LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN051820\
Data File : VN061468.D
Acq On : 19 May 2020 8:29
Operator : JC/MD
Sample : L2645-06
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-051320

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\82N051820W.M
Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.995	60	69	102	rBV	2533728	3962050	4.09%	0.367%
2	2.156	105	119	138	rVV	11563380	21171406	21.86%	1.964%
3	2.246	139	147	170	rVV2	1109375	2138608	2.21%	0.198%
4	2.352	172	180	212	rVB	813330	1476671	1.52%	0.137%
5	2.632	254	267	293	rBV	944209	1906015	1.97%	0.177%
6	2.770	293	310	353	rBV	23229061	60958437	62.95%	5.654%
7	3.024	375	389	399	rBV	2037516	4092499	4.23%	0.380%
8	3.101	399	413	450	rVB	10015872	23902612	24.68%	2.217%
9	3.285	452	470	493	rBV	7228825	16053611	16.58%	1.489%
10	3.439	495	518	531	rVV	3652920	8625131	8.91%	0.800%
11	3.526	531	545	580	rVB	4682474	11538174	11.92%	1.070%
12	3.764	597	619	663	rVB4	667549	2176728	2.25%	0.202%
13	4.384	777	812	827	rBV4	1690834	6072683	6.27%	0.563%
14	4.487	828	844	847	rBV3	1571654	3607874	3.73%	0.335%
15	4.998	967	1003	1042	rVB3	3712992	13737295	14.19%	1.274%
16	5.339	1080	1109	1137	rBV5	481244	2123143	2.19%	0.197%
17	5.519	1140	1165	1195	rBV	1743706	5491777	5.67%	0.509%
18	5.693	1197	1219	1234	rBV	730603	2230810	2.30%	0.207%
19	5.812	1235	1256	1265	rBV	1287263	4110150	4.24%	0.381%
20	6.021	1300	1321	1335	rBV	962701	3118020	3.22%	0.289%
21	6.114	1336	1350	1356	rBV	801344	2116745	2.19%	0.196%
22	6.326	1397	1416	1445	rVB	1180583	3646923	3.77%	0.338%
23	6.580	1474	1495	1515	rBV	7484139	23647641	24.42%	2.193%
24	7.339	1696	1731	1752	rBV2	1648494	5337608	5.51%	0.495%
25	7.616	1795	1817	1833	rBV5	2791398	8578784	8.86%	0.796%
26	7.706	1835	1845	1868	rVB4	819499	2442597	2.52%	0.227%
27	7.838	1868	1886	1899	rBV	668156	1649835	1.70%	0.153%
28	8.005	1917	1938	1955	rBV	18059280	46896741	48.43%	4.349%
29	8.098	1955	1967	1974	rBV2	2085076	4991781	5.15%	0.463%
30	8.265	2010	2019	2041	rVB2	584812	1484853	1.53%	0.138%
31	8.403	2041	2062	2088	rVB5	385919	1503819	1.55%	0.139%
32	8.548	2088	2107	2132	rBV2	1496863	4211607	4.35%	0.391%
33	9.011	2227	2251	2254	rBV3	1128563	2016618	2.08%	0.187%
34	9.047	2256	2262	2276	rVB	1912922	3817382	3.94%	0.354%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN051820\
Data File : VN061468.D
Acq On : 19 May 2020 8:29
Operator : JC/MD
Sample : L2645-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-051320

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\82N051820W.M
Title : SW846 8260

35	9.583	2416	2429	2438	rBV	1325502	2635246	2.72%	0.244%
36	9.825	2489	2504	2508	rBV	738906	1331362	1.37%	0.123%
37	9.863	2508	2516	2529	rVB	2231349	4135611	4.27%	0.384%
38	9.940	2529	2540	2564	rVB3	560010	1392913	1.44%	0.129%
39	10.063	2564	2578	2585	rBV2	1166432	2328506	2.40%	0.216%
40	10.130	2585	2599	2612	rBV	42270441	96834415	100.00%	8.981%
41	11.265	2935	2952	2973	rVB	1059649	2220362	2.29%	0.206%
42	11.397	2975	2993	3007	rBV2	1656560	4589982	4.74%	0.426%
43	11.490	3007	3022	3039	rBV	44058048	85793678	88.60%	7.957%
44	11.599	3039	3056	3058	rBV2	54488665	92653048	95.68%	8.593%
45	11.924	3133	3157	3174	rBV	42407256	79296028	81.89%	7.354%
46	12.223	3239	3250	3262	rVB	4128500	6660202	6.88%	0.618%
47	12.378	3285	3298	3319	rVB4	1397943	2610666	2.70%	0.242%
48	12.567	3343	3357	3366	rBV	9029171	15926998	16.45%	1.477%
49	12.632	3366	3377	3384	rBV	28609756	51081651	52.75%	4.738%
50	12.709	3394	3401	3421	rVB	16113651	27134111	28.02%	2.517%
51	12.866	3437	3450	3477	rVB	14972054	26234546	27.09%	2.433%
52	13.017	3479	3497	3510	rBV	48923967	87389512	90.25%	8.105%
53	13.143	3522	3536	3547	rVB3	676758	1778557	1.84%	0.165%
54	13.210	3547	3557	3566	rBV	954942	1468774	1.52%	0.136%
55	13.349	3581	3600	3627	rVB	16503588	30895785	31.91%	2.865%
56	13.519	3628	3653	3661	rVV	19734976	39235281	40.52%	3.639%
57	13.561	3661	3666	3684	rVV2	5982509	10693093	11.04%	0.992%
58	13.709	3700	3712	3723	rVV2	2373923	5213333	5.38%	0.484%
59	13.776	3723	3733	3738	rVV	3763766	5993494	6.19%	0.556%
60	13.805	3739	3742	3751	rVV	3038195	4103861	4.24%	0.381%
61	13.863	3751	3760	3770	rVV	5823486	9167056	9.47%	0.850%
62	13.921	3770	3778	3784	rVV	992404	1508823	1.56%	0.140%
63	13.969	3784	3793	3812	rVB2	3568147	5789757	5.98%	0.537%
64	14.101	3824	3834	3845	rVB	1480483	2336156	2.41%	0.217%
65	14.185	3845	3860	3866	rBV	3398831	5484820	5.66%	0.509%
66	14.223	3866	3872	3881	rVB	4174070	6238838	6.44%	0.579%
67	14.451	3933	3943	3953	rBV	3967011	6175382	6.38%	0.573%
68	14.567	3965	3979	3991	rBV3	6880978	13178210	13.61%	1.222%
69	14.632	3991	3999	4010	rVV3	745691	1235520	1.28%	0.115%
70	14.715	4014	4025	4041	rVV	1499706	2513622	2.60%	0.233%
71	14.828	4041	4060	4079	rVV4	606599	1915527	1.98%	0.178%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN051820\
Data File : VN061468.D
Acq On : 19 May 2020 8:29
Operator : JC/MD
Sample : L2645-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-051320

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\82N051820W.M
Title : SW846 8260

72	14.931	4080	4092	4106	rVB2	475508	1001047	1.03%	0.093%
73	15.104	4128	4146	4160	rBV	9307252	17549604	18.12%	1.628%
74	16.123	4450	4463	4487	rVB	1168531	2413242	2.49%	0.224%
75	16.313	4508	4522	4539	rVB	596367	1256584	1.30%	0.117%

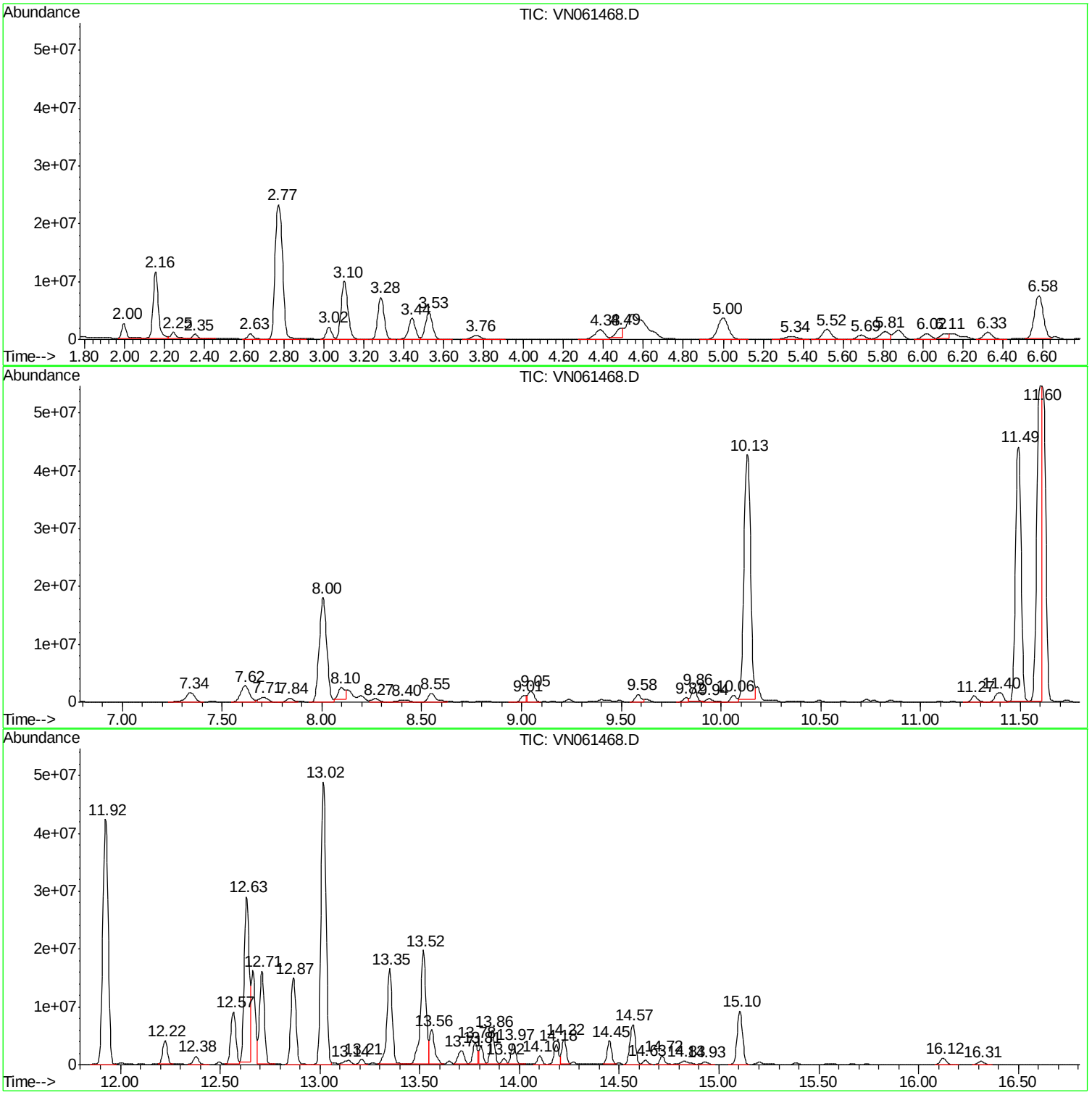
Sum of corrected areas: 1078231831

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN051820\
Data File : VN061468.D
Acq On : 19 May 2020 8:29
Operator : JC/MD
Sample : L2645-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-051320

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N051820W.M
Quant Title : SW846 8260

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



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN051820\
Data File : VN061468.D
Acq On : 19 May 2020 8:29
Operator : JC/MD
Sample : L2645-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-051320

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N051820W.M
Quant Title : SW846 8260

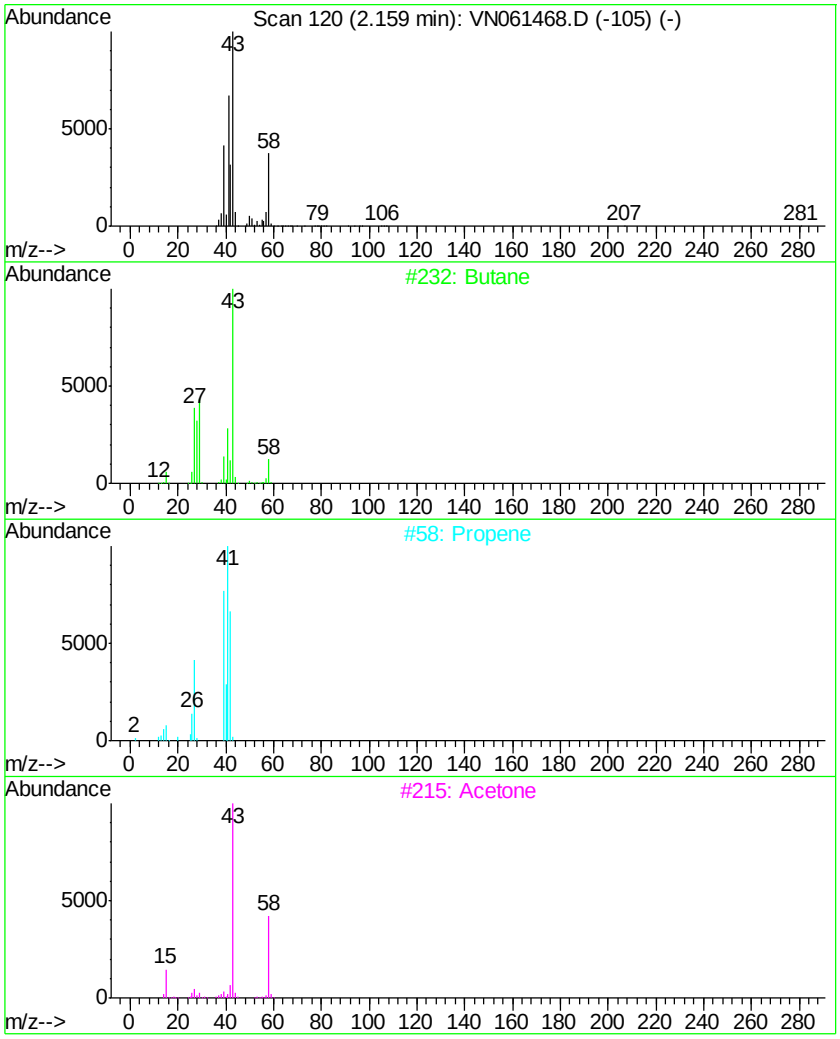
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

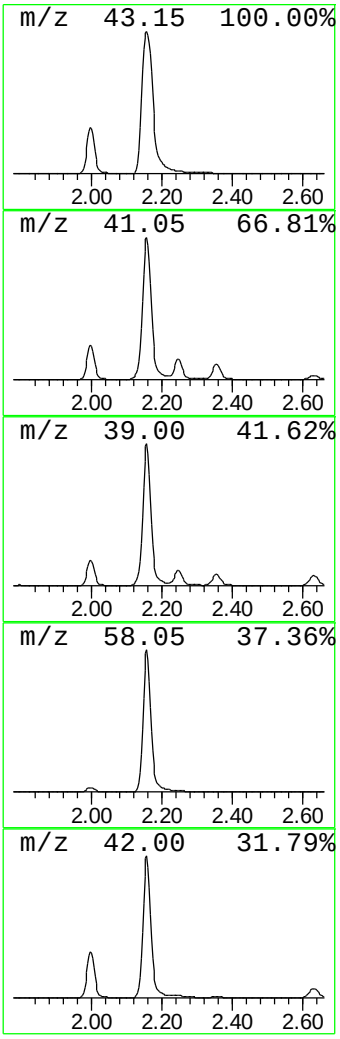
****

Peak Number 1 unknown2.16 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.16	123.39 ug/l	21171400	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane	58	C4H10	000106-97-8	45
2	Propene	42	C3H6	000115-07-1	9
3	Acetone	58	C3H6O	000067-64-1	5
4	Diazene, dimethyl-	58	C2H6N2	000503-28-6	5
5	Propylene oxide	58	C3H6O	000075-56-9	4





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN051820\
Data File : VN061468.D
Acq On : 19 May 2020 8:29
Operator : JC/MD
Sample : L2645-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-051320

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N051820W.M
Quant Title : SW846 8260

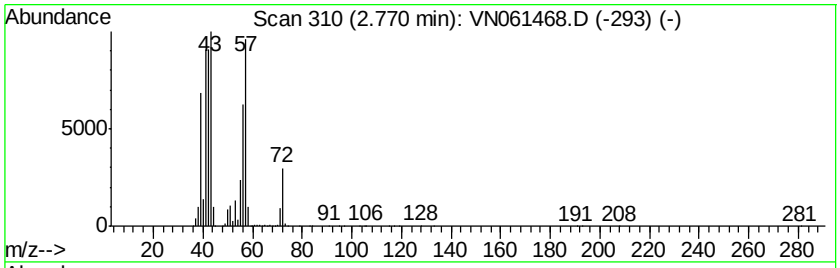
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

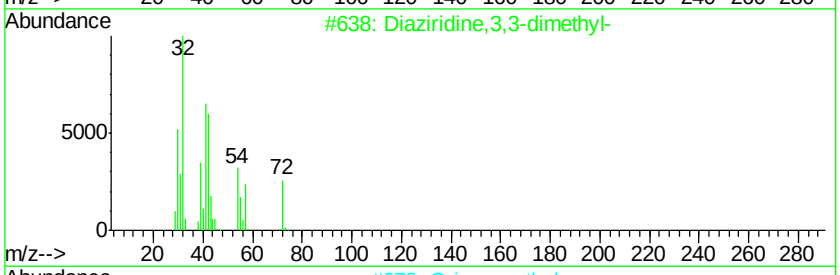
****

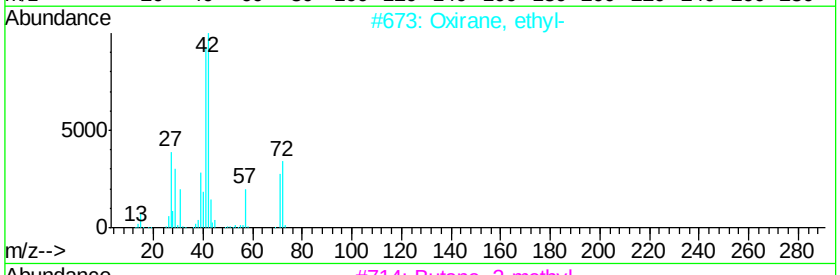
Peak Number 2 unknown2.77 Concentration Rank 1

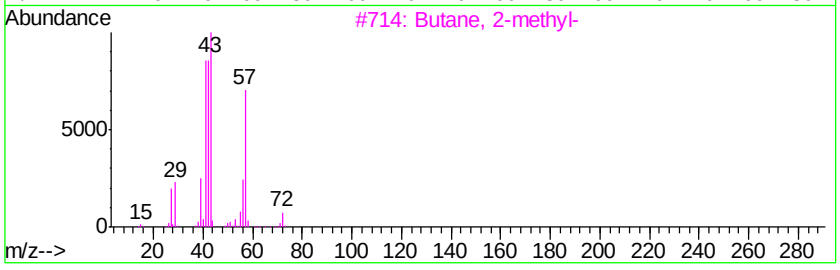
R.T.	EstConc	Area	Relative to ISTD	R.T.
2.77	355.29 ug/l	60958400	Pentafluorobenzene	7.63

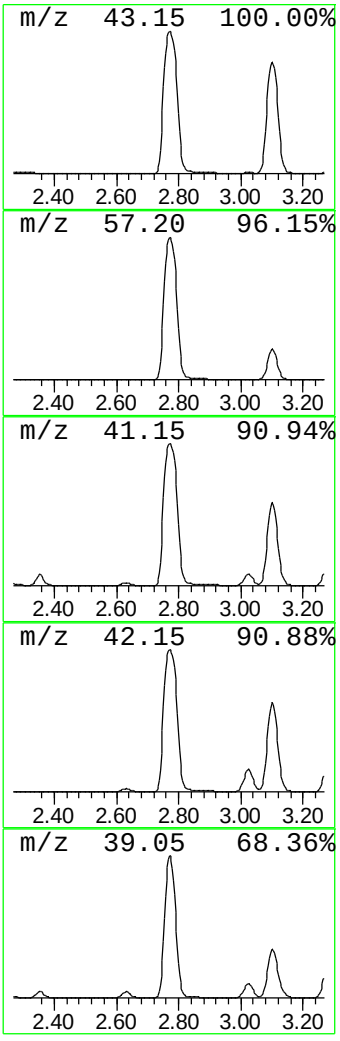
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	49
2	Oxirane, ethyl-	72	C4H8O	000106-88-7	35
3	Butane, 2-methyl-	72	C5H12	000078-78-4	32
4	2-Propen-1-ol, 2-methyl-	72	C4H8O	000513-42-8	30
5	2-Buten-1-ol	72	C4H8O	006117-91-5	27











Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN051820\
Data File : VN061468.D
Acq On : 19 May 2020 8:29
Operator : JC/MD
Sample : L2645-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-051320

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N051820W.M
Quant Title : SW846 8260

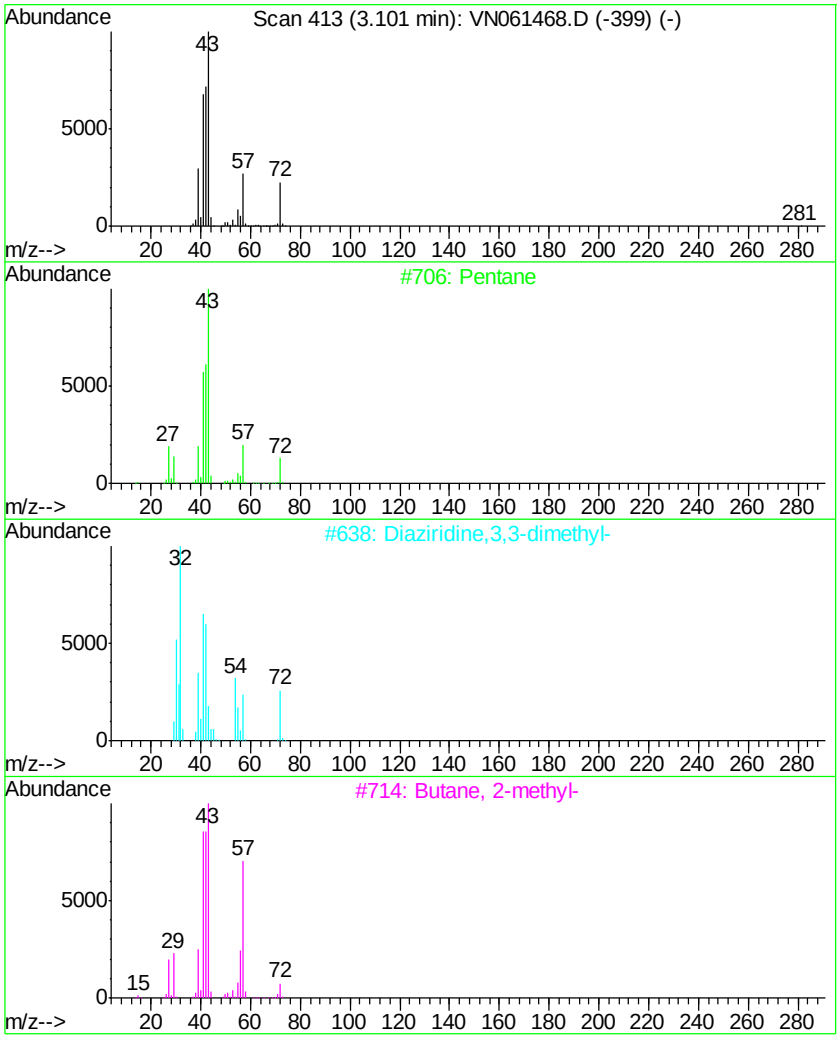
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

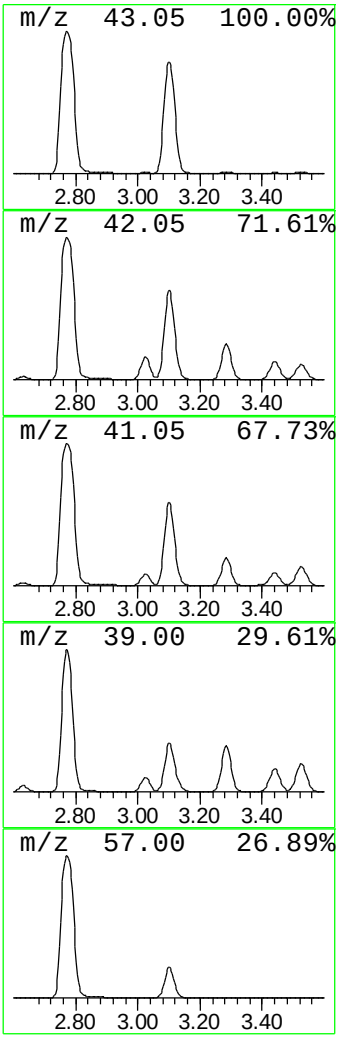
****

Peak Number 3 Pentane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.10	139.31 ug/l	23902600	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	91
2	Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	64
3	Butane, 2-methyl-	72	C5H12	000078-78-4	53
4	Oxirane, ethyl-	72	C4H8O	000106-88-7	53
5	Isobutylene epoxide	72	C4H8O	000558-30-5	50





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN051820\
Data File : VN061468.D
Acq On : 19 May 2020 8:29
Operator : JC/MD
Sample : L2645-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-051320

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N051820W.M
Quant Title : SW846 8260

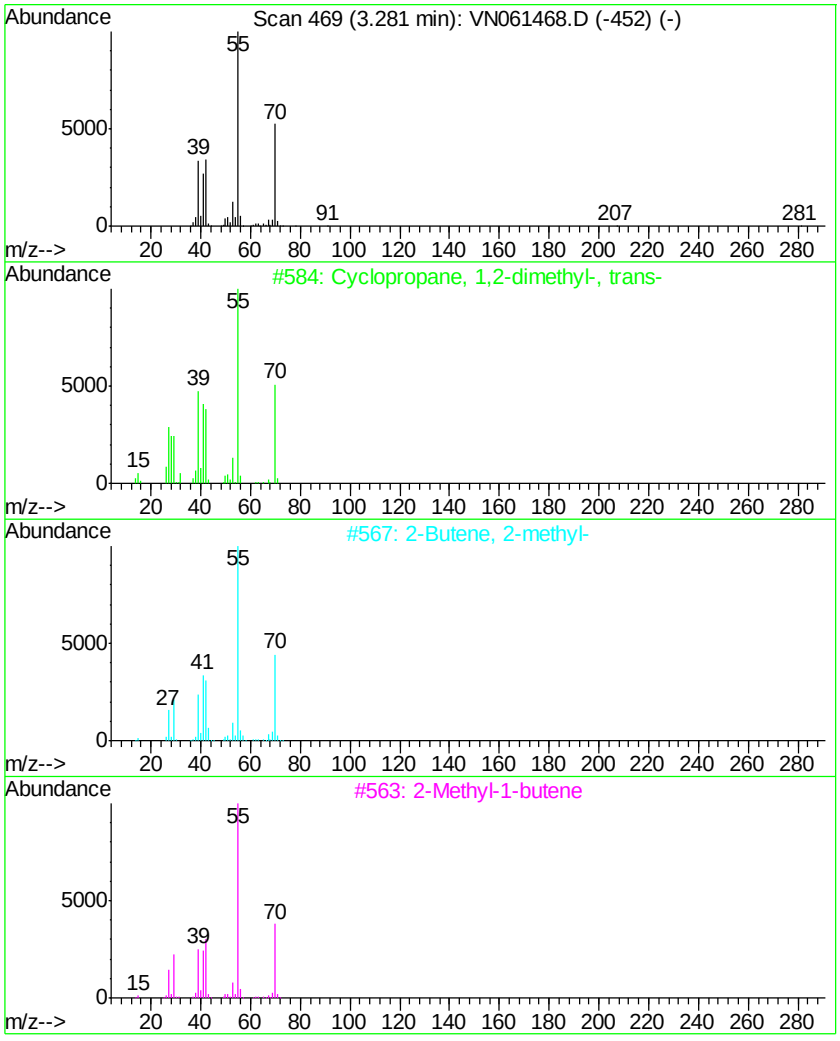
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

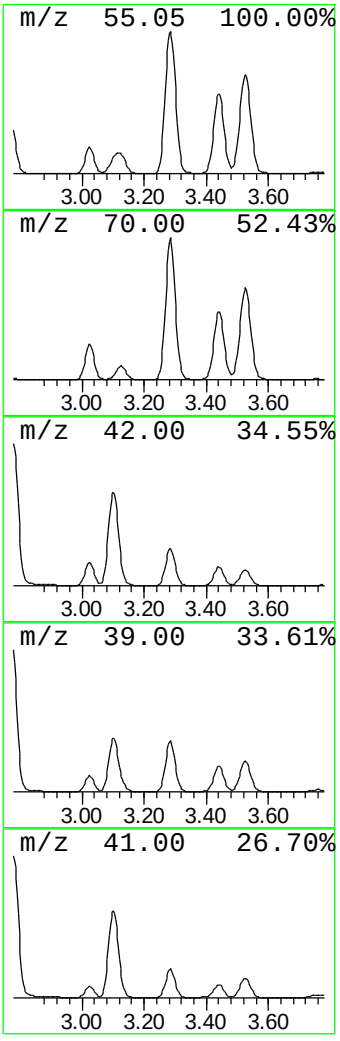
****

Peak Number 4 Cyclopropane, 1,2-dimethyl-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.28	93.57 ug/l	16053600	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	87
2	2-Butene, 2-methyl-	70	C5H10	000513-35-9	87
3	2-Methyl-1-butene	70	C5H10	000563-46-2	87
4	2-Pentene, (Z)-	70	C5H10	000627-20-3	87
5	2-Pentene, (E)-	70	C5H10	000646-04-8	87





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN051820\
Data File : VN061468.D
Acq On : 19 May 2020 8:29
Operator : JC/MD
Sample : L2645-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-051320

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N051820W.M
Quant Title : SW846 8260

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

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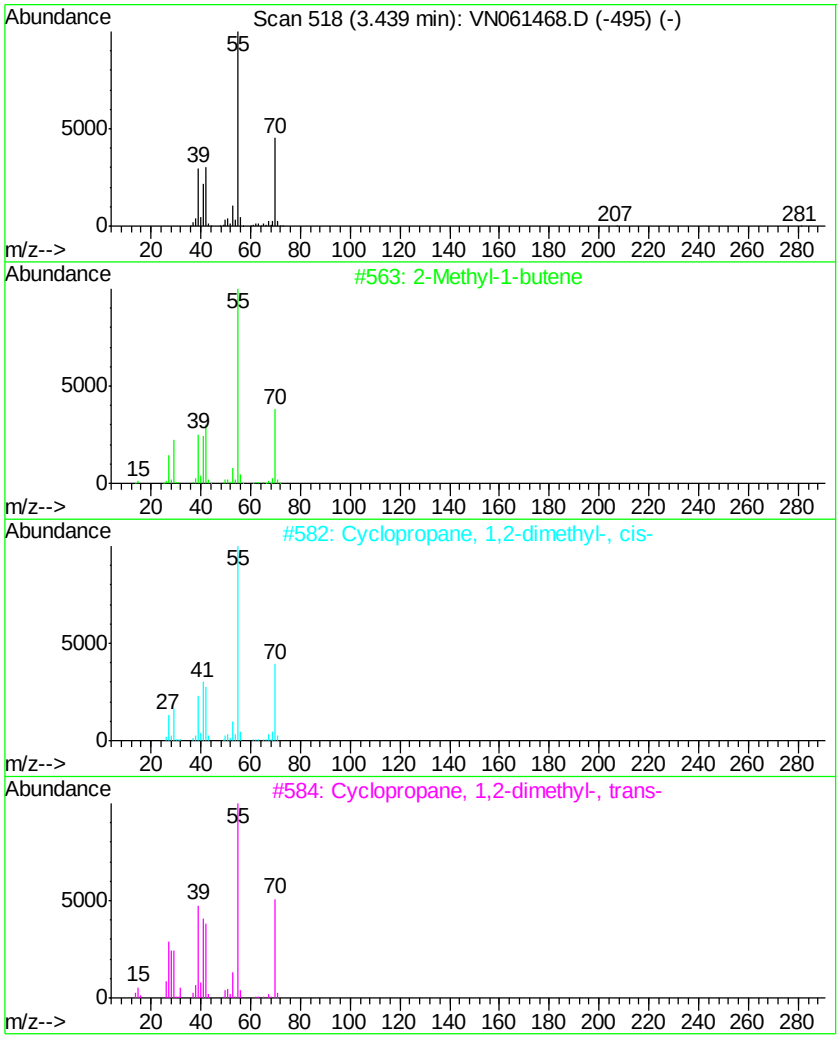
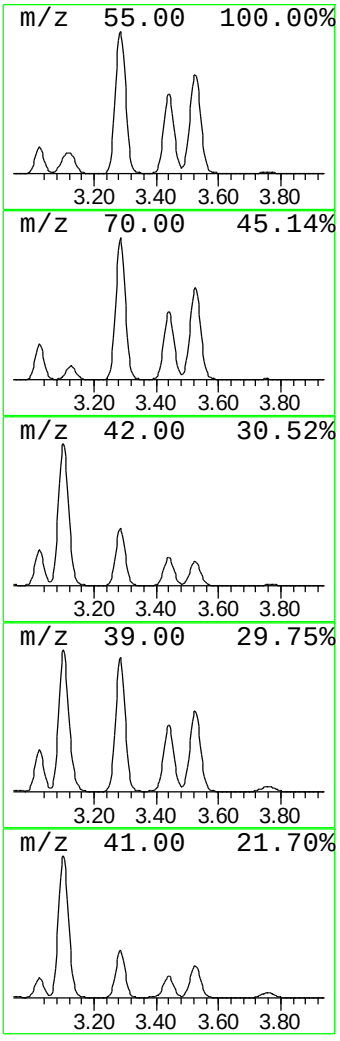
Peak Number 5 2-Methyl-1-butene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.44	50.27 ug/l	8625130	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-1-butene	70	C5H10	000563-46-2	91
2	Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	90
3	Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	87
4	2-Butene, 2-methyl-	70	C5H10	000513-35-9	87
5	2-Pentene, (Z)-	70	C5H10	000627-20-3	68





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN051820\
Data File : VN061468.D
Acq On : 19 May 2020 8:29
Operator : JC/MD
Sample : L2645-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-051320

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N051820W.M
Quant Title : SW846 8260

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

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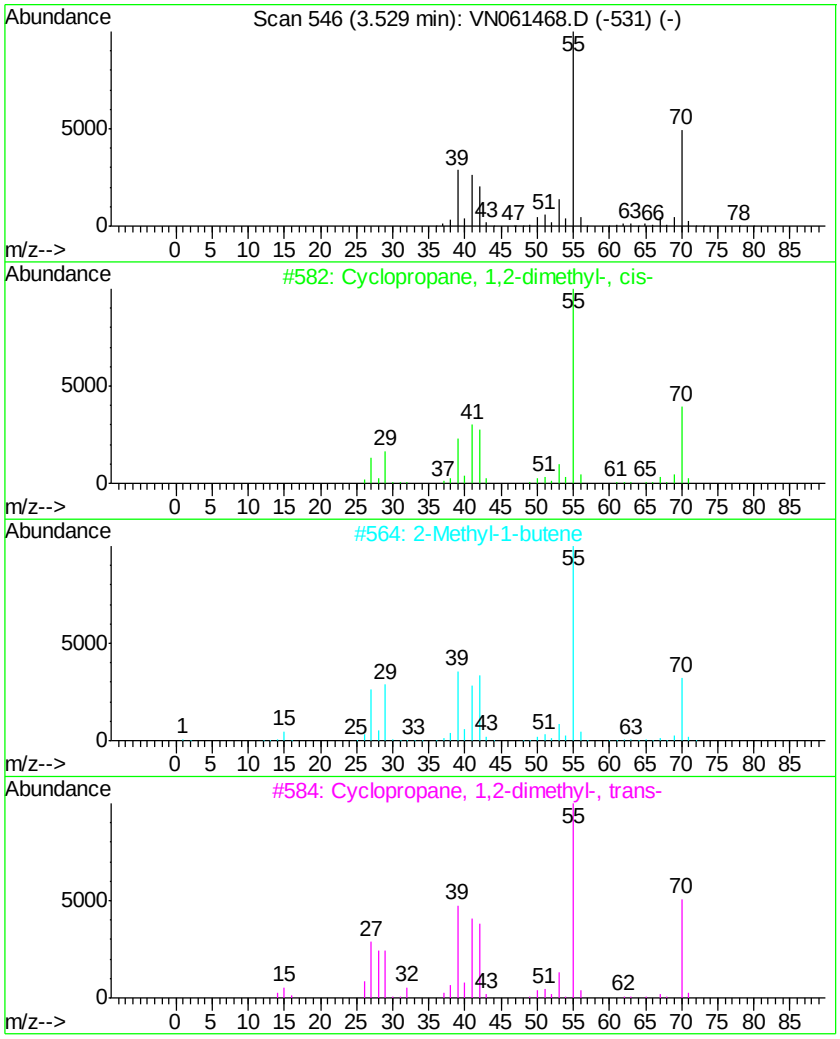
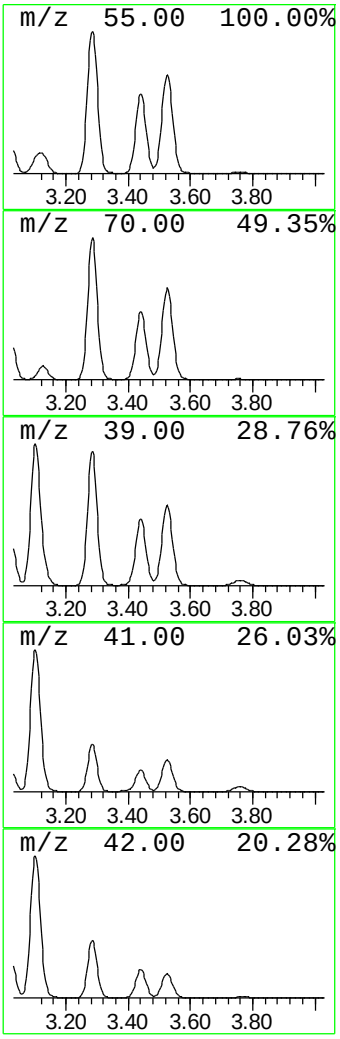
Peak Number 6 Cyclopropane, 1,2-dimethyl-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.53	67.25 ug/l	11538200	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	87
2	2-Methyl-1-butene	70	C5H10	000563-46-2	83
3	Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	83
4	Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	80
5	1-Butene, 3-methyl-	70	C5H10	000563-45-1	78





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN051820\
Data File : VN061468.D
Acq On : 19 May 2020 8:29
Operator : JC/MD
Sample : L2645-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-051320

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N051820W.M
Quant Title : SW846 8260

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

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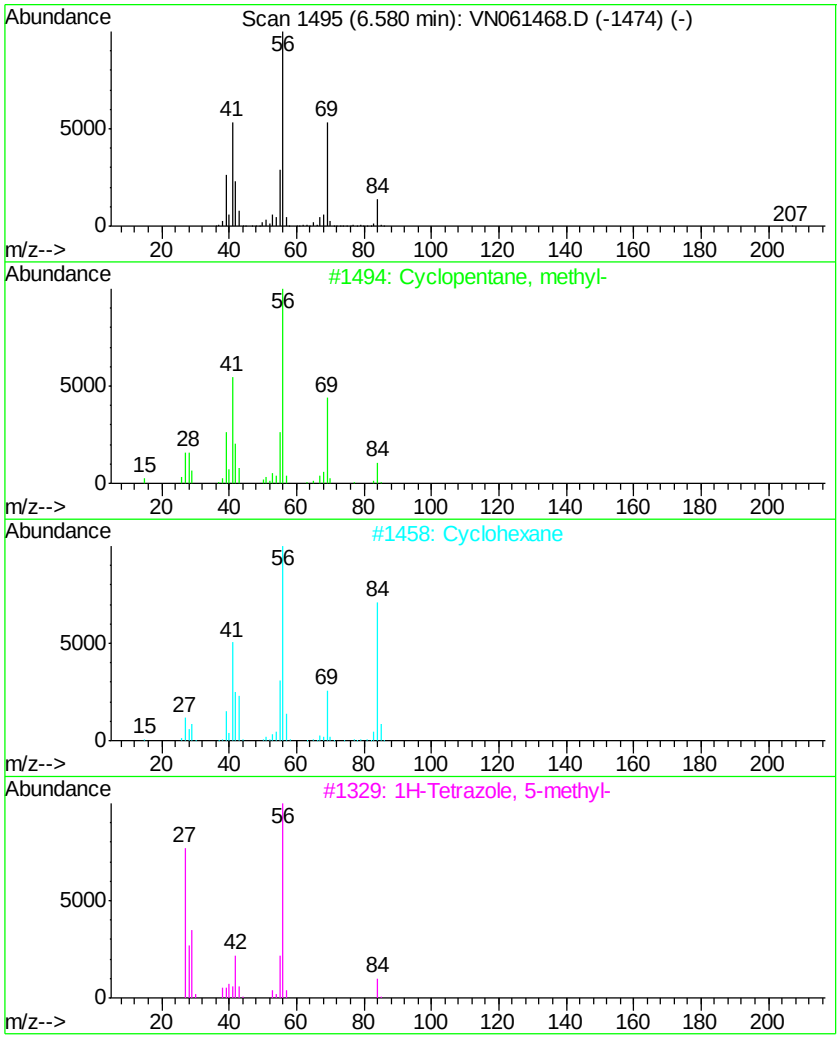
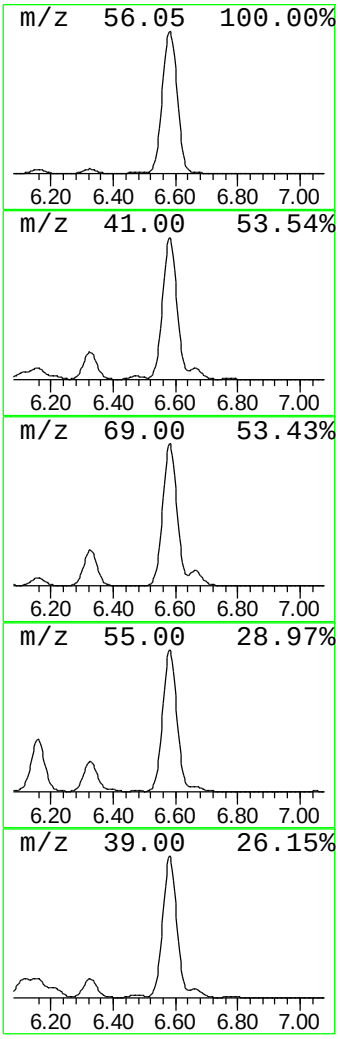
Peak Number 7 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	137.83 ug/l	23647600	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	Cyclohexane	84	C6H12	000110-82-7	78
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	72
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64





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Data File : VN061468.D
Acq On : 19 May 2020 8:29
Operator : JC/MD
Sample : L2645-06
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-051320

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N051820W.M
Quant Title : SW846 8260

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

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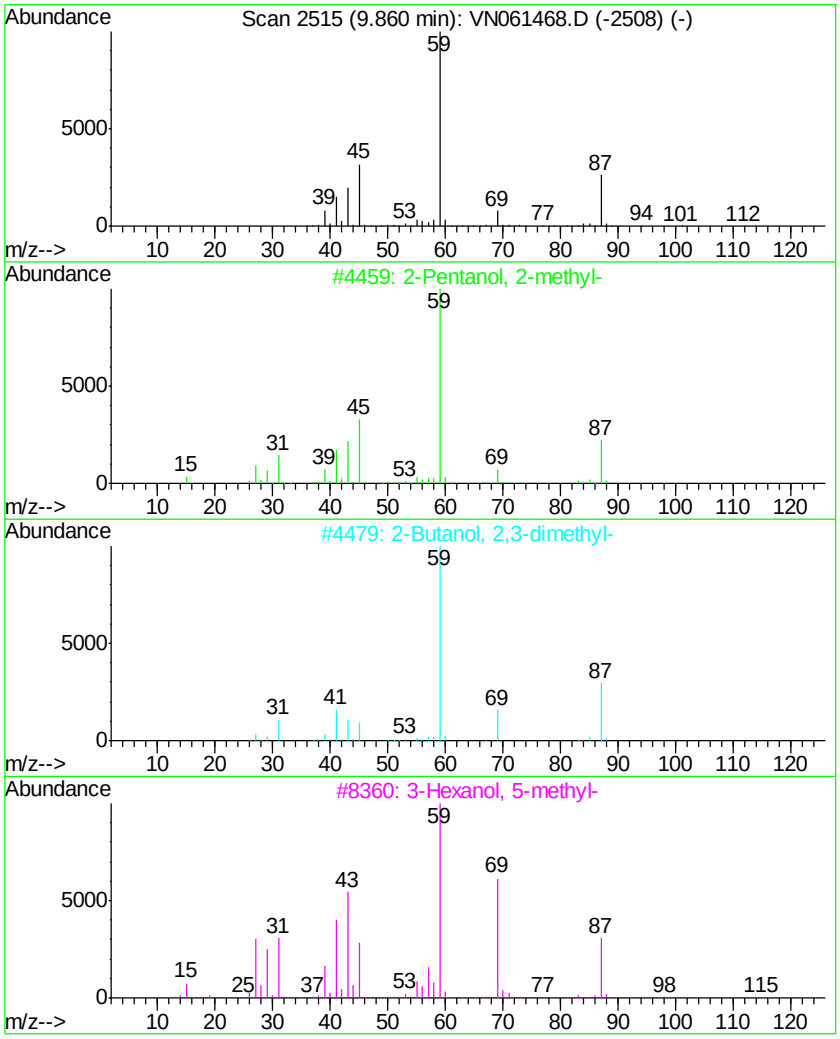
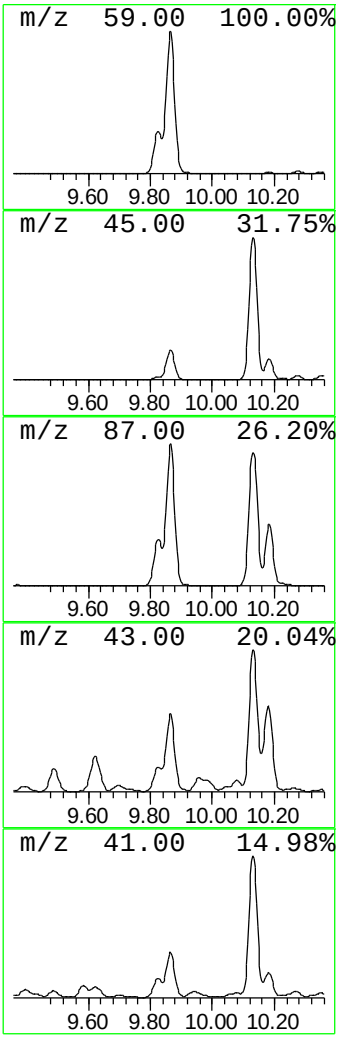
Peak Number 8 2-Pentanol, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.86	49.10 ug/l	4135610	1,4-Difluorobenzene	8.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	78
2	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	50
3	3-Hexanol, 5-methyl-	116	C7H16O	000623-55-2	43
4	Butane, 2-methoxy-	88	C5H12O	006795-87-5	40
5	1,2-Butanediol	90	C4H10O2	000584-03-2	38





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN051820\
Data File : VN061468.D
Acq On : 19 May 2020 8:29
Operator : JC/MD
Sample : L2645-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-051320

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N051820W.M
Quant Title : SW846 8260

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

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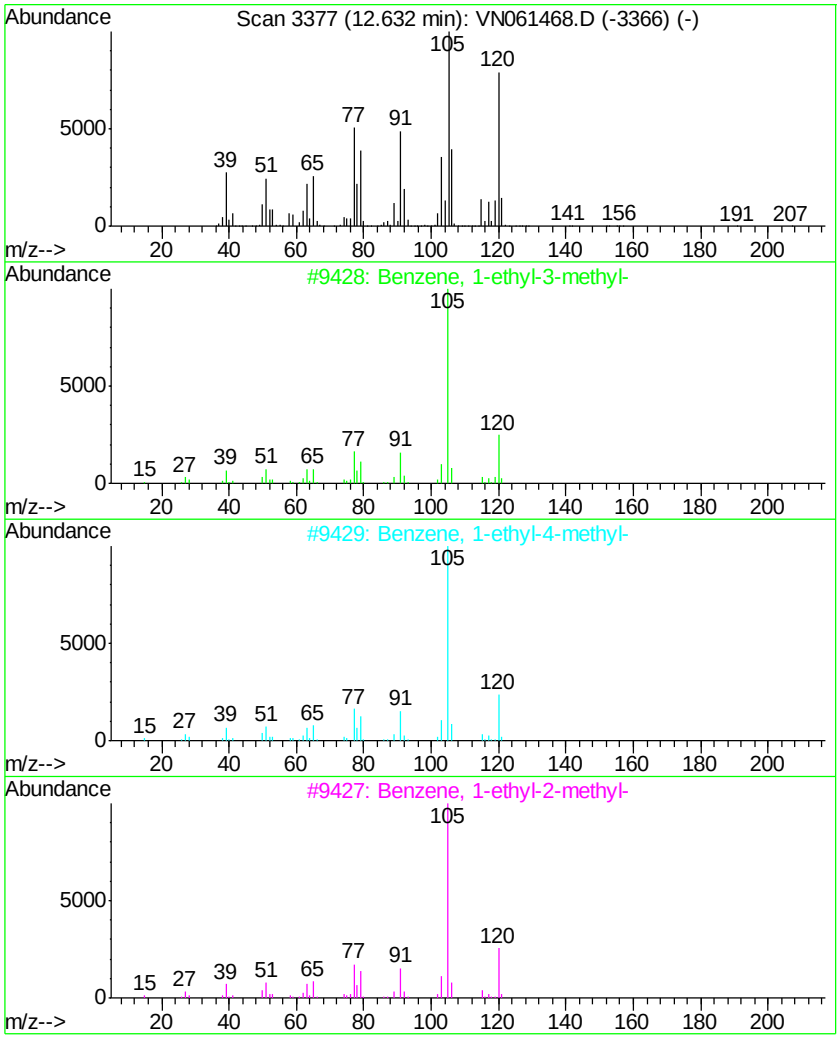
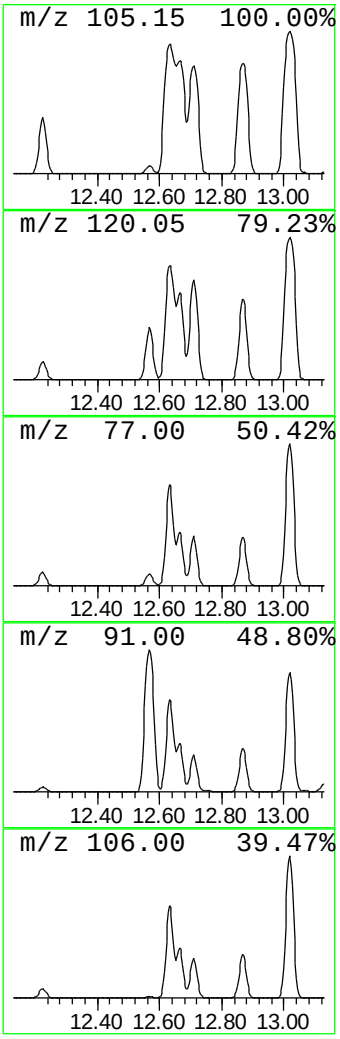
Peak Number 9 Benzene, 1-ethyl-3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.63	82.67 ug/l	51081700	1,4-Dichlorobenzene-d4	13.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	81
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	76
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	76
4	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	74
5	2,3-Heptadien-5-yne, 2,4-dimethyl-	120	C9H12	041898-89-9	72





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN051820\
Data File : VN061468.D
Acq On : 19 May 2020 8:29
Operator : JC/MD
Sample : L2645-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-051320

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N051820W.M
Quant Title : SW846 8260

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

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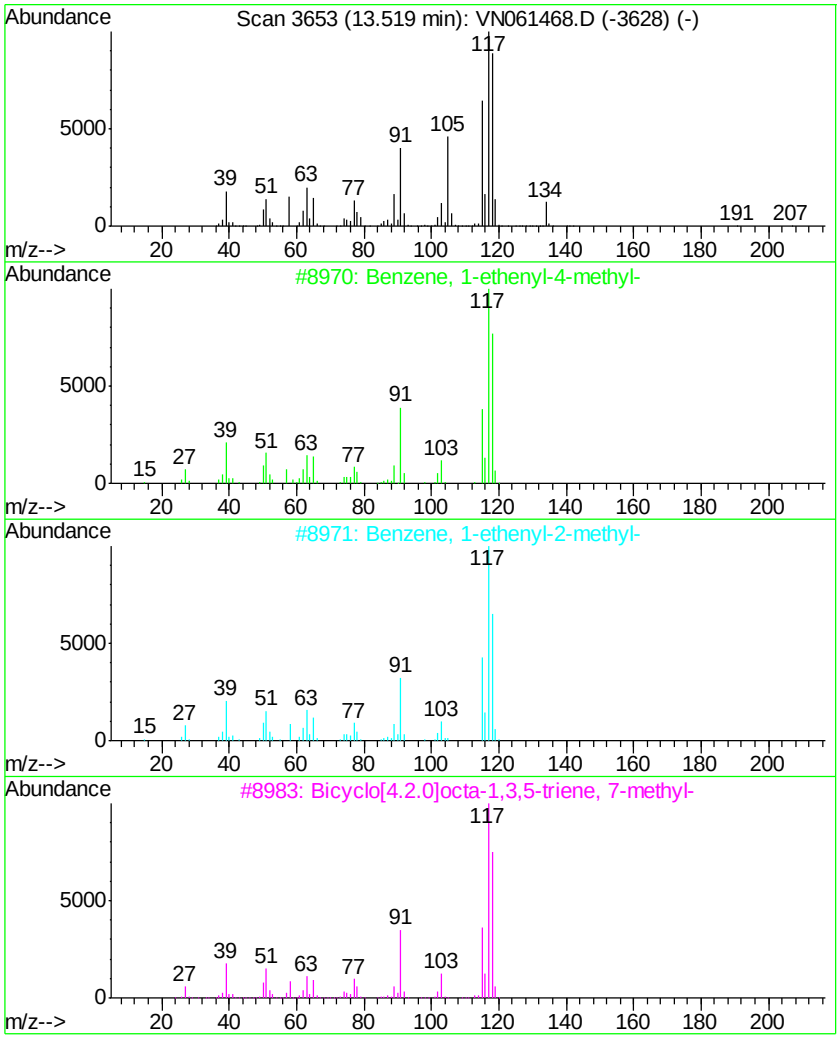
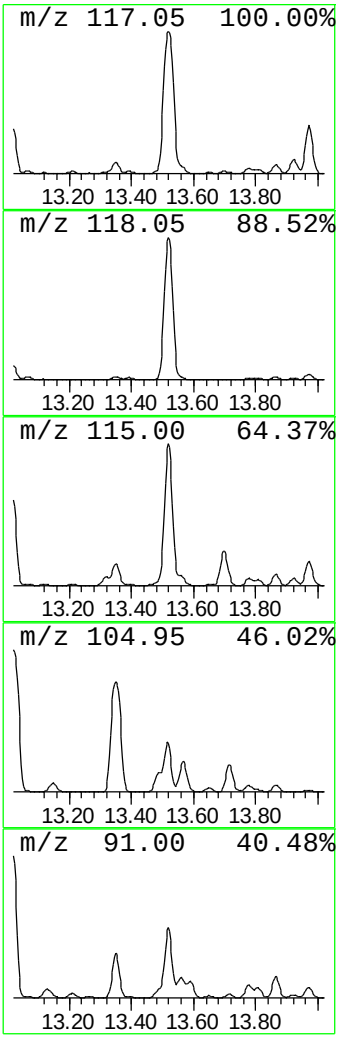
Peak Number 10 Benzene, 1-ethenyl-4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.52	63.50 ug/l	39235300	1,4-Dichlorobenzene-d4	13.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	96
2	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	96
3	Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	055337-80-9	94
4	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	94
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	94





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN051820\
Data File : VN061468.D
Acq On : 19 May 2020 8:29
Operator : JC/MD
Sample : L2645-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-051320

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N051820W.M
Quant Title : SW846 8260

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown2.16	2.16	123.4	ug/l	21171400	1	7.63	8578780	50.0
unknown2.77	2.77	355.3	ug/l	60958400	1	7.63	8578780	50.0
Pentane	3.10	139.3	ug/l	23902600	1	7.63	8578780	50.0
Cyclopropane, 1,2...	3.28	93.6	ug/l	16053600	1	7.63	8578780	50.0
2-Methyl-1-butene	3.44	50.3	ug/l	8625130	1	7.63	8578780	50.0
Cyclopropane, 1,2...	3.53	67.3	ug/l	11538200	1	7.63	8578780	50.0
Cyclopentane, met...	6.58	137.8	ug/l	23647600	1	7.63	8578780	50.0
2-Pentanol, 2-met...	9.86	49.1	ug/l	4135610	2	8.55	4211610	50.0
Benzene, 1-ethyl-...	12.63	82.7	ug/l	51081700	4	13.32	30895800	50.0
Benzene, 1-etheny...	13.52	63.5	ug/l	39235300	4	13.32	30895800	50.0