

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080122\
 Data File : BP011198.D
 Acq On : 01 Aug 2022 14:11
 Operator : CG/JU
 Sample : N3790-16
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBUL1

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_P\Methods\SFAM-EPA-BP072822.M
 Title : SVOA CALIBRATION

Signal : TIC: BP011198.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.564	78	81	95	rVB	234926	334099	4.24%	0.510%
2	6.828	630	636	648	rVB	1280343	1969168	24.97%	3.004%
3	6.993	658	664	674	rVB	1085822	1652999	20.96%	2.521%
4	7.181	690	696	705	rVB	1298188	1970816	24.99%	3.006%
5	7.646	769	775	784	rVB	1334538	2019382	25.61%	3.080%
6	8.363	891	897	913	rBV	1496806	2389980	30.31%	3.645%
7	8.810	966	973	985	rVB	1283859	2050533	26.01%	3.128%
8	9.528	1088	1095	1104	rBV	939408	1512966	19.19%	2.308%
9	10.063	1179	1186	1197	rBV	1309581	2181919	27.67%	3.328%
10	10.440	1243	1250	1259	rBV	1807226	2933537	37.20%	4.474%
11	10.587	1268	1275	1288	rBV	1328212	2211536	28.05%	3.373%
12	12.040	1516	1522	1530	rVB3	24779	50913	0.65%	0.078%
13	12.316	1565	1569	1576	rBV2	7007	11775	0.15%	0.018%
14	12.663	1623	1628	1637	rBV10	5481	15518	0.20%	0.024%
15	12.922	1666	1672	1676	rVB6	10284	17683	0.22%	0.027%
16	13.216	1718	1722	1728	rVV3	16014	27870	0.35%	0.043%
17	13.275	1729	1732	1739	rVB6	6166	10778	0.14%	0.016%
18	13.534	1771	1776	1782	rBV10	6199	10077	0.13%	0.015%
19	13.734	1803	1810	1823	rBV	2125942	2963824	37.59%	4.521%
20	14.004	1846	1856	1869	rBV	2515950	3595401	45.60%	5.484%
21	14.316	1902	1909	1916	rBV	2289043	3175918	40.28%	4.844%
22	14.381	1916	1920	1923	rVB6	4906	8408	0.11%	0.013%
23	14.539	1941	1947	1961	rBV	572120	972990	12.34%	1.484%
24	14.751	1979	1983	1990	rVB4	31816	47241	0.60%	0.072%
25	15.134	2044	2048	2050	rBV3	8259	11474	0.15%	0.018%
26	15.316	2073	2079	2095	rVB	2887721	4024580	51.04%	6.139%
27	15.445	2095	2101	2113	rBV	583403	817848	10.37%	1.247%
28	15.557	2115	2120	2124	rVB2	31771	45746	0.58%	0.070%
29	15.886	2172	2176	2181	rBV7	5785	10905	0.14%	0.017%
30	16.016	2194	2198	2202	rBV7	5883	12761	0.16%	0.019%
31	16.057	2202	2205	2208	rVV5	6675	10830	0.14%	0.017%
32	16.110	2211	2214	2221	rVB7	9734	16114	0.20%	0.025%
33	16.757	2321	2324	2328	rVB6	5426	8432	0.11%	0.013%
34	16.857	2335	2341	2346	rBV5	9123	16029	0.20%	0.024%
35	17.022	2365	2369	2373	rVB7	7110	11462	0.15%	0.017%
36	17.086	2373	2380	2390	rBV	2216383	3099413	39.31%	4.727%

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Peak Location: TOP

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072822.M
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37	17.186	2390	2397	2409	rVV	2739638	3871512	49.10%	5.905%
38	17.292	2411	2415	2420	rVB3	36018	49194	0.62%	0.075%
39	17.386	2429	2431	2436	rBV6	6831	9502	0.12%	0.014%
40	17.498	2441	2450	2458	rVV6	18181	43984	0.56%	0.067%
41	17.781	2493	2498	2507	rBV	84122	102674	1.30%	0.157%
42	17.998	2531	2535	2539	rBV6	14499	21861	0.28%	0.033%
43	18.069	2543	2547	2551	rBV2	10907	17934	0.23%	0.027%
44	18.922	2688	2692	2696	rBV	199528	213960	2.71%	0.326%
45	19.016	2704	2708	2711	rBV3	39130	58055	0.74%	0.089%
46	19.057	2711	2715	2717	rVV	72224	79664	1.01%	0.122%
47	19.086	2717	2720	2723	rVB	27231	37420	0.47%	0.057%
48	19.445	2775	2781	2789	rBV	3324319	4443364	56.35%	6.777%
49	20.563	2961	2971	2974	rBV4	481292	1363506	17.29%	2.080%
50	20.939	3032	3035	3042	rBV3	148260	252820	3.21%	0.386%
51	21.210	3076	3081	3087	rBV	2695210	3319175	42.10%	5.063%
52	23.404	3447	3454	3472	rVB2	3152768	7884869	100.00%	12.027%
53	23.557	3474	3480	3489	rVB2	1806618	3571746	45.30%	5.448%

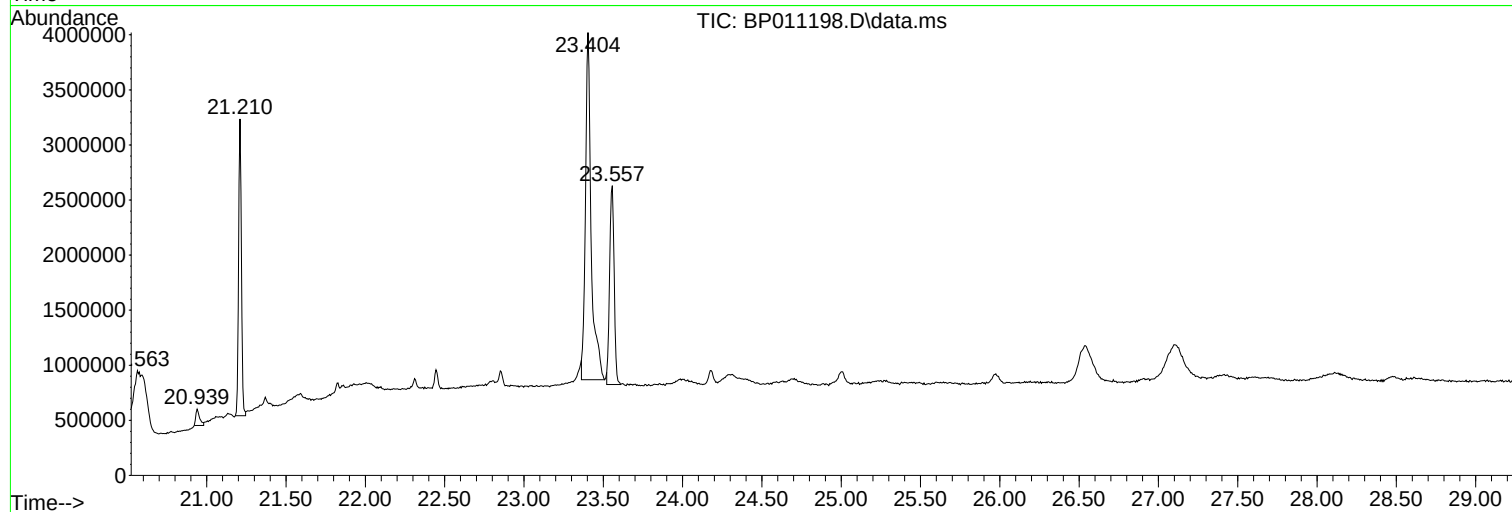
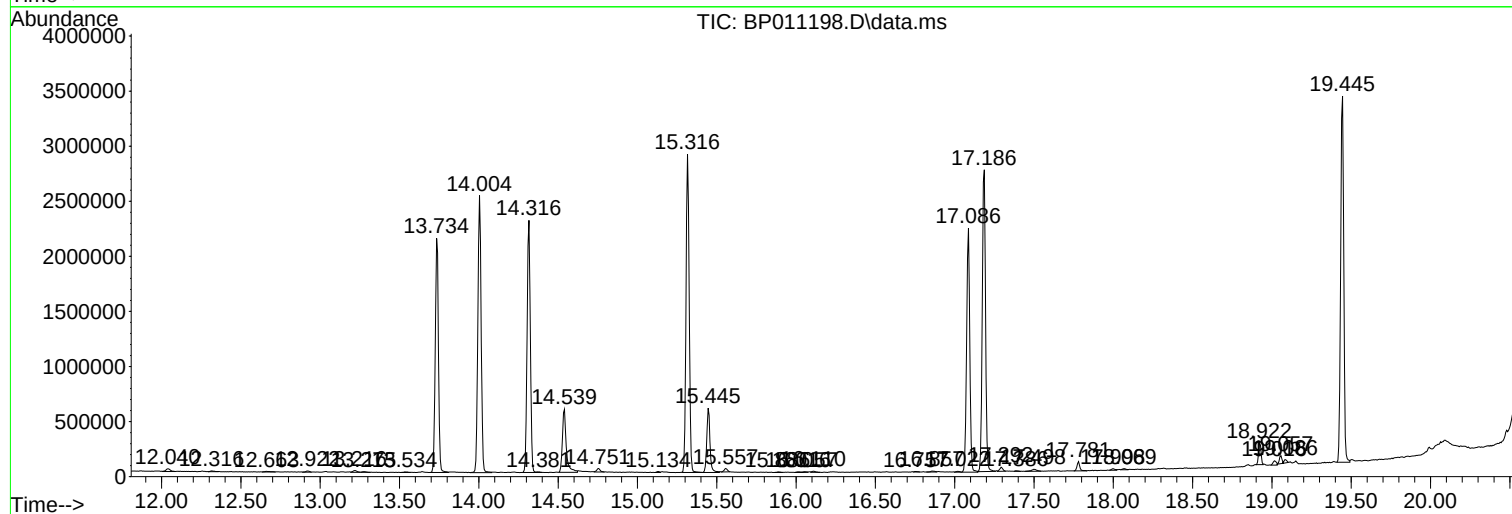
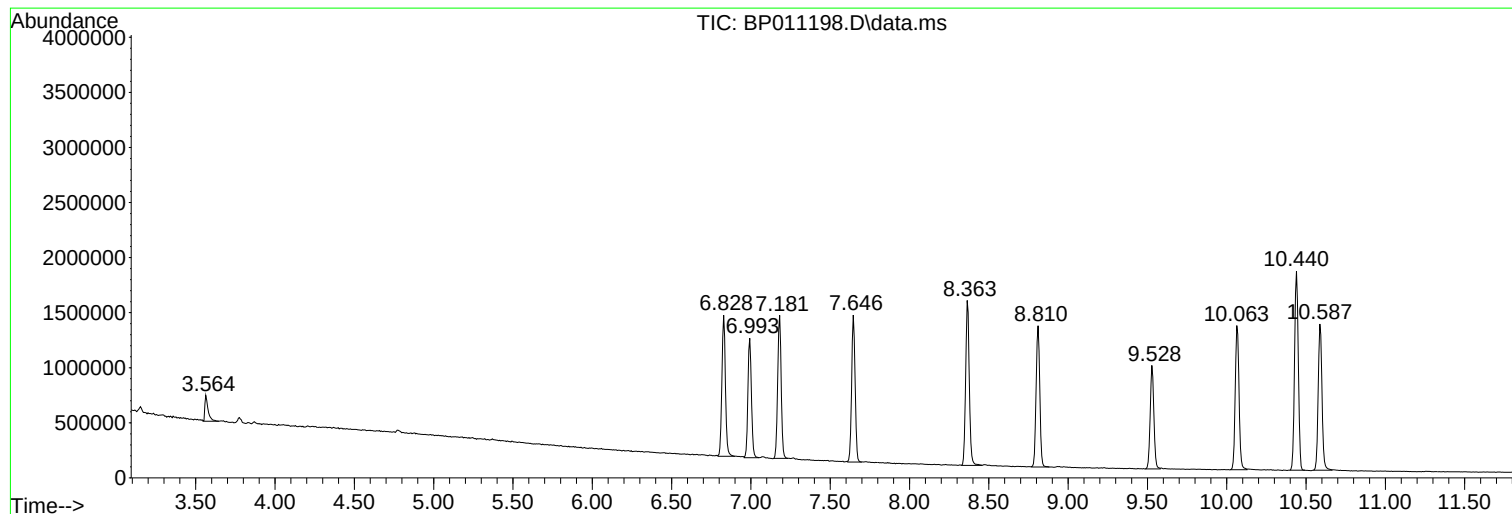
Sum of corrected areas: 65562165

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072822.M
Quant Title : SVOA CALIBRATION

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



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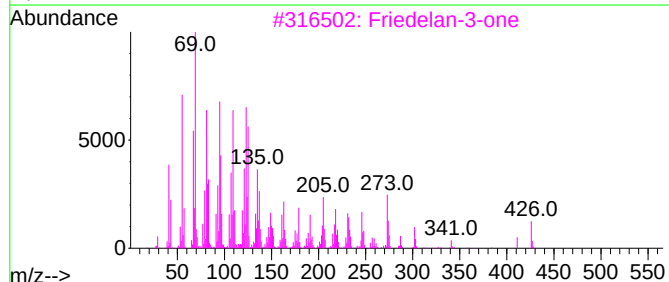
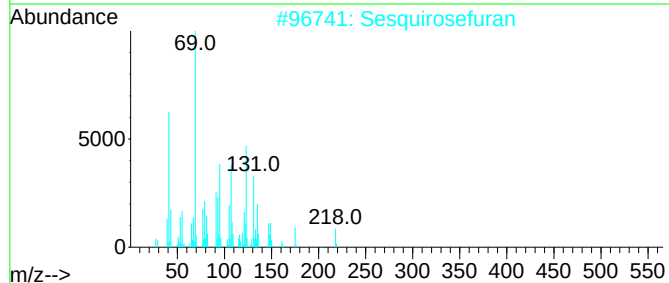
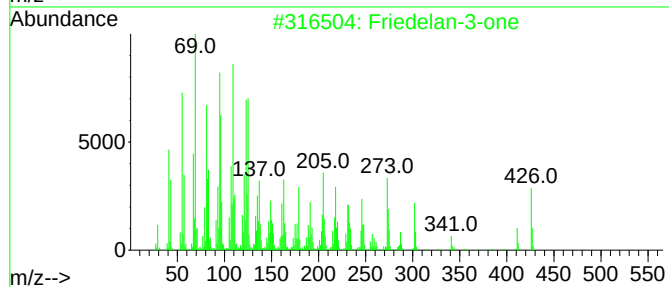
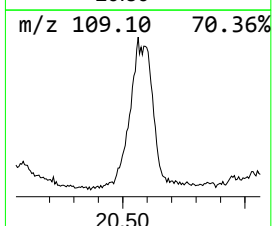
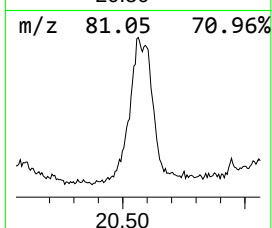
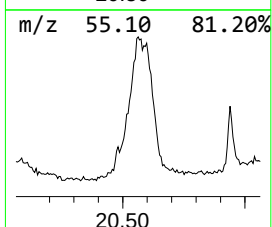
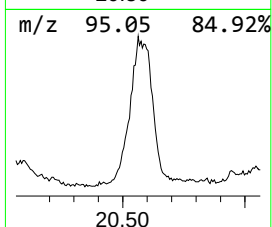
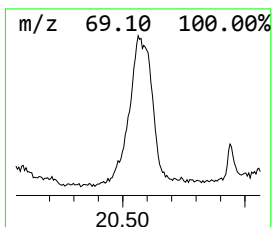
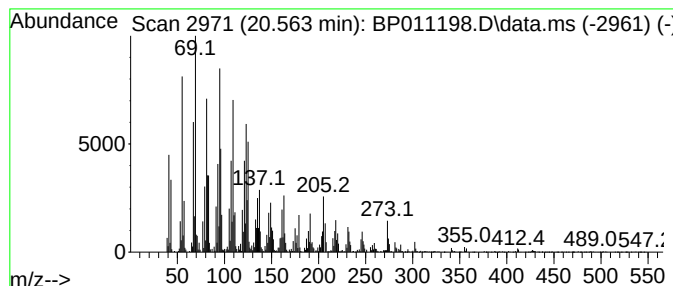
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072822.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Sesquirosefuran Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.563	8.22 ng/ul	1363510	Chrysene-d12	21.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Friedelan-3-one	426	C30H50O	000559-74-0	96
2			Sesquirosefuran	218	C15H22O	039007-93-7	86
3			Friedelan-3-one	426	C30H50O	000559-74-0	74
4			5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	53
5			Sesquirosefuran	218	C15H22O	039007-93-7	50



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Sesquirosefuran	20.563	8.2	ng/ul	1363510	5	21.210	3319180	20.0