

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
 Data File : BP009882.D
 Acq On : 16 Apr 2022 03:24
 Operator : CG/JU
 Sample : N2421-01
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J55

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: BP009882.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.122	3	6	18	rVB	361514	502994	8.78%	1.064%
2	3.381	46	50	61	rVB2	30866	47219	0.82%	0.100%
3	3.799	118	121	132	rBV	148433	252800	4.41%	0.535%
4	4.134	174	178	184	rBV2	11022	15622	0.27%	0.033%
5	5.263	366	370	377	rVB	15401	25177	0.44%	0.053%
6	7.110	677	684	701	rVB2	160981	297113	5.19%	0.629%
7	7.281	706	713	731	rVB	528624	858914	14.99%	1.817%
8	7.487	741	748	760	rBV	551321	877776	15.32%	1.857%
9	7.957	820	828	838	rBV	497480	833862	14.56%	1.764%
10	8.310	882	888	894	rBV3	12201	18676	0.33%	0.040%
11	8.657	938	947	957	rBV	474475	762943	13.32%	1.614%
12	9.116	1017	1025	1038	rBV	721811	1220836	21.31%	2.583%
13	9.298	1050	1056	1060	rBV6	5179	8715	0.15%	0.018%
14	9.575	1097	1103	1108	rVB4	10112	18569	0.32%	0.039%
15	9.840	1138	1148	1160	rBV	589136	976124	17.04%	2.065%
16	10.381	1231	1240	1255	rBV	794784	1376076	24.02%	2.911%
17	10.545	1263	1268	1276	rBV3	27147	53830	0.94%	0.114%
18	10.628	1276	1282	1287	rVB2	20180	35630	0.62%	0.075%
19	10.763	1294	1305	1317	rBV	731680	1233885	21.54%	2.610%
20	10.892	1319	1327	1337	rBV	775709	1319962	23.04%	2.792%
21	11.151	1363	1371	1378	rVB	5017	11712	0.20%	0.025%
22	12.345	1569	1574	1581	rVB3	18784	34626	0.60%	0.073%
23	12.563	1606	1611	1617	rVB9	2947	6538	0.11%	0.014%
24	12.963	1673	1679	1683	rVB7	5448	9414	0.16%	0.020%
25	13.110	1700	1704	1707	rVB6	4498	6320	0.11%	0.013%
26	13.210	1715	1721	1727	rVB6	8535	15930	0.28%	0.034%
27	13.398	1745	1753	1757	rBV4	25322	45589	0.80%	0.096%
28	13.428	1757	1758	1762	rVV4	7962	7664	0.13%	0.016%
29	13.563	1776	1781	1782	rVV5	5595	7523	0.13%	0.016%
30	13.592	1782	1786	1791	rVB3	9218	14939	0.26%	0.032%
31	13.992	1847	1854	1869	rBV	1862349	2638680	46.06%	5.582%
32	14.128	1872	1877	1883	rVB9	5620	10746	0.19%	0.023%
33	14.233	1890	1895	1897	rBV2	30445	41821	0.73%	0.088%
34	14.280	1897	1903	1913	rVB	1946937	2818332	49.20%	5.962%
35	14.492	1936	1939	1944	rVB4	6661	9400	0.16%	0.020%
36	14.586	1947	1955	1965	rBV	1205032	1792614	31.29%	3.792%

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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
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37	14.663	1965	1968	1974	rVB8	4481	6639	0.12%	0.014%
38	14.763	1975	1985	1996	rBV	181707	290188	5.07%	0.614%
39	14.998	2021	2025	2029	rVB6	4040	6273	0.11%	0.013%
40	15.222	2059	2063	2070	rVB3	14160	22378	0.39%	0.047%
41	15.398	2087	2093	2098	rBV2	35035	58878	1.03%	0.125%
42	15.445	2099	2101	2105	rVB4	9726	10330	0.18%	0.022%
43	15.580	2114	2124	2134	rBV	2825371	3970900	69.32%	8.401%
44	15.686	2136	2142	2159	rVV	1015106	1404984	24.53%	2.972%
45	15.822	2159	2165	2170	rVB2	29749	48606	0.85%	0.103%
46	15.945	2181	2186	2191	rBV6	10601	18512	0.32%	0.039%
47	16.004	2192	2196	2200	rBV7	6002	8296	0.14%	0.018%
48	16.222	2230	2233	2237	rBV6	5512	8510	0.15%	0.018%
49	16.269	2237	2241	2245	rVV4	17662	26178	0.46%	0.055%
50	16.310	2245	2248	2251	rVV4	12567	17582	0.31%	0.037%
51	16.363	2254	2257	2264	rVB7	15476	25736	0.45%	0.054%
52	16.480	2271	2277	2284	rBV2	14550	23227	0.41%	0.049%
53	16.586	2291	2295	2301	rBV8	3769	6946	0.12%	0.015%
54	16.651	2305	2306	2312	rVV6	3932	7094	0.12%	0.015%
55	16.721	2315	2318	2321	rBV5	4840	5927	0.10%	0.013%
56	16.816	2331	2334	2339	rVB6	8682	12352	0.22%	0.026%
57	17.016	2365	2368	2372	rVB5	7261	8977	0.16%	0.019%
58	17.057	2372	2375	2378	rBV4	6637	7371	0.13%	0.016%
59	17.110	2378	2384	2388	rVV5	16926	32473	0.57%	0.069%
60	17.157	2390	2392	2397	rVB5	9384	12210	0.21%	0.026%
61	17.263	2406	2410	2415	rVB4	32641	42843	0.75%	0.091%
62	17.339	2416	2423	2432	rBV	1516960	2235429	39.02%	4.729%
63	17.439	2433	2440	2451	rVV	3384573	4810095	83.97%	10.176%
64	17.527	2451	2455	2459	rVB7	7210	11886	0.21%	0.025%
65	18.010	2534	2537	2543	rVB4	21214	25148	0.44%	0.053%
66	18.221	2568	2573	2579	rBV	55607	83919	1.47%	0.178%
67	18.286	2581	2584	2591	rVB3	19803	27005	0.47%	0.057%
68	18.516	2620	2623	2624	rBV3	10266	10438	0.18%	0.022%
69	18.968	2696	2700	2702	rBV4	11512	17218	0.30%	0.036%
70	19.051	2711	2714	2718	rBV5	13848	21304	0.37%	0.045%
71	19.245	2743	2747	2750	rBV2	50284	63956	1.12%	0.135%
72	19.280	2751	2753	2755	rVV3	17337	19430	0.34%	0.041%
73	19.310	2755	2758	2765	rVB2	44097	61002	1.06%	0.129%
74	19.445	2779	2781	2788	rBV7	12446	15702	0.27%	0.033%
75	19.668	2812	2819	2829	rBV	4273789	5728217	100.00%	12.118%
76	21.127	3063	3067	3074	rBV	193835	272343	4.75%	0.576%

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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

77	21.415	3111	3116	3123	rBV	1820241	2444034	42.67%	5.170%
78	23.721	3500	3508	3516	rBV2	2450274	5007250	87.41%	10.593%
79	23.880	3528	3535	3544	rVB2	1044537	2123385	37.07%	4.492%

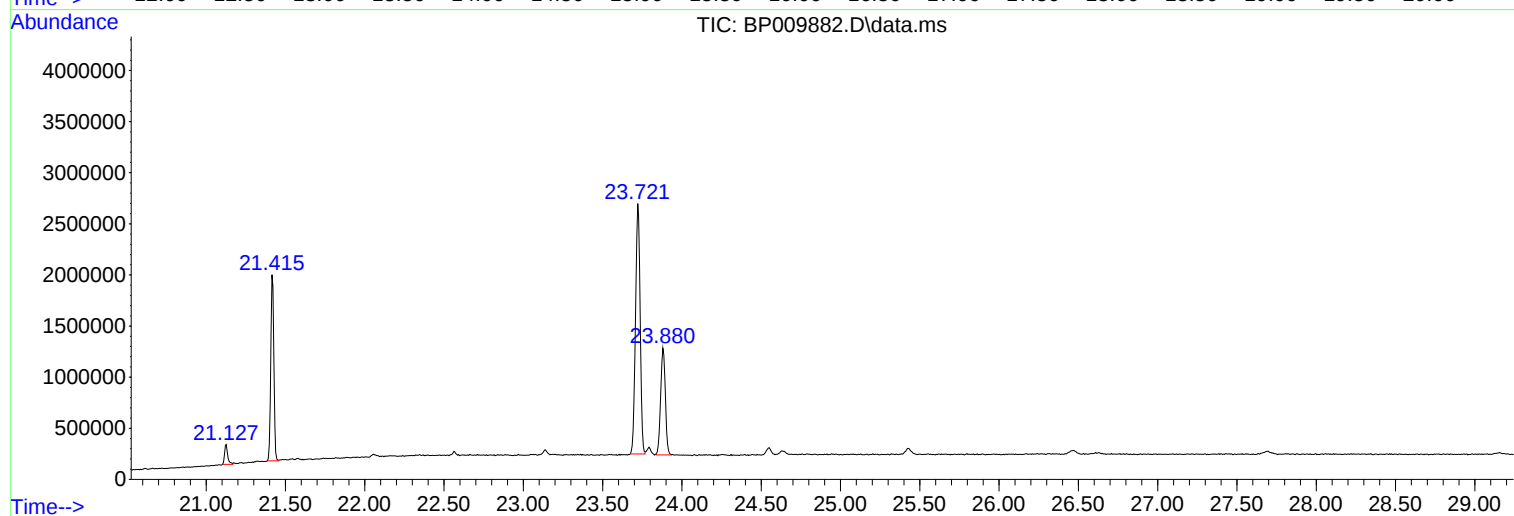
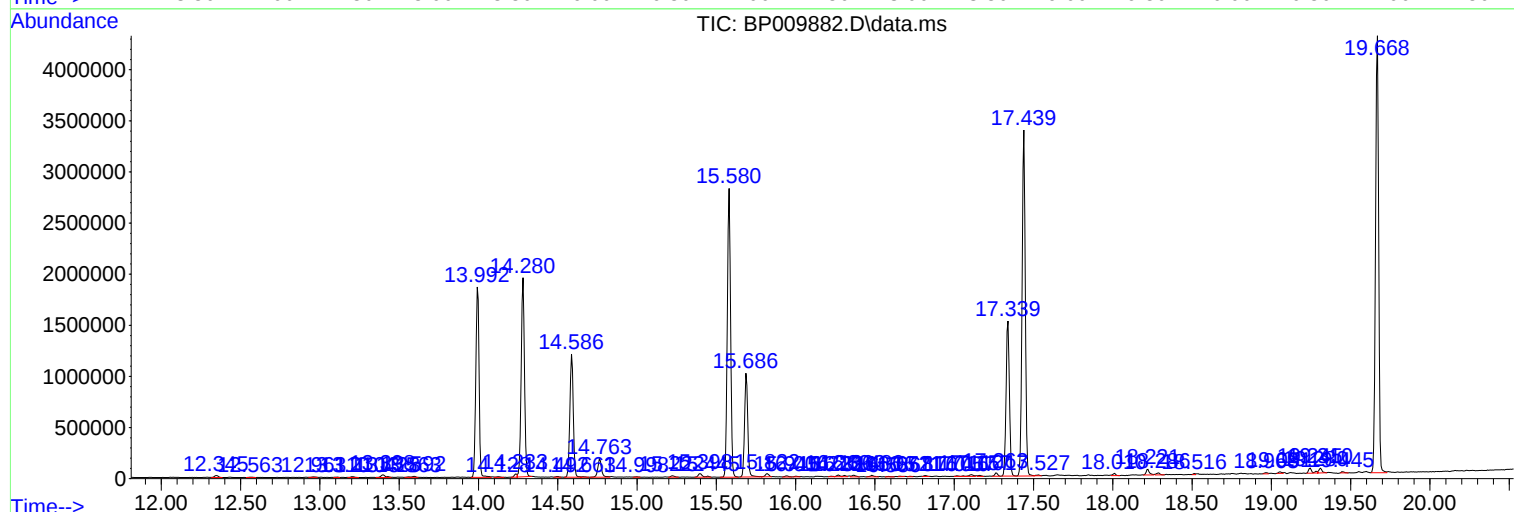
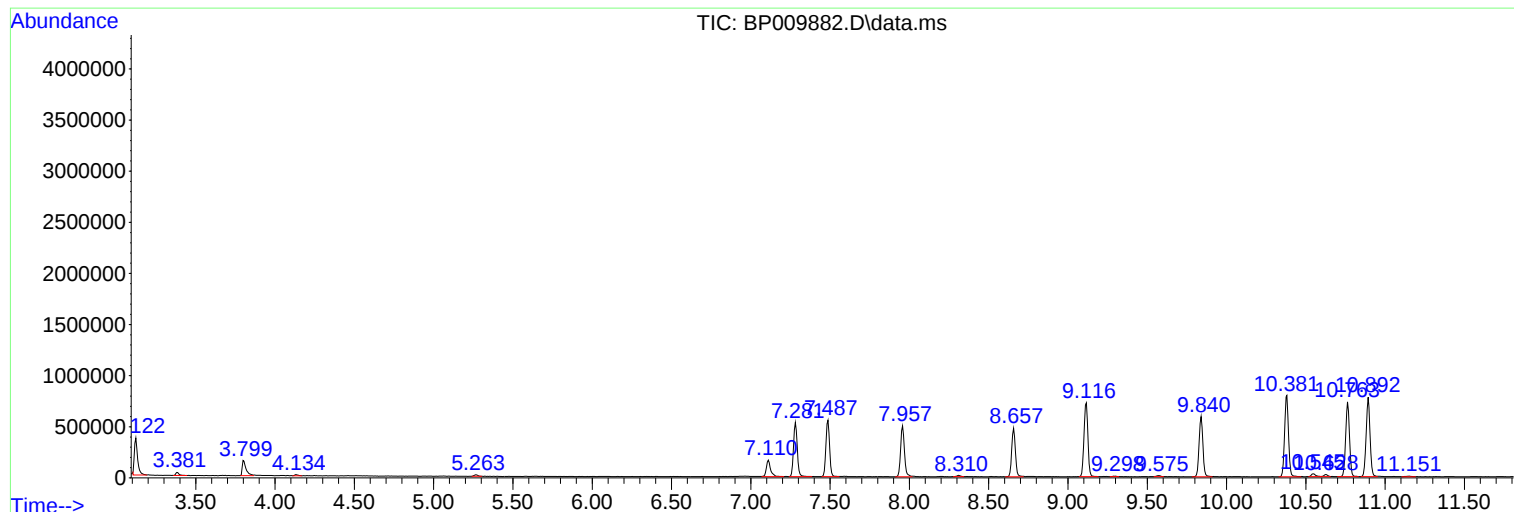
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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



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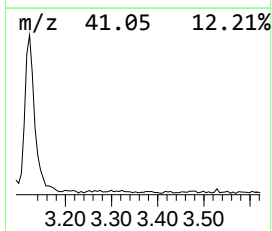
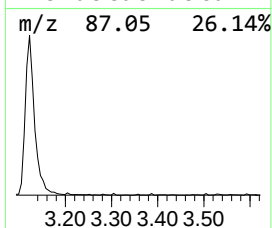
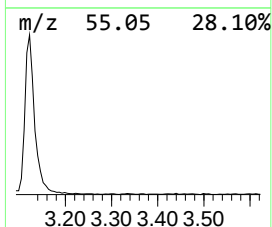
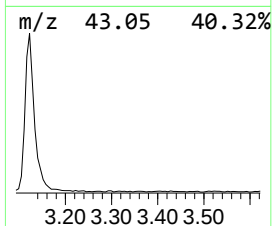
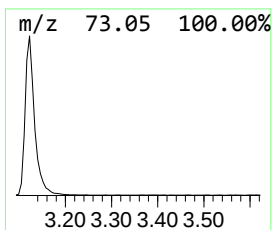
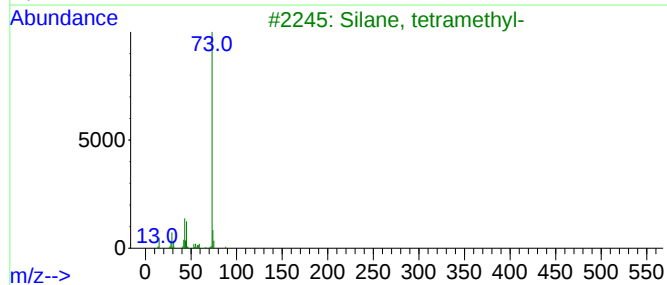
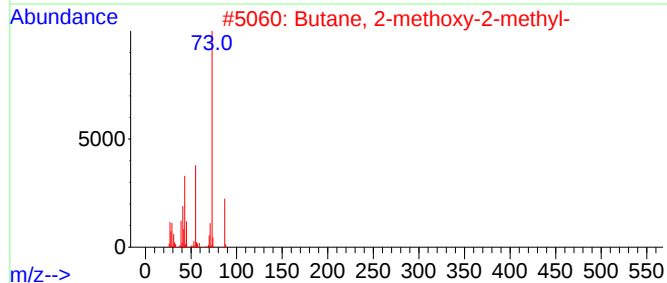
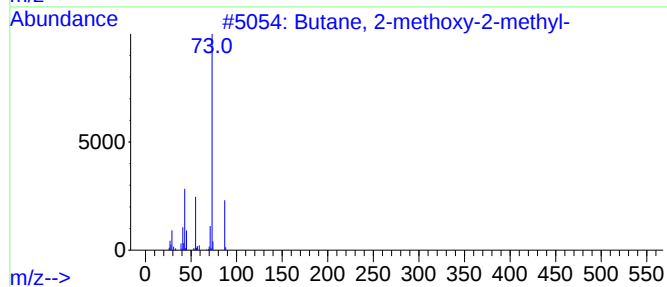
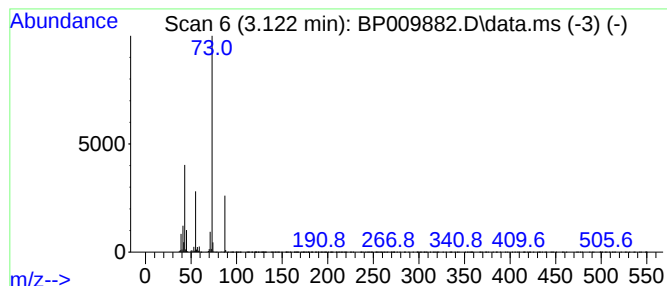
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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.
3.122	12.06 ng/ul	502994	1,4-Dichlorobenzene-d4	7.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	50		
3	Silane, tetramethyl-	88	C4H12Si	000075-76-3	27		
4	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25		
5	1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	23		



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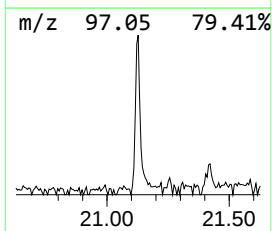
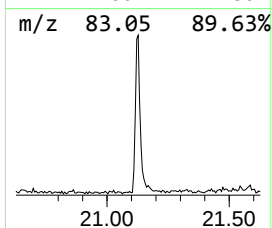
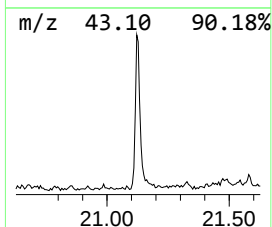
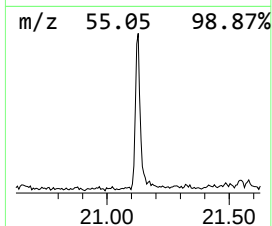
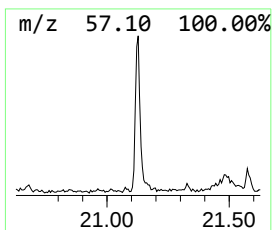
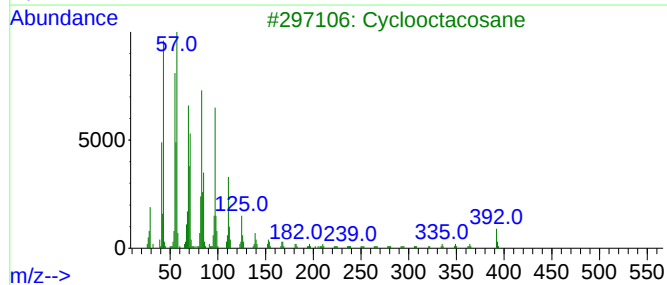
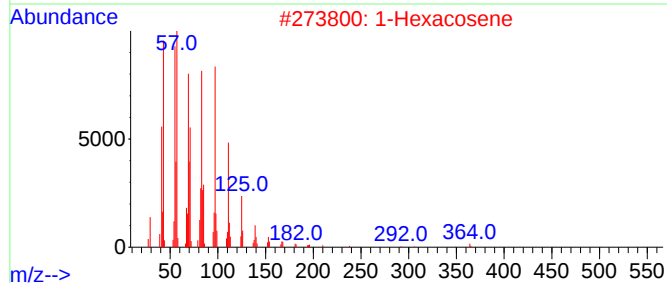
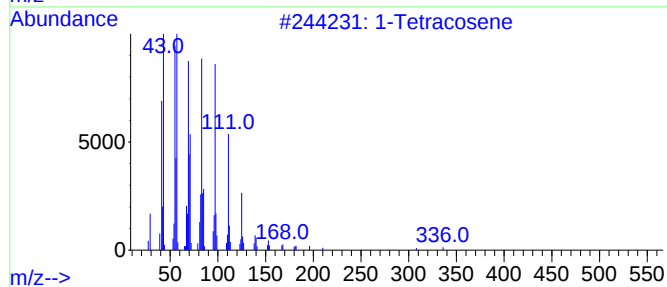
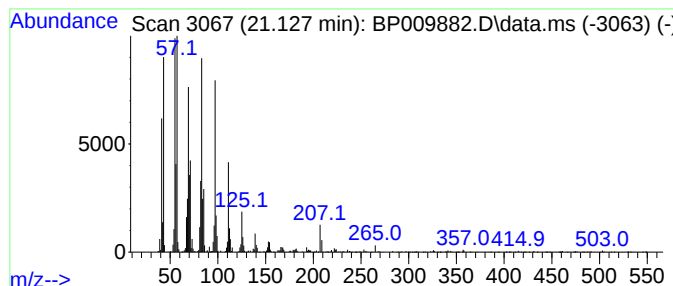
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.127	2.23 ng/ul	272343	Chrysene-d12	21.415

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene		336	C24H48	010192-32-2	94
2	1-Hexacosene		364	C26H52	018835-33-1	94
3	Cyclooctacosane		392	C28H56	000297-24-5	92
4	Behenic alcohol		326	C22H46O	000661-19-8	91
5	Pentafluoropropionic acid, penta...		374	C18H31F5O2	959092-08-1	91



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TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.122	12.1	ng/u1	502994	1	7.957	833862	20.0
(DEL) Alkane: S...	21.127	2.2	ng/u1	272343	5	21.415	2444030	20.0