

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100122\
 Data File : BP011853.D
 Acq On : 01 Oct 2022 10:10
 Operator : CG/JU
 Sample : PB147968BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK968

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: BP011853.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.040	5	9	18	rVB	137028	199384	2.13%	0.272%
2	3.305	50	54	63	rVB	49315	71098	0.76%	0.097%
3	3.728	123	126	144	rBV	555395	927481	9.91%	1.267%
4	4.058	177	182	190	rVB	12623	21125	0.23%	0.029%
5	7.052	684	691	704	rBV	810688	1329450	14.21%	1.817%
6	7.222	713	720	739	rVB	761180	1243224	13.29%	1.699%
7	7.422	746	754	767	rBV	860208	1424212	15.22%	1.946%
8	7.893	826	834	844	rBV	708438	1151938	12.31%	1.574%
9	8.599	945	954	974	rBV	1117080	1910883	20.43%	2.611%
10	9.058	1024	1032	1045	rBV	844601	1429859	15.28%	1.954%
11	9.781	1147	1155	1167	rBV	681263	1171710	12.52%	1.601%
12	10.322	1238	1247	1268	rBV	1060828	1931648	20.65%	2.640%
13	10.705	1304	1312	1322	rBV	1088269	1873565	20.03%	2.560%
14	10.840	1328	1335	1358	rBV	1107646	2072075	22.15%	2.832%
15	12.299	1576	1583	1591	rBV3	19618	43462	0.46%	0.059%
16	13.434	1771	1776	1782	rBV9	5979	10975	0.12%	0.015%
17	13.516	1782	1790	1795	rBV4	4571	9890	0.11%	0.014%
18	13.951	1855	1864	1881	rBV	2681955	3735922	39.93%	5.105%
19	14.234	1905	1912	1925	rBV	2827551	4242083	45.34%	5.797%
20	14.540	1957	1964	1973	rBV	1946301	2804311	29.98%	3.832%
21	14.734	1991	1997	2015	rBV	1074121	1693066	18.10%	2.314%
22	15.534	2126	2133	2142	rBV	3865037	5681088	60.73%	7.763%
23	15.651	2147	2153	2169	rVV	1182863	1704467	18.22%	2.329%
24	15.775	2170	2174	2183	rVB3	34478	57894	0.62%	0.079%
25	16.322	2264	2267	2275	rVB6	11168	18490	0.20%	0.025%
26	16.975	2374	2378	2382	rBV6	8991	14664	0.16%	0.020%
27	17.298	2426	2433	2442	rBV	2466588	3506757	37.48%	4.792%
28	17.398	2443	2450	2470	rVV2	4480730	6600909	70.56%	9.020%
29	19.210	2754	2758	2763	rBV	55668	71006	0.76%	0.097%
30	19.275	2765	2769	2776	rBV	55659	87006	0.93%	0.119%
31	19.628	2823	2829	2840	rBV	5861758	7924691	84.71%	10.829%
32	21.380	3122	3127	3136	rBV	3191198	4189764	44.78%	5.725%
33	23.663	3508	3515	3532	rVB2	4569895	9355398	100.00%	12.784%
34	23.821	3535	3542	3557	rVB2	2310381	4669231	49.91%	6.381%

Sum of corrected areas: 73178726

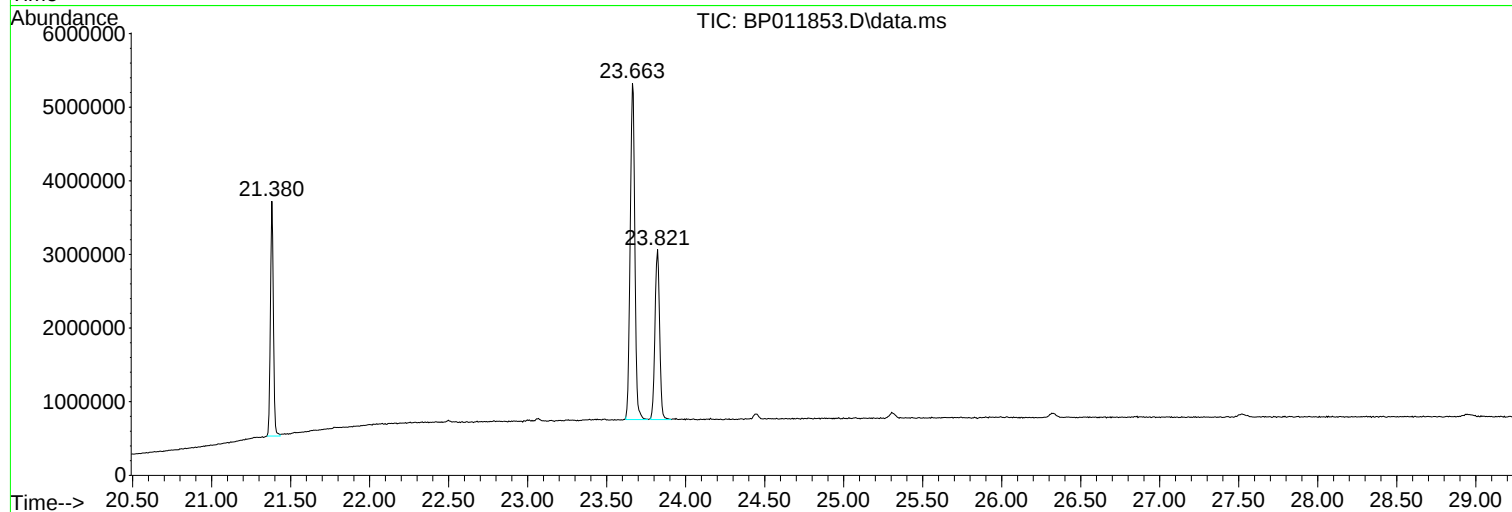
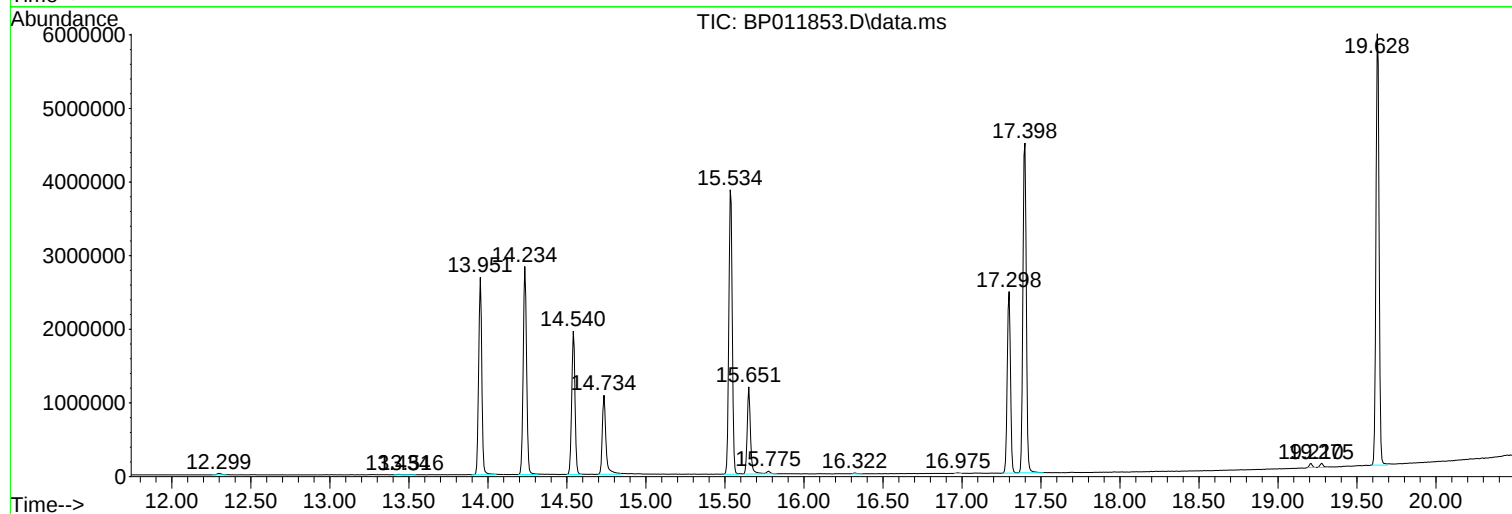
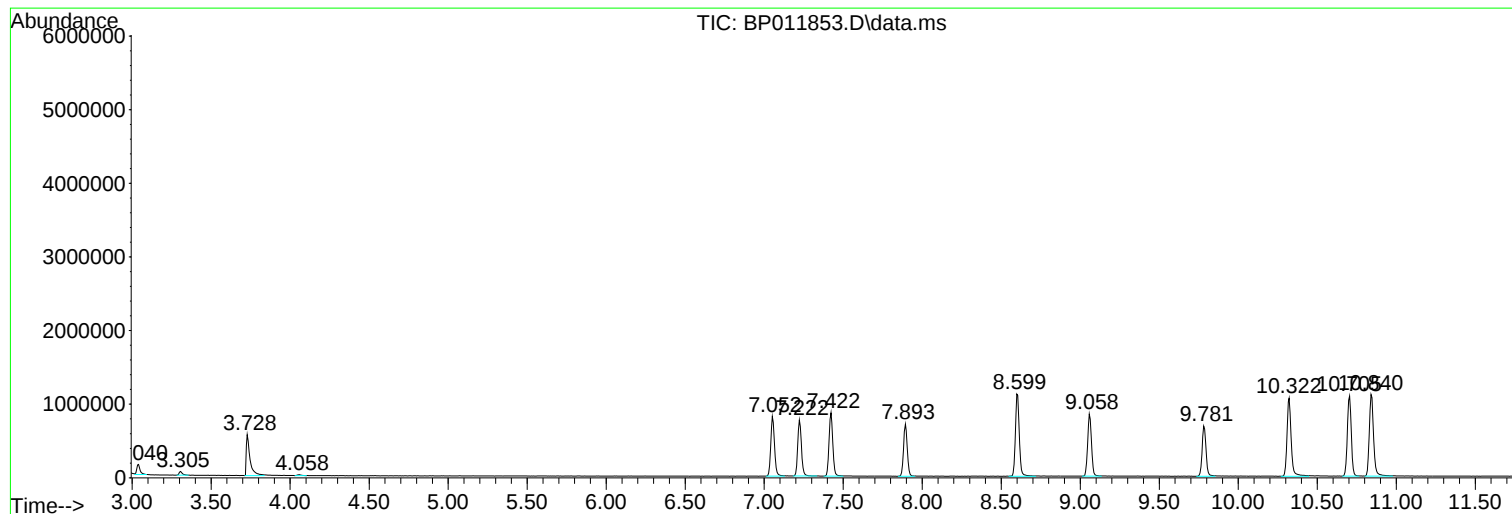
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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



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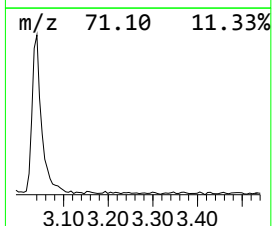
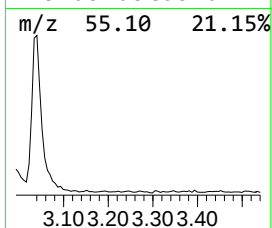
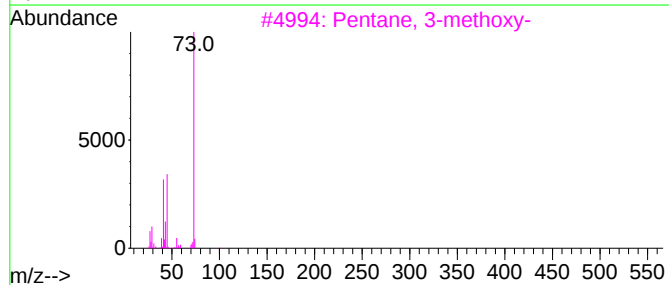
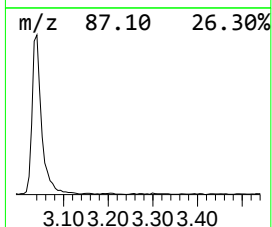
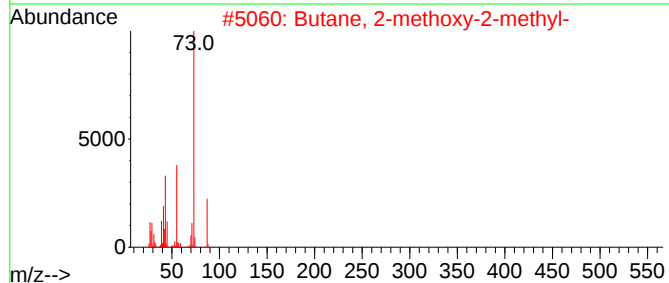
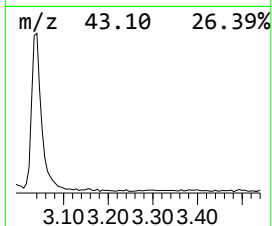
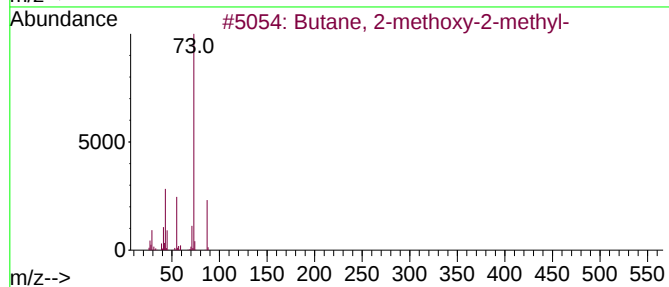
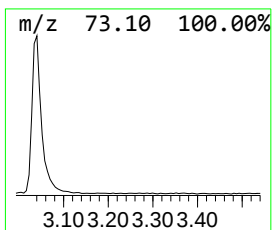
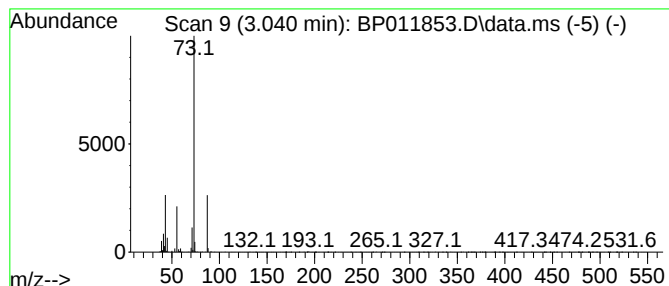
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.040	3.46 ng/ul	199384	1,4-Dichlorobenzene-d4	7.893

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	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12		
4	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9		
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.040	3.5	ng/u1	199384	1	7.893	1151940	20.0