

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
 Data File : BM033945.D
 Acq On : 01 Feb 2022 18:32
 Operator : CG/JU
 Sample : N1328-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COHM2

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
 Title : SVOA CALIBRATION

Signal : TIC: BM033945.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.325	12	15	35	rVB	1709594	2354597	6.79%	3.725%
2	4.160	152	157	166	rBV2	60683	92116	0.27%	0.146%
3	4.613	225	234	247	rBV	21963746	34663035	100.00%	54.842%
4	6.660	577	582	588	rBV	23673	37451	0.11%	0.059%
5	7.084	648	654	664	rBV	112054	186343	0.54%	0.295%
6	7.248	675	682	694	rBV	407822	662150	1.91%	1.048%
7	7.454	710	717	728	rBV	424327	660735	1.91%	1.045%
8	7.931	791	798	805	rBV	401318	652811	1.88%	1.033%
9	8.637	908	918	930	rBV	293900	478037	1.38%	0.756%
10	9.089	987	995	1006	rBV	531319	860919	2.48%	1.362%
11	9.542	1064	1072	1079	rVB	38150	61518	0.18%	0.097%
12	9.813	1108	1118	1131	rBV	412048	705044	2.03%	1.115%
13	10.354	1203	1210	1221	rBV	548960	931178	2.69%	1.473%
14	10.742	1268	1276	1284	rVB	545022	927782	2.68%	1.468%
15	10.872	1291	1298	1307	rBV	377127	637418	1.84%	1.008%
16	13.472	1735	1740	1746	rBV	52324	78864	0.23%	0.125%
17	13.977	1819	1826	1831	rBV	1106279	1597060	4.61%	2.527%
18	14.260	1868	1874	1888	rVB	1251237	1817407	5.24%	2.875%
19	14.566	1919	1926	1936	rBV	811743	1188760	3.43%	1.881%
20	14.760	1954	1959	1971	rBV2	41670	103133	0.30%	0.163%
21	15.554	2087	2094	2105	rBV	1602422	2272774	6.56%	3.596%
22	15.666	2107	2113	2127	rVB	526378	704594	2.03%	1.115%
23	17.301	2384	2391	2400	rBV	951930	1306379	3.77%	2.067%
24	17.401	2401	2408	2420	rVV	1755449	2445279	7.05%	3.869%
25	17.977	2500	2506	2512	rBV	57814	72957	0.21%	0.115%
26	18.195	2538	2543	2551	rBV2	45306	72845	0.21%	0.115%
27	19.142	2700	2704	2709	rBV2	150023	170502	0.49%	0.270%
28	19.277	2724	2727	2731	rVV	41149	48607	0.14%	0.077%
29	19.465	2756	2759	2765	rBV5	27743	42423	0.12%	0.067%
30	19.683	2790	2796	2803	rBV	1947073	2574660	7.43%	4.073%
31	21.465	3094	3099	3104	rBV	958302	1194230	3.45%	1.889%
32	23.700	3472	3479	3486	rBV2	1185683	2394067	6.91%	3.788%
33	23.859	3499	3506	3515	rVB2	600281	1209445	3.49%	1.914%

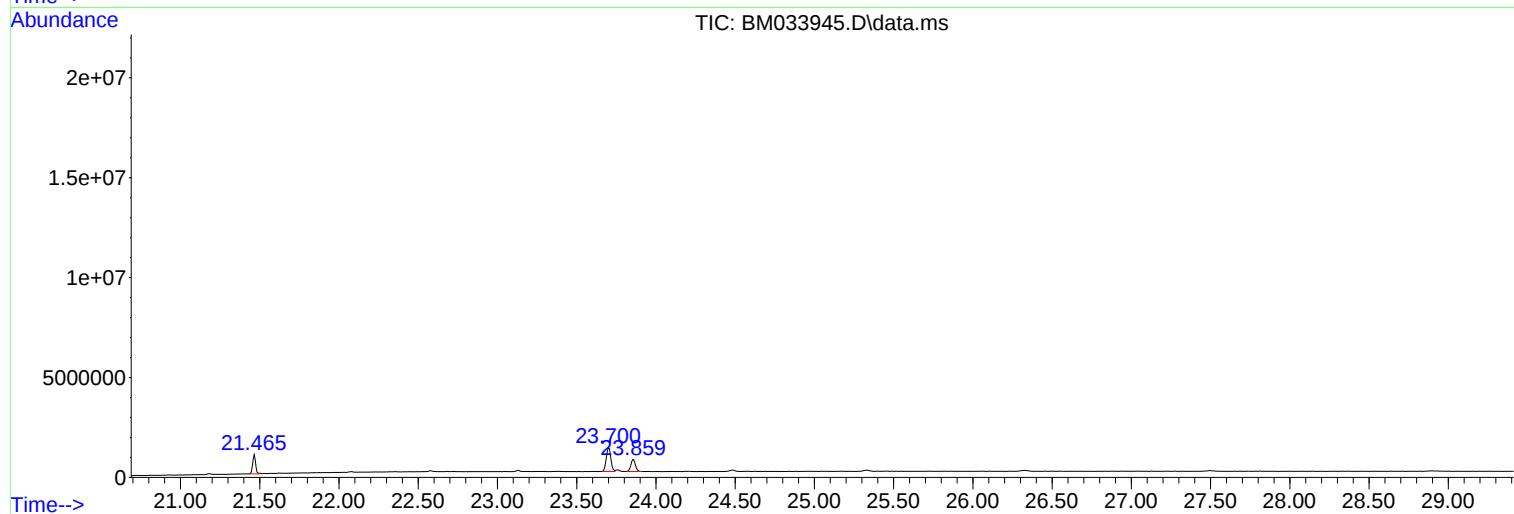
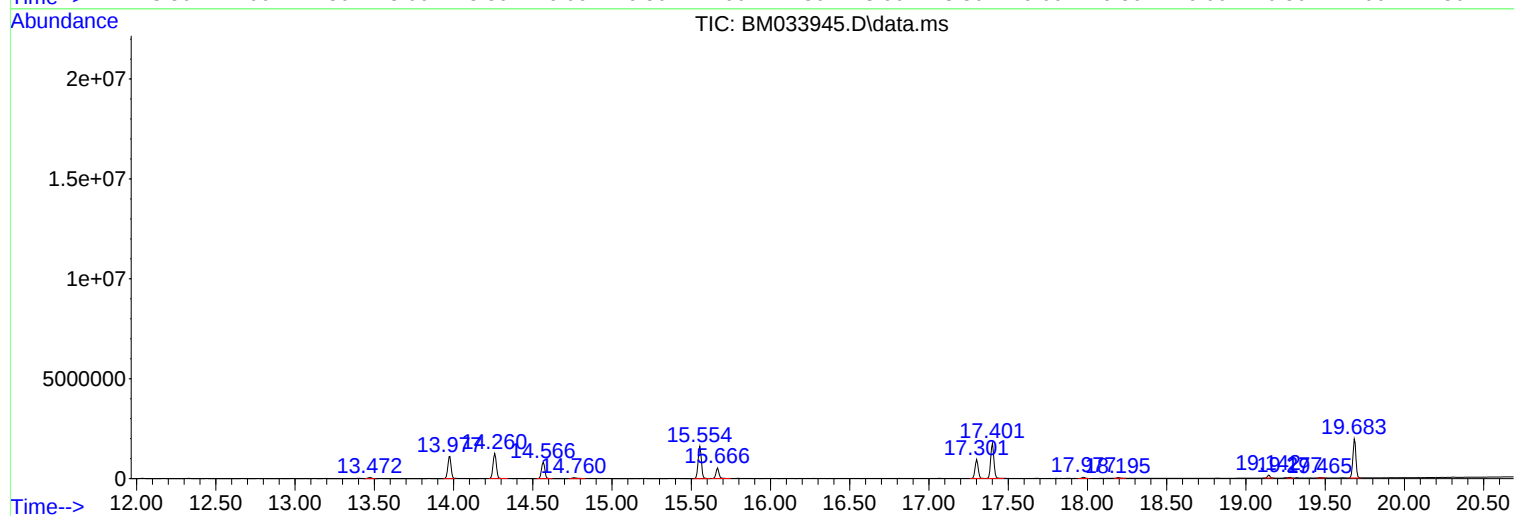
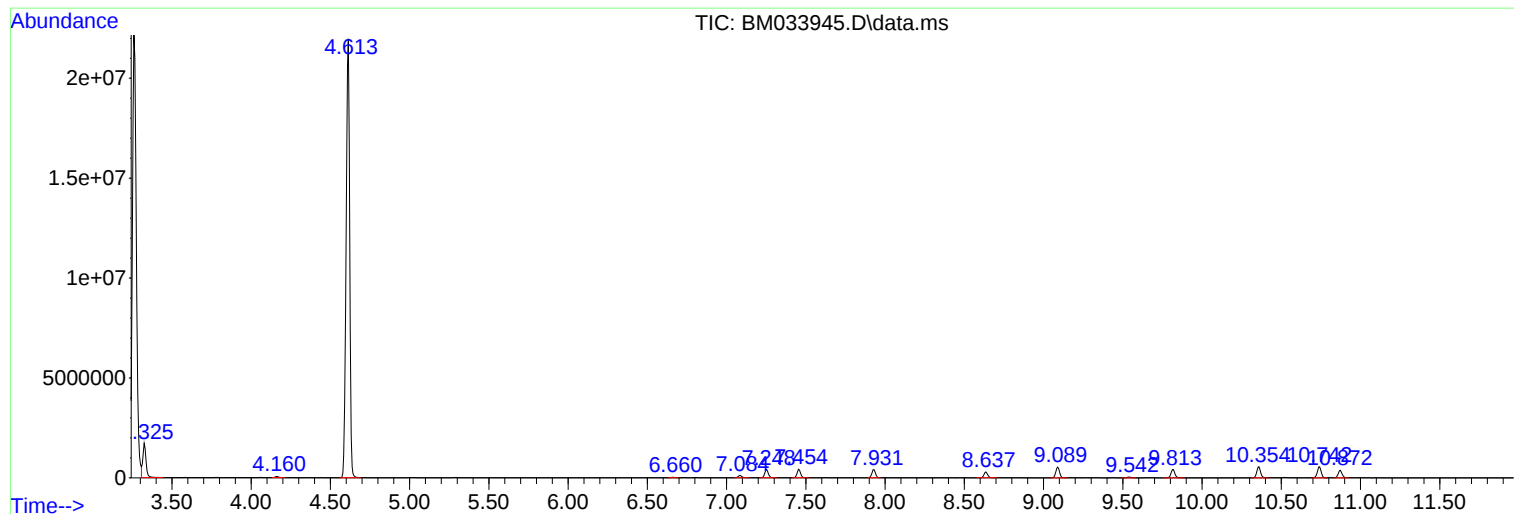
Sum of corrected areas: 63205120

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C0HM2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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



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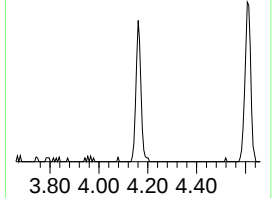
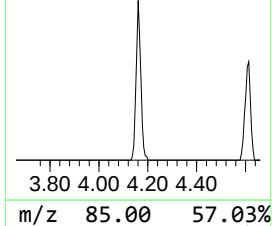
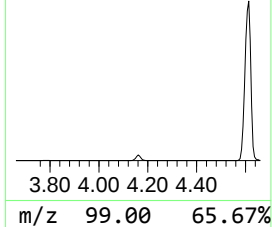
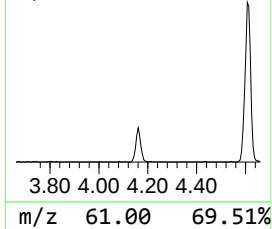
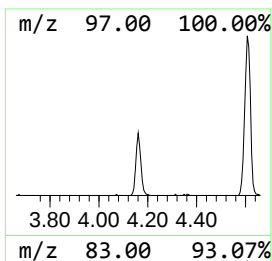
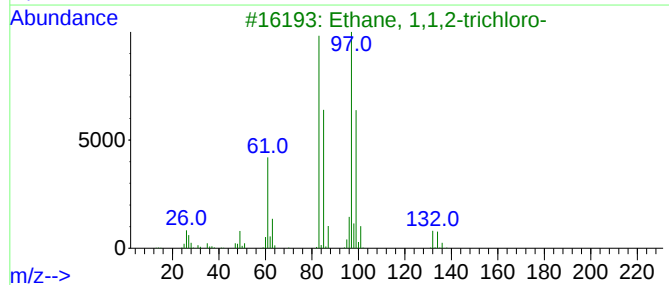
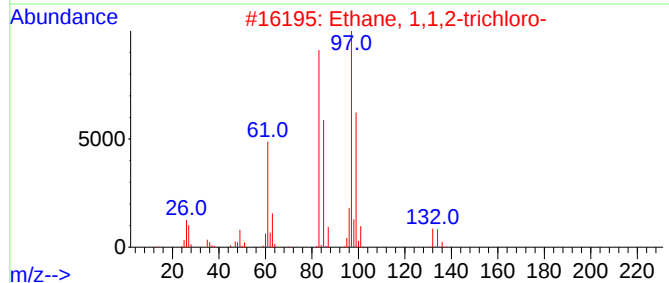
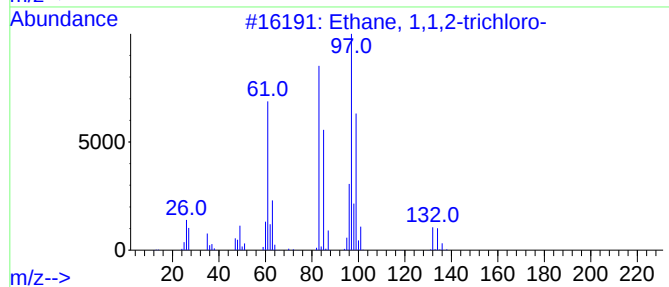
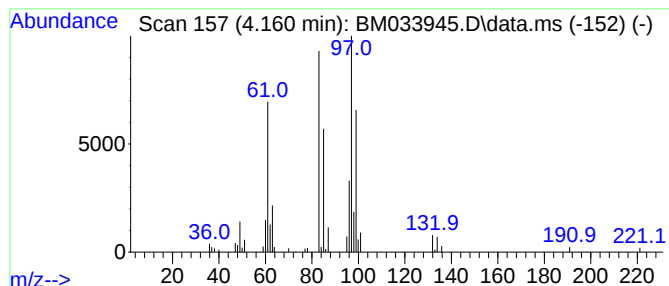
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethane, 1,1,2-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.160	2.82 ng/ul	92116	1,4-Dichlorobenzene-d4	7.931

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethane, 1,1,2-trichloro-	132	C2H3Cl3	000079-00-5	98
2			Ethane, 1,1,2-trichloro-	132	C2H3Cl3	000079-00-5	94
3			Ethane, 1,1,2-trichloro-	132	C2H3Cl3	000079-00-5	94
4			Ethane, 1,1,2-trichloro-	132	C2H3Cl3	000079-00-5	81
5			Ethane, 1,1,2-trichloro-	132	C2H3Cl3	000079-00-5	64



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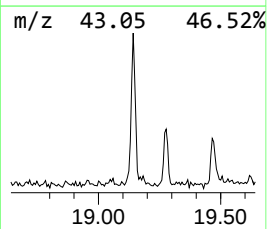
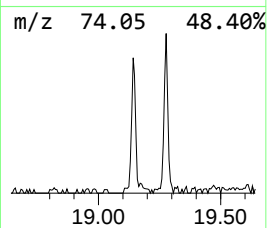
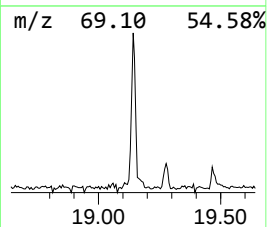
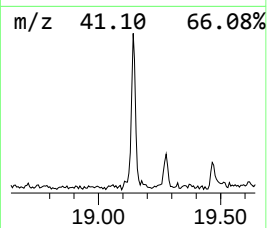
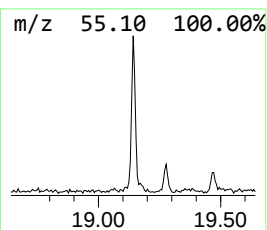
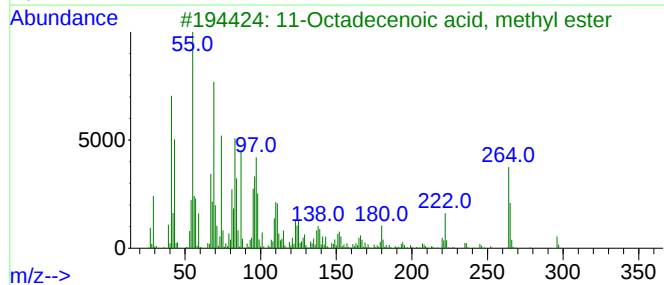
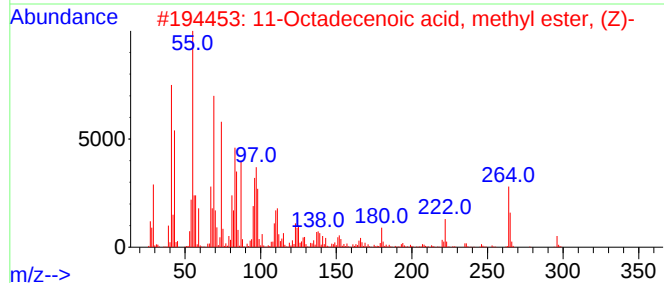
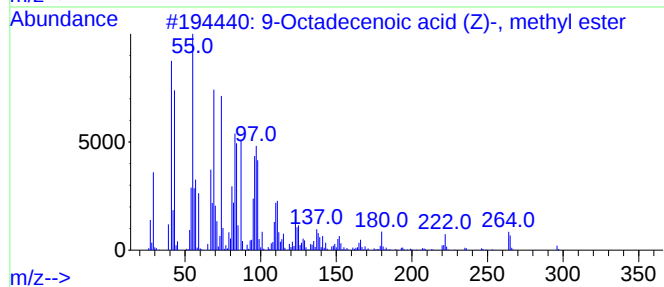
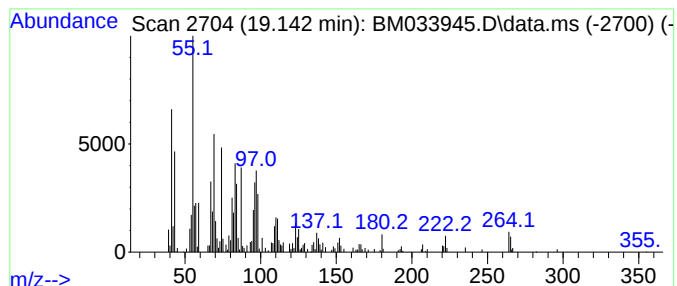
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 9-Octadecenoic acid (Z)-, m... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.142	2.61 ng/ul	170502	Phenanthrene-d10	17.301

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2			11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	99
3			11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	99
4			cis-13-Octadecenoic acid, methyl...	296	C19H36O2	1010333-58-3	99
5			trans-13-Octadecenoic acid, meth...	296	C19H36O2	1000333-61-3	99



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TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethane, 1,1,2-t...	4.160	2.8	ng/ul	92116	1	7.931	652811	20.0
9-Octadecenoic ...	19.142	2.6	ng/ul	170502	4	17.301	1306380	20.0