

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013136.D
 Acq On : 19 Dec 2022 20:38
 Operator : CG/JU
 Sample : N6006-14
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBYA0

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: BP013136.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.340	22	26	32	rVB	140459	186870	1.39%	0.116%
2	3.775	97	100	113	rBV	357242	557016	4.15%	0.345%
3	3.875	113	117	125	rVB	70791	100930	0.75%	0.062%
4	3.999	133	138	143	rVB	84419	126859	0.94%	0.078%
5	4.099	151	155	159	rBV	41347	51892	0.39%	0.032%
6	4.475	216	219	224	rVB4	22054	32329	0.24%	0.020%
7	5.040	308	315	322	rBV	403222	605906	4.51%	0.375%
8	6.628	579	585	593	rVB	87550	137223	1.02%	0.085%
9	7.111	658	667	679	rBV	2202481	3591466	26.73%	2.222%
10	7.281	689	696	708	rVV	2084165	3189866	23.75%	1.974%
11	7.481	723	730	738	rBV	2348584	3714424	27.65%	2.298%
12	7.569	740	745	752	rVB6	21368	40731	0.30%	0.025%
13	7.769	773	779	783	rBV2	10430	20080	0.15%	0.012%
14	7.958	802	811	818	rBV	2051646	3236530	24.09%	2.002%
15	8.016	818	821	826	rVB5	16497	25677	0.19%	0.016%
16	8.199	843	852	856	rBV5	7650	20025	0.15%	0.012%
17	8.663	920	931	942	rBV	2347147	3731698	27.78%	2.309%
18	8.787	946	952	960	rVB	104944	176703	1.32%	0.109%
19	9.122	1002	1009	1023	rBV	2416338	3926016	29.23%	2.429%
20	9.228	1023	1027	1031	rVV6	16060	28652	0.21%	0.018%
21	9.275	1031	1035	1042	rVB8	19033	35095	0.26%	0.022%
22	9.528	1069	1078	1086	rVV2	76606	123112	0.92%	0.076%
23	9.846	1122	1132	1146	rBV	1895194	3179921	23.67%	1.967%
24	9.981	1151	1155	1162	rVV6	22746	49606	0.37%	0.031%
25	10.075	1165	1171	1176	rVB10	20041	37878	0.28%	0.023%
26	10.387	1217	1224	1235	rBV	2972744	4904744	36.51%	3.035%
27	10.769	1279	1289	1296	rBV	2957159	4851522	36.11%	3.002%
28	10.905	1305	1312	1334	rVB	1494573	2683665	19.98%	1.660%
29	11.299	1375	1379	1387	rVB	13607	28237	0.21%	0.017%
30	11.557	1415	1423	1428	rBV10	18919	43100	0.32%	0.027%
31	11.675	1440	1443	1453	rVB10	8161	19260	0.14%	0.012%
32	11.787	1453	1462	1469	rBV9	12028	24014	0.18%	0.015%
33	12.181	1523	1529	1535	rBV	49939	76491	0.57%	0.047%
34	12.293	1538	1548	1553	rBV2	29516	61896	0.46%	0.038%
35	12.351	1553	1558	1564	rVB2	60156	89727	0.67%	0.056%
36	12.457	1567	1576	1586	rBV	92626	180297	1.34%	0.112%

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

37	12.634	1601	1606	1614	rVB	8825	22393	0.17%	0.014%
38	13.128	1685	1690	1693	rBV3	16880	21528	0.16%	0.013%
39	13.410	1734	1738	1744	rVB	91259	125569	0.93%	0.078%
40	13.487	1746	1751	1754	rBV7	13499	27746	0.21%	0.017%
41	13.557	1759	1763	1772	rVB9	19581	51055	0.38%	0.032%
42	13.916	1818	1824	1830	rBV10	24242	43726	0.33%	0.027%
43	14.004	1832	1839	1848	rBV	5638119	7667847	57.08%	4.744%
44	14.069	1848	1850	1854	rVB4	26173	26981	0.20%	0.017%
45	14.116	1854	1858	1864	rVB6	34596	58554	0.44%	0.036%
46	14.187	1864	1870	1874	rVB	25016	36724	0.27%	0.023%
47	14.240	1874	1879	1881	rBV5	16958	27582	0.21%	0.017%
48	14.293	1881	1888	1896	rBV	6259841	9285561	69.12%	5.745%
49	14.440	1907	1913	1916	rBV8	13321	28260	0.21%	0.017%
50	14.481	1916	1920	1925	rVB	75252	107772	0.80%	0.067%
51	14.598	1929	1940	1947	rBV	4569525	6785957	50.51%	4.199%
52	14.657	1947	1950	1964	rVB	38083	91652	0.68%	0.057%
53	14.787	1966	1972	1989	rBV	1453681	2411456	17.95%	1.492%
54	14.951	1996	2000	2003	rVB3	30166	41145	0.31%	0.025%
55	15.010	2005	2010	2020	rVB9	48094	112049	0.83%	0.069%
56	15.387	2068	2074	2077	rBV5	17890	35621	0.27%	0.022%
57	15.440	2079	2083	2087	rVB2	25176	35331	0.26%	0.022%
58	15.504	2090	2094	2101	rVB6	34012	52890	0.39%	0.033%
59	15.592	2102	2109	2116	rBV	8308709	11921897	88.75%	7.376%
60	15.704	2122	2128	2145	rBV	2265235	3204608	23.86%	1.983%
61	15.828	2145	2149	2155	rVB2	42085	71274	0.53%	0.044%
62	15.957	2165	2171	2180	rBV	2817752	3855972	28.70%	2.386%
63	16.163	2202	2206	2210	rVB7	30738	42929	0.32%	0.027%
64	16.304	2226	2230	2237	rVB8	25259	57041	0.42%	0.035%
65	16.481	2257	2260	2263	rVV3	23025	34015	0.25%	0.021%
66	16.522	2263	2267	2275	rVB2	65722	104248	0.78%	0.064%
67	16.739	2300	2304	2307	rBV5	23139	31196	0.23%	0.019%
68	16.887	2325	2329	2337	rVB4	32135	44244	0.33%	0.027%
69	17.139	2368	2372	2375	rBV3	16445	25762	0.19%	0.016%
70	17.234	2384	2388	2391	rBV4	42004	57611	0.43%	0.036%
71	17.298	2396	2399	2402	rBV5	25643	31349	0.23%	0.019%
72	17.351	2402	2408	2414	rVV	5608767	7995325	59.52%	4.947%
73	17.392	2414	2415	2419	rVV	103215	106246	0.79%	0.066%
74	17.451	2419	2425	2435	rVV2	8082670	11730613	87.32%	7.258%
75	17.534	2437	2439	2448	rVB6	62650	117065	0.87%	0.072%
76	17.634	2453	2456	2460	rVV4	21306	27493	0.20%	0.017%

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 Title : SVOA CALIBRATION

77	17.728	2469	2472	2475	rVV4	19087	21112	0.16%	0.013%
78	17.839	2488	2491	2495	rBV5	21445	30934	0.23%	0.019%
79	18.010	2517	2520	2524	rBV4	34185	44838	0.33%	0.028%
80	18.110	2532	2537	2542	rBV2	62447	92594	0.69%	0.057%
81	18.163	2542	2546	2551	rBV2	118555	170657	1.27%	0.106%
82	18.228	2551	2557	2566	rBV	1770332	2569783	19.13%	1.590%
83	18.504	2601	2604	2609	rBV4	48535	82484	0.61%	0.051%
84	18.557	2610	2613	2619	rVV7	38810	53448	0.40%	0.033%
85	18.634	2622	2626	2631	rBV2	168562	211610	1.58%	0.131%
86	18.704	2635	2638	2641	rVV3	48720	60429	0.45%	0.037%
87	19.110	2702	2707	2714	rBV6	110685	237763	1.77%	0.147%
88	19.263	2729	2733	2738	rVB2	133129	183197	1.36%	0.113%
89	19.345	2738	2747	2761	rBV3	1710013	3855631	28.70%	2.386%
90	19.457	2761	2766	2779	rVB	609122	1061538	7.90%	0.657%
91	19.686	2799	2805	2817	rVB	9953198	13433645	100.00%	8.312%
92	19.833	2826	2830	2835	rVB	421714	462023	3.44%	0.286%
93	20.192	2884	2891	2894	rBV4	251373	615756	4.58%	0.381%
94	21.127	3046	3050	3056	rBV	1941283	2874146	21.40%	1.778%
95	21.439	3098	3103	3108	rBV	4878599	6379981	47.49%	3.947%
96	23.133	3387	3391	3397	rBV	1165083	1936778	14.42%	1.198%
97	23.757	3491	3497	3512	rVB2	4289016	9223964	68.66%	5.707%
98	23.921	3518	3525	3540	rVB2	2741664	5899417	43.92%	3.650%
99	24.545	3625	3631	3639	rVB	1234405	2668307	19.86%	1.651%
100	26.457	3950	3956	3972	rVB	987282	3010391	22.41%	1.863%

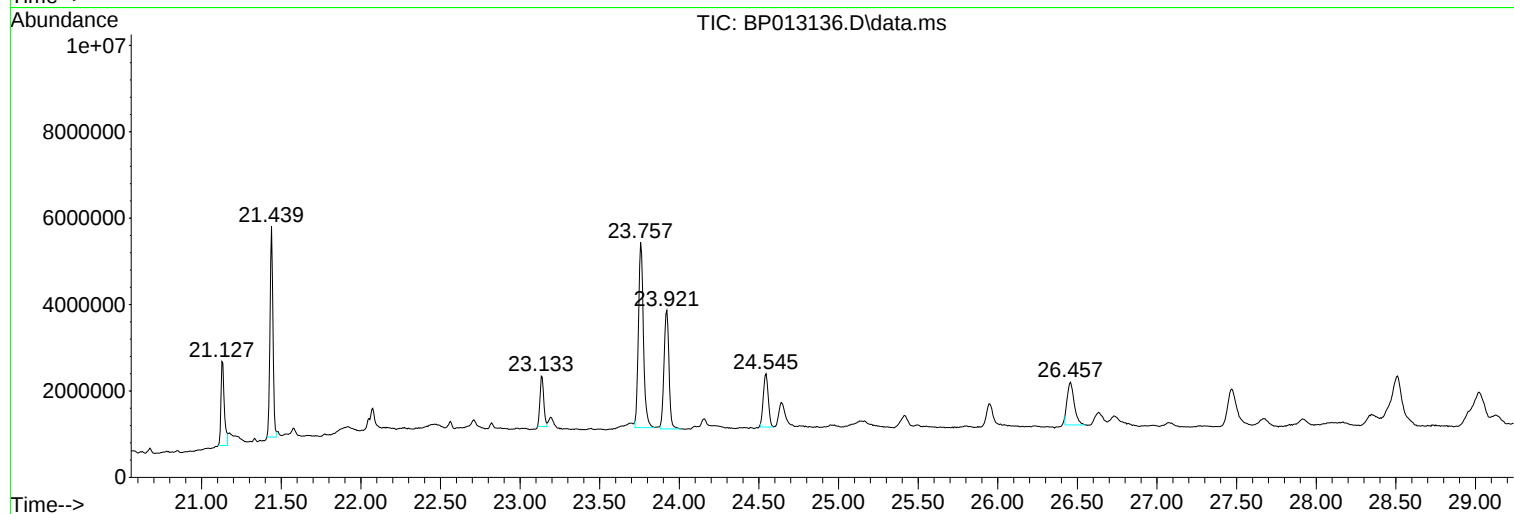
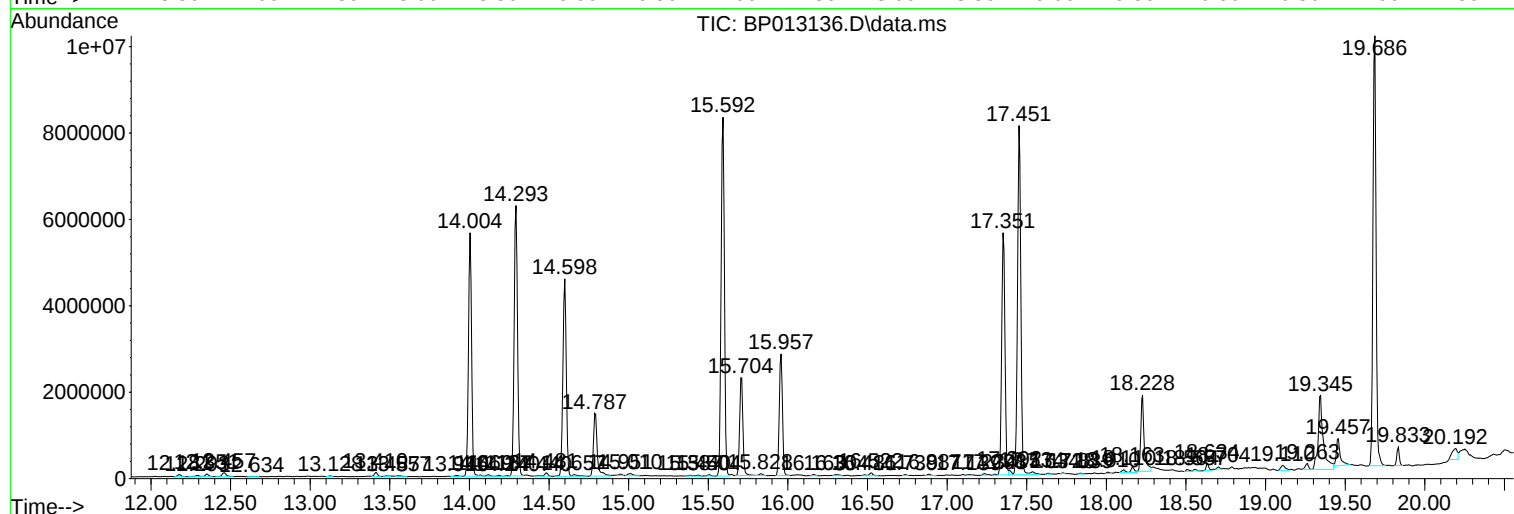
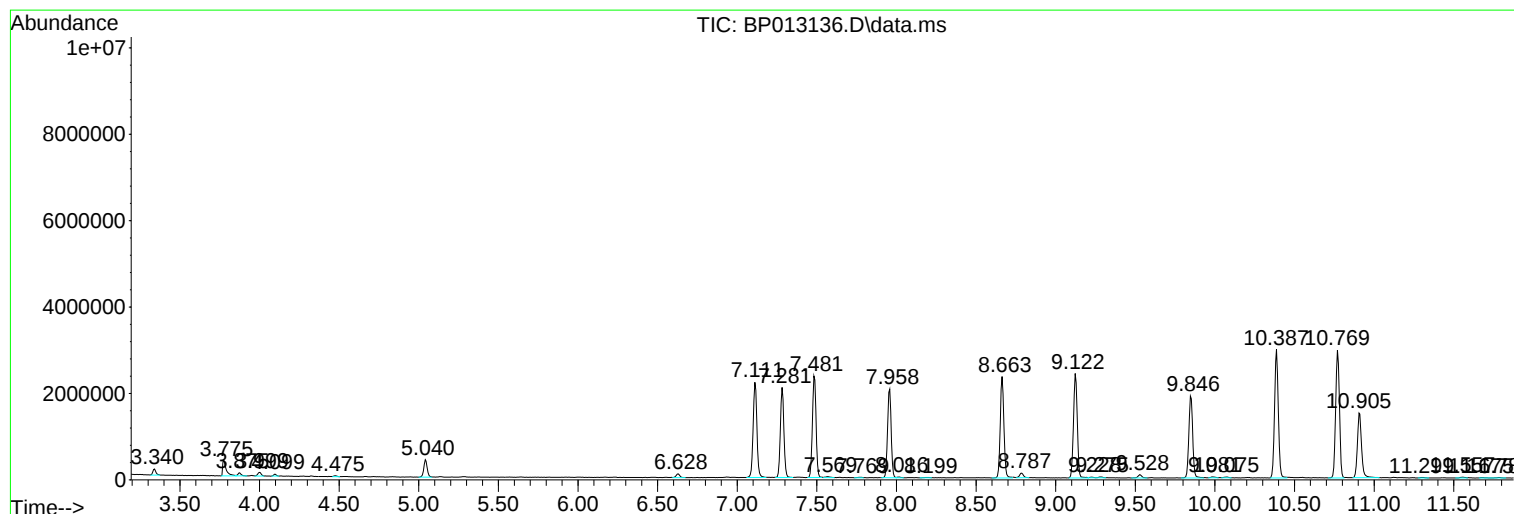
Sum of corrected areas: 161626171

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Instrument :
BNA_P
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



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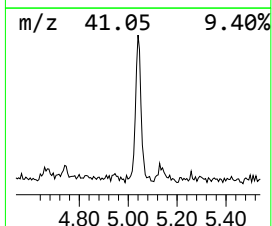
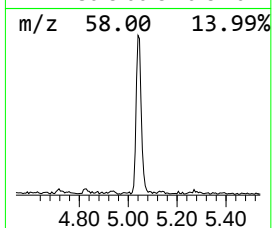
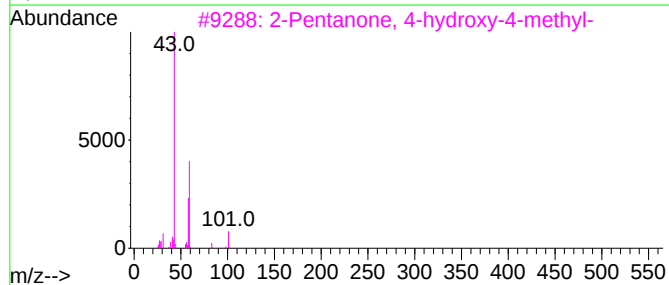
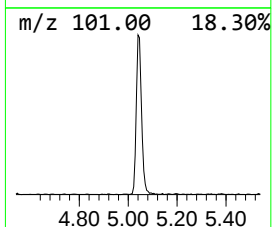
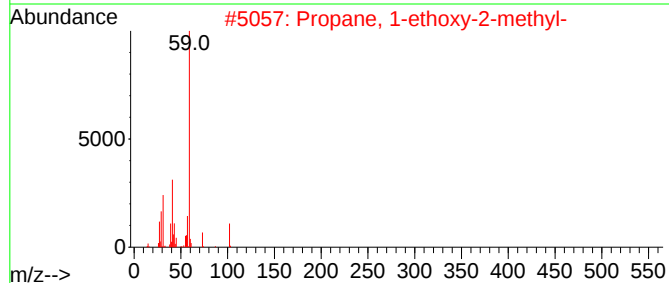
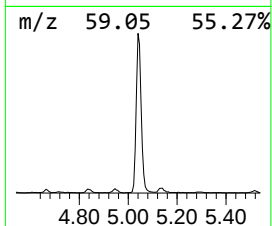
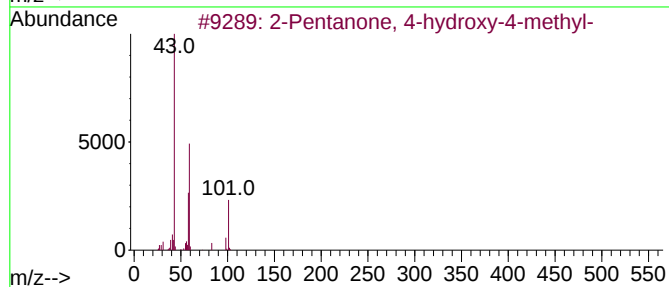
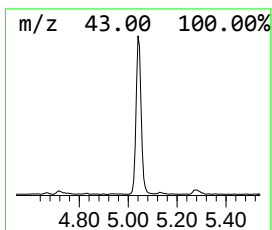
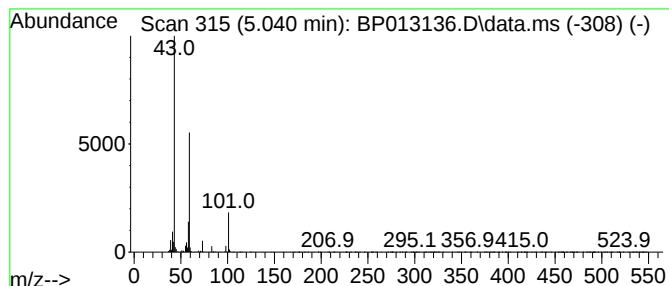
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.040	3.74 ng/ul	605906	1,4-Dichlorobenzene-d4	7.958

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	46		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	28		
5	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25		



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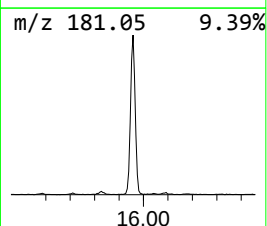
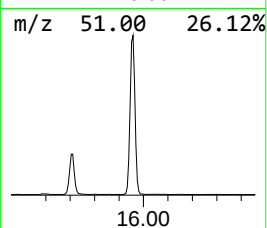
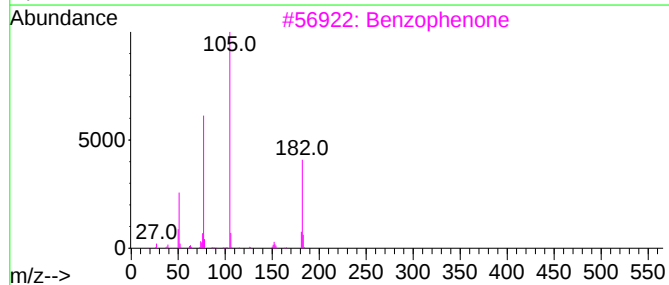
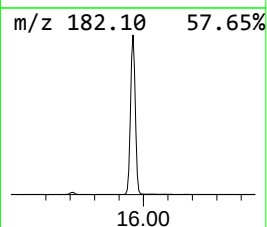
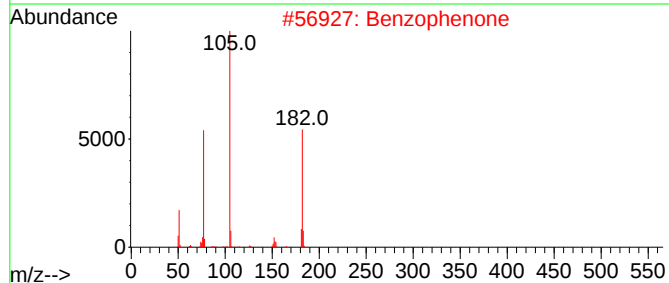
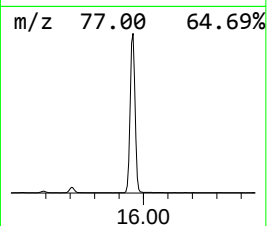
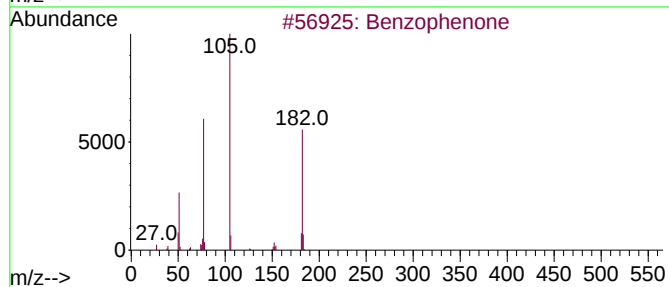
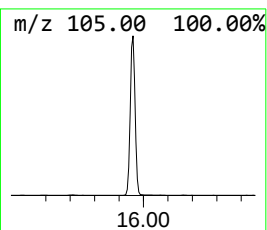
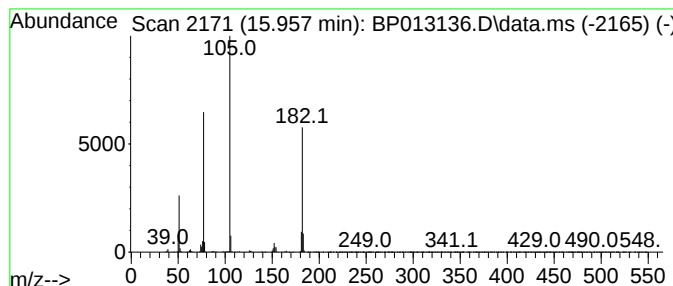
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.957	11.36 ng/ul	3855970	Acenaphthene-d10	14.598

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	94
4			Benzophenone	182	C13H10O	000119-61-9	93
5			Azobenzene	182	C12H10N2	000103-33-3	64



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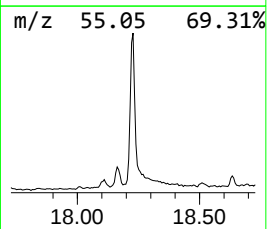
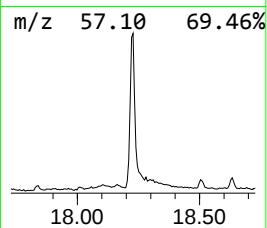
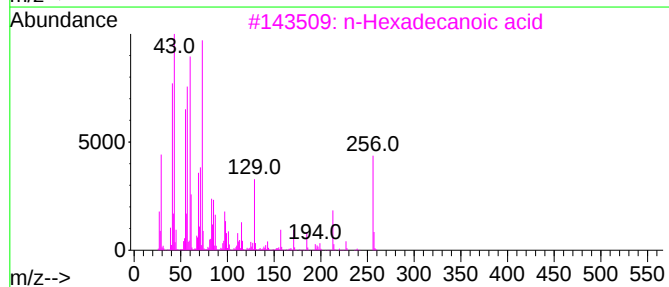
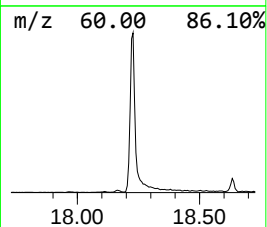
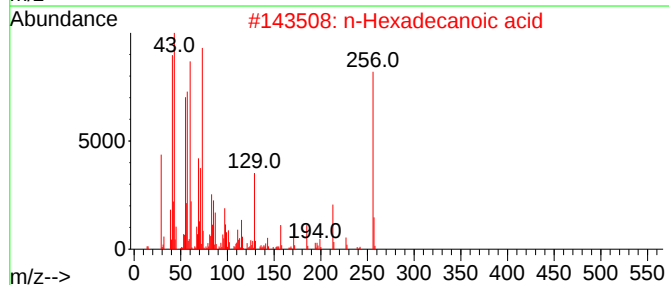
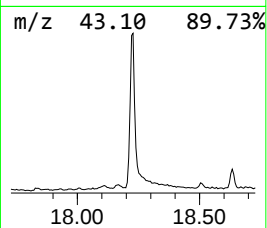
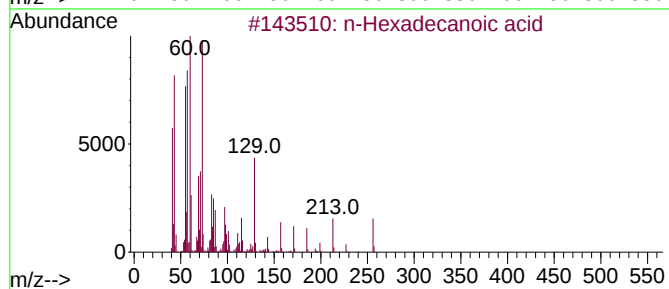
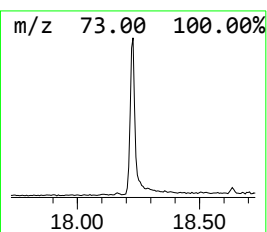
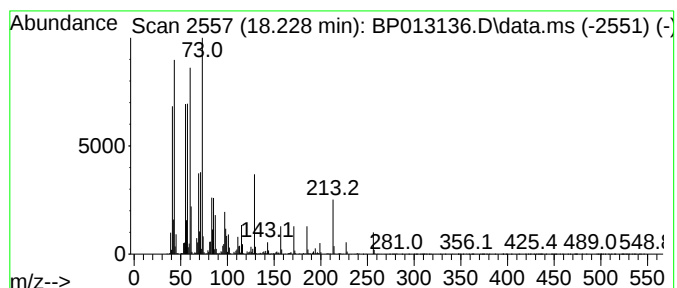
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.228	6.43 ng/ul	2569780	Phenanthrene-d10	17.351

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	97		
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013136.D
Acq On : 19 Dec 2022 20:38
Operator : CG/JU
Sample : N6006-14
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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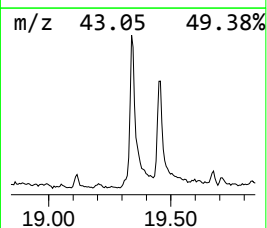
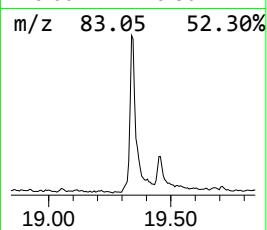
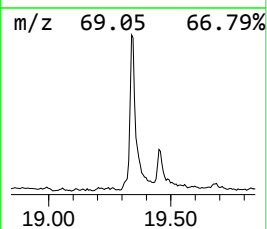
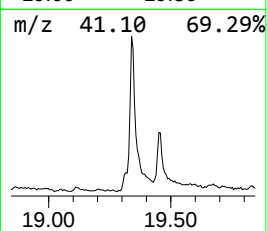
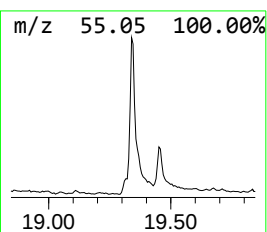
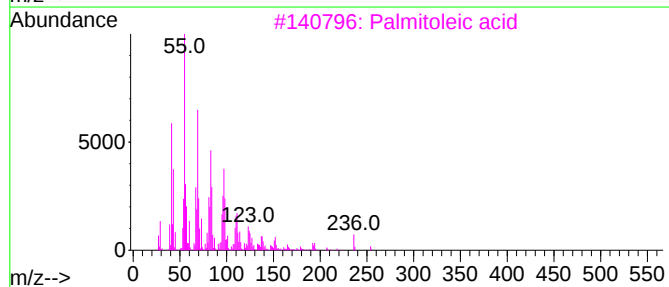
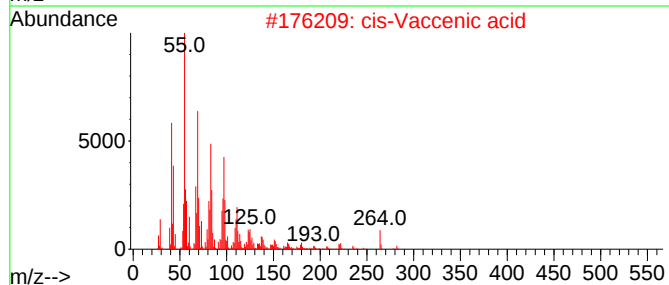
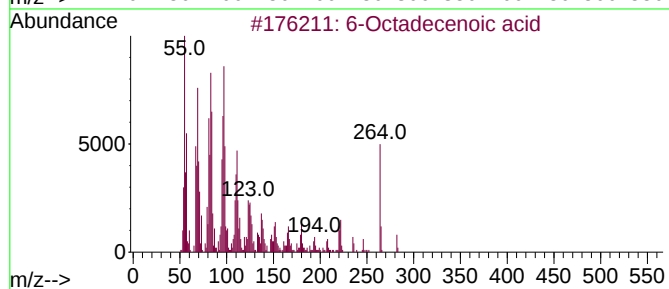
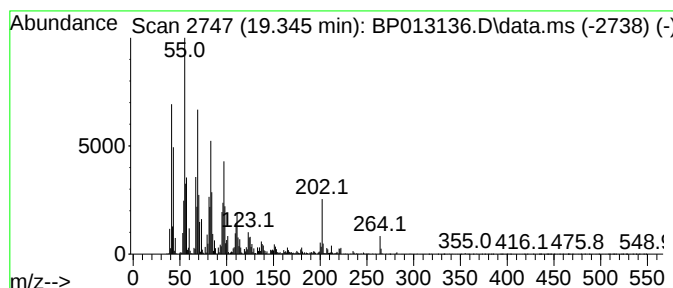
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 6-Octadecenoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.345	9.64 ng/ul	3855630	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Octadecenoic acid	282	C18H34O2	1000336-66-8	93
2			cis-Vaccenic acid	282	C18H34O2	000506-17-2	91
3			Palmitoleic acid	254	C16H30O2	000373-49-9	87
4			cis-13-Octadecenoic acid	282	C18H34O2	013126-39-1	87
5			Oleic Acid	282	C18H34O2	000112-80-1	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013136.D
Acq On : 19 Dec 2022 20:38
Operator : CG/JU
Sample : N6006-14
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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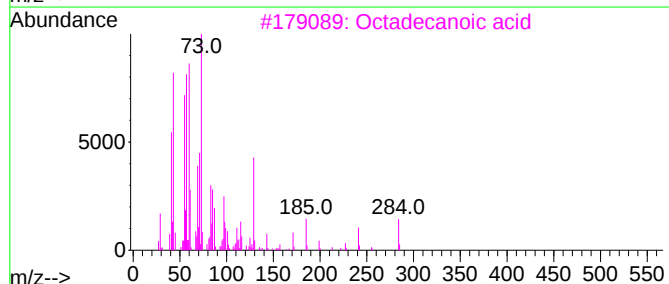
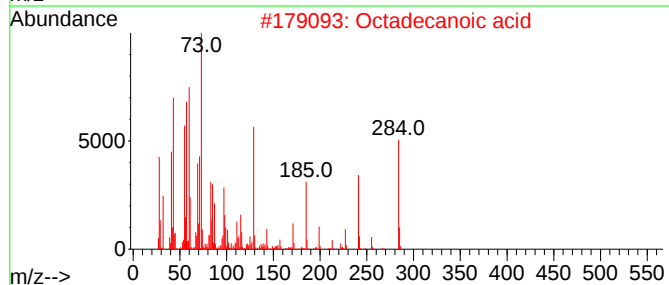
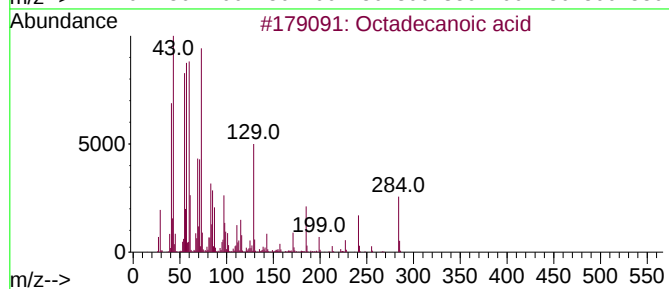
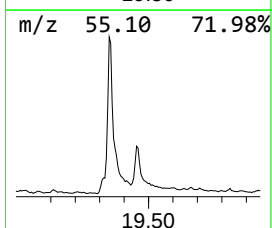
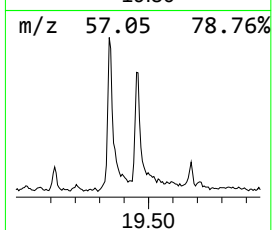
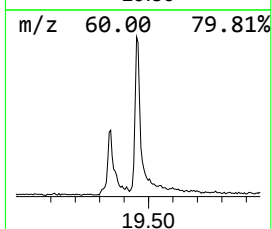
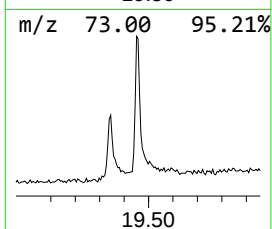
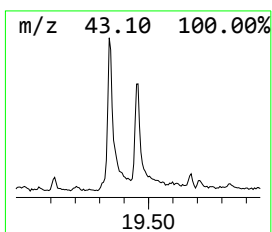
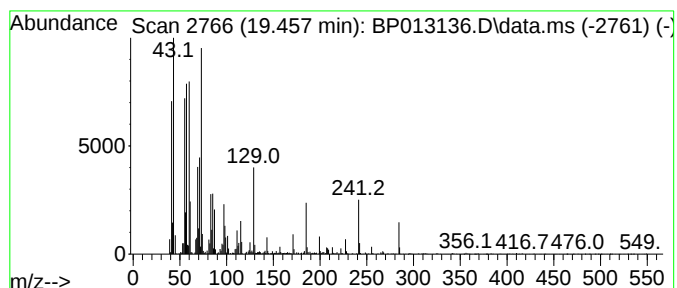
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.457	3.33 ng/ul	1061540	Chrysene-d12	21.439

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\BP121922\
Data File : BP013136.D
Acq On : 19 Dec 2022 20:38
Operator : CG/JU
Sample : N6006-14
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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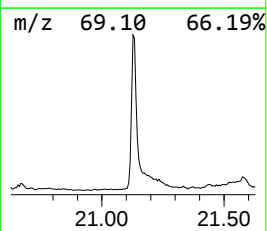
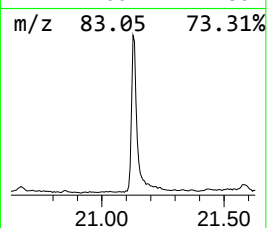
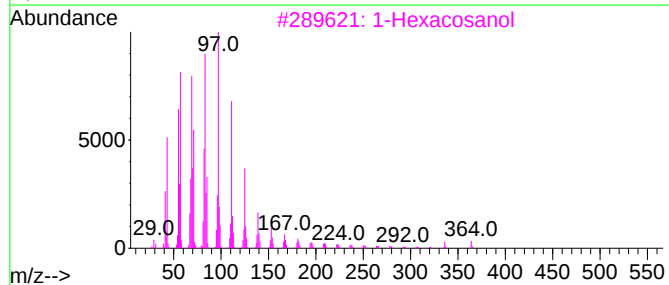
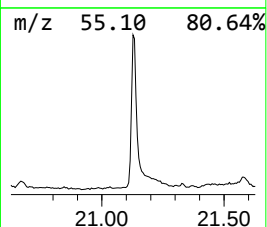
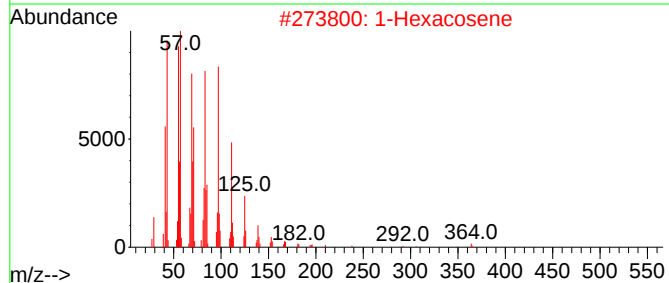
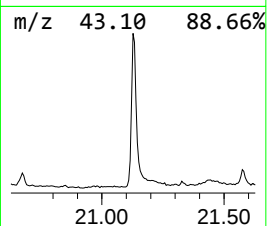
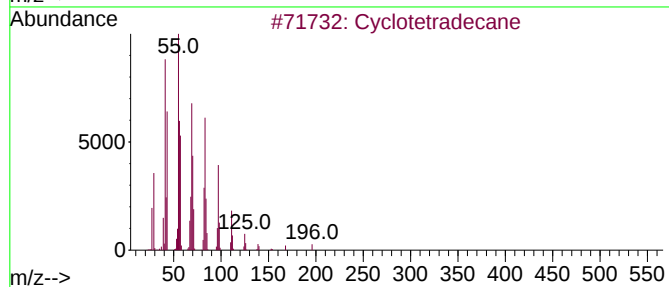
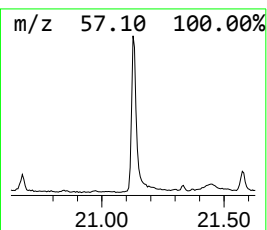
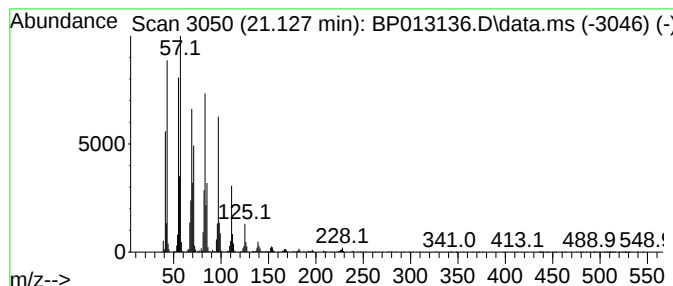
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic21.128 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.128	9.01 ng/ul	2874150	Chrysene-d12	21.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetradecane	196	C14H28	000295-17-0	96
2			1-Hexacosene	364	C26H52	018835-33-1	94
3			1-Hexacosanol	382	C26H54O	000506-52-5	91
4			1-Heptacosanol	396	C27H56O	002004-39-9	91
5			Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013136.D
Acq On : 19 Dec 2022 20:38
Operator : CG/JU
Sample : N6006-14
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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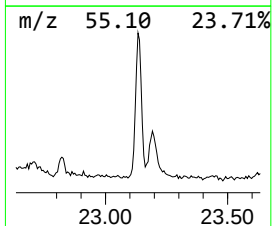
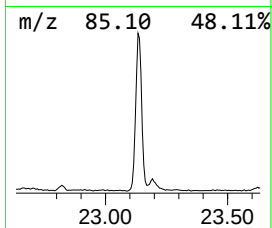
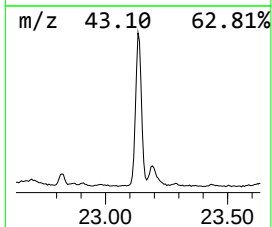
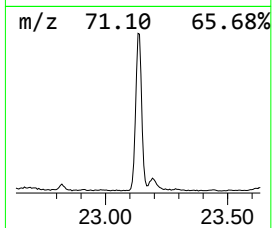
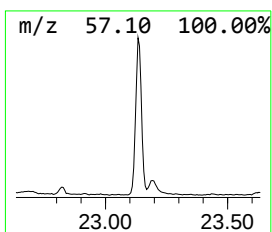
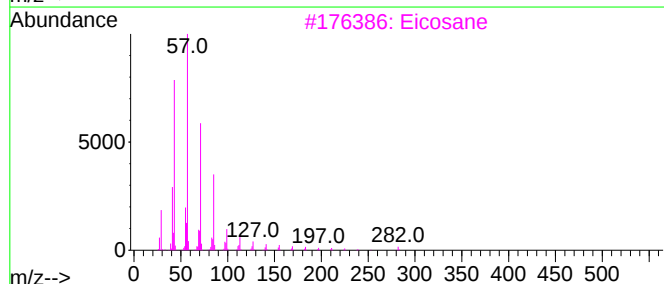
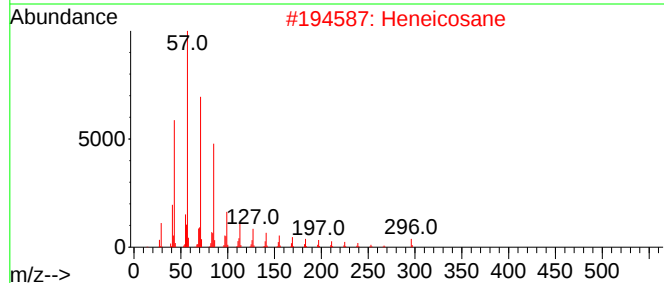
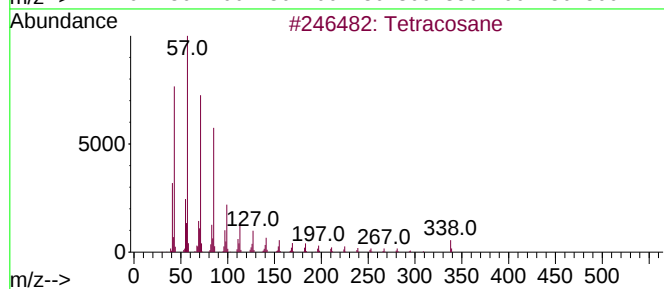
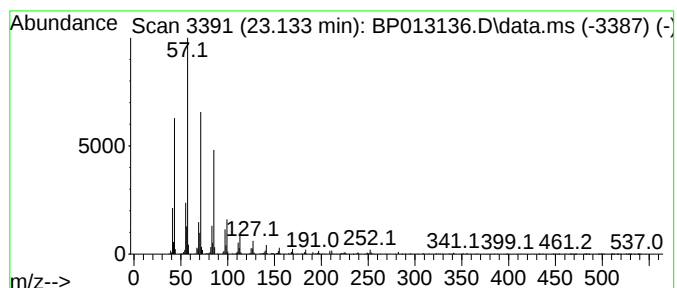
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.133	6.57 ng/ul	1936780	Perylene-d12	23.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	97
2			Heneicosane	296	C21H44	000629-94-7	97
3			Eicosane	282	C20H42	000112-95-8	96
4			Eicosane	282	C20H42	000112-95-8	94
5			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013136.D
Acq On : 19 Dec 2022 20:38
Operator : CG/JU
Sample : N6006-14
Misc :
ALS Vial : 15 Sample Multiplier: 1

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ClientSampleId :
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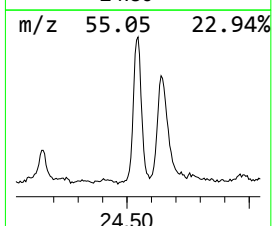
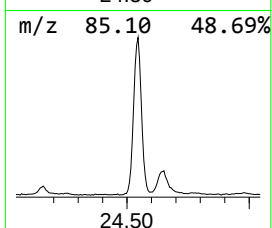
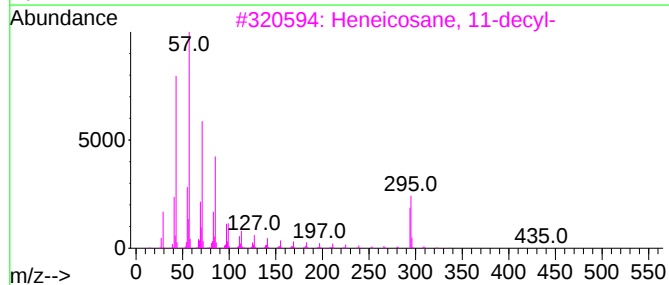
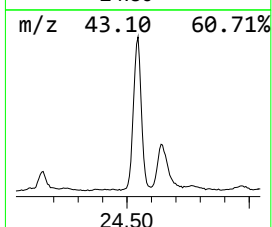
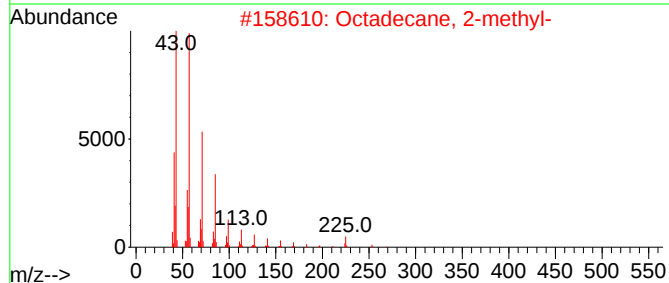
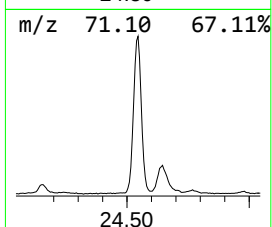
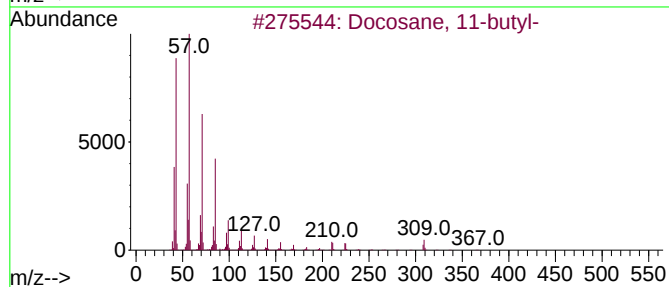
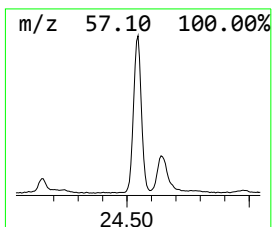
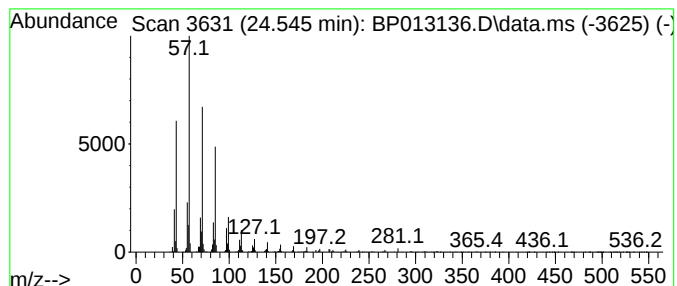
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.545	9.05 ng/ul	2668310	Perylene-d12	23.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-			366	C26H54	013475-76-8	95
2	Octadecane, 2-methyl-			268	C19H40	001560-88-9	94
3	Heneicosane, 11-decyl-			437	C31H64	055320-06-4	93
4	Heneicosane, 11-pentyl-			366	C26H54	014739-72-1	93
5	Heneicosane			296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013136.D
Acq On : 19 Dec 2022 20:38
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Sample : N6006-14
Misc :
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Instrument :
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ClientSampleId :
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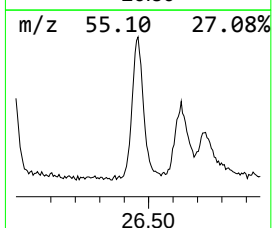
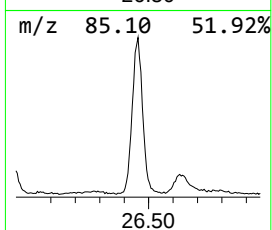
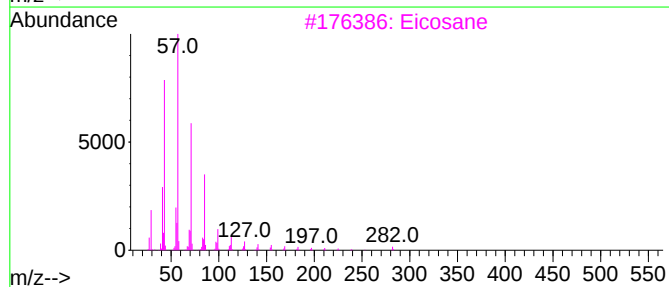
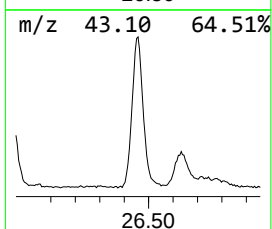
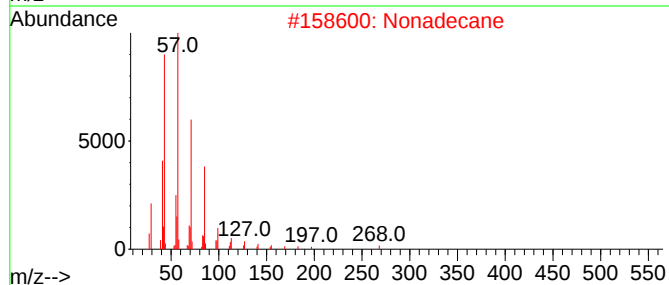
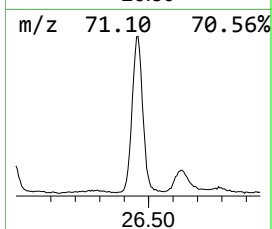
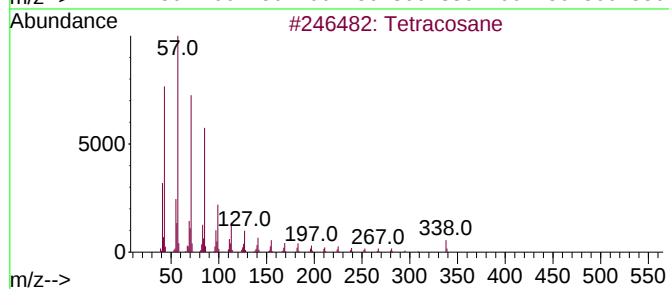
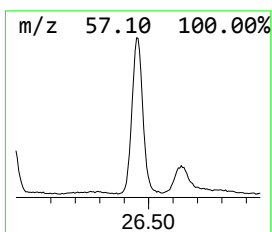
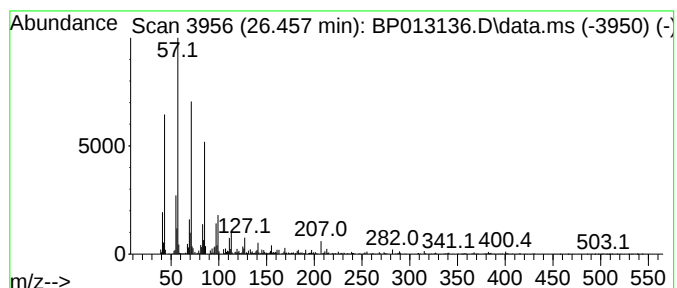
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.457	10.21 ng/ul	3010390	Perylene-d12	23.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	96
2			Nonadecane	268	C19H40	000629-92-5	94
3			Eicosane	282	C20H42	000112-95-8	93
4			Tetracosane	338	C24H50	000646-31-1	93
5			Heptacosane	380	C27H56	000593-49-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013136.D
Acq On : 19 Dec 2022 20:38
Operator : CG/JU
Sample : N6006-14
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYA0

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

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					#	RT	Resp	Conc
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Benzophenone	15.957	11.4	ng/u1	3855970	3	14.598	6785960	20.0
n-Hexadecanoic ...	18.228	6.4	ng/u1	2569780	4	17.351	7995330	20.0
6-Octadecenoic ...	19.345	9.6	ng/u1	3855630	4	17.351	7995330	20.0
Octadecanoic acid	19.457	3.3	ng/u1	1061540	5	21.439	6379980	20.0
(DEL) Alkane: C...	21.128	9.0	ng/u1	2874150	5	21.439	6379980	20.0
(DEL) Alkane: S...	23.133	6.6	ng/u1	1936780	6	23.922	5899420	20.0
(DEL) Alkane: S...	24.545	9.1	ng/u1	2668310	6	23.922	5899420	20.0
(DEL) Alkane: S...	26.457	10.2	ng/u1	3010390	6	23.922	5899420	20.0