

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
 Data File : BM036586.D
 Acq On : 19 Sep 2022 18:47
 Operator : CG/JU
 Sample : N4604-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A5767

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-BM091522.M
 Title : SVOA CALIBRATION

Signal : TIC: BM036586.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.128	3	7	31	rVB	14319463	20615961	100.00%	14.927%
2	3.393	49	52	61	rVB	95517	137603	0.67%	0.100%
3	3.828	124	126	139	rBV2	87091	148345	0.72%	0.107%
4	3.934	140	144	153	rVV	91723	130743	0.63%	0.095%
5	4.069	161	167	175	rVV	87444	146072	0.71%	0.106%
6	4.163	178	183	193	rVB	51510	81885	0.40%	0.059%
7	5.104	337	343	354	rBV	617298	880522	4.27%	0.638%
8	7.175	688	695	704	rVV	1522041	2466461	11.96%	1.786%
9	7.351	719	725	732	rVV	1345544	2078907	10.08%	1.505%
10	7.551	749	759	768	rBV	1525747	2391702	11.60%	1.732%
11	8.028	832	840	847	rBV	1697942	2707979	13.14%	1.961%
12	8.728	950	959	968	rVB	1361061	2128649	10.33%	1.541%
13	8.857	974	981	987	rBV	43593	86783	0.42%	0.063%
14	9.192	1031	1038	1047	rVV	1527138	2469134	11.98%	1.788%
15	9.269	1047	1051	1061	rVB3	38531	84368	0.41%	0.061%
16	9.916	1153	1161	1168	rBV	1132819	1803771	8.75%	1.306%
17	10.451	1244	1252	1263	rBV	1938141	3050748	14.80%	2.209%
18	10.616	1274	1280	1286	rBV	123240	207964	1.01%	0.151%
19	10.839	1308	1318	1324	rBV	2498426	4248066	20.61%	3.076%
20	10.892	1324	1327	1333	rVV	66148	101609	0.49%	0.074%
21	10.969	1333	1340	1354	rVV	781708	1441647	6.99%	1.044%
22	11.869	1488	1493	1499	rBV	85740	125481	0.61%	0.091%
23	12.257	1554	1559	1566	rBV	69455	112317	0.54%	0.081%
24	12.416	1578	1586	1591	rBV	40611	76342	0.37%	0.055%
25	12.492	1593	1599	1606	rVB	122241	200701	0.97%	0.145%
26	12.710	1630	1636	1643	rVV	104788	162464	0.79%	0.118%
27	13.204	1716	1720	1725	rVB	61418	87829	0.43%	0.064%
28	13.492	1761	1769	1775	rBV	142027	217242	1.05%	0.157%
29	13.975	1846	1851	1856	rVV	139098	232662	1.13%	0.168%
30	14.063	1856	1866	1875	rVV	3710899	5113762	24.80%	3.703%
31	14.139	1875	1879	1883	rVV2	99863	143309	0.70%	0.104%
32	14.210	1887	1891	1895	rVV	67285	107965	0.52%	0.078%
33	14.345	1908	1914	1926	rVB	4195937	6080432	29.49%	4.403%
34	14.551	1945	1949	1955	rVB	110020	146536	0.71%	0.106%
35	14.651	1956	1966	1973	rBV	4014694	5827450	28.27%	4.219%
36	14.827	1989	1996	2002	rBV	1360844	1990689	9.66%	1.441%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Title : SVOA CALIBRATION

37	14.986	2019	2023	2029	rVV4	44202	87351	0.42%	0.063%
38	15.051	2029	2034	2040	rVV3	140411	247868	1.20%	0.179%
39	15.122	2040	2046	2050	rVV	61625	96447	0.47%	0.070%
40	15.263	2063	2070	2075	rBV2	64514	122898	0.60%	0.089%

41	15.316	2075	2079	2083	rVB	60890	86072	0.42%	0.062%
42	15.439	2093	2100	2107	rBV5	86152	196434	0.95%	0.142%
43	15.498	2107	2110	2116	rVB	102393	126017	0.61%	0.091%
44	15.639	2126	2134	2139	rVV	5321039	7634069	37.03%	5.528%
45	15.680	2139	2141	2145	rVB4	65784	80143	0.39%	0.058%

46	15.745	2145	2152	2160	rBV	1196189	1621167	7.86%	1.174%
47	15.904	2176	2179	2184	rVB	72964	100576	0.49%	0.073%
48	15.980	2187	2192	2202	rVB4	72429	153523	0.74%	0.111%
49	16.133	2214	2218	2224	rVB3	61357	130968	0.64%	0.095%
50	16.216	2226	2232	2238	rVB4	44702	93736	0.45%	0.068%

51	16.363	2252	2257	2258	rBV	145818	186383	0.90%	0.135%
52	16.380	2258	2260	2265	rVB	201163	259827	1.26%	0.188%
53	16.698	2305	2314	2318	rBV6	39606	94703	0.46%	0.069%
54	16.927	2348	2353	2356	rBV3	97850	148538	0.72%	0.108%
55	17.039	2368	2372	2379	rBV2	63199	128724	0.62%	0.093%

56	17.121	2381	2386	2387	rVV2	46101	75341	0.37%	0.055%
57	17.151	2387	2391	2397	rVV2	122626	205947	1.00%	0.149%
58	17.204	2397	2400	2409	rVB5	76122	120738	0.59%	0.087%
59	17.386	2421	2431	2436	rBV	4791311	6669649	32.35%	4.829%
60	17.427	2436	2438	2442	rVV	340411	392823	1.91%	0.284%

61	17.486	2442	2448	2458	rVV	5499754	7556745	36.65%	5.472%
62	17.568	2458	2462	2468	rVV4	65471	121527	0.59%	0.088%
63	17.698	2481	2484	2491	rVB	76461	97711	0.47%	0.071%
64	17.886	2513	2516	2521	rVB	95727	118506	0.57%	0.086%
65	17.962	2526	2529	2533	rVB2	77070	89842	0.44%	0.065%

66	18.221	2569	2573	2576	rBV3	76556	105905	0.51%	0.077%
67	18.280	2576	2583	2586	rBV2	480248	882907	4.28%	0.639%
68	18.445	2606	2611	2615	rBV2	210080	314118	1.52%	0.227%
69	18.486	2615	2618	2624	rVB	142201	203486	0.99%	0.147%
70	18.580	2629	2634	2638	rBV	171897	263891	1.28%	0.191%

71	18.757	2660	2664	2668	rBV	144193	207783	1.01%	0.150%
72	19.051	2711	2714	2717	rVV	75783	102108	0.50%	0.074%
73	19.092	2718	2721	2724	rVB	68552	81141	0.39%	0.059%
74	19.168	2729	2734	2738	rVV2	322795	507882	2.46%	0.368%
75	19.209	2738	2741	2747	rVV2	182568	308777	1.50%	0.224%

76	19.262	2747	2750	2754	rVB	101823	116725	0.57%	0.085%
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77	19.327	2754	2761	2766	rBV4	112912	273917	1.33%	0.198%
78	19.433	2775	2779	2786	rVB	578829	782009	3.79%	0.566%
79	19.551	2795	2799	2802	rBV2	178342	218303	1.06%	0.158%
80	19.768	2830	2836	2848	rBV	6819945	10073121	48.86%	7.294%
81	19.868	2849	2853	2858	rVB2	121350	190055	0.92%	0.138%
82	20.109	2891	2894	2904	rBV4	85188	184028	0.89%	0.133%
83	20.280	2920	2923	2926	rVV2	122456	140669	0.68%	0.102%
84	20.315	2926	2929	2933	rVB	235258	244432	1.19%	0.177%
85	20.768	3002	3006	3009	rBV	195753	234616	1.14%	0.170%
86	20.809	3010	3013	3016	rVB	208602	211928	1.03%	0.153%
87	21.027	3047	3050	3056	rVB	171211	244486	1.19%	0.177%
88	21.262	3086	3090	3096	rBV2	1078807	1448670	7.03%	1.049%
89	21.550	3133	3139	3143	rBV	4964705	6718091	32.59%	4.864%
90	21.715	3164	3167	3171	rBV	307284	408570	1.98%	0.296%
91	22.198	3244	3249	3257	rBV2	727373	1327273	6.44%	0.961%
92	22.703	3331	3335	3346	rVB	248561	459564	2.23%	0.333%
93	22.956	3373	3378	3386	rVB	515952	804622	3.90%	0.583%
94	23.274	3428	3432	3435	rBV	484700	723826	3.51%	0.524%
95	23.315	3435	3439	3453	rVB	1055423	1875256	9.10%	1.358%
96	23.839	3521	3528	3537	rBV2	3175835	6021340	29.21%	4.360%
97	23.997	3548	3555	3564	rVB2	2351965	4760819	23.09%	3.447%
98	24.274	3598	3602	3610	rVB2	288165	586178	2.84%	0.424%
99	24.668	3663	3669	3677	rBV	670222	1429492	6.93%	1.035%
100	24.756	3681	3684	3696	rVB2	302945	573811	2.78%	0.415%

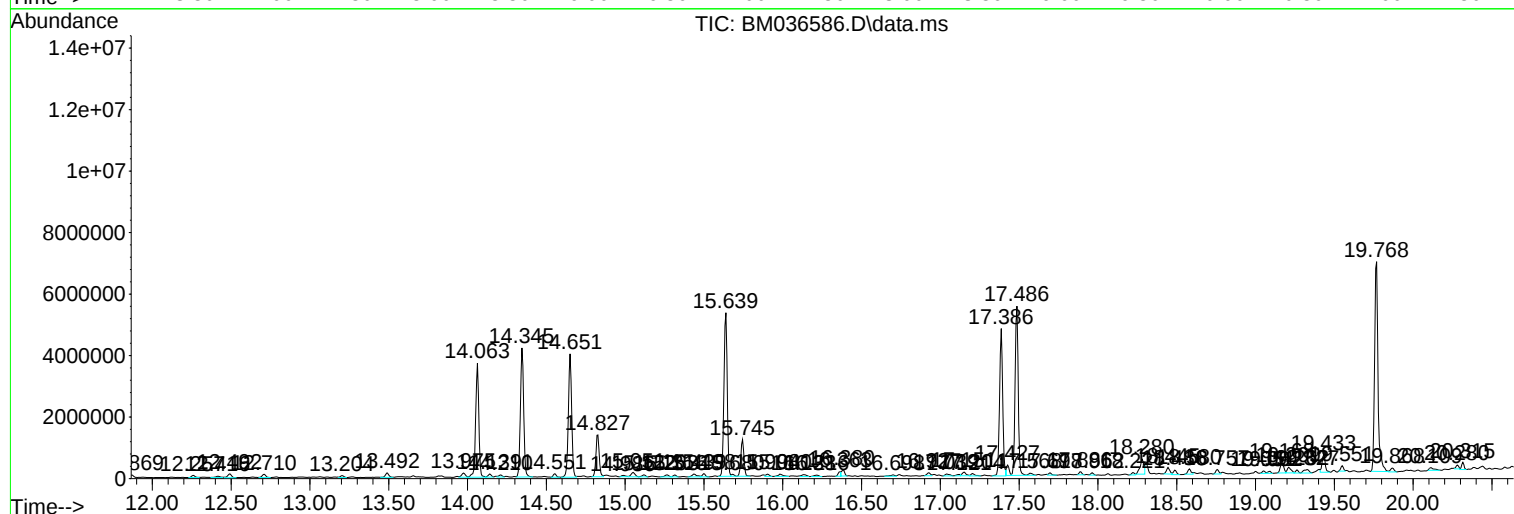
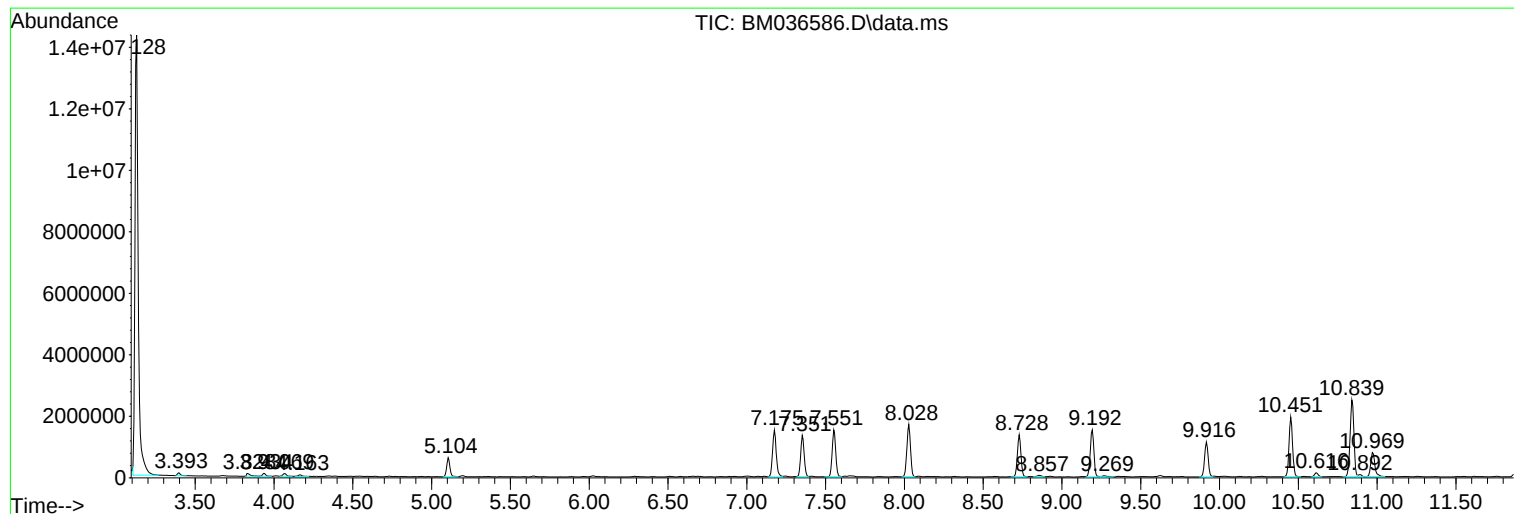
Sum of corrected areas: 138108172

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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



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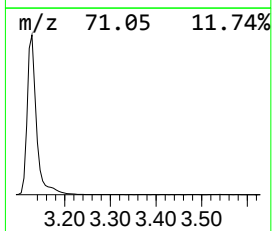
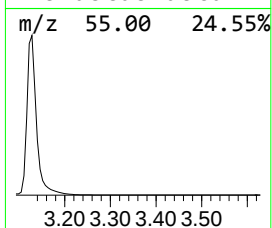
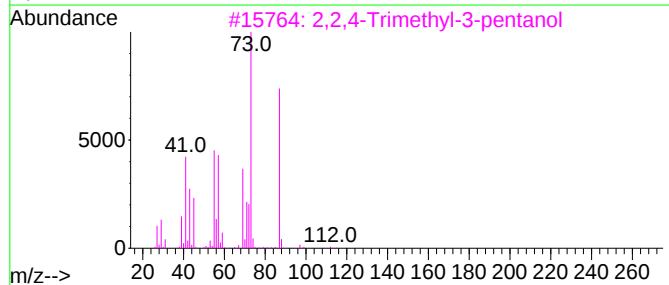
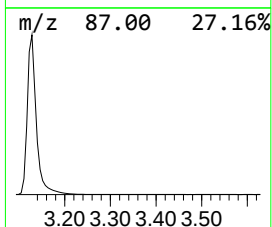
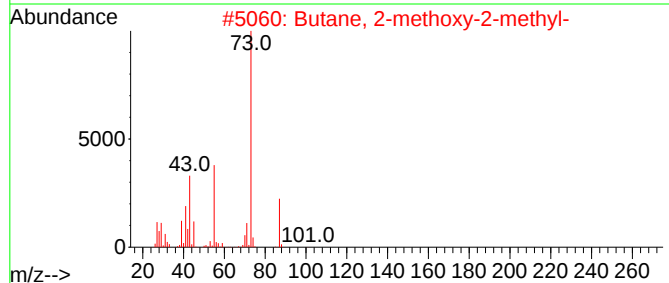
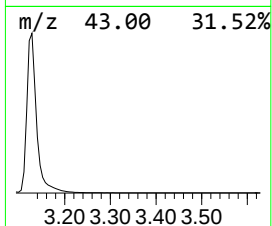
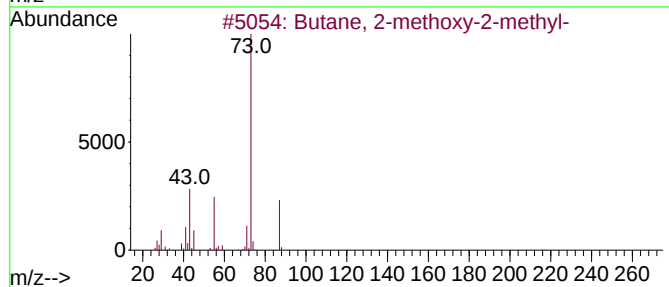
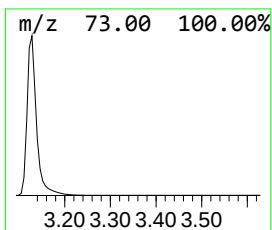
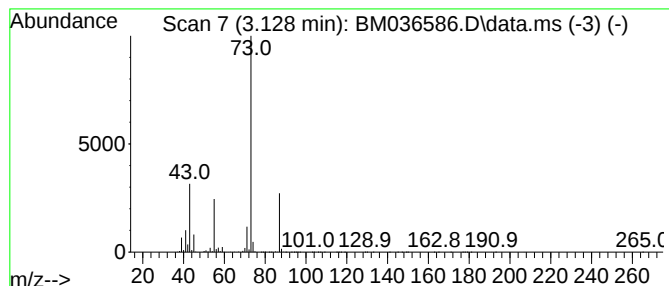
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.128	152.26 ng/ul	20616000	1,4-Dichlorobenzene-d4	8.028

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5			Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9



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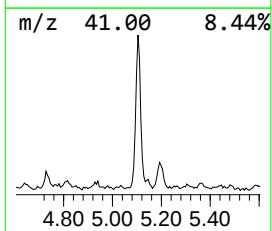
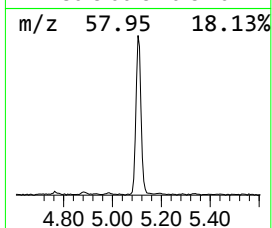
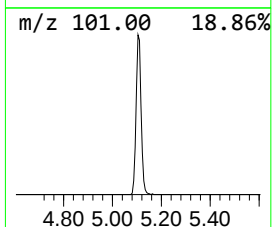
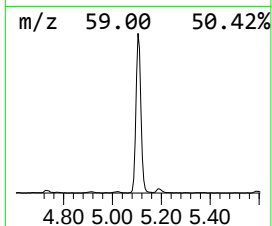
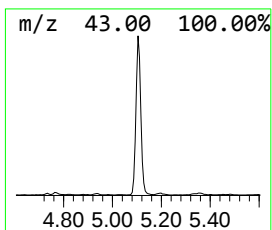
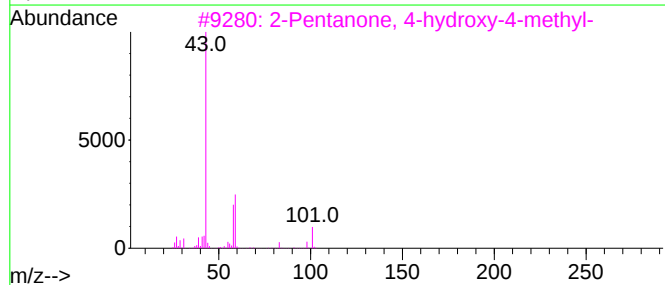
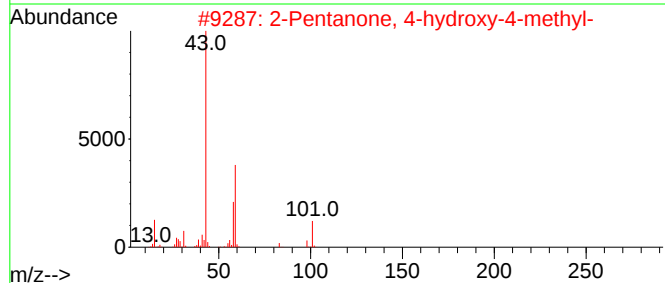
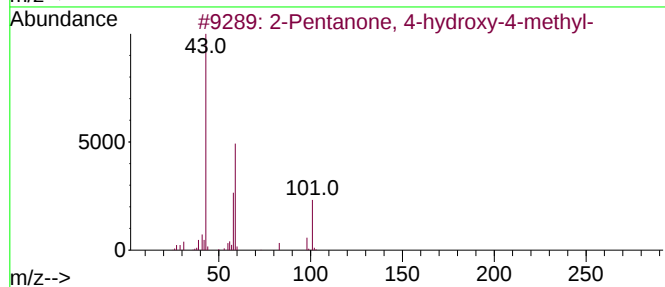
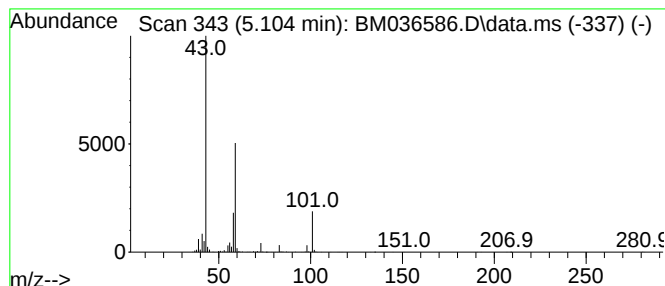
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.104	6.50 ng/ul	880522	1,4-Dichlorobenzene-d4	8.028

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42		
5	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38		



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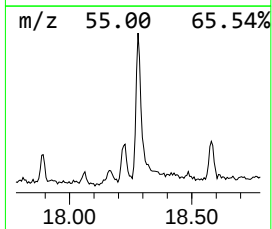
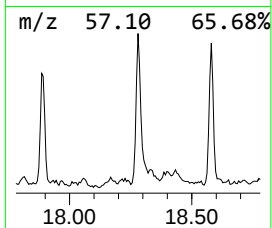
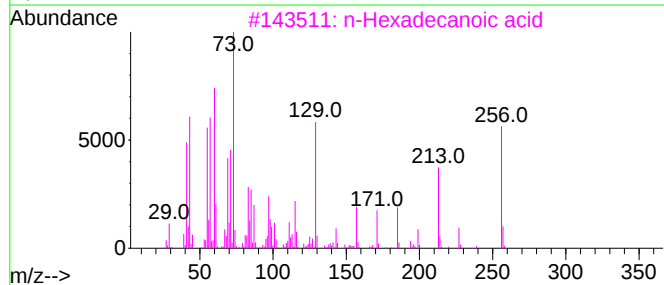
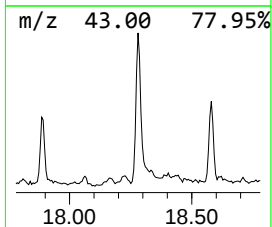
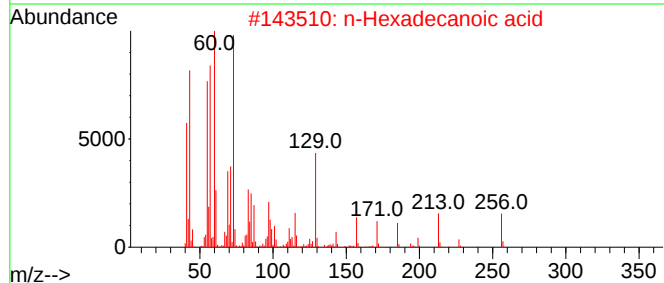
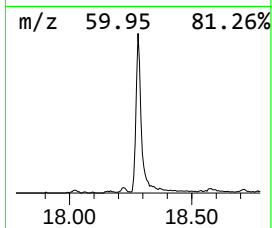
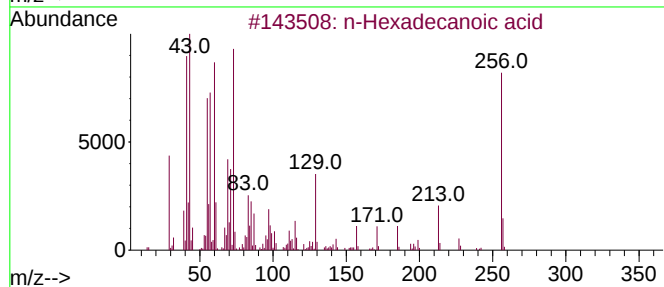
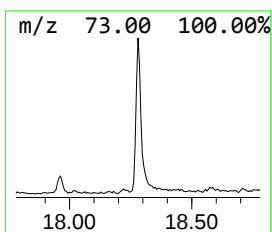
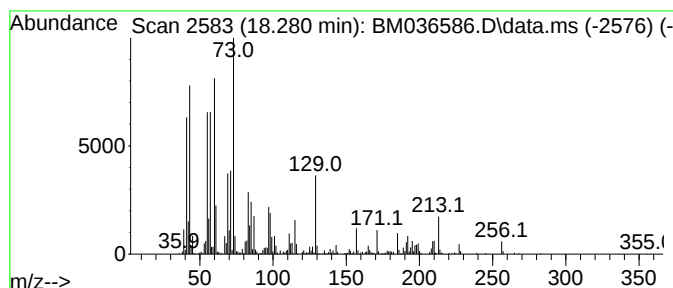
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.280	2.65 ng/ul	882907	Phenanthrene-d10	17.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
Data File : BM036586.D
Acq On : 19 Sep 2022 18:47
Operator : CG/JU
Sample : N4604-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

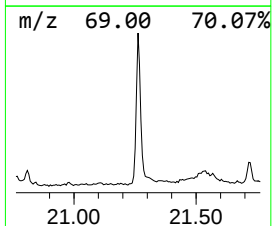
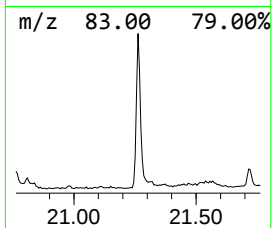
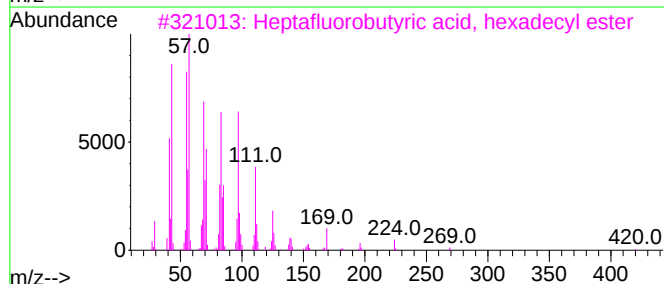
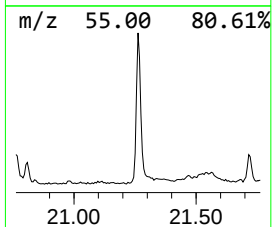
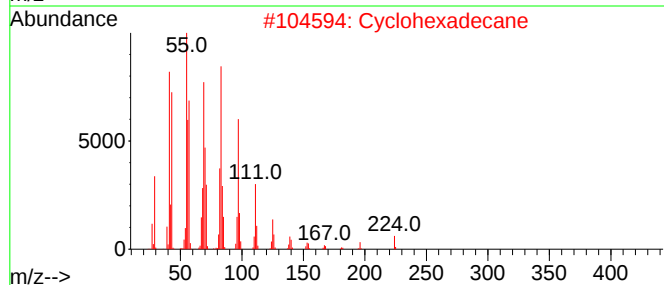
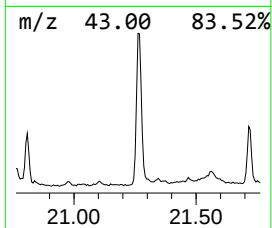
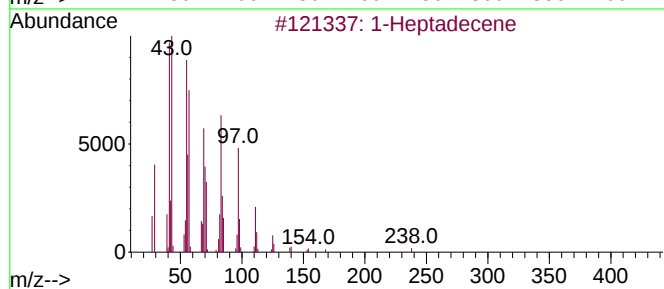
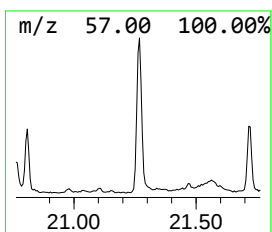
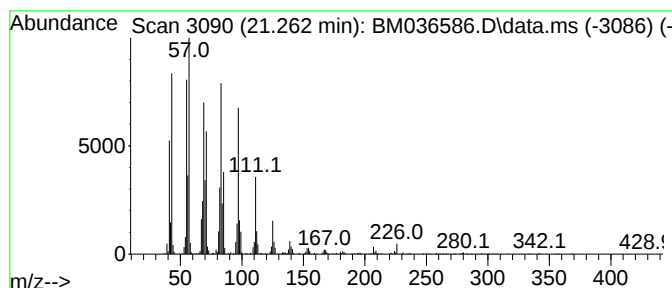
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.262	4.31 ng/ul	1448670	Chrysene-d12	21.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptadecene	238	C17H34	006765-39-5	95
2	Cyclohexadecane	224	C16H32	000295-65-8	94
3	Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	94
4	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
5	1-Hexacosene	364	C26H52	018835-33-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
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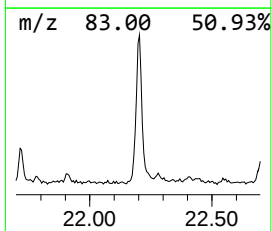
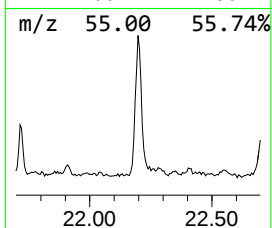
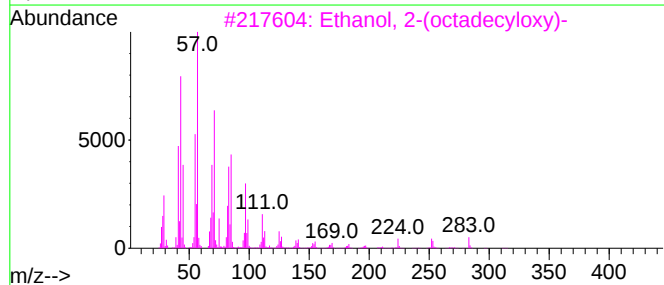
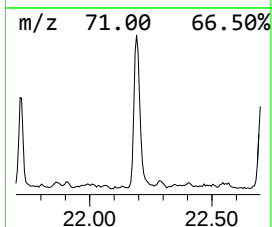
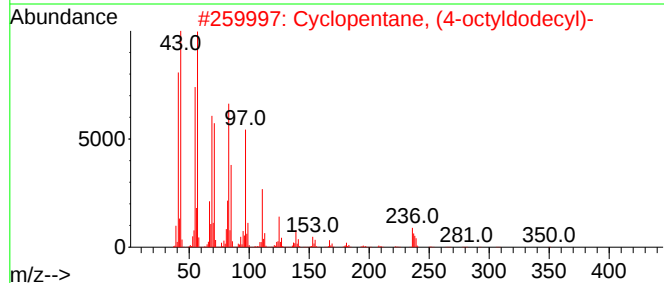
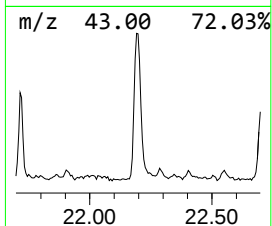
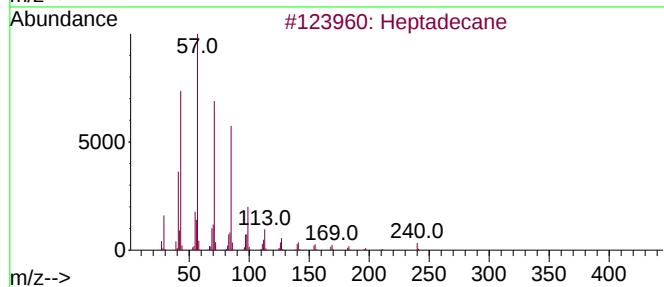
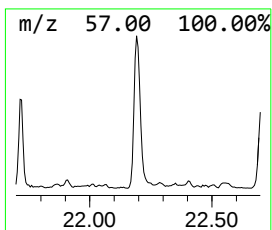
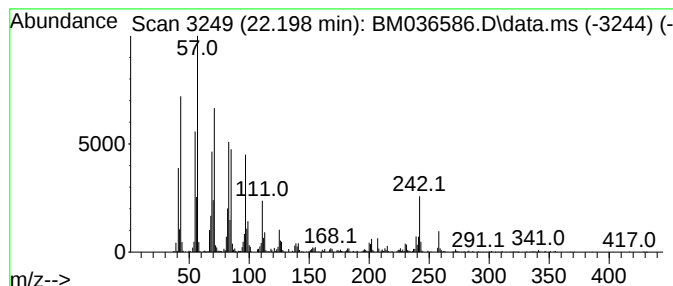
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.198	3.95 ng/ul	1327270	Chrysene-d12	21.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	86
2	Cyclopentane, (4-octyldodecyl)-	350	C25H50	005638-09-5	86
3	Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	78
4	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	70
5	Heneicosyl heptafluorobutyrate	508	C25H43F7O2	1000351-83-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
Data File : BM036586.D
Acq On : 19 Sep 2022 18:47
Operator : CG/JU
Sample : N4604-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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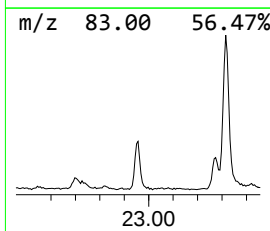
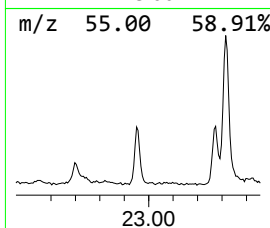
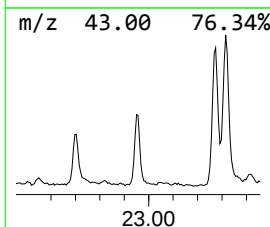
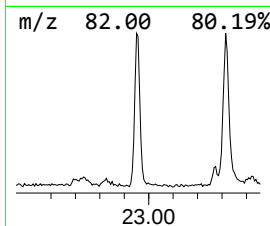
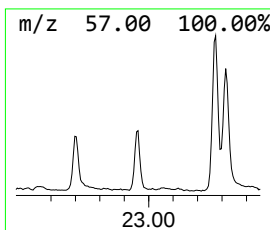
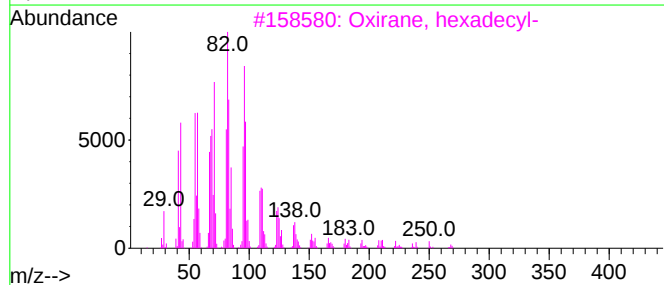
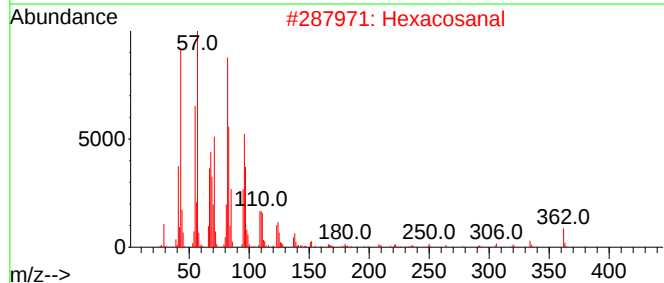
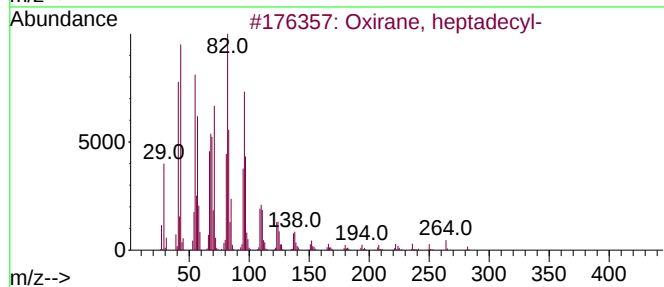
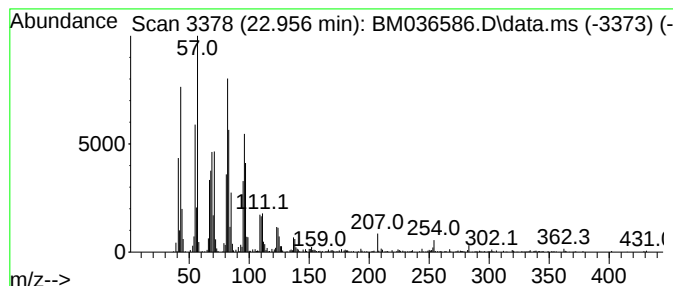
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Oxirane, heptadecyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.956	3.38 ng/ul	804622	Perylene-d12	23.997

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
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Operator : CG/JU
Sample : N4604-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

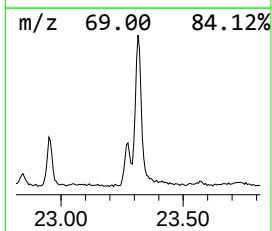
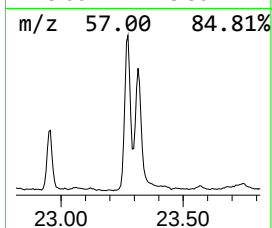
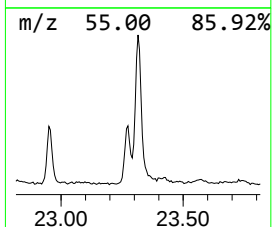
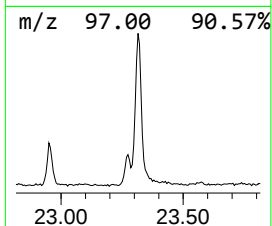
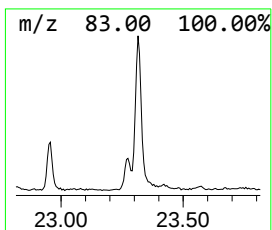
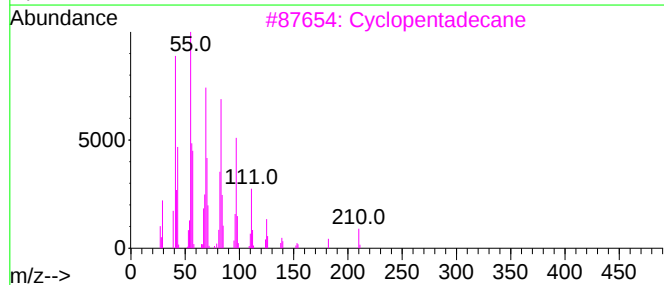
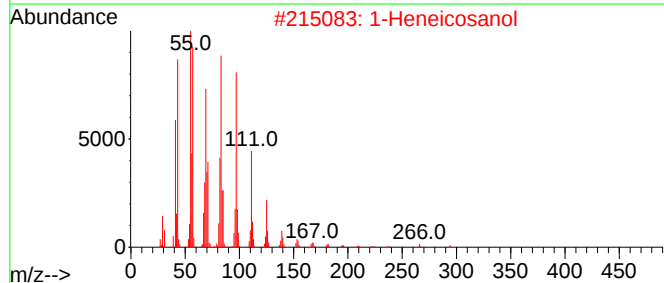
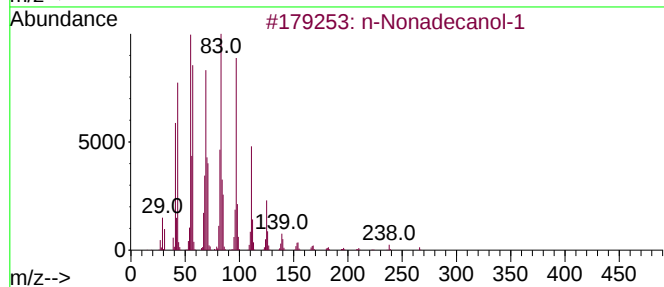
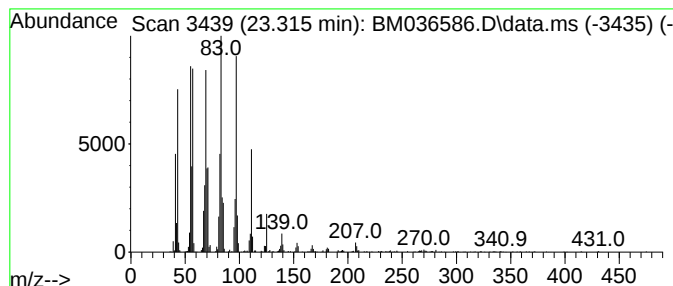
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Nonadecanol-1 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.315	7.88 ng/ul	1875260	Perylene-d12	23.997	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
2	1-Heneicosanol	312	C21H44O	015594-90-8	94
3	Cyclopentadecane	210	C15H30	000295-48-7	93
4	1-Octadecanol	270	C18H38O	000112-92-5	91
5	n-Nonadecanol-1	284	C19H40O	001454-84-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
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Sample : N4604-08
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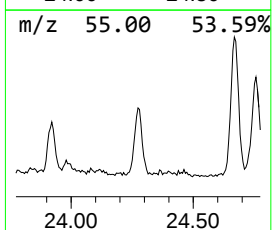
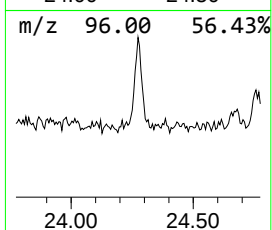
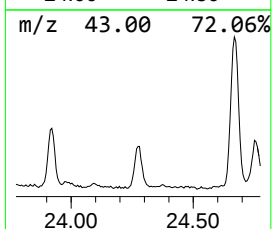
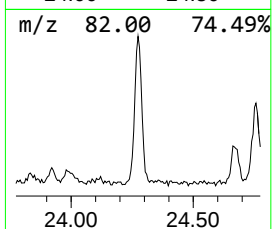
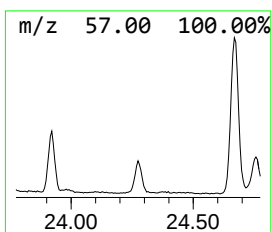
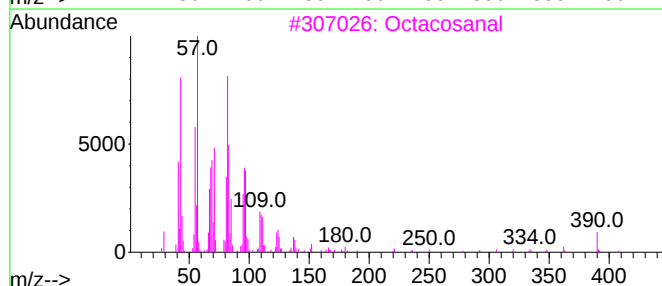
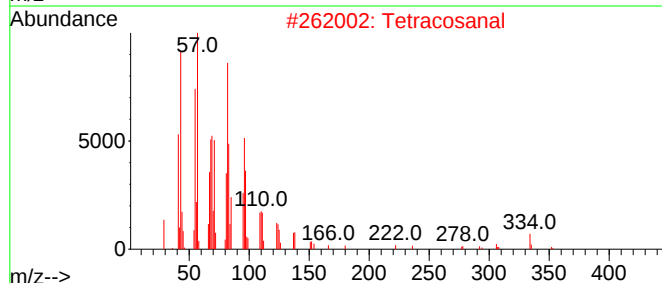
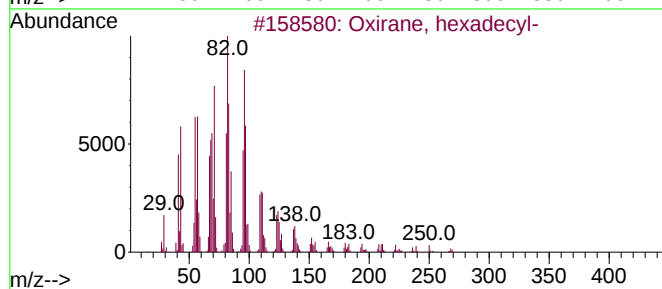
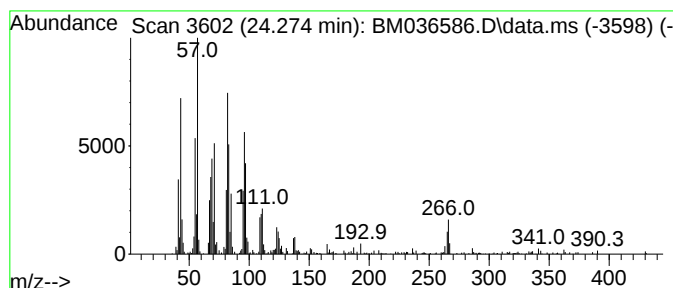
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Oxirane, hexadecyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.274	2.46 ng/ul	586178	Perylene-d12	23.997	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	90
2	Tetracosanal	352	C24H48O	057866-08-7	83
3	Octacosanal	408	C28H56O	022725-64-0	83
4	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	83
5	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
Data File : BM036586.D
Acq On : 19 Sep 2022 18:47
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Misc :
ALS Vial : 14 Sample Multiplier: 1

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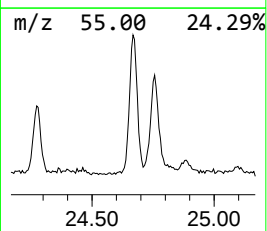
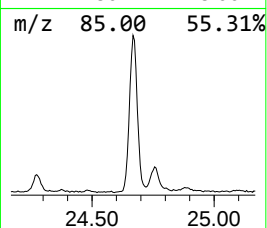
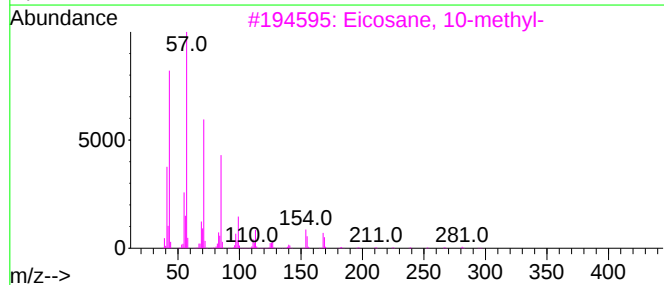
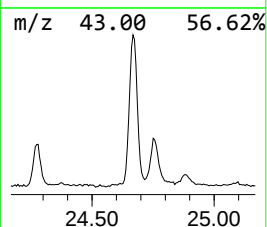
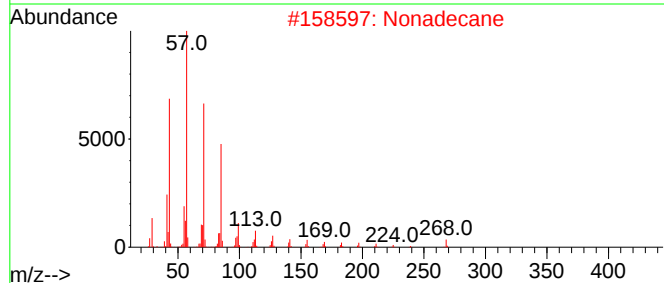
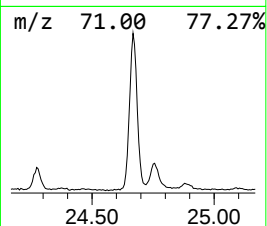
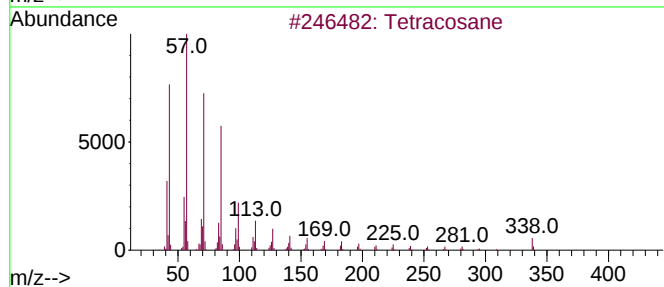
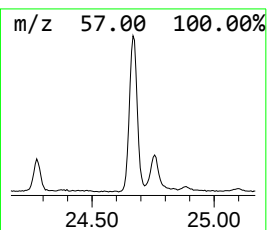
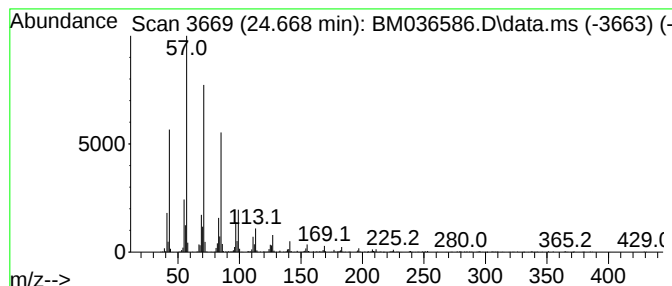
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.668	6.01 ng/ul	1429490	Perylene-d12	23.997

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	96
2			Nonadecane	268	C19H40	000629-92-5	94
3			Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4			Tetracosane	338	C24H50	000646-31-1	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
Data File : BM036586.D
Acq On : 19 Sep 2022 18:47
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Sample : N4604-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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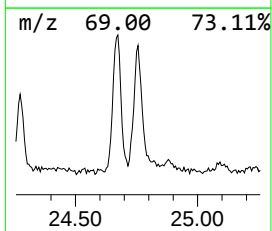
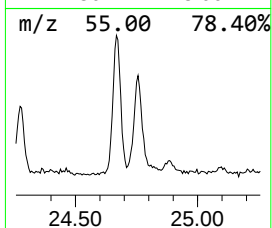
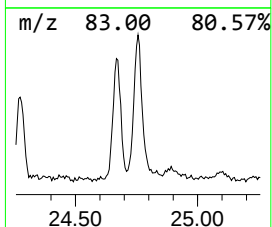
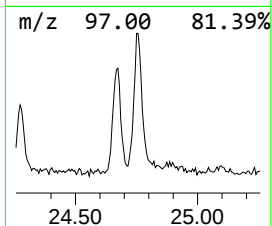
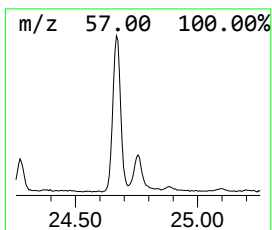
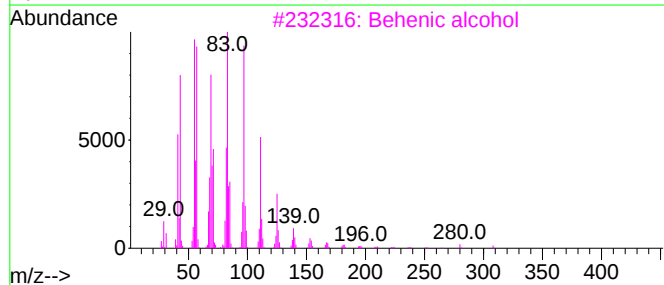
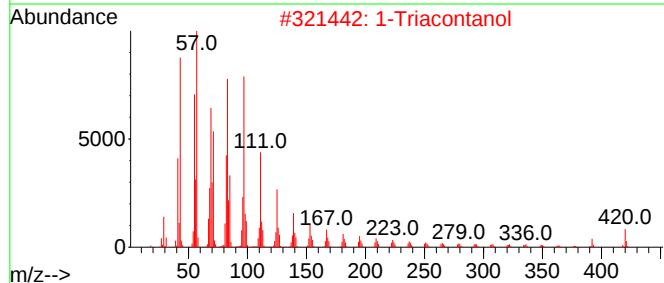
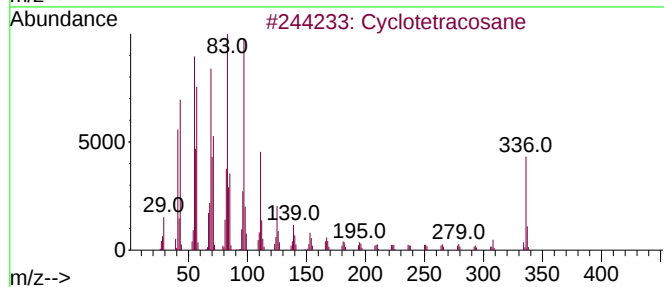
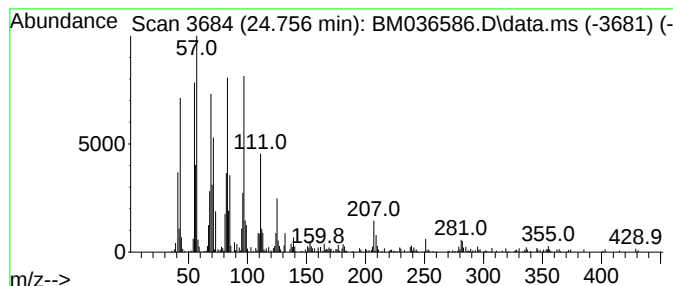
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Cyclic24.756 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.756	2.41 ng/ul	573811	Perylene-d12	23.997

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclotetracosane	336	C24H48	000297-03-0	95
2		1-Triacontanol	438	C30H62O	000593-50-0	94
3		Behenic alcohol	326	C22H46O	000661-19-8	93
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	93
5		1-Hexacosene	364	C26H52	018835-33-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
Data File : BM036586.D
Acq On : 19 Sep 2022 18:47
Operator : CG/JU
Sample : N4604-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A5767

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.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
Butane, 2-metho...	3.128	152.3	ng/ul	20616000	1	8.028	2707980	20.0
2-Pentanone, 4-...	5.104	6.5	ng/ul	880522	1	8.028	2707980	20.0
n-Hexadecanoic ...	18.280	2.6	ng/ul	882907	4	17.386	6669650	20.0
(DEL) Alkane: S...	21.262	4.3	ng/ul	1448670	5	21.551	6718090	20.0
(DEL) Alkane: S...	22.198	4.0	ng/ul	1327270	5	21.551	6718090	20.0
Oxirane, heptad...	22.956	3.4	ng/ul	804622	6	23.997	4760820	20.0
n-Nonadecanol-1	23.315	7.9	ng/ul	1875260	6	23.997	4760820	20.0
Oxirane, hexade...	24.274	2.5	ng/ul	586178	6	23.997	4760820	20.0
(DEL) Alkane: S...	24.668	6.0	ng/ul	1429490	6	23.997	4760820	20.0
(DEL) Alkane: C...	24.756	2.4	ng/ul	573811	6	23.997	4760820	20.0