

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011956.D
 Acq On : 06 Oct 2022 06:30
 Operator : CG/JU
 Sample : N4956-04
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB8S1

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: BP011956.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	29	rVB	5525260	7146931	100.00%	10.621%
2	3.299	49	53	58	rVB	50522	61897	0.87%	0.092%
3	3.728	123	126	139	rBV	190845	309092	4.32%	0.459%
4	3.946	159	163	168	rVB	30389	46211	0.65%	0.069%
5	4.046	176	180	186	rVB	18835	27619	0.39%	0.041%
6	4.269	215	218	223	rVB5	6493	9621	0.13%	0.014%
7	7.040	681	689	703	rVB	171617	303907	4.25%	0.452%
8	7.211	711	718	732	rBV	721252	1186228	16.60%	1.763%
9	7.405	744	751	762	rBV	669123	1129468	15.80%	1.679%
10	7.675	791	797	804	rBV	3811	7725	0.11%	0.011%
11	7.875	822	831	840	rBV	843708	1413406	19.78%	2.101%
12	7.946	840	843	852	rVB3	10004	19141	0.27%	0.028%
13	8.587	945	952	972	rBV	531607	938243	13.13%	1.394%
14	9.046	1021	1030	1043	rBV	781414	1342680	18.79%	1.995%
15	9.452	1094	1099	1104	rBV5	6493	8781	0.12%	0.013%
16	9.769	1146	1153	1162	rBV	559444	974225	13.63%	1.448%
17	10.304	1237	1244	1260	rBV	919109	1648068	23.06%	2.449%
18	10.687	1301	1309	1322	rBV	1231017	2097726	29.35%	3.118%
19	10.828	1326	1333	1359	rBV	742601	1446424	20.24%	2.150%
20	12.275	1573	1579	1589	rBV2	17158	36963	0.52%	0.055%
21	13.422	1767	1774	1779	rVB4	12722	20468	0.29%	0.030%
22	13.498	1781	1787	1791	rVB7	4953	9137	0.13%	0.014%
23	13.934	1854	1861	1878	rBV	1952435	2868810	40.14%	4.263%
24	14.216	1902	1909	1929	rBV	2256000	3484405	48.75%	5.178%
25	14.522	1954	1961	1968	rBV	1907863	2858985	40.00%	4.249%
26	14.575	1968	1970	1976	rVB3	14094	25719	0.36%	0.038%
27	14.734	1991	1997	2011	rBV2	84385	237095	3.32%	0.352%
28	14.951	2029	2034	2040	rVB10	10713	21052	0.29%	0.031%
29	15.345	2098	2101	2111	rVB	27440	38495	0.54%	0.057%
30	15.522	2124	2131	2137	rBV	3162465	4736270	66.27%	7.039%
31	15.639	2145	2151	2165	rVV	880463	1301951	18.22%	1.935%
32	15.757	2167	2171	2179	rVB3	16156	32857	0.46%	0.049%
33	17.281	2423	2430	2441	rBV	2412602	3465362	48.49%	5.150%
34	17.381	2441	2447	2467	rVB	3946028	5677905	79.45%	8.438%
35	18.163	2575	2580	2586	rBV	199383	291893	4.08%	0.434%
36	19.192	2751	2755	2761	rBV	52887	72705	1.02%	0.108%

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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

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37	19.263	2763	2767	2770	rBV	55832	83138	1.16%	0.124%
38	19.398	2786	2790	2797	rBV2	137549	213585	2.99%	0.317%
39	19.616	2821	2827	2844	rBV	5229427	7047821	98.61%	10.474%
40	21.369	3120	3125	3136	rBV	2986713	3861258	54.03%	5.738%
41	23.639	3504	3511	3524	rVB2	3488813	6989435	97.80%	10.387%
42	23.798	3532	3538	3550	rVB2	1868108	3795901	53.11%	5.641%

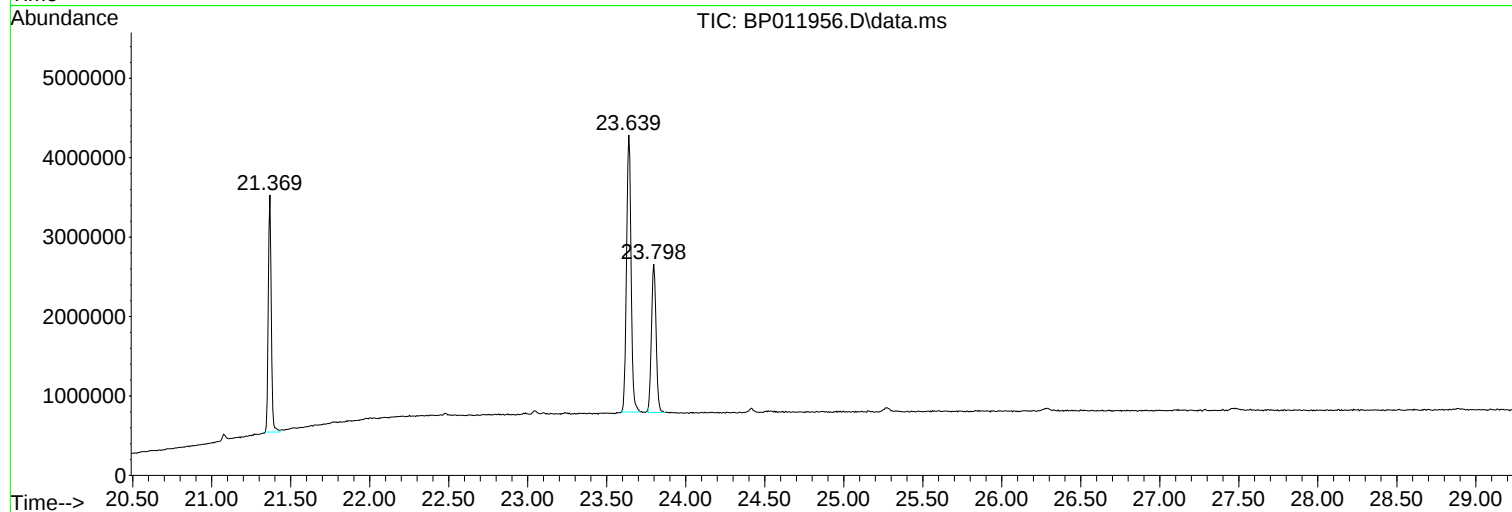
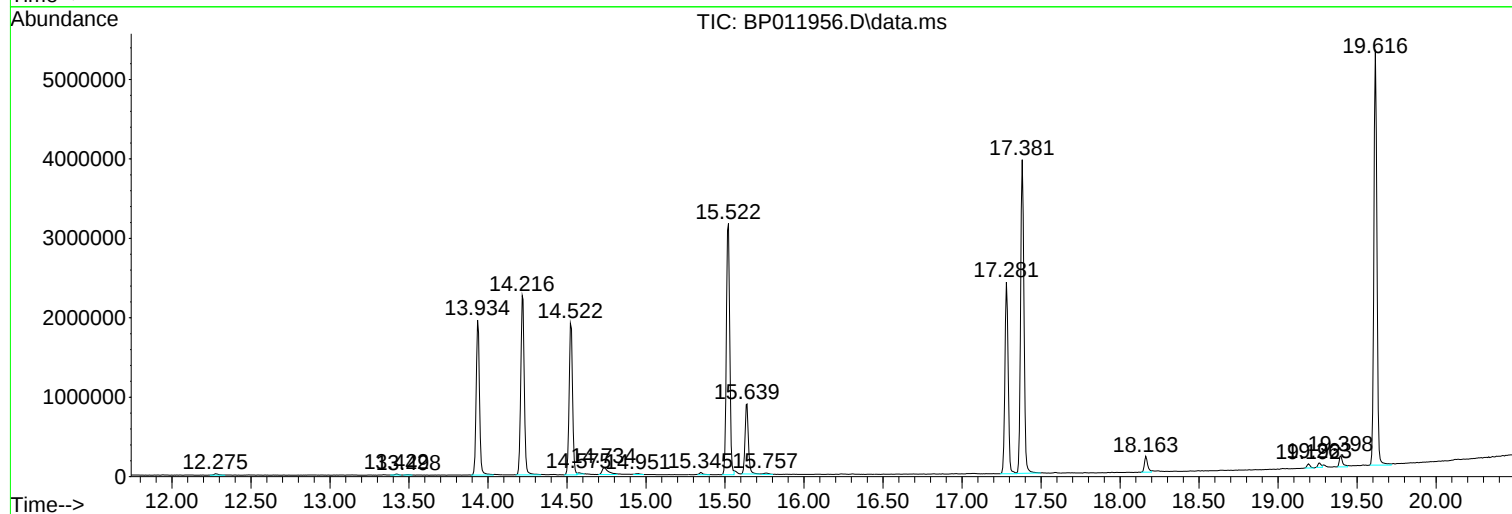
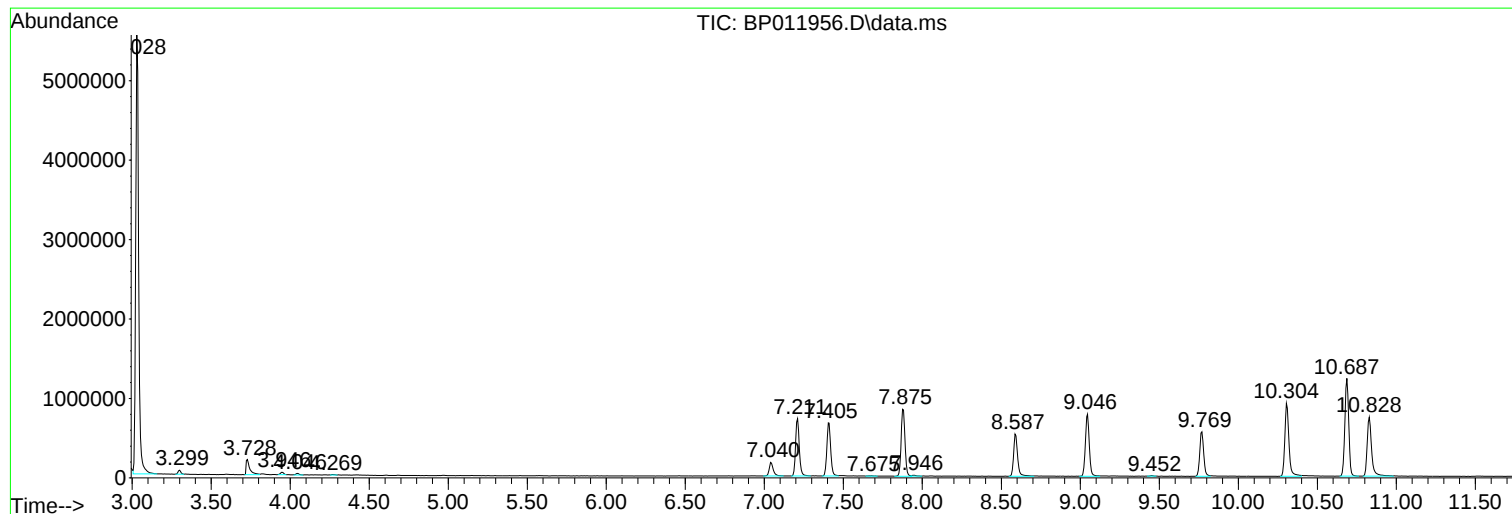
Sum of corrected areas: 67288603

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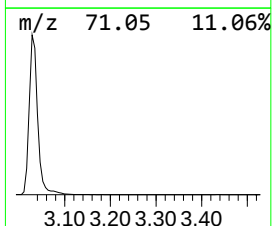
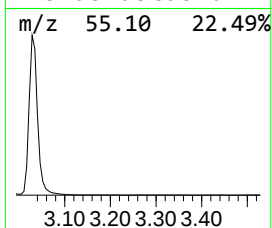
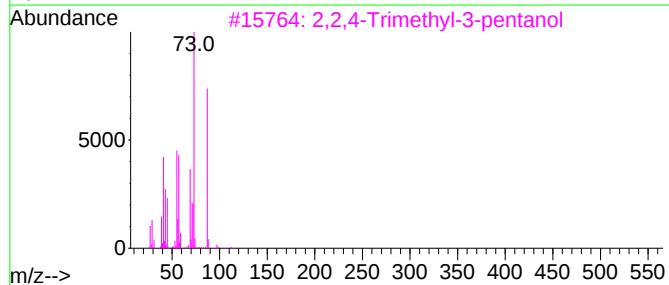
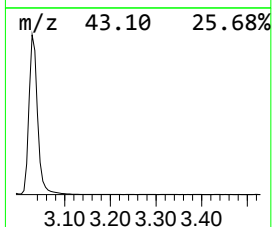
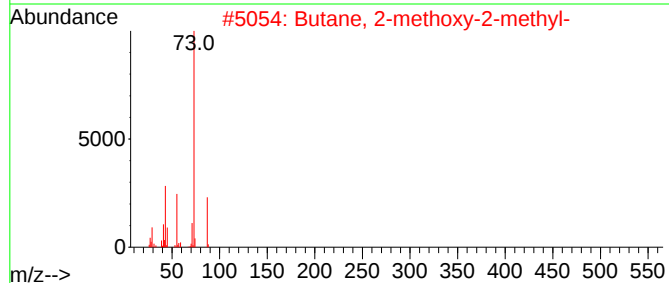
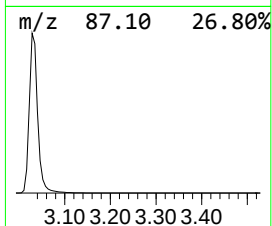
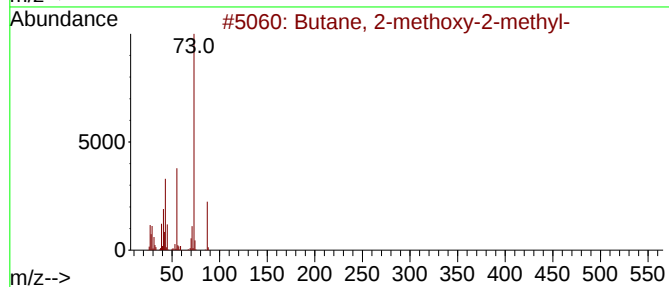
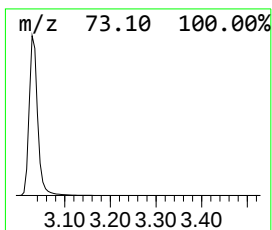
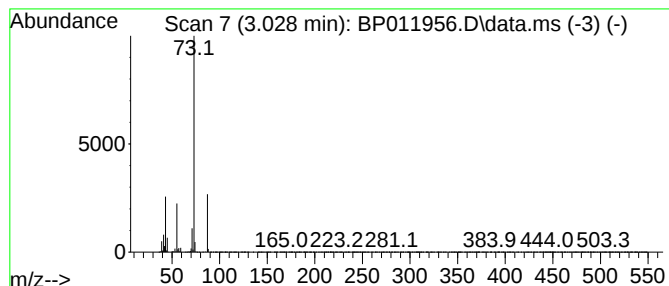
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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	101.13 ng/ul	7146930	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	72		
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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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.028	101.1	ng/ul	7146930	1	7.881	1413410	20.0