

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011960.D
 Acq On : 06 Oct 2022 09:20
 Operator : CG/JU
 Sample : PB148128BL
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK128

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: BP011960.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	3	7	18	rVB	274423	344601	4.74%	0.484%
2	3.299	49	53	60	rVB	81438	105633	1.45%	0.148%
3	3.717	121	124	143	rBV	919617	1376820	18.95%	1.933%
4	4.040	176	179	186	rVB	18698	28852	0.40%	0.041%
5	7.040	682	689	702	rBV	1116051	1797650	24.75%	2.524%
6	7.211	711	718	731	rVB	956533	1554089	21.39%	2.182%
7	7.405	744	751	764	rBV	1136692	1893169	26.06%	2.658%
8	7.875	824	831	844	rBV	858877	1426708	19.64%	2.003%
9	8.587	945	952	974	rBV	1395334	2327066	32.04%	3.267%
10	9.046	1022	1030	1046	rBV	1035017	1747897	24.06%	2.454%
11	9.769	1144	1153	1169	rBV	796792	1377277	18.96%	1.934%
12	10.305	1236	1244	1266	rBV	1299041	2358586	32.47%	3.311%
13	10.687	1299	1309	1325	rVB	1251475	2127746	29.29%	2.987%
14	10.828	1325	1333	1354	rBV	1276480	2373585	32.68%	3.332%
15	12.287	1575	1581	1592	rVB2	19630	45105	0.62%	0.063%
16	13.422	1766	1774	1779	rBV	21117	36043	0.50%	0.051%
17	13.934	1855	1861	1876	rBV	2290444	3306696	45.52%	4.642%
18	14.216	1902	1909	1924	rBV	2769197	4266922	58.74%	5.990%
19	14.522	1954	1961	1973	rBV	1850684	2835714	39.04%	3.981%
20	14.728	1988	1996	2020	rBV	807995	1591451	21.91%	2.234%
21	14.934	2028	2031	2037	rVB7	6642	12806	0.18%	0.018%
22	15.522	2124	2131	2144	rBV	3721946	5480500	75.45%	7.694%
23	15.640	2144	2151	2168	rVV	943449	1412326	19.44%	1.983%
24	15.757	2168	2171	2177	rVB4	17505	29129	0.40%	0.041%
25	16.304	2259	2264	2269	rBV8	5248	8804	0.12%	0.012%
26	16.957	2372	2375	2380	rBV7	5110	8793	0.12%	0.012%
27	17.281	2423	2430	2440	rBV	2340516	3356384	46.21%	4.712%
28	17.381	2440	2447	2464	rVV2	4203116	6066453	83.51%	8.517%
29	19.192	2751	2755	2759	rBV3	51503	70921	0.98%	0.100%
30	19.263	2764	2767	2772	rVB	44939	59085	0.81%	0.083%
31	19.616	2821	2827	2845	rBV	5433415	7193483	99.03%	10.099%
32	21.369	3120	3125	3135	rBV	2796072	3685597	50.74%	5.174%
33	23.639	3504	3511	3525	rVB2	3740064	7263948	100.00%	10.198%
34	23.798	3532	3538	3549	rVB2	1824784	3660537	50.39%	5.139%

Sum of corrected areas: 71230376

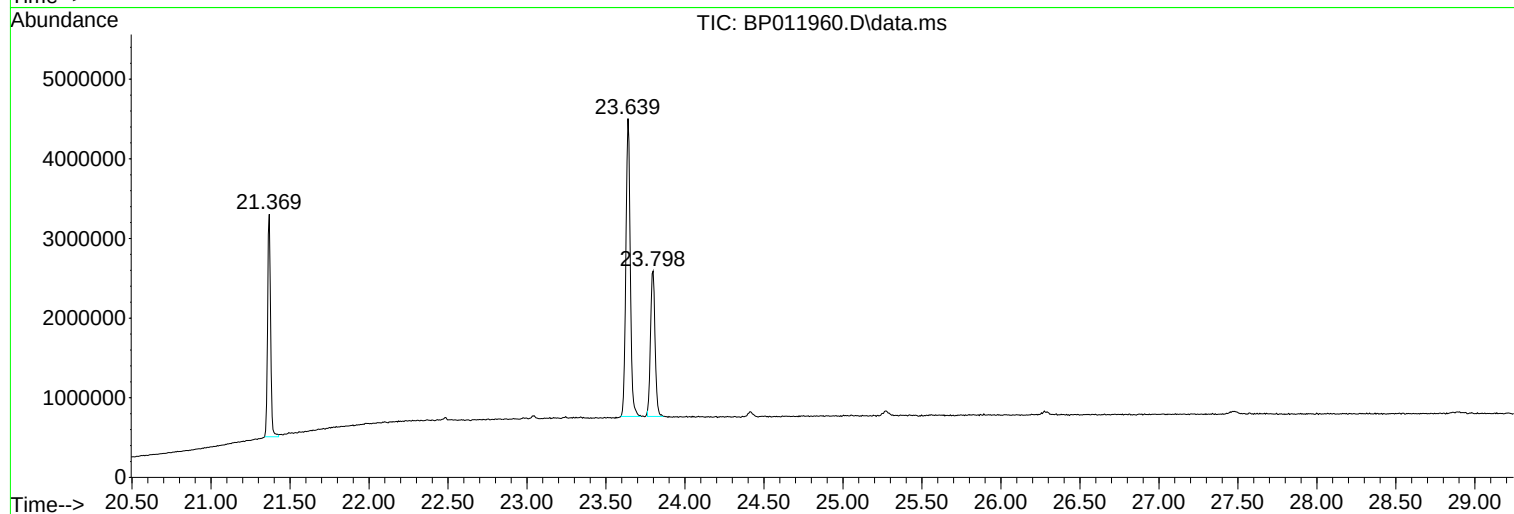
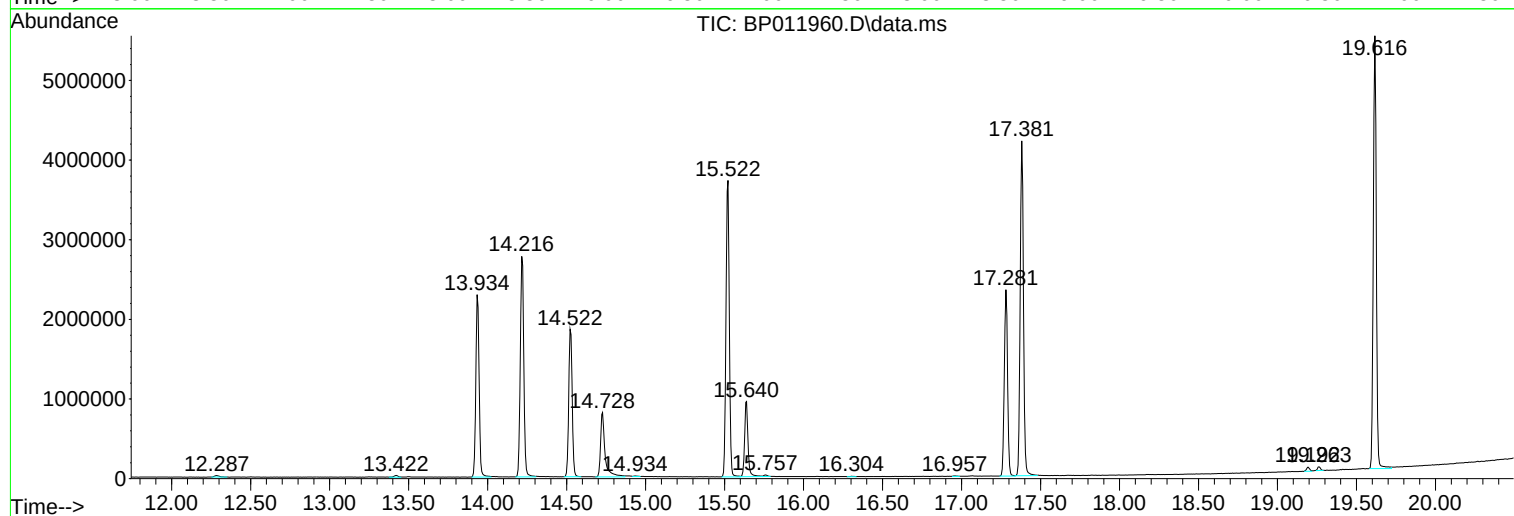
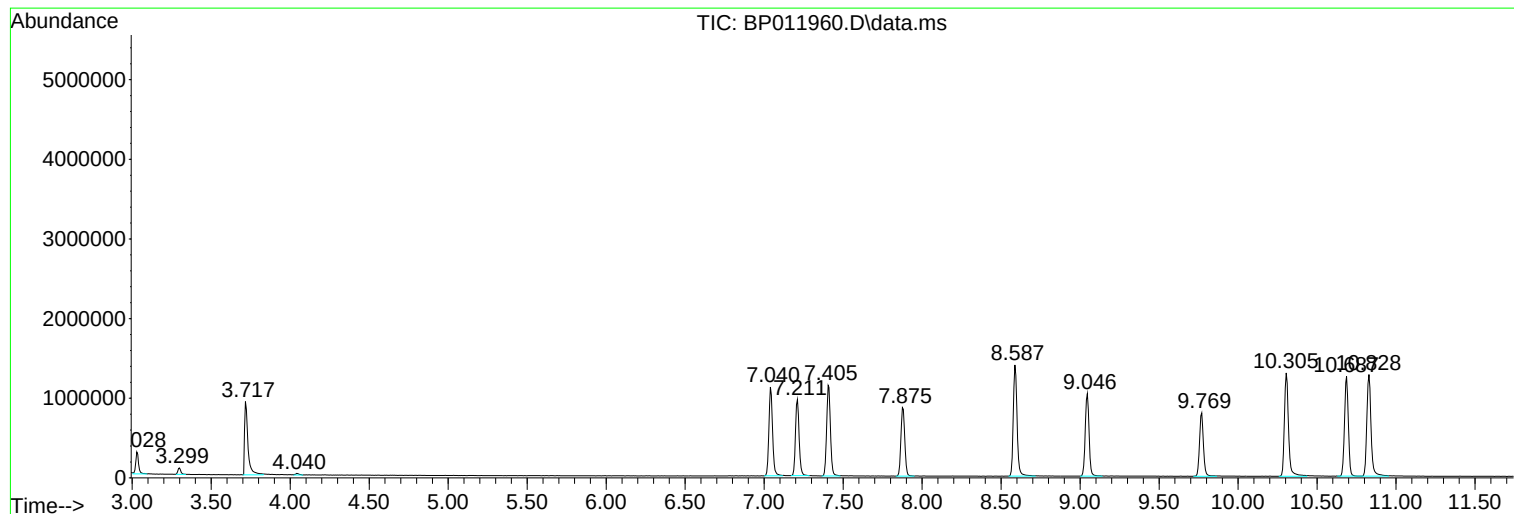
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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



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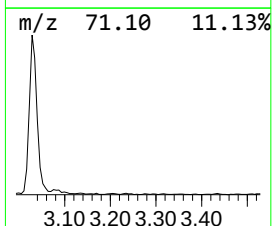
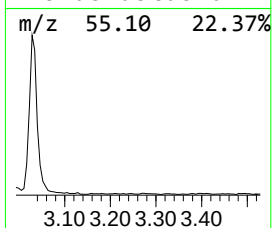
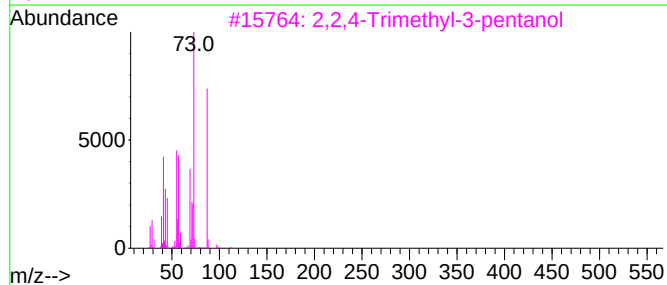
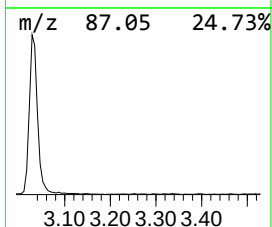
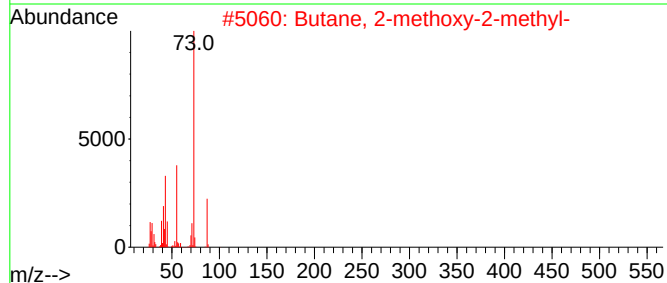
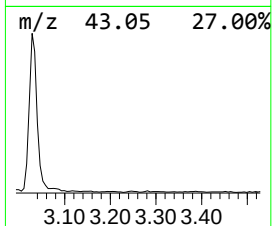
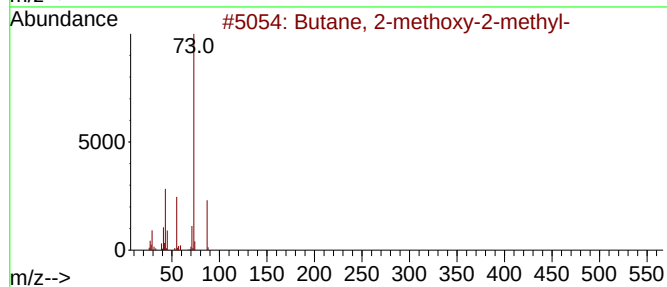
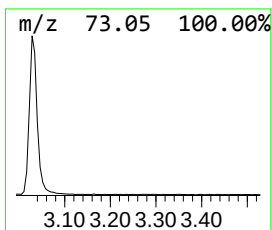
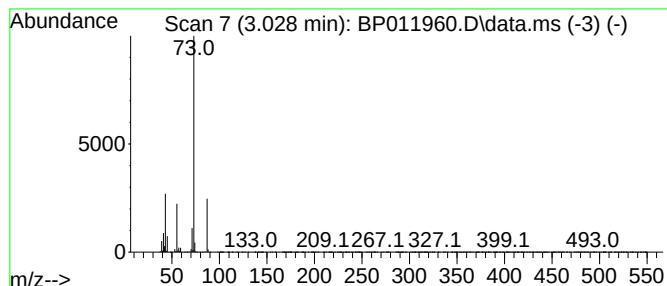
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Quant Title : SVOA CALIBRATION

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.
3.028	4.83 ng/ul	344601	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	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40		
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12		
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	4.8	ng/u1	344601	1	7.881	1426710	20.0