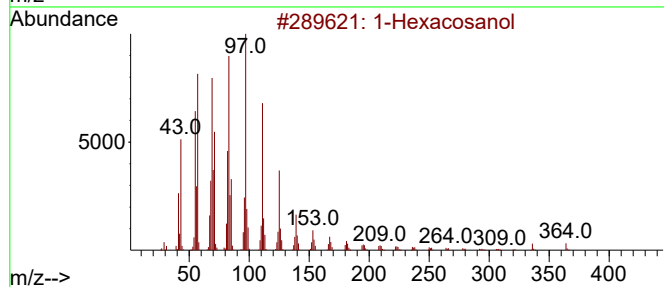
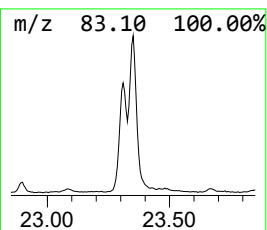
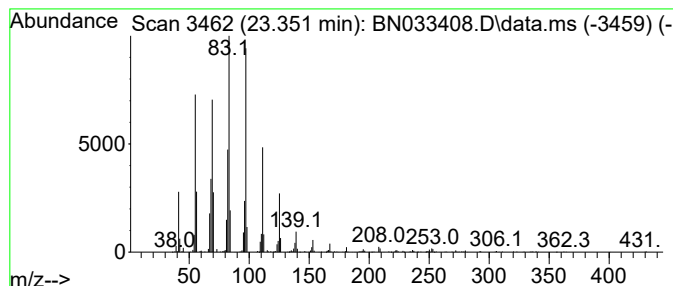
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.351	15.27 ng/ul	3827700	Perylene-d12	24.015

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosanol			382	C26H54O	000506-52-5	49
2	Cyclotetracosane			336	C24H48	000297-03-0	49
3	1-Docosanol, methyl ether			340	C23H48O	1000333-91-9	47
4	Triallylmethylsilane			166	C10H18Si	001112-91-0	43
5	1,7-Dimethyl-4-(1-methylethyl)cy...			210	C15H30	000645-10-3	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
Operator : MA/JU
Sample : P3502-16
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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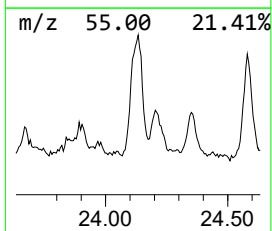
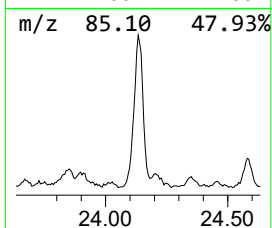
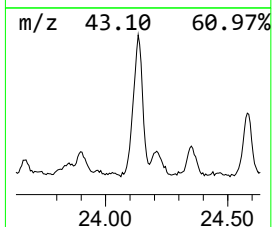
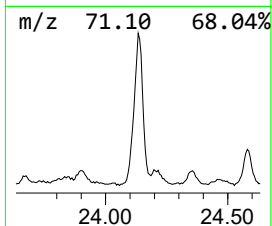
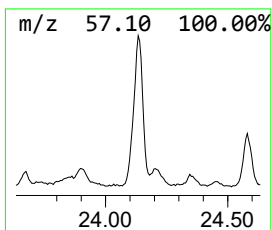
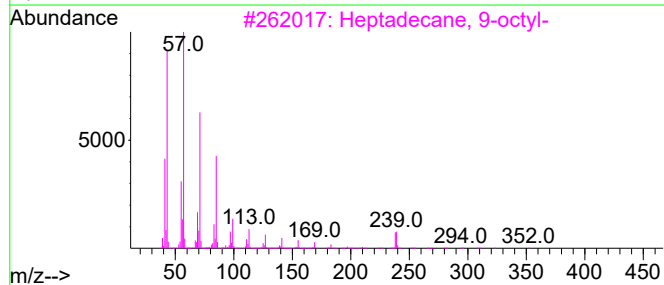
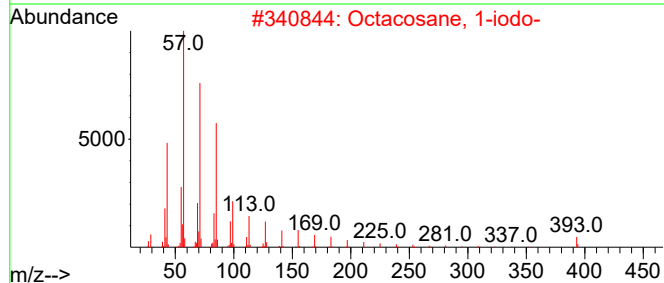
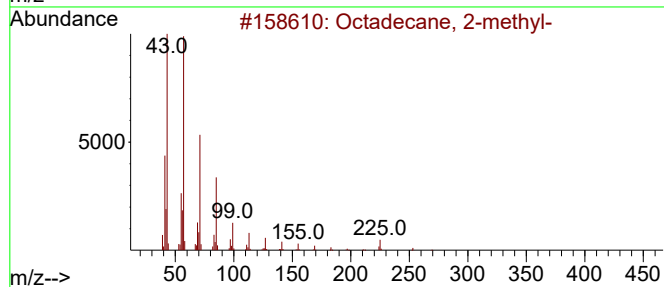
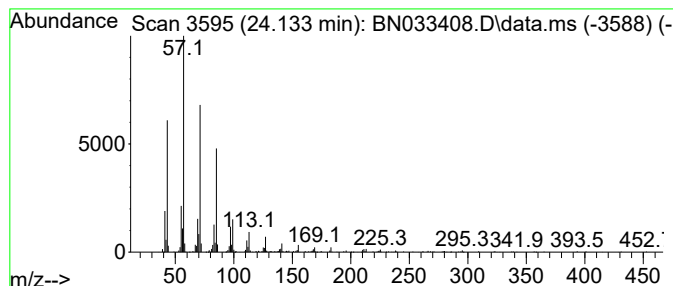
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.133	4.26 ng/ul	1068140	Perylene-d12	24.015

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane, 2-methyl-	268	C19H40	001560-88-9	93
2		Octacosane, 1-iodo-	520	C28H57I	1000406-32-2	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Triacontane, 1-iodo-	548	C30H61I	1000406-32-3	91
5		Hexacosane, 1-iodo-	492	C26H53I	1000406-32-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
Operator : MA/JU
Sample : P3502-16
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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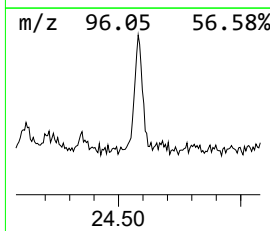
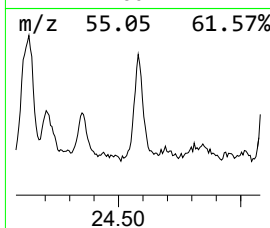
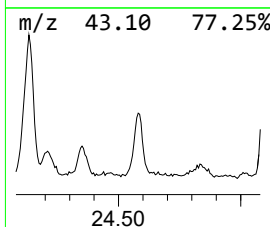
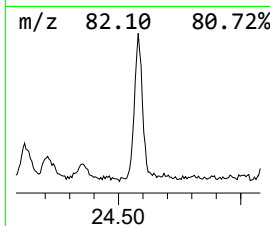
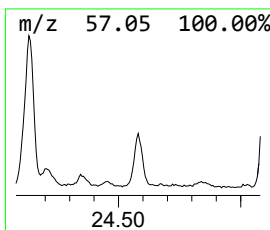
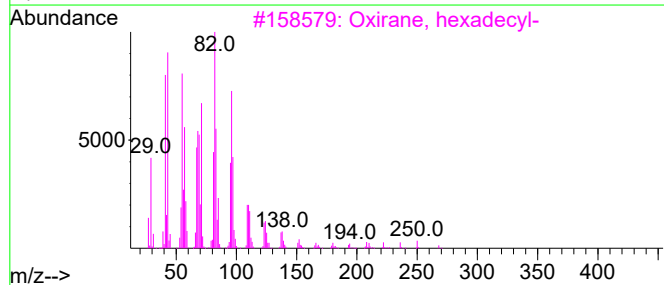
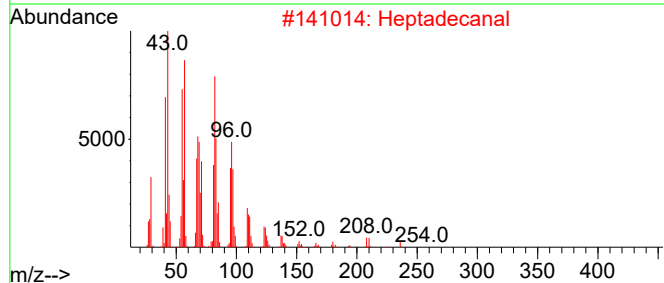
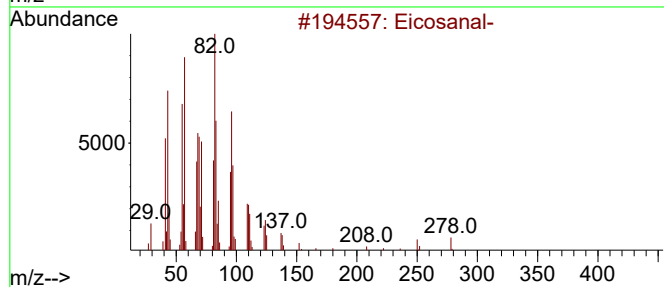
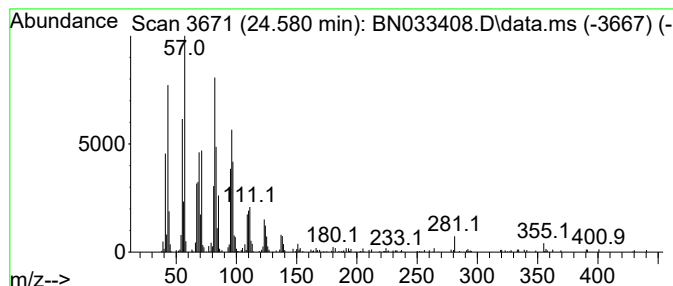
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Eicosanal- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.580	2.77 ng/ul	693331	Perylene-d12	24.015

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosanal-	296	C20H40O	002400-66-0	91
2			Heptadecanal	254	C17H34O	1000376-70-0	91
3			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
4			Oxirane, heptadecyl-	282	C19H38O	067860-04-2	87
5			Tetracosanal	352	C24H48O	057866-08-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
Operator : MA/JU
Sample : P3502-16
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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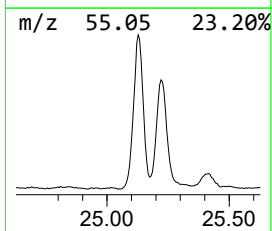
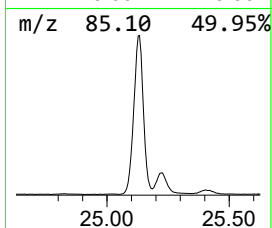
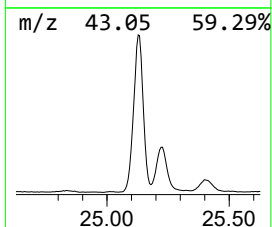
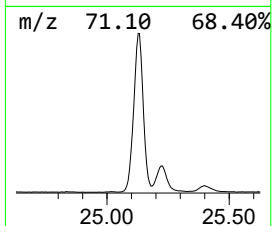
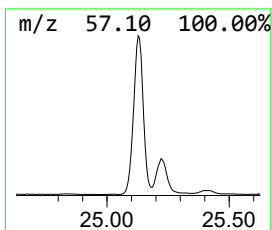
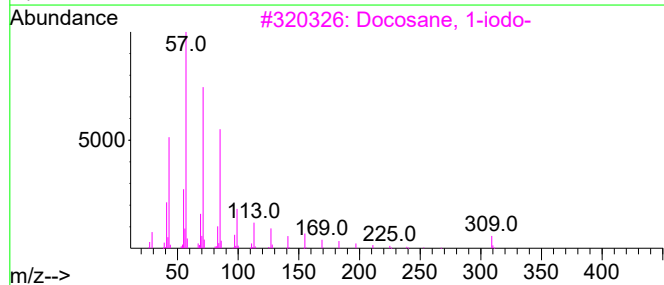
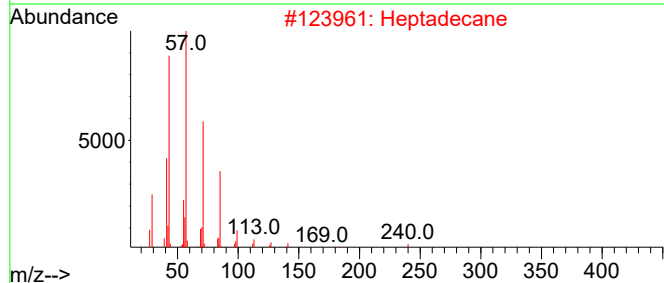
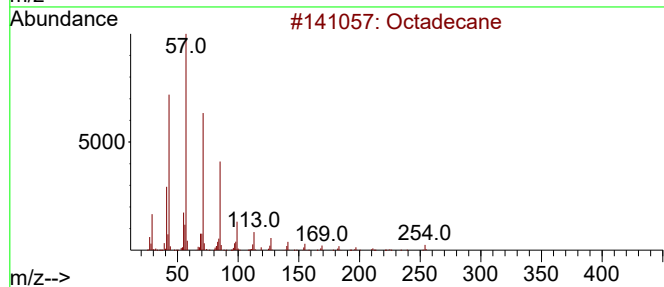
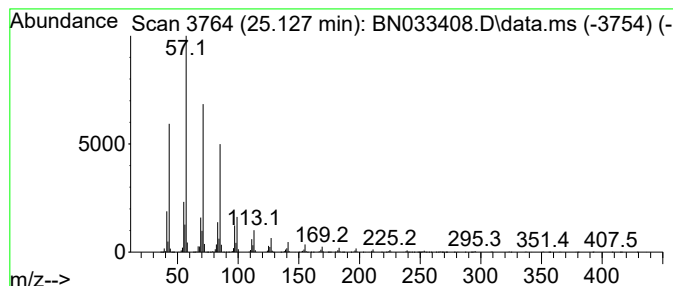
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.127	41.00 ng/ul	10278900	Perylene-d12	24.015

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane	254	C18H38	000593-45-3	95
2		Heptadecane	240	C17H36	000629-78-7	93
3		Docosane, 1-iodo-	436	C22H45I	1000406-31-9	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
Operator : MA/JU
Sample : P3502-16
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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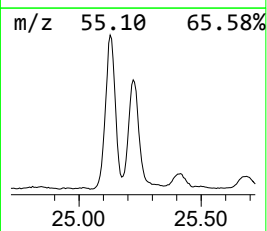
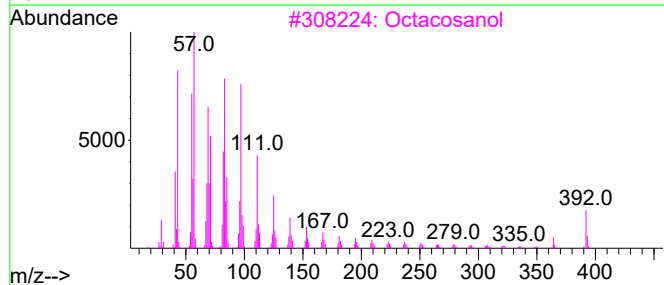
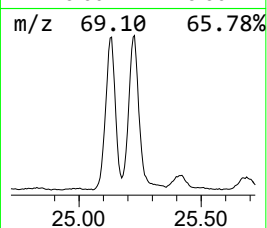
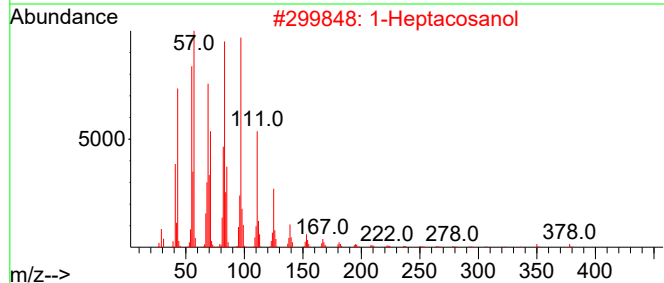
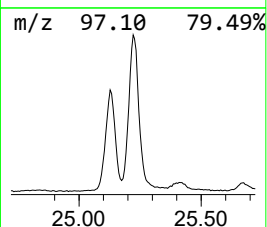
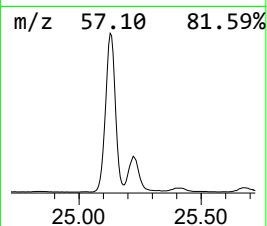
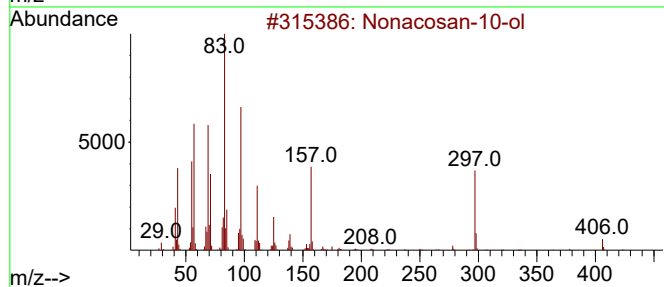
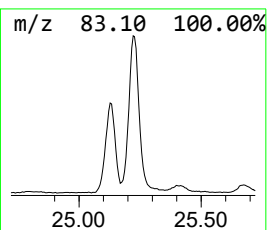
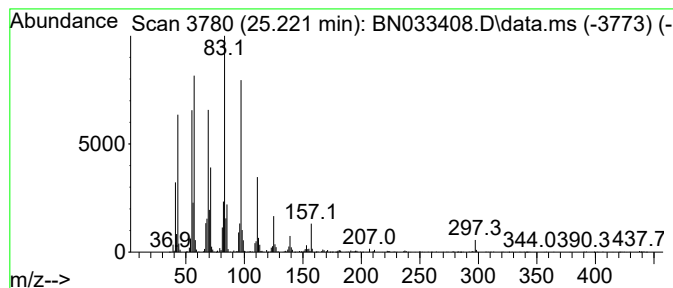
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Nonacosan-10-ol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.221	19.75 ng/ul	4952370	Perylene-d12	24.015

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonacosan-10-ol	424	C29H60O	000504-55-2	87
2			1-Heptacosanol	396	C27H56O	002004-39-9	64
3			Octacosanol	410	C28H58O	000557-61-9	64
4			Octacosyl acetate	452	C30H60O2	018206-97-8	64
5			Triacetyl acetate	480	C32H64O2	041755-58-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
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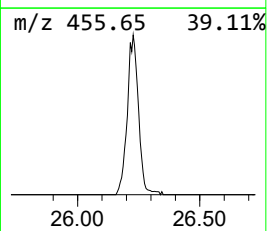
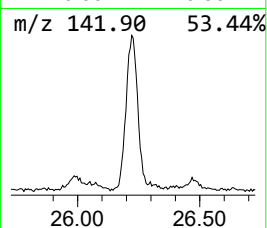
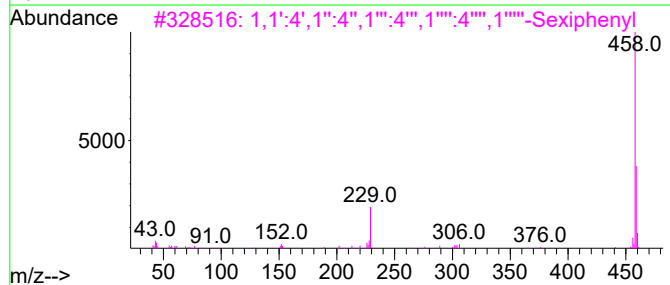
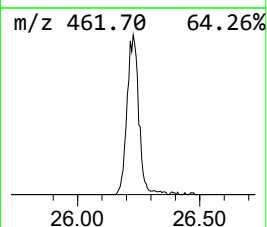
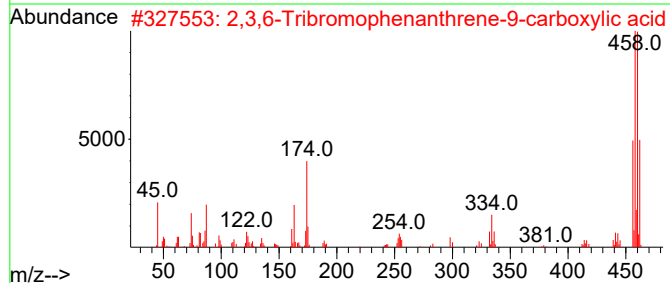
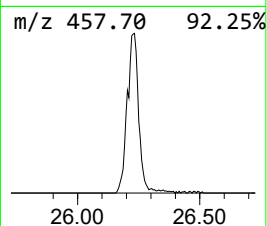
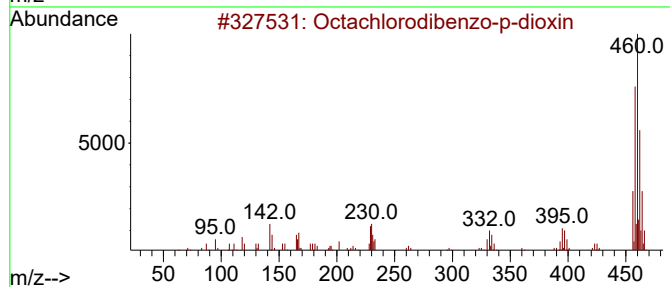
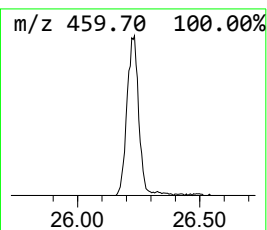
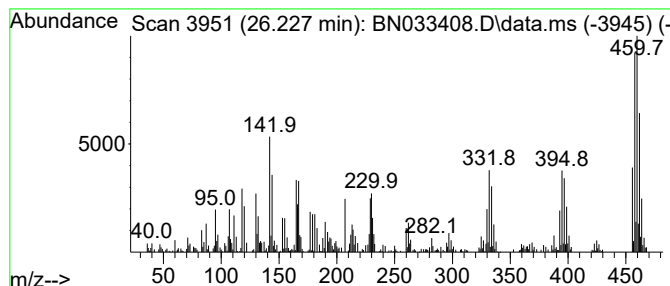
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Octachlorodibenzo-p-dioxin Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.	
26.227	2.49 ng/ul	623344	Perylene-d12	24.015	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octachlorodibenzo-p-dioxin	456	C12Cl8O2	003268-87-9	96
2	2,3,6-Tribromophenanthrene-9-car...	456	C15H7Br3O2	056988-60-4	46
3	1,1':4',1'':4'',1''':4''',1'''':4'''':...	458	C36H26	004499-83-6	10
4	1H-Isoindole-1,3(2H)-dione, 5,5'...	460	C28H16N2O5	033734-37-1	9
5	Silane, dimethyl(4-bromophenoxy)...	456	C23H41BrO2Si	1000347-11-7	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
Operator : MA/JU
Sample : P3502-16
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ALS Vial : 19 Sample Multiplier: 1

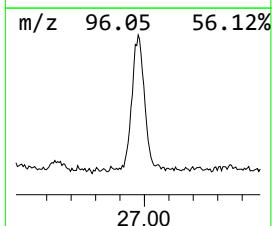
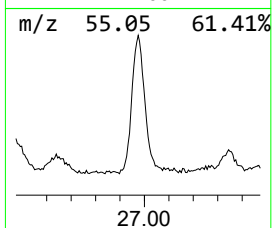
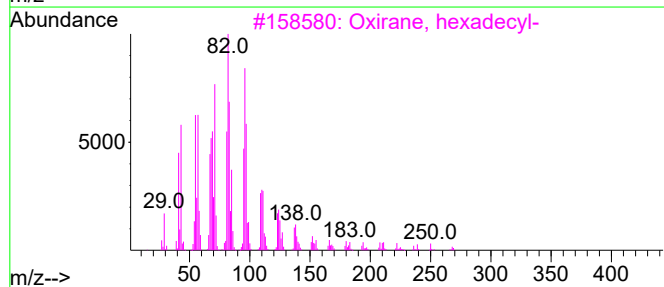
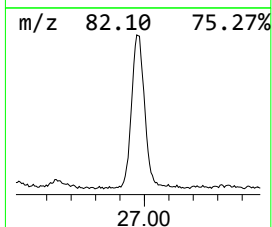
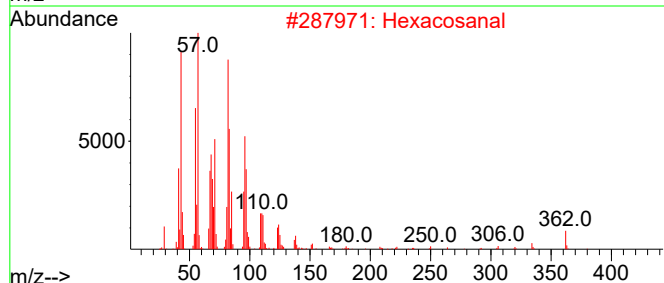
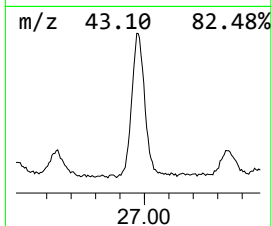
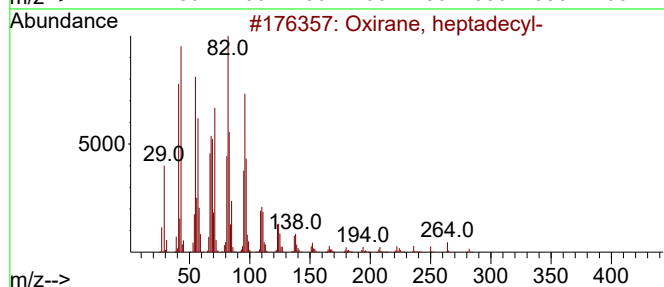
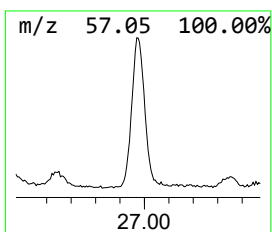
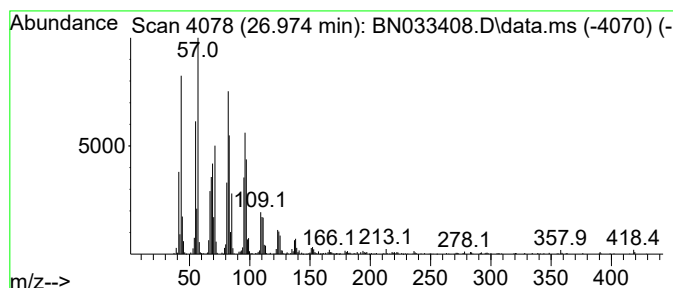
Instrument :
BNA_N
ClientSampleId :
DCXU9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 24 Oxirane, heptadecyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.	
26.974	8.96 ng/ul	2246890	Perylene-d12		24.015	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Oxirane, heptadecyl-		282	C19H38O	067860-04-2	91
2	Hexacosanal		380	C26H52O	026627-85-0	91
3	Oxirane, hexadecyl-		268	C18H36O	007390-81-0	91
4	Oxirane, hexadecyl-		268	C18H36O	007390-81-0	91
5	Oxirane, hexadecyl-		268	C18H36O	007390-81-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033408.D
Acq On : 15 Aug 2024 20:52
Operator : MA/JU
Sample : P3502-16
Misc :
ALS Vial : 19 Sample Multiplier: 1

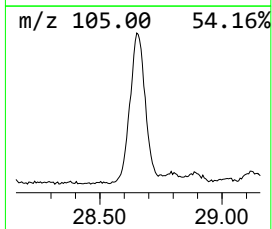
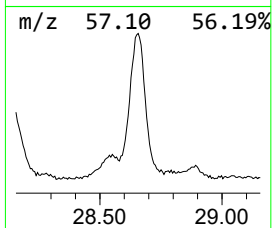
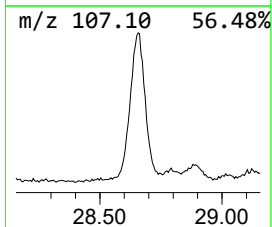
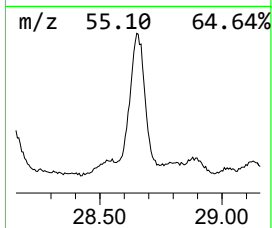
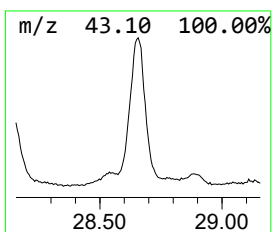
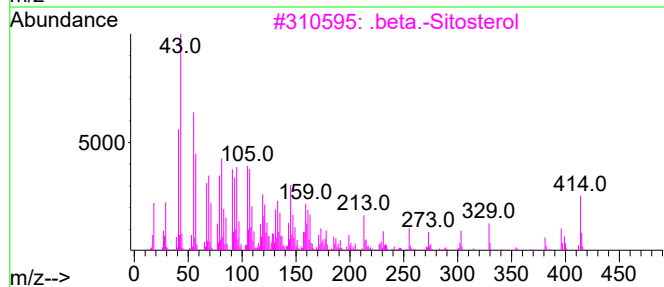
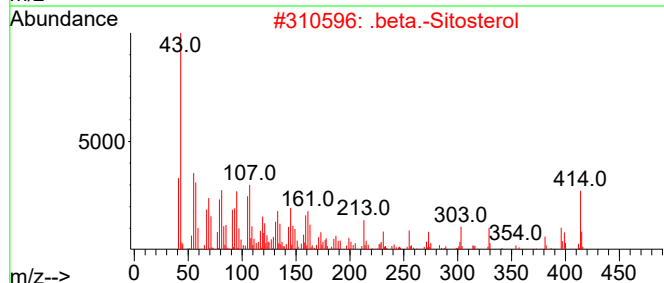
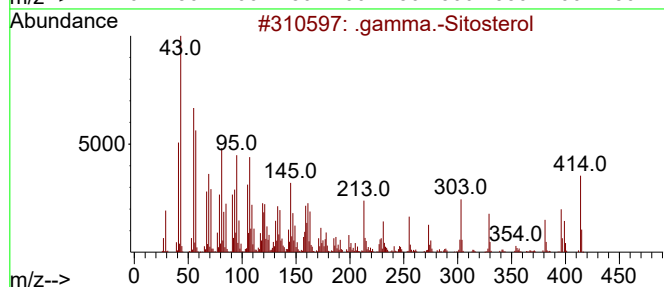
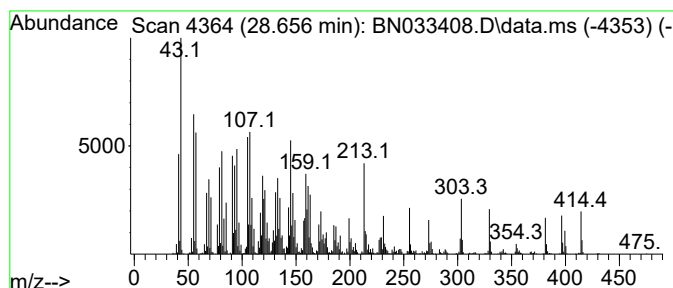
Instrument :
BNA_N
ClientSampleId :
DCXU9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 25 .gamma.-Sitosterol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
28.656	25.69 ng/ul	6441600	Perylene-d12		24.015	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	.gamma.-Sitosterol		414	C29H50O	000083-47-6	99
2	.beta.-Sitosterol		414	C29H50O	000083-46-5	95
3	.beta.-Sitosterol		414	C29H50O	000083-46-5	95
4	Pregn-5-en-3-ol, 21-bromo-20-met...		394	C22H35BrO	055103-80-5	64
5	Campesterol		400	C28H48O	000474-62-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
 Data File : BN033408.D
 Acq On : 15 Aug 2024 20:52
 Operator : MA/JU
 Sample : P3502-16
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DCXU9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propylene Glycol	3.411	2.6	ng/ul	403636	1	7.569	3143880	20.0
(1R)-2,6,6-Trim...	6.287	3.3	ng/ul	522992	1	7.569	3143880	20.0
1-Phenoxypropan...	11.081	2.9	ng/ul	674063	2	10.334	4686390	20.0
Phenol, 4-(1-me...	17.345	19.0	ng/ul	6552090	4	16.957	6908780	20.0
n-Hexadecanoic ...	17.892	9.8	ng/ul	3386860	4	16.957	6908780	20.0
Octadecanoic acid	19.192	9.3	ng/ul	2490300	5	21.186	5324060	20.0
Phenol, 4,4'-(1...	19.439	222.4	ng/ul	59212800	5	21.186	5324060	20.0
1-Heneicosanol	19.922	2.6	ng/ul	704465	5	21.186	5324060	20.0
Cinnamyl cinnamate	20.733	13.3	ng/ul	3548290	5	21.186	5324060	20.0
Dehydroabiatic ...	20.869	9.3	ng/ul	2481290	5	21.186	5324060	20.0
Trichloroacetic...	20.916	26.6	ng/ul	7081230	5	21.186	5324060	20.0
Docosanoic acid	21.257	3.6	ng/ul	960090	5	21.186	5324060	20.0
(DEL) Alkane: S...	21.427	3.1	ng/ul	828128	5	21.186	5324060	20.0
1-Heptacosanol	21.992	44.1	ng/ul	11749600	5	21.186	5324060	20.0
Tetracosanoic acid	22.404	4.6	ng/ul	1221340	5	21.186	5324060	20.0
Hexadecane, 1-i...	22.598	5.3	ng/ul	1414180	5	21.186	5324060	20.0
(DEL) Alkane: S...	23.310	37.9	ng/ul	9495450	6	24.015	5014040	20.0
unknown-01	23.351	15.3	ng/ul	3827700	6	24.015	5014040	20.0
(DEL) Alkane: S...	24.133	4.3	ng/ul	1068140	6	24.015	5014040	20.0
Eicosanal-	24.580	2.8	ng/ul	693331	6	24.015	5014040	20.0
(DEL) Alkane: S...	25.127	41.0	ng/ul	10278900	6	24.015	5014040	20.0
Nonacosan-10-ol	25.221	19.8	ng/ul	4952370	6	24.015	5014040	20.0
Octachlorodiben...	26.227	2.5	ng/ul	623344	6	24.015	5014040	20.0
Oxirane, heptad...	26.974	9.0	ng/ul	2246890	6	24.015	5014040	20.0
.gamma.-Sitosterol	28.656	25.7	ng/ul	6441600	6	24.015	5014040	20.0