

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011949.D
 Acq On : 06 Oct 2022 02:31
 Operator : CG/JU
 Sample : PB148042BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK042

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: BP011949.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.028	3	7	17	rVB	267808	329245	4.92%	0.501%
2	3.299	49	53	62	rVB	76579	97084	1.45%	0.148%
3	3.717	121	124	145	rBV	802910	1284452	19.21%	1.953%
4	4.046	176	180	186	rVB	14223	21676	0.32%	0.033%
5	7.040	683	689	704	rBV	1003354	1639801	24.52%	2.493%
6	7.211	711	718	730	rBV	877970	1409608	21.08%	2.143%
7	7.411	744	752	768	rBV	1081550	1761234	26.33%	2.678%
8	7.881	825	832	847	rBV	761433	1228863	18.37%	1.868%
9	8.593	945	953	968	rBV	1260709	2175771	32.53%	3.308%
10	9.046	1022	1030	1044	rBV	964585	1603452	23.98%	2.438%
11	9.769	1144	1153	1163	rBV	758489	1265959	18.93%	1.925%
12	10.310	1236	1245	1266	rBV	1184046	2212152	33.08%	3.363%
13	10.687	1302	1309	1326	rVB	1102392	1856931	27.77%	2.823%
14	10.828	1326	1333	1358	rBV	1244035	2303153	34.44%	3.502%
15	12.287	1572	1581	1588	rBV	19637	43231	0.65%	0.066%
16	12.499	1609	1617	1624	rBV	2982	6941	0.10%	0.011%
17	13.422	1767	1774	1780	rBV2	23026	38913	0.58%	0.059%
18	13.940	1855	1862	1873	rBV	2151401	3157335	47.21%	4.800%
19	14.222	1901	1910	1924	rBV	2701689	4089069	61.14%	6.217%
20	14.528	1955	1962	1975	rBV	1711210	2512419	37.57%	3.820%
21	14.728	1989	1996	2019	rBV	733497	1463065	21.88%	2.224%
22	14.940	2030	2032	2037	rVB6	6228	8563	0.13%	0.013%
23	15.522	2124	2131	2144	rBV	3665364	5300649	79.26%	8.059%
24	15.640	2144	2151	2166	rVV	881145	1269176	18.98%	1.930%
25	15.763	2168	2172	2179	rVB3	17519	30384	0.45%	0.046%
26	16.963	2371	2376	2384	rBV8	5702	11612	0.17%	0.018%
27	17.281	2424	2430	2440	rBV	2062165	2960934	44.27%	4.502%
28	17.381	2441	2447	2464	rVV	3910751	5719811	85.53%	8.696%
29	19.192	2752	2755	2760	rBV2	43117	56444	0.84%	0.086%
30	19.263	2763	2767	2771	rBV2	45998	62672	0.94%	0.095%
31	19.616	2821	2827	2836	rBV	5142394	6687855	100.00%	10.168%
32	21.369	3120	3125	3134	rBV	2542664	3277410	49.01%	4.983%
33	23.639	3504	3511	3525	rVB2	3416288	6681590	99.91%	10.158%
34	23.798	3532	3538	3547	rVB	1572585	3207418	47.96%	4.876%

Sum of corrected areas: 65774872

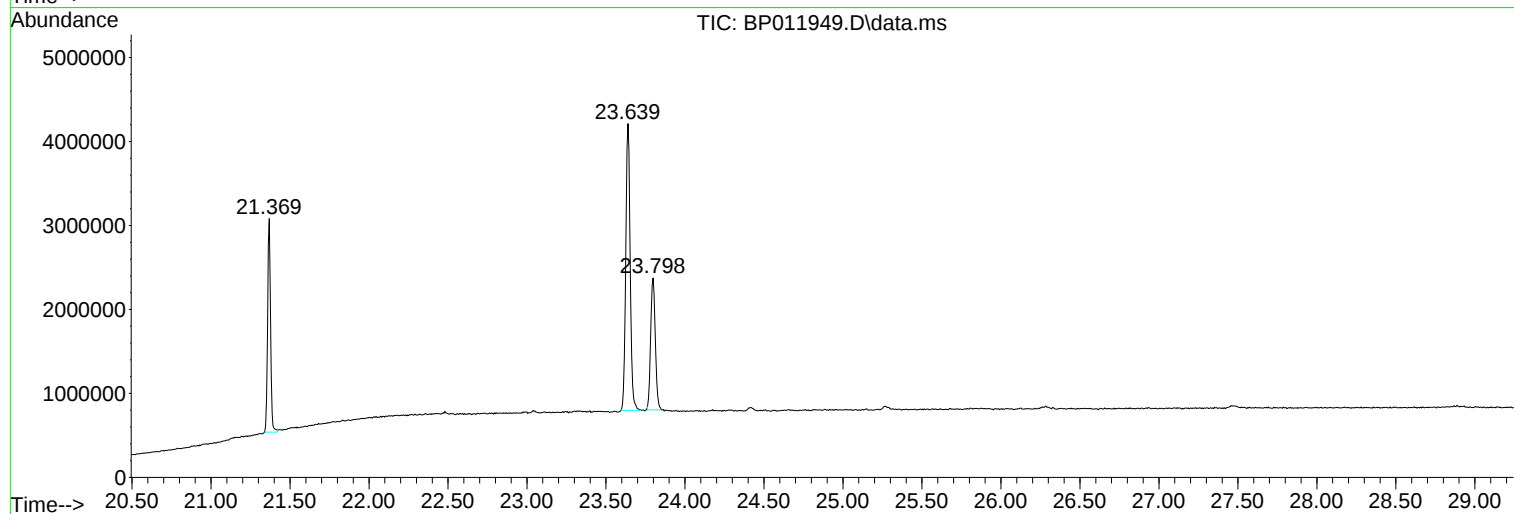
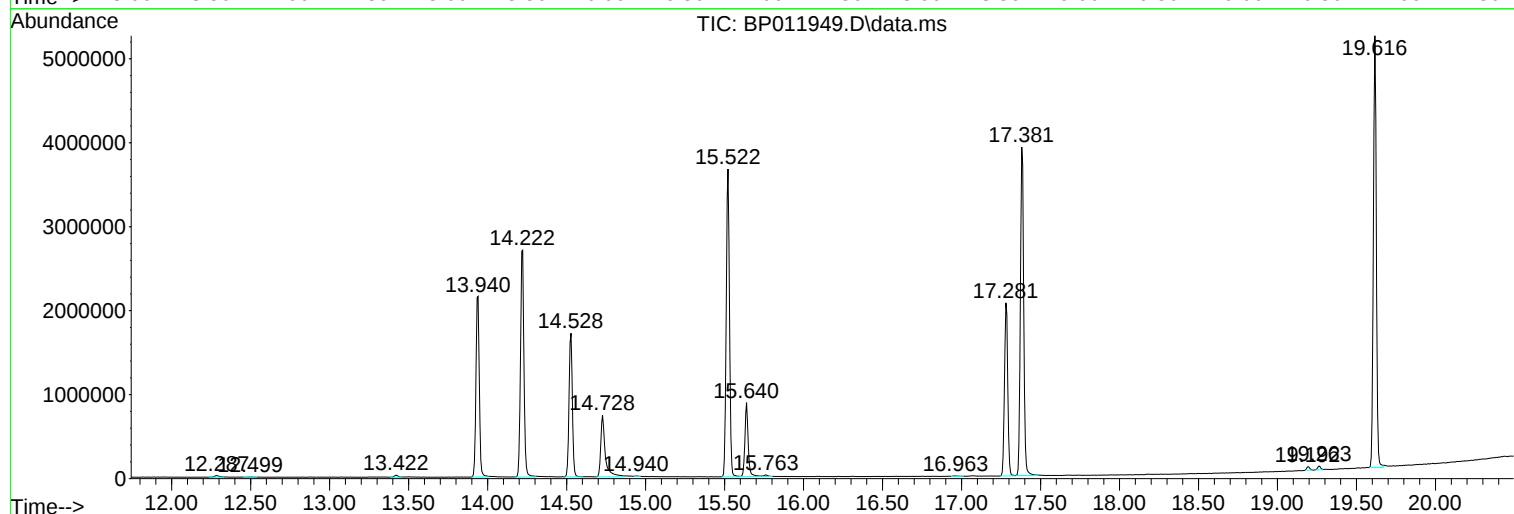
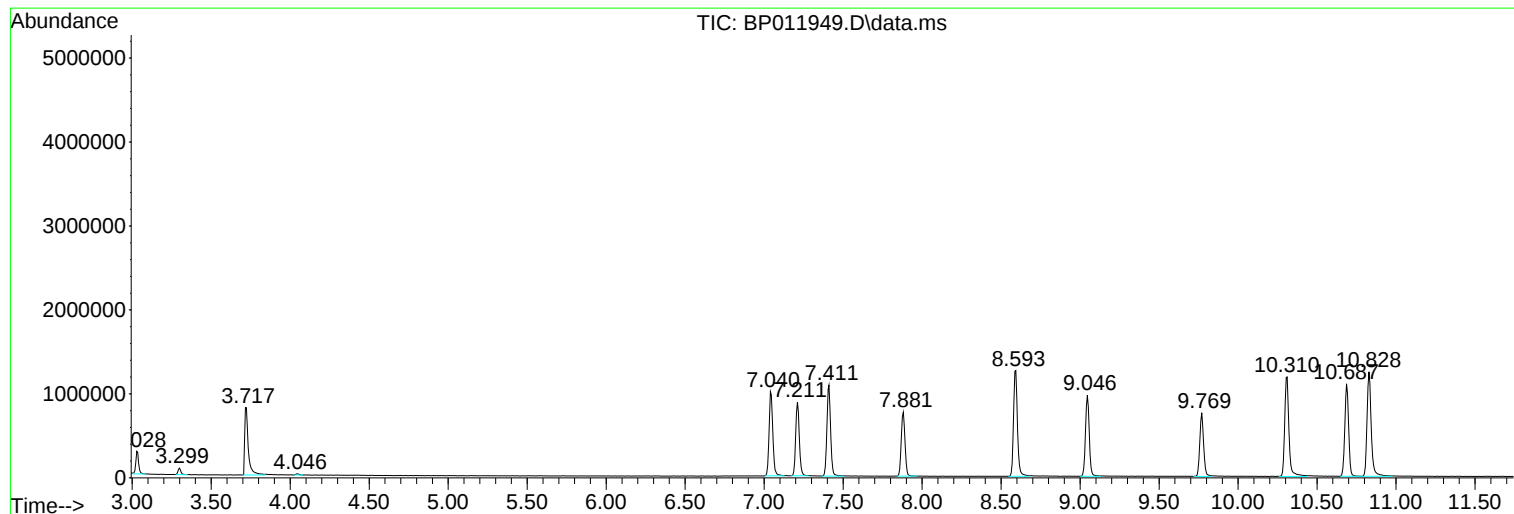
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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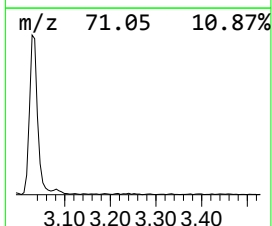
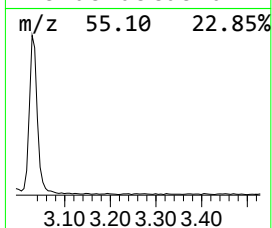
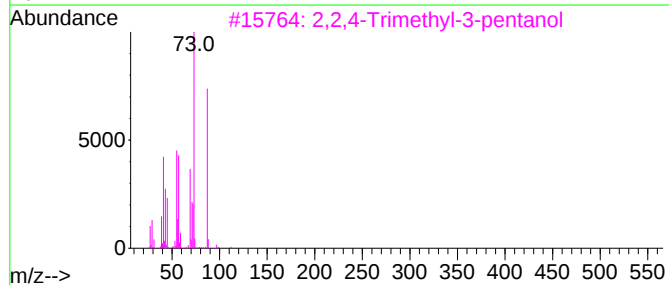
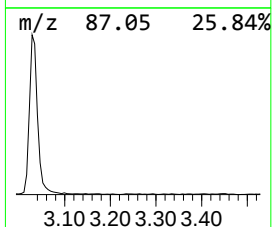
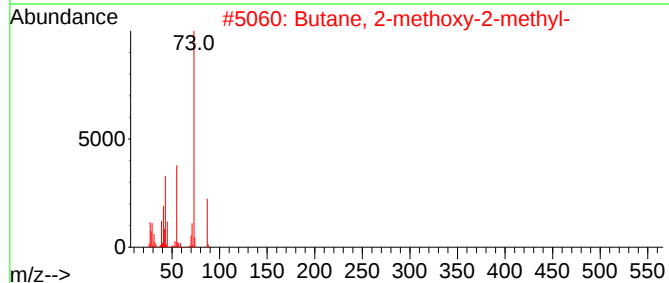
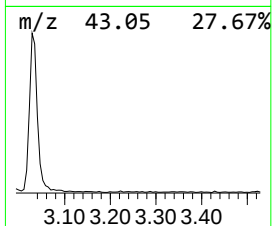
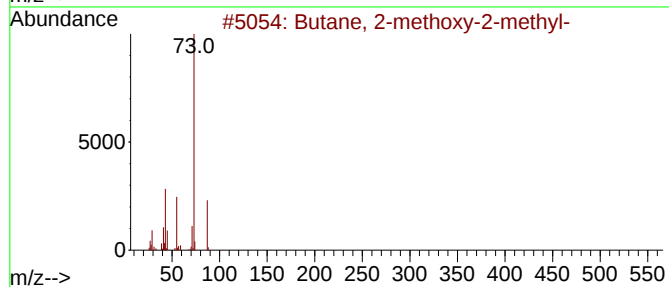
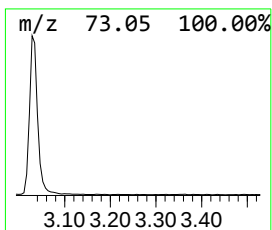
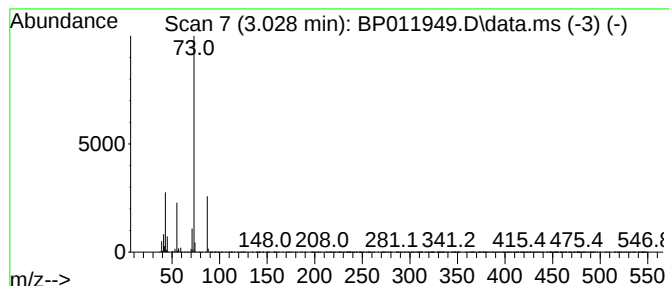
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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.
3.028	5.36 ng/ul	329245	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	72		
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	17		
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...	3.028	5.4	ng/u1	329245	1	7.881	1228860	20.0