

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022824\
 Data File : BM044406.D
 Acq On : 28 Feb 2024 16:17
 Operator : MA/JU
 Sample : PB159332BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK332

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-BM022324.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM044406.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.146	6	10	24	rVB	498022	598402	13.70%	1.214%
2	3.369	44	48	60	rVB	78728	112884	2.58%	0.229%
3	3.781	115	118	136	rBV	882304	1394345	31.92%	2.828%
4	4.105	169	173	183	rVB4	34265	61863	1.42%	0.125%
5	4.281	201	203	208	rVB6	5378	7210	0.17%	0.015%
6	4.334	209	212	217	rVB	22044	27970	0.64%	0.057%
7	4.487	236	238	243	rVB4	11657	13414	0.31%	0.027%
8	4.669	265	269	276	rVB	46556	68665	1.57%	0.139%
9	4.728	276	279	281	rBV2	16410	23172	0.53%	0.047%
10	5.840	462	468	474	rBV10	6578	11986	0.27%	0.024%
11	5.893	474	477	483	rBV8	2737	5486	0.13%	0.011%
12	6.010	496	497	503	rVB6	3193	4781	0.11%	0.010%
13	6.263	535	540	544	rVB8	5323	10646	0.24%	0.022%
14	6.734	617	620	623	rVB5	3529	5738	0.13%	0.012%
15	6.763	623	625	628	rBV4	4454	6749	0.15%	0.014%
16	7.110	677	684	693	rBV	913263	1457551	33.36%	2.956%
17	7.287	708	714	728	rVB	810579	1268257	29.03%	2.572%
18	7.469	737	745	766	rVB	954193	1494605	34.21%	3.031%
19	7.657	772	777	785	rVB	7312	16020	0.37%	0.032%
20	7.946	819	826	835	rBV	660901	1040465	23.82%	2.110%
21	8.187	860	867	871	rVB8	5877	11628	0.27%	0.024%
22	8.657	940	947	958	rBV	1073223	1742762	39.89%	3.535%
23	9.016	1005	1008	1014	rVB8	2640	4598	0.11%	0.009%
24	9.116	1018	1025	1034	rBV	843256	1372487	31.42%	2.784%
25	9.834	1139	1147	1159	rBV	654804	1112262	25.46%	2.256%
26	10.369	1230	1238	1253	rBV	973430	1687680	38.63%	3.423%
27	10.622	1278	1281	1284	rVB5	4559	6354	0.15%	0.013%
28	10.751	1295	1303	1315	rVB	897762	1511484	34.60%	3.066%
29	10.898	1321	1328	1351	rBV	968844	1836119	42.03%	3.724%
30	11.134	1365	1368	1371	rBV5	3915	4862	0.11%	0.010%
31	11.416	1408	1416	1417	rBV6	2935	7204	0.16%	0.015%
32	11.439	1417	1420	1426	rVB8	4592	8175	0.19%	0.017%
33	11.516	1431	1433	1438	rVB6	5015	5487	0.13%	0.011%
34	11.592	1443	1446	1449	rBV5	3274	4403	0.10%	0.009%
35	12.345	1567	1574	1579	rBV	19407	33296	0.76%	0.068%
36	13.492	1762	1769	1776	rBV3	19955	35124	0.80%	0.071%

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Peak Location: TOP

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37	13.933	1838	1844	1845	rVB6	2794	4443	0.10%	0.009%
38	14.010	1850	1857	1870	rBV	1754755	2413249	55.24%	4.895%
39	14.280	1896	1903	1920	rBV	2065714	3009433	68.89%	6.104%
40	14.586	1947	1955	1963	rBV	1277087	1910007	43.72%	3.874%
41	14.798	1984	1991	2013	rBV	379095	1015851	23.25%	2.060%
42	15.263	2063	2070	2073	rVB9	3197	6256	0.14%	0.013%
43	15.580	2116	2124	2135	rBV	2541702	3703294	84.77%	7.511%
44	15.704	2138	2145	2158	rVV	689084	1012742	23.18%	2.054%
45	15.821	2161	2165	2171	rVB4	13680	22837	0.52%	0.046%
46	16.363	2253	2257	2264	rBV10	8910	11864	0.27%	0.024%
47	17.010	2363	2367	2373	rBV8	4631	10150	0.23%	0.021%
48	17.333	2415	2422	2433	rBV	1540594	2195524	50.26%	4.453%
49	17.433	2433	2439	2454	rVV	2752997	3886762	88.97%	7.883%
50	18.251	2574	2578	2583	rBV7	7370	11312	0.26%	0.023%
51	19.298	2751	2756	2763	rBV2	34587	54891	1.26%	0.111%
52	19.368	2763	2768	2771	rBV2	27349	38757	0.89%	0.079%
53	19.680	2818	2821	2823	rBV4	7617	9139	0.21%	0.019%
54	19.733	2823	2830	2839	rVV	3171097	4368602	100.00%	8.861%
55	21.533	3131	3136	3146	rBV	1547358	2076707	47.54%	4.212%
56	23.797	3514	3521	3536	rBV2	2116533	4262713	97.58%	8.646%
57	23.956	3541	3548	3561	rVB	1056507	2264804	51.84%	4.594%

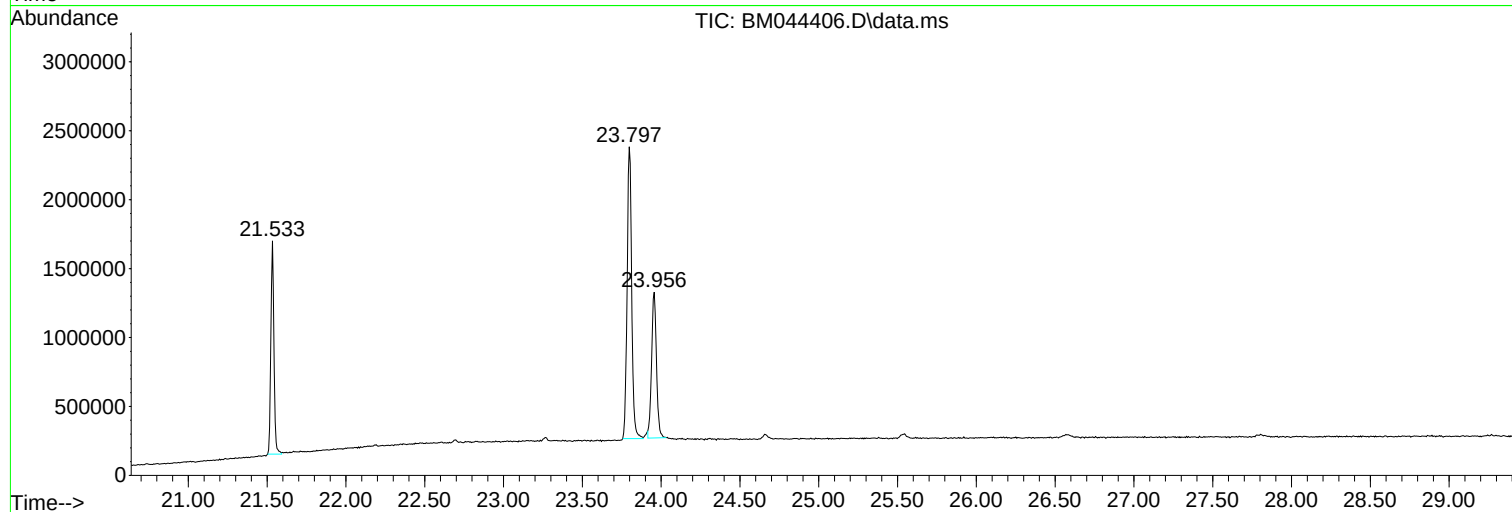
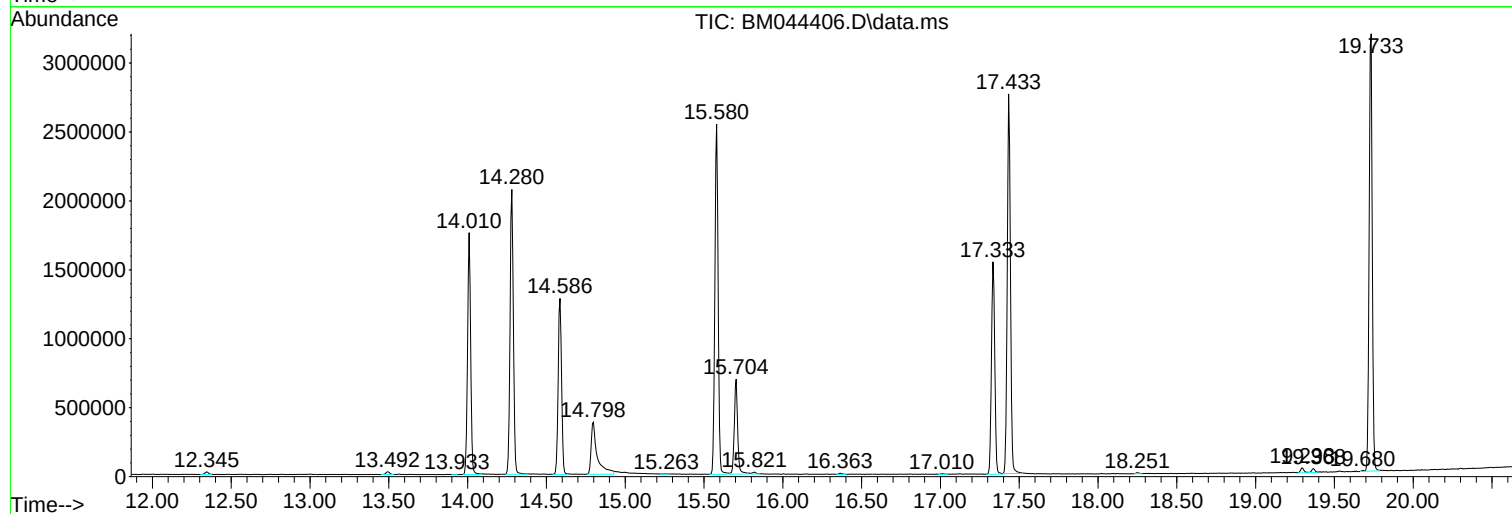
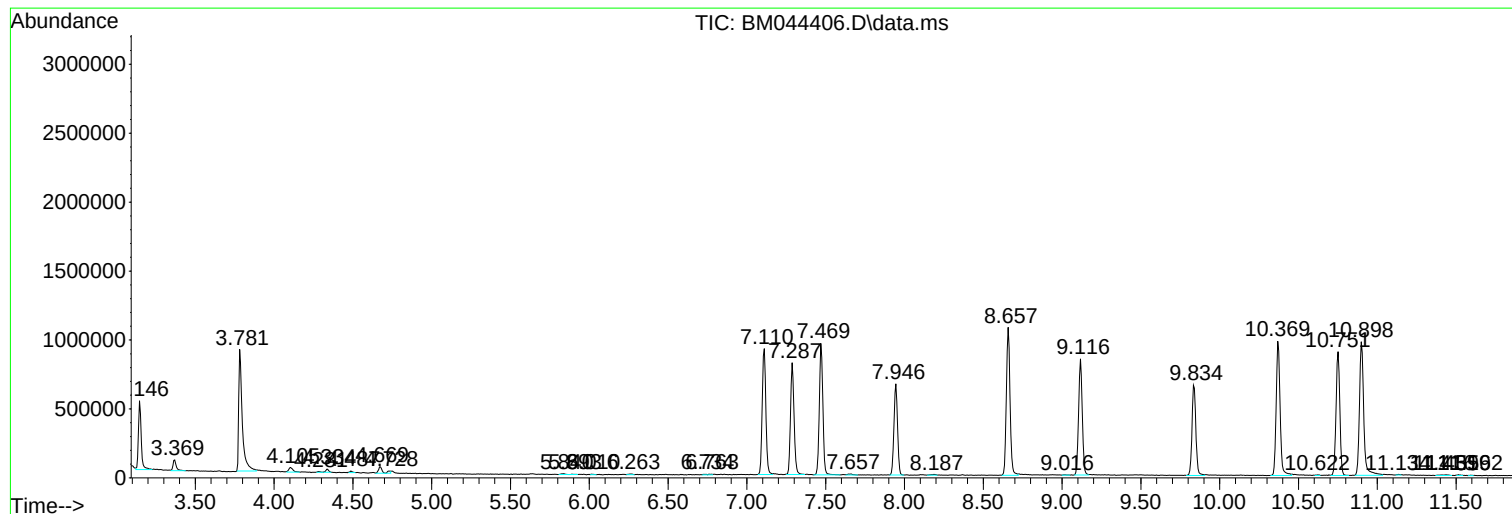
Sum of corrected areas: 49303471

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
Quant Title : SVOA CALIBRATION

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



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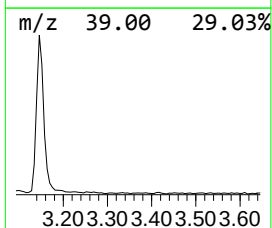
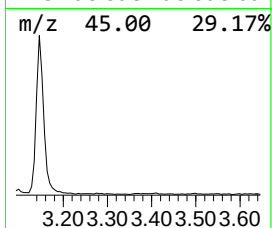
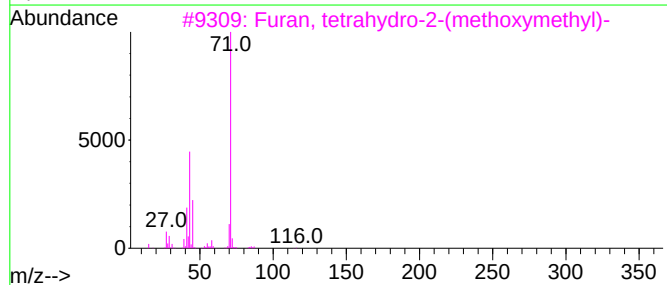
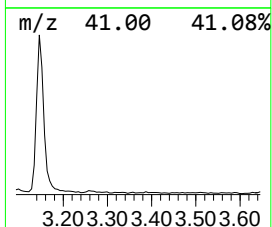
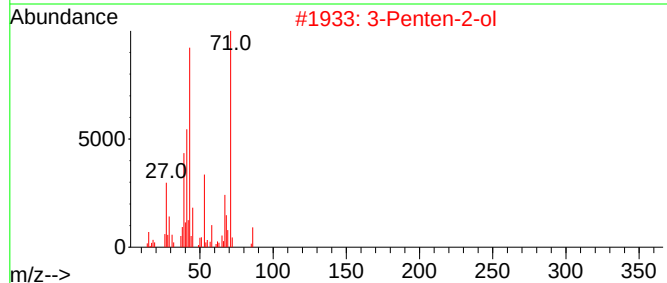
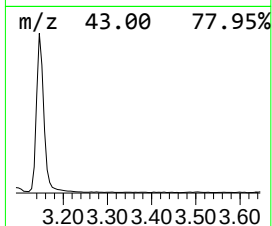
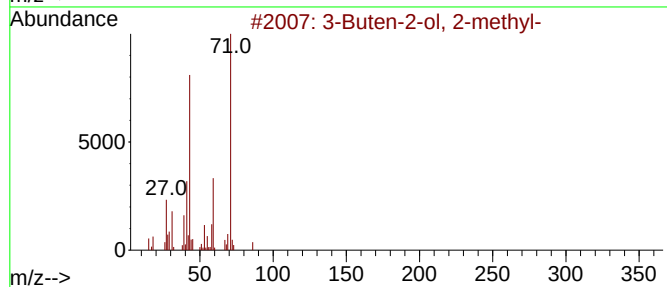
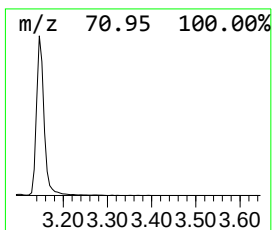
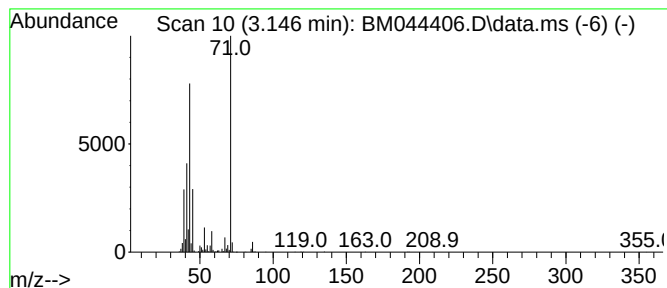
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 3-Buten-2-ol, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.146	11.50 ng/ul	598402	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	72
2			3-Penten-2-ol	86	C5H10O	001569-50-2	72
3			Furan, tetrahydro-2-(methoxymeth...	116	C6H12O2	019354-27-9	72
4			3-Penten-2-ol	86	C5H10O	001569-50-2	64
5			3-Penten-2-ol	86	C5H10O	001569-50-2	56



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Buten-2-ol, 2...	3.146	11.5	ng/ul	598402	1	7.946	1040470	20.0