

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101421\
 Data File : BP007415.D
 Acq On : 14 Oct 2021 14:04
 Operator : CG/JU
 Sample : PB139790BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK790

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-BP101421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP007415.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.128	3	7	18	rVB	200436	283977	6.90%	0.627%
2	3.387	46	51	56	rBV	46657	61858	1.50%	0.137%
3	3.793	117	120	136	rBV	597973	913153	22.20%	2.015%
4	4.140	175	179	185	rVB	13131	20815	0.51%	0.046%
5	4.646	260	265	268	rBV7	2954	6786	0.16%	0.015%
6	7.116	678	685	697	rBV	692363	1070109	26.01%	2.362%
7	7.287	708	714	725	rVB	517555	842827	20.49%	1.860%
8	7.493	742	749	758	rBV	674034	1071119	26.04%	2.364%
9	7.963	822	829	837	rBV	729240	1146536	27.87%	2.530%
10	8.663	941	948	962	rBV	858732	1329958	32.33%	2.935%
11	9.116	1017	1025	1035	rBV	629079	1003421	24.39%	2.214%
12	9.846	1141	1149	1158	rBV	420431	697944	16.96%	1.540%
13	10.381	1233	1240	1257	rBV	773489	1295740	31.50%	2.860%
14	10.769	1297	1306	1315	rVB	950381	1614719	39.25%	3.564%
15	10.898	1319	1328	1335	rBV	833165	1412603	34.34%	3.118%
16	12.351	1569	1575	1583	rBV2	18599	35299	0.86%	0.078%
17	13.569	1777	1782	1787	rVB8	5207	7533	0.18%	0.017%
18	14.004	1849	1856	1862	rBV	1473807	2093209	50.88%	4.620%
19	14.287	1895	1904	1915	rBV	1675383	2478883	60.25%	5.471%
20	14.598	1948	1957	1964	rBV	1475711	2224036	54.06%	4.908%
21	14.769	1980	1986	2002	rBV	663179	962203	23.39%	2.124%
22	15.004	2023	2026	2030	rVB6	4051	5211	0.13%	0.012%
23	15.586	2116	2125	2132	rBV	2100717	3064891	74.50%	6.764%
24	15.692	2137	2143	2153	rBV	429700	603358	14.67%	1.332%
25	15.828	2162	2166	2174	rVB3	28246	42769	1.04%	0.094%
26	16.151	2219	2221	2226	rVB6	3127	4182	0.10%	0.009%
27	16.251	2235	2238	2239	rBV3	4481	4126	0.10%	0.009%
28	16.375	2256	2259	2265	rBV8	5642	8527	0.21%	0.019%
29	17.128	2382	2387	2393	rVB2	11938	21607	0.53%	0.048%
30	17.345	2418	2424	2434	rBV	1898017	2753157	66.92%	6.076%
31	17.445	2434	2441	2453	rVV2	2346178	3375278	82.04%	7.449%
32	17.710	2484	2486	2490	rBV5	4401	7243	0.18%	0.016%
33	18.245	2574	2577	2581	rBV6	5849	7852	0.19%	0.017%
34	19.257	2745	2749	2753	rVB4	29279	36841	0.90%	0.081%
35	19.322	2756	2760	2763	rBV	37051	45205	1.10%	0.100%
36	19.674	2814	2820	2830	rBV	3110802	4058763	98.66%	8.957%

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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	21.427	3113	3118	3130	rVB	2322432	3225402	78.40%	7.118%
38	23.721	3501	3508	3519	rVB2	2036571	4114095	100.00%	9.080%
39	23.880	3528	3535	3548	rVB2	1583429	3360468	81.68%	7.416%

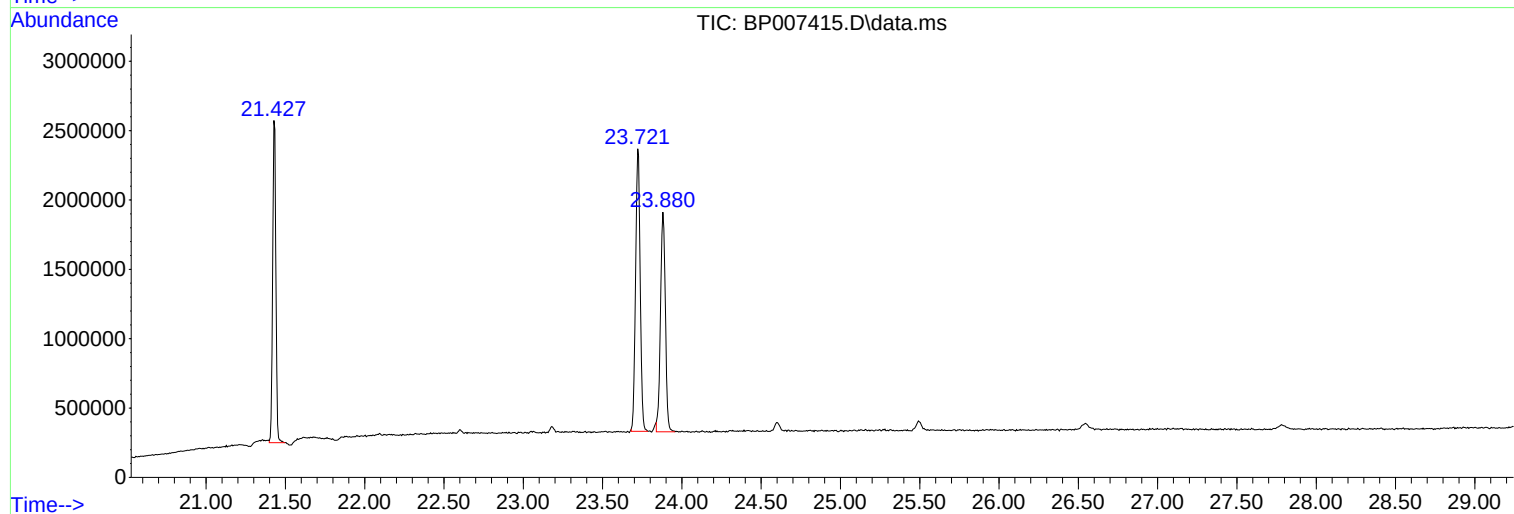
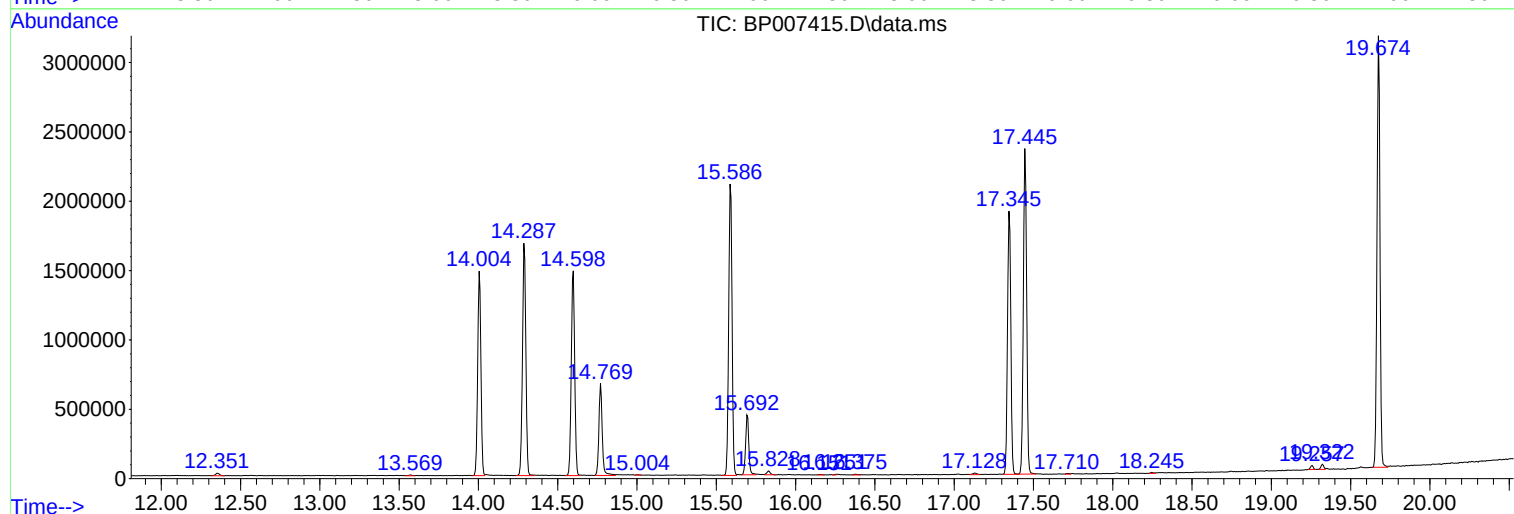
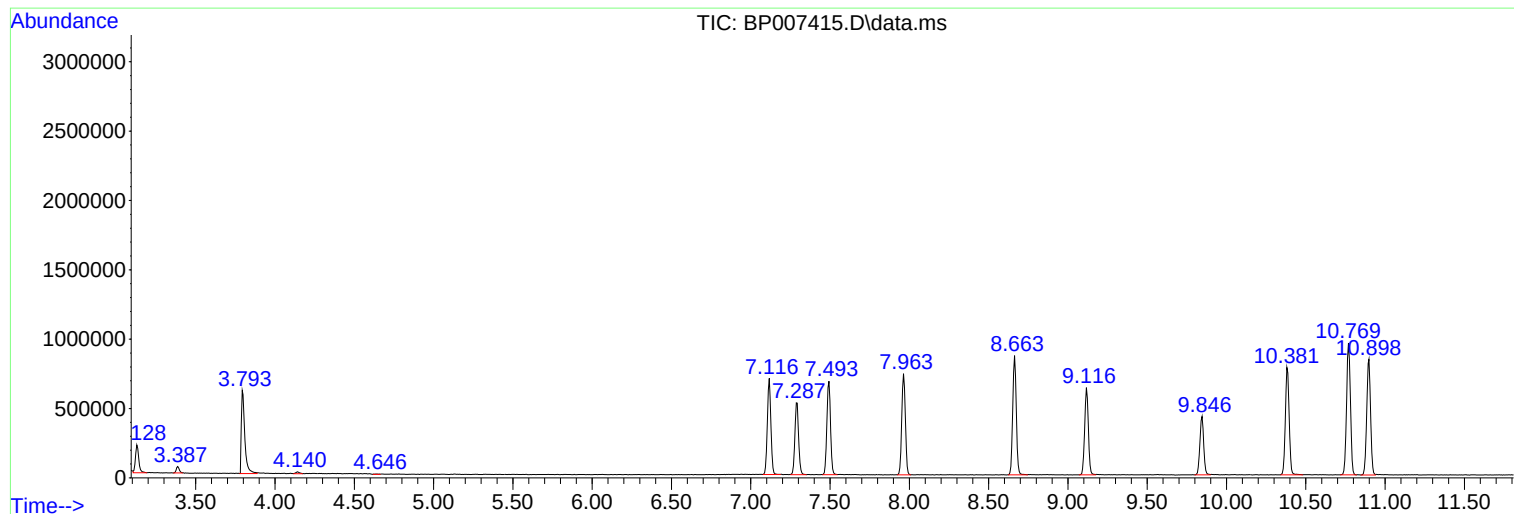
Sum of corrected areas: 45311703

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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



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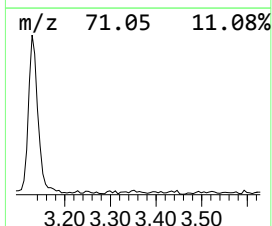
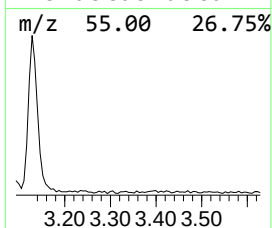
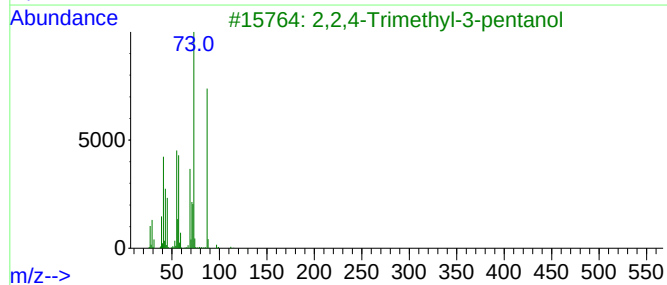
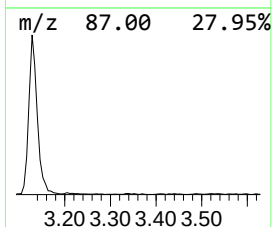
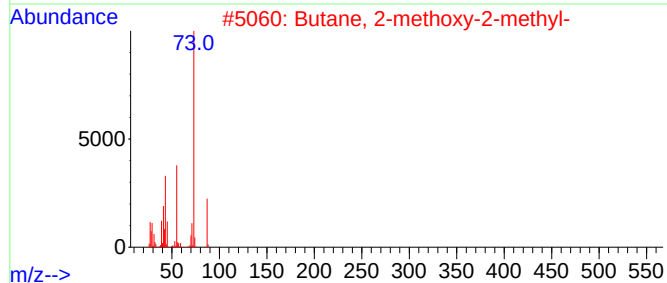
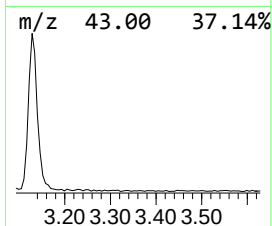
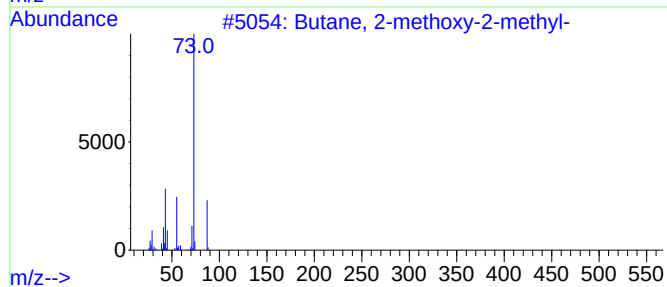
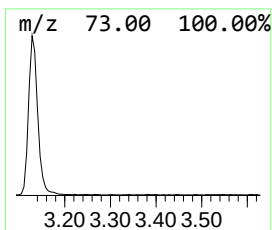
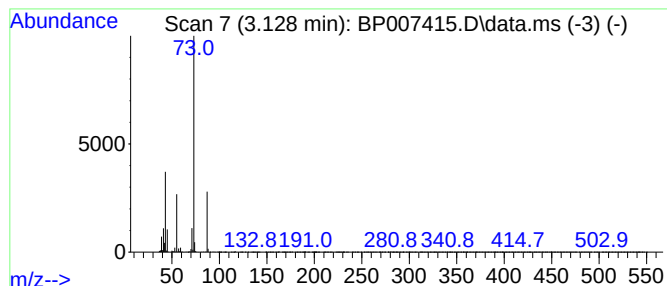
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.128	4.95 ng/ul	283977	1,4-Dichlorobenzene-d4	7.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
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, tetramethyl-	88	C4H12Si	000075-76-3	25
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23



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Quant Title : SVOA CALIBRATION

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...	3.128	5.0	ng/ul	283977	1	7.963	1146540	20.0