

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
 Data File : BP015873.D
 Acq On : 01 Jul 2023 08:05
 Operator : MA/JU
 Sample : 03239-06
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFY4

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-BP062623.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015873.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	29	32	39	rVB	32461	46938	1.28%	0.079%
2	3.822	104	108	116	rVB	28845	39841	1.08%	0.067%
3	3.875	116	117	119	rVV2	14061	14483	0.39%	0.024%
4	3.922	121	125	132	rVV	147853	208958	5.68%	0.352%
5	3.999	135	138	141	rVV	20606	28604	0.78%	0.048%
6	4.040	141	145	152	rVB	36204	60693	1.65%	0.102%
7	4.140	158	162	168	rVB	17341	29298	0.80%	0.049%
8	4.540	226	230	235	rBV7	7806	14177	0.39%	0.024%
9	4.722	257	261	265	rBV	20990	28894	0.79%	0.049%
10	4.911	287	293	296	rBV8	7054	14539	0.40%	0.025%
11	5.093	322	324	329	rVB2	10106	11945	0.32%	0.020%
12	5.222	342	346	354	rBV8	11820	28590	0.78%	0.048%
13	6.011	474	480	486	rBV10	9683	19253	0.52%	0.032%
14	6.693	592	596	600	rVB5	9315	13845	0.38%	0.023%
15	6.911	627	633	638	rBV10	5580	11286	0.31%	0.019%
16	7.240	681	689	707	rBV	286231	855282	23.25%	1.442%
17	7.381	707	713	731	rVB	342628	683839	18.59%	1.153%
18	7.587	740	748	764	rBV	417351	849317	23.09%	1.432%
19	8.052	819	827	842	rBV	929875	1828463	49.70%	3.082%
20	8.610	918	922	927	rVB8	6599	10937	0.30%	0.018%
21	8.799	946	954	967	rBV2	216821	671813	18.26%	1.132%
22	8.905	967	972	1001	rVB	235163	604371	16.43%	1.019%
23	9.246	1023	1030	1047	rBV	355366	817521	22.22%	1.378%
24	9.393	1052	1055	1063	rVB10	8963	18761	0.51%	0.032%
25	9.587	1083	1088	1093	rVB4	7563	13808	0.38%	0.023%
26	9.975	1145	1154	1171	rBV	211589	652900	17.75%	1.101%
27	10.228	1192	1197	1198	rBV4	17870	20322	0.55%	0.034%
28	10.375	1215	1222	1236	rVB7	18534	67487	1.83%	0.114%
29	10.557	1245	1253	1279	rBV2	272013	967171	26.29%	1.630%
30	10.881	1293	1308	1325	rBV	1299450	2792377	75.90%	4.707%
31	11.087	1337	1343	1358	rBV2	63466	208703	5.67%	0.352%
32	11.516	1408	1416	1418	rBV9	9958	22614	0.61%	0.038%
33	11.557	1418	1423	1427	rVV7	7935	16555	0.45%	0.028%
34	11.781	1456	1461	1467	rVB8	7699	19586	0.53%	0.033%
35	11.875	1472	1477	1481	rVV7	10533	16648	0.45%	0.028%
36	12.281	1541	1546	1550	rVB5	8393	14549	0.40%	0.025%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
 Data File : BP015873.D
 Acq On : 01 Jul 2023 08:05
 Operator : MA/JU
 Sample : 03239-06
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFY4

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-BP062623.MA.M
 Title : SVOA CALIBRATION

37	12.481	1576	1580	1589	rVB9	8972	16189	0.44%	0.027%
38	12.563	1589	1594	1604	rBV	15012	36238	0.99%	0.061%
39	12.769	1625	1629	1635	rVB2	10672	21869	0.59%	0.037%
40	13.063	1673	1679	1687	rVB9	13686	41347	1.12%	0.070%

41	13.210	1699	1704	1711	rBV6	11752	20990	0.57%	0.035%
42	13.298	1713	1719	1723	rBV4	15649	25588	0.70%	0.043%
43	13.446	1739	1744	1748	rBV8	11562	16335	0.44%	0.028%
44	13.498	1749	1753	1758	rVB2	17255	24966	0.68%	0.042%
45	13.575	1758	1766	1772	rBV3	41031	92440	2.51%	0.156%

46	13.881	1814	1818	1819	rBV4	9271	11682	0.32%	0.020%
47	14.016	1835	1841	1847	rBV7	18119	44579	1.21%	0.075%
48	14.116	1847	1858	1871	rBV	866348	1605830	43.65%	2.707%
49	14.204	1871	1873	1877	rVB5	7404	10538	0.29%	0.018%
50	14.293	1886	1888	1893	rVB6	13351	14916	0.41%	0.025%

51	14.398	1897	1906	1921	rBV2	1128133	2006179	54.53%	3.382%
52	14.522	1923	1927	1931	rVB6	25488	39324	1.07%	0.066%
53	14.563	1931	1934	1939	rVB2	34109	45968	1.25%	0.077%
54	14.645	1943	1948	1951	rBV7	7292	15173	0.41%	0.026%
55	14.698	1951	1957	1966	rBV	2056466	3383947	91.99%	5.704%

56	14.922	1988	1995	2001	rBV	19308	36434	0.99%	0.061%
57	15.016	2005	2011	2023	rBV4	63395	264137	7.18%	0.445%
58	15.116	2025	2028	2036	rVB5	38045	68798	1.87%	0.116%
59	15.469	2085	2088	2092	rBV3	16219	20288	0.55%	0.034%
60	15.516	2093	2096	2102	rVB2	29804	45040	1.22%	0.076%

61	15.604	2106	2111	2118	rVB2	8816	20458	0.56%	0.034%
62	15.704	2121	2128	2135	rBV	1268296	2210343	60.08%	3.726%
63	15.881	2152	2158	2171	rBV3	135115	399790	10.87%	0.674%
64	16.016	2178	2181	2185	rVB6	12221	15438	0.42%	0.026%
65	16.069	2185	2190	2198	rBV2	162056	293734	7.98%	0.495%

66	16.151	2201	2204	2209	rVB6	23059	36727	1.00%	0.062%
67	16.257	2218	2222	2226	rVB5	31445	49000	1.33%	0.083%
68	16.404	2240	2247	2253	rBV4	68138	158438	4.31%	0.267%
69	16.551	2268	2272	2278	rBV6	27082	44843	1.22%	0.076%
70	16.616	2278	2283	2289	rVV6	39538	79402	2.16%	0.134%

71	16.763	2306	2308	2314	rVB6	13823	18395	0.50%	0.031%
72	17.175	2375	2378	2382	rBV2	36355	47797	1.30%	0.081%
73	17.281	2394	2396	2402	rVB	23543	30602	0.83%	0.052%
74	17.339	2402	2406	2413	rVB8	25653	44826	1.22%	0.076%
75	17.463	2421	2427	2431	rBV	2219354	3287692	89.37%	5.542%

76	17.504	2431	2434	2440	rVB	493146	723184	19.66%	1.219%
----	--------	------	------	------	-----	--------	--------	--------	--------

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
 Data File : BP015873.D
 Acq On : 01 Jul 2023 08:05
 Operator : MA/JU
 Sample : 03239-06
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFY4

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-BP062623.MA.M
 Title : SVOA CALIBRATION

77	17.569	2440	2445	2450	rBV2	1026709	1734796	47.16%	2.924%
78	17.898	2497	2501	2502	rBV2	34429	49744	1.35%	0.084%
79	18.092	2532	2534	2539	rVB	45801	51064	1.39%	0.086%
80	18.245	2556	2560	2567	rBV2	195125	261259	7.10%	0.440%
81	18.316	2567	2572	2579	rBV2	953560	1446357	39.32%	2.438%
82	18.381	2580	2583	2588	rVB2	110155	186840	5.08%	0.315%
83	18.522	2603	2607	2610	rBV2	84248	116045	3.15%	0.196%
84	18.581	2615	2617	2620	rVV2	48500	49195	1.34%	0.083%
85	18.886	2666	2669	2674	rVB3	85163	116453	3.17%	0.196%
86	19.192	2718	2721	2728	rBV3	101646	184711	5.02%	0.311%
87	19.463	2763	2767	2775	rVB	1180541	1710965	46.51%	2.884%
88	19.539	2776	2780	2786	rBV	426821	578130	15.72%	0.974%
89	19.792	2818	2823	2825	rBV2	1205082	1713304	46.57%	2.888%
90	19.816	2825	2827	2836	rVB	762368	1018205	27.68%	1.716%
91	20.263	2900	2903	2907	rBV3	199631	281562	7.65%	0.475%
92	21.222	3059	3066	3072	rBV4	750618	1596571	43.40%	2.691%
93	21.551	3116	3122	3126	rBV2	2079514	3525099	95.82%	5.942%
94	21.927	3183	3186	3203	rVB3	1273189	3144340	85.47%	5.300%
95	22.127	3217	3220	3223	rBV	301696	381038	10.36%	0.642%
96	22.174	3225	3228	3240	rVB	632273	1196025	32.51%	2.016%
97	23.233	3402	3408	3415	rVB	1966382	3331937	90.57%	5.616%
98	23.316	3417	3422	3436	rVB	920149	2644060	71.87%	4.457%
99	24.092	3548	3554	3571	rVB	1382632	3235977	87.96%	5.455%
100	24.663	3645	3651	3659	rVB	1719638	3678791	100.00%	6.201%

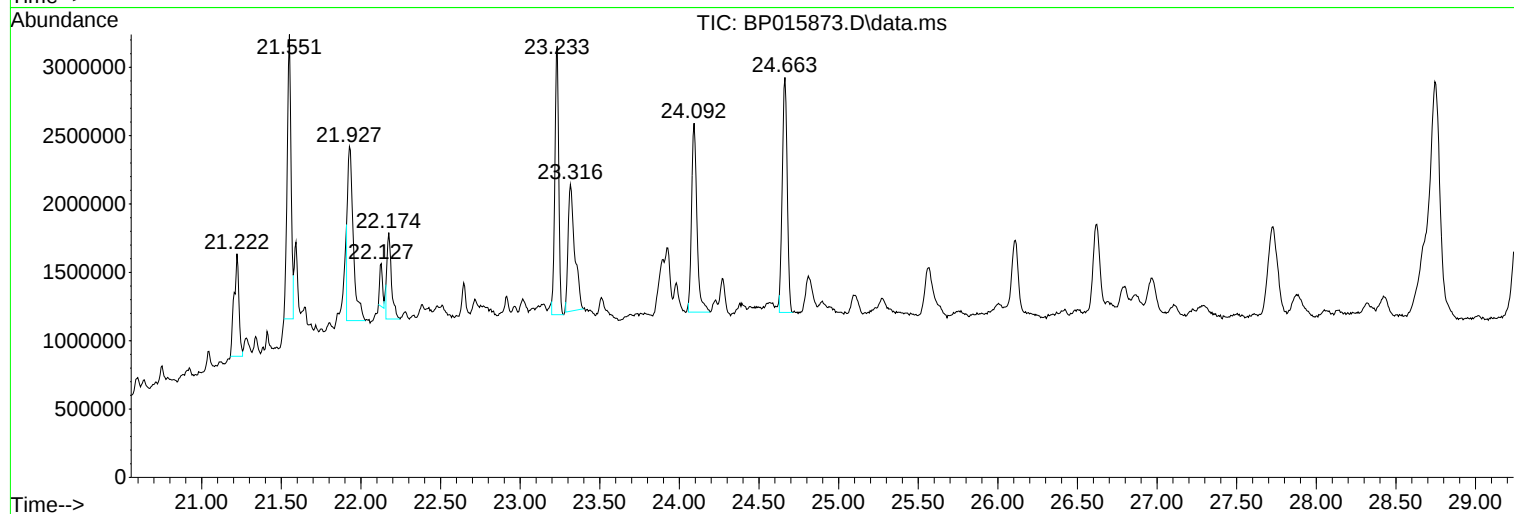
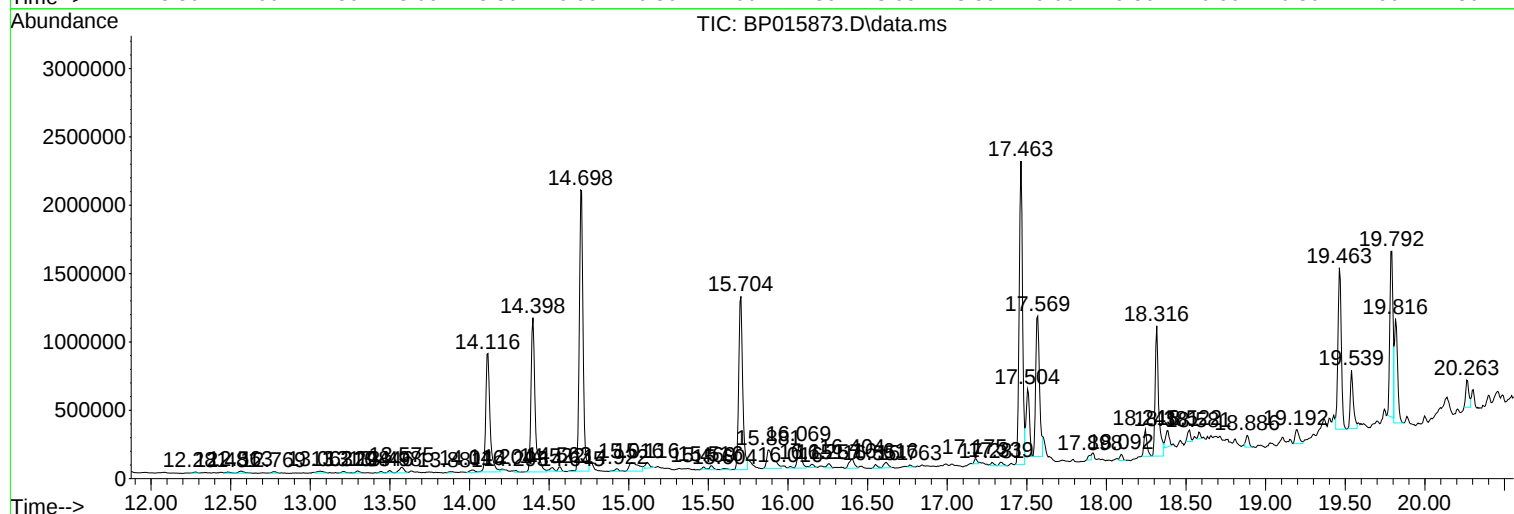
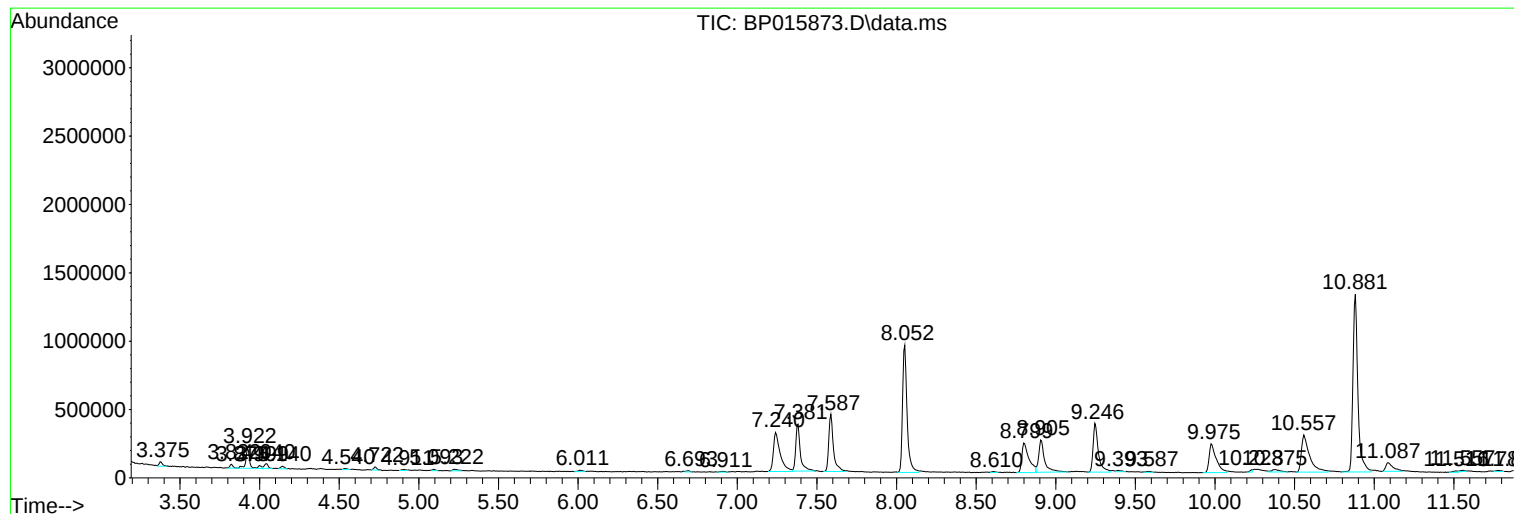
Sum of corrected areas: 59326300

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015873.D
Acq On : 01 Jul 2023 08:05
Operator : MA/JU
Sample : 03239-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFY4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015873.D
Acq On : 01 Jul 2023 08:05
Operator : MA/JU
Sample : 03239-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFY4

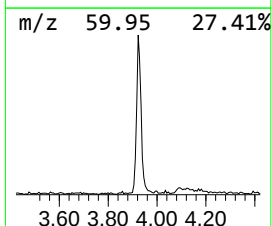
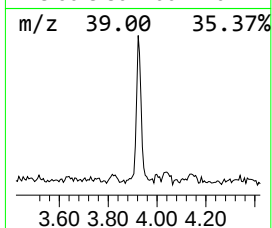
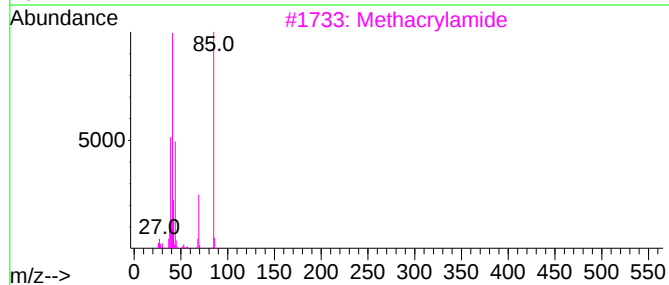
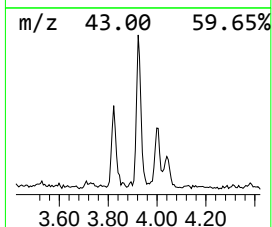
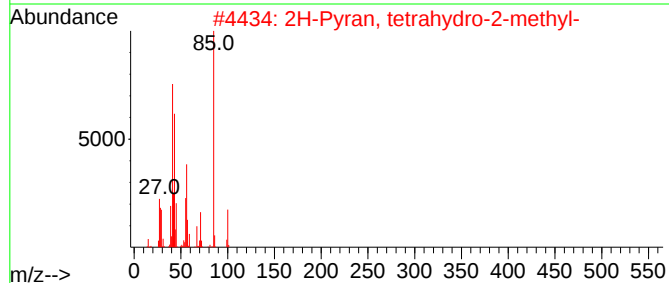
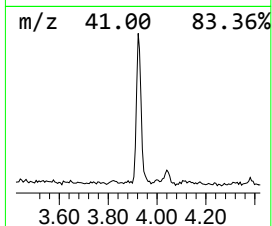
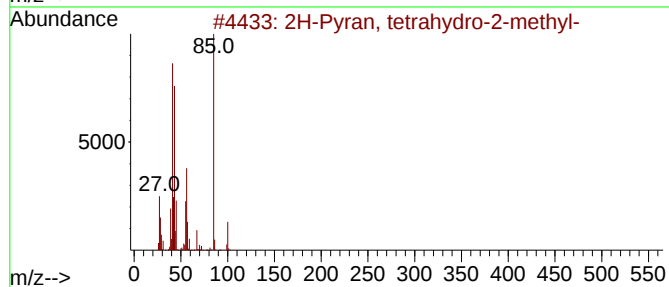
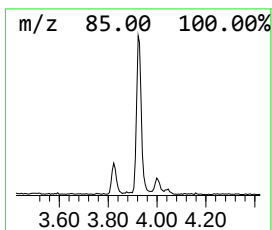
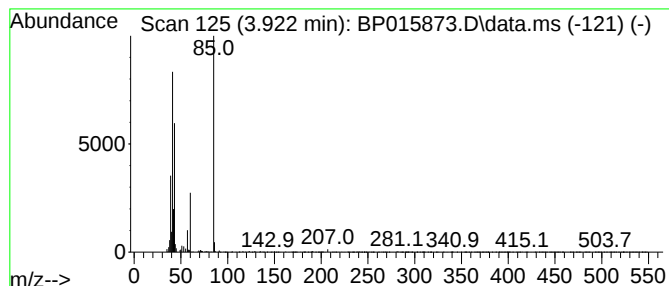
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.922	2.29 ng/ul	208958	1,4-Dichlorobenzene-d4	8.052

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	50
2	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	39
3	Methacrylamide			85	C4H7NO	000079-39-0	38
4	5-(Bromomethyl)oxolan-2-one			178	C5H7BrO2	099393-06-3	9
5	10-(Tetrahydro-pyran-2-yloxy)-tr...			252	C15H24O3	1010192-28-8	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015873.D
Acq On : 01 Jul 2023 08:05
Operator : MA/JU
Sample : 03239-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFY4

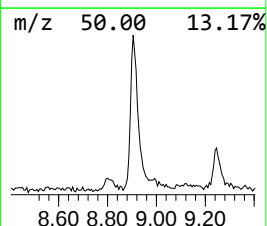
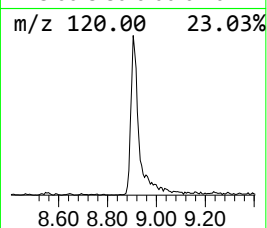
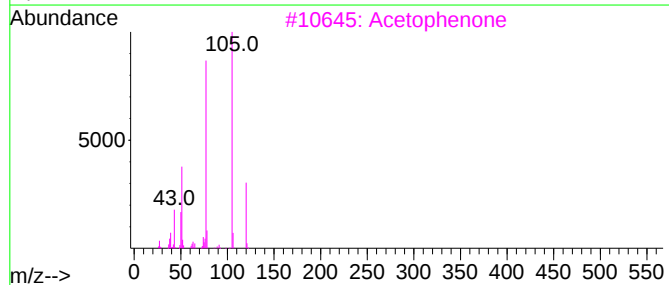
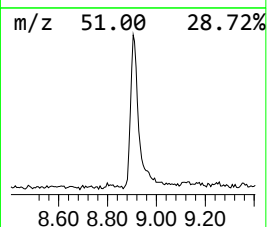
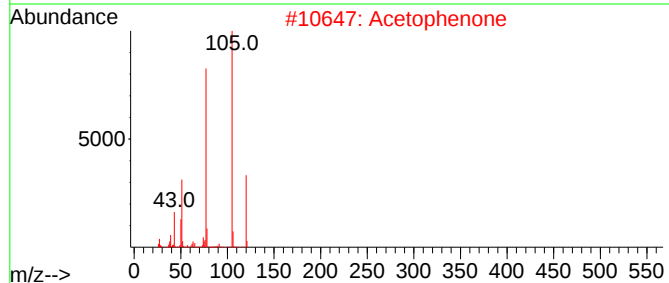
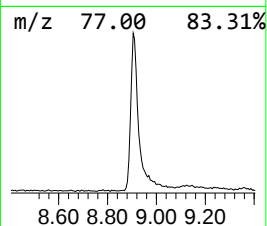
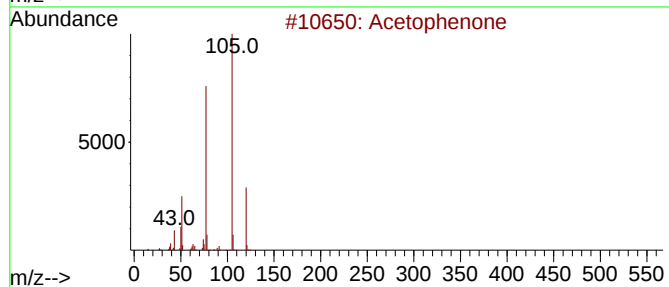
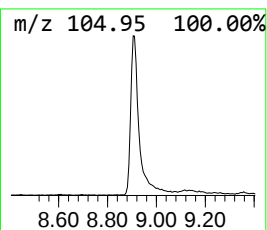
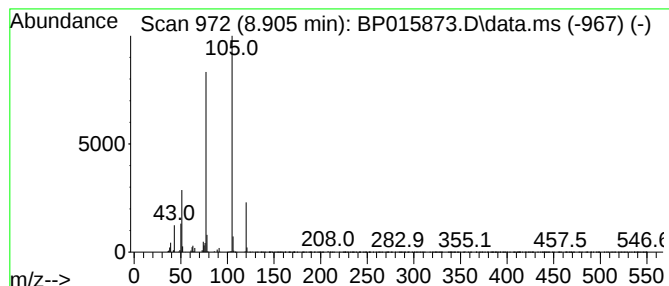
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Aceto phenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.905	6.61 ng/ul	604371	1,4-Dichlorobenzene-d4	8.052

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone			120	C8H8O	000098-86-2	95
2	Acetophenone			120	C8H8O	000098-86-2	94
3	Acetophenone			120	C8H8O	000098-86-2	94
4	Acetophenone			120	C8H8O	000098-86-2	91
5	Acetophenone			120	C8H8O	000098-86-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015873.D
Acq On : 01 Jul 2023 08:05
Operator : MA/JU
Sample : 03239-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFY4

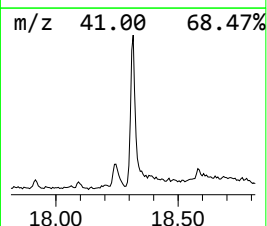
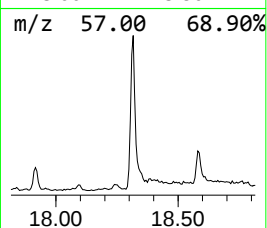
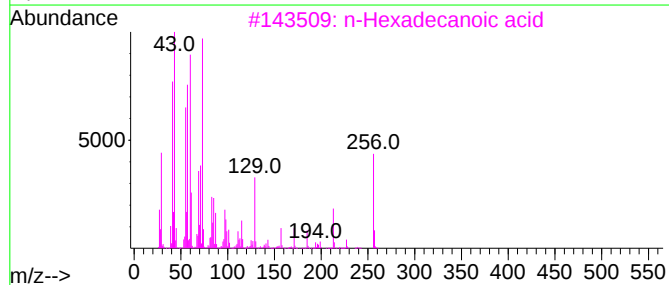
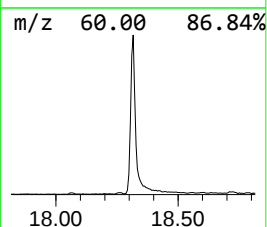
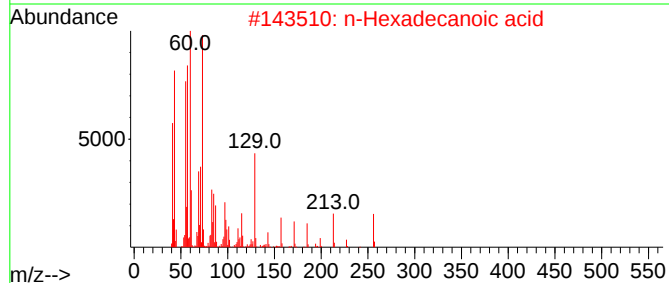
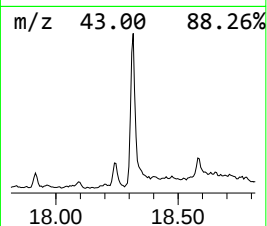
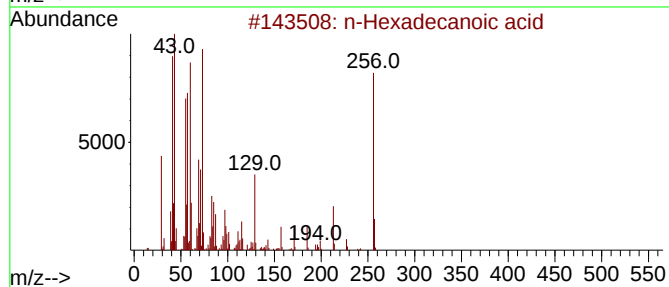
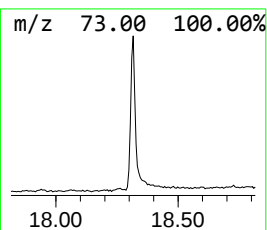
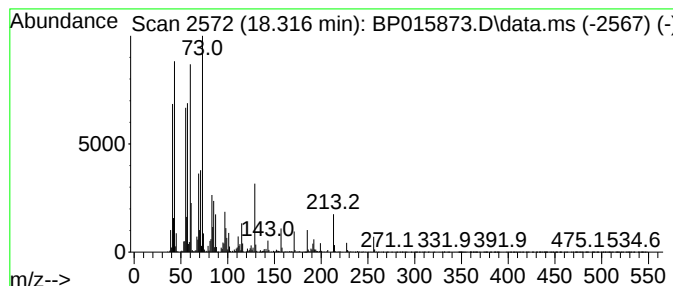
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.316	8.80 ng/ul	1446360	Phenanthrene-d10	17.463

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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015873.D
Acq On : 01 Jul 2023 08:05
Operator : MA/JU
Sample : 03239-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFY4

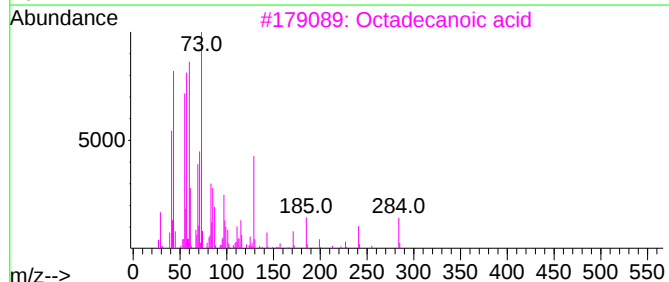
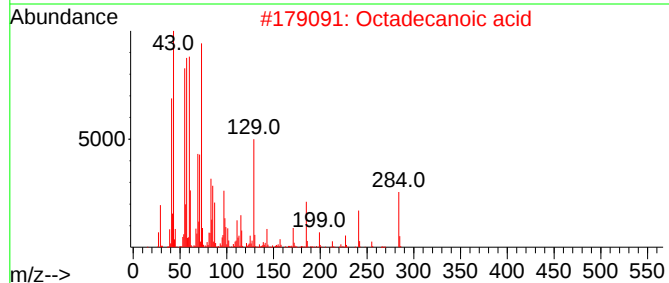
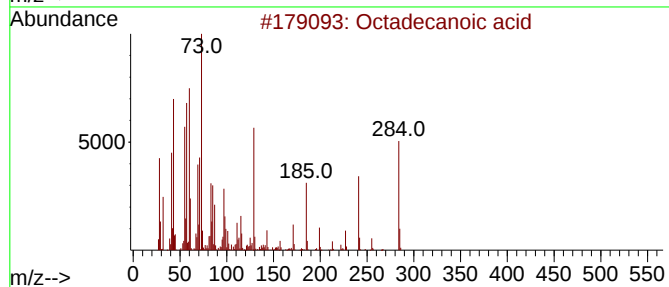
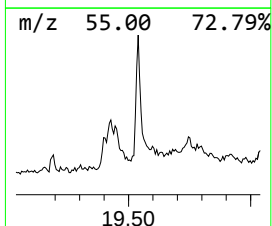
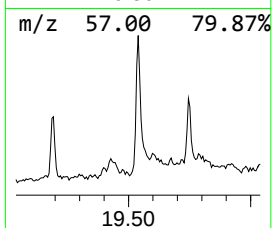
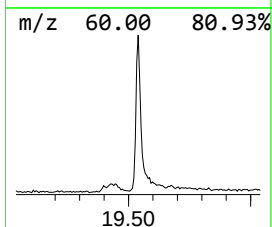
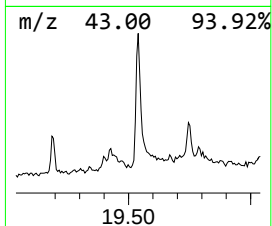
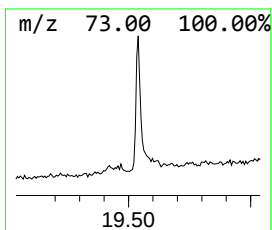
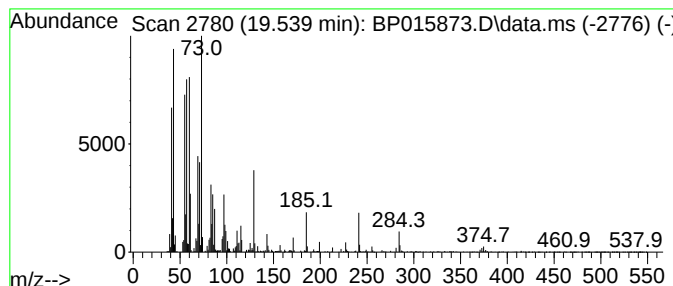
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.539	3.28 ng/ul	578130	Chrysene-d12	21.551

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	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015873.D
Acq On : 01 Jul 2023 08:05
Operator : MA/JU
Sample : 03239-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFY4

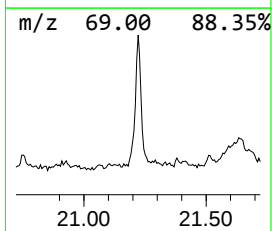
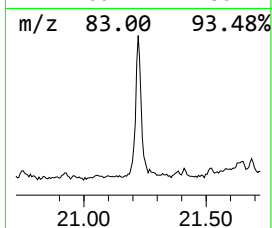
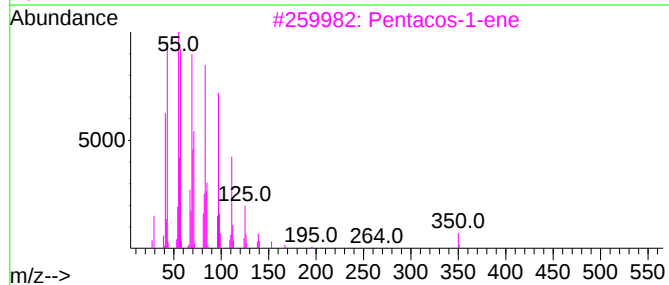
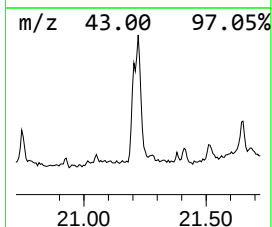
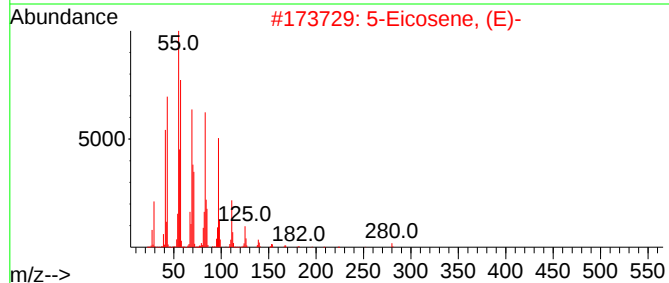
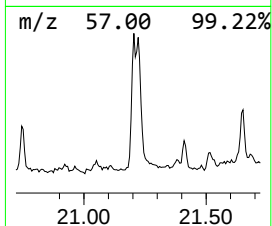
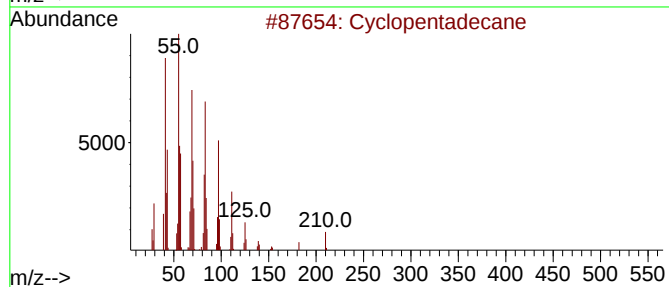
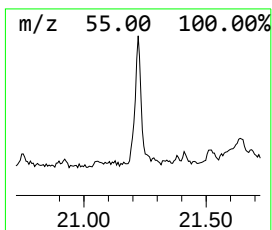
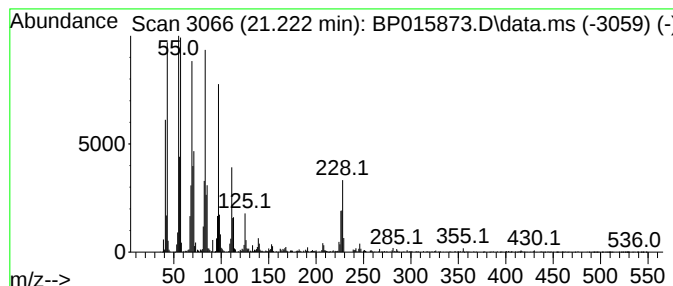
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Cyclic21.222 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.222	9.06 ng/ul	1596570	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentadecane	210	C15H30	000295-48-7	96
2			5-Eicosene, (E)-	280	C20H40	074685-30-6	95
3			Pentacos-1-ene	350	C25H50	016980-85-1	94
4			Cyclohexadecane	224	C16H32	000295-65-8	94
5			n-Tetracosanol-1	354	C24H50O	000506-51-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015873.D
Acq On : 01 Jul 2023 08:05
Operator : MA/JU
Sample : 03239-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFY4

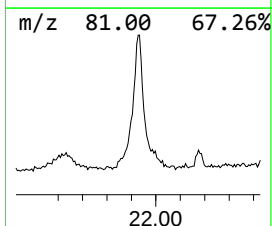
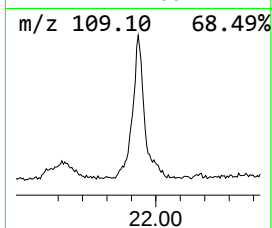
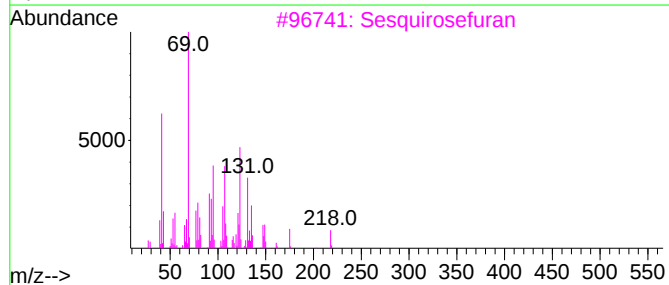
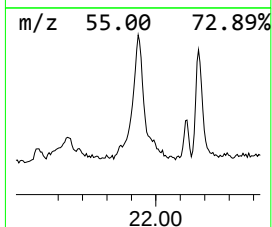
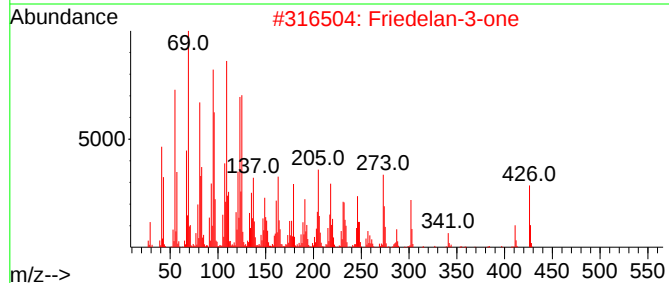
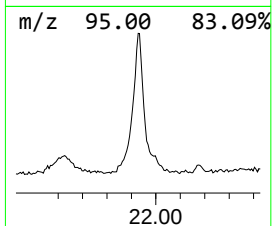
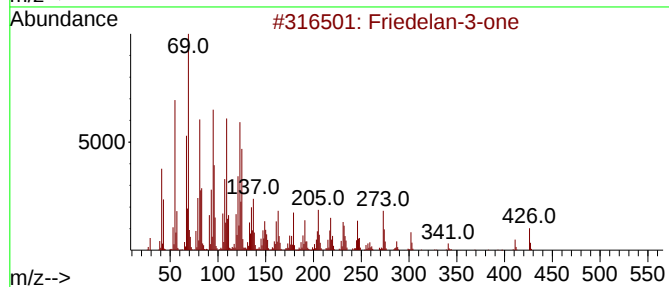
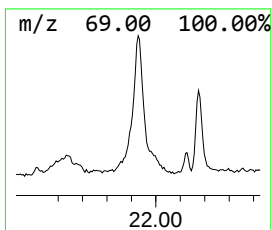
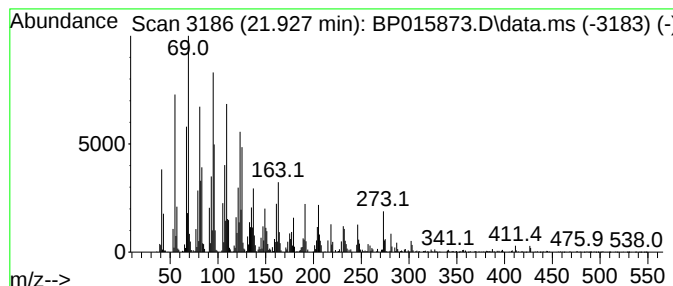
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Sesquirosefuran Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.927	17.84 ng/ul	3144340	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Friedelan-3-one	426	C30H50O	000559-74-0	93
2			Friedelan-3-one	426	C30H50O	000559-74-0	90
3			Sesquirosefuran	218	C15H22O	039007-93-7	84
4			Sesquirosefuran	218	C15H22O	039007-93-7	50
5			(-)-Neoclovene-(I), dihydro-	206	C15H26	1000152-82-1	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015873.D
Acq On : 01 Jul 2023 08:05
Operator : MA/JU
Sample : 03239-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFY4

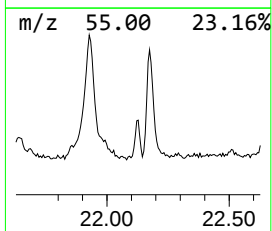
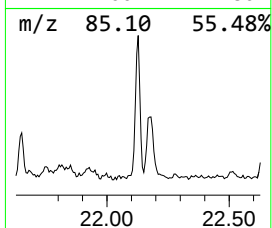
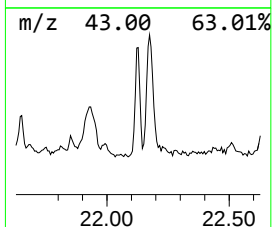
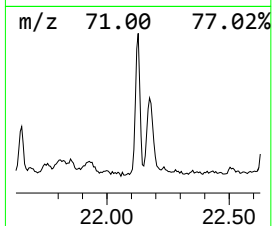
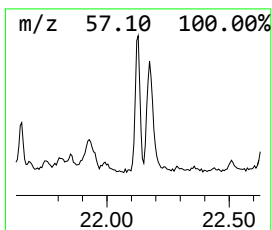
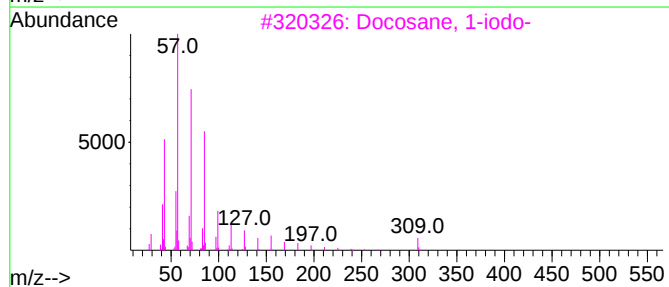
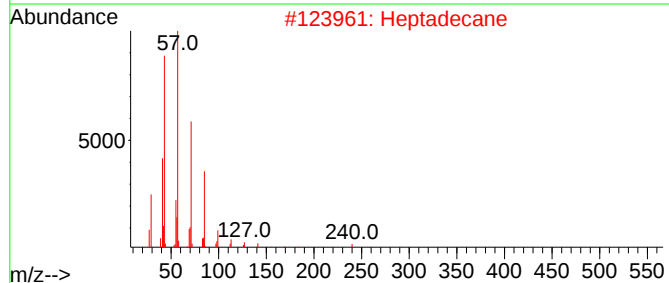
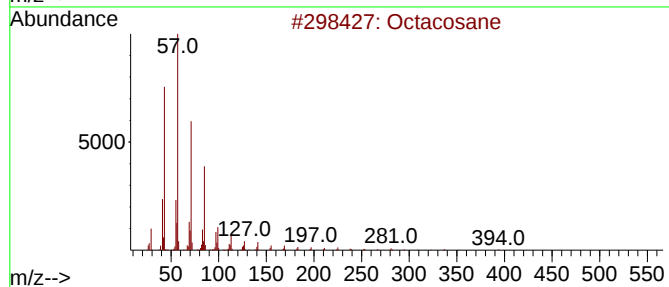
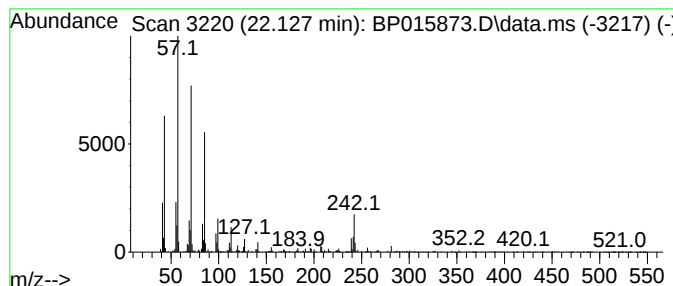
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.127	2.16 ng/ul	381038	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	95
2			Heptadecane	240	C17H36	000629-78-7	93
3			Docosane, 1-iodo-	436	C22H45I	1000406-31-9	83
4			Tetracosane, 1-iodo-	464	C24H49I	1000406-32-0	83
5			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015873.D
Acq On : 01 Jul 2023 08:05
Operator : MA/JU
Sample : 03239-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

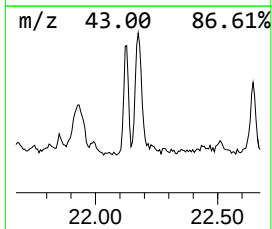
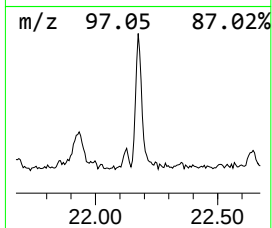
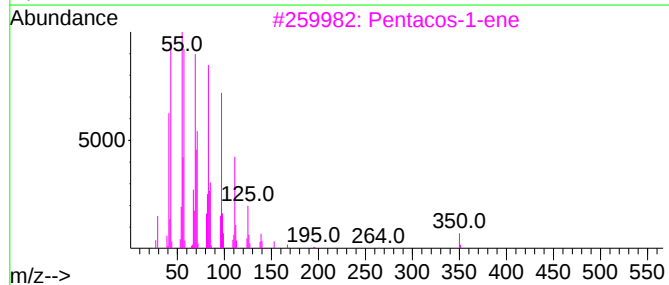
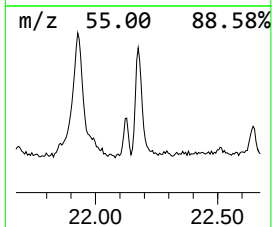
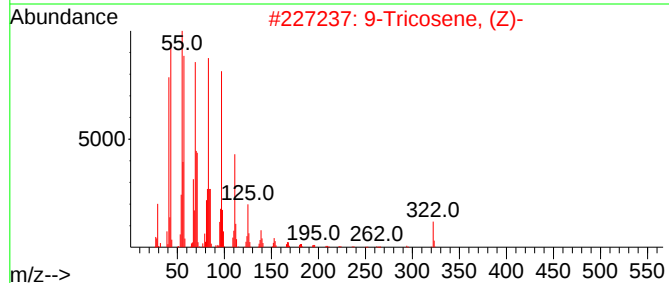
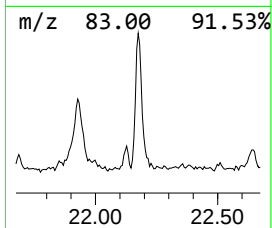
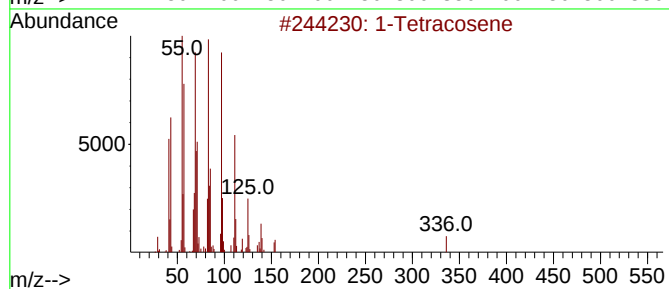
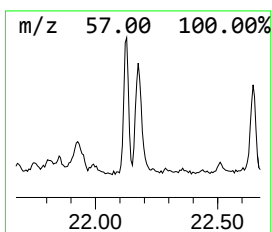
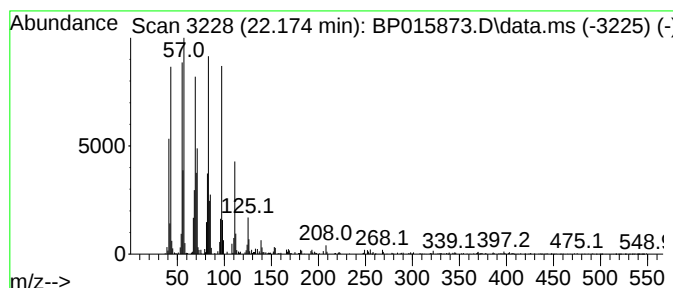
Instrument :
BNA_P
ClientSampleId :
DCFY4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.174	6.79 ng/ul	1196030	Chrysene-d12	21.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene	336	C24H48	010192-32-2	97
2	9-Tricosene, (Z)-	322	C23H46	027519-02-4	94
3	Pentacos-1-ene	350	C25H50	016980-85-1	94
4	Cyclotetracosane	336	C24H48	000297-03-0	94
5	2-Methyl-2-docosene	322	C23H46	1000131-16-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015873.D
Acq On : 01 Jul 2023 08:05
Operator : MA/JU
Sample : 03239-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

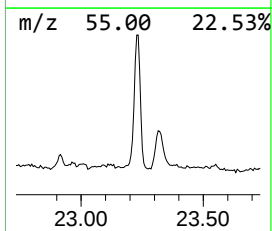
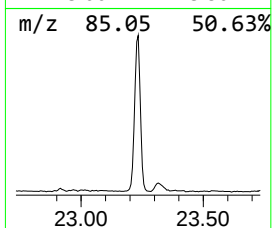
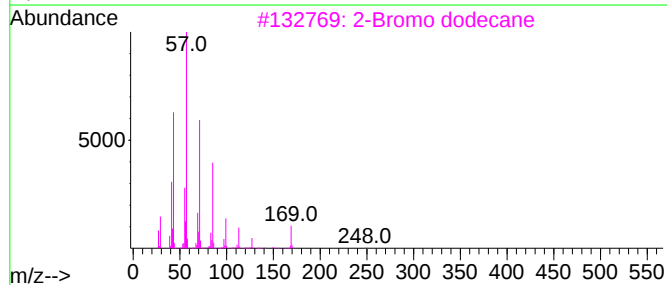
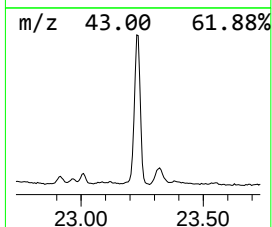
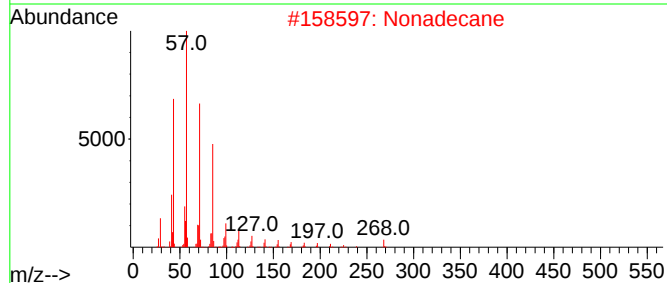
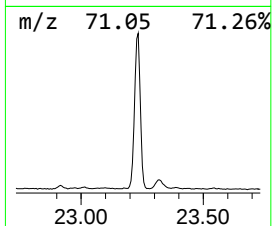
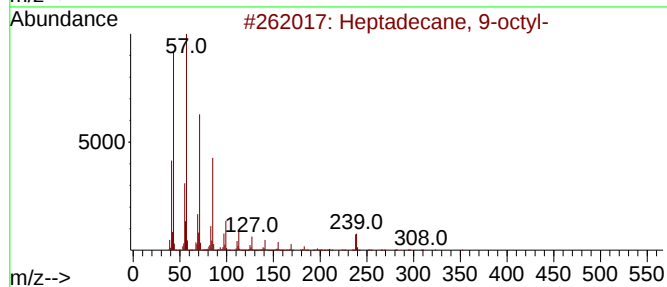
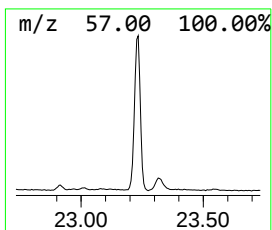
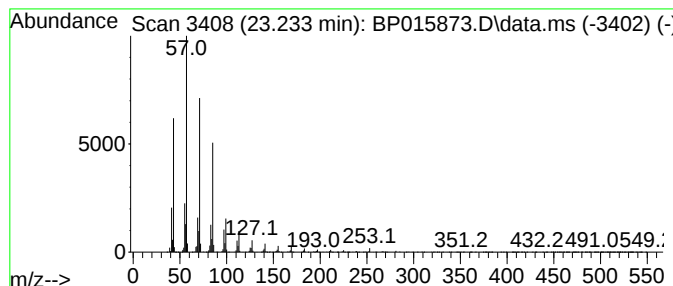
Instrument :
BNA_P
ClientSampleId :
DCFY4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.233	20.59 ng/ul	3331940	Perylene-d12	24.092	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	96
2	Nonadecane	268	C19H40	000629-92-5	93
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	93
4	Heneicosane	296	C21H44	000629-94-7	91
5	Hexatriacontane	507	C36H74	000630-06-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015873.D
Acq On : 01 Jul 2023 08:05
Operator : MA/JU
Sample : 03239-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFY4

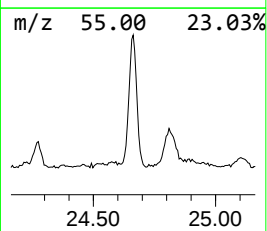
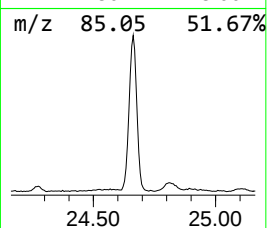
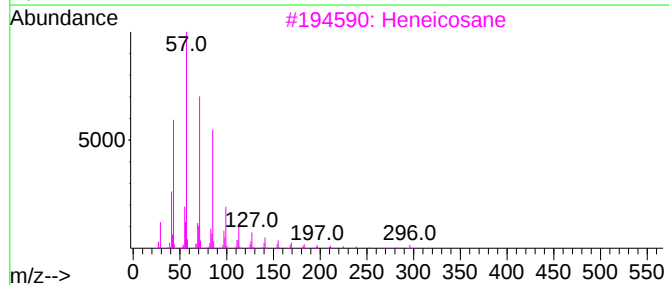
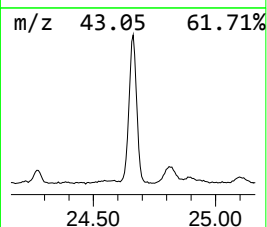
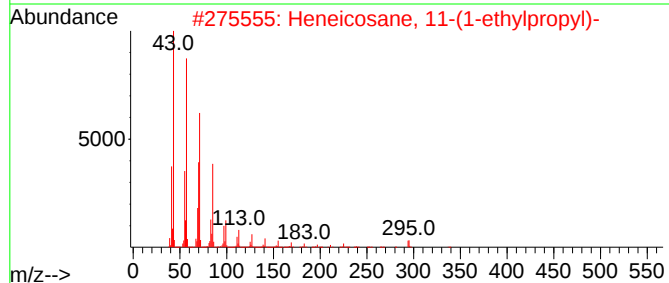
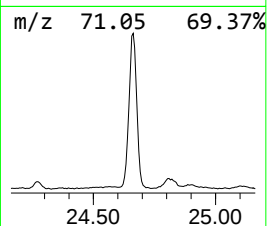
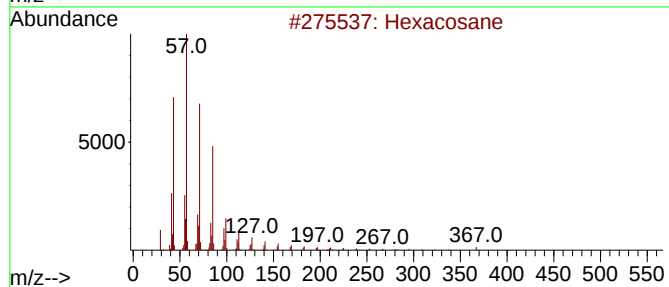
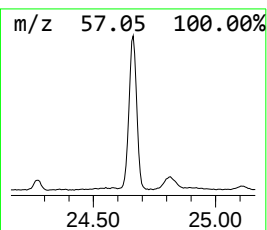
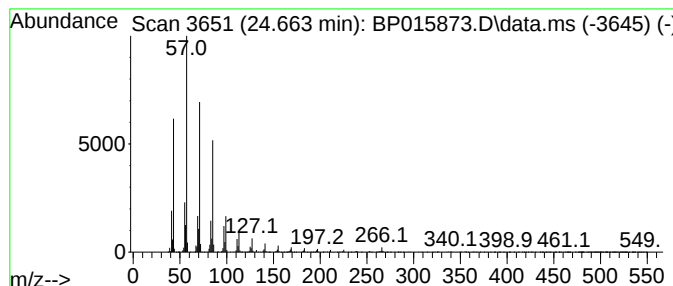
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.663	22.74 ng/ul	3678790	Perylene-d12	24.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	91
2			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
5			Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015873.D
Acq On : 01 Jul 2023 08:05
Operator : MA/JU
Sample : 03239-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFY4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2H-Pyran, tetra...	3.922	2.3	ng/ul	208958	1	8.052	1828460	20.0
Aceto phenone	8.905	6.6	ng/ul	604371	1	8.052	1828460	20.0
n-Hexadecanoic ...	18.316	8.8	ng/ul	1446360	4	17.463	3287690	20.0
Octadecanoic acid	19.539	3.3	ng/ul	578130	5	21.551	3525100	20.0
(DEL) Alkane: C...	21.222	9.1	ng/ul	1596570	5	21.551	3525100	20.0
Sesquirosefuran	21.927	17.8	ng/ul	3144340	5	21.551	3525100	20.0
(DEL) Alkane: S...	22.127	2.2	ng/ul	381038	5	21.551	3525100	20.0
(DEL) Alkane: S...	22.174	6.8	ng/ul	1196030	5	21.551	3525100	20.0
(DEL) Alkane: S...	23.233	20.6	ng/ul	3331940	6	24.092	3235980	20.0
(DEL) Alkane: S...	24.663	22.7	ng/ul	3678790	6	24.092	3235980	20.0