LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE8

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 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_V\Method\SFAMVTR110421WMA.M
Title : TRACE VOA SFAM1.0

Signal : TIC: VV023346.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.304	74	82	92	rBV2	50132	53985	3.04%	0.544%
2	1.564	156	163	176	rBV	40524	45588	2.57%	0.460%
3	2.108	321	332	343	rBV	75939	107488	6.05%	1.084%
4	3.191	651	669	686	rBV	12847	28287	1.59%	0.285%
5	3.905	872	891	918	rBV2	79751	290370	16.35%	2.927%
6	4.352	1010	1030	1059	rBV3	78170	216613	12.20%	2.184%
7	4.609	1093	1110	1128	rBV3	30427	76257	4.29%	0.769%
8	5.047	1229	1246	1281	rBV2	149386	387046	21.80%	3.902%
9	5.619	1412	1424	1444	rBV	152199	307503	17.32%	3.100%
10	5.915	1502	1516	1545	rBV	273467	538934	30.35%	5.433%
11	6.072	1552	1565	1579	rBV	87806	183888	10.36%	1.854%
12	6.989	1838	1850	1871	rBV2	43038	80544	4.54%	0.812%
13	7.140	1886	1897	1911	rVB3	9124	18147	1.02%	0.183%
14	7.317	1941	1952	1965	rBV	172129	298395	16.80%	3.008%
15	7.391	1965	1975	1997	rVB	130242	230901	13.00%	2.328%
16	7.625	2039	2048	2071	rBV2	26426	50721	2.86%	0.511%
17	7.979	2149	2158	2172	rVB4	13453	24997	1.41%	0.252%
18	8.088	2181	2192	2224	rBV	230967	413540	23.29%	4.169%
19	8.854	2419	2430	2435	rBV	251908	409920	23.08%	4.132%
20	9.146	2509	2521	2533	rVB2	11173	22089	1.24%	0.223%
21	9.481	2609	2625	2637	rBV9	6655	19133	1.08%	0.193%
22	9.548	2637	2646	2661	rVB	36318	62335	3.51%	0.628%
23	9.706	2677	2695	2705	rVB4	12301	26582	1.50%	0.268%
24	9.799	2713	2724	2741	rBV4	9000	21299	1.20%	0.215%
25	9.934	2756	2766	2789	rBV	66763	110666	6.23%	1.116%
26	10.217	2842	2854	2874	rBV	75576	140668	7.92%	1.418%
27	10.355	2886	2897	2916	rBV	193603	315040	17.74%	3.176%
28	10.448	2916	2926	2944	rVV	82734	188753	10.63%	1.903%
29	10.542	2948	2955	2969	rVB3	10992	22348	1.26%	0.225%
30	10.738	3004	3016	3034	rBV	372281	573680	32.31%	5.783%
31	10.915	3059	3071	3094	rVB	1186417	1775770	100.00%	17.900%
32	11.056	3094	3115	3120	rBV	39271	70307	3.96%	0.709%
33	11.088	3120	3125	3137	rVB	52506	78635	4.43%	0.793%
34	11.194	3151	3158	3165	rVB3	17571	24908	1.40%	0.251%
35	11.249	3165	3175	3192	rBV	284672	486285	27.38%	4.902%
36	11.336	3192	3202	3218	rVB	359065	532017	29.96%	5.363%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE8

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 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_V\Method\SFAMVTR110421WMA.M
Title : TRACE VOA SFAM1.0

37	11.455	3231	3239	3249	rVB3	14788	22496	1.27%	0.227%
38	11.529	3249	3262	3282	rBV3	213662	515437	29.03%	5.196%
39	11.625	3282	3292	3312	rVB4	270858	548047	30.86%	5.524%
40	11.744	3320	3329	3340	rVB2	16648	26352	1.48%	0.266%
41	11.821	3340	3353	3366	rVB	34136	53279	3.00%	0.537%
42	11.911	3370	3381	3386	rBV	48363	72755	4.10%	0.733%
43	11.940	3387	3390	3399	rVB	32579	40337	2.27%	0.407%
44	12.017	3400	3414	3420	rBV	56946	87484	4.93%	0.882%
45	12.117	3434	3445	3453	rBV2	44988	82635	4.65%	0.833%
46	12.255	3475	3488	3496	rBV6	10410	24716	1.39%	0.249%
47	12.307	3496	3504	3513	rVV7	12479	22590	1.27%	0.228%
48	12.413	3531	3537	3546	rVB2	21528	30455	1.72%	0.307%
49	12.468	3546	3554	3570	rVB2	26352	44140	2.49%	0.445%
50	12.728	3626	3635	3650	rVB2	13636	24450	1.38%	0.246%
51	12.882	3674	3683	3694	rVB3	31781	50674	2.85%	0.511%
52	13.050	3725	3735	3747	rVB5	10830	19065	1.07%	0.192%
53	13.519	3864	3881	3898	rBV8	7594	21785	1.23%	0.220%

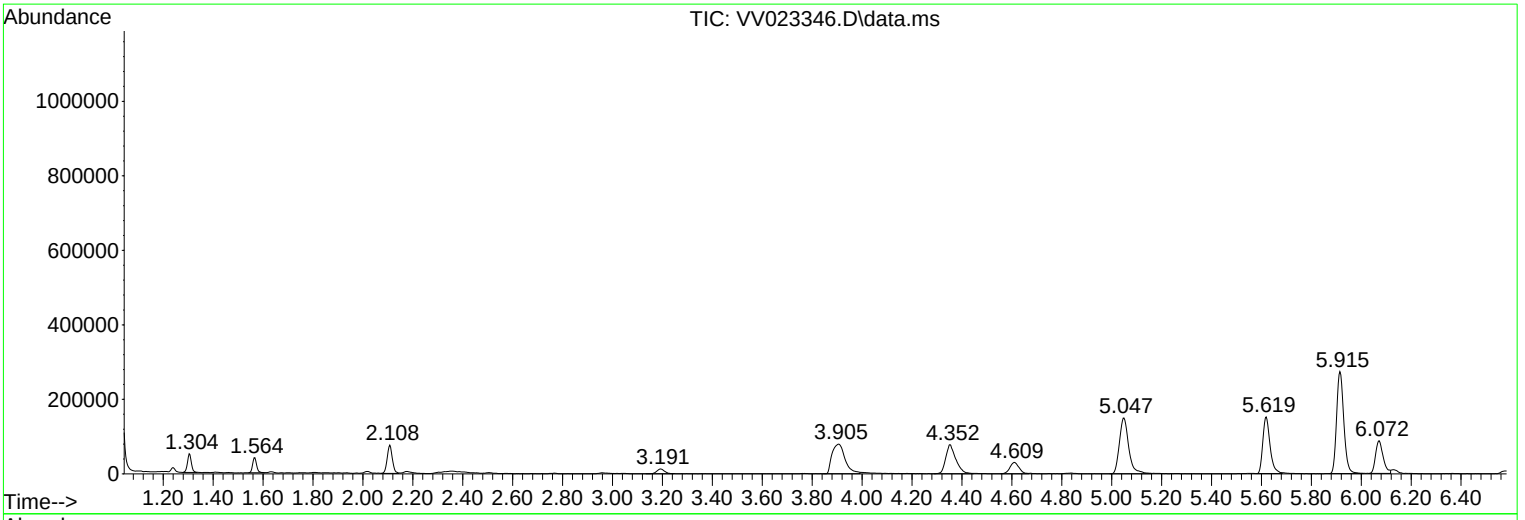
Sum of corrected areas: 9920336

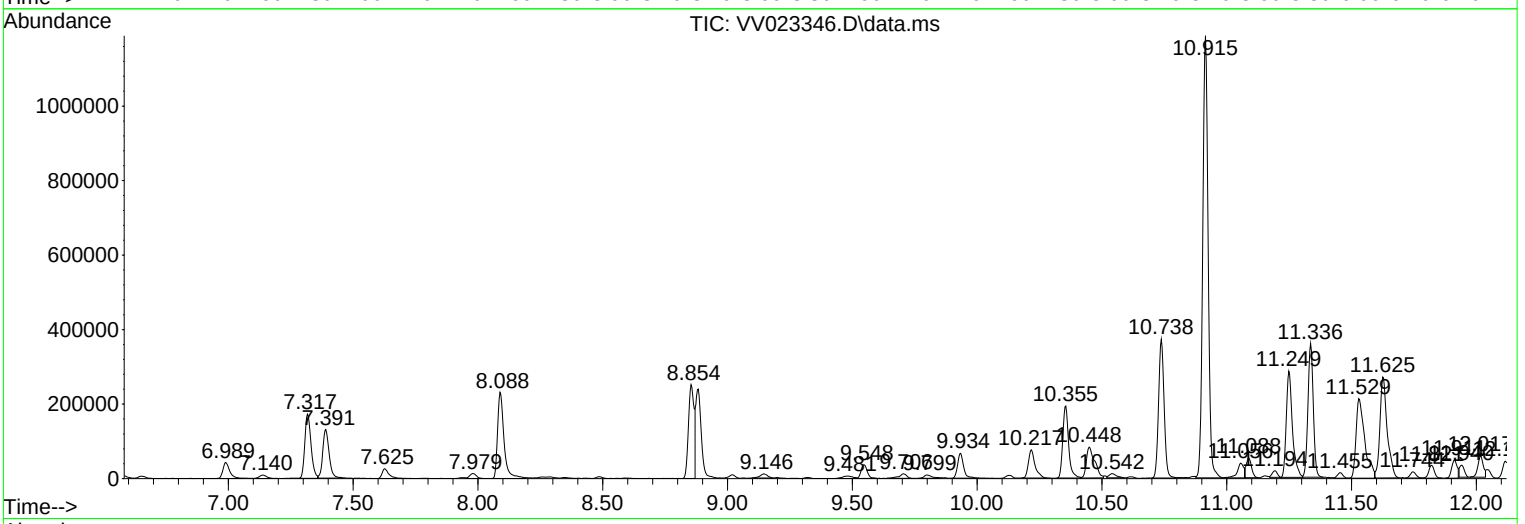
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

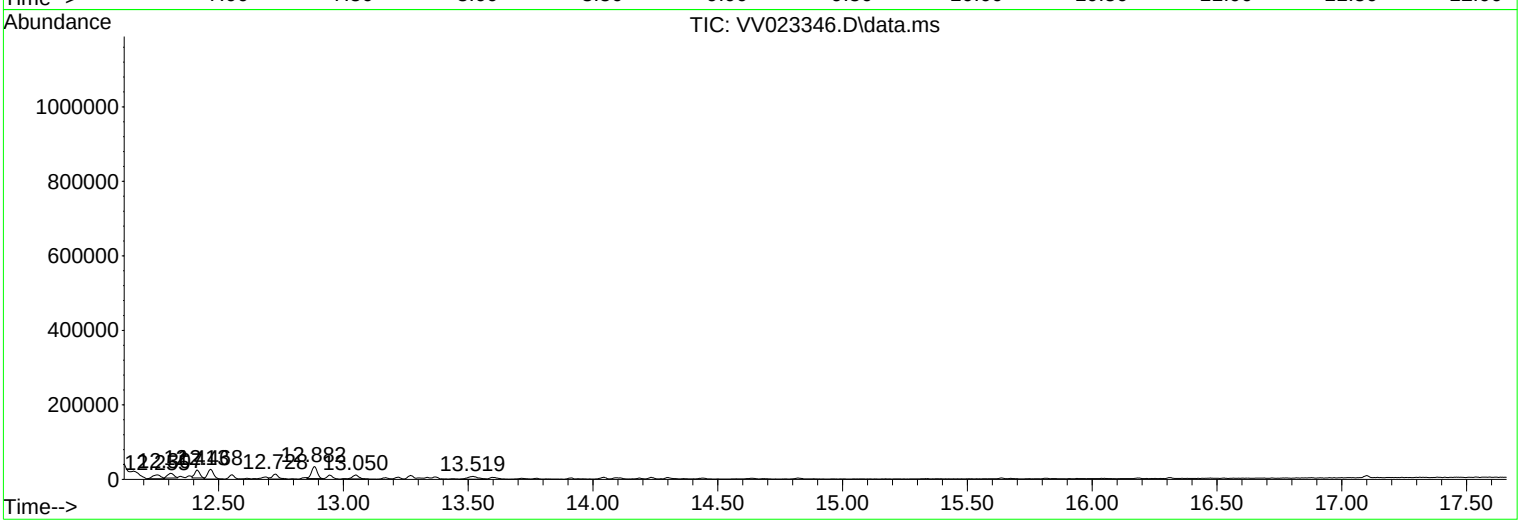
Instrument :
MSVOA_V
ClientSampleId :
BGGE8

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

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







Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE8

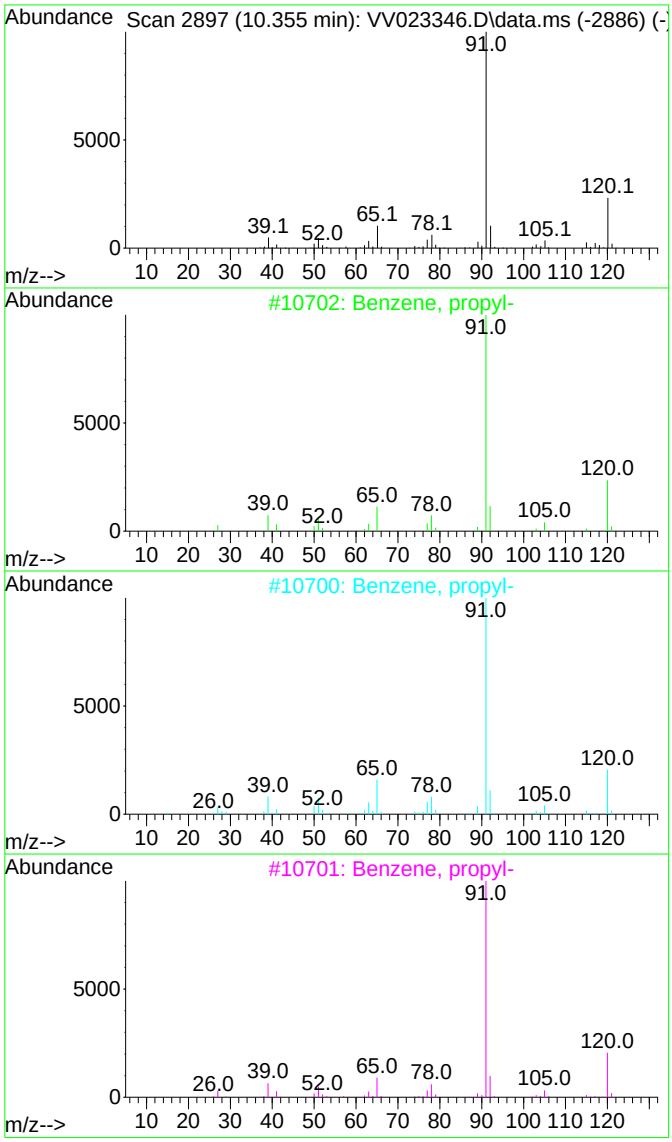
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

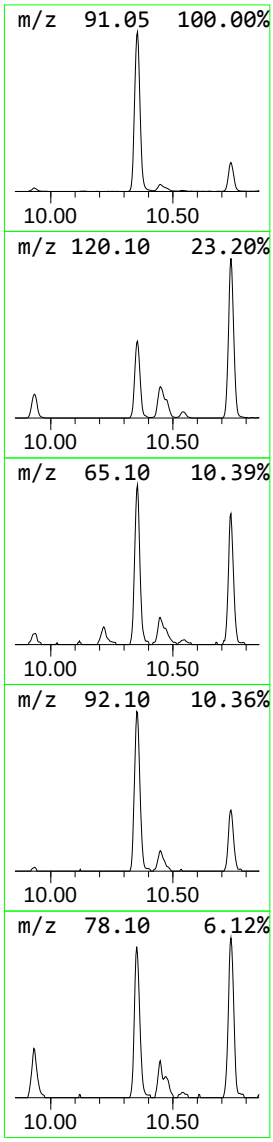
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, propyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.355	3.24 ug/L	315040	1,4-Dichlorobenzene-d4	11.249

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	90
3	Benzene, propyl-	120	C9H12	000103-65-1	87
4	Benzene, propyl-	120	C9H12	000103-65-1	87
5	1,2-Ethanediamine, N-(phenylmeth...	150	C9H14N2	004152-09-4	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE8

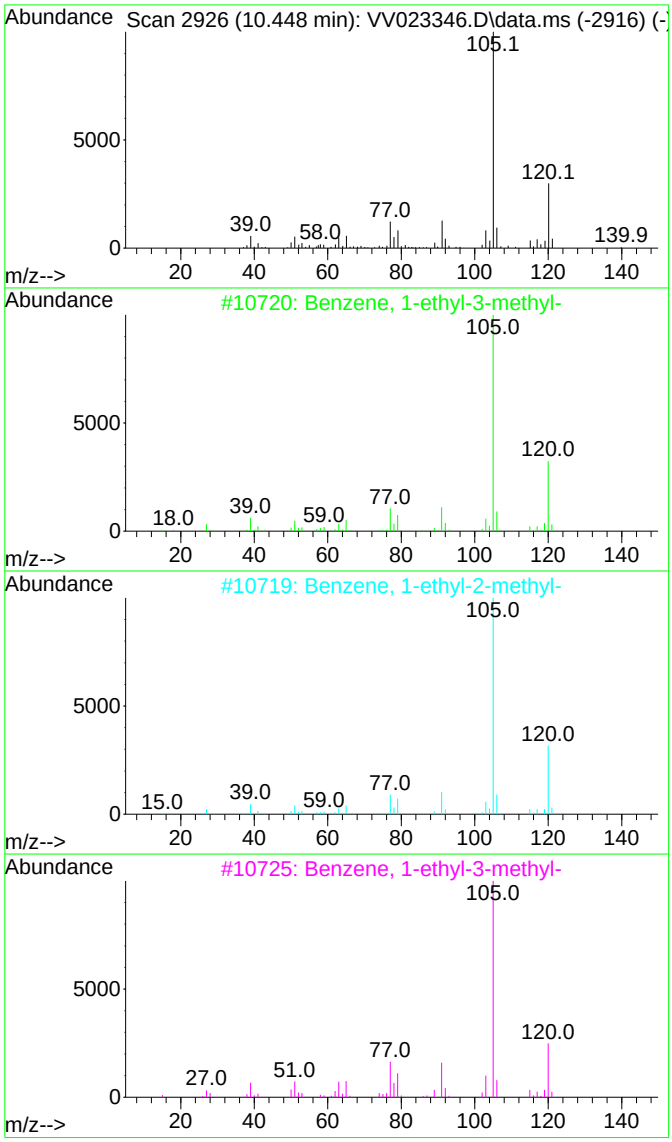
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

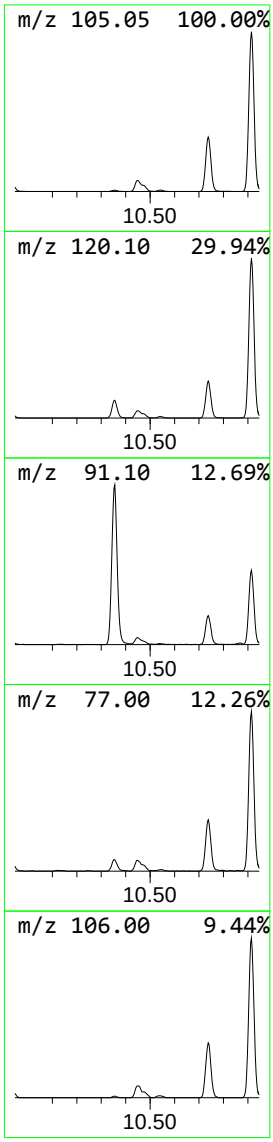
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.448	1.94 ug/L	188753	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE8

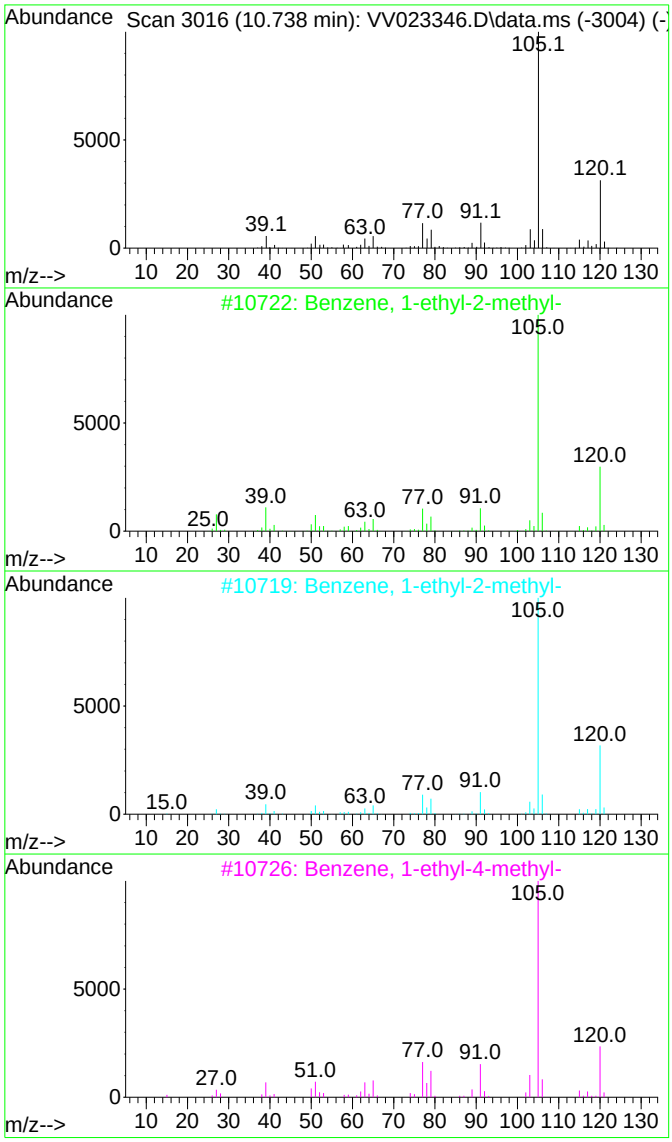
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

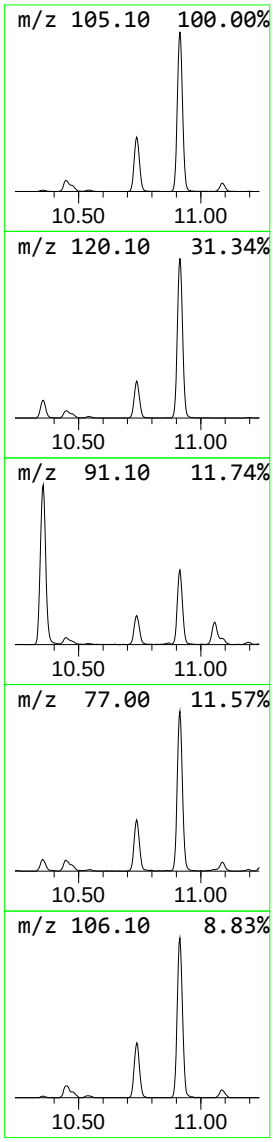
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.738	5.90 ug/L	573680	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE8

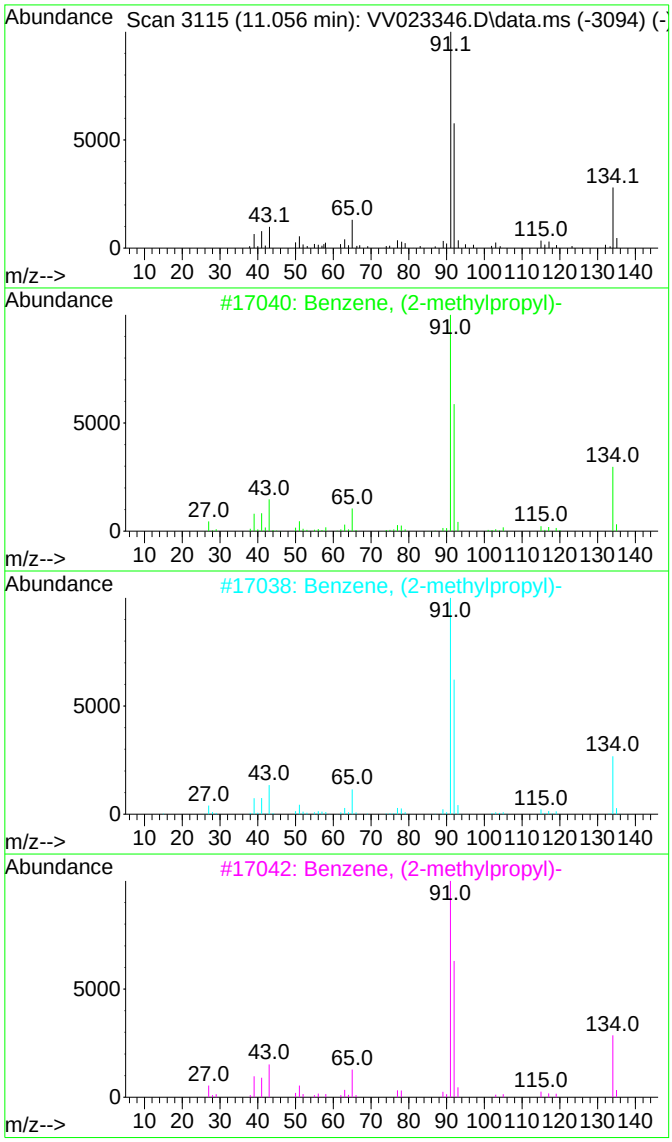
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

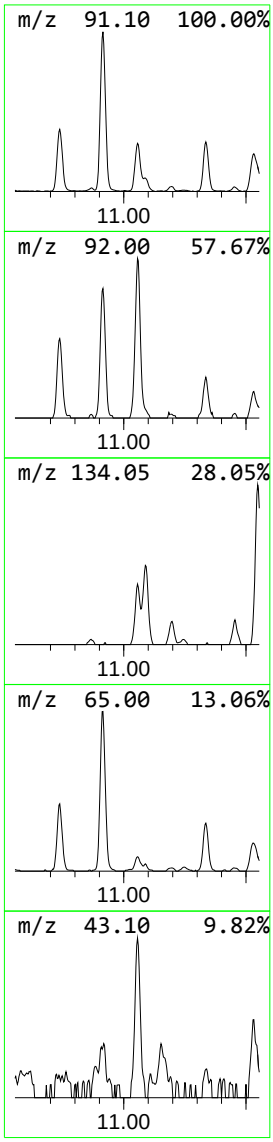
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, (2-methylpropyl)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.056	0.72 ug/L	70307	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
2	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
3	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
4	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	87
5	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE8

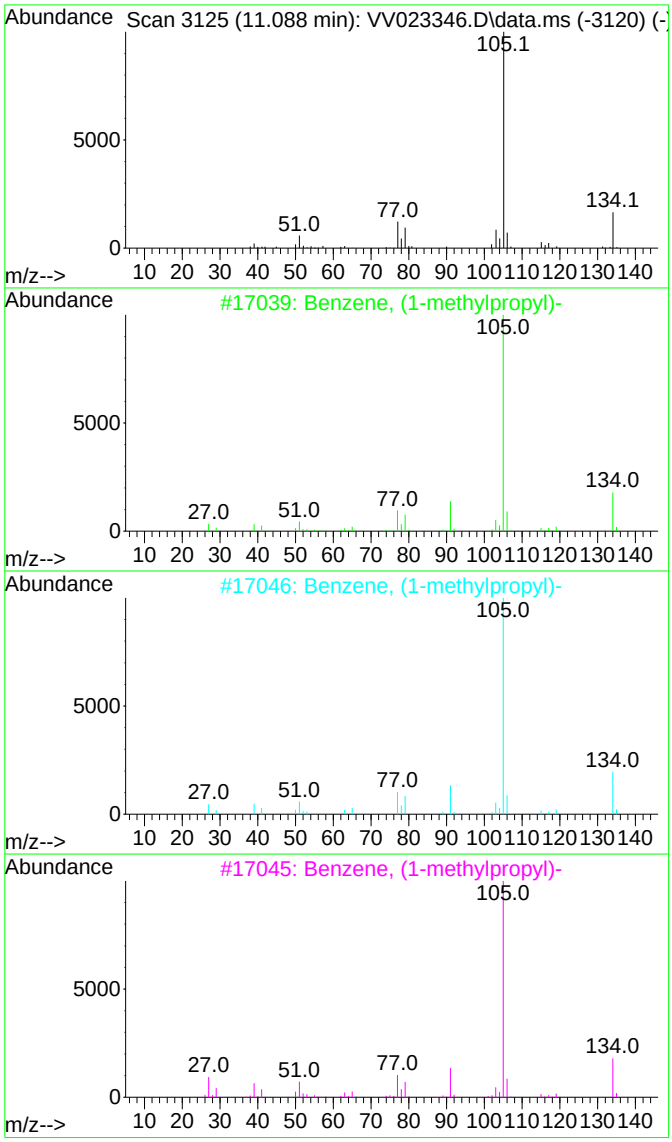
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

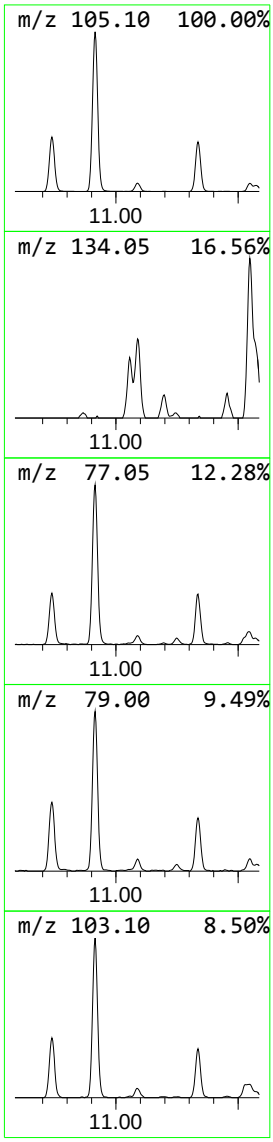
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, (1-methylpropyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.088	0.81 ug/L	78635	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	86
2	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	86
3	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	78
4	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	78
5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE8

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

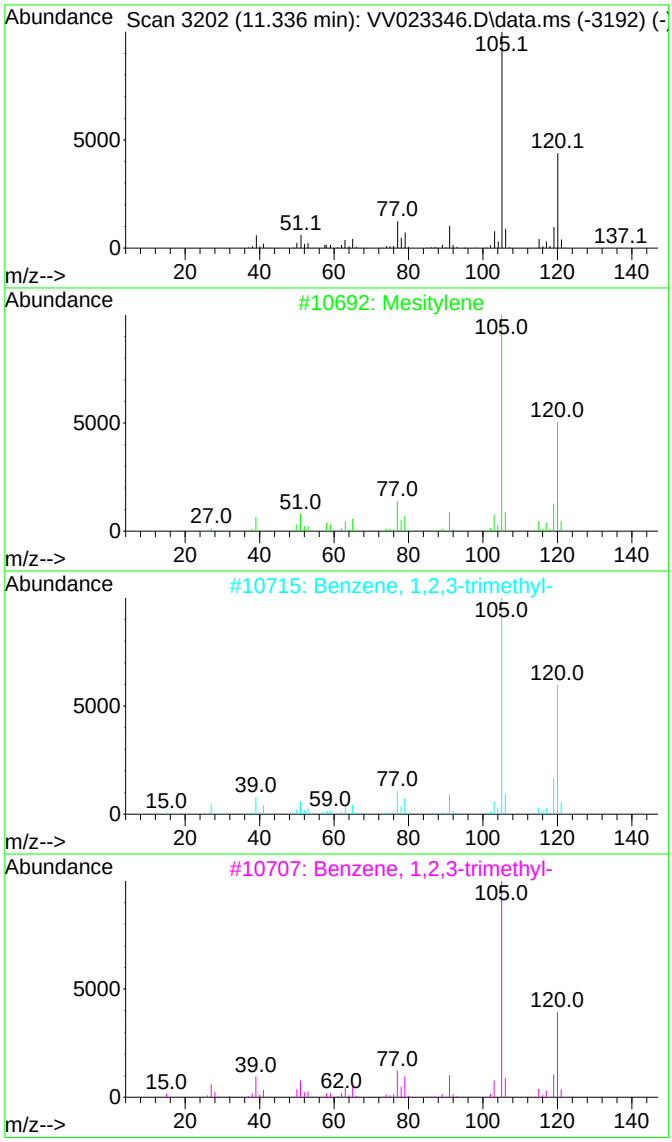
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

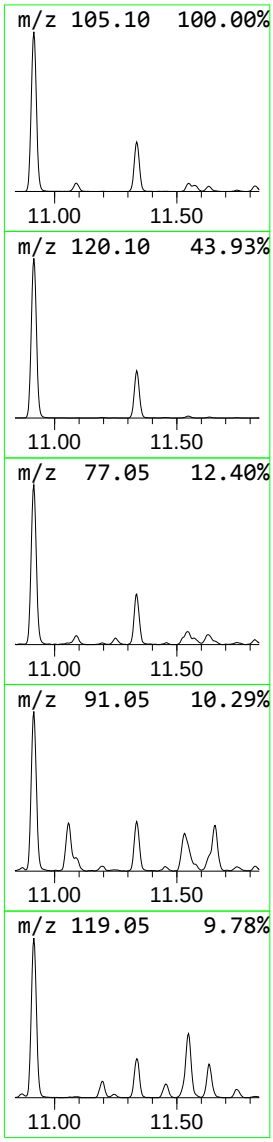
****

Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.336	5.47 ug/L	532017	1,4-Dichlorobenzene-d4	11.249

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	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4	Mesitylene	120	C9H12	000108-67-8	93
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE8

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

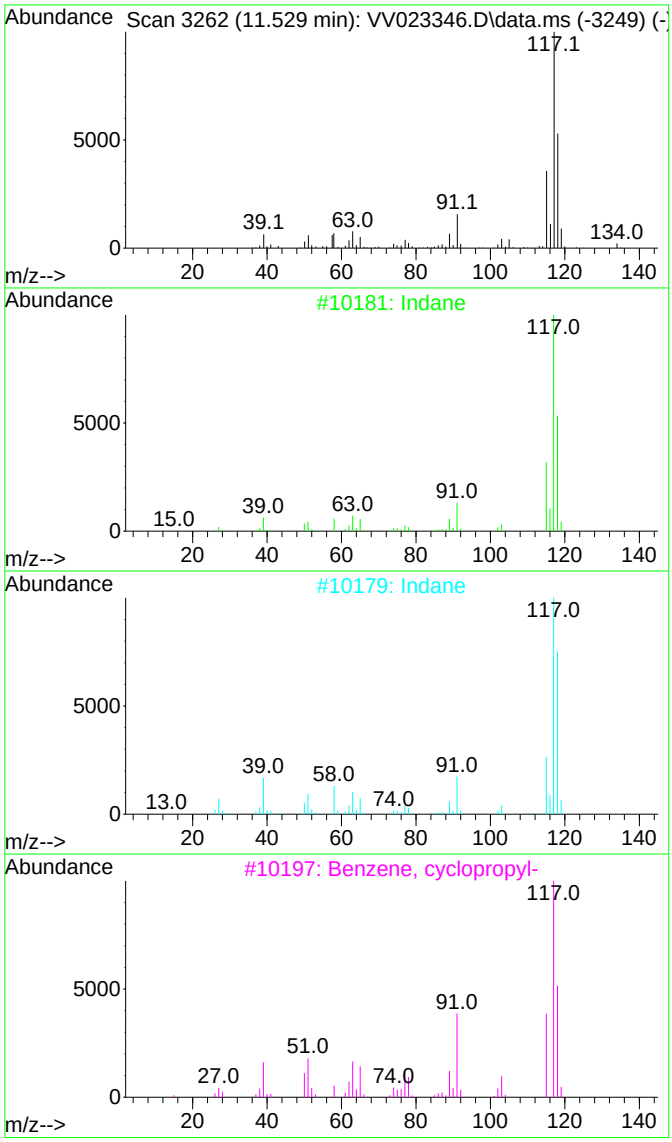
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

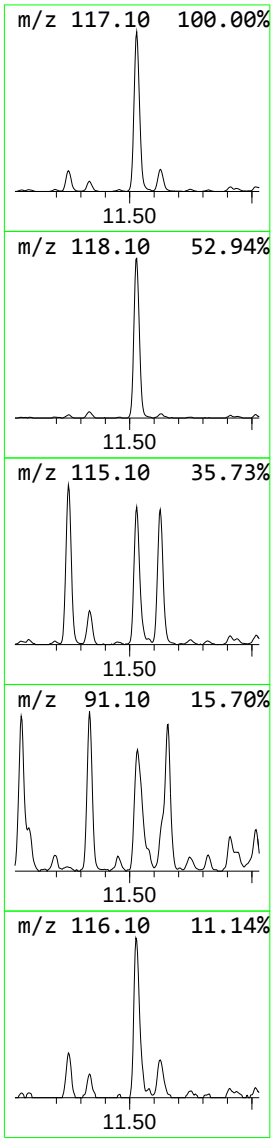
****

Peak Number 8 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	5.30 ug/L	515437	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	94
2	Indane	118	C9H10	000496-11-7	91
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
4	Benzene, 2-propenyl-	118	C9H10	000300-57-2	74
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	74





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE8

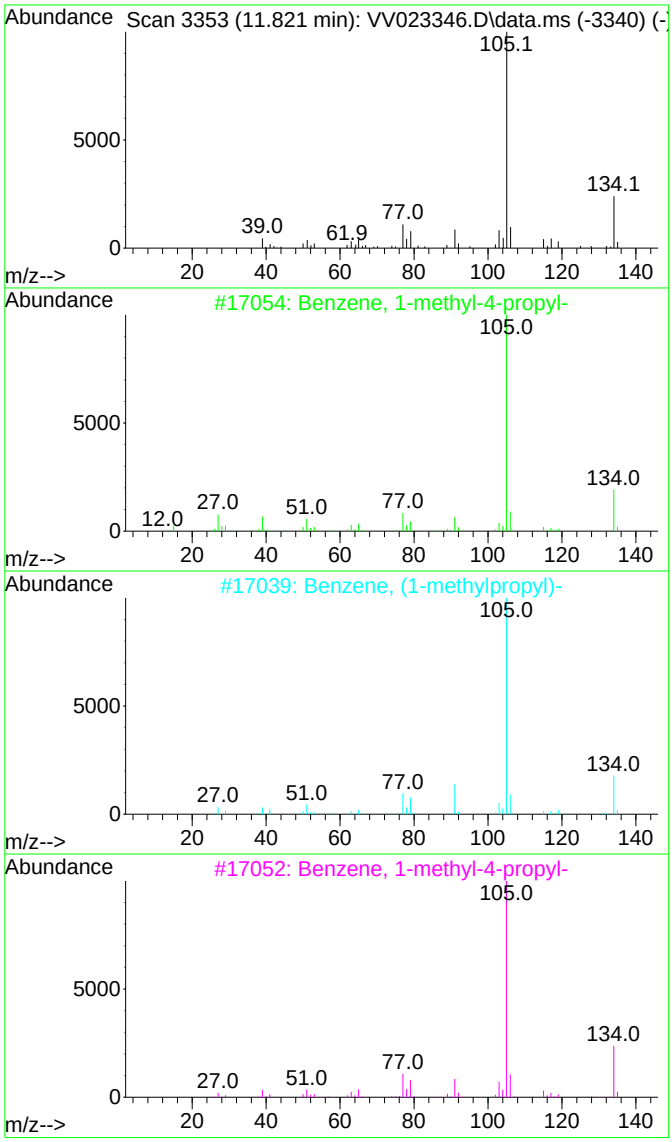
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

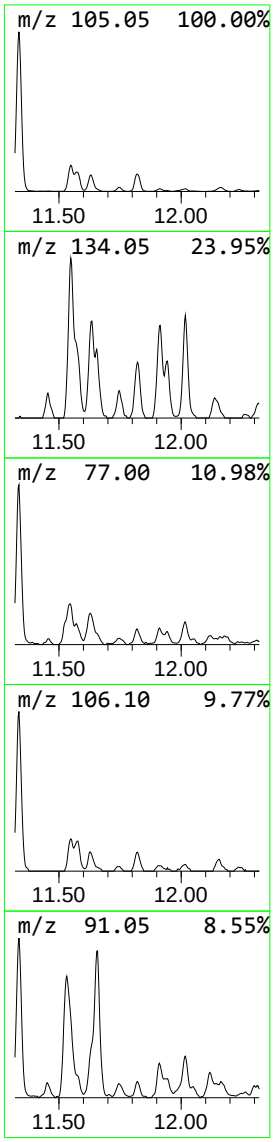
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-methyl-4-propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.821	0.55 ug/L	53279	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	90
2	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	90
3	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	90
4	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	87
5	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE8

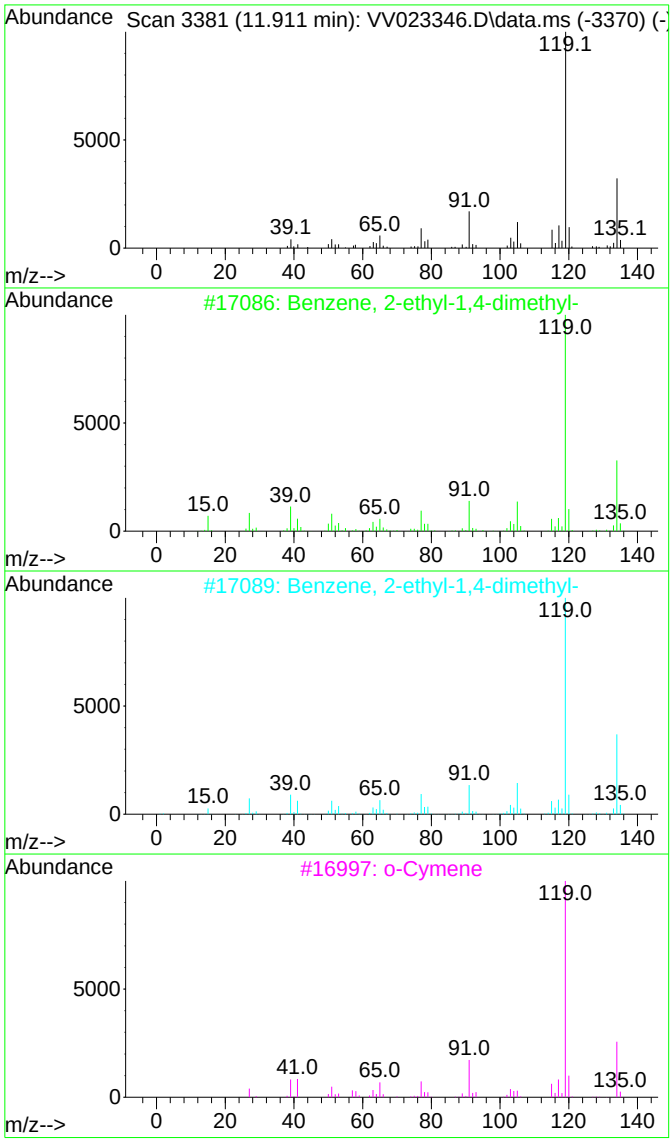
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

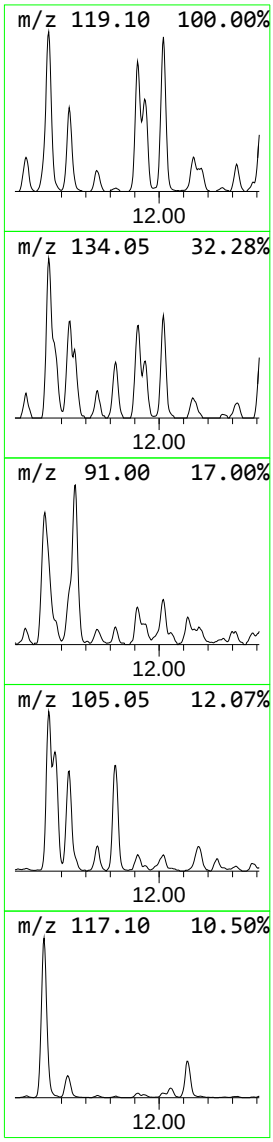
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.911	0.75 ug/L	72755	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
5	o-Cymene	134	C10H14	000527-84-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE8

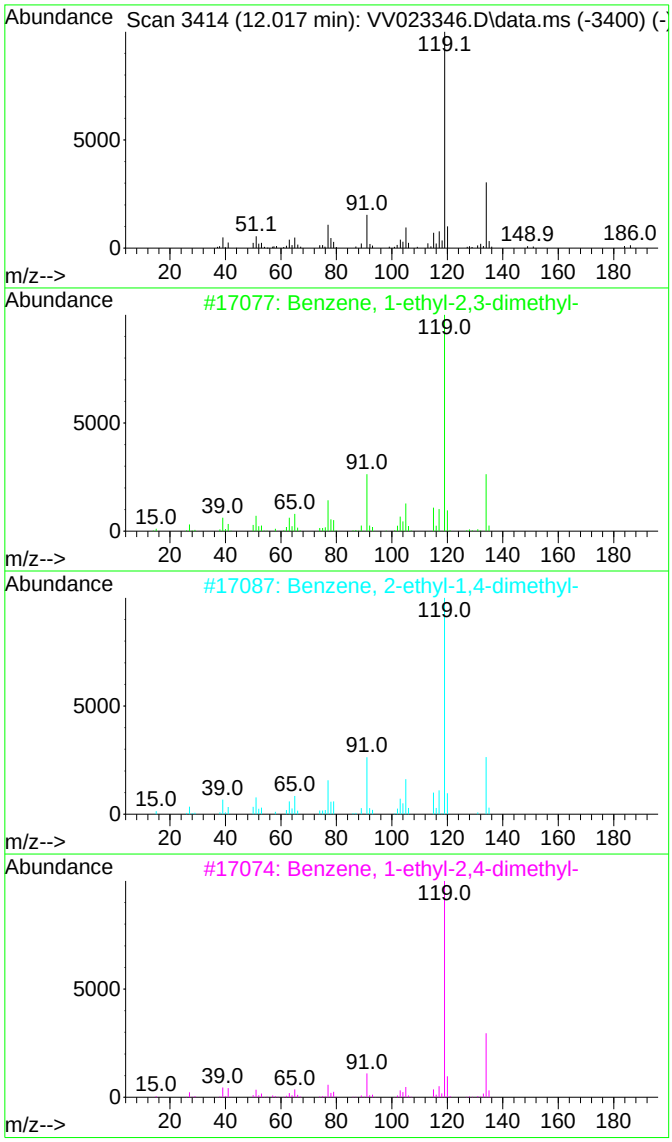
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

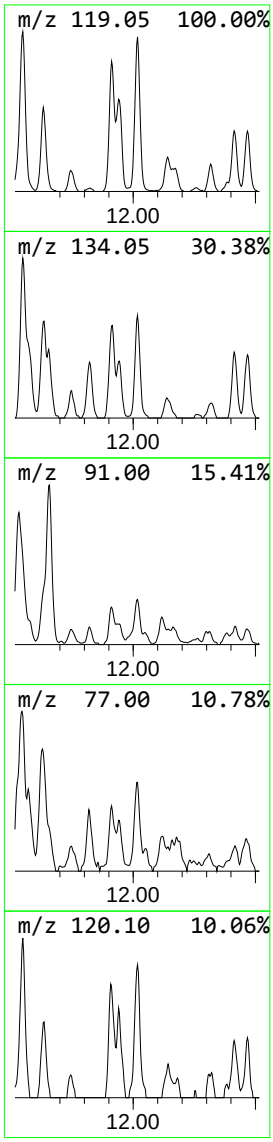
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.017	0.90 ug/L	87484	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	p-Cymene	134	C10H14	000099-87-6	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE8

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

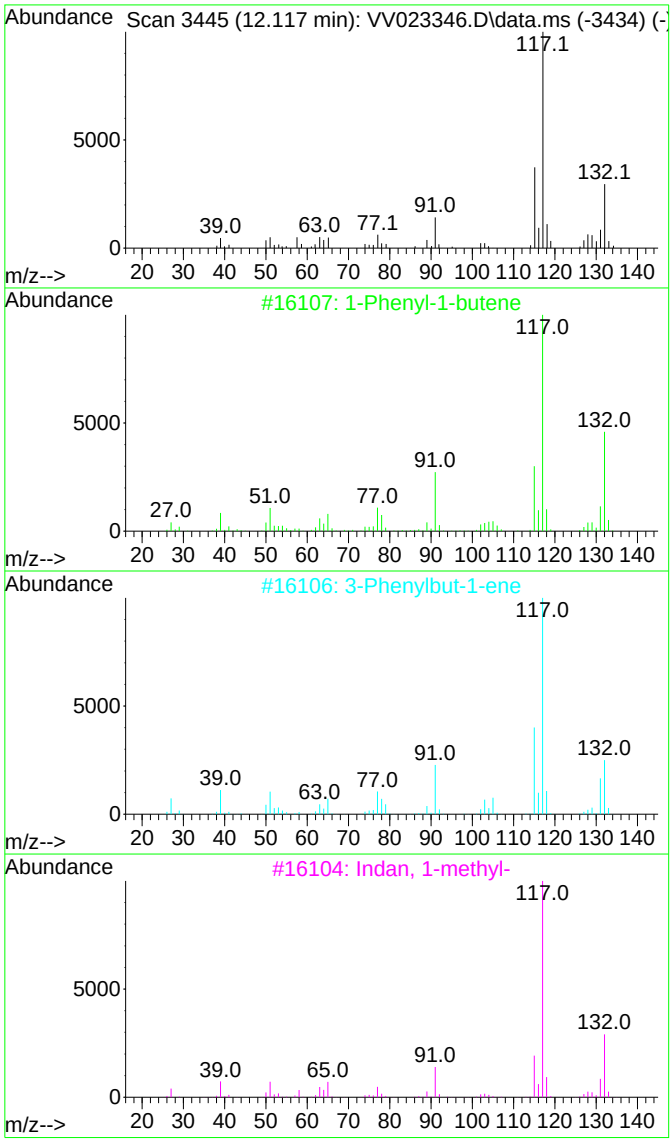
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

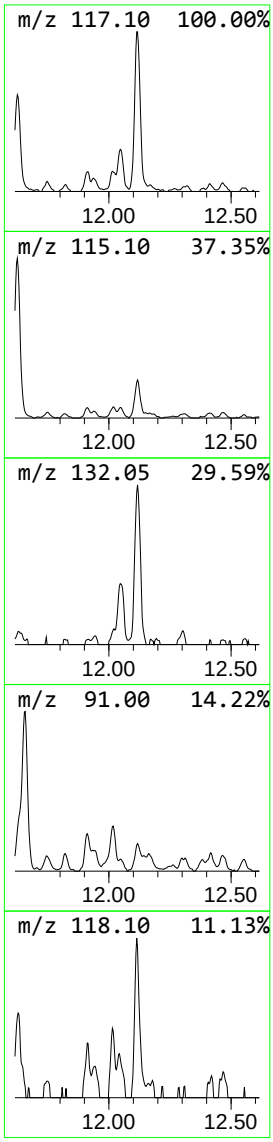
****

Peak Number 12 1-Phenyl-1-butene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.117	0.85 ug/L	82635	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
3	Indan, 1-methyl-	132	C10H12	000767-58-8	87
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
5	Indan, 1-methyl-	132	C10H12	000767-58-8	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE8

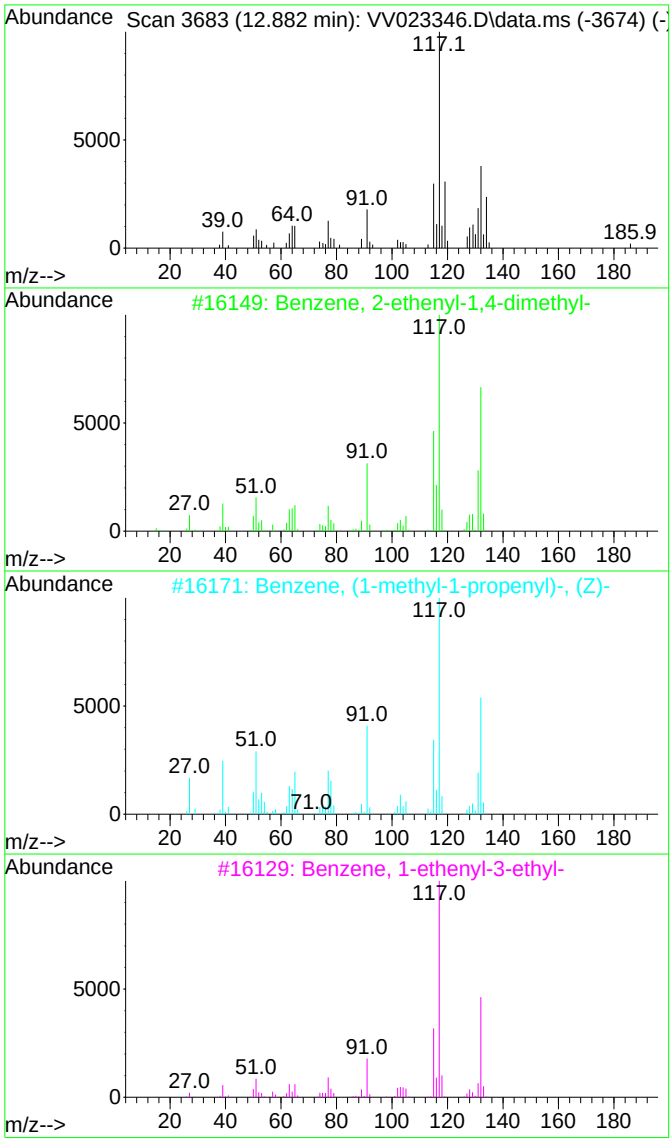
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

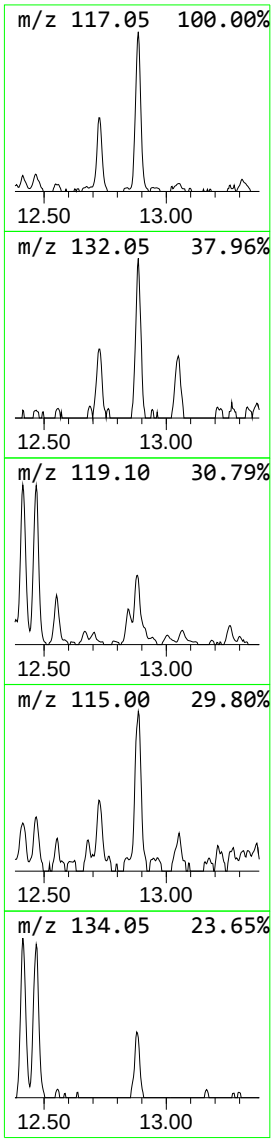
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.882	0.52 ug/L	50674	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	70
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	70
4	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	70
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	70





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023346.D
Acq On : 10 Nov 2021 20:47
Operator : SY/MD
Sample : M4542-02
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE8

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, propyl-	10.355	3.2	ug/L	315040	3	11.249	486285	5.0
Benzene, 1-ethy...	10.448	1.9	ug/L	188753	3	11.249	486285	5.0
Benzene, 1-ethy...	10.738	5.9	ug/L	573680	3	11.249	486285	5.0
Benzene, (2-met...	11.056	0.7	ug/L	70307	3	11.249	486285	5.0
Benzene, (1-met...	11.088	0.8	ug/L	78635	3	11.249	486285	5.0
Benzene, 1,2,3-...	11.336	5.5	ug/L	532017	3	11.249	486285	5.0
Indane	11.529	5.3	ug/L	515437	3	11.249	486285	5.0
Benzene, 1-meth...	11.821	0.6	ug/L	53279	3	11.249	486285	5.0
Benzene, 2-ethy...	11.911	0.8	ug/L	72755	3	11.249	486285	5.0
Benzene, 1-ethy...	12.017	0.9	ug/L	87484	3	11.249	486285	5.0
1-Phenyl-1-butene	12.117	0.8	ug/L	82635	3	11.249	486285	5.0
Benzene, 2-ethe...	12.882	0.5	ug/L	50674	3	11.249	486285	5.0