

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091324\
 Data File : BP021905.D
 Acq On : 13 Sep 2024 12:26
 Operator : RC/JU
 Sample : PB163298TB 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLEB298

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-BP091124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021905.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.010	3	4	11	rVB	68200	70687	1.62%	0.187%
2	3.110	18	21	26	rVB	10983	14275	0.33%	0.038%
3	3.269	44	48	57	rVB	40181	52541	1.20%	0.139%
4	3.681	114	118	143	rBV	413009	651067	14.90%	1.719%
5	4.010	170	174	179	rVB3	7063	9865	0.23%	0.026%
6	6.934	663	671	684	rBV	387413	647229	14.81%	1.708%
7	7.099	690	699	709	rBV	303132	486882	11.14%	1.285%
8	7.299	724	733	742	rBV	417175	646797	14.80%	1.707%
9	7.763	804	812	821	rBV	345870	544594	12.46%	1.437%
10	8.457	922	930	948	rBV	525533	817987	18.72%	2.159%
11	8.904	997	1006	1013	rBV	394921	662074	15.15%	1.748%
12	9.616	1120	1127	1142	rBV	391190	684606	15.67%	1.807%
13	10.151	1211	1218	1231	rBV	669976	1130945	25.89%	2.985%
14	10.534	1275	1283	1290	rBV	585657	1028444	23.54%	2.715%
15	10.657	1294	1304	1320	rBV	622996	1105436	25.30%	2.918%
16	12.116	1546	1552	1560	rBV5	8828	17524	0.40%	0.046%
17	13.769	1826	1833	1848	rBV	1264330	1939020	44.38%	5.118%
18	14.057	1874	1882	1898	rBV	1433890	2126321	48.67%	5.612%
19	14.369	1927	1935	1945	rBV	996357	1536292	35.16%	4.055%
20	14.575	1963	1970	1985	rBV	539908	931950	21.33%	2.460%
21	15.375	2098	2106	2113	rBV	1894285	2872620	65.75%	7.582%
22	15.492	2120	2126	2142	rBV	541731	807228	18.48%	2.131%
23	15.610	2142	2146	2155	rVB5	9109	19232	0.44%	0.051%
24	16.169	2237	2241	2245	rBV5	7777	11650	0.27%	0.031%
25	17.157	2403	2409	2418	rBV	1307722	2064982	47.26%	5.451%
26	17.263	2420	2427	2446	rVB	2409916	3380275	77.37%	8.922%
27	19.186	2749	2754	2759	rBV2	27901	37820	0.87%	0.100%
28	19.263	2763	2767	2770	rVB	20808	29578	0.68%	0.078%
29	19.633	2824	2830	2840	rBV	3044313	4220081	96.59%	11.139%
30	21.598	3157	3164	3170	rBV2	1526999	2477550	56.71%	6.540%
31	24.686	3680	3689	3707	rVB2	1575900	4369018	100.00%	11.532%
32	24.921	3721	3729	3741	rVB	1008687	2491099	57.02%	6.575%

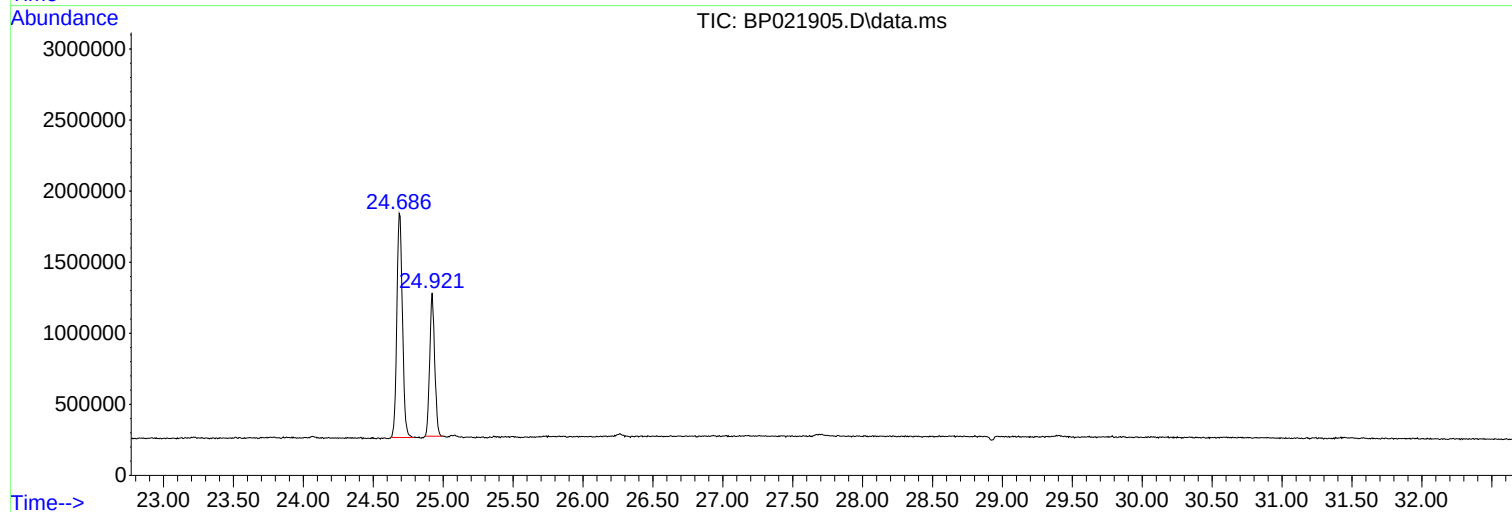
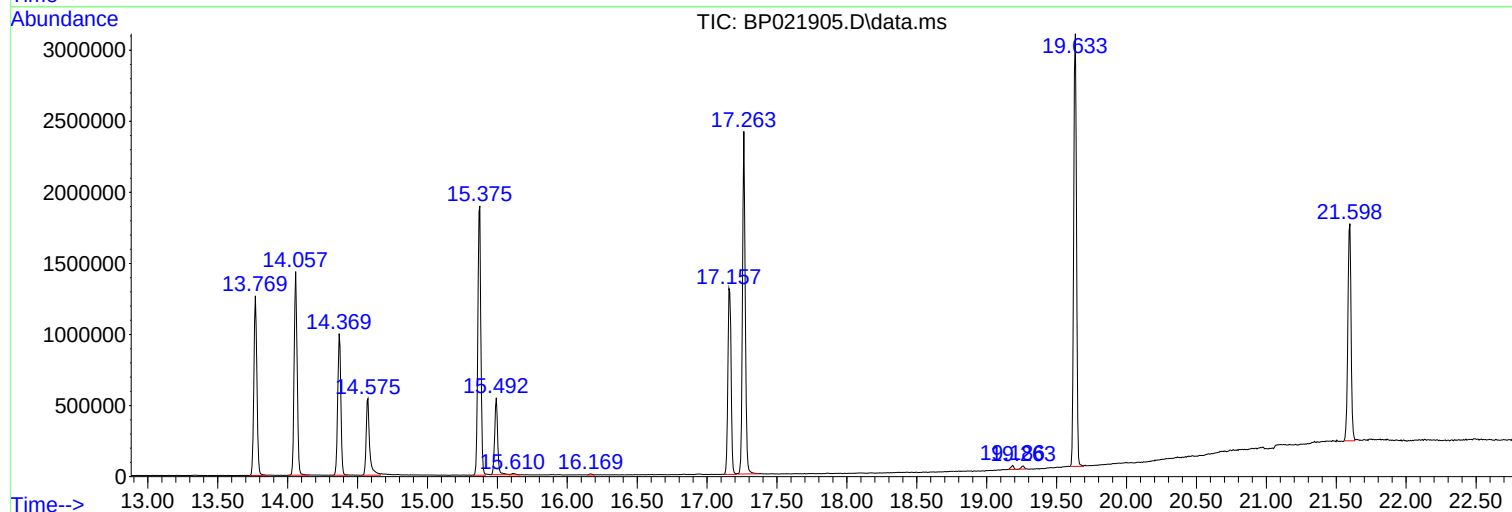
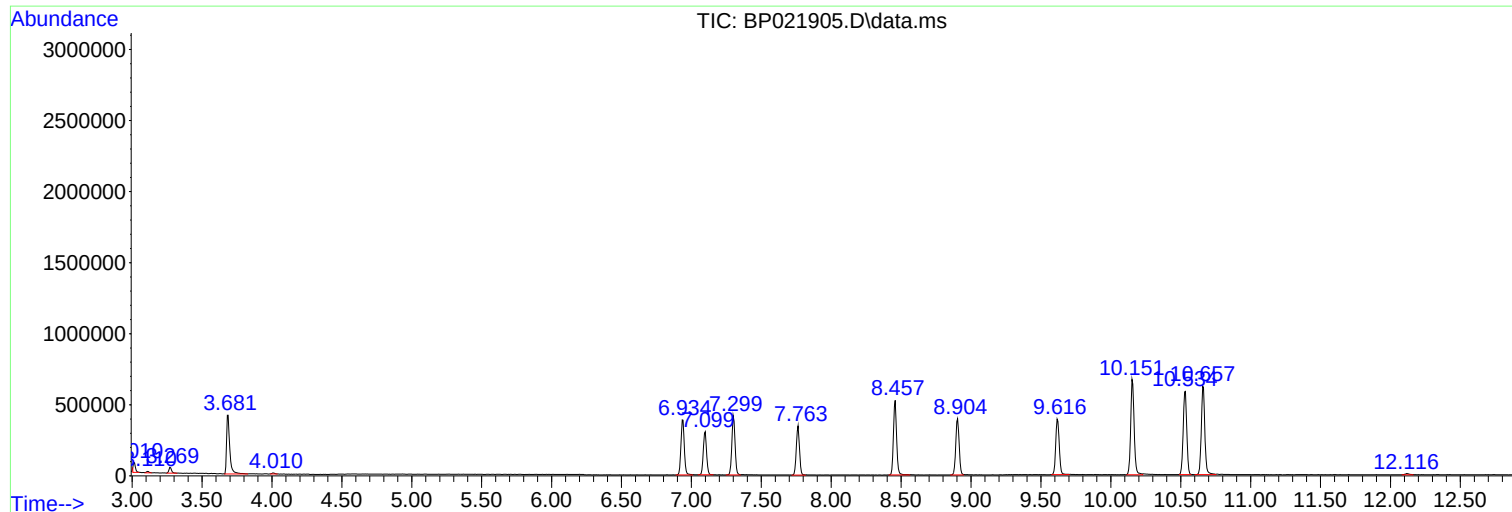
Sum of corrected areas: 37885669

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091124.MA.M
Quant Title : SVOA CALIBRATION

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



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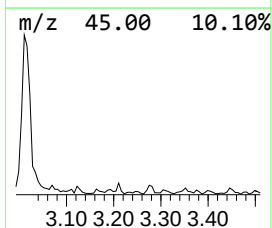
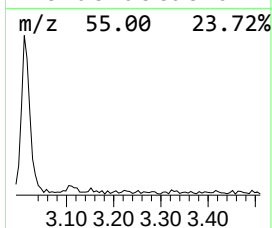
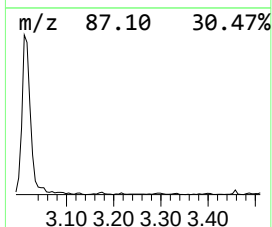
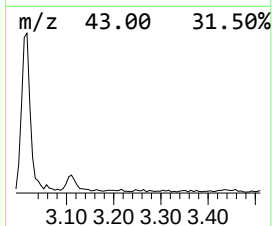
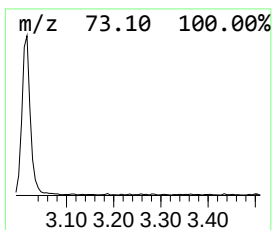
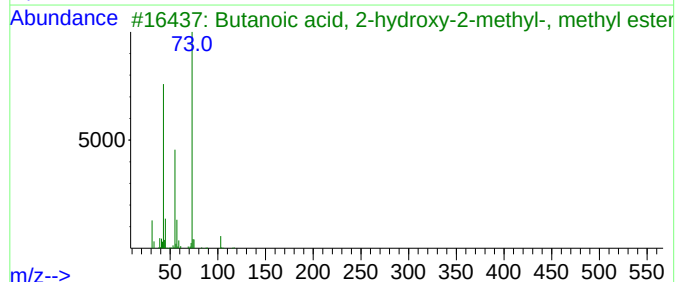
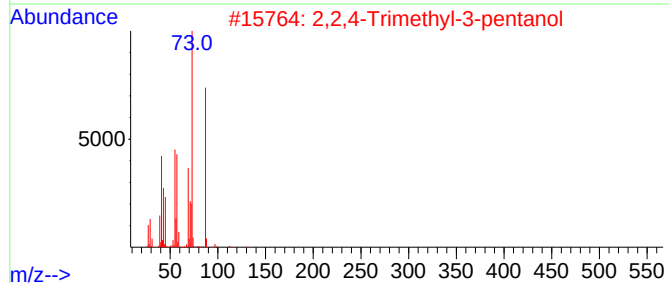
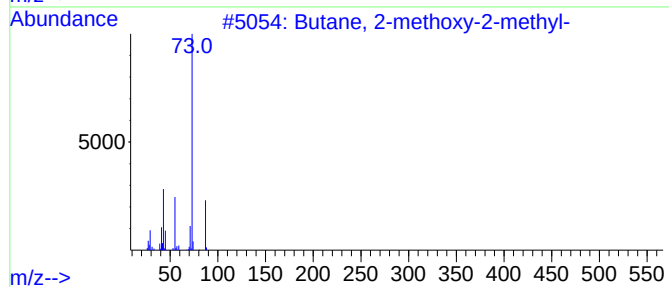
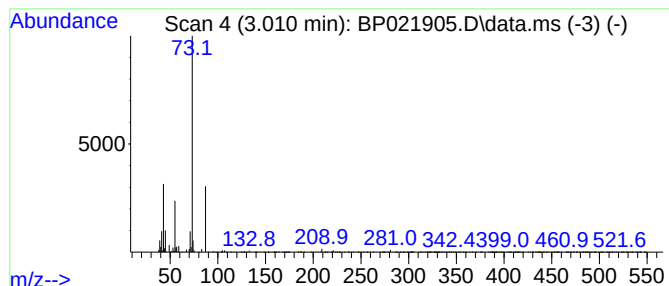
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091124.MA.M
Quant Title : SVOA CALIBRATION

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.
3.010	2.60 ng/ul	70687	1,4-Dichlorobenzene-d4	7.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Butanoic acid, 2-hydroxy-2-methyl...	132	C6H12O3	032793-34-3	25
4			1-(2-Methoxyethoxy)-2-methyl-2-p...	162	C8H18O3	1000364-24-4	23
5			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	17



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.010	2.6	ng/ul	70687	1	7.763	544594	20.0