LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062922\
Data File : VU049548.D
Acq On : 30 Jun 2022 06:16
Operator : SY/MD
Sample : N3524-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
31466

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\82U062722W.M
Title : SW846 8260

Signal : TIC: VU049548.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.359	3	6	21	rVB	114266	105093	2.84%	0.261%
2	1.584	68	76	87	rBV2	74790	90744	2.45%	0.225%
3	1.687	101	108	118	rVB	108884	130675	3.53%	0.324%
4	2.543	364	374	386	rBV	36486	60973	1.65%	0.151%
5	2.633	389	402	440	rVB	379267	737872	19.95%	1.831%
6	3.198	560	578	624	rBV2	269167	1027077	27.77%	2.549%
7	3.578	683	696	714	rBV3	29047	67362	1.82%	0.167%
8	4.446	947	966	990	rBV4	45867	120344	3.25%	0.299%
9	4.706	1034	1047	1082	rVB	103036	245144	6.63%	0.608%
10	5.288	1211	1228	1241	rBV	208438	443527	11.99%	1.101%
11	5.372	1241	1254	1274	rVB	349366	714637	19.32%	1.773%
12	5.700	1339	1356	1375	rBV	225055	467484	12.64%	1.160%
13	5.838	1384	1399	1429	rBV2	142785	434798	11.76%	1.079%
14	5.964	1429	1438	1454	rVB5	23879	49838	1.35%	0.124%
15	6.150	1477	1496	1514	rBV	621252	1263048	34.15%	3.134%
16	6.247	1514	1526	1556	rVB	529634	1018074	27.53%	2.527%
17	6.385	1556	1569	1589	rBV	105427	213429	5.77%	0.530%
18	6.507	1595	1607	1628	rVB2	49938	108852	2.94%	0.270%
19	6.668	1638	1657	1672	rBV4	65129	187023	5.06%	0.464%
20	6.767	1674	1688	1714	rVB2	235944	502352	13.58%	1.247%
21	6.877	1714	1722	1732	rBV3	30252	52029	1.41%	0.129%
22	7.163	1786	1811	1835	rBV	578371	1242749	33.60%	3.084%
23	7.629	1944	1956	1966	rBV	132665	261910	7.08%	0.650%
24	7.693	1966	1976	1994	rVB	367400	711906	19.25%	1.767%
25	7.793	1994	2007	2011	rBV	428995	751805	20.33%	1.866%
26	7.896	2028	2039	2056	rVB	773633	1364830	36.91%	3.387%
27	8.028	2069	2080	2095	rBV	264116	506587	13.70%	1.257%
28	8.230	2133	2143	2163	rVB3	67411	165971	4.49%	0.412%
29	8.404	2186	2197	2210	rBV5	35461	95765	2.59%	0.238%
30	8.488	2210	2223	2236	rVV	1117946	1919311	51.90%	4.763%
31	8.549	2237	2242	2255	rVB2	25765	45289	1.22%	0.112%
32	8.684	2272	2284	2301	rBV	793735	1395590	37.74%	3.463%
33	8.848	2310	2335	2343	rBV6	83356	279934	7.57%	0.695%
34	8.893	2343	2349	2358	rVV4	55164	112666	3.05%	0.280%
35	8.951	2358	2367	2377	rVV	279401	500813	13.54%	1.243%
36	9.005	2378	2384	2399	rVB	79139	139213	3.76%	0.345%

LSC Area Percent Report

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Operator : SY/MD
Sample : N3524-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
31466

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\82U062722W.M
Title : SW846 8260

37	9.102	2402	2414	2430	rBV	208614	373002	10.09%	0.926%
38	9.269	2453	2466	2477	rBV	375263	630104	17.04%	1.564%
39	9.336	2477	2487	2502	rBV3	194276	420481	11.37%	1.043%
40	9.417	2502	2512	2526	rBV	668558	1132711	30.63%	2.811%
41	9.484	2527	2533	2542	rVB	80235	125708	3.40%	0.312%
42	9.574	2543	2561	2569	rBV	569446	1015407	27.46%	2.520%
43	9.758	2608	2618	2630	rBV2	431542	783186	21.18%	1.944%
44	9.819	2630	2637	2652	rVB5	85424	166101	4.49%	0.412%
45	9.906	2653	2664	2676	rVB8	57265	145022	3.92%	0.360%
46	9.980	2677	2687	2694	rBV2	87602	147193	3.98%	0.365%
47	10.028	2694	2702	2712	rVB3	112221	188655	5.10%	0.468%
48	10.086	2712	2720	2727	rBV4	48092	73706	1.99%	0.183%
49	10.131	2727	2734	2744	rVB3	60226	92046	2.49%	0.228%
50	10.195	2744	2754	2760	rBV4	95874	166067	4.49%	0.412%
51	10.234	2760	2766	2776	rVB2	116112	176812	4.78%	0.439%
52	10.285	2777	2782	2792	rVB3	67633	101358	2.74%	0.252%
53	10.372	2793	2809	2822	rBV4	429865	1279143	34.59%	3.174%
54	10.430	2822	2827	2854	rVB2	166281	369571	9.99%	0.917%
55	10.545	2855	2863	2875	rBV4	30794	82705	2.24%	0.205%
56	10.639	2877	2892	2907	rBV2	917232	1859778	50.29%	4.615%
57	10.738	2916	2923	2936	rVB7	170205	367052	9.93%	0.911%
58	10.809	2937	2945	2952	rBV2	90278	157629	4.26%	0.391%
59	10.854	2952	2959	2964	rBV	97752	125673	3.40%	0.312%
60	10.886	2964	2969	2976	rVB2	64550	73353	1.98%	0.182%
61	10.931	2976	2983	2988	rBV	151039	205608	5.56%	0.510%
62	11.057	3013	3022	3026	rBV3	63291	98376	2.66%	0.244%
63	11.105	3026	3037	3056	rVB5	311348	752344	20.34%	1.867%
64	11.211	3065	3070	3078	rBV4	25248	41760	1.13%	0.104%
65	11.391	3120	3126	3135	rVV4	34974	55673	1.51%	0.138%
66	11.459	3137	3147	3159	rVB	304464	490799	13.27%	1.218%
67	11.571	3171	3182	3191	rBV4	54919	106539	2.88%	0.264%
68	11.742	3224	3235	3247	rBV	2367597	3698169	100.00%	9.178%
69	11.809	3248	3256	3264	rBV	664739	1019124	27.56%	2.529%
70	11.934	3286	3295	3304	rBV3	133888	238030	6.44%	0.591%
71	12.018	3313	3321	3327	rBV2	147734	255095	6.90%	0.633%
72	12.208	3371	3380	3385	rBV	262080	439840	11.89%	1.092%
73	12.320	3407	3415	3422	rBV2	131838	223253	6.04%	0.554%
74	12.442	3446	3453	3459	rBV3	62464	95956	2.59%	0.238%
75	12.635	3506	3513	3530	rVB6	69798	161014	4.35%	0.400%
76	12.860	3577	3583	3601	rVB4	59720	109758	2.97%	0.272%

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ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
31466

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\82U062722W.M
Title : SW846 8260

77	13.012	3623	3630	3641	rVB3	77149	128162	3.47%	0.318%
78	13.086	3644	3653	3663	rVV4	142937	291196	7.87%	0.723%
79	13.143	3664	3671	3689	rVB2	295843	499709	13.51%	1.240%
80	13.285	3709	3715	3723	rVB	134164	183718	4.97%	0.456%
81	13.333	3724	3730	3737	rVB4	85154	116188	3.14%	0.288%
82	13.381	3738	3745	3757	rVB	135585	197069	5.33%	0.489%
83	13.629	3815	3822	3835	rVB2	92387	149470	4.04%	0.371%
84	13.841	3883	3888	3894	rBV2	27467	39025	1.06%	0.097%
85	13.979	3927	3931	3943	rVB3	29344	39915	1.08%	0.099%
86	16.191	4612	4619	4632	rVB	602520	947701	25.63%	2.352%
87	16.478	4703	4708	4718	rVV4	227519	360876	9.76%	0.896%
88	16.542	4721	4728	4738	rVB2	695985	1088237	29.43%	2.701%
89	17.423	4995	5002	5012	rVB2	442177	675858	18.28%	1.677%

Sum of corrected areas: 40295639

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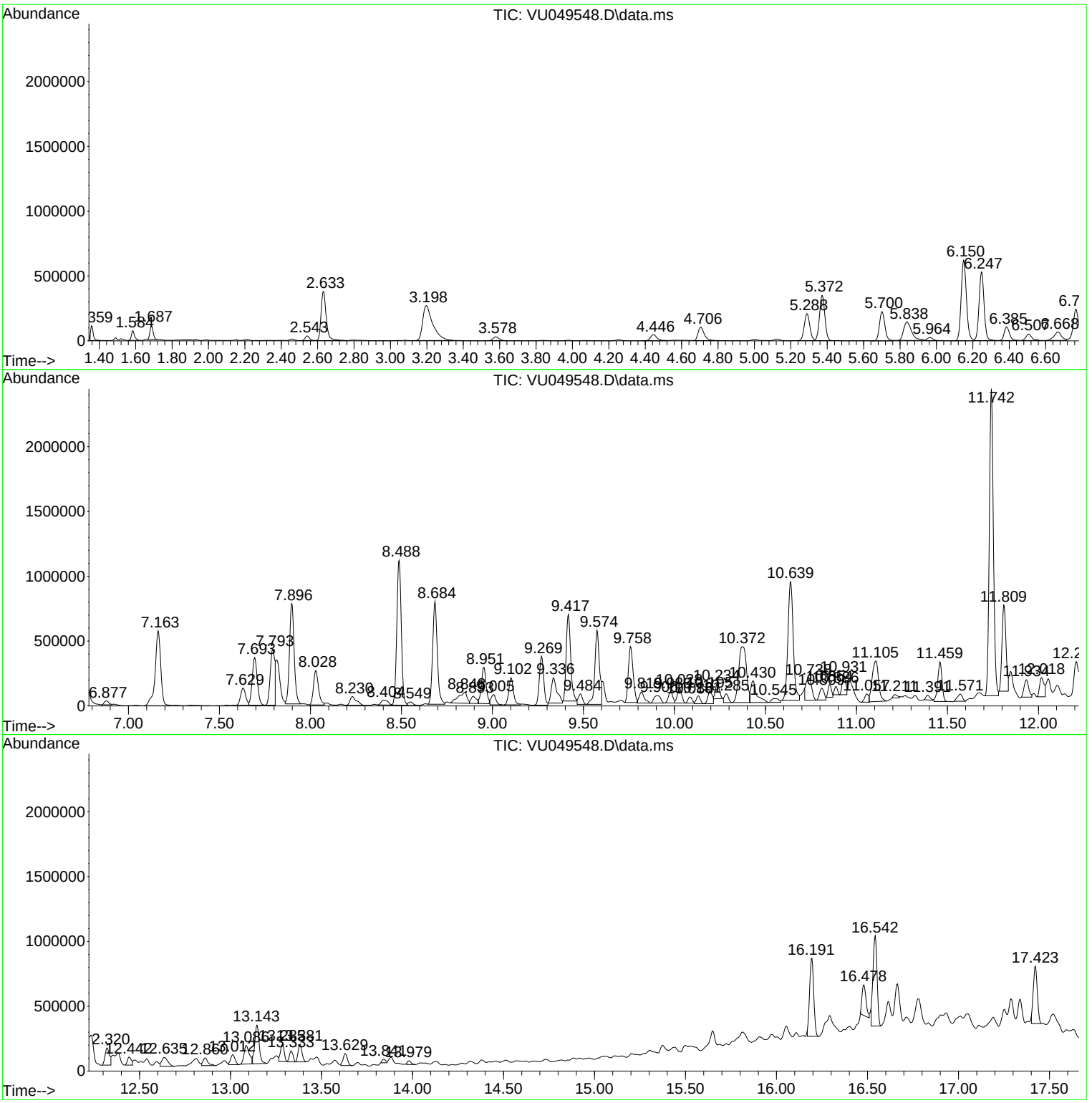
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
31466

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062922\
Data File : VU049548.D
Acq On : 30 Jun 2022 06:16
Operator : SY/MD
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

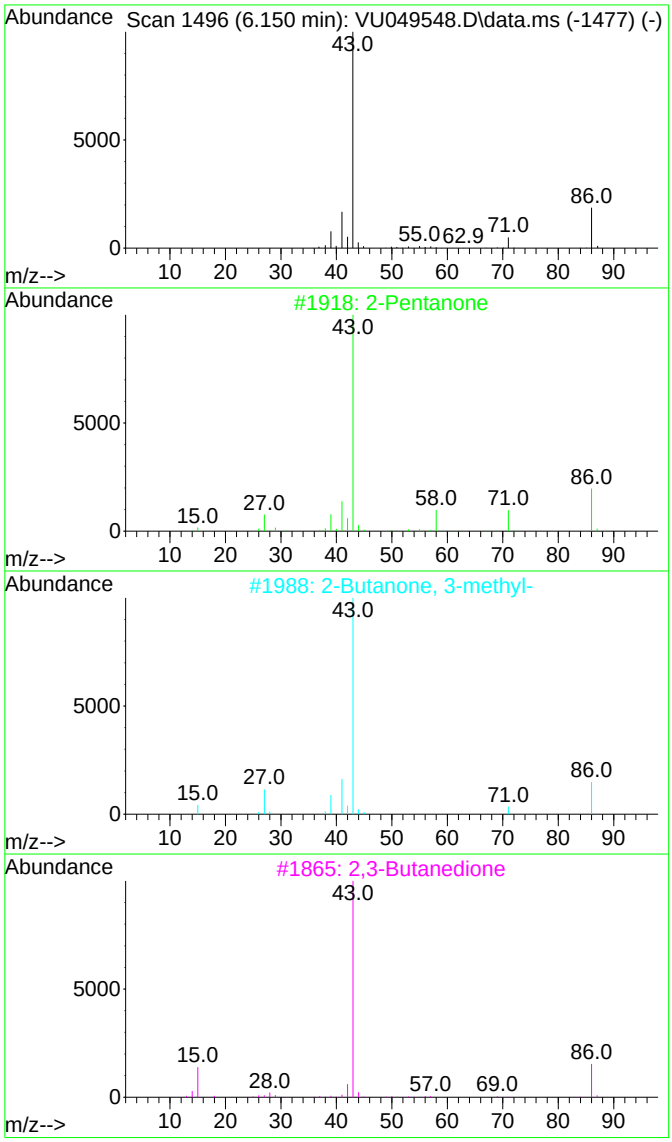
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

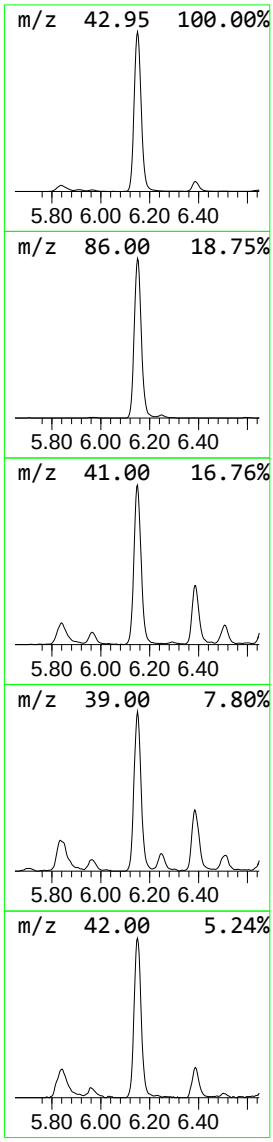
****

Peak Number 1 2-Pentanone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.150	62.03 ug/l	1263050	1,4-Difluorobenzene	6.250

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone	86	C5H10O	000107-87-9	72
2	2-Butanone, 3-methyl-	86	C5H10O	000563-80-4	72
3	2,3-Butanedione	86	C4H6O2	000431-03-8	50
4	3,3-Dimethyl-2,4-pentane dione	128	C7H12O2	003142-58-3	36
5	2-Pentanone, 3-ethyl-	114	C7H14O	006137-03-7	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062922\
Data File : VU049548.D
Acq On : 30 Jun 2022 06:16
Operator : SY/MD
Sample : N3524-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
31466

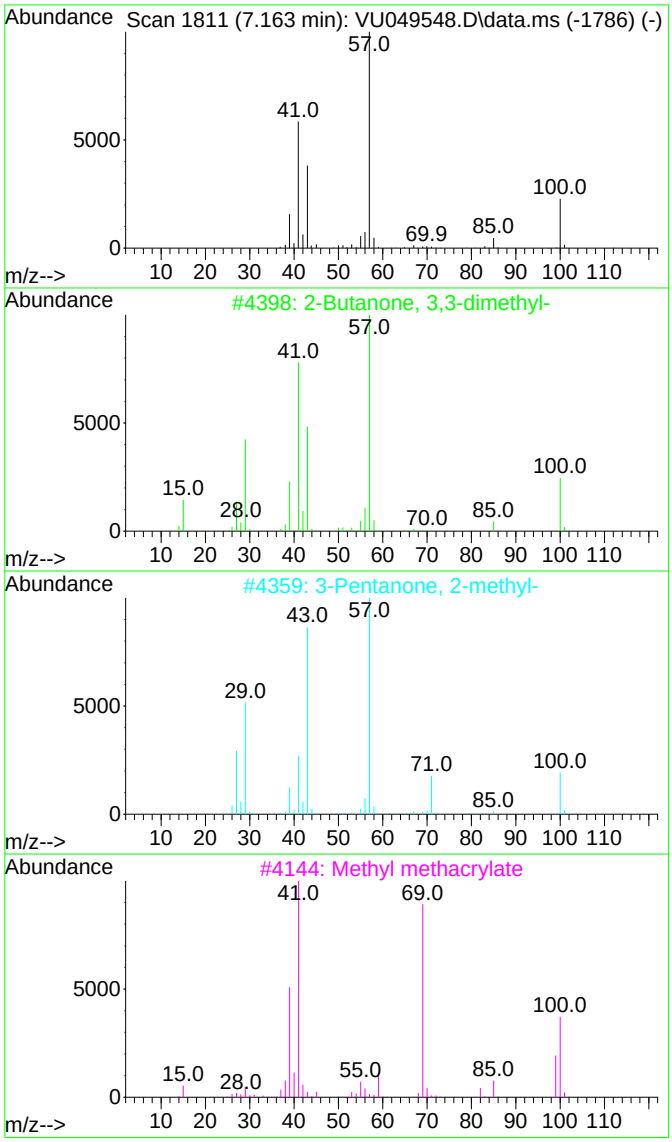
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

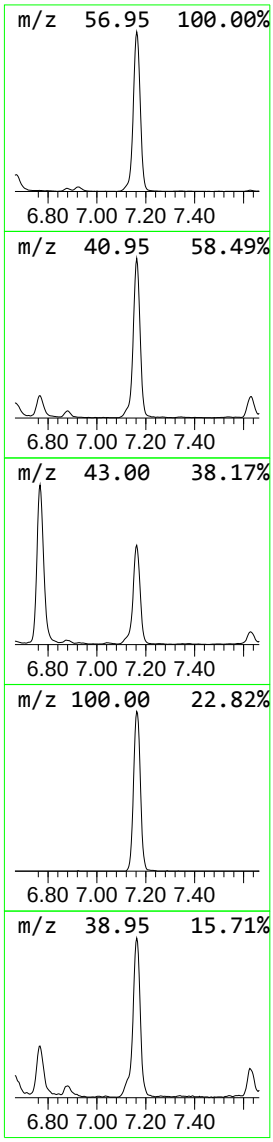
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2-Butanone, 3,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.163	61.03 ug/l	1242750	1,4-Difluorobenzene	6.250

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanone, 3,3-dimethyl-	100	C6H12O	000075-97-8	91
2	3-Pentanone, 2-methyl-	100	C6H12O	000565-69-5	64
3	Methyl methacrylate	100	C5H8O2	000080-62-6	43
4	Propane, 2-methyl-2-nitro-	103	C4H9NO2	000594-70-7	43
5	Propane, 1-(ethenyloxy)-2-methyl-	100	C6H12O	000109-53-5	42





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Data File : VU049548.D
Acq On : 30 Jun 2022 06:16
Operator : SY/MD
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Instrument :
MSVOA_U
ClientSampleId :
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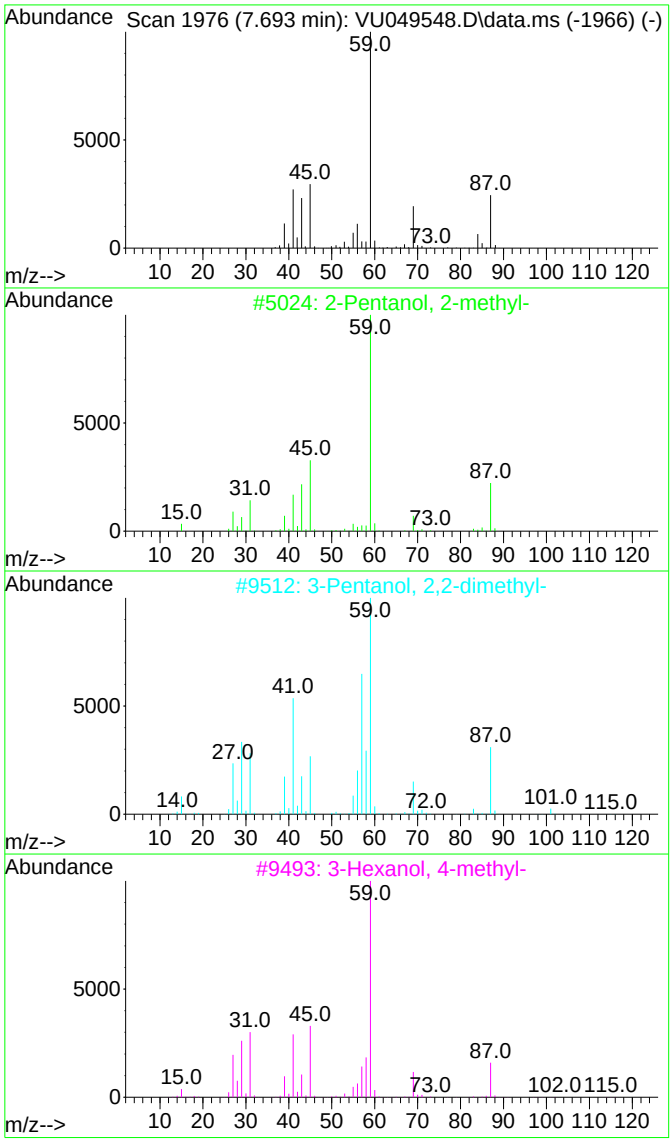
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

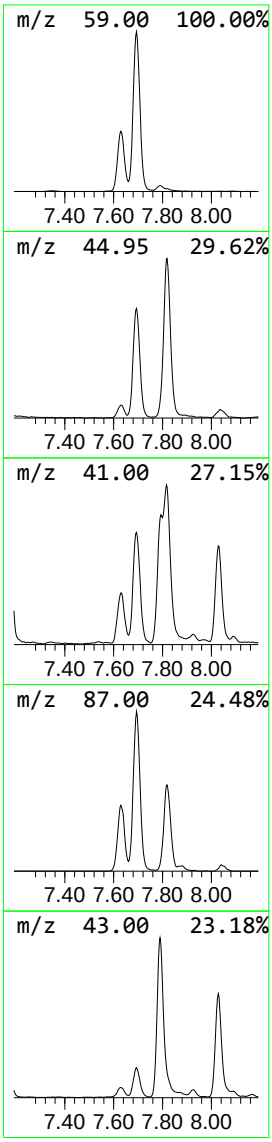
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-Pentanol, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.693	34.96 ug/l	711906	1,4-Difluorobenzene	6.250

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	78
2	3-Pentanol, 2,2-dimethyl-	116	C7H16O	003970-62-5	64
3	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	64
4	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	64
5	3-Hexanol, 5-methyl-	116	C7H16O	000623-55-2	40





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ALS Vial : 51 Sample Multiplier: 1

Instrument :
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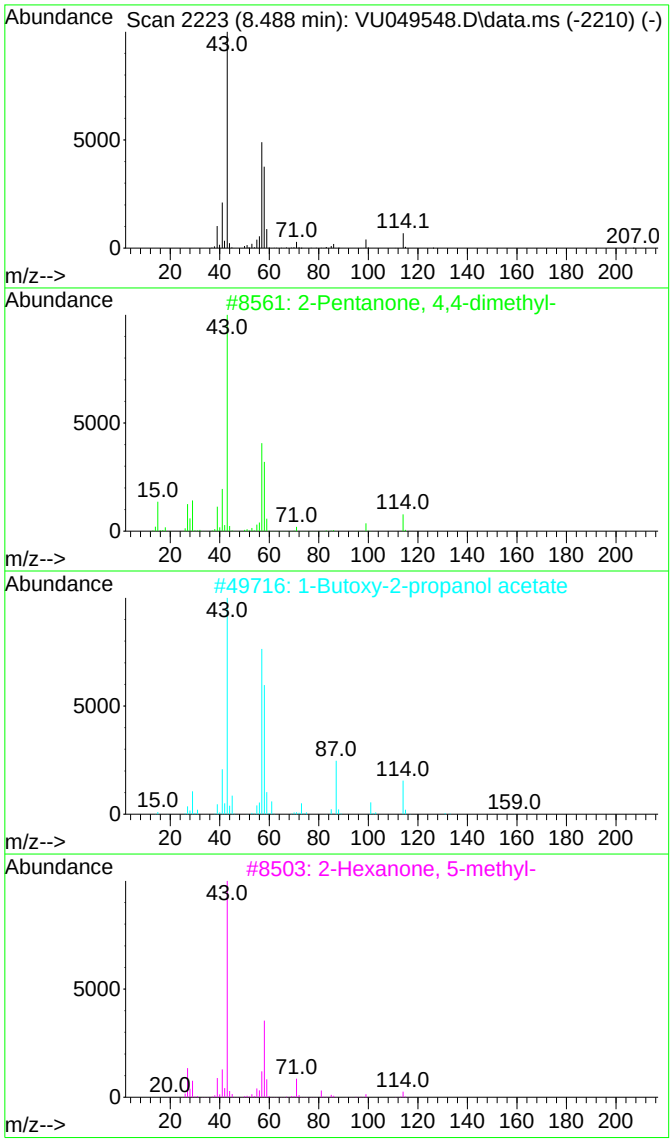
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

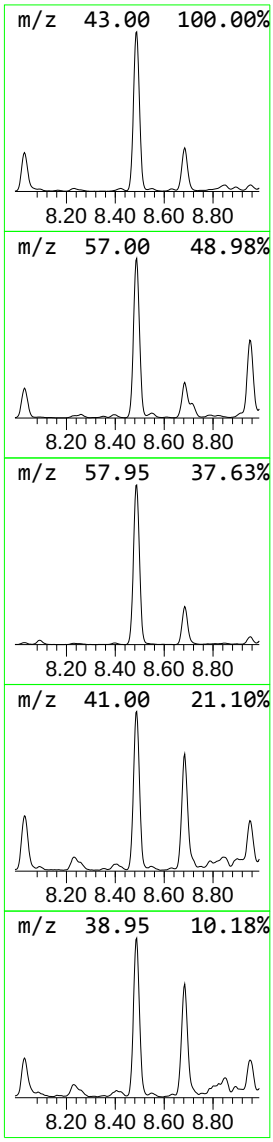
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 2-Pentanone, 4,4-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.488	84.72 ug/l	1919310	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4,4-dimethyl-	114	C7H14O	000590-50-1	90
2	1-Butoxy-2-propanol acetate	174	C9H18O3	085409-76-3	72
3	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	64
4	2-Heptanone	114	C7H14O	000110-43-0	59
5	Butanal, 2-methyl-	86	C5H10O	000096-17-3	25





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Data File : VU049548.D
Acq On : 30 Jun 2022 06:16
Operator : SY/MD
Sample : N3524-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
31466

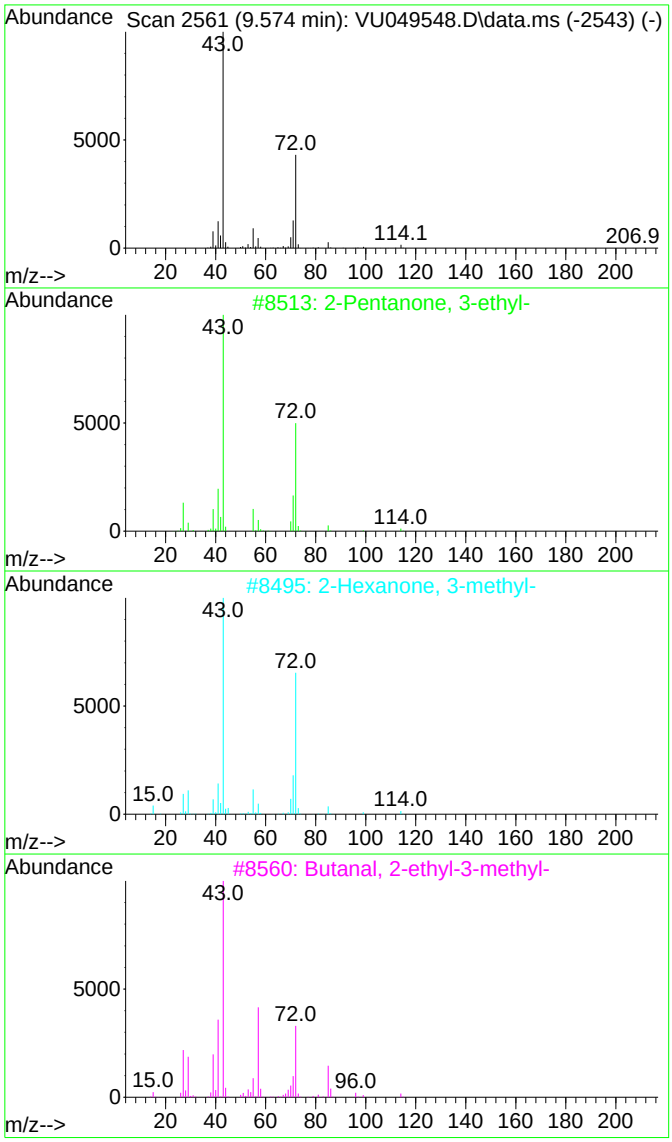
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

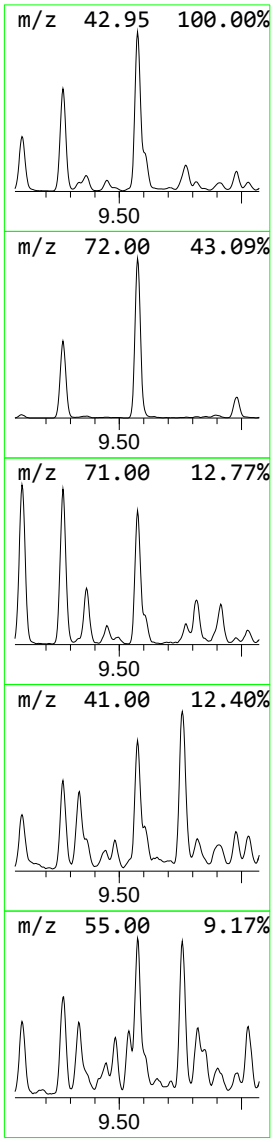
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 2-Pentanone, 3-ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.574	44.82 ug/l	1015410	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 3-ethyl-	114	C7H14O	006137-03-7	80
2	2-Hexanone, 3-methyl-	114	C7H14O	002550-21-2	80
3	Butanal, 2-ethyl-3-methyl-	114	C7H14O	026254-92-2	72
4	Propanal, 2-methyl-	72	C4H8O	000078-84-2	40
5	2-Heptanone, 3-methyl-	128	C8H16O	002371-19-9	36





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062922\
Data File : VU049548.D
Acq On : 30 Jun 2022 06:16
Operator : SY/MD
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ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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

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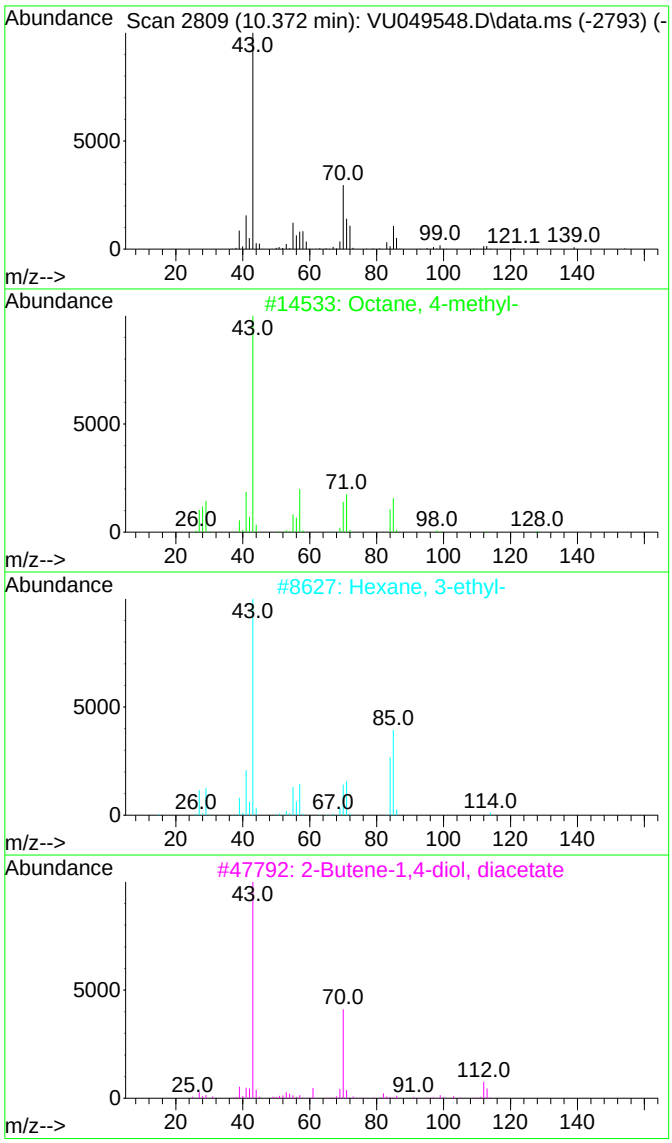
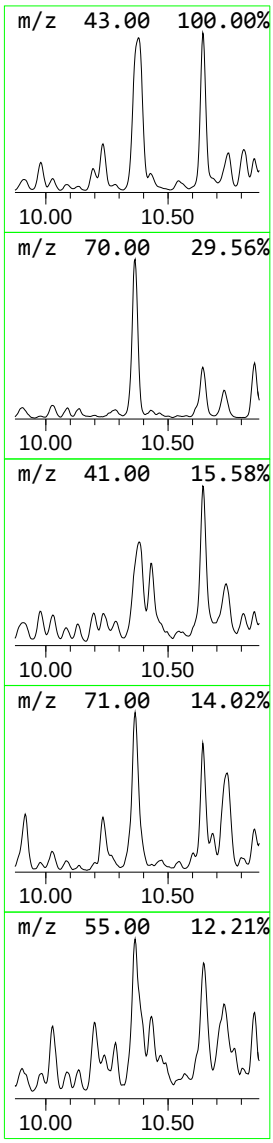
Peak Number 6 unknown10.372 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.372	56.46 ug/l	1279140	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 4-methyl-	128	C9H20	002216-34-4	37
2	Hexane, 3-ethyl-	114	C8H18	000619-99-8	37
3	2-Butene-1,4-diol, diacetate	172	C8H12O4	018621-75-5	37
4	Heptane, 2,5,5-trimethyl-	142	C10H22	001189-99-7	36
5	1-Butanol, 2,2-dimethyl-	102	C6H14O	001185-33-7	33





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062922\
Data File : VU049548.D
Acq On : 30 Jun 2022 06:16
Operator : SY/MD
Sample : N3524-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
31466

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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

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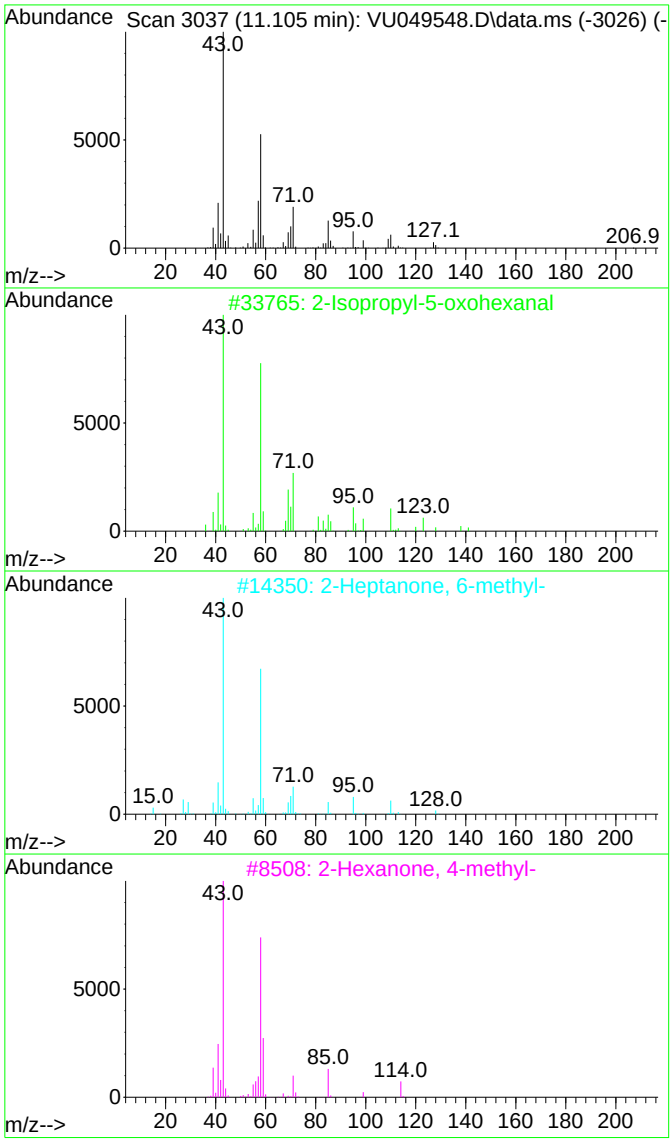
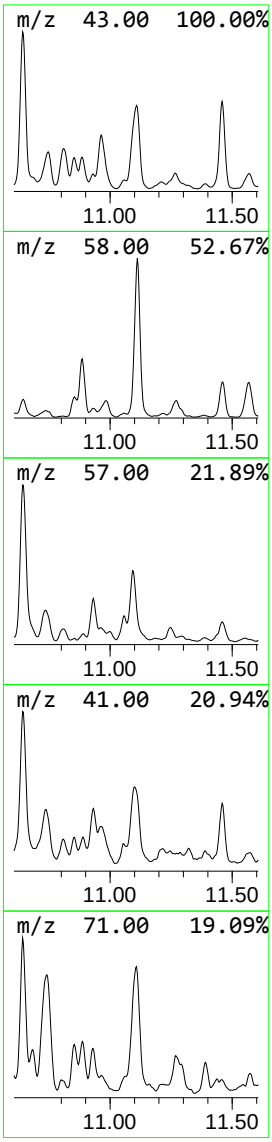
Peak Number 7 2-Isopropyl-5-oxohexanal Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.105	36.91 ug/l	752344	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Isopropyl-5-oxohexanal	156	C9H16O2	015303-46-5	78
2	2-Heptanone, 6-methyl-	128	C8H16O	000928-68-7	62
3	2-Hexanone, 4-methyl-	114	C7H14O	000105-42-0	43
4	2-Heptanone	114	C7H14O	000110-43-0	38
5	2-Hexanone	100	C6H12O	000591-78-6	37





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Data File : VU049548.D
Acq On : 30 Jun 2022 06:16
Operator : SY/MD
Sample : N3524-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
31466

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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

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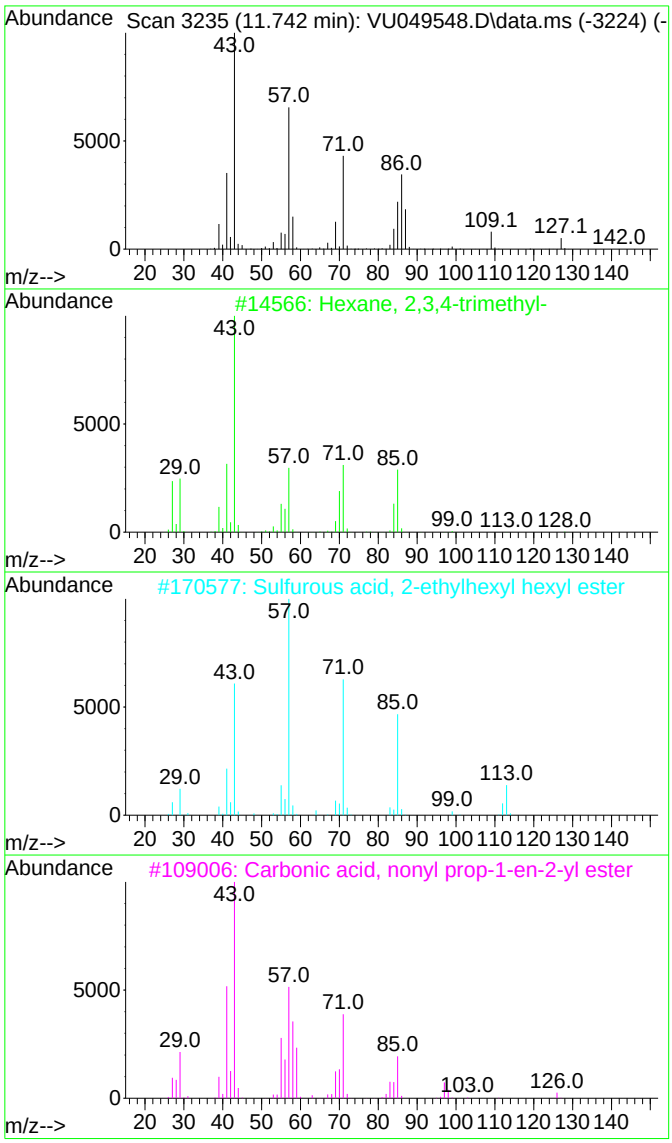
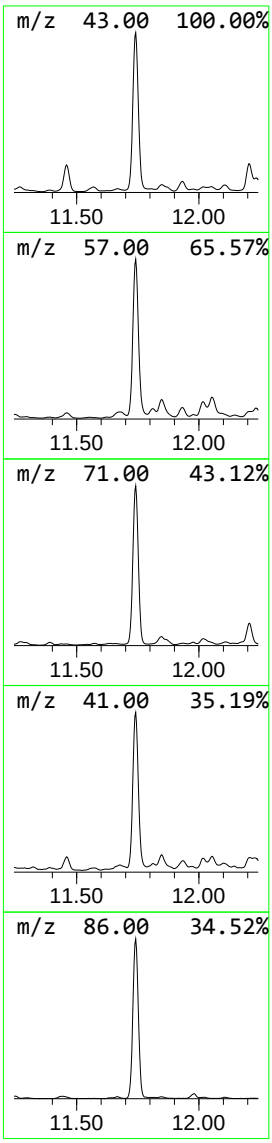
Peak Number 8 Hexane, 2,3,4-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.742	181.44 ug/l	3698170	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	50
2	Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	43
3	Carbonic acid, nonyl prop-1-en-2...	228	C13H24O3	1000382-53-9	40
4	Carbonic acid, nonyl vinyl ester	214	C12H22O3	1000383-25-6	40
5	Sulfurous acid, hexyl pentyl ester	236	C11H24O3S	1000309-14-1	38





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Data File : VU049548.D
Acq On : 30 Jun 2022 06:16
Operator : SY/MD
Sample : N3524-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
31466

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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

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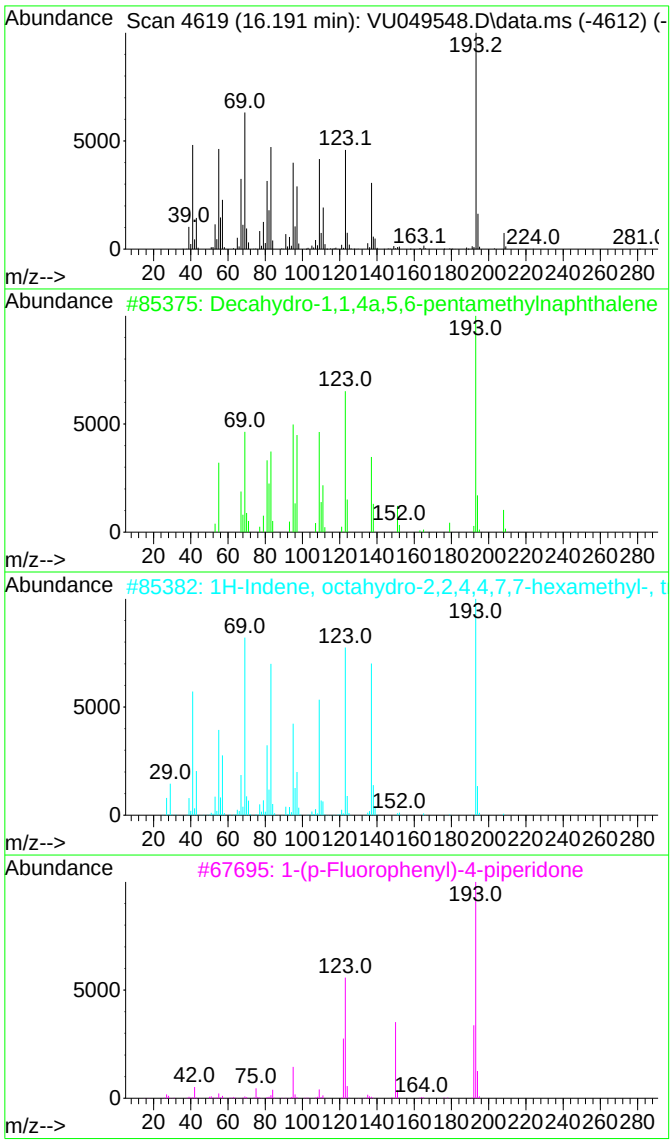
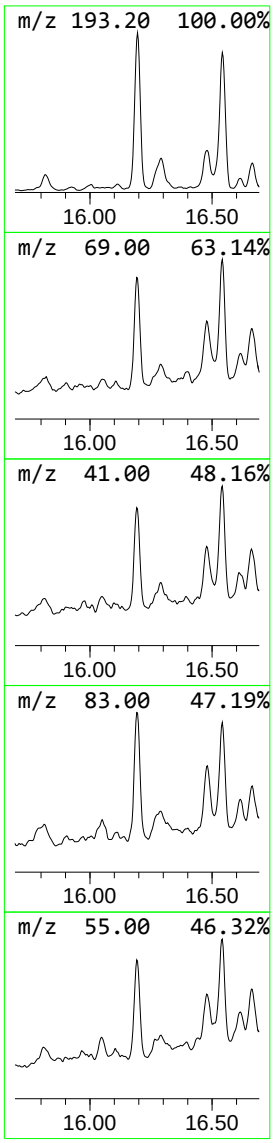
Peak Number 9 unknown16.192 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.192	46.50 ug/l	947701	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	43
2	1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	38
3	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	25
4	2-Anthracenamine	193	C14H11N	000613-13-8	22
5	Butan-2-one, 1-[3-(3,3-dimethyl-...	278	C16H26N2O2	1000303-34-2	22





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Data File : VU049548.D
Acq On : 30 Jun 2022 06:16
Operator : SY/MD
Sample : N3524-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
31466

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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

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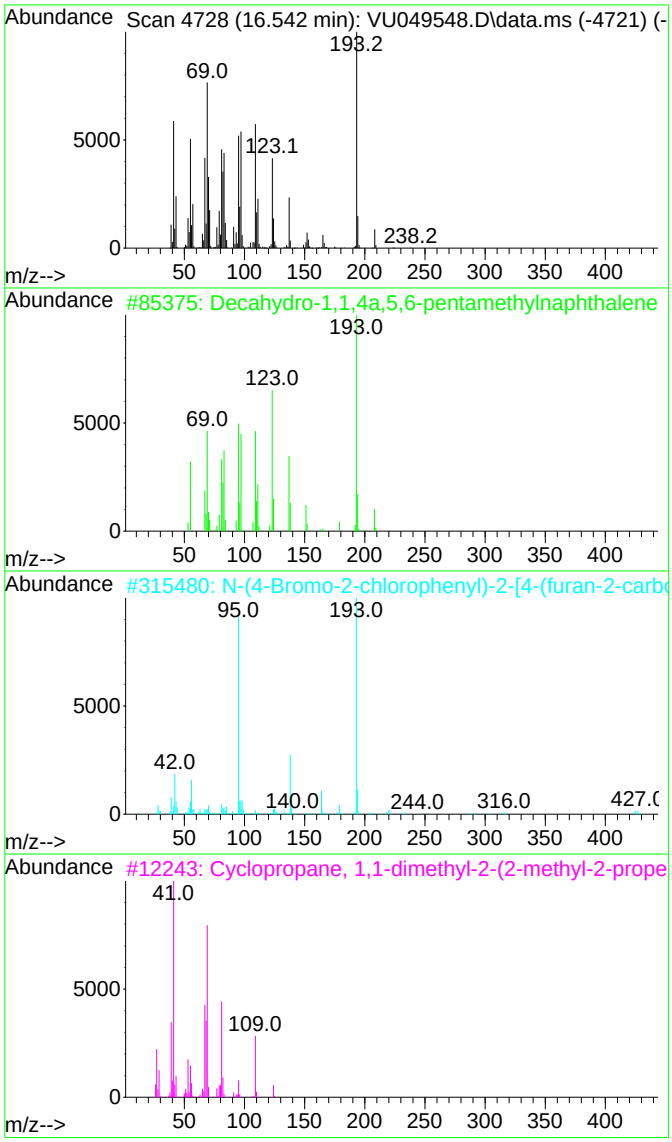
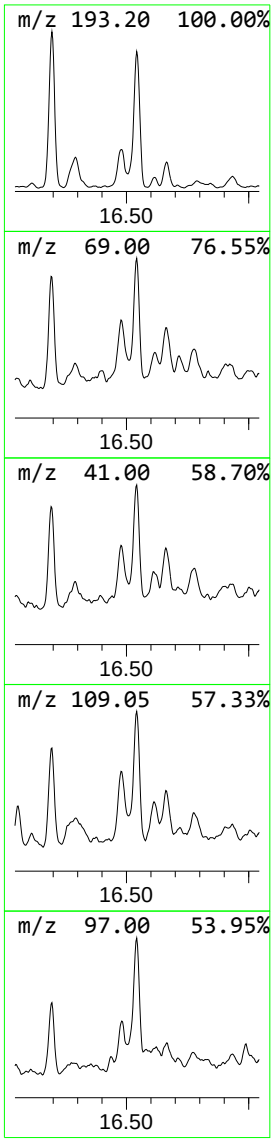
Peak Number 10 unknown16.542 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.542	53.39 ug/l	1088240	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	38
2	N-(4-Bromo-2-chlorophenyl)-2-[4-...	425	C17H17BrClN3O3	1000482-48-9	22
3	Cyclopropane, 1,1-dimethyl-2-(2-...	124	C9H16	069147-03-1	20
4	Bicyclo[3.1.1]heptane, 2,6,6-tri...	138	C10H18	000473-55-2	15
5	(-)-trans-Pinane	138	C10H18	033626-25-4	15





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062922\
Data File : VU049548.D
Acq On : 30 Jun 2022 06:16
Operator : SY/MD
Sample : N3524-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
31466

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Pentanone	6.150	62.0	ug/l	1263050	2	6.250	1018070	50.0
2-Butanone, 3,3...	7.163	61.0	ug/l	1242750	2	6.250	1018070	50.0
2-Pentanol, 2-m...	7.693	35.0	ug/l	711906	2	6.250	1018070	50.0
2-Pentanone, 4,...	8.488	84.7	ug/l	1919310	3	9.417	1132710	50.0
2-Pentanone, 3-...	9.574	44.8	ug/l	1015410	3	9.417	1132710	50.0
unknown10.372	10.372	56.5	ug/l	1279140	3	9.417	1132710	50.0
2-Isopropyl-5-o...	11.105	36.9	ug/l	752344	4	11.809	1019120	50.0
Hexane, 2,3,4-t...	11.742	181.4	ug/l	3698170	4	11.809	1019120	50.0
unknown16.192	16.192	46.5	ug/l	947701	4	11.809	1019120	50.0
unknown16.542	16.542	53.4	ug/l	1088240	4	11.809	1019120	50.0