

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052324\
 Data File : VN082317.D
 Acq On : 23 May 2024 13:29
 Operator : JC\MD
 Sample : P2556-01 20X
 Misc : 5.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 COMP-1

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_N\methods\82N051524W.M

Title : SW846 8260

Signal : TIC: VN082317.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.377	65	72	74	rBV6	5819	13236	1.22%	0.206%
2	2.824	139	148	156	rVB6	20025	64323	5.94%	1.003%
3	8.171	1047	1057	1061	rBV	134269	318382	29.40%	4.965%
4	8.224	1061	1066	1077	rVB	230073	521631	48.17%	8.135%
5	8.583	1117	1127	1136	rBV	166369	369365	34.11%	5.760%
6	8.683	1136	1144	1145	rBV4	7619	15628	1.44%	0.244%
7	9.106	1207	1216	1228	rBV	354138	740160	68.36%	11.543%
8	10.018	1362	1371	1372	rBV6	5401	11703	1.08%	0.183%
9	10.571	1457	1465	1473	rVV	581172	1082811	100.00%	16.887%
10	10.629	1473	1475	1482	rVB3	15300	25142	2.32%	0.392%
11	11.865	1677	1685	1695	rBV	520694	948682	87.61%	14.795%
12	12.770	1830	1839	1840	rBV5	10471	19752	1.82%	0.308%
13	12.853	1846	1853	1861	rVB	418836	734253	67.81%	11.451%
14	13.170	1904	1907	1914	rVB6	7241	11738	1.08%	0.183%
15	13.423	1939	1950	1951	rBV8	11551	24583	2.27%	0.383%
16	13.482	1958	1960	1966	rVB4	13177	15993	1.48%	0.249%
17	13.706	1994	1998	2003	rVB7	9836	15829	1.46%	0.247%
18	13.794	2003	2013	2023	rBV	504500	897013	82.84%	13.989%
19	14.117	2062	2068	2074	rBV7	13497	29311	2.71%	0.457%
20	14.435	2119	2122	2126	rBV5	6958	11053	1.02%	0.172%
21	14.506	2130	2134	2140	rBV9	5681	12043	1.11%	0.188%
22	14.629	2151	2155	2159	rBV7	16006	29591	2.73%	0.461%
23	14.900	2198	2201	2207	rVB8	8631	15845	1.46%	0.247%
24	15.064	2220	2229	2233	rBV10	22976	59466	5.49%	0.927%
25	15.212	2249	2254	2259	rBV5	21217	38944	3.60%	0.607%
26	15.264	2261	2263	2268	rVB6	9648	16095	1.49%	0.251%
27	15.435	2288	2292	2298	rVB8	12204	25617	2.37%	0.400%
28	15.700	2333	2337	2343	rBV5	18576	36253	3.35%	0.565%
29	15.788	2348	2352	2362	rVB5	18798	44592	4.12%	0.695%
30	16.017	2385	2391	2398	rBV5	39309	87375	8.07%	1.363%
31	16.170	2412	2417	2422	rBV6	33692	64719	5.98%	1.009%
32	16.711	2503	2509	2515	rBV4	49361	111089	10.26%	1.732%

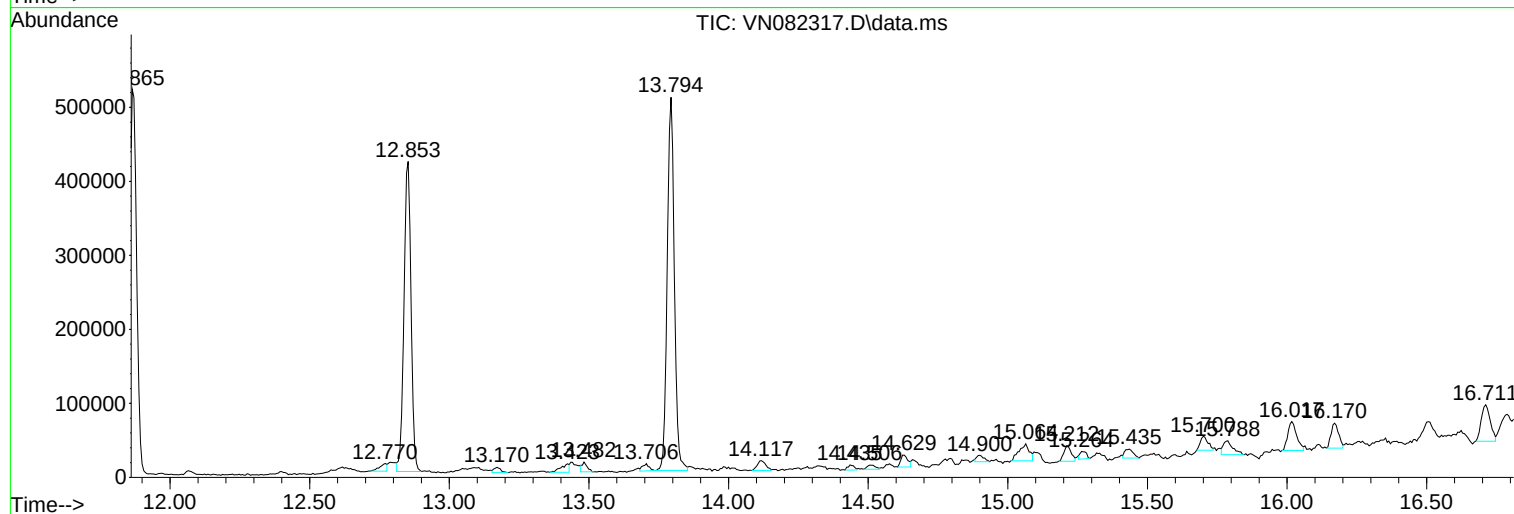
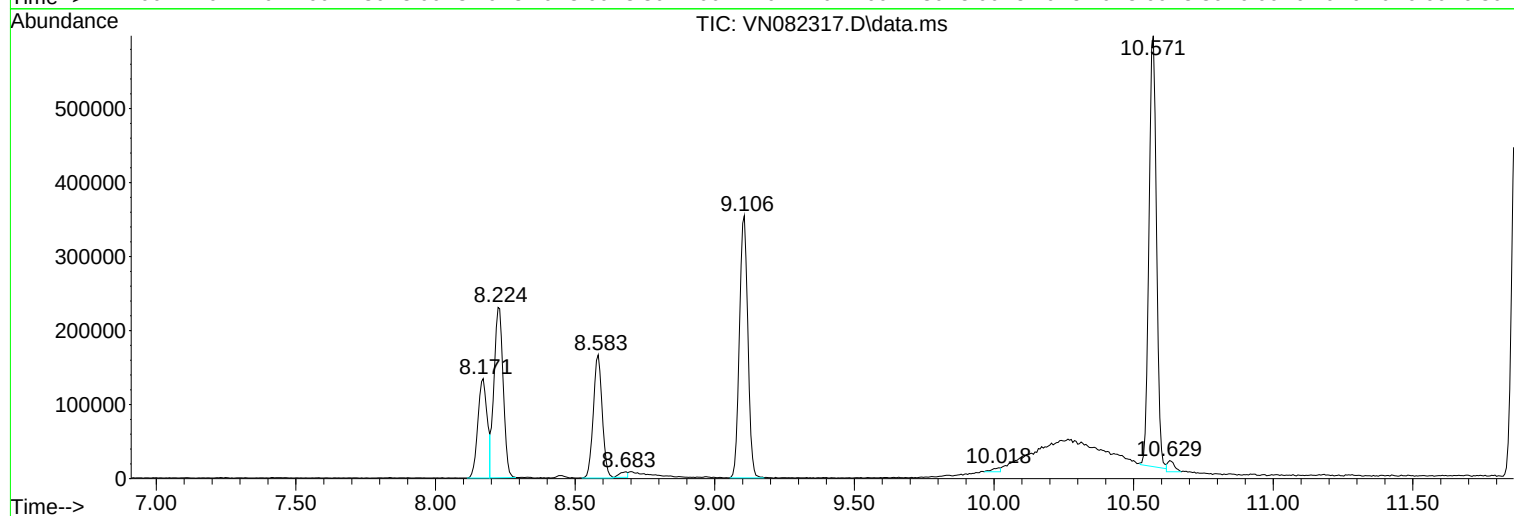
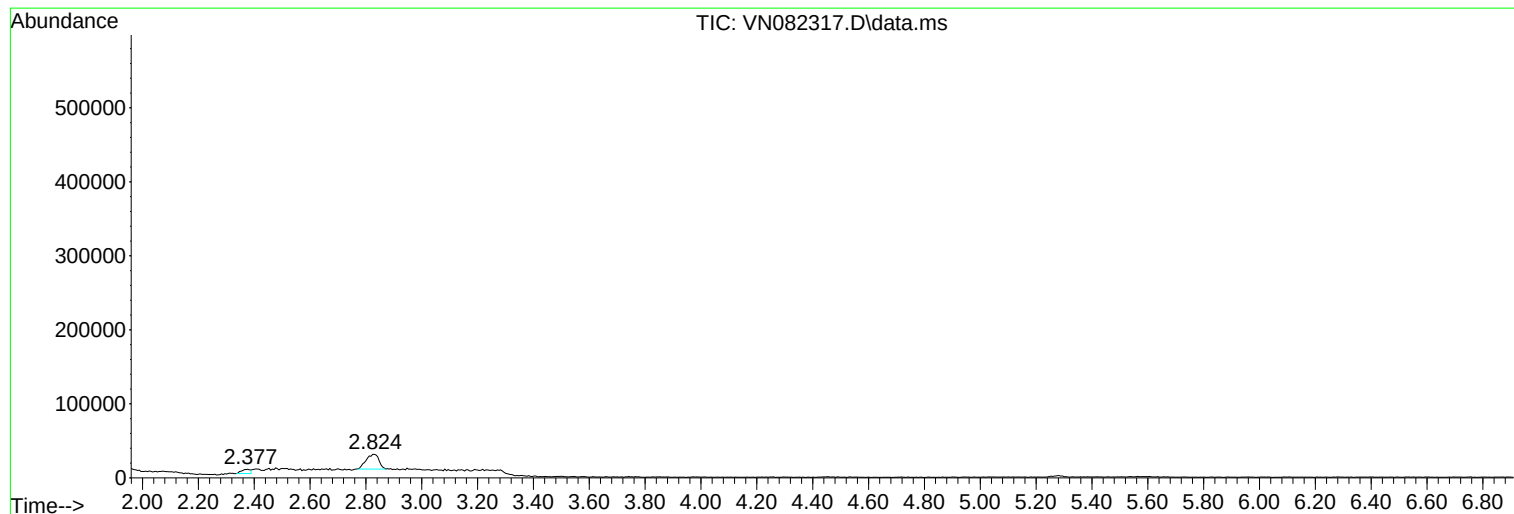
Sum of corrected areas: 6412217

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Instrument :
MSVOA_N
ClientSampleId :
COMP-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051524W.M
Quant Title : SW846 8260

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



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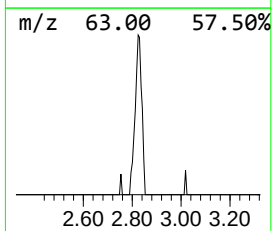
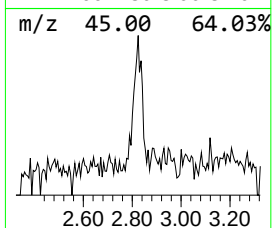
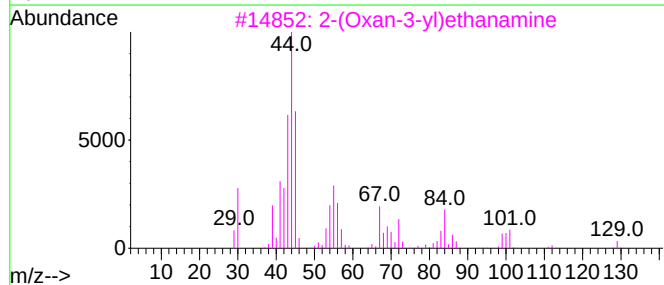
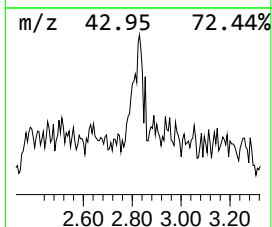
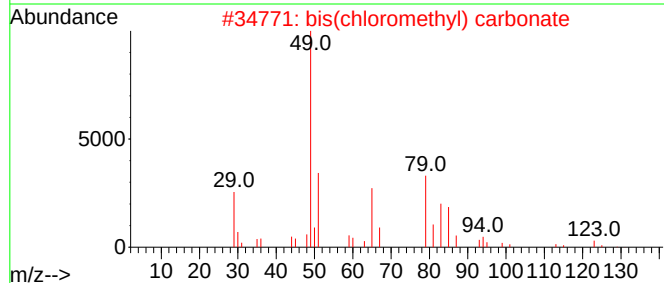
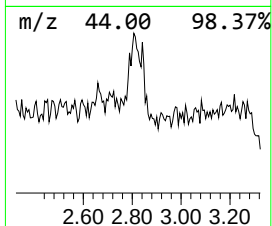
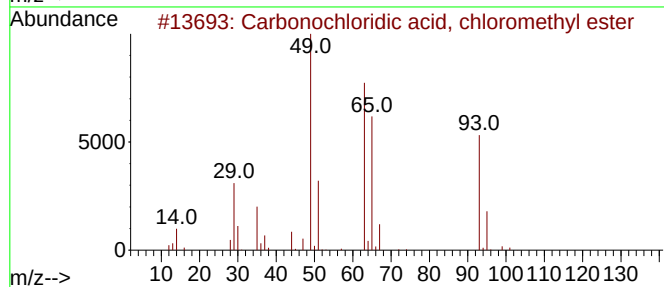
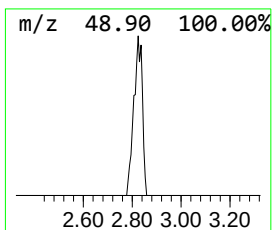
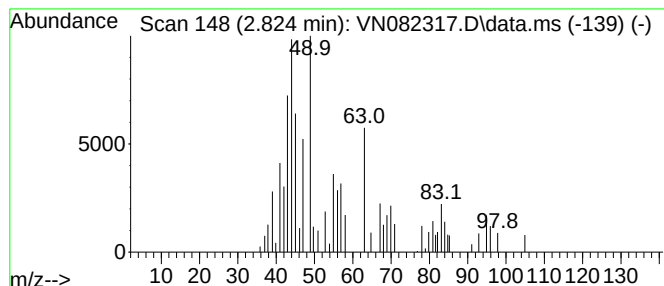
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051524W.M
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown2.824 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.824	6.17 ug/l	64323	Pentafluorobenzene	8.230

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Carbonochloridic acid, chloromet...	128	C2H2Cl2O2	022128-62-7	36
2		bis(chloromethyl) carbonate	158	C3H4Cl2O3	1000404-37-0	17
3		2-(Oxan-3-yl)ethanamine	129	C7H15NO	1000411-07-5	10
4		Silane, methyl-	46	CH6Si	000992-94-9	9
5		Phosgene	98	CCl2O	000075-44-5	9



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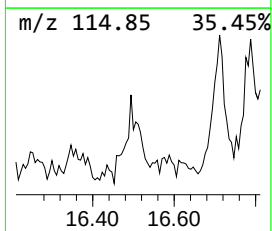
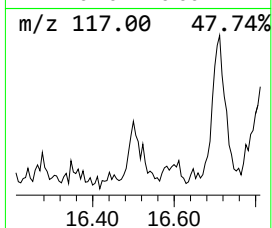
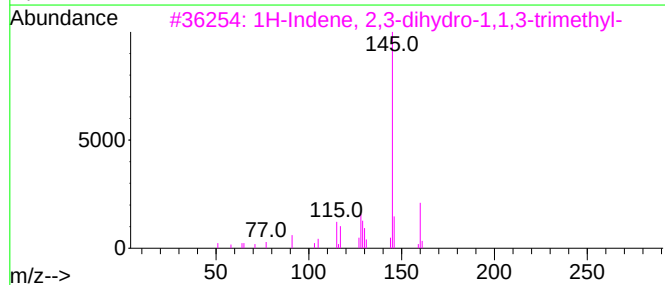
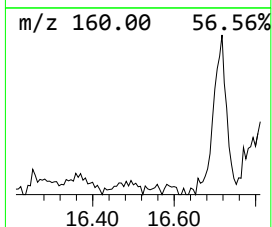
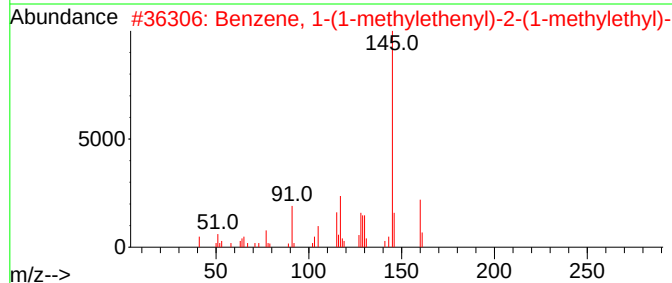
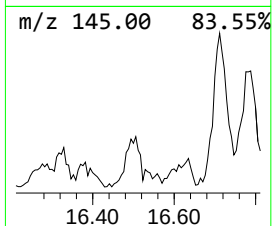
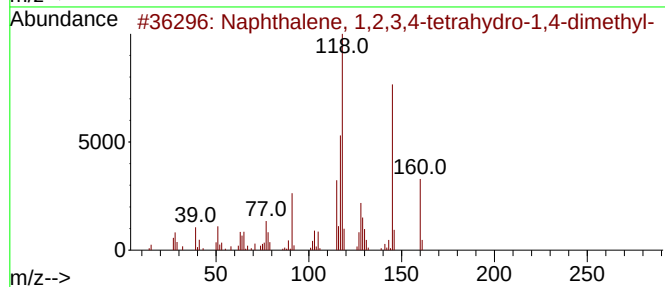
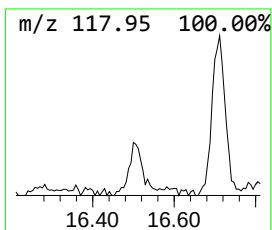
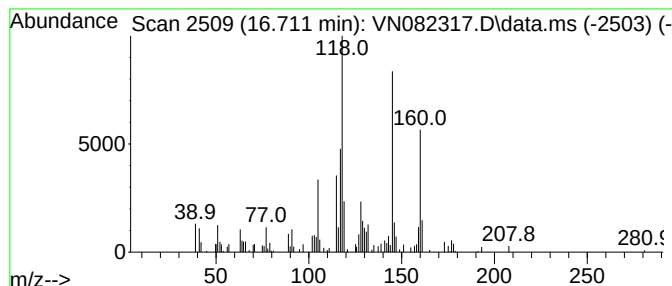
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Peak Number 2 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.711	6.19 ug/l	111089	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	97
2			Benzene, 1-(1-methylethenyl)-2-(...	160	C12H16	005557-93-7	90
3			1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	89
4			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	87
5			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	78



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown2.824	2.824	6.2	ug/l	64323	1	8.230	521631	50.0
Naphthalene, 1,...	16.711	6.2	ug/l	111089	4	13.794	897013	50.0