

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041122\
 Data File : BM034613.D
 Acq On : 11 Apr 2022 17:04
 Operator : CG/JU
 Sample : N2338-05
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0J90

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_M\Methods\SFAM-EPA-BM040922.M
 Title : SVOA CALIBRATION

Signal : TIC: BM034613.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.255	34	37	42	rVB	28610	31683	1.20%	0.137%
2	3.684	107	110	128	rBV	96156	161810	6.13%	0.698%
3	4.019	164	167	172	rVB4	9158	14275	0.54%	0.062%
4	4.966	324	328	335	rVB	14920	21250	0.81%	0.092%
5	7.048	676	682	692	rBV	65440	138765	5.26%	0.598%
6	7.201	702	708	722	rBV	316326	547740	20.75%	2.361%
7	7.407	737	743	758	rBV	336868	574316	21.76%	2.476%
8	7.878	817	823	833	rBV	348186	579000	21.94%	2.496%
9	8.590	938	944	958	rBV	220563	417725	15.83%	1.801%
10	9.042	1014	1021	1043	rBV	375554	690231	26.15%	2.976%
11	9.213	1047	1050	1053	rBV5	2469	3440	0.13%	0.015%
12	9.489	1092	1097	1107	rVB	12071	19809	0.75%	0.085%
13	9.766	1137	1144	1160	rBV	273311	508109	19.25%	2.191%
14	10.313	1230	1237	1252	rBV	357651	697636	26.43%	3.008%
15	10.560	1274	1279	1283	rVB6	2800	4260	0.16%	0.018%
16	10.689	1294	1301	1315	rVB	458055	781003	29.59%	3.367%
17	10.825	1318	1324	1352	rBV	235856	590067	22.36%	2.544%
18	11.978	1517	1520	1528	rVB6	1557	3269	0.12%	0.014%
19	12.125	1541	1545	1551	rVB5	4095	7143	0.27%	0.031%
20	12.225	1554	1562	1571	rBV2	55081	127856	4.84%	0.551%
21	12.936	1679	1683	1687	rVB4	3038	4159	0.16%	0.018%
22	13.083	1703	1708	1710	rBV5	2444	3078	0.12%	0.013%
23	13.148	1716	1719	1722	rVB3	2914	3386	0.13%	0.015%
24	13.366	1751	1756	1761	rBV2	9924	15897	0.60%	0.069%
25	13.777	1821	1826	1828	rBV5	1627	2949	0.11%	0.013%
26	13.936	1846	1853	1865	rBV	820136	1235771	46.83%	5.328%
27	14.013	1865	1866	1870	rVB3	3833	4500	0.17%	0.019%
28	14.219	1891	1901	1915	rBV	938912	1433462	54.32%	6.180%
29	14.436	1934	1938	1943	rVB5	3518	6079	0.23%	0.026%
30	14.524	1945	1953	1963	rVV	641565	954088	36.15%	4.113%
31	14.607	1965	1967	1972	rVB6	2340	2727	0.10%	0.012%
32	14.766	1988	1994	1997	rBV7	5317	11329	0.43%	0.049%
33	15.324	2085	2089	2096	rVB7	3500	6590	0.25%	0.028%
34	15.513	2115	2121	2136	rBV	1187454	1873879	71.00%	8.079%
35	15.630	2136	2141	2155	rVV	307757	549506	20.82%	2.369%
36	15.754	2159	2162	2171	rVB6	12885	24329	0.92%	0.105%

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37	16.301	2251	2255	2261	rBV8	3957	7331	0.28%	0.032%
38	16.695	2318	2322	2324	rBV4	1932	2679	0.10%	0.012%
39	17.142	2397	2398	2402	rBV3	2793	2756	0.10%	0.012%
40	17.195	2404	2407	2413	rBV6	2875	4209	0.16%	0.018%
41	17.265	2413	2419	2427	rBV	751890	1069120	40.51%	4.609%
42	17.365	2430	2436	2452	rVV	1396911	2137265	80.98%	9.214%
43	17.948	2531	2535	2541	rVB	35044	45072	1.71%	0.194%
44	18.165	2568	2572	2580	rBV2	56566	101192	3.83%	0.436%
45	18.318	2597	2598	2602	rVB4	2590	2851	0.11%	0.012%
46	18.377	2605	2608	2615	rBV4	11645	19873	0.75%	0.086%
47	19.224	2749	2752	2756	rVB3	18308	21125	0.80%	0.091%
48	19.295	2760	2764	2768	rBV	15421	27477	1.04%	0.118%
49	19.448	2787	2790	2796	rBV6	12098	25202	0.95%	0.109%
50	19.653	2820	2825	2832	rBV	1868555	2639115	100.00%	11.378%
51	21.153	3076	3080	3088	rBV2	194893	312572	11.84%	1.348%
52	21.442	3125	3129	3135	rBV	597770	955991	36.22%	4.121%
53	23.671	3501	3508	3526	rBV2	1060761	2503272	94.85%	10.792%
54	23.824	3528	3534	3549	rVB2	571170	1267386	48.02%	5.464%

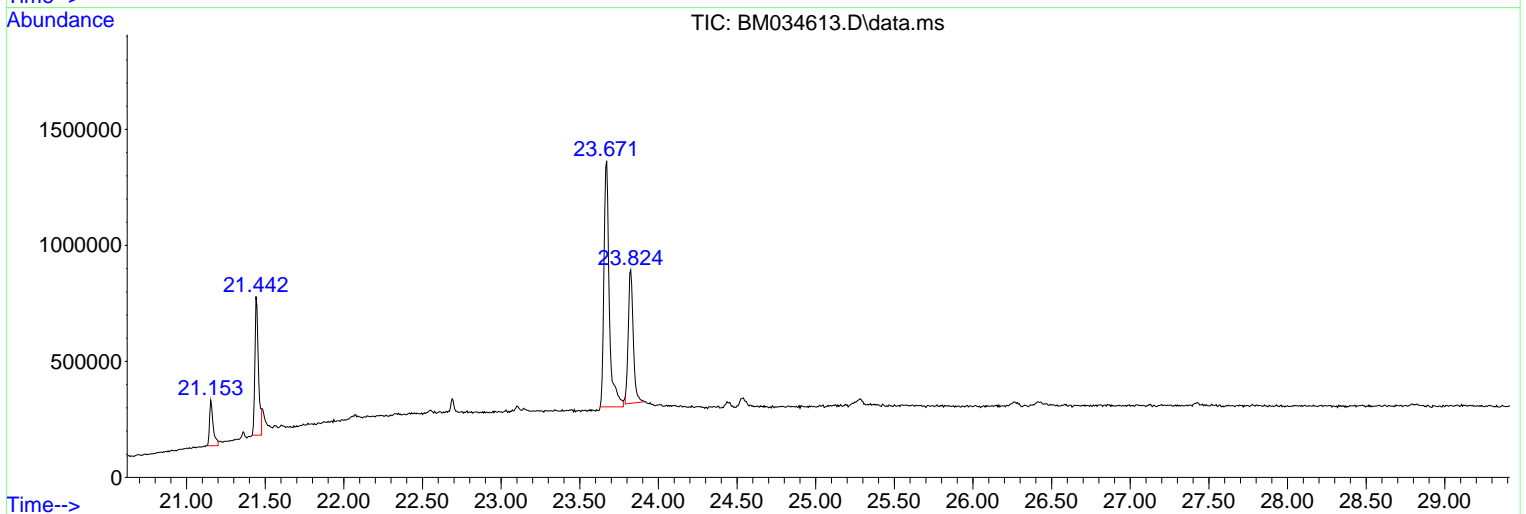
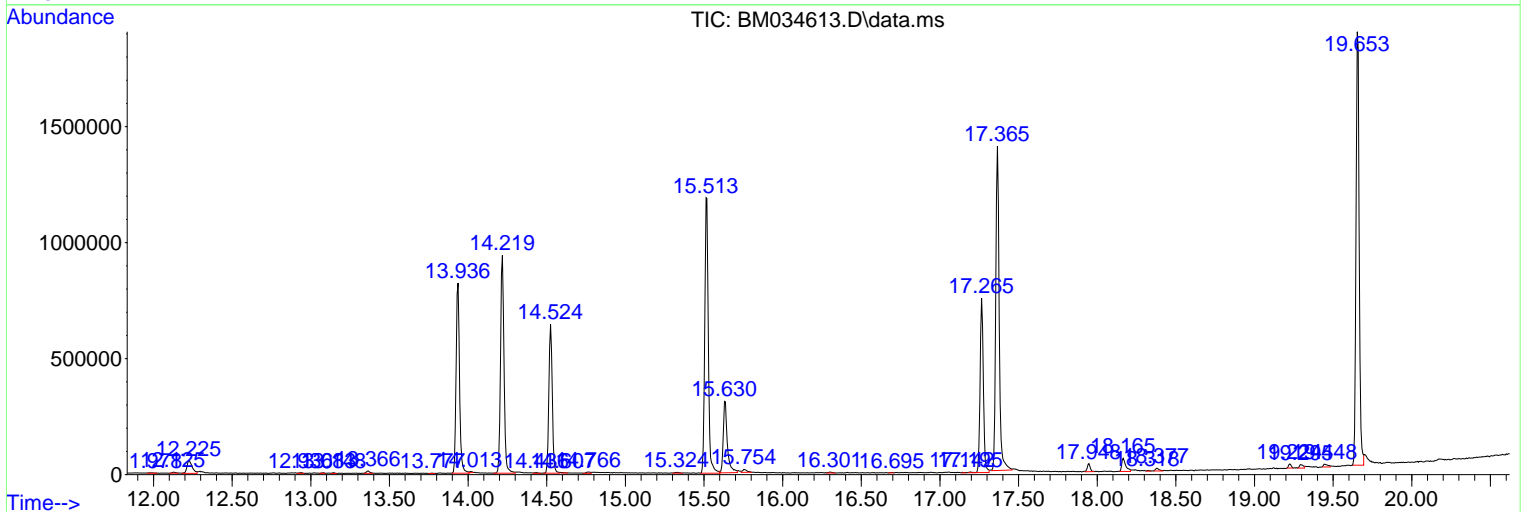
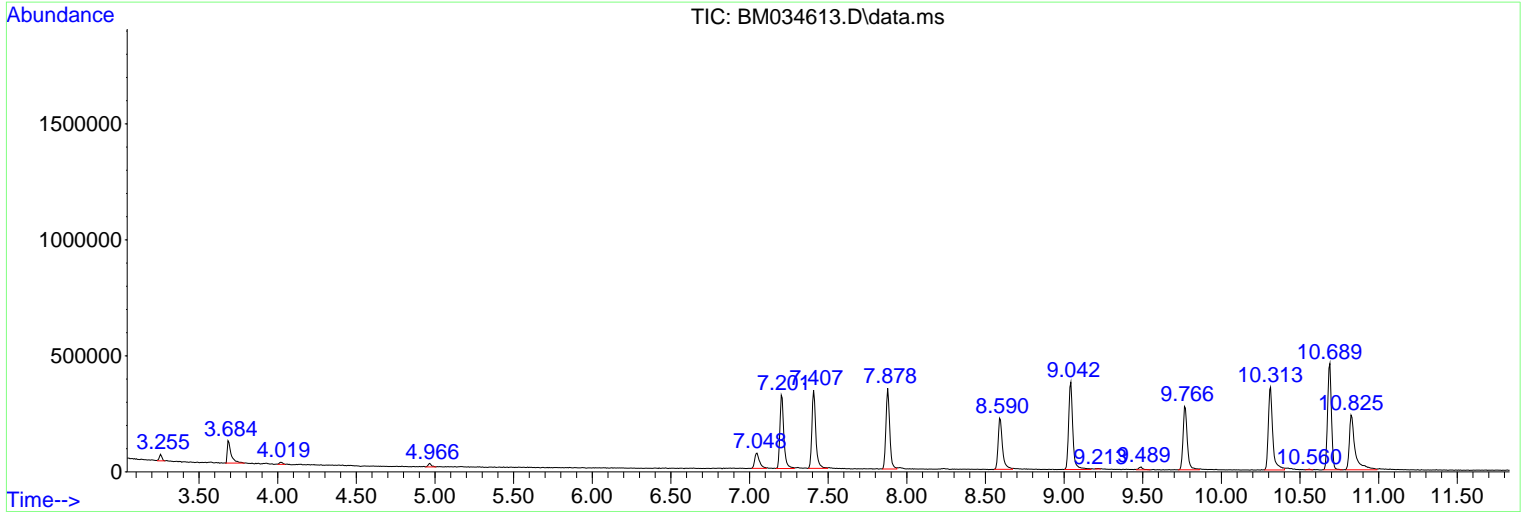
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TIC Library : C:\Database\NIST20.L
 TIC Integration Parameters: LSCINT.P



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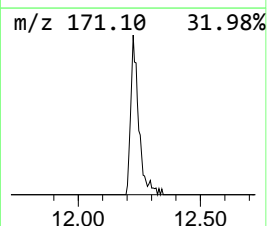
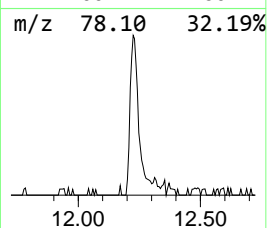
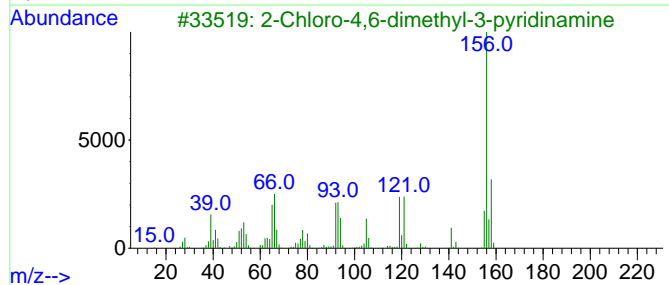
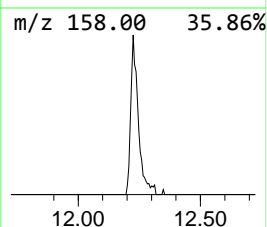
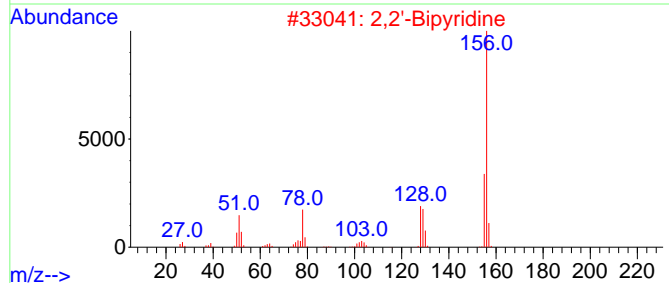
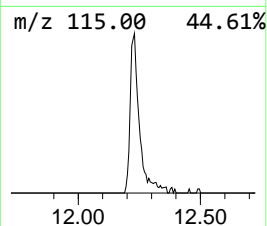
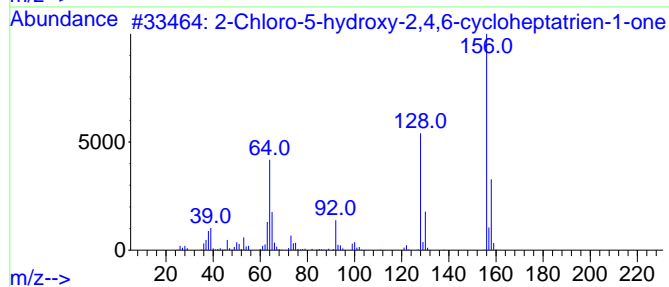
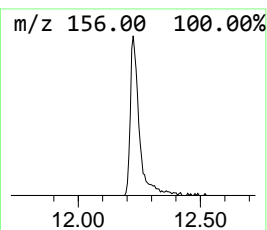
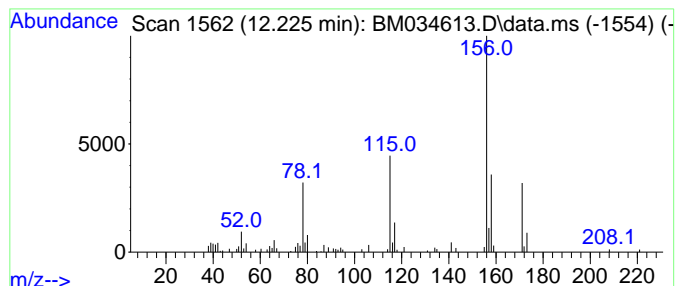
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
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TIC Library : C:\Database\NIST20.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 1 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.225	3.27 ng/ul	127856	Naphthalene-d8	10.689

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Chloro-5-hydroxy-2,4,6-cyclohe...	156	C7H5ClO2	007009-16-7	43
2		2,2'-Bipyridine	156	C10H8N2	000366-18-7	35
3		2-Chloro-4,6-dimethyl-3-pyridina...	156	C7H9ClN2	140413-40-7	32
4		Benzene, 2,5-cyclohexadien-1-yl-	156	C12H12	004794-05-2	28
5		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	25



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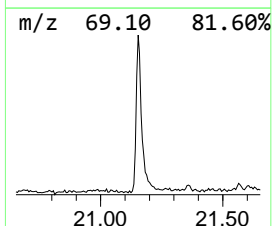
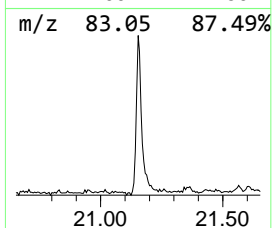
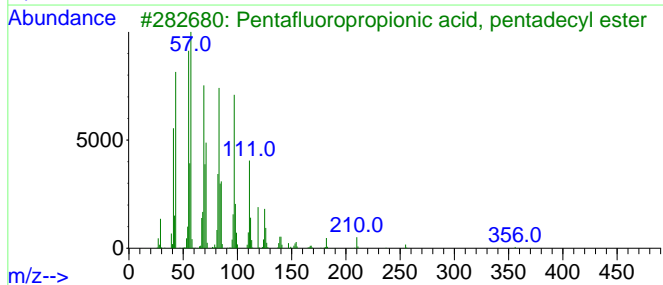
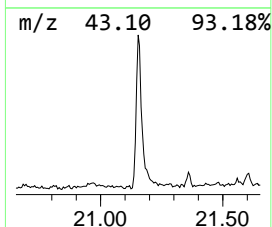
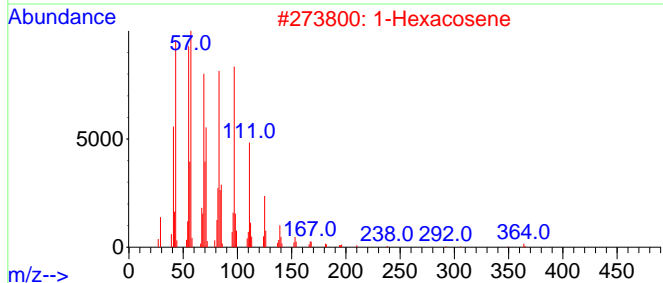
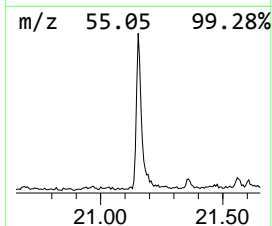
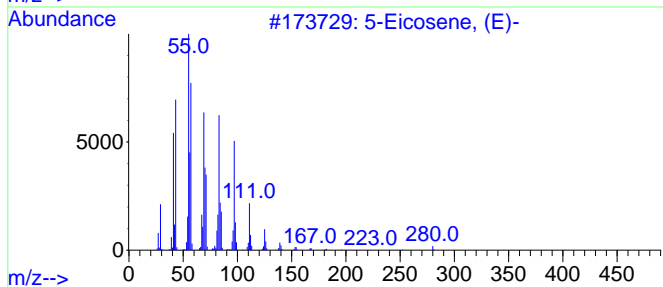
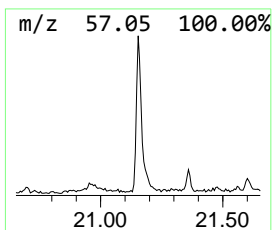
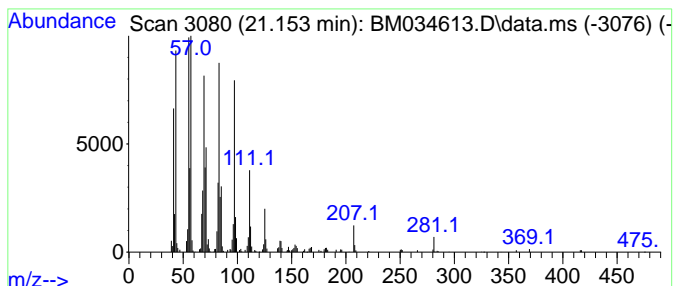
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 Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.153	6.54 ng/ul	312572	Chrysene-d12	21.442

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		5-Eicosene, (E)-	280	C20H40	074685-30-6	97
2		1-Hexacosene	364	C26H52	018835-33-1	94
3		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
4		Pentacos-1-ene	350	C25H50	016980-85-1	94
5		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	12.225	3.3	ng/u1	127856	2	10.689	781003	20.0
(DEL) Alkane: S...	21.153	6.5	ng/u1	312572	5	21.442	955991	20.0