

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP009979.D
 Acq On : 20 Apr 2022 16:43
 Operator : CG/JU
 Sample : N2374-01
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFFD4

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-BP040722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP009979.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	45	49	53	rVB	11660	15730	0.51%	0.042%
2	3.799	118	121	132	rBV	42293	72471	2.35%	0.192%
3	3.893	134	137	143	rVB2	25775	37053	1.20%	0.098%
4	4.122	173	176	180	rVB3	10412	12110	0.39%	0.032%
5	4.675	266	270	275	rBV7	5796	8460	0.27%	0.022%
6	5.040	328	332	338	rVB3	8499	14462	0.47%	0.038%
7	5.128	345	347	356	rVB9	4761	9801	0.32%	0.026%
8	7.099	675	682	692	rBV	242245	418496	13.59%	1.107%
9	7.269	705	711	721	rBV	200877	329952	10.72%	0.873%
10	7.475	739	746	758	rBV	263324	426667	13.86%	1.129%
11	7.581	760	764	770	rVB7	4141	6976	0.23%	0.018%
12	7.946	816	826	835	rBV	478773	820564	26.65%	2.171%
13	8.646	936	945	956	rBV	323563	554235	18.00%	1.466%
14	8.763	960	965	973	rVB	42848	72638	2.36%	0.192%
15	9.099	1013	1022	1031	rBV	260771	449963	14.61%	1.191%
16	9.552	1090	1099	1108	rVB2	18165	35454	1.15%	0.094%
17	9.828	1138	1146	1155	rBV	206153	368178	11.96%	0.974%
18	9.981	1163	1172	1176	rBV10	5320	11385	0.37%	0.030%
19	10.181	1201	1206	1211	rBV7	5722	10279	0.33%	0.027%
20	10.363	1230	1237	1251	rBV	337068	599458	19.47%	1.586%
21	10.751	1295	1303	1310	rBV	744473	1274899	41.41%	3.373%
22	10.881	1317	1325	1343	rBV	240865	460622	14.96%	1.219%
23	11.281	1387	1393	1401	rVB10	7458	15772	0.51%	0.042%
24	11.551	1435	1439	1442	rVV6	4788	8407	0.27%	0.022%
25	12.181	1541	1546	1552	rVB5	10310	17239	0.56%	0.046%
26	12.334	1568	1572	1578	rBV5	7151	11226	0.36%	0.030%
27	12.616	1615	1620	1628	rBV8	5691	13649	0.44%	0.036%
28	13.416	1751	1756	1763	rVB6	7717	13267	0.43%	0.035%
29	13.498	1763	1770	1773	rBV8	3837	6383	0.21%	0.017%
30	13.751	1808	1813	1819	rVB10	4074	9369	0.30%	0.025%
31	13.898	1833	1838	1843	rVB4	12393	20601	0.67%	0.055%
32	13.981	1843	1852	1865	rBV	726900	1031486	33.50%	2.729%
33	14.128	1873	1877	1881	rBV7	4467	6981	0.23%	0.018%
34	14.263	1891	1900	1912	rBV	717338	1167133	37.91%	3.088%
35	14.575	1943	1953	1960	rBV	1274716	1875858	60.92%	4.963%
36	14.639	1960	1964	1972	rVB6	15883	29342	0.95%	0.078%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Title : SVOA CALIBRATION

37	14.757	1978	1984	1995	rBV	206696	374125	12.15%	0.990%
38	14.910	2004	2010	2014	rBV9	5853	13398	0.44%	0.035%
39	14.975	2017	2021	2027	rVV4	11813	20424	0.66%	0.054%
40	15.051	2030	2034	2039	rVB6	7527	11115	0.36%	0.029%
41	15.239	2062	2066	2072	rVB9	6283	13838	0.45%	0.037%
42	15.381	2085	2090	2096	rVV4	11256	23452	0.76%	0.062%
43	15.569	2115	2122	2127	rBV	1058868	1568574	50.94%	4.150%
44	15.622	2127	2131	2136	rVV3	27960	43039	1.40%	0.114%
45	15.681	2136	2141	2148	rVV	173862	256089	8.32%	0.678%
46	15.804	2155	2162	2171	rBV8	12905	30742	1.00%	0.081%
47	15.928	2177	2183	2189	rVB10	11051	25866	0.84%	0.068%
48	16.145	2217	2220	2227	rVB8	7720	13510	0.44%	0.036%
49	16.298	2241	2246	2247	rBV5	14437	20307	0.66%	0.054%
50	16.316	2247	2249	2253	rVB	19639	21424	0.70%	0.057%
51	16.357	2253	2256	2260	rVB6	6545	8634	0.28%	0.023%
52	16.404	2260	2264	2268	rBV	31544	47137	1.53%	0.125%
53	16.463	2270	2274	2281	rVB3	33076	58579	1.90%	0.155%
54	16.528	2281	2285	2289	rBV2	34774	46800	1.52%	0.124%
55	16.575	2289	2293	2296	rBV2	21194	25416	0.83%	0.067%
56	16.628	2298	2302	2303	rBV4	12004	13966	0.45%	0.037%
57	16.675	2308	2310	2314	rVB5	18147	20895	0.68%	0.055%
58	16.728	2314	2319	2325	rBV4	34756	59648	1.94%	0.158%
59	16.792	2325	2330	2334	rBV	33720	50648	1.64%	0.134%
60	16.828	2334	2336	2339	rVV3	10377	13155	0.43%	0.035%
61	16.869	2339	2343	2347	rVB3	21912	27692	0.90%	0.073%
62	16.981	2357	2362	2364	rBV2	13105	19036	0.62%	0.050%
63	17.086	2370	2380	2382	rBV10	13296	34421	1.12%	0.091%
64	17.139	2386	2389	2399	rVB	21580	36432	1.18%	0.096%
65	17.328	2414	2421	2425	rBV	1561542	2330387	75.69%	6.166%
66	17.369	2425	2428	2432	rVV	363648	509739	16.56%	1.349%
67	17.422	2432	2437	2449	rVV2	1184775	1852619	60.17%	4.902%
68	17.516	2449	2453	2461	rVB3	22944	43688	1.42%	0.116%
69	17.728	2485	2489	2497	rVB	24837	38730	1.26%	0.102%
70	17.833	2503	2507	2514	rBV7	15080	27759	0.90%	0.073%
71	17.904	2516	2519	2523	rVB2	19329	23502	0.76%	0.062%
72	18.204	2564	2570	2573	rBV2	123076	223724	7.27%	0.592%
73	18.239	2573	2576	2581	rVB	90813	116253	3.78%	0.308%
74	18.375	2594	2599	2603	rBV2	105802	183679	5.97%	0.486%
75	18.416	2603	2606	2611	rVB	56095	74663	2.42%	0.198%
76	18.669	2646	2649	2652	rBV	49865	64235	2.09%	0.170%

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77	18.710	2652	2656	2661	rVB3	87584	122433	3.98%	0.324%
78	18.910	2687	2690	2694	rBV5	27974	33252	1.08%	0.088%
79	18.957	2695	2698	2702	rBV4	20417	33124	1.08%	0.088%
80	19.075	2714	2718	2721	rBV	65438	94204	3.06%	0.249%
81	19.210	2737	2741	2743	rBV3	41843	60560	1.97%	0.160%
82	19.327	2756	2761	2767	rVB	1001023	1323838	43.00%	3.503%
83	19.569	2799	2802	2808	rVB2	74201	103417	3.36%	0.274%
84	19.657	2811	2817	2819	rVV	1620137	2413344	78.38%	6.385%
85	19.680	2819	2821	2829	rVB	938850	1189119	38.62%	3.146%
86	19.992	2870	2874	2881	rBV5	61420	161970	5.26%	0.429%
87	20.163	2900	2903	2910	rVB2	155567	228826	7.43%	0.605%
88	20.263	2916	2920	2923	rBV	90817	114981	3.73%	0.304%
89	21.063	3053	3056	3059	rBV2	142206	166698	5.41%	0.441%
90	21.110	3061	3064	3072	rVB3	598497	863378	28.04%	2.284%
91	21.316	3096	3099	3103	rVB	173332	174380	5.66%	0.461%
92	21.404	3108	3114	3118	rVV2	1794061	3078986	100.00%	8.146%
93	21.445	3118	3121	3132	rVB	480118	683245	22.19%	1.808%
94	22.045	3217	3223	3228	rBV3	524574	1047784	34.03%	2.772%
95	23.110	3399	3404	3409	rBV	963026	1815811	58.97%	4.804%
96	23.157	3409	3412	3421	rVB	473616	794108	25.79%	2.101%
97	23.698	3499	3504	3509	rBV2	665371	1220941	39.65%	3.230%
98	23.857	3525	3531	3546	rVB2	983569	2145164	69.67%	5.676%
99	24.510	3637	3642	3650	rVB	390221	803062	26.08%	2.125%
100	24.604	3654	3658	3668	rVB3	224161	499277	16.22%	1.321%

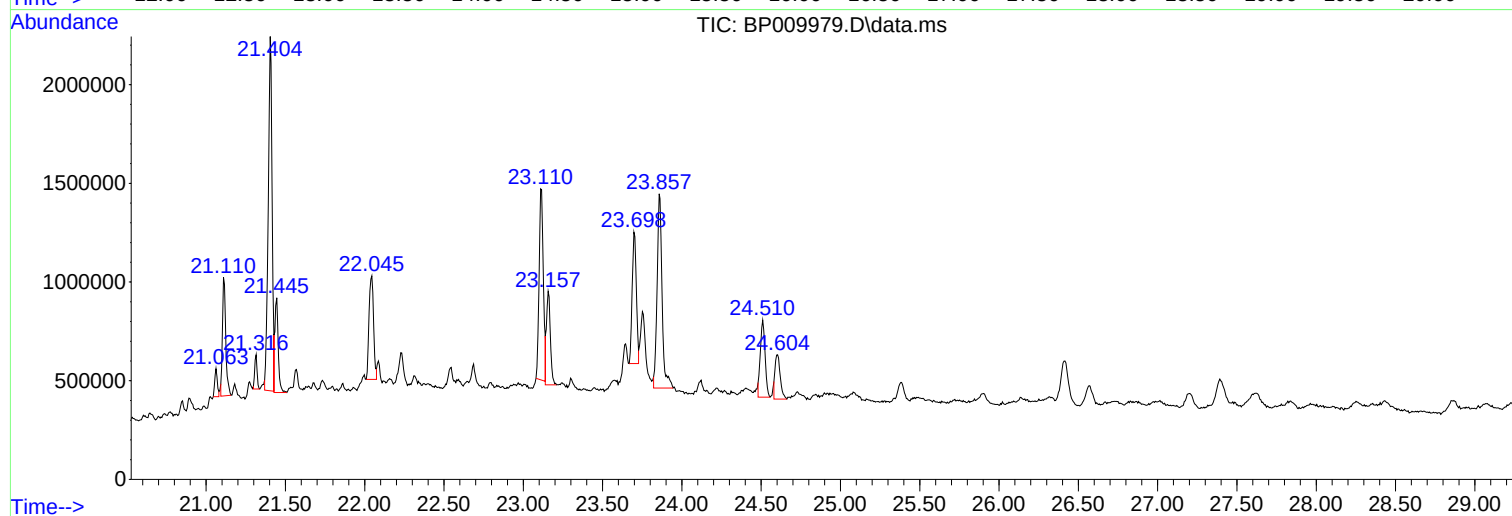
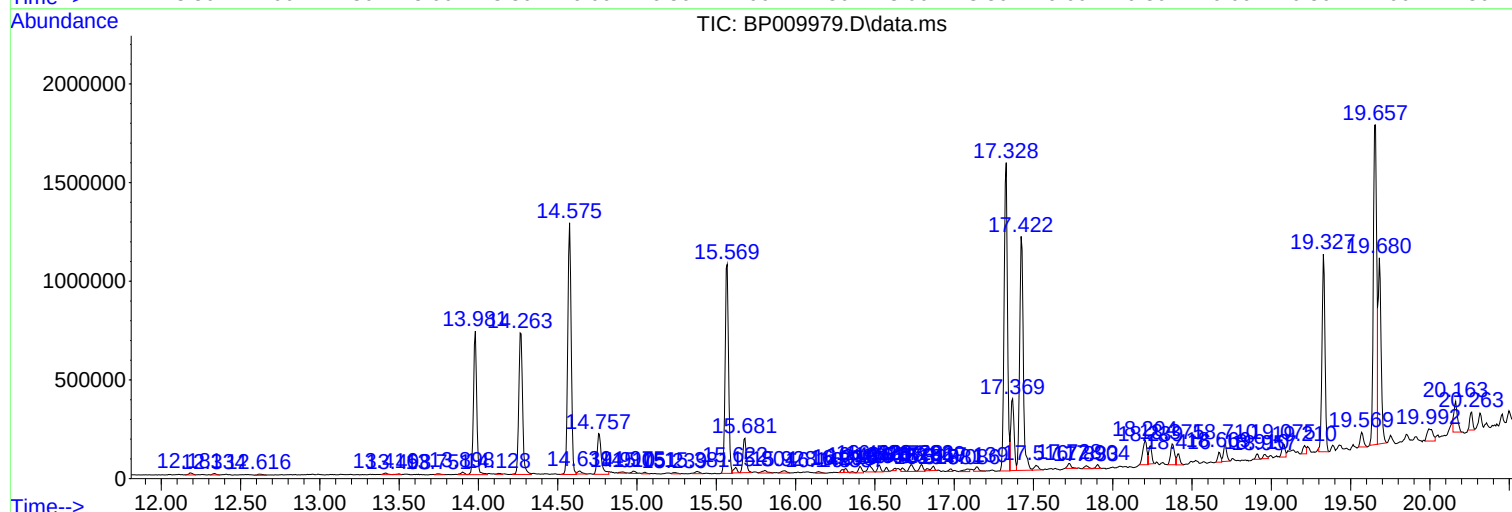
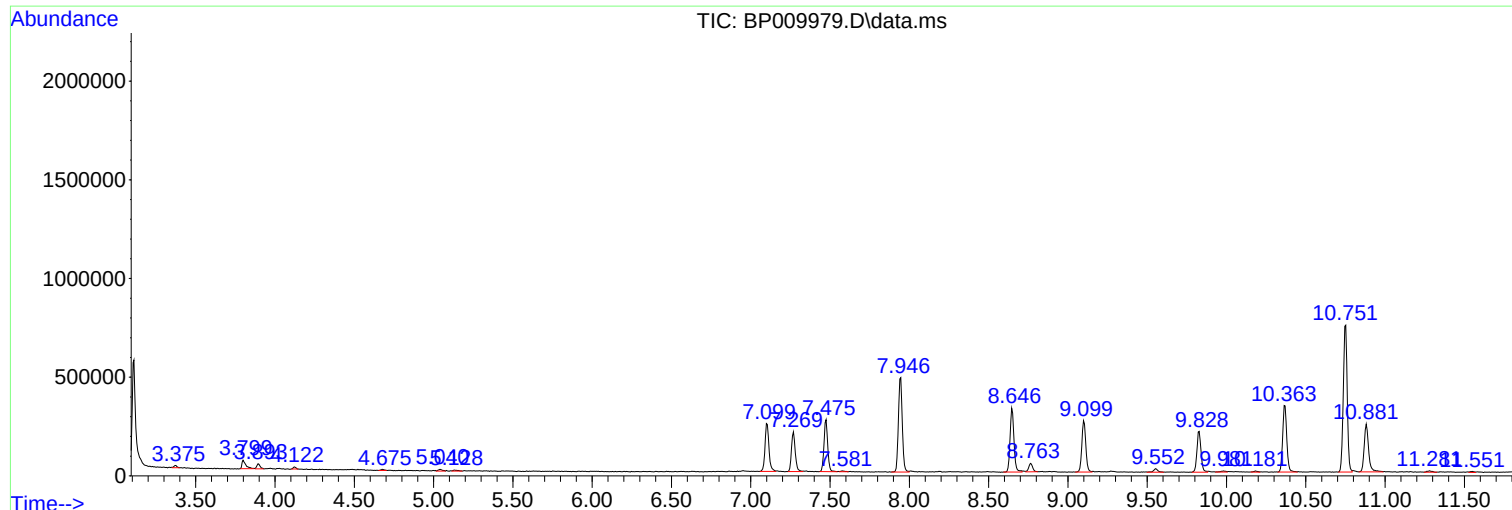
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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



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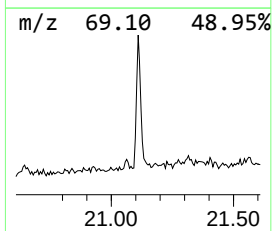
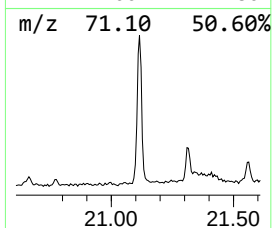
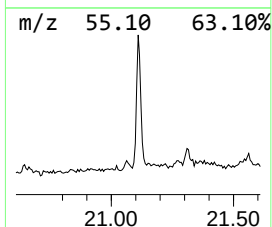
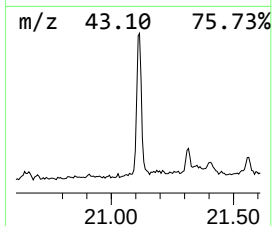
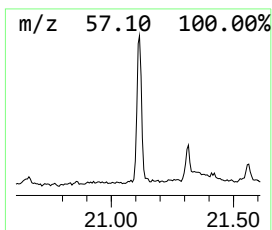
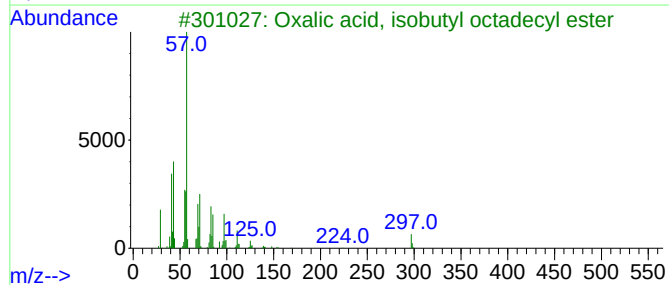
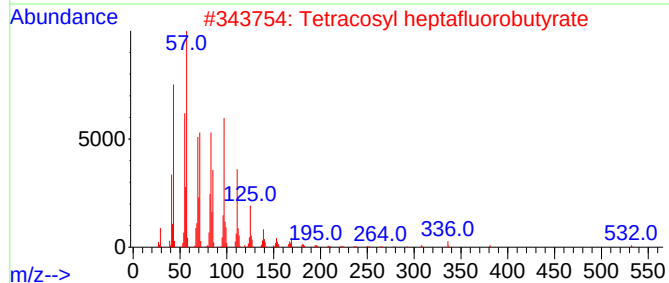
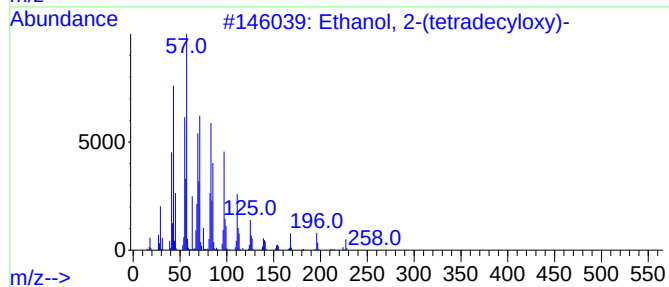
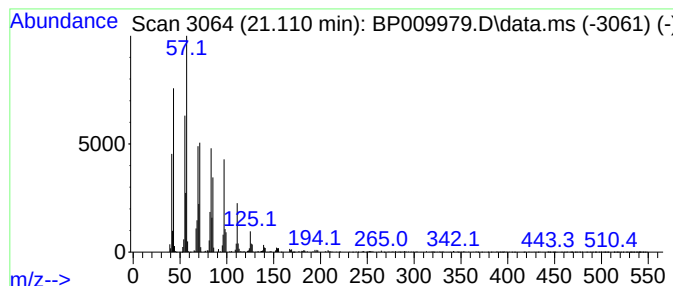
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 2-(tetradecyloxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.110	5.61 ng/ul	863378	Chrysene-d12	21.404

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
2			Tetracosyl heptafluorobutyrate	550	C28H49F7O2	1000351-83-7	91
3			Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	91
4			1-Dodecanol, 2-octyl-	298	C20H42O	005333-42-6	91
5			17-Pentatriacontene	491	C35H70	006971-40-0	91



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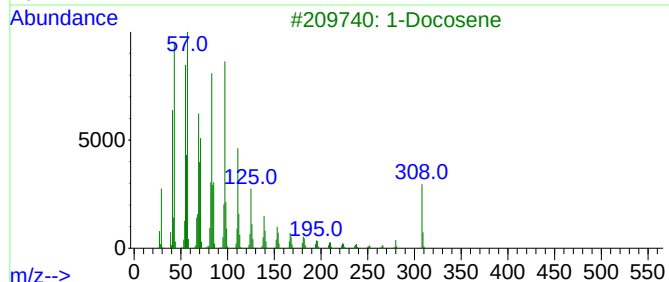
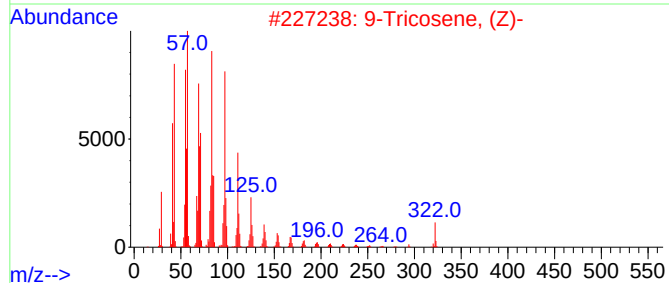
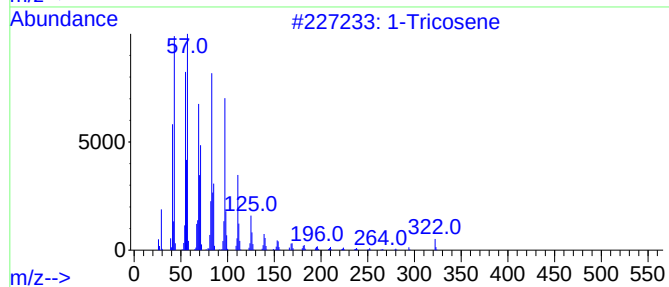
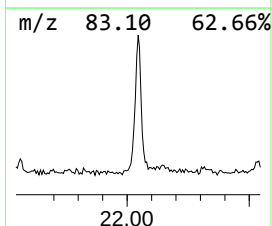
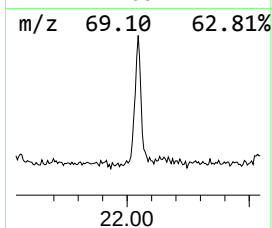
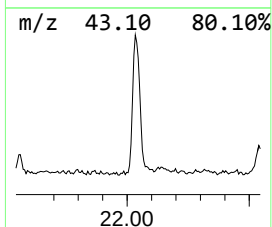
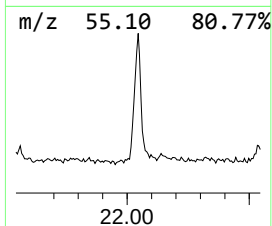
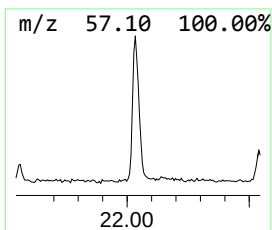
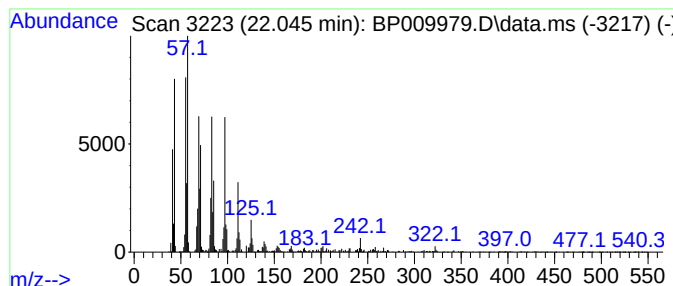
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.045	6.81 ng/ul	1047780	Chrysene-d12	21.404

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tricosene			322	C23H46	018835-32-0	98
2	9-Tricosene, (Z)-			322	C23H46	027519-02-4	96
3	1-Docosene			308	C22H44	001599-67-3	94
4	Heneicosyl heptafluorobutyrate			508	C25H43F7O2	1000351-83-8	94
5	Pentafluoropropionic acid, hepta...			402	C20H35F5O2	959218-78-1	94



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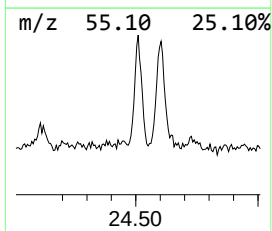
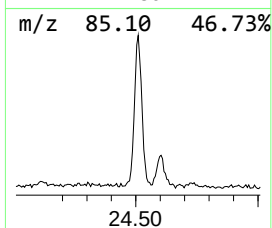
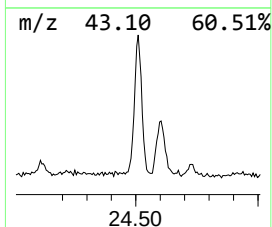
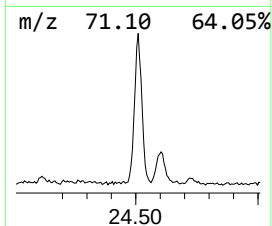
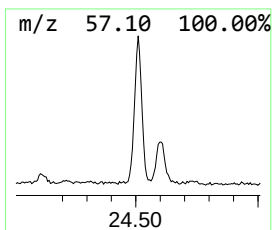
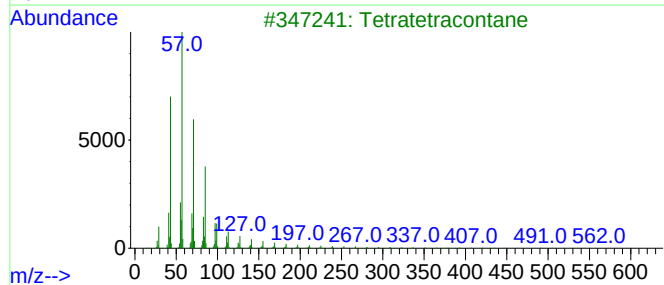
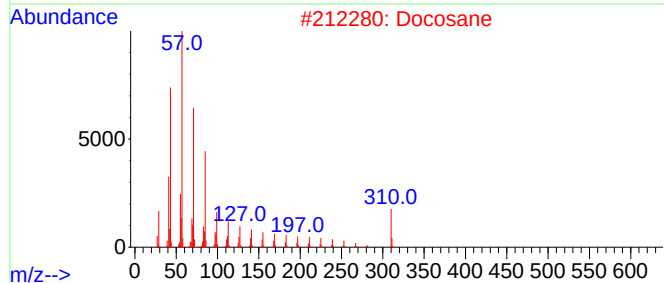
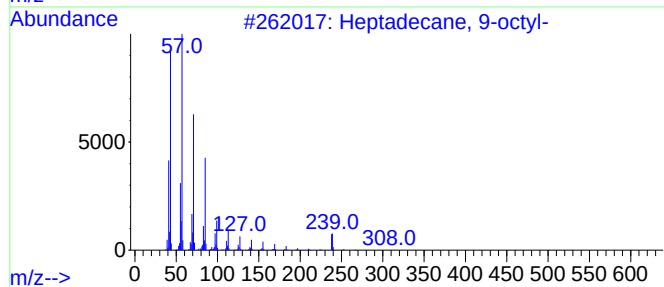
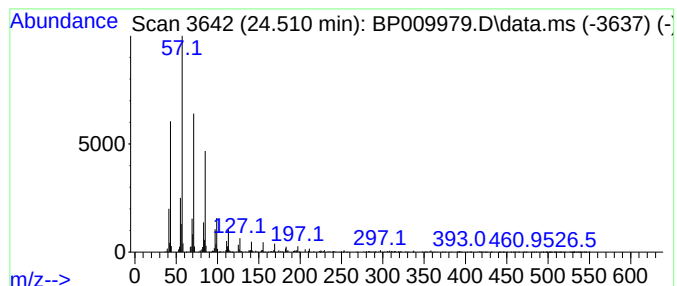
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.509	7.49 ng/ul	803062	Perylene-d12	23.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2			Docosane	310	C22H46	000629-97-0	93
3			Tetratetracontane	619	C44H90	007098-22-8	91
4			Hentriacontane	437	C31H64	000630-04-6	91
5			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009979.D
Acq On : 20 Apr 2022 16:43
Operator : CG/JU
Sample : N2374-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFD4

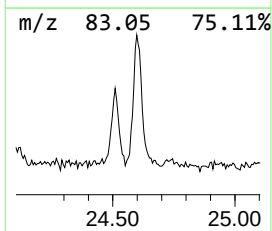
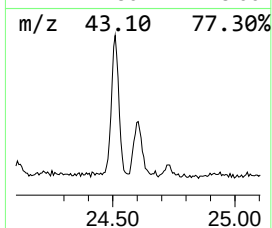
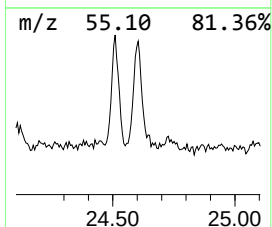
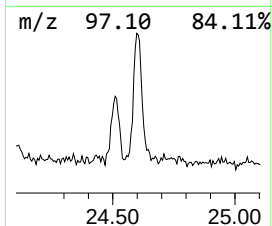
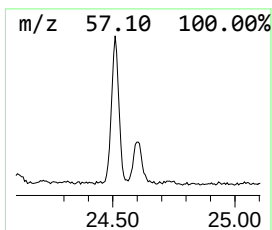
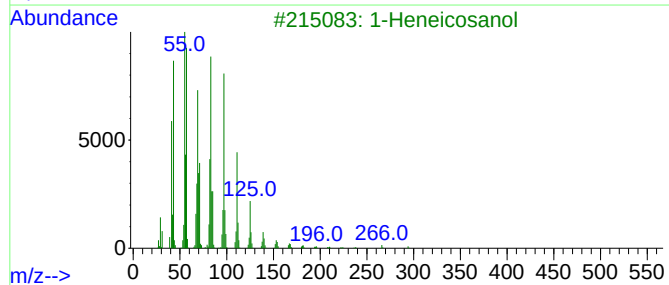
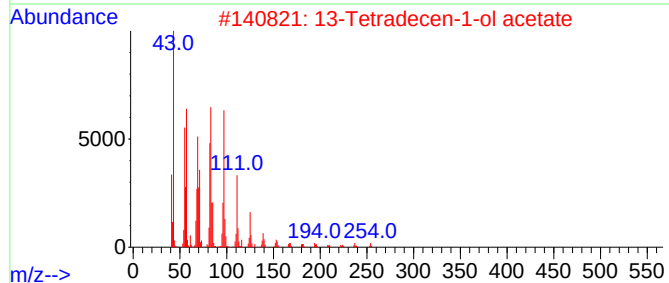
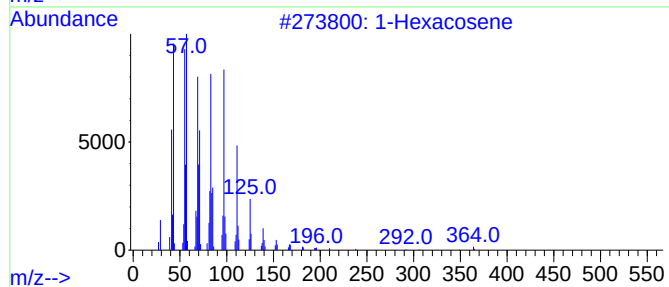
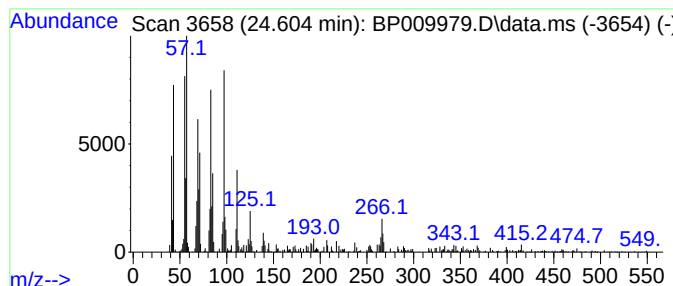
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.604	4.65 ng/ul	499277	Perylene-d12	23.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene			364	C26H52	018835-33-1	95
2	13-Tetradecen-1-ol acetate			254	C16H30O2	056221-91-1	93
3	1-Heneicosanol			312	C21H44O	015594-90-8	93
4	10-Heneicosene (c,t)			294	C21H42	095008-11-0	91
5	Heptafluorobutyric acid, n-octad...			466	C22H37F7O2	000400-57-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009979.D
Acq On : 20 Apr 2022 16:43
Operator : CG/JU
Sample : N2374-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(tet...	21.110	5.6	ng/ul	863378	5	21.404	3078990	20.0
(DEL) Alkane: S...	22.045	6.8	ng/ul	1047780	5	21.404	3078990	20.0
(DEL) Alkane: S...	24.509	7.5	ng/ul	803062	6	23.857	2145160	20.0
(DEL) Alkane: S...	24.604	4.7	ng/ul	499277	6	23.857	2145160	20.0