

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011883.D
 Acq On : 03 Oct 2022 21:22
 Operator : CG/JU
 Sample : N4847-07
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB8R1

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: BP011883.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.034	4	8	30	rVB	4590363	5764619	80.61%	9.479%
2	3.305	50	54	63	rVB	51013	71157	0.99%	0.117%
3	3.734	124	127	138	rBV	158663	243545	3.41%	0.400%
4	3.958	159	165	173	rVB	29482	51865	0.73%	0.085%
5	4.052	177	181	187	rVB	17377	23080	0.32%	0.038%
6	4.987	335	340	347	rVB7	5162	10101	0.14%	0.017%
7	7.052	686	691	705	rVB	169709	305669	4.27%	0.503%
8	7.216	713	719	734	rBV	706186	1172673	16.40%	1.928%
9	7.416	746	753	764	rBV	675104	1109984	15.52%	1.825%
10	7.887	825	833	842	rBV	829015	1353821	18.93%	2.226%
11	8.599	947	954	971	rBV	499536	878376	12.28%	1.444%
12	9.057	1024	1032	1045	rBV	751649	1291791	18.06%	2.124%
13	9.463	1097	1101	1105	rBV2	9037	13438	0.19%	0.022%
14	9.781	1147	1155	1168	rBV	562143	964648	13.49%	1.586%
15	10.316	1239	1246	1268	rBV	851978	1564559	21.88%	2.573%
16	10.698	1303	1311	1325	rBV	1126879	1936672	27.08%	3.184%
17	10.840	1328	1335	1352	rBV	618859	1153497	16.13%	1.897%
18	12.240	1567	1573	1576	rBV3	4964	9261	0.13%	0.015%
19	12.287	1576	1581	1587	rVV2	17791	30159	0.42%	0.050%
20	12.516	1613	1620	1627	rBV2	5636	12521	0.18%	0.021%
21	13.357	1757	1763	1767	rVV6	8156	15473	0.22%	0.025%
22	13.428	1770	1775	1784	rVV	32715	54884	0.77%	0.090%
23	13.510	1784	1789	1795	rVB9	4760	9246	0.13%	0.015%
24	13.945	1856	1863	1879	rBV	1753646	2488531	34.80%	4.092%
25	14.228	1904	1911	1931	rBV	2085870	3087557	43.17%	5.077%
26	14.534	1956	1963	1972	rBV	1632643	2422658	33.88%	3.984%
27	14.745	1993	1999	2016	rBV	82506	220347	3.08%	0.362%
28	15.281	2084	2090	2097	rBV	128056	182846	2.56%	0.301%
29	15.381	2104	2107	2113	rVB5	6392	10817	0.15%	0.018%
30	15.528	2125	2132	2146	rBV	2944528	4194470	58.65%	6.897%
31	15.645	2146	2152	2167	rVV	819691	1130747	15.81%	1.859%
32	15.769	2169	2173	2182	rVB2	14776	27718	0.39%	0.046%
33	17.292	2425	2432	2441	rBV	1998423	2857659	39.96%	4.699%
34	17.392	2442	2449	2474	rVB	3415867	5026274	70.28%	8.265%
35	18.169	2577	2581	2591	rBV2	54876	94972	1.33%	0.156%
36	19.198	2753	2756	2763	rVB3	41220	54309	0.76%	0.089%

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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

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
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37	19.269	2764	2768	2772	rBV	53460	80841	1.13%	0.133%
38	19.404	2789	2791	2797	rVB5	27540	29239	0.41%	0.048%
39	19.622	2822	2828	2842	rBV	4691321	6342629	88.69%	10.429%
40	21.374	3121	3126	3133	rBV	2767807	3579446	50.05%	5.886%
41	23.651	3506	3513	3527	rVB2	3470665	7151621	100.00%	11.759%
42	23.810	3533	3540	3551	rVB2	1844914	3792155	53.03%	6.235%

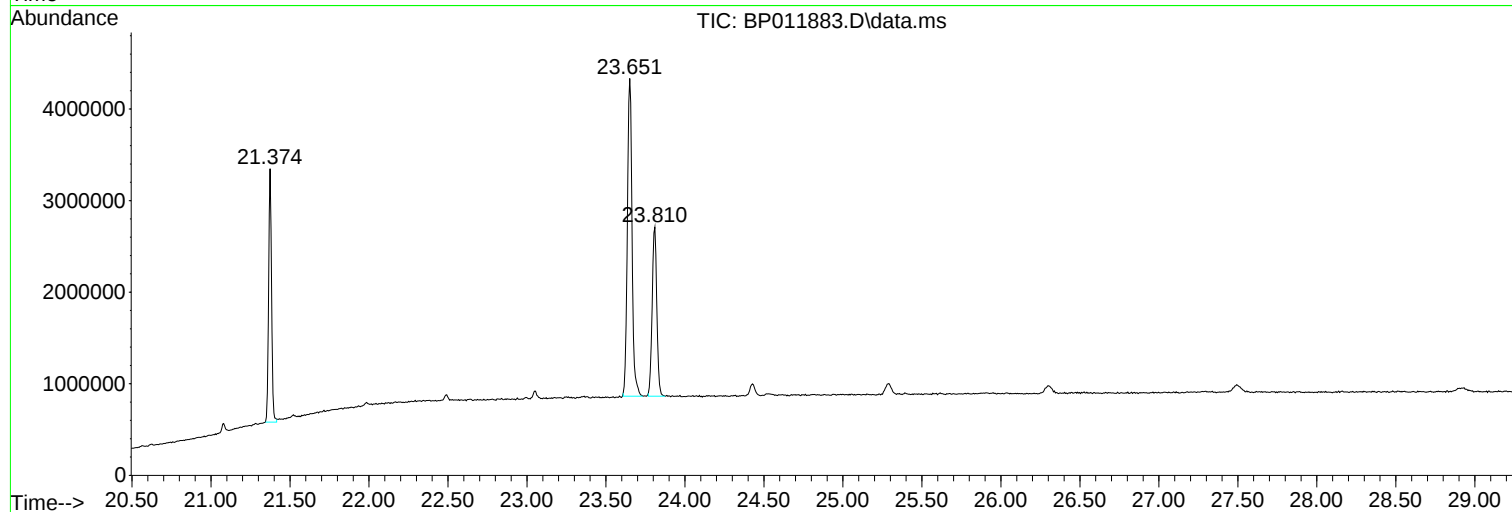
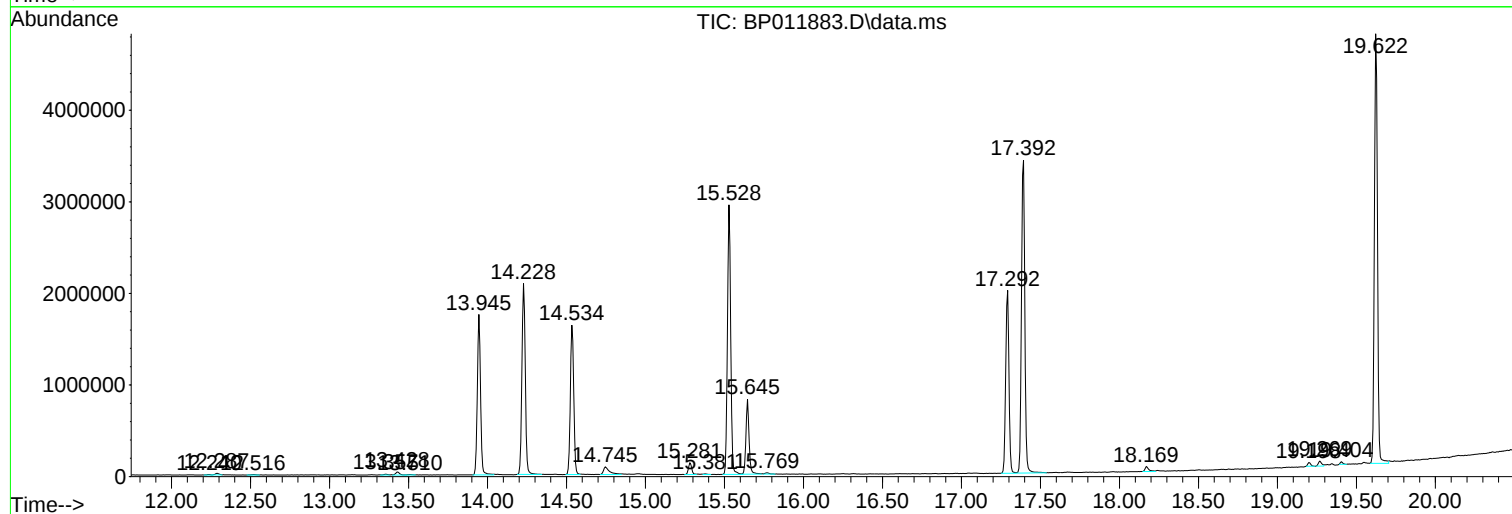
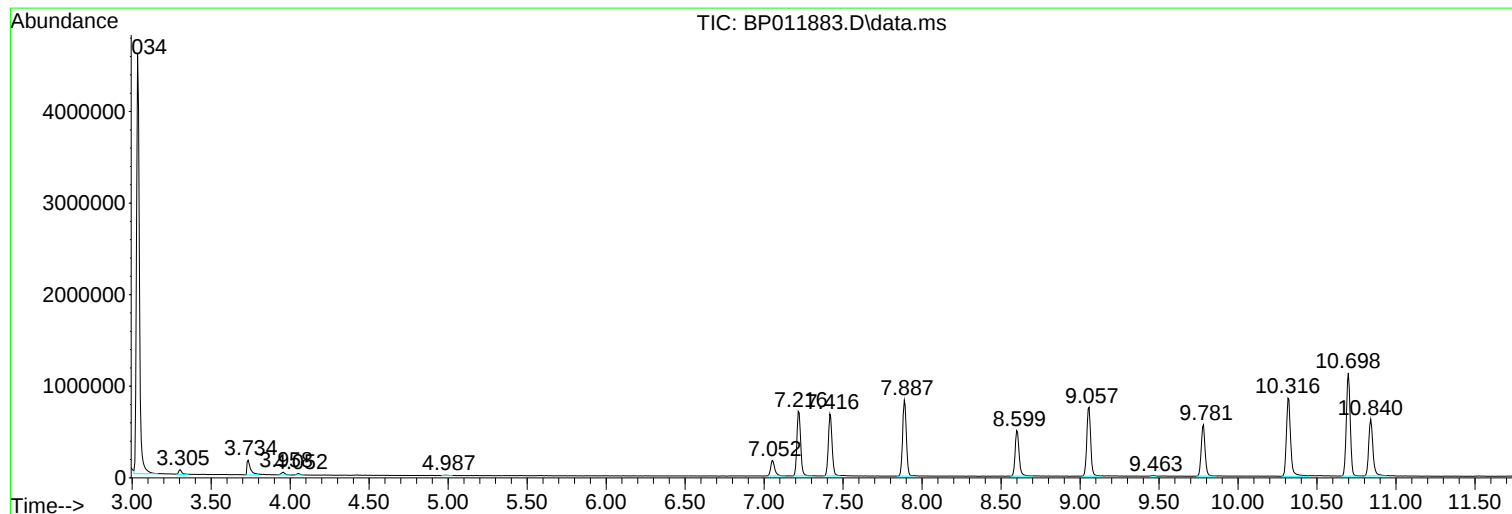
Sum of corrected areas: 60815875

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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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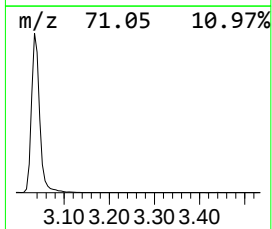
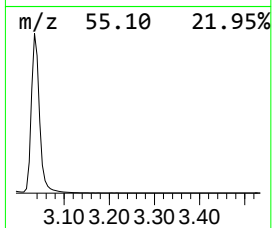
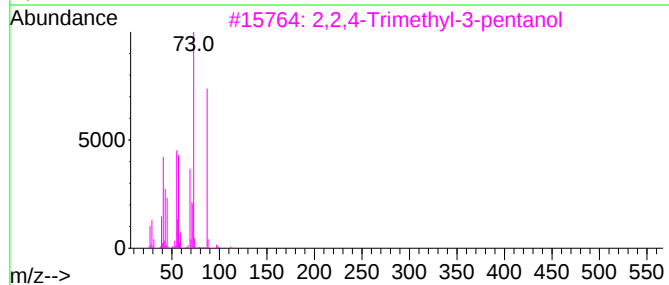
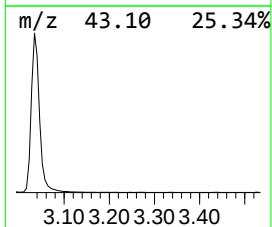
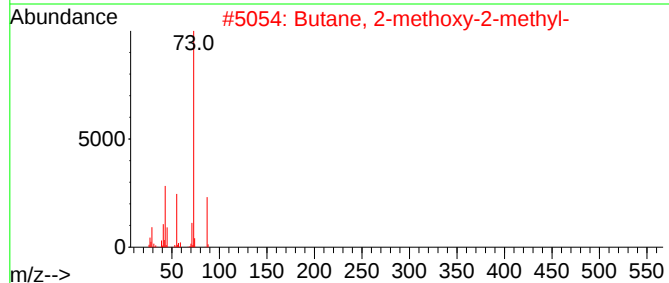
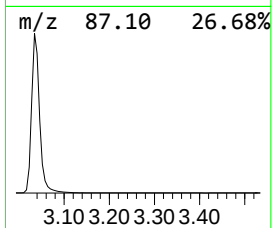
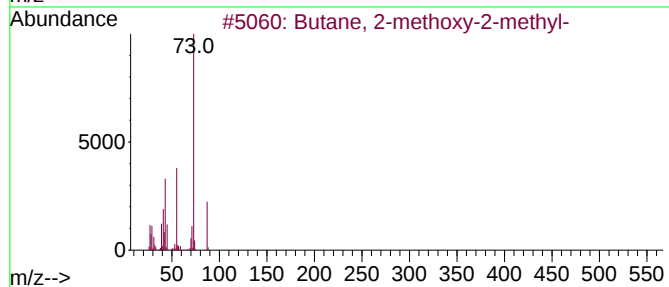
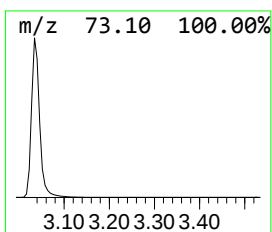
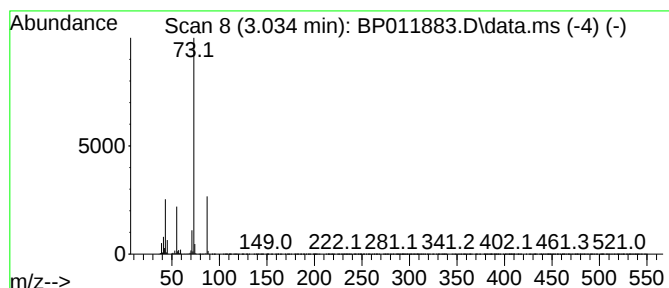
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.034	85.16 ng/ul	5764620	1,4-Dichlorobenzene-d4	7.887

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72		
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43		
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17		
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9		



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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.034	85.2	ng/u1	5764620	1	7.887	1353820	20.0