

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
 Data File : BP015747.D
 Acq On : 27 Jun 2023 15:13
 Operator : MA/JU
 Sample : 03085-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFN9

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: BP015747.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.387	30	34	40	rVB	29461	37406	0.53%	0.094%
2	3.834	108	110	123	rBV	180990	286871	4.06%	0.722%
3	3.934	123	127	135	rVV2	40034	60390	0.86%	0.152%
4	4.052	143	147	151	rVB	18129	27490	0.39%	0.069%
5	4.158	161	165	170	rVB	13386	19168	0.27%	0.048%
6	4.734	261	263	269	rVB7	6849	8838	0.13%	0.022%
7	5.105	323	326	331	rBV	4777	7211	0.10%	0.018%
8	6.711	595	599	604	rVB4	15782	25039	0.35%	0.063%
9	7.234	680	688	705	rBV	262165	625563	8.86%	1.575%
10	7.387	707	714	726	rBV	280716	488677	6.92%	1.230%
11	7.593	743	749	768	rVB	381509	692450	9.81%	1.743%
12	8.063	820	829	841	rBV	377570	692463	9.81%	1.743%
13	8.193	847	851	859	rVB10	4478	9286	0.13%	0.023%
14	8.269	859	864	869	rVV8	4024	7263	0.10%	0.018%
15	8.799	947	954	974	rBV2	146454	395357	5.60%	0.995%
16	9.252	1023	1031	1049	rBV	316112	672586	9.53%	1.693%
17	9.599	1087	1090	1096	rVB7	4859	8657	0.12%	0.022%
18	9.981	1148	1155	1168	rBV	256027	564657	8.00%	1.422%
19	10.546	1243	1251	1274	rBV	231879	736533	10.43%	1.854%
20	10.893	1300	1310	1329	rBV	463509	906213	12.84%	2.281%
21	11.069	1333	1340	1362	rBV2	97087	360050	5.10%	0.906%
22	12.510	1578	1585	1591	rBV8	4807	10707	0.15%	0.027%
23	13.063	1673	1679	1690	rVB8	11040	24901	0.35%	0.063%
24	13.316	1718	1722	1727	rVB8	4473	7239	0.10%	0.018%
25	13.575	1761	1766	1776	rVV7	15650	31881	0.45%	0.080%
26	13.698	1782	1787	1791	rBV8	5614	10346	0.15%	0.026%
27	14.022	1838	1842	1848	rVB8	14009	21052	0.30%	0.053%
28	14.122	1851	1859	1874	rVV	702661	1143265	16.19%	2.878%
29	14.222	1874	1876	1880	rVB5	6537	7854	0.11%	0.020%
30	14.304	1886	1890	1894	rVB	8183	10781	0.15%	0.027%
31	14.410	1901	1908	1921	rBV	875258	1405504	19.91%	3.538%
32	14.651	1945	1949	1953	rVV4	16920	23630	0.33%	0.059%
33	14.710	1953	1959	1980	rVB	699875	1130745	16.02%	2.847%
34	14.940	1995	1998	2000	rBV4	9910	13144	0.19%	0.033%
35	14.981	2000	2005	2024	rVV2	88806	328469	4.65%	0.827%
36	15.128	2026	2030	2039	rVB2	12660	36980	0.52%	0.093%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
 Data File : BP015747.D
 Acq On : 27 Jun 2023 15:13
 Operator : MA/JU
 Sample : 03085-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFN9

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	15.534	2095	2099	2103	rBV7	7412	10332	0.15%	0.026%
38	15.710	2122	2129	2137	rBV	1028528	1680421	23.80%	4.230%
39	15.869	2149	2156	2167	rBV2	181633	426159	6.04%	1.073%
40	16.075	2186	2191	2203	rBV	476872	704909	9.98%	1.775%
41	16.175	2205	2208	2212	rVB5	14467	20524	0.29%	0.052%
42	16.416	2243	2249	2255	rVB10	9195	23602	0.33%	0.059%
43	16.557	2270	2273	2281	rBV6	22444	39834	0.56%	0.100%
44	16.639	2281	2287	2294	rVB	107729	162061	2.30%	0.408%
45	16.857	2321	2324	2328	rBV5	9605	10557	0.15%	0.027%
46	17.010	2346	2350	2354	rVB3	12019	18724	0.27%	0.047%
47	17.198	2379	2382	2384	rBV3	9792	11484	0.16%	0.029%
48	17.345	2403	2407	2414	rBV2	48968	75153	1.06%	0.189%
49	17.416	2415	2419	2422	rBV4	33117	47762	0.68%	0.120%
50	17.475	2423	2429	2435	rBV	894565	1333240	18.88%	3.356%
51	17.575	2440	2446	2459	rVB	1025552	1629024	23.07%	4.101%
52	18.216	2553	2555	2560	rVB5	27807	35189	0.50%	0.089%
53	18.275	2561	2565	2569	rBV	70705	91263	1.29%	0.230%
54	18.328	2569	2574	2583	rBV	838213	1162623	16.47%	2.927%
55	18.610	2618	2622	2629	rBV4	52780	117099	1.66%	0.295%
56	18.804	2648	2655	2656	rBV7	63384	127251	1.80%	0.320%
57	19.210	2721	2724	2730	rBV7	31908	46461	0.66%	0.117%
58	19.416	2756	2759	2761	rBV2	59962	74031	1.05%	0.186%
59	19.439	2761	2763	2765	rBV3	52770	59866	0.85%	0.151%
60	19.557	2779	2783	2791	rBV	250875	372968	5.28%	0.939%
61	19.798	2819	2824	2834	rBV	1562221	2447248	34.66%	6.161%
62	21.092	3040	3044	3049	rBV	243418	321101	4.55%	0.808%
63	21.233	3063	3068	3077	rVB2	522976	916665	12.98%	2.308%
64	21.427	3099	3101	3105	rVB	164426	188137	2.66%	0.474%
65	21.563	3120	3124	3135	rVB	1046902	1535130	21.74%	3.865%
66	22.192	3228	3231	3239	rVB2	412664	682657	9.67%	1.719%
67	23.257	3408	3412	3419	rVB	546901	877318	12.43%	2.209%
68	23.345	3423	3427	3436	rVB5	315392	679662	9.63%	1.711%
69	23.945	3522	3529	3543	rVB2	1138074	2736596	38.76%	6.889%
70	24.110	3552	3557	3565	rVB	712035	1515773	21.47%	3.816%
71	24.704	3652	3658	3666	rVB	782619	1642695	23.27%	4.135%
72	28.809	4346	4356	4374	rVB3	1784012	7060221	100.00%	17.774%

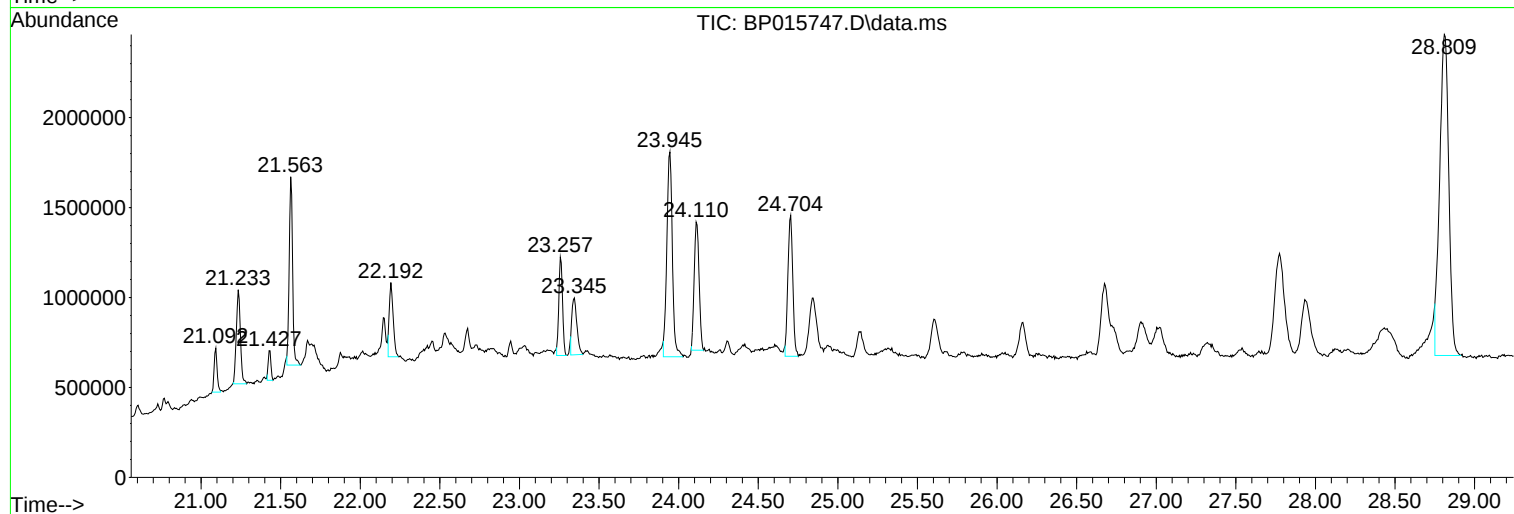
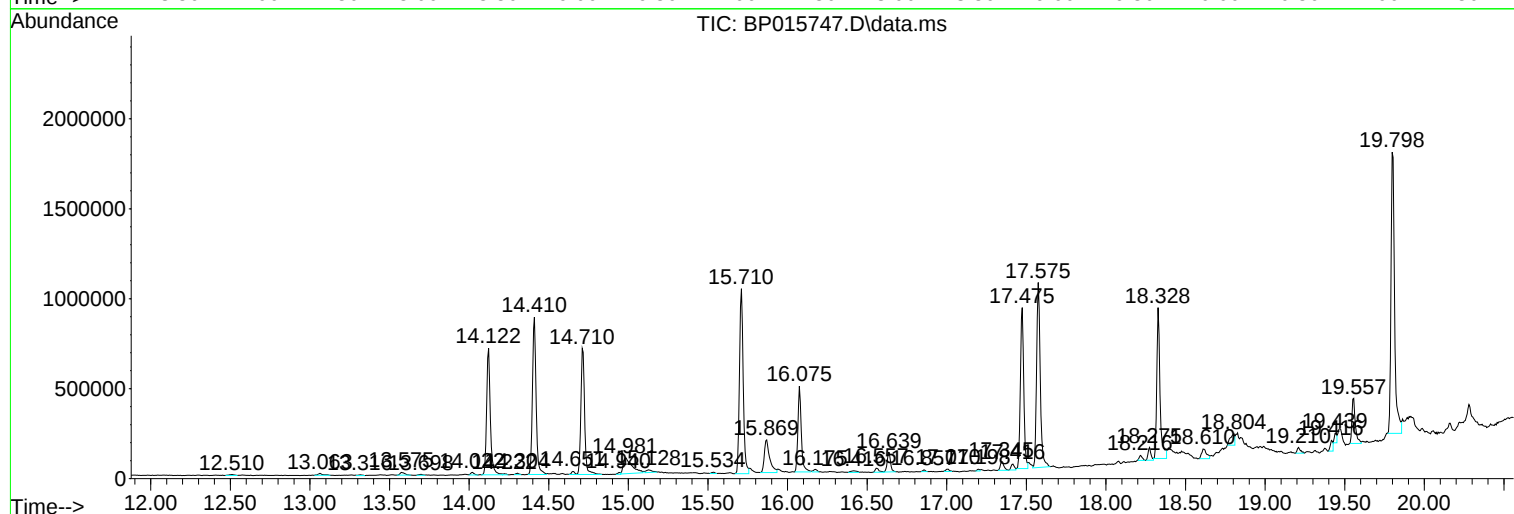
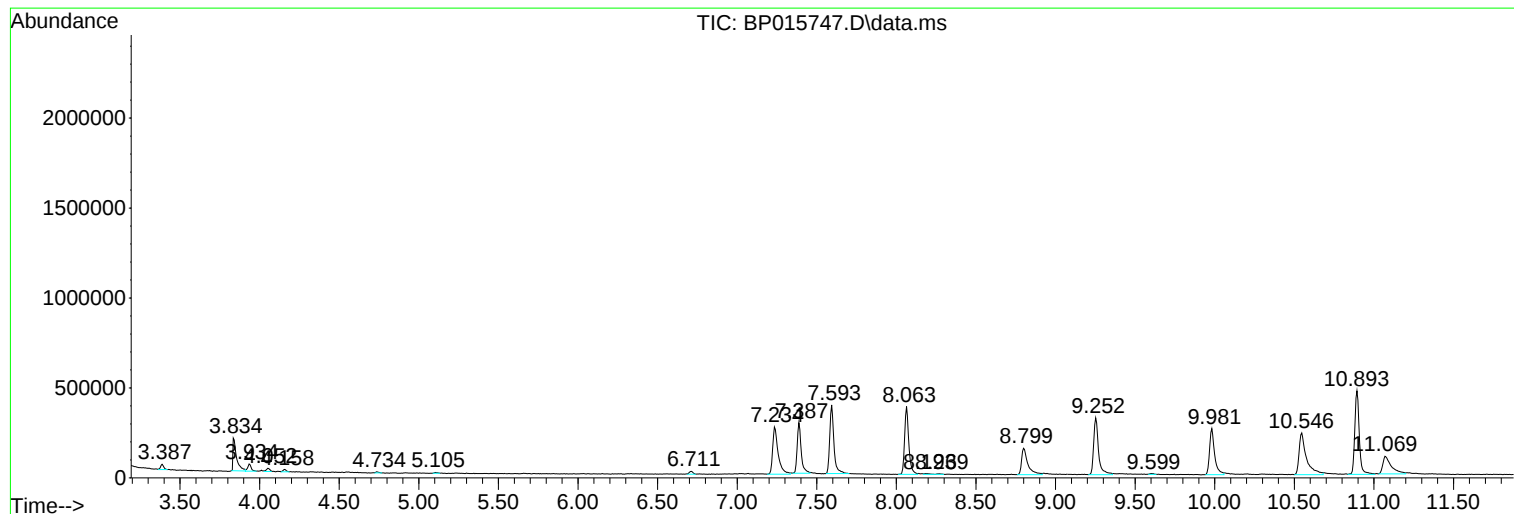
Sum of corrected areas: 39722406

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
Data File : BP015747.D
Acq On : 27 Jun 2023 15:13
Operator : MA/JU
Sample : 03085-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN9

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\BP062723\
Data File : BP015747.D
Acq On : 27 Jun 2023 15:13
Operator : MA/JU
Sample : 03085-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN9

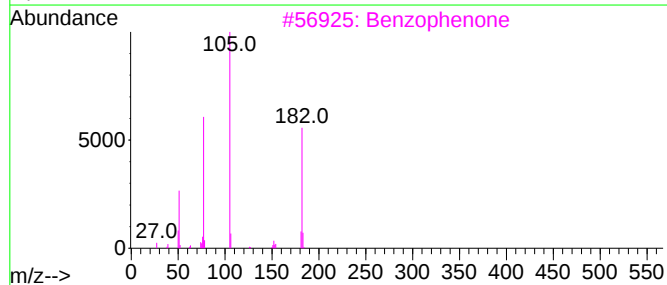
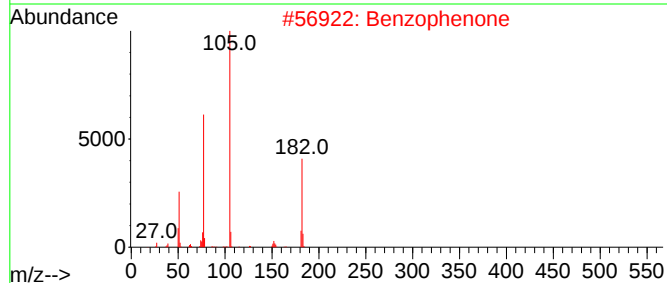
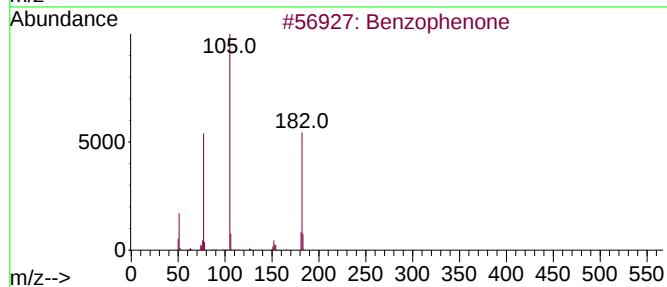
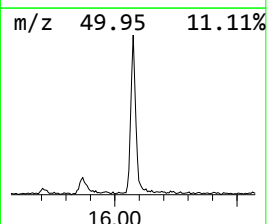
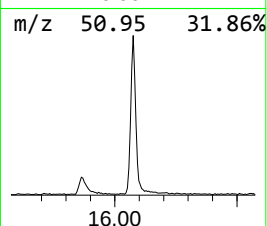
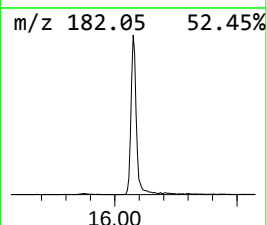
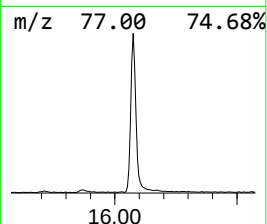
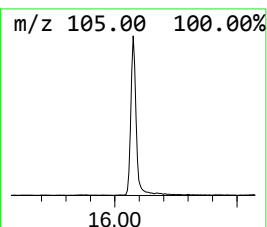
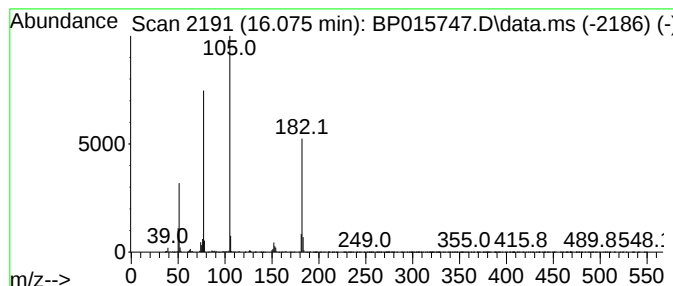
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.075	12.47 ng/ul	704909	Acenaphthene-d10	14.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	96
4			Benzophenone	182	C13H10O	000119-61-9	94
5			Benzophenone	182	C13H10O	000119-61-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
Data File : BP015747.D
Acq On : 27 Jun 2023 15:13
Operator : MA/JU
Sample : 03085-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN9

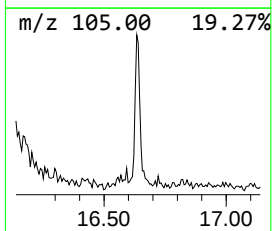
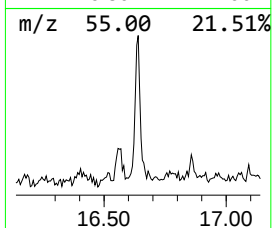
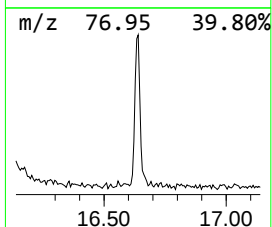
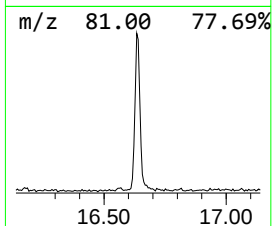
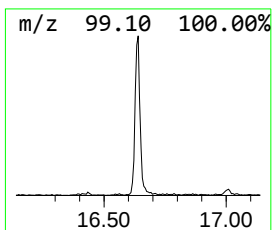
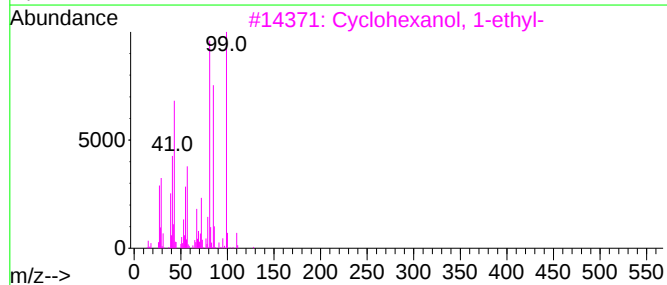
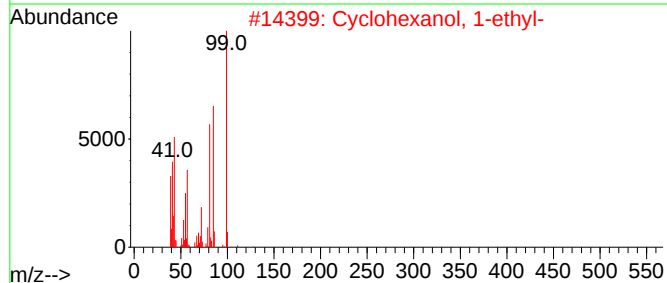
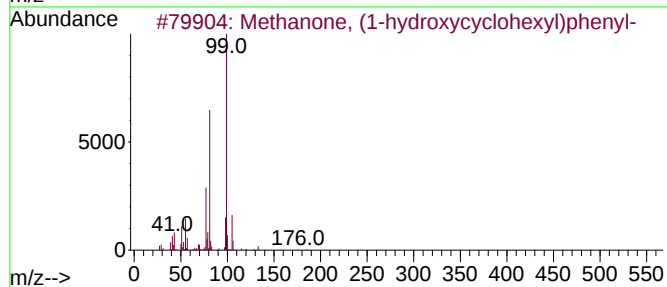
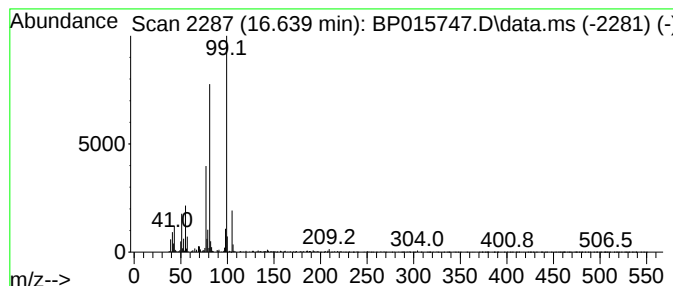
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Methanone, (1-hydroxycyclohexyl) Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.640	2.43 ng/ul	162061	Phenanthrene-d10	17.475

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methanone, (1-hydroxycyclohexyl)...	204	C13H16O2	000947-19-3	91
2			Cyclohexanol, 1-ethyl-	128	C8H16O	001940-18-7	50
3			Cyclohexanol, 1-ethyl-	128	C8H16O	001940-18-7	40
4			3-Methylcyclohexyl methylphospho...	194	C8H16FO2P	113548-86-0	40
5			Cyclohexanol, 1-(aminomethyl)-	129	C7H15NO	004000-72-0	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
Data File : BP015747.D
Acq On : 27 Jun 2023 15:13
Operator : MA/JU
Sample : 03085-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN9

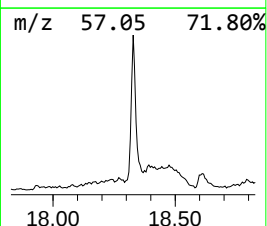
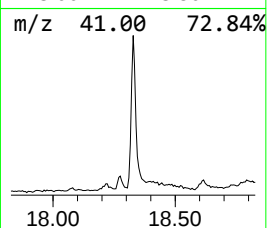
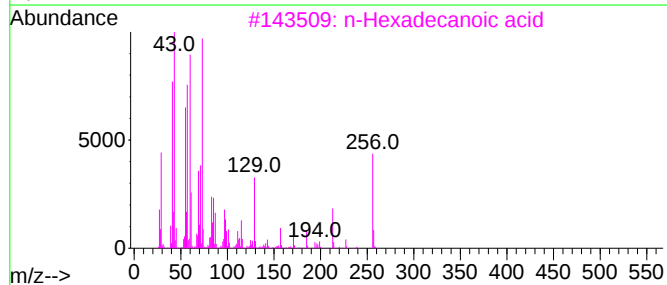
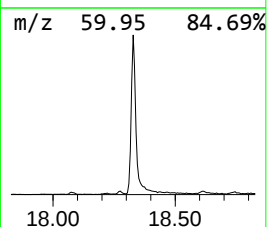
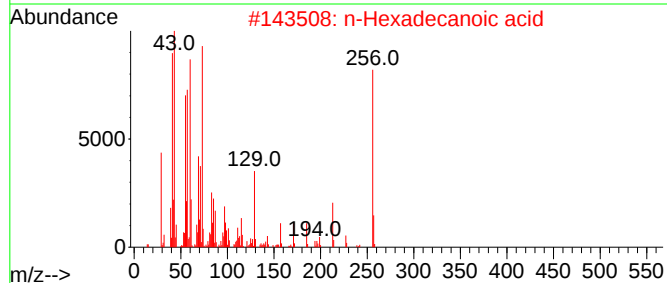
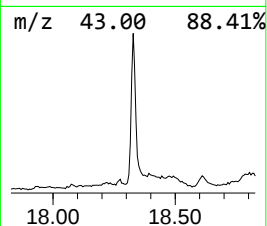
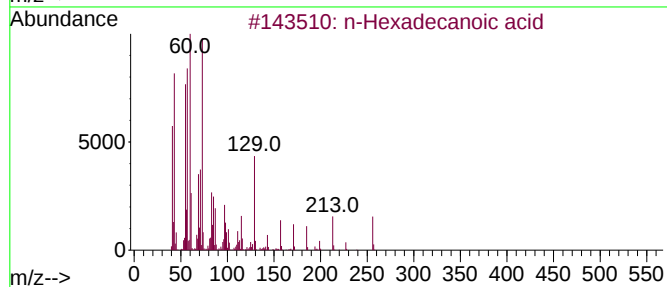
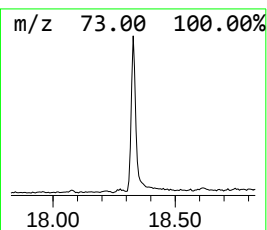
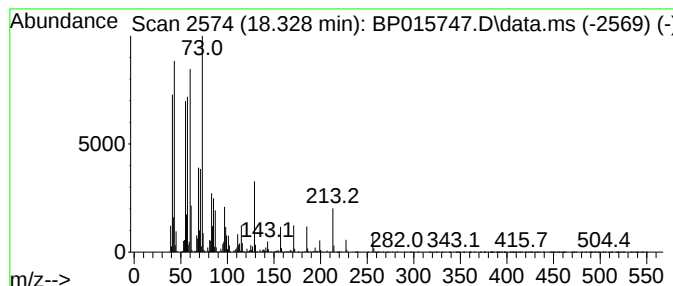
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.328	17.44 ng/ul	1162620	Phenanthrene-d10	17.475

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	95		
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
Data File : BP015747.D
Acq On : 27 Jun 2023 15:13
Operator : MA/JU
Sample : 03085-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN9

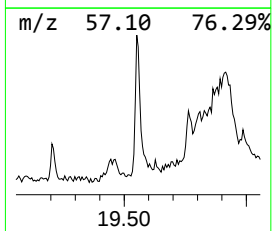
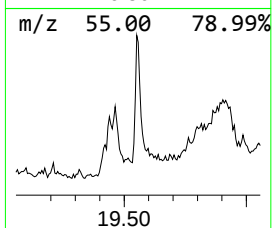
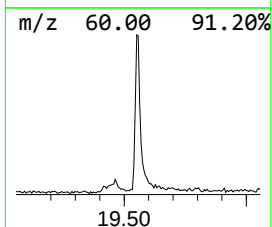
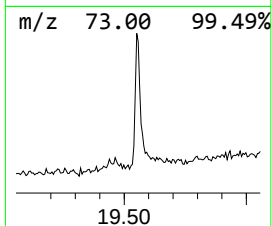
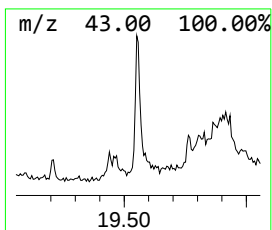
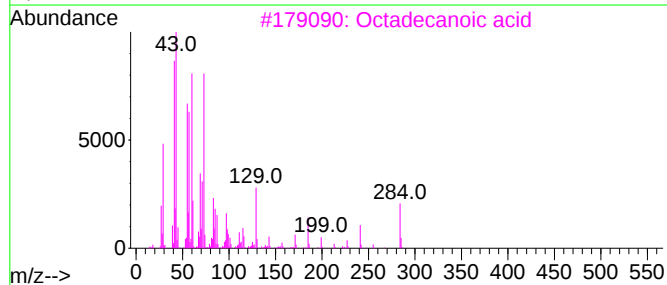
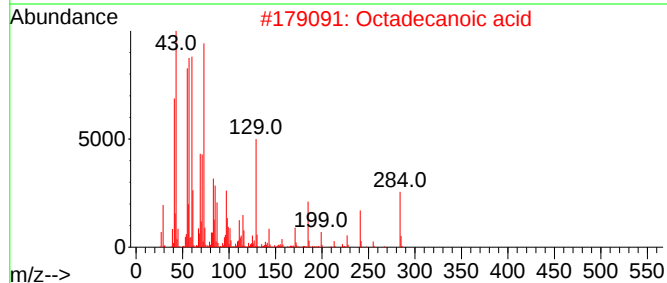
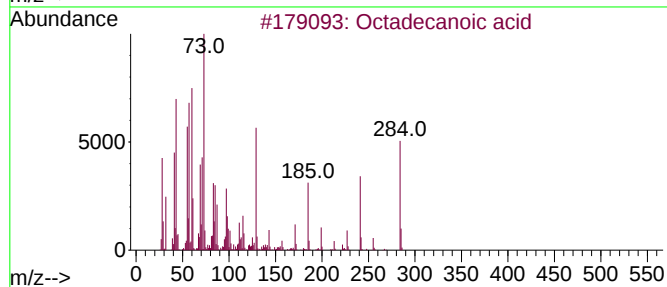
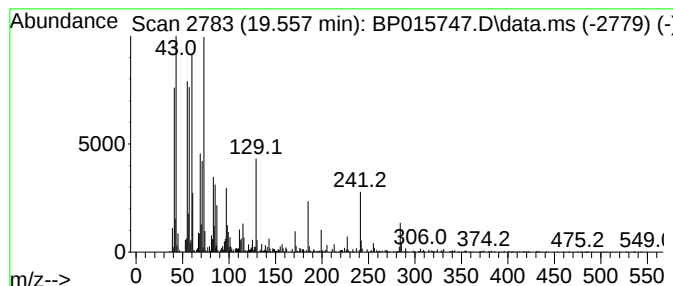
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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.
19.557	4.86 ng/ul	372968	Chrysene-d12	21.563

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	98
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
Data File : BP015747.D
Acq On : 27 Jun 2023 15:13
Operator : MA/JU
Sample : 03085-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN9

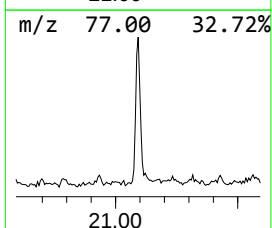
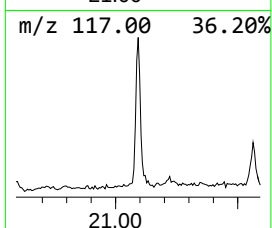
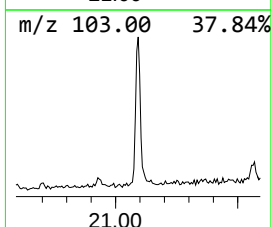
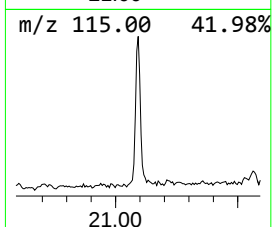
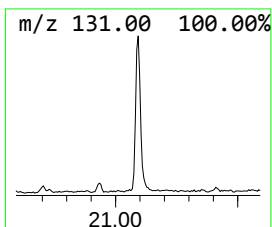
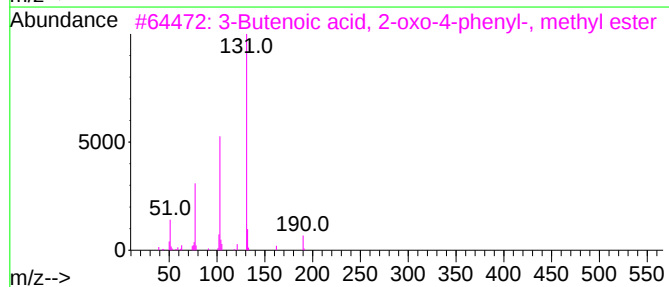
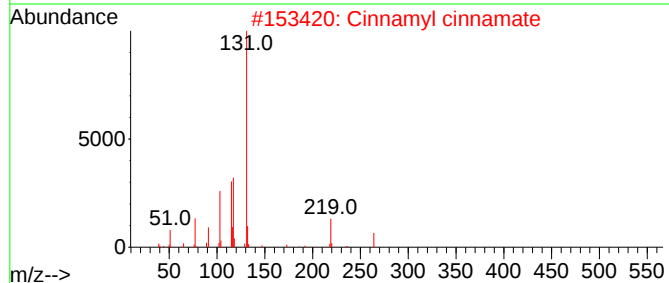
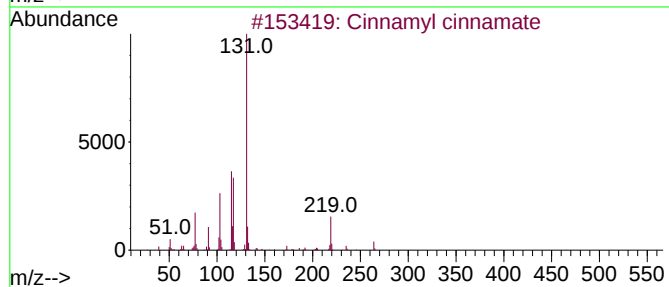
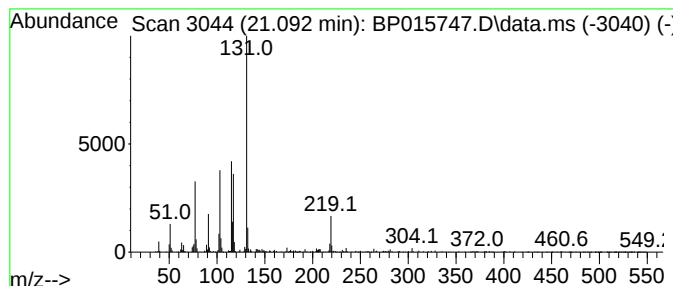
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Cinnamyl cinnamate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.092	4.18 ng/ul	321101	Chrysene-d12	21.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cinnamyl cinnamate	264	C18H16O2	000122-69-0	91
2			Cinnamyl cinnamate	264	C18H16O2	000122-69-0	86
3			3-Butenoic acid, 2-oxo-4-phenyl-...	190	C11H10O3	1000142-22-4	49
4			1-(4-Ethynylphenyl)-2-methylprop...	174	C12H14O	1000466-39-4	49
5			3-phenylprop-2-enoic anhydride	278	C18H14O3	1010401-87-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
Data File : BP015747.D
Acq On : 27 Jun 2023 15:13
Operator : MA/JU
Sample : 03085-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

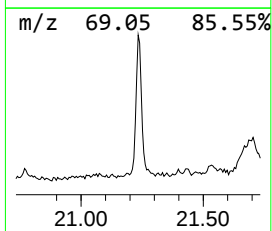
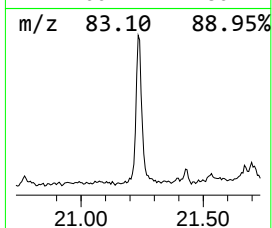
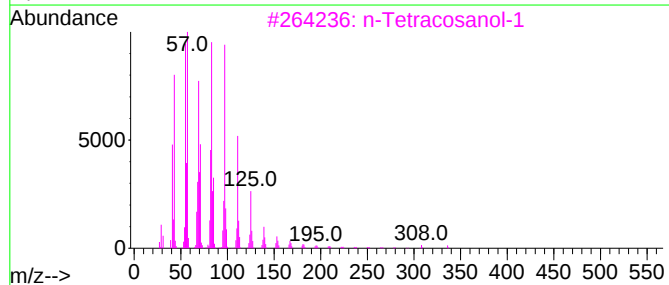
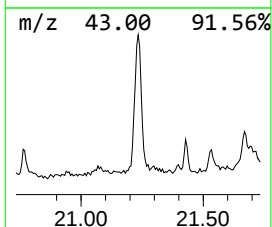
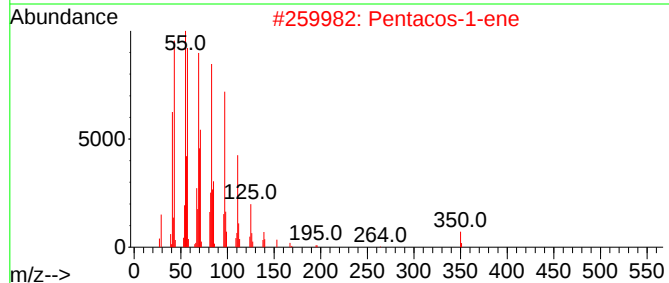
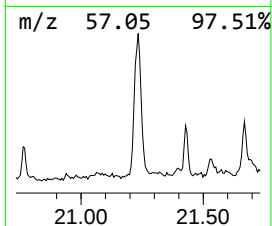
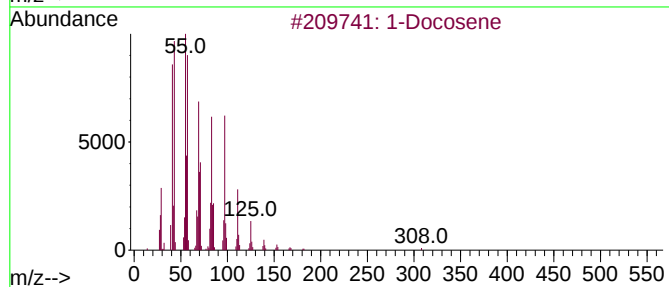
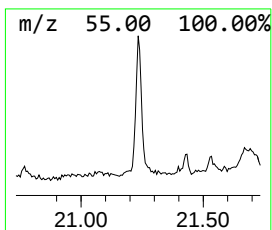
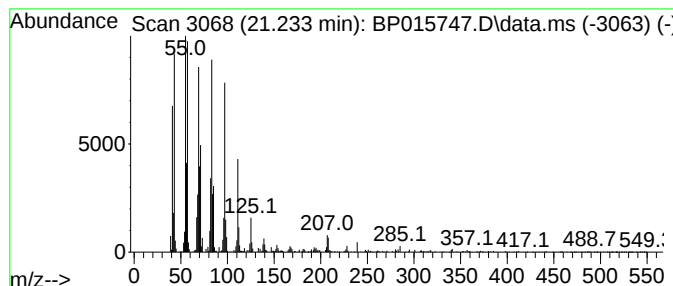
Instrument :
BNA_P
ClientSampleId :
DCFN9

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.233	11.94 ng/ul	916665	Chrysene-d12		21.563	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Docosene		308	C22H44	001599-67-3	96
2	Pentacos-1-ene		350	C25H50	016980-85-1	94
3	n-Tetracosanol-1		354	C24H50O	000506-51-4	94
4	1-Hexacosene		364	C26H52	018835-33-1	93
5	1-Nonadecene		266	C19H38	018435-45-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
Data File : BP015747.D
Acq On : 27 Jun 2023 15:13
Operator : MA/JU
Sample : 03085-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN9

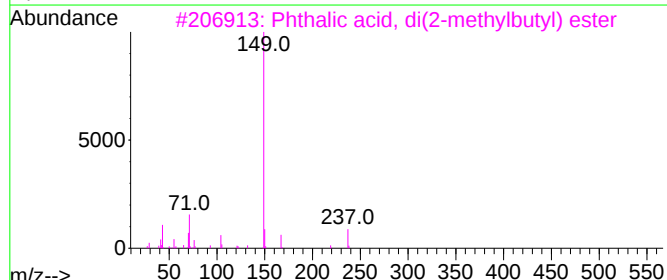
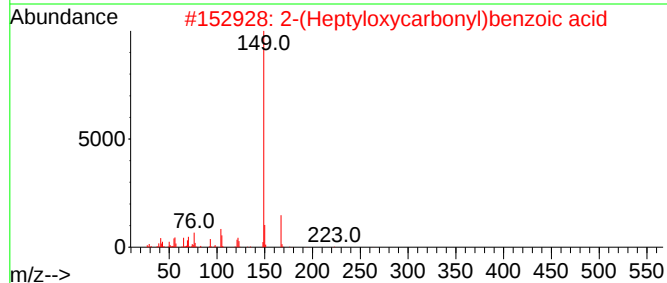
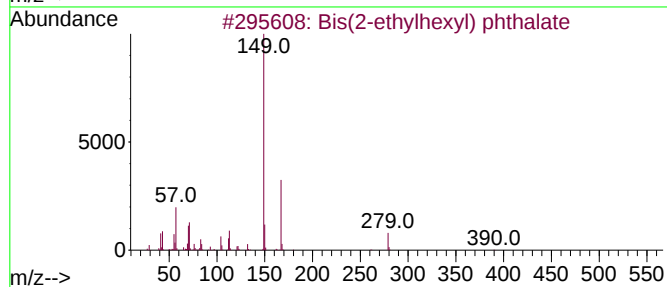
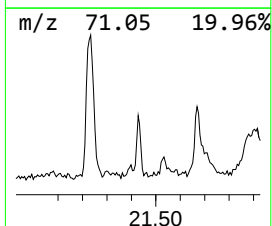
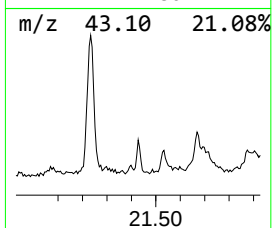
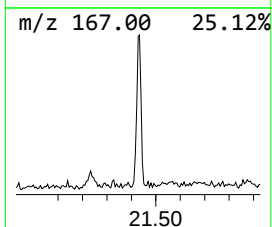
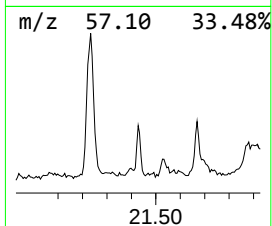
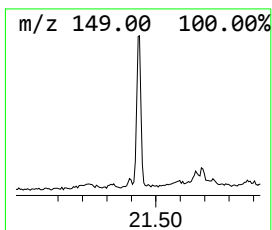
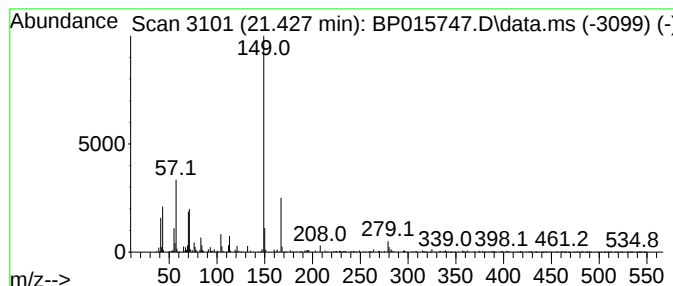
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Bis(2-ethylhexyl) phthalate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.427	2.45 ng/ul	188137	Chrysene-d12	21.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	68
2			2-(Heptyloxycarbonyl)benzoic acid	264	C15H20O4	024539-58-0	64
3			Phthalic acid, di(2-methylbutyl)...	306	C18H26O4	1000322-82-5	59
4			Phthalic acid, cyclohexyl 2-pent...	318	C19H26O4	1000315-55-3	59
5			Phthalic acid, cyclohexylmethyl ...	332	C20H28O4	1000315-55-5	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
Data File : BP015747.D
Acq On : 27 Jun 2023 15:13
Operator : MA/JU
Sample : 03085-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

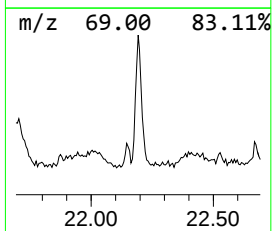
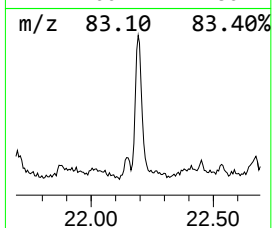
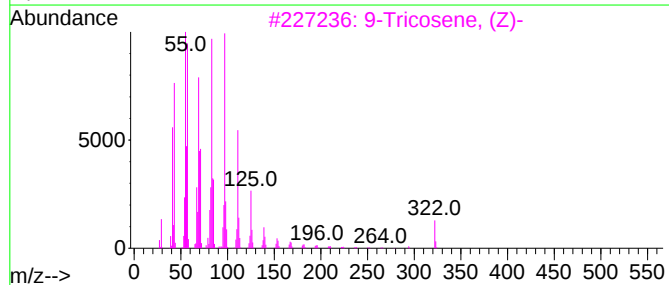
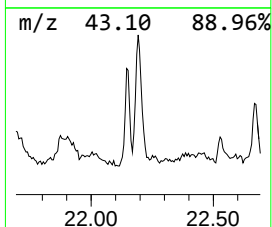
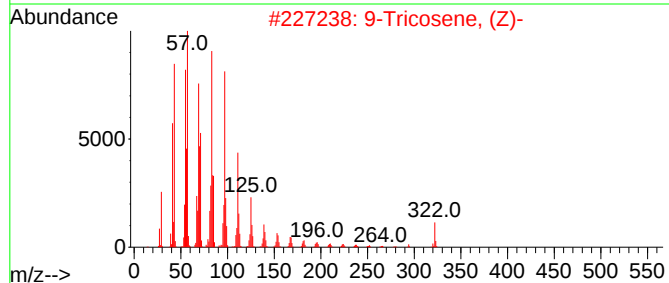
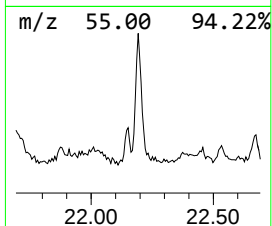
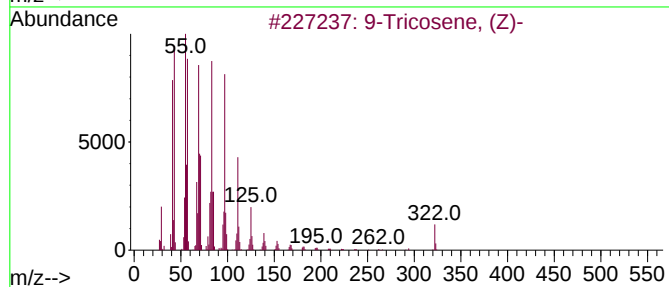
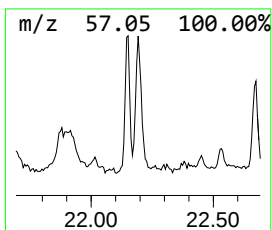
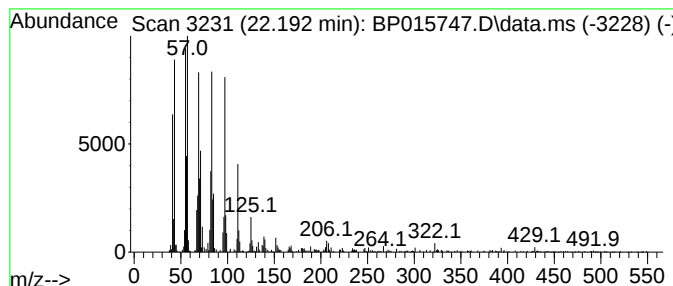
Instrument :
BNA_P
ClientSampleId :
DCFN9

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 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
22.192	8.89 ng/ul	682657	Chrysene-d12		21.563	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	9-Tricosene, (Z)-		322	C23H46	027519-02-4	96
2	9-Tricosene, (Z)-		322	C23H46	027519-02-4	95
3	9-Tricosene, (Z)-		322	C23H46	027519-02-4	95
4	2-Methyl-2-docosene		322	C23H46	1000131-16-9	94
5	Trifluoroacetoxy hexadecane		338	C18H33F3O2	006222-03-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
Data File : BP015747.D
Acq On : 27 Jun 2023 15:13
Operator : MA/JU
Sample : 03085-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

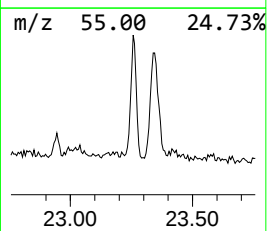
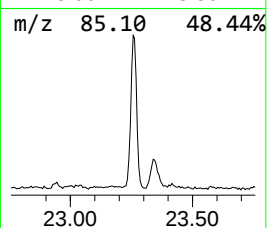
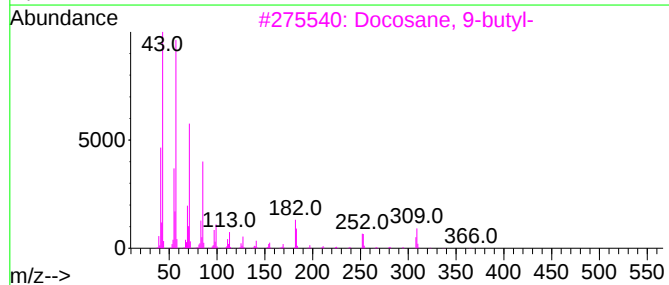
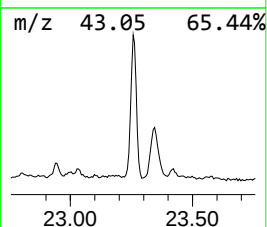
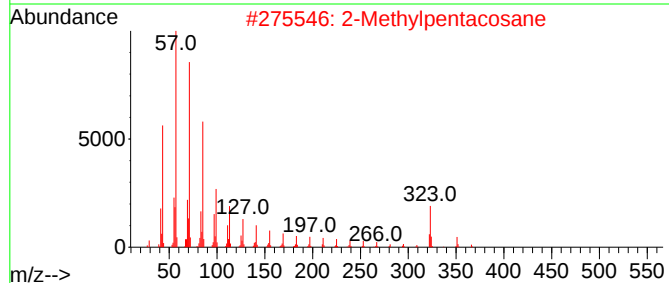
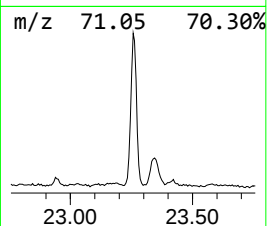
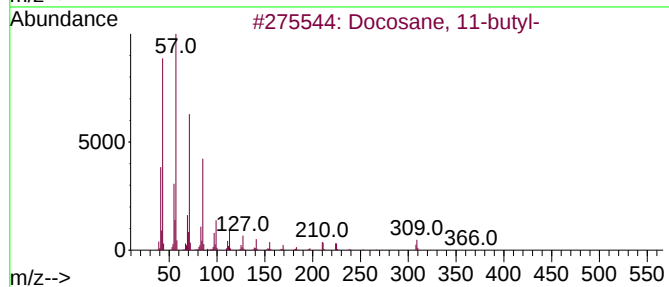
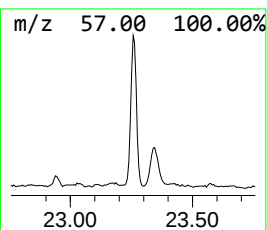
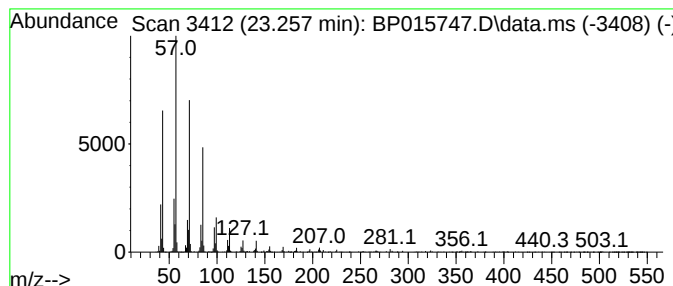
Instrument :
BNA_P
ClientSampleId :
DCFN9

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 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.257	11.58 ng/ul	877318	Perylene-d12	24.110	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-	366	C26H54	013475-76-8	93
2	2-Methylpentacosane	366	C26H54	000629-87-8	93
3	Docosane, 9-butyl-	366	C26H54	055282-14-9	93
4	Heneicosane	296	C21H44	000629-94-7	90
5	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
Data File : BP015747.D
Acq On : 27 Jun 2023 15:13
Operator : MA/JU
Sample : 03085-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

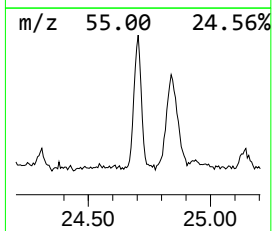
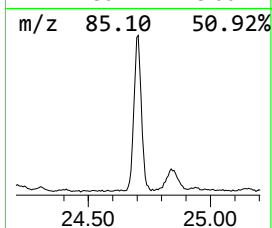
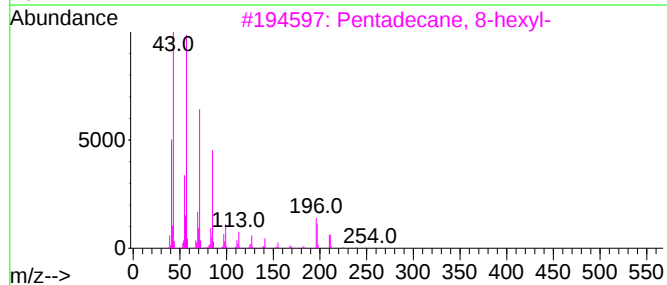
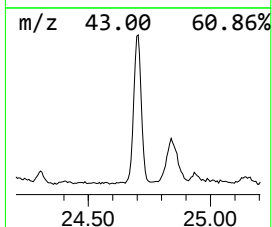
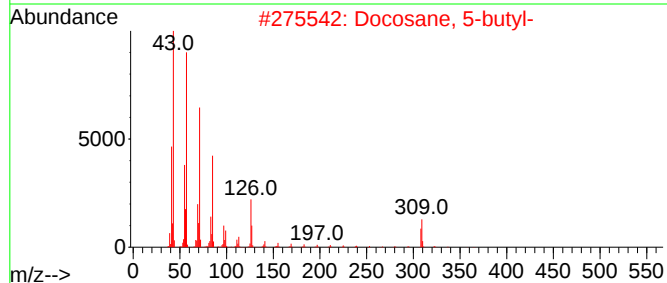
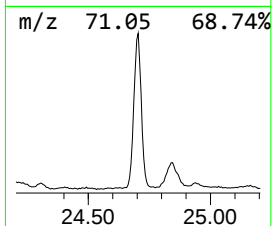
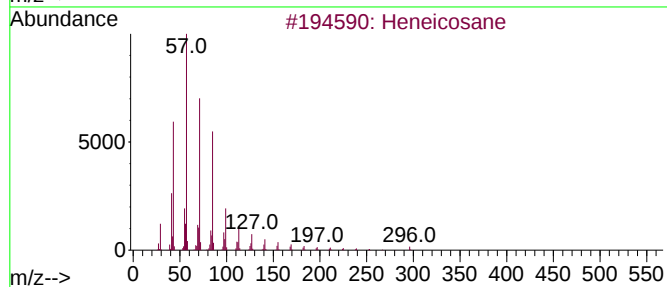
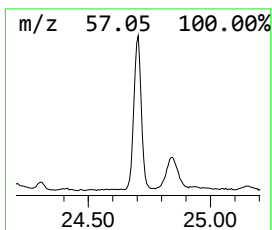
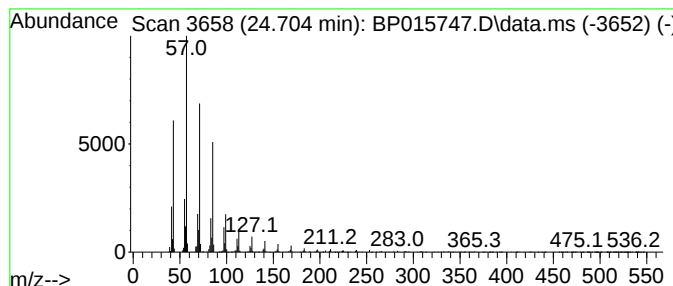
Instrument :
BNA_P
ClientSampleId :
DCFN9

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 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.704	21.67 ng/ul	1642700	Perylene-d12	24.110	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Docosane, 5-butyl-	366	C26H54	055282-16-1	91
3	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
4	Tetracosane	338	C24H50	000646-31-1	91
5	Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
Data File : BP015747.D
Acq On : 27 Jun 2023 15:13
Operator : MA/JU
Sample : 03085-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN9

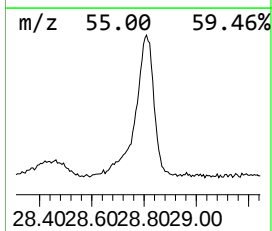
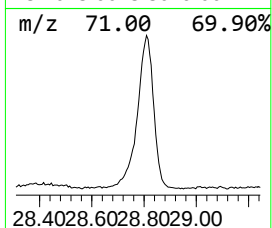
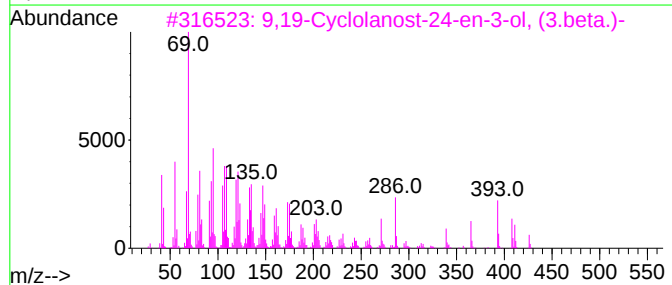
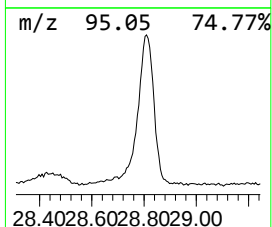
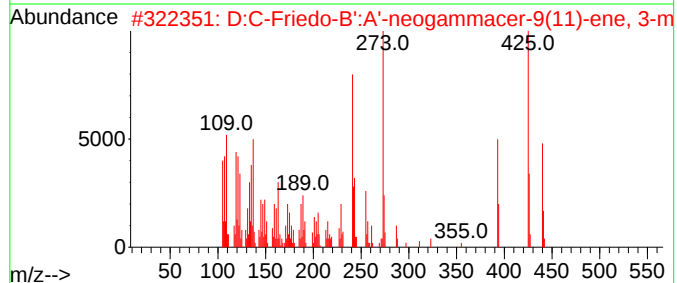
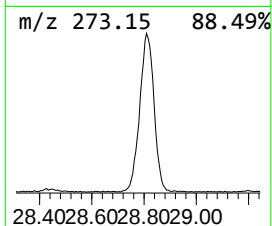
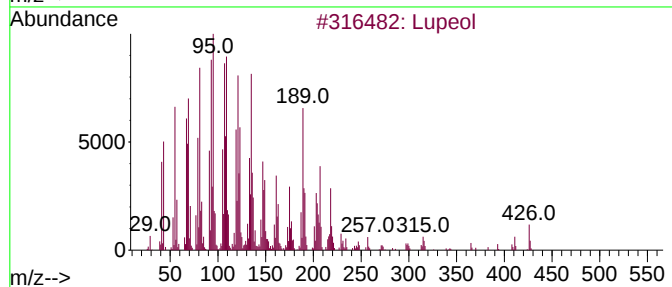
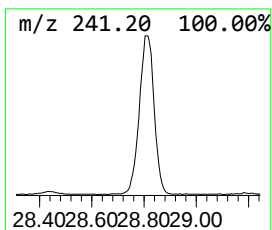
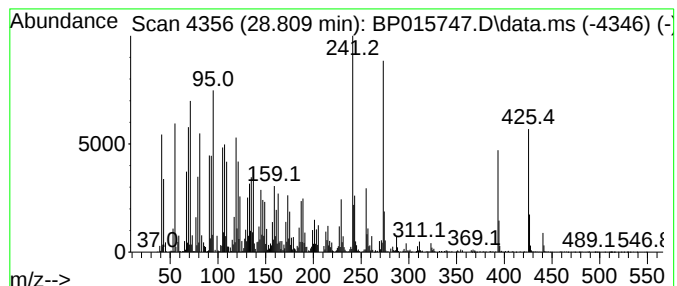
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.809	93.16 ng/ul	7060220	Perylene-d12	24.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Lupeol	426	C30H50O	000545-47-1	43
2			D:C-Friedo-B':A'-neogammacer-9(1...	440	C31H52O	004555-56-0	40
3			9,19-Cyclolanost-24-en-3-ol, (3...	426	C30H50O	000469-38-5	30
4			9,19-Cyclolanost-25-en-3-ol, 24-...	440	C31H52O	000511-61-5	30
5			2-(2-Benzothiazolyl)-5-methylphenol	241	C14H11NOS	056048-54-5	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
Data File : BP015747.D
Acq On : 27 Jun 2023 15:13
Operator : MA/JU
Sample : 03085-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN9

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
Benzophenone	16.075	12.5	ng/u1	704909	3	14.710	1130750	20.0
Methanone, (1-h...	16.640	2.4	ng/u1	162061	4	17.475	1333240	20.0
n-Hexadecanoic ...	18.328	17.4	ng/u1	1162620	4	17.475	1333240	20.0
Octadecanoic acid	19.557	4.9	ng/u1	372968	5	21.563	1535130	20.0
Cinnamyl cinnamate	21.092	4.2	ng/u1	321101	5	21.563	1535130	20.0
(DEL) Alkane: S...	21.233	11.9	ng/u1	916665	5	21.563	1535130	20.0
Bis(2-ethylhexy...	21.427	2.5	ng/u1	188137	5	21.563	1535130	20.0
(DEL) Alkane: S...	22.192	8.9	ng/u1	682657	5	21.563	1535130	20.0
(DEL) Alkane: S...	23.257	11.6	ng/u1	877318	6	24.110	1515770	20.0
(DEL) Alkane: S...	24.704	21.7	ng/u1	1642700	6	24.110	1515770	20.0
unknown-01	28.809	93.2	ng/u1	7060220	6	24.110	1515770	20.0