

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011923.D
 Acq On : 05 Oct 2022 11:04
 Operator : CG/JU
 Sample : PB148074BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK074

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: BP011923.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	16	rVB	190026	224558	3.13%	0.373%
2	3.299	49	53	62	rVB	64195	82016	1.14%	0.136%
3	3.722	122	125	146	rBV	642158	938276	13.08%	1.559%
4	4.052	176	181	186	rBV	16331	25087	0.35%	0.042%
5	7.046	683	690	704	rBV	762684	1239711	17.28%	2.060%
6	7.210	712	718	729	rBV	669745	1099746	15.33%	1.828%
7	7.410	745	752	769	rVB	830247	1332444	18.57%	2.214%
8	7.881	824	832	844	rBV	787865	1256344	17.51%	2.088%
9	8.593	944	953	967	rBV	949930	1594002	22.21%	2.649%
10	9.052	1023	1031	1044	rBV	671318	1166086	16.25%	1.938%
11	9.769	1146	1153	1167	rBV	533102	915031	12.75%	1.521%
12	10.310	1237	1245	1265	rBV	954806	1760629	24.54%	2.926%
13	10.693	1301	1310	1323	rVB	1157741	1994730	27.80%	3.315%
14	10.834	1326	1334	1358	rBV	928228	1770909	24.68%	2.943%
15	12.287	1574	1581	1589	rBV3	15150	35408	0.49%	0.059%
16	13.428	1769	1775	1780	rBV	15139	24442	0.34%	0.041%
17	13.504	1782	1788	1794	rBV5	4220	8738	0.12%	0.015%
18	13.940	1854	1862	1876	rBV	1724205	2495027	34.77%	4.147%
19	14.222	1901	1910	1932	rBV	2206067	3252609	45.33%	5.406%
20	14.528	1955	1962	1972	rBV	1810232	2613875	36.43%	4.344%
21	14.734	1990	1997	2019	rBV	491127	1080179	15.05%	1.795%
22	15.522	2124	2131	2145	rBV	2891758	4224938	58.88%	7.022%
23	15.639	2145	2151	2168	rVV	663336	972271	13.55%	1.616%
24	15.763	2168	2172	2180	rVB3	14290	26237	0.37%	0.044%
25	17.063	2391	2393	2398	rBV5	5766	9668	0.13%	0.016%
26	17.286	2424	2431	2441	rBV	2202708	3187618	44.42%	5.298%
27	17.386	2441	2448	2468	rVV	3365765	4886466	68.10%	8.121%
28	19.198	2753	2756	2762	rVB	37902	52532	0.73%	0.087%
29	19.263	2764	2767	2773	rBV	42538	62434	0.87%	0.104%
30	19.622	2822	2828	2841	rBV	4507708	6264756	87.30%	10.412%
31	21.369	3121	3125	3135	rBV	3165504	4033136	56.21%	6.703%
32	23.645	3505	3512	3524	rVB	3594599	7175737	100.00%	11.926%
33	23.804	3533	3539	3551	rVB2	2194186	4364054	60.82%	7.253%

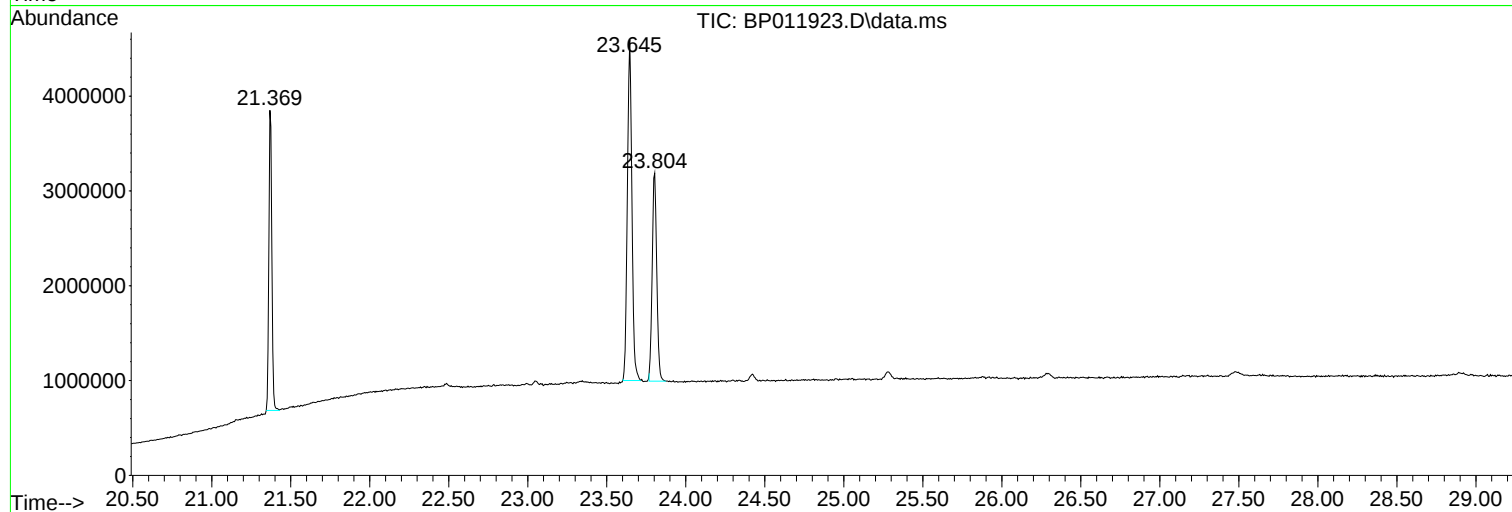
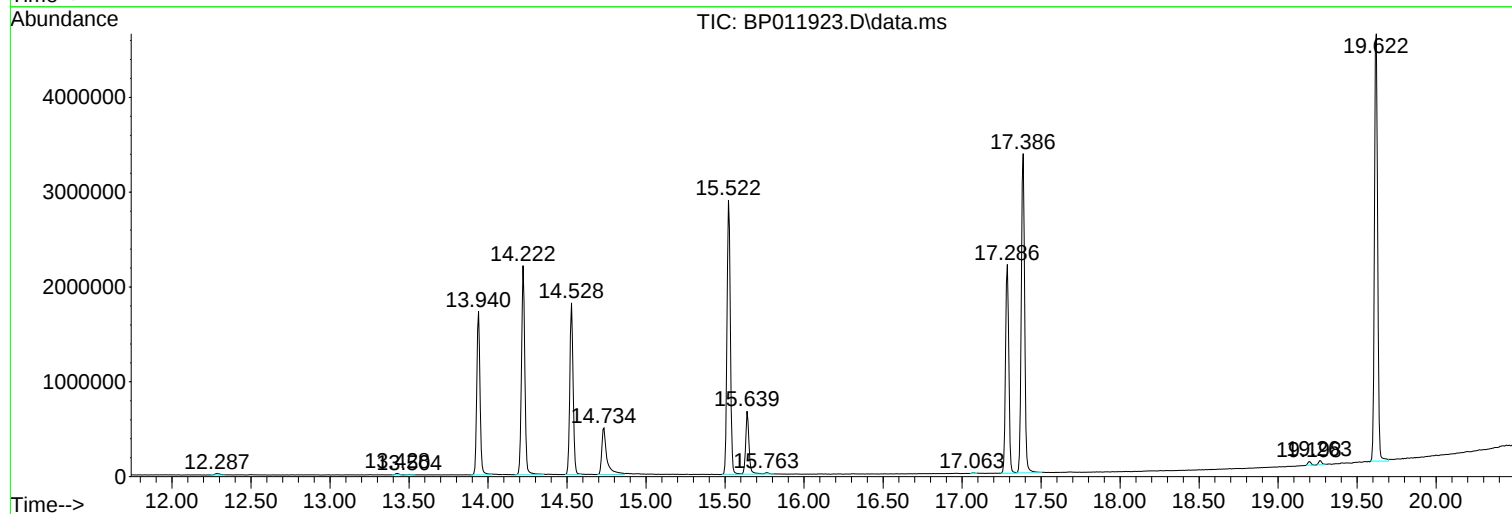
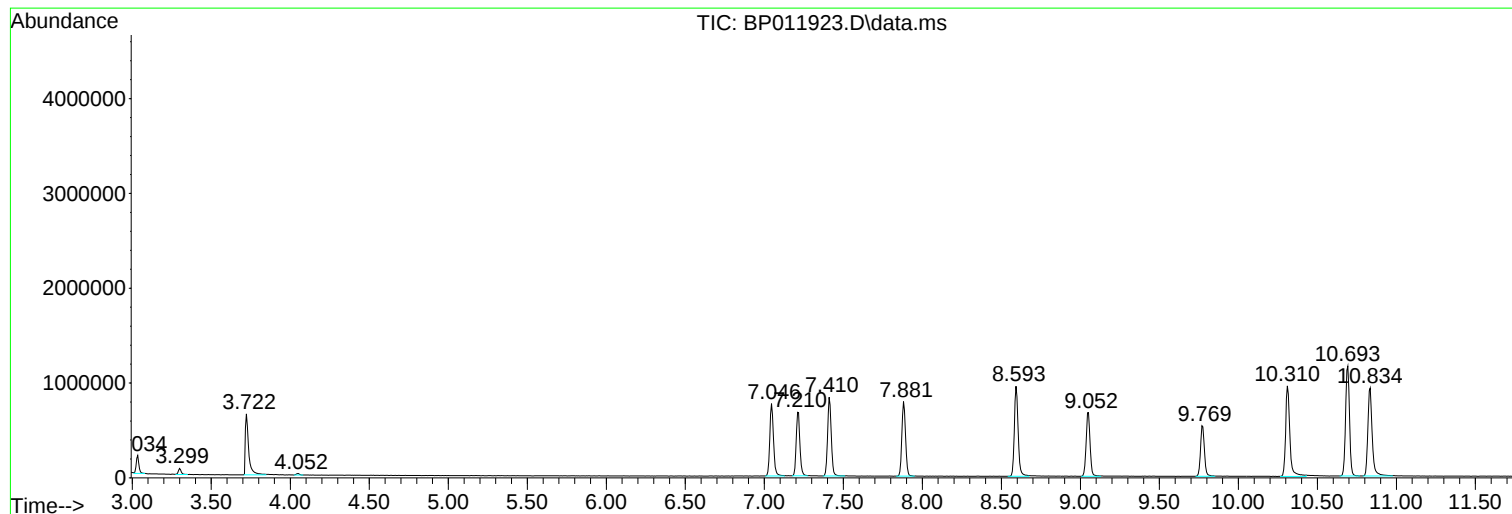
Sum of corrected areas: 60169694

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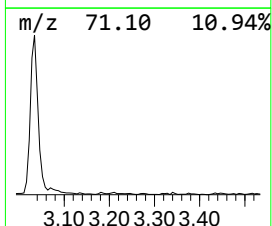
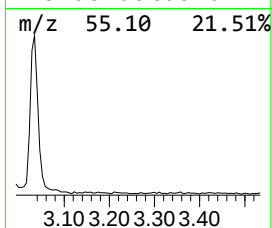
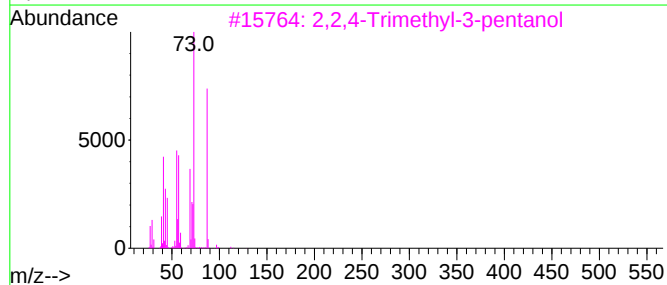
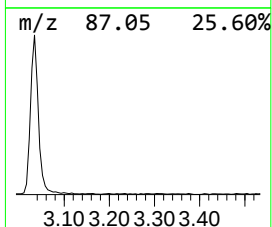
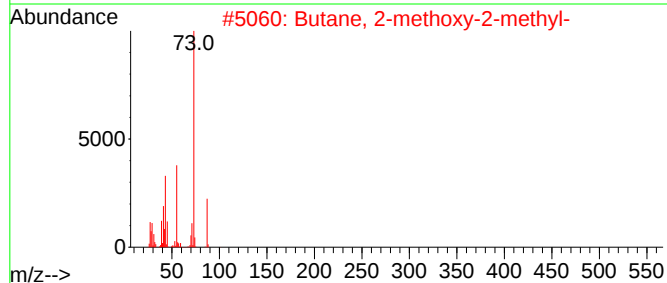
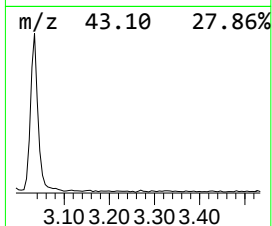
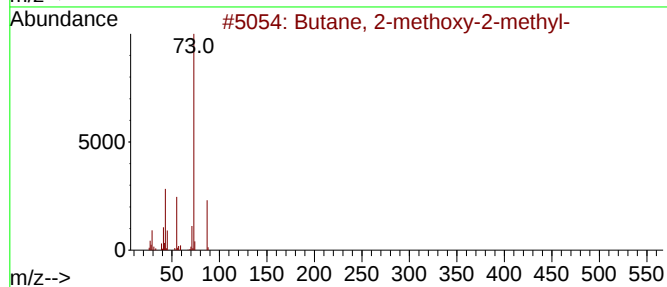
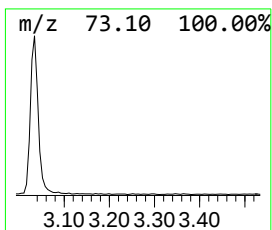
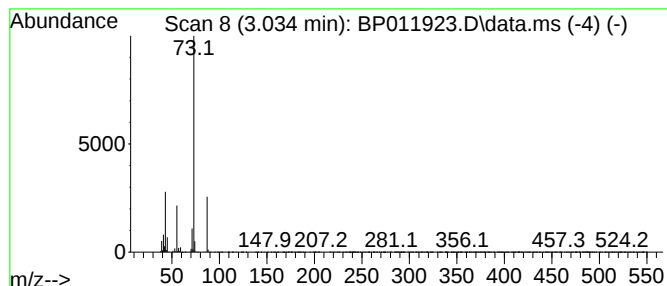
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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.
3.034	3.57 ng/ul	224558	1,4-Dichlorobenzene-d4	7.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	90		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64		
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43		
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9		
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9		



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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	3.6	ng/u1	224558	1	7.881	1256340	20.0