

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
 Data File : VU043618.D
 Acq On : 07 May 2021 17:26
 Operator : SY/MD
 Sample : M2303-01
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-18

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_U\Method\82U050621W.M
 Title : SW846 8260

Signal : TIC: VU043618.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.794	127	141	147	rBV6	7010	13373	1.54%	0.251%
2	2.630	393	401	413	rBV	5889	9466	1.09%	0.178%
3	2.761	431	442	459	rBV	15144	29907	3.43%	0.562%
4	5.298	1214	1231	1244	rBV	137430	290416	33.34%	5.454%
5	5.382	1244	1257	1278	rVB	195669	397821	45.67%	7.471%
6	5.710	1343	1359	1377	rBV	164448	333207	38.25%	6.257%
7	6.256	1515	1529	1548	rBV	308672	573902	65.88%	10.778%
8	7.906	2027	2042	2062	rBV	514040	871091	100.00%	16.359%
9	9.423	2496	2514	2533	rBV	440822	724180	83.13%	13.600%
10	10.639	2880	2892	2912	rBV	364946	576578	66.19%	10.828%
11	11.819	3247	3259	3275	rBV	439853	682644	78.37%	12.820%
12	12.102	3330	3347	3359	rBV3	53605	84659	9.72%	1.590%
13	12.192	3367	3375	3386	rVB7	8106	12328	1.42%	0.232%
14	12.700	3520	3533	3546	rBV7	14301	36514	4.19%	0.686%
15	12.873	3580	3587	3599	rVB2	8977	14596	1.68%	0.274%
16	12.980	3604	3620	3631	rBV2	21100	37506	4.31%	0.704%
17	13.118	3654	3663	3674	rVB6	7836	13038	1.50%	0.245%
18	13.259	3696	3707	3713	rBV2	9762	13975	1.60%	0.262%
19	13.301	3713	3720	3726	rBV	6533	8900	1.02%	0.167%
20	13.462	3761	3770	3783	rVB3	30396	50285	5.77%	0.944%
21	13.841	3880	3888	3897	rBV4	9372	14416	1.65%	0.271%
22	13.970	3922	3928	3936	rVB8	6057	9614	1.10%	0.181%
23	14.070	3946	3959	3967	rBV4	5998	11680	1.34%	0.219%
24	14.198	3989	3999	4009	rVV3	18653	28939	3.32%	0.543%
25	14.295	4013	4029	4040	rBV3	23147	42703	4.90%	0.802%
26	14.816	4182	4191	4199	rBV3	12488	20340	2.34%	0.382%
27	14.889	4205	4214	4226	rVB3	27619	46673	5.36%	0.877%
28	15.034	4250	4259	4269	rBV10	7080	11667	1.34%	0.219%
29	15.179	4293	4304	4311	rBV	24098	43218	4.96%	0.812%
30	15.282	4327	4336	4347	rVB6	9971	16849	1.93%	0.316%
31	15.420	4365	4379	4398	rBV3	79804	159434	18.30%	2.994%
32	15.934	4527	4539	4548	rBV4	9806	19090	2.19%	0.359%
33	16.198	4616	4621	4630	rVB9	6603	10268	1.18%	0.193%
34	16.410	4679	4687	4696	rBV10	5446	9486	1.09%	0.178%
35	16.619	4739	4752	4757	rBV4	14362	24539	2.82%	0.461%
36	16.655	4758	4763	4776	rVB3	16278	25572	2.94%	0.480%

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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

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Title : SW846 8260

37	16.793	4797	4806	4816	rBV2	19930	32329	3.71%	0.607%
38	17.105	4894	4903	4916	rVB4	6840	13535	1.55%	0.254%
39	17.549	5034	5041	5051	rVB5	6431	10188	1.17%	0.191%

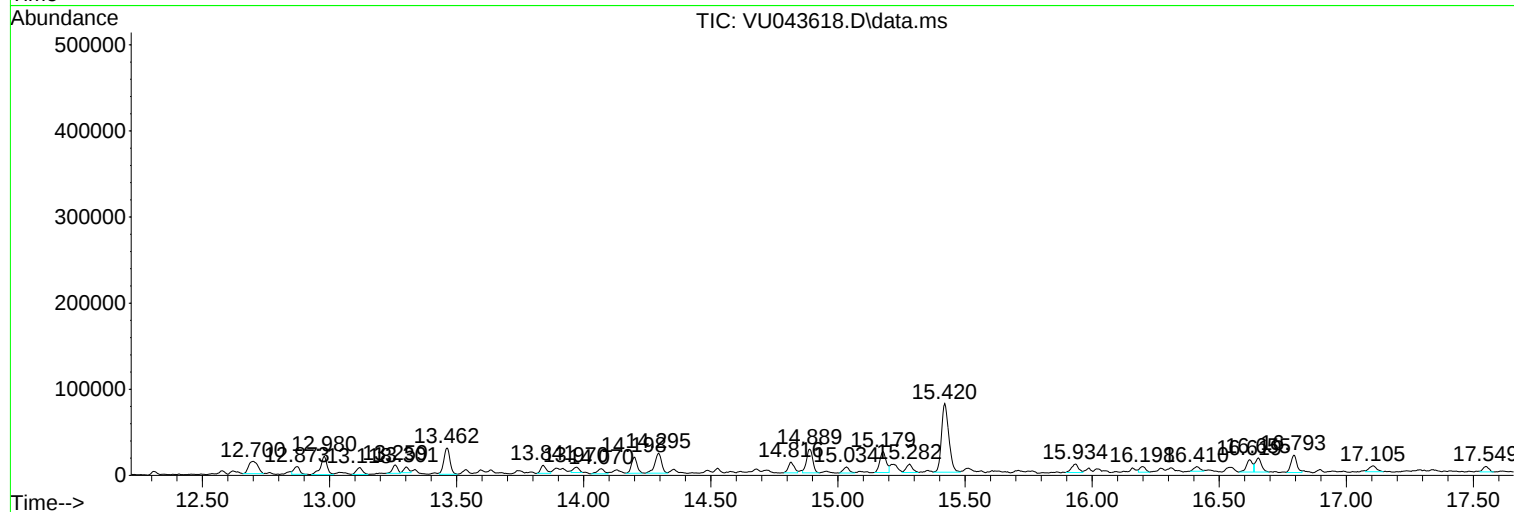
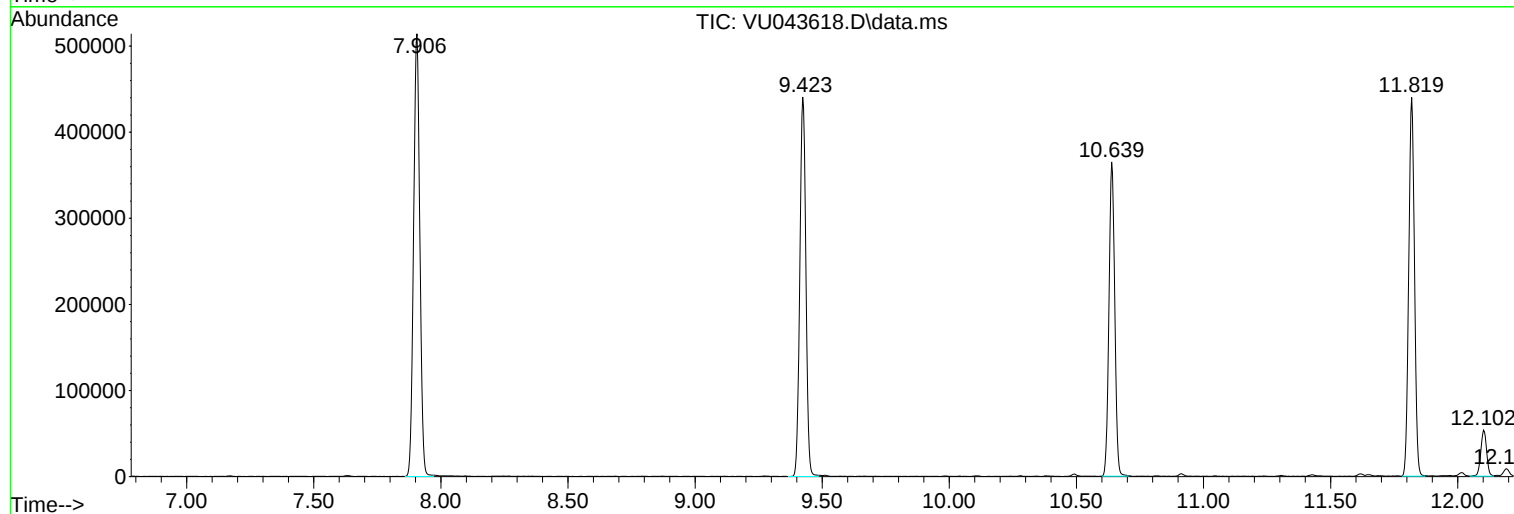
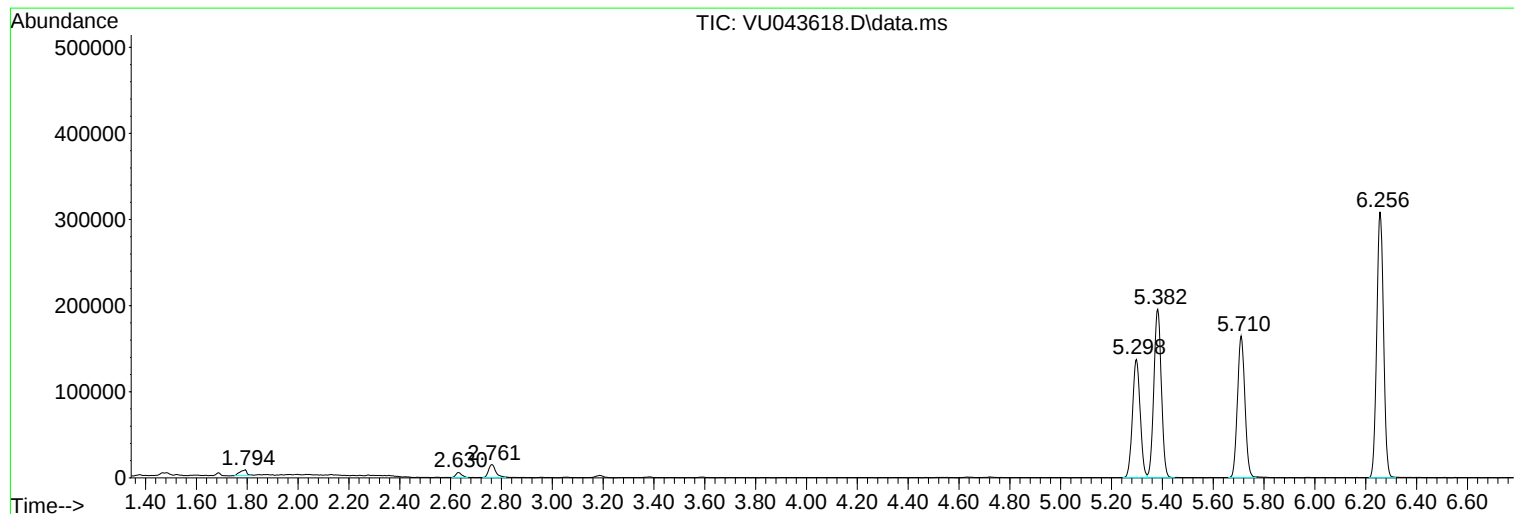
Sum of corrected areas: 5324926

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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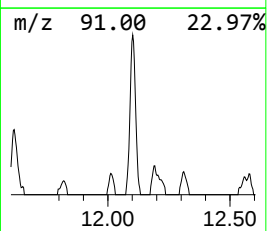
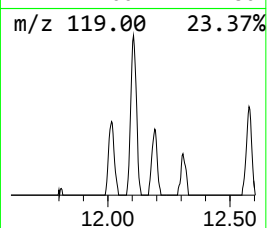
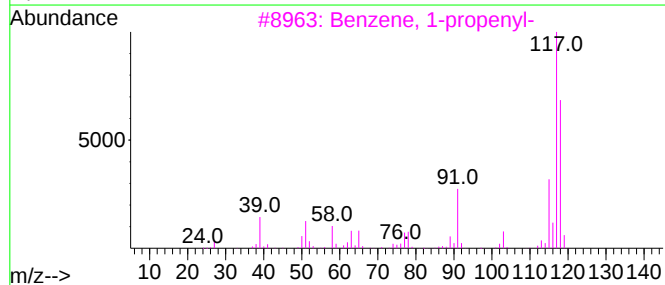
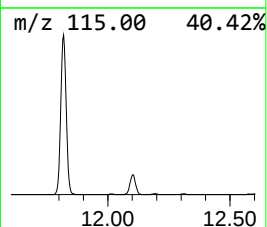
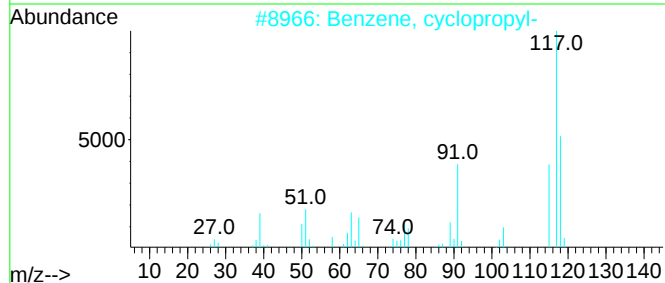
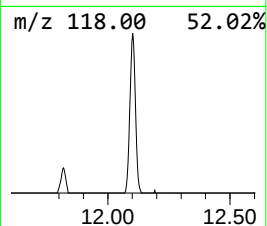
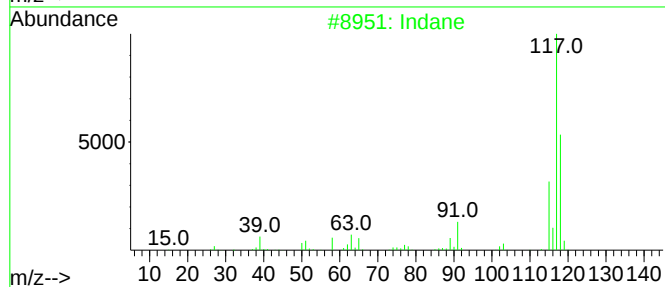
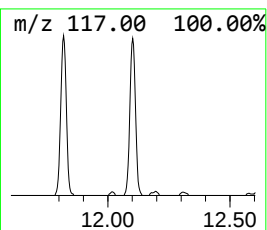
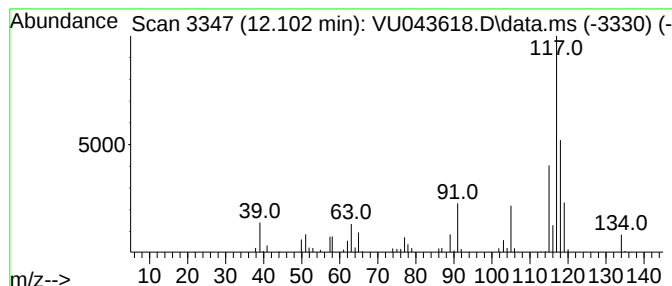
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.102	6.20 ug/l	84659	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	64
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	60
3			Benzene, 1-propenyl-	118	C9H10	000637-50-3	53
4			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	49
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	46



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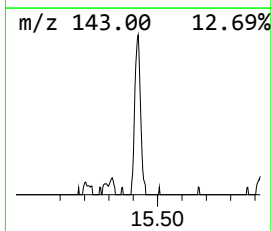
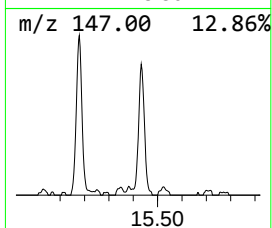
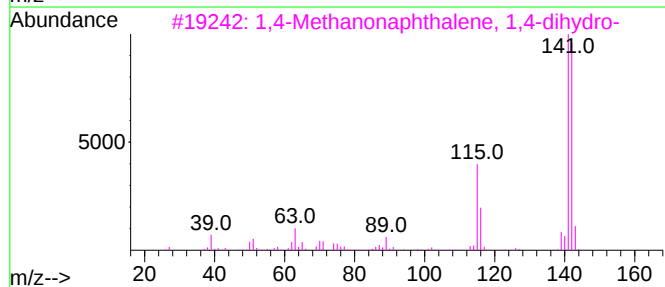
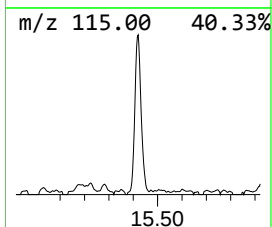
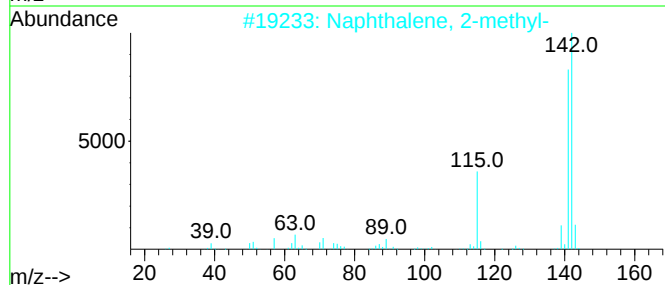
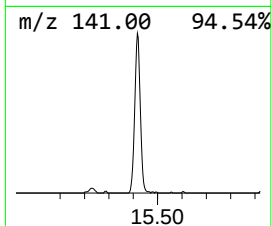
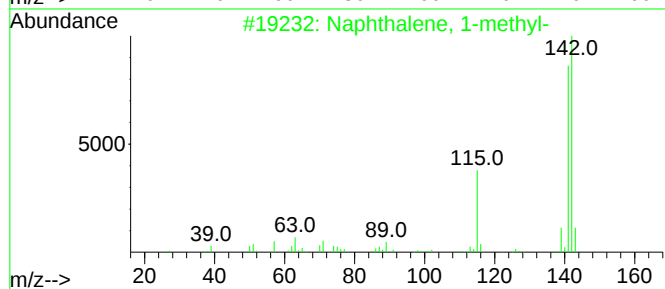
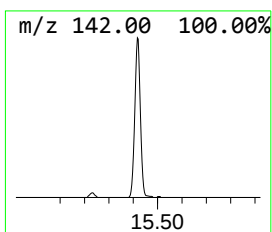
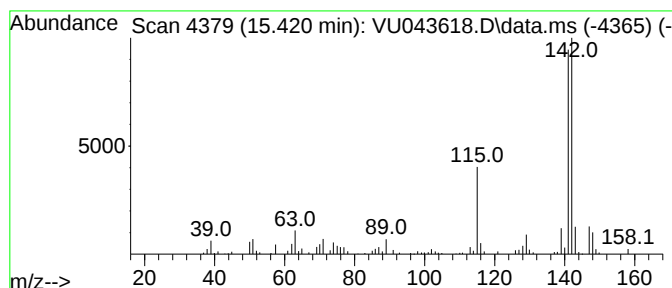
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Peak Number 2 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.420	11.68 ug/l	159434	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91
5			Benzocycloheptatriene	142	C11H10	000264-09-5	90



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Indane	12.102	6.2	ug/l	84659	4	11.819	682644	50.0
Naphthalene, 1-...	15.420	11.7	ug/l	159434	4	11.819	682644	50.0