

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
 Data File : BP021656.D
 Acq On : 27 Aug 2024 04:03
 Operator : RC/JU
 Sample : PB162959BL
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK959

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-BP082324.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021656.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.022	3	6	18	rVB	159134	202664	2.16%	0.215%
2	3.122	19	23	28	rVB	20545	28013	0.30%	0.030%
3	3.275	45	49	57	rVB	98831	130205	1.39%	0.138%
4	3.687	115	119	135	rBV	1213385	1760135	18.80%	1.864%
5	4.022	172	176	183	rVB	24118	34741	0.37%	0.037%
6	6.958	668	675	686	rBV	1345418	2294747	24.51%	2.430%
7	7.128	696	704	713	rVB	1236834	1902513	20.32%	2.014%
8	7.328	731	738	750	rBV	1289713	2158591	23.05%	2.286%
9	7.793	810	817	825	rBV	1182028	1928040	20.59%	2.041%
10	8.487	927	935	953	rBV	1755297	2985808	31.89%	3.161%
11	8.934	1004	1011	1024	rBV	1325076	2259901	24.14%	2.393%
12	9.657	1125	1134	1146	rBV	1018075	1773755	18.94%	1.878%
13	10.199	1217	1226	1246	rBV	1529468	2842396	30.36%	3.010%
14	10.581	1283	1291	1304	rVB	1635524	2847133	30.41%	3.015%
15	10.710	1304	1313	1332	rBV	1544325	2777207	29.66%	2.941%
16	11.381	1420	1427	1431	rBV9	6624	15422	0.16%	0.016%
17	12.175	1552	1562	1572	rBV2	21544	50971	0.54%	0.054%
18	12.410	1595	1602	1608	rVB	9433	17631	0.19%	0.019%
19	13.398	1764	1770	1775	rBV9	5660	13370	0.14%	0.014%
20	13.828	1837	1843	1858	rBV	3007507	4494961	48.00%	4.759%
21	14.116	1885	1892	1910	rBV2	3212078	5200746	55.54%	5.507%
22	14.428	1937	1945	1956	rBV	2554025	4053286	43.29%	4.292%
23	14.622	1971	1978	1997	rBV	1433958	2504269	26.74%	2.652%
24	14.851	2013	2017	2024	rVB10	10989	21834	0.23%	0.023%
25	15.439	2109	2117	2125	rBV	4580976	6652494	71.05%	7.044%
26	15.551	2130	2136	2152	rVV	1370905	1958031	20.91%	2.073%
27	15.687	2154	2159	2163	rVB3	18888	33274	0.36%	0.035%
28	15.951	2199	2204	2210	rVB10	6107	9399	0.10%	0.010%
29	16.239	2249	2253	2258	rVB2	16155	24453	0.26%	0.026%
30	16.904	2362	2366	2369	rVB6	9643	13051	0.14%	0.014%
31	17.010	2378	2384	2393	rBV4	9957	26106	0.28%	0.028%
32	17.228	2414	2421	2430	rBV	3511464	5141784	54.91%	5.444%
33	17.333	2431	2439	2449	rBV2	4572722	7535955	80.48%	7.979%
34	19.251	2761	2765	2771	rBV3	67106	102195	1.09%	0.108%
35	19.328	2774	2778	2784	rVB	55142	83439	0.89%	0.088%
36	19.704	2836	2842	2852	rVB	6788775	9198082	98.23%	9.739%

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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

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Title : SVOA CALIBRATION

37	21.686	3173	3179	3192	rVB	3614955	5863756	62.62%	6.209%
38	24.874	3711	3721	3738	rVB3	3467648	9363546	100.00%	9.914%
39	25.104	3749	3760	3774	rVB	2097759	6142599	65.60%	6.504%

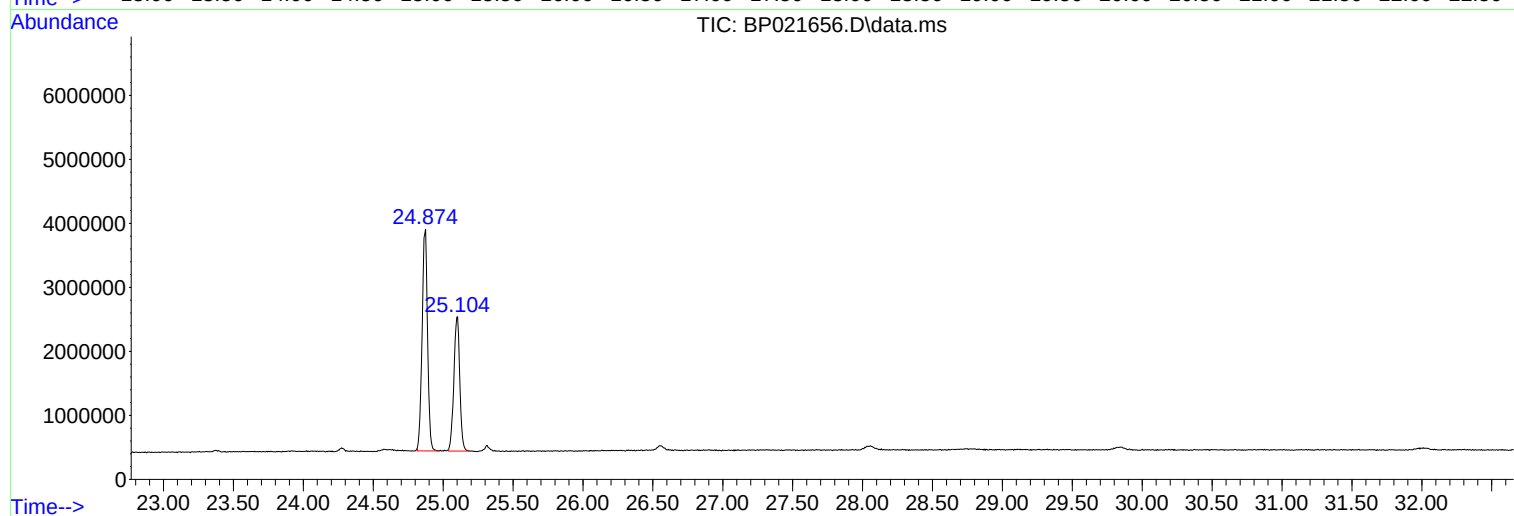
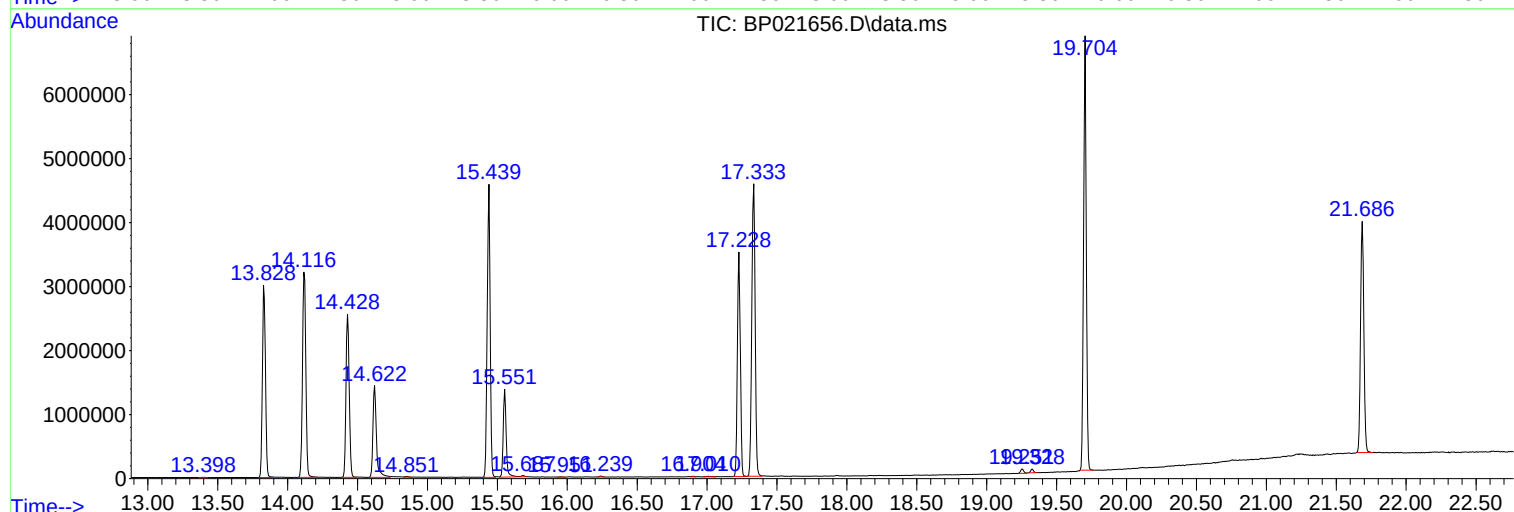
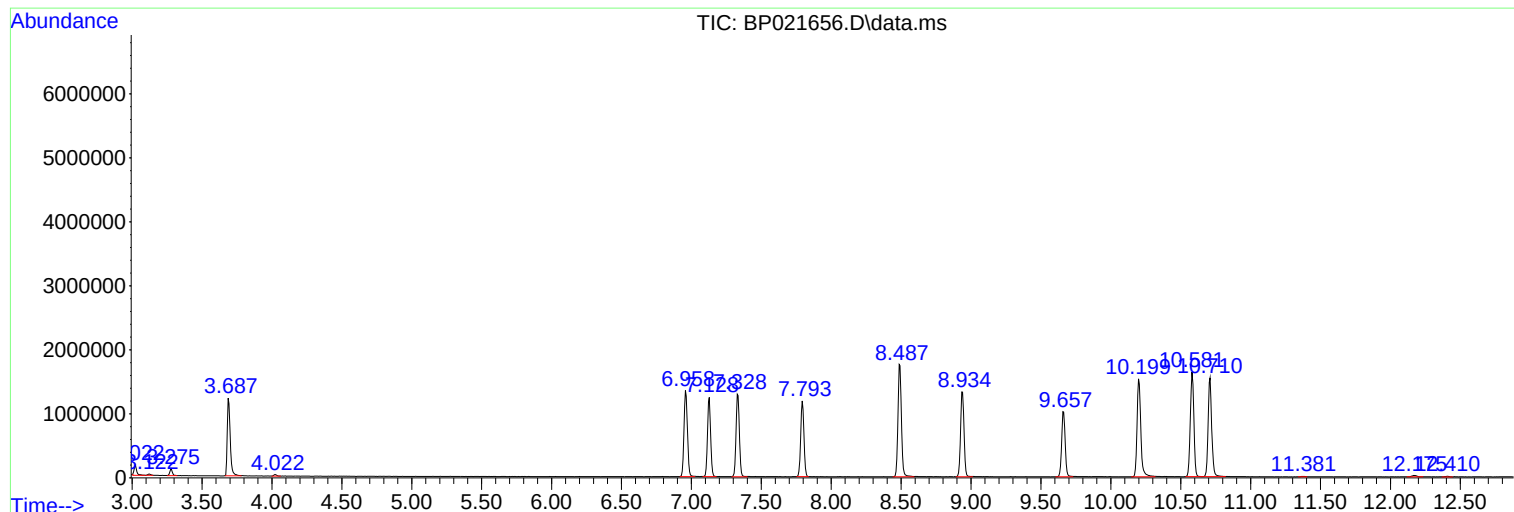
Sum of corrected areas: 94446503

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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



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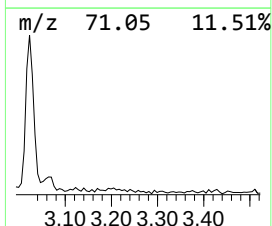
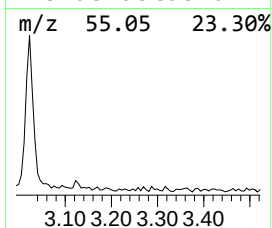
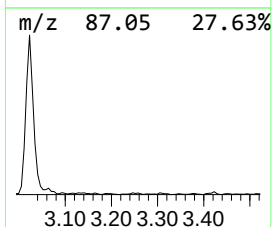
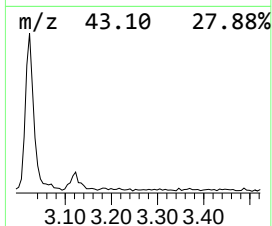
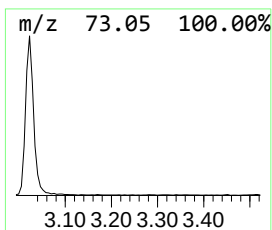
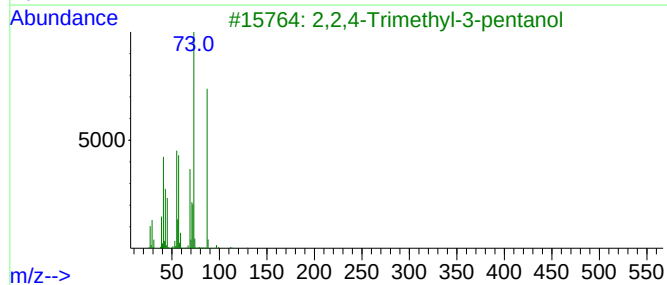
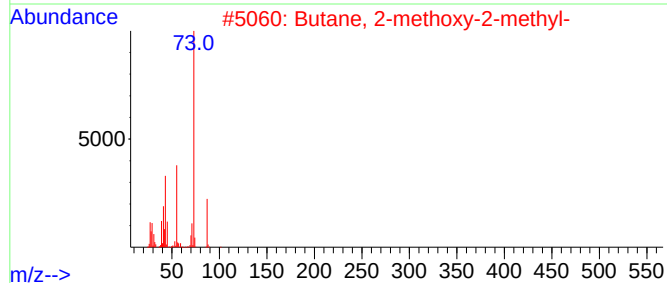
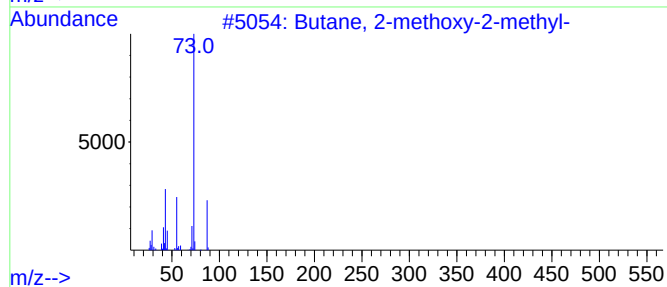
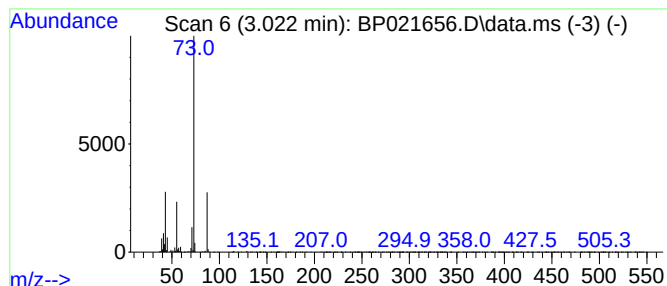
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.022	2.10 ng/ul	202664	1,4-Dichlorobenzene-d4	7.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
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	17
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.022	2.1	ng/u1	202664	1	7.793	1928040	20.0