

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040922\
 Data File : BM034573.D
 Acq On : 08 Apr 2022 22:47
 Operator : CG/JU
 Sample : N2319-06
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 YB2H3

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-BM040922.M
 Title : SVOA CALIBRATION

Signal : TIC: BM034573.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.225	27	32	42	rBV2	192066	328201	8.53%	0.957%
2	3.690	108	111	127	rBV	63289	120451	3.13%	0.351%
3	4.019	165	167	170	rVB2	11529	12458	0.32%	0.036%
4	4.560	254	259	270	rVB	142882	220908	5.74%	0.644%
5	5.907	482	488	496	rBV	69098	111446	2.90%	0.325%
6	6.696	618	622	624	rBV3	2506	4183	0.11%	0.012%
7	6.878	649	653	657	rVB2	6347	8159	0.21%	0.024%
8	7.043	674	681	695	rBV	91831	189152	4.92%	0.551%
9	7.201	702	708	722	rBV	536060	876269	22.78%	2.554%
10	7.407	736	743	758	rBV	495317	851059	22.12%	2.481%
11	7.878	816	823	833	rBV	470031	776769	20.19%	2.264%
12	8.119	861	864	871	rVB5	3494	7185	0.19%	0.021%
13	8.590	936	944	957	rBV	321311	584842	15.20%	1.705%
14	9.037	1014	1020	1035	rBV	625591	1097144	28.52%	3.198%
15	9.484	1091	1096	1102	rVB2	13567	22841	0.59%	0.067%
16	9.766	1136	1144	1156	rBV	494454	859392	22.34%	2.505%
17	10.307	1229	1236	1252	rBV	614582	1125757	29.26%	3.281%
18	10.683	1293	1300	1315	rBV	628053	1073613	27.91%	3.129%
19	10.825	1317	1324	1349	rBV	368691	858608	22.32%	2.503%
20	12.130	1544	1546	1553	rVB8	3610	6579	0.17%	0.019%
21	12.301	1570	1575	1583	rVB3	7176	17305	0.45%	0.050%
22	13.366	1750	1756	1762	rBV4	8307	12347	0.32%	0.036%
23	13.936	1846	1853	1865	rBV	1395329	1990150	51.73%	5.801%
24	14.219	1894	1901	1915	rBV	1560711	2298820	59.76%	6.700%
25	14.395	1927	1931	1935	rBV7	3929	6710	0.17%	0.020%
26	14.524	1946	1953	1964	rBV	916063	1328420	34.53%	3.872%
27	14.754	1985	1992	1994	rBV7	11135	24343	0.63%	0.071%
28	15.518	2115	2122	2136	rBV	1979475	2968501	77.17%	8.652%
29	15.630	2136	2141	2157	rVV	599347	980995	25.50%	2.859%
30	15.760	2159	2163	2168	rVB	18885	28422	0.74%	0.083%
31	16.307	2252	2256	2259	rVB3	6195	8343	0.22%	0.024%
32	17.054	2376	2383	2389	rBV7	6616	16113	0.42%	0.047%
33	17.265	2413	2419	2429	rBV	1011147	1471067	38.24%	4.288%
34	17.365	2429	2436	2461	rVB	2214614	3293281	85.61%	9.599%
35	18.171	2568	2573	2574	rBV2	27456	35007	0.91%	0.102%
36	18.189	2574	2576	2581	rVB4	31185	39752	1.03%	0.116%

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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

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37	18.595	2642	2645	2650	rVB7	4677	5170	0.13%	0.015%
38	19.224	2748	2752	2757	rBV4	26740	37753	0.98%	0.110%
39	19.295	2760	2764	2769	rVV2	20649	30238	0.79%	0.088%
40	19.454	2787	2791	2793	rBV4	11756	17917	0.47%	0.052%
41	19.659	2819	2826	2832	rBV	2759666	3846838	100.00%	11.212%
42	21.153	3077	3080	3087	rBV	83248	123099	3.20%	0.359%
43	21.442	3124	3129	3135	rBV	825379	1282039	33.33%	3.737%
44	23.671	3501	3508	3525	rBV2	1650473	3689874	95.92%	10.755%
45	23.824	3528	3534	3545	rVB	765280	1621081	42.14%	4.725%

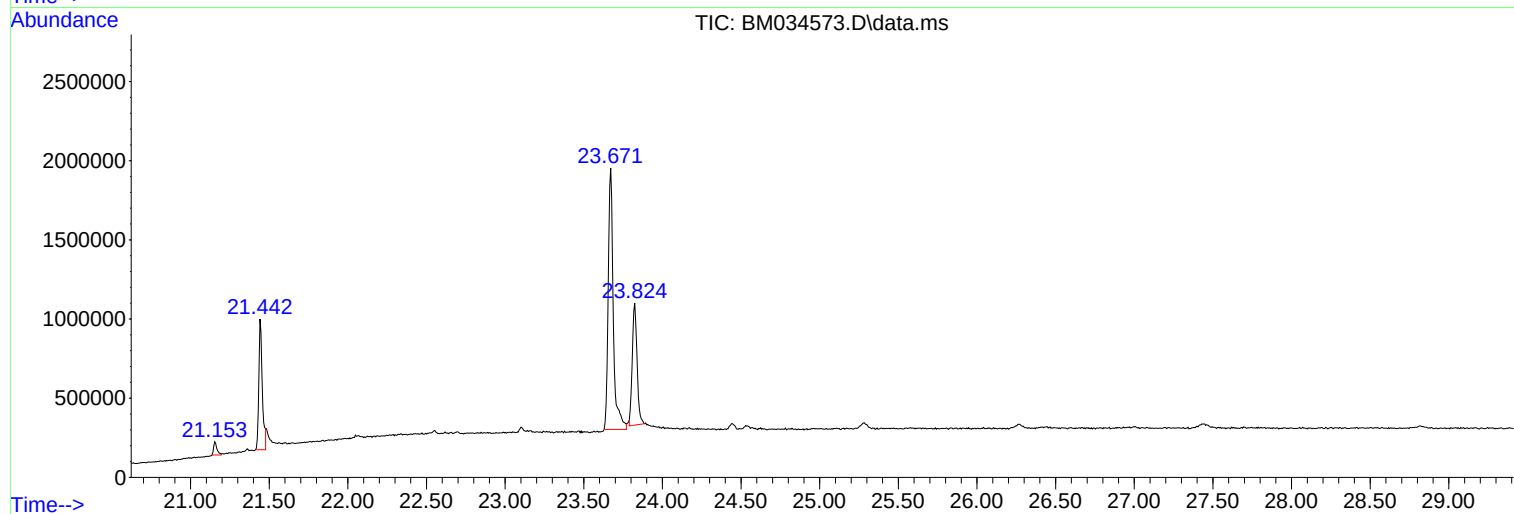
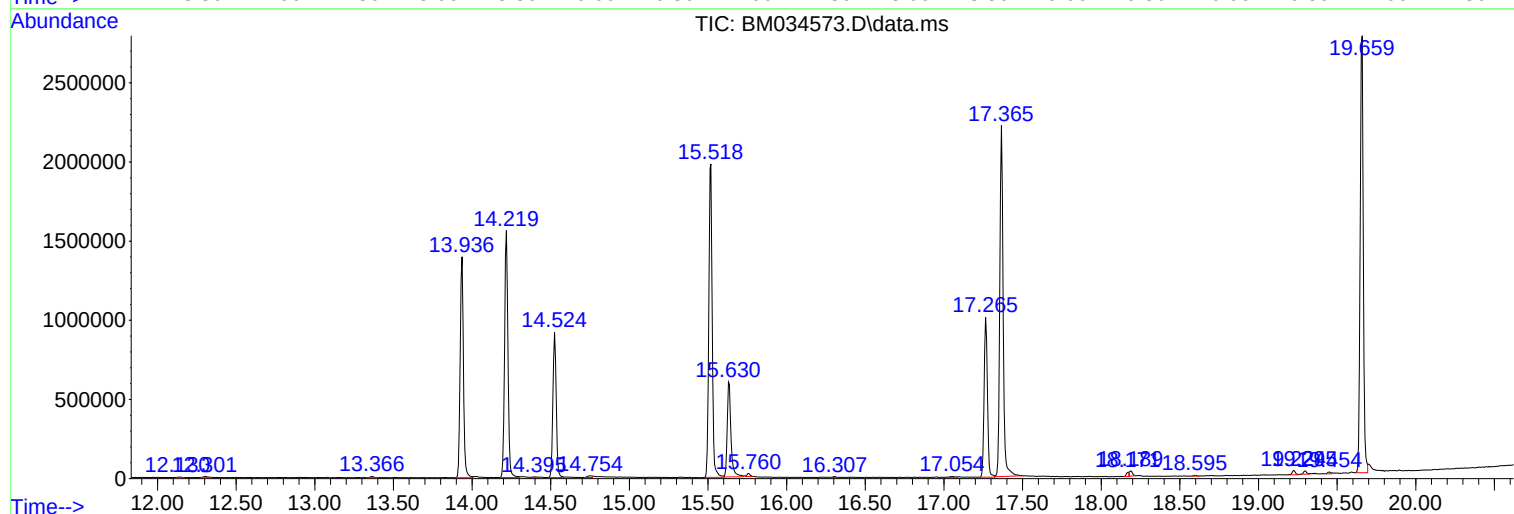
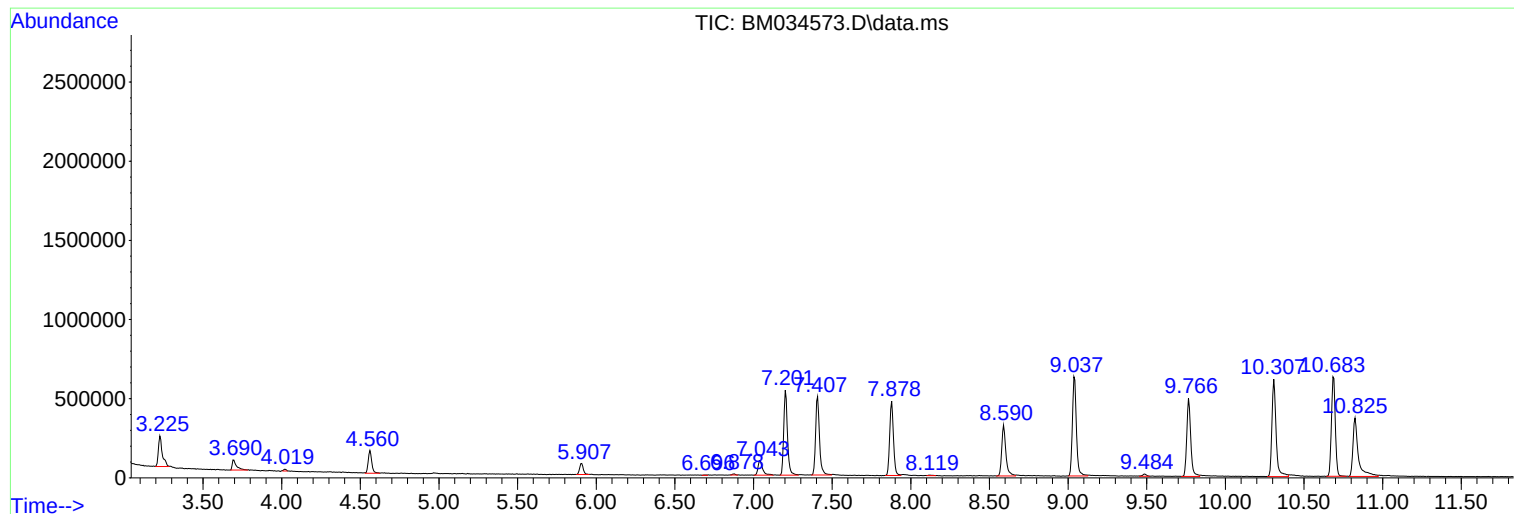
Sum of corrected areas: 34308601

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

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



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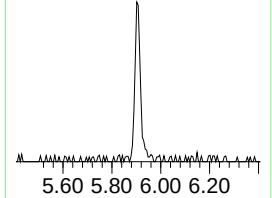
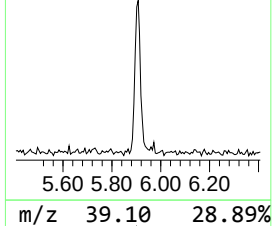
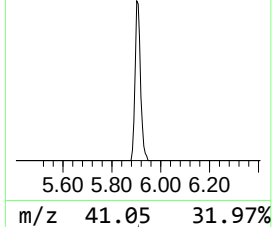
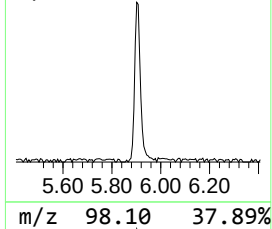
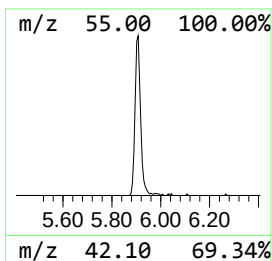
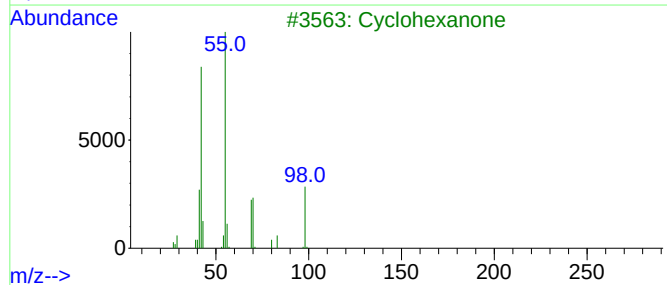
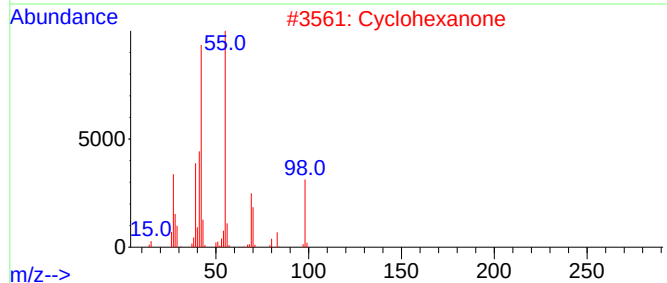
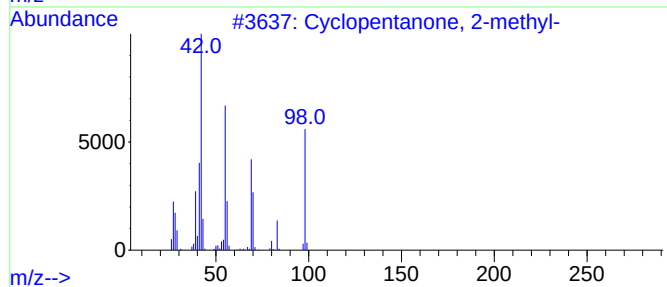
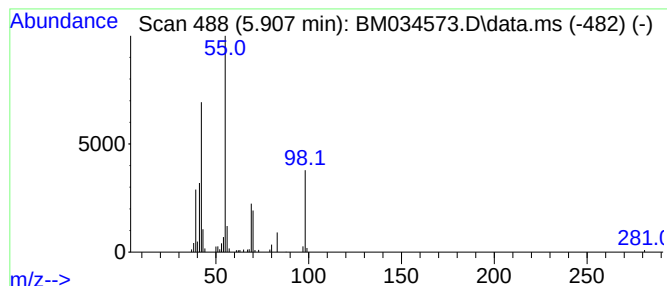
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Cyclopentanone, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.907	2.87 ng/ul	111446	1,4-Dichlorobenzene-d4	7.878

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentanone, 2-methyl-	98	C6H10O	001120-72-5	93
2			Cyclohexanone	98	C6H10O	000108-94-1	91
3			Cyclohexanone	98	C6H10O	000108-94-1	90
4			Cyclohexanone	98	C6H10O	000108-94-1	90
5			Cyclopentanone, 2-methyl-	98	C6H10O	001120-72-5	80



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclopentanone,...	5.907	2.9	ng/ul	111446	1	7.878	776769	20.0