

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
 Data File : BM029814.D
 Acq On : 05 May 2021 07:17
 Operator : CG/JU
 Sample : M2127-15
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E0AB6

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 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-BM050421.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.757	653	658	660	rBV	93405	162932	4.94%	0.607%
2	6.910	679	684	697	rBV	302022	541080	16.40%	2.015%
3	7.104	711	717	730	rBV	382117	660348	20.01%	2.459%
4	7.298	746	750	758	rVB4	15237	28731	0.87%	0.107%
5	7.563	789	795	804	rBV	356318	616631	18.68%	2.297%
6	7.657	806	811	824	rVB	59295	132916	4.03%	0.495%
7	8.081	877	883	895	rVB2	40408	81974	2.48%	0.305%
8	8.286	911	918	933	rBV	206322	440187	13.34%	1.639%
9	8.722	983	992	1010	rBV	376425	732448	22.19%	2.728%
10	9.139	1059	1063	1070	rVB2	14377	22109	0.67%	0.082%
11	9.439	1107	1114	1131	rBV	314243	594607	18.02%	2.215%
12	9.763	1166	1169	1171	rBV4	2920	3730	0.11%	0.014%
13	9.980	1199	1206	1234	rBV	320916	822342	24.92%	3.063%
14	10.339	1259	1267	1278	rBV	475065	846979	25.66%	3.154%
15	10.733	1330	1334	1335	rBV4	3402	4347	0.13%	0.016%
16	11.180	1405	1410	1413	rBV6	5162	8957	0.27%	0.033%
17	11.204	1413	1414	1420	rVV5	4027	5688	0.17%	0.021%
18	11.292	1424	1429	1432	rVV6	2312	4112	0.12%	0.015%
19	11.422	1446	1451	1453	rBV6	2980	3615	0.11%	0.013%
20	11.951	1535	1541	1550	rVB5	6970	17905	0.54%	0.067%
21	12.251	1586	1592	1601	rBV	34325	96573	2.93%	0.360%
22	12.510	1632	1636	1639	rBV6	2094	3618	0.11%	0.013%
23	12.821	1685	1689	1690	rBV3	3617	3693	0.11%	0.014%
24	13.045	1722	1727	1734	rBV7	6027	11580	0.35%	0.043%
25	13.633	1820	1827	1832	rBV	849720	1230980	37.30%	4.585%
26	13.680	1832	1835	1845	rVB	39768	64413	1.95%	0.240%
27	13.904	1863	1873	1894	rBV	1052206	1620231	49.09%	6.034%
28	14.127	1908	1911	1914	rVB4	6015	6997	0.21%	0.026%
29	14.216	1918	1926	1934	rBV	691125	1076975	32.63%	4.011%
30	14.480	1965	1971	1972	rBV5	19183	29375	0.89%	0.109%
31	14.568	1984	1986	1991	rVB6	4060	5743	0.17%	0.021%
32	14.621	1991	1995	1998	rVV5	2364	4485	0.14%	0.017%
33	14.663	2000	2002	2004	rVV3	4514	4349	0.13%	0.016%
34	14.704	2007	2009	2014	rVB5	3702	4854	0.15%	0.018%

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35	14.815	2025	2028	2032	rBV5	3464	6251	0.19%	0.023%
36	14.851	2032	2034	2042	rVB3	7769	12541	0.38%	0.047%
37	14.968	2050	2054	2061	rBV	90714	134998	4.09%	0.503%
38	15.027	2061	2064	2067	rVV2	8796	11579	0.35%	0.043%
39	15.080	2070	2073	2081	rVB7	6895	11391	0.35%	0.042%
40	15.210	2088	2095	2109	rBV	1336083	1922755	58.26%	7.161%
41	15.345	2112	2118	2132	rVV	379697	599349	18.16%	2.232%
42	15.451	2132	2136	2141	rVB3	13479	22348	0.68%	0.083%
43	15.645	2165	2169	2171	rBV3	2536	3455	0.10%	0.013%
44	15.757	2184	2188	2189	rBV3	2705	3568	0.11%	0.013%
45	15.786	2189	2193	2200	rVB8	6188	10414	0.32%	0.039%
46	15.851	2200	2204	2207	rBV5	5203	9327	0.28%	0.035%
47	15.898	2209	2212	2216	rVV6	8462	11419	0.35%	0.043%
48	15.945	2216	2220	2226	rVV4	11234	20453	0.62%	0.076%
49	15.998	2226	2229	2235	rVB8	12070	18461	0.56%	0.069%
50	16.110	2243	2248	2254	rVV8	5464	11304	0.34%	0.042%
51	16.198	2258	2263	2268	rVB5	9803	17311	0.52%	0.064%
52	16.292	2274	2279	2285	rBV10	5318	11005	0.33%	0.041%
53	16.615	2329	2334	2335	rBV5	2809	3459	0.10%	0.013%
54	16.633	2335	2337	2342	rVB6	5194	7371	0.22%	0.027%
55	16.704	2344	2349	2352	rBV5	9489	14977	0.45%	0.056%
56	16.739	2352	2355	2357	rBV4	6332	7589	0.23%	0.028%
57	16.904	2380	2383	2386	rVB5	4386	4922	0.15%	0.018%
58	16.956	2386	2392	2399	rBV	812310	1214756	36.81%	4.524%
59	17.015	2400	2402	2403	rVV2	13061	11206	0.34%	0.042%
60	17.057	2403	2409	2423	rVB	1359490	2005674	60.77%	7.470%
61	17.286	2446	2448	2452	rVB5	6292	5996	0.18%	0.022%
62	17.339	2452	2457	2462	rBV8	12529	18708	0.57%	0.070%
63	17.509	2484	2486	2488	rBV3	3827	4750	0.14%	0.018%
64	17.633	2503	2507	2512	rVB7	16867	24282	0.74%	0.090%
65	17.745	2523	2526	2527	rBV3	4941	5884	0.18%	0.022%
66	17.774	2527	2531	2536	rVB2	18689	28788	0.87%	0.107%
67	17.868	2542	2547	2565	rBV	239079	435089	13.18%	1.620%
68	18.356	2627	2630	2637	rVB3	11544	19034	0.58%	0.071%
69	18.627	2671	2676	2681	rVB	35283	45994	1.39%	0.171%
70	18.709	2686	2690	2696	rBV	21569	33980	1.03%	0.127%
71	18.992	2734	2738	2743	rBV2	14961	28459	0.86%	0.106%

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72	19.156	2757	2766	2777	rBV4	231994	610243	18.49%	2.273%
73	19.239	2778	2780	2786	rVB2	22778	35366	1.07%	0.132%
74	19.356	2794	2800	2808	rBV	1918201	2690866	81.53%	10.022%
75	20.880	3055	3059	3066	rBV	126144	197959	6.00%	0.737%
76	21.144	3099	3104	3116	rBV	1191290	1601968	48.54%	5.966%
77	23.168	3442	3448	3462	rVB	1726608	3300267	100.00%	12.291%
78	23.309	3465	3472	3481	rVV	921082	1759035	53.30%	6.551%

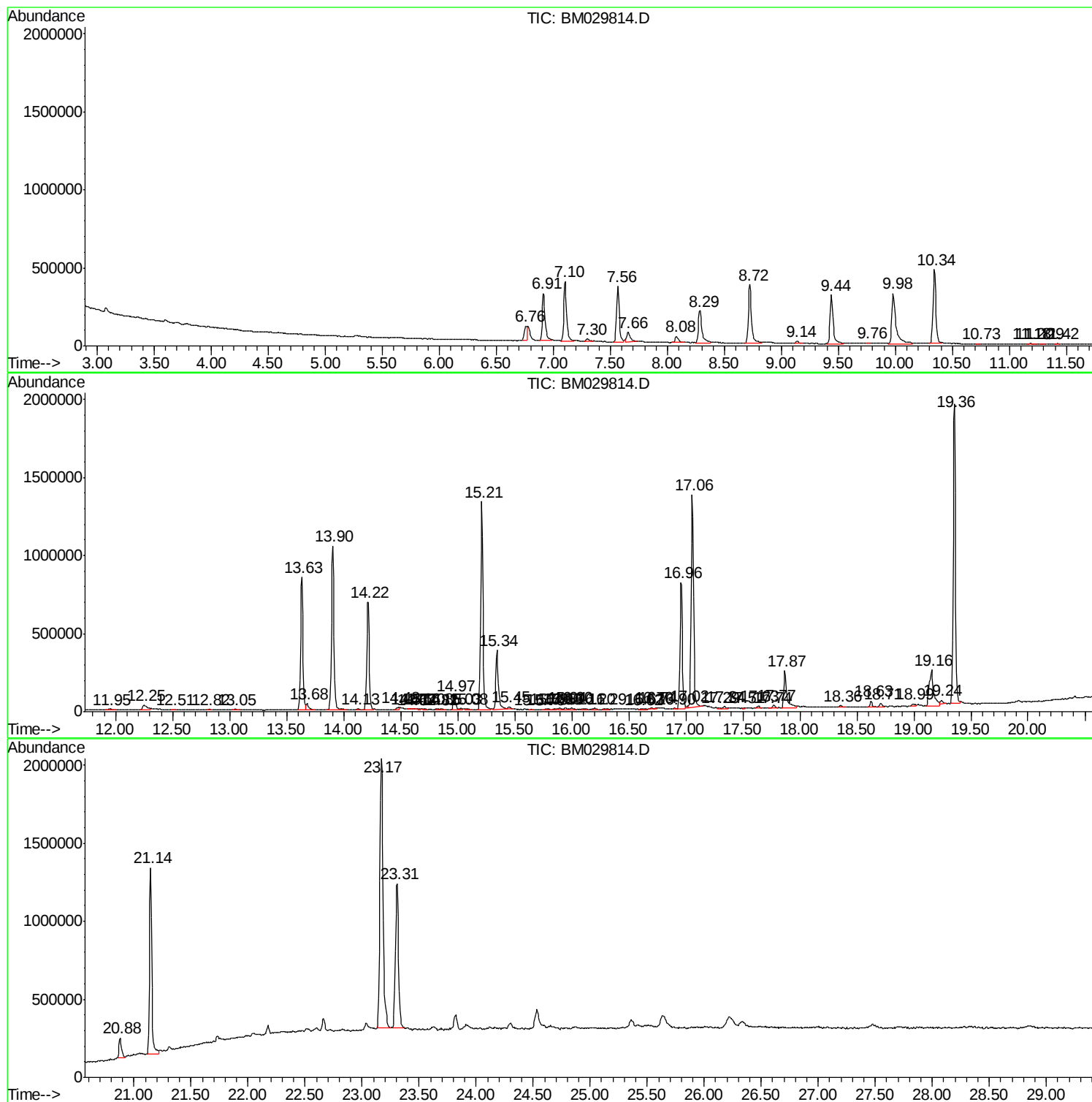
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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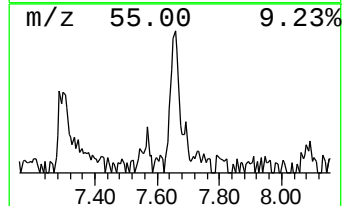
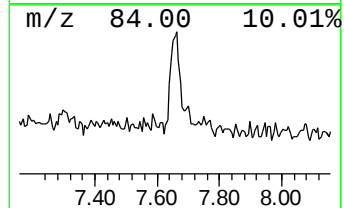
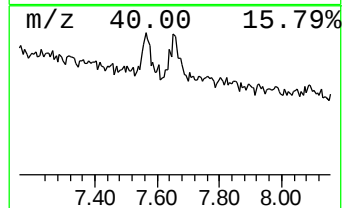
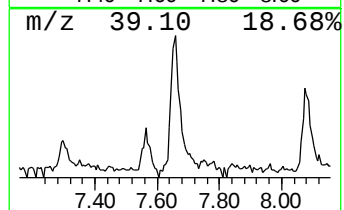
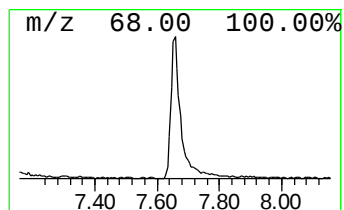
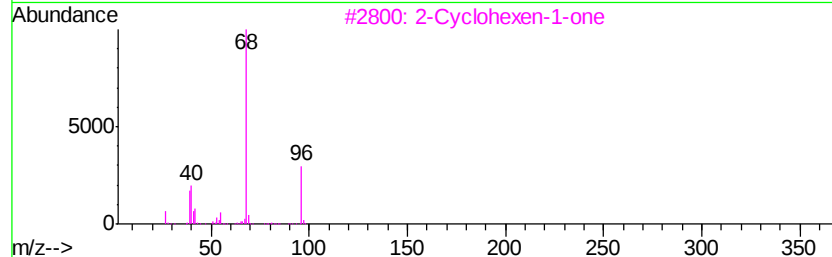
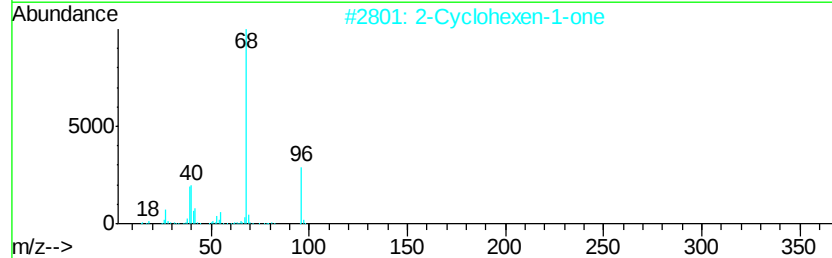
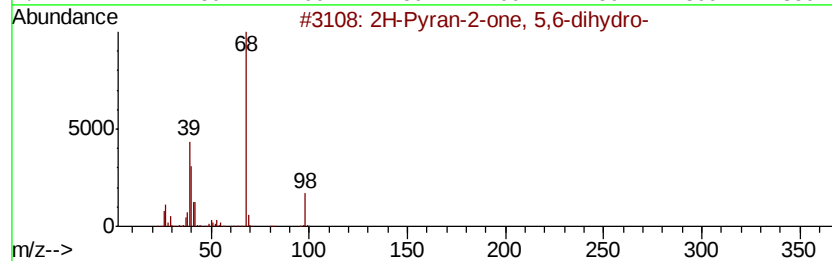
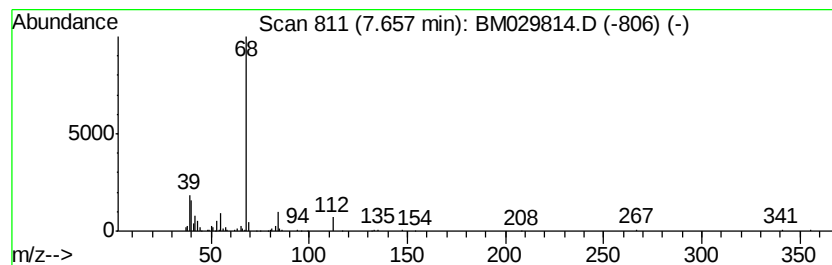
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2H-Pyran-2-one, 5,6-dihydro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	4.31 ng/ul	132916	1,4-Dichlorobenzene-d4	7.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2H-Pyran-2-one, 5,6-dihydro-	98	C5H6O2	003393-45-1	59
2			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	45
3			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	45
4			Furan	68	C4H4O	000110-00-9	42
5			Tetrahydrocyclopenta[1,3]dioxin-...	142	C7H10O3	124898-85-7	39



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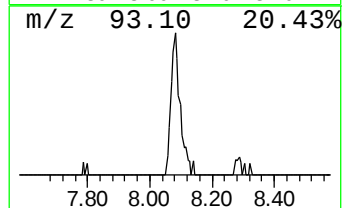
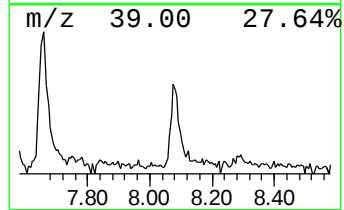
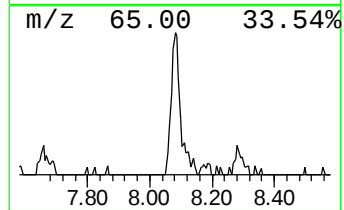
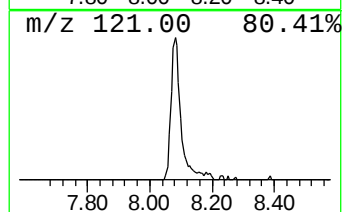
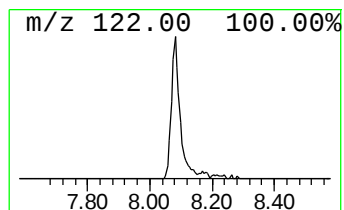
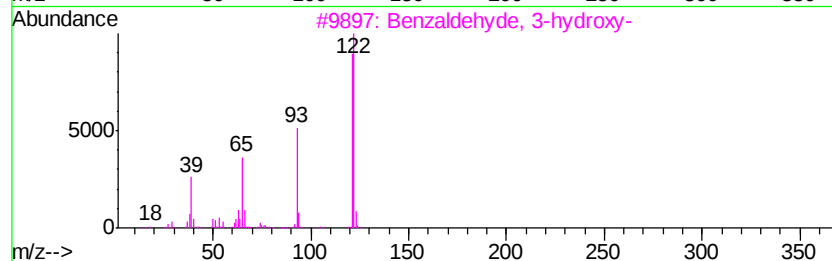
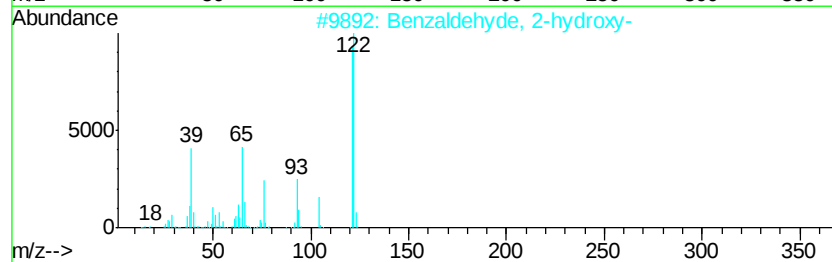
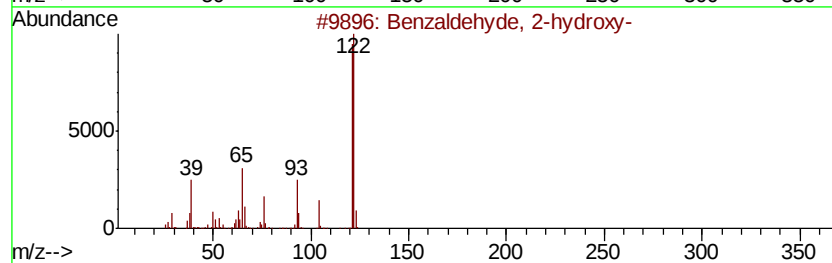
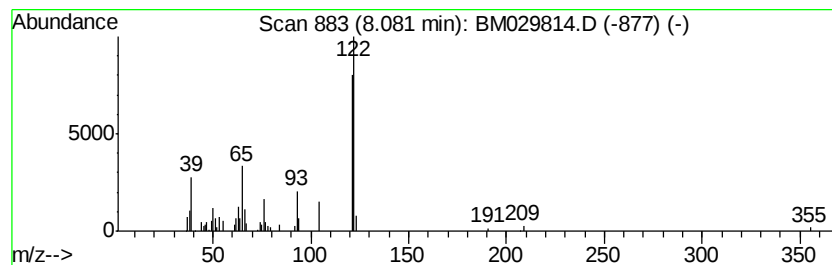
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzaldehyde, 2-hydroxy- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.08	2.66 ng/ul	81974	1,4-Dichlorobenzene-d4	7.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyd, 2-hydroxy-	122	C7H6O2	000090-02-8	96
2			Benzaldehyd, 2-hydroxy-	122	C7H6O2	000090-02-8	95
3			Benzaldehyd, 3-hydroxy-	122	C7H6O2	000100-83-4	89
4			Benzaldehyd, 3-hydroxy-	122	C7H6O2	000100-83-4	89
5			Propanoic acid, 3-chloro-, 4-for...	212	C10H9ClO3	1000142-41-5	78



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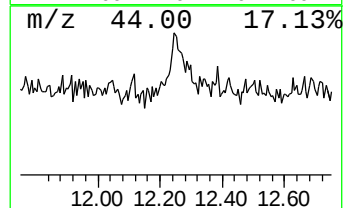
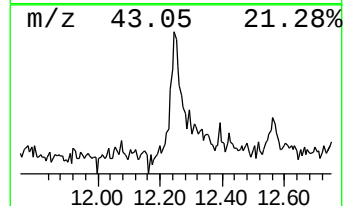
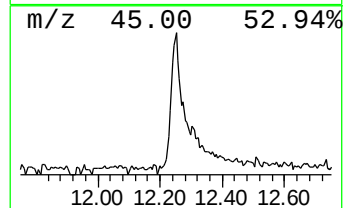
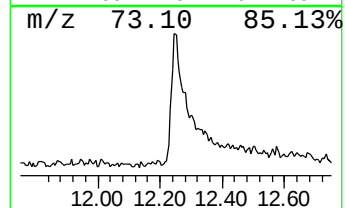
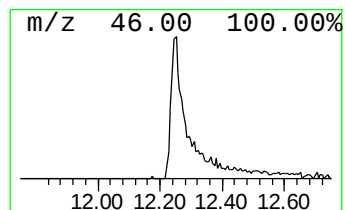
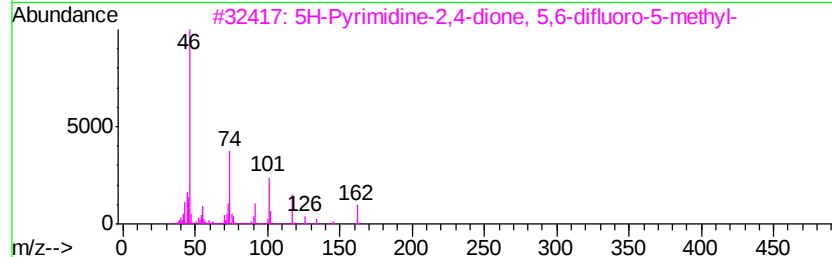
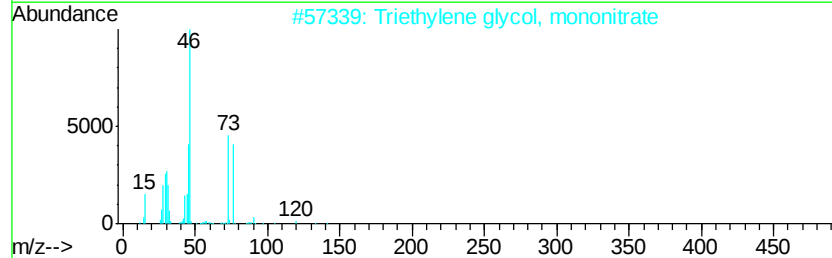
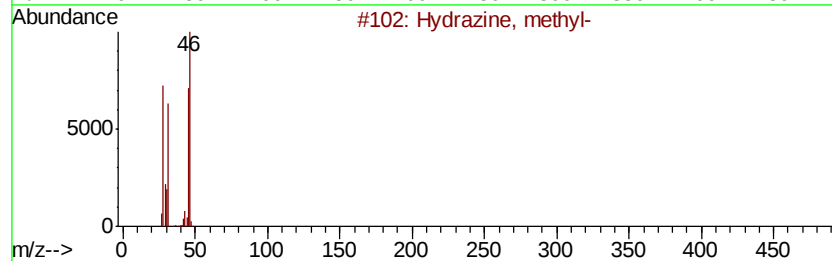
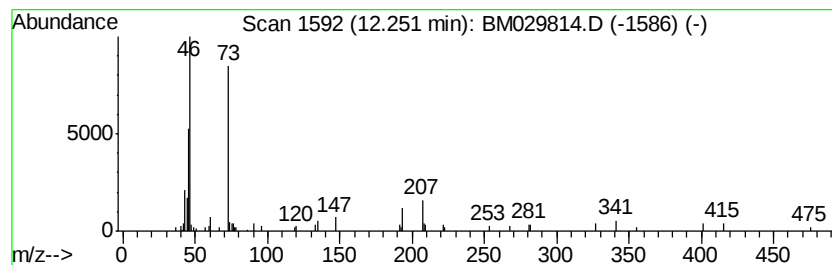
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Peak Number 3 Hydrazine, methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.25	2.28 ng/ul	96573	Naphthalene-d8	10.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hydrazine, methyl-	46	CH6N2	000060-34-4	53
2		Triethylene glycol, mononitrate	195	C6H13N06	020600-87-7	38
3		5H-Pyrimidine-2,4-dione, 5,6-dif...	162	C5H4F2N2O2	1000302-97-4	25
4		Diethylhydroxylamine	89	C4H11NO	003710-84-7	17
5		6-Oxa-3-thiaoctanoic acid	164	C6H12O3S	162514-01-4	10



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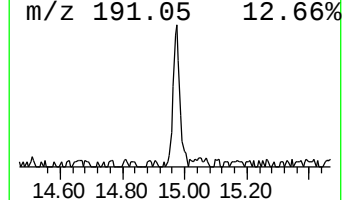
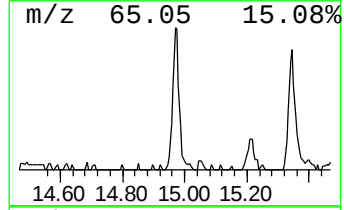
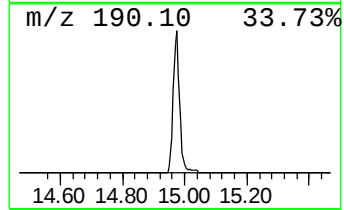
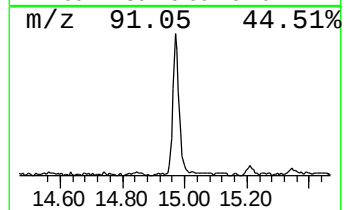
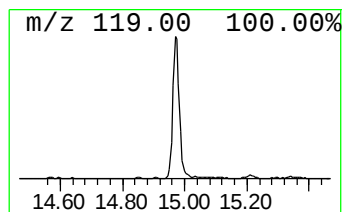
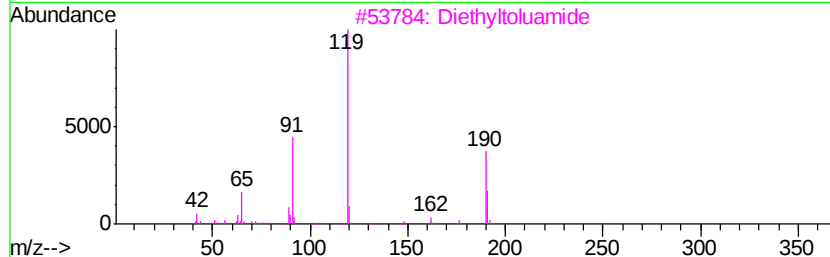
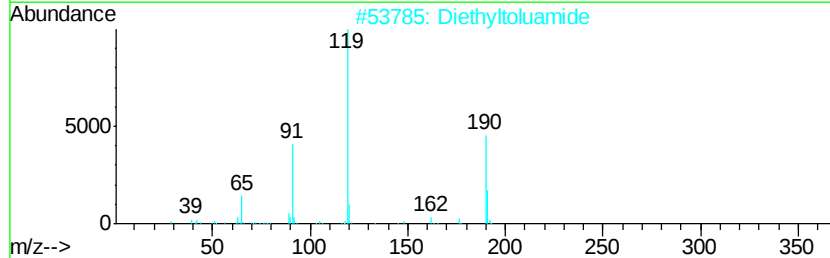
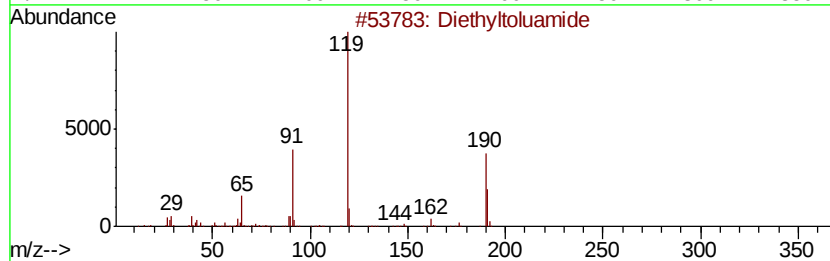
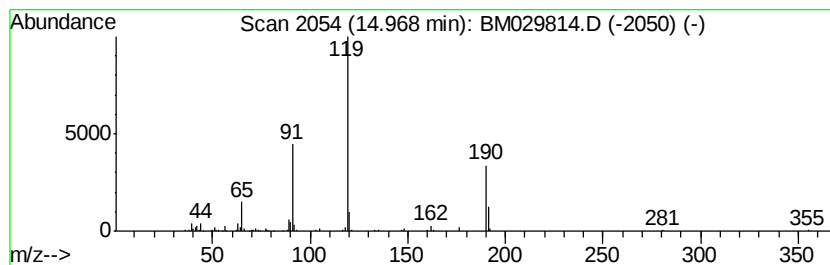
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Diethyltoluamide Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.97	2.51 ng/ul	134998	Acenaphthene-d10	14.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diethyltoluamide	191	C12H17NO	000134-62-3	97
2		Diethyltoluamide	191	C12H17NO	000134-62-3	96
3		Diethyltoluamide	191	C12H17NO	000134-62-3	93
4		Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	90
5		3-(4-Methylbenzoyl)acrylic acid	190	C11H10O3	020972-36-5	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029814.D
Acq On : 05 May 2021 07:17
Operator : CG/JU
Sample : M2127-15
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E0AB6

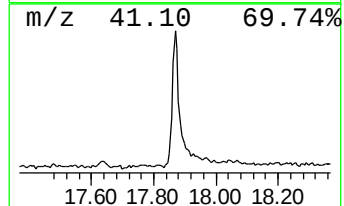
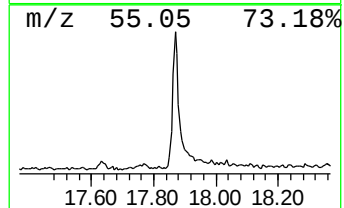
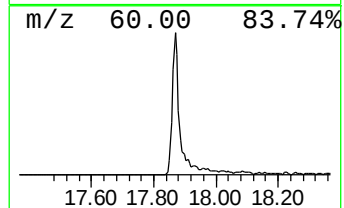
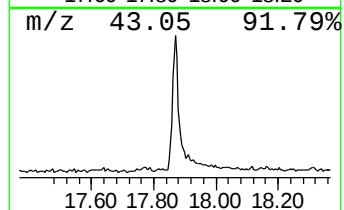
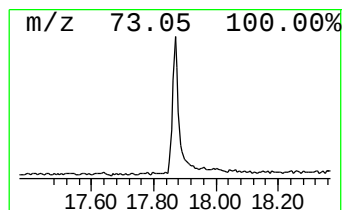
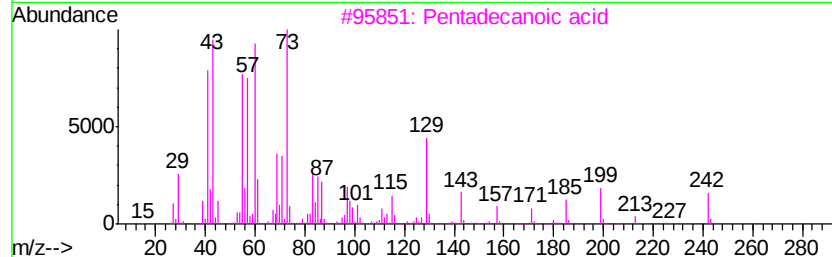
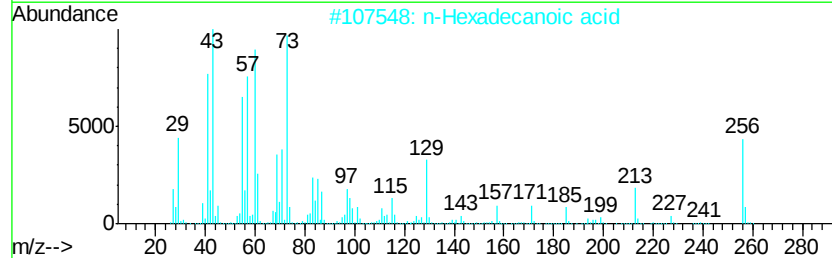
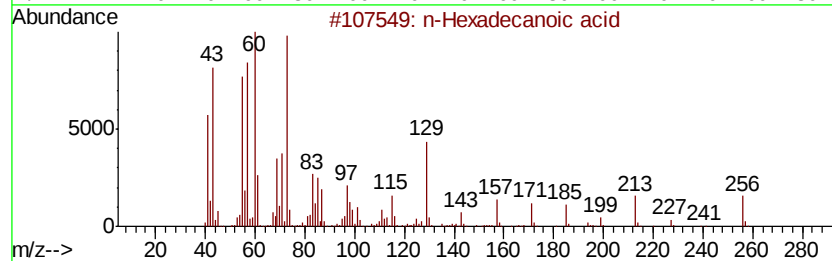
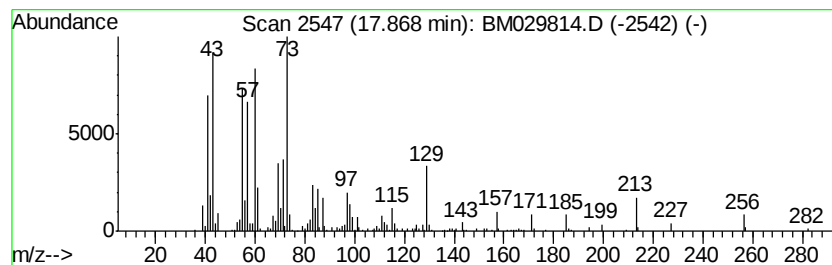
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.87	7.16 ng/ul	435089	Phenanthrene-d10	16.96

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4		Tridecanoic acid	214	C13H26O2	000638-53-9	87
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029814.D
Acq On : 05 May 2021 07:17
Operator : CG/JU
Sample : M2127-15
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E0AB6

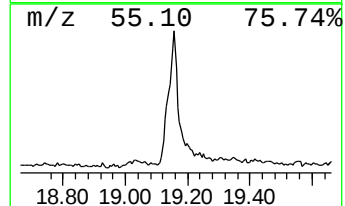
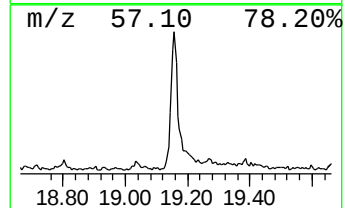
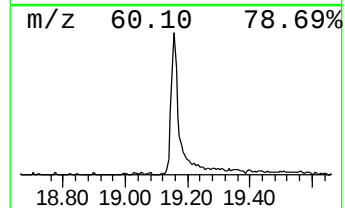
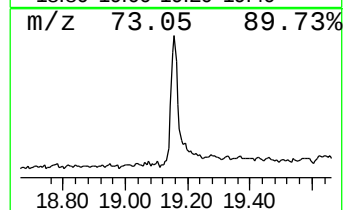
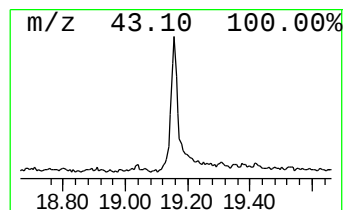
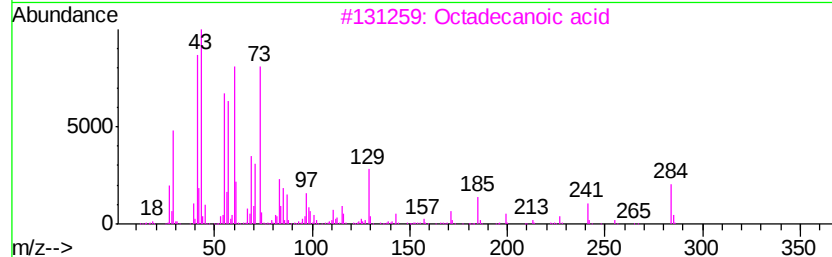
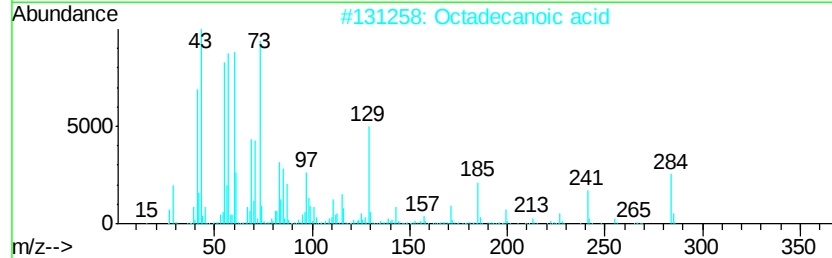
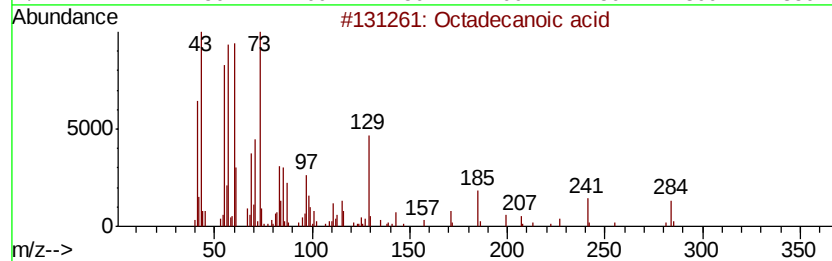
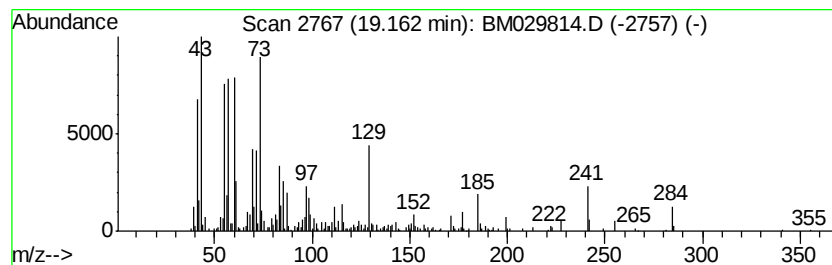
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Octadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.16	7.62 ng/ul	610243	Chrysene-d12	21.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	99
5			Octadecanoic acid	284	C18H36O2	000057-11-4	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029814.D
Acq On : 05 May 2021 07:17
Operator : CG/JU
Sample : M2127-15
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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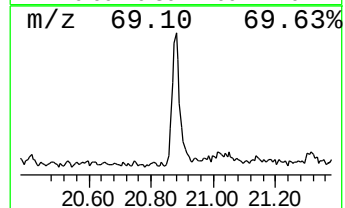
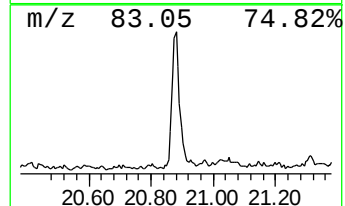
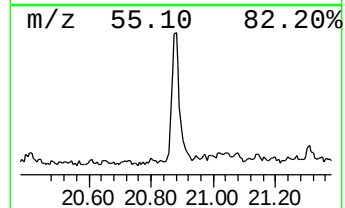
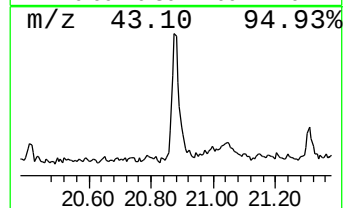
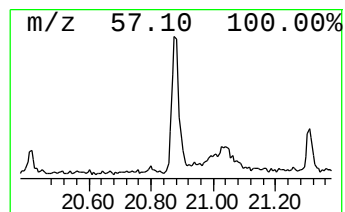
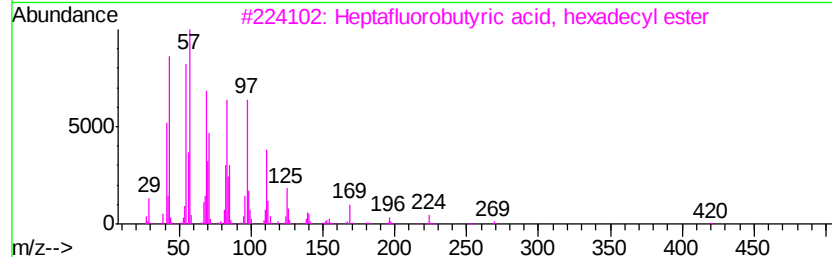
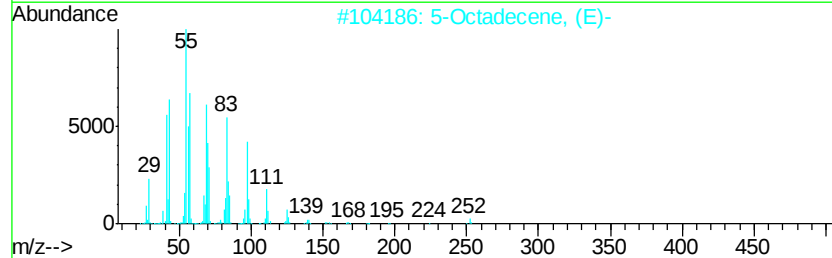
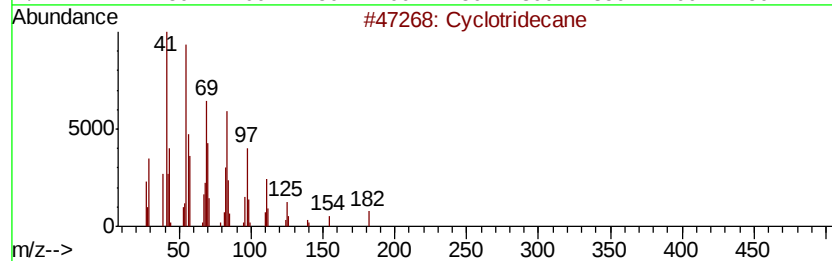
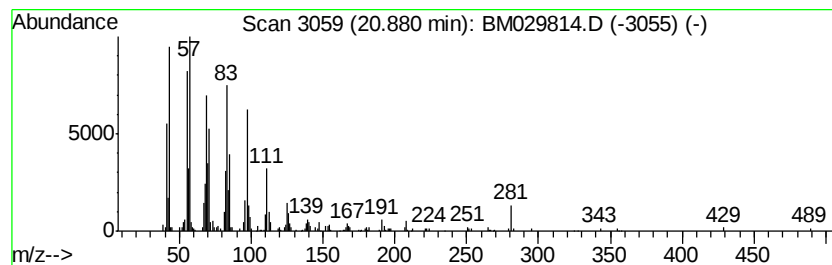
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Cyclic20.88 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.88	2.47 ng/ul	197959	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotridecane	182	C13H26	000295-02-3	95
2		5-Octadecene, (E)-	252	C18H36	007206-21-5	95
3		Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91
4		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
5		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM050421\
Data File : BM029814.D
Acq On : 05 May 2021 07:17
Operator : CG/JU
Sample : M2127-15
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2H-Pyran-2-one, 5...	7.66	4.3	ng/ul	132916	1	7.57	616631	20.0
Benzaldehyde, 2-h...	8.08	2.7	ng/ul	81974	1	7.57	616631	20.0
Hydrazine, methyl-	12.25	2.3	ng/ul	96573	2	10.34	846979	20.0
Diethyltoluamide	14.97	2.5	ng/ul	134998	3	14.22	1076980	20.0
n-Hexadecanoic acid	17.87	7.2	ng/ul	435089	4	16.96	1214760	20.0
Octadecanoic acid	19.16	7.6	ng/ul	610243	5	21.14	1601970	20.0
(DEL) Alkane: Cyc...	20.88	2.5	ng/ul	197959	5	21.14	1601970	20.0