

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
 Data File : BP021657.D
 Acq On : 27 Aug 2024 04:46
 Operator : RC/JU
 Sample : PB162966BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK966

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: BP021657.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	160926	220592	2.51%	0.256%
2	3.122	20	23	27	rVB2	24986	33318	0.38%	0.039%
3	3.281	46	50	62	rVB	106651	147281	1.68%	0.171%
4	3.693	115	120	137	rBV	1191125	1776974	20.22%	2.062%
5	4.022	172	176	183	rVB	25335	38677	0.44%	0.045%
6	6.958	668	675	684	rBV	1410735	2248567	25.58%	2.609%
7	7.128	697	704	713	rVB	1203219	1856160	21.12%	2.154%
8	7.334	731	739	750	rBV	1281405	2139686	24.35%	2.483%
9	7.793	810	817	827	rBV	1171637	1904606	21.67%	2.210%
10	8.493	928	936	952	rBV	1678285	2813306	32.01%	3.264%
11	8.946	1005	1013	1023	rBV	1247659	2115465	24.07%	2.455%
12	9.663	1126	1135	1149	rBV	954637	1660284	18.89%	1.926%
13	10.199	1217	1226	1243	rBV	1391566	2662803	30.30%	3.090%
14	10.581	1283	1291	1304	rBV	1503494	2620807	29.82%	3.041%
15	10.710	1305	1313	1333	rVB	1463291	2537036	28.87%	2.944%
16	12.175	1554	1562	1571	rBV3	23648	51086	0.58%	0.059%
17	12.404	1592	1601	1606	rBV	19565	31266	0.36%	0.036%
18	13.840	1836	1845	1861	rBV	2858389	4067464	46.28%	4.720%
19	14.128	1886	1894	1907	rBV	3044227	4708710	53.58%	5.464%
20	14.439	1939	1947	1959	rBV	2208091	3565770	40.57%	4.137%
21	14.528	1959	1962	1968	rVB8	6555	9694	0.11%	0.011%
22	14.628	1972	1979	1996	rBV	1275904	2140808	24.36%	2.484%
23	14.863	2014	2019	2024	rVB9	8664	15067	0.17%	0.017%
24	15.445	2111	2118	2130	rBV	3959713	5990541	68.16%	6.951%
25	15.557	2131	2137	2154	rVV	1189963	1731037	19.70%	2.009%
26	15.686	2154	2159	2166	rVB2	18251	36768	0.42%	0.043%
27	16.239	2250	2253	2258	rVB5	13493	18688	0.21%	0.022%
28	16.904	2363	2366	2372	rVB8	8999	15138	0.17%	0.018%
29	17.022	2382	2386	2391	rVB4	7037	11726	0.13%	0.014%
30	17.233	2415	2422	2432	rBV	2852671	4420812	50.30%	5.130%
31	17.339	2433	2440	2453	rVV2	4300829	6672139	75.91%	7.742%
32	18.169	2579	2581	2586	rBV6	6008	10253	0.12%	0.012%
33	19.257	2762	2766	2772	rVB2	59519	78105	0.89%	0.091%
34	19.327	2775	2778	2782	rVB	45711	63572	0.72%	0.074%
35	19.704	2835	2842	2854	rBV	5630829	8351987	95.03%	9.691%
36	21.686	3173	3179	3190	rBV2	2928746	5145629	58.55%	5.971%

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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-BP082324.MA.M
Title : SVOA CALIBRATION

37	24.868	3710	3720	3737	rVB	3305893	8788986	100.00%	10.198%
38	25.098	3750	3759	3774	rVB	1864586	5482296	62.38%	6.361%

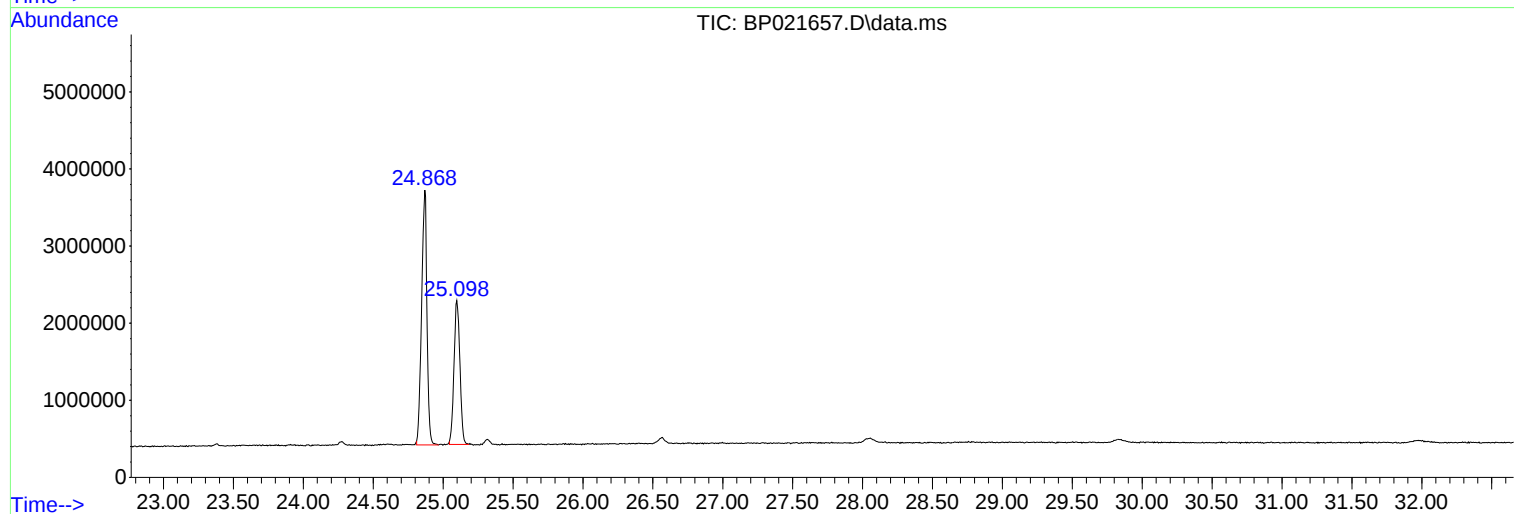
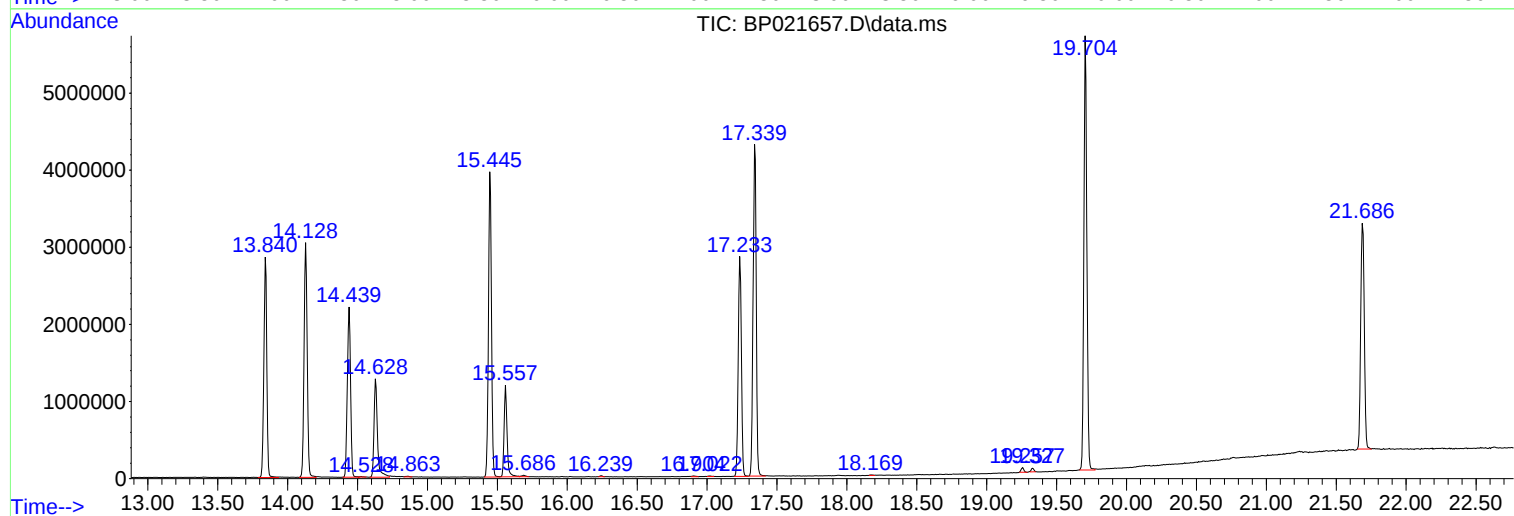
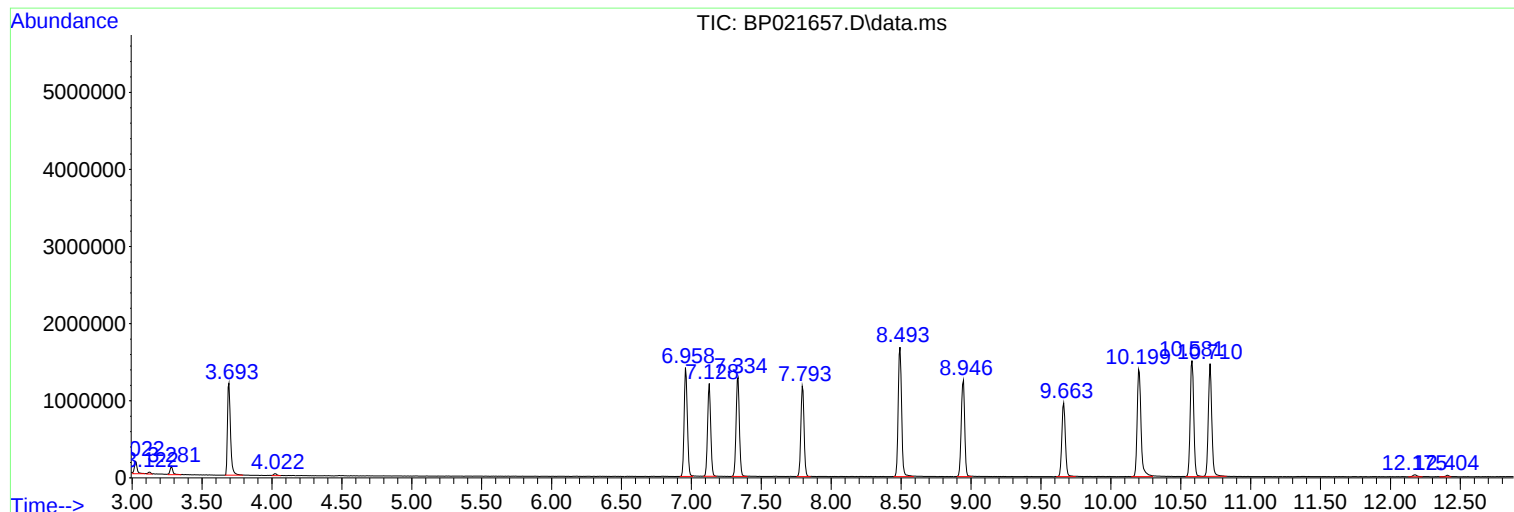
Sum of corrected areas: 86183104

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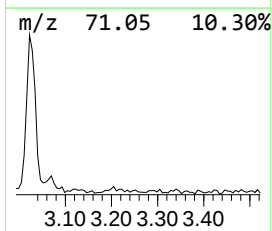
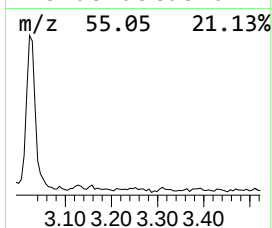
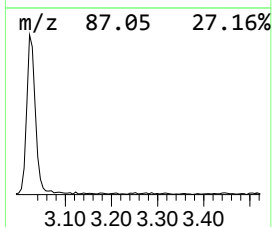
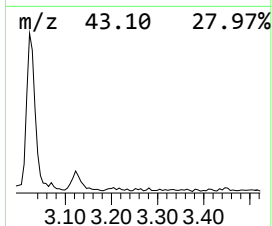
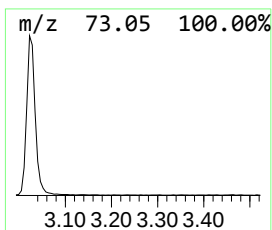
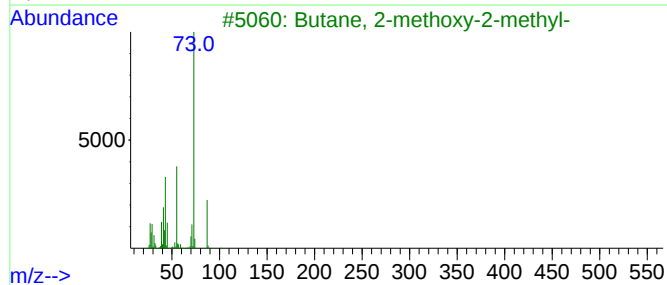
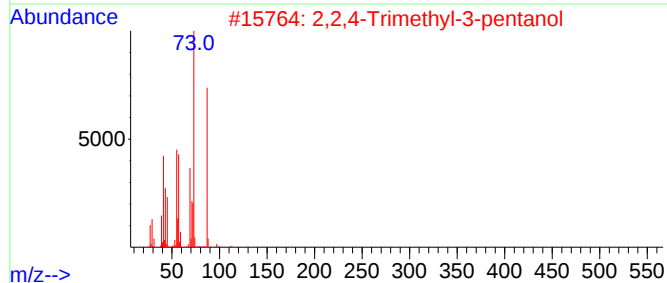
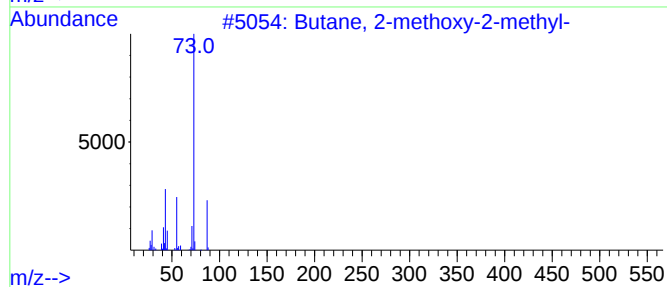
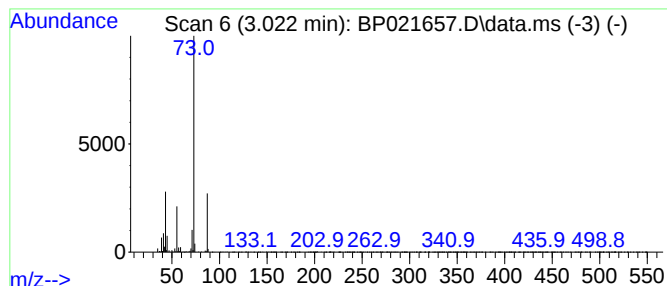
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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.
3.022	2.32 ng/ul	220592	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	86
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	64
3			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
4			Myo-Inositol, 4-C-methyl-	194	C7H14O6	000472-95-7	36
5			Myo-Inositol, 2-C-methyl-	194	C7H14O6	000472-96-8	36



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.022	2.3	ng/u1	220592	1	7.793	1904610	20.0