

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011892.D
 Acq On : 04 Oct 2022 03:02
 Operator : CG/JU
 Sample : PB147998BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK998

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: BP011892.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.034	4	8	14	rVB	219553	246947	3.72%	0.387%
2	3.299	49	53	63	rVB	69252	94891	1.43%	0.149%
3	3.722	123	125	145	rBV	763962	1122498	16.90%	1.757%
4	4.052	177	181	187	rVB	18279	25637	0.39%	0.040%
5	7.052	683	691	706	rBV	889242	1504161	22.64%	2.355%
6	7.216	712	719	733	rBV	824619	1317670	19.83%	2.063%
7	7.416	745	753	767	rBV	984980	1600273	24.09%	2.505%
8	7.887	824	833	848	rBV	990590	1566306	23.58%	2.452%
9	8.599	945	954	969	rBV	1139473	1939705	29.20%	3.037%
10	9.052	1023	1031	1044	rBV	859289	1429780	21.52%	2.238%
11	9.775	1144	1154	1168	rBV	659996	1127267	16.97%	1.765%
12	10.316	1238	1246	1265	rBV	1054969	1928554	29.03%	3.019%
13	10.699	1302	1311	1325	rBV	1284963	2263924	34.08%	3.544%
14	10.834	1327	1334	1358	rBV	1054005	1978504	29.78%	3.097%
15	12.293	1575	1582	1590	rBV2	17151	37749	0.57%	0.059%
16	13.428	1769	1775	1780	rBV2	26727	43024	0.65%	0.067%
17	13.510	1783	1789	1796	rVB6	3673	7566	0.11%	0.012%
18	13.945	1856	1863	1876	rBV	1806776	2650941	39.90%	4.150%
19	14.228	1903	1911	1927	rBV	2324857	3462197	52.11%	5.420%
20	14.534	1956	1963	1974	rBV	1959755	2895480	43.58%	4.533%
21	14.734	1990	1997	2022	rBV	636326	1239791	18.66%	1.941%
22	15.528	2125	2132	2146	rBV	3036663	4407723	66.34%	6.900%
23	15.645	2146	2152	2167	rVV	731564	1080173	16.26%	1.691%
24	15.769	2169	2173	2179	rVB2	14431	24027	0.36%	0.038%
25	17.081	2392	2396	2402	rVB4	5849	8900	0.13%	0.014%
26	17.286	2425	2431	2441	rBV	2291856	3441755	51.80%	5.388%
27	17.386	2442	2448	2466	rVV2	3255281	4799750	72.24%	7.514%
28	19.198	2753	2756	2762	rBV2	40410	55482	0.84%	0.087%
29	19.269	2765	2768	2773	rVB	40599	51712	0.78%	0.081%
30	19.622	2822	2828	2839	rBV	4548890	6096226	91.76%	9.544%
31	21.374	3121	3126	3133	rBV	3354672	4279414	64.41%	6.700%
32	23.651	3506	3513	3527	rVB2	3223417	6643852	100.00%	10.401%
33	23.804	3533	3539	3552	rVB2	2131889	4503543	67.79%	7.051%

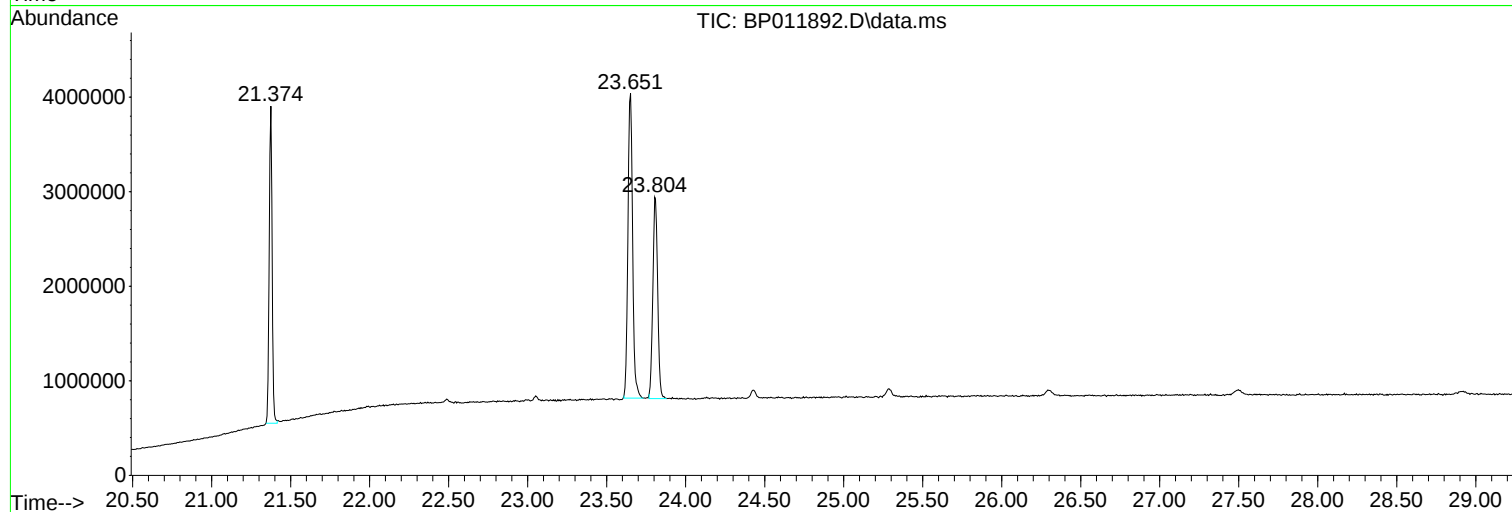
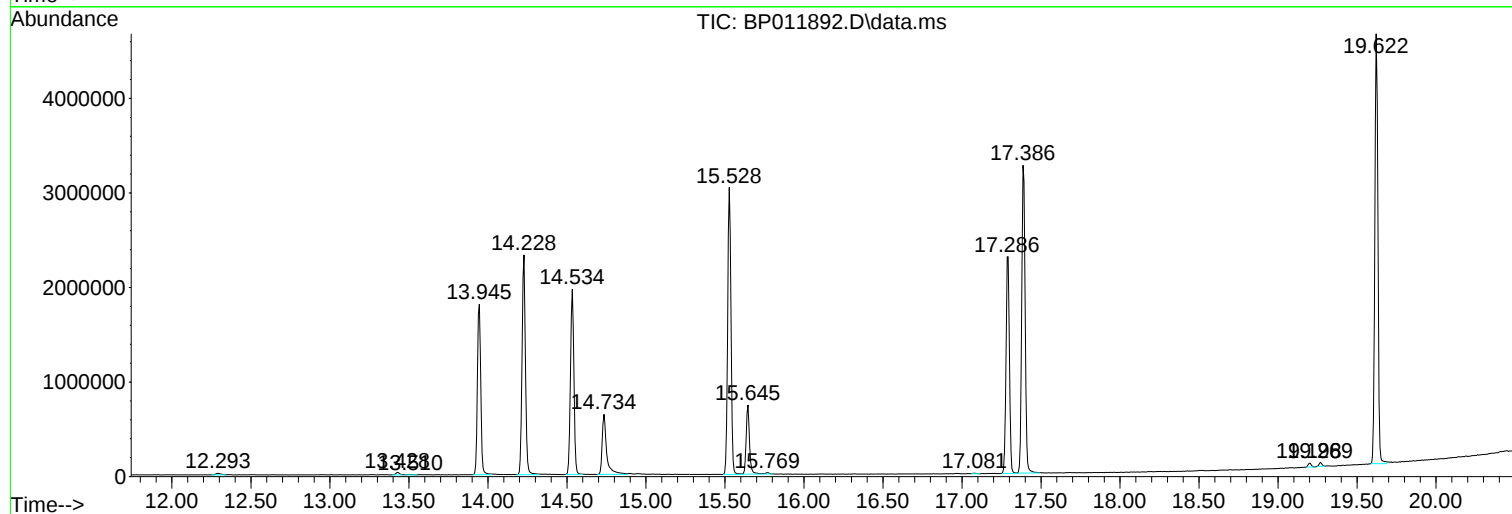
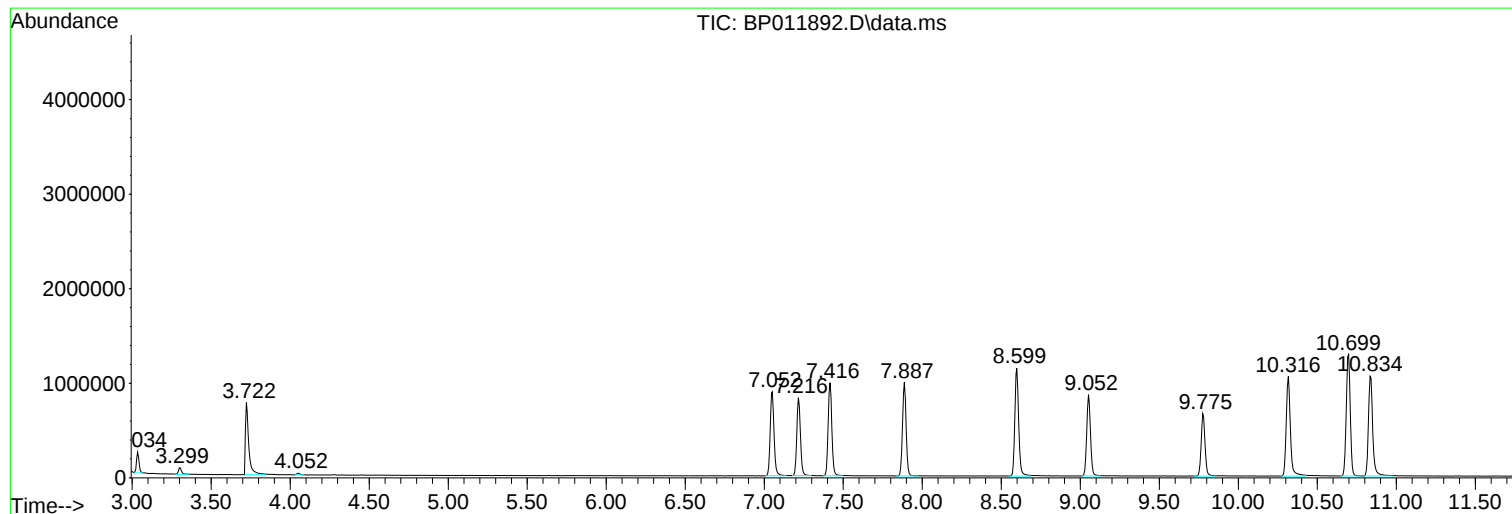
Sum of corrected areas: 63875422

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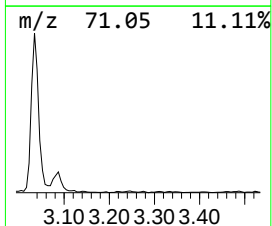
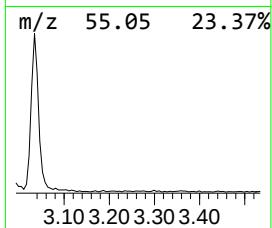
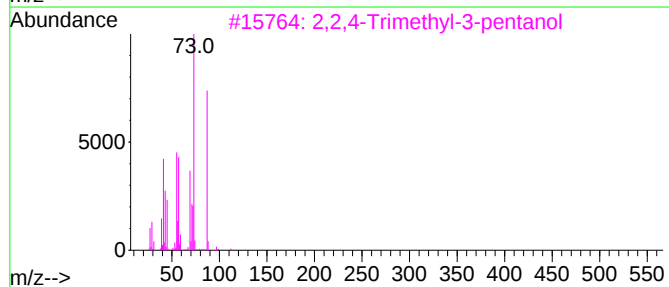
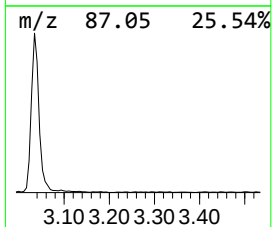
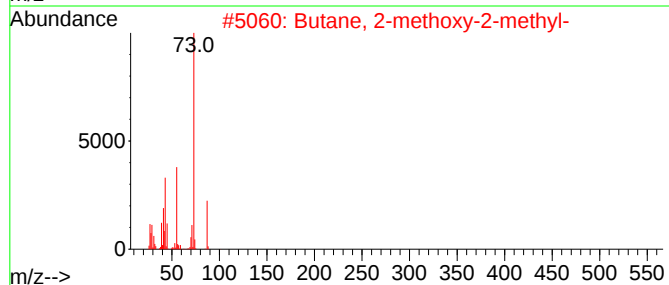
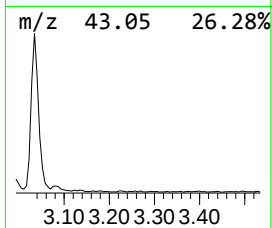
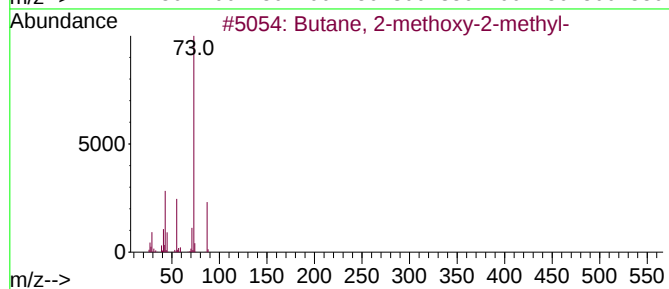
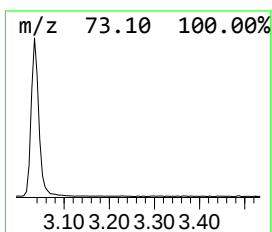
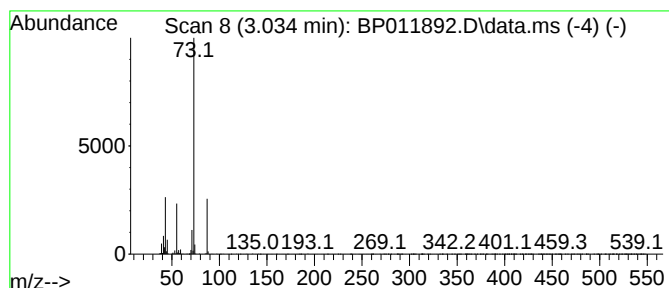
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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.
3.034	3.15 ng/ul	246947	1,4-Dichlorobenzene-d4	7.887

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	40
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	12



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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...	3.034	3.1	ng/u1	246947	1	7.887	1566310	20.0