

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011886.D
 Acq On : 03 Oct 2022 23:38
 Operator : CG/JU
 Sample : PB148023BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK023

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: BP011886.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	15	rVB	202852	232222	3.60%	0.399%
2	3.305	49	54	66	rVB	68403	96308	1.49%	0.166%
3	3.722	123	125	153	rBV	637702	1001798	15.51%	1.722%
4	4.052	177	181	186	rVB	18572	23574	0.37%	0.041%
5	4.381	235	237	244	rVB6	4867	7221	0.11%	0.012%
6	7.052	684	691	708	rBV	865362	1410924	21.85%	2.425%
7	7.216	713	719	732	rVB	773889	1243517	19.25%	2.137%
8	7.416	746	753	768	rBV	945400	1525319	23.62%	2.622%
9	7.887	826	833	843	rBV	837656	1350655	20.91%	2.322%
10	8.599	946	954	970	rBV	1106378	1849576	28.64%	3.179%
11	9.052	1023	1031	1048	rBV	795702	1356574	21.00%	2.332%
12	9.775	1145	1154	1167	rBV	612961	1063920	16.47%	1.829%
13	10.316	1238	1246	1272	rBV	1006745	1814992	28.10%	3.120%
14	10.699	1299	1311	1327	rBV	1112411	1936650	29.99%	3.329%
15	10.840	1327	1335	1353	rBV	957556	1769982	27.41%	3.042%
16	12.293	1574	1582	1589	rBV2	16158	35787	0.55%	0.062%
17	13.428	1769	1775	1780	rBV9	3919	8010	0.12%	0.014%
18	13.504	1785	1788	1795	rVB7	4086	7772	0.12%	0.013%
19	13.945	1856	1863	1876	rBV	1828114	2522126	39.05%	4.335%
20	14.228	1903	1911	1925	rBV	2145042	3198083	49.52%	5.497%
21	14.534	1956	1963	1979	rBV	1693510	2455611	38.02%	4.221%
22	14.734	1990	1997	2021	rBV	595283	1162382	18.00%	1.998%
23	15.528	2125	2132	2145	rBV	2861552	4091407	63.35%	7.033%
24	15.645	2146	2152	2167	rVV	749802	1050795	16.27%	1.806%
25	15.769	2168	2173	2180	rVB2	24243	41927	0.65%	0.072%
26	16.316	2262	2266	2270	rBV4	7994	10867	0.17%	0.019%
27	16.969	2372	2377	2382	rBV9	4837	7234	0.11%	0.012%
28	17.292	2425	2432	2442	rBV	1964338	2833477	43.87%	4.870%
29	17.386	2442	2448	2465	rVV	3144052	4467368	69.17%	7.679%
30	19.198	2753	2756	2762	rVB2	35677	50141	0.78%	0.086%
31	19.269	2765	2768	2772	rVV	36468	46263	0.72%	0.080%
32	19.622	2822	2828	2840	rBV	4343576	5776351	89.44%	9.929%
33	21.374	3121	3126	3134	rBV	2823699	3587408	55.55%	6.166%
34	23.651	3506	3513	3526	rVB2	3271881	6458443	100.00%	11.101%
35	23.810	3533	3540	3548	rVB2	1893849	3683524	57.03%	6.331%

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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

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

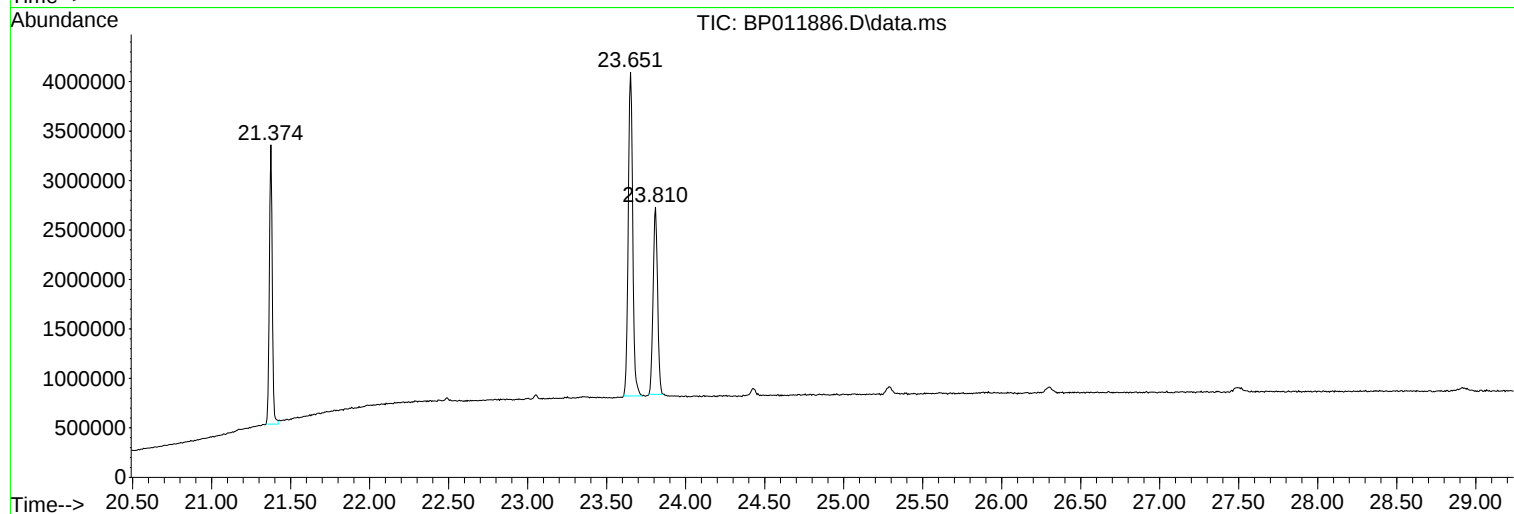
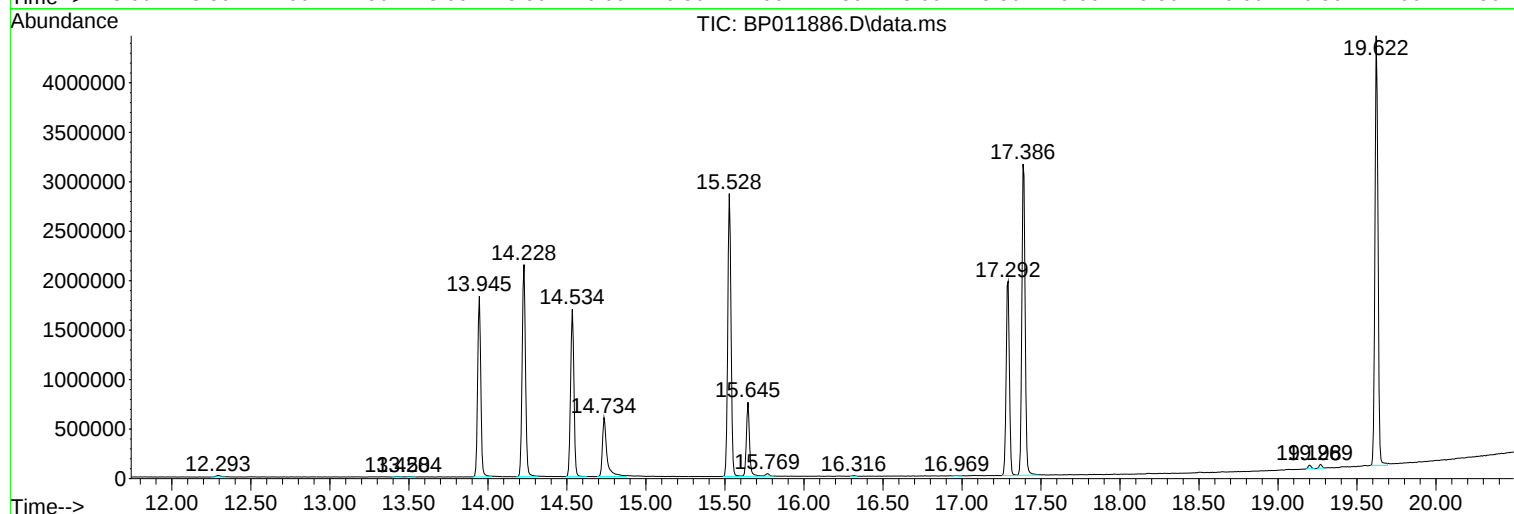
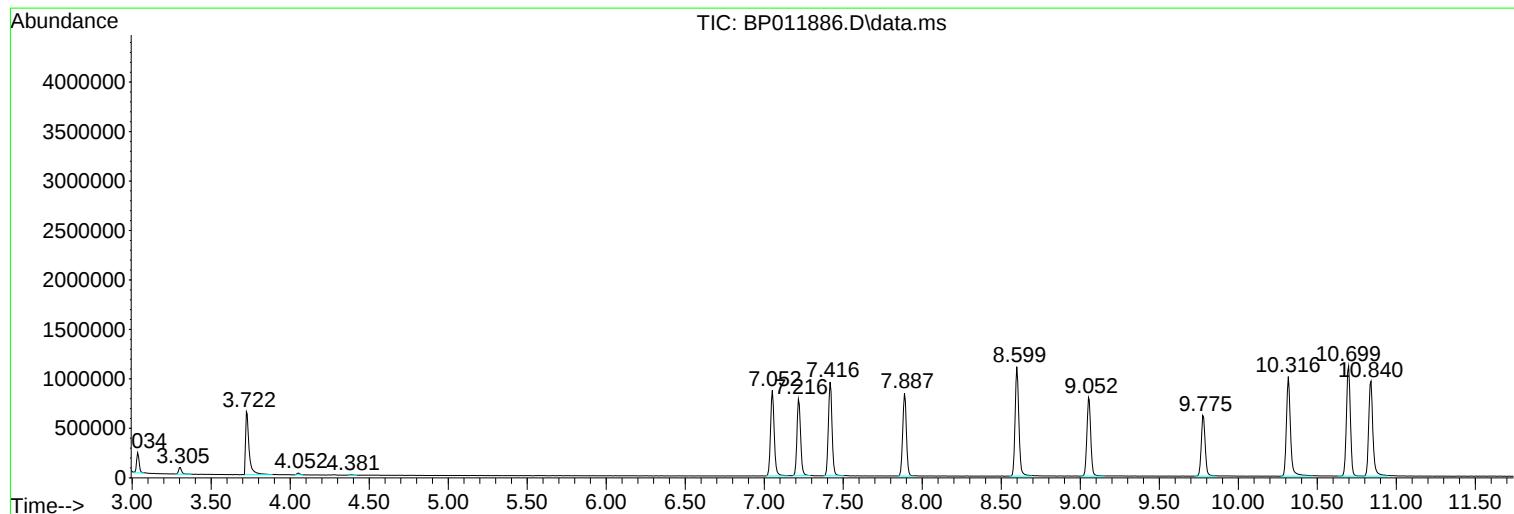
Sum of corrected areas: 58178208

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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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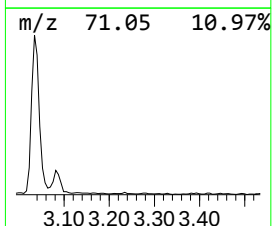
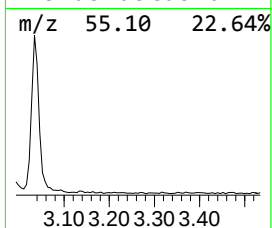
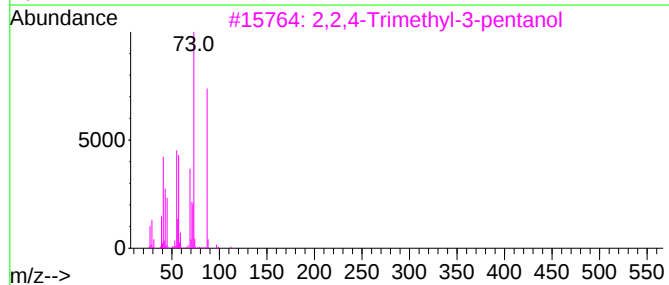
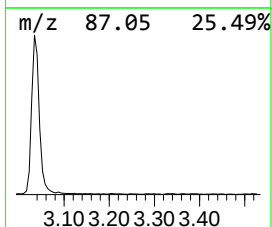
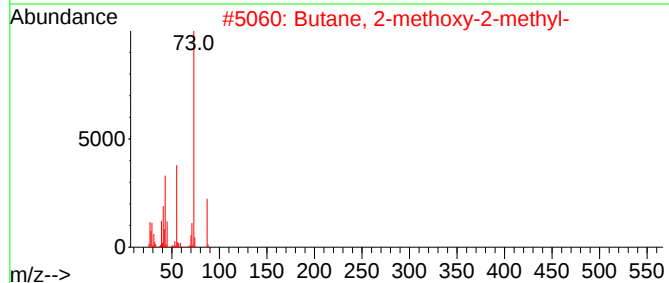
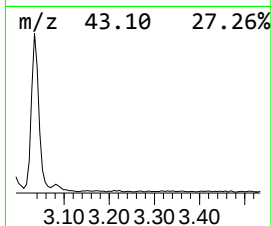
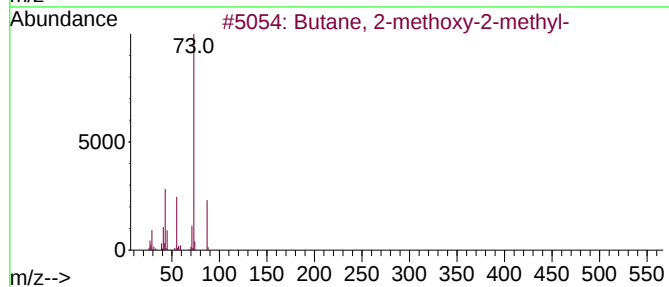
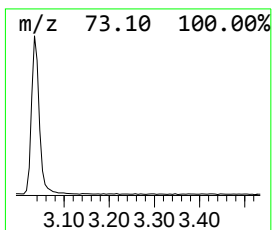
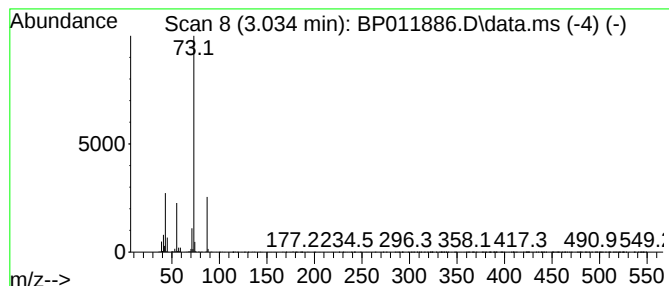
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Quant Title : SVOA CALIBRATION

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.
3.034	3.44 ng/ul	232222	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	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			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...	3.034	3.4	ng/u1	232222	1	7.887	1350660	20.0