

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
 Data File : BP017205.D
 Acq On : 18 Sep 2023 15:07
 Operator : MA/JU
 Sample : 04332-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AC6

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-BP091223.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017205.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.310	35	38	45	rVB	27476	36601	0.35%	0.044%
2	3.757	111	114	121	rBV	28283	48554	0.47%	0.058%
3	3.940	140	145	150	rBV	75589	107739	1.04%	0.129%
4	4.281	199	203	209	rVB2	18797	28079	0.27%	0.034%
5	4.440	227	230	238	rVB4	19943	34368	0.33%	0.041%
6	4.993	319	324	328	rBV	42681	57323	0.55%	0.069%
7	6.898	641	648	657	rVB6	14929	31964	0.31%	0.038%
8	7.098	675	682	696	rBV	364664	641080	6.18%	0.767%
9	7.287	707	714	725	rBV	441208	666619	6.43%	0.797%
10	7.463	738	744	752	rBV	465041	731837	7.05%	0.875%
11	7.540	752	757	764	rVV	34054	55972	0.54%	0.067%
12	7.728	784	789	796	rVB3	30444	60592	0.58%	0.072%
13	7.945	819	826	834	rBV	829731	1273161	12.27%	1.523%
14	8.057	837	845	853	rBV	235868	402743	3.88%	0.482%
15	8.204	864	870	877	rBV	47680	93075	0.90%	0.111%
16	8.481	913	917	922	rBV3	18532	30661	0.30%	0.037%
17	8.657	938	947	959	rBV	260186	453030	4.37%	0.542%
18	8.798	962	971	977	rVV	199295	331486	3.20%	0.397%
19	9.145	1020	1030	1039	rBV	484640	817419	7.88%	0.978%
20	9.251	1040	1048	1059	rVV4	56641	155838	1.50%	0.186%
21	9.351	1061	1065	1075	rVB7	8469	20199	0.19%	0.024%
22	9.857	1145	1151	1161	rBV	404324	646419	6.23%	0.773%
23	10.022	1168	1179	1202	rVV2	201720	532338	5.13%	0.637%
24	10.222	1208	1213	1219	rVB4	16963	28393	0.27%	0.034%
25	10.381	1233	1240	1252	rBV	572551	932097	8.98%	1.115%
26	10.545	1261	1268	1280	rBV	25896	66262	0.64%	0.079%
27	10.769	1298	1306	1321	rBV	1140204	1888896	18.21%	2.260%
28	10.934	1326	1334	1341	rBV	65798	125386	1.21%	0.150%
29	11.316	1392	1399	1416	rBV5	19027	67243	0.65%	0.080%
30	11.528	1429	1435	1440	rVB4	21587	35835	0.35%	0.043%
31	11.592	1440	1446	1457	rVB	13710	33160	0.32%	0.040%
32	11.786	1475	1479	1486	rVV7	8401	20067	0.19%	0.024%
33	11.881	1490	1495	1504	rVB	35352	57939	0.56%	0.069%
34	12.351	1568	1575	1582	rVB2	13753	23044	0.22%	0.028%
35	12.451	1582	1592	1601	rBV10	8562	28759	0.28%	0.034%
36	12.716	1630	1637	1645	rBV	86081	135836	1.31%	0.162%

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Peak Location: TOP

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
 Title : SVOA CALIBRATION

37	12.845	1653	1659	1666	rBV2	36343	56980	0.55%	0.068%
38	13.451	1754	1762	1768	rBV	112336	192085	1.85%	0.230%
39	13.551	1773	1779	1789	rBV	108861	178930	1.72%	0.214%
40	13.639	1789	1794	1799	rVB	36860	55946	0.54%	0.067%
41	13.786	1811	1819	1826	rBV3	22507	64498	0.62%	0.077%
42	14.022	1852	1859	1865	rBV	1082063	1464608	14.12%	1.752%
43	14.304	1900	1907	1922	rBV	1230073	1761128	16.98%	2.107%
44	14.475	1932	1936	1942	rVB5	15071	24823	0.24%	0.030%
45	14.604	1951	1958	1967	rBV	1711399	2460880	23.72%	2.944%
46	14.827	1988	1996	2002	rBV	163855	291271	2.81%	0.348%
47	14.998	2019	2025	2030	rBV7	17211	36943	0.36%	0.044%
48	15.422	2093	2097	2103	rVB2	14189	22019	0.21%	0.026%
49	15.604	2117	2128	2139	rBV	1579088	2307776	22.24%	2.761%
50	15.727	2143	2149	2156	rBV	348269	482732	4.65%	0.577%
51	16.427	2264	2268	2274	rBV5	29332	47452	0.46%	0.057%
52	16.727	2315	2319	2322	rBV3	25796	37379	0.36%	0.045%
53	17.092	2369	2381	2386	rBV9	42444	89628	0.86%	0.107%
54	17.157	2389	2392	2399	rVB9	9345	20964	0.20%	0.025%
55	17.227	2399	2404	2411	rBV	77454	135229	1.30%	0.162%
56	17.292	2411	2415	2422	rVB2	76930	96542	0.93%	0.115%
57	17.374	2423	2429	2435	rBV	2189329	3008992	29.00%	3.599%
58	17.474	2440	2446	2456	rBV	1600148	2256876	21.75%	2.700%
59	17.804	2499	2502	2508	rVB6	14219	21609	0.21%	0.026%
60	17.863	2509	2512	2521	rVB5	15451	26692	0.26%	0.032%
61	17.963	2526	2529	2532	rBV4	17590	21220	0.20%	0.025%
62	17.998	2532	2535	2539	rVB	41360	47960	0.46%	0.057%
63	18.110	2547	2554	2560	rBV3	100341	196520	1.89%	0.235%
64	18.168	2560	2564	2568	rVV	96808	126732	1.22%	0.152%
65	18.221	2568	2573	2586	rBV	989367	1431746	13.80%	1.713%
66	18.510	2615	2622	2633	rBV8	51044	115835	1.12%	0.139%
67	18.815	2669	2674	2678	rVB2	48656	77699	0.75%	0.093%
68	19.039	2708	2712	2715	rBV	40125	55652	0.54%	0.067%
69	19.098	2718	2722	2726	rVV	138663	184525	1.78%	0.221%
70	19.133	2726	2728	2735	rVB4	52046	70618	0.68%	0.084%
71	19.286	2745	2754	2757	rBV6	94665	248636	2.40%	0.297%
72	19.321	2757	2760	2762	rVV	82693	116518	1.12%	0.139%
73	19.368	2762	2768	2776	rVV4	223385	496128	4.78%	0.593%
74	19.457	2779	2783	2799	rVB	305573	453467	4.37%	0.542%
75	19.657	2814	2817	2821	rVB	62048	65976	0.64%	0.079%
76	19.715	2822	2827	2839	rVB	2283156	2938297	28.32%	3.515%

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77	20.180	2900	2906	2918	rBV	1075793	1743172	16.80%	2.085%
78	20.668	2983	2989	2993	rBV2	175150	275096	2.65%	0.329%
79	20.768	3004	3006	3014	rVB5	68584	81911	0.79%	0.098%
80	20.857	3018	3021	3024	rVB2	70041	76409	0.74%	0.091%
81	21.127	3062	3067	3069	rBV	1672520	1977436	19.06%	2.365%
82	21.292	3092	3095	3102	rVB3	251892	410464	3.96%	0.491%
83	21.392	3109	3112	3116	rBV	196174	233236	2.25%	0.279%
84	21.445	3118	3121	3122	rVV2	176698	192562	1.86%	0.230%
85	21.480	3122	3127	3132	rVB	2268291	2884814	27.81%	3.451%
86	21.568	3137	3142	3146	rVV3	253358	413703	3.99%	0.495%
87	21.774	3172	3177	3185	rBV2	2347584	3151316	30.38%	3.770%
88	22.045	3218	3223	3227	rBV	2717850	3399127	32.76%	4.066%
89	22.092	3227	3231	3248	rVB2	696185	1967832	18.97%	2.354%
90	22.280	3260	3263	3268	rVB	343073	418683	4.04%	0.501%
91	22.839	3352	3358	3368	rBV	4184271	6374587	61.44%	7.625%
92	23.139	3403	3409	3417	rVB	1610641	2716822	26.19%	3.250%
93	23.233	3418	3425	3436	rBV2	1021954	3542530	34.15%	4.238%
94	23.462	3460	3464	3474	rVB	347915	603054	5.81%	0.721%
95	23.833	3522	3527	3535	rVB2	1007620	1918158	18.49%	2.295%
96	24.003	3550	3556	3568	rVB	1238750	2580734	24.88%	3.087%
97	24.192	3581	3588	3607	rVB	5184424	10374613	100.00%	12.410%
98	24.562	3645	3651	3661	rVB	583393	1269712	12.24%	1.519%
99	24.721	3671	3678	3690	rBV3	693121	2890208	27.86%	3.457%
100	26.033	3895	3901	3915	rVB2	611988	1610142	15.52%	1.926%

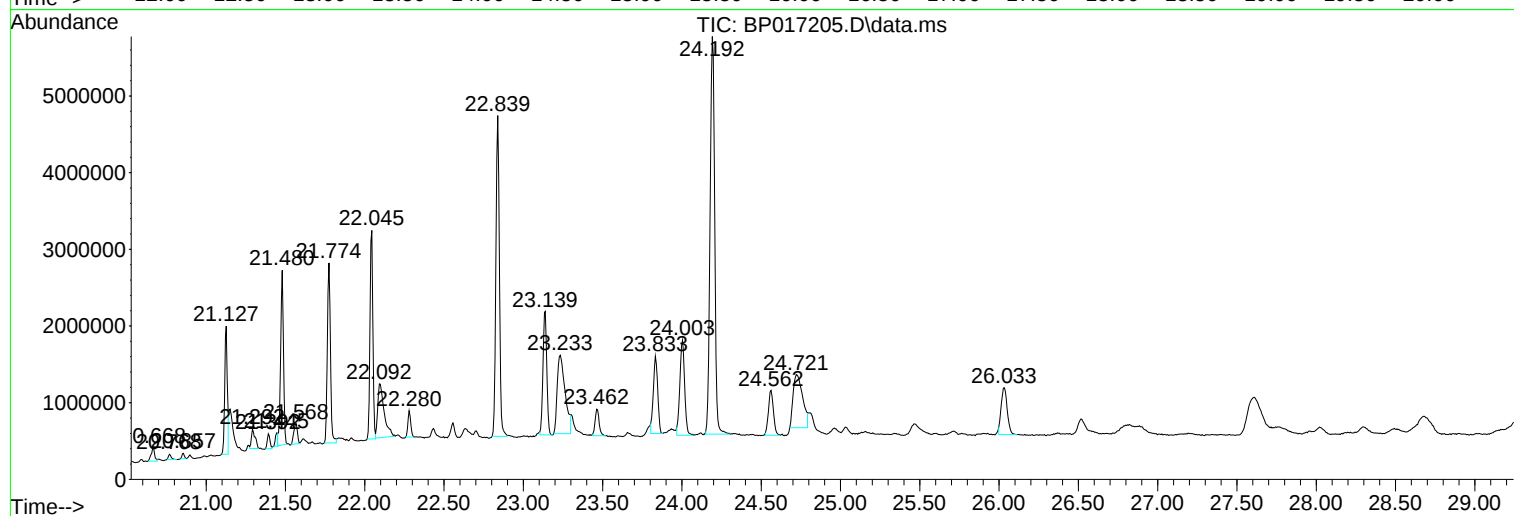
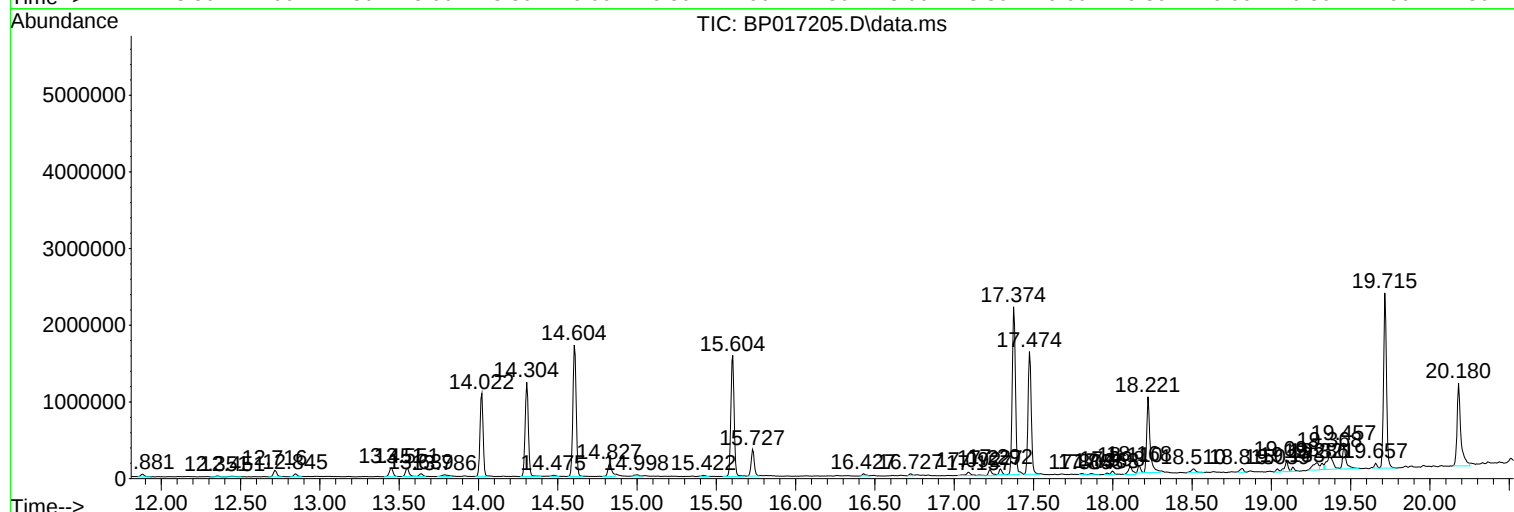
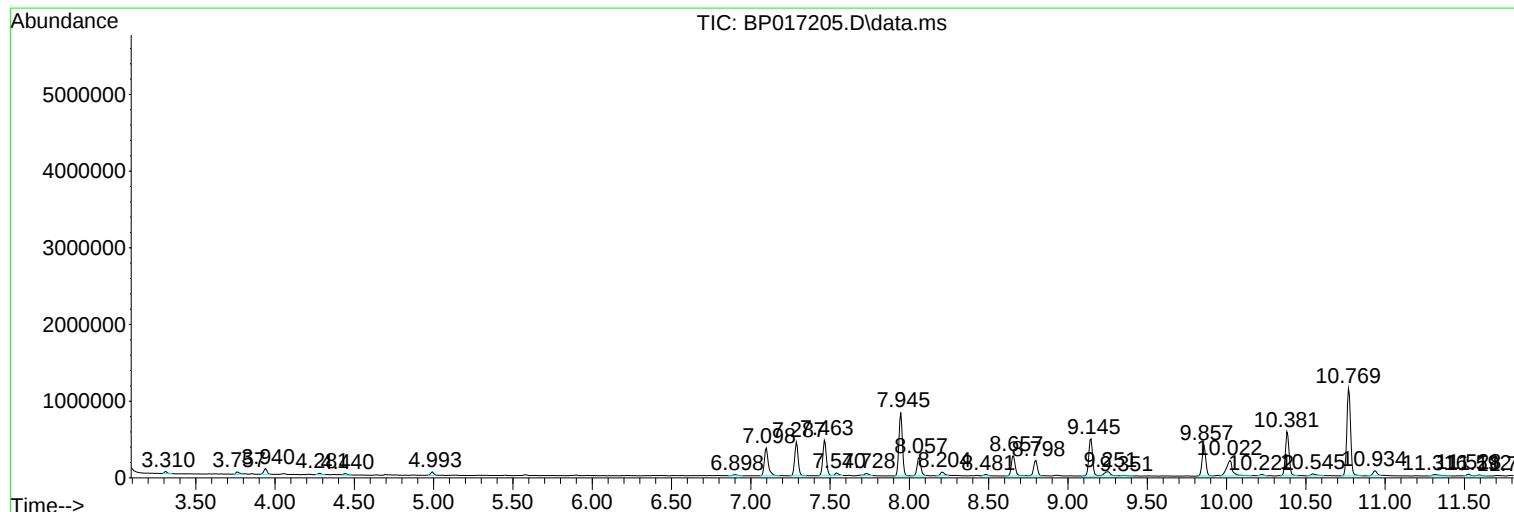
Sum of corrected areas: 83595846

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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



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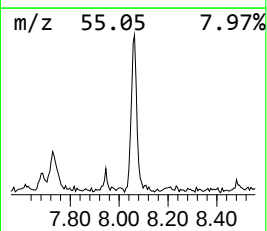
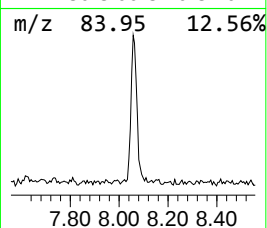
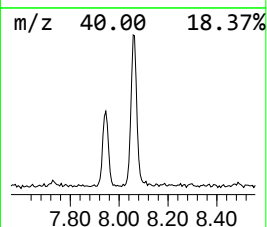
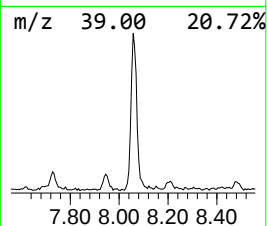
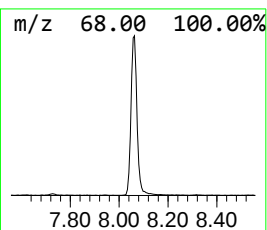
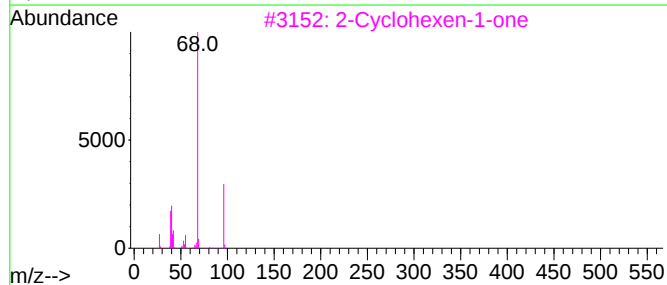
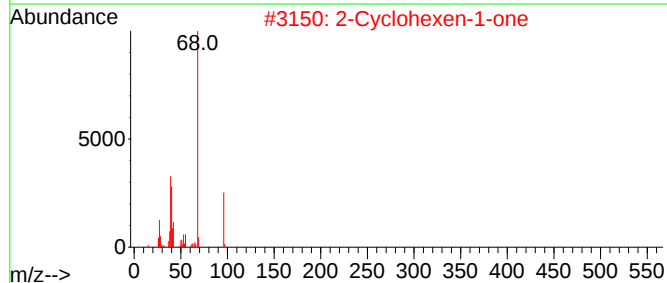
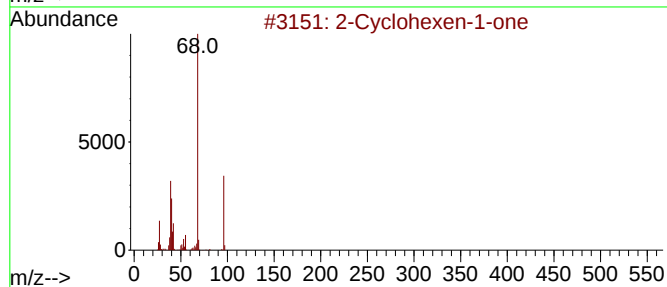
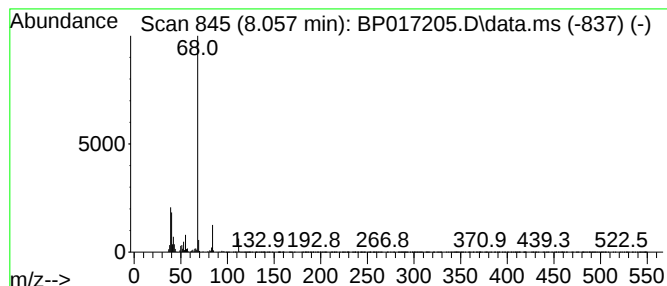
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Cyclohexen-1-one Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.057	6.33 ng/ul	402743	1,4-Dichlorobenzene-d4	7.945

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	80
2			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	72
3			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	64
4			Proline, 3,4-didehydro-	113	C5H7NO2	003395-35-5	59
5			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	50



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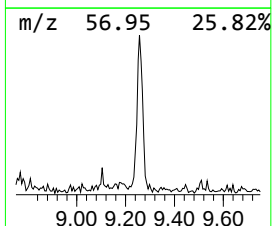
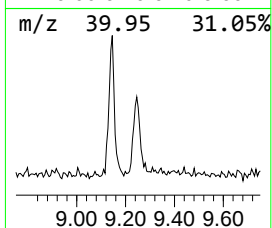
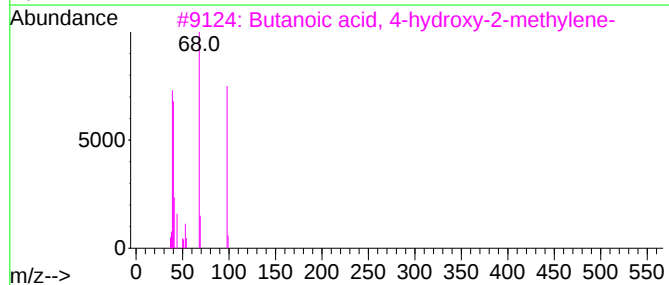
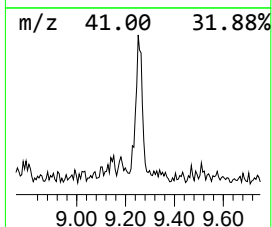
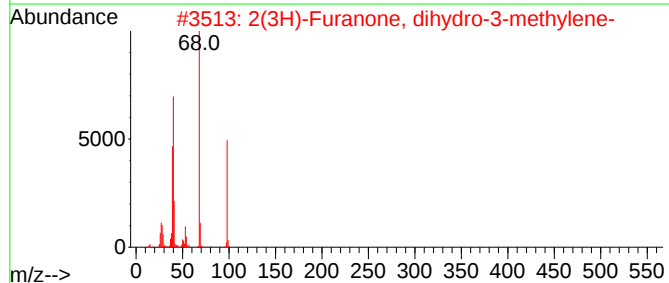
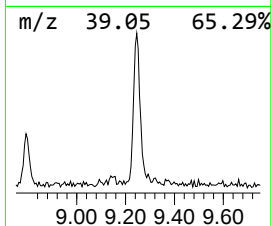
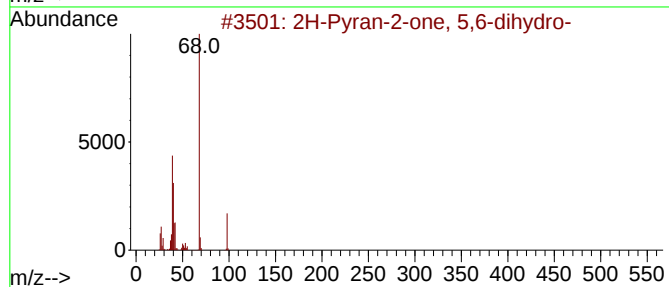
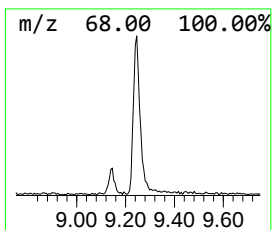
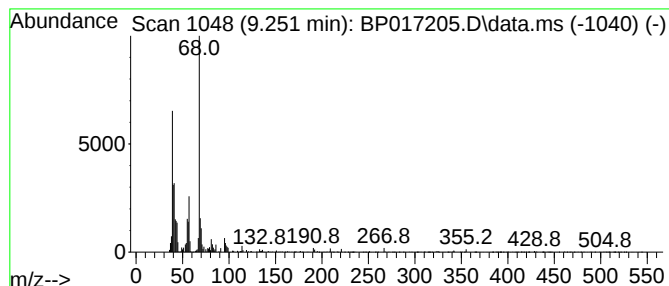
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2H-Pyran-2-one, 5,6-dihydro- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.251	2.45 ng/ul	155838	1,4-Dichlorobenzene-d4	7.945

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2H-Pyran-2-one, 5,6-dihydro-	98	C5H6O2	003393-45-1	64
2			2(3H)-Furanone, dihydro-3-methyl...	98	C5H6O2	000547-65-9	50
3			Butanoic acid, 4-hydroxy-2-methy...	116	C5H8O3	024923-76-0	50
4			1,5-Cyclooctadien-4-one	122	C8H10O	001460-21-5	50
5			(+)-3,4-Dehydroproline amide	112	C5H8N2O	1000132-73-7	45



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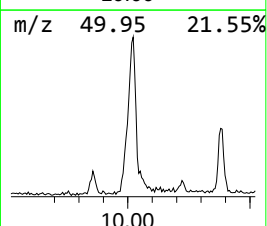
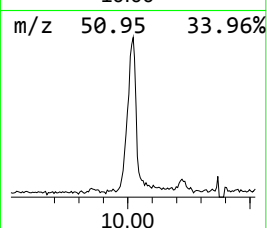
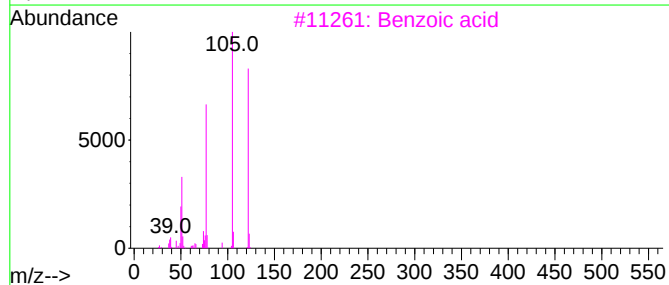
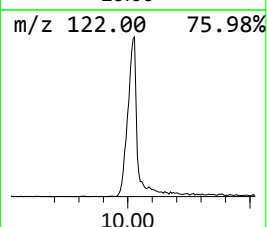
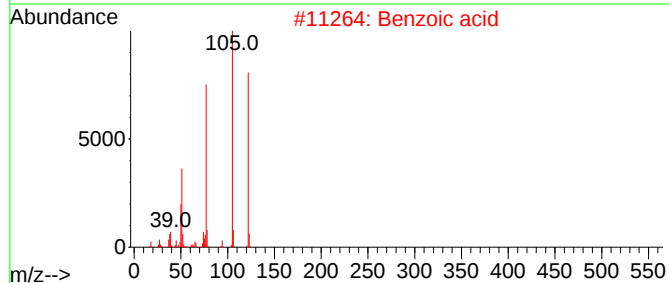
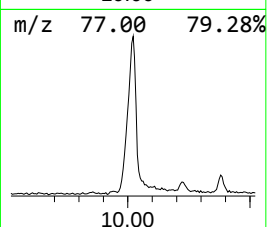
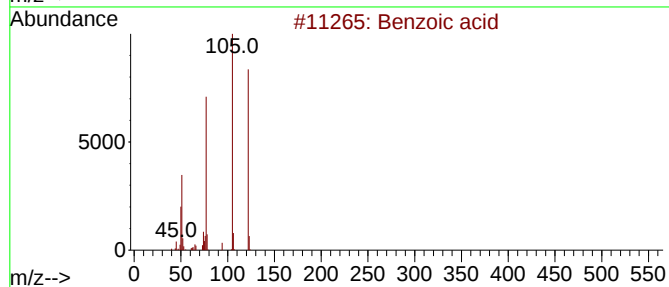
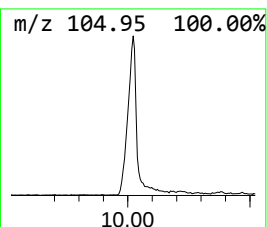
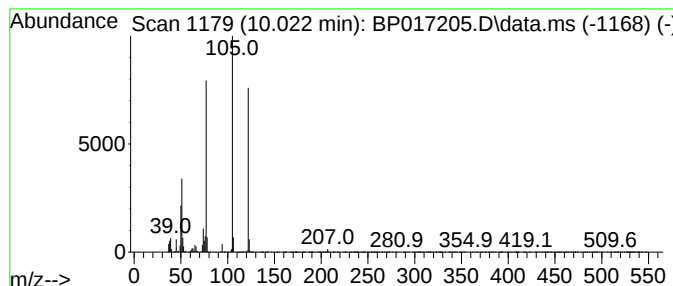
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.022	5.64 ng/ul	532338	Naphthalene-d8	10.769	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid	122	C7H6O2	000065-85-0	96
2	Benzoic acid	122	C7H6O2	000065-85-0	94
3	Benzoic acid	122	C7H6O2	000065-85-0	94
4	Cyclobutane-1,1-dicarboxamide, N...	382	C20H18N2O6	1000253-25-3	91
5	Benzoic acid, silver(1+) salt	228	C7H5AgO2	000532-31-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
Operator : MA/JU
Sample : 04332-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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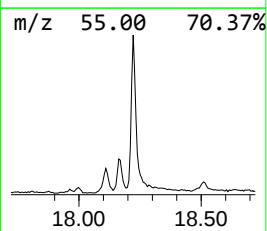
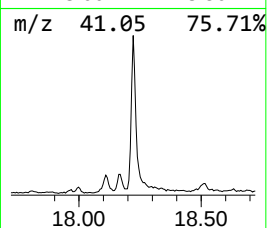
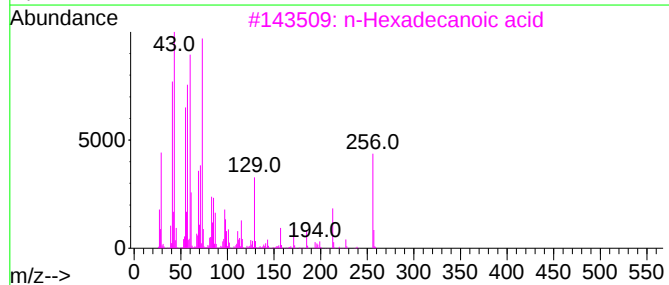
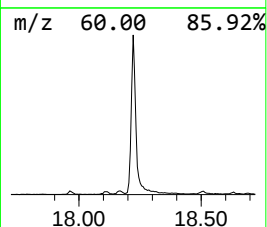
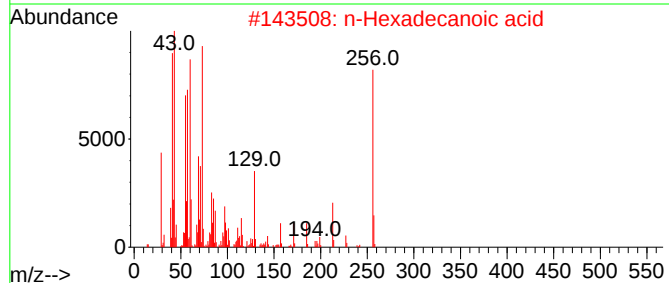
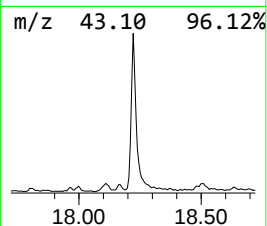
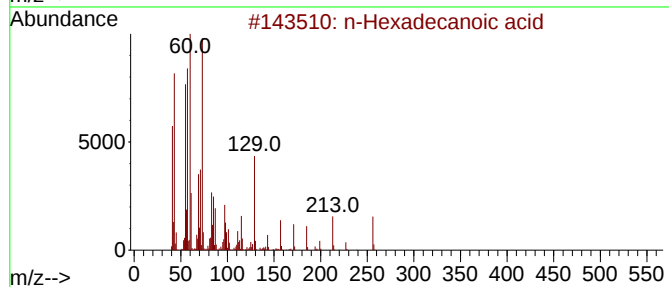
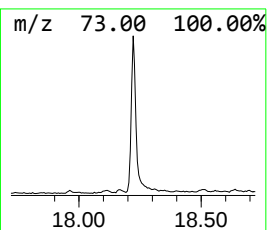
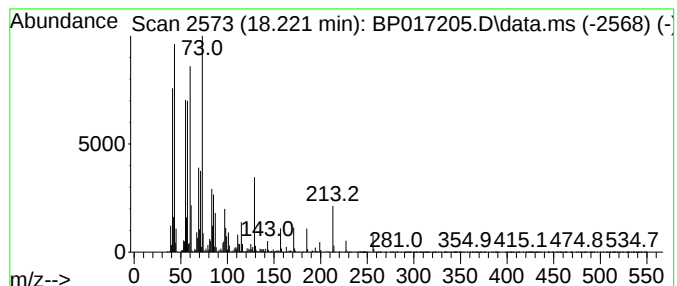
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.221	9.52 ng/ul	1431750	Phenanthrene-d10	17.374

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	96
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
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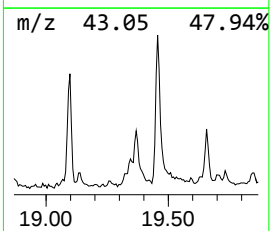
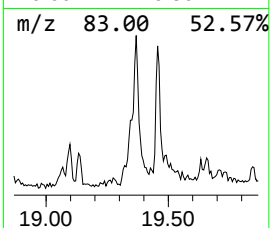
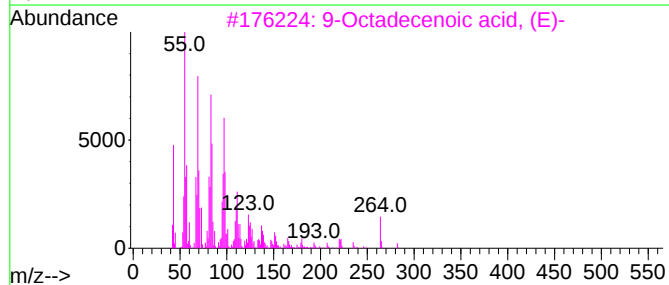
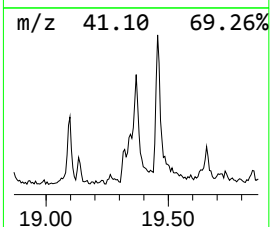
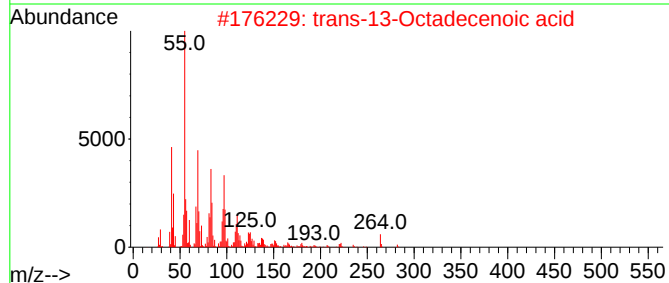
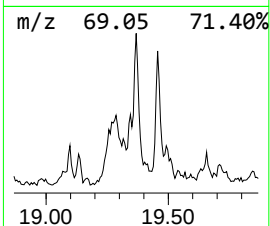
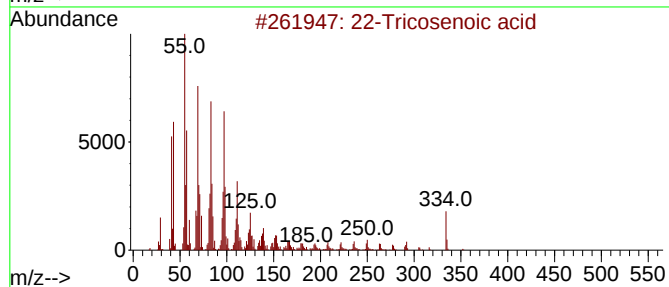
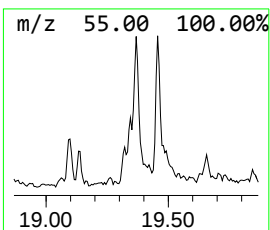
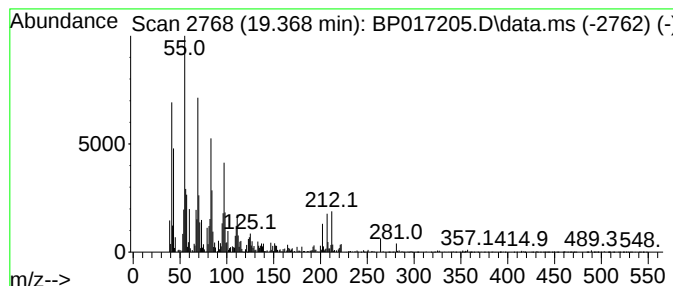
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 22-Tricosenoic acid Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.368	3.30 ng/ul	496128	Phenanthrene-d10		17.374	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	22-Tricosenoic acid		352	C23H44O2	065119-95-1	86
2	trans-13-Octadecenoic acid		282	C18H34O2	000693-71-0	83
3	9-Octadecenoic acid, (E)-		282	C18H34O2	000112-79-8	72
4	cis-13-Octadecenoic acid		282	C18H34O2	013126-39-1	70
5	Erucic acid		338	C22H42O2	000112-86-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
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Sample : 04332-06
Misc :
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Instrument :
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ClientSampleId :
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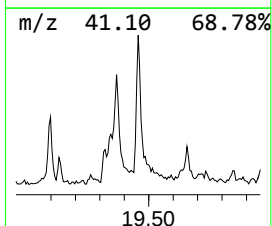
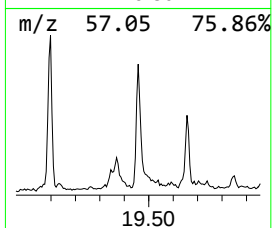
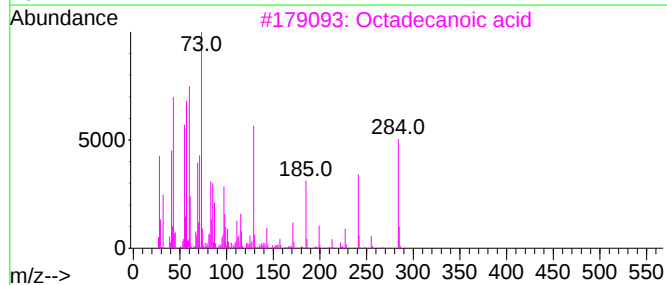
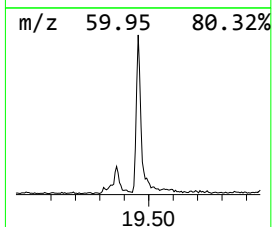
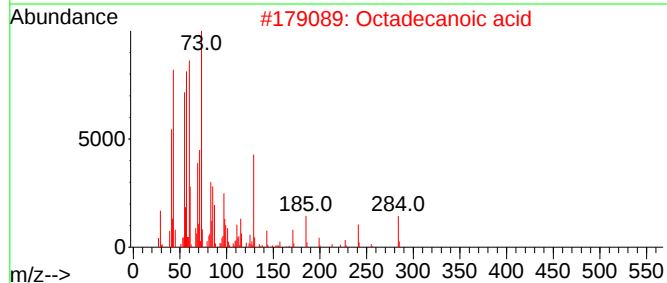
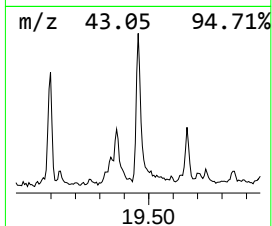
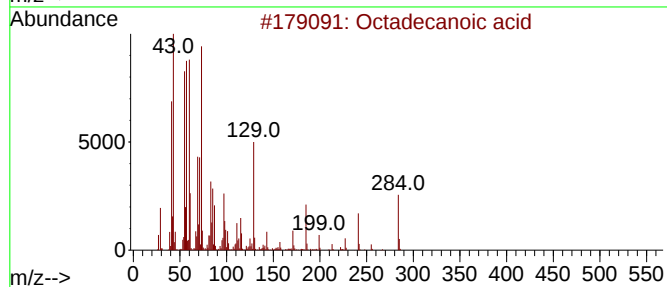
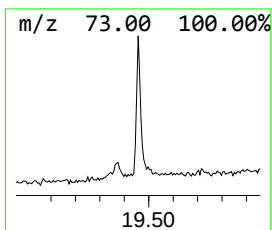
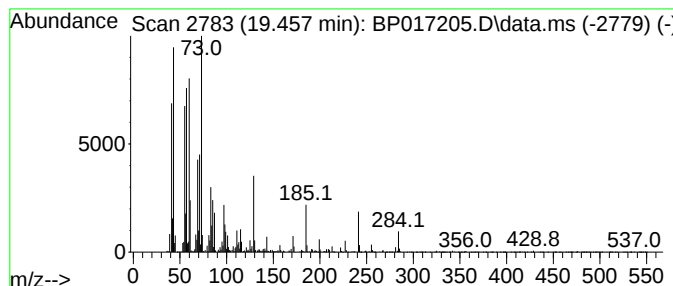
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.457	3.14 ng/ul	453467	Chrysene-d12	21.480

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	98
3			Octadecanoic acid	284	C18H36O2	000057-11-4	92
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
Operator : MA/JU
Sample : 04332-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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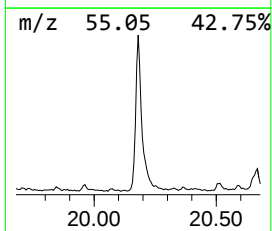
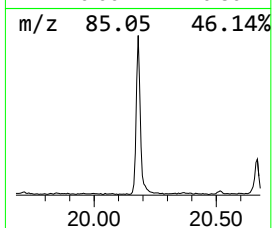
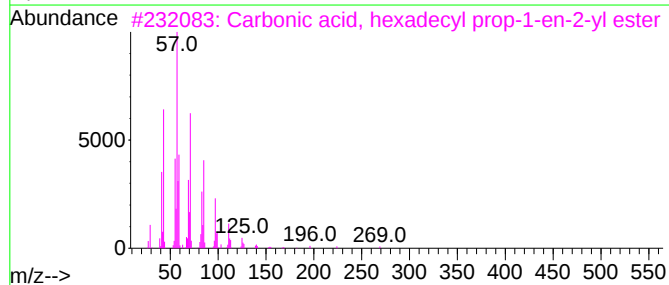
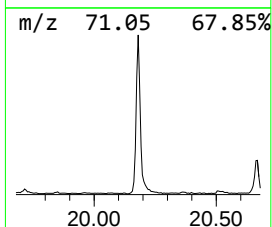
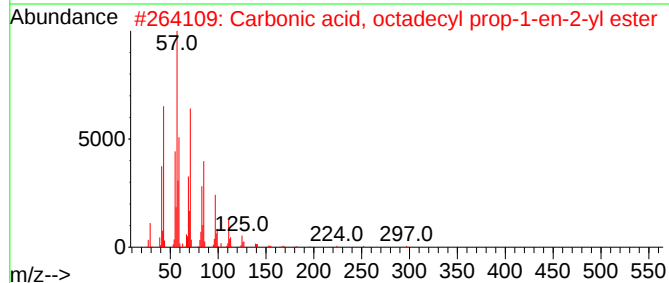
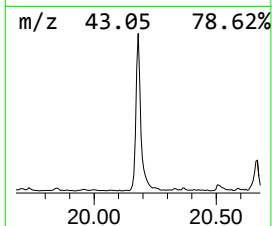
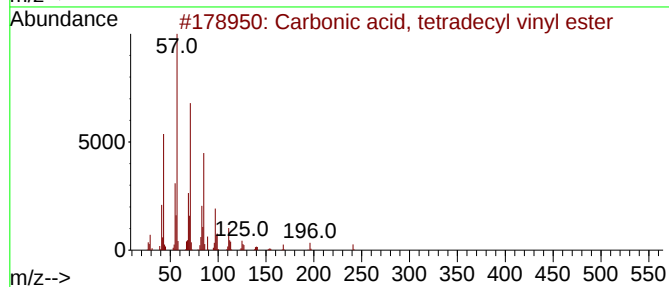
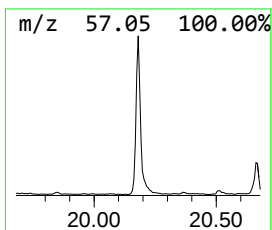
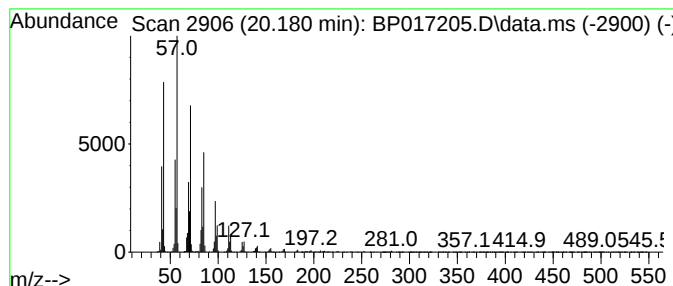
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Carbonic acid, tetradecyl v... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.180	12.09 ng/ul	1743170	Chrysene-d12	21.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, tetradecyl vinyl ...	284	C17H32O3	1000382-54-5	91
2			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	91
3			Carbonic acid, hexadecyl prop-1-...	326	C20H38O3	1000382-90-3	91
4			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	91
5			Carbonic acid, tridecyl vinyl ester	270	C16H30O3	1000382-54-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
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Sample : 04332-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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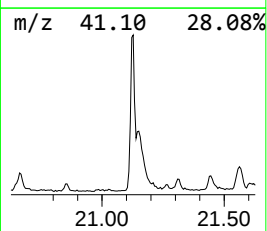
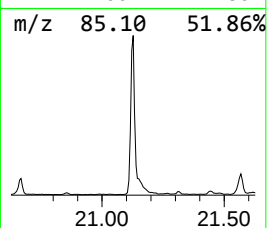
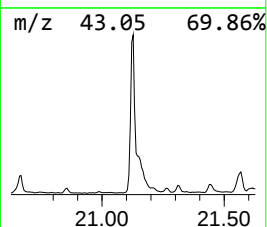
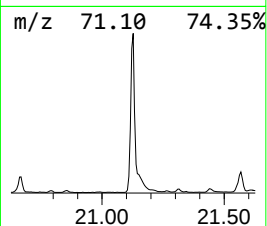
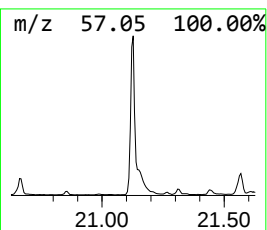
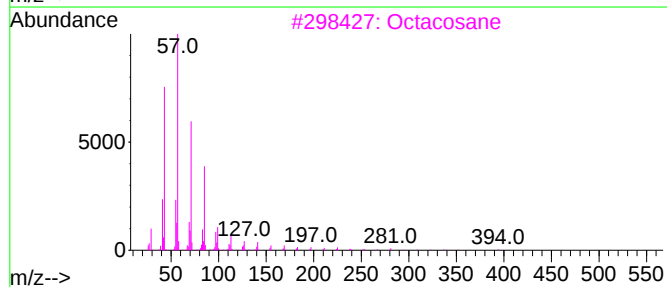
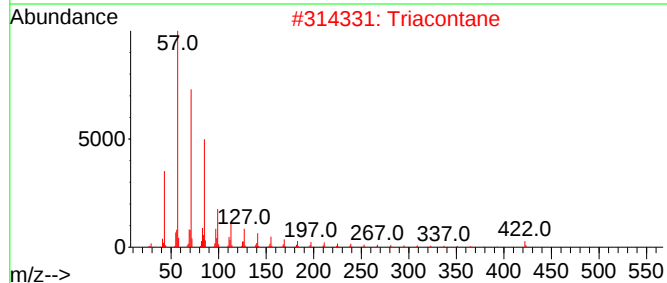
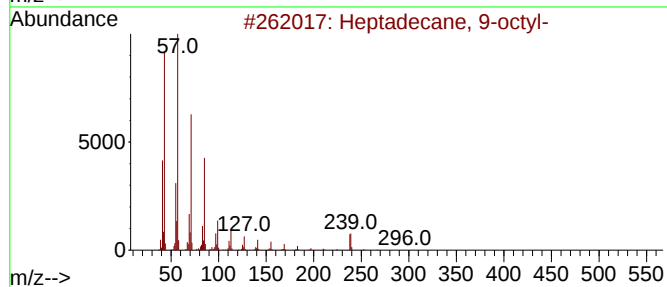
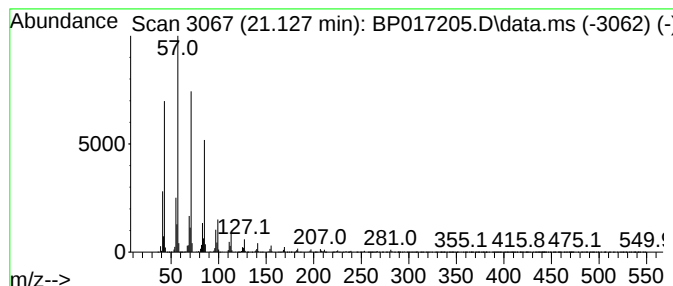
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.127	13.71 ng/ul	1977440	Chrysene-d12	21.480

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	96
2		Triacontane	422	C30H62	000638-68-6	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
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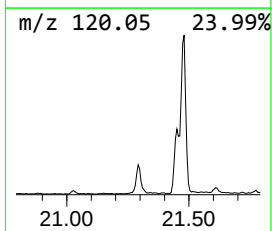
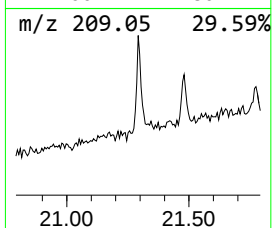
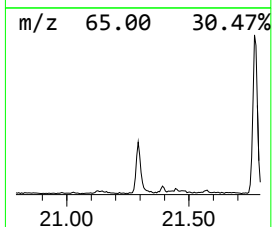
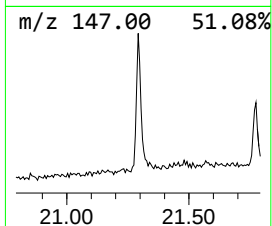
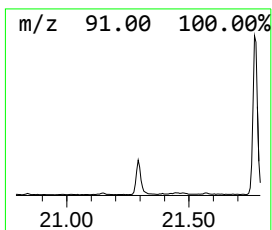
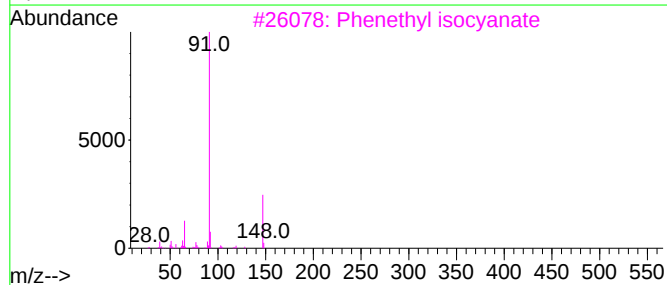
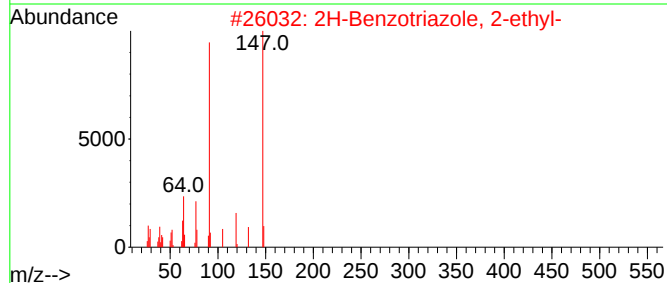
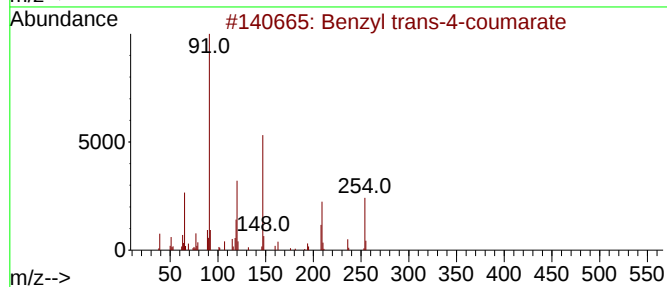
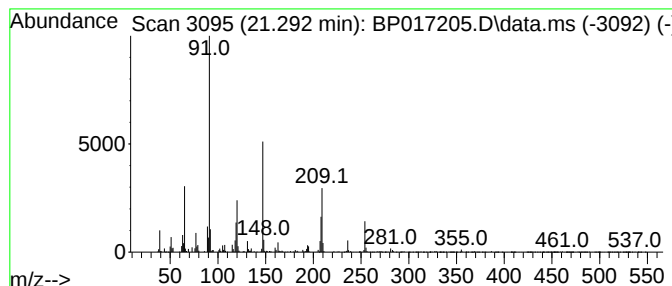
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzyl trans-4-coumarate Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.292	2.85 ng/ul	410464	Chrysene-d12	21.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzyl trans-4-coumarate	254	C16H14O3	061844-62-0	95
2			2H-Benzotriazole, 2-ethyl-	147	C8H9N3	016584-04-6	43
3			Phenethyl isocyanate	147	C9H9NO	001943-82-4	38
4			Pyrido[2,3-d]pyrimidin-4(1H)-one	147	C7H5N3O	024410-19-3	35
5			1-Methyl-1H-indazol-5-amine	147	C8H9N3	050593-24-3	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
Operator : MA/JU
Sample : 04332-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

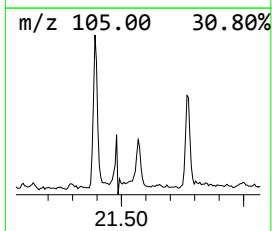
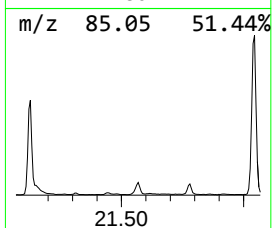
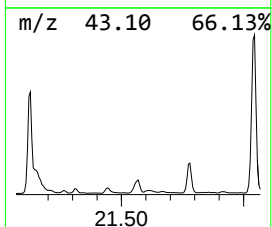
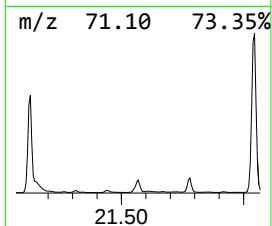
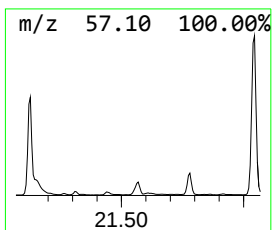
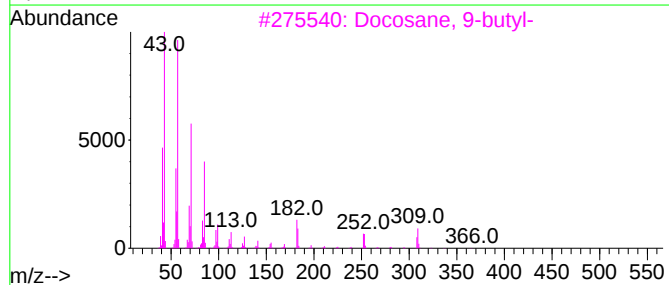
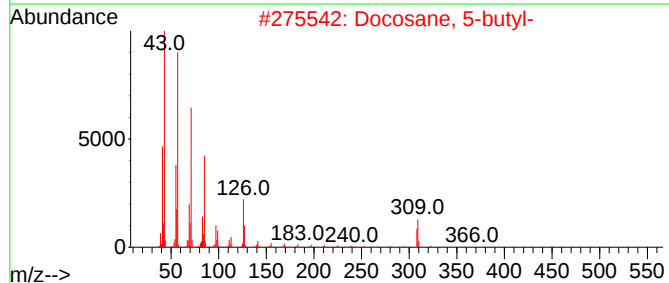
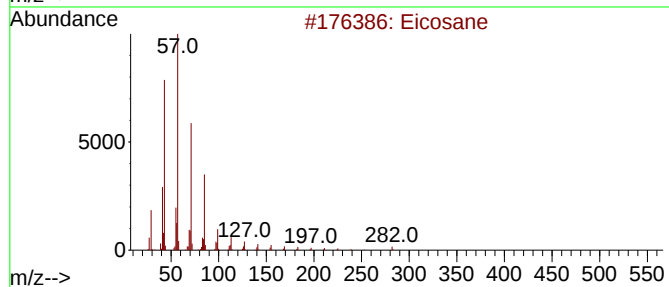
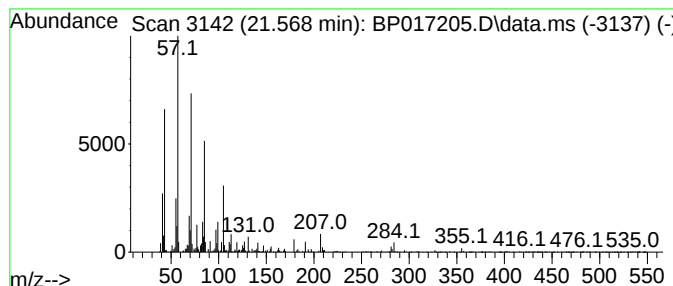
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.568	2.87 ng/ul	413703	Chrysene-d12		21.480	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	97
2	Docosane, 5-butyl-		366	C26H54	055282-16-1	90
3	Docosane, 9-butyl-		366	C26H54	055282-14-9	90
4	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	90
5	Pentadecane, 8-hexyl-		296	C21H44	013475-75-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
Operator : MA/JU
Sample : 04332-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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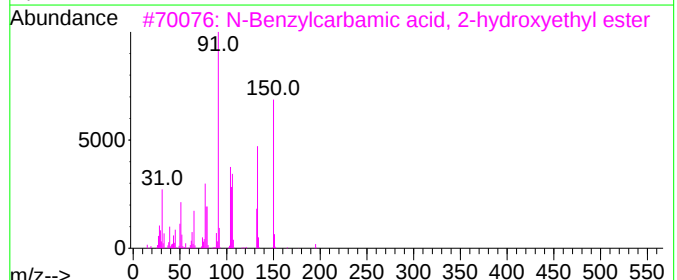
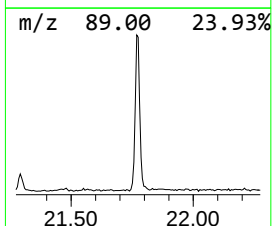
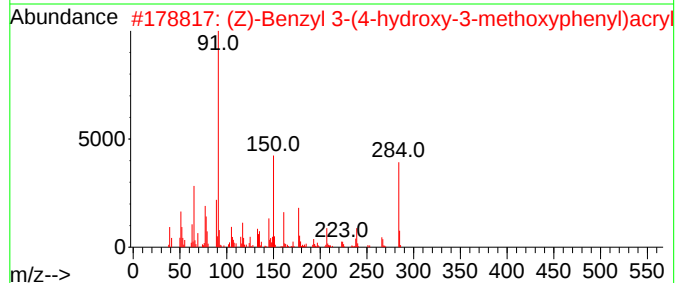
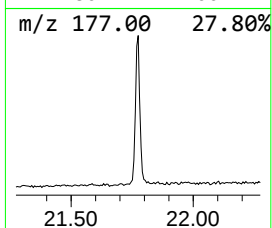
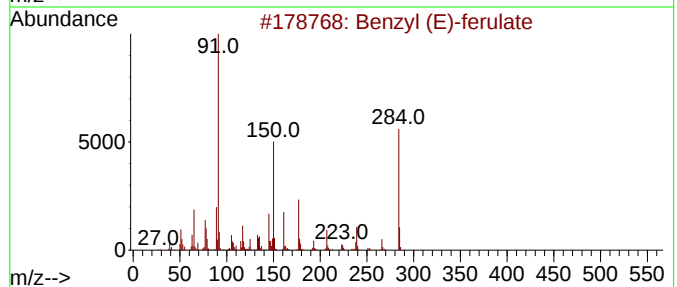
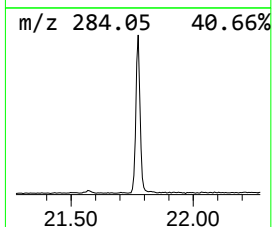
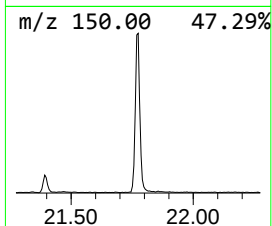
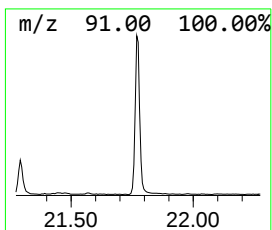
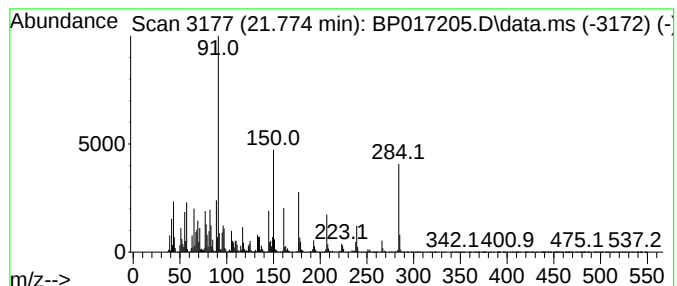
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzyl (E)-ferulate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.774	21.85 ng/ul	3151320	Chrysene-d12	21.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzyl (E)-ferulate	284	C17H16O4	132335-97-8	99
2			(Z)-Benzyl 3-(4-hydroxy-3-methox...	284	C17H16O4	132336-01-7	96
3			N-Benzylcarbamic acid, 2-hydroxy...	195	C10H13NO3	004663-83-6	11
4			Benzeneacetic acid, methyl ester	150	C9H10O2	000101-41-7	11
5			Hydrocinnamic acid	150	C9H10O2	000501-52-0	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
Operator : MA/JU
Sample : 04332-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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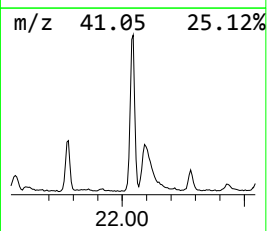
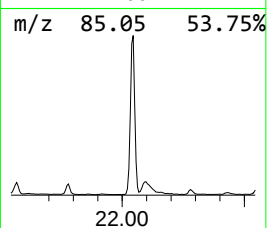
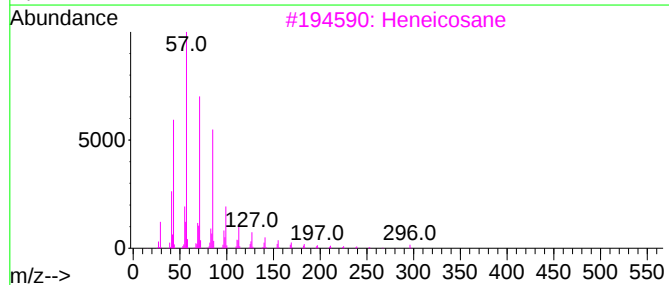
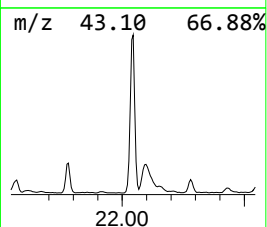
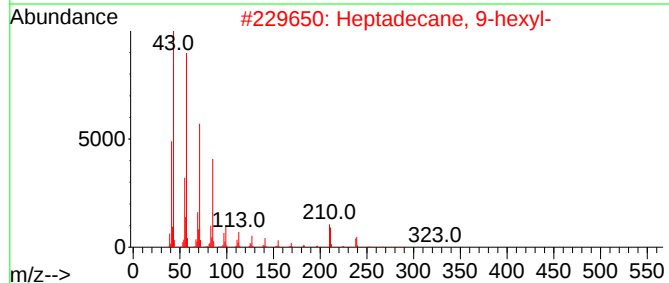
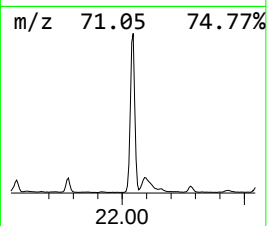
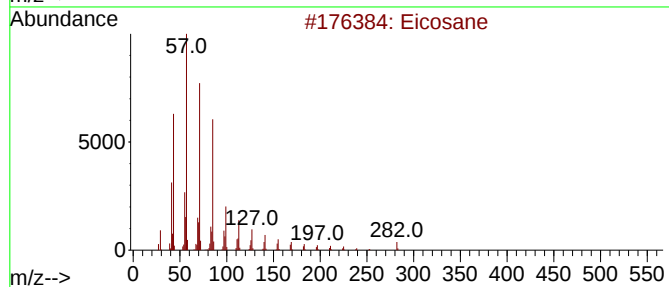
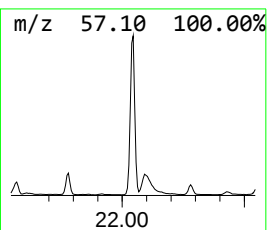
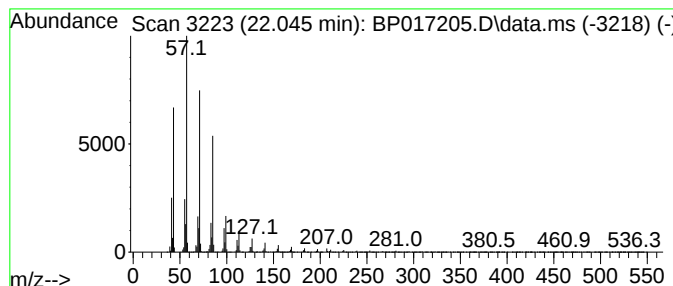
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.045	23.57 ng/ul	3399130	Chrysene-d12	21.480

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	97
2		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
3		Heneicosane	296	C21H44	000629-94-7	91
4		Hexacosane	366	C26H54	000630-01-3	91
5		Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
Operator : MA/JU
Sample : 04332-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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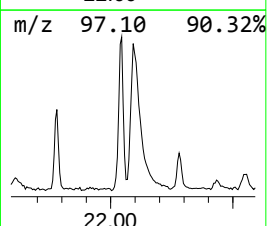
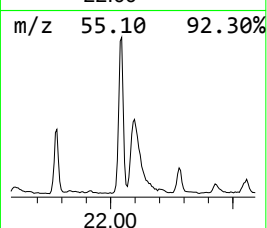
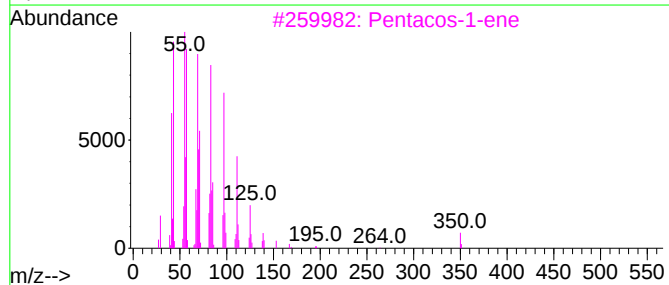
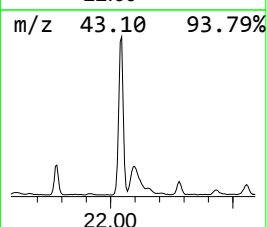
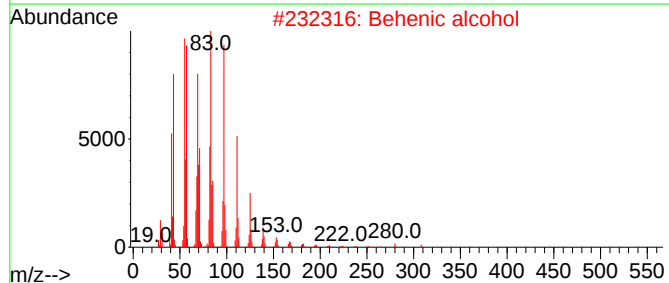
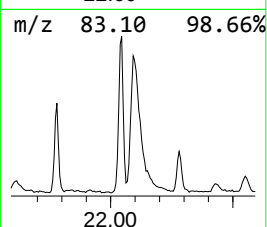
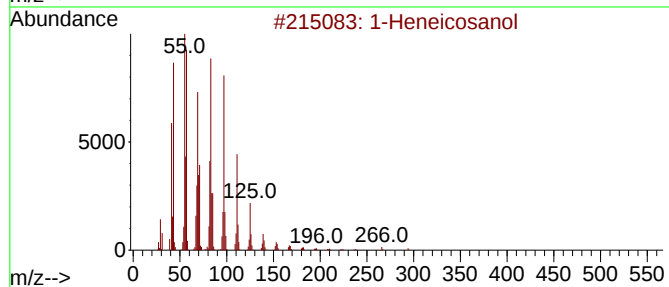
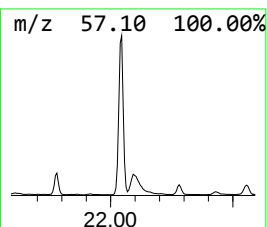
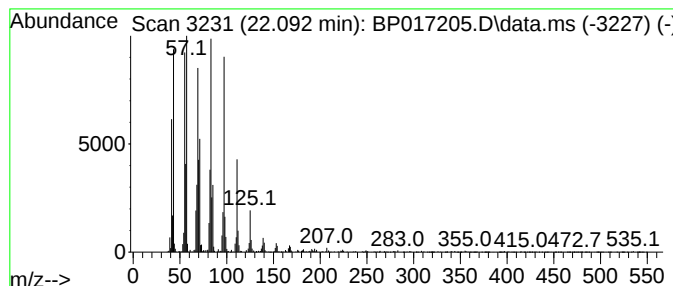
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 1-Heneicosanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.092	13.64 ng/ul	1967830	Chrysene-d12	21.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	95
2			Behenic alcohol	326	C22H46O	000661-19-8	94
3			Pentacos-1-ene	350	C25H50	016980-85-1	94
4			Heptacos-1-ene	378	C27H54	015306-27-1	94
5			1-Heptacosanol	396	C27H56O	002004-39-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
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Sample : 04332-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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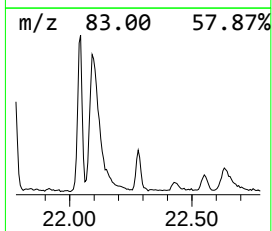
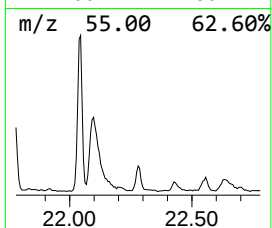
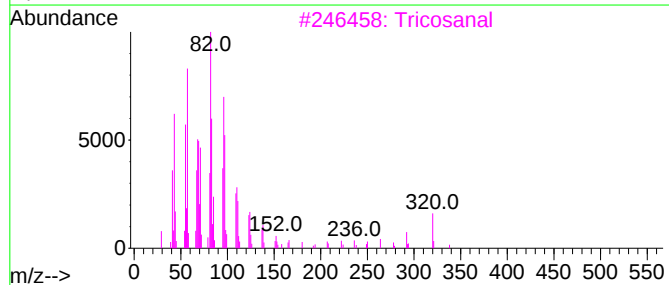
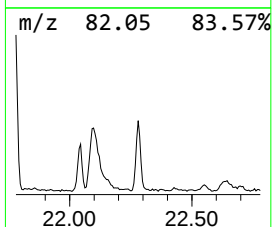
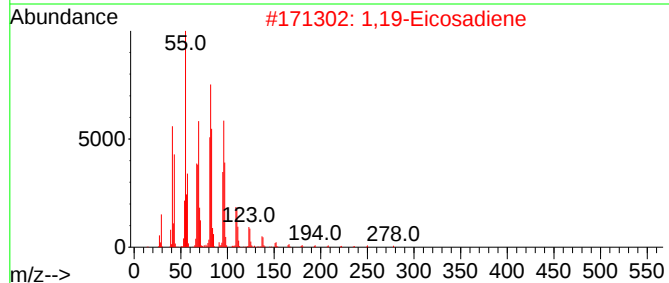
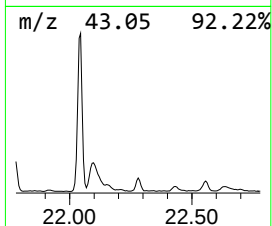
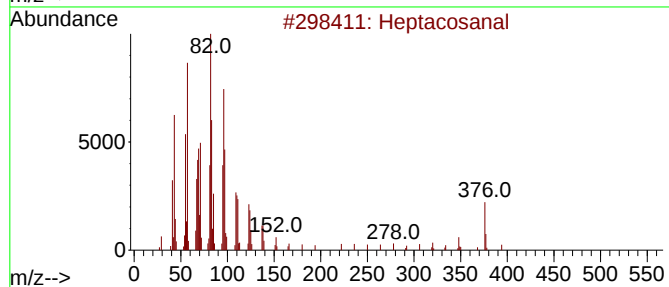
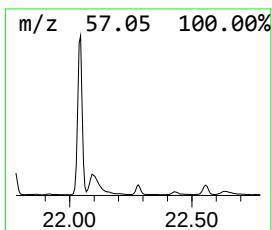
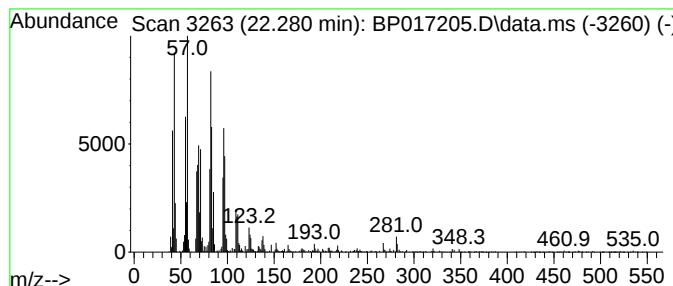
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Heptacosanal Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.280	2.90 ng/ul	418683	Chrysene-d12	21.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosanal	394	C27H54O	072934-03-3	95
2			1,19-Eicosadiene	278	C20H38	014811-95-1	95
3			Tricosanal	338	C23H46O	072934-02-2	94
4			Hexacosanal	380	C26H52O	026627-85-0	94
5			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
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Operator : MA/JU
Sample : 04332-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

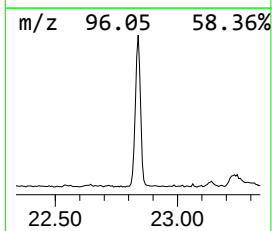
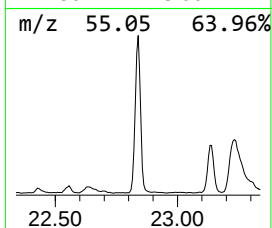
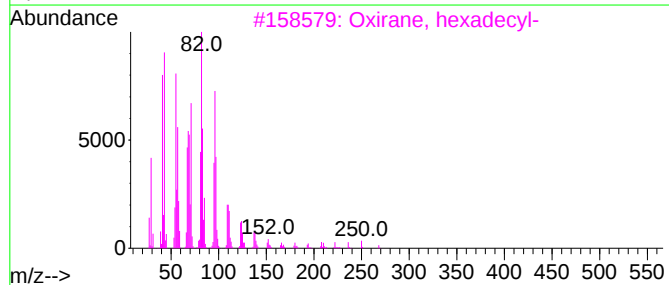
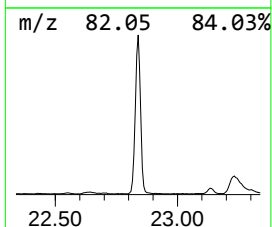
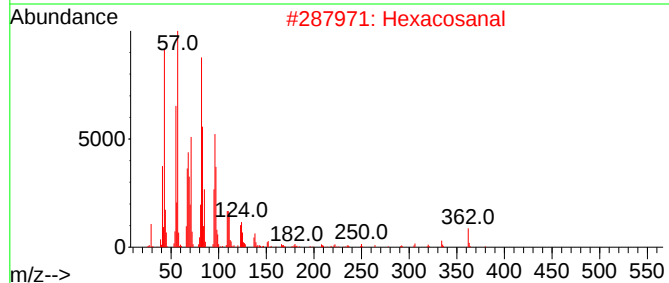
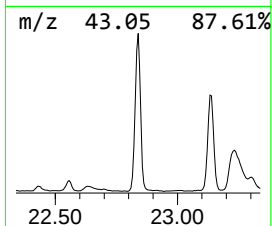
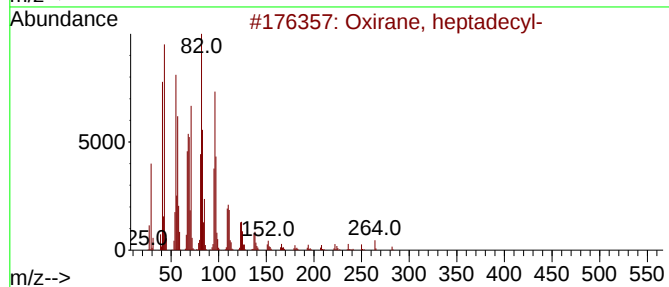
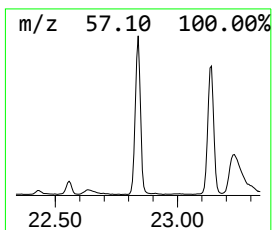
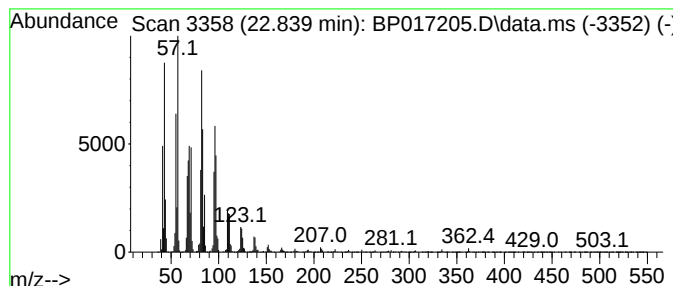
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 15 Oxirane, heptadecyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.839	49.40 ng/ul	6374590	Perylene-d12	24.003	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	96
2	Hexacosanal	380	C26H52O	026627-85-0	94
3	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	94
4	Octadecanal	268	C18H36O	000638-66-4	91
5	Octadecanal	268	C18H36O	000638-66-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
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Misc :
ALS Vial : 10 Sample Multiplier: 1

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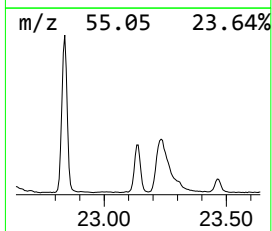
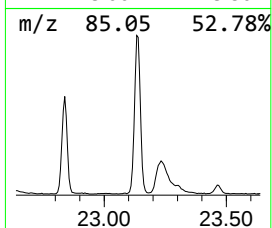
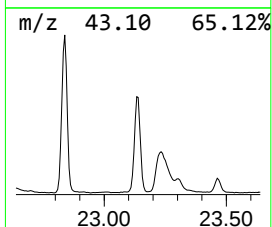
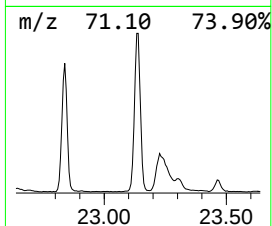
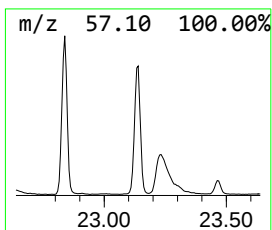
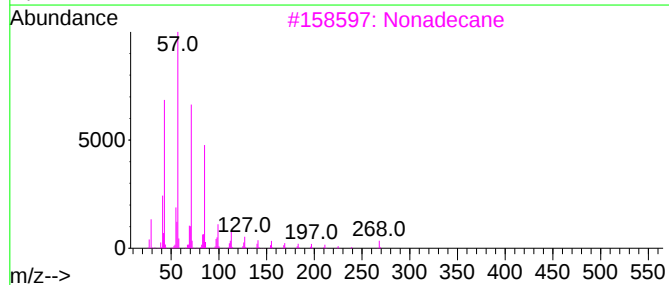
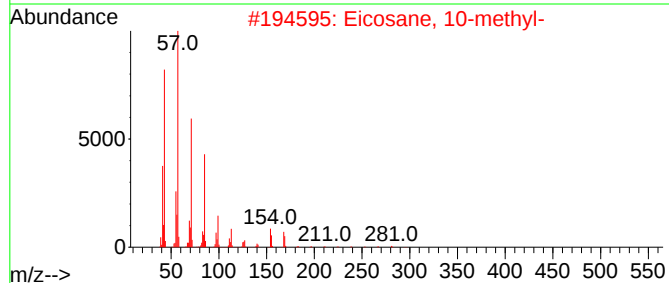
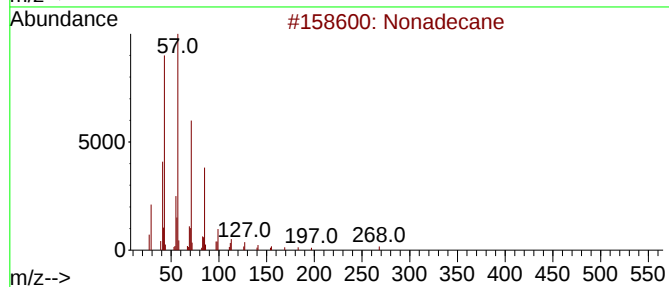
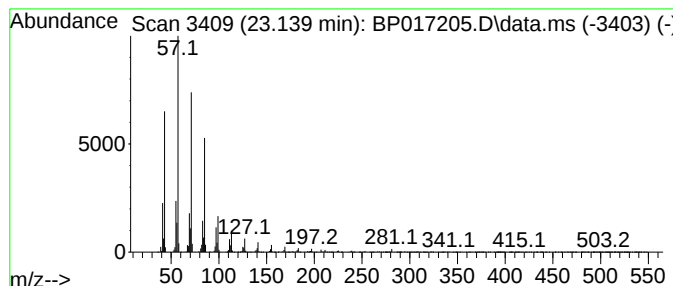
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.139	21.05 ng/ul	2716820	Perylene-d12	24.003

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	95
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
3		Nonadecane	268	C19H40	000629-92-5	93
4		Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
Operator : MA/JU
Sample : 04332-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AC6

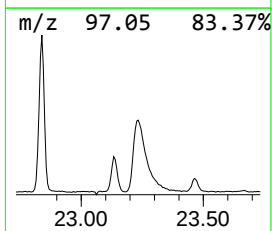
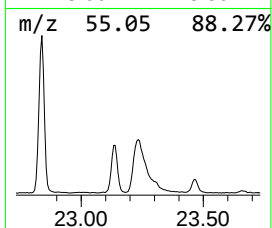
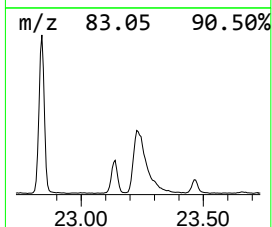
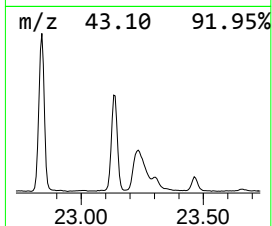
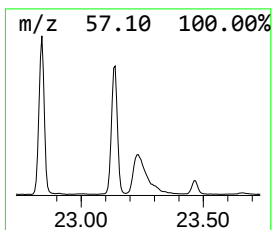
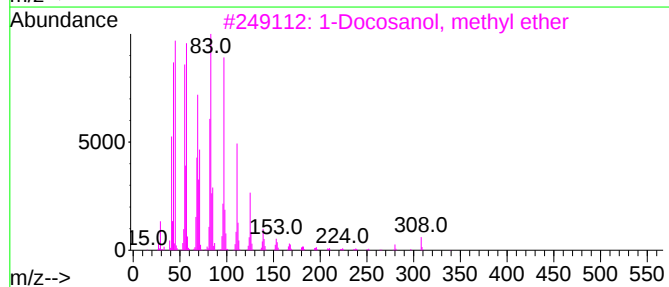
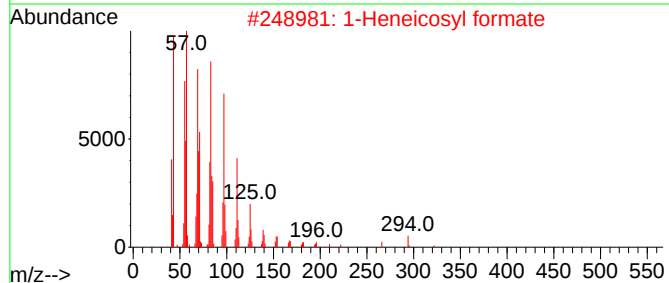
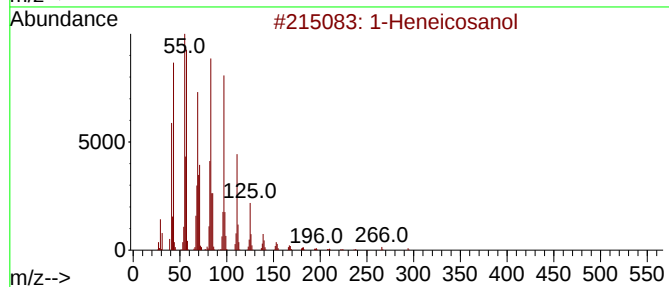
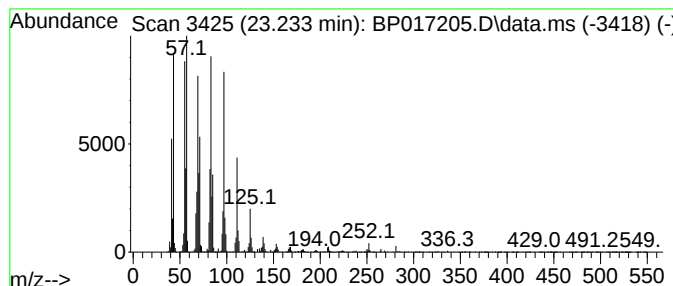
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 1-Heneicosyl formate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.233	27.45 ng/ul	3542530	Perylene-d12	24.003

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	95
2			1-Heneicosyl formate	340	C22H44O2	077899-03-7	95
3			1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	95
4			Heptacos-1-ene	378	C27H54	015306-27-1	94
5			Nonacos-1-ene	406	C29H58	018835-35-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
Operator : MA/JU
Sample : 04332-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

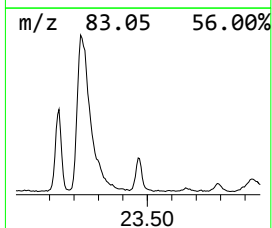
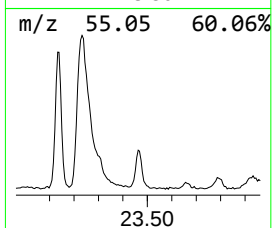
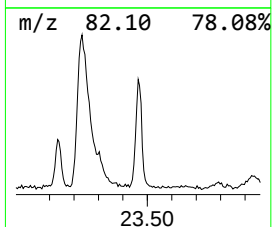
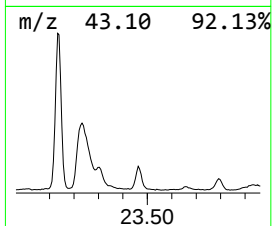
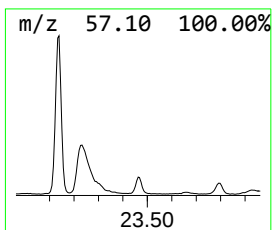
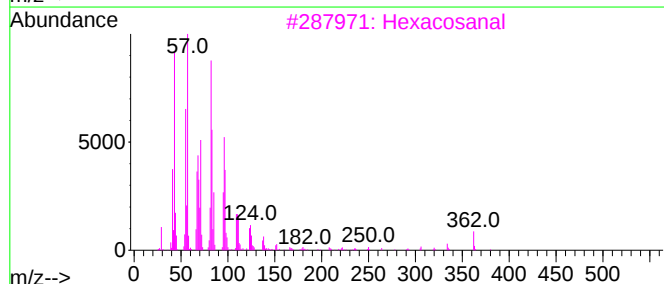
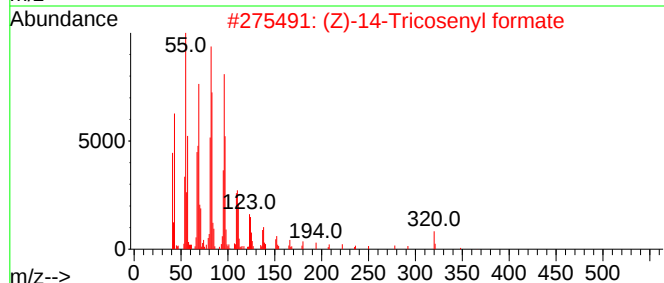
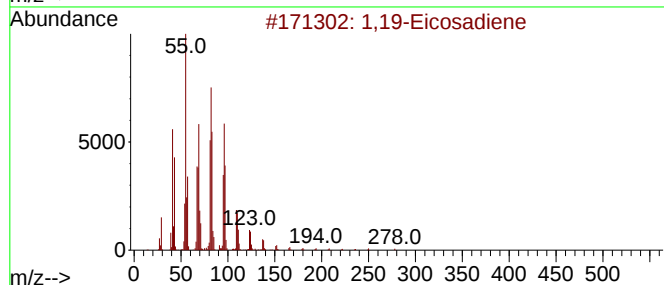
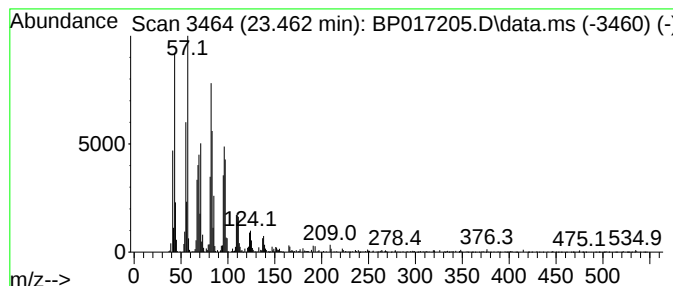
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 1,19-Eicosadiene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.462	4.67 ng/ul	603054	Perylene-d12	24.003	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,19-Eicosadiene	278	C20H38	014811-95-1	95
2	(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	95
3	Hexacosanal	380	C26H52O	026627-85-0	94
4	Octadecanal	268	C18H36O	000638-66-4	93
5	Octacosanal	408	C28H56O	022725-64-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
Operator : MA/JU
Sample : 04332-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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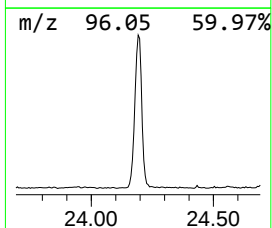
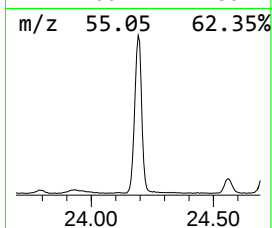
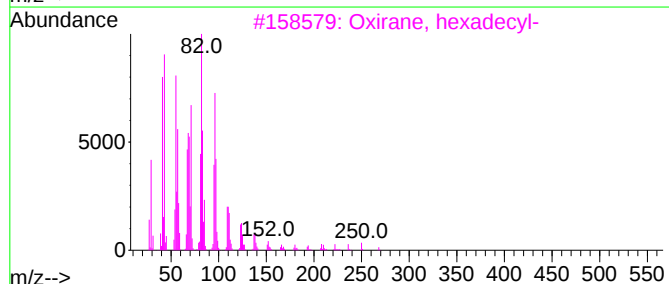
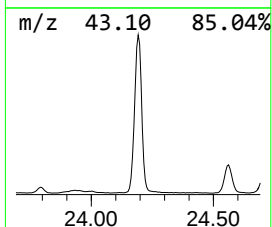
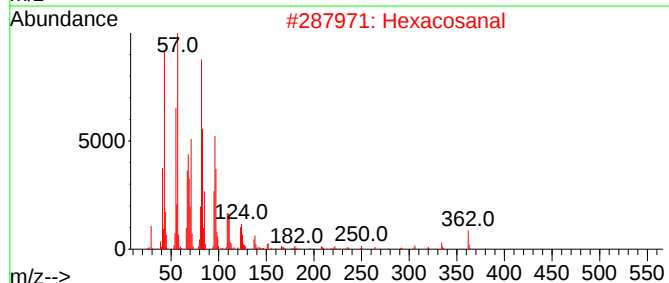
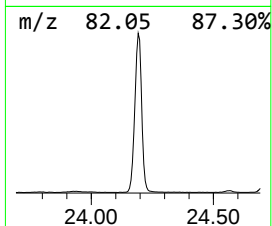
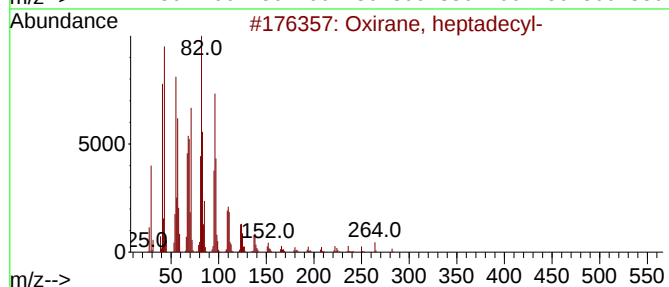
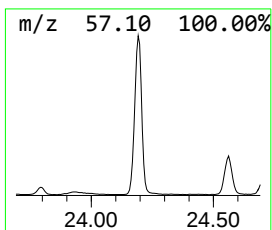
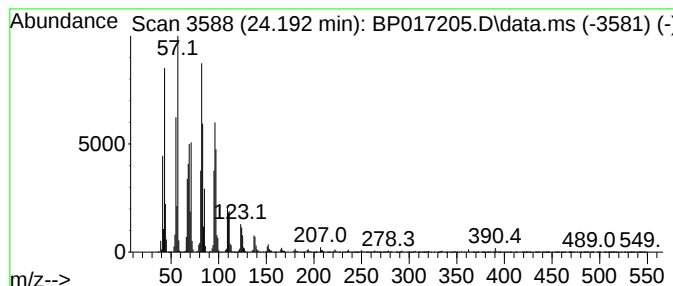
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Hexacosanal Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.192	80.40 ng/ul	10374600	Perylene-d12	24.003

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	94	
2	Hexacosanal	380	C26H52O	026627-85-0	94	
3	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91	
4	Octacosanal	408	C28H56O	022725-64-0	91	
5	Octadecanal	268	C18H36O	000638-66-4	90	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
Operator : MA/JU
Sample : 04332-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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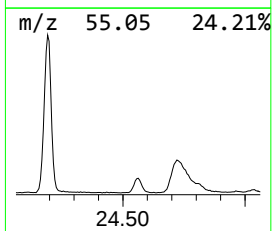
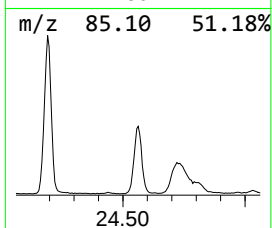
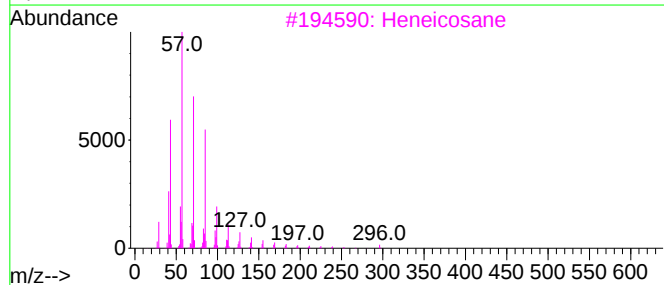
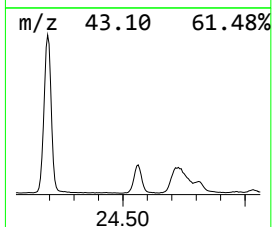
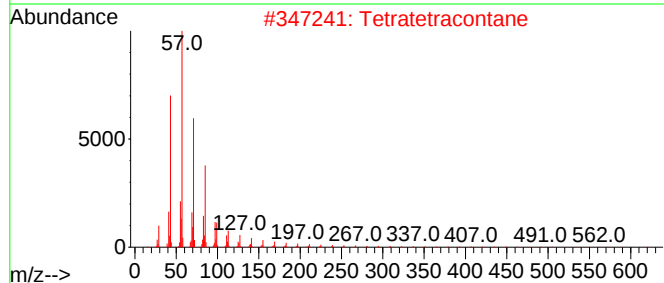
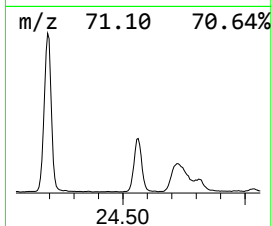
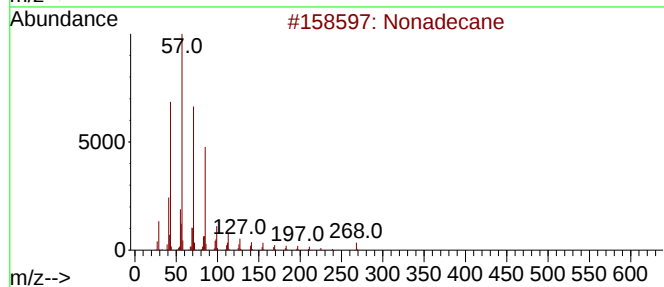
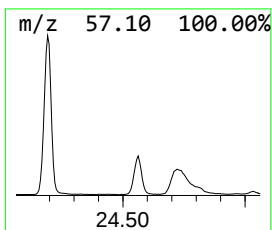
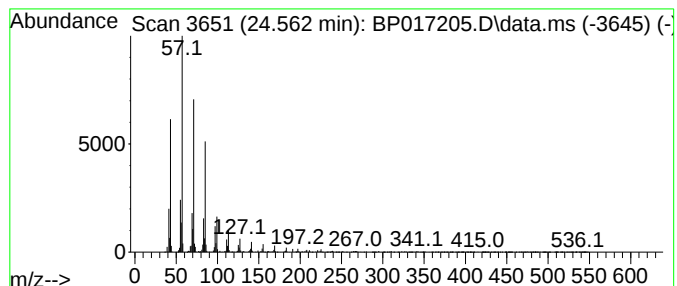
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.562	9.84 ng/ul	1269710	Perylene-d12	24.003

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	93
2			Tetratetracontane	619	C44H90	007098-22-8	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
Operator : MA/JU
Sample : 04332-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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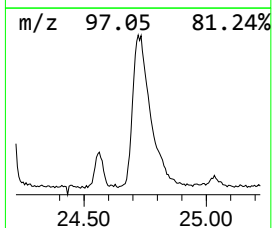
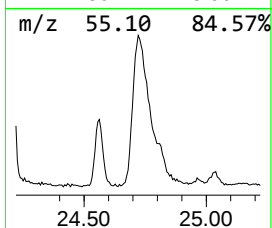
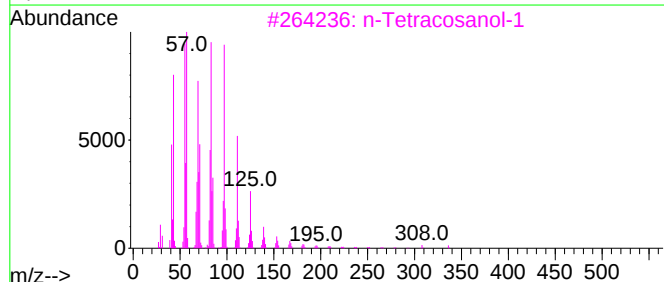
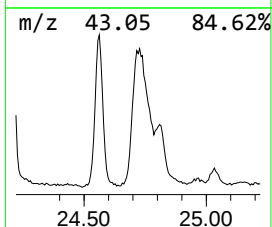
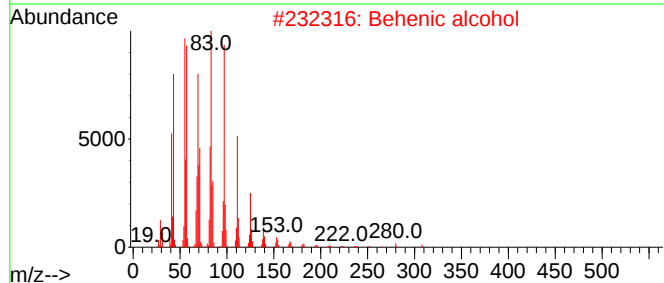
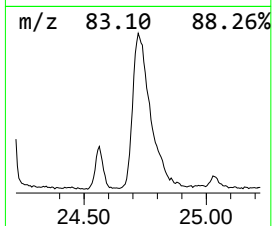
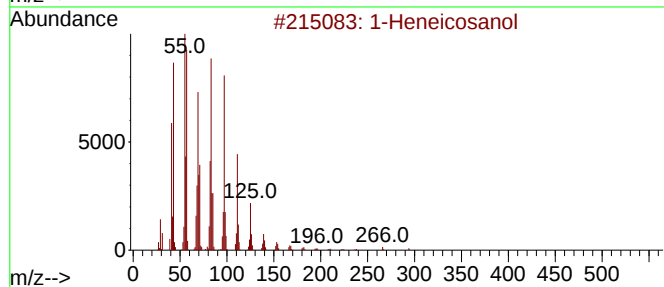
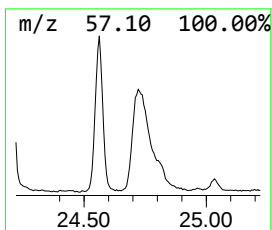
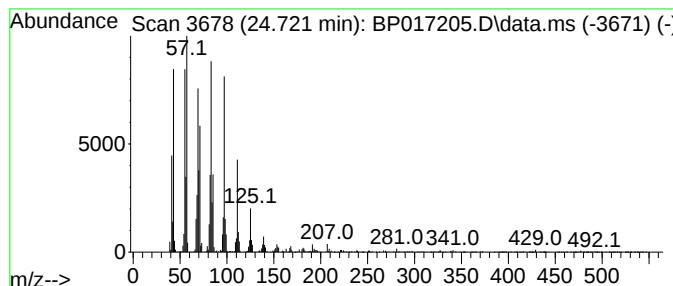
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Behenic alcohol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.721	22.40 ng/ul	2890210	Perylene-d12	24.003

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	95
2			Behenic alcohol	326	C22H46O	000661-19-8	94
3			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4			Heptacos-1-ene	378	C27H54	015306-27-1	94
5			Pentacos-1-ene	350	C25H50	016980-85-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017205.D
Acq On : 18 Sep 2023 15:07
Operator : MA/JU
Sample : 04332-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

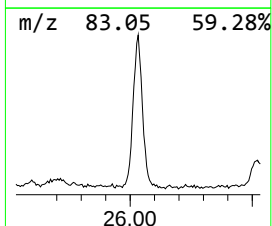
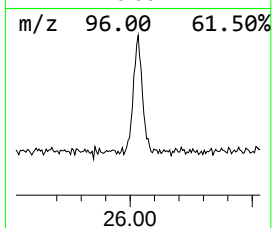
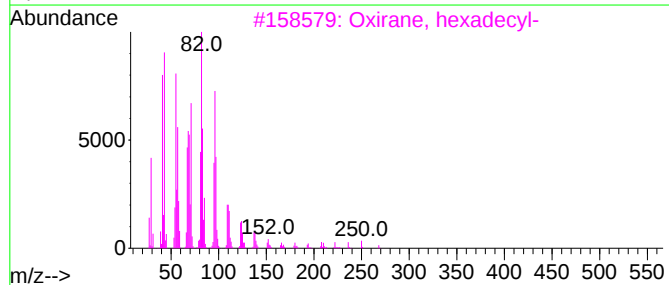
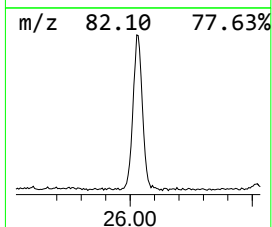
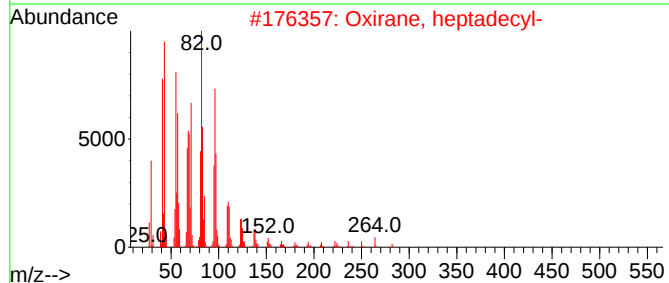
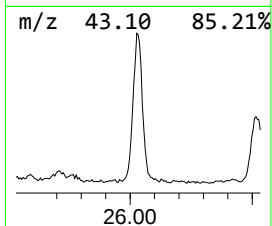
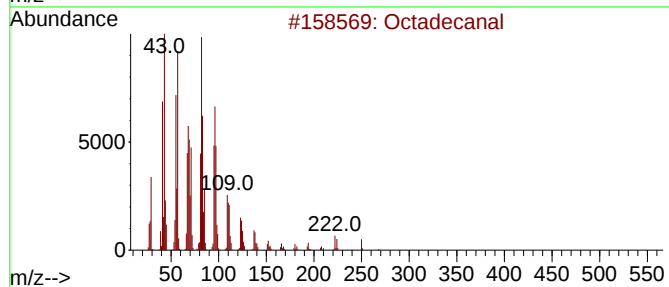
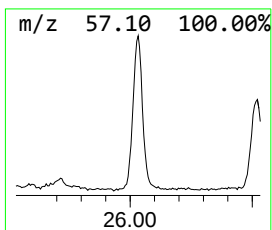
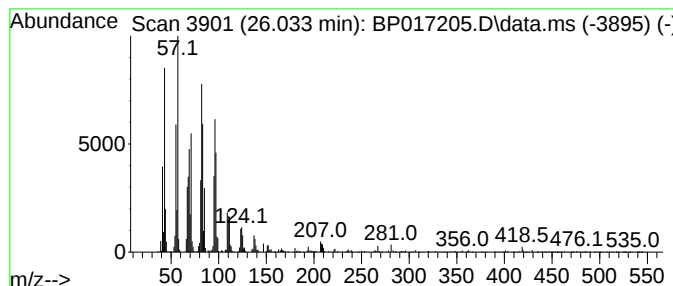
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 22 Octadecanal Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
26.033	12.48 ng/ul	1610140	Perylene-d12		24.003	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octadecanal		268	C18H36O	000638-66-4	92
2	Oxirane, heptadecyl-		282	C19H38O	067860-04-2	90
3	Oxirane, hexadecyl-		268	C18H36O	007390-81-0	90
4	13-Methyltetradecanal		226	C15H30O	075853-51-9	90
5	1,19-Eicosadiene		278	C20H38	014811-95-1	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
 Data File : BP017205.D
 Acq On : 18 Sep 2023 15:07
 Operator : MA/JU
 Sample : 04332-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.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
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2H-Pyran-2-one,...	9.251	2.5	ng/ul	155838	1	7.945	1273160	20.0
Benzoic acid	10.022	5.6	ng/ul	532338	2	10.769	1888900	20.0
n-Hexadecanoic ...	18.221	9.5	ng/ul	1431750	4	17.374	3008990	20.0
22-Tricosenoic ...	19.368	3.3	ng/ul	496128	4	17.374	3008990	20.0
Octadecanoic acid	19.457	3.1	ng/ul	453467	5	21.480	2884810	20.0
Carbonic acid, ...	20.180	12.1	ng/ul	1743170	5	21.480	2884810	20.0
(DEL) Alkane: S...	21.127	13.7	ng/ul	1977440	5	21.480	2884810	20.0
Benzyl trans-4-...	21.292	2.9	ng/ul	410464	5	21.480	2884810	20.0
(DEL) Alkane: S...	21.568	2.9	ng/ul	413703	5	21.480	2884810	20.0
Benzyl (E)-feru...	21.774	21.9	ng/ul	3151320	5	21.480	2884810	20.0
(DEL) Alkane: S...	22.045	23.6	ng/ul	3399130	5	21.480	2884810	20.0
1-Heneicosanol	22.092	13.6	ng/ul	1967830	5	21.480	2884810	20.0
Heptacosanal	22.280	2.9	ng/ul	418683	5	21.480	2884810	20.0
Oxirane, heptad...	22.839	49.4	ng/ul	6374590	6	24.003	2580730	20.0
(DEL) Alkane: S...	23.139	21.1	ng/ul	2716820	6	24.003	2580730	20.0
1-Heneicosyl fo...	23.233	27.4	ng/ul	3542530	6	24.003	2580730	20.0
1,19-Eicosadiene	23.462	4.7	ng/ul	603054	6	24.003	2580730	20.0
Hexacosanal	24.192	80.4	ng/ul	10374600	6	24.003	2580730	20.0
(DEL) Alkane: S...	24.562	9.8	ng/ul	1269710	6	24.003	2580730	20.0
Behenic alcohol	24.721	22.4	ng/ul	2890210	6	24.003	2580730	20.0
Octadecanal	26.033	12.5	ng/ul	1610140	6	24.003	2580730	20.0