

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011869.D
 Acq On : 03 Oct 2022 13:24
 Operator : CG/JU
 Sample : PB147928BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK928

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: BP011869.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	23	rVB	209554	307561	3.89%	0.440%
2	3.305	51	54	69	rVB	71416	104263	1.32%	0.149%
3	3.728	123	126	151	rBV	765158	1132001	14.31%	1.619%
4	4.052	178	181	191	rVB	16456	25638	0.32%	0.037%
5	4.281	216	220	225	rVB3	7277	11735	0.15%	0.017%
6	7.058	682	692	702	rBV	944379	1580541	19.98%	2.261%
7	7.222	714	720	733	rVB	861291	1353731	17.12%	1.937%
8	7.422	746	754	770	rBV	1016733	1663819	21.04%	2.380%
9	7.893	825	834	847	rBV	914362	1494210	18.89%	2.138%
10	8.604	944	955	973	rBV	1224694	2076507	26.25%	2.971%
11	9.057	1024	1032	1048	rBV	890098	1513878	19.14%	2.166%
12	9.781	1148	1155	1174	rBV	707002	1219951	15.42%	1.745%
13	10.322	1239	1247	1263	rBV	1137490	2039415	25.78%	2.918%
14	10.699	1301	1311	1322	rBV	1264677	2205760	27.89%	3.156%
15	10.840	1328	1335	1359	rBV	1151029	2151106	27.20%	3.077%
16	12.293	1576	1582	1592	rBV2	19211	40319	0.51%	0.058%
17	13.428	1769	1775	1783	rBV6	6901	12514	0.16%	0.018%
18	13.951	1857	1864	1885	rBV	2085333	3071941	38.84%	4.395%
19	14.228	1904	1911	1930	rBV2	2557752	3910952	49.45%	5.595%
20	14.539	1956	1964	1978	rBV	1935648	2965120	37.49%	4.242%
21	14.739	1990	1998	2022	rBV	799344	1468158	18.56%	2.100%
22	14.945	2031	2033	2045	rVB9	8255	16751	0.21%	0.024%
23	15.534	2125	2133	2143	rBV	3375596	4999687	63.21%	7.153%
24	15.651	2146	2153	2169	rVV	889621	1326715	16.77%	1.898%
25	15.775	2169	2174	2181	rVB	15284	30679	0.39%	0.044%
26	17.086	2393	2397	2402	rVB4	5089	9160	0.12%	0.013%
27	17.292	2425	2432	2442	rBV	2560432	3594135	45.44%	5.142%
28	17.392	2442	2449	2465	rBV2	3832196	5596343	70.75%	8.006%
29	19.204	2753	2757	2763	rBV2	51049	78211	0.99%	0.112%
30	19.275	2765	2769	2773	rVV	51691	75884	0.96%	0.109%
31	19.627	2823	2829	2837	rBV	5192443	6927924	87.59%	9.911%
32	21.374	3121	3126	3136	rBV	3199951	4334775	54.80%	6.202%
33	23.651	3507	3513	3530	rVB2	3763168	7909511	100.00%	11.316%
34	23.810	3534	3540	3554	rVB2	2297303	4649829	58.79%	6.652%

Sum of corrected areas: 69898724

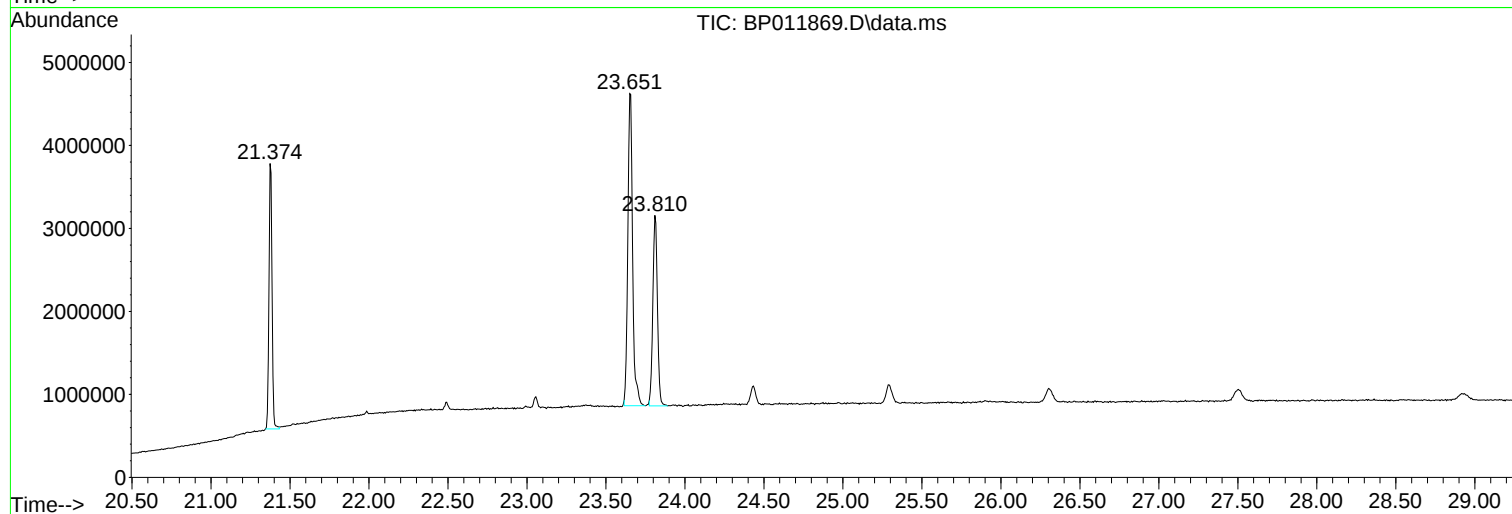
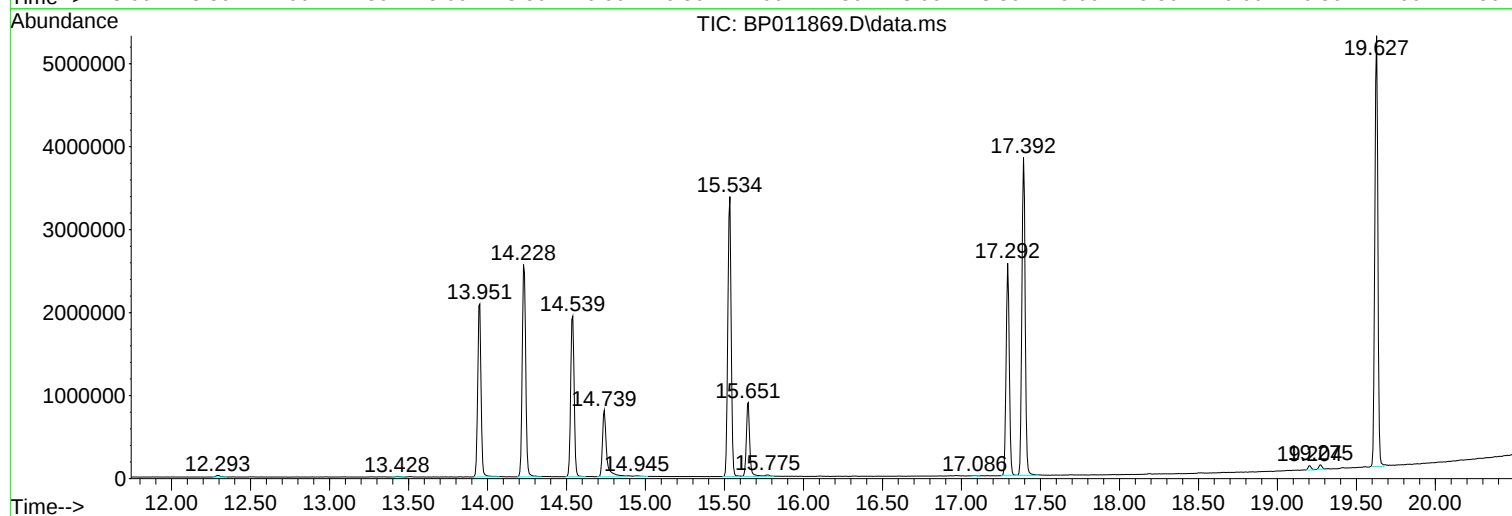
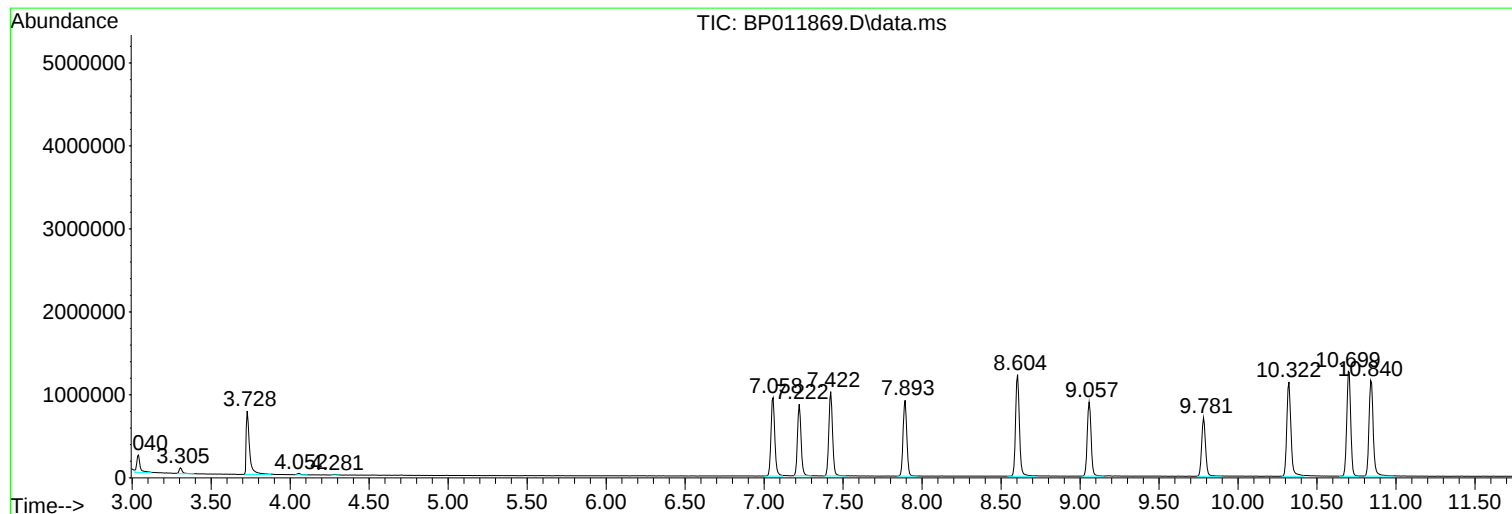
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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



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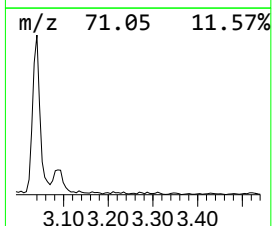
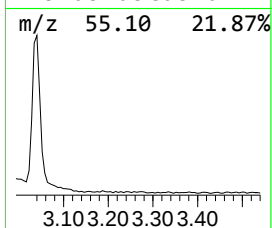
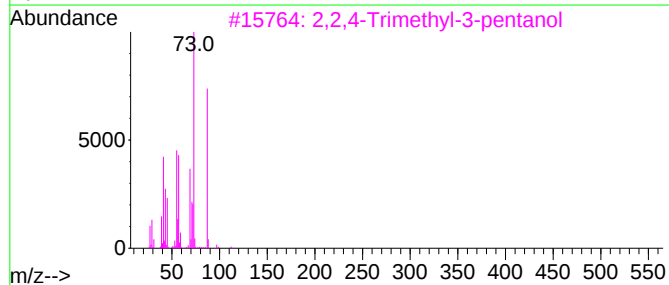
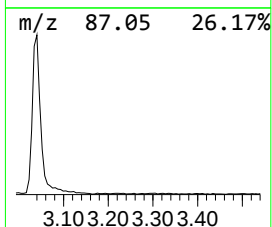
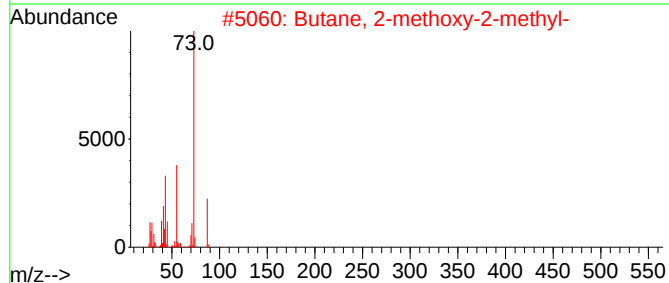
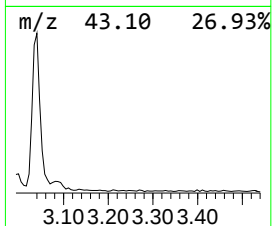
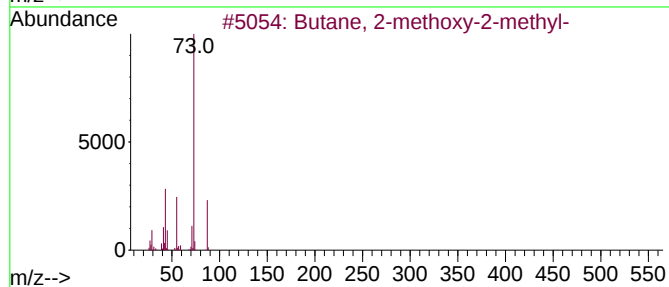
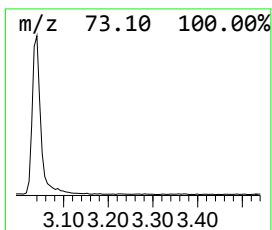
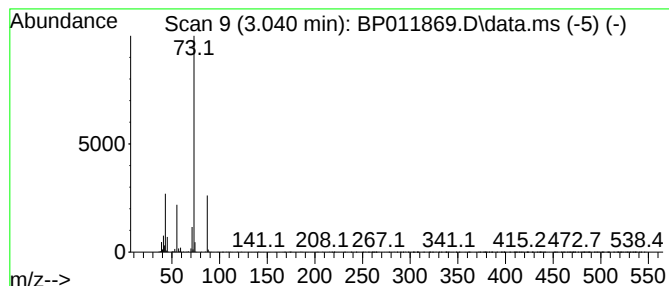
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Quant Title : SVOA CALIBRATION

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.
3.040	4.12 ng/ul	307561	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	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	43		
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23		
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.040	4.1	ng/u1	307561	1	7.893	1494210	20.0