

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF120318\
 Data File : VF060898.D
 Acq On : 3 Dec 2018 18:24
 Operator : VA/AP
 Sample : J5775-03
 Misc : 7.23g/5mL/MSVOA-F/SOIL
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 TR-04-113018

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % 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_F\METHODS\82F112618S.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.135	46	51	54	rBV5	3204	12319	1.29%	0.193%
2	2.191	153	158	160	rBV4	4071	9630	1.01%	0.151%
3	2.250	160	164	171	rVB	18693	55510	5.81%	0.870%
4	3.088	247	249	256	rVB2	7157	12915	1.35%	0.202%
5	4.045	340	346	349	rBV2	9384	32908	3.45%	0.515%
6	4.134	349	355	367	rVB2	138172	425441	44.55%	6.664%
7	4.883	422	431	443	rBV2	273729	955009	100.00%	14.960%
8	5.623	496	506	515	rBV	299415	830710	86.98%	13.013%
9	6.195	562	564	569	rVB	16755	14646	1.53%	0.229%
10	6.264	569	571	573	rBV	10597	9974	1.04%	0.156%
11	7.576	698	704	709	rBV	339823	807869	84.59%	12.655%
12	7.645	709	711	717	rVB	21579	47550	4.98%	0.745%
13	8.652	808	813	820	rBV7	3322	12182	1.28%	0.191%
14	9.796	924	929	937	rBV	407943	882786	92.44%	13.829%
15	9.924	937	942	945	rVB5	5269	14406	1.51%	0.226%
16	10.151	962	965	968	rVB	13818	25063	2.62%	0.393%
17	10.733	1015	1024	1030	rBV	13325	34070	3.57%	0.534%
18	10.871	1034	1038	1043	rVB2	4821	10266	1.07%	0.161%
19	11.433	1090	1095	1102	rBV	310990	587379	61.51%	9.201%
20	11.739	1122	1126	1129	rVB3	6825	14390	1.51%	0.225%
21	11.838	1134	1136	1140	rBV3	6887	10225	1.07%	0.160%
22	12.035	1154	1156	1165	rVB	13146	20116	2.11%	0.315%
23	12.222	1172	1175	1178	rVB	14821	22039	2.31%	0.345%
24	12.568	1206	1210	1214	rVV	473835	750598	78.60%	11.758%
25	12.617	1214	1215	1219	rVV	24515	34802	3.64%	0.545%
26	12.735	1222	1227	1231	rBV	69325	110895	11.61%	1.737%
27	12.824	1234	1236	1242	rVB5	14142	32022	3.35%	0.502%
28	13.021	1252	1256	1259	rVB3	10470	16408	1.72%	0.257%
29	13.091	1261	1263	1268	rBV4	4698	11713	1.23%	0.183%
30	13.169	1268	1271	1276	rBV3	12673	22800	2.39%	0.357%
31	13.298	1282	1284	1287	rVB2	7397	9860	1.03%	0.154%
32	13.436	1296	1298	1302	rVB5	5224	9671	1.01%	0.151%
33	13.505	1302	1305	1308	rBV2	9421	15243	1.60%	0.239%
34	13.554	1308	1310	1312	rBV	8962	10591	1.11%	0.166%

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35	13.722	1324	1327	1331	rBV3	22092	30803	3.23%	0.483%
36	13.870	1339	1342	1346	rBV3	21079	51727	5.42%	0.810%
37	13.978	1350	1353	1359	rVB7	8954	18689	1.96%	0.293%
38	14.067	1359	1362	1363	rBV3	4904	9603	1.01%	0.150%
39	14.097	1363	1365	1368	rVB3	11678	15741	1.65%	0.247%
40	14.166	1368	1372	1376	rBV6	9704	24854	2.60%	0.389%
41	14.225	1376	1378	1381	rVV3	6124	12247	1.28%	0.192%
42	14.472	1399	1403	1409	rVB	59408	80675	8.45%	1.264%
43	14.610	1415	1417	1424	rVB7	6619	15848	1.66%	0.248%
44	14.758	1428	1432	1435	rVB3	9656	17520	1.83%	0.274%
45	15.024	1455	1459	1463	rBV5	5632	15222	1.59%	0.238%
46	15.132	1467	1470	1473	rBV3	7831	13725	1.44%	0.215%
47	15.300	1484	1487	1490	rVB	50179	70520	7.38%	1.105%
48	15.428	1497	1500	1505	rVB	65010	111253	11.65%	1.743%
49	15.803	1534	1538	1540	rBV4	6375	12292	1.29%	0.193%
50	16.020	1557	1560	1563	rVB2	6119	11023	1.15%	0.173%

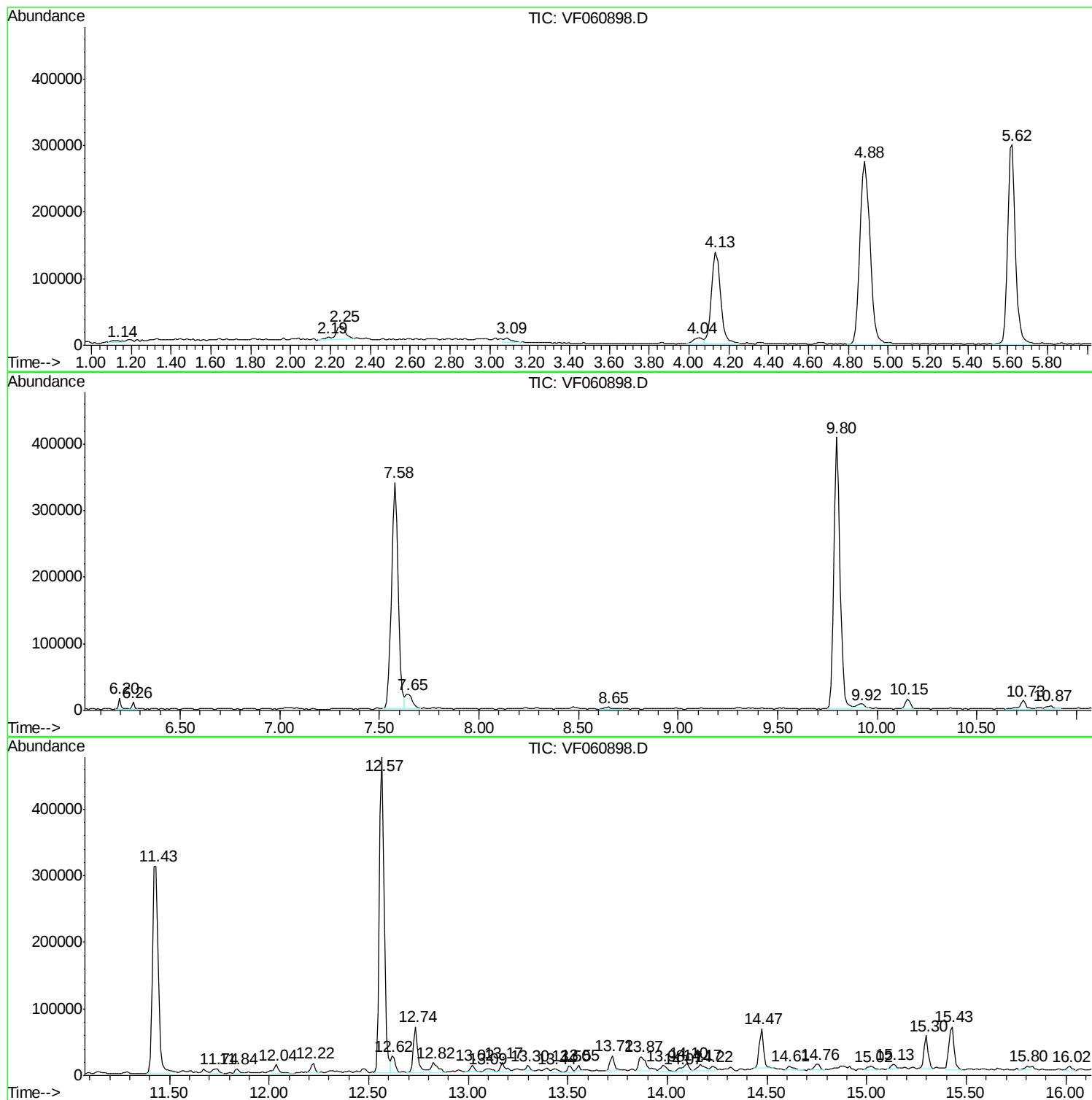
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F112618S.M
Quant Title : SW846 8260

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



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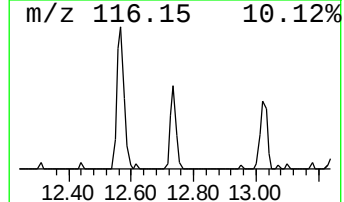
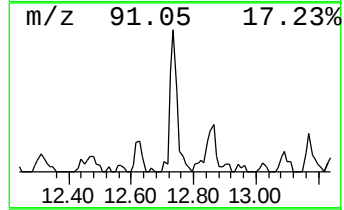
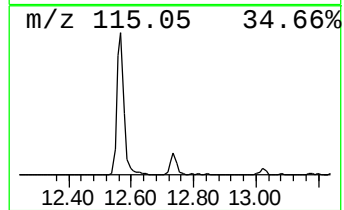
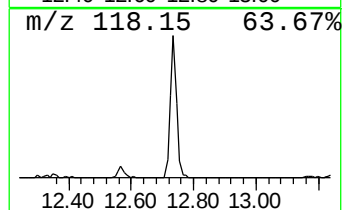
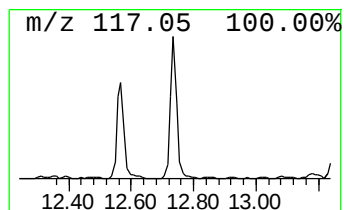
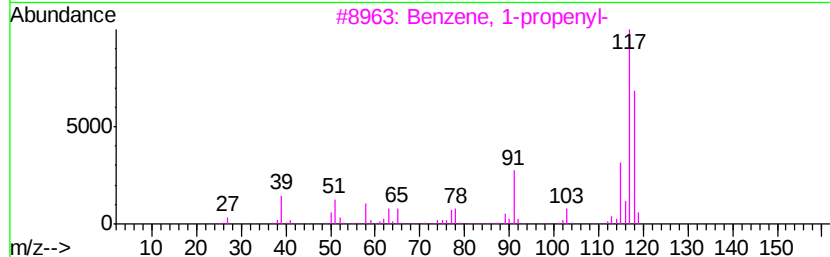
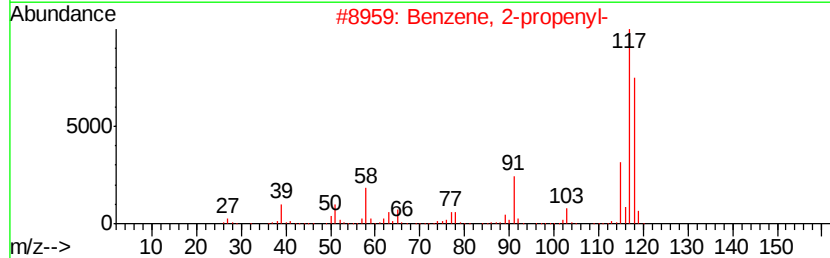
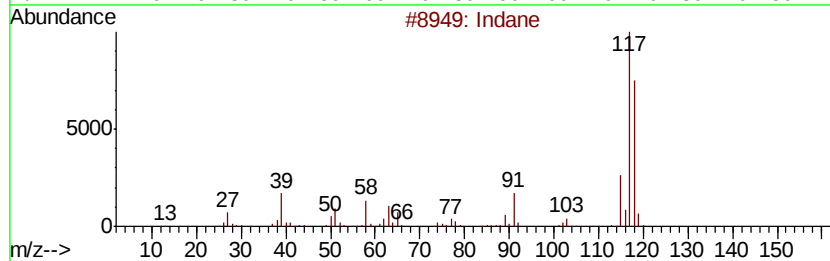
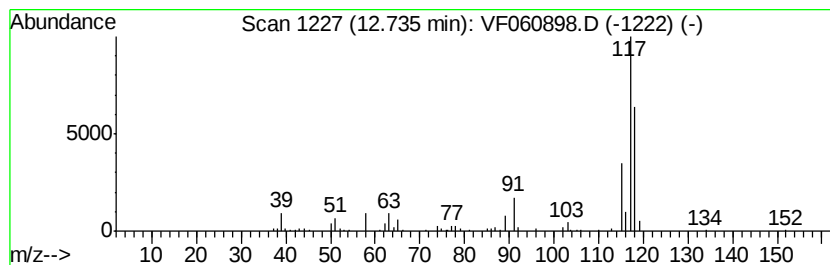
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Quant Title : SW846 8260

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

Peak Number 1 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	7.39 ug/l	110895	1,4-Dichlorobenzene-d4	12.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	95
2		Benzene, 2-propenyl-	118	C9H10	000300-57-2	81
3		Benzene, 1-propenyl-	118	C9H10	000637-50-3	72
4		Benzene, cyclopropyl-	118	C9H10	000873-49-4	72
5		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	68



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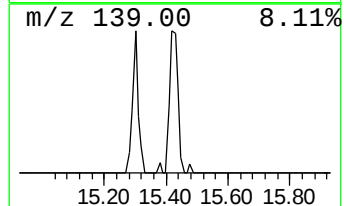
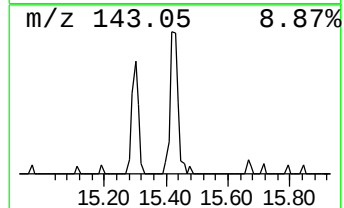
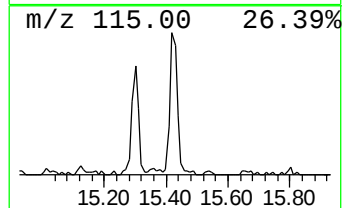
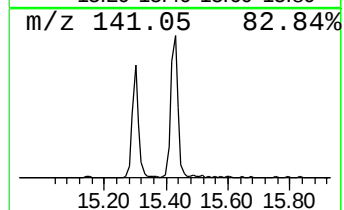
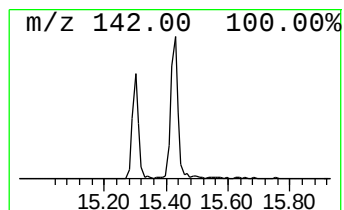
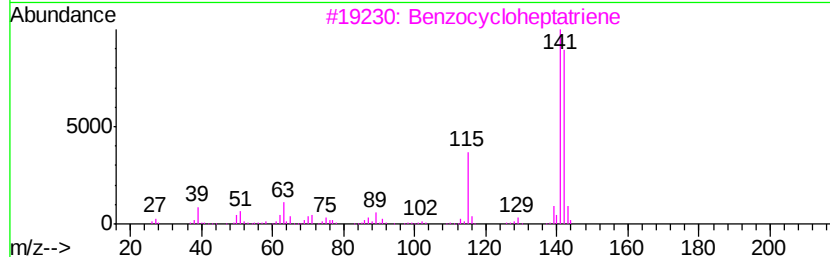
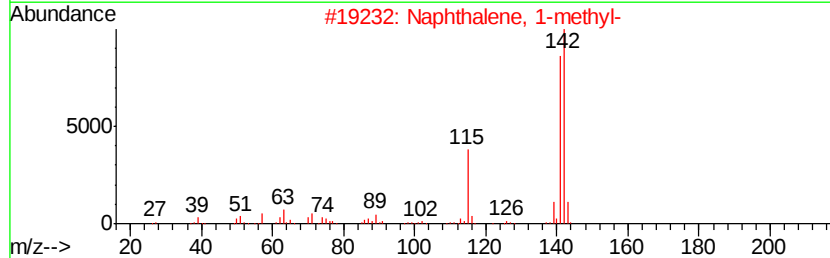
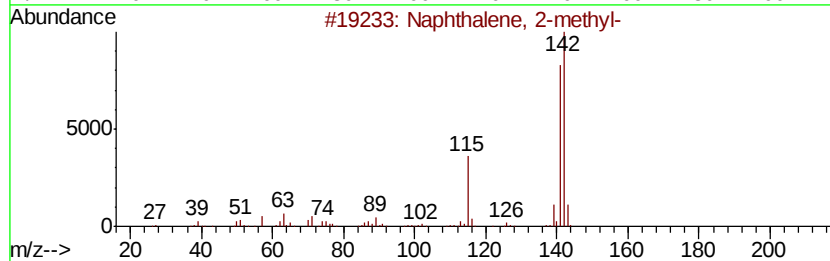
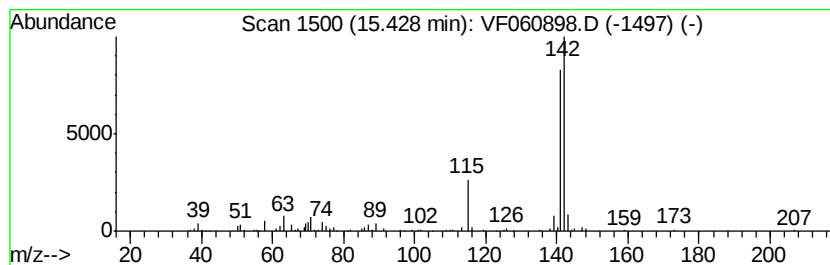
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Peak Number 2 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.43	7.41 ug/l	111253	1,4-Dichlorobenzene-d4	12.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
3		Benzocycloheptatriene	142	C11H10	000264-09-5	91
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	90
5		Naphthalene, 1-[(2-propenyloxy)m...	198	C14H14O	074685-39-5	64



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Indane	12.74	7.4	ug/l	110895	4	12.57	750598	50.0
Naphthalene, 1-me...	15.43	7.4	ug/l	111253	4	12.57	750598	50.0