

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
 Data File : BP015996.D
 Acq On : 05 Jul 2023 02:59
 Operator : MA/JU
 Sample : 03244-14
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGF3

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: BP015996.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.370	28	31	40	rVB	62463	85943	0.39%	0.084%
2	3.823	105	108	121	rBV	512076	848285	3.85%	0.829%
3	3.917	121	124	134	rVB	60866	94207	0.43%	0.092%
4	4.034	139	144	150	rVB	78269	126084	0.57%	0.123%
5	4.134	158	161	170	rVB2	27125	47875	0.22%	0.047%
6	5.081	319	322	333	rVB	15614	30977	0.14%	0.030%
7	5.628	410	415	422	rBV6	12854	32253	0.15%	0.032%
8	6.005	473	479	487	rBV7	15452	33503	0.15%	0.033%
9	6.675	586	593	599	rVB5	27194	50218	0.23%	0.049%
10	7.117	661	668	675	rBV9	15587	51054	0.23%	0.050%
11	7.234	680	688	705	rBV	595164	1543733	7.01%	1.508%
12	7.369	705	711	727	rVB	662975	1261862	5.73%	1.233%
13	7.575	739	746	759	rBV	862857	1631954	7.41%	1.594%
14	8.040	818	825	838	rBV	894543	1650109	7.49%	1.612%
15	8.593	914	919	924	rVB9	10997	22080	0.10%	0.022%
16	8.793	945	953	979	rBV2	624960	1825601	8.29%	1.783%
17	9.234	1021	1028	1043	rBV	738499	1562028	7.09%	1.526%
18	9.375	1049	1052	1058	rVB6	17156	28457	0.13%	0.028%
19	9.963	1144	1152	1173	rBV	485294	1330117	6.04%	1.299%
20	10.540	1242	1250	1278	rBV	558920	1888439	8.57%	1.845%
21	10.805	1290	1295	1298	rBV3	16549	26843	0.12%	0.026%
22	10.869	1298	1306	1323	rVV	1171631	2417802	10.97%	2.362%
23	11.052	1330	1337	1361	rBV	283323	1003555	4.56%	0.980%
24	11.846	1467	1472	1478	rBV	27093	48257	0.22%	0.047%
25	12.252	1537	1541	1547	rVV2	19767	35608	0.16%	0.035%
26	12.481	1575	1580	1586	rBV6	12898	28109	0.13%	0.027%
27	12.552	1588	1592	1602	rVB3	17403	38592	0.18%	0.038%
28	12.763	1623	1628	1633	rBV3	16132	33133	0.15%	0.032%
29	13.193	1697	1701	1707	rVB3	24304	36207	0.16%	0.035%
30	13.481	1747	1750	1757	rVB2	34246	52596	0.24%	0.051%
31	13.563	1757	1764	1771	rBV4	70050	157869	0.72%	0.154%
32	13.834	1807	1810	1814	rBV4	29309	32223	0.15%	0.031%
33	13.987	1832	1836	1845	rBV7	57990	112982	0.51%	0.110%
34	14.104	1848	1856	1866	rBV	1605006	2684653	12.19%	2.622%
35	14.387	1898	1904	1917	rBV2	1964635	3393886	15.40%	3.315%
36	14.551	1928	1932	1937	rBV	51244	79901	0.36%	0.078%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
 Data File : BP015996.D
 Acq On : 05 Jul 2023 02:59
 Operator : MA/JU
 Sample : 03244-14
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGF3

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	14.628	1941	1945	1949	rBV2	69903	112588	0.51%	0.110%
38	14.693	1950	1956	1964	rBV	1707189	2627955	11.93%	2.567%
39	14.898	1986	1991	2005	rBV	261779	608979	2.76%	0.595%
40	15.016	2006	2011	2020	rBV	143452	495049	2.25%	0.484%
41	15.504	2091	2094	2101	rVB3	59946	87650	0.40%	0.086%
42	15.693	2120	2126	2142	rBV	2056655	3587400	16.28%	3.504%
43	15.881	2154	2158	2177	rVB4	191997	662484	3.01%	0.647%
44	16.063	2184	2189	2198	rBV	302808	631273	2.87%	0.617%
45	16.151	2201	2204	2208	rVB2	134746	185020	0.84%	0.181%
46	16.387	2238	2244	2248	rBV2	111409	248910	1.13%	0.243%
47	17.163	2374	2376	2380	rVB3	73025	72338	0.33%	0.071%
48	17.328	2400	2404	2410	rVB2	125112	181776	0.83%	0.178%
49	17.392	2412	2415	2420	rVB3	87311	120674	0.55%	0.118%
50	17.457	2421	2426	2430	rBV	1543924	2272032	10.31%	2.219%
51	17.498	2430	2433	2438	rVB	301911	442063	2.01%	0.432%
52	17.557	2438	2443	2449	rBV2	1643572	2726927	12.38%	2.664%
53	17.904	2499	2502	2509	rVB3	90599	115034	0.52%	0.112%
54	18.251	2557	2561	2566	rBV	695141	922780	4.19%	0.901%
55	18.310	2566	2571	2577	rBV	2844347	3742365	16.99%	3.655%
56	19.392	2753	2755	2757	rBV	183319	195803	0.89%	0.191%
57	19.416	2757	2759	2761	rVV	424741	505413	2.29%	0.494%
58	19.463	2763	2767	2771	rVV	997292	1687906	7.66%	1.649%
59	19.528	2775	2778	2789	rVB	835470	1503120	6.82%	1.468%
60	19.786	2817	2822	2825	rBV	2369988	3529913	16.02%	3.448%
61	21.228	3064	3067	3072	rVB2	638437	833174	3.78%	0.814%
62	21.398	3093	3096	3099	rBV	1625436	2278551	10.34%	2.226%
63	21.545	3116	3121	3125	rBV2	1830360	3252158	14.76%	3.177%
64	21.639	3132	3137	3171	rVB5	3812837	22031815	100.00%	21.520%
65	21.928	3178	3186	3193	rBV4	3451746	10385209	47.14%	10.144%
66	23.210	3400	3404	3415	rVB	1605336	2901639	13.17%	2.834%
67	23.922	3520	3525	3531	rVB2	1288223	2405775	10.92%	2.350%
68	24.086	3549	3553	3561	rVB	1051532	2130291	9.67%	2.081%
69	24.639	3641	3647	3655	rVB	2151018	4565230	20.72%	4.459%

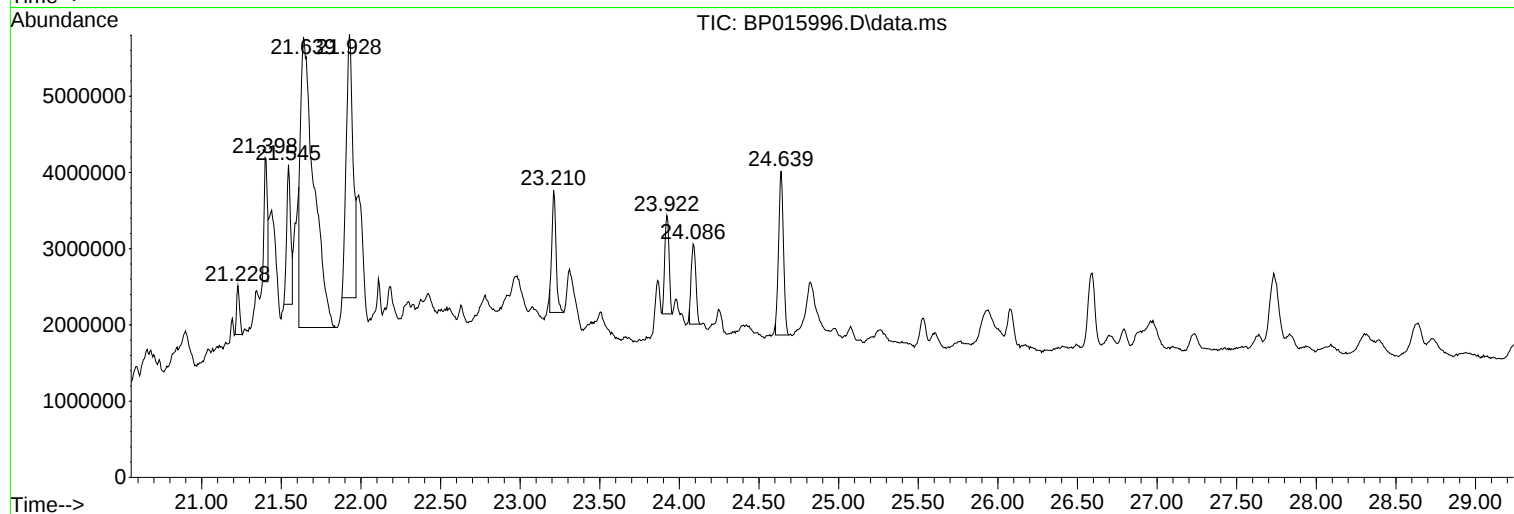
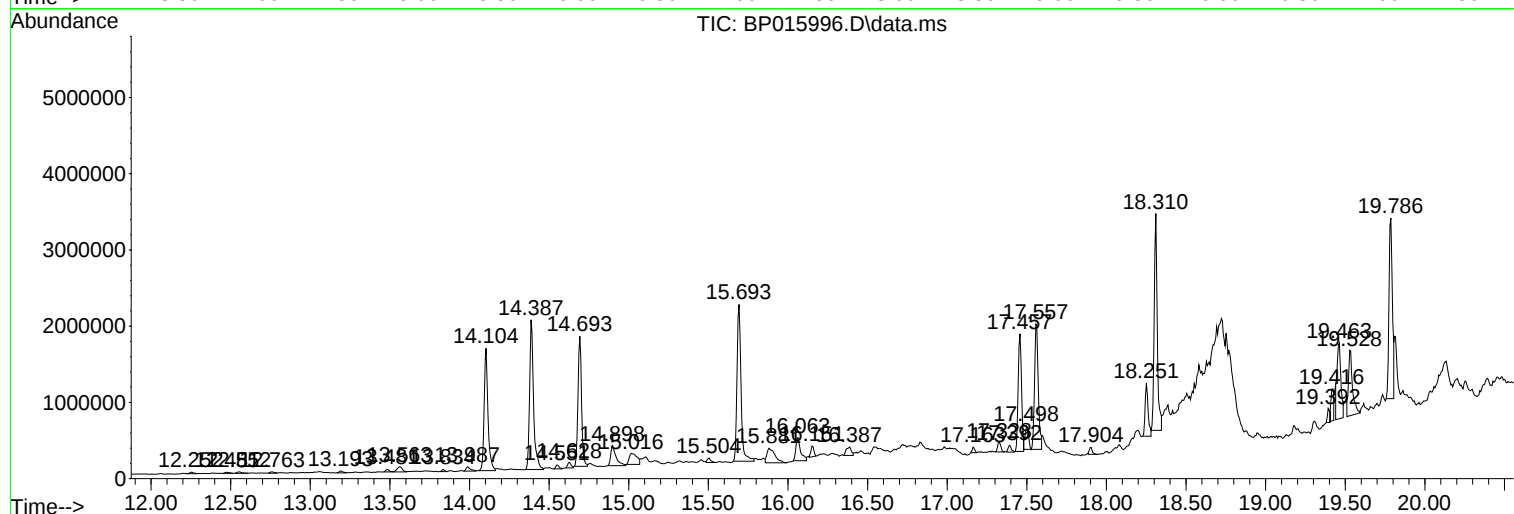
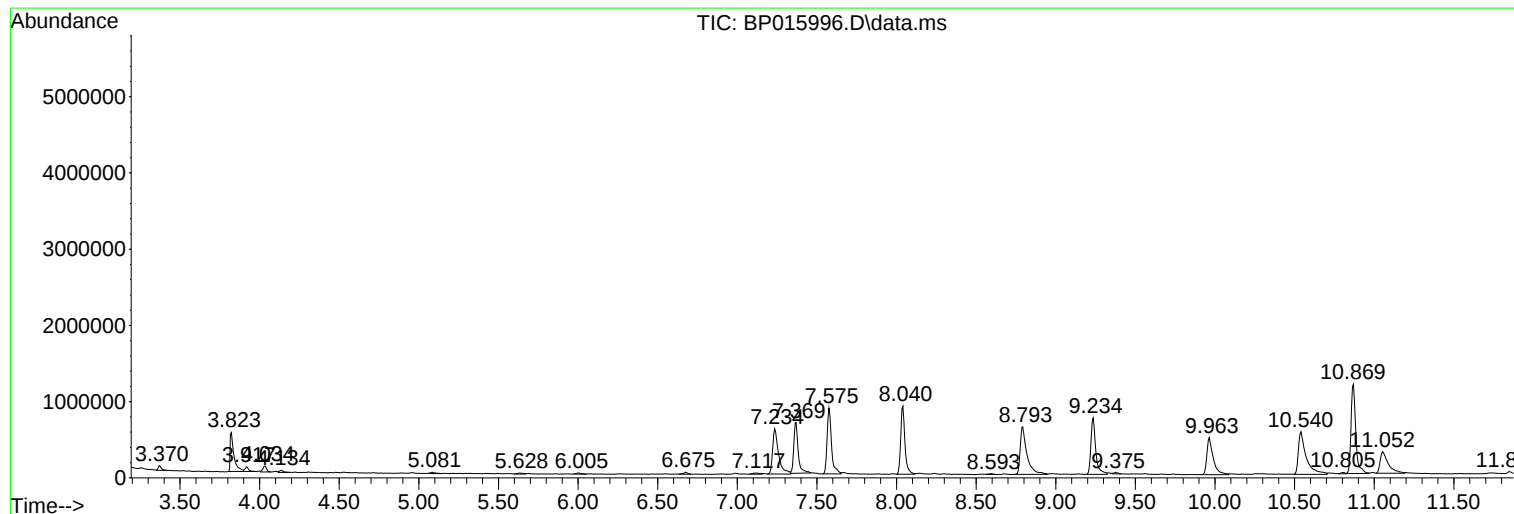
Sum of corrected areas: 102376289

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015996.D
Acq On : 05 Jul 2023 02:59
Operator : MA/JU
Sample : 03244-14
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGF3

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\BP070423\
Data File : BP015996.D
Acq On : 05 Jul 2023 02:59
Operator : MA/JU
Sample : 03244-14
Misc :
ALS Vial : 42 Sample Multiplier: 1

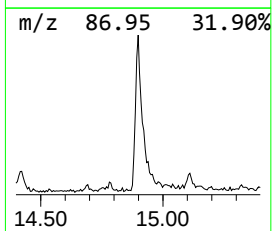
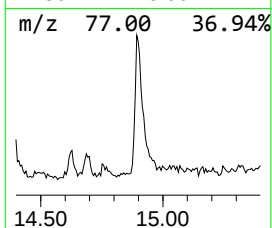
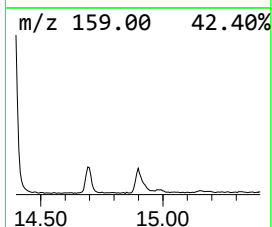
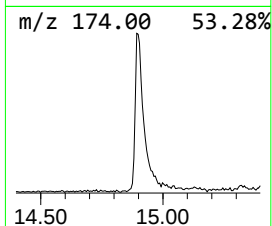
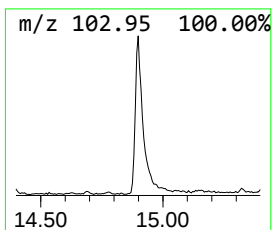
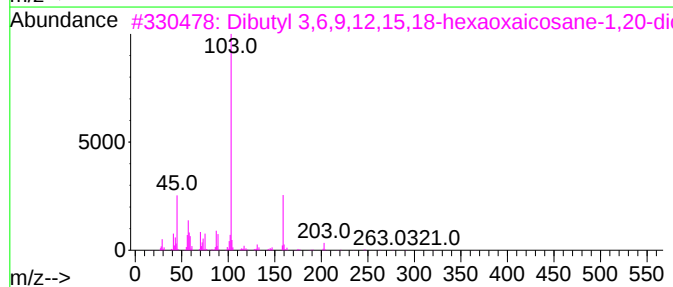
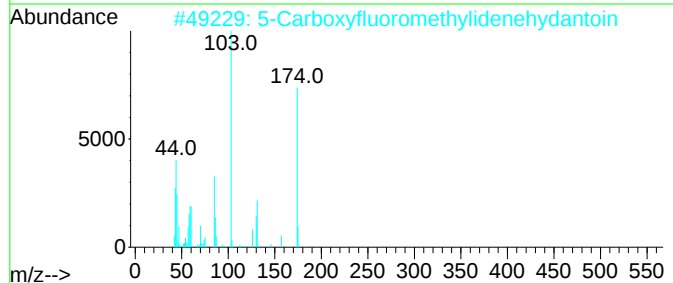
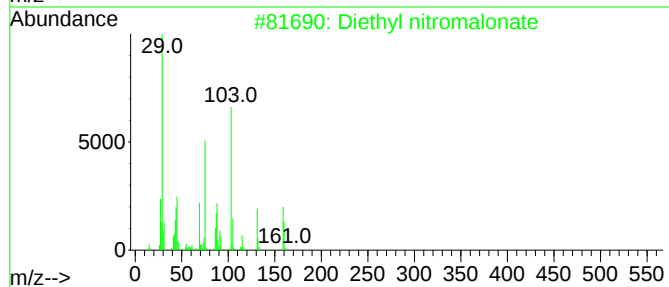
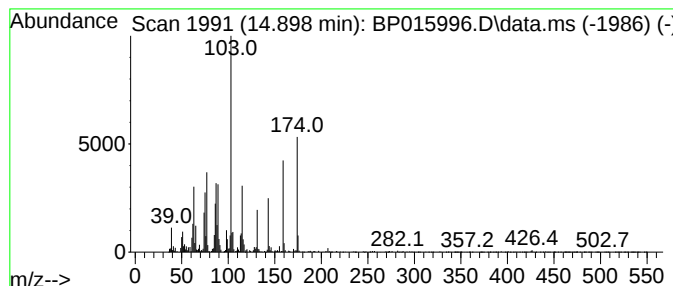
Instrument :
BNA_P
ClientSampleId :
DCGF3

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 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.899	4.63 ng/ul	608979	Acenaphthene-d10	14.693	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyl nitromalonate	205	C7H11NO6	000603-67-8	38
2	5-Carboxyfluoromethylidenehydantoin	174	C5H3FN2O4	057646-16-9	35
3	Dibutyl 3,6,9,12,15,18-hexaoxaic...	466	C22H42O10	1000366-80-0	16
4	5-Fluoro-2-phenylpyrimidine	174	C10H7FN2	054376-51-1	16
5	8-Quinolinol, 5-nitroso-	174	C9H6N2O2	003565-26-2	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015996.D
Acq On : 05 Jul 2023 02:59
Operator : MA/JU
Sample : 03244-14
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGF3

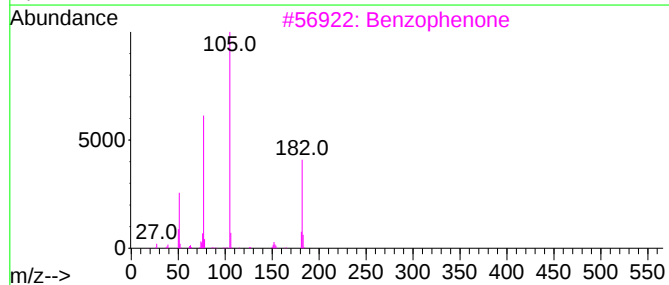
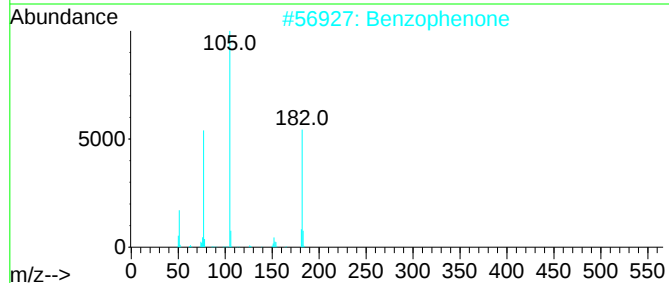
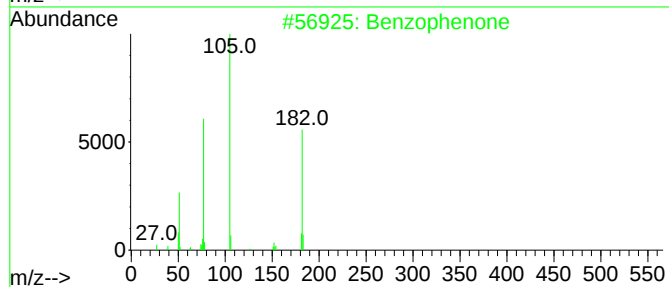
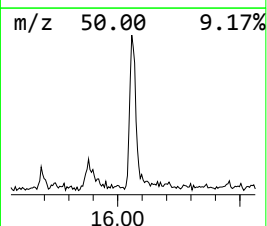
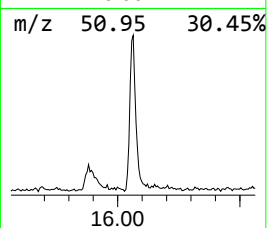
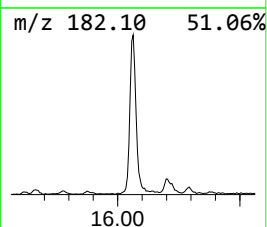
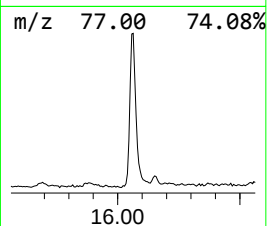
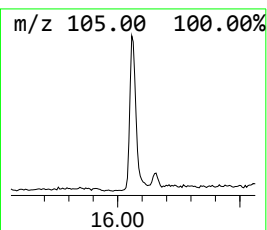
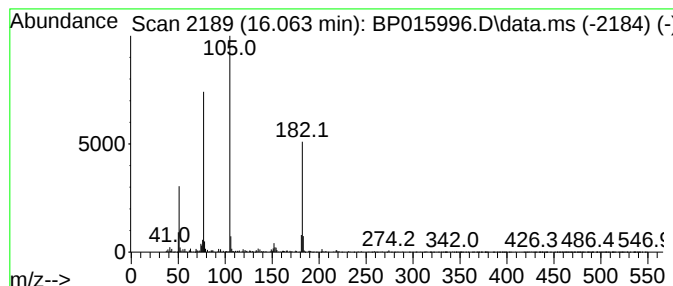
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.063	4.80 ng/ul	631273	Acenaphthene-d10	14.693

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015996.D
Acq On : 05 Jul 2023 02:59
Operator : MA/JU
Sample : 03244-14
Misc :
ALS Vial : 42 Sample Multiplier: 1

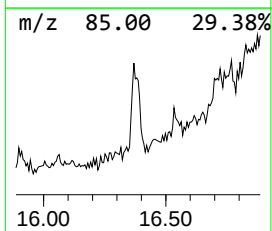
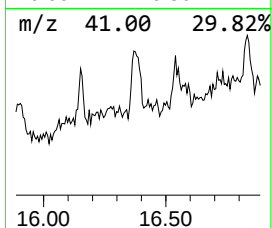
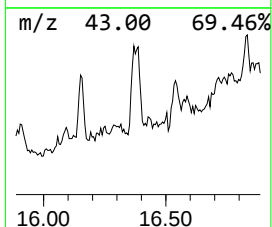
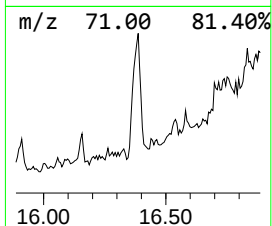
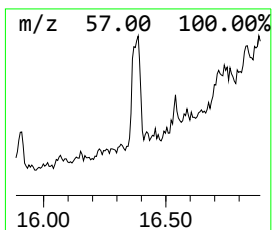
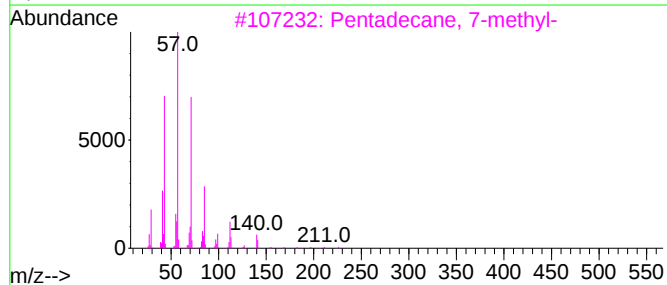
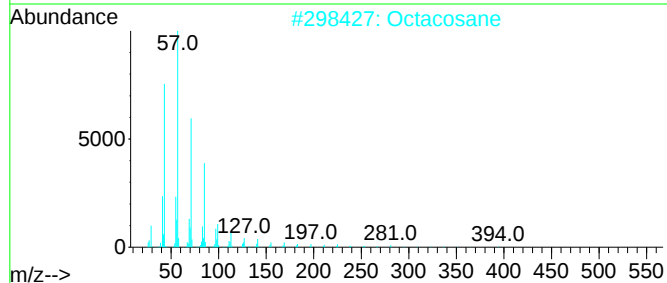
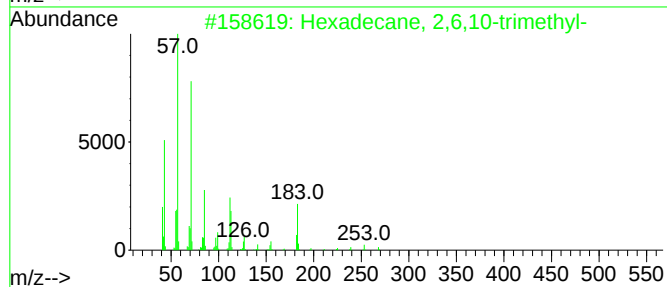
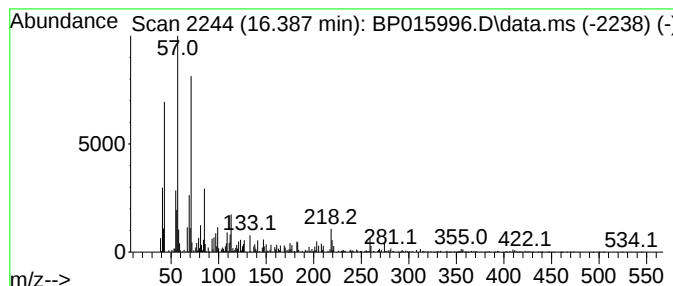
Instrument :
BNA_P
ClientSampleId :
DCGF3

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 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.387	2.19 ng/ul	248910	Phenanthrene-d10	17.457	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 2,6,10-trimethyl-	268	C19H40	055000-52-7	90
2	Octacosane	394	C28H58	000630-02-4	90
3	Pentadecane, 7-methyl-	226	C16H34	006165-40-8	89
4	Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	81
5	Octacosane	394	C28H58	000630-02-4	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015996.D
Acq On : 05 Jul 2023 02:59
Operator : MA/JU
Sample : 03244-14
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGF3

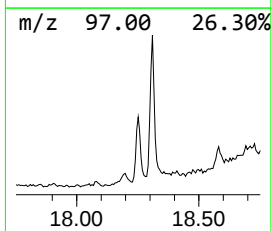
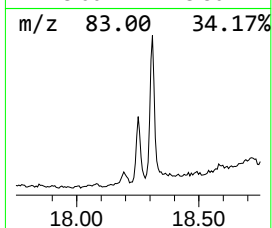
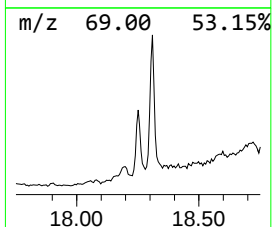
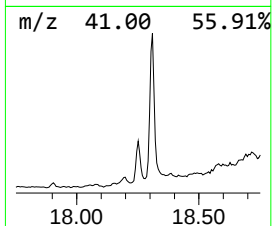
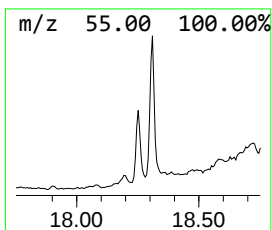
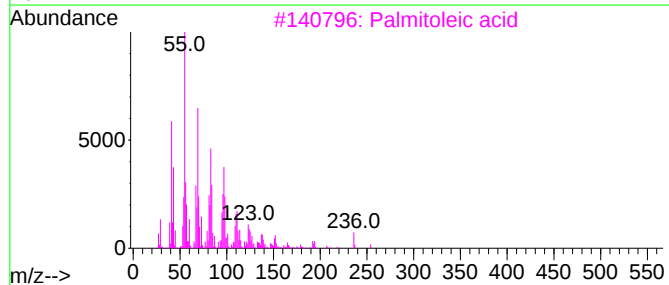
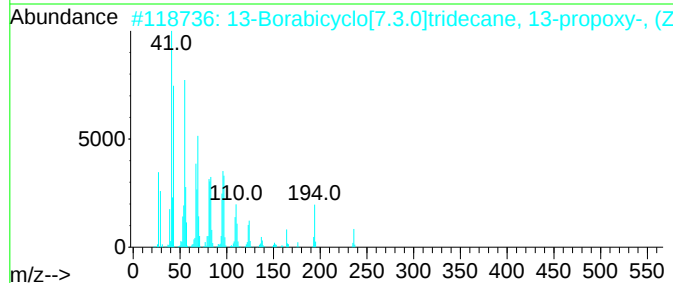
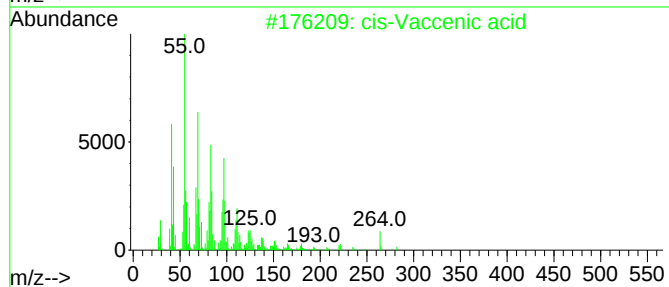
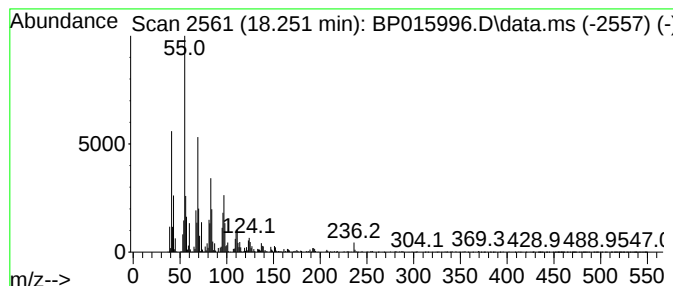
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 cis-Vaccenic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.251	8.12 ng/ul	922780	Phenanthrene-d10	17.457

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-Vaccenic acid	282	C18H34O2	000506-17-2	95
2			13-Borabicyclo[7.3.0]tridecane, ...	236	C15H29BO	1000156-41-7	91
3			Palmitoleic acid	254	C16H30O2	000373-49-9	91
4			cis-10-Heptadecenoic acid	268	C17H32O2	029743-97-3	91
5			trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015996.D
Acq On : 05 Jul 2023 02:59
Operator : MA/JU
Sample : 03244-14
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGF3

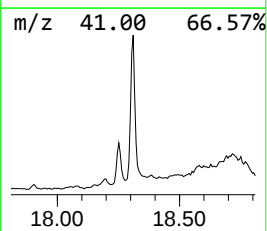
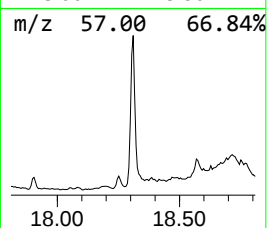
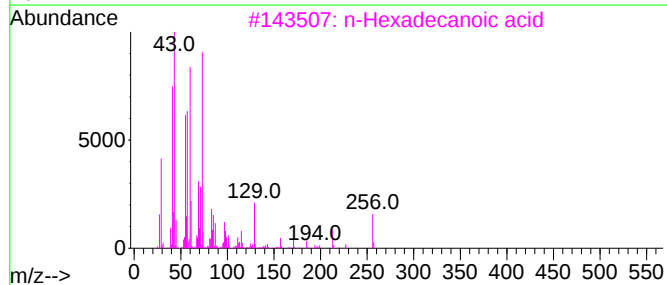
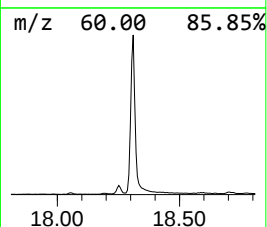
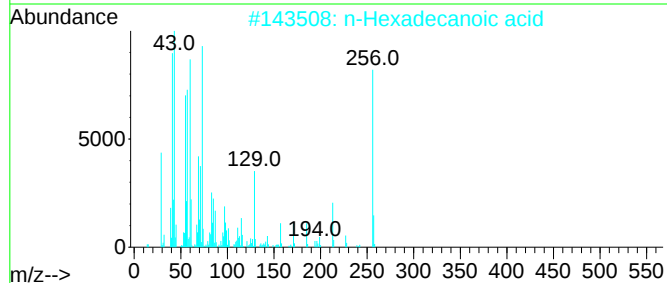
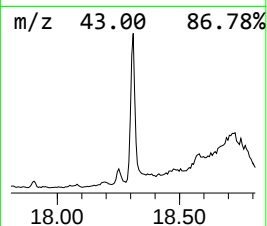
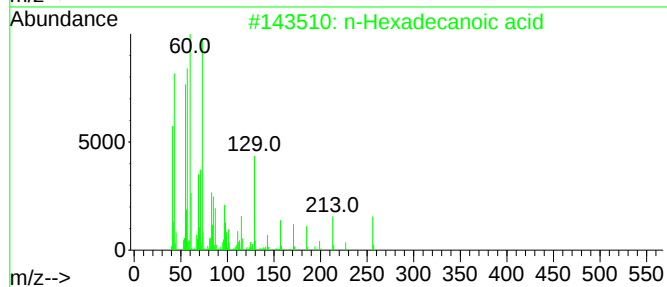
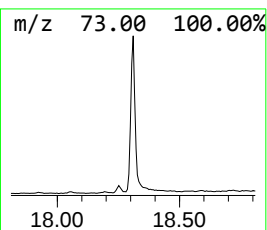
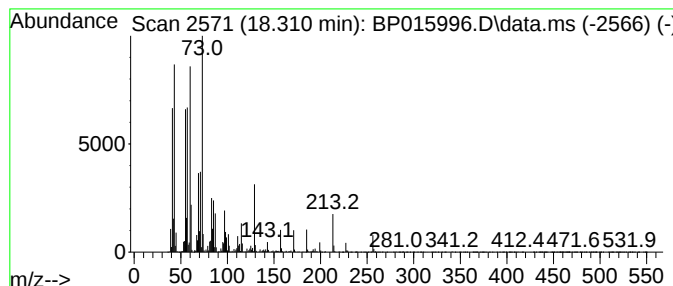
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.310	32.94 ng/ul	3742370	Phenanthrene-d10	17.457

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	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Tridecanoic acid	214	C13H26O2	000638-53-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015996.D
Acq On : 05 Jul 2023 02:59
Operator : MA/JU
Sample : 03244-14
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGF3

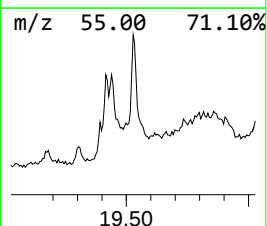
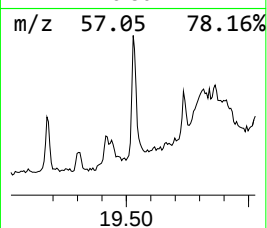
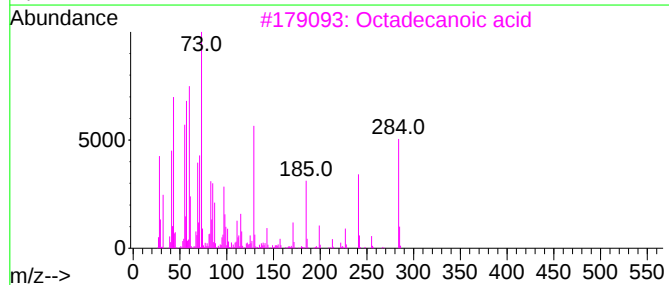
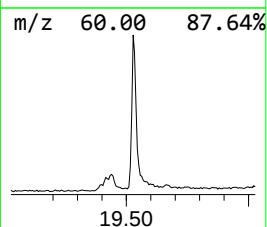
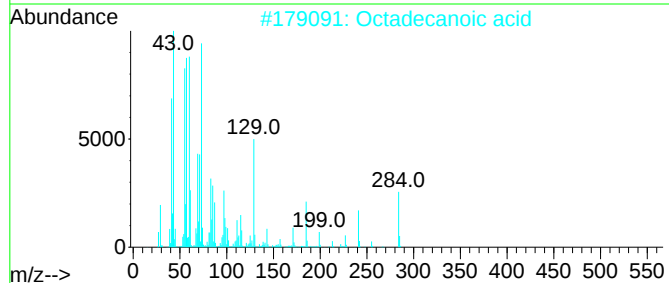
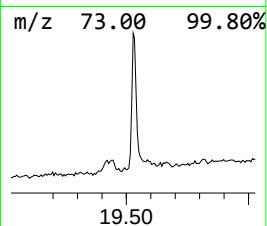
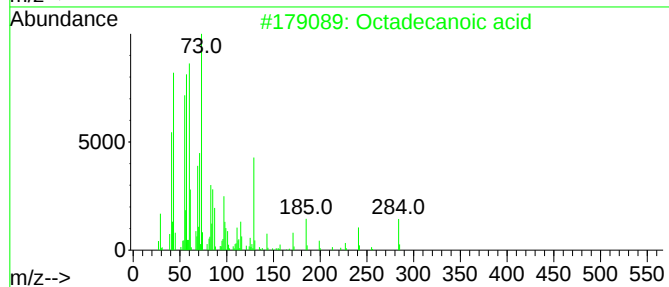
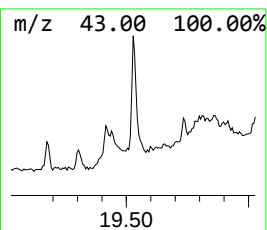
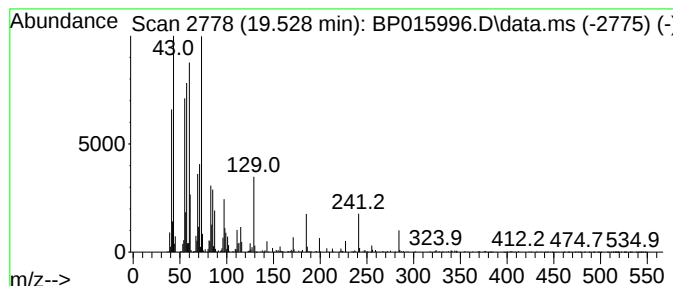
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.528	9.24 ng/ul	1503120	Chrysene-d12	21.545

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	96
5			Octadecanoic acid	284	C18H36O2	000057-11-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015996.D
Acq On : 05 Jul 2023 02:59
Operator : MA/JU
Sample : 03244-14
Misc :
ALS Vial : 42 Sample Multiplier: 1

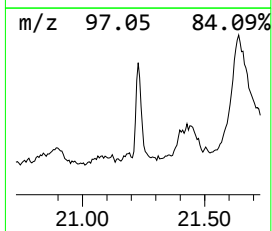
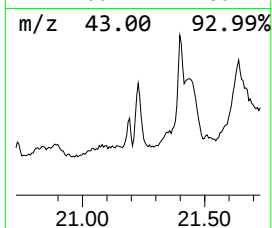
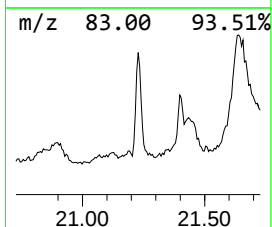
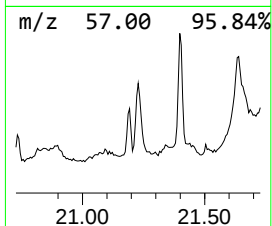
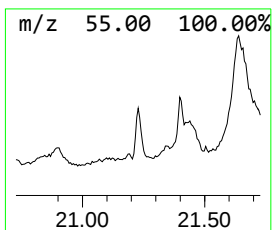
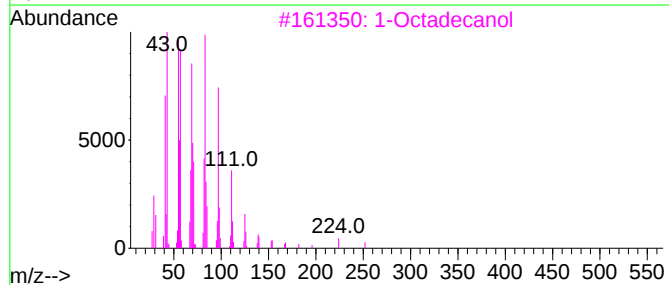
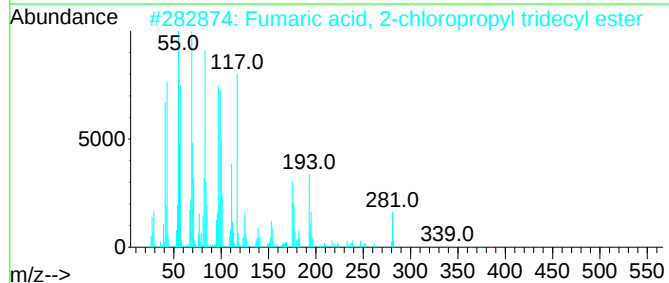
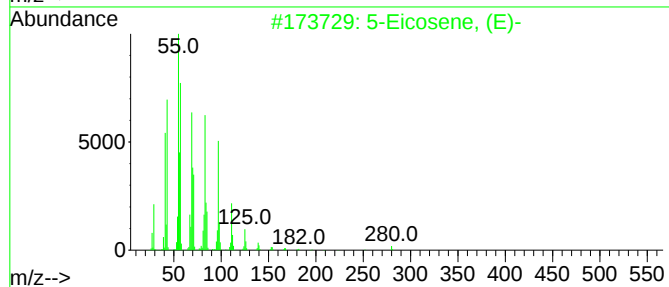
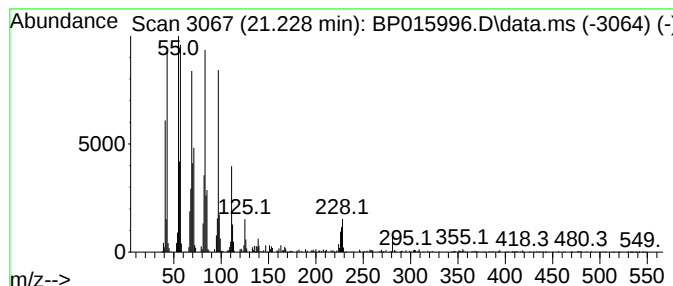
Instrument :
BNA_P
ClientSampleId :
DCGF3

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 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.228	5.12 ng/ul	833174	Chrysene-d12		21.545	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-		280	C20H40	074685-30-6	99
2	Fumaric acid, 2-chloropropyl tri...		374	C20H35ClO4	1010348-57-1	96
3	1-Octadecanol		270	C18H38O	000112-92-5	95
4	Cyclohexadecane		224	C16H32	000295-65-8	94
5	Trifluoroacetoxy hexadecane		338	C18H33F3O2	006222-03-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015996.D
Acq On : 05 Jul 2023 02:59
Operator : MA/JU
Sample : 03244-14
Misc :
ALS Vial : 42 Sample Multiplier: 1

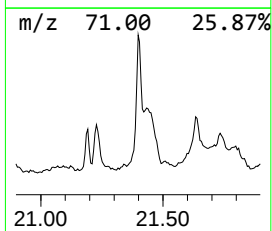
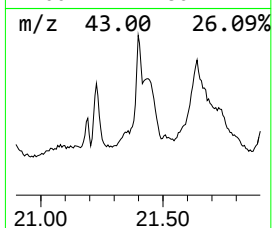
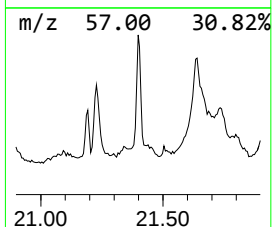
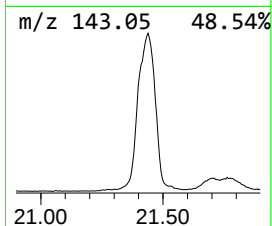
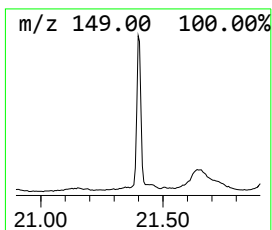
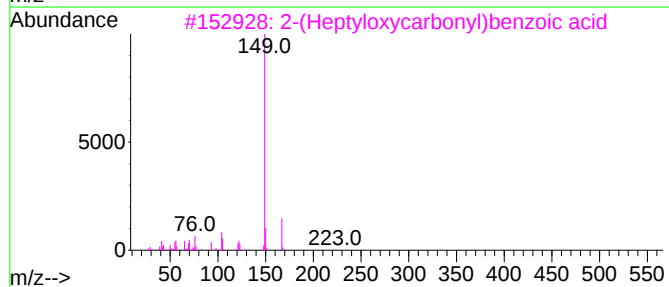
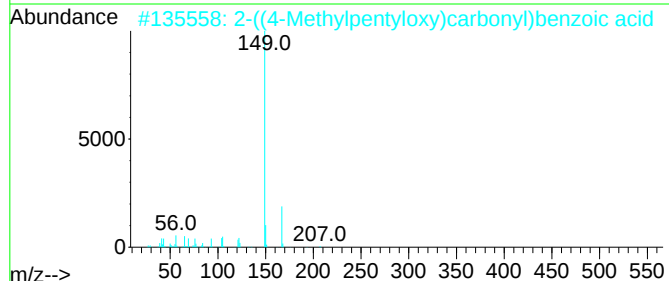
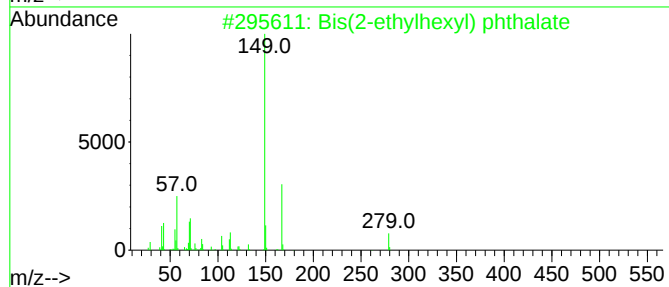
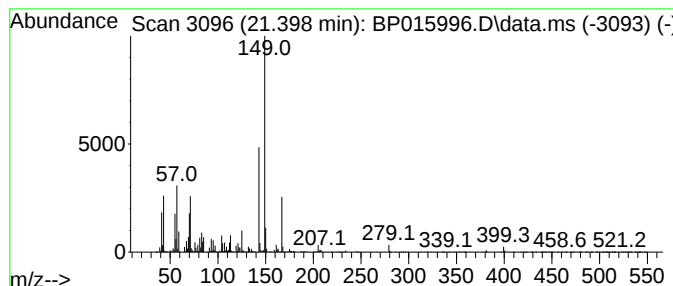
Instrument :
BNA_P
ClientSampleId :
DCGF3

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 Bis(2-ethylhexyl) phthalate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.398	14.01 ng/ul	2278550	Chrysene-d12	21.545	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	60
2	2-((4-Methylpentyloxy)carbonyl)b...	250	C14H18O4	848131-14-6	46
3	2-(Heptyloxy)carbonyl)benzoic acid	264	C15H20O4	024539-58-0	45
4	Phthalic acid, 4,4-dimethylpent...	516	C33H56O4	1000371-12-8	43
5	Phthalic acid, 2-ethylhexyl isoh...	362	C22H34O4	1000308-98-5	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015996.D
Acq On : 05 Jul 2023 02:59
Operator : MA/JU
Sample : 03244-14
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGF3

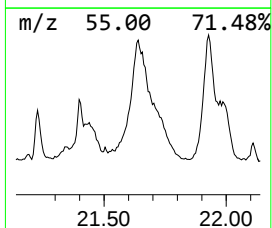
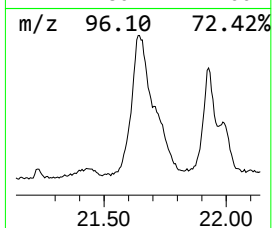
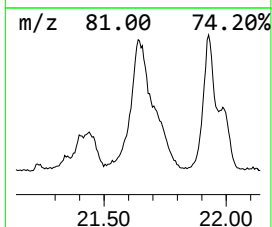
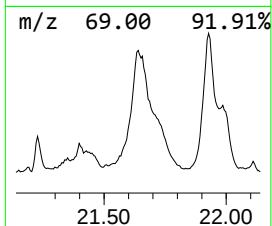
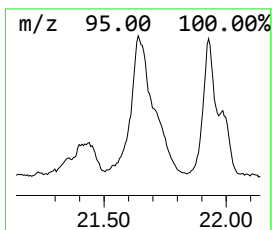
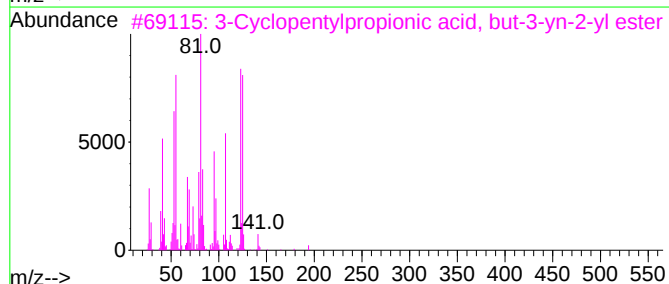
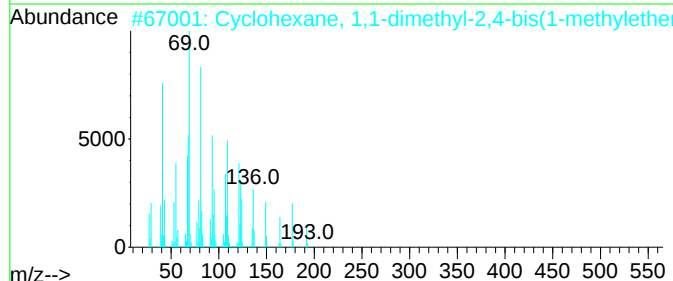
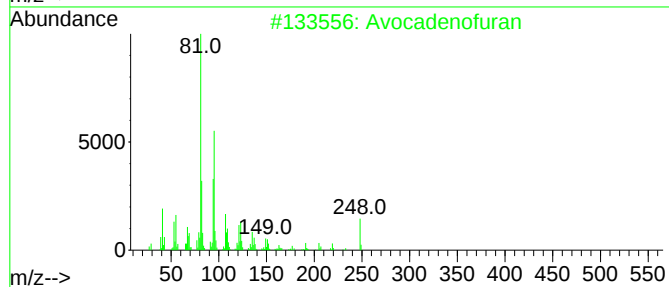
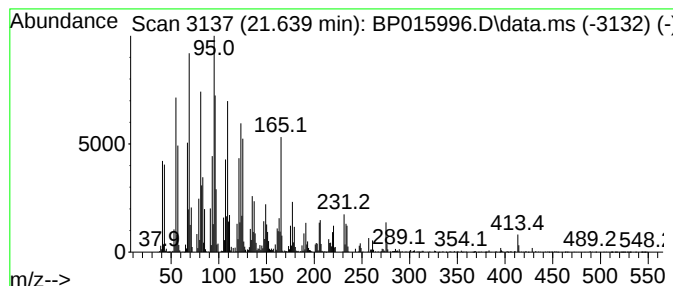
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 Avocadenofuran Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.639	135.49 ng/ul	22031800	Chrysene-d12	21.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Avocadenofuran	248	C17H28O	025346-24-1	64
2			Cyclohexane, 1,1-dimethyl-2,4-bi...	192	C14H24	062337-98-8	46
3			3-Cyclopentylpropionic acid, but...	194	C12H18O2	1000292-46-4	46
4			Cyclohexane, 3,4-bis(1-methyleth...	192	C14H24	061142-74-3	46
5			1,1,6-trimethyl-3-methylene-2-(3...	452	C33H56	1000373-94-5	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015996.D
Acq On : 05 Jul 2023 02:59
Operator : MA/JU
Sample : 03244-14
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGF3

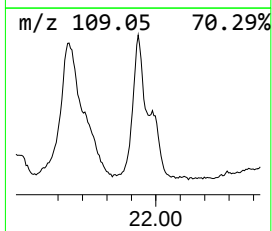
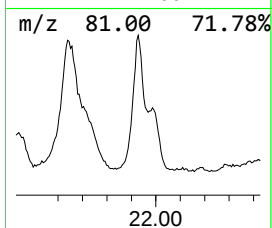
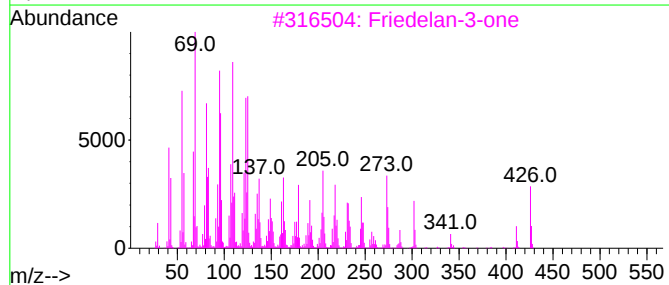
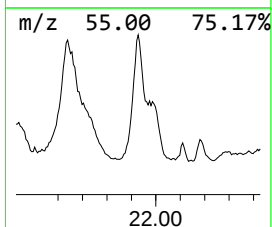
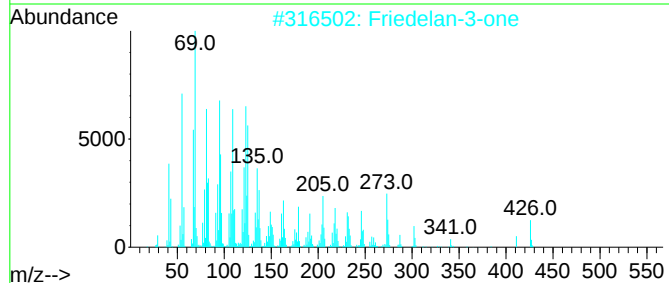
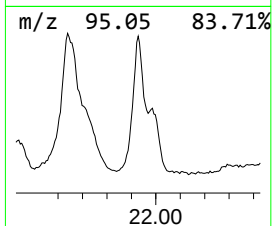
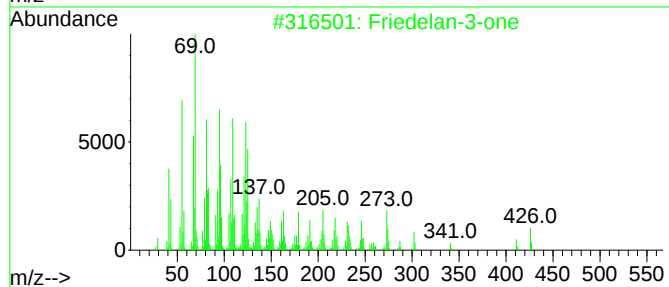
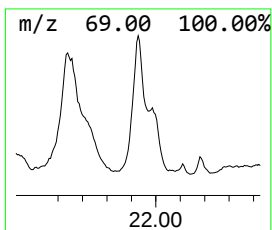
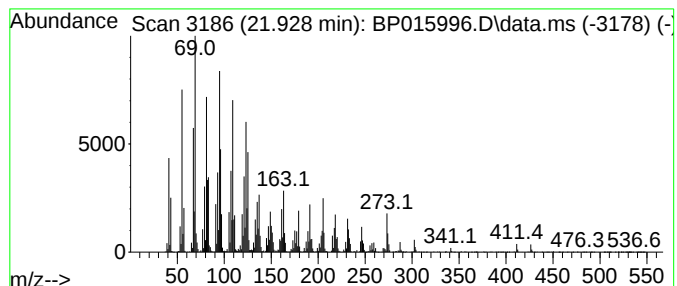
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 Sesquirosefuran Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.927	63.87 ng/ul	10385200	Chrysene-d12	21.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Friedelan-3-one	426	C30H50O	000559-74-0	99
2			Friedelan-3-one	426	C30H50O	000559-74-0	97
3			Friedelan-3-one	426	C30H50O	000559-74-0	95
4			Sesquirosefuran	218	C15H22O	039007-93-7	80
5			Friedelan-3-one	426	C30H50O	000559-74-0	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015996.D
Acq On : 05 Jul 2023 02:59
Operator : MA/JU
Sample : 03244-14
Misc :
ALS Vial : 42 Sample Multiplier: 1

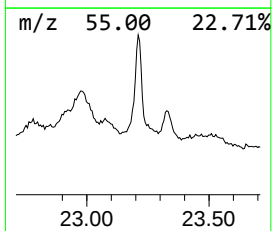
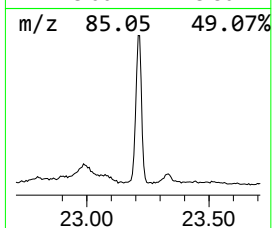
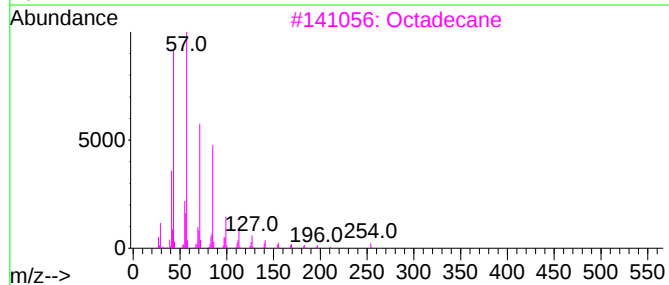
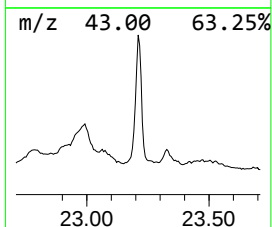
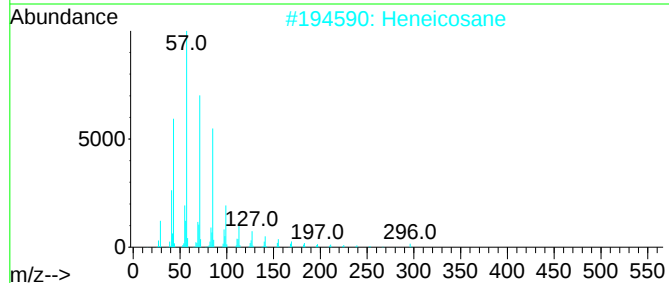
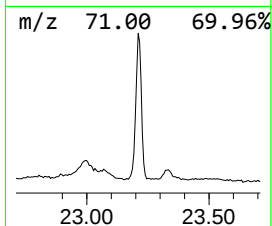
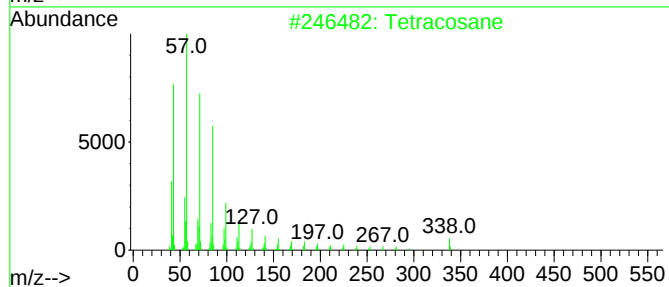
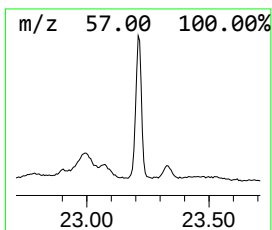
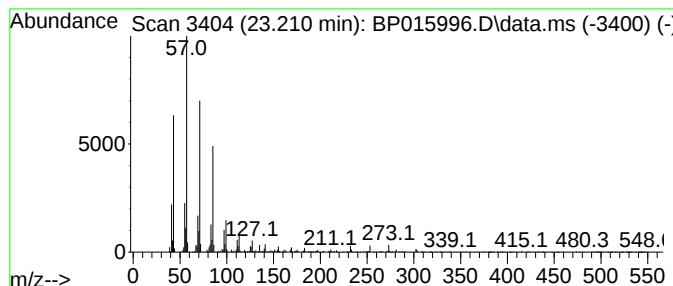
Instrument :
BNA_P
ClientSampleId :
DCGF3

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 11 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
23.210	27.24 ng/ul	2901640	Perylene-d12		24.086	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetracosane		338	C24H50	000646-31-1	95
2	Heneicosane		296	C21H44	000629-94-7	87
3	Octadecane		254	C18H38	000593-45-3	83
4	Heneicosane		296	C21H44	000629-94-7	83
5	Octacosane		394	C28H58	000630-02-4	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015996.D
Acq On : 05 Jul 2023 02:59
Operator : MA/JU
Sample : 03244-14
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGF3

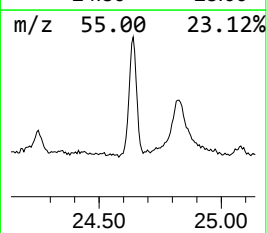
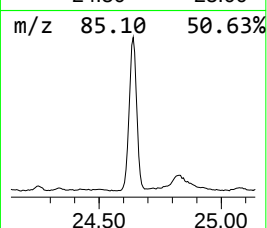
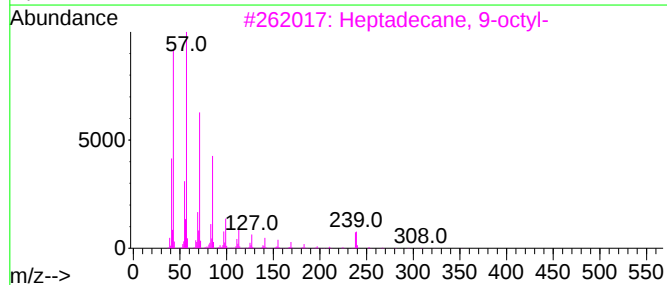
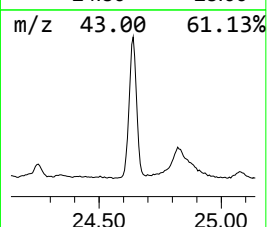
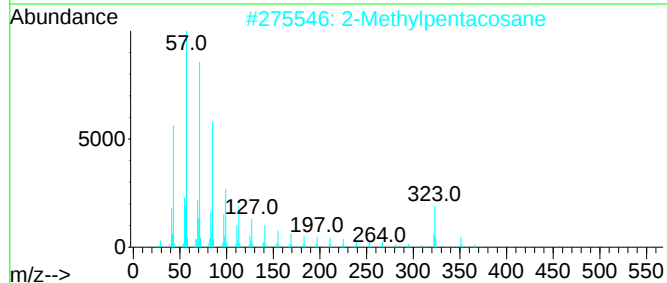
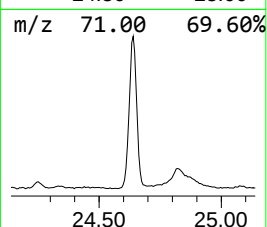
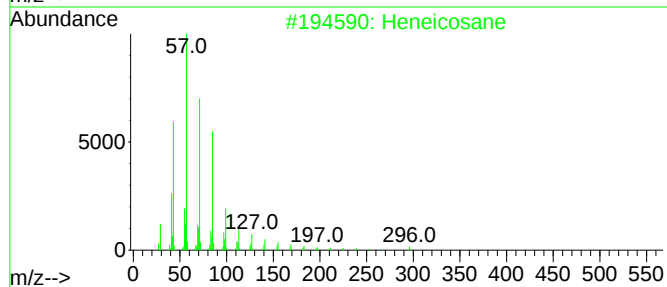
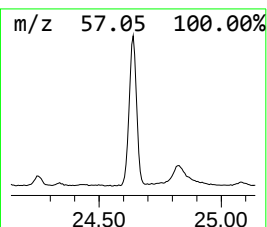
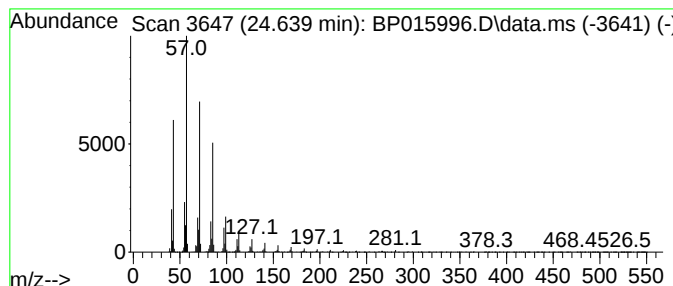
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.639	42.86 ng/ul	4565230	Perylene-d12	24.086

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	97
2		2-Methylpentacosane	366	C26H54	000629-87-8	95
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
4		Octacosane	394	C28H58	000630-02-4	91
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
 Data File : BP015996.D
 Acq On : 05 Jul 2023 02:59
 Operator : MA/JU
 Sample : 03244-14
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGF3

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
unknown-01	14.899	4.6	ng/u1	608979	3	14.693	2627960	20.0
Benzophenone	16.063	4.8	ng/u1	631273	3	14.693	2627960	20.0
(DEL) Alkane: S...	16.387	2.2	ng/u1	248910	4	17.457	2272030	20.0
cis-Vaccenic acid	18.251	8.1	ng/u1	922780	4	17.457	2272030	20.0
n-Hexadecanoic ...	18.310	32.9	ng/u1	3742370	4	17.457	2272030	20.0
Octadecanoic acid	19.528	9.2	ng/u1	1503120	5	21.545	3252160	20.0
(DEL) Alkane: S...	21.228	5.1	ng/u1	833174	5	21.545	3252160	20.0
Bis(2-ethylhexy...	21.398	14.0	ng/u1	2278550	5	21.545	3252160	20.0
Avocadenofuran	21.639	135.5	ng/u1	22031800	5	21.545	3252160	20.0
Sesquirosefuran	21.927	63.9	ng/u1	10385200	5	21.545	3252160	20.0
(DEL) Alkane: S...	23.210	27.2	ng/u1	2901640	6	24.086	2130290	20.0
(DEL) Alkane: S...	24.639	42.9	ng/u1	4565230	6	24.086	2130290	20.0