

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011974.D
 Acq On : 06 Oct 2022 17:25
 Operator : CG/JU
 Sample : PB148078BL
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK078

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: BP011974.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	4	7	19	rVB	193543	241352	4.57%	0.485%
2	3.299	49	53	58	rBV	61103	76398	1.45%	0.154%
3	3.722	121	125	148	rBV	660137	1009329	19.12%	2.029%
4	4.046	177	180	185	rVB	16100	21549	0.41%	0.043%
5	7.046	683	690	703	rBV	788949	1280316	24.26%	2.574%
6	7.211	712	718	734	rVB	702188	1128678	21.38%	2.269%
7	7.411	745	752	767	rBV	856338	1357072	25.71%	2.728%
8	7.881	824	832	843	rBV	705847	1153548	21.86%	2.319%
9	8.593	946	953	973	rBV	952389	1624830	30.79%	3.266%
10	9.046	1023	1030	1043	rBV	722455	1227371	23.25%	2.467%
11	9.769	1142	1153	1168	rBV	569429	964887	18.28%	1.939%
12	10.310	1235	1245	1268	rBV	858535	1604820	30.41%	3.226%
13	10.687	1300	1309	1324	rVB	943931	1622137	30.73%	3.261%
14	10.828	1326	1333	1351	rBV	866509	1618458	30.66%	3.253%
15	12.287	1574	1581	1588	rBV2	13420	30140	0.57%	0.061%
16	13.428	1770	1775	1780	rVB9	3800	8481	0.16%	0.017%
17	13.498	1780	1787	1791	rBV10	3644	7600	0.14%	0.015%
18	13.940	1855	1862	1878	rBV	1438179	2124065	40.24%	4.270%
19	14.222	1900	1910	1930	rBV	1832298	2775709	52.59%	5.579%
20	14.528	1955	1962	1975	rBV	1312584	2008008	38.05%	4.036%
21	14.734	1990	1997	2022	rBV	457650	1025656	19.43%	2.062%
22	15.522	2124	2131	2144	rBV	2377405	3523251	66.75%	7.082%
23	15.639	2144	2151	2166	rVV	596271	870722	16.50%	1.750%
24	15.757	2168	2171	2176	rVB6	9974	15945	0.30%	0.032%
25	17.281	2424	2430	2440	rBV	1694110	2381862	45.13%	4.788%
26	17.381	2441	2447	2466	rVV	2773519	4012322	76.02%	8.065%
27	19.198	2752	2756	2761	rBV4	31354	43871	0.83%	0.088%
28	19.263	2764	2767	2771	rVB	29922	40521	0.77%	0.081%
29	19.616	2821	2827	2835	rBV	3616059	4906934	92.97%	9.863%
30	21.369	3120	3125	3133	rBV	2121796	2845582	53.91%	5.720%
31	23.639	3505	3511	3526	rVB2	2728327	5277920	100.00%	10.609%
32	23.798	3532	3538	3548	rVB	1443557	2920142	55.33%	5.870%

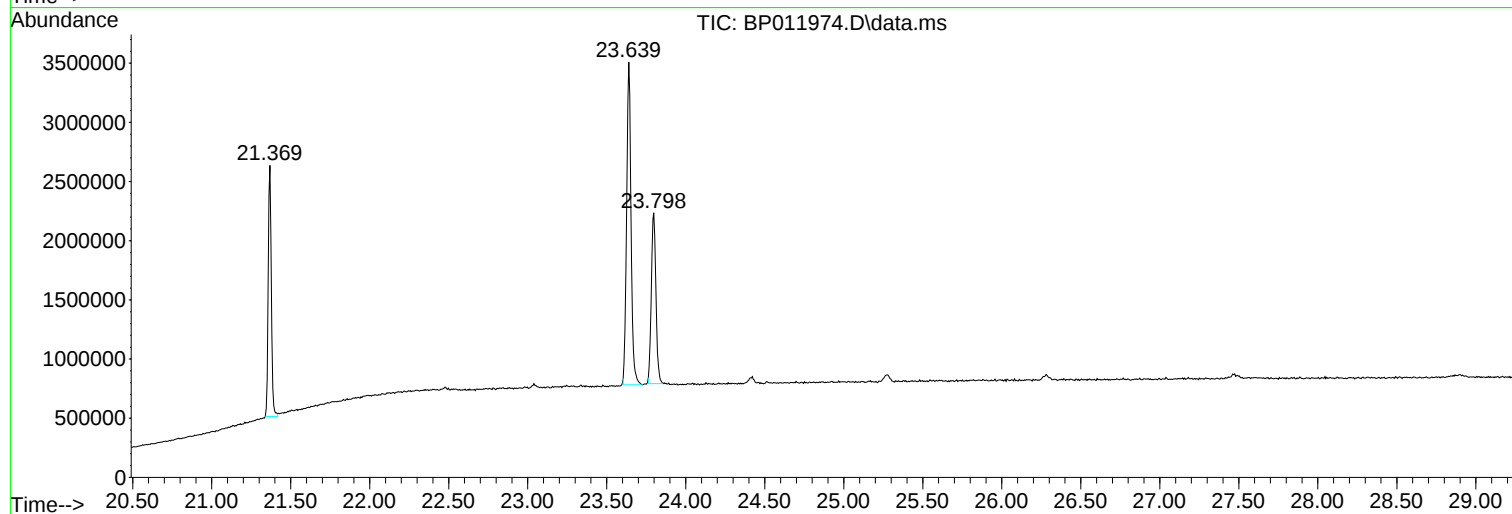
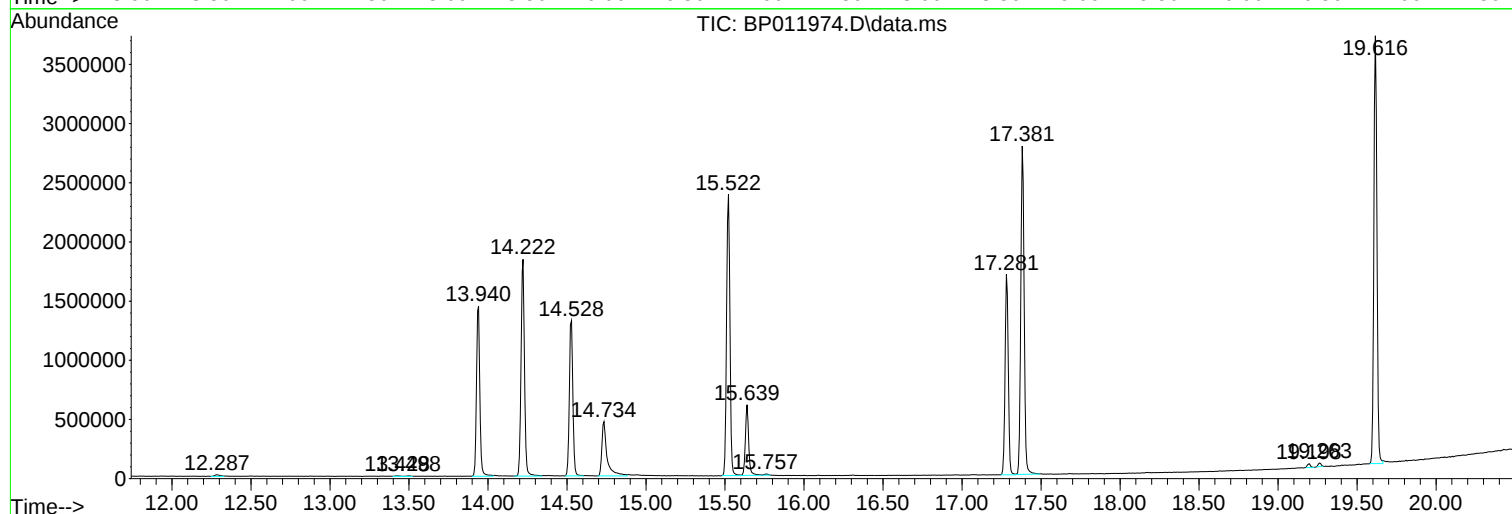
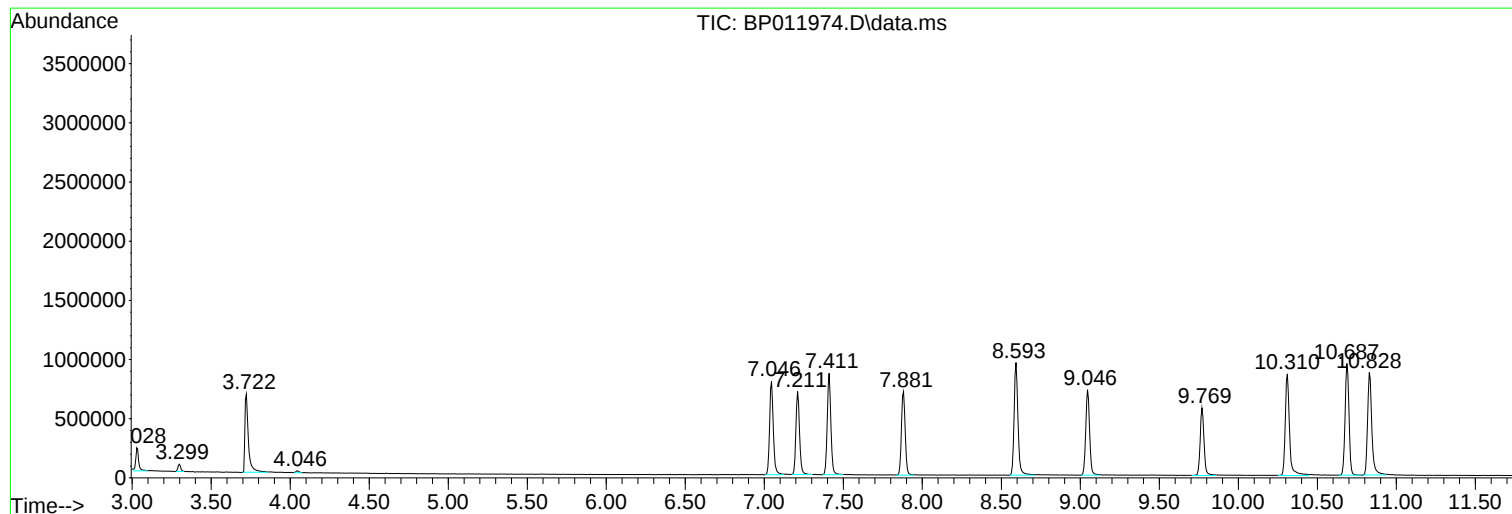
Sum of corrected areas: 49749476

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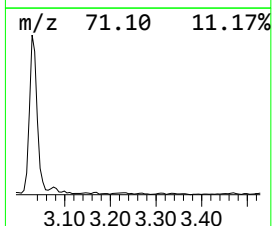
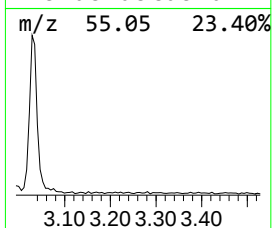
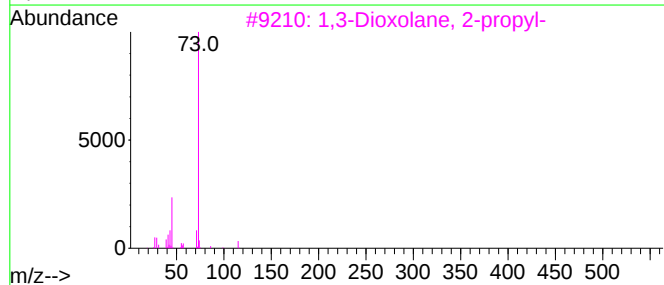
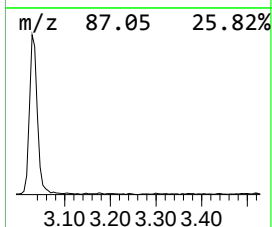
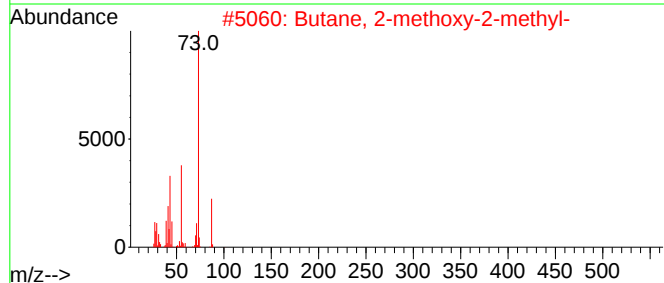
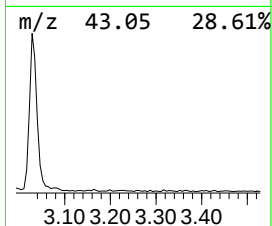
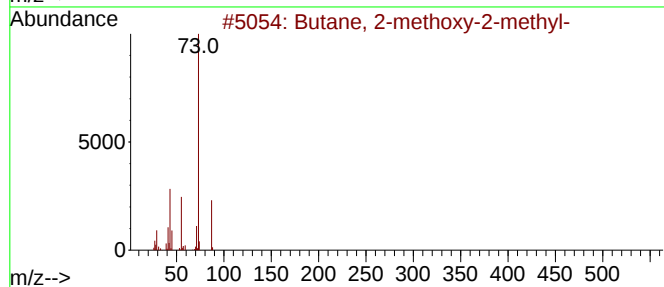
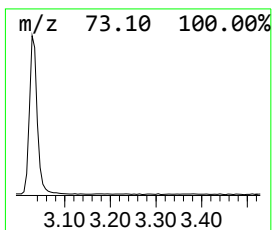
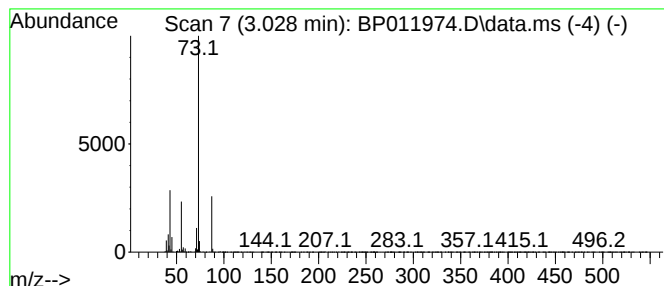
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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	4.18 ng/ul	241352	1,4-Dichlorobenzene-d4	7.881

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			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	23
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
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...	3.028	4.2	ng/u1	241352	1	7.881	1153550	20.0