

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011868.D
 Acq On : 03 Oct 2022 12:50
 Operator : CG/JU
 Sample : PB147991BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK991

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: BP011868.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	14	rVB	191397	234049	3.50%	0.373%
2	3.305	50	54	62	rVB	69156	86010	1.28%	0.137%
3	3.728	123	126	151	rBV	742835	1119383	16.72%	1.784%
4	4.052	177	181	187	rBV	18234	25919	0.39%	0.041%
5	4.281	216	220	224	rVB3	6075	8719	0.13%	0.014%
6	7.058	681	692	708	rBV	907738	1513024	22.60%	2.411%
7	7.222	713	720	732	rVV	827043	1319700	19.71%	2.103%
8	7.422	746	754	770	rBV	996416	1604248	23.96%	2.557%
9	7.893	827	834	844	rBV	888423	1445095	21.58%	2.303%
10	8.605	947	955	972	rBV	1164131	1959261	29.26%	3.122%
11	9.058	1022	1032	1047	rBV	871915	1458608	21.79%	2.325%
12	9.781	1146	1155	1168	rBV	687689	1158656	17.31%	1.847%
13	10.322	1237	1247	1273	rBV	1076496	1967094	29.38%	3.135%
14	10.705	1304	1312	1326	rBV	1203931	2096434	31.31%	3.341%
15	10.840	1328	1335	1355	rBV	1099726	2014947	30.10%	3.211%
16	12.293	1576	1582	1592	rBV2	17379	37146	0.55%	0.059%
17	13.434	1771	1776	1781	rVB9	5793	8950	0.13%	0.014%
18	13.951	1855	1864	1881	rBV	1862624	2796053	41.76%	4.456%
19	14.228	1903	1911	1928	rBV	2370633	3638274	54.34%	5.798%
20	14.540	1955	1964	1978	rBV	1846027	2754705	41.15%	4.390%
21	14.740	1991	1998	2026	rBV	659303	1256588	18.77%	2.003%
22	14.945	2031	2033	2040	rVB8	5368	10108	0.15%	0.016%
23	15.534	2126	2133	2143	rBV	3199259	4534486	67.73%	7.226%
24	15.651	2146	2153	2168	rVV	758133	1134315	16.94%	1.808%
25	15.775	2170	2174	2183	rVB3	16215	30790	0.46%	0.049%
26	16.316	2262	2266	2270	rBV7	4136	7068	0.11%	0.011%
27	16.969	2374	2377	2380	rBV5	5311	7810	0.12%	0.012%
28	17.081	2392	2396	2400	rBV3	6552	9330	0.14%	0.015%
29	17.292	2425	2432	2440	rBV	2232111	3185121	47.57%	5.076%
30	17.392	2443	2449	2464	rVV	3421378	4927041	73.59%	7.852%
31	19.204	2753	2757	2762	rBV2	42131	60344	0.90%	0.096%
32	19.269	2765	2768	2773	rBV	43909	68135	1.02%	0.109%
33	19.628	2823	2829	2837	rBV	4466720	5916953	88.38%	9.430%
34	21.375	3121	3126	3134	rBV	2787162	3718421	55.54%	5.926%
35	23.657	3507	3514	3529	rVB2	3139522	6694970	100.00%	10.670%
36	23.810	3534	3540	3552	rVB2	1902217	3940670	58.86%	6.280%

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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

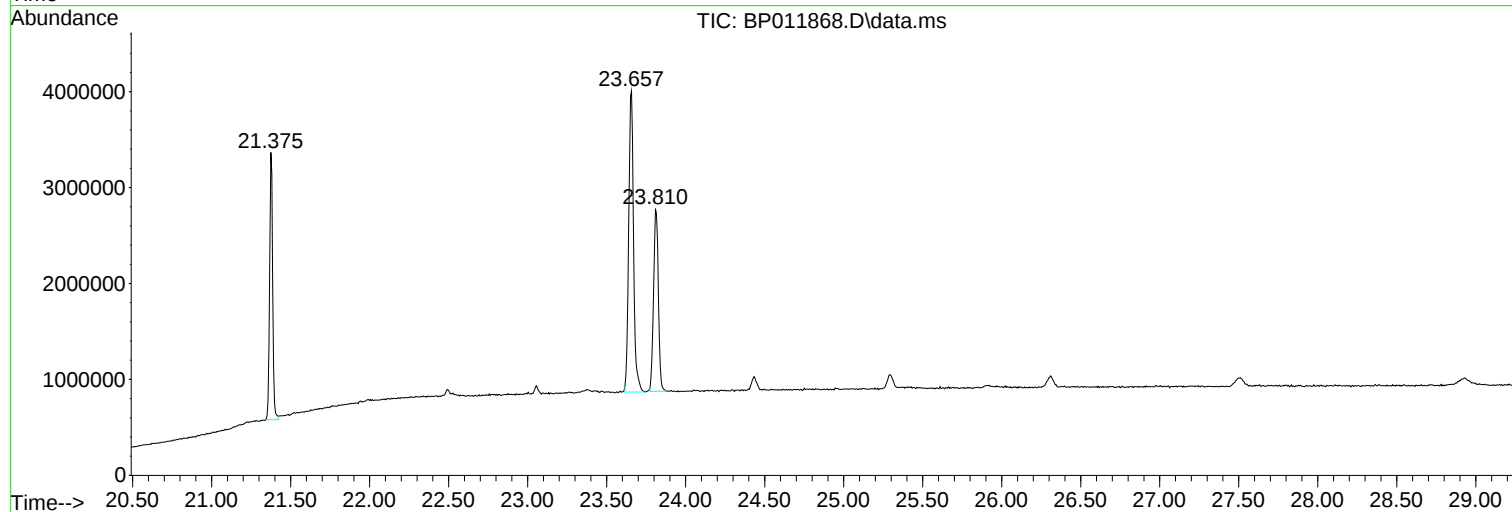
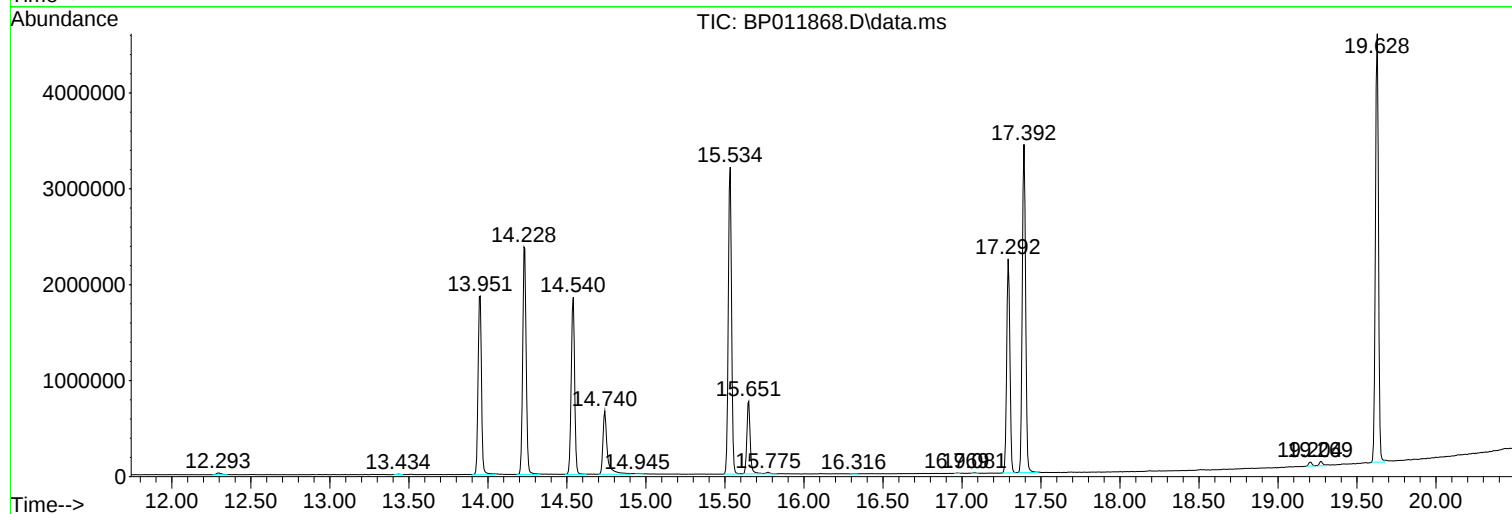
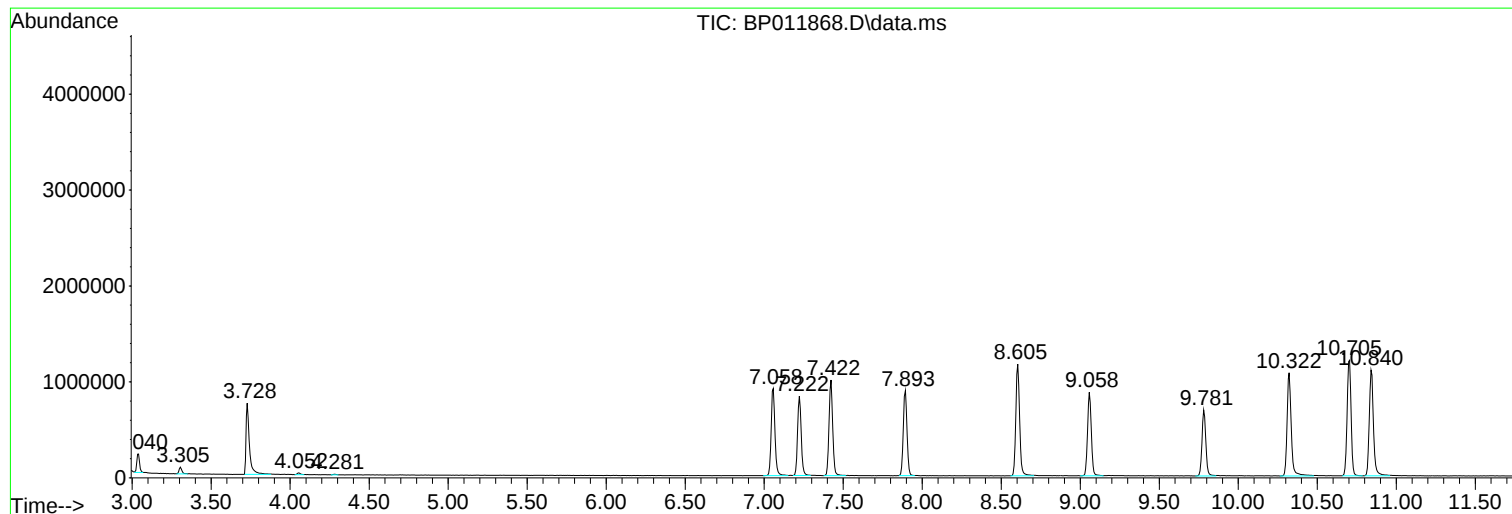
Sum of corrected areas: 62748425

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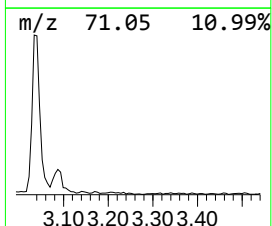
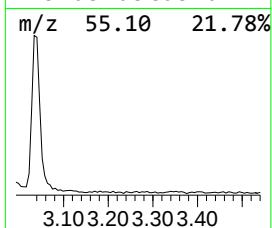
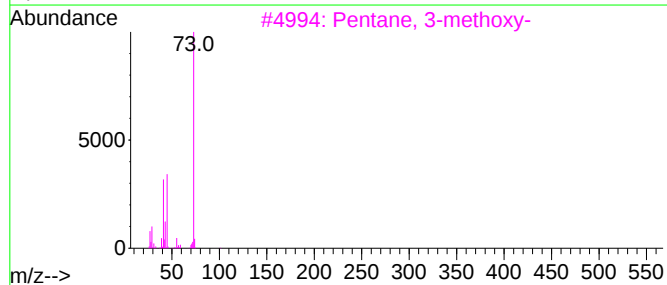
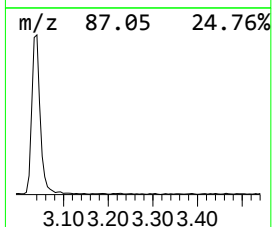
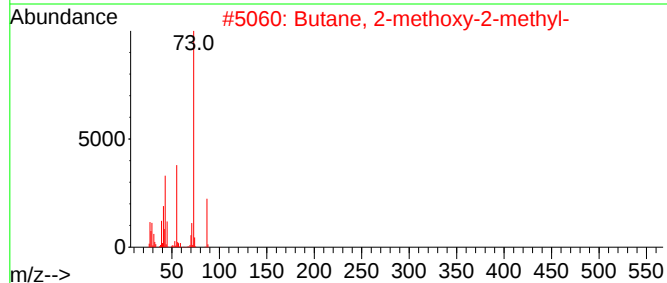
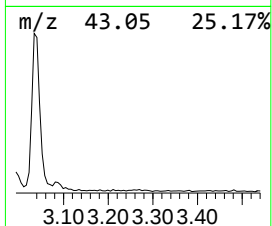
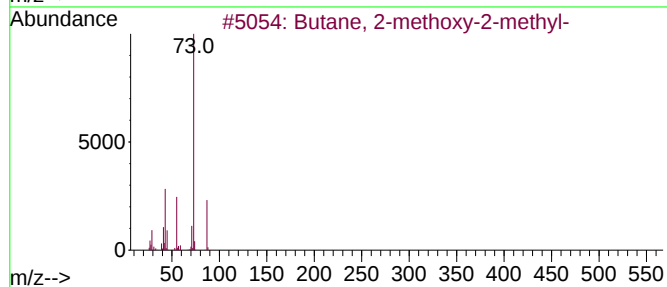
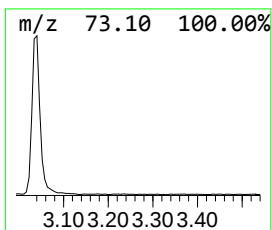
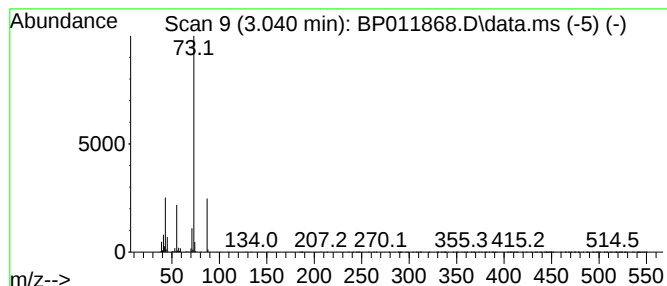
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.040	3.24 ng/ul	234049	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	83
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4			N-Ethylformamide	73	C3H7NO	000627-45-2	9
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.040	3.2	ng/u1	234049	1	7.893	1445100	20.0