

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022122\
 Data File : VD072040.D
 Acq On : 21 Feb 2022 18:51
 Operator : VA/SY
 Sample : N1626-11
 Misc : 6.83G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 CC-A6

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % 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:\voasrv\HPCHEM1\MSVOA_D\Method\82D021822S.M
 Title : SW846 8260

Signal : TIC: VD072040.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.321	61	68	75	rBV4	16193	29437	1.16%	0.188%
2	2.491	89	97	105	rBV	16585	31579	1.24%	0.202%
3	4.156	372	380	391	rBV	39288	110675	4.34%	0.708%
4	4.968	506	518	532	rBV3	32684	109528	4.30%	0.701%
5	7.214	890	900	907	rBV3	12850	38751	1.52%	0.248%
6	7.914	1005	1019	1023	rBV2	207480	489546	19.22%	3.131%
7	7.979	1023	1030	1041	rVB	601944	1358057	53.31%	8.687%
8	8.332	1076	1090	1100	rBV	303393	695048	27.28%	4.446%
9	8.861	1172	1180	1193	rBV	953941	1899917	74.58%	12.153%
10	9.303	1249	1255	1262	rBV2	15428	31138	1.22%	0.199%
11	9.426	1268	1276	1286	rBV3	15256	32909	1.29%	0.210%
12	10.338	1419	1431	1439	rBV	1384599	2547438	100.00%	16.294%
13	10.979	1534	1540	1547	rBV	17555	28792	1.13%	0.184%
14	11.079	1551	1557	1562	rBV4	26374	45777	1.80%	0.293%
15	11.638	1643	1652	1663	rBV	1250514	2179903	85.57%	13.944%
16	12.167	1737	1742	1746	rBV3	15254	29486	1.16%	0.189%
17	12.244	1749	1755	1762	rVV	43838	72864	2.86%	0.466%
18	12.349	1766	1773	1781	rVV7	16305	34291	1.35%	0.219%
19	12.620	1812	1819	1830	rBV	1409276	2313846	90.83%	14.800%
20	12.944	1869	1874	1880	rVV2	123476	203432	7.99%	1.301%
21	13.079	1890	1897	1905	rBV5	17434	48462	1.90%	0.310%
22	13.296	1929	1934	1947	rVB2	44659	96800	3.80%	0.619%
23	13.561	1971	1979	1987	rBV	986851	1668921	65.51%	10.675%
24	13.655	1989	1995	2002	rVV	293531	473880	18.60%	3.031%
25	13.720	2002	2006	2010	rVB6	14759	28339	1.11%	0.181%
26	13.885	2028	2034	2045	rBV4	36804	113012	4.44%	0.723%
27	13.979	2045	2050	2054	rVV2	48428	81464	3.20%	0.521%
28	14.026	2054	2058	2064	rVV6	30454	63389	2.49%	0.405%
29	14.085	2064	2068	2075	rVB2	29949	56334	2.21%	0.360%
30	14.232	2087	2093	2098	rBV2	44892	82434	3.24%	0.527%
31	14.338	2106	2111	2116	rVB4	26017	37442	1.47%	0.239%
32	14.638	2156	2162	2168	rVB2	48377	81624	3.20%	0.522%
33	14.749	2177	2181	2185	rVB2	27013	38972	1.53%	0.249%
34	14.855	2193	2199	2201	rBV5	20162	30917	1.21%	0.198%
35	14.879	2201	2203	2211	rVB8	17418	26889	1.06%	0.172%
36	15.020	2224	2227	2231	rVV6	23444	48445	1.90%	0.310%

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37	15.091	2234	2239	2245	rVV4	64957	136469	5.36%	0.873%
38	15.173	2249	2253	2257	rVB6	20692	34931	1.37%	0.223%
39	15.314	2270	2277	2284	rVB10	12982	36476	1.43%	0.233%
40	15.379	2284	2288	2292	rBV2	15023	27616	1.08%	0.177%
41	15.761	2344	2353	2358	rBV10	11608	29852	1.17%	0.191%
42	16.473	2469	2474	2483	rBV4	27001	63000	2.47%	0.403%
43	16.685	2504	2510	2517	rVB4	20734	45713	1.79%	0.292%

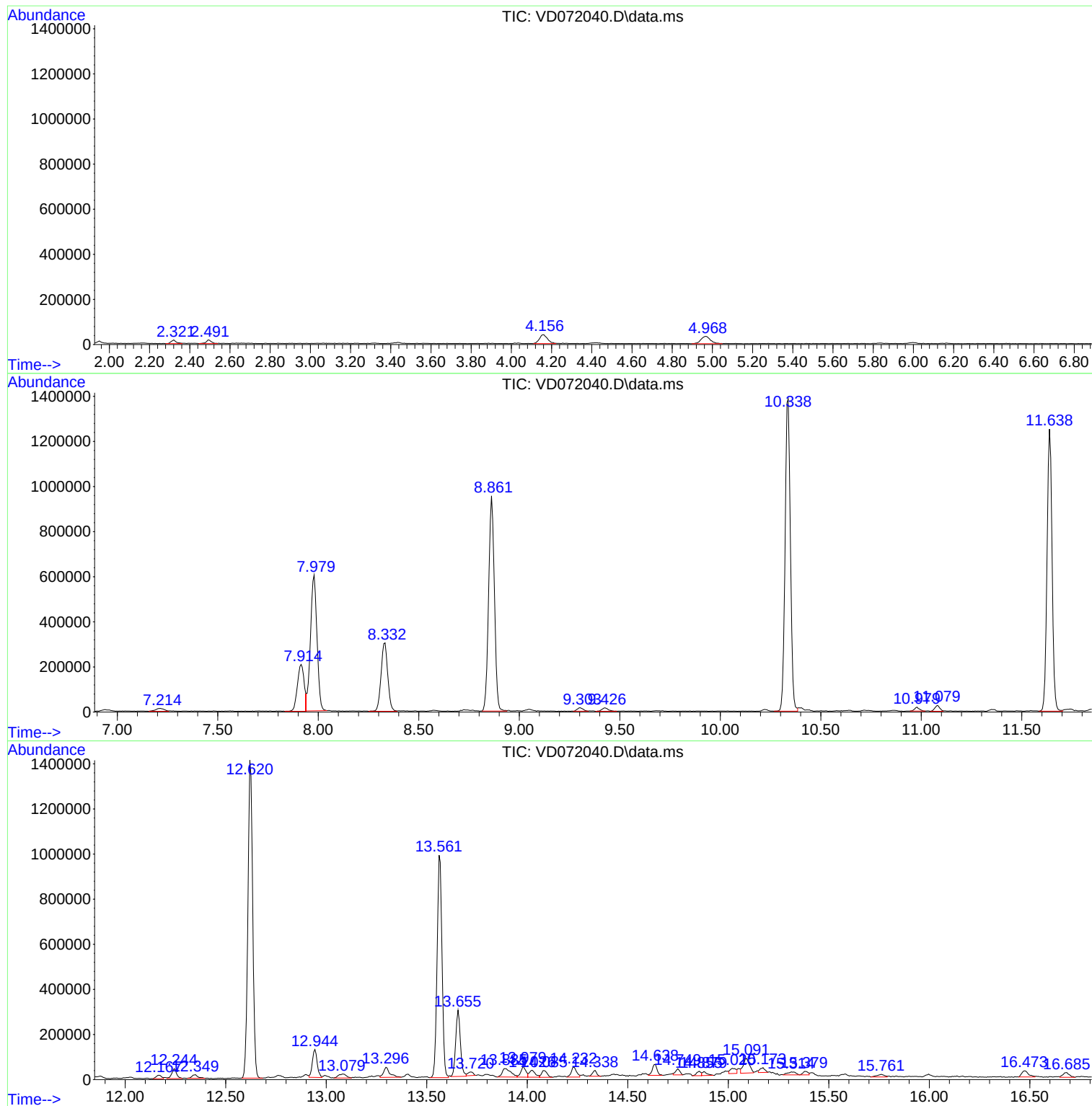
Sum of corrected areas: 15633795

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021822S.M
Quant Title : SW846 8260

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



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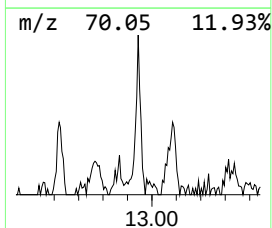
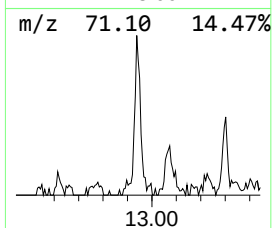
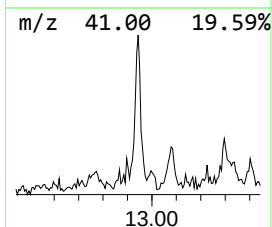
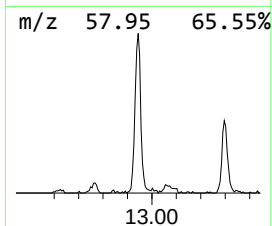
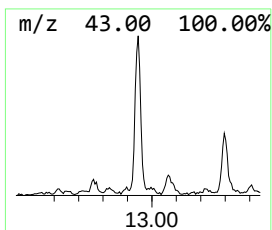
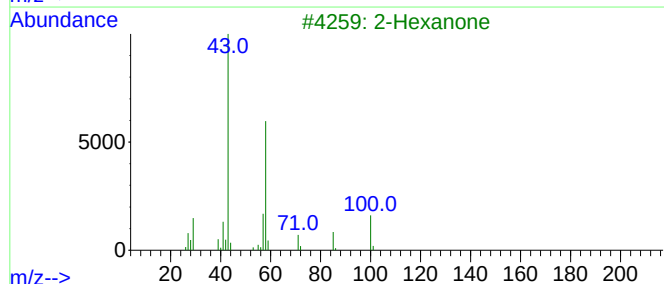
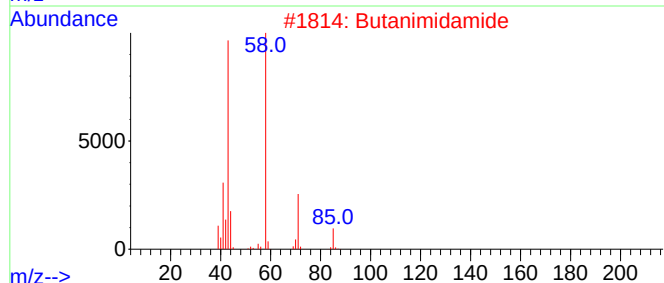
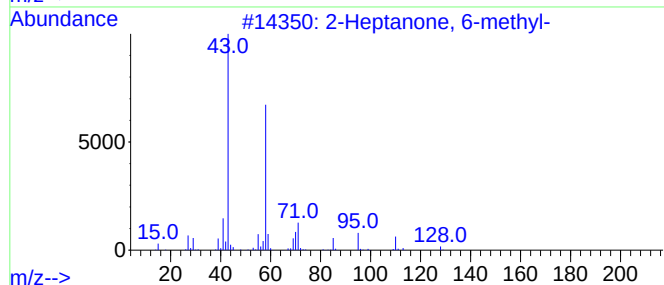
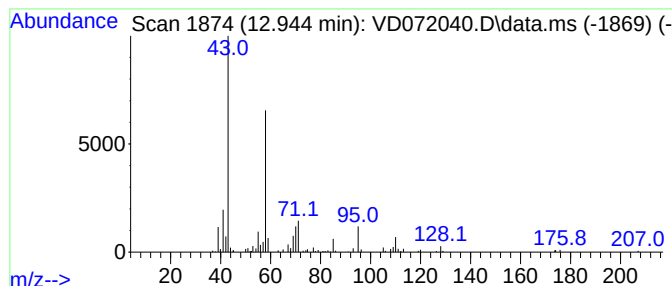
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021822S.M
Quant Title : SW846 8260

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

Peak Number 1 2-Heptanone, 6-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.944	6.09 ug/l	203432	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Heptanone, 6-methyl-	128	C8H16O	000928-68-7	91
2			Butanimidamide	86	C4H10N2	000107-90-4	53
3			2-Hexanone	100	C6H12O	000591-78-6	53
4			2-Octanone	128	C8H16O	000111-13-7	42
5			Pentanal, 2-methyl-	100	C6H12O	000123-15-9	42



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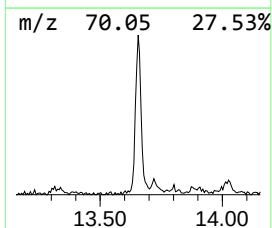
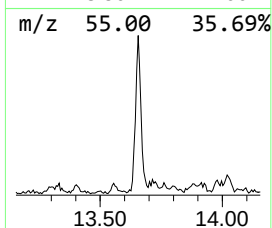
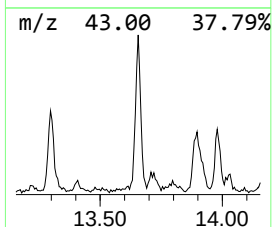
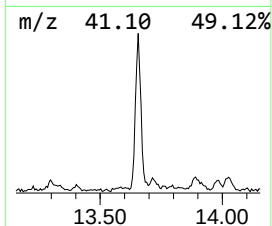
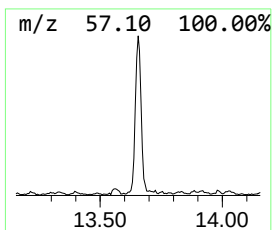
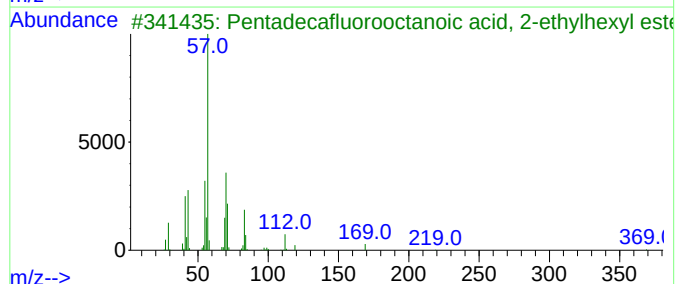
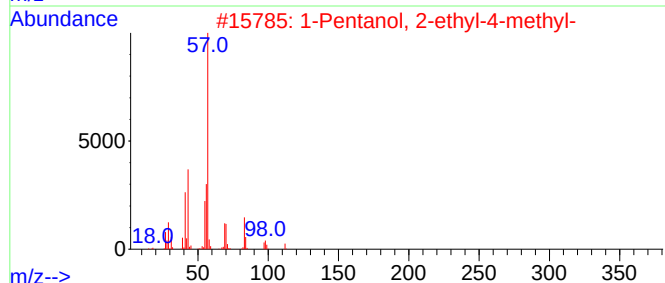
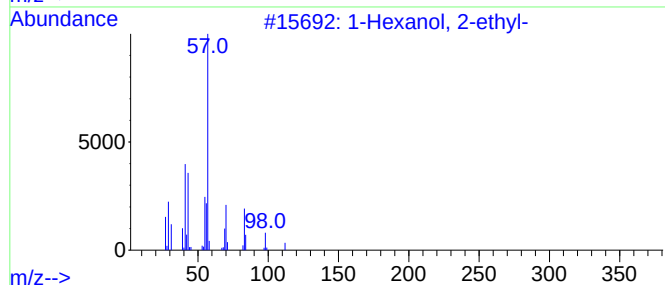
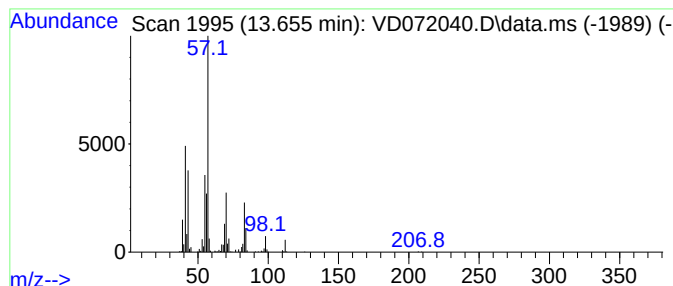
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Peak Number 2 1-Hexanol, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.655	14.20 ug/l	473880	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	90
2			1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	72
3			Pentadecafluorooctanoic acid, 2-...	526	C16H17F15O2	1000406-80-9	59
4			Decanal	156	C10H20O	000112-31-2	50
5			1-Decene, 2,4-dimethyl-	168	C12H24	055170-80-4	47



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Heptanone, 6-...	12.944	6.1	ug/l	203432	4	13.561	1668920	50.0
1-Hexanol, 2-et...	13.655	14.2	ug/l	473880	4	13.561	1668920	50.0