

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011941.D
 Acq On : 05 Oct 2022 21:57
 Operator : CG/JU
 Sample : PB148075BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK075

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: BP011941.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	19	rVB	195533	248061	4.30%	0.426%
2	3.299	49	53	65	rVB	62737	81529	1.41%	0.140%
3	3.722	122	125	143	rBV	663436	985373	17.08%	1.692%
4	4.046	176	180	186	rVB	17535	23843	0.41%	0.041%
5	7.046	683	690	702	rBV	798918	1326438	23.00%	2.278%
6	7.210	712	718	732	rVB	713663	1140394	19.77%	1.959%
7	7.410	744	752	762	rBV	878127	1411975	24.48%	2.425%
8	7.881	824	832	843	rBV	867775	1376848	23.87%	2.365%
9	8.593	944	953	968	rBV	1038463	1740968	30.19%	2.990%
10	9.046	1021	1030	1043	rBV	743463	1272348	22.06%	2.185%
11	9.769	1146	1153	1166	rBV	590089	1009845	17.51%	1.734%
12	10.310	1237	1245	1267	rBV	945341	1762911	30.57%	3.028%
13	10.687	1302	1309	1324	rVB	1187675	2050505	35.55%	3.522%
14	10.828	1324	1333	1349	rBV	945445	1789553	31.03%	3.074%
15	12.287	1574	1581	1588	rBV	14552	33644	0.58%	0.058%
16	13.422	1769	1774	1780	rBV5	9446	16906	0.29%	0.029%
17	13.504	1783	1788	1795	rVB8	3859	7910	0.14%	0.014%
18	13.939	1855	1862	1877	rBV	1781417	2534828	43.95%	4.354%
19	14.222	1903	1910	1925	rBV	2138637	3212660	55.70%	5.518%
20	14.528	1955	1962	1973	rBV	1899157	2761535	47.88%	4.743%
21	14.728	1990	1996	2029	rBV	542953	1196160	20.74%	2.054%
22	15.522	2124	2131	2144	rBV	2988324	4221119	73.19%	7.250%
23	15.639	2145	2151	2168	rVV	715751	1047484	18.16%	1.799%
24	15.763	2168	2172	2182	rVB2	14278	26876	0.47%	0.046%
25	17.069	2390	2394	2398	rBV4	5212	10693	0.19%	0.018%
26	17.286	2424	2431	2441	rBV	2255503	3308624	57.37%	5.683%
27	17.386	2441	2448	2466	rVV	3182435	4727091	81.96%	8.119%
28	19.198	2752	2756	2759	rBV2	35893	45274	0.78%	0.078%
29	19.263	2764	2767	2774	rVV	35107	50854	0.88%	0.087%
30	19.616	2822	2827	2839	rBV	4191976	5602908	97.15%	9.623%
31	21.368	3120	3125	3132	rBV	2922216	3677228	63.76%	6.316%
32	23.639	3505	3511	3527	rVB2	2775787	5767544	100.00%	9.906%
33	23.798	3532	3538	3551	rVB2	1832993	3753469	65.08%	6.447%

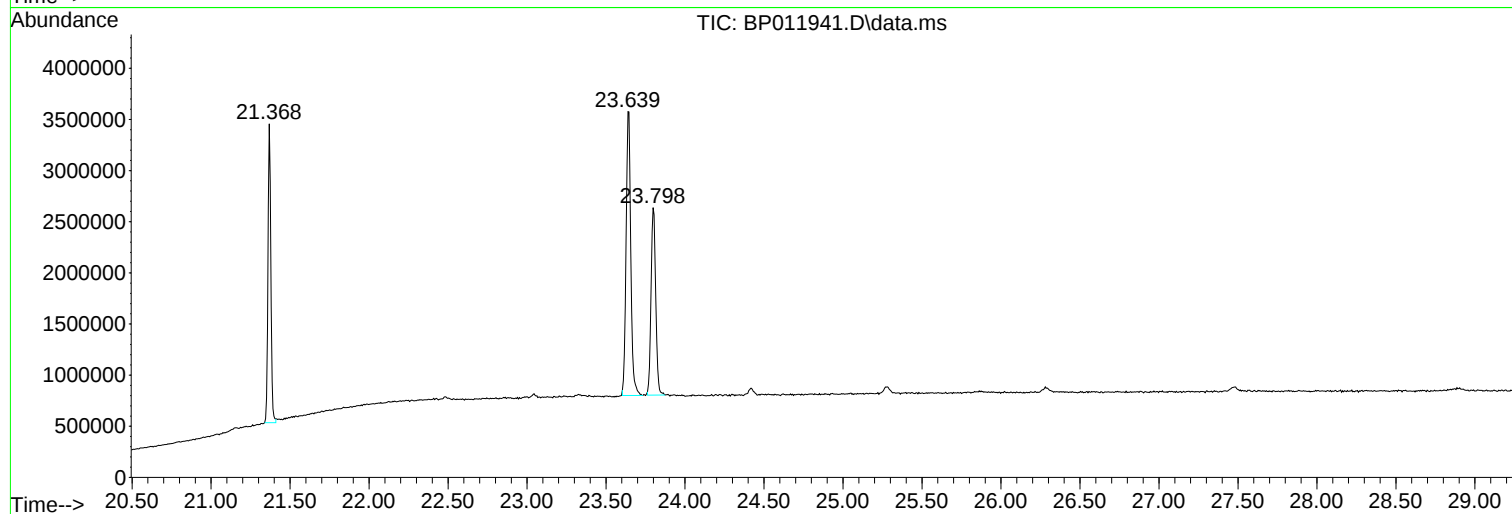
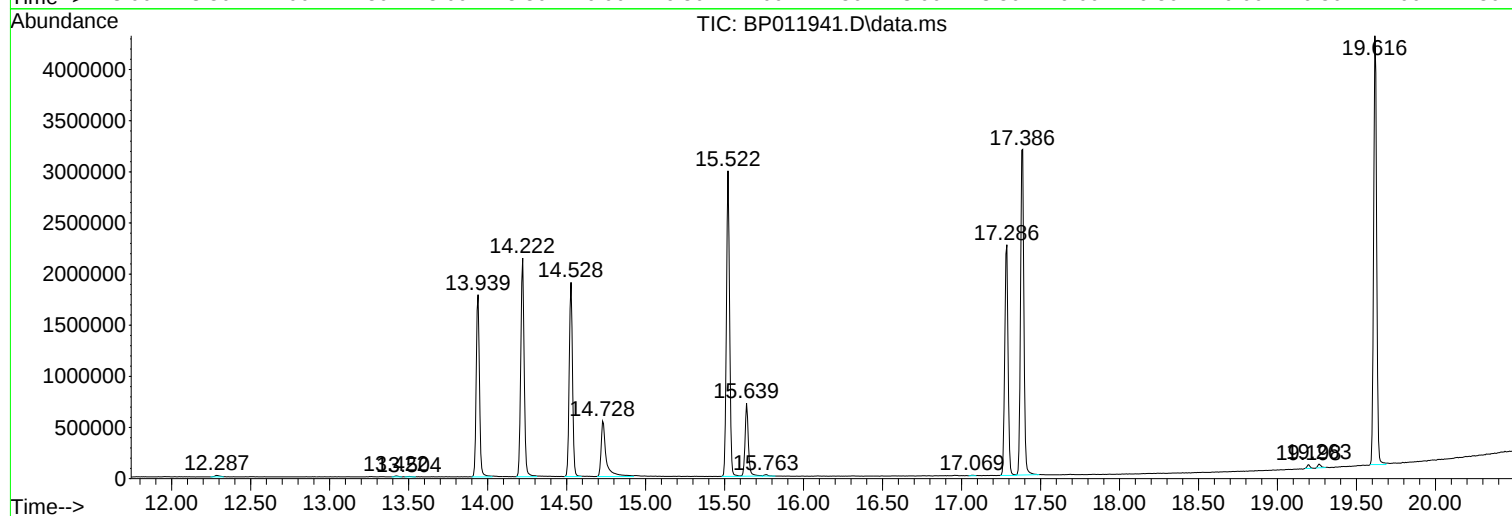
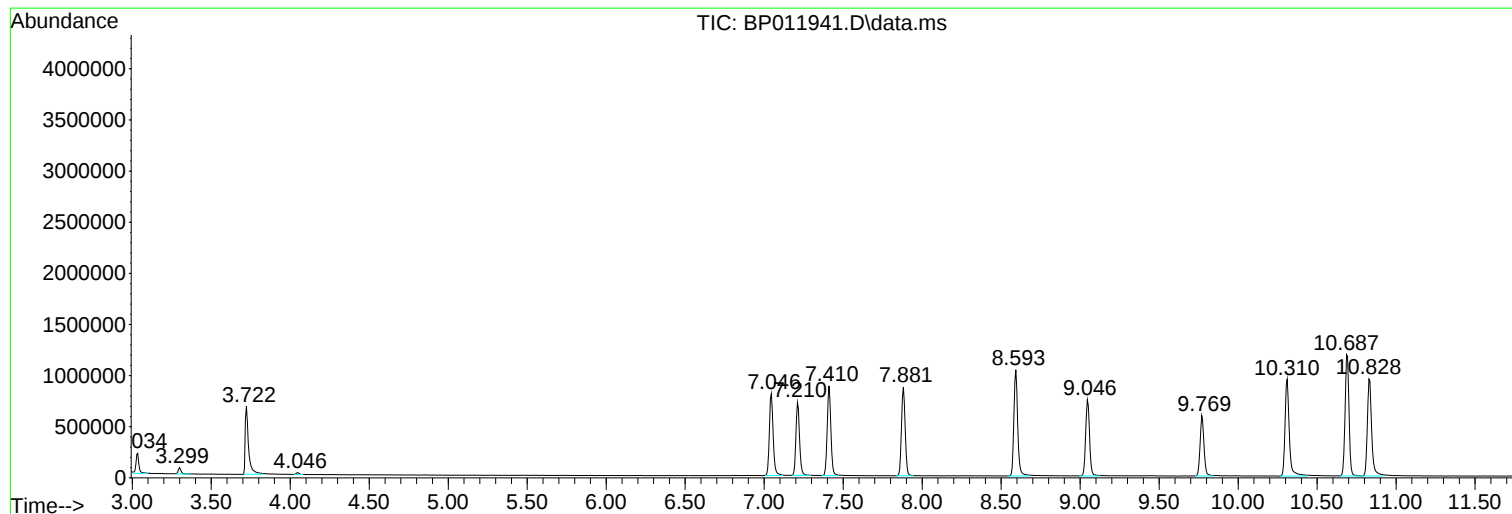
Sum of corrected areas: 58223398

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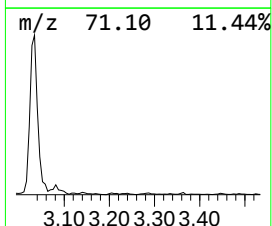
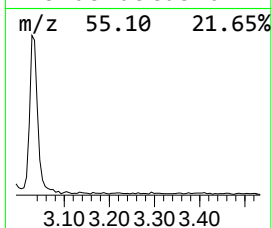
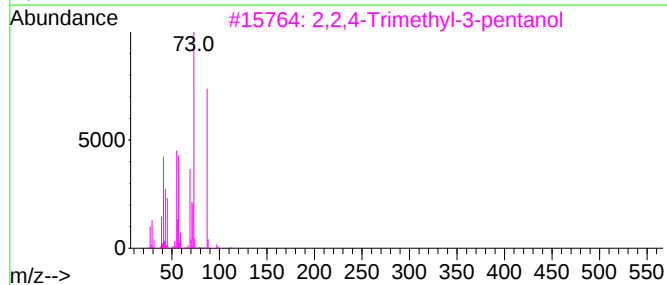
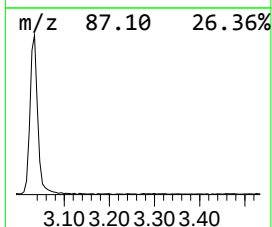
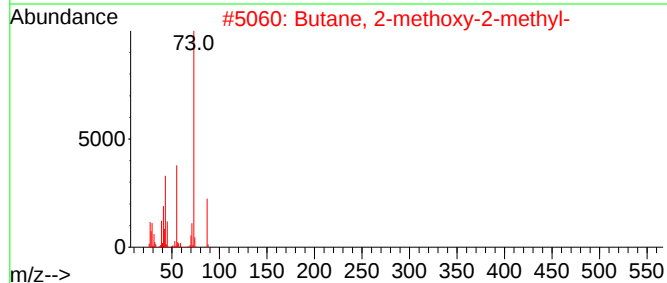
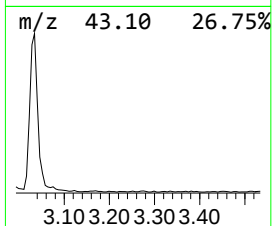
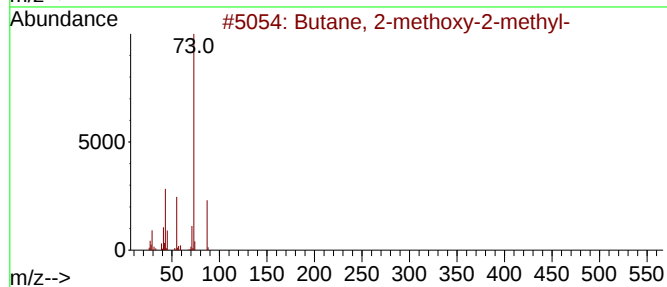
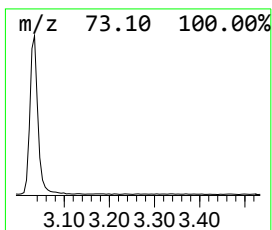
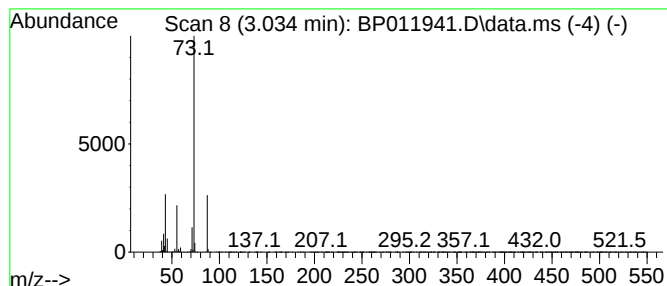
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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	3.60 ng/ul	248061	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	40
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	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/ul	248061	1	7.881	1376850	20.0