

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011980.D
 Acq On : 06 Oct 2022 20:53
 Operator : CG/JU
 Sample : N4923-23
 Misc :
 ALS Vial : 68 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E10029

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-BP093022.M
 Title : SVOA CALIBRATION

Signal : TIC: BP011980.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.028	3	7	29	rVB	5317457	6414054	97.18%	9.846%
2	3.299	49	53	63	rVB	46697	68520	1.04%	0.105%
3	3.393	67	69	73	rVB2	11586	13915	0.21%	0.021%
4	3.646	109	112	116	rVB	15333	18005	0.27%	0.028%
5	3.728	124	126	140	rBV	130680	220157	3.34%	0.338%
6	3.946	159	163	169	rVB	29607	43826	0.66%	0.067%
7	4.040	176	179	185	rVB	26035	34145	0.52%	0.052%
8	4.987	335	340	346	rVB3	13996	23050	0.35%	0.035%
9	5.375	401	406	411	rBV	18170	26679	0.40%	0.041%
10	5.505	423	428	436	rBV	37605	71773	1.09%	0.110%
11	5.875	485	491	498	rVB2	25199	44374	0.67%	0.068%
12	7.046	680	690	704	rBV	133277	281343	4.26%	0.432%
13	7.210	711	718	734	rBV	761323	1267997	19.21%	1.946%
14	7.405	745	751	764	rBV	663207	1089837	16.51%	1.673%
15	7.493	764	766	774	rVB9	6610	8020	0.12%	0.012%
16	7.875	822	831	841	rBV	961569	1559222	23.62%	2.394%
17	8.140	872	876	882	rBV8	4397	7858	0.12%	0.012%
18	8.587	946	952	967	rBV	430816	758655	11.49%	1.165%
19	9.046	1022	1030	1045	rBV	825984	1433420	21.72%	2.200%
20	9.204	1054	1057	1061	rVB5	8840	11752	0.18%	0.018%
21	9.763	1143	1152	1166	rBV	595000	1042895	15.80%	1.601%
22	10.304	1237	1244	1268	rBV	883610	1623416	24.60%	2.492%
23	10.687	1300	1309	1322	rBV	1312158	2286614	34.65%	3.510%
24	10.828	1325	1333	1355	rBV	843314	1607711	24.36%	2.468%
25	12.275	1573	1579	1589	rBV	17656	37330	0.57%	0.057%
26	13.422	1768	1774	1778	rBV2	18212	27871	0.42%	0.043%
27	13.934	1854	1861	1877	rBV	1878118	2694235	40.82%	4.136%
28	14.216	1902	1909	1924	rBV	2234351	3388406	51.34%	5.202%
29	14.522	1954	1961	1977	rBV	1987631	2994799	45.38%	4.597%
30	14.745	1992	1999	2013	rBV3	41227	149500	2.27%	0.229%
31	14.951	2029	2034	2044	rVB3	8346	21642	0.33%	0.033%
32	15.210	2072	2078	2081	rBV7	4322	6985	0.11%	0.011%
33	15.516	2123	2130	2144	rBV	3069317	4530693	68.65%	6.955%
34	15.639	2144	2151	2166	rVV	717747	1103187	16.72%	1.693%
35	15.763	2168	2172	2178	rVB3	15939	27867	0.42%	0.043%
36	16.310	2261	2265	2269	rVB7	4557	6846	0.10%	0.011%

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Max Peaks: 100

Peak Location: TOP

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37	17.281	2424	2430	2440	rBV	2566343	3556107	53.88%	5.459%
38	17.381	2440	2447	2462	rVV	3830904	5362978	81.26%	8.233%
39	18.169	2577	2581	2586	rBV5	22866	38893	0.59%	0.060%
40	19.192	2752	2755	2761	rVB2	41162	52331	0.79%	0.080%
41	19.263	2763	2767	2773	rBV2	44275	72232	1.09%	0.111%
42	19.616	2821	2827	2839	rBV	4915656	6392099	96.85%	9.812%
43	21.369	3120	3125	3132	rBV	3178402	4050220	61.37%	6.217%
44	23.639	3504	3511	3525	rVB2	3282468	6599981	100.00%	10.132%
45	23.792	3531	3537	3548	rVB2	2036156	4071154	61.68%	6.250%

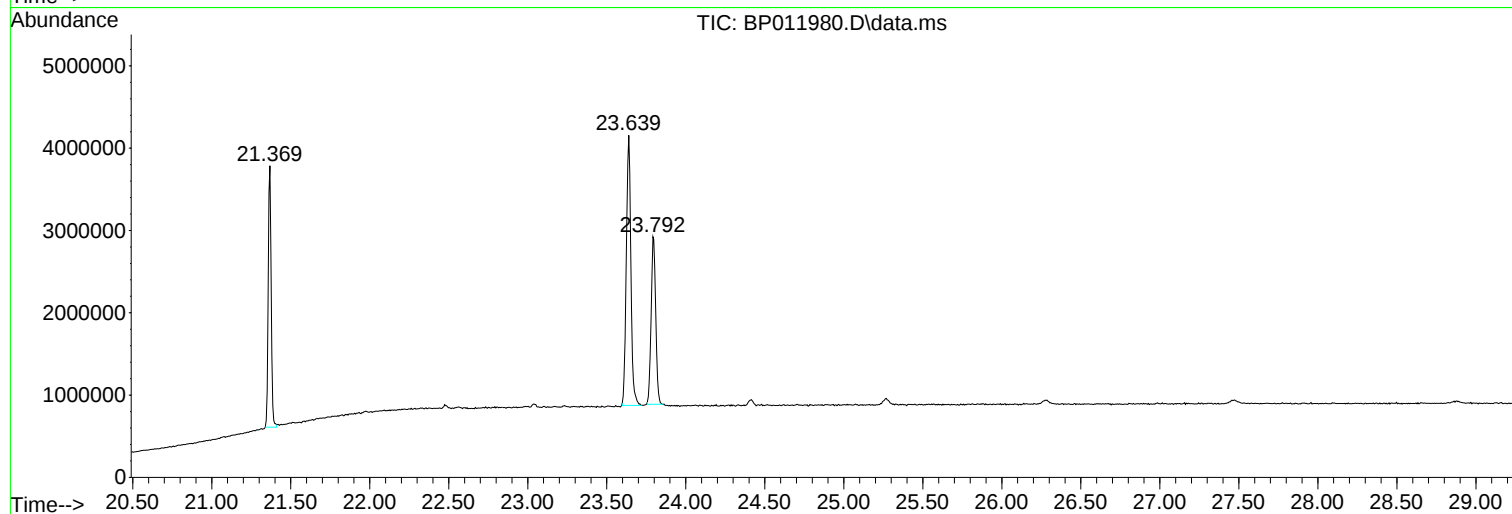
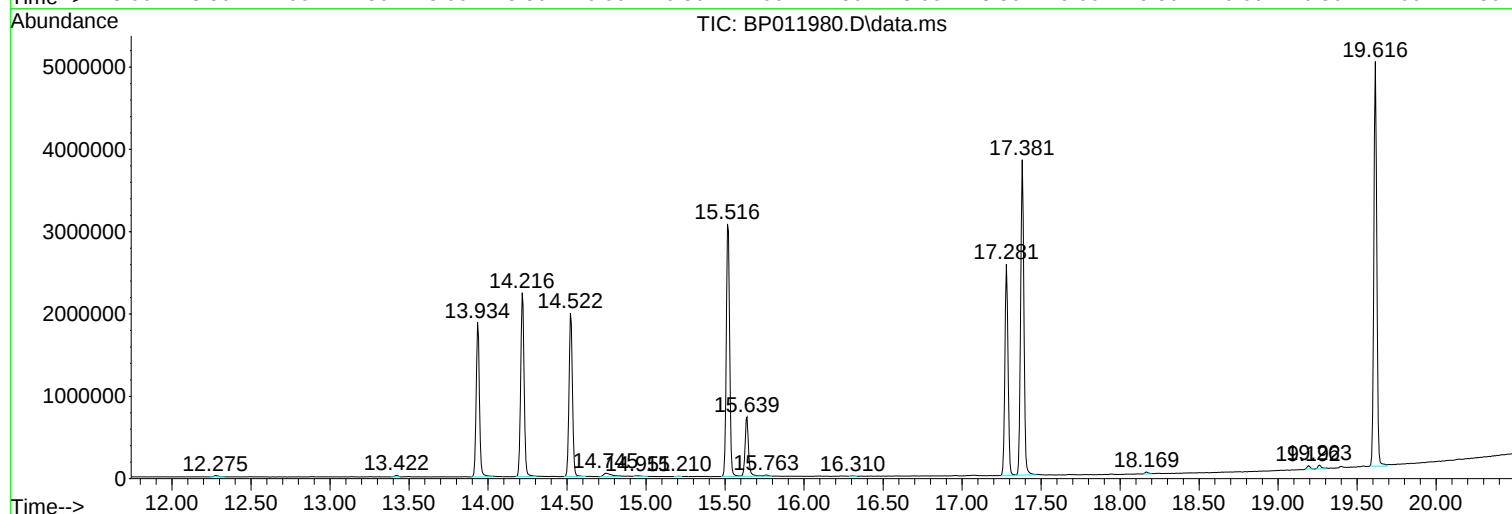
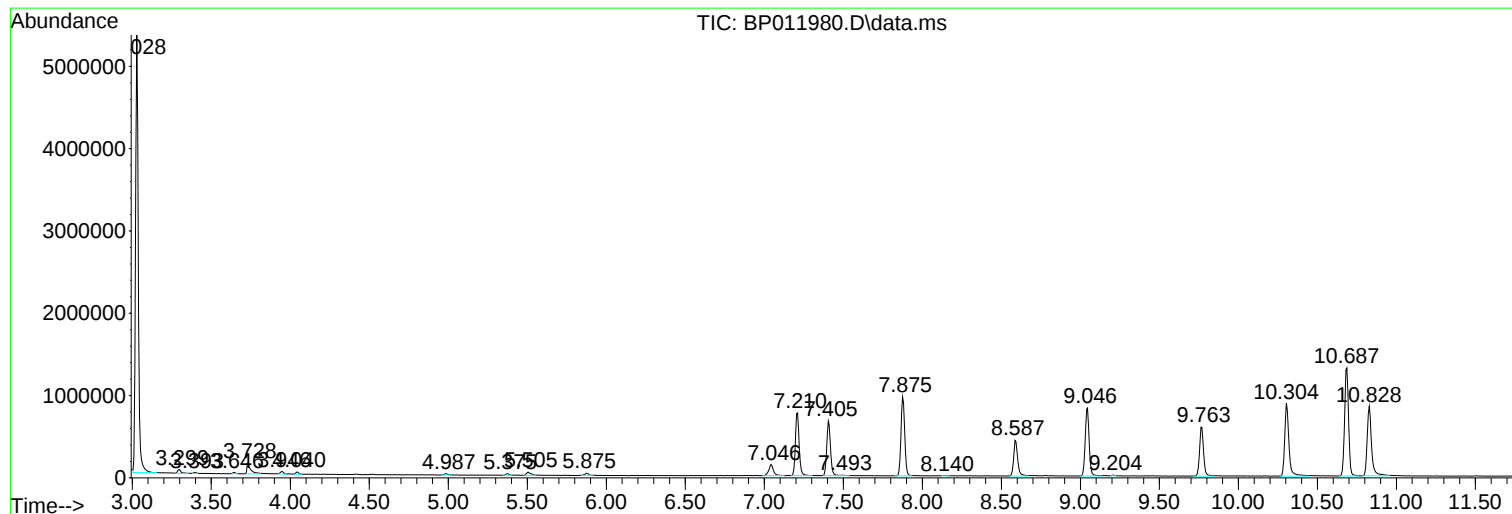
Sum of corrected areas: 65142594

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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



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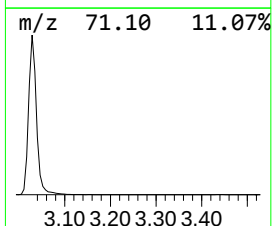
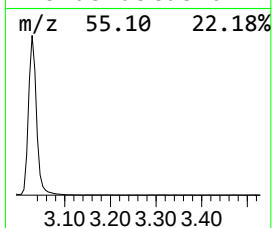
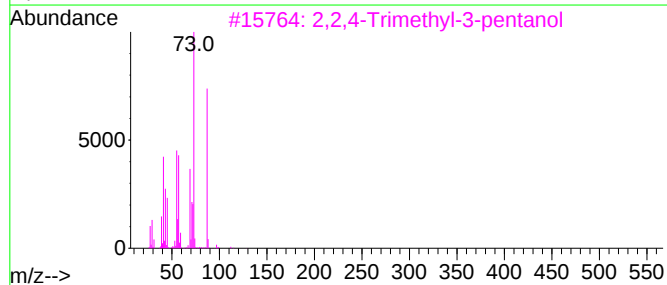
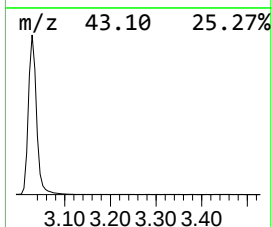
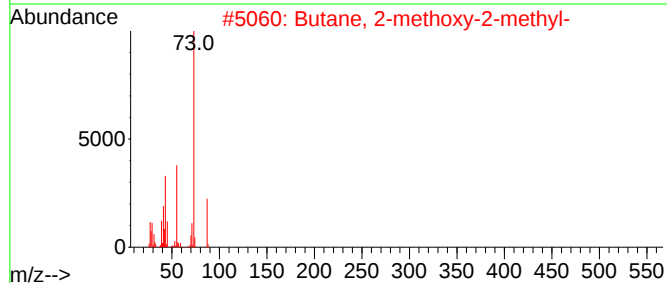
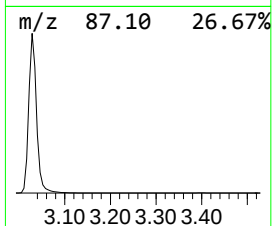
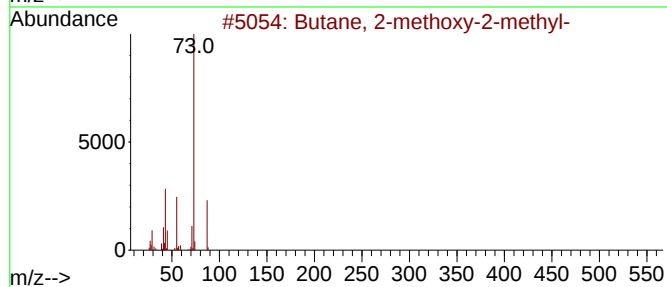
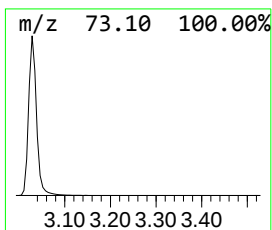
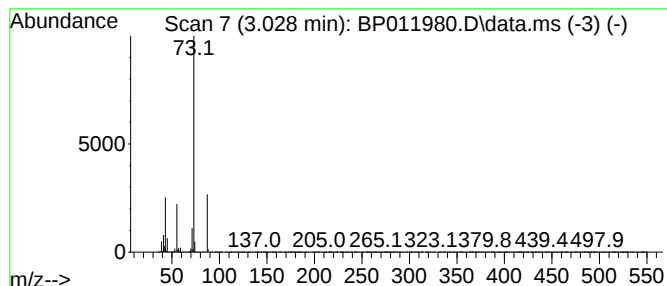
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.028	82.27 ng/ul	6414050	1,4-Dichlorobenzene-d4	7.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	86		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78		
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43		
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17		
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9		



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.028	82.3	ng/ul	6414050	1	7.875	1559220	20.0