

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011885.D
 Acq On : 03 Oct 2022 23:04
 Operator : CG/JU
 Sample : PB148012BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK012

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: BP011885.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	139876	162880	2.20%	0.241%
2	3.305	49	54	67	rVB	75343	103732	1.40%	0.153%
3	3.728	122	126	151	rBV	618562	997399	13.45%	1.474%
4	4.052	177	181	186	rBV	21211	26362	0.36%	0.039%
5	4.281	215	220	224	rVB7	6378	10889	0.15%	0.016%
6	7.052	684	691	711	rBV	1020465	1645913	22.19%	2.433%
7	7.216	713	719	733	rVB	884891	1439330	19.41%	2.127%
8	7.416	746	753	764	rBV	1080933	1769828	23.86%	2.616%
9	7.887	826	833	845	rBV	846275	1365435	18.41%	2.018%
10	8.599	946	954	974	rBV	1322482	2201422	29.68%	3.254%
11	9.057	1021	1032	1046	rBV	933052	1606872	21.67%	2.375%
12	9.775	1146	1154	1169	rBV	739044	1298318	17.51%	1.919%
13	10.316	1236	1246	1268	rBV	1231486	2218615	29.91%	3.279%
14	10.699	1302	1311	1324	rVB	1166648	2031131	27.39%	3.002%
15	10.840	1328	1335	1350	rBV	689105	1289330	17.38%	1.906%
16	12.293	1573	1582	1590	rBV	20393	42067	0.57%	0.062%
17	13.428	1771	1775	1780	rBV6	7338	10879	0.15%	0.016%
18	13.504	1784	1788	1795	rVB4	4747	9311	0.13%	0.014%
19	13.945	1856	1863	1875	rBV	2281986	3159446	42.60%	4.670%
20	14.228	1903	1911	1928	rBV	2828348	4085821	55.09%	6.039%
21	14.534	1953	1963	1974	rBV	1827361	2729715	36.81%	4.035%
22	14.734	1991	1997	2022	rBV	832538	1555866	20.98%	2.300%
23	15.528	2125	2132	2145	rBV	3625352	5225771	70.46%	7.724%
24	15.645	2146	2152	2167	rVV	972318	1389318	18.73%	2.054%
25	15.769	2169	2173	2182	rVB2	30769	52117	0.70%	0.077%
26	16.310	2261	2265	2271	rBV4	9205	14980	0.20%	0.022%
27	16.969	2374	2377	2381	rVB5	6417	8856	0.12%	0.013%
28	17.286	2425	2431	2441	rBV	2205340	3283560	44.27%	4.853%
29	17.392	2442	2449	2467	rVV	3839154	5620136	75.78%	8.307%
30	17.957	2541	2545	2549	rBV6	13415	16542	0.22%	0.024%
31	19.198	2753	2756	2762	rBV2	47924	61464	0.83%	0.091%
32	19.269	2764	2768	2773	rBV	51308	74401	1.00%	0.110%
33	19.622	2822	2828	2840	rBV	5163114	6989338	94.24%	10.331%
34	21.374	3121	3126	3133	rBV	2913792	3787124	51.06%	5.598%
35	23.651	3506	3513	3529	rVB2	3674964	7416582	100.00%	10.962%
36	23.810	3533	3540	3553	rVB2	1930449	3954634	53.32%	5.845%

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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

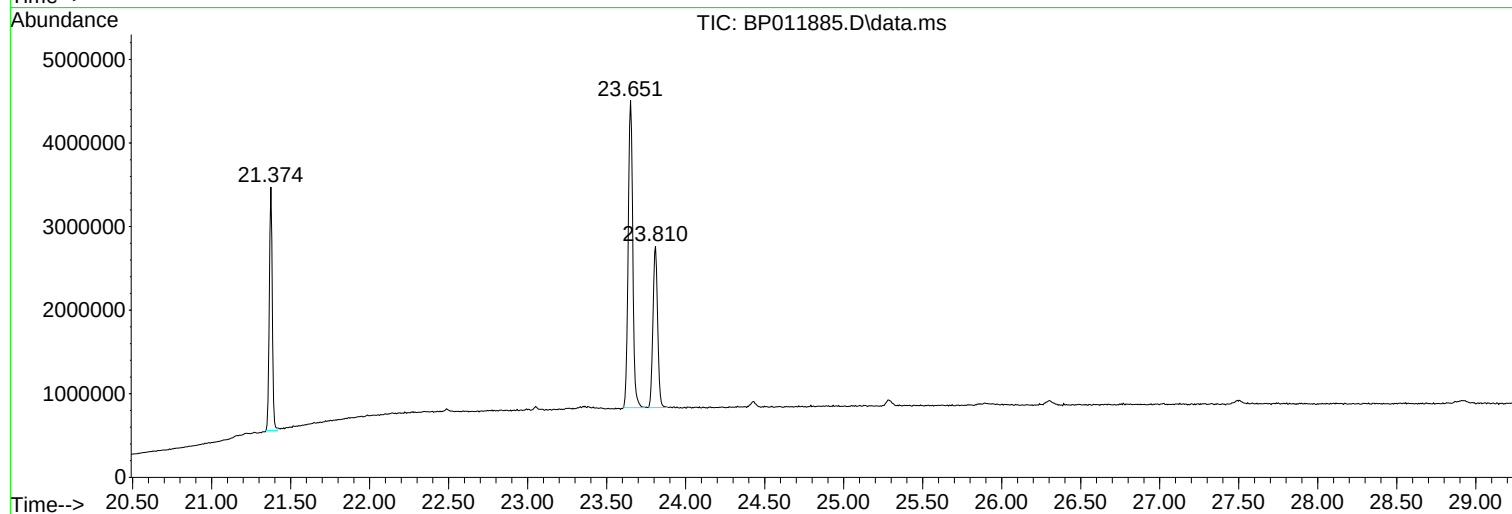
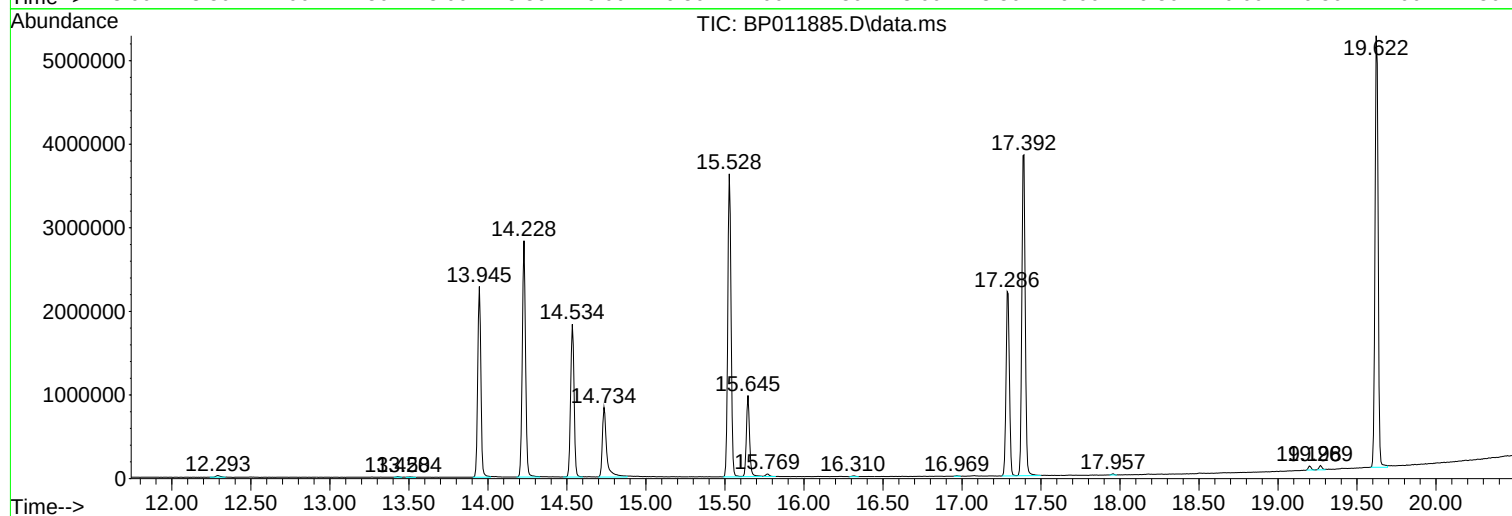
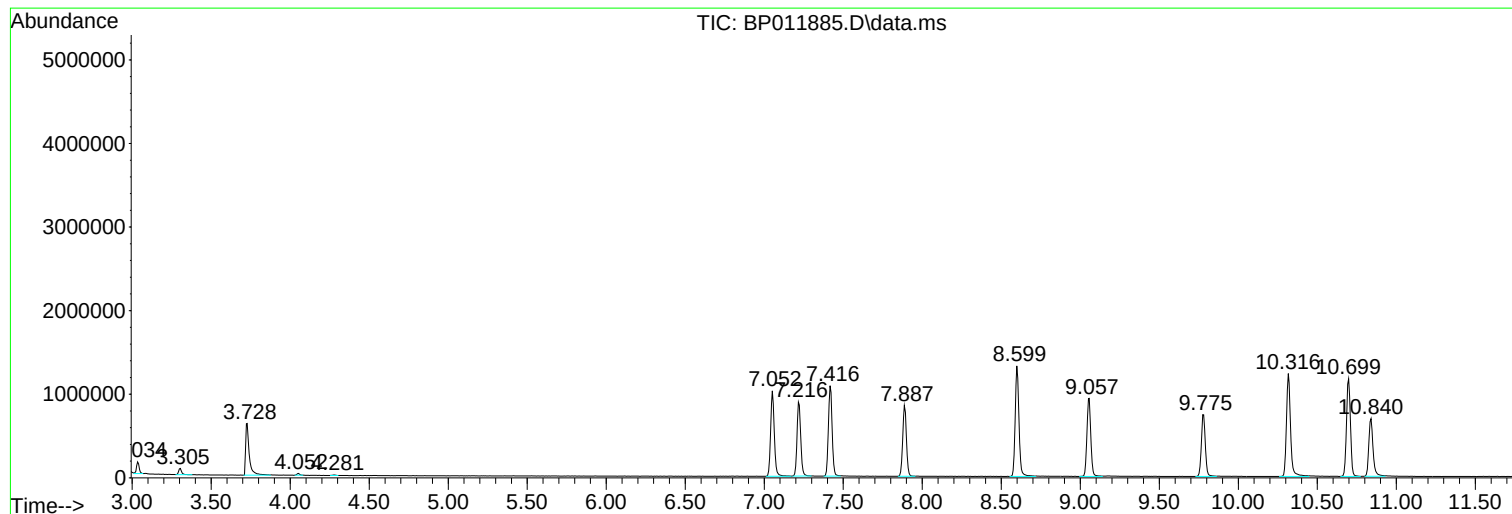
Sum of corrected areas: 67655384

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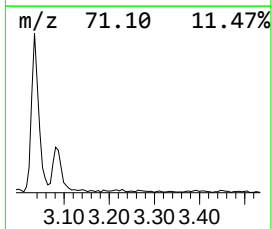
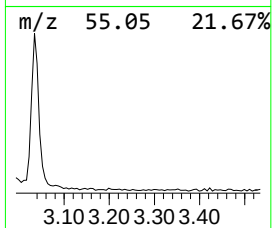
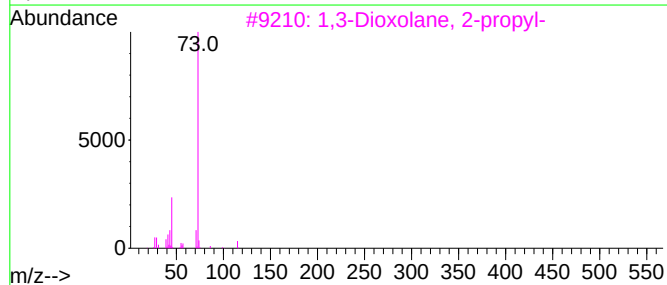
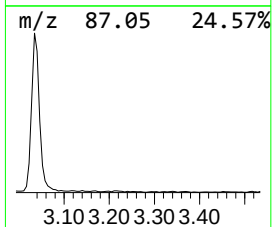
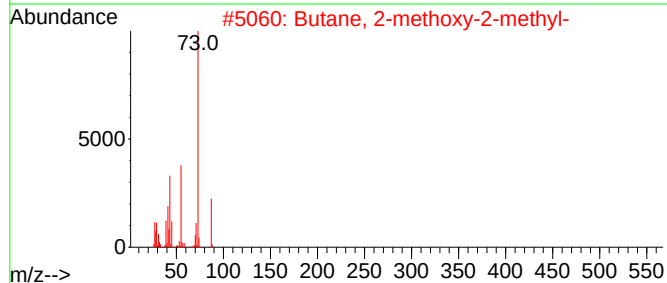
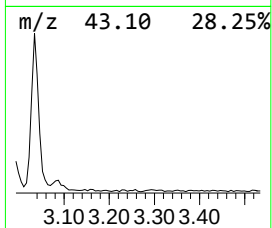
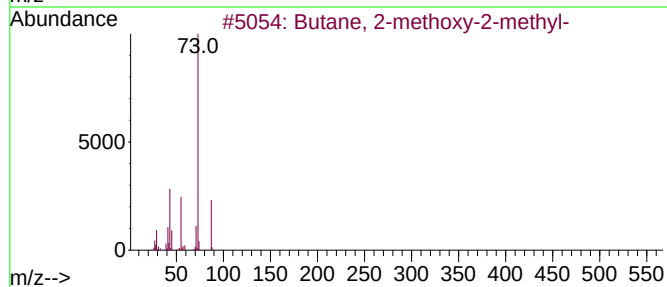
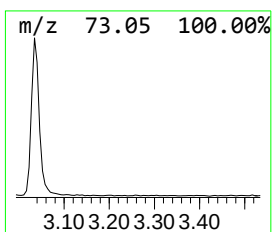
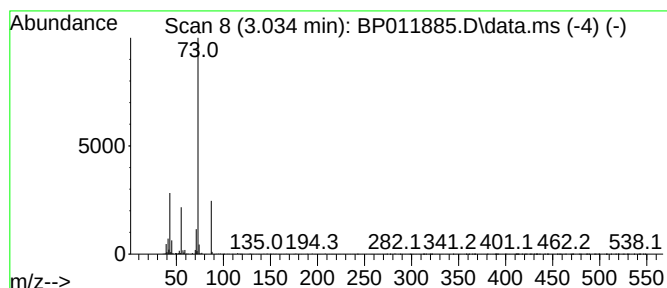
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.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	2.39 ng/ul	162880	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	28		
3	1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	23		
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9		
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...	3.034	2.4	ng/u1	162880	1	7.887	1365440	20.0