

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011959.D
 Acq On : 06 Oct 2022 08:46
 Operator : CG/JU
 Sample : PB148127BL
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK127

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: BP011959.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	15	rVB	158693	203507	3.32%	0.320%
2	3.299	49	53	60	rVB	71231	88980	1.45%	0.140%
3	3.716	122	124	140	rBV	729628	1147701	18.73%	1.807%
4	4.046	176	180	187	rVB	17555	26014	0.42%	0.041%
5	7.040	682	689	703	rBV	937717	1502165	24.52%	2.365%
6	7.210	711	718	733	rVB	808978	1307119	21.34%	2.058%
7	7.404	744	751	768	rBV	972395	1591232	25.97%	2.505%
8	7.875	822	831	842	rBV	869802	1401092	22.87%	2.206%
9	8.587	945	952	975	rBV	1201379	1987321	32.44%	3.128%
10	9.045	1021	1030	1045	rBV	842638	1478780	24.14%	2.328%
11	9.769	1143	1153	1168	rBV	675592	1156907	18.88%	1.821%
12	10.304	1237	1244	1259	rBV	1098212	1961593	32.02%	3.088%
13	10.686	1300	1309	1321	rBV	1216981	2089776	34.11%	3.290%
14	10.828	1325	1333	1359	rBV	1107893	2063310	33.68%	3.248%
15	12.280	1572	1580	1590	rBV	16727	38369	0.63%	0.060%
16	13.422	1767	1774	1779	rBV	30651	52752	0.86%	0.083%
17	13.492	1782	1786	1793	rVB6	4097	8118	0.13%	0.013%
18	13.933	1854	1861	1880	rBV	2052870	2929692	47.82%	4.612%
19	14.216	1902	1909	1926	rBV	2444798	3742201	61.09%	5.891%
20	14.527	1955	1962	1970	rBV	1897873	2840217	46.36%	4.471%
21	14.727	1989	1996	2029	rBV	649930	1343300	21.93%	2.115%
22	15.521	2124	2131	2144	rBV	3223772	4864578	79.41%	7.658%
23	15.639	2145	2151	2166	rVV	771565	1163345	18.99%	1.831%
24	15.757	2168	2171	2179	rVB2	15856	26541	0.43%	0.042%
25	17.068	2390	2394	2398	rBV5	5036	9284	0.15%	0.015%
26	17.280	2424	2430	2440	rBV	2423521	3411804	55.69%	5.371%
27	17.380	2441	2447	2464	rVV	3782233	5295486	86.44%	8.336%
28	19.192	2752	2755	2762	rBV3	40240	55908	0.91%	0.088%
29	19.262	2763	2767	2774	rBV	43089	69943	1.14%	0.110%
30	19.615	2821	2827	2842	rBV	4742638	6126154	100.00%	9.644%
31	21.368	3120	3125	3134	rBV	2897699	3718215	60.69%	5.853%
32	23.639	3504	3511	3526	rVB2	3042017	6069647	99.08%	9.555%
33	23.797	3531	3538	3552	rVB2	1899085	3754578	61.29%	5.910%

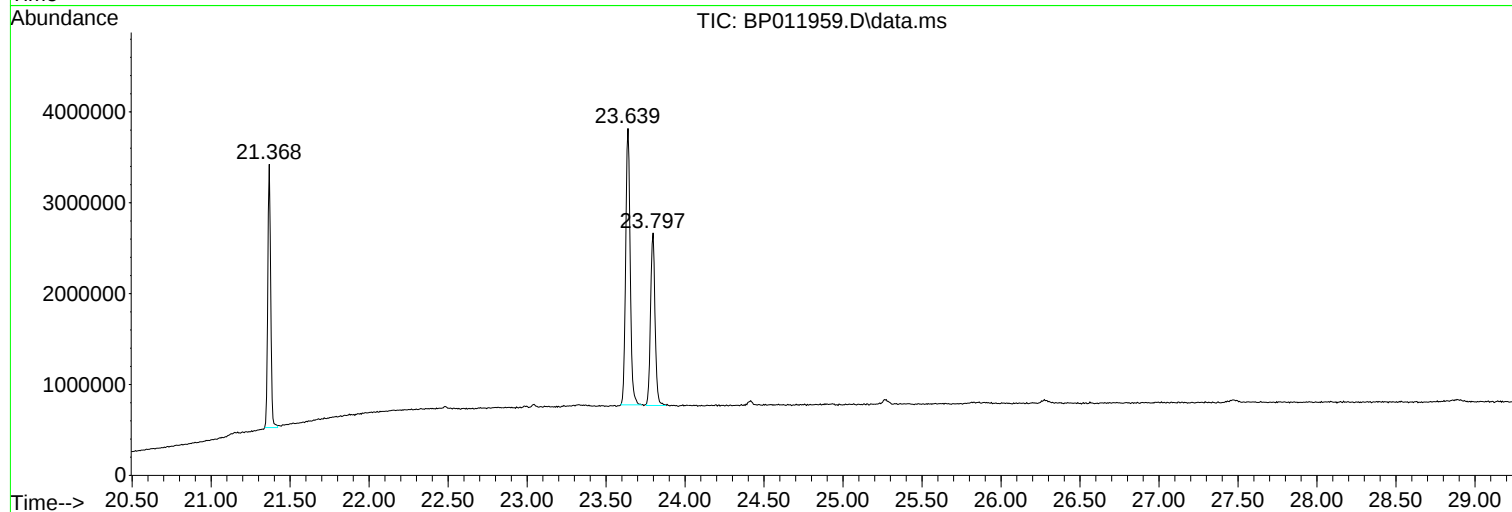
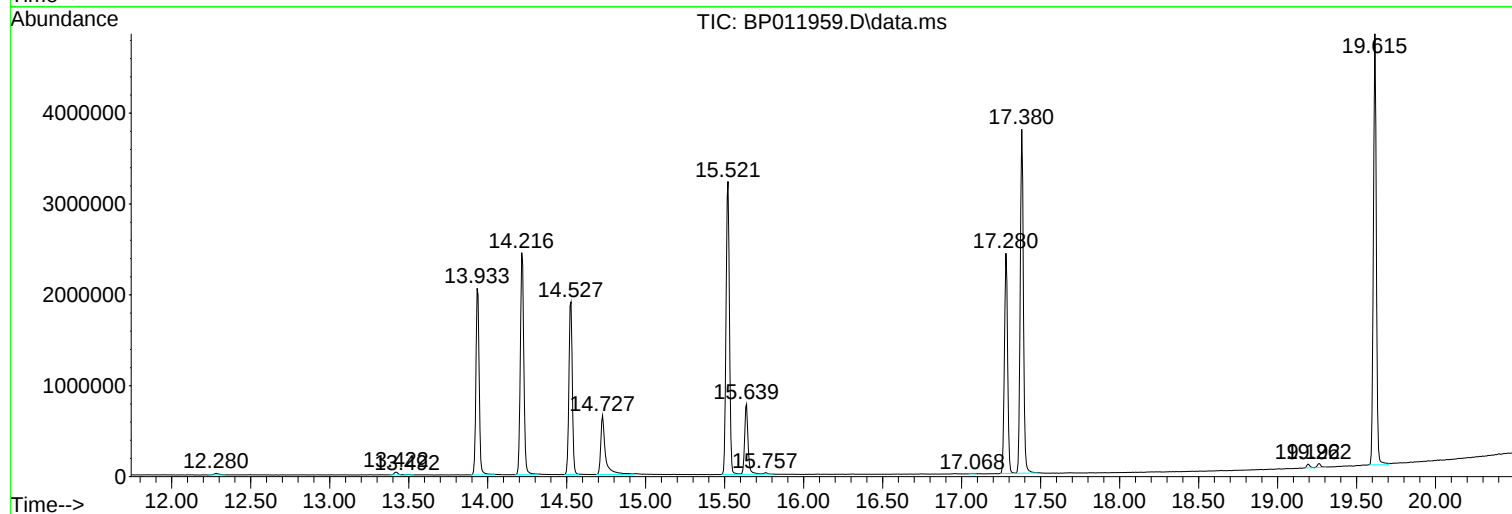
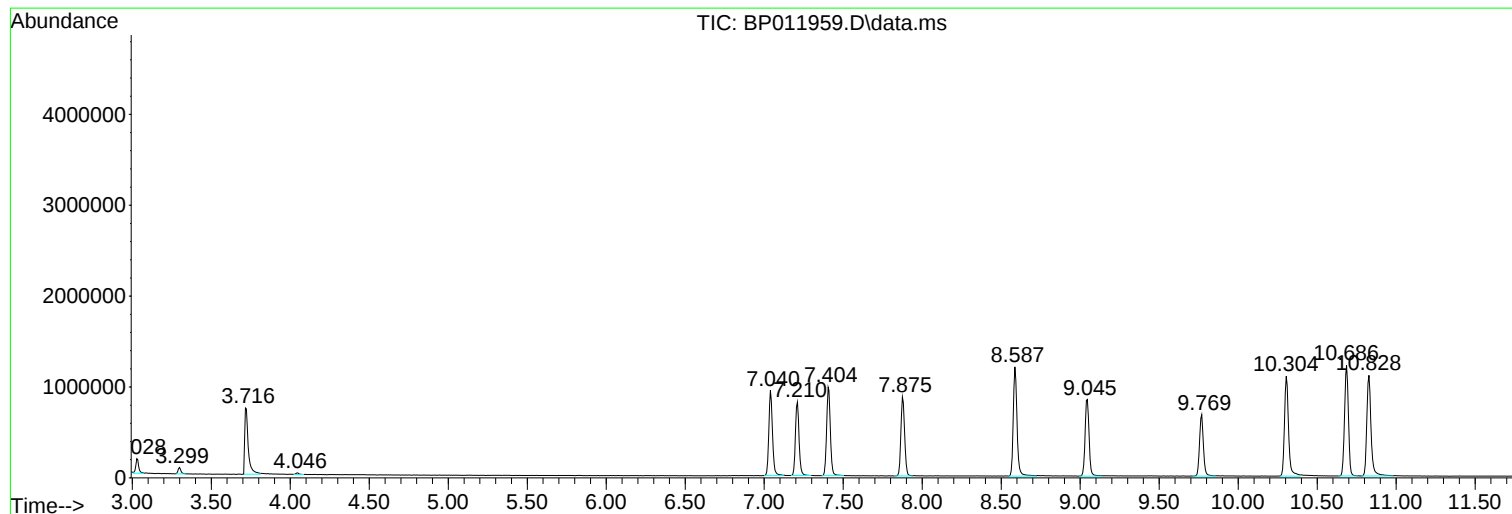
Sum of corrected areas: 63525629

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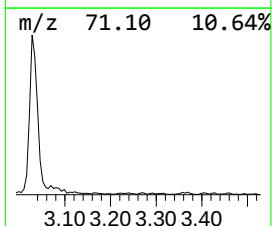
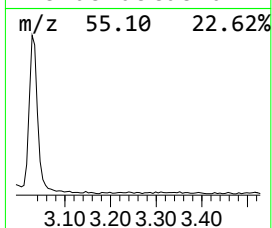
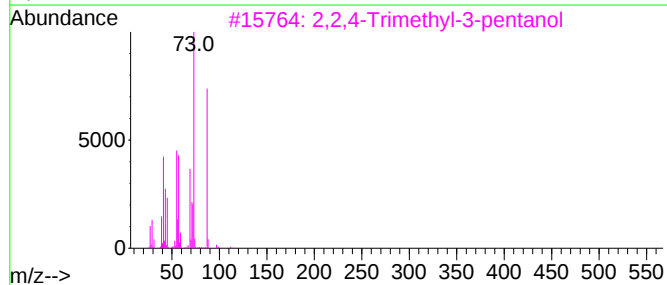
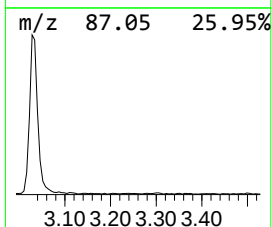
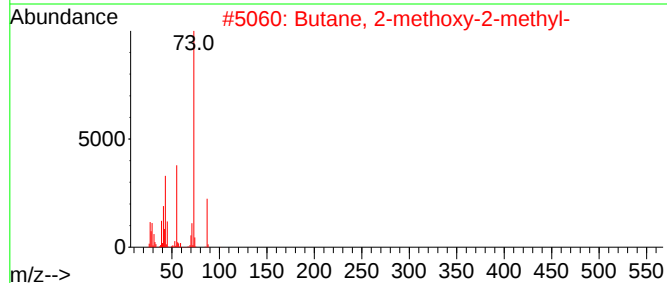
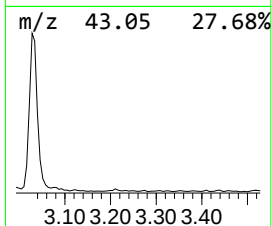
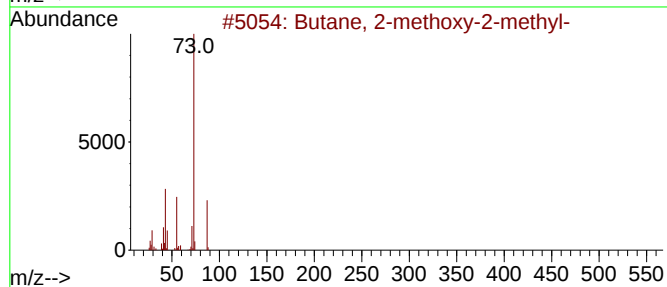
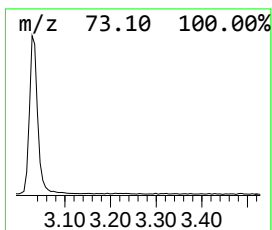
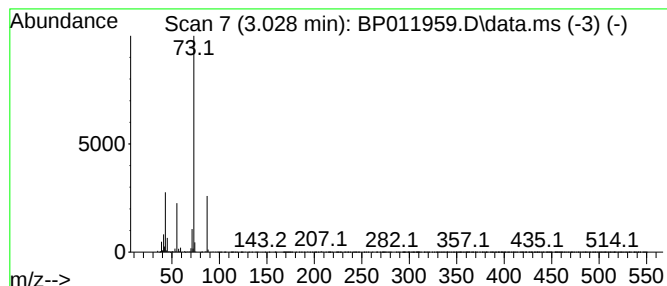
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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	2.90 ng/ul	203507	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	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	39		
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	35		
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	2.9	ng/u1	203507	1	7.875	1401090	20.0