

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
 Data File : BP023546.D
 Acq On : 20 Dec 2024 04:07
 Operator : RC/JU
 Sample : P5265-10
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E28Z5

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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

Signal : TIC: BP023546.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.093	14	18	30	rVB	46210	76328	1.42%	0.099%
2	3.511	85	89	107	rBV	382632	703638	13.08%	0.914%
3	5.575	436	440	445	rVB3	22898	36355	0.68%	0.047%
4	5.822	475	482	489	rBV4	23138	62550	1.16%	0.081%
5	6.028	511	517	524	rBV8	16432	43761	0.81%	0.057%
6	6.093	524	528	533	rBV	24278	41253	0.77%	0.054%
7	6.211	539	548	556	rVB5	36260	95007	1.77%	0.123%
8	6.416	576	583	590	rBV4	15585	41345	0.77%	0.054%
9	6.622	613	618	621	rBV5	34523	56872	1.06%	0.074%
10	6.664	621	625	632	rVB3	65182	131767	2.45%	0.171%
11	6.740	632	638	653	rBV	533330	1250944	23.26%	1.625%
12	6.875	655	661	670	rVB	552721	903047	16.79%	1.173%
13	7.005	679	683	688	rBV4	19933	38760	0.72%	0.050%
14	7.075	689	695	703	rVV	670318	1102568	20.50%	1.432%
15	7.169	708	711	717	rVB2	27375	40721	0.76%	0.053%
16	7.428	746	755	759	rBV5	36534	108517	2.02%	0.141%
17	7.487	759	765	766	rVV2	72140	99116	1.84%	0.129%
18	7.522	766	771	778	rVB	894132	1456270	27.08%	1.891%
19	7.628	784	789	794	rBV7	33364	65894	1.23%	0.086%
20	7.687	794	799	802	rBV5	34389	47694	0.89%	0.062%
21	7.728	802	806	810	rBV	56568	81517	1.52%	0.106%
22	8.028	845	857	864	rBV4	77296	209115	3.89%	0.272%
23	8.158	871	879	887	rBV2	65644	158478	2.95%	0.206%
24	8.240	887	893	903	rBV	686602	1425576	26.51%	1.852%
25	8.358	909	913	917	rVB4	26290	40484	0.75%	0.053%
26	8.587	943	952	959	rVB6	37443	99365	1.85%	0.129%
27	8.663	959	965	976	rBV	600452	1198946	22.29%	1.557%
28	8.793	982	987	988	rBV4	27680	39471	0.73%	0.051%
29	8.828	989	993	1000	rVB7	56915	119189	2.22%	0.155%
30	8.934	1007	1011	1017	rVB6	20733	46447	0.86%	0.060%
31	9.040	1024	1029	1035	rVB3	50672	92877	1.73%	0.121%
32	9.222	1057	1060	1067	rVB2	46595	70781	1.32%	0.092%
33	9.375	1075	1086	1094	rBV	481313	984549	18.31%	1.279%
34	9.440	1094	1097	1105	rVB5	39095	82757	1.54%	0.107%
35	9.534	1107	1113	1118	rVB6	22574	44030	0.82%	0.057%
36	9.663	1129	1135	1143	rBV5	47515	112107	2.08%	0.146%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
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37	9.775	1150	1154	1159	rVB3	20616	37442	0.70%	0.049%
38	9.922	1172	1179	1194	rBV	614354	1511619	28.10%	1.963%
39	10.263	1224	1237	1243	rBV	1275209	2311241	42.97%	3.002%
40	10.310	1243	1245	1252	rVB3	40992	65015	1.21%	0.084%
41	10.428	1258	1265	1281	rBV	467608	1064945	19.80%	1.383%
42	10.822	1324	1332	1343	rVB	16217	52009	0.97%	0.068%
43	11.099	1372	1379	1385	rBV5	19215	44586	0.83%	0.058%
44	11.628	1461	1469	1474	rBV2	38502	77662	1.44%	0.101%
45	12.934	1685	1691	1698	rVV	42092	68175	1.27%	0.089%
46	13.169	1726	1731	1739	rBV3	25444	58475	1.09%	0.076%
47	13.551	1786	1796	1803	rBV	1521317	2224177	41.35%	2.889%
48	13.828	1833	1843	1861	rBV	1526804	2577342	47.92%	3.348%
49	14.028	1871	1877	1883	rVB	32950	58831	1.09%	0.076%
50	14.134	1888	1895	1904	rBV	1933840	3426810	63.71%	4.451%
51	14.451	1943	1949	1966	rBV2	217610	865633	16.09%	1.124%
52	14.587	1969	1972	1978	rVB3	36553	60741	1.13%	0.079%
53	15.004	2037	2043	2046	rVB2	37830	50924	0.95%	0.066%
54	15.151	2060	2068	2075	rBV	1945670	3061682	56.92%	3.977%
55	15.322	2091	2097	2106	rBV	254041	473835	8.81%	0.615%
56	15.534	2125	2133	2138	rBV	364622	575516	10.70%	0.747%
57	15.575	2138	2140	2147	rVB3	26963	39081	0.73%	0.051%
58	15.887	2190	2193	2195	rBV2	36427	42739	0.79%	0.056%
59	16.075	2217	2225	2228	rBV2	30364	69748	1.30%	0.091%
60	16.145	2232	2237	2241	rVB	122730	168483	3.13%	0.219%
61	16.222	2246	2250	2259	rBV5	21542	53596	1.00%	0.070%
62	16.351	2267	2272	2276	rBV3	41910	65708	1.22%	0.085%
63	16.698	2327	2331	2337	rBV3	39540	81232	1.51%	0.106%
64	16.857	2353	2358	2365	rVB3	44593	74934	1.39%	0.097%
65	16.939	2365	2372	2378	rBV	2852176	4624408	85.98%	6.006%
66	16.987	2378	2380	2384	rVV	303642	371876	6.91%	0.483%
67	17.045	2384	2390	2401	rVV2	2170200	3494230	64.97%	4.538%
68	17.128	2401	2404	2409	rVB7	34348	51835	0.96%	0.067%
69	17.375	2440	2446	2453	rBV4	38785	93708	1.74%	0.122%
70	17.516	2468	2470	2475	rVB	45216	58370	1.09%	0.076%
71	17.557	2475	2477	2483	rVB2	27360	42654	0.79%	0.055%
72	17.751	2505	2510	2516	rBV2	67847	144144	2.68%	0.187%
73	17.816	2516	2521	2527	rVV2	306918	631751	11.75%	0.821%
74	17.881	2527	2532	2548	rVB	1886438	3095959	57.56%	4.021%
75	18.051	2558	2561	2565	rVB3	53099	67025	1.25%	0.087%
76	18.169	2578	2581	2587	rBV3	53124	96075	1.79%	0.125%

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77	18.375	2614	2616	2623	rVB3	37162	60119	1.12%	0.078%
78	18.451	2625	2629	2635	rBV7	33912	79446	1.48%	0.103%
79	18.663	2661	2665	2672	rBV	133406	208806	3.88%	0.271%
80	18.810	2687	2690	2692	rBV3	50728	56175	1.04%	0.073%
81	18.892	2701	2704	2710	rVB	225385	262601	4.88%	0.341%
82	19.081	2729	2736	2746	rBV	817976	2014584	37.46%	2.617%
83	19.216	2755	2759	2772	rBV	781843	1272976	23.67%	1.653%
84	19.386	2785	2788	2790	rVB	88964	79592	1.48%	0.103%
85	19.428	2790	2795	2799	rBV	2589623	4004550	74.45%	5.201%
86	19.498	2803	2807	2821	rVB	3018114	4812468	89.48%	6.251%
87	19.851	2864	2867	2873	rVB	329909	416206	7.74%	0.541%
88	20.151	2915	2918	2921	rBV4	121182	139351	2.59%	0.181%
89	20.198	2922	2926	2931	rVV2	548080	621657	11.56%	0.807%
90	20.322	2944	2947	2950	rBV	170795	187923	3.49%	0.244%
91	20.410	2960	2962	2978	rVB2	378143	701306	13.04%	0.911%
92	20.528	2978	2982	2987	rBV	622935	791264	14.71%	1.028%
93	20.933	3047	3051	3058	rBV2	369416	542593	10.09%	0.705%
94	21.045	3065	3070	3079	rBV3	718659	1249523	23.23%	1.623%
95	21.269	3105	3108	3115	rVB	652614	962749	17.90%	1.250%
96	21.363	3118	3124	3130	rBV2	3264216	5378490	100.00%	6.986%
97	21.733	3184	3187	3195	rVB2	366225	473483	8.80%	0.615%
98	22.739	3353	3358	3371	rVB	1300333	2687395	49.97%	3.490%
99	24.321	3621	3627	3635	rVB	1150975	2343490	43.57%	3.044%
100	24.527	3656	3662	3671	rBV	1891748	4403181	81.87%	5.719%

Sum of corrected areas: 76992240

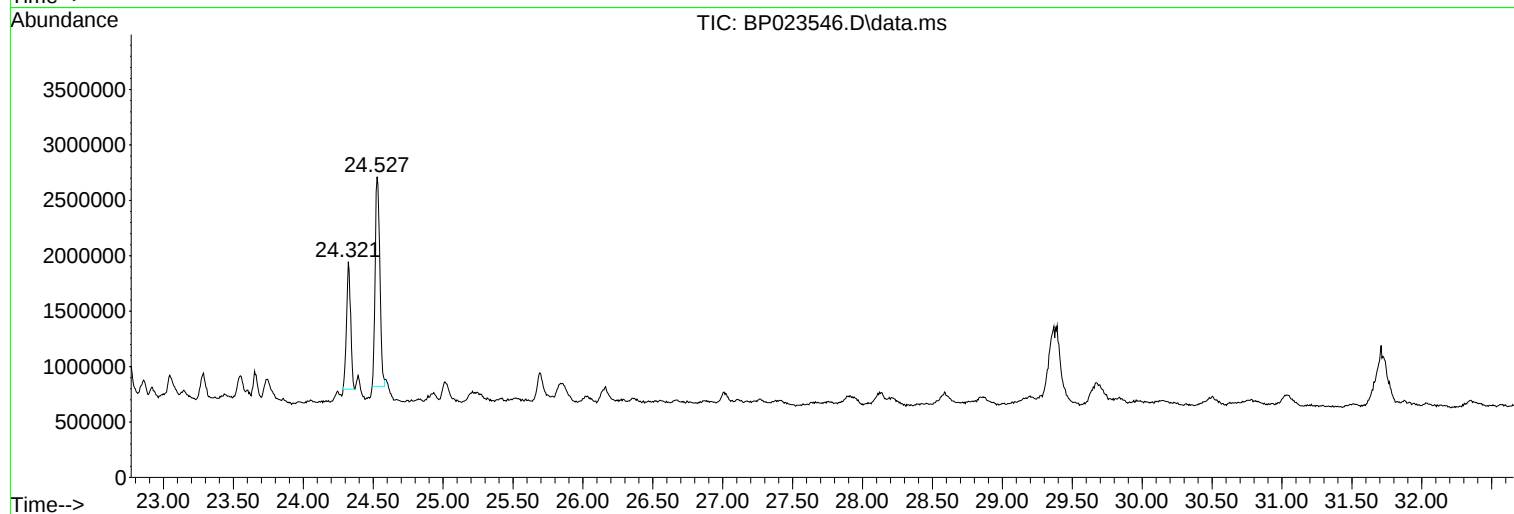
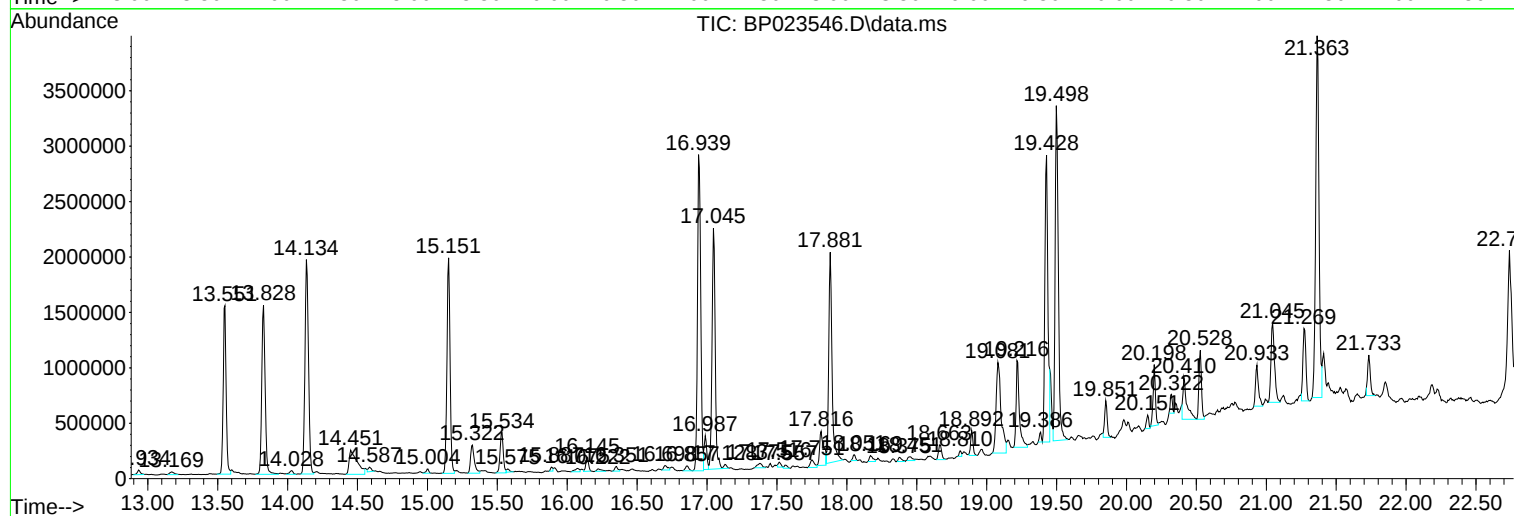
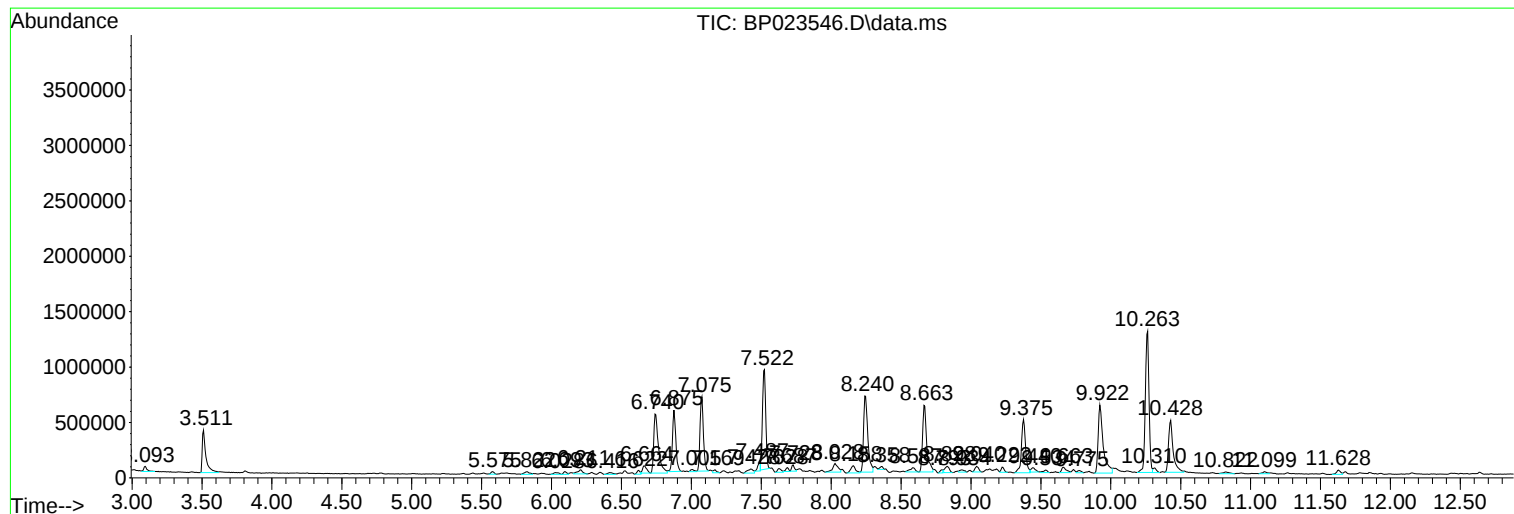
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P



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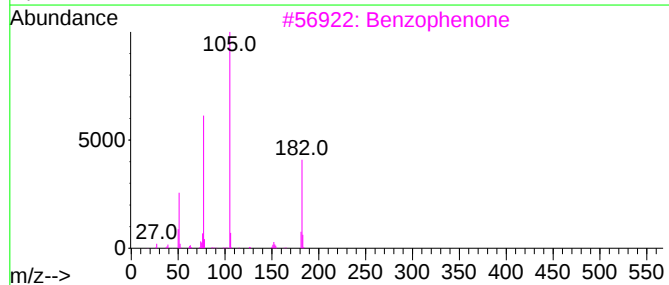
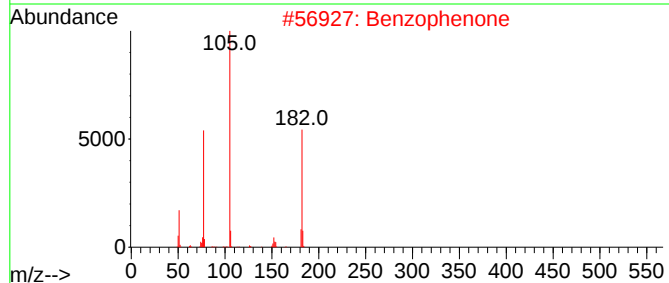
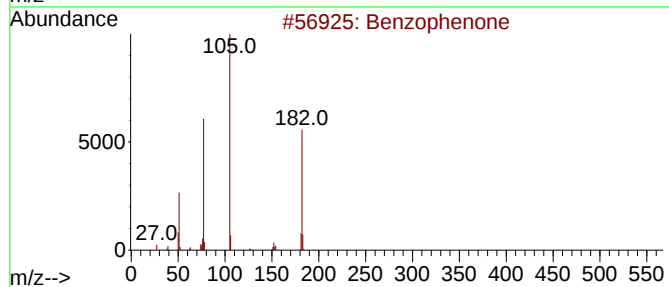
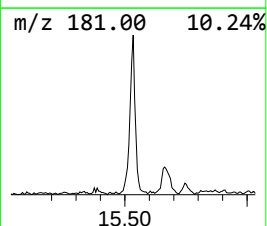
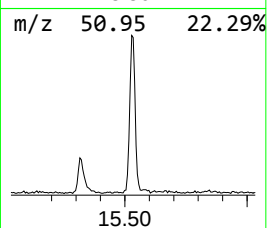
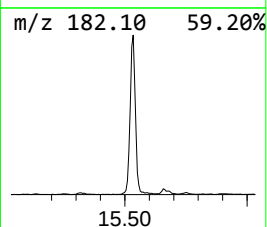
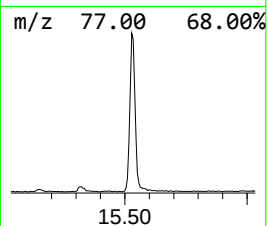
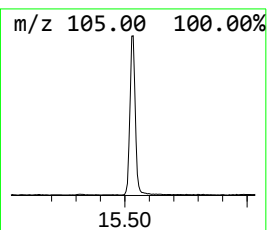
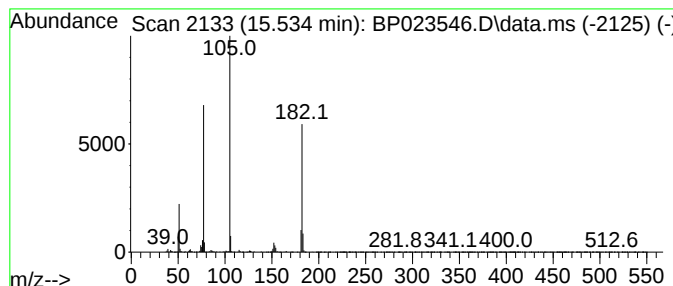
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.534	3.36 ng/ul	575516	Acenaphthene-d10	14.134

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



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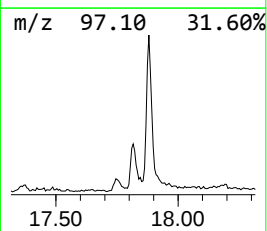
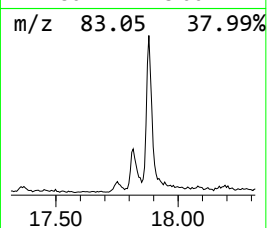
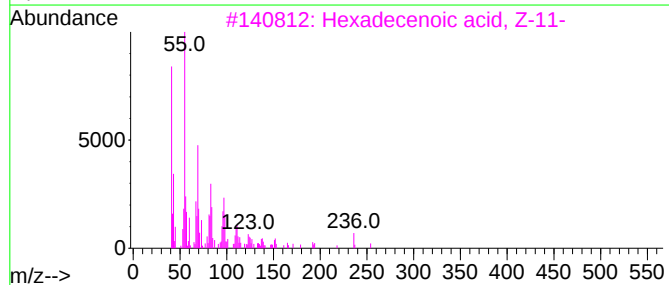
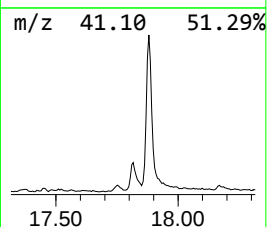
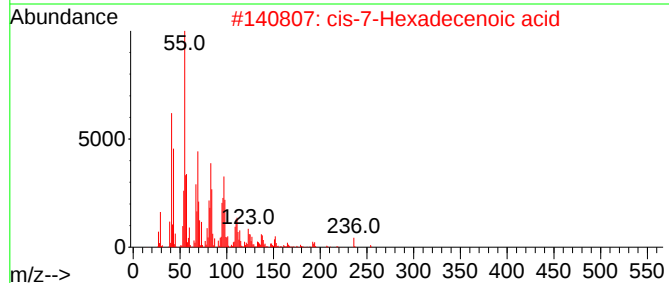
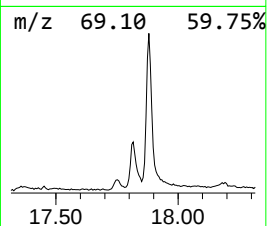
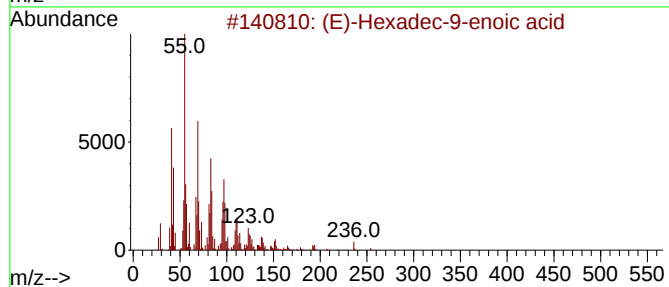
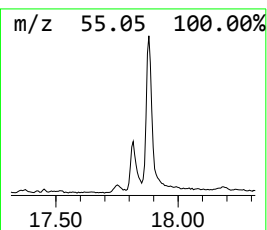
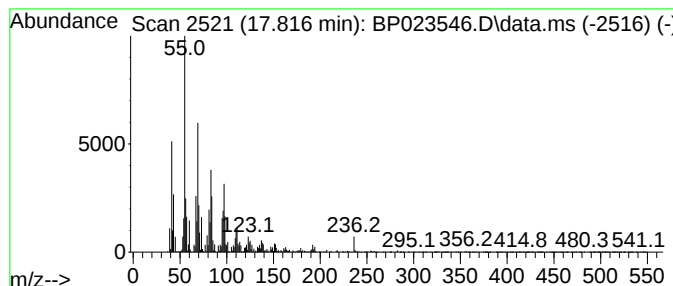
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (E)-Hexadec-9-enoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.816	2.73 ng/ul	631751	Phenanthrene-d10	16.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	96
2			cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	94
3			Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	91
4			Palmitoleic acid	254	C16H30O2	000373-49-9	91
5			10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	91



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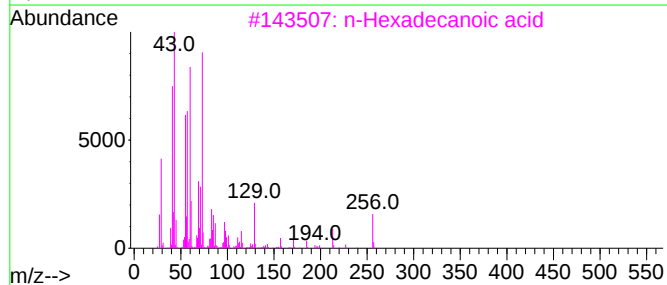
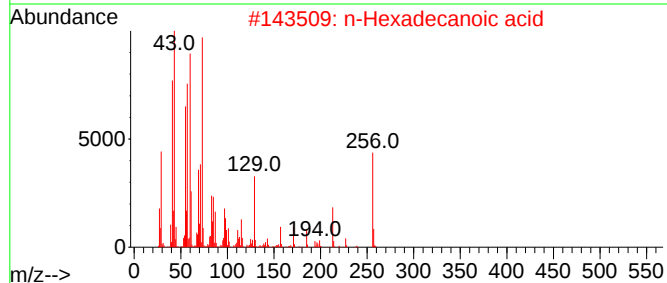
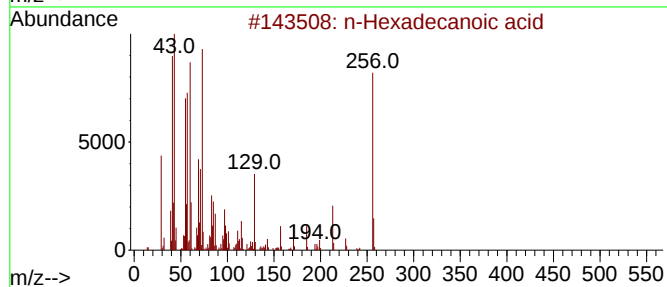
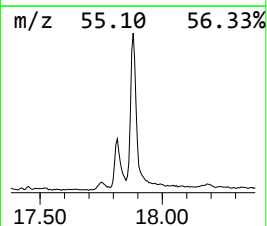
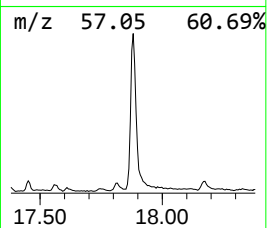
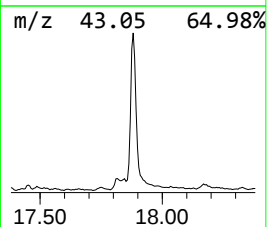
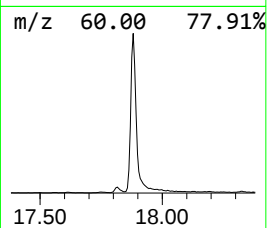
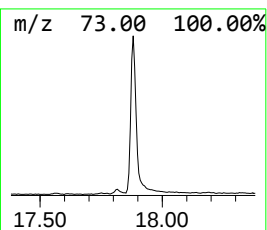
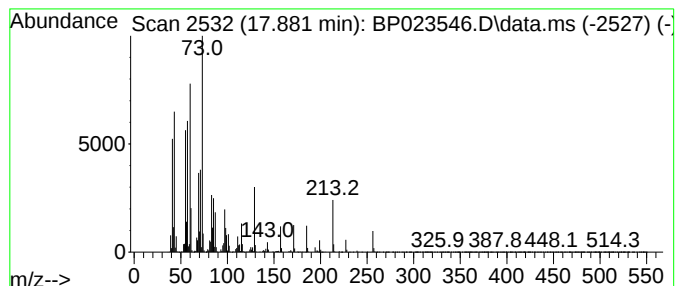
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.881	13.39 ng/ul	3095960	Phenanthrene-d10	16.939

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	96
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023546.D
Acq On : 20 Dec 2024 04:07
Operator : RC/JU
Sample : P5265-10
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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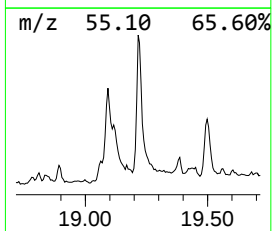
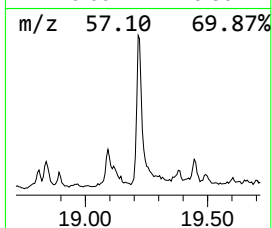
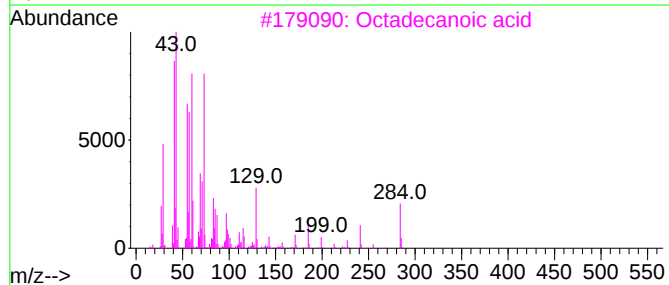
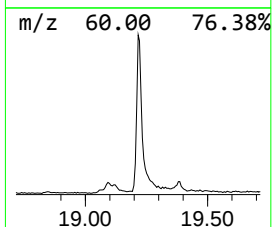
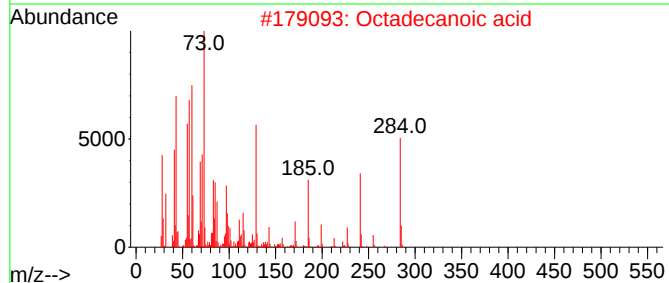
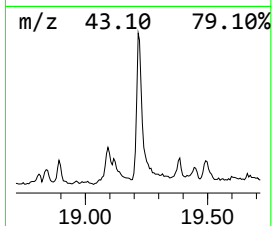
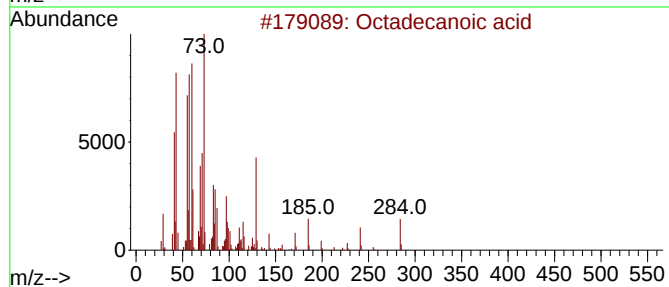
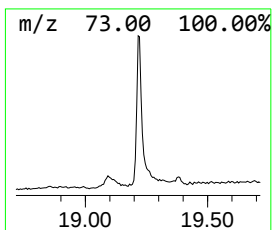
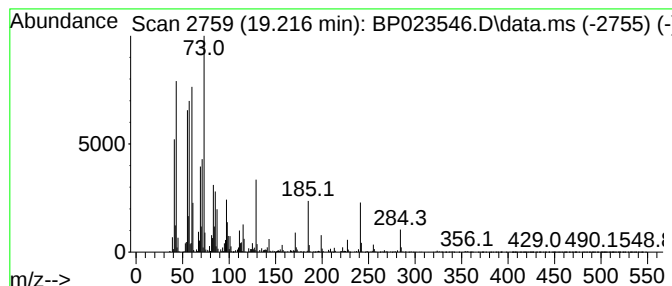
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.216	4.73 ng/ul	1272980	Chrysene-d12	21.363

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	96
4			Octadecanoic acid	284	C18H36O2	000057-11-4	96
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023546.D
Acq On : 20 Dec 2024 04:07
Operator : RC/JU
Sample : P5265-10
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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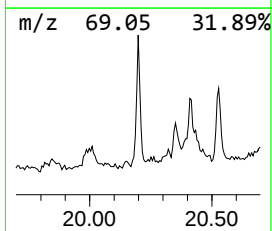
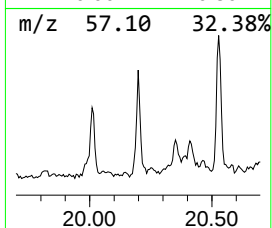
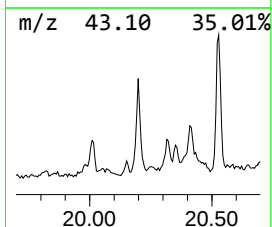
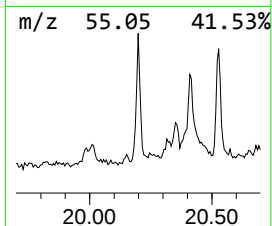
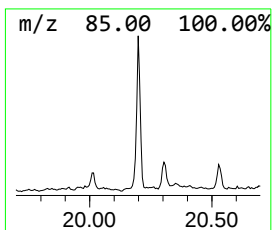
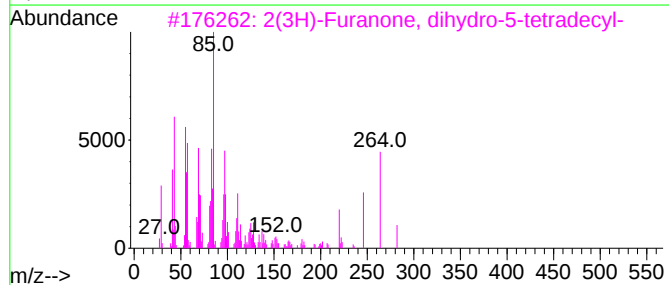
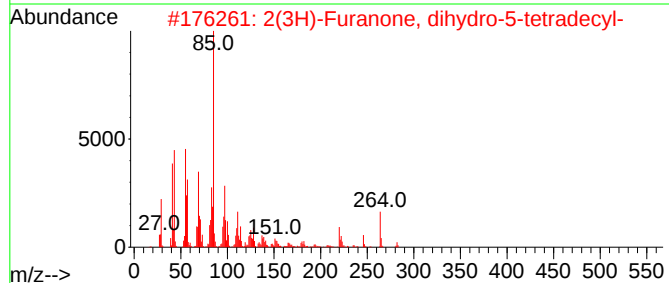
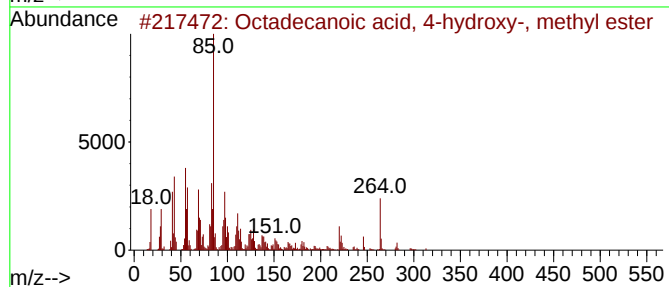
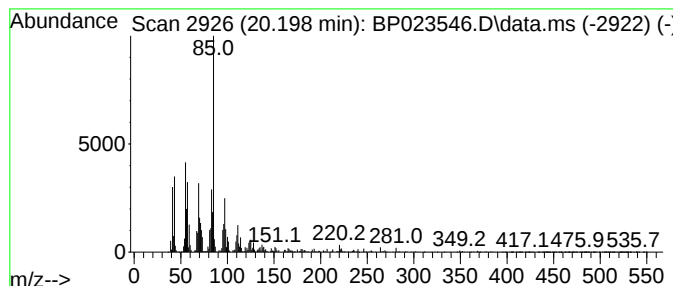
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octadecanoic acid, 4-hydrox... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.198	2.31 ng/ul	621657	Chrysene-d12	21.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid, 4-hydroxy-, m...	314	C19H38O3	002420-38-4	90
2			2(3H)-Furanone, dihydro-5-tetrad...	282	C18H34O2	000502-26-1	90
3			2(3H)-Furanone, dihydro-5-tetrad...	282	C18H34O2	000502-26-1	53
4			.gamma.-Tetradecalactone	226	C14H26O2	002721-23-5	52
5			10-Methylundecan-4-olide	198	C12H22O2	1000370-40-5	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023546.D
Acq On : 20 Dec 2024 04:07
Operator : RC/JU
Sample : P5265-10
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E28Z5

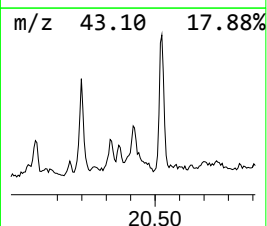
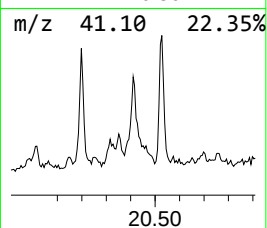
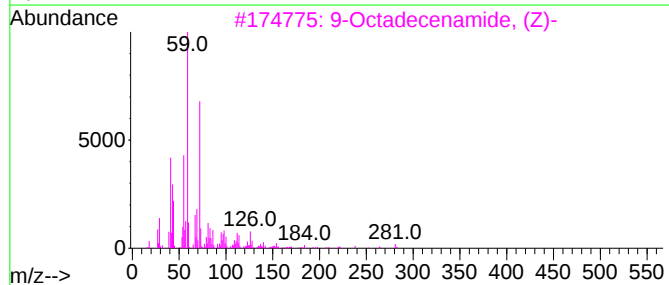
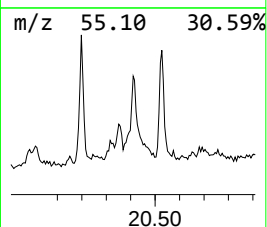
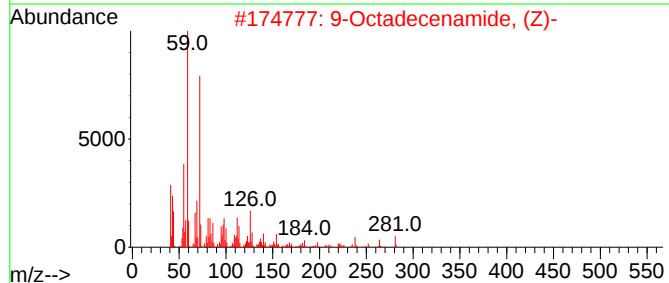
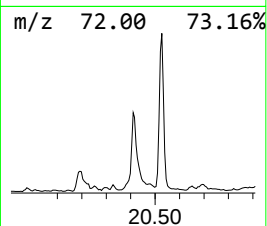
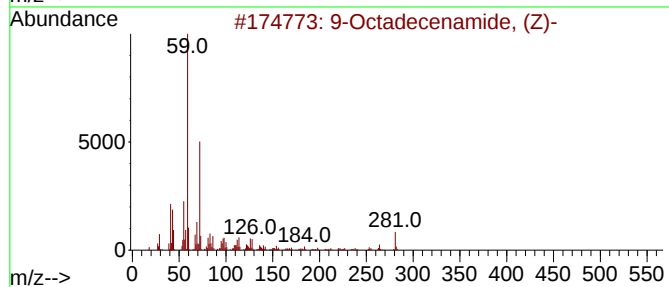
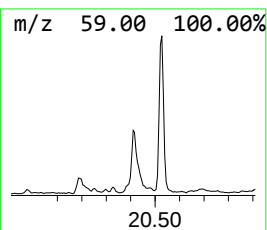
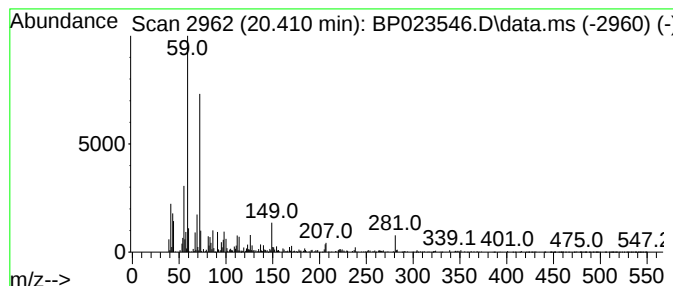
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 9-Octadecenamide, (Z)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.410	2.61 ng/ul	701306	Chrysene-d12	21.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	93
2	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	93
3	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	87
4	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	78
5	9-Octadecenamide			281	C18H35NO	003322-62-1	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023546.D
Acq On : 20 Dec 2024 04:07
Operator : RC/JU
Sample : P5265-10
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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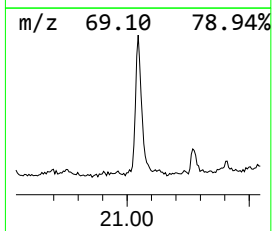
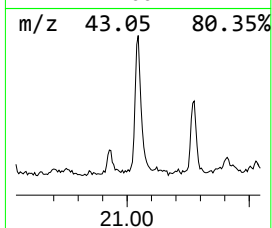
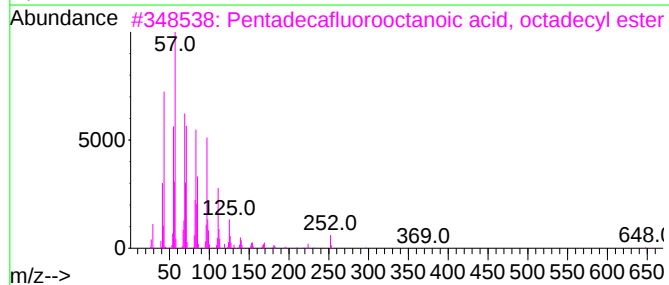
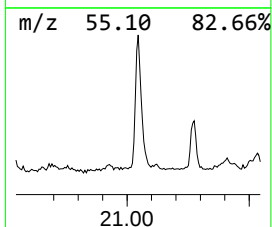
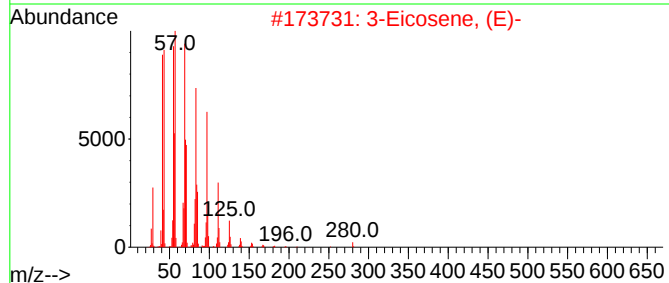
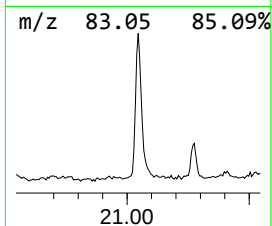
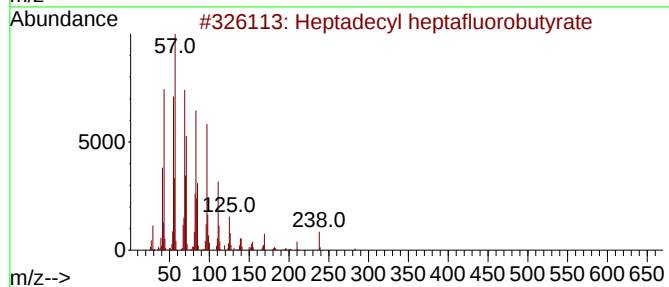
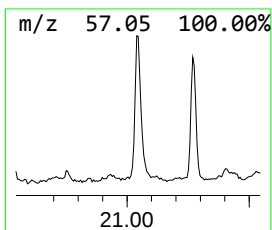
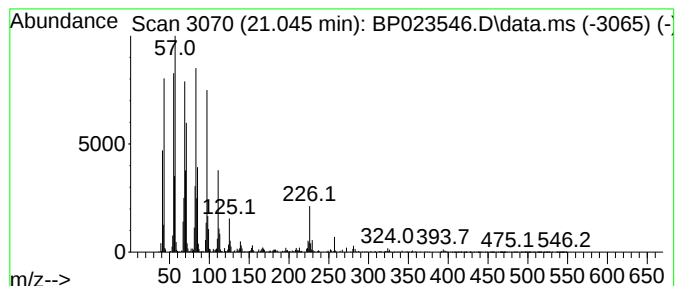
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Heptadecyl heptafluorobutyrate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.045	4.65 ng/ul	1249520	Chrysene-d12	21.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93
2			3-Eicosene, (E)-	280	C20H40	074685-33-9	92
3			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91
4			Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	90
5			9-Eicosene, (E)-	280	C20H40	074685-29-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023546.D
Acq On : 20 Dec 2024 04:07
Operator : RC/JU
Sample : P5265-10
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E28Z5

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

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	15.534	3.4	ng/ul	575516	3	14.134	3426810	20.0
(E)-Hexadec-9-e...	17.816	2.7	ng/ul	631751	4	16.939	4624410	20.0
n-Hexadecanoic ...	17.881	13.4	ng/ul	3095960	4	16.939	4624410	20.0
Octadecanoic acid	19.216	4.7	ng/ul	1272980	5	21.363	5378490	20.0
Octadecanoic ac...	20.198	2.3	ng/ul	621657	5	21.363	5378490	20.0
9-Octadecenamid...	20.410	2.6	ng/ul	701306	5	21.363	5378490	20.0
Heptadecyl hept...	21.045	4.7	ng/ul	1249520	5	21.363	5378490	20.0