

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011989.D
 Acq On : 07 Oct 2022 02:37
 Operator : CG/JU
 Sample : PB148077BL
 Misc :
 ALS Vial : 78 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK077

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: BP011989.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	20	rVB	214710	276634	4.24%	0.427%
2	3.299	49	53	58	rVB	67092	87361	1.34%	0.135%
3	3.716	122	124	142	rBV	705215	1098493	16.85%	1.696%
4	4.040	176	179	187	rVB	18200	25779	0.40%	0.040%
5	7.040	682	689	701	rBV	944330	1532883	23.51%	2.366%
6	7.205	711	717	737	rVB	819412	1335576	20.49%	2.062%
7	7.405	744	751	764	rBV	1013621	1597703	24.51%	2.466%
8	7.875	824	831	844	rBV	846136	1395528	21.41%	2.154%
9	8.587	944	952	969	rBV	1167184	2001746	30.71%	3.090%
10	9.040	1020	1029	1044	rVV	899725	1513604	23.22%	2.337%
11	9.763	1144	1152	1170	rBV	707108	1187181	18.21%	1.833%
12	10.304	1236	1244	1266	rBV	1092336	2002180	30.71%	3.091%
13	10.681	1300	1308	1320	rBV	1247614	2101680	32.24%	3.244%
14	10.822	1325	1332	1354	rBV	1055381	1990339	30.53%	3.073%
15	12.281	1572	1580	1585	rBV3	17271	36349	0.56%	0.056%
16	13.416	1767	1773	1782	rBV3	17245	32849	0.50%	0.051%
17	13.492	1782	1786	1792	rVB7	4391	8283	0.13%	0.013%
18	13.934	1854	1861	1877	rBV	2039425	2940063	45.10%	4.539%
19	14.216	1901	1909	1923	rBV	2436680	3759664	57.67%	5.804%
20	14.522	1952	1961	1975	rBV	1925404	2888564	44.31%	4.459%
21	14.722	1989	1995	2018	rBV	675188	1373571	21.07%	2.120%
22	14.934	2029	2031	2040	rVB10	6687	14425	0.22%	0.022%
23	15.516	2121	2130	2142	rBV	3387348	4860747	74.56%	7.504%
24	15.633	2144	2150	2166	rVV	832801	1189179	18.24%	1.836%
25	15.757	2166	2171	2177	rVB3	16418	30371	0.47%	0.047%
26	16.957	2372	2375	2380	rVB7	7130	9962	0.15%	0.015%
27	17.280	2423	2430	2440	rBV	2338111	3482960	53.43%	5.377%
28	17.375	2440	2446	2462	rVV	3759897	5408655	82.97%	8.350%
29	19.192	2751	2755	2761	rBV4	43153	59488	0.91%	0.092%
30	19.257	2763	2766	2770	rBV	44089	57228	0.88%	0.088%
31	19.616	2821	2827	2840	rBV	4595721	6321494	96.97%	9.759%
32	21.363	3119	3124	3134	rBV	2845366	3843472	58.96%	5.933%
33	23.633	3503	3510	3527	rVB2	3230438	6518871	100.00%	10.063%
34	23.792	3531	3537	3551	rVB2	1875689	3795276	58.22%	5.859%

Sum of corrected areas: 64778158

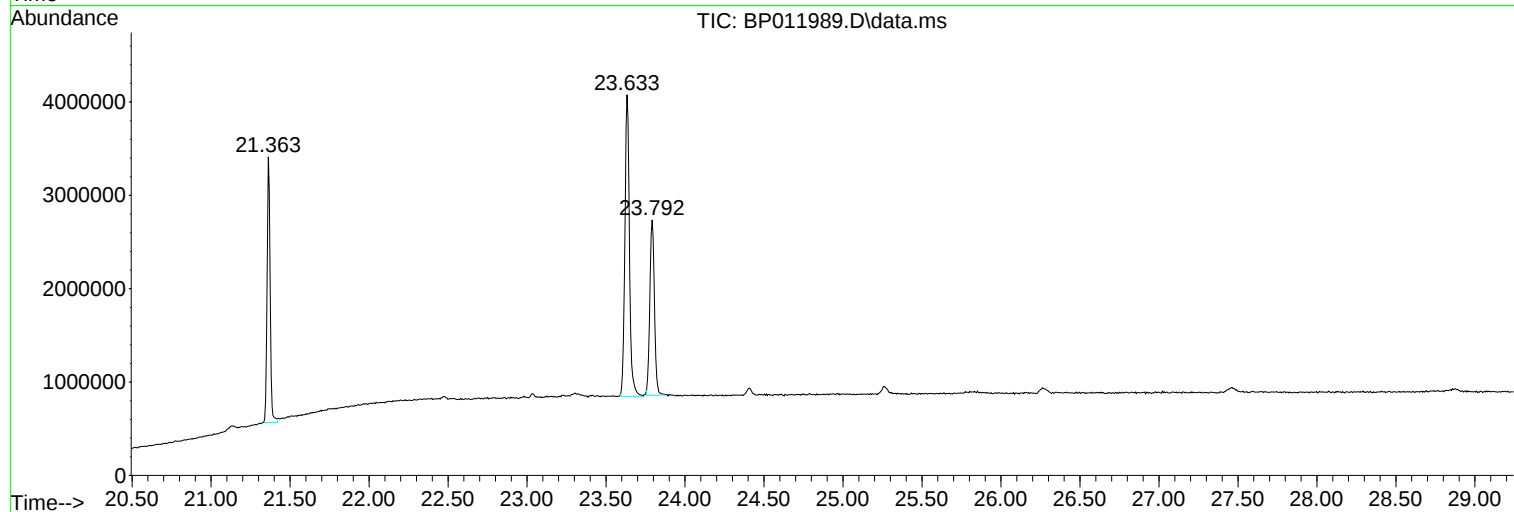
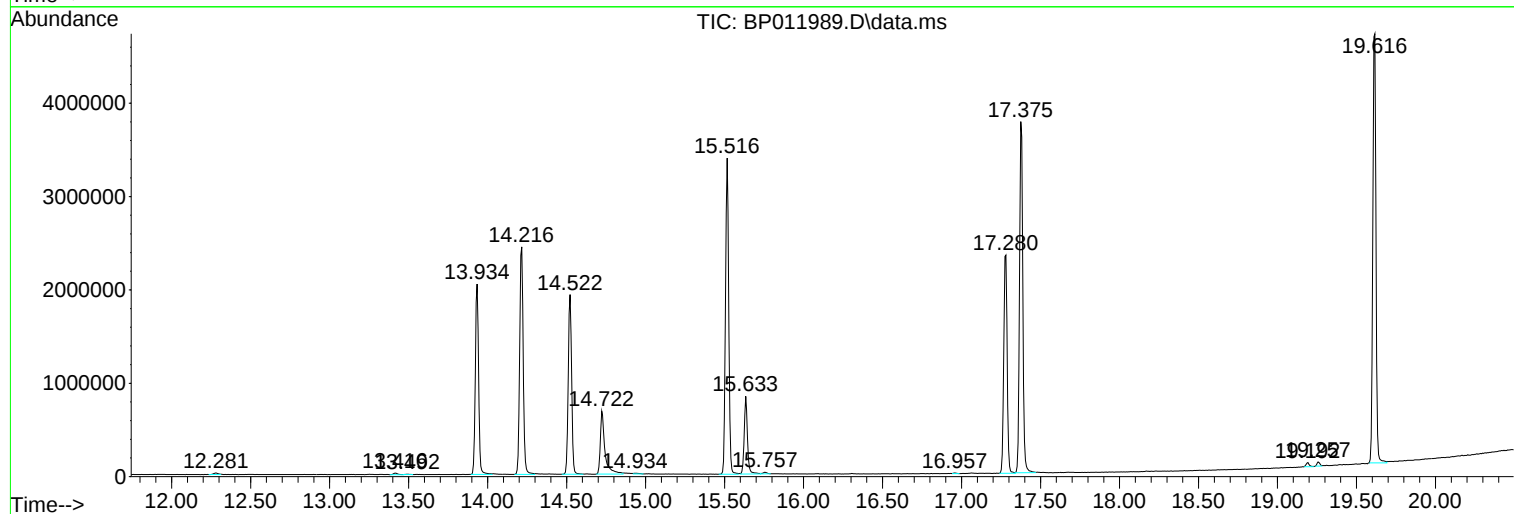
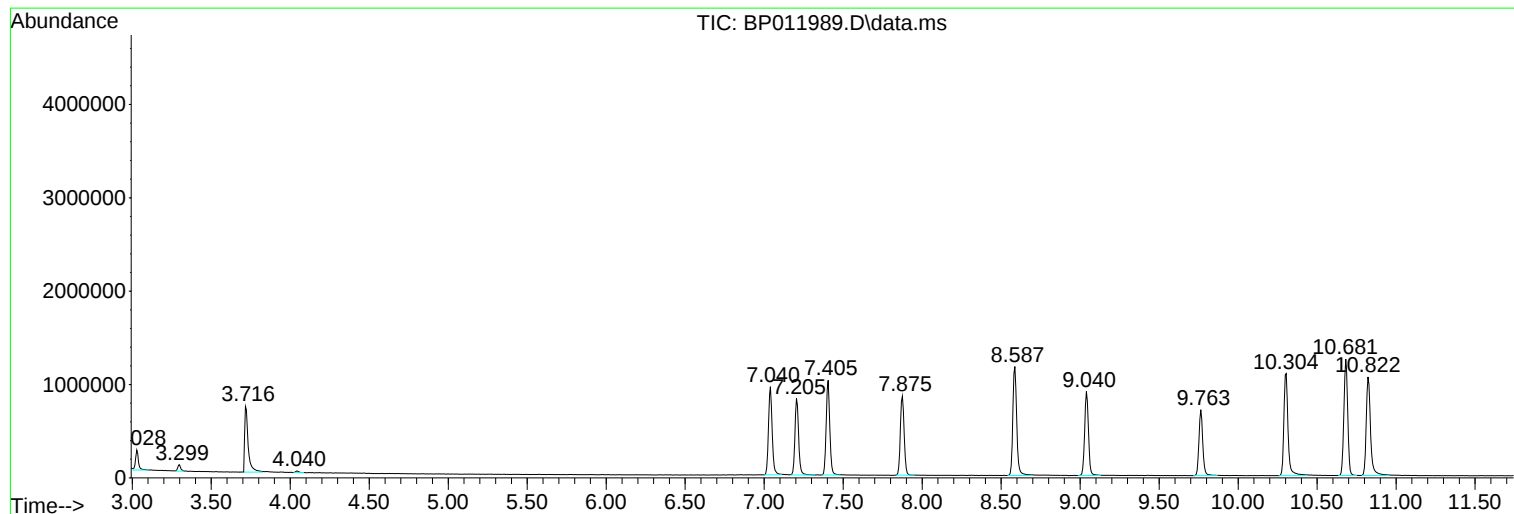
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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



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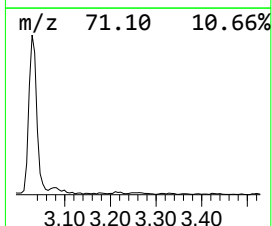
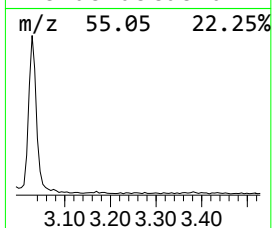
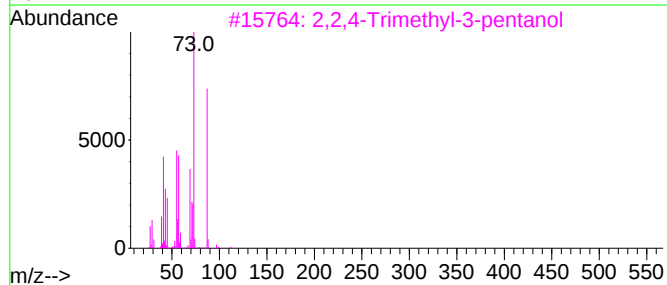
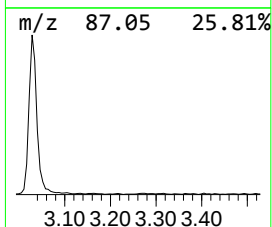
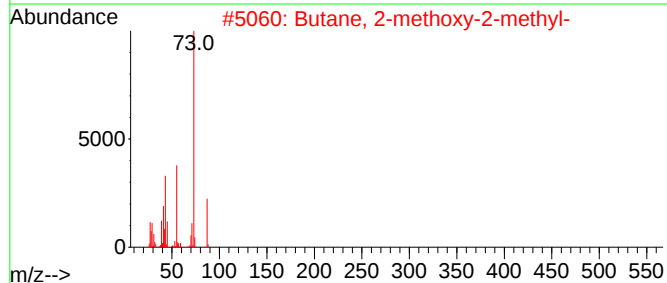
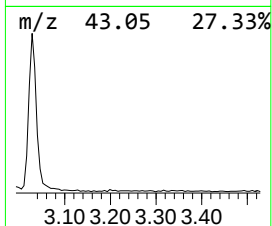
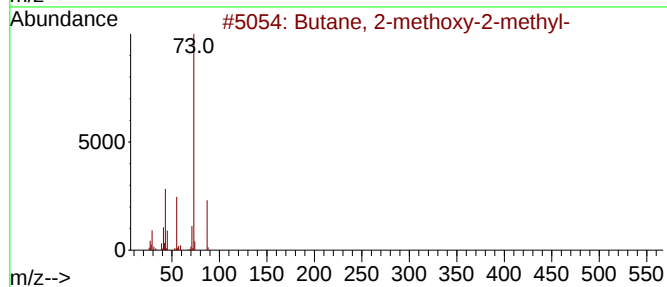
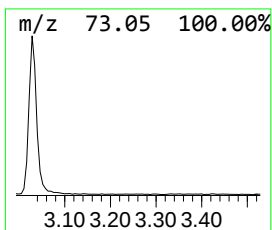
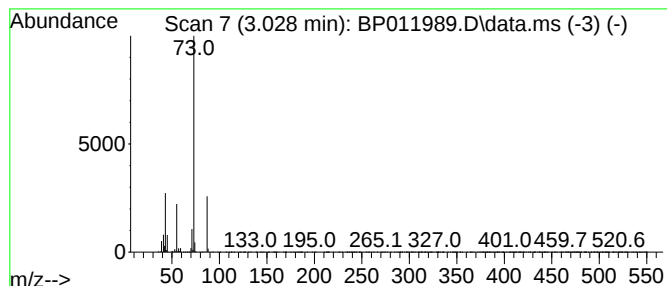
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Quant Title : SVOA CALIBRATION

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	3.96 ng/ul	276634	1,4-Dichlorobenzene-d4	7.875

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	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...	3.028	4.0	ng/u1	276634	1	7.875	1395530	20.0