

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
 Data File : BP009943.D
 Acq On : 19 Apr 2022 07:17
 Operator : CG/JU
 Sample : N2421-10
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COHY7

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: BP009943.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.116	3	5	21	rVB	354195	496022	10.23%	1.127%
2	3.381	46	50	52	rBV	28269	38895	0.80%	0.088%
3	3.410	52	55	65	rVB4	33591	78563	1.62%	0.179%
4	3.893	135	137	144	rBV8	4700	9125	0.19%	0.021%
5	3.963	144	149	156	rBV8	13885	40671	0.84%	0.092%
6	4.128	175	177	184	rVB3	10342	13530	0.28%	0.031%
7	4.487	234	238	245	rVB3	38040	61182	1.26%	0.139%
8	4.599	252	257	266	rVB	27663	60004	1.24%	0.136%
9	5.257	365	369	372	rBV6	6026	7767	0.16%	0.018%
10	5.640	428	434	442	rBV3	17560	42832	0.88%	0.097%
11	5.999	491	495	499	rBV7	8568	11968	0.25%	0.027%
12	6.493	574	579	585	rVB2	18202	33812	0.70%	0.077%
13	7.110	677	684	695	rVB2	226638	429860	8.86%	0.977%
14	7.275	705	712	721	rBV	542550	857257	17.67%	1.948%
15	7.481	738	747	761	rBV	587265	971251	20.03%	2.207%
16	7.575	761	763	768	rVV5	4436	7795	0.16%	0.018%
17	7.640	771	774	778	rVB6	8725	11563	0.24%	0.026%
18	7.951	819	827	836	rBV	618783	1006920	20.76%	2.288%
19	8.310	882	888	889	rBV6	8414	11123	0.23%	0.025%
20	8.369	892	898	903	rVV4	10436	22079	0.46%	0.050%
21	8.504	917	921	927	rBV9	4939	7978	0.16%	0.018%
22	8.616	934	940	941	rBV6	12487	17584	0.36%	0.040%
23	8.651	941	946	958	rVB	436715	742512	15.31%	1.687%
24	8.875	982	984	992	rVB2	25699	44476	0.92%	0.101%
25	9.110	1016	1024	1034	rBV	702383	1161149	23.94%	2.639%
26	9.228	1037	1044	1047	rVV2	28066	46497	0.96%	0.106%
27	9.281	1047	1053	1061	rVB	264232	407205	8.40%	0.925%
28	9.563	1092	1101	1108	rVB	55800	102842	2.12%	0.234%
29	9.669	1113	1119	1123	rBV9	6054	8500	0.18%	0.019%
30	9.834	1140	1147	1162	rBV	328078	555321	11.45%	1.262%
31	9.951	1162	1167	1177	rBV	69705	109487	2.26%	0.249%
32	10.063	1184	1186	1190	rVB5	8004	7966	0.16%	0.018%
33	10.281	1220	1223	1226	rVV5	5615	8284	0.17%	0.019%
34	10.375	1229	1239	1254	rBV	860061	1512539	31.19%	3.437%
35	10.539	1261	1267	1277	rBV	29890	66733	1.38%	0.152%
36	10.704	1289	1295	1297	rBV4	10838	18094	0.37%	0.041%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
 Data File : BP009943.D
 Acq On : 19 Apr 2022 07:17
 Operator : CG/JU
 Sample : N2421-10
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0HY7

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

37	10.757	1297	1304	1314	rVB	886936	1476189	30.44%	3.355%
38	10.934	1325	1334	1350	rBV	116975	350108	7.22%	0.796%
39	11.175	1374	1375	1383	rVB8	9230	8642	0.18%	0.020%
40	11.434	1416	1419	1424	rVB7	7464	11931	0.25%	0.027%
41	11.551	1432	1439	1448	rBV	74049	150439	3.10%	0.342%
42	11.769	1470	1476	1480	rBV7	7421	12697	0.26%	0.029%
43	12.057	1518	1525	1532	rBV2	24099	46636	0.96%	0.106%
44	12.151	1537	1541	1543	rBV5	5598	7855	0.16%	0.018%
45	12.339	1567	1573	1581	rBV	18374	32904	0.68%	0.075%
46	12.516	1598	1603	1607	rBV6	10393	20001	0.41%	0.045%
47	12.551	1607	1609	1616	rVB8	5061	7622	0.16%	0.017%
48	12.781	1643	1648	1652	rBV7	7523	15260	0.31%	0.035%
49	12.851	1659	1660	1664	rVB4	7919	7831	0.16%	0.018%
50	12.969	1673	1680	1684	rBV6	15433	35181	0.73%	0.080%
51	13.198	1714	1719	1724	rVB7	11769	21531	0.44%	0.049%
52	13.322	1733	1740	1744	rBV10	5768	13266	0.27%	0.030%
53	13.392	1746	1752	1754	rVV7	5150	10356	0.21%	0.024%
54	13.422	1754	1757	1762	rVB5	11270	15360	0.32%	0.035%
55	13.557	1775	1780	1785	rBV9	8409	14064	0.29%	0.032%
56	13.986	1847	1853	1859	rBV	1765224	2519022	51.94%	5.724%
57	14.233	1885	1895	1896	rBV7	30306	41017	0.85%	0.093%
58	14.275	1896	1902	1916	rVB	1772680	2583271	53.26%	5.870%
59	14.492	1935	1939	1943	rVB5	10545	14937	0.31%	0.034%
60	14.580	1945	1954	1964	rBV	1321244	1979817	40.82%	4.499%
61	14.775	1981	1987	1996	rBV	64599	129159	2.66%	0.294%
62	15.004	2020	2026	2033	rBV	329745	524139	10.81%	1.191%
63	15.239	2059	2066	2069	rBV4	19874	40641	0.84%	0.092%
64	15.386	2086	2091	2097	rBV2	45152	76715	1.58%	0.174%
65	15.439	2097	2100	2105	rVB2	12631	16039	0.33%	0.036%
66	15.575	2116	2123	2136	rBV	2348663	3598921	74.20%	8.178%
67	15.686	2136	2142	2150	rVB	231418	334540	6.90%	0.760%
68	15.816	2160	2164	2169	rBV	27922	40935	0.84%	0.093%
69	15.939	2181	2185	2189	rVB7	10786	17515	0.36%	0.040%
70	15.992	2191	2194	2197	rBV5	8552	11193	0.23%	0.025%
71	16.133	2214	2218	2225	rBV10	10812	24162	0.50%	0.055%
72	16.222	2229	2233	2236	rBV5	13063	16445	0.34%	0.037%
73	16.263	2237	2240	2243	rVB4	11777	13336	0.27%	0.030%
74	16.304	2243	2247	2250	rBV5	11636	15330	0.32%	0.035%
75	16.474	2272	2276	2280	rBV4	17900	22567	0.47%	0.051%
76	16.633	2299	2303	2304	rBV4	7011	8862	0.18%	0.020%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
 Data File : BP009943.D
 Acq On : 19 Apr 2022 07:17
 Operator : CG/JU
 Sample : N2421-10
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0HY7

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

77	16.733	2313	2320	2327	rBV2	87777	146566	3.02%	0.333%
78	16.816	2331	2334	2337	rVV5	10455	16331	0.34%	0.037%
79	17.051	2372	2374	2377	rVB3	7589	7613	0.16%	0.017%
80	17.104	2378	2383	2397	rVB2	105453	212805	4.39%	0.484%
81	17.263	2405	2410	2415	rVB3	37111	52326	1.08%	0.119%
82	17.333	2416	2422	2433	rBV	1571318	2299769	47.42%	5.226%
83	17.433	2433	2439	2450	rBV2	2552204	3732471	76.96%	8.482%
84	17.621	2468	2471	2476	rBV	37397	50993	1.05%	0.116%
85	17.727	2485	2489	2499	rVB5	32945	59534	1.23%	0.135%
86	17.839	2506	2508	2511	rBV3	9044	9085	0.19%	0.021%
87	18.004	2534	2536	2541	rVB5	23747	28171	0.58%	0.064%
88	18.221	2568	2573	2581	rBV	773217	1205353	24.85%	2.739%
89	18.674	2648	2650	2654	rBV4	14015	19861	0.41%	0.045%
90	19.051	2709	2714	2723	rBV	632609	918669	18.94%	2.088%
91	19.239	2743	2746	2750	rBV2	34892	44943	0.93%	0.102%
92	19.310	2755	2758	2762	rVV	30581	31632	0.65%	0.072%
93	19.451	2778	2782	2792	rBV	202882	264346	5.45%	0.601%
94	19.663	2812	2818	2828	rBV	3612522	4850113	100.00%	11.022%
95	20.292	2921	2925	2931	rVB	180643	215400	4.44%	0.489%
96	21.127	3064	3067	3075	rBV3	96780	133560	2.75%	0.304%
97	21.327	3098	3101	3105	rVB	125306	135145	2.79%	0.307%
98	21.415	3111	3116	3128	rBV2	1498372	2051919	42.31%	4.663%
99	23.721	3501	3508	3515	rBV2	1298002	2652663	54.69%	6.028%
100	23.880	3529	3535	3548	rVB3	720567	1549730	31.95%	3.522%

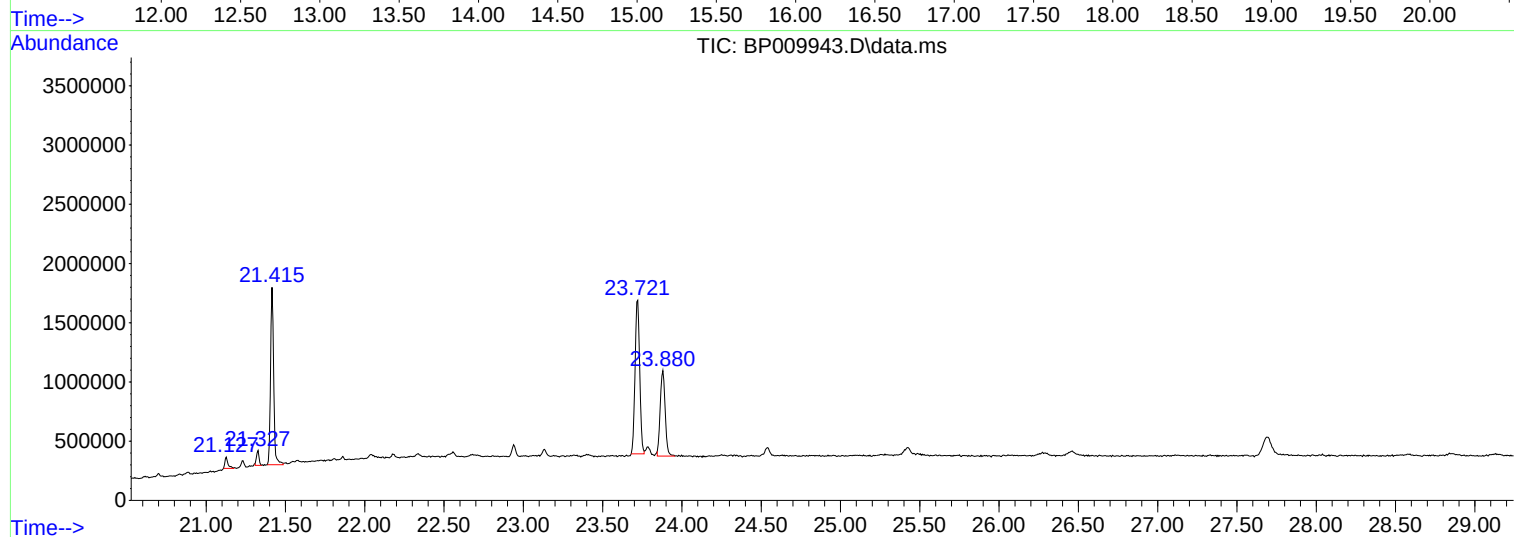
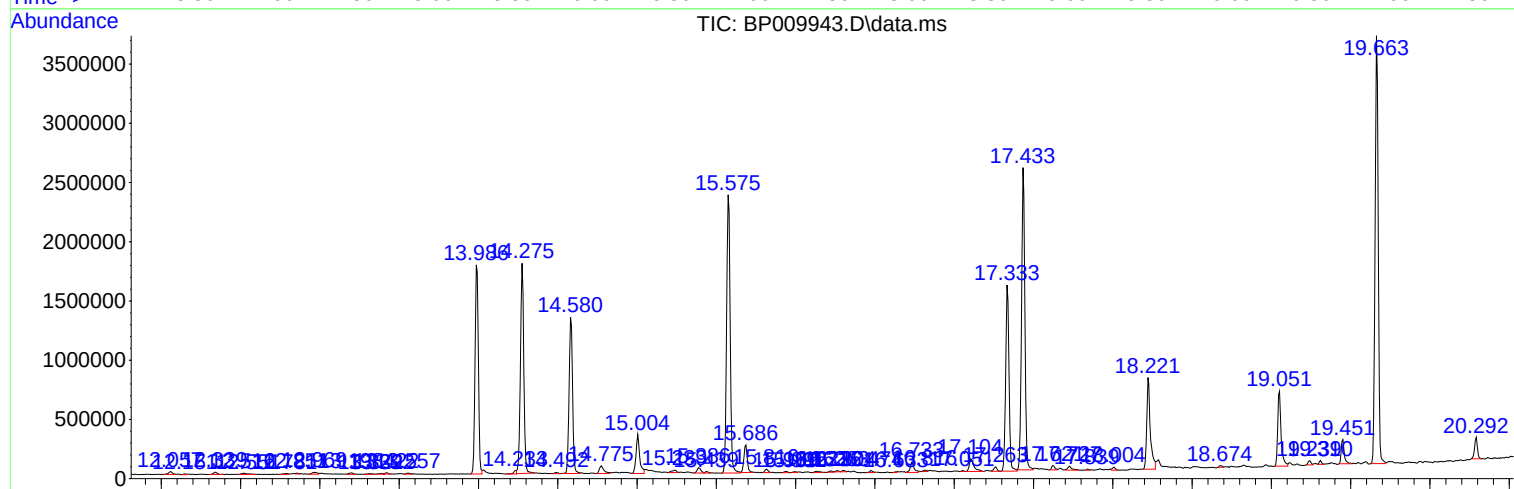
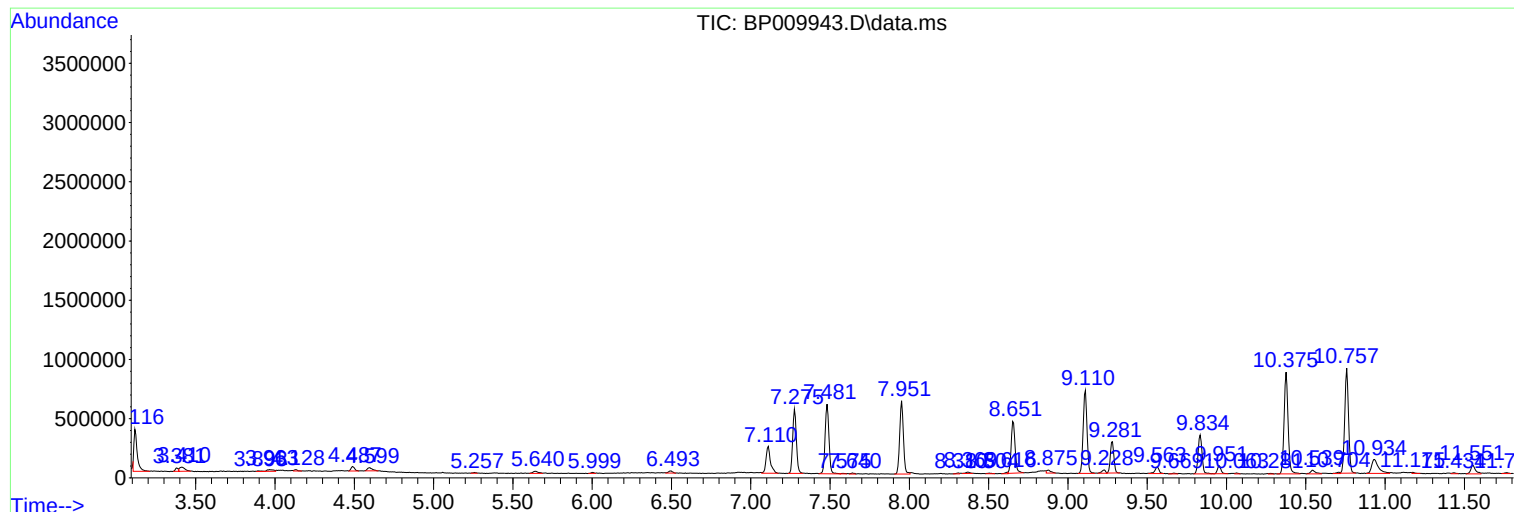
Sum of corrected areas: 44004820

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
Data File : BP009943.D
Acq On : 19 Apr 2022 07:17
Operator : CG/JU
Sample : N2421-10
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HY7

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
Data File : BP009943.D
Acq On : 19 Apr 2022 07:17
Operator : CG/JU
Sample : N2421-10
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HY7

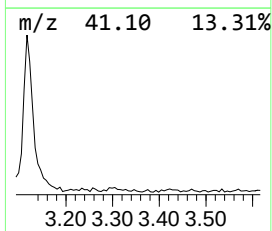
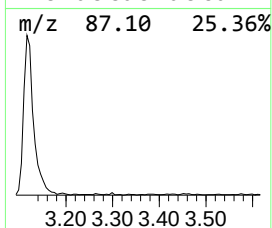
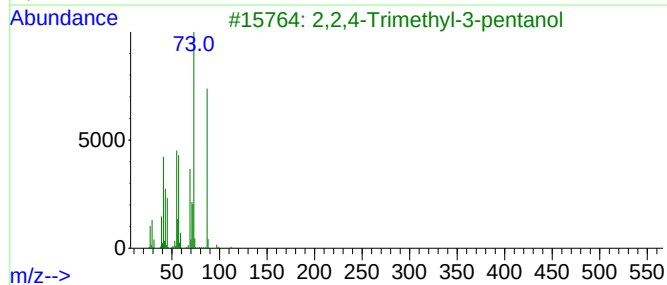
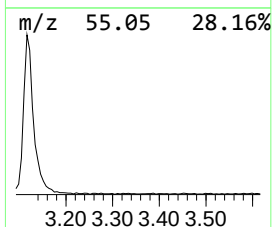
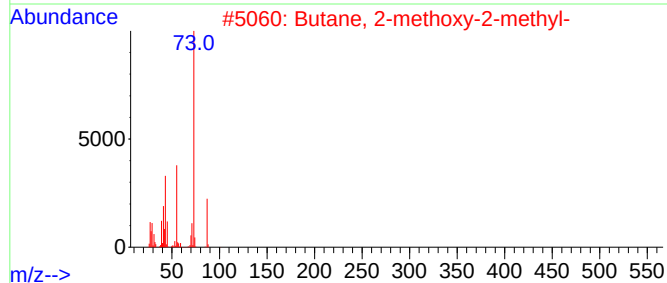
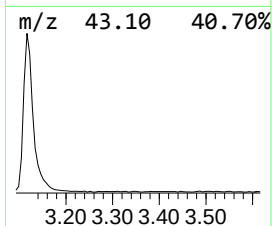
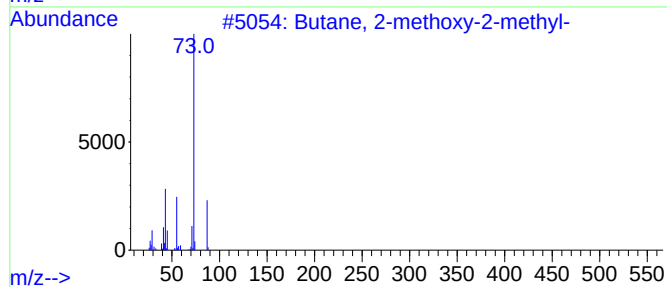
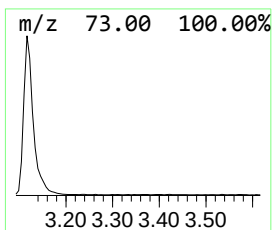
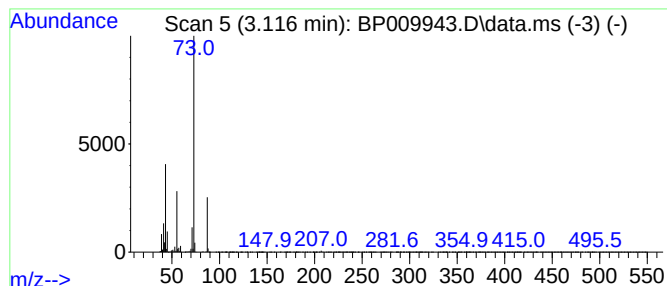
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.116	9.85 ng/ul	496022	1,4-Dichlorobenzene-d4	7.951

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	64
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	36
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
Data File : BP009943.D
Acq On : 19 Apr 2022 07:17
Operator : CG/JU
Sample : N2421-10
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HY7

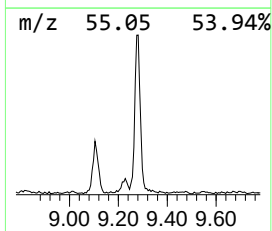
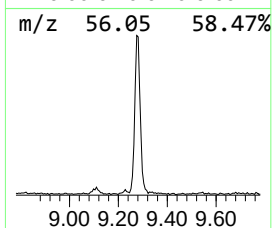
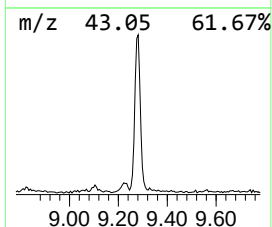
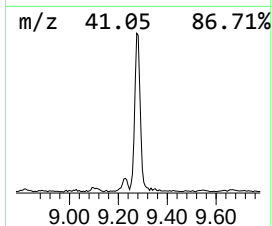
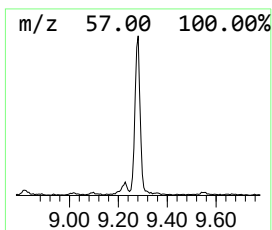
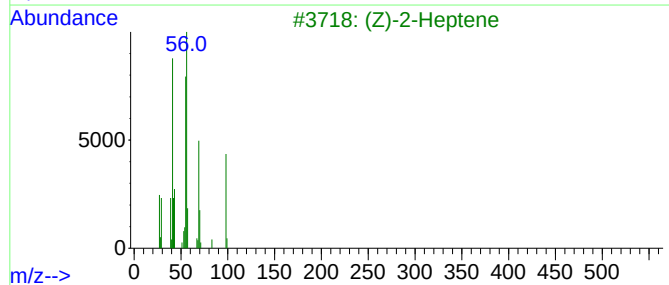
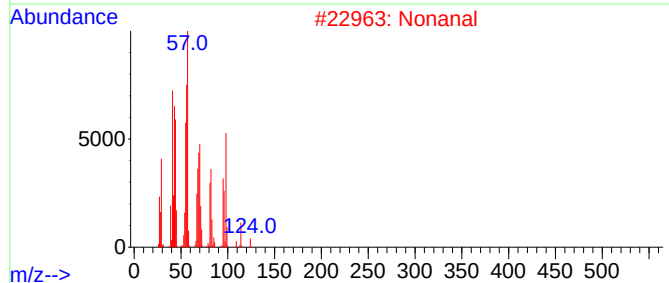
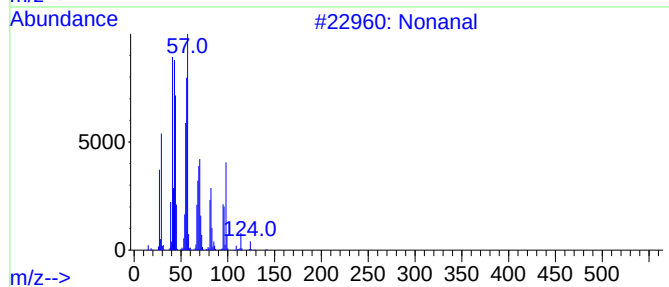
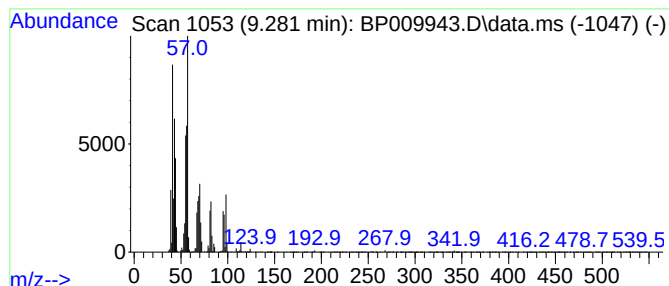
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Nonanal Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.281	8.09 ng/ul	407205	1,4-Dichlorobenzene-d4	7.951

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonanal	142	C9H18O	000124-19-6	83
2			Nonanal	142	C9H18O	000124-19-6	78
3			(Z)-2-Heptene	98	C7H14	006443-92-1	30
4			2-Hexene, 5-methyl-, (E)-	98	C7H14	007385-82-2	27
5			5-Methyl-2-hexene,c&t	98	C7H14	003404-62-4	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
Data File : BP009943.D
Acq On : 19 Apr 2022 07:17
Operator : CG/JU
Sample : N2421-10
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HY7

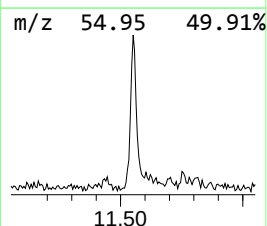
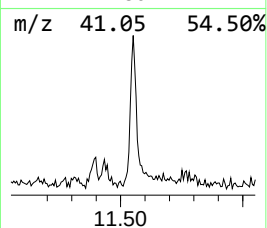
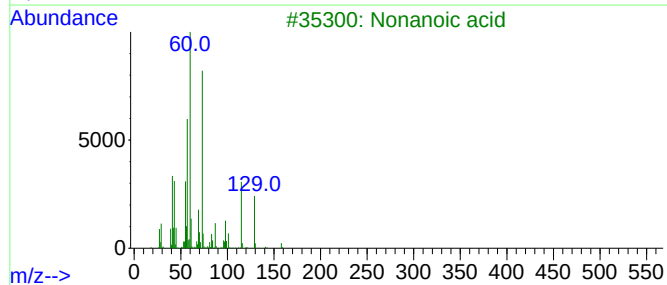
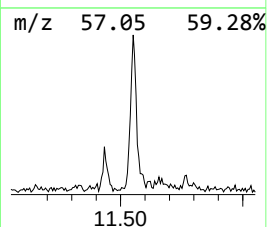
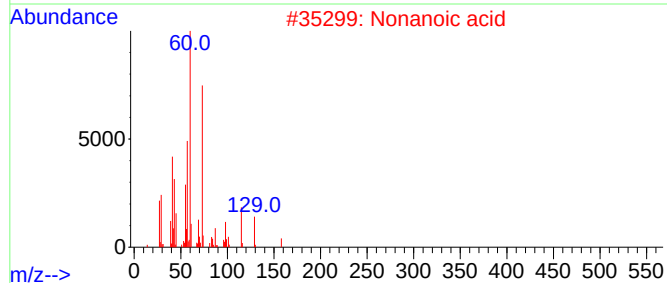
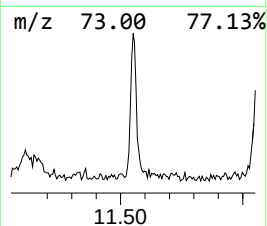
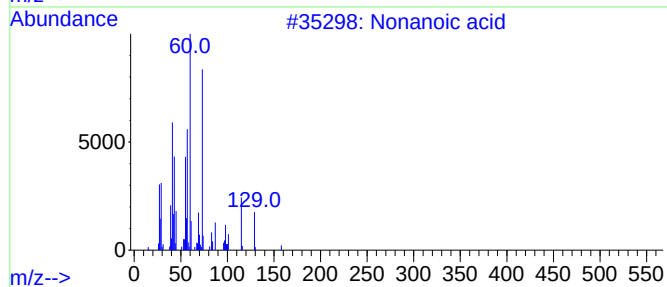
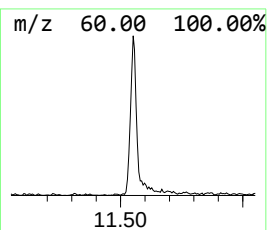
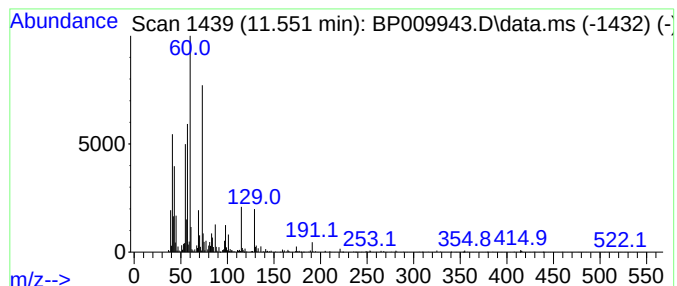
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Nonanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.551	2.04 ng/ul	150439	Naphthalene-d8	10.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonanoic acid	158	C9H18O2	000112-05-0	91
2			Nonanoic acid	158	C9H18O2	000112-05-0	90
3			Nonanoic acid	158	C9H18O2	000112-05-0	90
4			Nonanoic acid	158	C9H18O2	000112-05-0	90
5			n-Decanoic acid	172	C10H20O2	000334-48-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
Data File : BP009943.D
Acq On : 19 Apr 2022 07:17
Operator : CG/JU
Sample : N2421-10
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HY7

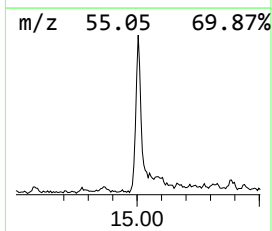
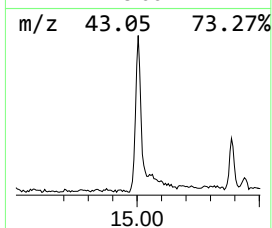
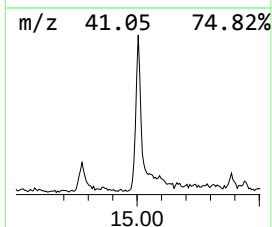
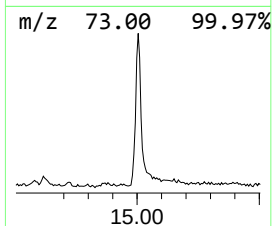
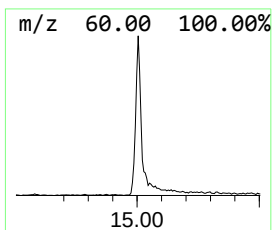
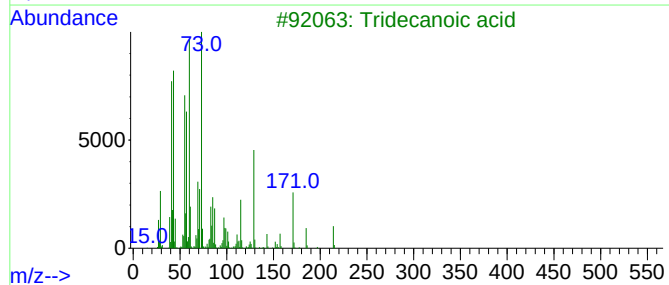
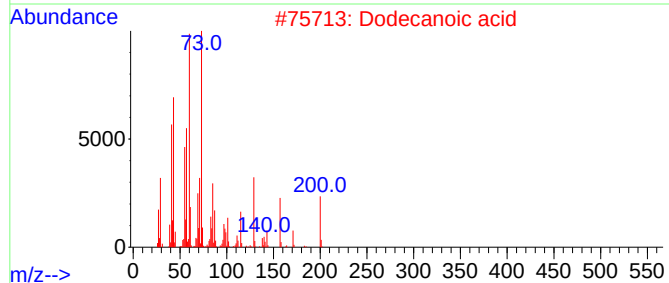
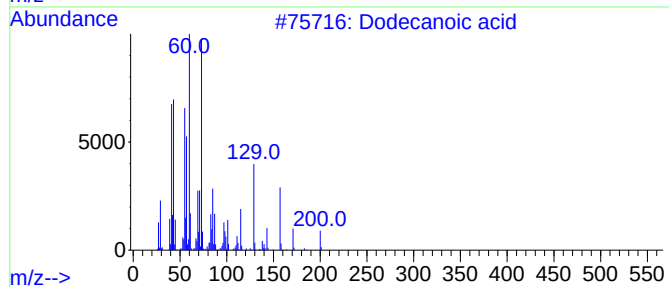
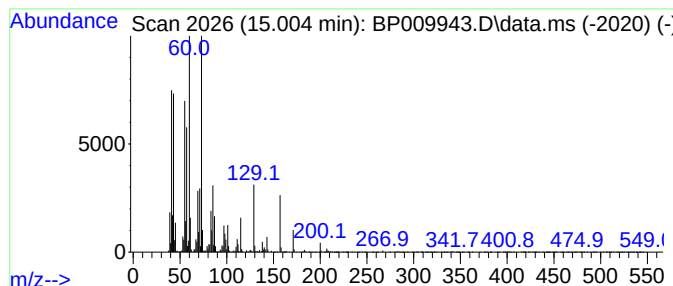
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Dodecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.004	5.29 ng/ul	524139	Acenaphthene-d10	14.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecanoic acid	200	C12H24O2	000143-07-7	98
2			Dodecanoic acid	200	C12H24O2	000143-07-7	91
3			Tridecanoic acid	214	C13H26O2	000638-53-9	91
4			Dodecanoic acid	200	C12H24O2	000143-07-7	91
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
Data File : BP009943.D
Acq On : 19 Apr 2022 07:17
Operator : CG/JU
Sample : N2421-10
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HY7

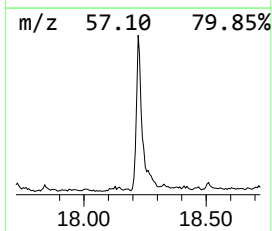
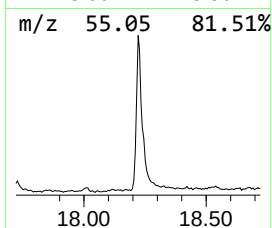
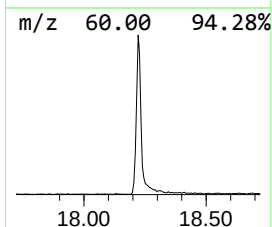
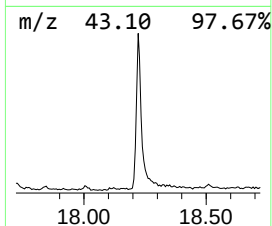
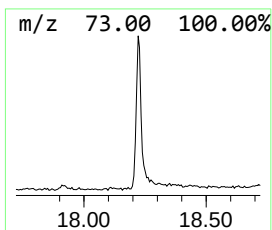
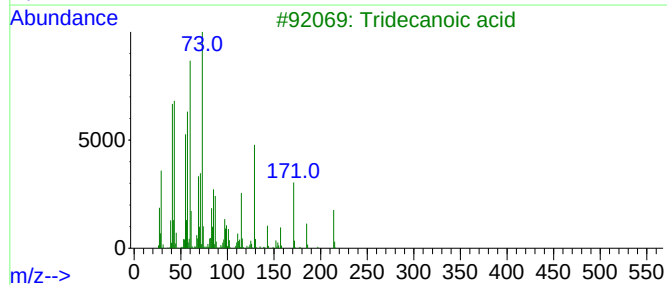
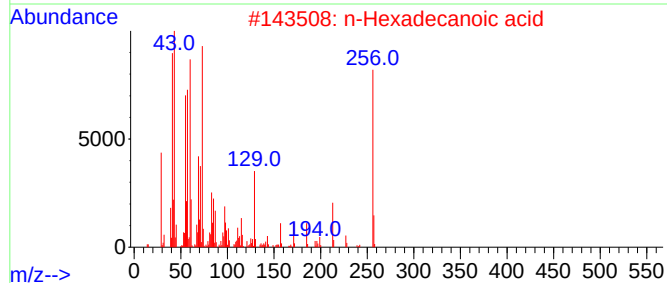
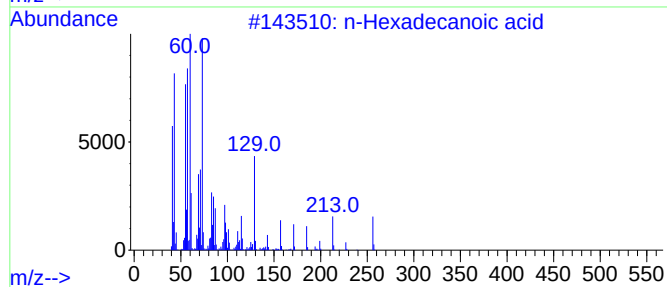
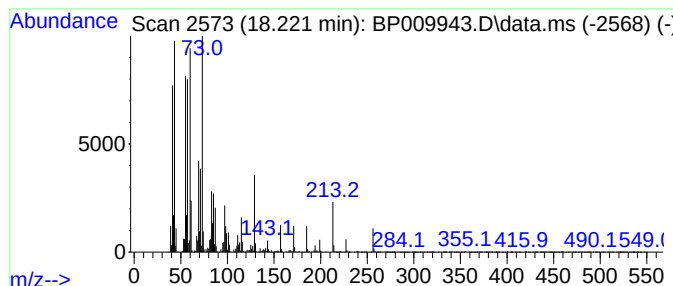
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.221	10.48 ng/ul	1205350	Phenanthrene-d10	17.333

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			Tridecanoic acid	214	C13H26O2	000638-53-9	96
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
Data File : BP009943.D
Acq On : 19 Apr 2022 07:17
Operator : CG/JU
Sample : N2421-10
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HY7

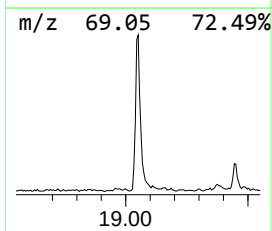
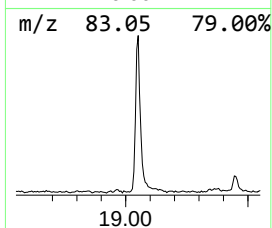
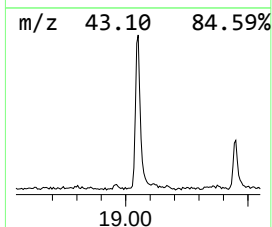
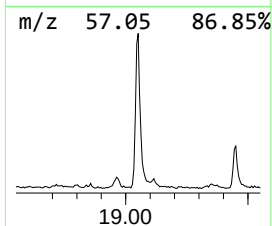
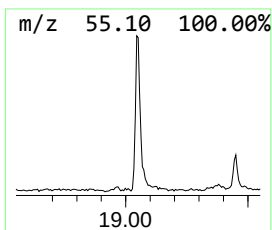
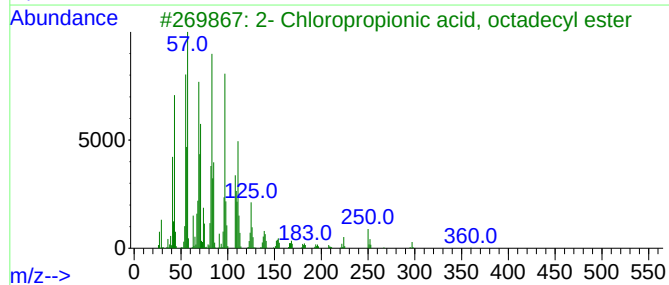
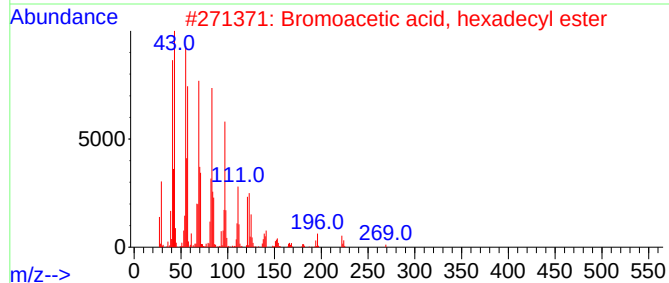
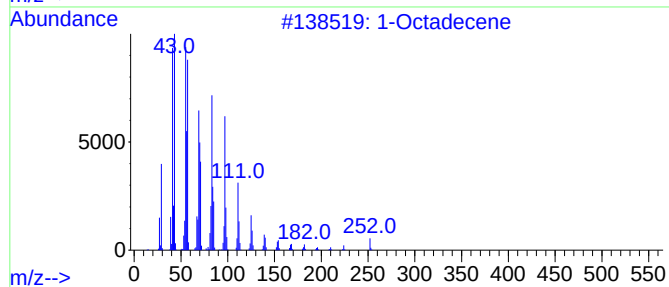
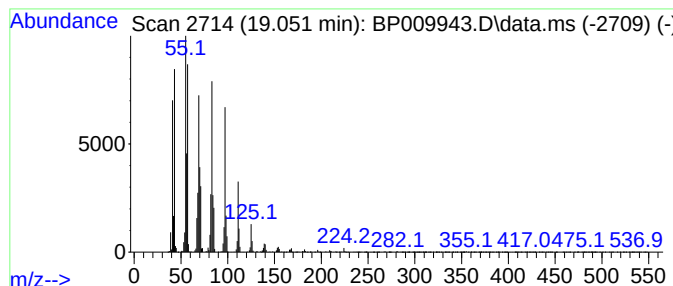
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.051	7.99 ng/ul	918669	Phenanthrene-d10	17.333

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Octadecene	252	C18H36	000112-88-9	94
2			Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	94
3			2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	93
4			1-Nonadecene	266	C19H38	018435-45-5	91
5			E-14-Hexadecenal	238	C16H30O	330207-53-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
Data File : BP009943.D
Acq On : 19 Apr 2022 07:17
Operator : CG/JU
Sample : N2421-10
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HY7

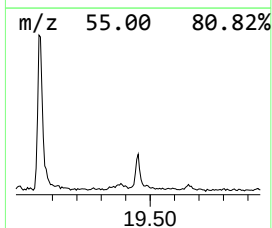
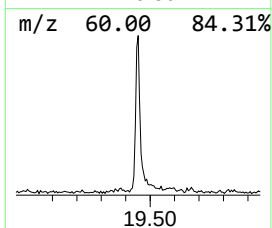
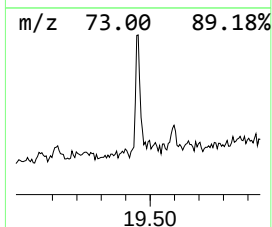
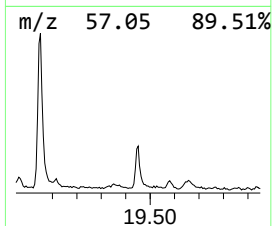
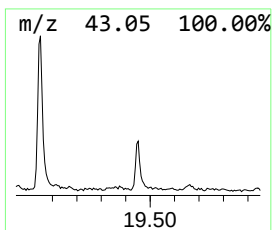
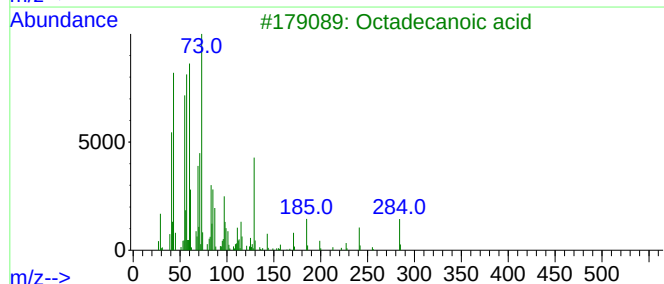
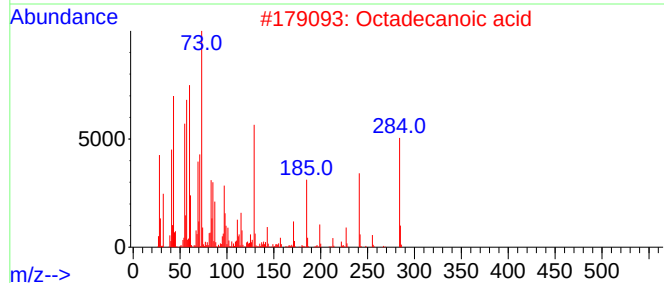
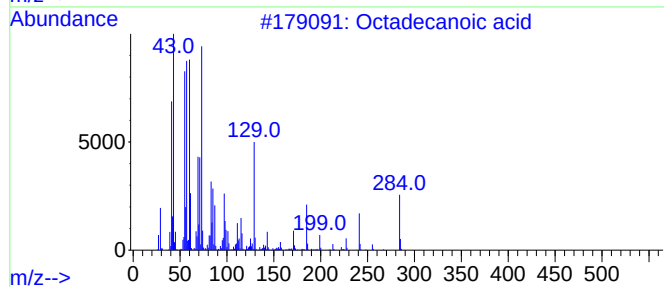
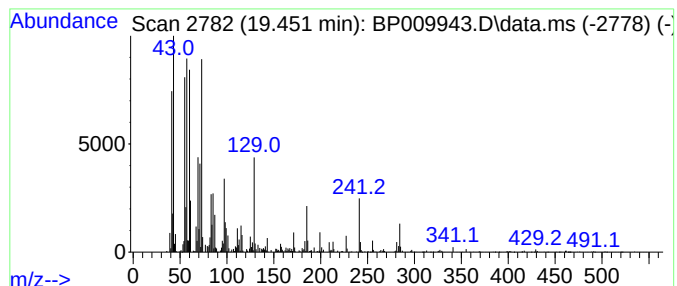
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.451	2.58 ng/ul	264346	Chrysene-d12	21.415

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	94
4			Octadecanoic acid	284	C18H36O2	000057-11-4	94
5			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
Data File : BP009943.D
Acq On : 19 Apr 2022 07:17
Operator : CG/JU
Sample : N2421-10
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HY7

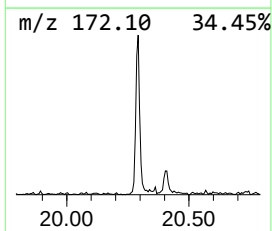
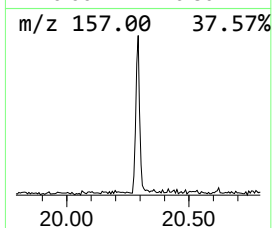
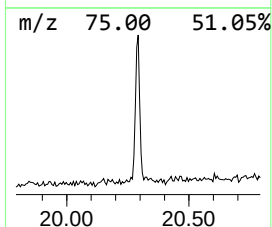
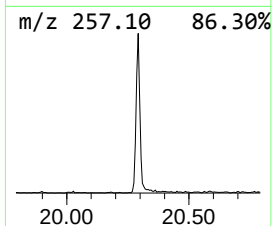
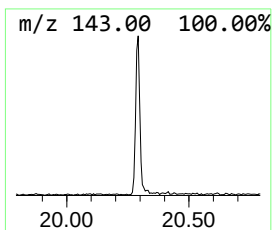
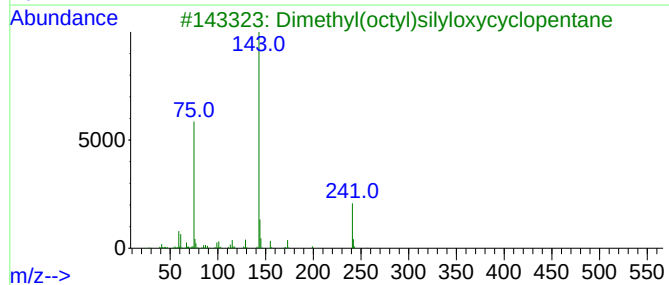
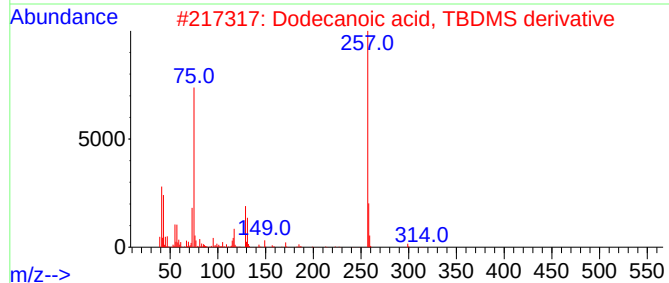
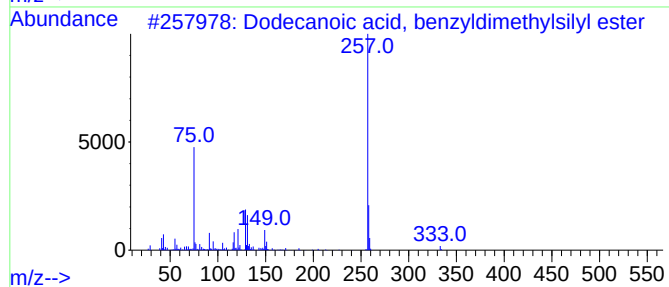
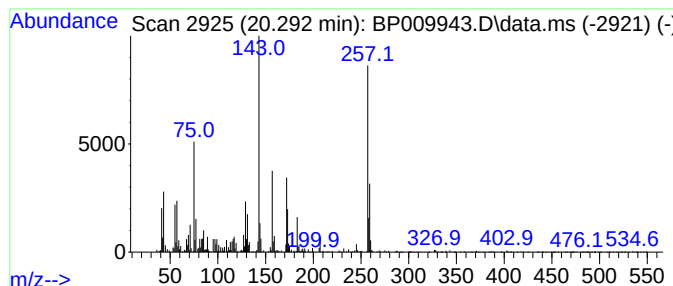
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.292	2.10 ng/ul	215400	Chrysene-d12	21.415

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecanoic acid, benzyldimethyls...	348	C21H36O2Si	1000375-51-2	43
2			Dodecanoic acid, TBDMS derivative	314	C18H38O2Si	104255-77-8	35
3			Dimethyl(octyl)silyloxycyclopentane	256	C15H32OSi	1010282-50-7	27
4			4-(4-Chloro-3-trifluoromethyl)py...	257	C12H7ClF3N	1000280-83-9	25
5			12-Tricosanol, TMS	412	C26H56OSi	1000506-69-9	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
Data File : BP009943.D
Acq On : 19 Apr 2022 07:17
Operator : CG/JU
Sample : N2421-10
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HY7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.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
Butane, 2-metho...	3.116	9.8	ng/ul	496022	1	7.951	1006920	20.0
Nonanal	9.281	8.1	ng/ul	407205	1	7.951	1006920	20.0
Nonanoic acid	11.551	2.0	ng/ul	150439	2	10.757	1476190	20.0
Dodecanoic acid	15.004	5.3	ng/ul	524139	3	14.581	1979820	20.0
n-Hexadecanoic ...	18.221	10.5	ng/ul	1205350	4	17.333	2299770	20.0
(DEL) Alkane: S...	19.051	8.0	ng/ul	918669	4	17.333	2299770	20.0
Octadecanoic acid	19.451	2.6	ng/ul	264346	5	21.415	2051920	20.0
unknown-01	20.292	2.1	ng/ul	215400	5	21.415	2051920	20.0