

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121324\
 Data File : BP023422.D
 Acq On : 13 Dec 2024 16:11
 Operator : RC/JU
 Sample : P5232-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E28S3

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-BP112624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023422.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.099	16	19	26	rVB	17116	20422	0.83%	0.057%
2	3.517	86	90	102	rBV	170764	291044	11.80%	0.814%
3	6.752	634	640	650	rBV	225514	456864	18.52%	1.277%
4	6.887	658	663	675	rVV	183191	308685	12.51%	0.863%
5	7.087	691	697	711	rBV	271582	437033	17.71%	1.222%
6	7.534	763	773	783	rBV	461777	777764	31.52%	2.174%
7	8.263	890	897	918	rBV2	188557	423439	17.16%	1.184%
8	8.681	961	968	984	rBV	229716	435755	17.66%	1.218%
9	9.246	1058	1064	1077	rVB	22051	44684	1.81%	0.125%
10	9.393	1082	1089	1102	rBV	176390	354402	14.36%	0.991%
11	9.940	1176	1182	1204	rBV	275535	656846	26.62%	1.836%
12	10.281	1232	1240	1247	rBV	666617	1082220	43.86%	3.025%
13	10.440	1259	1267	1280	rBV	208352	454574	18.42%	1.271%
14	12.957	1690	1695	1703	rVV8	8398	16028	0.65%	0.045%
15	13.163	1725	1730	1741	rVB3	19715	49088	1.99%	0.137%
16	13.551	1787	1796	1804	rBV	712416	1026078	41.59%	2.868%
17	13.834	1839	1844	1858	rVV	682748	1003992	40.69%	2.807%
18	14.151	1891	1898	1912	rBV	1172724	1656299	67.13%	4.630%
19	14.463	1944	1951	1970	rBV2	111275	402823	16.33%	1.126%
20	15.163	2064	2070	2081	rBV	810232	1237285	50.15%	3.459%
21	15.334	2095	2099	2107	rVB8	13821	25052	1.02%	0.070%
22	15.540	2127	2134	2140	rBV	350432	491426	19.92%	1.374%
23	15.898	2192	2195	2196	rBV3	12965	14158	0.57%	0.040%
24	15.916	2196	2198	2204	rVB7	16628	21269	0.86%	0.059%
25	16.057	2218	2222	2227	rBV7	15501	25421	1.03%	0.071%
26	16.357	2269	2273	2279	rVB5	23591	35864	1.45%	0.100%
27	16.604	2313	2315	2323	rVB9	13650	21818	0.88%	0.061%
28	16.869	2356	2360	2366	rBV3	52292	82992	3.36%	0.232%
29	16.957	2366	2375	2380	rBV2	1444962	2142289	86.83%	5.989%
30	17.051	2386	2391	2402	rVB	932999	1384675	56.12%	3.871%
31	17.151	2404	2408	2412	rBV5	23102	37060	1.50%	0.104%
32	17.522	2467	2471	2474	rBV6	13611	25397	1.03%	0.071%
33	17.763	2507	2512	2518	rBV6	79806	184450	7.48%	0.516%
34	17.834	2519	2524	2529	rBV	284778	461468	18.70%	1.290%
35	17.892	2529	2534	2541	rBV	862563	1285019	52.08%	3.592%
36	18.192	2583	2585	2596	rVB6	62108	112854	4.57%	0.315%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
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37	18.486	2628	2635	2637	rBV7	59432	110875	4.49%	0.310%
38	18.516	2637	2640	2642	rBV3	85764	77936	3.16%	0.218%
39	18.745	2674	2679	2680	rBV4	55731	83168	3.37%	0.232%
40	18.816	2688	2691	2693	rBV2	67971	79169	3.21%	0.221%

41	18.886	2701	2703	2706	rVB	113607	95719	3.88%	0.268%
42	18.922	2707	2709	2711	rBV2	38771	32424	1.31%	0.091%
43	18.981	2717	2719	2722	rVB3	72748	62828	2.55%	0.176%
44	19.104	2731	2740	2748	rBV4	195636	621236	25.18%	1.737%
45	19.163	2748	2750	2751	rVV	74749	58990	2.39%	0.165%

46	19.210	2755	2758	2759	rVV2	94900	96966	3.93%	0.271%
47	19.233	2759	2762	2766	rVV	208027	317164	12.86%	0.887%
48	19.281	2768	2770	2774	rVV3	163350	228137	9.25%	0.638%
49	19.316	2774	2776	2778	rVV3	126115	132128	5.36%	0.369%
50	19.357	2778	2783	2787	rVV2	217475	408767	16.57%	1.143%

51	19.392	2787	2789	2791	rVV2	114896	124498	5.05%	0.348%
52	19.439	2791	2797	2806	rVB2	873938	1633277	66.20%	4.566%
53	19.545	2812	2815	2817	rBV4	89859	99976	4.05%	0.279%
54	19.581	2819	2821	2824	rVB3	99446	92553	3.75%	0.259%
55	19.616	2824	2827	2828	rBV2	135320	106067	4.30%	0.296%

56	19.657	2831	2834	2837	rVV4	116999	183649	7.44%	0.513%
57	19.686	2837	2839	2841	rVV2	123673	79301	3.21%	0.222%
58	19.775	2852	2854	2855	rBV	50418	29444	1.19%	0.082%
59	19.792	2855	2857	2859	rVB	64494	51006	2.07%	0.143%
60	19.845	2864	2866	2868	rBV2	53757	53443	2.17%	0.149%

61	19.939	2879	2882	2887	rBV5	86660	155847	6.32%	0.436%
62	19.992	2887	2891	2892	rBV2	191673	260733	10.57%	0.729%
63	20.057	2900	2902	2905	rBV3	85456	83286	3.38%	0.233%
64	20.210	2926	2928	2929	rBV	107245	72154	2.92%	0.202%
65	20.269	2936	2938	2939	rBV	124892	82613	3.35%	0.231%

66	20.404	2957	2961	2963	rBV3	189394	253568	10.28%	0.709%
67	20.422	2963	2964	2969	rVB5	120041	169954	6.89%	0.475%
68	20.480	2971	2974	2976	rBV4	209097	259191	10.51%	0.725%
69	20.533	2981	2983	2990	rVB	158137	182672	7.40%	0.511%
70	20.586	2990	2992	2993	rBV	96913	55245	2.24%	0.154%

71	20.698	3009	3011	3013	rBV2	132211	104795	4.25%	0.293%
72	20.728	3015	3016	3018	rBV2	124257	76609	3.11%	0.214%
73	20.757	3018	3021	3023	rBV2	245630	286006	11.59%	0.799%
74	20.845	3034	3036	3037	rBV	89313	76021	3.08%	0.213%
75	20.863	3037	3039	3046	rVB3	189983	254577	10.32%	0.712%

76	20.945	3051	3053	3054	rBV	156563	114158	4.63%	0.319%
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77	20.963	3054	3056	3058	rVV2	220395	214466	8.69%	0.600%
78	20.986	3058	3060	3064	rVV3	330527	515453	20.89%	1.441%
79	21.022	3064	3066	3067	rVV	315701	258488	10.48%	0.723%
80	21.045	3067	3070	3078	rVB3	356496	758256	30.73%	2.120%
81	21.145	3084	3087	3089	rBV3	110853	142526	5.78%	0.398%
82	21.245	3102	3104	3107	rVB	171690	175929	7.13%	0.492%
83	21.375	3120	3126	3132	rBV2	1206751	2269287	91.98%	6.344%
84	21.480	3143	3144	3146	rBV2	251034	219767	8.91%	0.614%
85	21.804	3197	3199	3201	rBV	240360	178273	7.23%	0.498%
86	21.845	3204	3206	3210	rVB4	154255	198998	8.07%	0.556%
87	22.098	3247	3249	3250	rBV	93084	49993	2.03%	0.140%
88	22.127	3251	3254	3256	rBV	214585	185602	7.52%	0.519%
89	22.192	3263	3265	3269	rVB	139792	119124	4.83%	0.333%
90	22.404	3298	3301	3302	rBV	241622	221098	8.96%	0.618%
91	23.010	3402	3404	3406	rBV2	76438	68641	2.78%	0.192%
92	23.051	3406	3411	3418	rVV	216188	502036	20.35%	1.403%
93	23.216	3433	3439	3442	rBV3	145522	249288	10.10%	0.697%
94	23.633	3507	3510	3519	rVB4	272426	732819	29.70%	2.049%
95	24.016	3573	3575	3577	rBV2	179055	159378	6.46%	0.446%
96	24.321	3623	3627	3635	rVB	431379	1000314	40.54%	2.796%
97	24.445	3646	3648	3652	rBV4	86928	119218	4.83%	0.333%
98	24.533	3656	3663	3674	rVB	850318	2467173	100.00%	6.897%
99	26.674	4025	4027	4028	rBV2	89594	78582	3.19%	0.220%
100	28.051	4259	4261	4263	rBV3	69616	55884	2.27%	0.156%

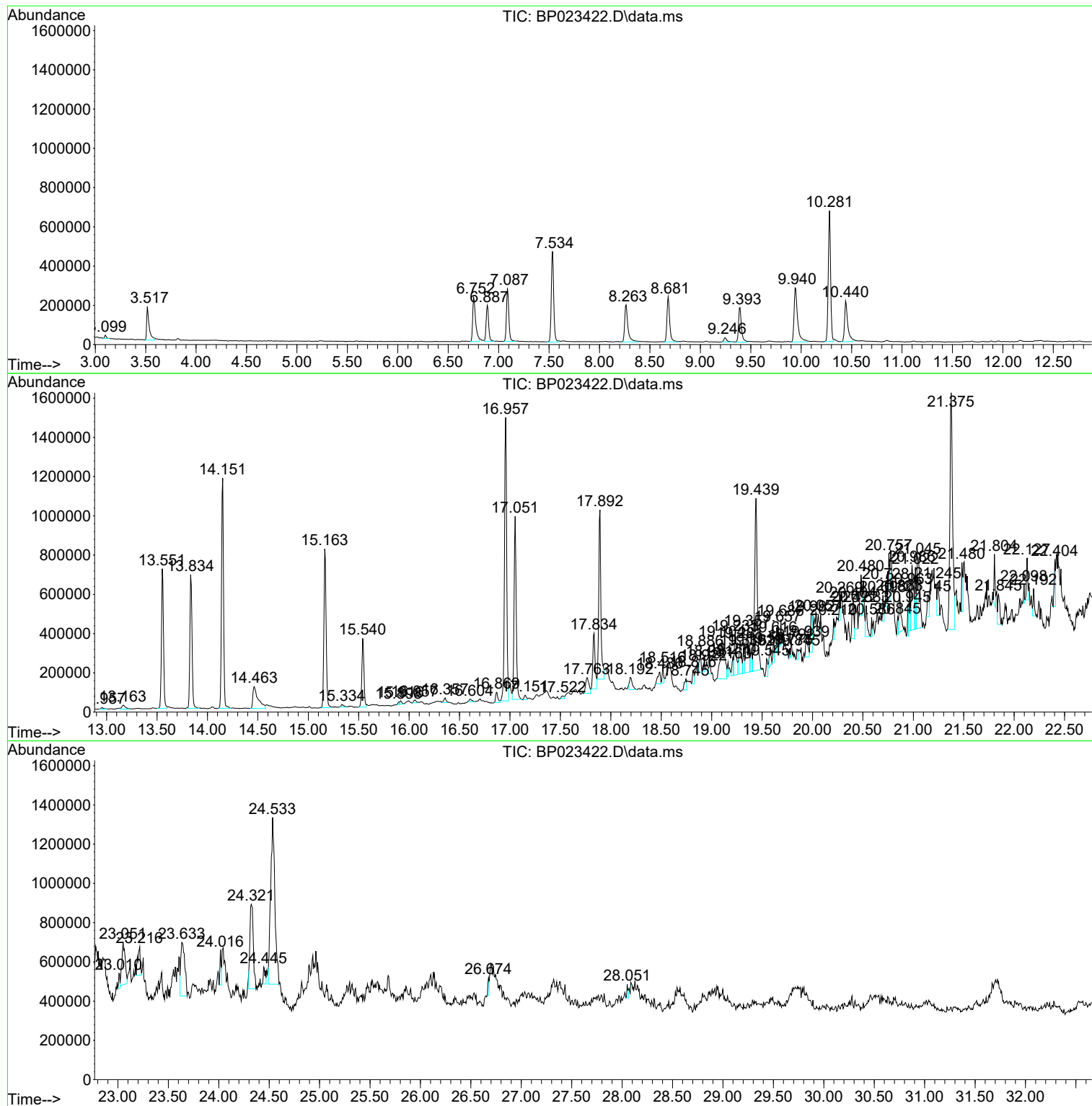
Sum of corrected areas: 35773280

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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



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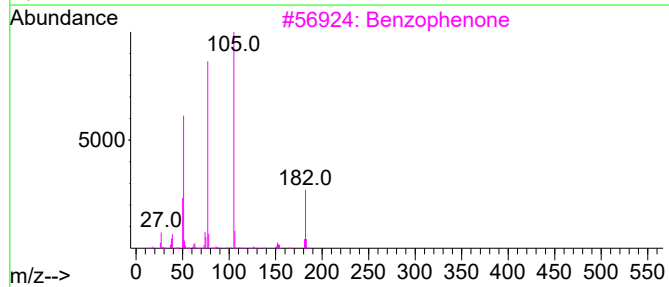
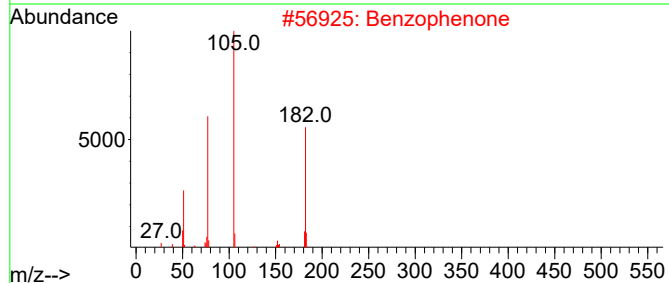
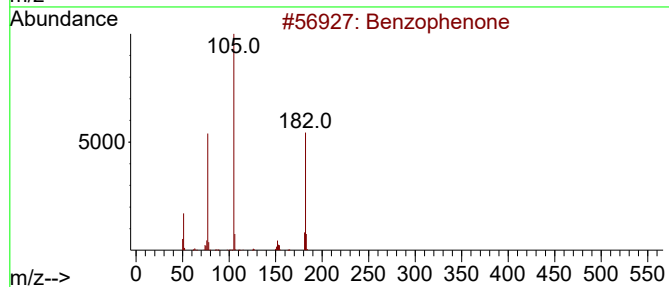
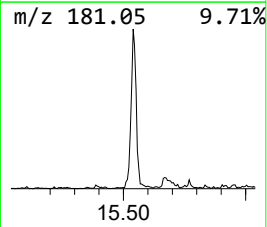
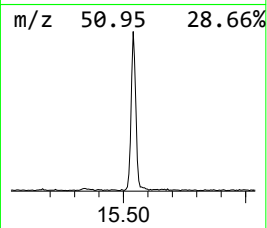
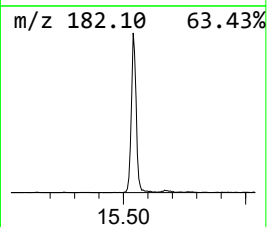
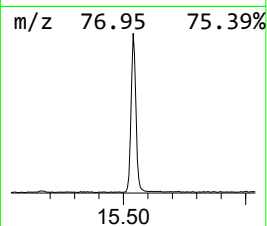
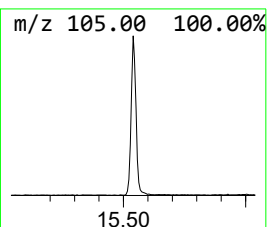
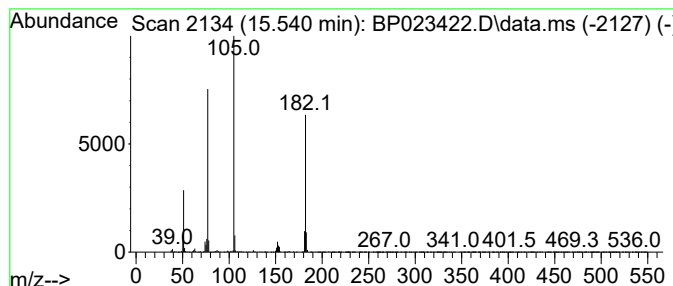
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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.
15.540	5.93 ng/ul	491426	Acenaphthene-d10	14.151

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	94
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	91



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ClientSampleId :
E28S3

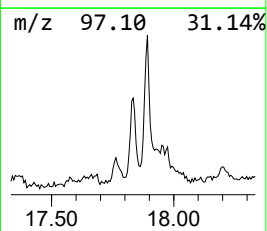
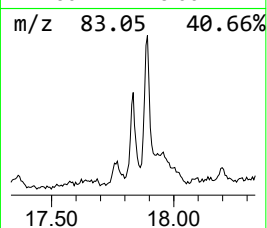
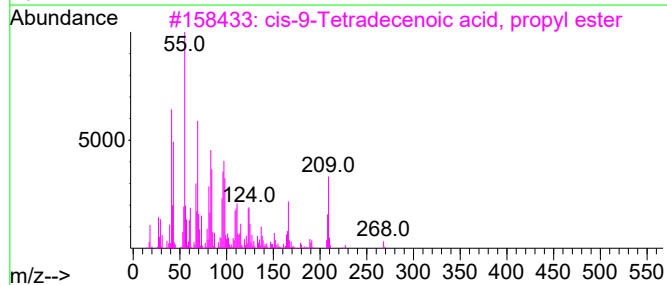
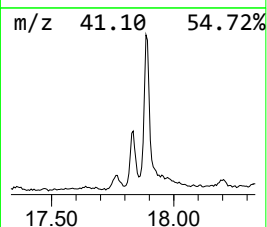
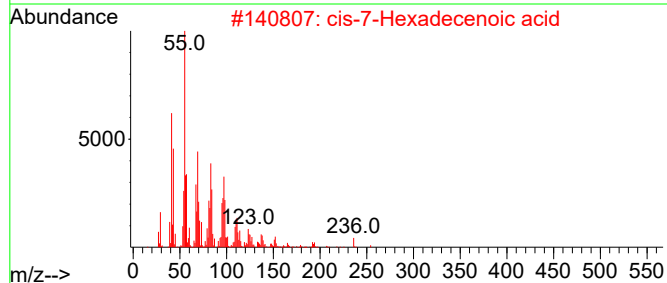
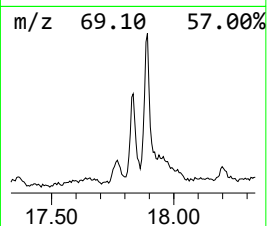
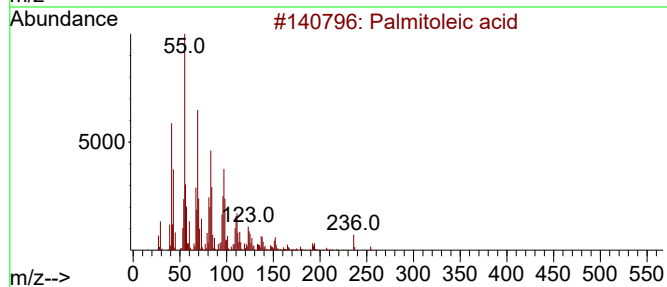
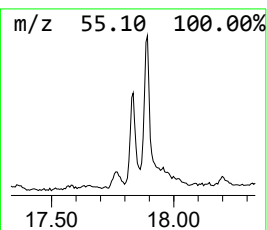
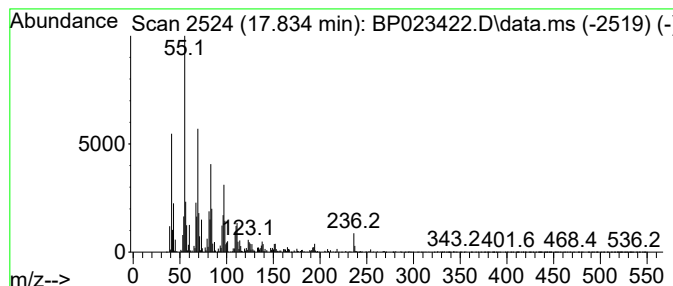
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Palmitoleic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.834	4.31 ng/ul	461468	Phenanthrene-d10	16.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Palmitoleic acid	254	C16H30O2	000373-49-9	98
2			cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	97
3			cis-9-Tetradecenoic acid, propyl...	268	C17H32O2	1000405-14-2	91
4			9-Hexadecenoic acid	254	C16H30O2	002091-29-4	91
5			trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	90



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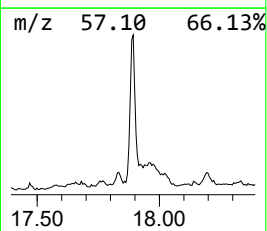
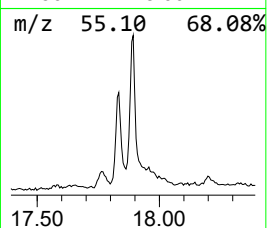
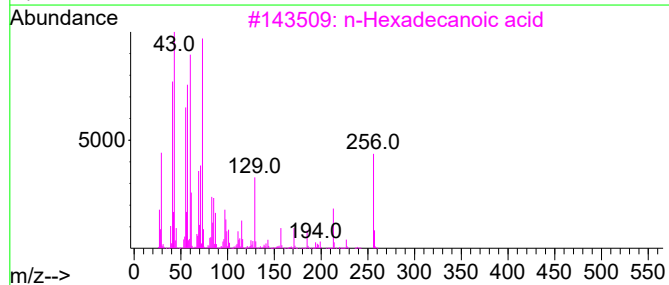
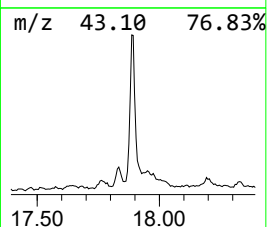
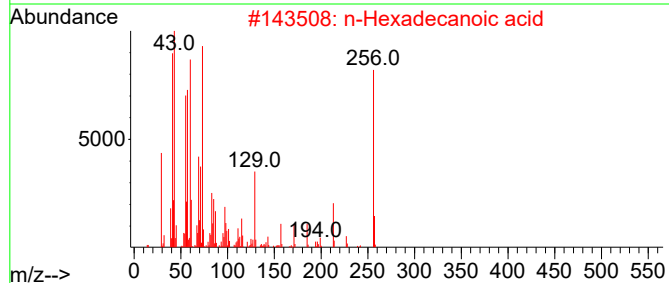
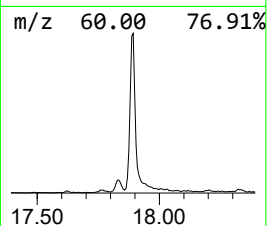
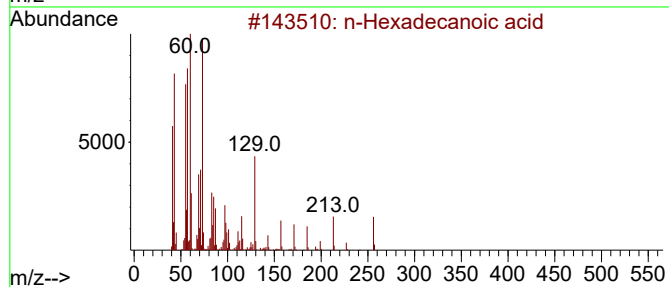
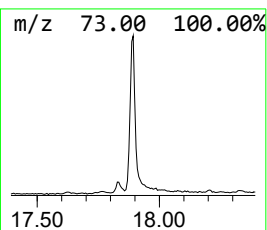
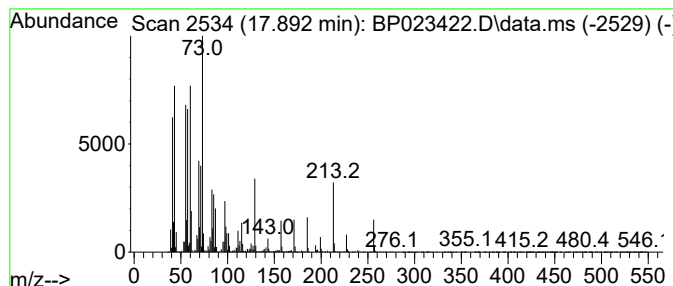
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.892	12.00 ng/ul	1285020	Phenanthrene-d10	16.957

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	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5			Tridecanoic acid	214	C13H26O2	000638-53-9	89



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Data File : BP023422.D
Acq On : 13 Dec 2024 16:11
Operator : RC/JU
Sample : P5232-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E28S3

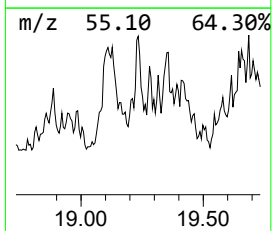
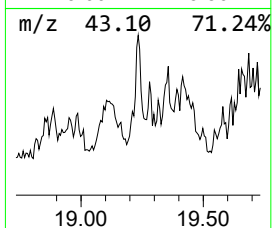
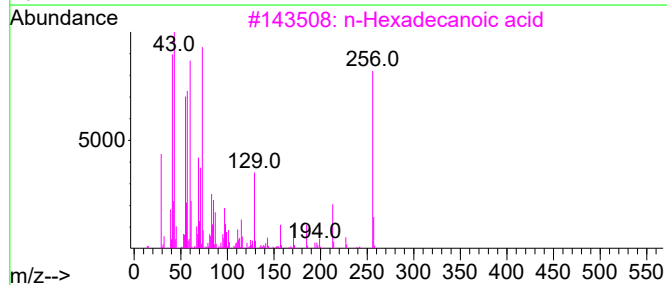
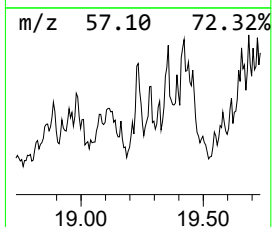
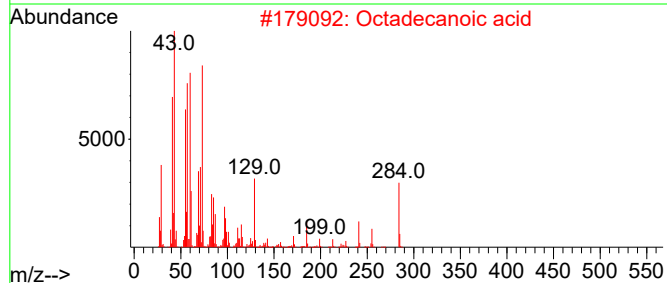
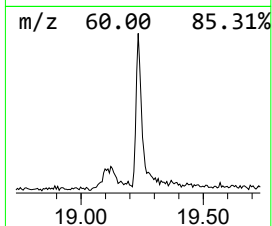
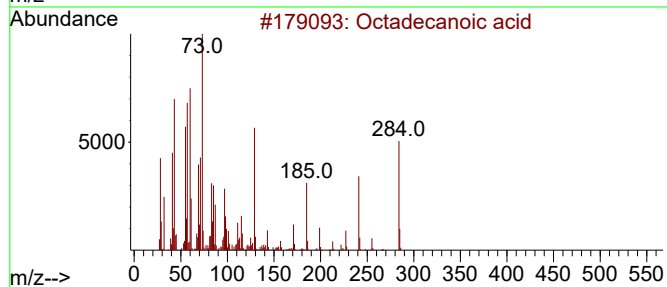
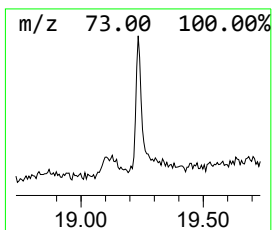
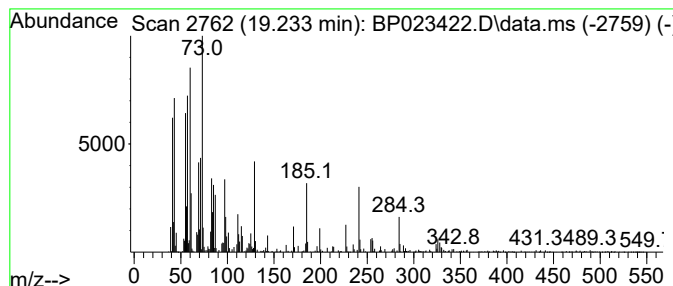
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.233	2.80 ng/ul	317164	Chrysene-d12	21.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	97
2			Octadecanoic acid	284	C18H36O2	000057-11-4	95
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
4			Octadecanoic acid	284	C18H36O2	000057-11-4	93
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121324\
Data File : BP023422.D
Acq On : 13 Dec 2024 16:11
Operator : RC/JU
Sample : P5232-09
Misc :
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Instrument :
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ClientSampleId :
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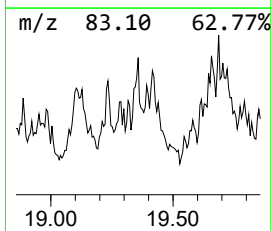
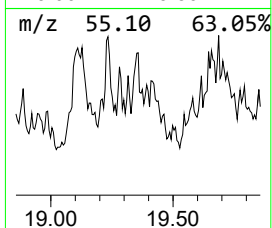
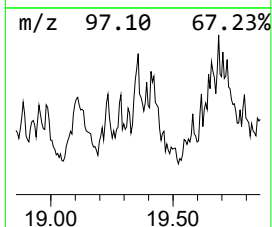
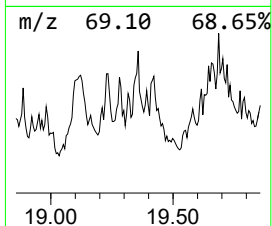
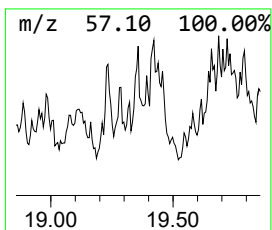
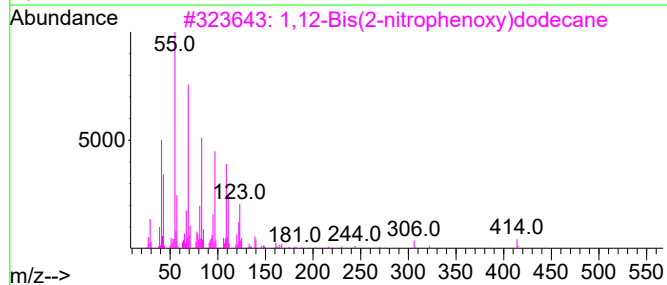
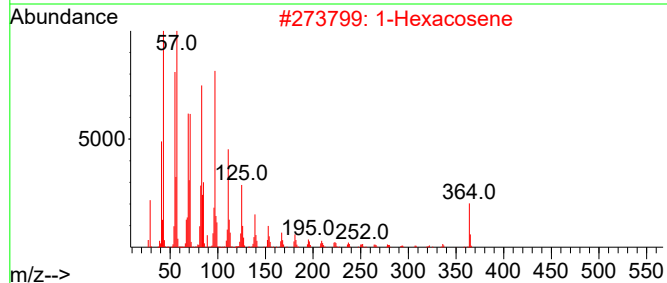
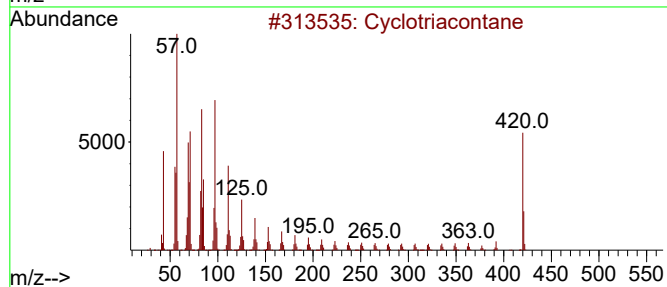
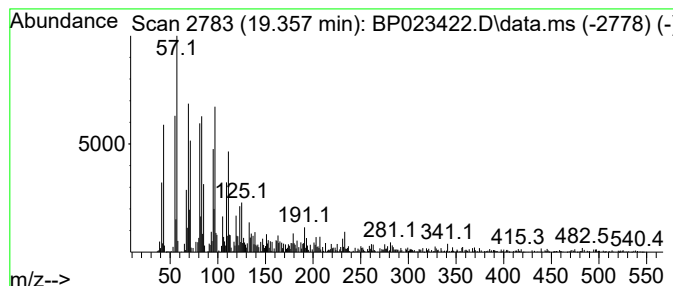
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic19.357 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.357	3.60 ng/ul	408767	Chrysene-d12	21.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotriacontane	420	C30H60	000297-35-8	91
2			1-Hexacosene	364	C26H52	018835-33-1	64
3			1,12-Bis(2-nitrophenoxy)dodecane	444	C24H32N2O6	155056-69-2	64
4			Decane, 5,6-bis(2,2-dimethylprop...	278	C20H38	073002-85-4	64
5			Cyclotetradecane, 1,7,11-trimeth...	280	C20H40	001786-12-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121324\
Data File : BP023422.D
Acq On : 13 Dec 2024 16:11
Operator : RC/JU
Sample : P5232-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

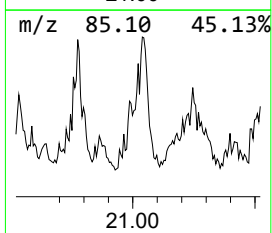
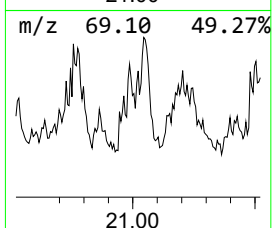
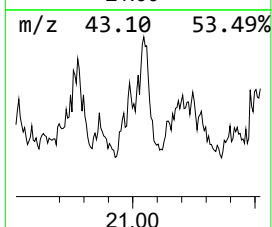
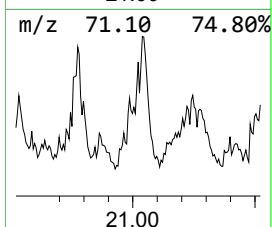
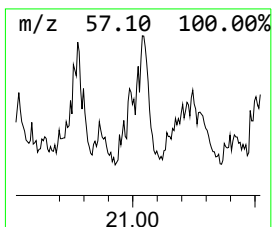
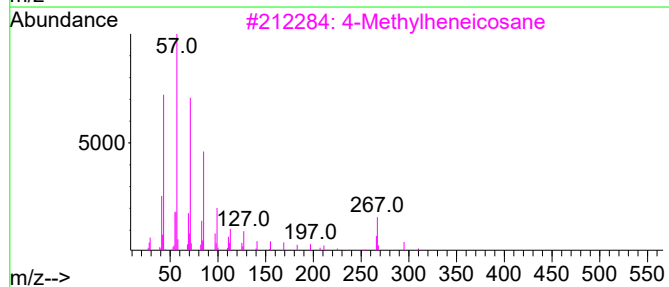
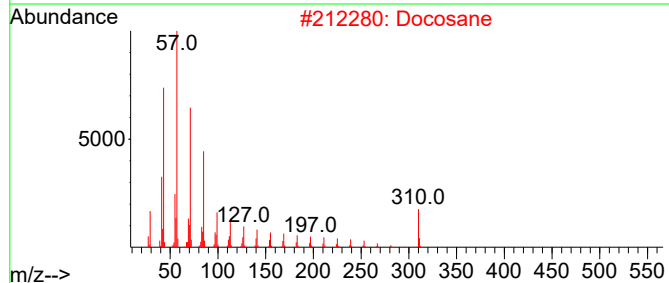
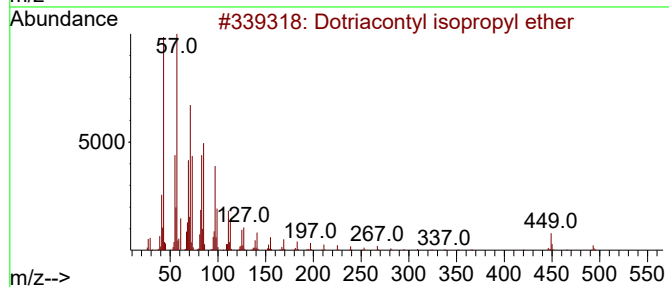
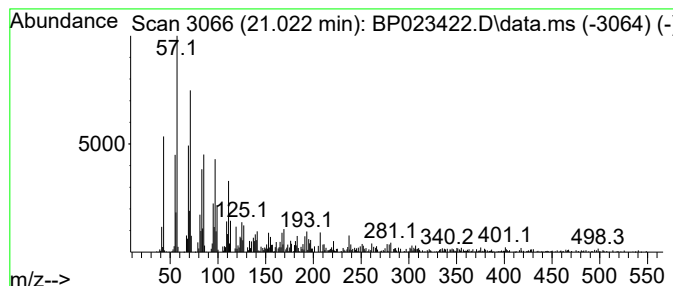
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 13 Dotriacontyl isopropyl ether Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.022	2.28 ng/ul	258488	Chrysene-d12	21.375	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dotriacontyl isopropyl ether	509	C35H72O	1000406-34-9	87
2	Docosane	310	C22H46	000629-97-0	86
3	4-Methylheneicosane	310	C22H46	025117-29-7	78
4	Tetratriacontane, 2-methyl-	493	C35H72	014167-65-8	78
5	Hexatriacontyl pentafluoropropio...	669	C39H73F5O2	1000351-89-0	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121324\
Data File : BP023422.D
Acq On : 13 Dec 2024 16:11
Operator : RC/JU
Sample : P5232-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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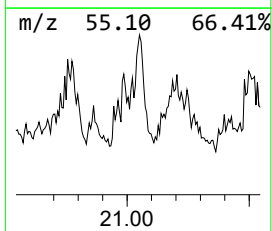
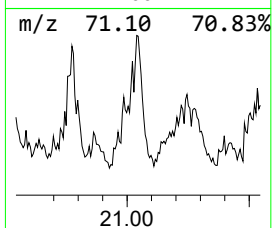
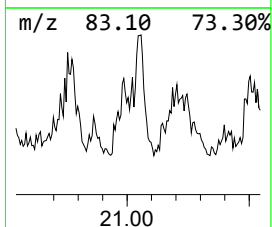
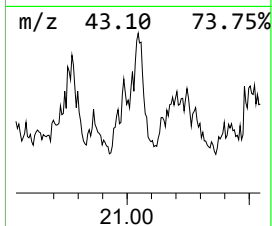
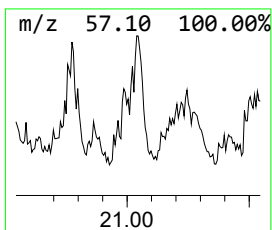
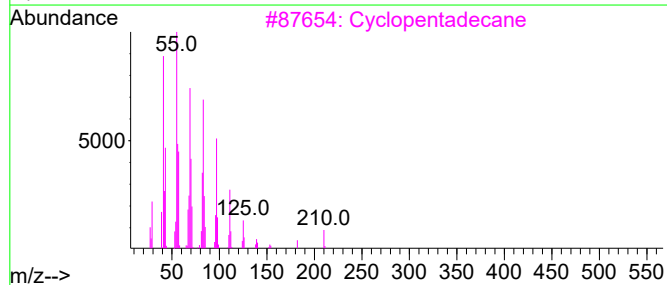
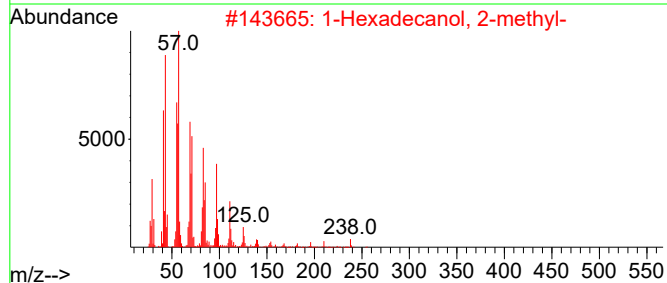
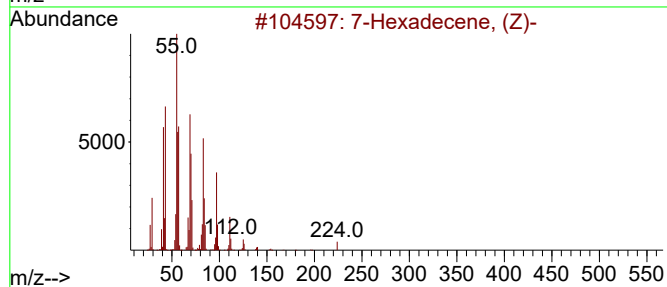
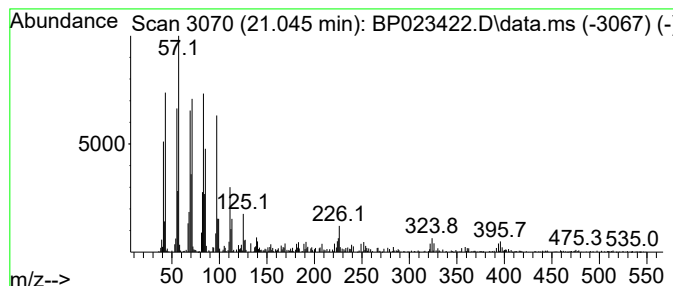
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.045	6.68 ng/ul	758256	Chrysene-d12	21.375

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		7-Hexadecene, (Z)-	224	C16H32	035507-09-6	95
2		1-Hexadecanol, 2-methyl-	256	C17H36O	002490-48-4	94
3		Cyclopentadecane	210	C15H30	000295-48-7	93
4		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
5		Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121324\
Data File : BP023422.D
Acq On : 13 Dec 2024 16:11
Operator : RC/JU
Sample : P5232-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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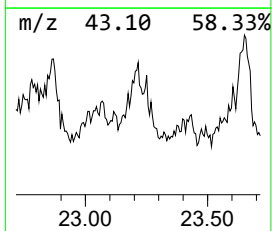
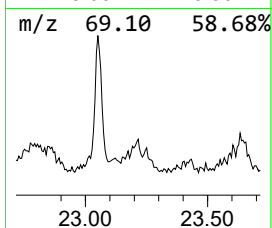
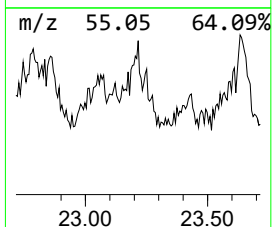
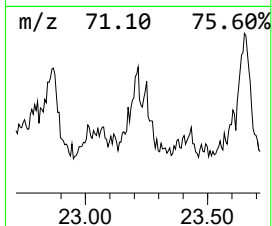
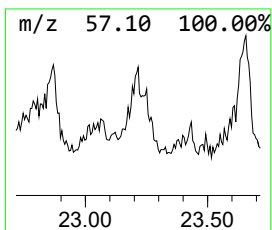
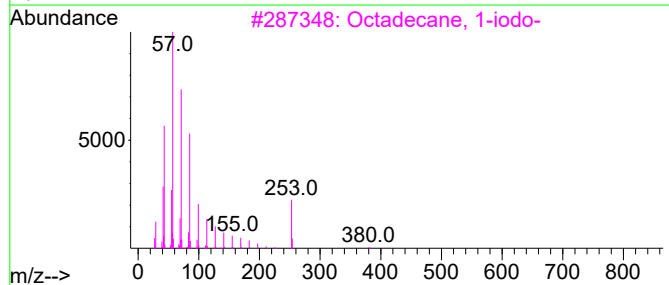
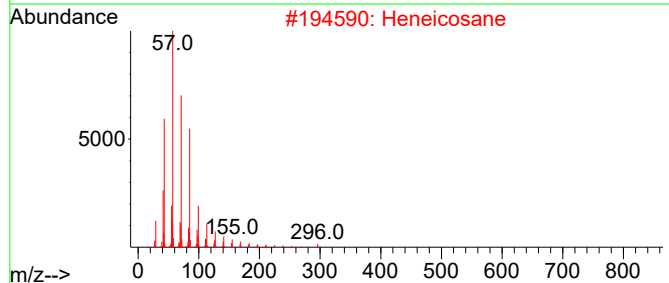
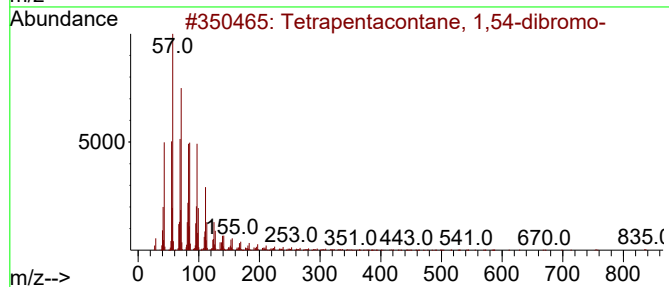
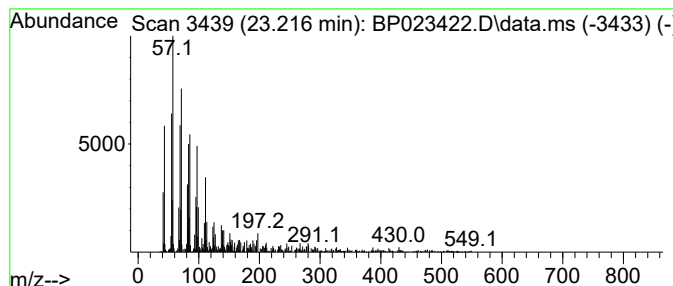
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Tetrapentacontane, 1,54-dib... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.216	2.02 ng/ul	249288	Perylene-d12	24.533

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetrapentacontane, 1,54-dibromo-	915	C54H108Br2	1000156-09-4	91
2		Heneicosane	296	C21H44	000629-94-7	90
3		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	86
4		1-Chloroeicosane	316	C20H41Cl	042217-02-7	86
5		Docosane, 1,22-dibromo-	466	C22H44Br2	034540-49-3	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121324\
Data File : BP023422.D
Acq On : 13 Dec 2024 16:11
Operator : RC/JU
Sample : P5232-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.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	15.540	5.9	ng/ul	491426	3	14.151	1656300	20.0
Palmitoleic acid	17.834	4.3	ng/ul	461468	4	16.957	2142290	20.0
n-Hexadecanoic ...	17.892	12.0	ng/ul	1285020	4	16.957	2142290	20.0
Octadecanoic acid	19.233	2.8	ng/ul	317164	5	21.375	2269290	20.0
(DEL) Alkane: C...	19.357	3.6	ng/ul	408767	5	21.375	2269290	20.0
Dotriacontyl is...	21.022	2.3	ng/ul	258488	5	21.375	2269290	20.0
(DEL) Alkane: S...	21.045	6.7	ng/ul	758256	5	21.375	2269290	20.0
2,6,10,14,18-Pe...	23.051	4.1	ng/ul	502036	6	24.533	2467170	20.0
Tetrapentaconta...	23.216	2.0	ng/ul	249288	6	24.533	2467170	20.0