

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011888.D
 Acq On : 04 Oct 2022 00:46
 Operator : CG/JU
 Sample : PB148000BL
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK000

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: BP011888.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	14	rVB	212971	239784	3.31%	0.361%
2	3.299	49	53	61	rVB	65902	86594	1.20%	0.130%
3	3.722	122	125	147	rBV	755835	1132062	15.63%	1.703%
4	4.052	177	181	188	rVB	18317	28467	0.39%	0.043%
5	4.281	217	220	225	rVB3	7834	10222	0.14%	0.015%
6	7.052	678	691	711	rBV	958115	1560618	21.54%	2.348%
7	7.216	712	719	730	rVV	831400	1340827	18.51%	2.017%
8	7.416	746	753	767	rBV	1018501	1635172	22.57%	2.460%
9	7.887	824	833	843	rBV	967469	1539232	21.25%	2.316%
10	8.598	946	954	974	rBV	1194126	2021111	27.90%	3.041%
11	9.051	1023	1031	1047	rBV	864647	1468748	20.27%	2.210%
12	9.775	1147	1154	1168	rBV	672741	1171743	16.17%	1.763%
13	10.316	1236	1246	1276	rBV	1138252	2021666	27.90%	3.041%
14	10.698	1303	1311	1322	rBV	1308102	2228465	30.76%	3.352%
15	10.840	1327	1335	1359	rBV	1102930	2060176	28.44%	3.099%
16	12.292	1573	1582	1590	rBV2	18394	41821	0.58%	0.063%
17	13.434	1770	1776	1781	rBV4	7151	12287	0.17%	0.018%
18	13.504	1785	1788	1795	rVB7	4662	8617	0.12%	0.013%
19	13.945	1856	1863	1877	rBV	2023423	2804503	38.71%	4.219%
20	14.228	1902	1911	1929	rBV	2395784	3655555	50.46%	5.499%
21	14.533	1956	1963	1974	rBV	2035102	2881563	39.77%	4.335%
22	14.733	1989	1997	2019	rBV	714550	1362607	18.81%	2.050%
23	15.528	2125	2132	2146	rBV	3257251	4667157	64.42%	7.021%
24	15.645	2146	2152	2169	rVV	838579	1207881	16.67%	1.817%
25	15.769	2169	2173	2179	rVB4	16294	26350	0.36%	0.040%
26	17.292	2425	2432	2441	rBV	2333118	3473159	47.94%	5.225%
27	17.392	2442	2449	2462	rVV	3502394	5167800	71.33%	7.774%
28	19.204	2753	2757	2761	rBV	46365	61014	0.84%	0.092%
29	19.269	2765	2768	2773	rVV2	43160	59142	0.82%	0.089%
30	19.621	2822	2828	2835	rBV	4859664	6535800	90.21%	9.832%
31	21.374	3121	3126	3135	rBV	3466289	4293577	59.26%	6.459%
32	23.651	3506	3513	3530	rVB2	3528554	7244876	100.00%	10.899%
33	23.809	3533	3540	3552	rVB2	2134196	4423361	61.06%	6.654%

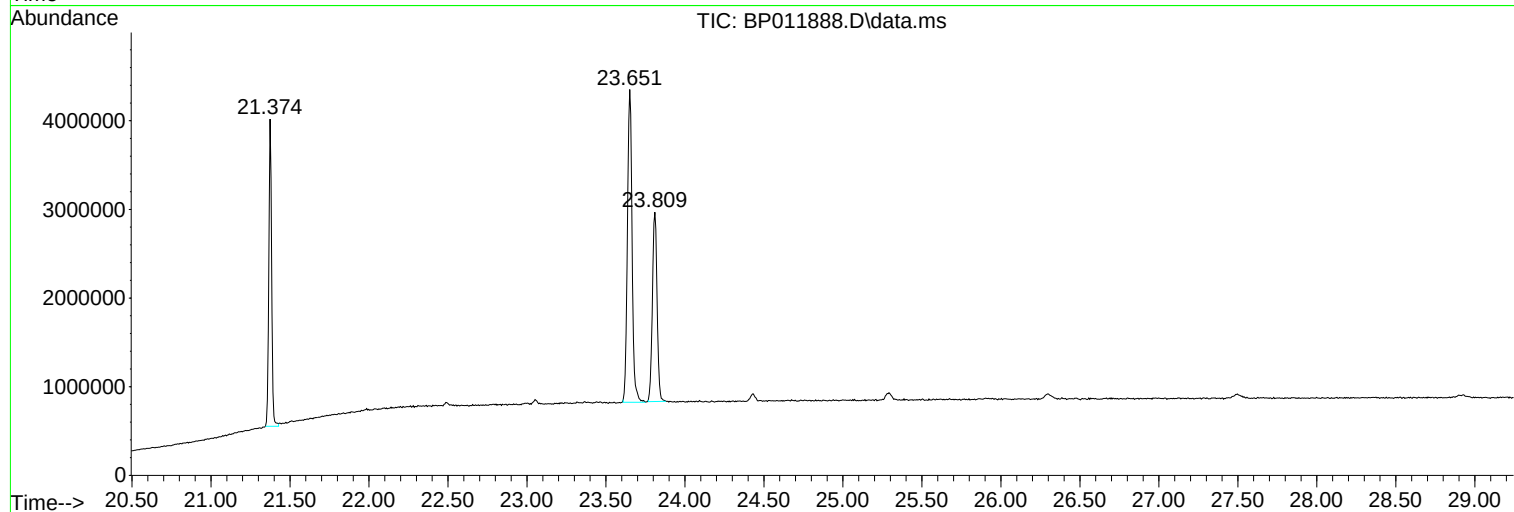
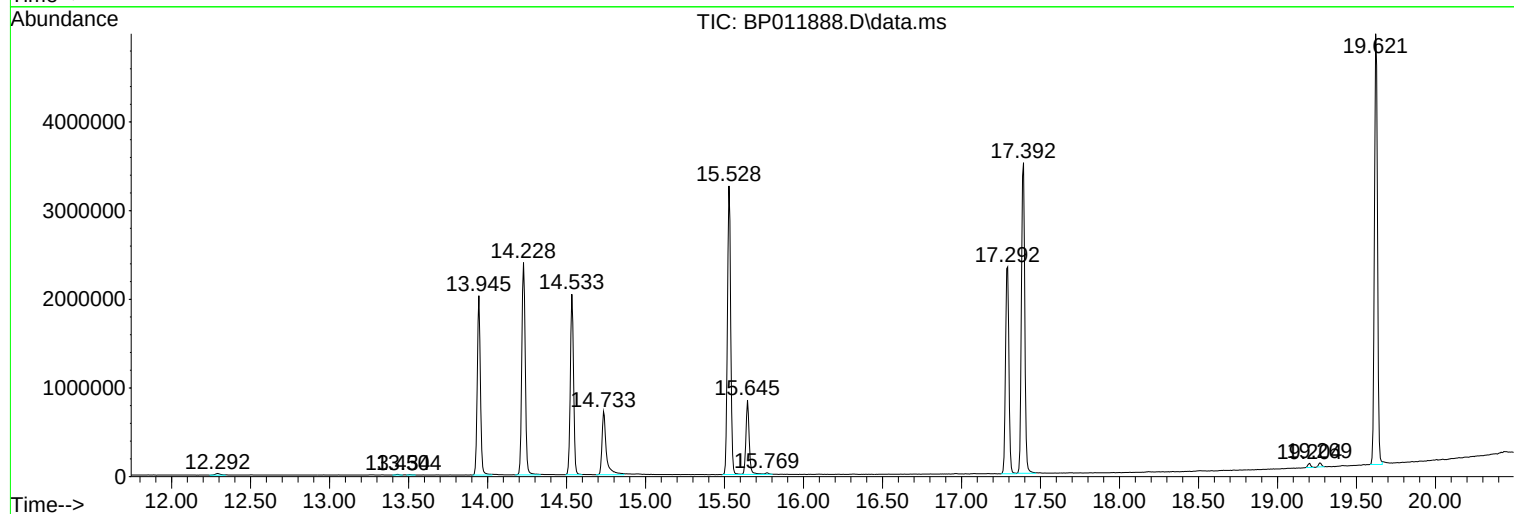
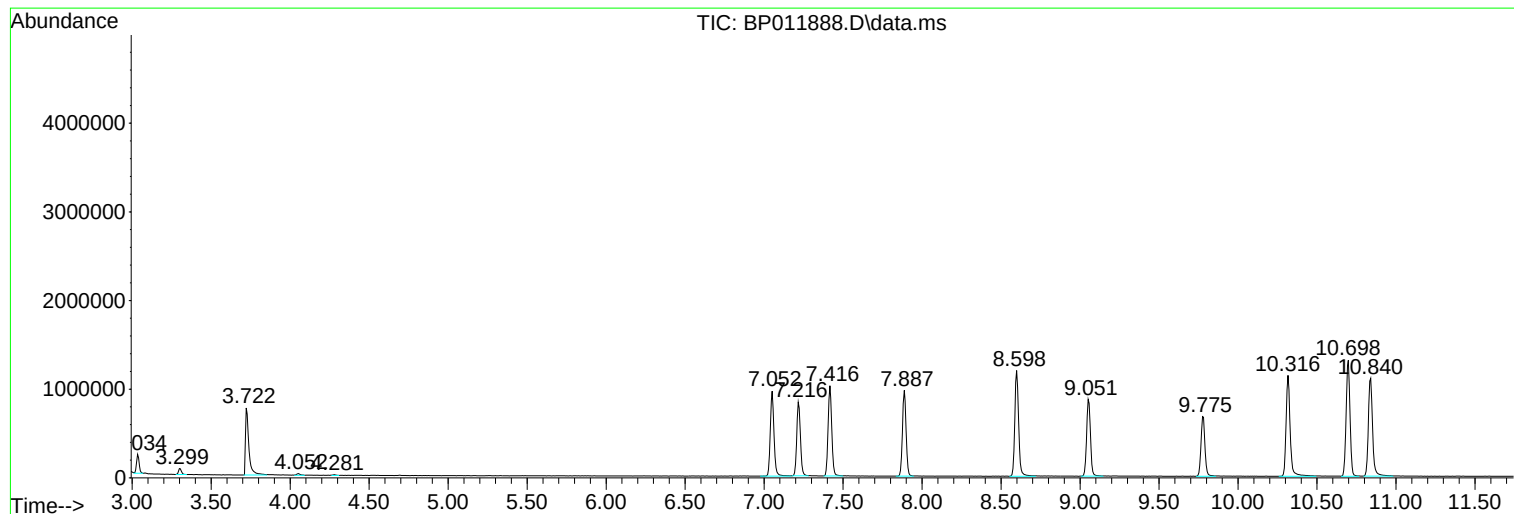
Sum of corrected areas: 66471957

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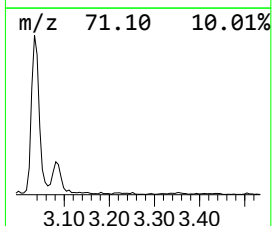
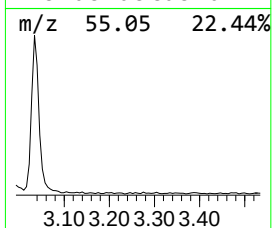
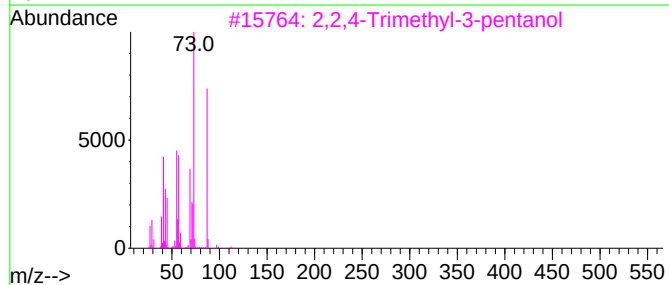
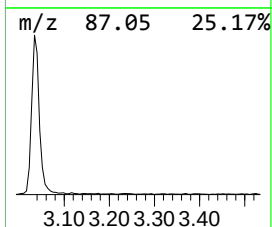
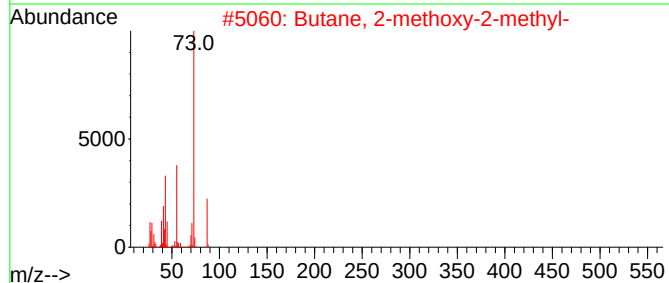
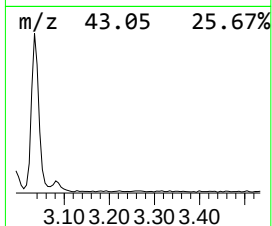
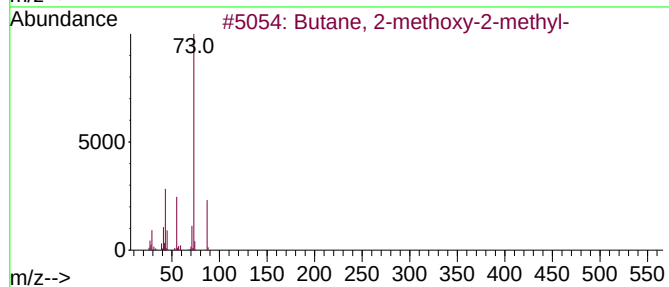
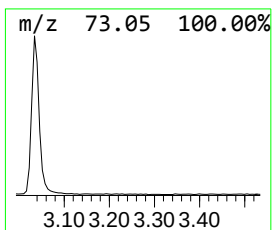
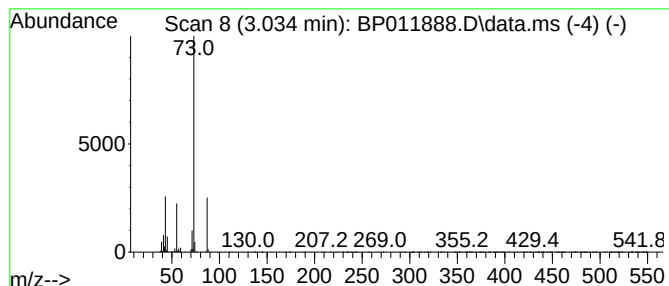
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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	3.12 ng/ul	239784	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	78
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	25
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	12



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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	3.1	ng/u1	239784	1	7.887	1539230	20.0