

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022522\
 Data File : BN018725.D
 Acq On : 25 Feb 2022 11:42
 Operator : CG/JU
 Sample : PB142908BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK908

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_N\Methods\SFAM-EPA-BN021622.M
 Title : SVOA CALIBRATION

Signal : TIC: BN018725.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.023	3	6	20	rVB	117380	189342	2.52%	0.228%
2	3.287	47	51	62	rVB	116054	159507	2.13%	0.192%
3	3.699	117	121	145	rBV	1377021	2152142	28.70%	2.596%
4	4.058	177	182	189	rVB2	21758	35610	0.47%	0.043%
5	5.070	351	354	359	rVB5	5900	8333	0.11%	0.010%
6	7.064	686	693	704	rBV	1616812	2607922	34.78%	3.146%
7	7.234	715	722	733	rVB	1304138	2055126	27.41%	2.479%
8	7.434	749	756	768	rBV	1601117	2601103	34.69%	3.137%
9	7.911	830	837	844	rBV	1104206	1755505	23.41%	2.117%
10	8.611	949	956	975	rBV	1984551	3143565	41.92%	3.792%
11	9.069	1026	1034	1045	rBV	1538380	2466837	32.90%	2.975%
12	9.793	1149	1157	1166	rBV	1264691	2035303	27.14%	2.455%
13	10.334	1239	1249	1266	rBV	1907619	3153361	42.05%	3.803%
14	10.716	1306	1314	1324	rVB	1484954	2461995	32.83%	2.970%
15	10.846	1326	1336	1349	rBV	1857838	3144216	41.93%	3.792%
16	12.304	1576	1584	1590	rBV2	32276	53630	0.72%	0.065%
17	13.522	1781	1791	1795	rVB2	5554	10989	0.15%	0.013%
18	13.951	1857	1864	1870	rBV	3298501	4460971	59.49%	5.381%
19	13.993	1870	1871	1880	rVB7	10006	12344	0.16%	0.015%
20	14.240	1905	1913	1924	rBV	3710646	5367871	71.58%	6.474%
21	14.546	1958	1965	1977	rBV	2152237	3127292	41.70%	3.772%
22	14.722	1988	1995	2008	rBV	1446956	2015068	26.87%	2.430%
23	15.534	2126	2133	2145	rBV	4439888	6406681	85.43%	7.727%
24	15.645	2145	2152	2165	rVV	1430619	1874322	24.99%	2.261%
25	15.775	2169	2174	2180	rVB	25546	37070	0.49%	0.045%
26	17.075	2388	2395	2401	rBV2	9628	17017	0.23%	0.021%
27	17.287	2424	2431	2440	rBV	2420295	3527008	47.03%	4.254%
28	17.387	2440	2448	2460	rBV2	4649755	6857200	91.44%	8.271%
29	19.239	2758	2763	2769	rVB2	60260	77134	1.03%	0.093%
30	19.310	2770	2775	2780	rVV	57009	76356	1.02%	0.092%
31	19.675	2831	2837	2848	rBV	5589414	7498981	100.00%	9.045%
32	20.269	2934	2938	2941	rBV5	5792	9404	0.13%	0.011%
33	21.463	3135	3141	3155	rBV	2621748	3338598	44.52%	4.027%
34	23.710	3514	3523	3540	rBV2	3412637	6916666	92.23%	8.343%
35	23.863	3541	3549	3560	rVB2	1527500	3254108	43.39%	3.925%

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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_N\Methods\SFAM-EPA-BN021622.M
Title : SVOA CALIBRATION

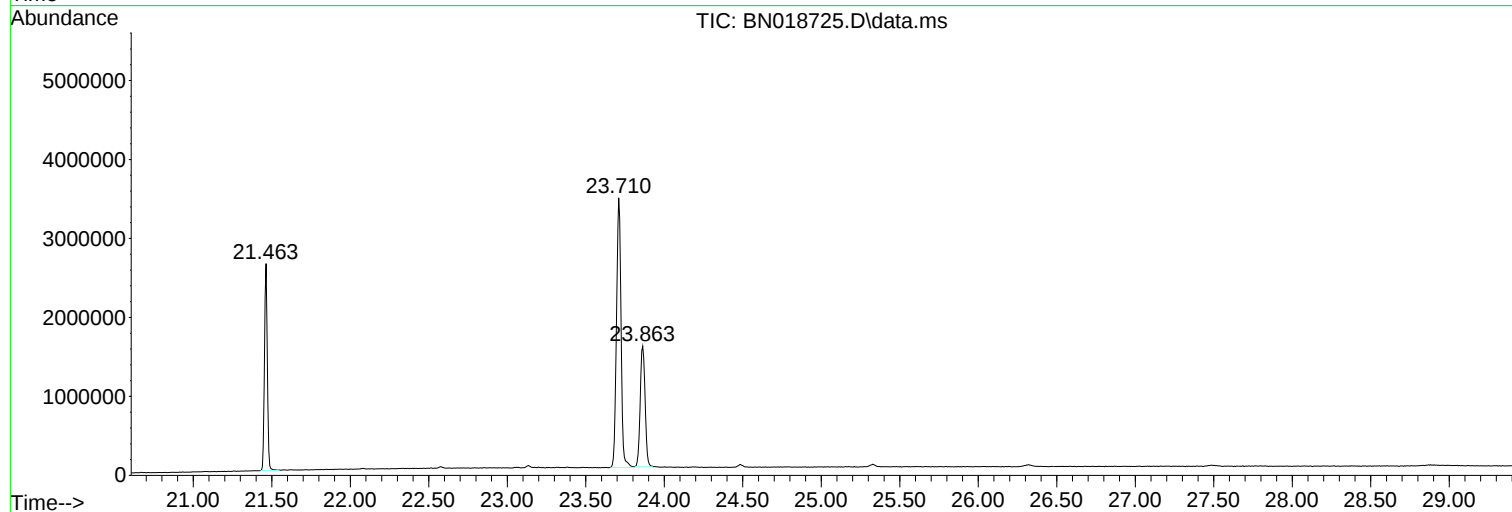
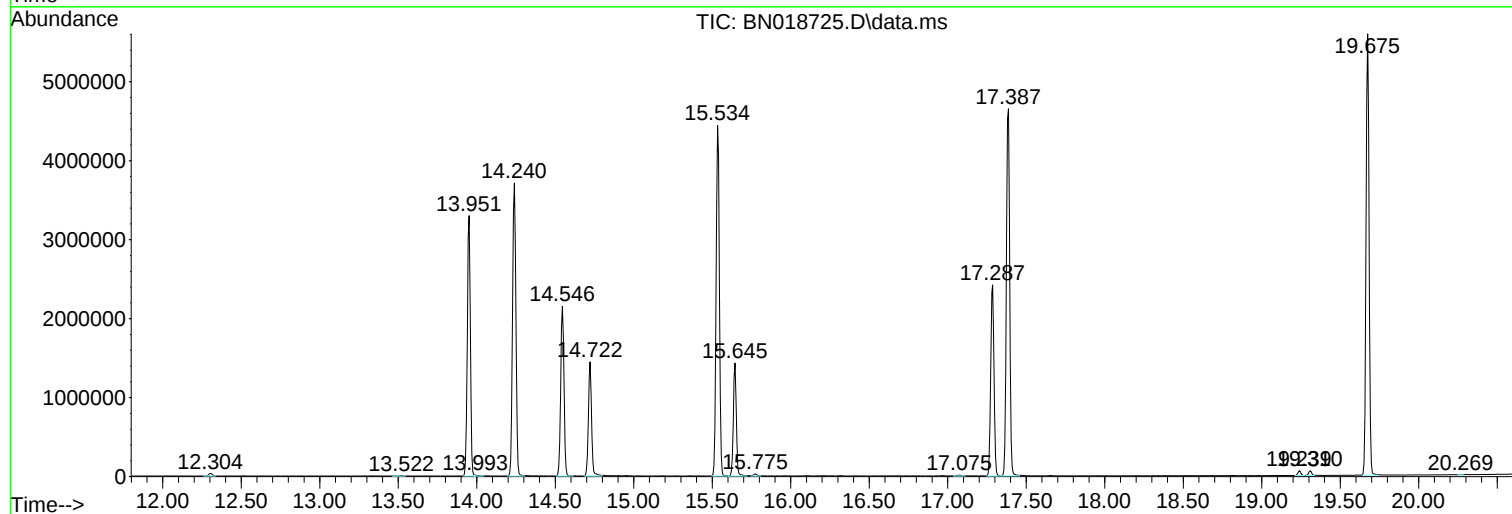
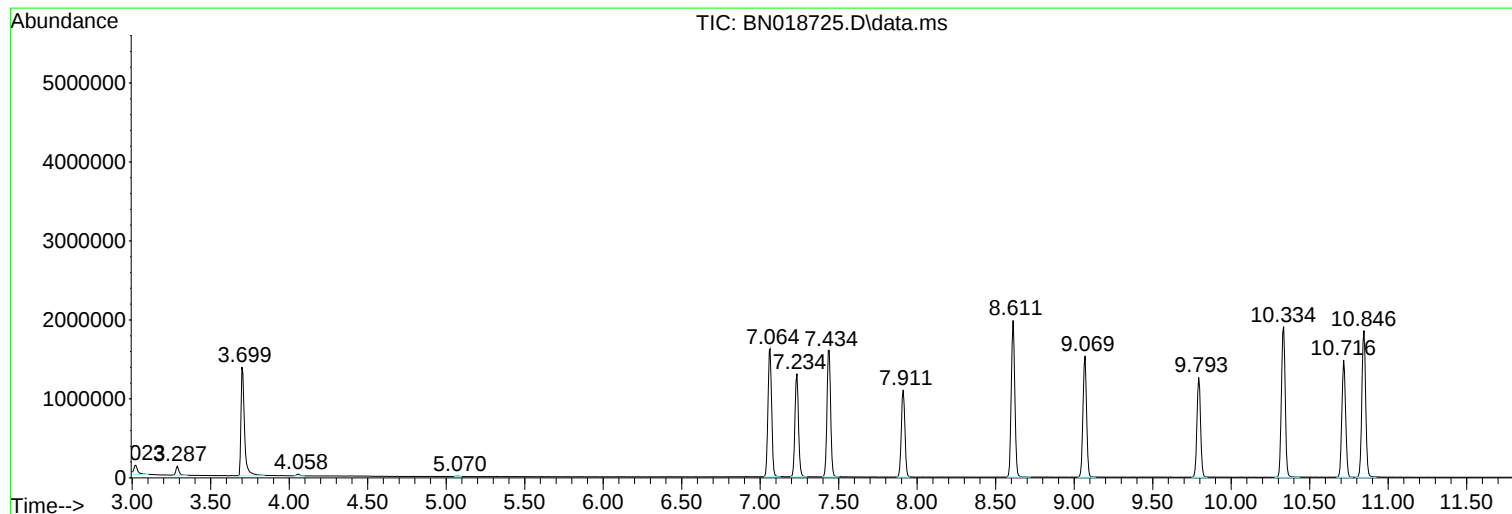
Sum of corrected areas: 82908577

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION

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



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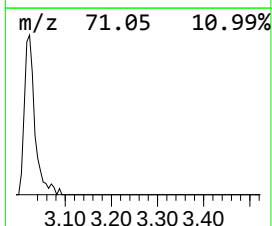
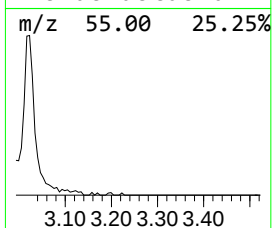
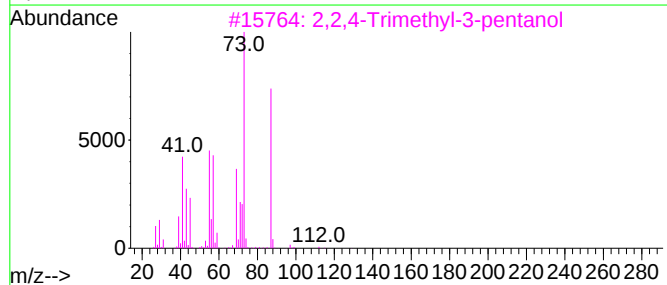
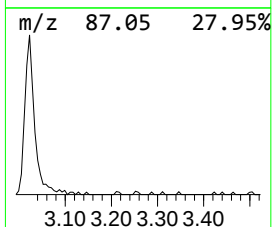
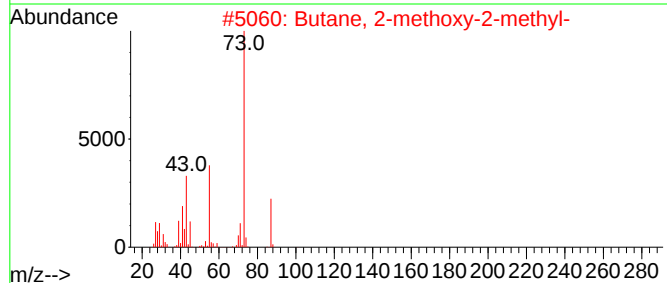
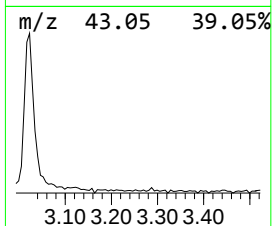
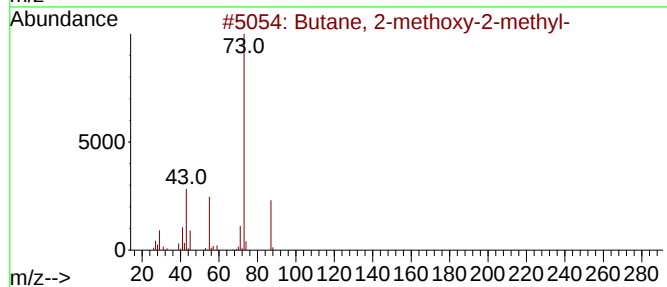
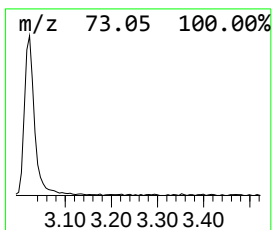
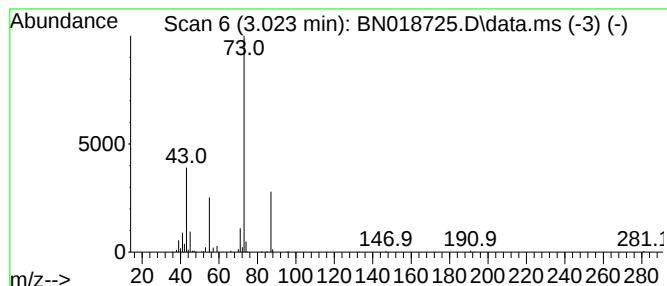
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.023	2.16 ng/ul	189342	1,4-Dichlorobenzene-d4	7.911

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
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			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	33
5			Myo-Inositol, 2-C-methyl-	194	C7H14O6	000472-96-8	33



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.023	2.2	ng/u1	189342	1	7.911	1755510	20.0