

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120924\
 Data File : VX044190.D
 Acq On : 09 Dec 2024 16:04
 Operator : JC/MD
 Sample : P5145-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 286085

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
 Title : SW846 8260

Signal : TIC: VX044190.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.142	5	9	17	rVB	105750	122956	18.59%	2.983%
2	1.563	68	78	79	rBV	15310	33717	5.10%	0.818%
3	2.392	208	214	220	rBV	14980	25979	3.93%	0.630%
4	2.709	259	266	277	rBV	105545	189558	28.67%	4.599%
5	4.587	565	574	586	rBV2	5537	17739	2.68%	0.430%
6	5.032	641	647	652	rBV4	3326	8227	1.24%	0.200%
7	5.385	694	705	721	rBV	73900	221352	33.47%	5.370%
8	5.544	721	731	747	rVB	121017	344160	52.05%	8.349%
9	5.952	788	798	813	rBV	89809	251181	37.99%	6.094%
10	6.757	920	930	943	rBV	198721	485338	73.40%	11.774%
11	8.647	1233	1240	1252	rBV	417148	661260	100.00%	16.042%
12	9.531	1379	1385	1392	rBV5	3880	6780	1.03%	0.164%
13	10.055	1464	1471	1479	rBV	389622	572918	86.64%	13.899%
14	11.079	1633	1639	1651	rBV	336330	432779	65.45%	10.499%
15	11.305	1670	1676	1683	rVB10	2961	7280	1.10%	0.177%
16	11.915	1771	1776	1785	rBV10	4025	11163	1.69%	0.271%
17	12.018	1788	1793	1802	rVB	381973	498450	75.38%	12.092%
18	12.128	1807	1811	1820	rVV8	5936	16939	2.56%	0.411%
19	12.219	1820	1826	1829	rVV3	9212	15633	2.36%	0.379%
20	12.280	1832	1836	1846	rVB9	8224	19401	2.93%	0.471%
21	12.670	1895	1900	1906	rBV8	6303	14764	2.23%	0.358%
22	12.859	1927	1931	1937	rVB7	7643	12891	1.95%	0.313%
23	12.920	1937	1941	1944	rBV4	8560	13850	2.09%	0.336%
24	13.597	2048	2052	2056	rBV5	5733	8975	1.36%	0.218%
25	13.640	2056	2059	2065	rVB8	3568	6813	1.03%	0.165%
26	13.853	2090	2094	2097	rBV6	4831	7458	1.13%	0.181%
27	14.195	2145	2150	2154	rBV2	26186	41912	6.34%	1.017%
28	14.323	2167	2171	2174	rBV6	4211	6969	1.05%	0.169%
29	14.359	2175	2177	2185	rVB10	6342	11841	1.79%	0.287%
30	14.445	2187	2191	2193	rBV4	4428	6711	1.01%	0.163%
31	14.896	2260	2265	2272	rBV4	5969	14043	2.12%	0.341%
32	15.213	2315	2317	2325	rVB8	4755	9321	1.41%	0.226%
33	15.316	2331	2334	2348	rBV8	8544	23670	3.58%	0.574%

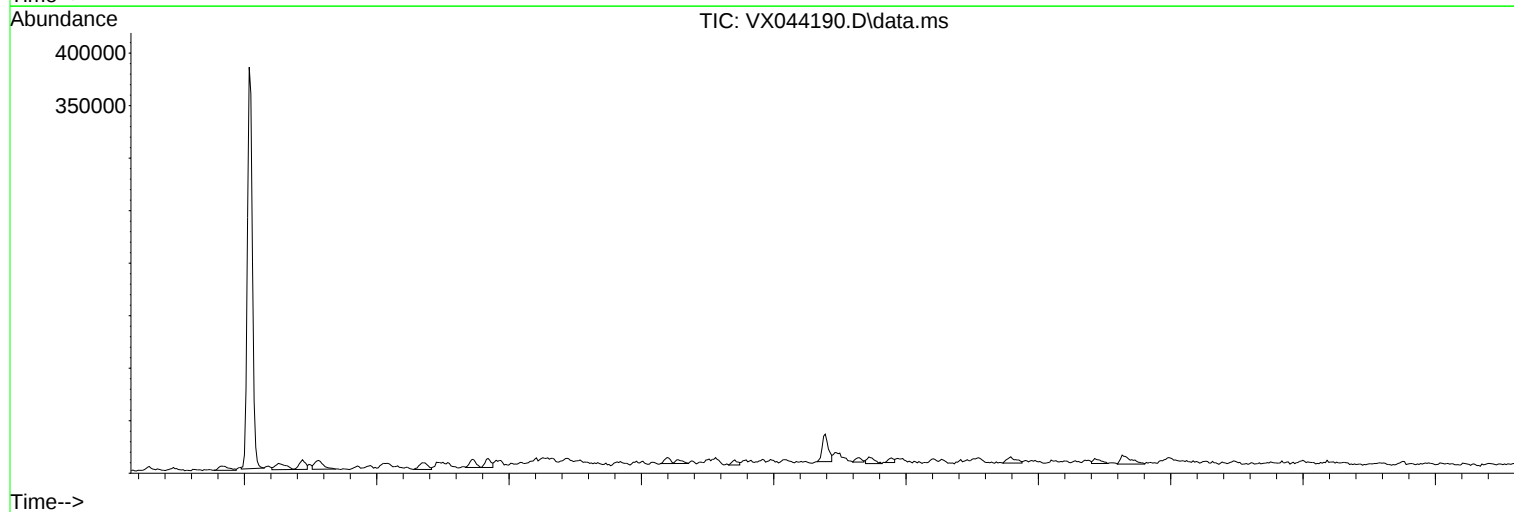
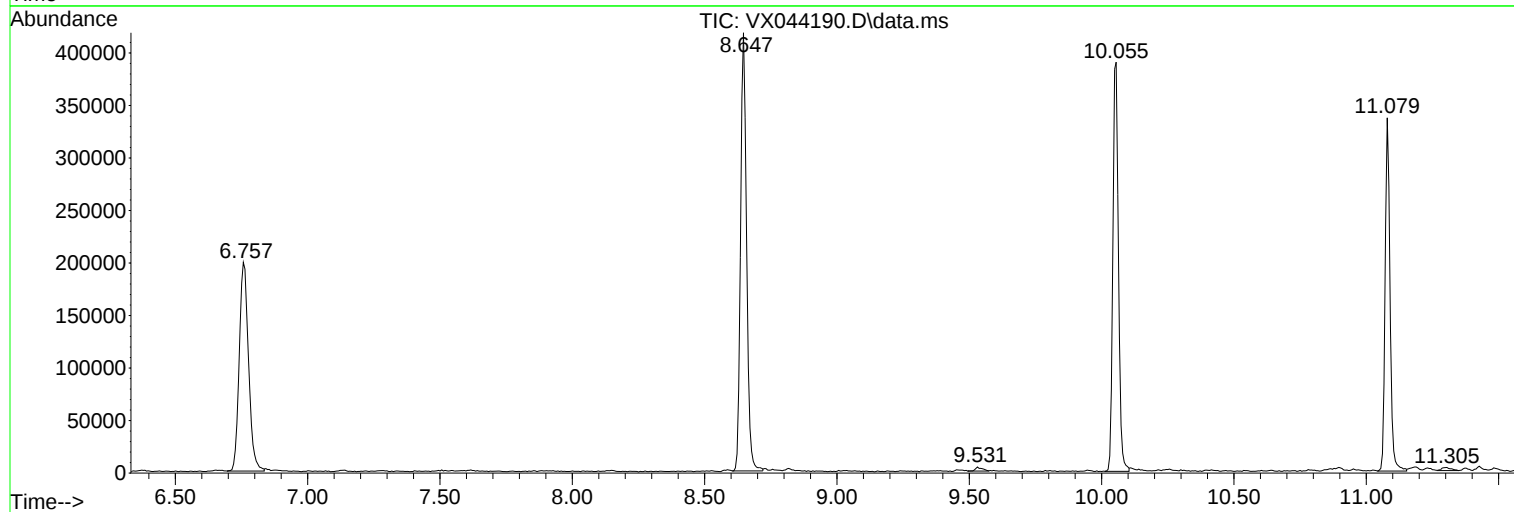
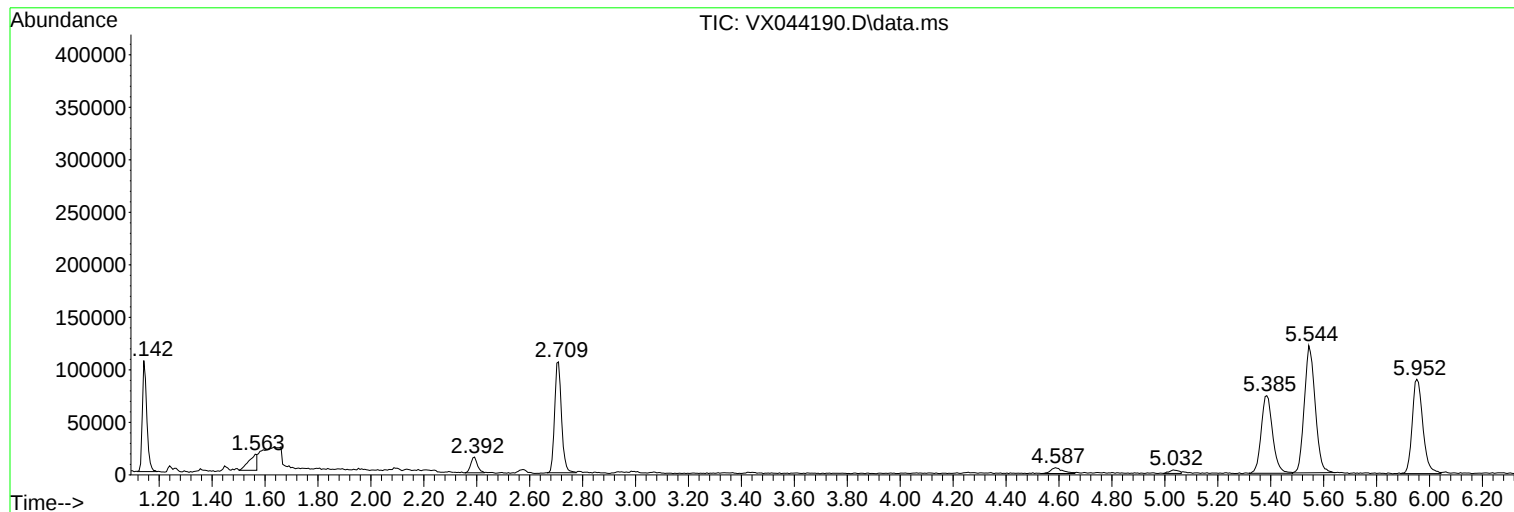
Sum of corrected areas: 4122028

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
Quant Title : SW846 8260

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



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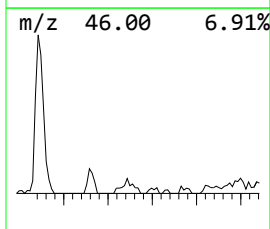
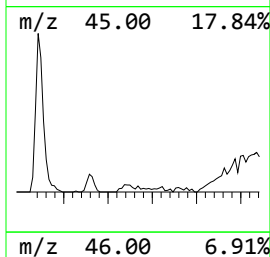
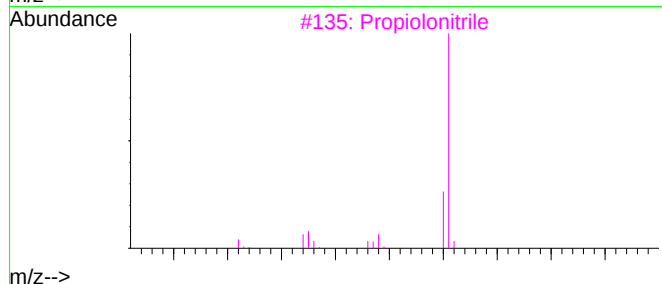
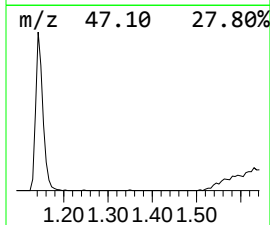
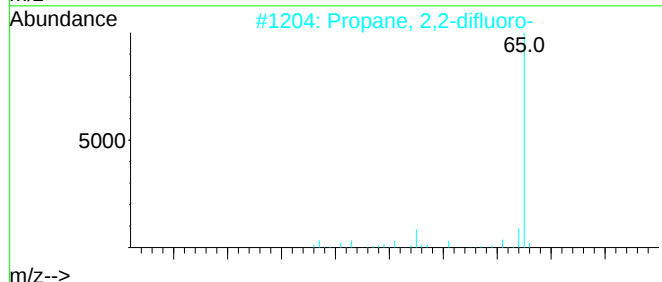
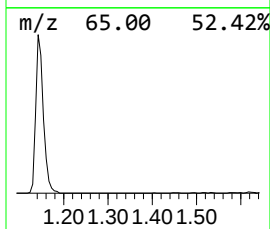
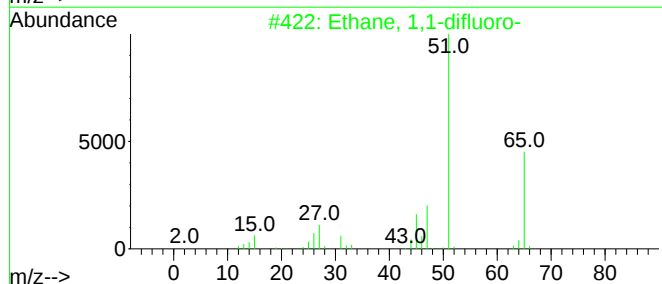
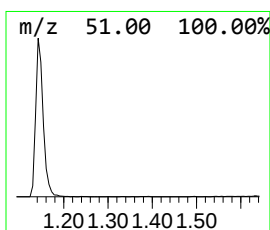
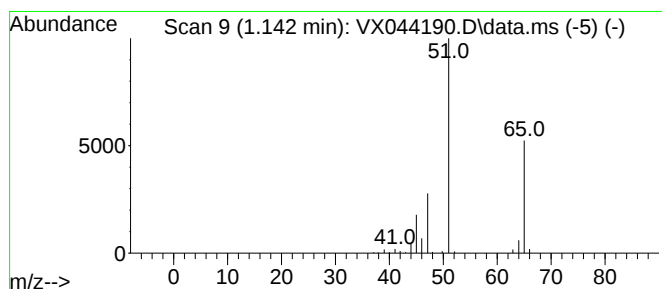
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethane, 1,1-difluoro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.142	17.86 ug/l	122956	Pentafluorobenzene	5.544

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethane, 1,1-difluoro-	66	C2H4F2	000075-37-6	91
2	Propane, 2,2-difluoro-	80	C3H6F2	000420-45-1	4
3	Propiolonitrile	51	C3HN	001070-71-9	3
4	Difluoromethane	52	CH2F2	000075-10-5	1
5	2-Chloroethanol	80	C2H5ClO	000107-07-3	1



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TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Ethane, 1,1-dif...	1.142	17.9	ug/l	122956	1	5.544	344160	50.0