

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011874.D
 Acq On : 03 Oct 2022 16:15
 Operator : CG/JU
 Sample : N4859-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E10019

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: BP011874.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.034	4	8	30	rVB	5276084	6827817	90.70%	10.648%
2	3.305	50	54	64	rVB	62296	79467	1.06%	0.124%
3	3.405	68	71	75	rVB3	10051	11792	0.16%	0.018%
4	3.652	110	113	118	rVB	17615	22456	0.30%	0.035%
5	3.734	125	127	143	rBV	114502	181781	2.41%	0.283%
6	3.958	160	165	170	rVB	33791	48393	0.64%	0.075%
7	4.052	177	181	187	rVB	20671	26518	0.35%	0.041%
8	5.387	402	408	412	rBV	22063	34617	0.46%	0.054%
9	5.517	425	430	439	rBV	48230	93449	1.24%	0.146%
10	5.893	485	494	500	rBV2	31255	58522	0.78%	0.091%
11	7.052	682	691	701	rBV	187255	357107	4.74%	0.557%
12	7.222	711	720	733	rBV	839593	1403843	18.65%	2.189%
13	7.422	746	754	764	rBV	773968	1280748	17.01%	1.997%
14	7.893	825	834	842	rBV	823885	1361286	18.08%	2.123%
15	8.146	872	877	885	rBV5	8937	19756	0.26%	0.031%
16	8.599	948	954	970	rBV	548571	975578	12.96%	1.521%
17	9.057	1024	1032	1048	rBV	917226	1540626	20.47%	2.403%
18	9.222	1055	1060	1066	rVB6	9557	17063	0.23%	0.027%
19	9.781	1148	1155	1169	rBV	601237	1023136	13.59%	1.596%
20	10.316	1238	1246	1266	rBV	958841	1776377	23.60%	2.770%
21	10.699	1303	1311	1327	rVB	1148345	1953419	25.95%	3.046%
22	10.840	1328	1335	1353	rBV	774111	1409416	18.72%	2.198%
23	12.293	1575	1582	1588	rBV2	19146	36401	0.48%	0.057%
24	12.504	1611	1618	1621	rBV9	3715	8895	0.12%	0.014%
25	13.434	1770	1776	1782	rVB2	13576	23536	0.31%	0.037%
26	13.504	1782	1788	1793	rBV5	5312	9997	0.13%	0.016%
27	13.945	1856	1863	1878	rBV	1922129	2724592	36.19%	4.249%
28	14.228	1901	1911	1927	rBV	2335583	3596997	47.78%	5.610%
29	14.534	1956	1963	1979	rBV	1601184	2440249	32.42%	3.806%
30	14.751	1994	2000	2013	rBV2	46765	145978	1.94%	0.228%
31	14.969	2030	2037	2041	rVB10	6234	13878	0.18%	0.022%
32	15.228	2077	2081	2085	rVB6	4917	7638	0.10%	0.012%
33	15.528	2125	2132	2146	rBV	3036518	4500245	59.78%	7.018%
34	15.645	2146	2152	2166	rVV	515824	770740	10.24%	1.202%
35	15.769	2169	2173	2179	rVB2	14994	25704	0.34%	0.040%
36	17.292	2425	2432	2442	rBV	1958487	2791484	37.08%	4.353%

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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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37	17.392	2442	2449	2468	rVV	3592073	5148049	68.39%	8.029%
38	18.175	2578	2582	2588	rBV4	34541	57706	0.77%	0.090%
39	19.204	2754	2757	2761	rVB	46753	55461	0.74%	0.086%
40	19.269	2765	2768	2774	rVV	49457	68540	0.91%	0.107%
41	19.404	2789	2791	2797	rBV5	27581	41173	0.55%	0.064%
42	19.628	2823	2829	2839	rBV	4791127	6429687	85.41%	10.027%
43	21.374	3121	3126	3134	rBV	2747945	3534477	46.95%	5.512%
44	23.651	3506	3513	3528	rVB2	3632863	7527670	100.00%	11.740%
45	23.810	3534	3540	3549	rVB2	1801650	3658824	48.60%	5.706%

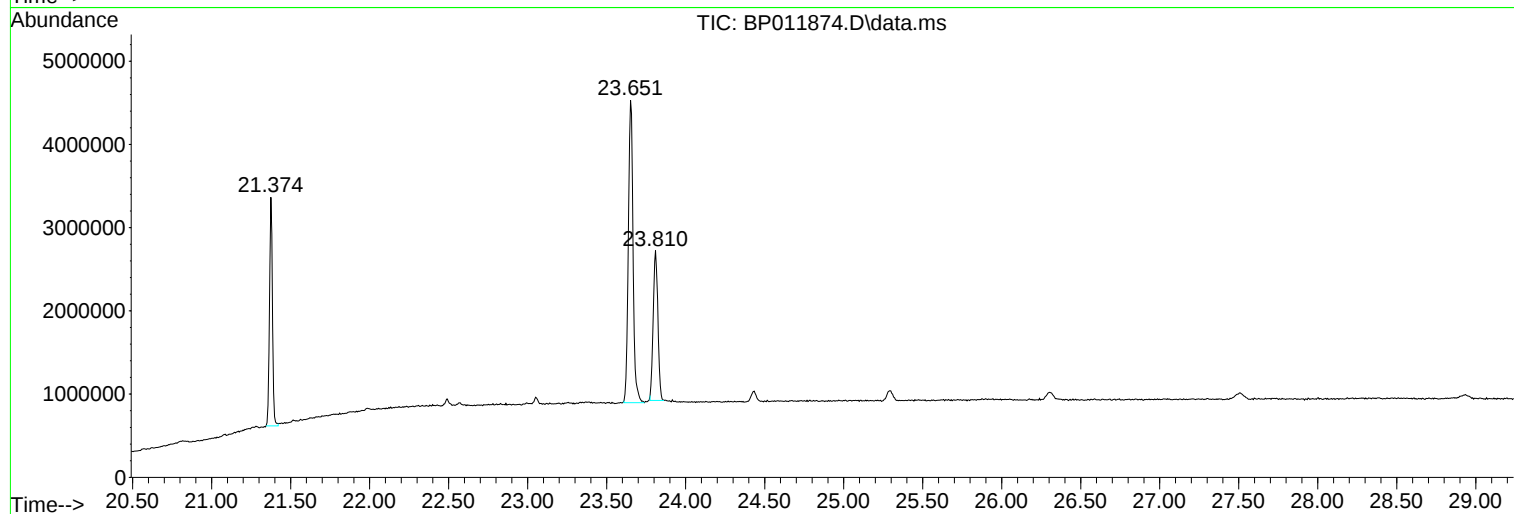
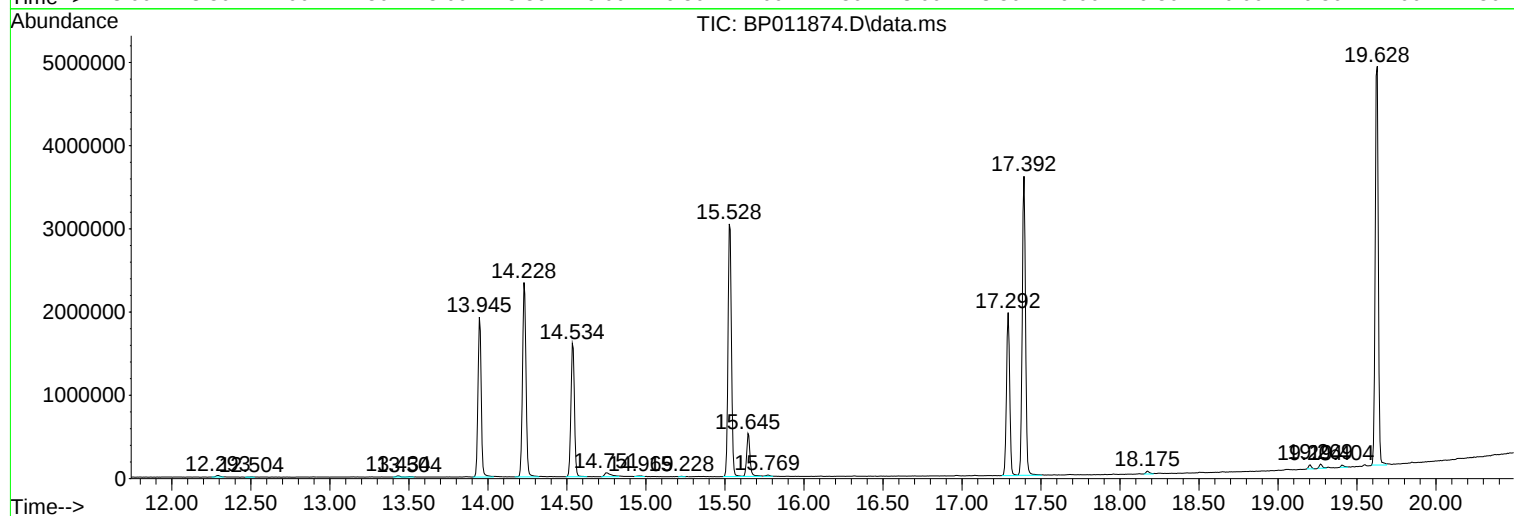
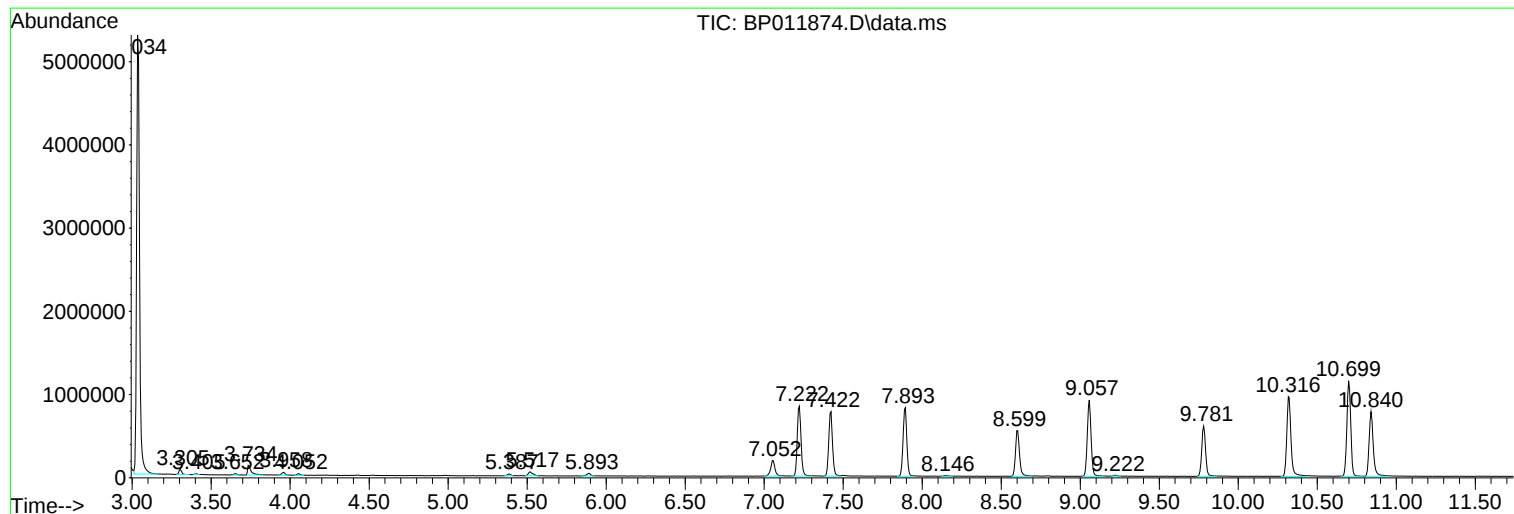
Sum of corrected areas: 64121088

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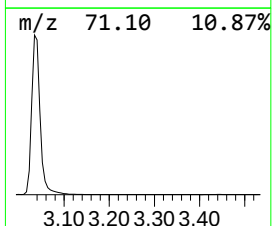
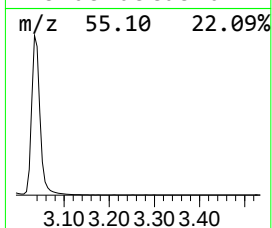
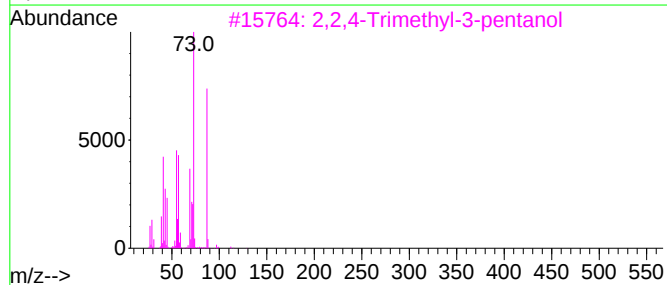
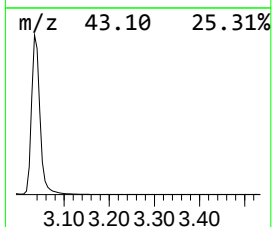
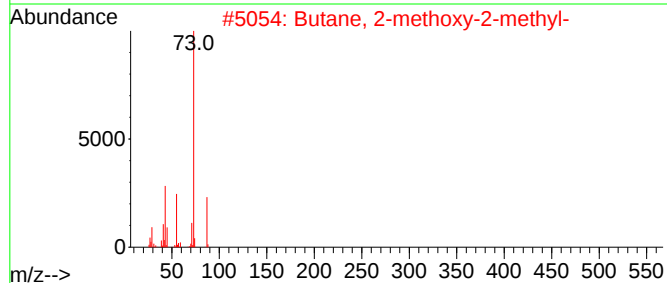
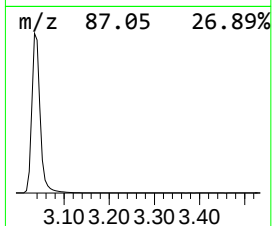
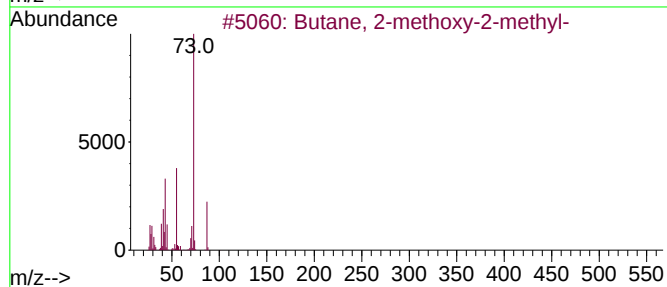
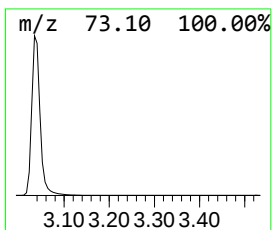
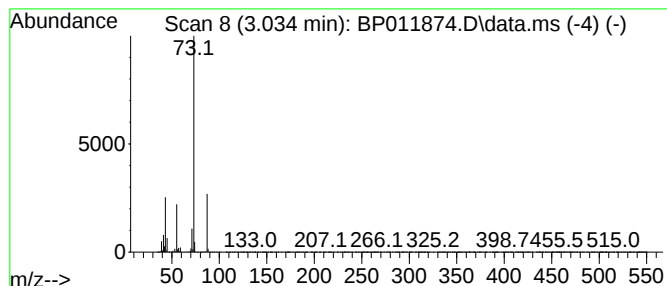
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.034	100.31 ng/ul	6827820	1,4-Dichlorobenzene-d4	7.893

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	72
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.034	100.3	ng/ul	6827820	1	7.893	1361290	20.0