

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
 Data File : BP013316.D
 Acq On : 25 Dec 2022 00:50
 Operator : CG/JU
 Sample : N6016-20
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBYE3

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-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP013316.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.328	21	24	32	rVB	119090	165971	0.92%	0.106%
2	3.781	99	101	110	rVV	52310	89901	0.50%	0.057%
3	3.864	111	115	124	rVV	139663	185597	1.03%	0.118%
4	3.987	131	136	141	rVB	154304	225908	1.25%	0.144%
5	4.087	149	153	160	rVB	50146	67208	0.37%	0.043%
6	4.464	214	217	223	rVB3	35034	51489	0.28%	0.033%
7	4.652	245	249	253	rBV	65660	88128	0.49%	0.056%
8	4.699	253	257	263	rVB	174213	237154	1.31%	0.151%
9	5.022	305	312	323	rBV2	585278	1145371	6.33%	0.730%
10	5.116	325	328	336	rVB2	25539	43211	0.24%	0.028%
11	6.540	565	570	575	rBV	57083	92713	0.51%	0.059%
12	6.610	577	582	587	rVB	44038	66528	0.37%	0.042%
13	6.916	629	634	640	rVB3	54771	88946	0.49%	0.057%
14	7.099	658	665	675	rBV	1878559	3083157	17.05%	1.966%
15	7.263	687	693	703	rVB	1601280	2454460	13.57%	1.565%
16	7.369	705	711	717	rVB	107023	166492	0.92%	0.106%
17	7.469	721	728	736	rBV	1981488	3108505	17.19%	1.982%
18	7.546	736	741	746	rVB	33710	48936	0.27%	0.031%
19	7.940	800	808	815	rBV	2586016	4026086	22.26%	2.567%
20	8.652	921	929	938	rBV	1949297	3060567	16.92%	1.951%
21	8.769	943	949	959	rVB	217355	340133	1.88%	0.217%
22	9.104	998	1006	1017	rBV	1827162	3043158	16.83%	1.940%
23	9.504	1066	1074	1080	rBV	57079	91581	0.51%	0.058%
24	9.699	1099	1107	1113	rVB	40713	72777	0.40%	0.046%
25	9.828	1122	1129	1137	rBV	1497489	2484663	13.74%	1.584%
26	9.981	1151	1155	1162	rBV2	26959	54259	0.30%	0.035%
27	10.051	1162	1167	1175	rVB2	21657	48349	0.27%	0.031%
28	10.187	1182	1190	1196	rVB3	21127	44857	0.25%	0.029%
29	10.369	1213	1221	1233	rBV	2247938	3754067	20.76%	2.393%
30	10.534	1243	1249	1256	rVB5	24619	48011	0.27%	0.031%
31	10.751	1278	1286	1297	rBV	3387875	5643784	31.21%	3.598%
32	10.893	1302	1310	1323	rBV	1104504	2011112	11.12%	1.282%
33	11.287	1370	1377	1383	rBV9	26231	49112	0.27%	0.031%
34	11.540	1409	1420	1425	rBV8	24941	71094	0.39%	0.045%
35	12.110	1513	1517	1522	rVV5	31877	49333	0.27%	0.031%
36	12.334	1551	1555	1561	rVB2	40008	61146	0.34%	0.039%

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Integrator: RTE

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Stop Thrs : 0

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Title : SVOA CALIBRATION

37	12.863	1638	1645	1654	rBV	572592	917228	5.07%	0.585%
38	13.104	1682	1686	1690	rBV3	31085	43408	0.24%	0.028%
39	13.340	1723	1726	1729	rVV3	27555	37724	0.21%	0.024%
40	13.392	1729	1735	1739	rVB2	43858	75379	0.42%	0.048%
41	13.451	1739	1745	1751	rBV	539752	736178	4.07%	0.469%
42	13.904	1817	1822	1829	rVB5	54114	89043	0.49%	0.057%
43	13.987	1829	1836	1843	rBV	3526300	4929641	27.26%	3.143%
44	14.104	1852	1856	1858	rBV3	64963	90064	0.50%	0.057%
45	14.134	1858	1861	1864	rVV2	124344	156763	0.87%	0.100%
46	14.169	1864	1867	1871	rVB	86725	118340	0.65%	0.075%
47	14.275	1879	1885	1894	rVB	3896016	5771498	31.91%	3.680%
48	14.416	1905	1909	1914	rVB	94995	125902	0.70%	0.080%
49	14.469	1914	1918	1924	rVB2	58009	88256	0.49%	0.056%
50	14.539	1924	1930	1931	rBV5	28030	38133	0.21%	0.024%
51	14.581	1931	1937	1943	rBV	4180171	6315933	34.92%	4.027%
52	14.639	1943	1947	1952	rVB	254376	335069	1.85%	0.214%
53	14.692	1952	1956	1962	rVB4	126562	194193	1.07%	0.124%
54	14.787	1966	1972	1977	rBV	825625	1343136	7.43%	0.856%
55	14.887	1984	1989	1994	rBV2	216969	310334	1.72%	0.198%
56	14.945	1994	1999	2003	rVV5	140254	273822	1.51%	0.175%
57	14.992	2003	2007	2014	rVB2	235510	362836	2.01%	0.231%
58	15.492	2084	2092	2097	rBV9	42443	98046	0.54%	0.063%
59	15.575	2100	2106	2113	rBV	4549939	6690931	37.00%	4.266%
60	15.698	2122	2127	2138	rVV	733014	989821	5.47%	0.631%
61	15.816	2143	2147	2155	rVB9	27700	54271	0.30%	0.035%
62	15.945	2160	2169	2178	rBV	3875776	5261253	29.09%	3.354%
63	16.057	2185	2188	2195	rVB2	52266	70054	0.39%	0.045%
64	16.228	2211	2217	2223	rVB10	24891	49194	0.27%	0.031%
65	16.304	2223	2230	2235	rBV6	47393	107452	0.59%	0.069%
66	16.510	2260	2265	2271	rVV	293662	396809	2.19%	0.253%
67	16.728	2297	2302	2306	rBV	166725	223388	1.24%	0.142%
68	17.086	2359	2363	2366	rBV5	25620	40754	0.23%	0.026%
69	17.222	2383	2386	2394	rVB5	43891	68713	0.38%	0.044%
70	17.292	2394	2398	2400	rBV5	31240	45229	0.25%	0.029%
71	17.345	2400	2407	2416	rVV	4183853	6094491	33.70%	3.885%
72	17.439	2417	2423	2431	rBV	4296599	6041881	33.41%	3.852%
73	17.998	2514	2518	2523	rBV3	122410	161048	0.89%	0.103%
74	18.157	2542	2545	2550	rBV3	82951	112605	0.62%	0.072%
75	18.216	2550	2555	2564	rBV	1735335	2359888	13.05%	1.505%
76	18.463	2594	2597	2600	rVB	104394	122771	0.68%	0.078%

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 Title : SVOA CALIBRATION

77	18.622	2622	2624	2632	rVB3	144747	198526	1.10%	0.127%
78	18.722	2638	2641	2646	rVB2	159846	201417	1.11%	0.128%
79	18.780	2647	2651	2654	rVV	388970	449180	2.48%	0.286%
80	18.816	2654	2657	2661	rVB3	128742	158584	0.88%	0.101%
81	18.992	2684	2687	2693	rBV2	47865	72681	0.40%	0.046%
82	19.116	2704	2708	2712	rBV2	186447	230956	1.28%	0.147%
83	19.251	2727	2731	2735	rVB2	213845	273399	1.51%	0.174%
84	19.333	2737	2745	2758	rBV3	758622	1685515	9.32%	1.075%
85	19.457	2760	2766	2784	rBV	12607530	18085575	100.00%	11.530%
86	19.674	2797	2803	2813	rVB	6058025	8209472	45.39%	5.234%
87	19.822	2824	2828	2838	rBV	3449138	4762311	26.33%	3.036%
88	19.898	2839	2841	2849	rVB7	128289	194291	1.07%	0.124%
89	20.080	2869	2872	2876	rBV5	111151	152067	0.84%	0.097%
90	20.133	2879	2881	2883	rVV2	99911	103547	0.57%	0.066%
91	20.169	2884	2887	2888	rVV	500459	505771	2.80%	0.322%
92	20.198	2888	2892	2897	rVB	4202434	4912470	27.16%	3.132%
93	20.316	2909	2912	2918	rBV6	125819	190704	1.05%	0.122%
94	20.445	2929	2934	2940	rBV	1141445	1359173	7.52%	0.867%
95	20.522	2944	2947	2953	rVB4	217380	323386	1.79%	0.206%
96	20.674	2969	2973	2979	rVB3	271155	364593	2.02%	0.232%
97	21.127	3046	3050	3060	rBV	1342920	2046266	11.31%	1.305%
98	21.433	3097	3102	3107	rVB	5312567	7110324	39.31%	4.533%
99	23.751	3489	3496	3512	rVB	4574878	9627900	53.24%	6.138%
100	23.915	3517	3524	3532	rVB2	3875414	7929263	43.84%	5.055%

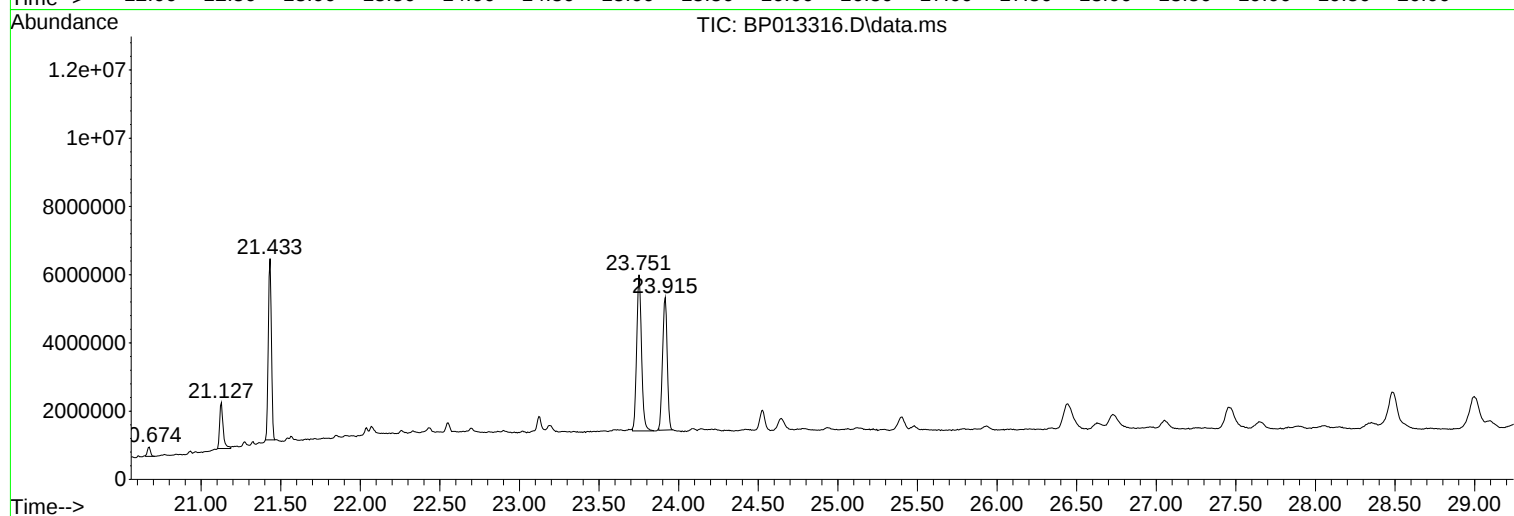
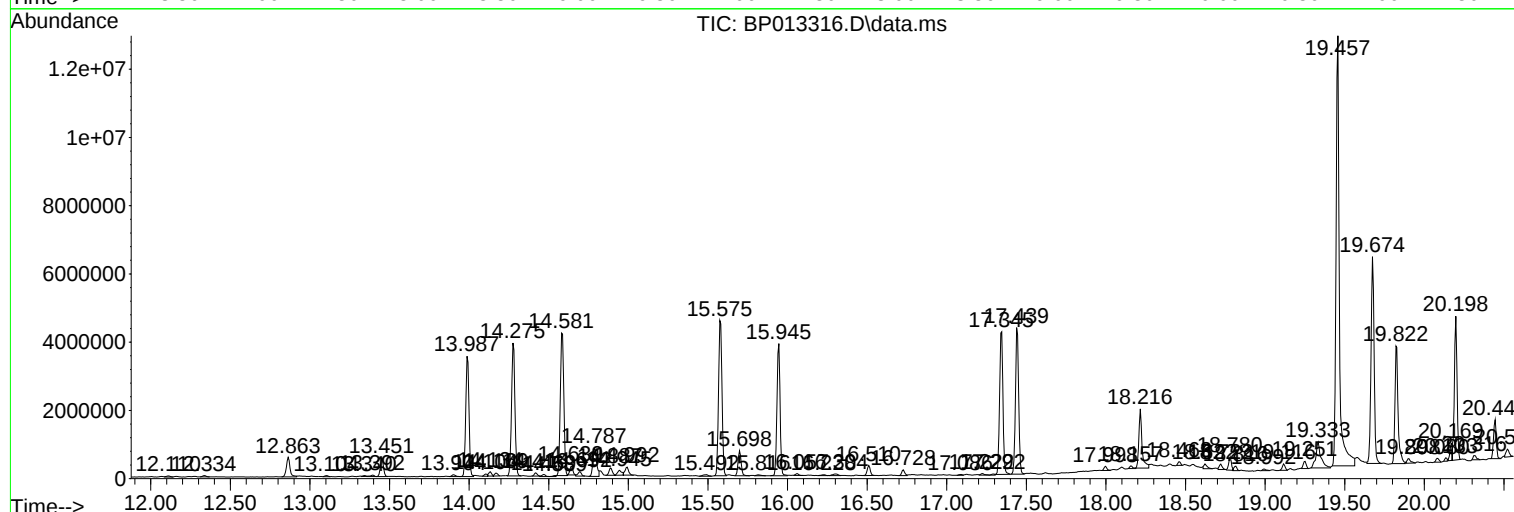
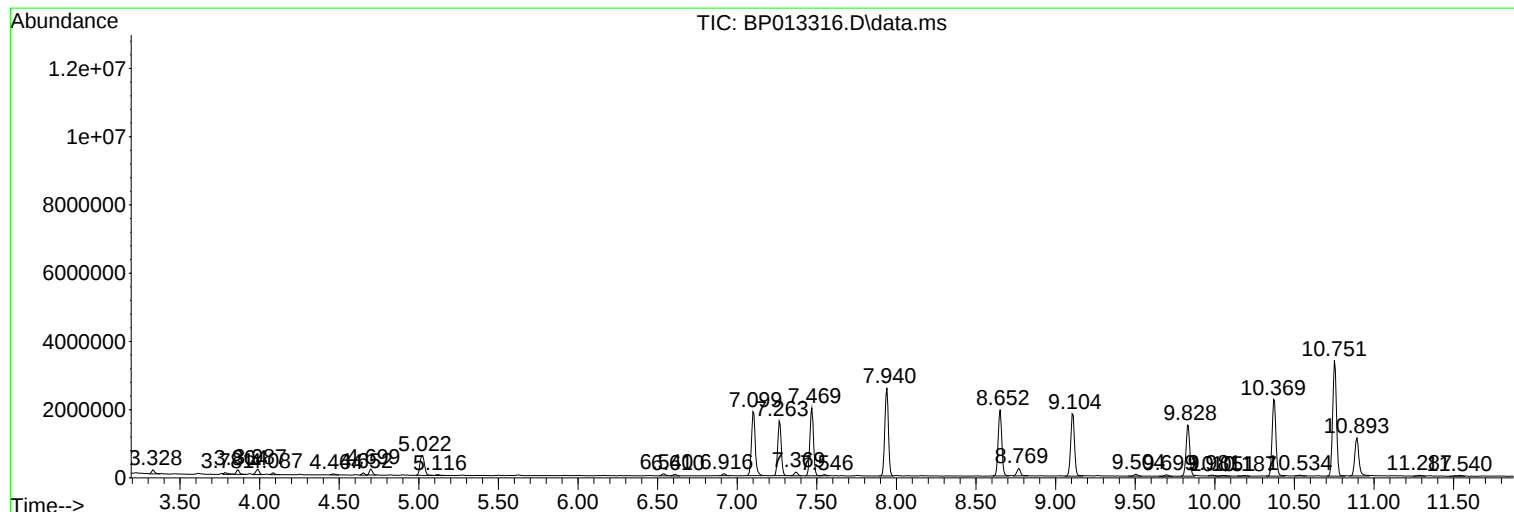
Sum of corrected areas: 156852614

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Instrument :
BNA_P
ClientSampleId :
DBYE3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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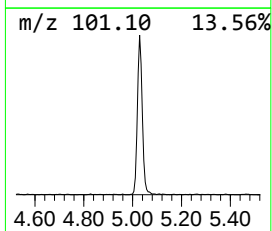
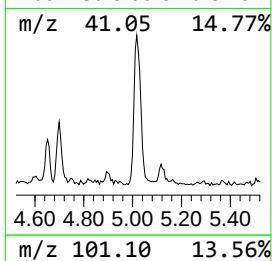
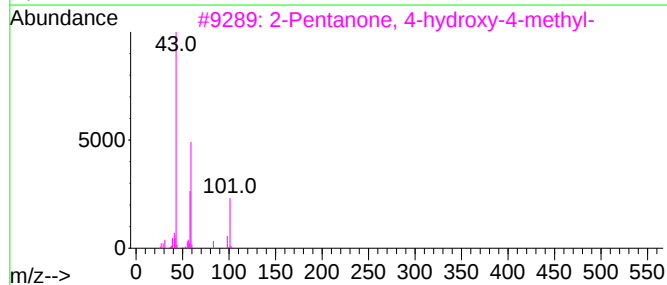
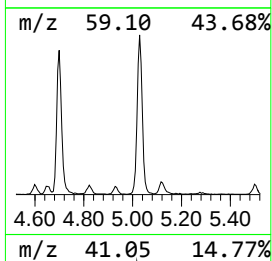
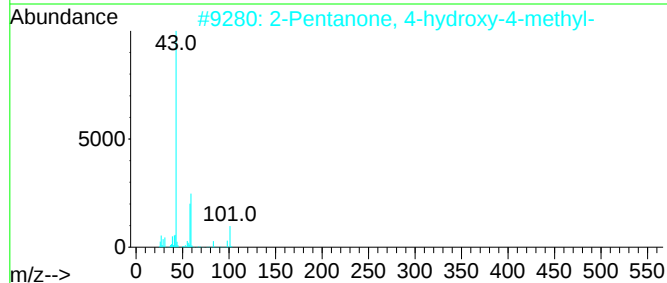
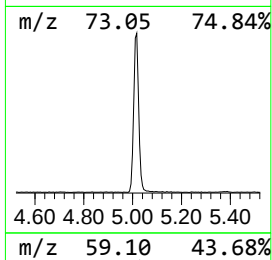
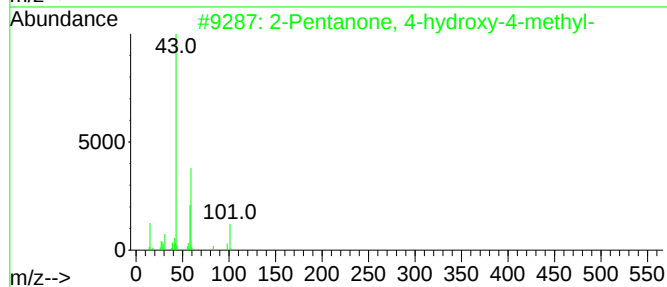
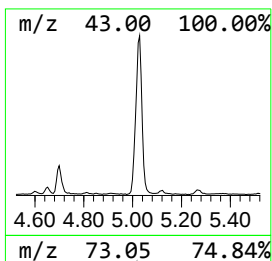
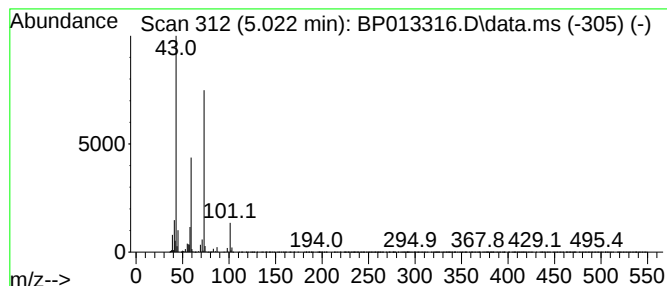
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.022	5.69 ng/ul	1145370	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	37		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	37		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	37		
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28		
5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	25		



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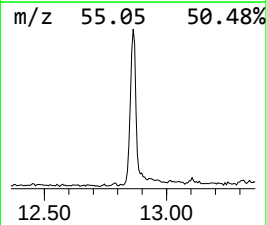
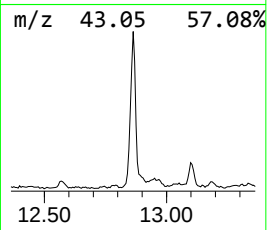
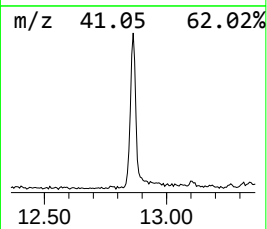
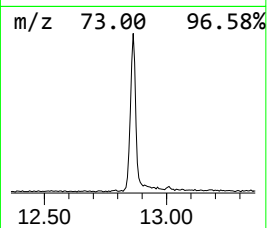
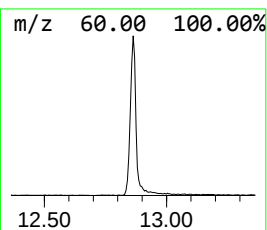
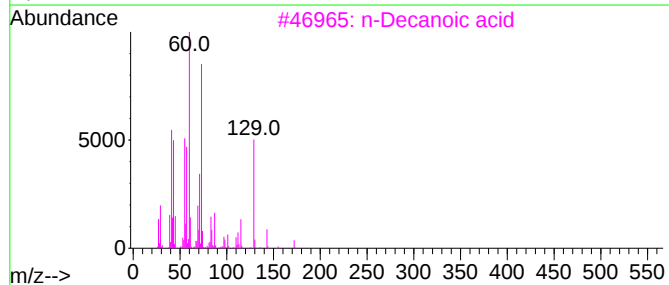
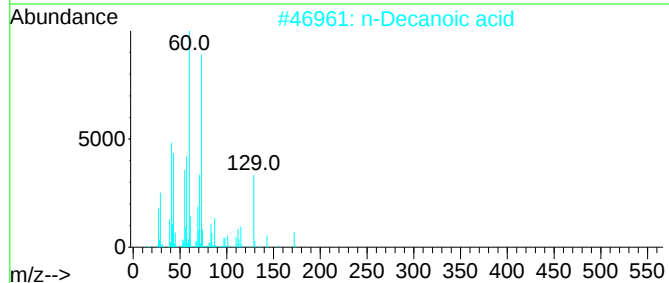
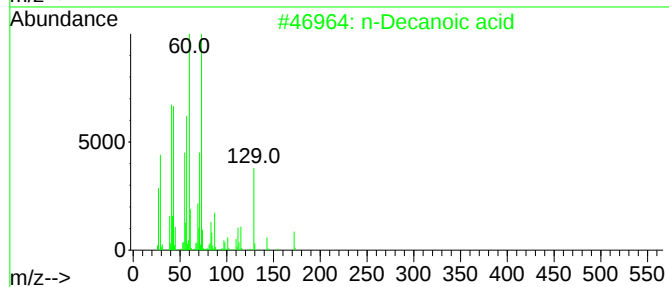
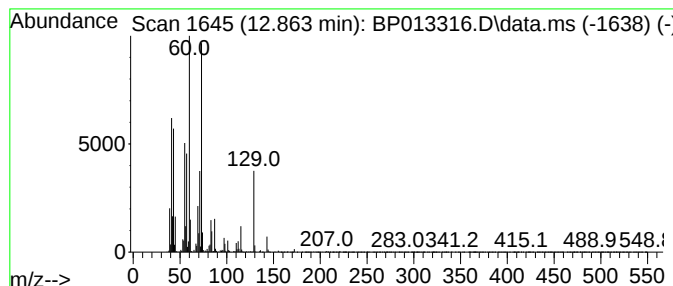
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Decanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.863	2.90 ng/ul	917228	Acenaphthene-d10	14.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Decanoic acid	172	C10H20O2	000334-48-5	91
2			n-Decanoic acid	172	C10H20O2	000334-48-5	90
3			n-Decanoic acid	172	C10H20O2	000334-48-5	86
4			n-Decanoic acid	172	C10H20O2	000334-48-5	86
5			n-Decanoic acid	172	C10H20O2	000334-48-5	78



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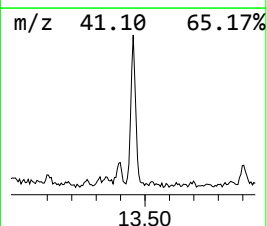
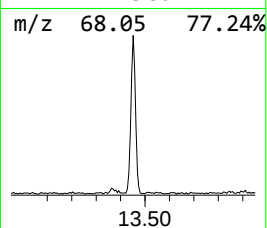
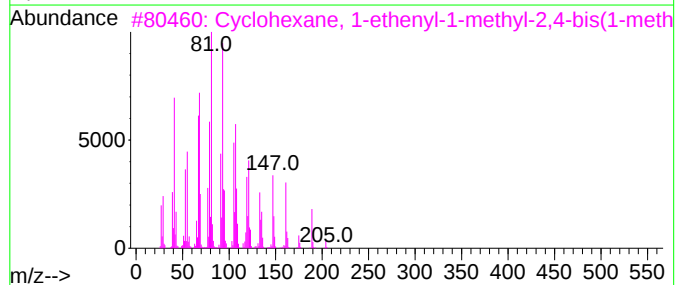
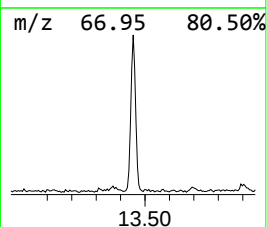
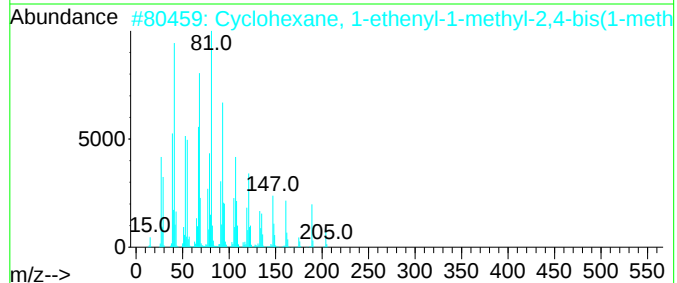
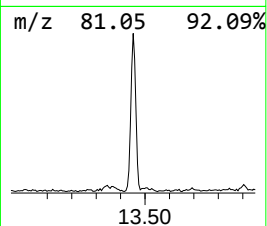
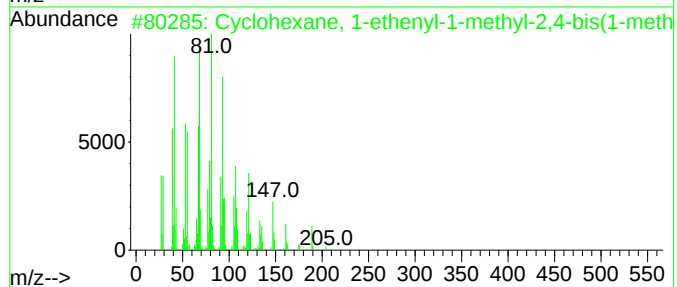
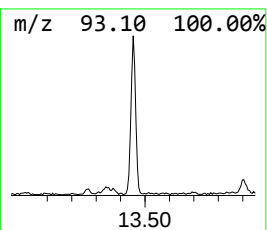
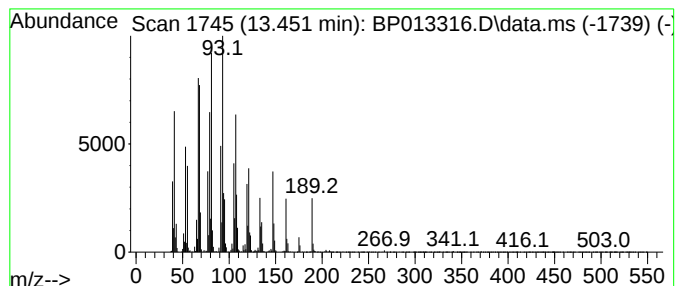
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Cyclohexane, 1-ethenyl-1-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.451	2.33 ng/ul	736178	Acenaphthene-d10	14.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-ethenyl-1-methyl-...	204	C15H24	110823-68-2	86
2			Cyclohexane, 1-ethenyl-1-methyl-...	204	C15H24	000515-13-9	72
3			Cyclohexane, 1-ethenyl-1-methyl-...	204	C15H24	000515-13-9	64
4			Cyclohexane, 1-ethenyl-1-methyl-...	204	C15H24	000515-13-9	58
5			1,3-Pentadiene	68	C5H8	000504-60-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013316.D
Acq On : 25 Dec 2022 00:50
Operator : CG/JU
Sample : N6016-20
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYE3

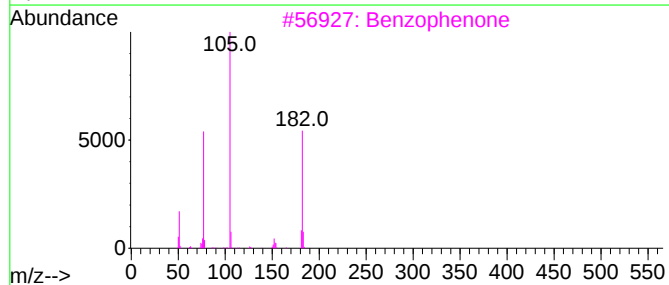
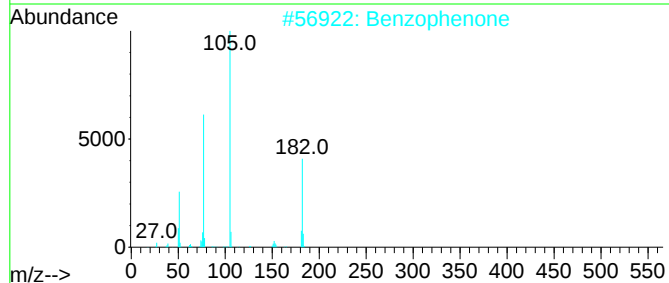
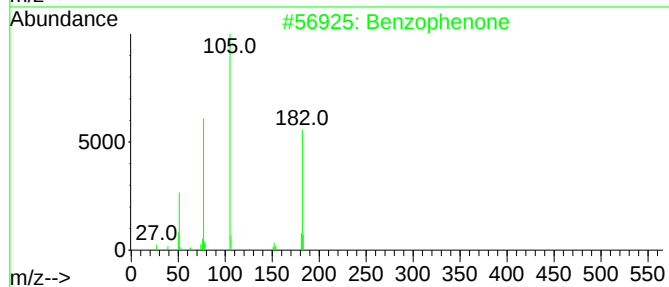
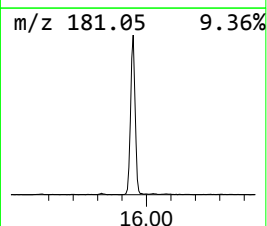
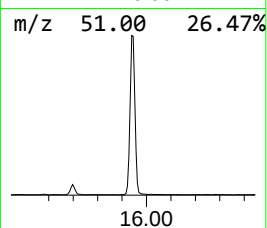
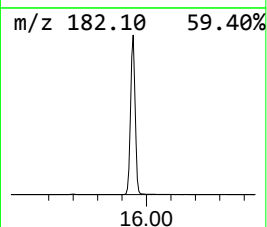
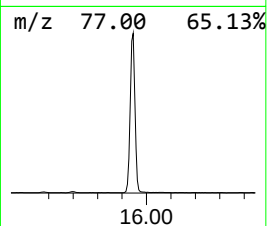
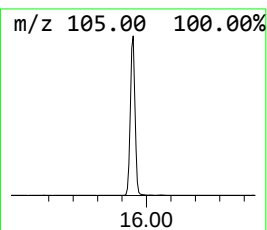
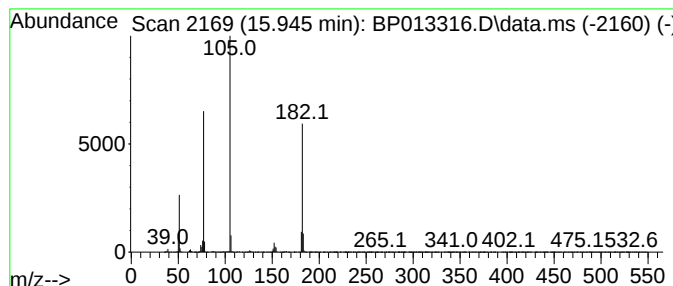
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.945	16.66 ng/ul	5261250	Acenaphthene-d10	14.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
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Operator : CG/JU
Sample : N6016-20
Misc :
ALS Vial : 28 Sample Multiplier: 1

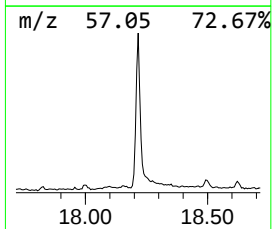
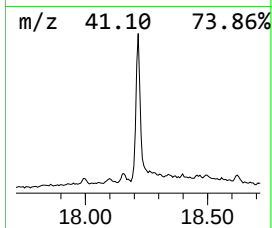
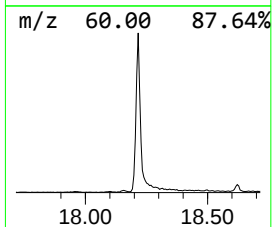
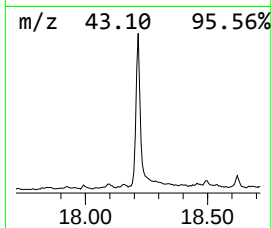
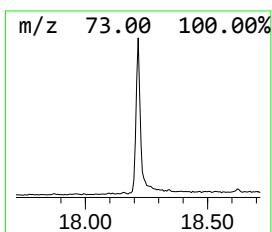
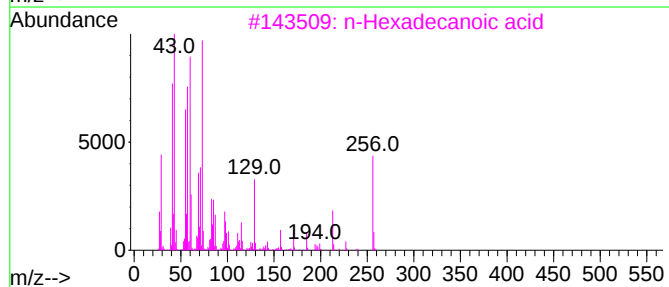
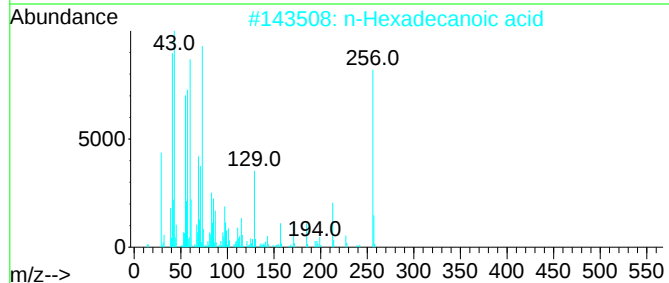
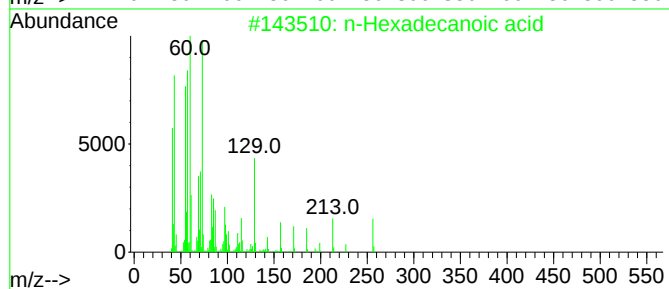
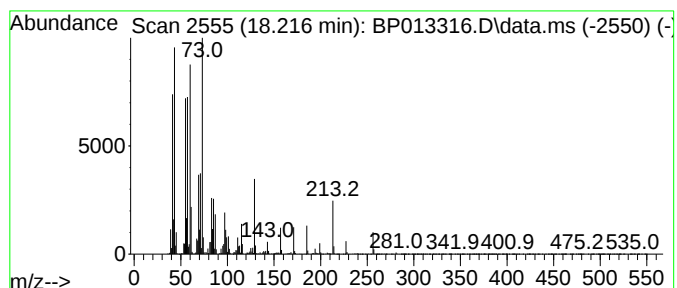
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.216	7.74 ng/ul	2359890	Phenanthrene-d10		17.345	
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	98
3	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	95
5	Tridecanoic acid		214	C13H26O2	000638-53-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013316.D
Acq On : 25 Dec 2022 00:50
Operator : CG/JU
Sample : N6016-20
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DBYE3

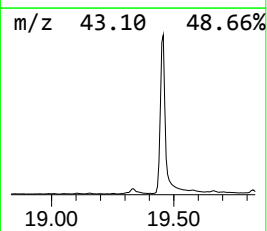
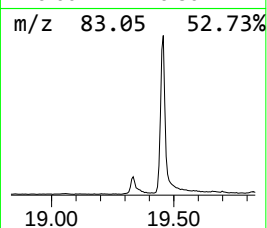
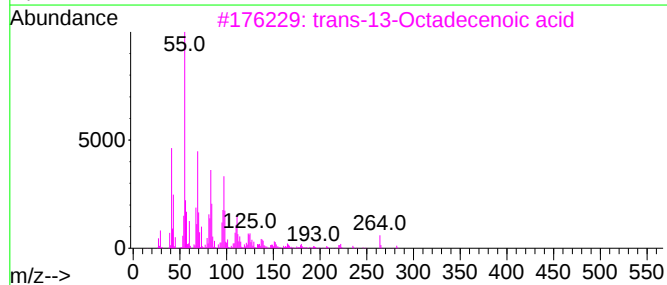
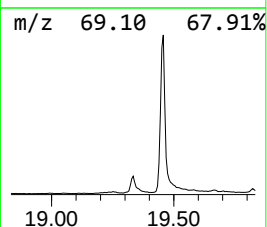
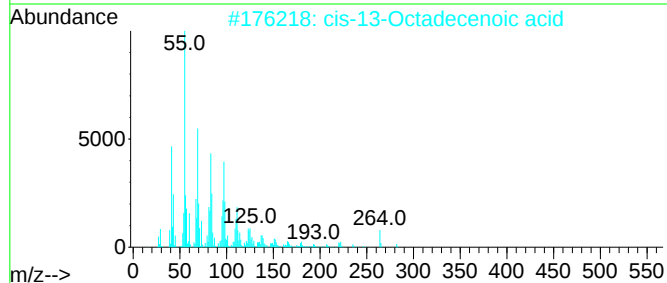
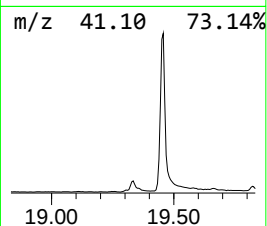
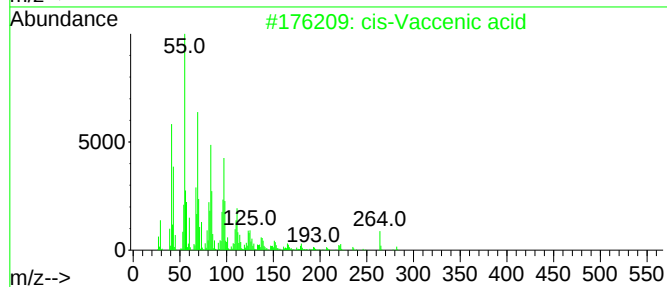
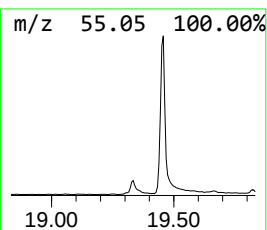
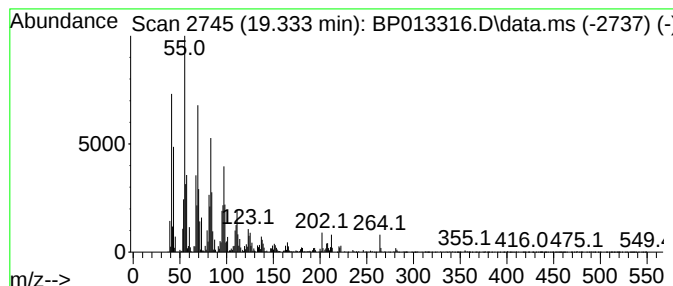
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 cis-Vaccenic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.333	5.53 ng/ul	1685520	Phenanthrene-d10	17.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-Vaccenic acid	282	C18H34O2	000506-17-2	98
2			cis-13-Octadecenoic acid	282	C18H34O2	013126-39-1	98
3			trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	98
4			9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	97
5			Oleic Acid	282	C18H34O2	000112-80-1	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013316.D
Acq On : 25 Dec 2022 00:50
Operator : CG/JU
Sample : N6016-20
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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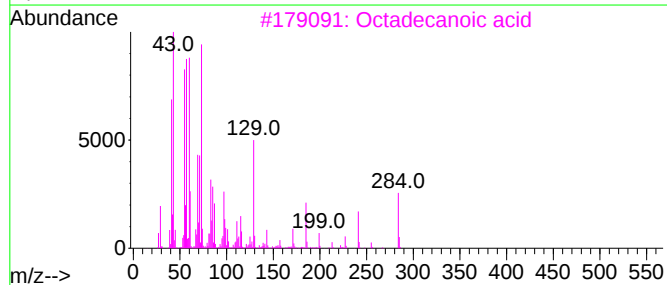
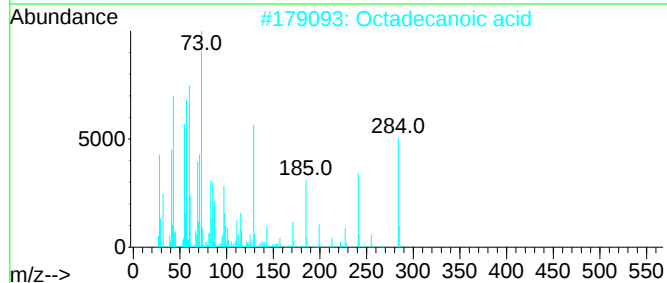
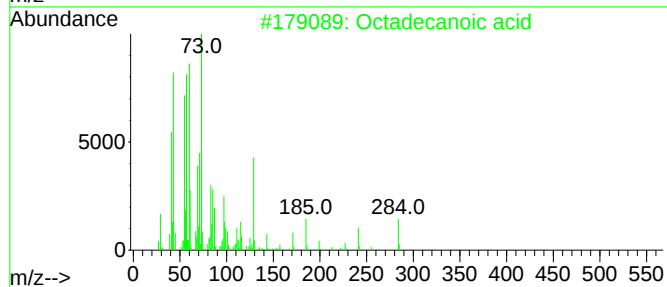
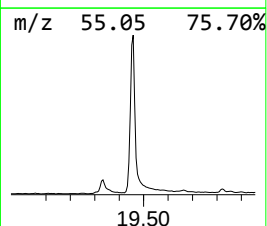
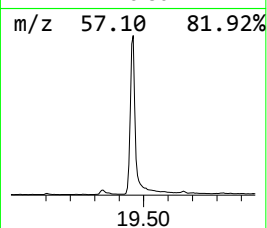
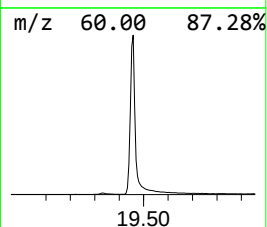
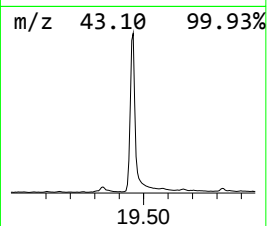
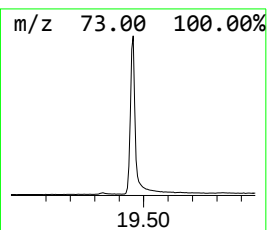
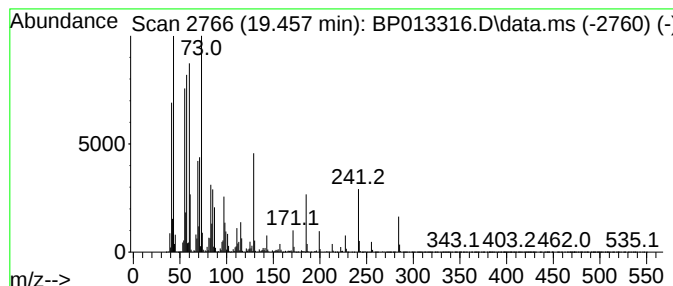
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.457	50.87 ng/ul	18085600	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	98
5			Octadecanoic acid	284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013316.D
Acq On : 25 Dec 2022 00:50
Operator : CG/JU
Sample : N6016-20
Misc :
ALS Vial : 28 Sample Multiplier: 1

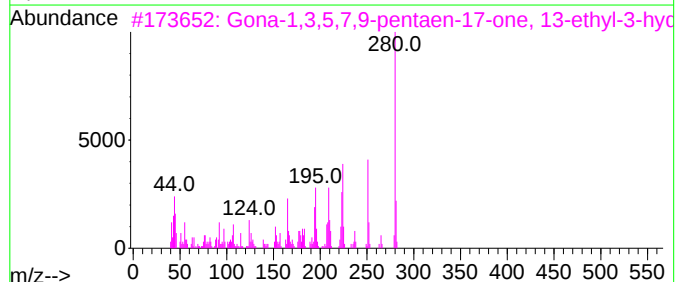
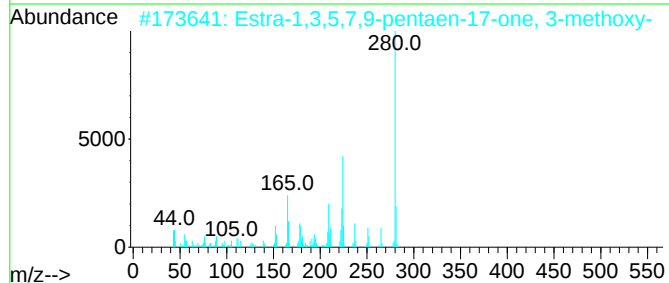
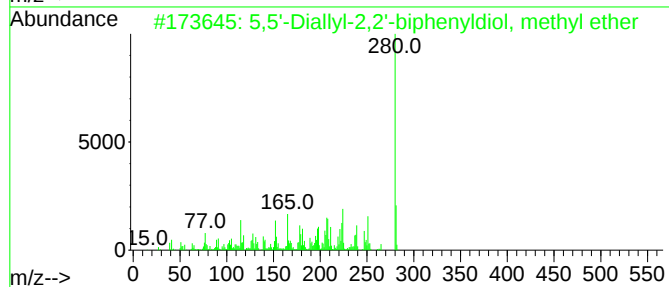
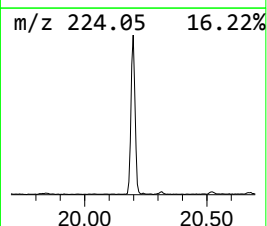
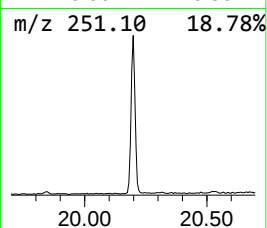
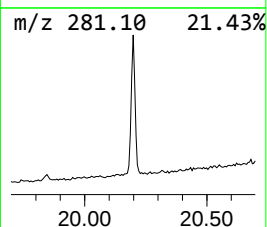
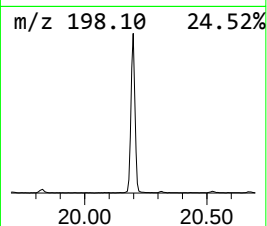
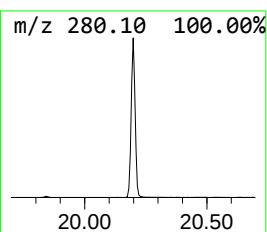
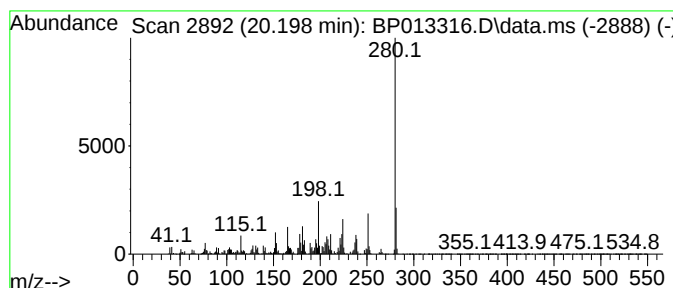
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 5,5'-Diallyl-2,2'-biphenyld... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.198	13.82 ng/ul	4912470	Chrysene-d12	21.433	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,5'-Diallyl-2,2'-biphenyldiol, ...	280	C19H20O2	1000384-18-6	83
2	Estra-1,3,5,7,9-pentaen-17-one, ...	280	C19H20O2	003907-67-3	74
3	Gona-1,3,5,7,9-pentaen-17-one, 1...	280	C19H20O2	030788-51-3	72
4	Gona-1,3,5,7,9-pentaen-17-one, 1...	280	C19H20O2	069853-73-2	64
5	6-Ethoxy-4-methyl-3-phenylcoumarin	280	C18H16O3	263365-04-4	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013316.D
Acq On : 25 Dec 2022 00:50
Operator : CG/JU
Sample : N6016-20
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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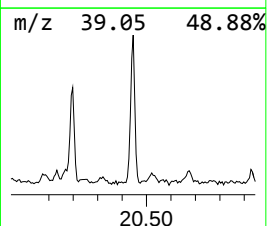
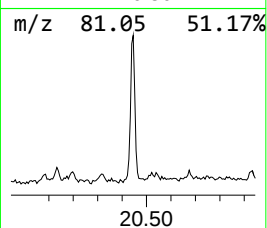
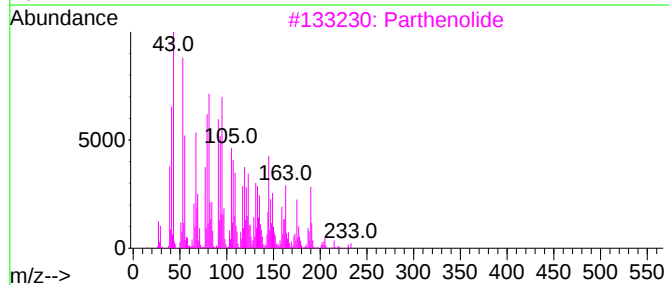
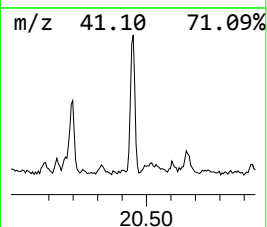
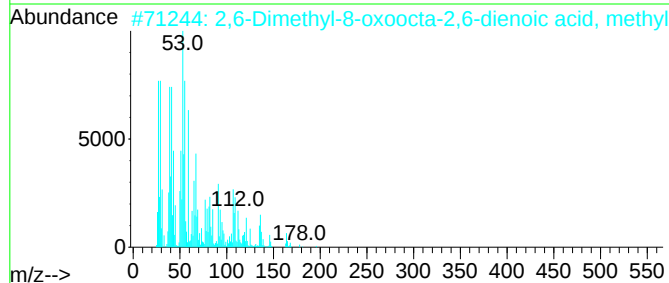
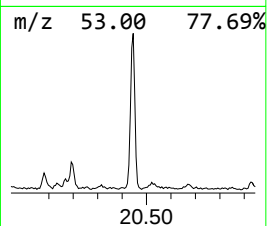
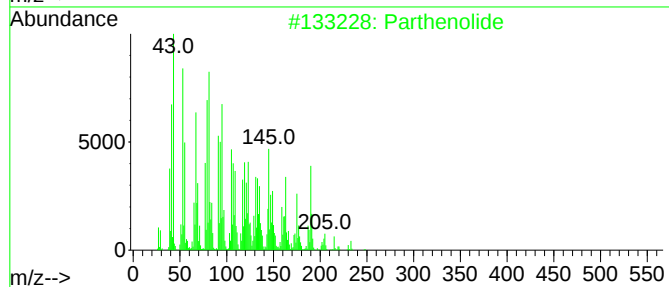
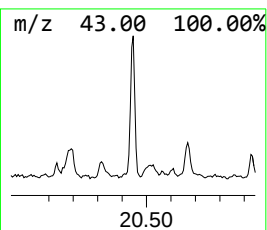
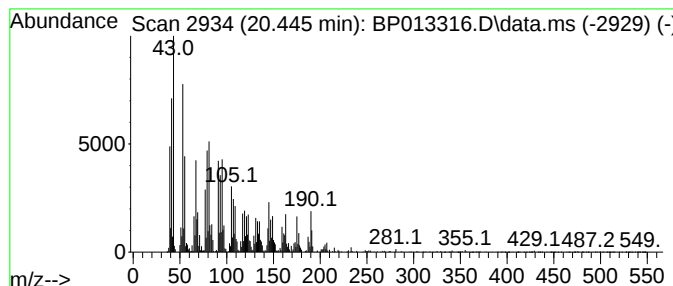
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Parthenolide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.445	3.82 ng/ul	1359170	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Parthenolide	248	C15H20O3	020554-84-1	99
2			2,6-Dimethyl-8-oxoocta-2,6-dieno...	196	C11H16O3	1000195-67-9	59
3			Parthenolide	248	C15H20O3	020554-84-1	50
4			Furan, 2-ethyl-	96	C6H8O	003208-16-0	49
5			2,4-Pentadienamide, N-phenyl-	173	C11H11NO	1000156-33-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013316.D
Acq On : 25 Dec 2022 00:50
Operator : CG/JU
Sample : N6016-20
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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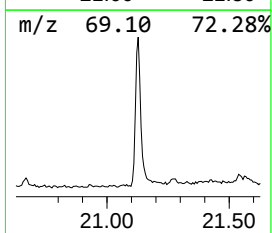
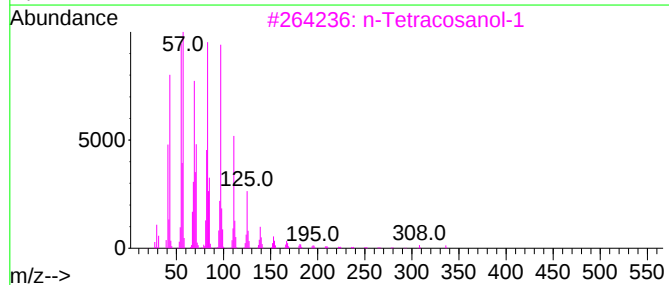
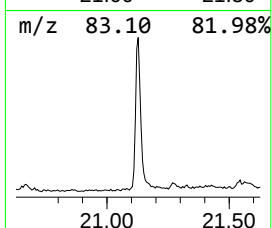
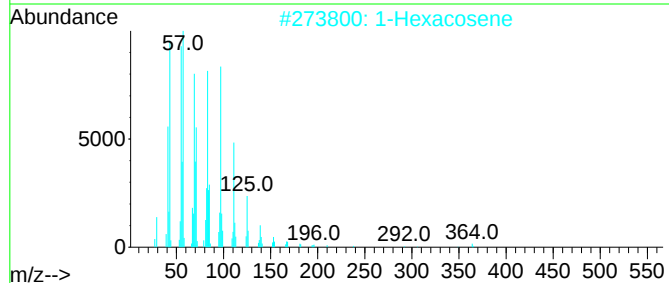
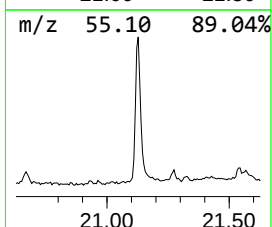
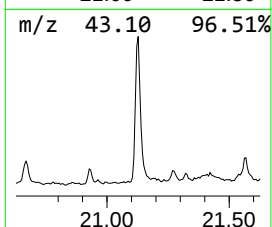
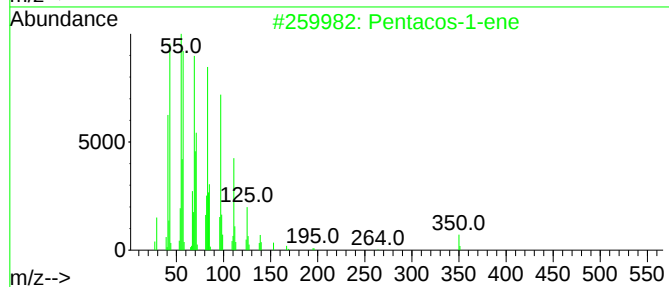
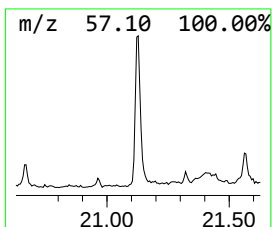
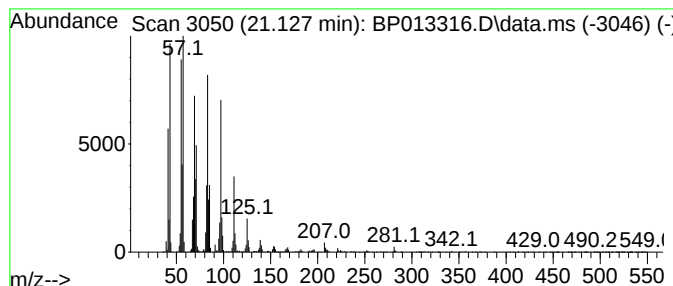
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.127	5.76 ng/ul	2046270	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentacos-1-ene	350	C25H50	016980-85-1	94
2			1-Hexacosene	364	C26H52	018835-33-1	94
3			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	94
5			Pentadecafluorooctanoic acid, pe...	624	C23H31F15O2	1000406-04-5	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013316.D
Acq On : 25 Dec 2022 00:50
Operator : CG/JU
Sample : N6016-20
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYE3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.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
2-Pentanone, 4-...	5.022	5.7	ng/ul	1145370	1	7.940	4026090	20.0
n-Decanoic acid	12.863	2.9	ng/ul	917228	3	14.581	6315930	20.0
Cyclohexane, 1-...	13.451	2.3	ng/ul	736178	3	14.581	6315930	20.0
Benzophenone	15.945	16.7	ng/ul	5261250	3	14.581	6315930	20.0
n-Hexadecanoic ...	18.216	7.7	ng/ul	2359890	4	17.345	6094490	20.0
cis-Vaccenic acid	19.333	5.5	ng/ul	1685520	4	17.345	6094490	20.0
Octadecanoic acid	19.457	50.9	ng/ul	18085600	5	21.433	7110320	20.0
4-(3-Fluoro-8-n...	19.822	13.4	ng/ul	4762310	5	21.433	7110320	20.0
5,5'-Diallyl-2,...	20.198	13.8	ng/ul	4912470	5	21.433	7110320	20.0
Parthenolide	20.445	3.8	ng/ul	1359170	5	21.433	7110320	20.0
(DEL) Alkane: S...	21.127	5.8	ng/ul	2046270	5	21.433	7110320	20.0