

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102422\
 Data File : BP012291.D
 Acq On : 24 Oct 2022 18:30
 Operator : CG/JU
 Sample : N5070-08
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BQ6

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012291.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.105	16	20	25	rVB	146630	173319	2.72%	0.269%
2	3.687	116	119	127	rBV2	73027	108362	1.70%	0.168%
3	3.781	132	135	140	rVB2	73477	86358	1.36%	0.134%
4	4.928	327	330	337	rVB	185472	273721	4.30%	0.425%
5	6.987	674	680	692	rBV	929184	1606971	25.23%	2.497%
6	7.152	703	708	719	rVB	822127	1320236	20.73%	2.052%
7	7.346	735	741	751	rBV	956209	1559768	24.49%	2.424%
8	7.816	815	821	828	rBV	1300455	2069125	32.49%	3.215%
9	8.534	937	943	959	rBV	538218	951400	14.94%	1.478%
10	8.987	1013	1020	1030	rBV	994117	1687510	26.50%	2.622%
11	9.375	1082	1086	1091	rVB3	25421	43335	0.68%	0.067%
12	9.705	1135	1142	1158	rBV	741587	1297754	20.38%	2.017%
13	10.246	1227	1234	1246	rBV	1062702	1844494	28.96%	2.866%
14	10.404	1256	1261	1275	rBV	157764	377196	5.92%	0.586%
15	10.622	1290	1298	1305	rBV	1795081	3094853	48.59%	4.809%
16	10.769	1316	1323	1337	rBV	760663	1432083	22.49%	2.225%
17	11.451	1432	1439	1442	rBV6	8875			

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37	14.163	1888	1900	1915	rBV2	2228715	3527376	55.39%	5.482%
38	14.357	1926	1933	1937	rBV2	23135	29624	0.47%	0.046%
39	14.469	1941	1952	1965	rBV	2410926	3605259	56.61%	5.603%
40	14.634	1976	1980	1983	rBV4	5259	6653	0.10%	0.010%

41	14.687	1983	1989	2004	rBV	545480	1061669	16.67%	1.650%
42	14.863	2014	2019	2020	rBV4	11077	16218	0.25%	0.025%
43	14.898	2021	2025	2031	rVB3	44132	72706	1.14%	0.113%
44	14.975	2037	2038	2047	rVB8	7984	13822	0.22%	0.021%
45	15.087	2054	2057	2061	rVB6	6921	10324	0.16%	0.016%

46	15.145	2061	2067	2069	rBV6	4979	8155	0.13%	0.013%
47	15.257	2080	2086	2090	rBV3	10924	21233	0.33%	0.033%
48	15.316	2092	2096	2101	rVB3	22936	32269	0.51%	0.050%
49	15.463	2113	2121	2128	rBV	2944775	4136179	64.94%	6.428%
50	15.592	2138	2143	2154	rBV	349514			

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77	18.392	2616	2619	2623	rBV4	21561	34716	0.55%	0.054%
78	18.445	2625	2628	2633	rVB4	20959	34938	0.55%	0.054%
79	18.586	2649	2652	2657	rBV6	35511	54386	0.85%	0.085%
80	19.139	2744	2746	2751	rVB3	39540	49026	0.77%	0.076%
81	19.210	2754	2758	2760	rBV	57014	80466	1.26%	0.125%
82	19.239	2760	2763	2770	rVB	203090	273068	4.29%	0.424%
83	19.345	2777	2781	2787	rBV	140877	203812	3.20%	0.317%
84	19.569	2813	2819	2831	rBV	4113189	5704262	89.57%	8.864%
85	20.569	2986	2989	2993	rBV2	66352	78136	1.23%	0.121%
86	21.022	3063	3066	3076	rBV2	221821	403850	6.34%	0.628%
87	21.316	3111	3116	3121	rBV	3038609	4015630	63.05%	6.240%
88	23.563	3491	3498	3514	rVB2	3045770	6368749	100.00%	9.897%
89	23.716	3518	3524	3538	rVB2	2101839	4228960	66.40%	6.572%

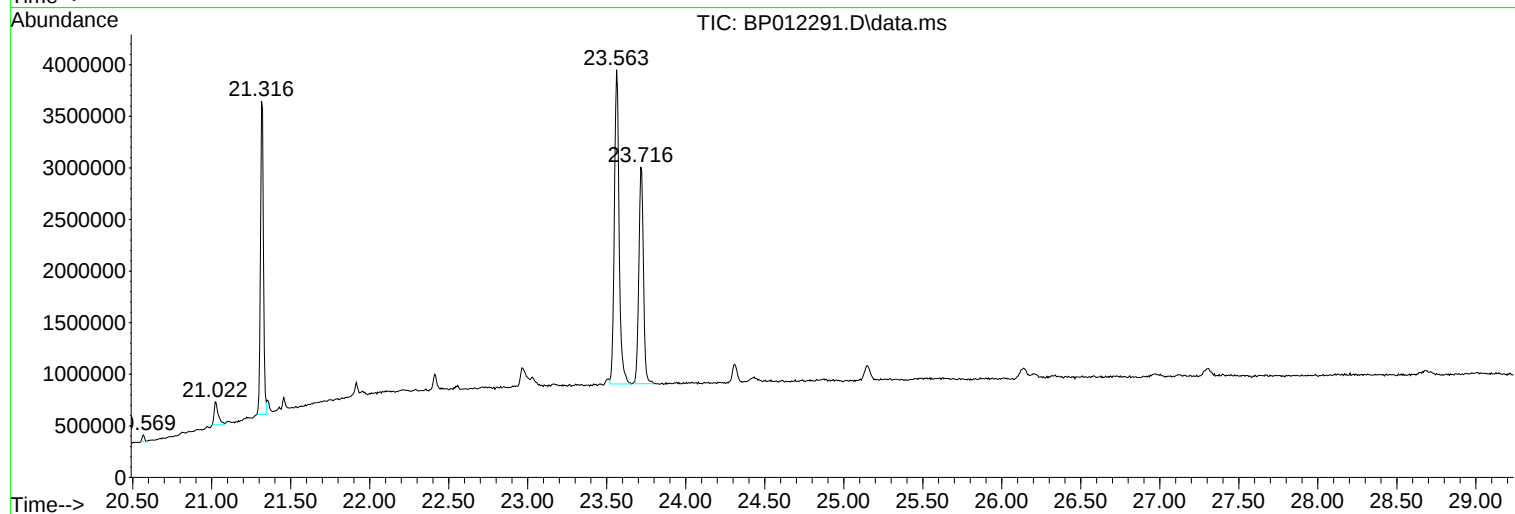
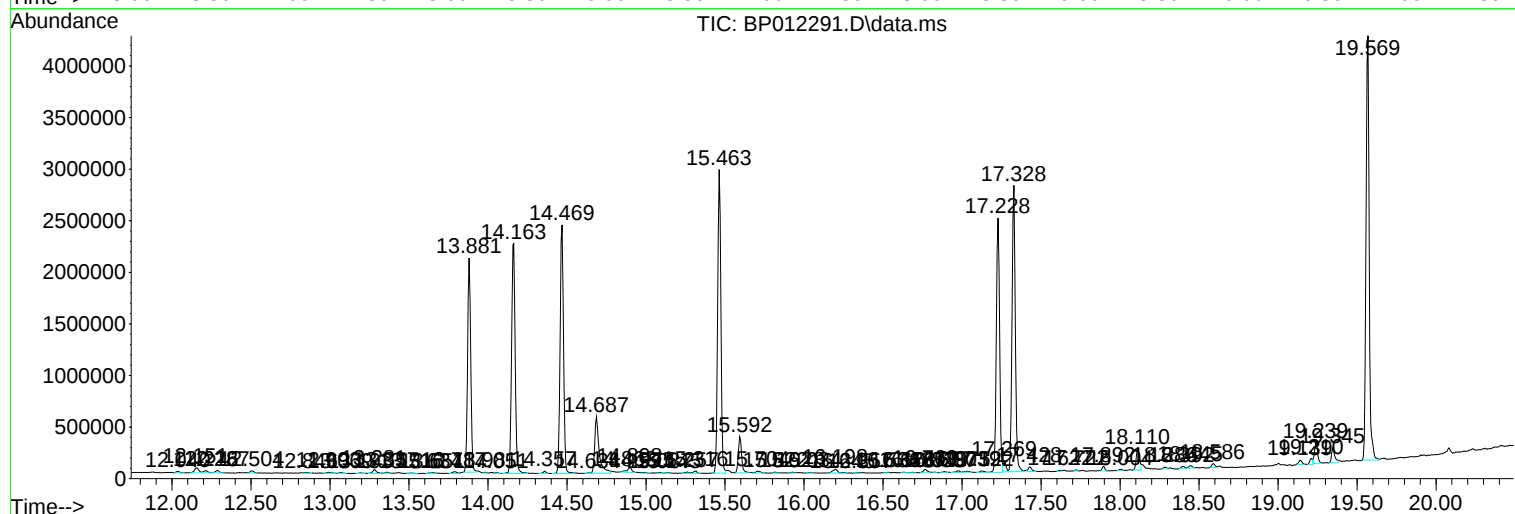
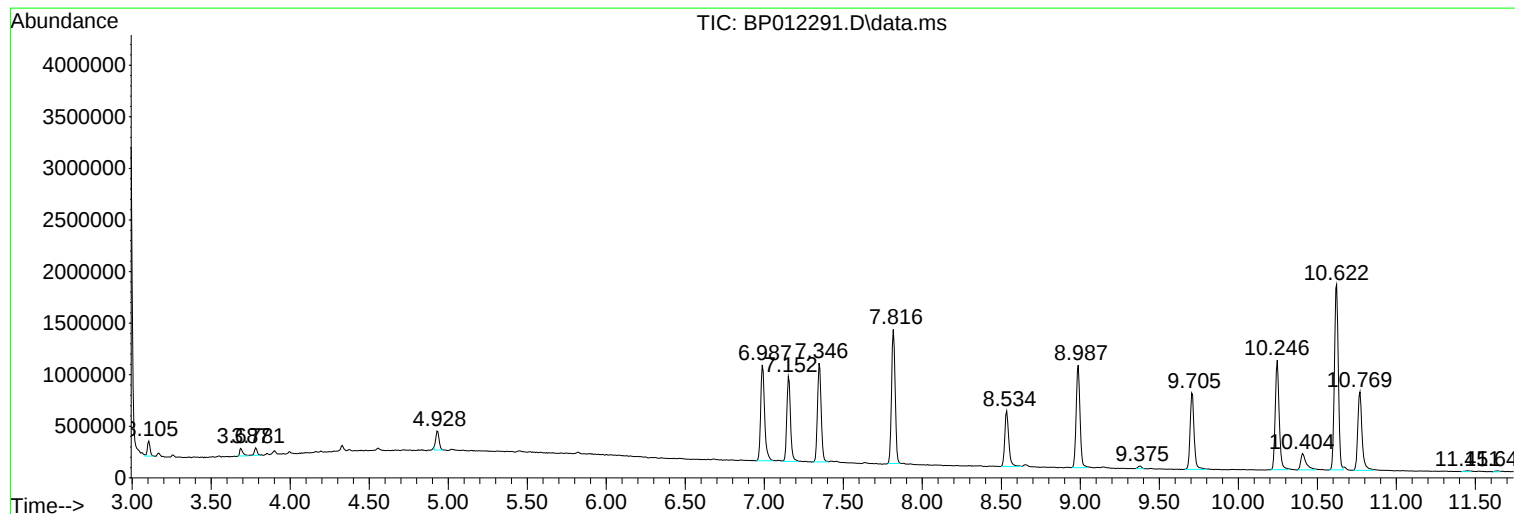
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION

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



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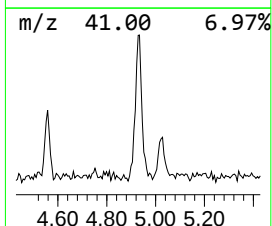
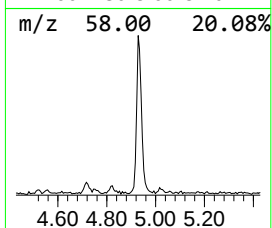
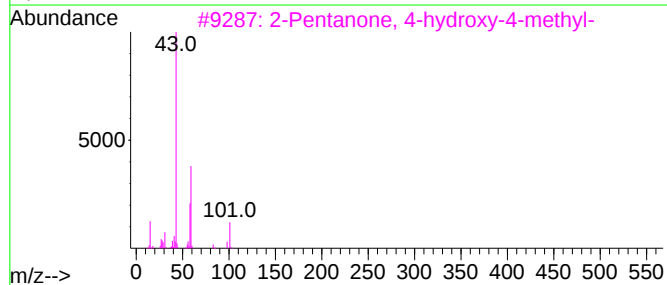
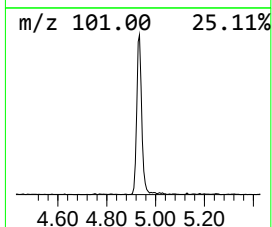
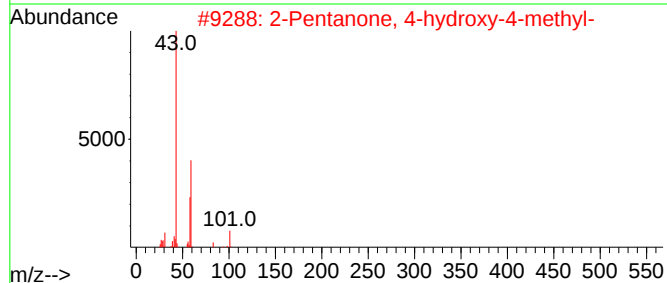
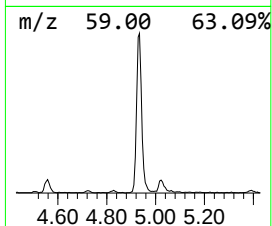
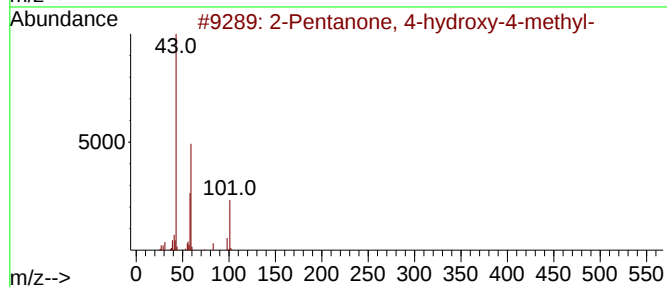
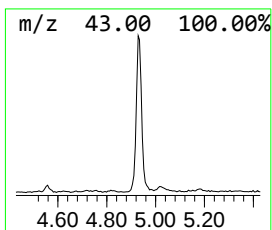
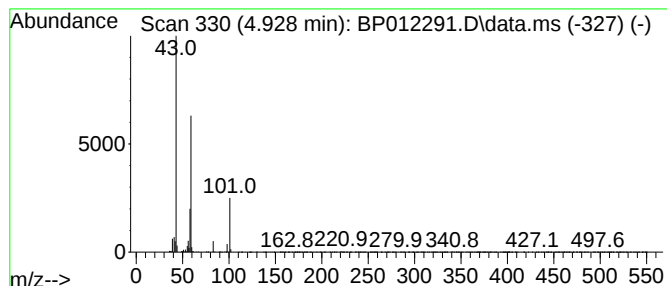
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.928	2.65 ng/ul	273721	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	47		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	37		



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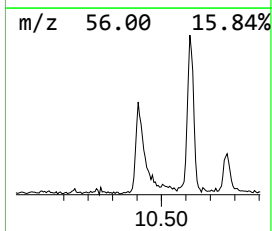
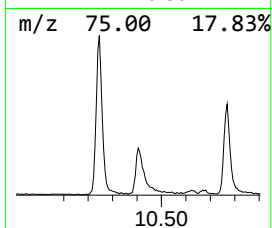
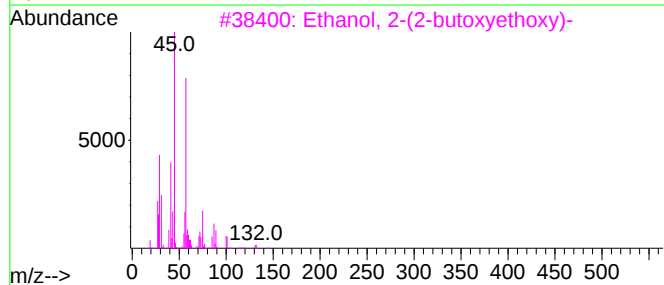
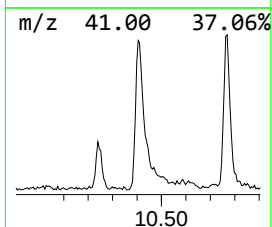
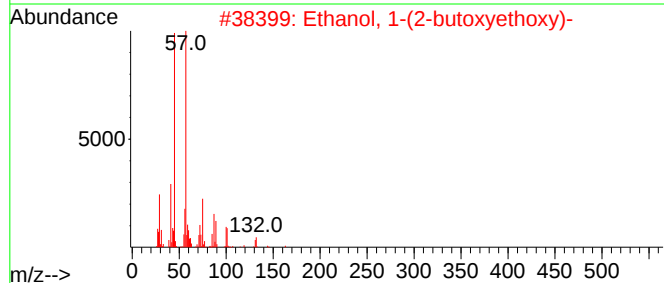
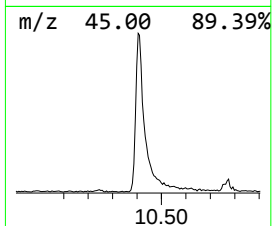
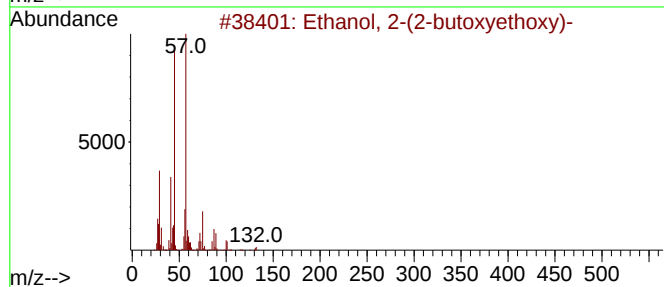
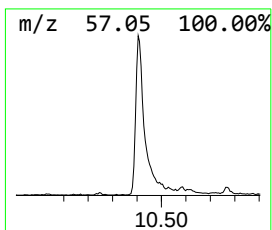
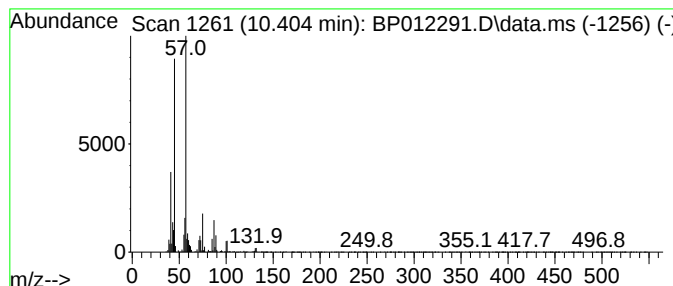
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.405	2.44 ng/ul	377196	Naphthalene-d8	10.622	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	86
3	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
4	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	78
5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72



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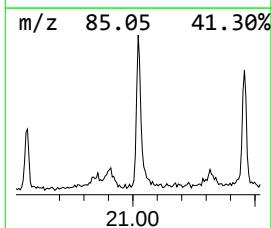
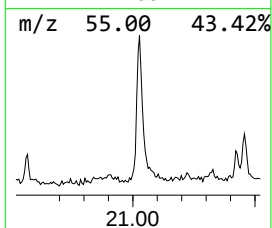
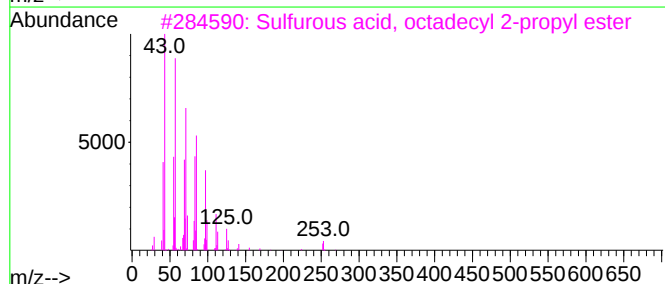
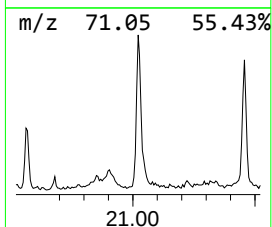
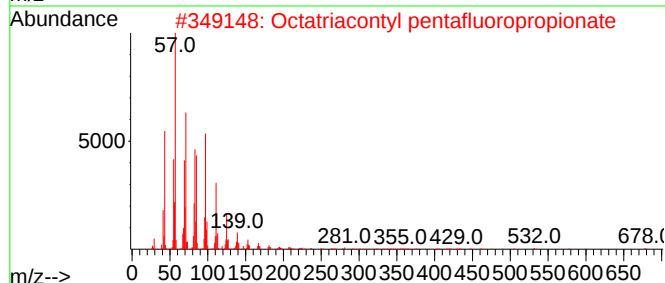
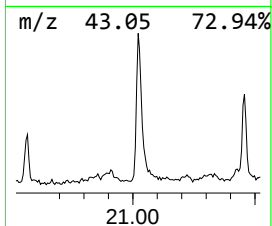
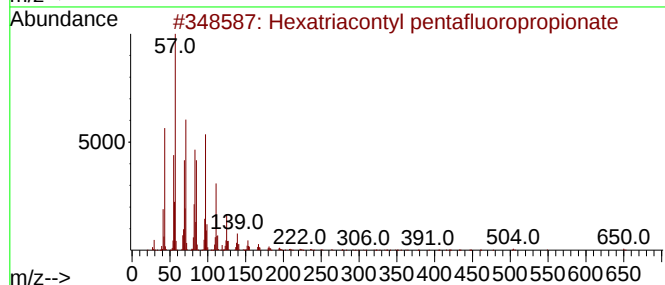
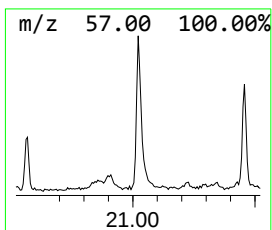
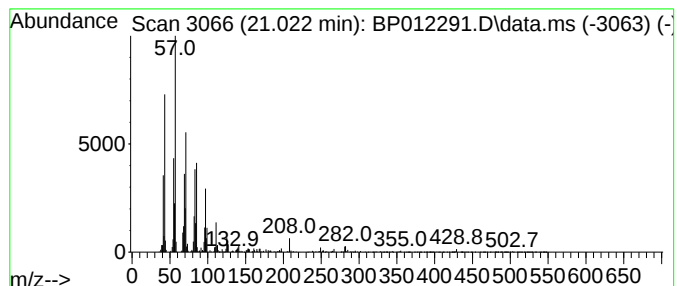
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TIC Library : C:\DATABASE\NIST20.L
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Peak Number 3 Hexatriacontyl pentafluorop... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.022	2.01 ng/ul	403850	Chrysene-d12	21.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexatriacontyl pentafluoropropio...	669	C39H73F5O2	1000351-89-0	91
2			Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	91
3			Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	87
4			Hexacosyl nonyl ether	509	C35H72O	1010406-38-1	86
5			Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	74



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TIC Library : C:\DATABASE\NIST20.L
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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.928	2.6	ng/ul	273721	1	7.816	2069130	20.0
Ethanol, 2-(2-b...	10.405	2.4	ng/ul	377196	2	10.622	3094850	20.0
Hexatriacontyl ...	21.022	2.0	ng/ul	403850	5	21.316	4015630	20.0