

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
 Data File : BP009728.D
 Acq On : 08 Apr 2022 21:35
 Operator : CG/JU
 Sample : N2182-12
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0R19

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-BP040722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP009728.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.129	3	7	20	rVB	169555	254389	7.71%	0.740%
2	3.393	48	52	61	rVB	22549	37273	1.13%	0.108%
3	3.805	118	122	137	rBV	301853	495942	15.04%	1.443%
4	4.146	177	180	184	rVB4	6874	10266	0.31%	0.030%
5	5.075	335	338	344	rBV7	3402	4752	0.14%	0.014%
6	7.122	679	686	698	rBV	440522	726205	22.02%	2.114%
7	7.299	704	716	726	rBV	341506	559932	16.98%	1.630%
8	7.505	741	751	764	rBV	435217	725603	22.00%	2.112%
9	7.975	822	831	839	rBV	388475	628984	19.07%	1.831%
10	8.040	839	842	848	rVB2	5665	9337	0.28%	0.027%
11	8.675	942	950	959	rBV	591074	977297	29.64%	2.844%
12	9.128	1019	1027	1036	rBV	449425	766023	23.23%	2.229%
13	9.305	1055	1057	1060	rBV3	3091	3730	0.11%	0.011%
14	9.599	1103	1107	1114	rVB3	9182	14809	0.45%	0.043%
15	9.858	1143	1151	1160	rBV	371355	629870	19.10%	1.833%
16	10.399	1234	1243	1253	rBV	580018	1007605	30.55%	2.933%
17	10.575	1268	1273	1275	rBV6	2834	3899	0.12%	0.011%
18	10.781	1298	1308	1318	rBV	559902	976999	29.63%	2.844%
19	10.910	1322	1330	1341	rBV	623554	1074989	32.60%	3.129%
20	11.346	1397	1404	1410	rVB	4038	7841	0.24%	0.023%
21	11.428	1415	1418	1423	rVB7	2262	3571	0.11%	0.010%
22	11.693	1457	1463	1473	rBV7	3164	9730	0.30%	0.028%
23	12.222	1546	1553	1561	rBV5	7664	13745	0.42%	0.040%
24	12.310	1561	1568	1572	rBV8	4013	7082	0.21%	0.021%
25	12.363	1572	1577	1583	rVB2	12916	20105	0.61%	0.059%
26	13.034	1684	1691	1695	rVB9	4103	6546	0.20%	0.019%
27	13.110	1700	1704	1709	rVV7	2805	5712	0.17%	0.017%
28	13.169	1709	1714	1718	rVV6	4750	6970	0.21%	0.020%
29	13.346	1738	1744	1749	rBV9	1989	4598	0.14%	0.013%
30	13.451	1756	1762	1768	rBV	28601	44841	1.36%	0.131%
31	13.575	1779	1783	1786	rBV6	2332	3636	0.11%	0.011%
32	14.010	1851	1857	1864	rBV	1256818	1730773	52.48%	5.037%
33	14.104	1872	1873	1878	rVB5	3776	4457	0.14%	0.013%
34	14.146	1878	1880	1886	rVB6	4090	6446	0.20%	0.019%
35	14.299	1896	1906	1918	rBV	1303060	1923214	58.32%	5.597%
36	14.522	1936	1944	1948	rBV5	11172	18190	0.55%	0.053%

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Integrator: RTE

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Filtering: 5

Sampling : 1

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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-BP040722.M
 Title : SVOA CALIBRATION

37	14.604	1951	1958	1973	rBV	988063	1470958	44.60%	4.281%
38	14.781	1981	1988	1998	rBV	509198	753170	22.84%	2.192%
39	14.963	2017	2019	2024	rVB6	2925	4169	0.13%	0.012%
40	15.016	2024	2028	2035	rBV10	5199	10521	0.32%	0.031%
41	15.157	2045	2052	2064	rBV	31267	98913	3.00%	0.288%
42	15.416	2092	2096	2100	rBV6	4087	7655	0.23%	0.022%
43	15.469	2102	2105	2110	rVB7	5384	8005	0.24%	0.023%
44	15.598	2118	2127	2136	rBV	1741876	2487314	75.42%	7.239%
45	15.704	2138	2145	2161	rVV	368522	537908	16.31%	1.566%
46	15.840	2163	2168	2173	rVB2	17054	24216	0.73%	0.070%
47	16.034	2197	2201	2203	rBV5	2326	3370	0.10%	0.010%
48	16.175	2220	2225	2226	rBV5	2902	4617	0.14%	0.013%
49	16.287	2237	2244	2249	rBV5	2920	8617	0.26%	0.025%
50	16.387	2257	2261	2267	rVB6	7157	13669	0.41%	0.040%
51	16.457	2270	2273	2276	rBV5	2361	3644	0.11%	0.011%
52	16.604	2296	2298	2303	rBV6	3353	5297	0.16%	0.015%
53	16.757	2321	2324	2334	rVB6	5387	9384	0.28%	0.027%
54	16.828	2334	2336	2340	rBV5	2500	3849	0.12%	0.011%
55	17.040	2366	2372	2376	rBV9	4607	8601	0.26%	0.025%
56	17.110	2378	2384	2387	rBV3	13597	22198	0.67%	0.065%
57	17.357	2419	2426	2435	rBV	1265188	1830275	55.50%	5.327%
58	17.457	2436	2443	2453	rVV	1881914	2752611	83.47%	8.011%
59	17.540	2453	2457	2463	rVB6	13741	21884	0.66%	0.064%
60	18.034	2538	2541	2546	rVB7	6366	8388	0.25%	0.024%
61	18.239	2572	2576	2584	rBV	77021	110230	3.34%	0.321%
62	18.539	2624	2627	2635	rVB9	6119	11983	0.36%	0.035%
63	18.663	2645	2648	2651	rBV5	12570	14850	0.45%	0.043%
64	18.757	2660	2664	2667	rBV4	10678	15712	0.48%	0.046%
65	19.092	2718	2721	2726	rVB7	18425	24560	0.74%	0.071%
66	19.263	2746	2750	2755	rBV3	29792	40125	1.22%	0.117%
67	19.334	2758	2762	2764	rVV	27532	33479	1.02%	0.097%
68	19.475	2783	2786	2790	rBV5	20200	23141	0.70%	0.067%
69	19.681	2816	2821	2828	rBV	2389486	3297772	100.00%	9.598%
70	21.151	3067	3071	3076	rBV2	326238	445387	13.51%	1.296%
71	21.433	3114	3119	3126	rBV	1551165	2082258	63.14%	6.060%
72	21.904	3195	3199	3206	rVB	101044	136862	4.15%	0.398%
73	22.086	3227	3230	3236	rBV2	91710	142427	4.32%	0.415%
74	23.751	3506	3513	3523	rVB2	1555057	3161428	95.87%	9.201%
75	23.910	3533	3540	3549	rVB2	964980	1994207	60.47%	5.804%

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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-BP040722.M
Title : SVOA CALIBRATION

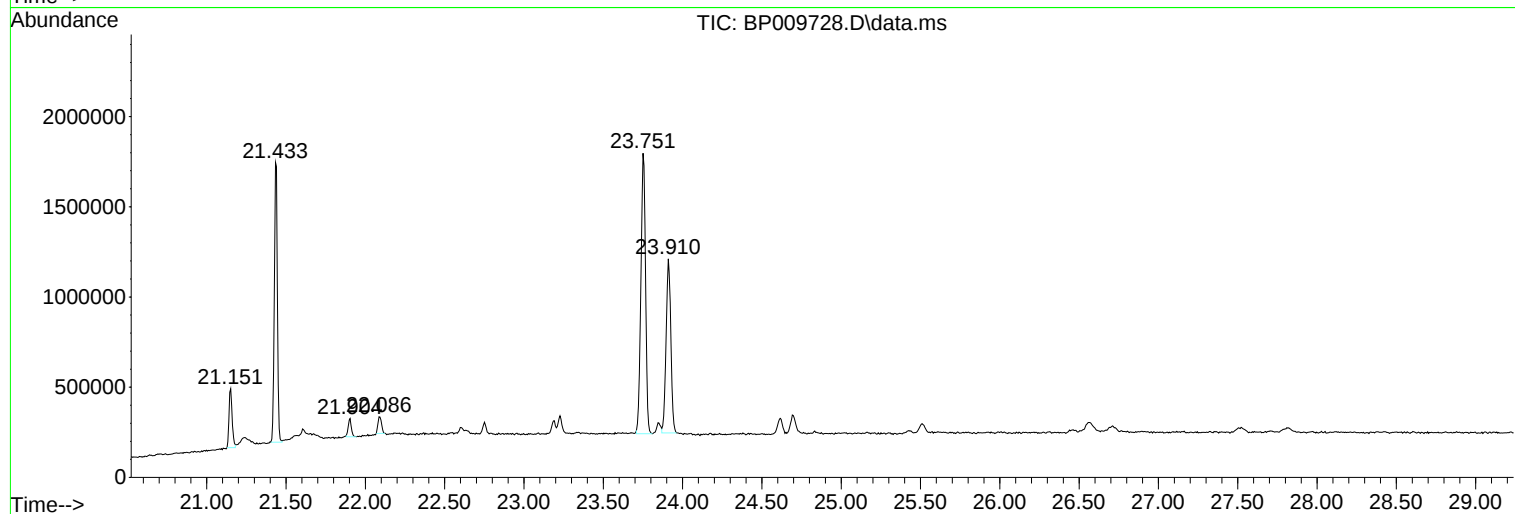
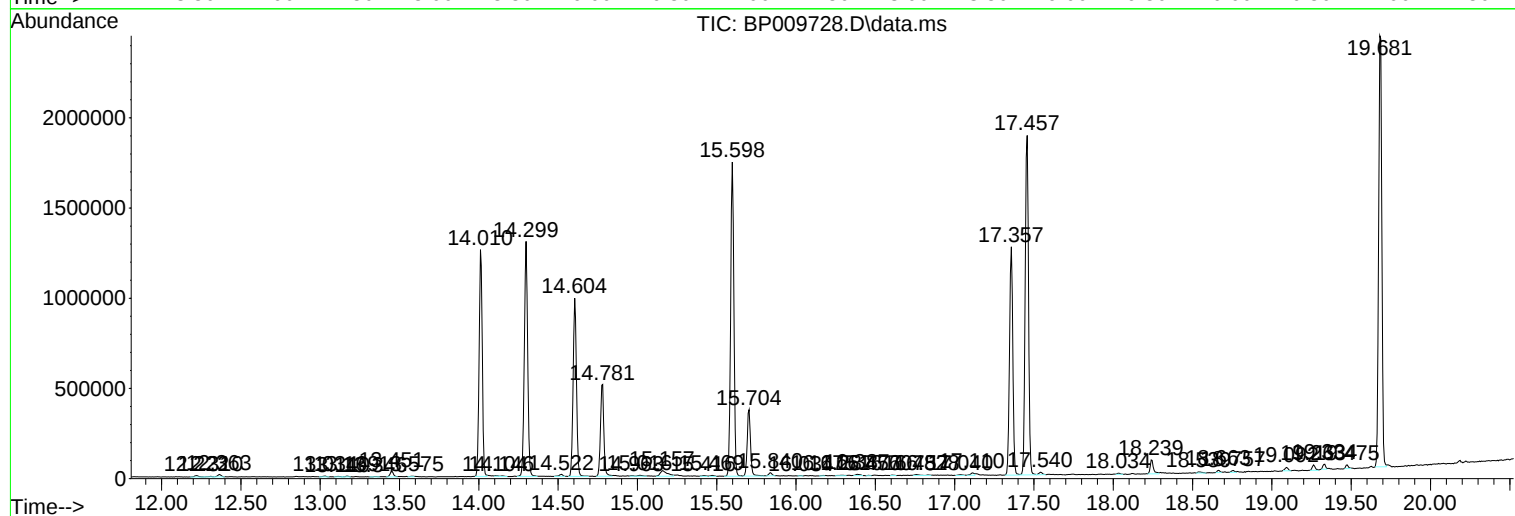
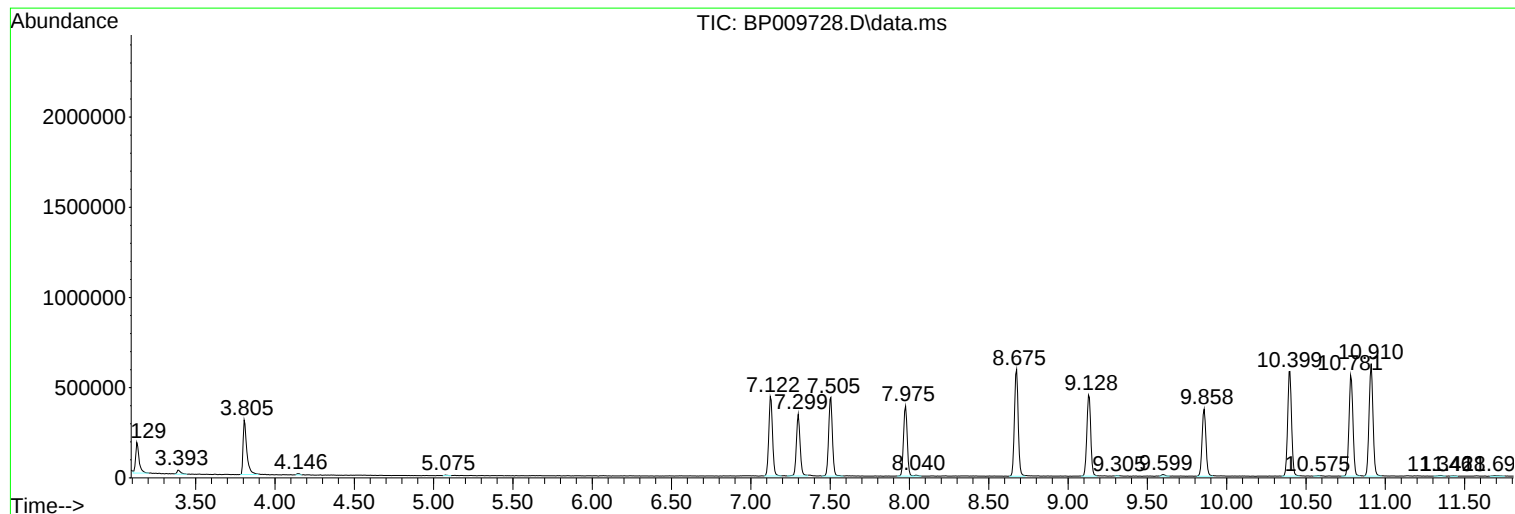
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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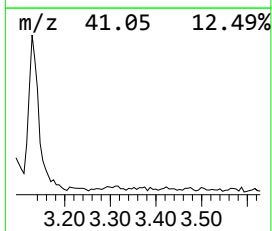
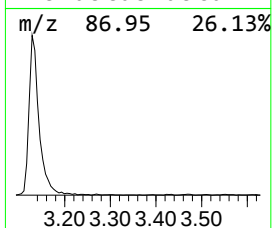
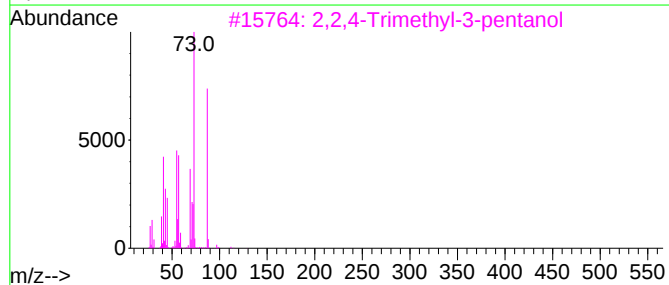
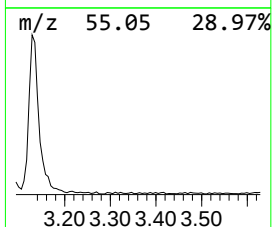
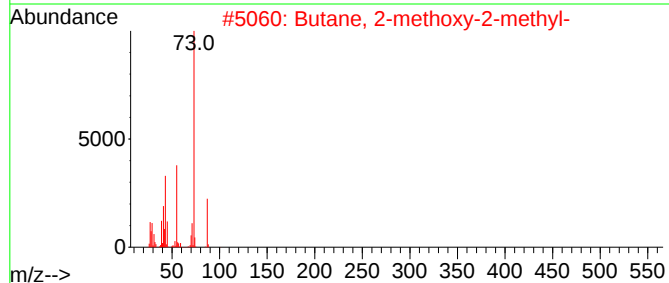
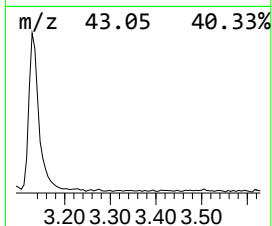
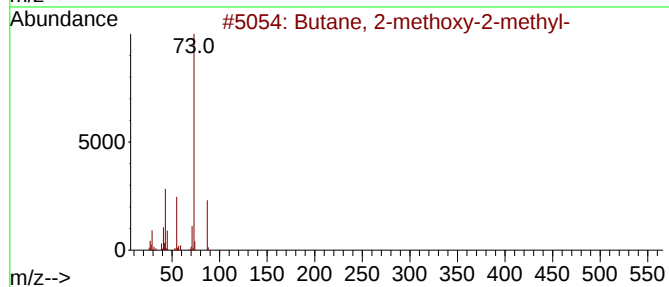
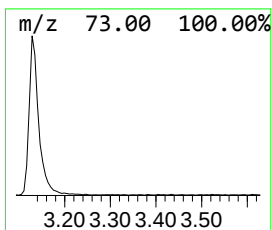
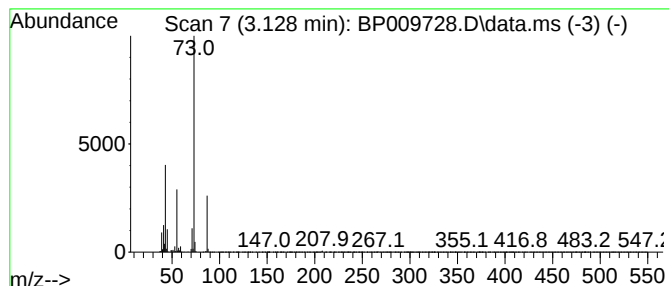
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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.
3.128	8.09 ng/ul	254389	1,4-Dichlorobenzene-d4	7.975

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	78		
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40		
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	35		
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	32		



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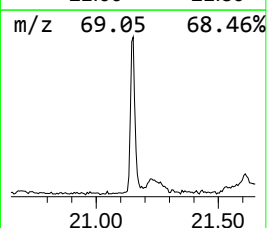
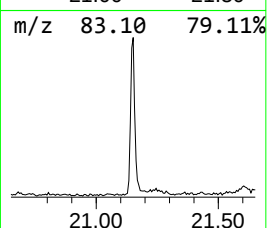
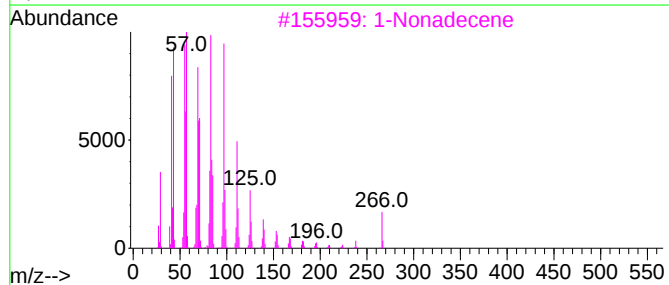
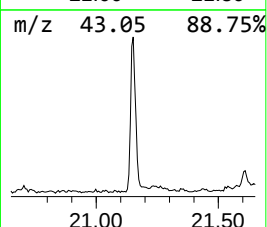
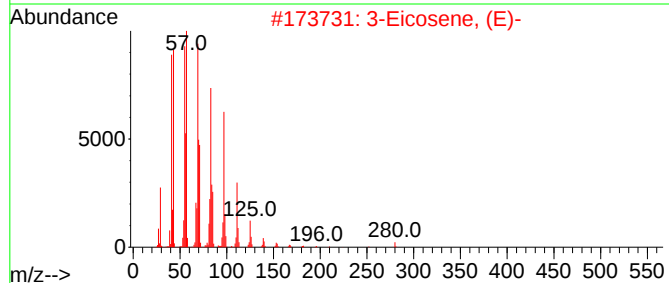
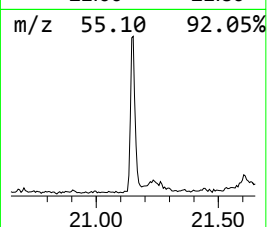
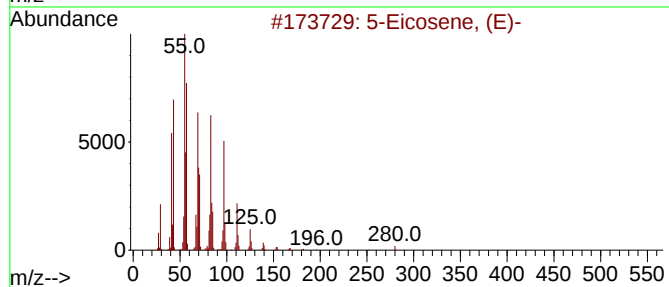
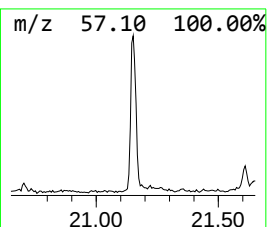
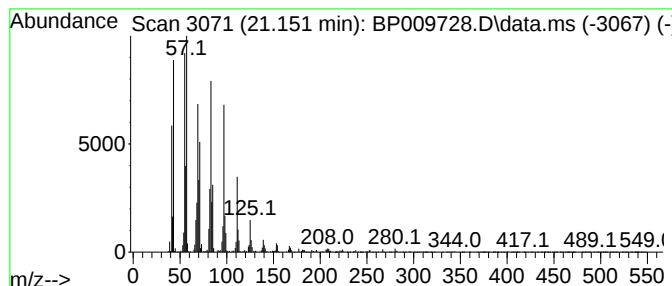
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.151	4.28 ng/ul	445387	Chrysene-d12	21.433

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-		280	C20H40	074685-30-6	97
2	3-Eicosene, (E)-		280	C20H40	074685-33-9	96
3	1-Nonadecene		266	C19H38	018435-45-5	95
4	Trichloroacetic acid, pentadecyl...		372	C17H31Cl3O2	074339-53-0	94
5	Pentadecafluorooctanoic acid, he...		638	C24H33F15O2	1000406-04-6	94



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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...	3.128	8.1	ng/ul	254389	1	7.975	628984	20.0
(DEL) Alkane: S...	21.151	4.3	ng/ul	445387	5	21.433	2082260	20.0