

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011942.D
 Acq On : 05 Oct 2022 22:31
 Operator : CG/JU
 Sample : PB148103BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK103

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: BP011942.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	3	8	18	rVB	253782	319402	4.78%	0.494%
2	3.299	49	53	59	rVB	75527	88400	1.32%	0.137%
3	3.722	122	125	144	rBV	789202	1210282	18.12%	1.870%
4	4.046	177	180	187	rVB2	15934	22555	0.34%	0.035%
5	7.046	682	690	706	rBV	977267	1614375	24.17%	2.494%
6	7.210	712	718	734	rVB	851678	1377489	20.63%	2.128%
7	7.410	745	752	764	rBV	1059942	1690506	25.31%	2.612%
8	7.881	823	832	844	rBV	803087	1298179	19.44%	2.006%
9	8.593	946	953	970	rBV	1253889	2090852	31.31%	3.231%
10	9.046	1023	1030	1046	rBV	899698	1522515	22.80%	2.352%
11	9.769	1146	1153	1170	rBV	718924	1216362	18.21%	1.879%
12	10.310	1236	1245	1265	rBV	1185053	2114823	31.67%	3.268%
13	10.687	1299	1309	1322	rBV	1107368	1930409	28.91%	2.983%
14	10.828	1326	1333	1353	rBV	1133163	2098884	31.43%	3.243%
15	12.281	1575	1580	1590	rBV2	17105	36236	0.54%	0.056%
16	13.422	1767	1774	1781	rBV	23787	40508	0.61%	0.063%
17	13.498	1781	1787	1795	rVB8	3830	9087	0.14%	0.014%
18	13.940	1855	1862	1876	rBV	2056273	2999293	44.91%	4.634%
19	14.222	1902	1910	1923	rBV	2564515	3841075	57.52%	5.935%
20	14.528	1955	1962	1975	rBV	1779954	2598066	38.90%	4.014%
21	14.728	1989	1996	2021	rBV	768688	1438408	21.54%	2.222%
22	14.939	2029	2032	2043	rVB	6992	15814	0.24%	0.024%
23	15.522	2124	2131	2144	rBV	3446209	4974548	74.49%	7.686%
24	15.639	2145	2151	2168	rVV	876817	1294827	19.39%	2.001%
25	15.763	2168	2172	2180	rVB3	16670	31512	0.47%	0.049%
26	17.286	2424	2431	2441	rBV	2084615	3093598	46.32%	4.780%
27	17.386	2441	2448	2470	rVV	3803096	5565092	83.33%	8.598%
28	19.198	2752	2756	2760	rBV3	44452	60305	0.90%	0.093%
29	19.263	2764	2767	2772	rVV	44868	61440	0.92%	0.095%
30	19.616	2822	2827	2838	rBV	4807760	6580739	98.54%	10.168%
31	21.369	3120	3125	3131	rBV	2674595	3363318	50.36%	5.197%
32	23.645	3505	3512	3524	rVB2	3355453	6678194	100.00%	10.318%
33	23.798	3532	3538	3551	rVB2	1669077	3444674	51.58%	5.322%

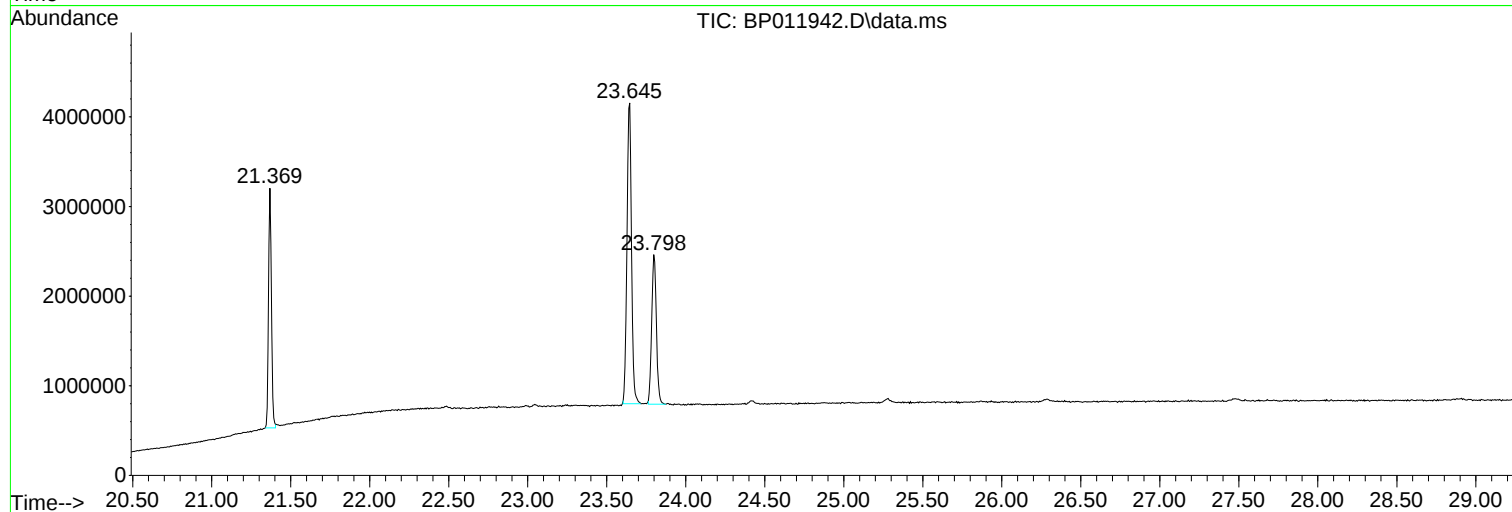
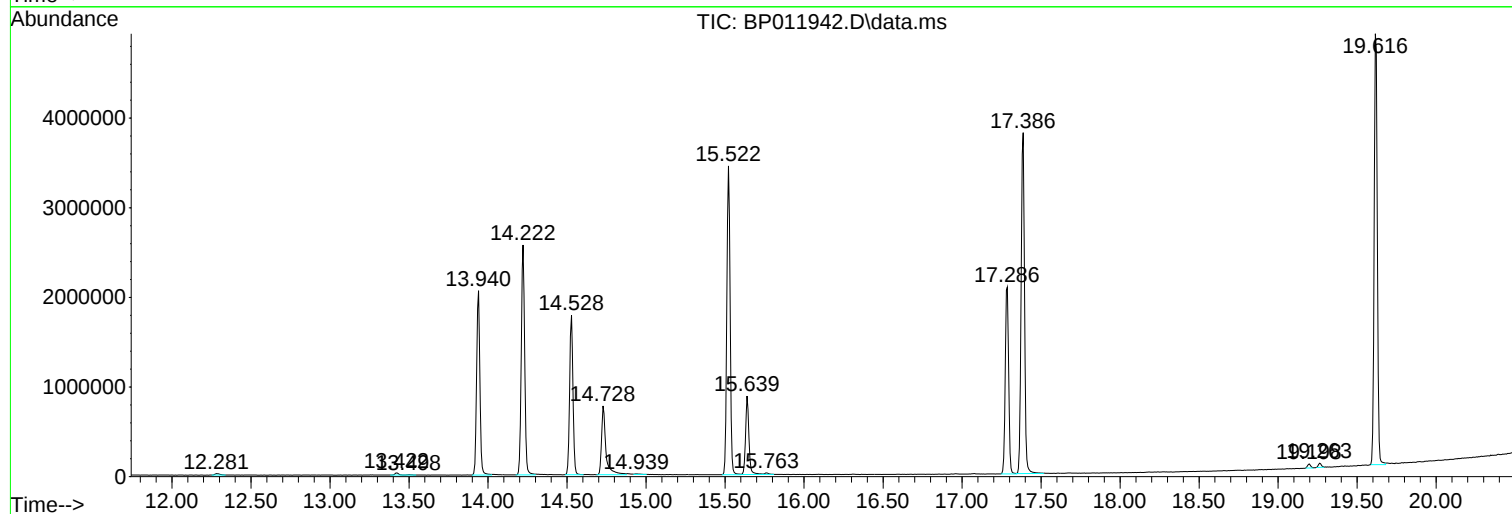
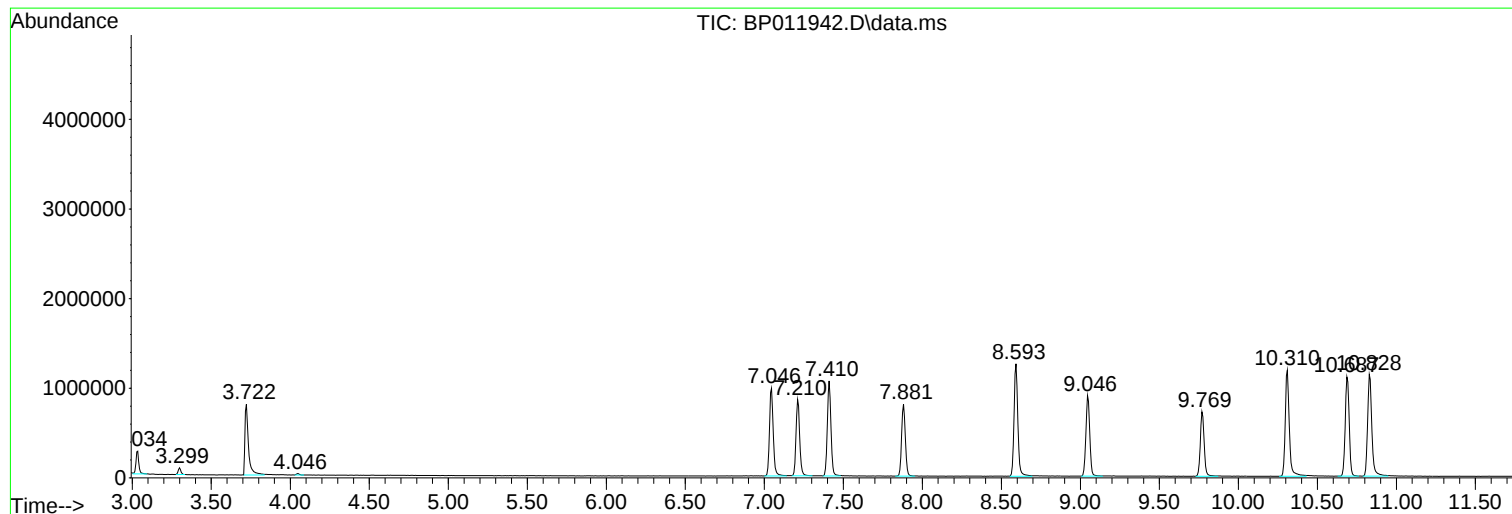
Sum of corrected areas: 64721767

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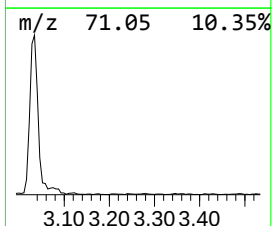
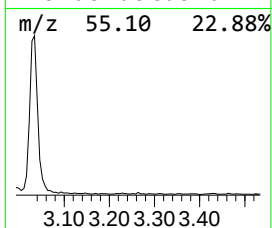
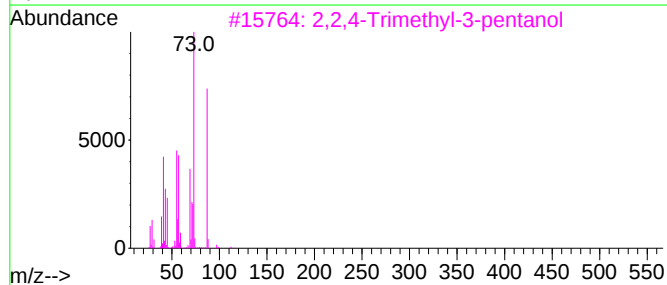
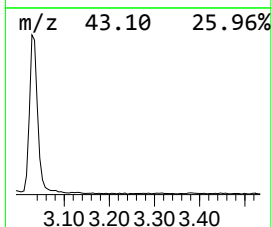
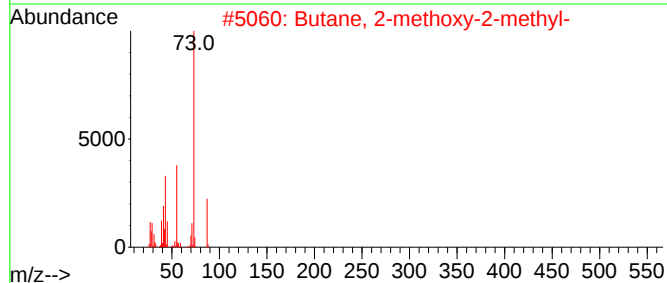
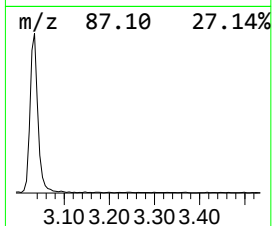
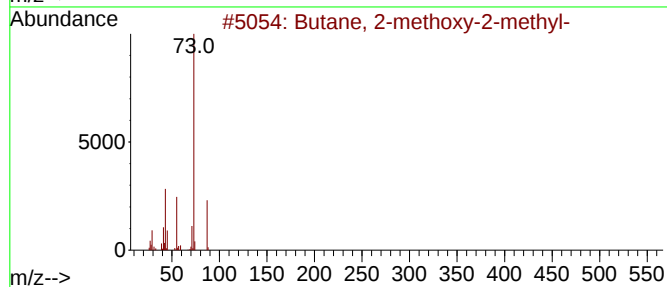
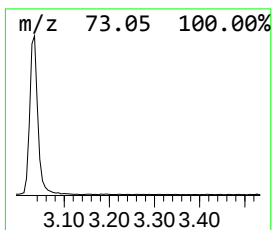
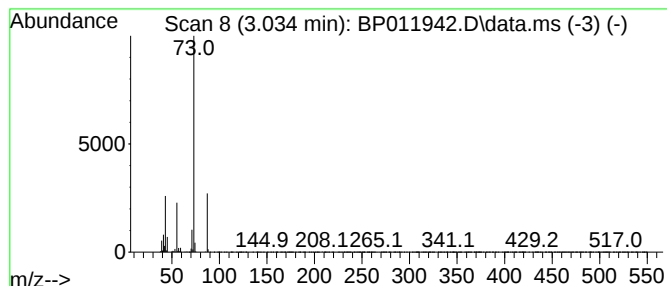
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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	4.92 ng/ul	319402	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	78		
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	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23		
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	4.9	ng/u1	319402	1	7.881	1298180	20.0