

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072921\
 Data File : BP006414.D
 Acq On : 29 Jul 2021 11:21
 Operator : CG/JU
 Sample : M3123-22
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0083

Integration Parameters: LSCINT.P

Integrator: RTE

Soothing : 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-BP072421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP006414.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.710	103	106	119	rBV	102183	188504	4.06%	0.336%
2	5.816	456	464	468	rVV	50134	79983	1.72%	0.143%
3	6.775	619	627	633	rBV7	37802	78807	1.70%	0.141%
4	6.951	647	657	664	rVV	452861	858973	18.51%	1.532%
5	7.122	680	686	692	rVV	363185	581792	12.54%	1.038%
6	7.310	704	718	727	rVV	441522	752296	16.21%	1.342%
7	7.410	730	735	742	rVV3	80704	143292	3.09%	0.256%
8	7.781	788	798	805	rVV	1384605	2200660	47.43%	3.925%
9	8.487	912	918	925	rVV	476317	769866	16.59%	1.373%
10	8.604	933	938	944	rVV	56555	102250	2.20%	0.182%
11	8.934	987	994	1002	rVV	417276	687496	14.82%	1.226%
12	9.010	1002	1007	1012	rVV	92580	137880	2.97%	0.246%
13	9.116	1017	1025	1032	rBV	22667	67736	1.46%	0.121%
14	9.651	1108	1116	1121	rBV	306156	511687	11.03%	0.913%
15	10.186	1198	1207	1215	rVB	480605	824227	17.76%	1.470%
16	10.351	1226	1235	1245	rBV	514030	827612	17.84%	1.476%
17	10.563	1262	1271	1278	rBV	1962070	3335135	71.88%	5.948%
18	10.704	1288	1295	1306	rVB2	301914	633004	13.64%	1.129%
19	11.275	1385	1392	1396	rVV	38579	72607	1.56%	0.129%
20	11.604	1442	1448	1453	rBV	118516	192620	4.15%	0.344%
21	11.998	1509	1515	1525	rVV	144926	244187	5.26%	0.435%
22	12.222	1546	1553	1559	rVB2	75572	133558	2.88%	0.238%
23	12.445	1583	1591	1595	rBV	51466	74632	1.61%	0.133%
24	12.692	1627	1633	1638	rBV4	48872	86323	1.86%	0.154%
25	12.957	1674	1678	1682	rVV	80800	120173	2.59%	0.214%
26	13.245	1719	1727	1733	rBV	205449	309578	6.67%	0.552%
27	13.727	1805	1809	1814	rVV2	68615	112333	2.42%	0.200%
28	13.780	1814	1818	1821	rVV	70458	99960	2.15%	0.178%
29	13.827	1821	1826	1831	rVV	910084	1281199	27.61%	2.285%
30	13.904	1833	1839	1843	rVV2	110047	164063	3.54%	0.293%
31	13.951	1843	1847	1852	rVB4	42814	91105	1.96%	0.162%
32	14.098	1862	1872	1884	rBV	965749	1458999	31.44%	2.602%
33	14.322	1904	1910	1915	rVB	190715	259386	5.59%	0.463%
34	14.404	1915	1924	1931	rBV	2743423	4187921	90.26%	7.469%
35	14.604	1950	1958	1972	rBV2	275648	485984	10.47%	0.867%
36	14.810	1988	1993	1995	rBV	47680	73198	1.58%	0.131%

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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-BP072421.M
 Title : SVOA CALIBRATION

37	14.880	2000	2005	2010	rVB	48196	76909	1.66%	0.137%
38	14.945	2010	2016	2023	rBV3	55378	91959	1.98%	0.164%
39	15.027	2024	2030	2035	rBV2	53391	97122	2.09%	0.173%
40	15.080	2035	2039	2043	rVB	66042	92756	2.00%	0.165%
41	15.198	2053	2059	2062	rBV3	55016	82837	1.79%	0.148%
42	15.274	2068	2072	2077	rVB	202780	237143	5.11%	0.423%
43	15.398	2087	2093	2099	rVV	1221072	1812634	39.06%	3.233%
44	15.516	2106	2113	2120	rBV2	294832	462868	9.98%	0.826%
45	15.680	2136	2141	2146	rVB3	59486	99222	2.14%	0.177%
46	15.751	2146	2153	2163	rVB5	56650	161189	3.47%	0.287%
47	15.892	2174	2177	2186	rVB8	50088	117054	2.52%	0.209%
48	15.986	2186	2193	2199	rVB4	44466	101392	2.19%	0.181%
49	16.139	2214	2219	2227	rBV	258571	457058	9.85%	0.815%
50	16.468	2264	2275	2279	rBV7	47736	149771	3.23%	0.267%
51	16.692	2309	2313	2316	rBV2	49324	68128	1.47%	0.122%
52	16.739	2316	2321	2324	rBV5	57079	85480	1.84%	0.152%
53	16.815	2330	2334	2341	rVB5	79173	144869	3.12%	0.258%
54	16.892	2342	2347	2351	rVV	88479	169927	3.66%	0.303%
55	16.939	2351	2355	2359	rVV2	248214	338571	7.30%	0.604%
56	16.986	2359	2363	2374	rVB5	67046	137897	2.97%	0.246%
57	17.163	2386	2393	2397	rBV	3069265	4544185	97.93%	8.104%
58	17.198	2397	2399	2404	rVV	201795	258744	5.58%	0.461%
59	17.257	2404	2409	2417	rVB2	1267754	1783269	38.43%	3.180%
60	17.351	2421	2425	2430	rBV2	63510	117732	2.54%	0.210%
61	17.474	2443	2446	2450	rVB	102690	129397	2.79%	0.231%
62	17.680	2477	2481	2484	rVV	210032	245301	5.29%	0.437%
63	17.739	2488	2491	2495	rVB2	79994	103668	2.23%	0.185%
64	17.851	2505	2510	2513	rBV2	109384	159386	3.44%	0.284%
65	18.033	2537	2541	2543	rBV	105327	135016	2.91%	0.241%
66	18.068	2543	2547	2553	rVB2	581195	901430	19.43%	1.608%
67	18.210	2566	2571	2575	rVV	223154	314334	6.77%	0.561%
68	18.251	2575	2578	2585	rVB2	127247	172214	3.71%	0.307%
69	18.357	2590	2596	2601	rBV	237963	346028	7.46%	0.617%
70	18.521	2619	2624	2628	rBV	183852	263408	5.68%	0.470%
71	18.657	2641	2647	2653	rBV7	35728	76785	1.65%	0.137%
72	18.751	2660	2663	2668	rVB3	78133	100047	2.16%	0.178%
73	18.804	2668	2672	2675	rBV	87863	103806	2.24%	0.185%
74	18.839	2675	2678	2683	rVB	76513	107804	2.32%	0.192%
75	18.921	2687	2692	2696	rBV	256337	367832	7.93%	0.656%
76	18.968	2696	2700	2705	rVV2	347071	520051	11.21%	0.927%

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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-BP072421.M
 Title : SVOA CALIBRATION

77	19.009	2705	2707	2711	rVB	105981	110591	2.38%	0.197%
78	19.068	2711	2717	2723	rVB4	70620	178460	3.85%	0.318%
79	19.121	2723	2726	2730	rBV5	41258	79646	1.72%	0.142%
80	19.174	2731	2735	2743	rVB2	289275	450440	9.71%	0.803%
81	19.245	2743	2747	2750	rBV	108327	130340	2.81%	0.232%
82	19.304	2753	2757	2763	rVB	230781	310344	6.69%	0.553%
83	19.498	2785	2790	2794	rBV	1625708	2274645	49.02%	4.057%
84	19.527	2794	2795	2804	rVV2	494635	496040	10.69%	0.885%
85	19.604	2805	2808	2813	rVB2	98687	146674	3.16%	0.262%
86	19.845	2845	2849	2859	rVB4	87822	211126	4.55%	0.377%
87	19.962	2866	2869	2874	rBV5	42271	81647	1.76%	0.146%
88	20.009	2875	2877	2880	rVV2	82862	88473	1.91%	0.158%
89	20.051	2880	2884	2888	rVB	359490	381618	8.22%	0.681%
90	20.533	2963	2966	2971	rBV	287899	333339	7.18%	0.594%
91	20.598	2974	2977	2979	rBV4	73805	98244	2.12%	0.175%
92	20.745	2999	3002	3009	rVB	162018	274313	5.91%	0.489%
93	20.992	3039	3044	3048	rBV	442713	608844	13.12%	1.086%
94	21.256	3084	3089	3094	rBV	3557529	4640049	100.00%	8.275%
95	21.427	3114	3118	3123	rBV	263430	325771	7.02%	0.581%
96	21.880	3191	3195	3201	rBV3	307267	468960	10.11%	0.836%
97	22.915	3368	3371	3376	rVB	363358	498854	10.75%	0.890%
98	23.450	3456	3462	3467	rBV	797912	1429954	30.82%	2.550%
99	23.603	3481	3488	3500	rVB	1978990	3890006	83.84%	6.938%
100	24.239	3592	3596	3604	rVB	310351	598589	12.90%	1.068%

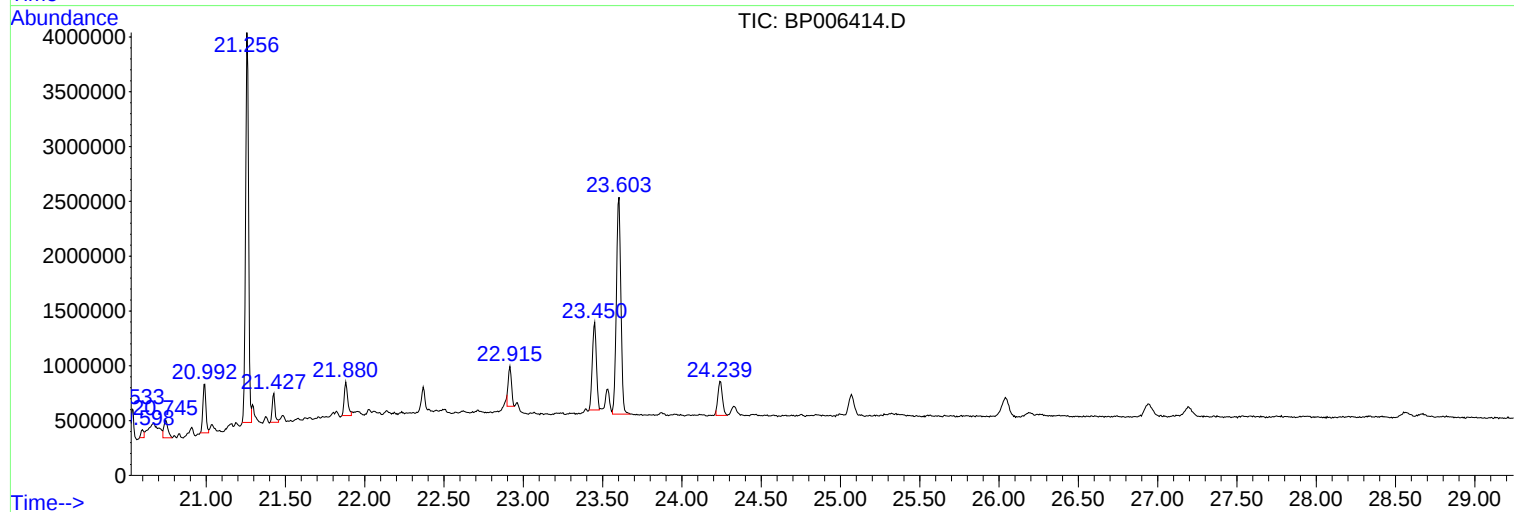
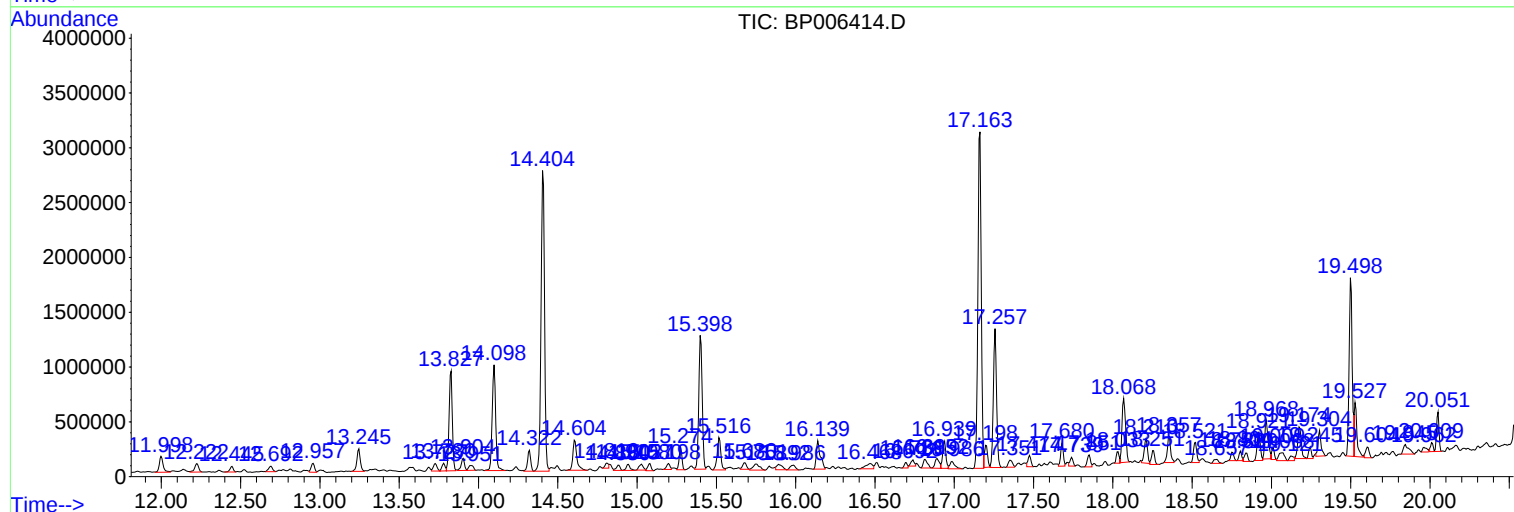
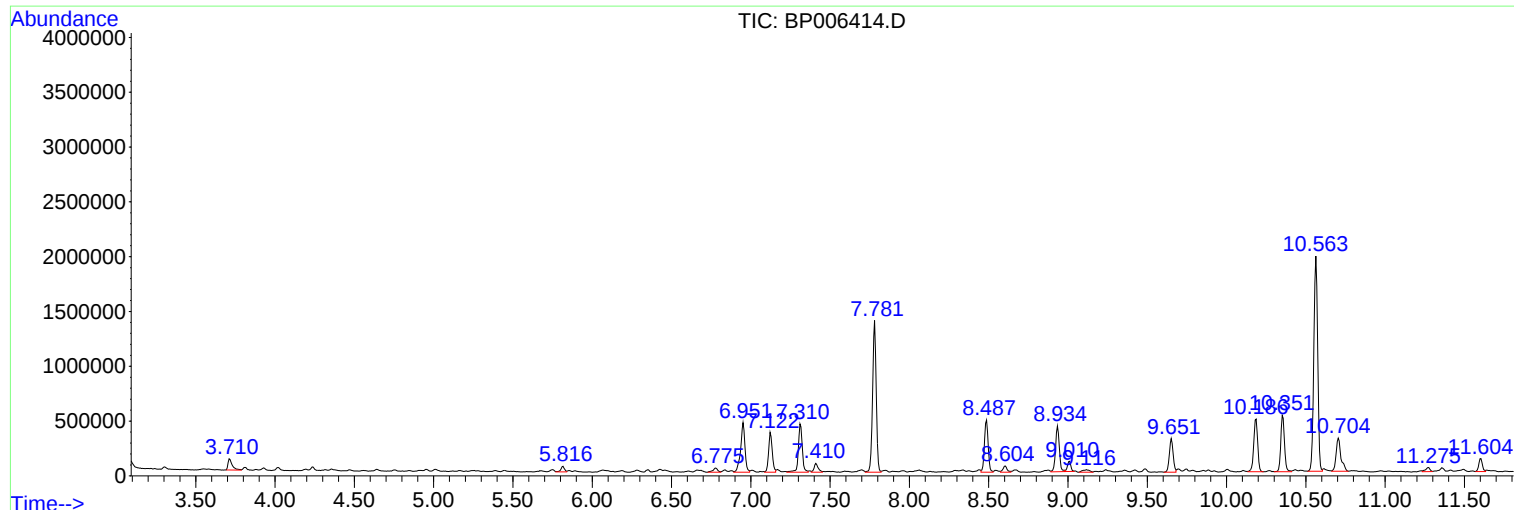
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Instrument :
BNA_P
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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



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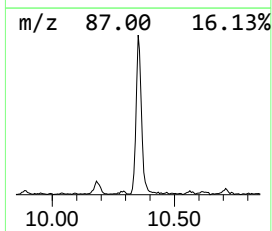
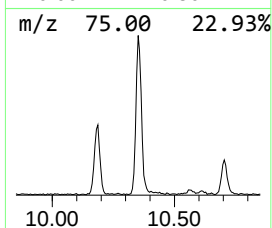
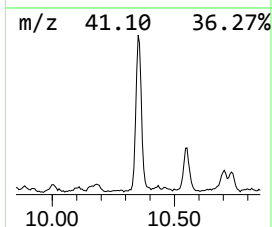
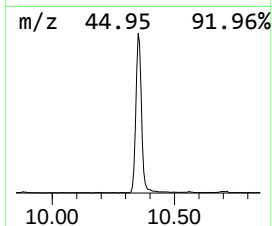
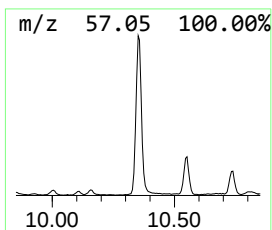
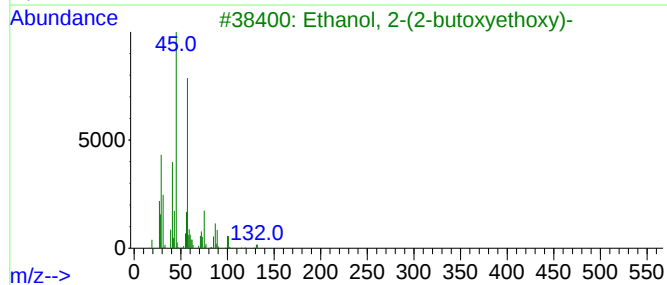
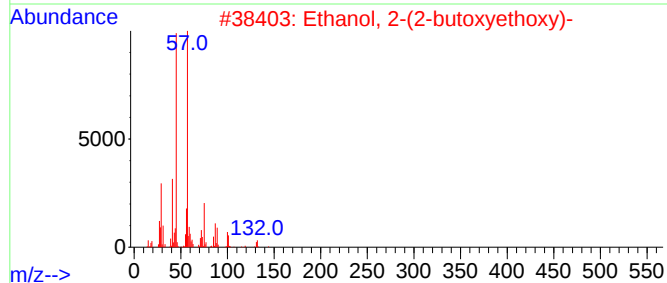
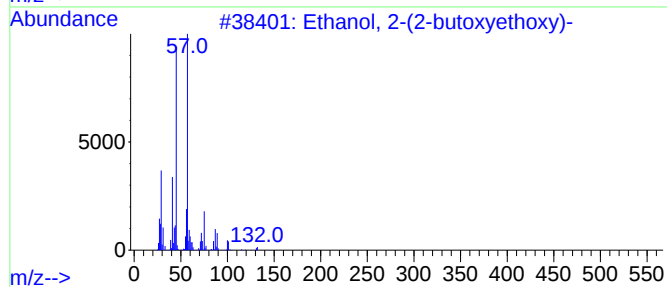
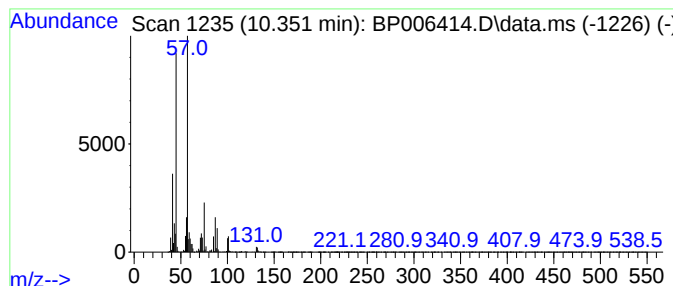
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.350	4.96 ng/ul	827612	Naphthalene-d8	10.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72



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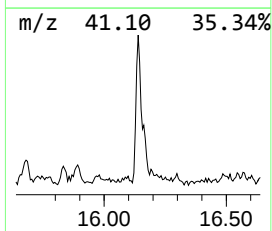
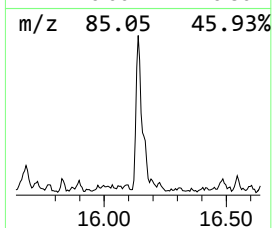
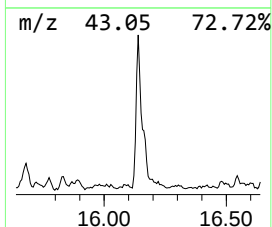
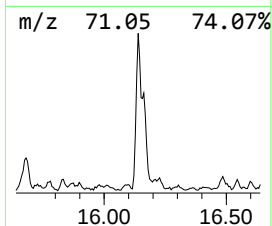
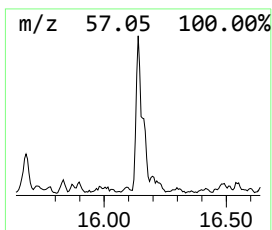
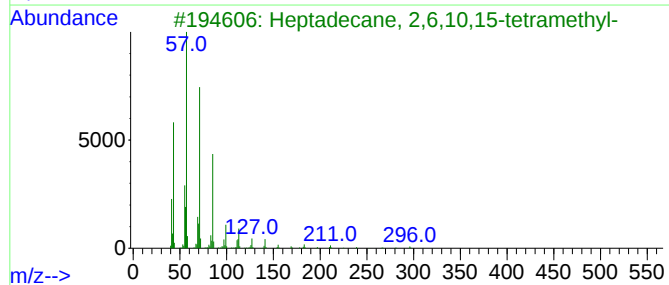
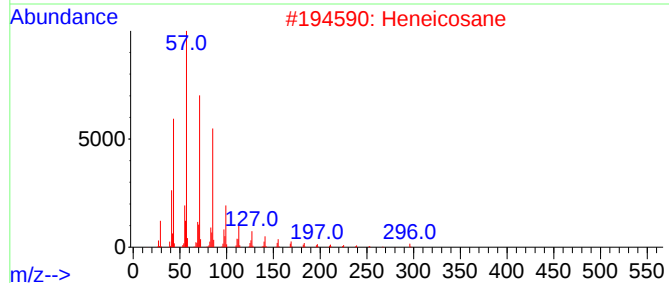
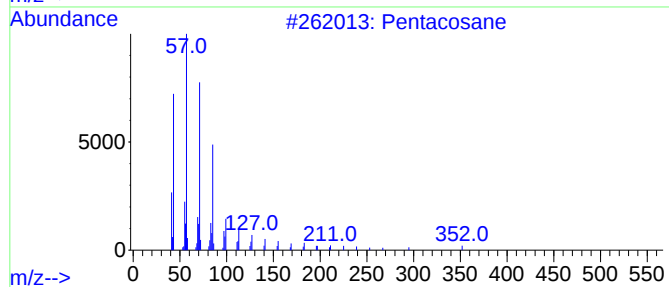
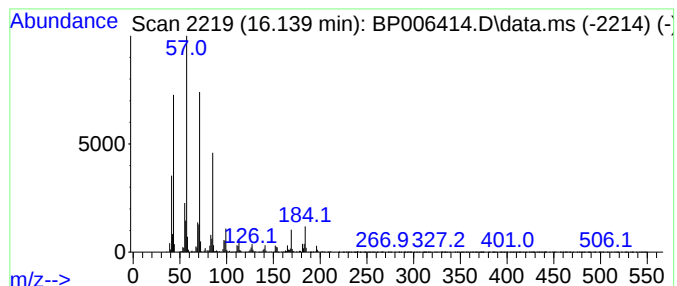
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.140	2.01 ng/ul	457058	Phenanthrene-d10	17.163	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane	352	C25H52	000629-99-2	96
2	Heneicosane	296	C21H44	000629-94-7	95
3	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
4	Tridecane	184	C13H28	000629-50-5	83
5	Tridecane	184	C13H28	000629-50-5	80



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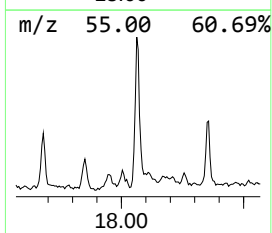
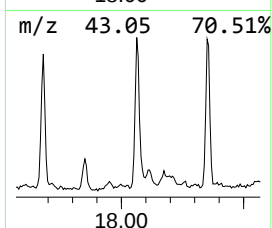
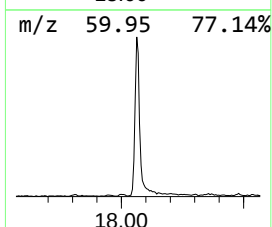
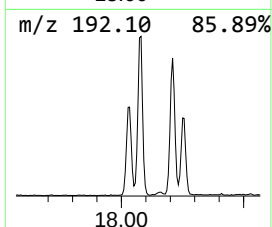
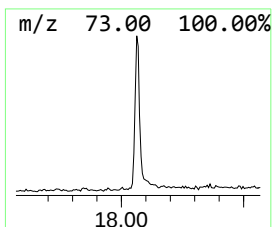
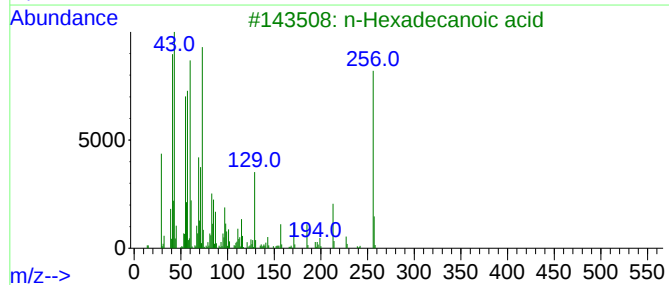
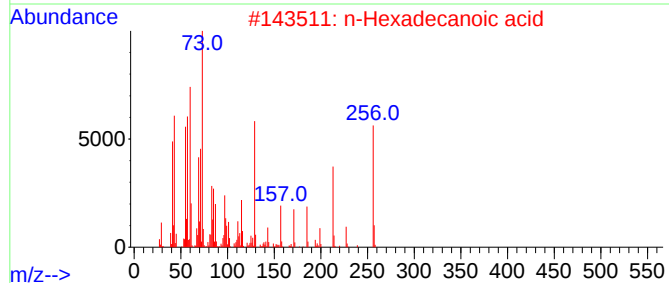
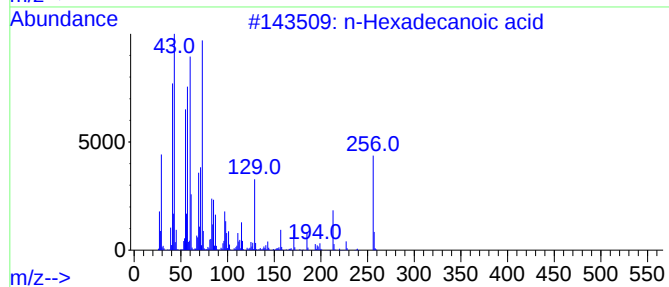
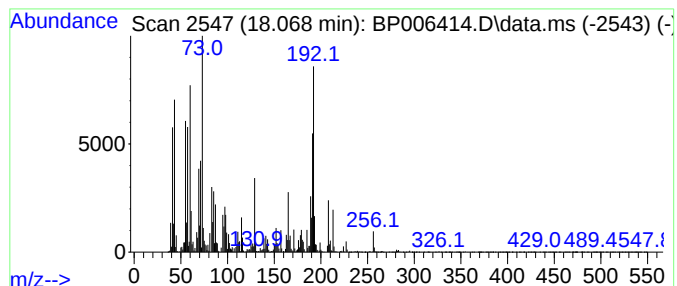
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.070	3.97 ng/ul	901430	Phenanthrene-d10	17.163

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86
5			Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072921\
Data File : BP006414.D
Acq On : 29 Jul 2021 11:21
Operator : CG/JU
Sample : M3123-22
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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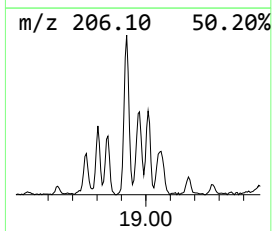
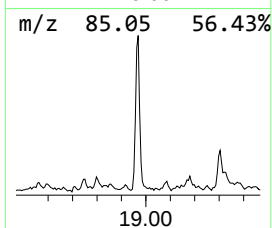
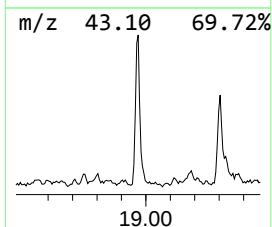
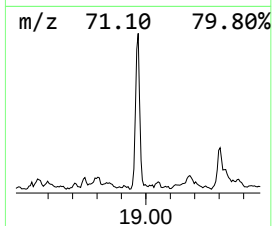
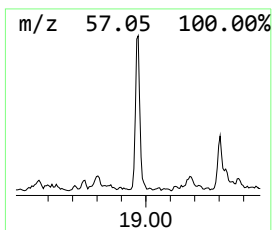
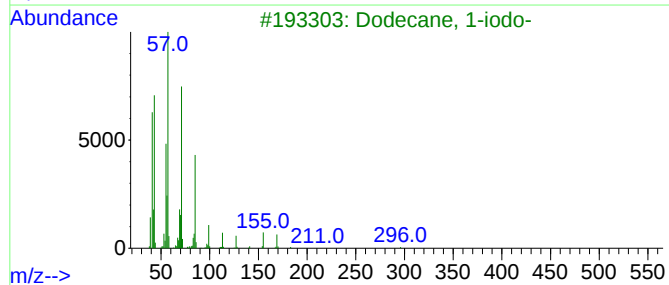
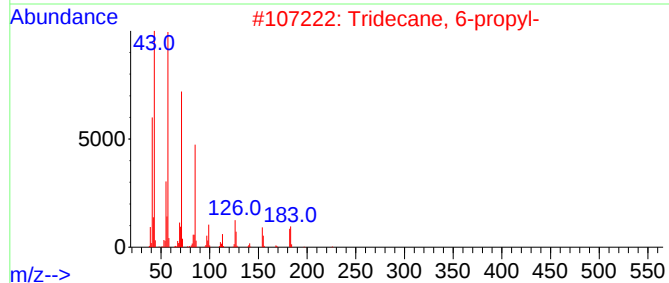
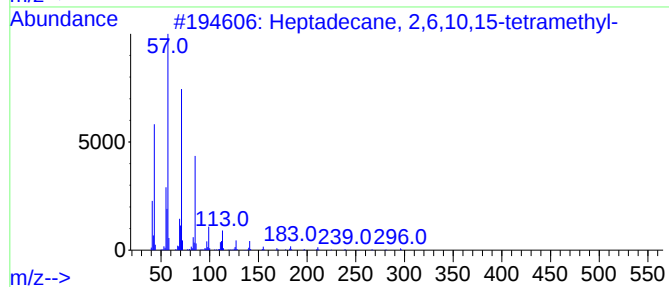
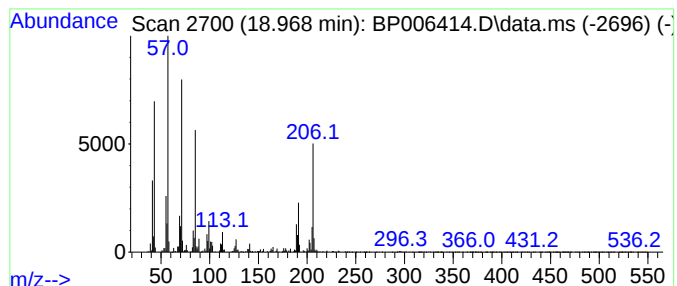
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.970	2.29 ng/ul	520051	Phenanthrene-d10	17.163

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	83
2			Tridecane, 6-propyl-	226	C16H34	055045-10-8	64
3			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	64
4			Docosane, 11-butyl-	366	C26H54	013475-76-8	59
5			Nonadecane	268	C19H40	000629-92-5	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072921\
Data File : BP006414.D
Acq On : 29 Jul 2021 11:21
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Sample : M3123-22
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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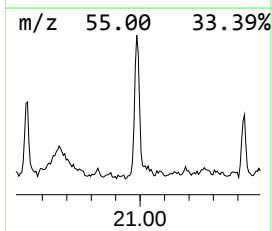
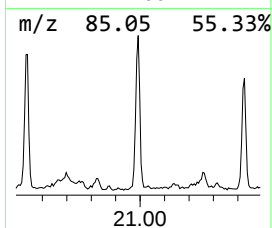
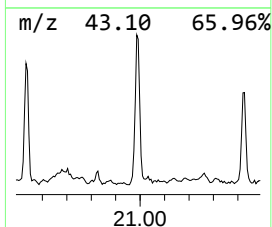
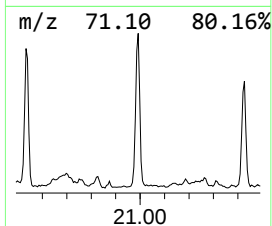
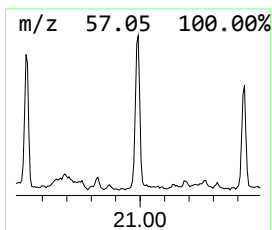
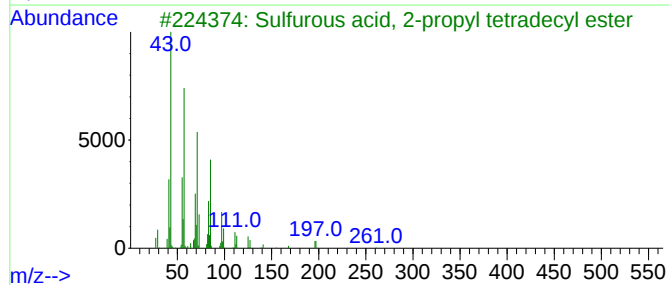
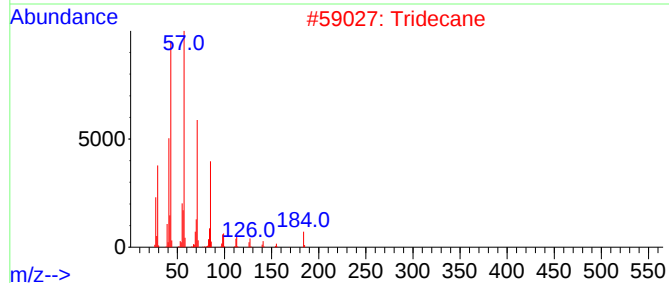
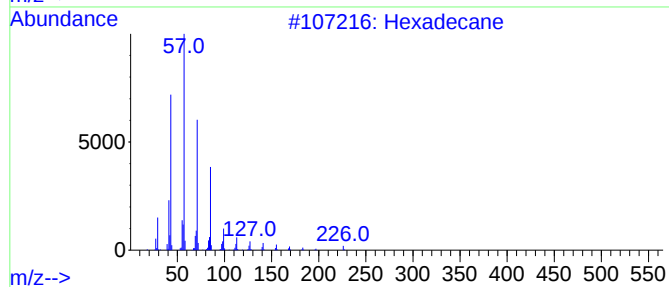
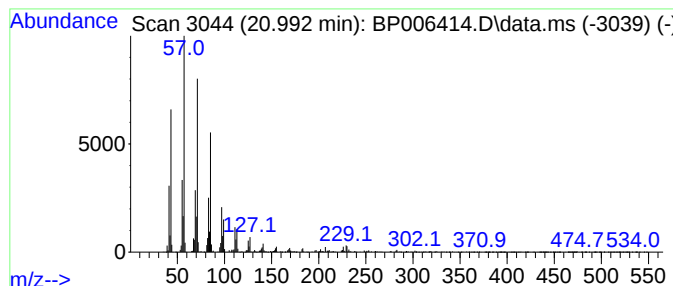
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.990	2.62 ng/ul	608844	Chrysene-d12	21.256

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	96
2		Tridecane	184	C13H28	000629-50-5	93
3		Sulfurous acid, 2-propyl tetradecyl...	320	C17H36O3S	1010309-12-5	91
4		Tritetracontane	605	C43H88	007098-21-7	91
5		Sulfurous acid, dodecyl 2-propyl...	292	C15H32O3S	1000309-12-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072921\
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Operator : CG/JU
Sample : M3123-22
Misc :
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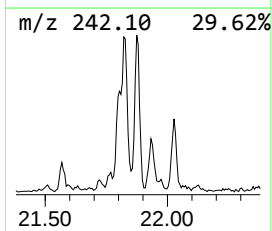
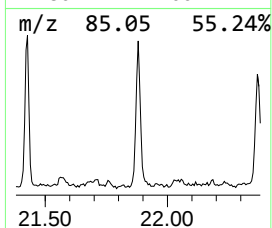
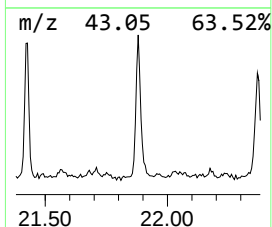
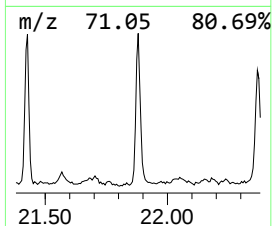
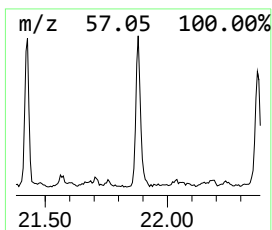
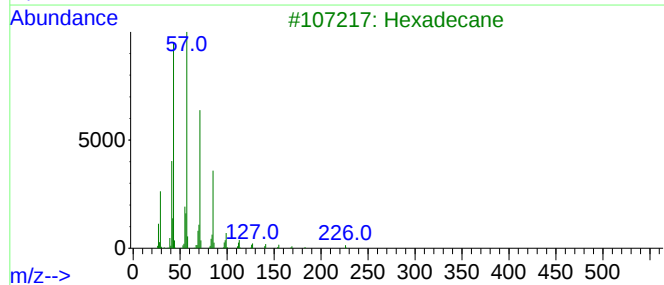
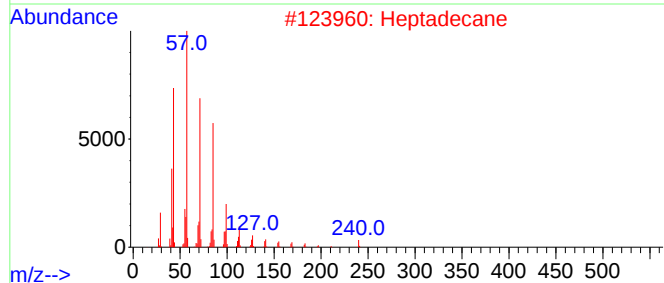
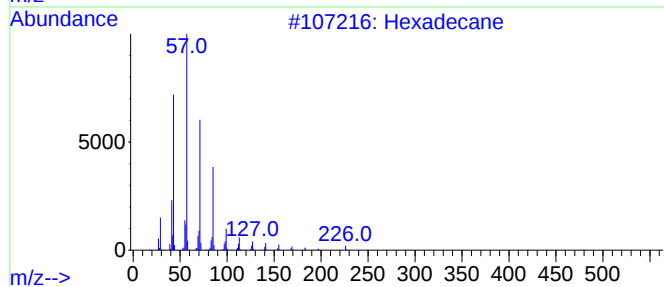
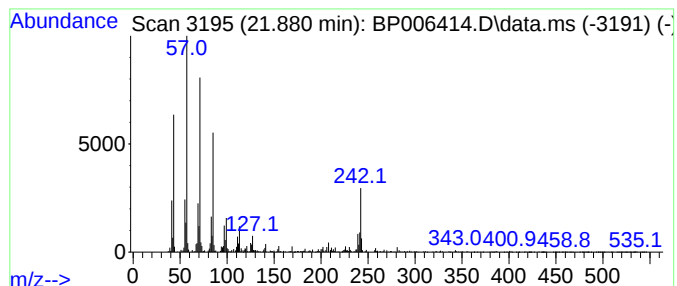
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.880	2.02 ng/ul	468960	Chrysene-d12		21.256	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	96
2	Heptadecane		240	C17H36	000629-78-7	95
3	Hexadecane		226	C16H34	000544-76-3	93
4	Heptadecane		240	C17H36	000629-78-7	89
5	Hexadecane		226	C16H34	000544-76-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072921\
Data File : BP006414.D
Acq On : 29 Jul 2021 11:21
Operator : CG/JU
Sample : M3123-22
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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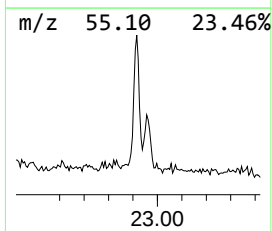
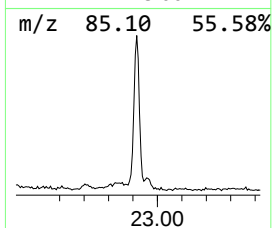
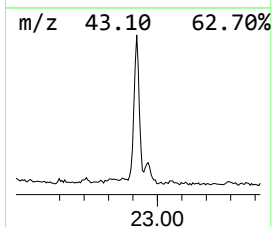
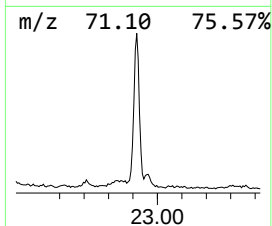
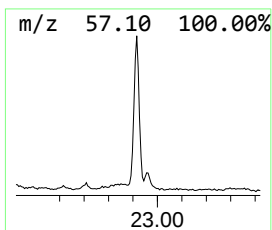
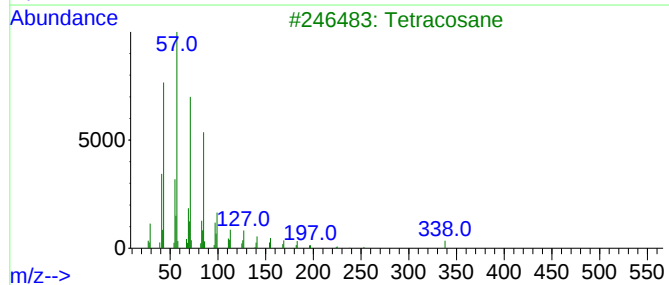
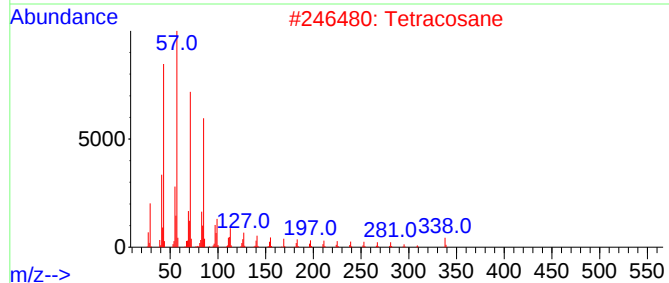
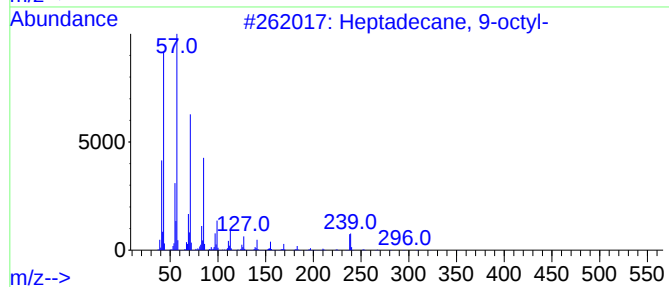
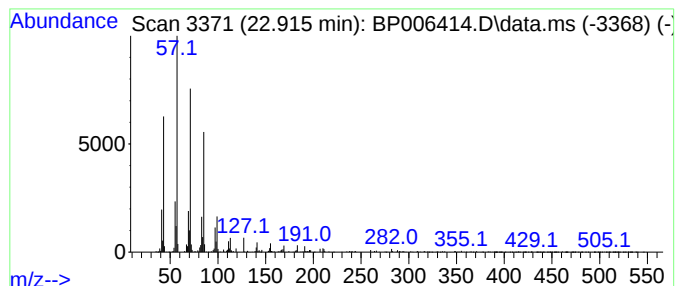
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.920	2.56 ng/ul	498854	Perylene-d12	23.603

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
2		Tetracosane	338	C24H50	000646-31-1	93
3		Tetracosane	338	C24H50	000646-31-1	90
4		Nonacosane	408	C29H60	000630-03-5	90
5		Eicosane	282	C20H42	000112-95-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072921\
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Acq On : 29 Jul 2021 11:21
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Misc :
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Instrument :
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ClientSampleId :
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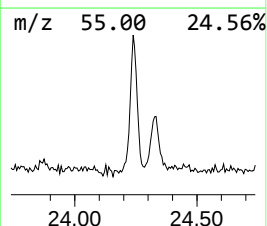
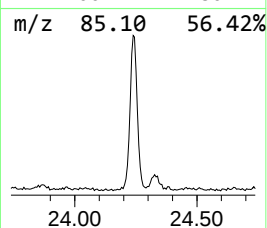
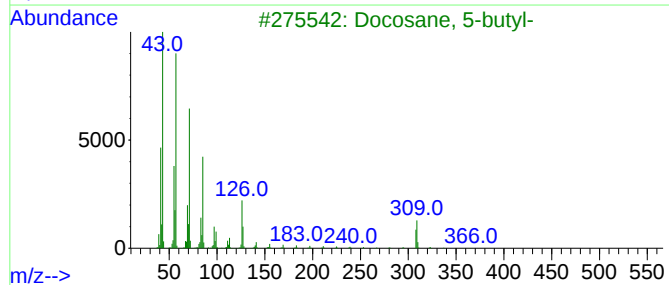
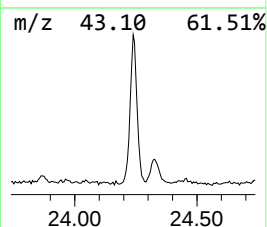
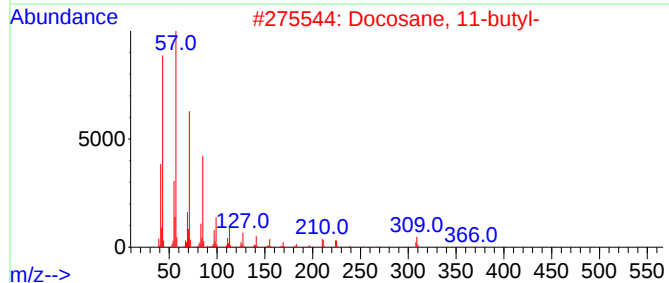
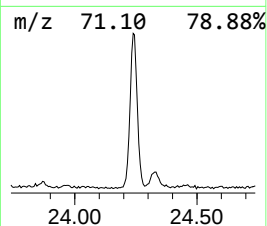
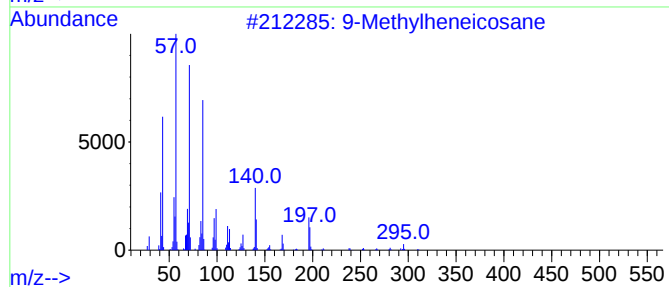
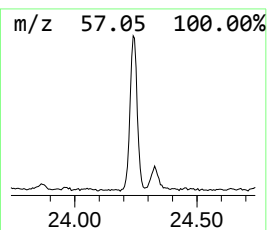
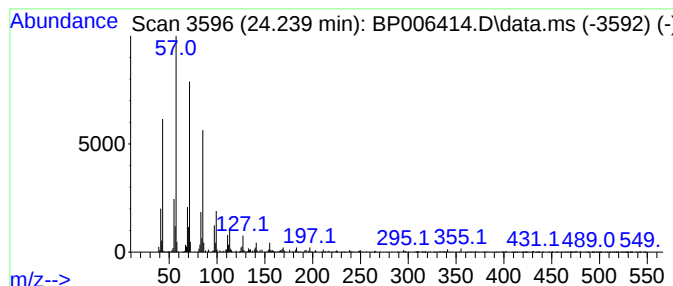
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.240	3.08 ng/ul	598589	Perylene-d12	23.603

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Methylheneicosane	310	C22H46	070475-51-3	94
2			Docosane, 11-butyl-	366	C26H54	013475-76-8	94
3			Docosane, 5-butyl-	366	C26H54	055282-16-1	93
4			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5			2-Methylheptacosane	394	C28H58	001561-00-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072921\
 Data File : BP006414.D
 Acq On : 29 Jul 2021 11:21
 Operator : CG/JU
 Sample : M3123-22
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.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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(DEL) Alkane: S...	16.140	2.0	ng/ul	457058	4	17.163	4544190	20.0
n-Hexadecanoic ...	18.070	4.0	ng/ul	901430	4	17.163	4544190	20.0
(DEL) Alkane: S...	18.970	2.3	ng/ul	520051	4	17.163	4544190	20.0
(DEL) Alkane: S...	20.990	2.6	ng/ul	608844	5	21.256	4640050	20.0
(DEL) Alkane: S...	21.880	2.0	ng/ul	468960	5	21.256	4640050	20.0
(DEL) Alkane: S...	22.920	2.6	ng/ul	498854	6	23.603	3890010	20.0
(DEL) Alkane: S...	24.240	3.1	ng/ul	598589	6	23.603	3890010	20.0