

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
 Data File : VV037562.D
 Acq On : 26 Sep 2024 20:26
 Operator : SY/MD
 Sample : P4184-14
 Misc : 5mL/MSVOA_V/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 EY040

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0 % 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_V\Method\SFAMVLM092624WMA.M
 Title : VOC Analysis

Signal : TIC: VV037562.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.175	32	42	52	rBV	30640	56454	0.04%	0.015%
2	1.285	66	76	95	rBV2	2615656	2943125	1.83%	0.778%
3	1.429	113	121	132	rVB4	8070	13258	0.01%	0.004%
4	1.548	145	158	180	rVB2	1158692	1690510	1.05%	0.447%
5	1.667	188	195	204	rBV	84767	102458	0.06%	0.027%
6	1.715	204	210	212	rVV	12903	14248	0.01%	0.004%
7	1.735	213	216	224	rVB	20556	24631	0.02%	0.007%
8	1.799	229	236	247	rVB	56103	69359	0.04%	0.018%
9	1.989	286	295	305	rBV2	18330	30641	0.02%	0.008%
10	2.066	305	319	340	rBV4	1841408	3037091	1.89%	0.803%
11	2.153	340	346	357	rVB2	22299	31512	0.02%	0.008%
12	2.445	426	437	454	rBV	50664	88804	0.06%	0.023%
13	2.690	500	513	540	rVB	674903	1147544	0.71%	0.303%
14	2.963	587	598	611	rVB3	11382	25540	0.02%	0.007%
15	3.108	623	643	687	rVB2	32190151	68573506	42.64%	18.126%
16	3.661	801	815	829	rBV5	11876	30107	0.02%	0.008%
17	3.809	841	861	921	rBV4	62954681	160834629	100.00%	42.514%
18	4.243	982	996	1026	rVB	415955	1060347	0.66%	0.280%
19	4.506	1053	1078	1117	rBV	15929086	39865054	24.79%	10.538%
20	4.950	1200	1216	1229	rBV2	1030248	2585633	1.61%	0.683%
21	5.230	1295	1303	1316	rVB2	6456	15701	0.01%	0.004%
22	5.529	1384	1396	1430	rBV	775388	1567138	0.97%	0.414%
23	5.828	1472	1489	1521	rBV	749062	1508440	0.94%	0.399%
24	5.985	1525	1538	1551	rBV	574079	1163034	0.72%	0.307%
25	6.908	1815	1825	1850	rVV	313686	581140	0.36%	0.154%
26	7.178	1898	1909	1917	rVV8	38348	89105	0.06%	0.024%
27	7.239	1917	1928	1939	rVV	1221571	2143061	1.33%	0.566%
28	7.307	1939	1949	1975	rVV	3935062	6781232	4.22%	1.793%
29	7.410	1975	1981	2001	rVV3	32496	95649	0.06%	0.025%
30	7.497	2001	2008	2015	rVV7	10634	20482	0.01%	0.005%
31	7.548	2015	2024	2052	rVB	217739	438859	0.27%	0.116%
32	7.709	2066	2074	2084	rVB5	11624	20823	0.01%	0.006%
33	7.773	2085	2094	2113	rVB2	38109	71877	0.04%	0.019%
34	7.898	2118	2133	2156	rBV	406113	706790	0.44%	0.187%
35	8.024	2162	2172	2205	rBV	377255	774914	0.48%	0.205%
36	8.217	2222	2232	2243	rVV2	43818	68361	0.04%	0.018%

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Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM092624WMA.M
 Title : VOC Analysis

37	8.278	2243	2251	2259	rVB3	17544	29290	0.02%	0.008%
38	8.519	2311	2326	2336	rBV3	5977	13945	0.01%	0.004%
39	8.651	2358	2367	2378	rBV5	11962	22311	0.01%	0.006%
40	8.779	2396	2407	2426	rBV	1207640	1968628	1.22%	0.520%
41	8.860	2427	2432	2444	rVV4	25989	46215	0.03%	0.012%
42	8.940	2444	2457	2480	rVV	2510433	3921163	2.44%	1.036%
43	9.062	2484	2495	2535	rVB	5913577	9599082	5.97%	2.537%
44	9.242	2545	2551	2567	rVB6	14213	27843	0.02%	0.007%
45	9.406	2590	2602	2610	rBV9	17058	39696	0.02%	0.010%
46	9.471	2610	2622	2643	rVV	4804207	7160314	4.45%	1.893%
47	9.632	2665	2672	2685	rVB3	49774	87052	0.05%	0.023%
48	9.728	2693	2702	2717	rVB6	37987	71533	0.04%	0.019%
49	9.863	2733	2744	2762	rBV	293705	475076	0.30%	0.126%
50	9.950	2763	2771	2783	rVB2	14470	25730	0.02%	0.007%
51	10.021	2783	2793	2798	rBV3	21688	36838	0.02%	0.010%
52	10.146	2820	2832	2852	rBV	711572	1150947	0.72%	0.304%
53	10.284	2864	2875	2893	rBV	1114521	1732106	1.08%	0.458%
54	10.377	2893	2904	2923	rVV	3234791	6625117	4.12%	1.751%
55	10.467	2923	2932	2950	rVB	2304032	3412603	2.12%	0.902%
56	10.667	2972	2994	3018	rBV	2785303	4348931	2.70%	1.150%
57	10.844	3038	3049	3074	rVB	10278459	14857141	9.24%	3.927%
58	10.988	3087	3094	3098	rBV2	78082	111005	0.07%	0.029%
59	11.021	3098	3104	3117	rVB	175284	270547	0.17%	0.072%
60	11.127	3126	3137	3144	rBV	264832	399188	0.25%	0.106%
61	11.178	3144	3153	3169	rVV	1496668	2308641	1.44%	0.610%
62	11.265	3169	3180	3202	rVB	3820381	5610939	3.49%	1.483%
63	11.384	3209	3217	3228	rVB	48139	75504	0.05%	0.020%
64	11.458	3228	3240	3250	rBV	1367721	2464089	1.53%	0.651%
65	11.506	3250	3255	3262	rVV	541821	786011	0.49%	0.208%
66	11.554	3262	3270	3297	rVB4	1581957	3567734	2.22%	0.943%
67	11.676	3300	3308	3320	rVB2	93481	147793	0.09%	0.039%
68	11.750	3320	3331	3347	rBV	285364	438659	0.27%	0.116%
69	11.850	3347	3362	3378	rBV3	1185169	2615568	1.63%	0.691%
70	11.947	3384	3392	3412	rVB2	583825	1063595	0.66%	0.281%
71	12.046	3413	3423	3429	rBV	203061	342879	0.21%	0.091%
72	12.191	3456	3468	3474	rBV5	37249	69066	0.04%	0.018%
73	12.246	3474	3485	3498	rVB2	167299	296253	0.18%	0.078%
74	12.316	3498	3507	3508	rBV3	50347	59403	0.04%	0.016%
75	12.345	3509	3516	3524	rVV	251432	378172	0.24%	0.100%
76	12.397	3524	3532	3550	rVB	463235	708345	0.44%	0.187%

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Integration Parameters: LSCINT.P

Integrator: RTE

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77	12.487	3551	3560	3565	rBV4	45311	68052	0.04%	0.018%
78	12.622	3595	3602	3605	rBV2	16373	21614	0.01%	0.006%
79	12.657	3606	3613	3627	rVB3	111134	184834	0.11%	0.049%
80	12.728	3629	3635	3642	rBV4	18245	22978	0.01%	0.006%
81	12.811	3642	3661	3677	rBV2	511207	867743	0.54%	0.229%
82	12.934	3693	3699	3704	rBV	20844	28572	0.02%	0.008%
83	12.975	3704	3712	3734	rVB	119522	214102	0.13%	0.057%
84	13.098	3739	3750	3759	rBV7	31075	58984	0.04%	0.016%
85	13.149	3759	3766	3773	rVB5	22106	29860	0.02%	0.008%
86	13.200	3773	3782	3789	rBV3	55327	90345	0.06%	0.024%
87	13.300	3808	3813	3819	rVV2	17506	23944	0.01%	0.006%
88	13.332	3819	3823	3832	rVB2	20014	28152	0.02%	0.007%
89	13.432	3842	3854	3879	rBV	360957	630785	0.39%	0.167%
90	13.545	3882	3889	3899	rVB	10276	17105	0.01%	0.005%
91	13.651	3911	3922	3923	rBV7	9379	13577	0.01%	0.004%
92	13.844	3972	3982	3996	rVB6	11212	24290	0.02%	0.006%
93	14.033	4028	4041	4052	rBV6	20803	47689	0.03%	0.013%
94	14.165	4075	4082	4093	rVV5	10007	16593	0.01%	0.004%
95	14.226	4093	4101	4112	rVB7	7643	15498	0.01%	0.004%
96	14.361	4133	4143	4157	rBV9	13992	27838	0.02%	0.007%
97	14.567	4196	4207	4223	rBV2	70476	123716	0.08%	0.033%
98	14.750	4253	4264	4276	rBV	44063	78112	0.05%	0.021%
99	15.747	4565	4574	4579	rBV8	12373	18108	0.01%	0.005%
100	16.110	4679	4687	4698	rBV9	12759	22303	0.01%	0.006%

Sum of corrected areas: 378310803

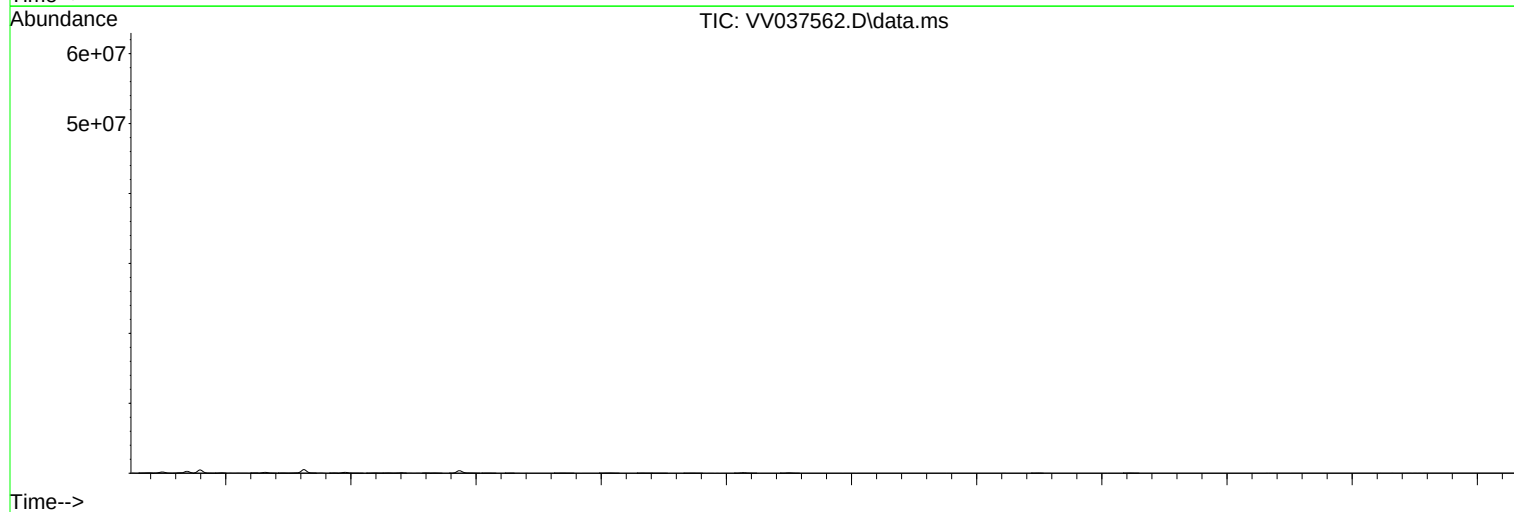
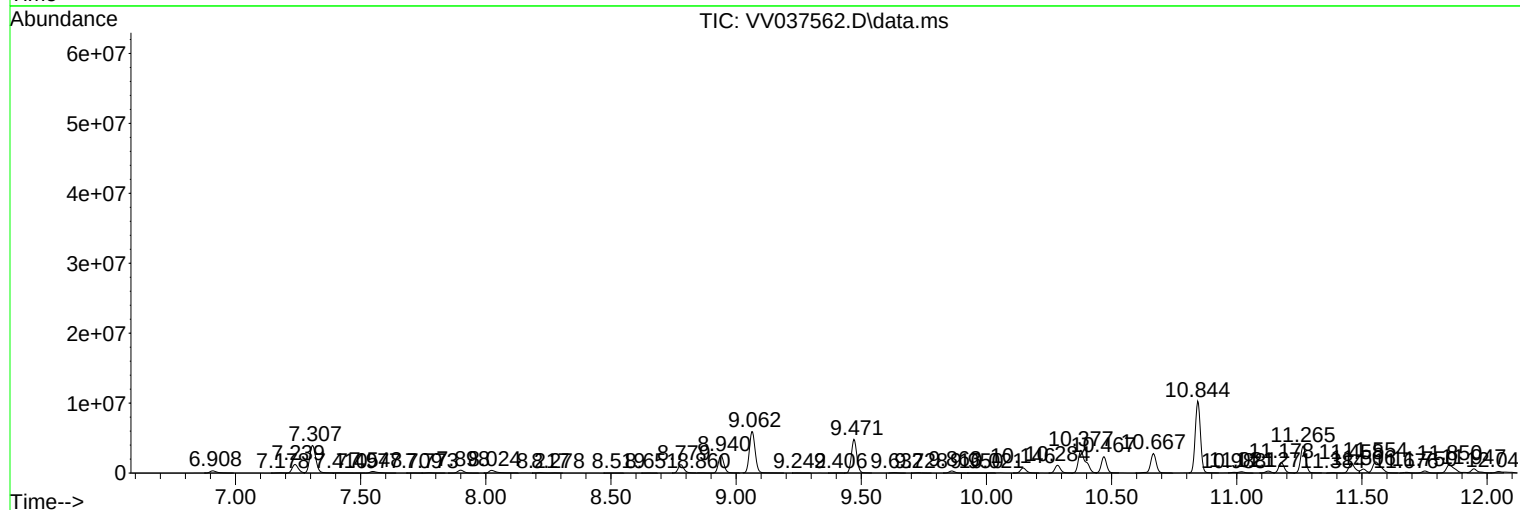
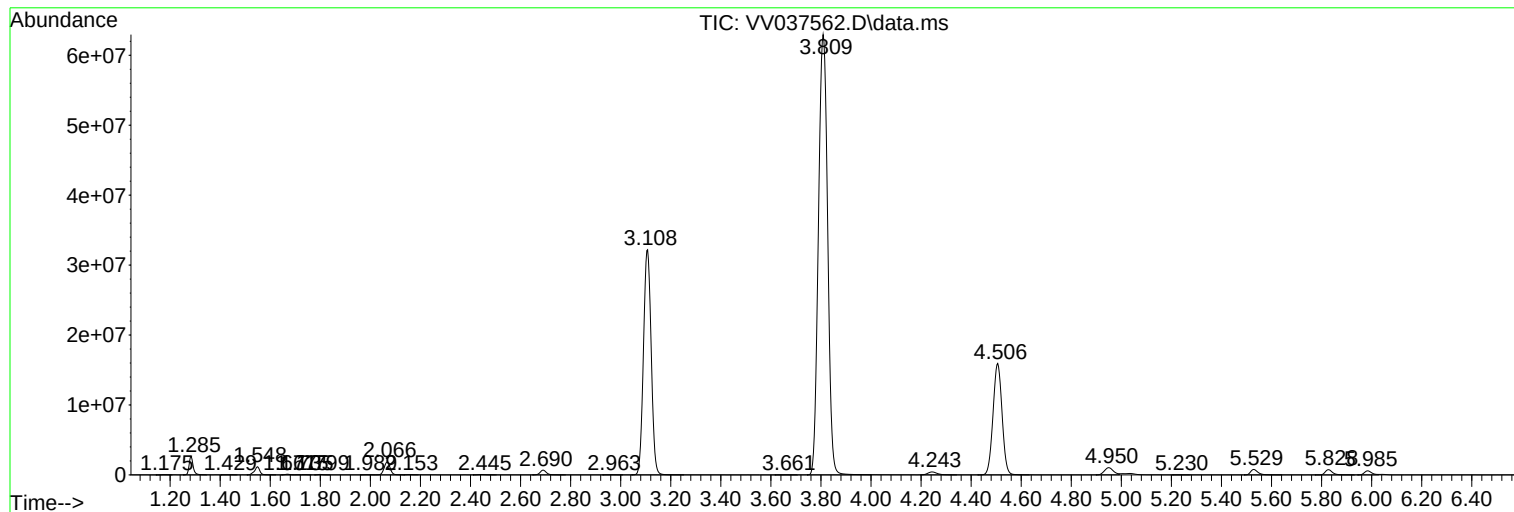
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Instrument :
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ClientSampleId :
EY040

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM092624WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P



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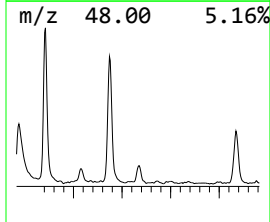
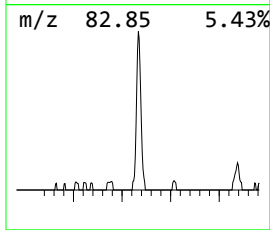
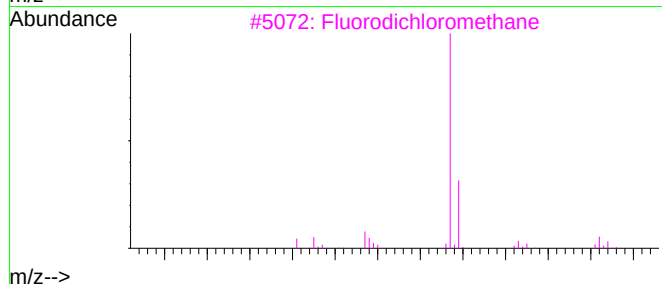
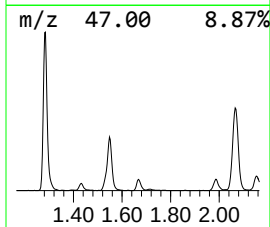
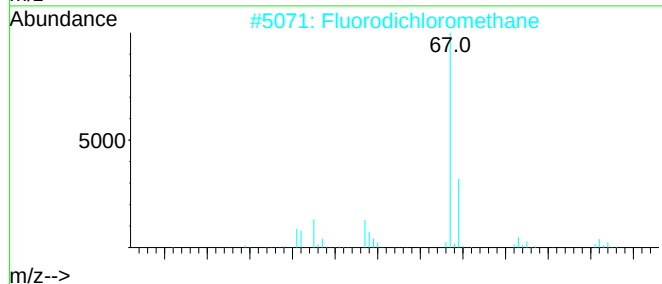
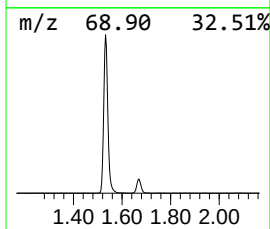
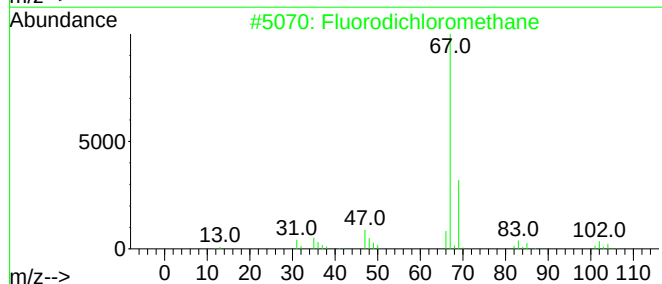
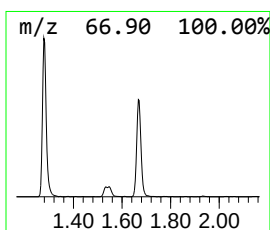
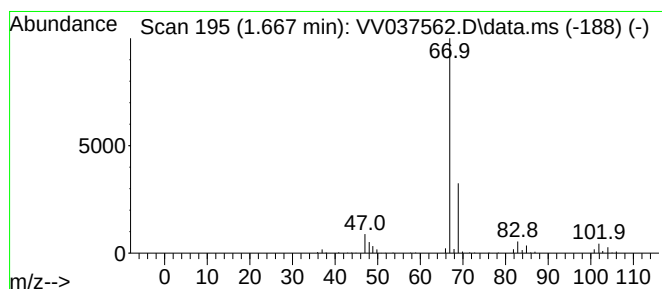
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 1 Fluorodichloromethane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.667	3.27 ug/L	102458	1,4-Difluorobenzene	5.529

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluorodichloromethane	102	CHCl2F	000075-43-4	91
2	Fluorodichloromethane	102	CHCl2F	000075-43-4	91
3	Fluorodichloromethane	102	CHCl2F	000075-43-4	91
4	Cyclopentene, 1-chloro-	102	C5H7Cl	000930-29-0	7
5	Methanamine, N,N-difluoro-	67	CH3F2N	000753-58-2	4



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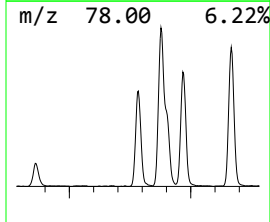
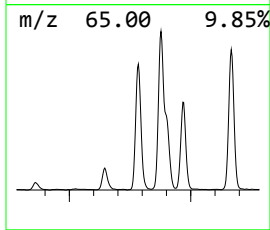
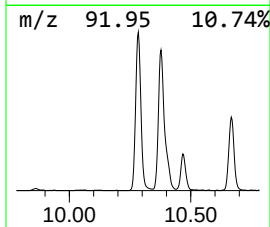
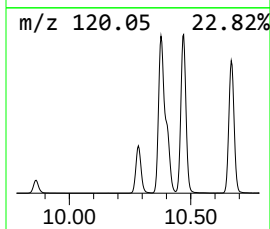
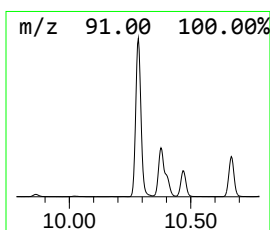
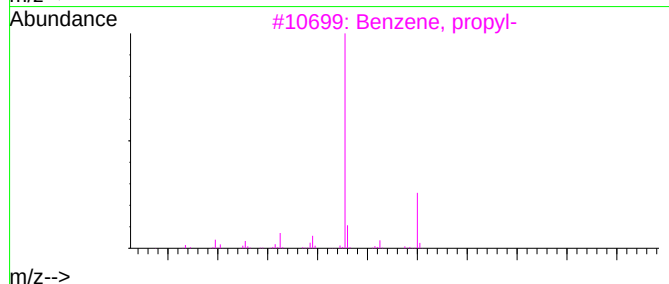
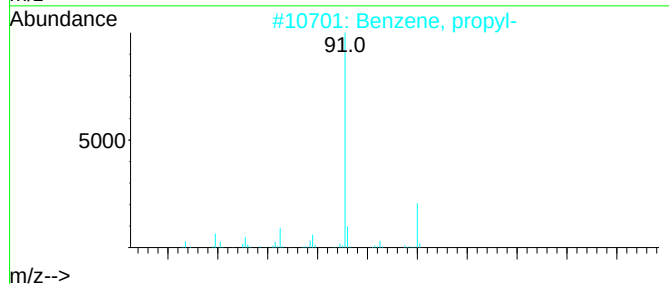
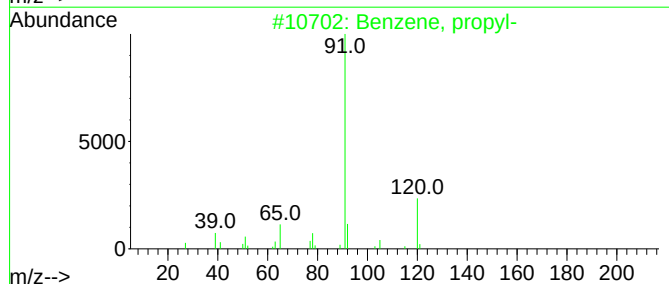
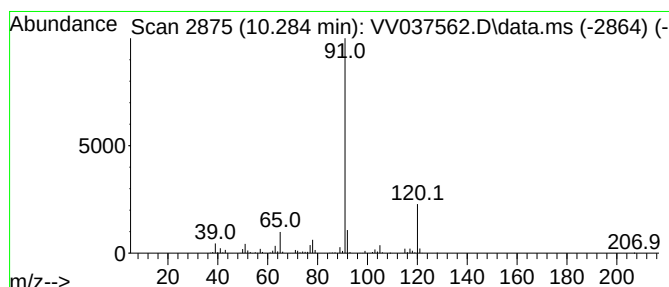
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM092624WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD			R.T.
10.284	37.51 ug/L	1732110	1,4-Dichlorobenzene-d4			11.178
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, propyl-		120	C9H12	000103-65-1	93
2	Benzene, propyl-		120	C9H12	000103-65-1	87
3	Benzene, propyl-		120	C9H12	000103-65-1	87
4	N-Benzyl-2-phenethylamine		211	C15H17N	003647-71-0	78
5	Benzene, propyl-		120	C9H12	000103-65-1	76



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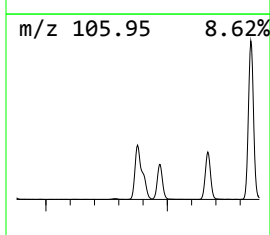
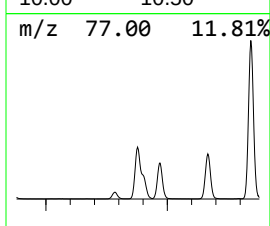
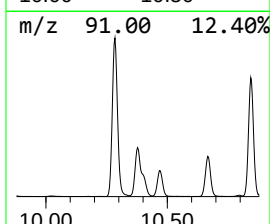
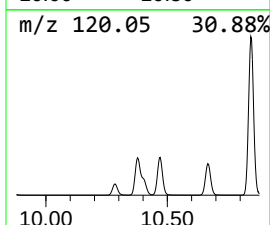
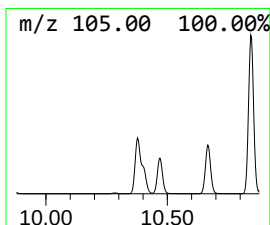
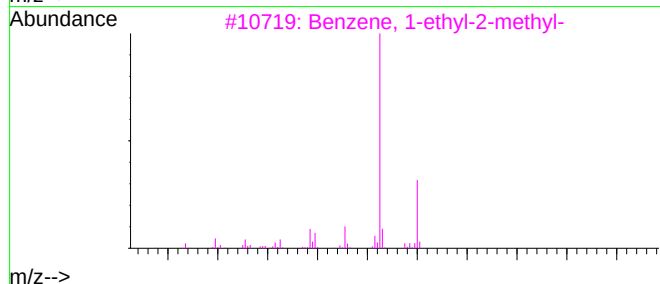
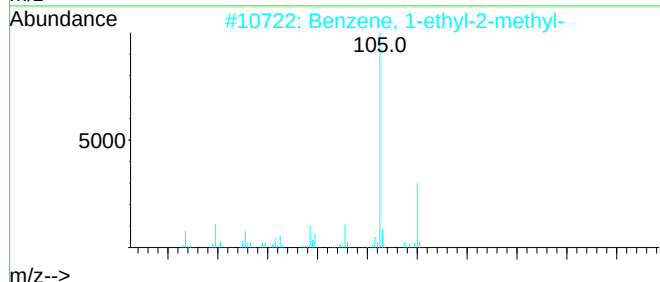
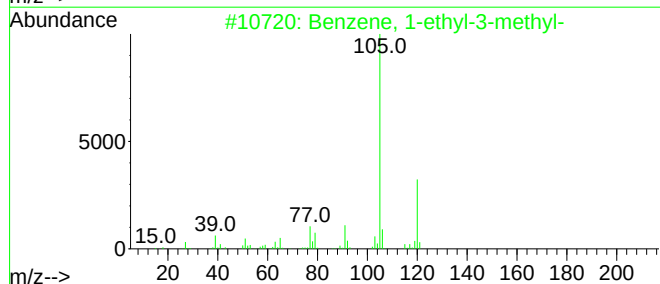
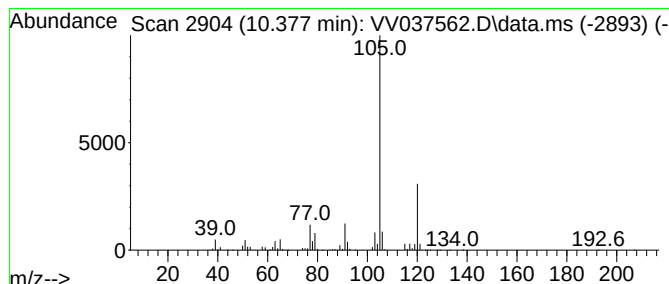
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM092624WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1-ethyl-3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
10.378	143.49 ug/L	6625120	1,4-Dichlorobenzene-d4			11.178
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-		120	C9H12	000620-14-4	95
2	Benzene, 1-ethyl-2-methyl-		120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-2-methyl-		120	C9H12	000611-14-3	95
4	Benzene, 1-ethyl-3-methyl-		120	C9H12	000620-14-4	94
5	Benzene, 1,2,4-trimethyl-		120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY040

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM092624WMA.M
Quant Title : VOC Analysis

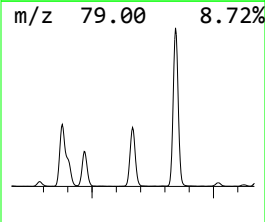
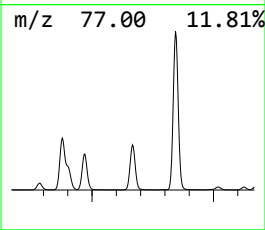
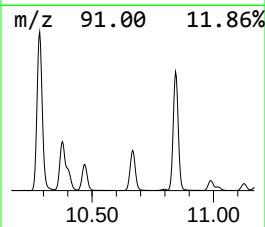
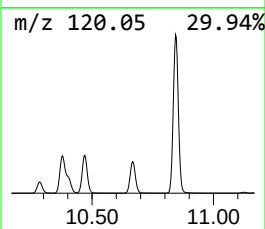
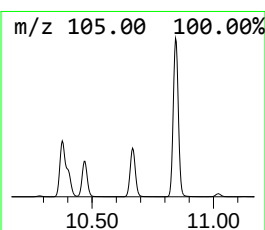
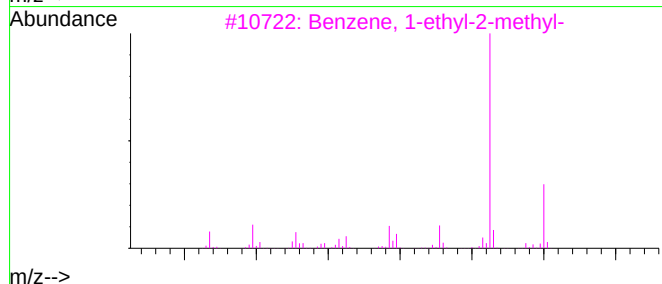
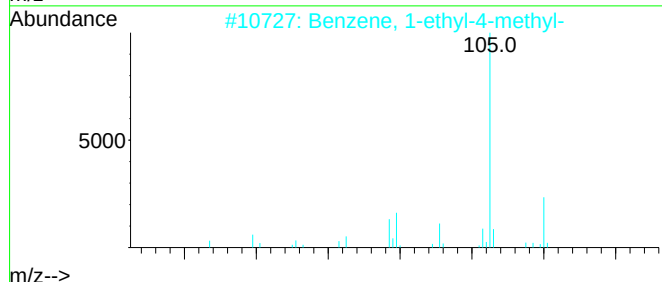
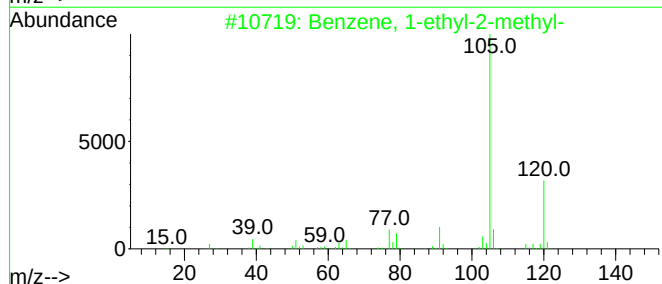
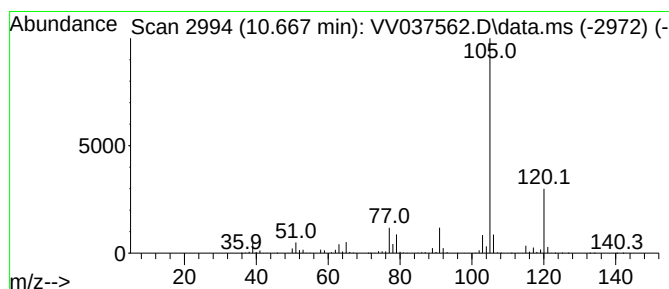
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.667	94.19 ug/L	4348930	1,4-Dichlorobenzene-d4	11.178

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
4	Mesitylene	120	C9H12	000108-67-8	90
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY040

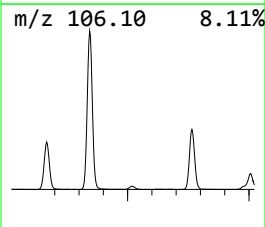
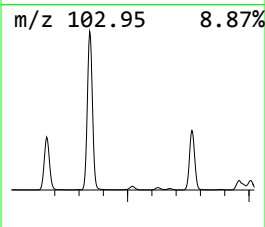
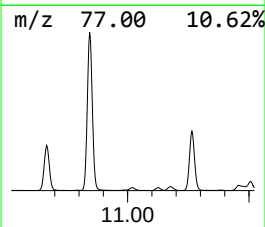
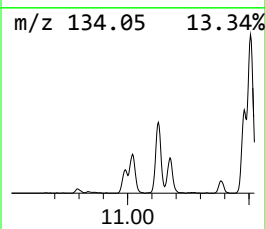
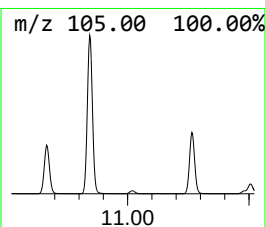
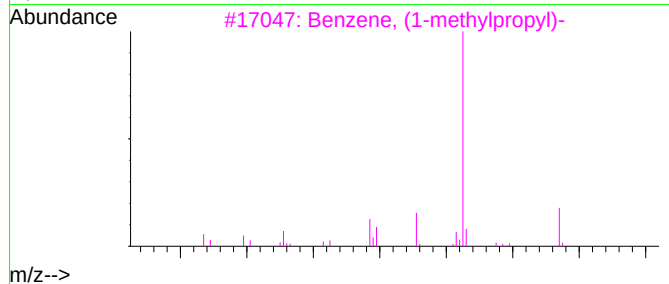
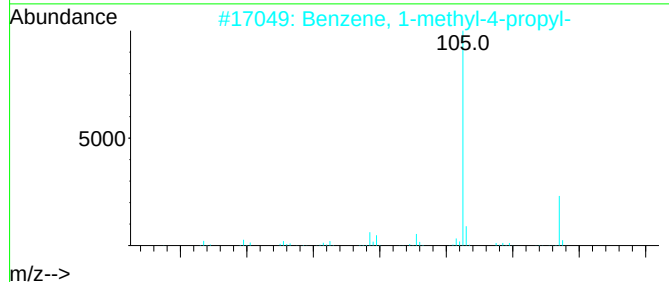
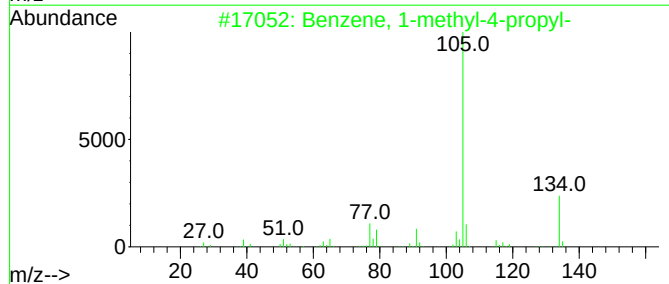
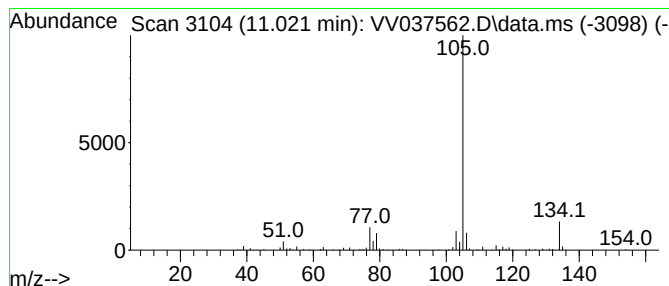
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-methyl-4-propyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.021	5.86 ug/L	270547	1,4-Dichlorobenzene-d4		11.178	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-		134	C10H14	001074-55-1	91
2	Benzene, 1-methyl-4-propyl-		134	C10H14	001074-55-1	86
3	Benzene, (1-methylpropyl)-		134	C10H14	000135-98-8	86
4	Benzeneacetaldehyde, .alpha.-met...		134	C9H10O	000093-53-8	83
5	Benzene, 1-methyl-4-propyl-		134	C10H14	001074-55-1	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY040

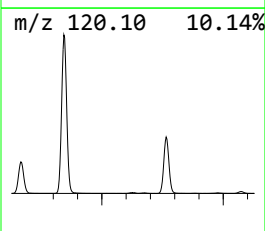
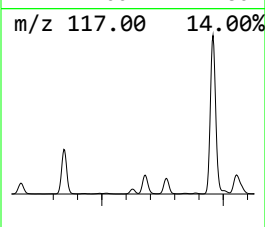
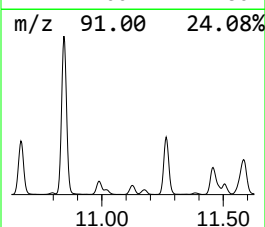
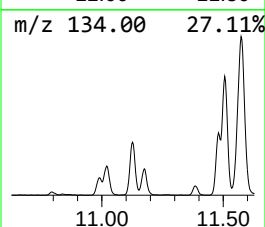
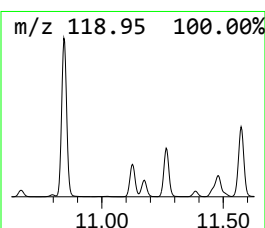
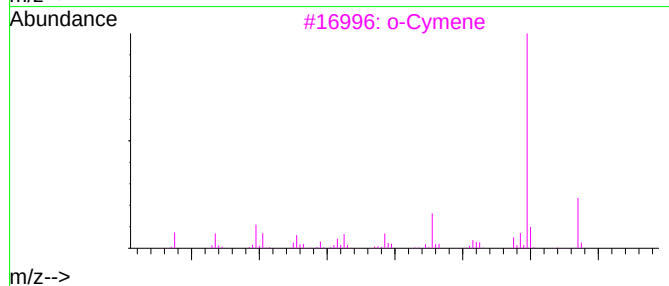
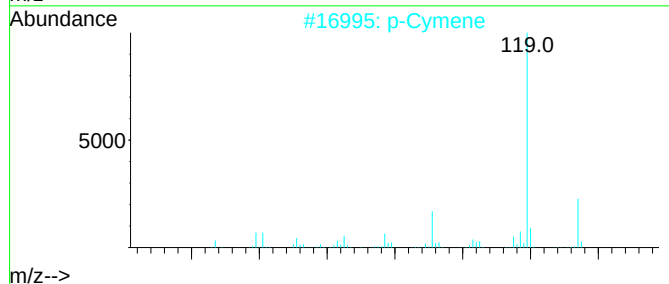
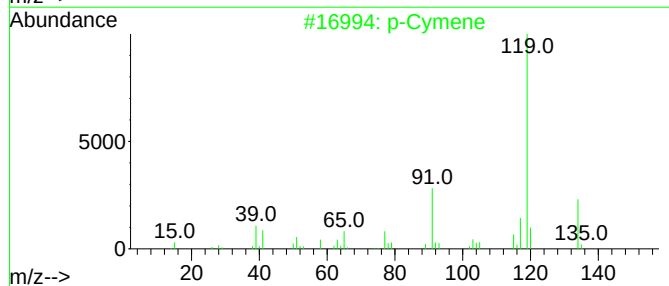
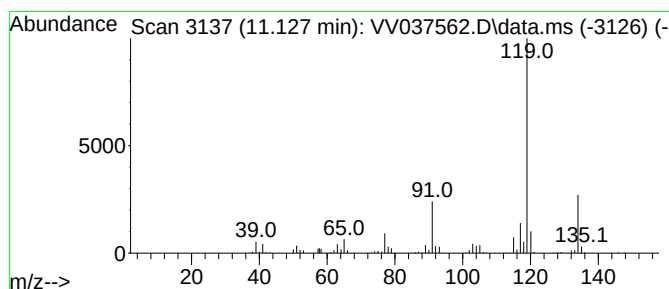
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 7 p-Cymene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.127	8.65 ug/L	399188	1,4-Dichlorobenzene-d4		11.178	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	p-Cymene		134	C10H14	000099-87-6	97
2	p-Cymene		134	C10H14	000099-87-6	95
3	o-Cymene		134	C10H14	000527-84-4	95
4	p-Cymene		134	C10H14	000099-87-6	95
5	o-Cymene		134	C10H14	000527-84-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY040

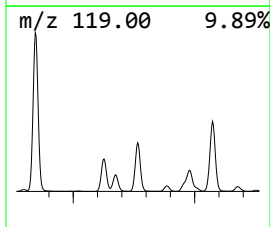
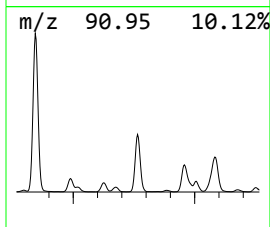
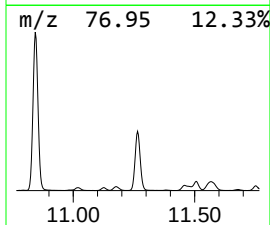
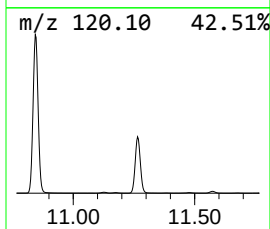
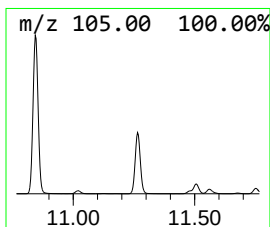
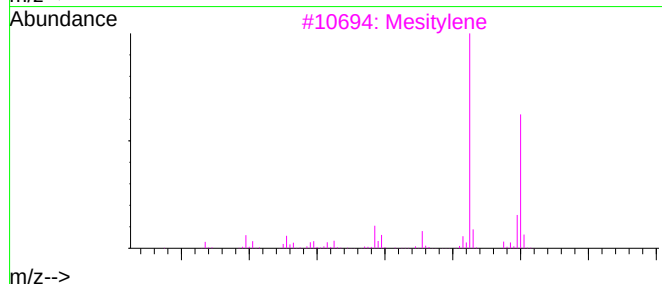
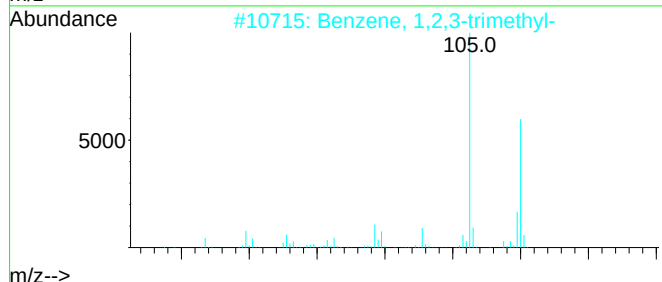
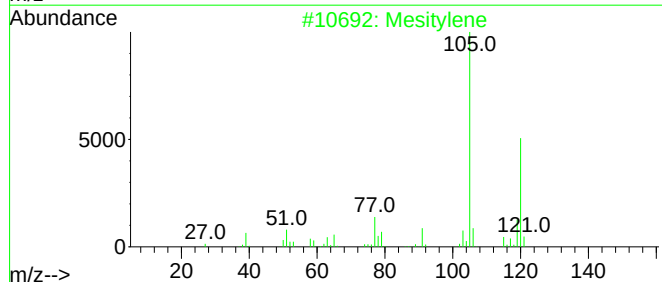
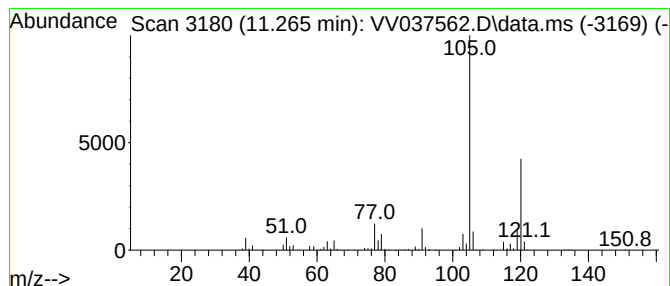
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.265	121.52 ug/L	5610940	1,4-Dichlorobenzene-d4	11.178	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Mesitylene	120	C9H12	000108-67-8	94
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5	Mesitylene	120	C9H12	000108-67-8	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
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ALS Vial : 1 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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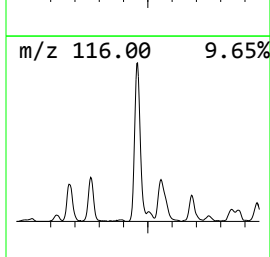
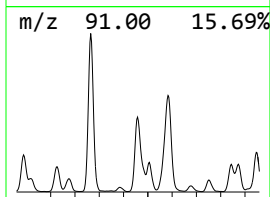
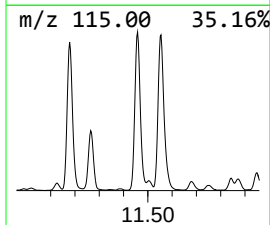
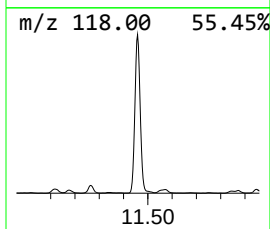
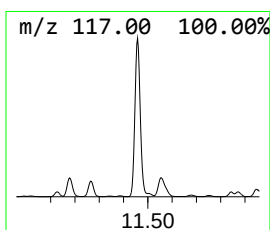
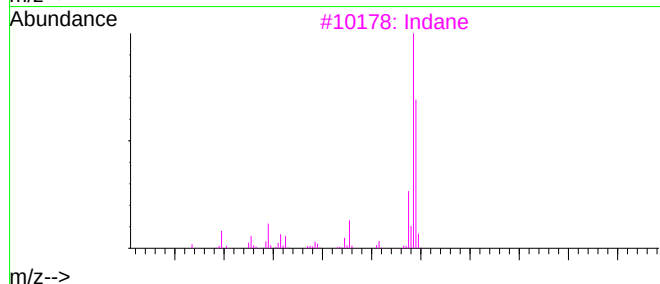
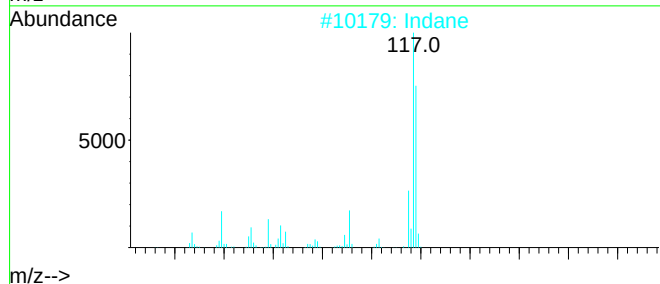
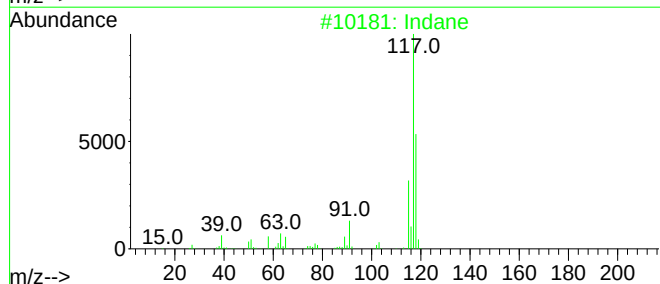
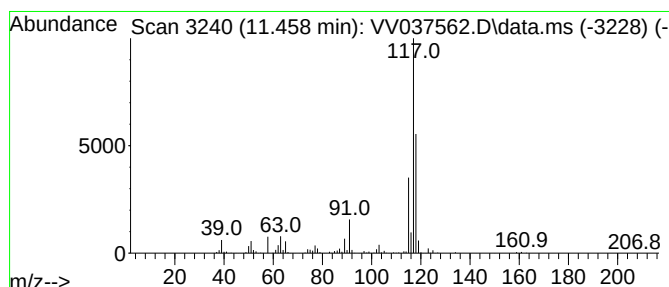
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 9 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.458	53.37 ug/L	2464090	1,4-Dichlorobenzene-d4		11.178	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	95
2	Indane		118	C9H10	000496-11-7	91
3	Indane		118	C9H10	000496-11-7	90
4	Indane		118	C9H10	000496-11-7	90
5	Benzene, 1-ethenyl-2-methyl-		118	C9H10	000611-15-4	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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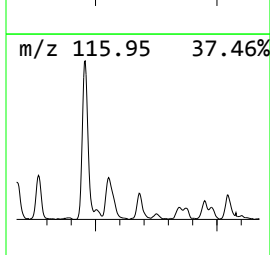
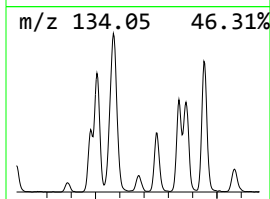
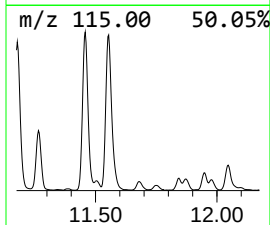
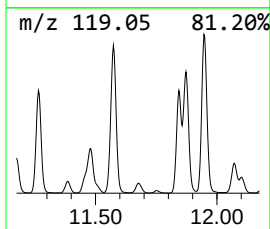
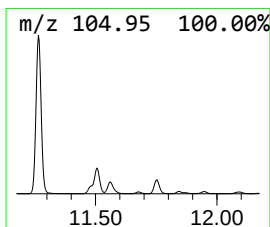
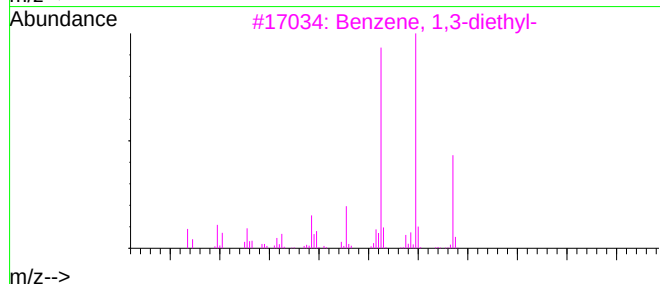
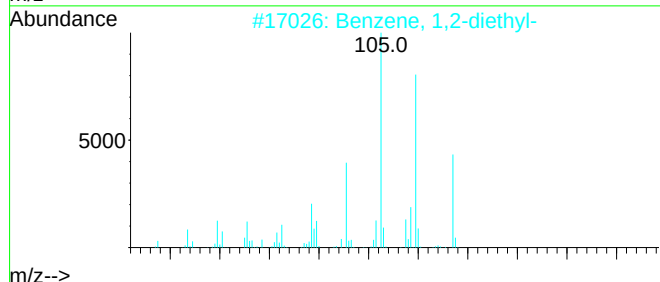
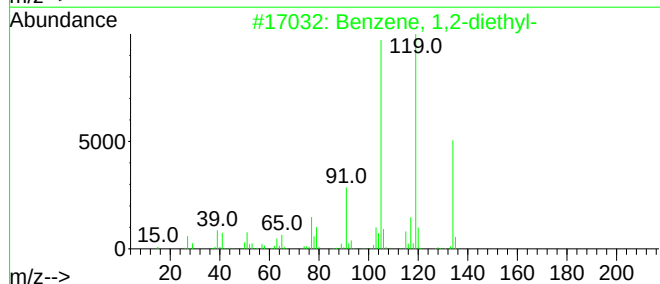
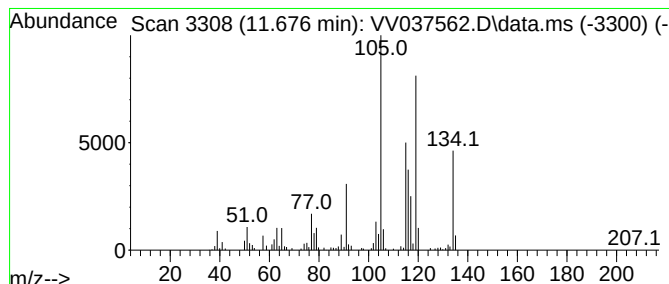
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 1,2-diethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.676	3.20 ug/L	147793	1,4-Dichlorobenzene-d4	11.178

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	92
4	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	76
5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
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Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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Quant Title : VOC Analysis

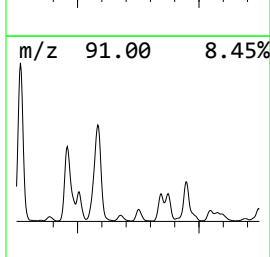
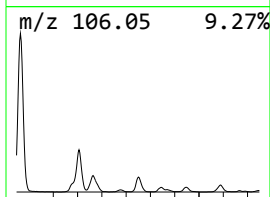
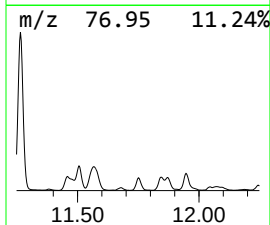
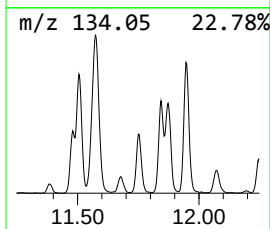
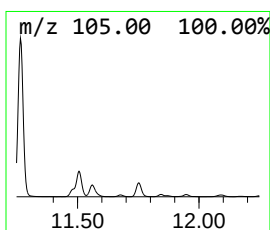
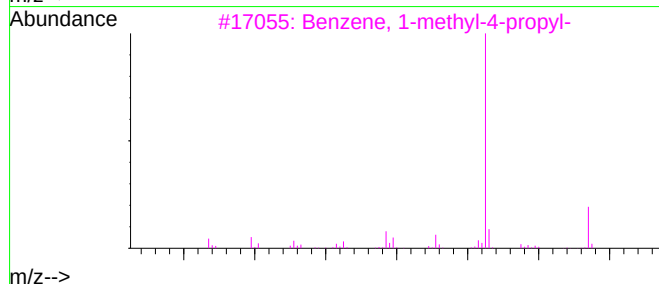
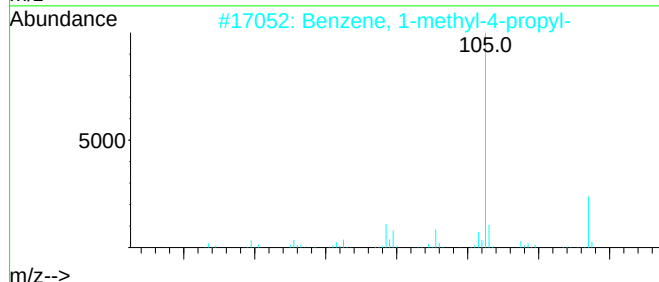
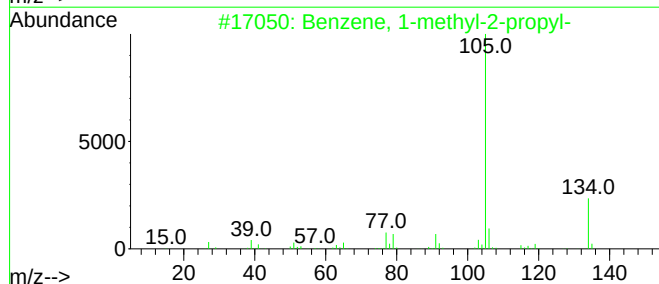
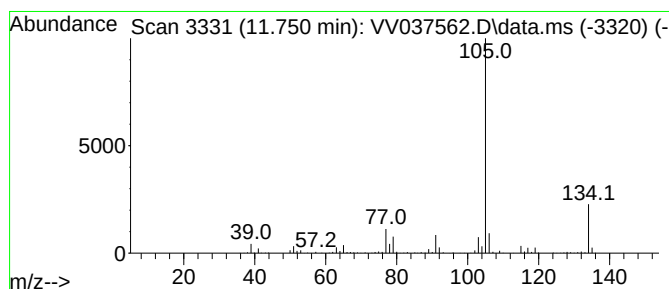
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 11 Benzene, 1-methyl-2-propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.750	9.50 ug/L	438659	1,4-Dichlorobenzene-d4	11.178

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	94
2	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	94
3	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
4	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY040

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM092624WMA.M
Quant Title : VOC Analysis

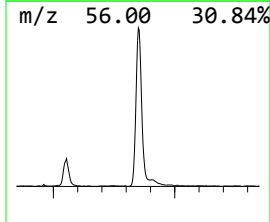
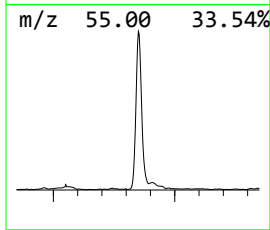
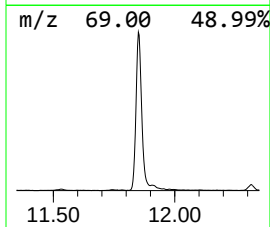
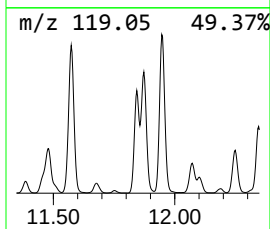
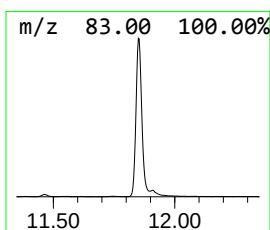
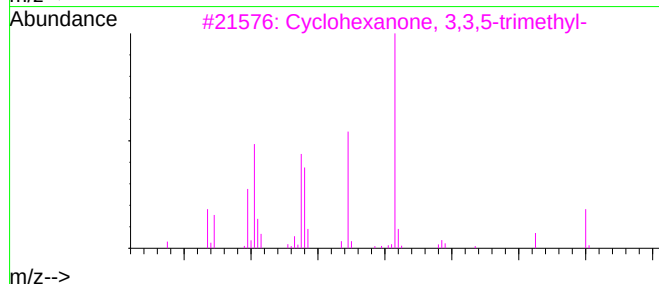
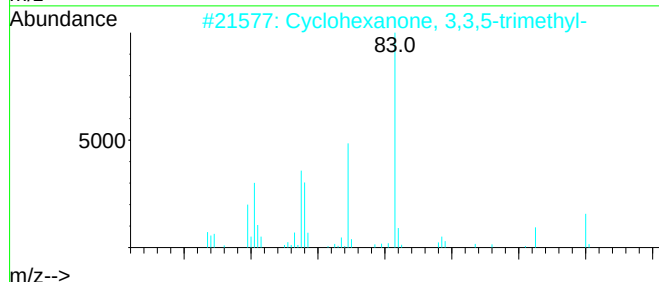
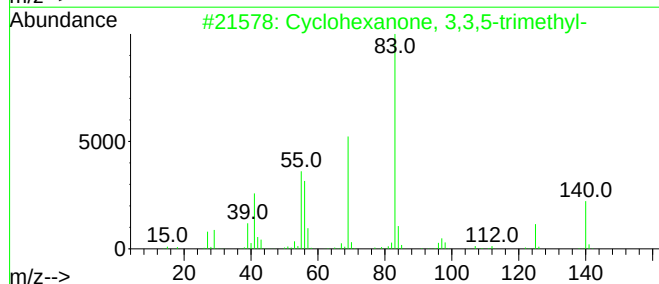
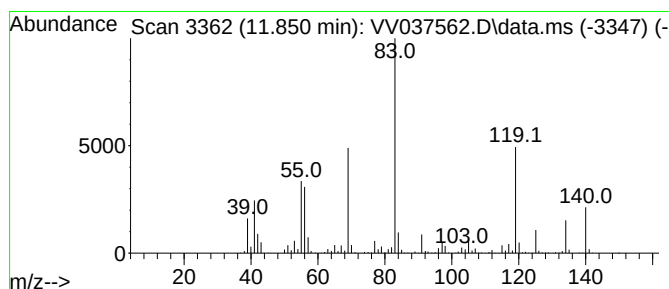
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 12 Cyclohexanone, 3,3,5-trimet... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.850	56.65 ug/L	2615570	1,4-Dichlorobenzene-d4	11.178

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	96
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	94
3	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	70
4	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	70
5	Verbenyl angelate, cis-	234	C15H22O2	1000383-56-3	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY040

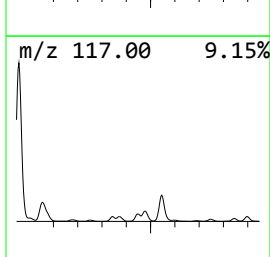
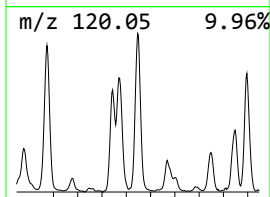
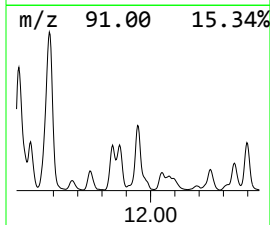
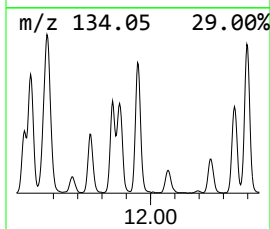
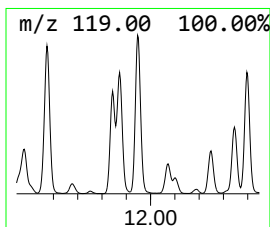
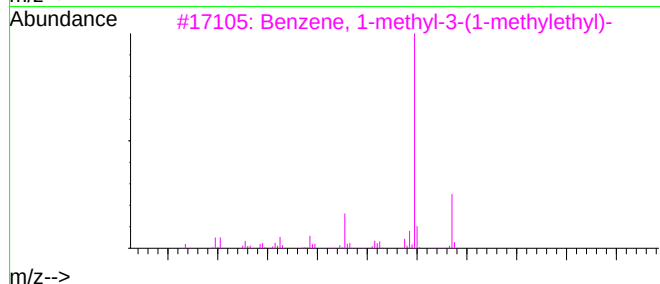
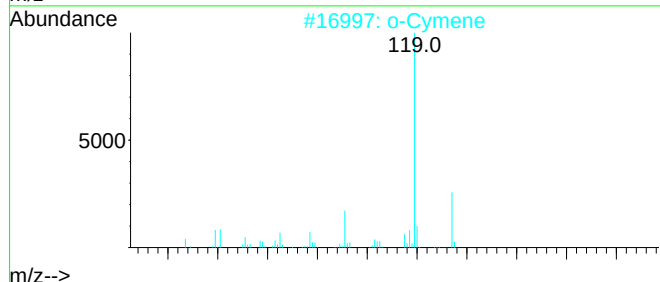
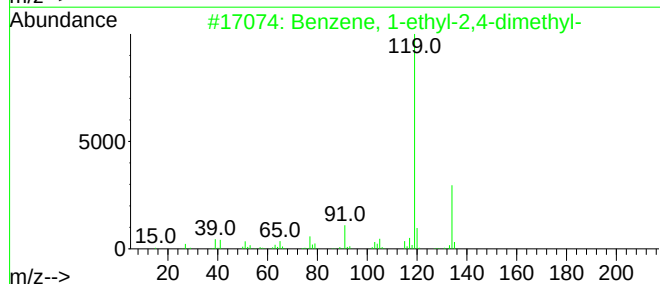
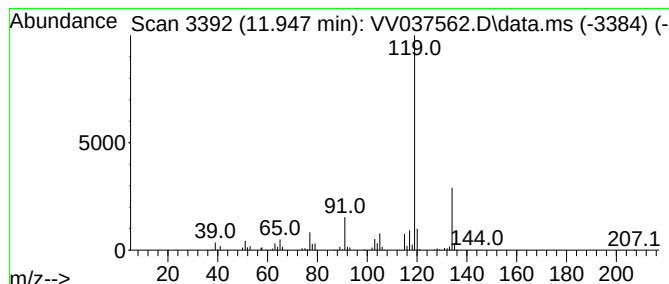
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 13 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.947	23.04 ug/L	1063600	1,4-Dichlorobenzene-d4	11.178	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5	o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY040

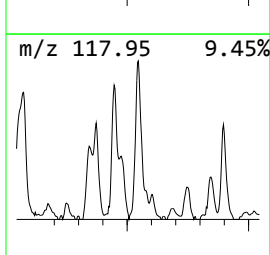
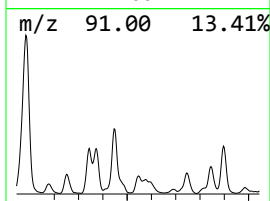
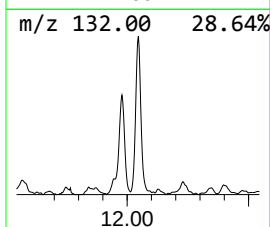
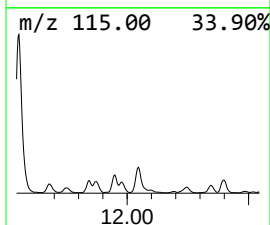
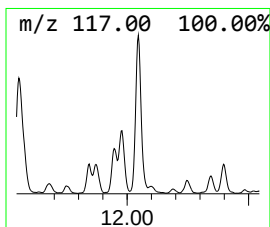
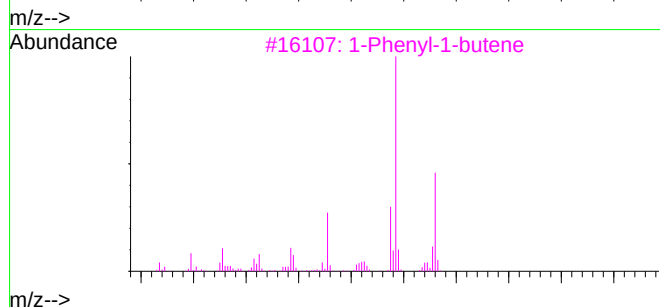
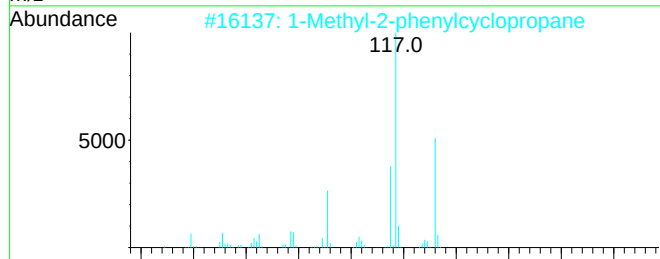
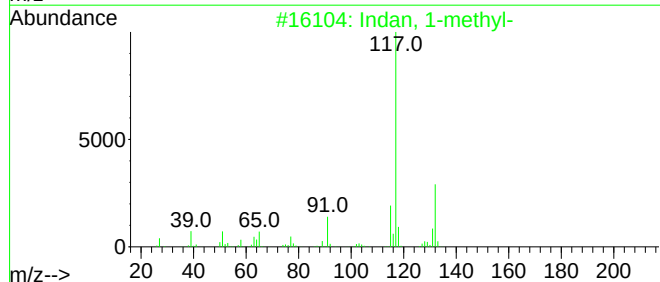
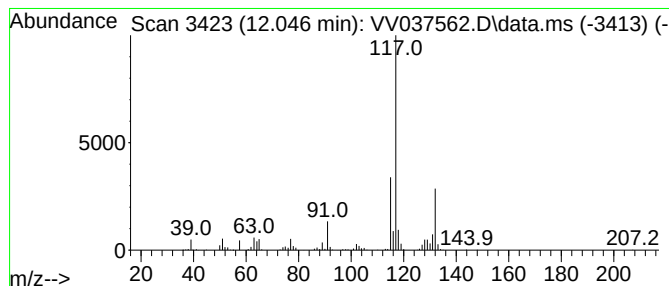
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 14 Indan, 1-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.046	7.43 ug/L	342879	1,4-Dichlorobenzene-d4		11.178	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-		132	C10H12	000767-58-8	90
2	1-Methyl-2-phenylcyclopropane		132	C10H12	003145-76-4	90
3	1-Phenyl-1-butene		132	C10H12	000824-90-8	87
4	Benzene, 1-ethenyl-4-ethyl-		132	C10H12	003454-07-7	83
5	Indan, 1-methyl-		132	C10H12	000767-58-8	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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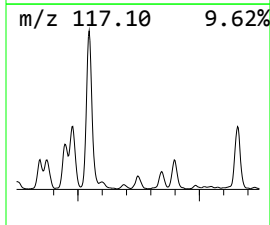
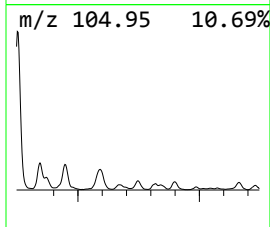
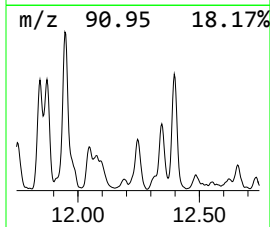
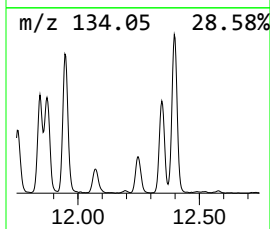
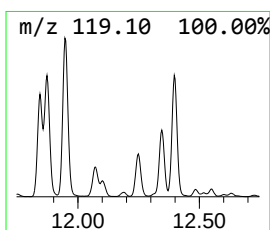
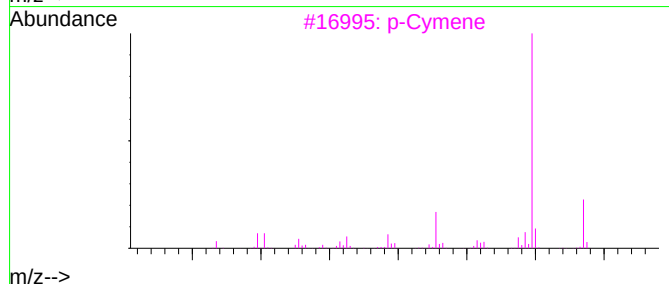
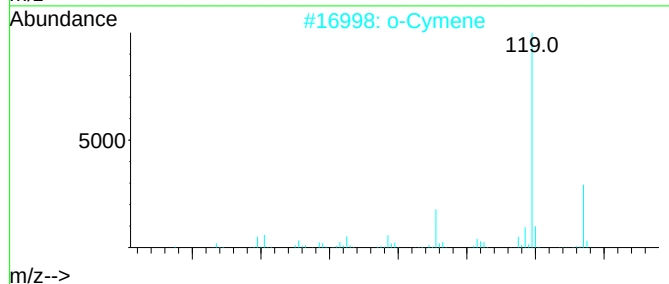
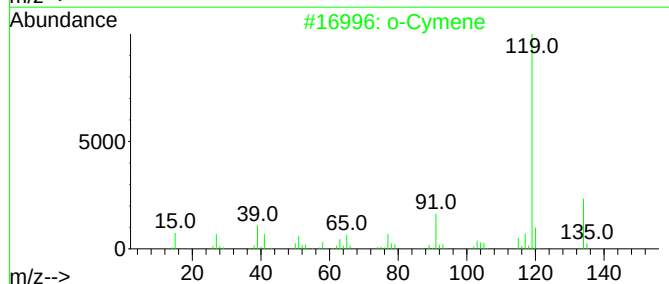
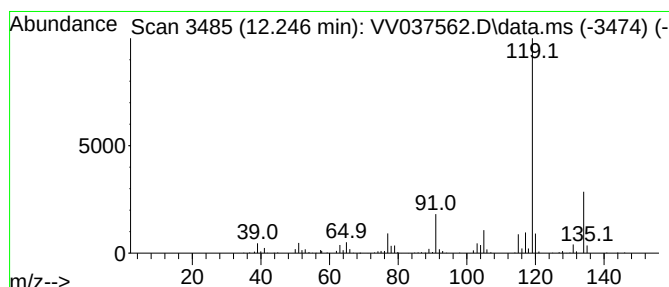
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 15 o-Cymene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.246	6.42 ug/L	296253	1,4-Dichlorobenzene-d4	11.178

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	95
2	o-Cymene	134	C10H14	000527-84-4	95
3	p-Cymene	134	C10H14	000099-87-6	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY040

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM092624WMA.M
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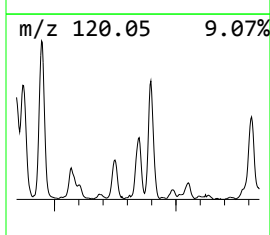
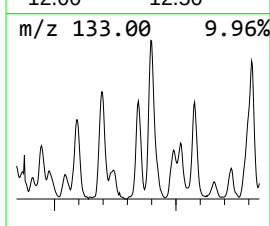
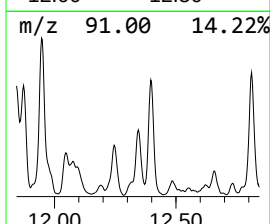
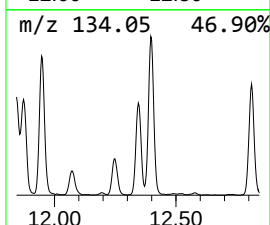
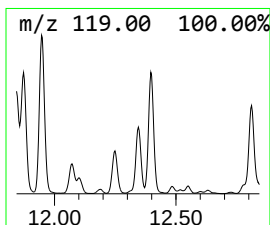
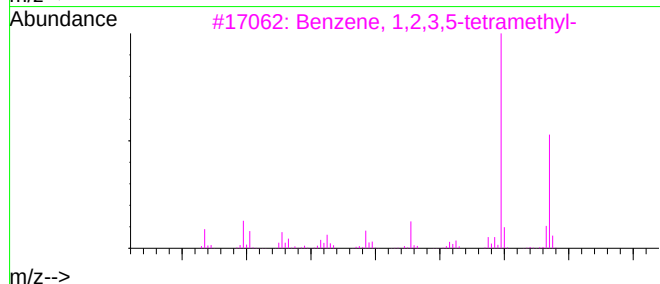
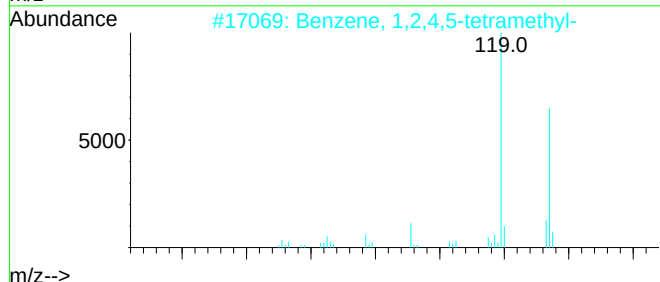
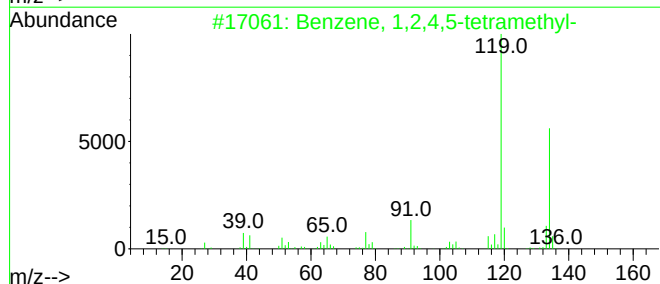
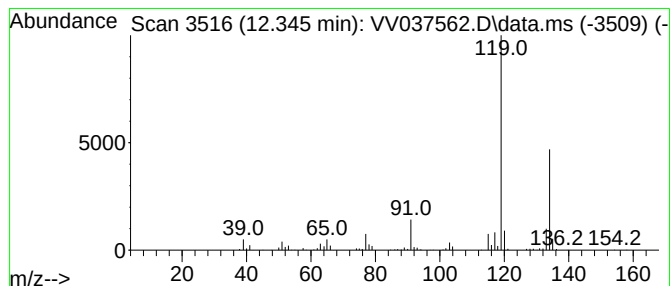
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 16 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.345	8.19 ug/L	378172	1,4-Dichlorobenzene-d4	11.178

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY040

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM092624WMA.M
Quant Title : VOC Analysis

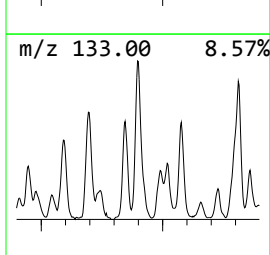
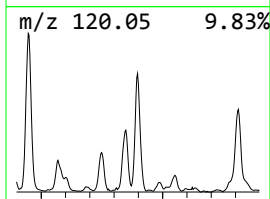
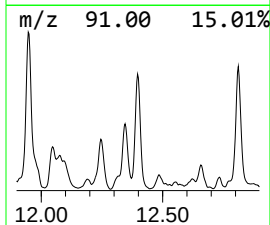
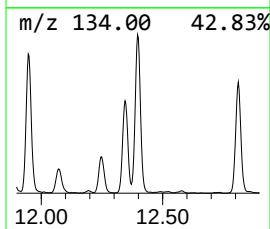
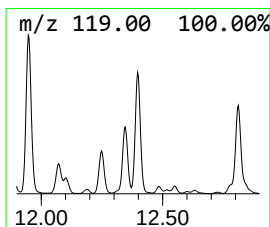
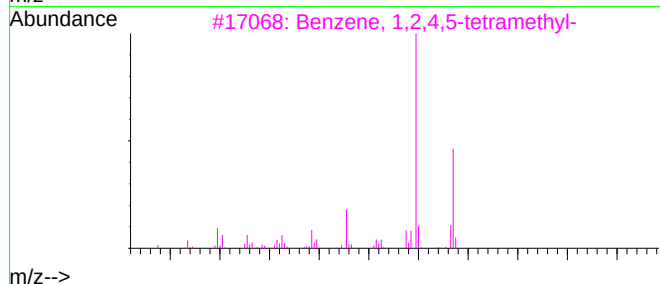
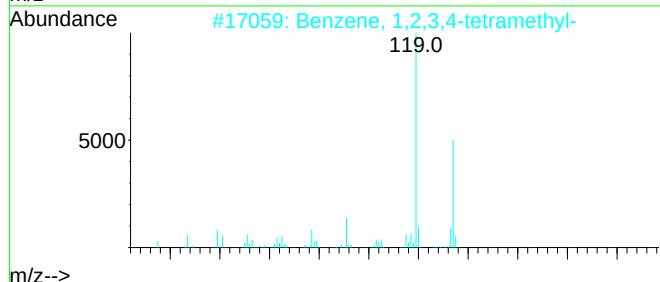
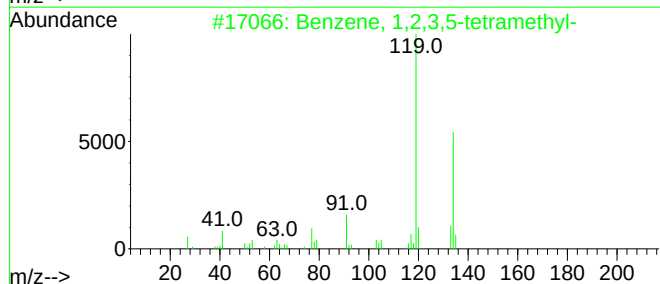
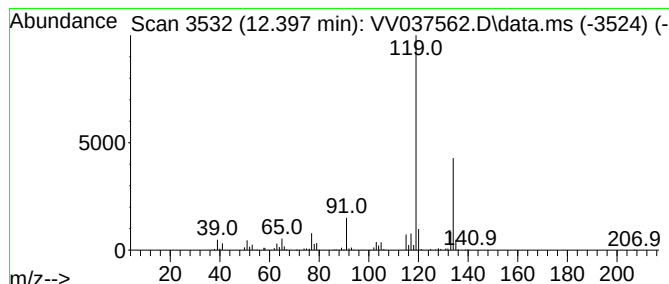
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 17 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.397	15.34 ug/L	708345	1,4-Dichlorobenzene-d4	11.178

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY040

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM092624WMA.M
Quant Title : VOC Analysis

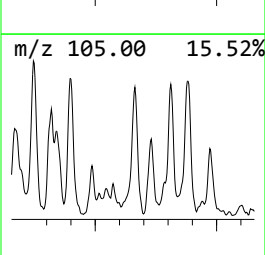
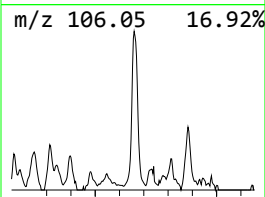
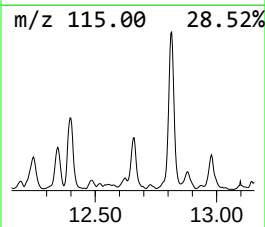
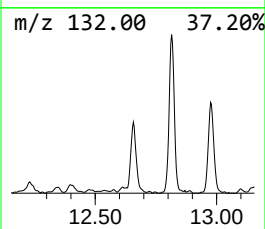
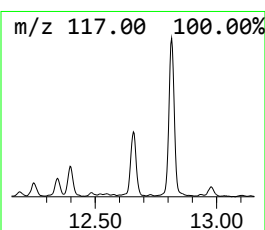
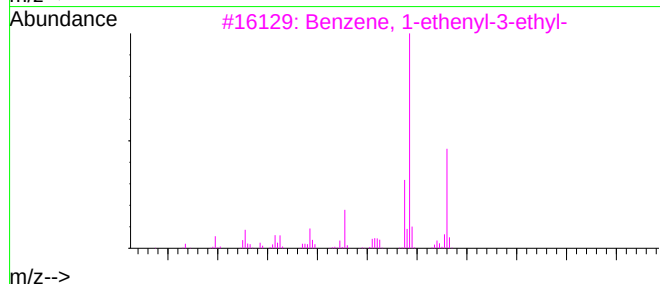
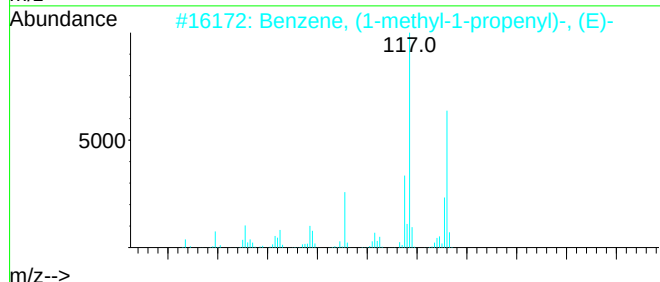
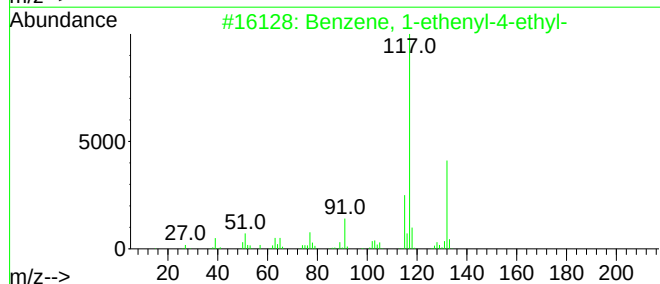
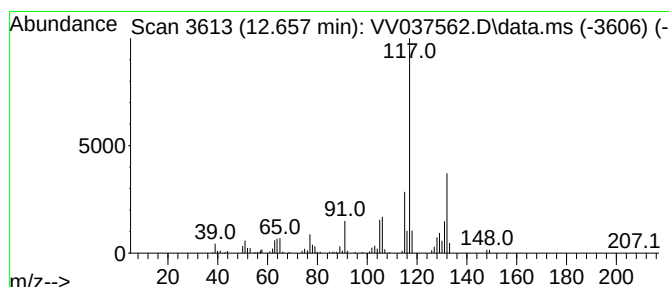
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 18 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.657	4.00 ug/L	184834	1,4-Dichlorobenzene-d4	11.178

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4	Indan, 1-methyl-	132	C10H12	000767-58-8	87
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY040

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM092624WMA.M
Quant Title : VOC Analysis

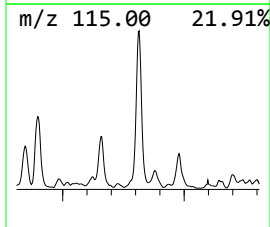
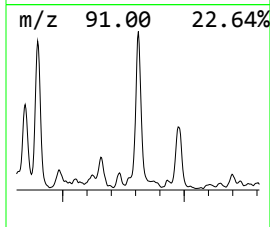
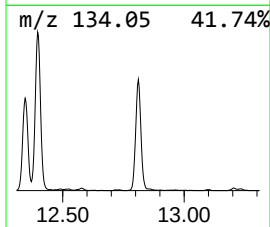
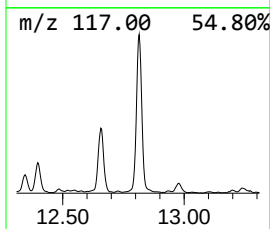
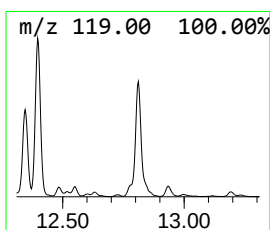
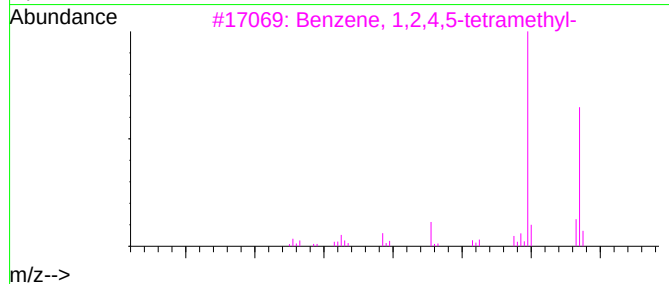
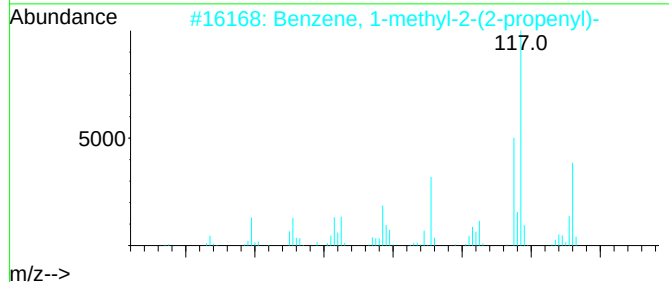
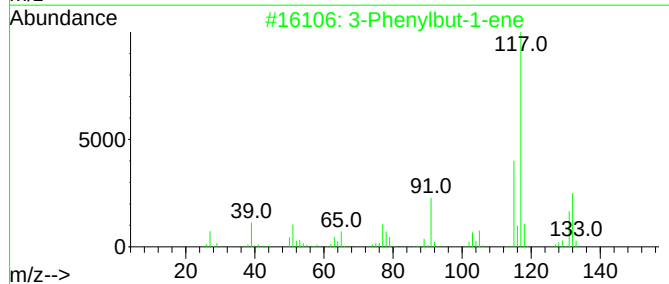
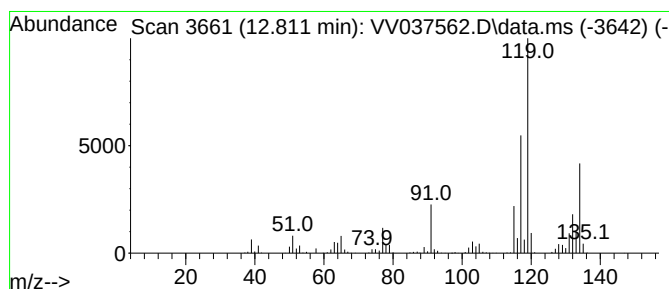
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 19 3-Phenylbut-1-ene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.811	18.79 ug/L	867743	1,4-Dichlorobenzene-d4	11.178

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY040

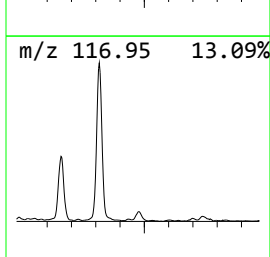
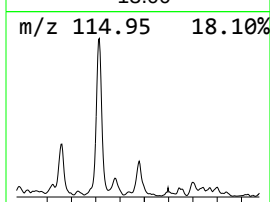
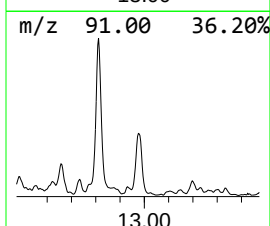
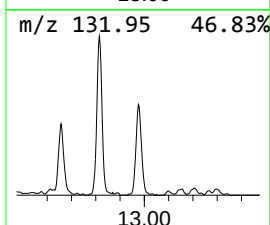
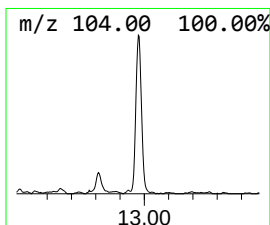
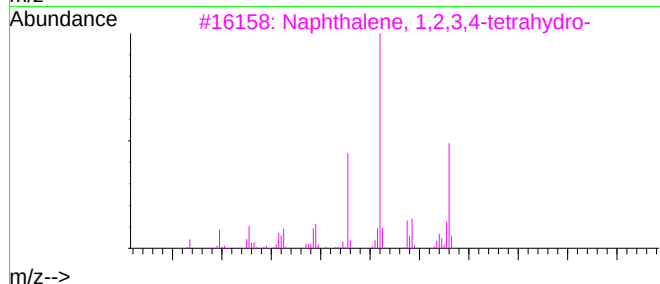
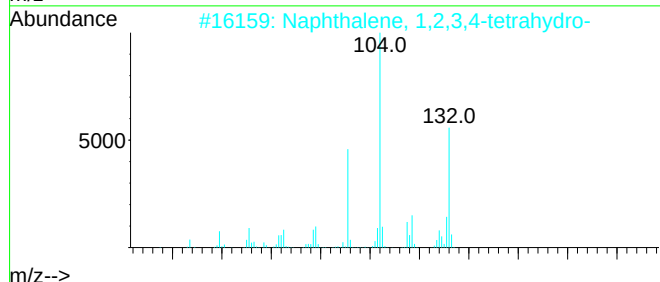
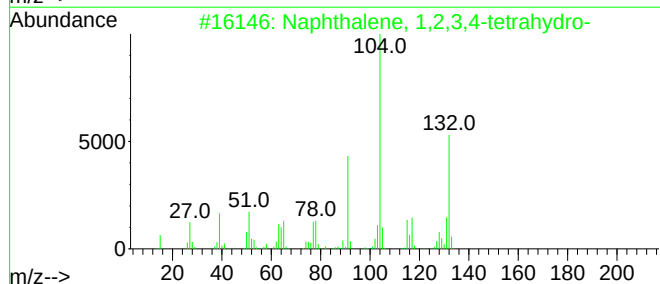
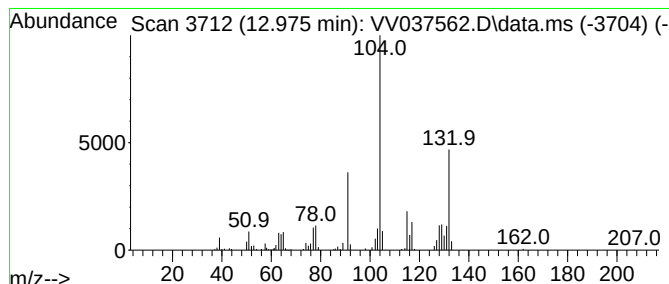
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 20 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.975	4.64 ug/L	214102	1,4-Dichlorobenzene-d4	11.178	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
3	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
4	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY040

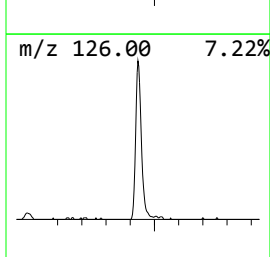
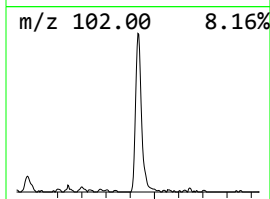
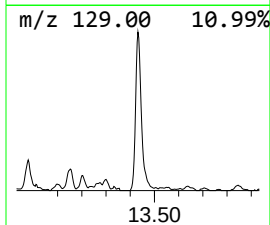
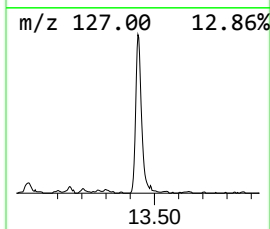
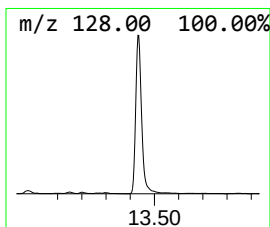
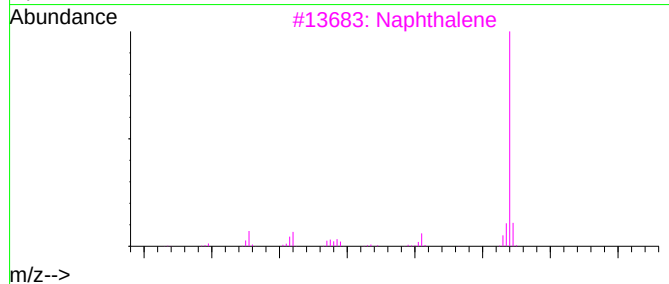
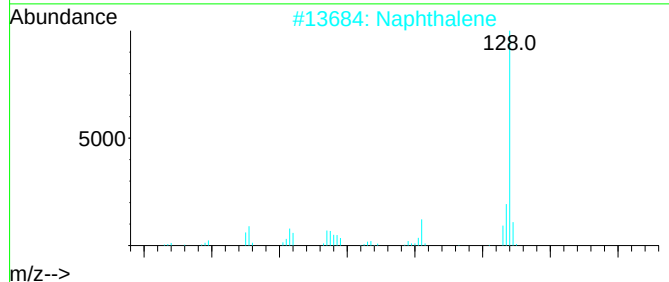
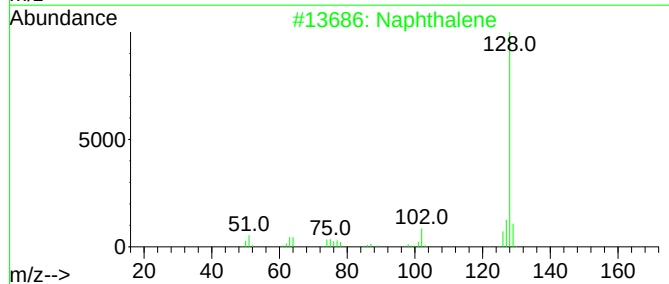
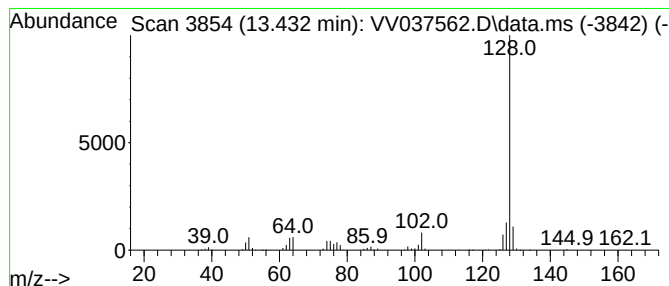
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 21 Azulene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.432	13.66 ug/L	630785	1,4-Dichlorobenzene-d4		11.178	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene		128	C10H8	000091-20-3	95
2	Naphthalene		128	C10H8	000091-20-3	94
3	Naphthalene		128	C10H8	000091-20-3	93
4	Azulene		128	C10H8	000275-51-4	91
5	Azulene		128	C10H8	000275-51-4	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY040

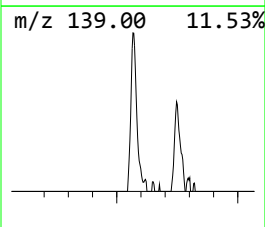
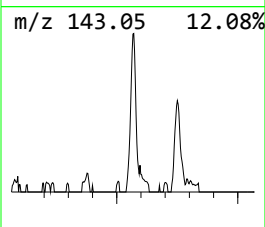
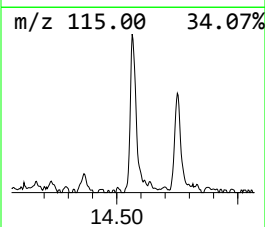
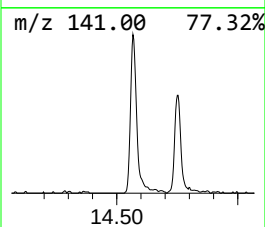
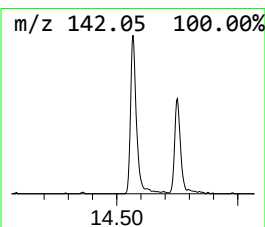
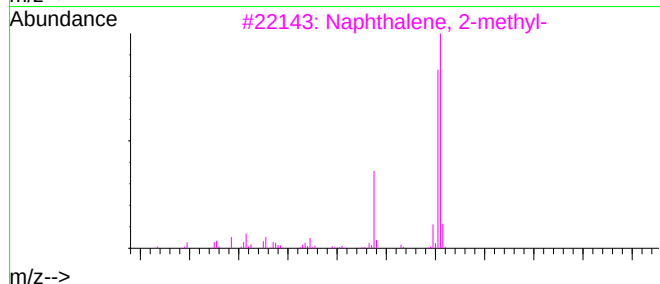
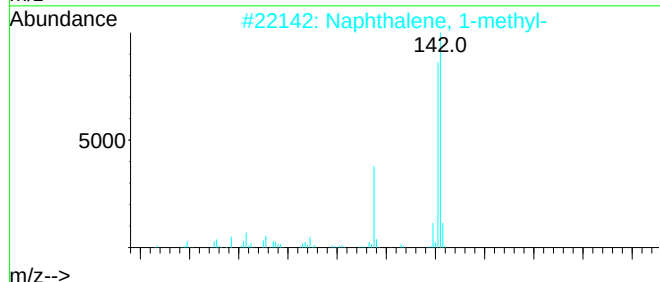
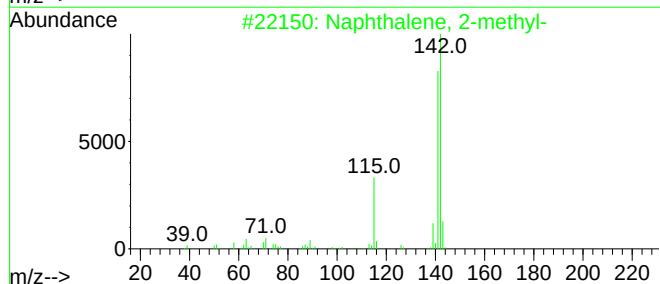
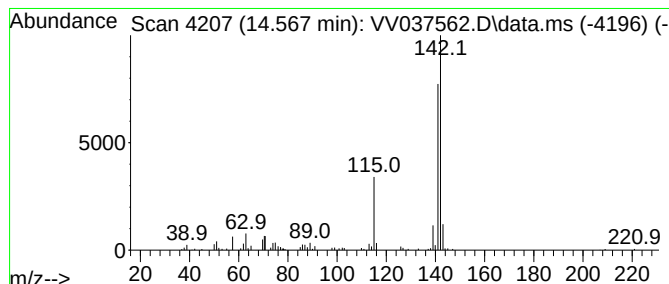
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 22 Naphthalene, 2-methyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.567	2.68 ug/L	123716	1,4-Dichlorobenzene-d4		11.178	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-		142	C11H10	000091-57-6	97
2	Naphthalene, 1-methyl-		142	C11H10	000090-12-0	96
3	Naphthalene, 2-methyl-		142	C11H10	000091-57-6	96
4	Naphthalene, 2-methyl-		142	C11H10	000091-57-6	95
5	Naphthalene, 1-methyl-		142	C11H10	000090-12-0	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092624\
Data File : VV037562.D
Acq On : 26 Sep 2024 20:26
Operator : SY/MD
Sample : P4184-14
Misc : 5mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY040

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM092624WMA.M
Quant Title : VOC Analysis

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Fluorodichlorom...	1.667	3.3	ug/L	102458	1	5.529	1567140	50.0
Benzene, propyl-	10.284	37.5	ug/L	1732110	3	11.178	2308640	50.0
Benzene, 1-ethy...	10.378	143.5	ug/L	6625120	3	11.178	2308640	50.0
Benzene, 1-ethy...	10.667	94.2	ug/L	4348930	3	11.178	2308640	50.0
Benzene, 1-meth...	11.021	5.9	ug/L	270547	3	11.178	2308640	50.0
p-Cymene	11.127	8.7	ug/L	399188	3	11.178	2308640	50.0
Benzene, 1,2,3-...	11.265	121.5	ug/L	5610940	3	11.178	2308640	50.0
Indane	11.458	53.4	ug/L	2464090	3	11.178	2308640	50.0
Benzene, 1,2-di...	11.676	3.2	ug/L	147793	3	11.178	2308640	50.0
Benzene, 1-meth...	11.750	9.5	ug/L	438659	3	11.178	2308640	50.0
Cyclohexanone, ...	11.850	56.6	ug/L	2615570	3	11.178	2308640	50.0
Benzene, 1-ethy...	11.947	23.0	ug/L	1063600	3	11.178	2308640	50.0
Indan, 1-methyl-	12.046	7.4	ug/L	342879	3	11.178	2308640	50.0
o-Cymene	12.246	6.4	ug/L	296253	3	11.178	2308640	50.0
Benzene, 1,2,4,...	12.345	8.2	ug/L	378172	3	11.178	2308640	50.0
Benzene, 1,2,3,...	12.397	15.3	ug/L	708345	3	11.178	2308640	50.0
Benzene, 1-ethe...	12.657	4.0	ug/L	184834	3	11.178	2308640	50.0
3-Phenylbut-1-ene	12.811	18.8	ug/L	867743	3	11.178	2308640	50.0
Naphthalene, 1,...	12.975	4.6	ug/L	214102	3	11.178	2308640	50.0
Azulene	13.432	13.7	ug/L	630785	3	11.178	2308640	50.0
Naphthalene, 2-...	14.567	2.7	ug/L	123716	3	11.178	2308640	50.0