

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
 Data File : BP023434.D
 Acq On : 16 Dec 2024 14:27
 Operator : RC/JU
 Sample : P5153-21
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E28R4

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: BP023434.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.105	17	20	27	rVB	18183	25714	0.65%	0.063%
2	3.611	102	106	115	rVV	24745	39978	1.01%	0.097%
3	3.823	138	142	146	rVB	9122	13036	0.33%	0.032%
4	4.370	231	235	240	rVB7	7377	11525	0.29%	0.028%
5	4.717	288	294	295	rBV	12041	16353	0.41%	0.040%
6	5.593	435	443	448	rBV9	8060	15430	0.39%	0.038%
7	6.758	632	641	658	rVV	262104	592451	14.97%	1.443%
8	6.887	658	663	676	rVB	233556	384421	9.71%	0.937%
9	7.087	691	697	705	rBV	311759	517209	13.07%	1.260%
10	7.152	706	708	711	rVV4	11306	13782	0.35%	0.034%
11	7.534	761	773	783	rBV	512148	833311	21.05%	2.030%
12	7.811	814	820	827	rBV2	19570	53461	1.35%	0.130%
13	8.063	858	863	871	rVB2	7680	15000	0.38%	0.037%
14	8.258	889	896	908	rBV	312450	612603	15.48%	1.492%
15	8.358	908	913	922	rVB	77180	147219	3.72%	0.359%
16	8.599	951	954	961	rVB8	5421	10099	0.26%	0.025%
17	8.681	961	968	981	rBV	272491	526219	13.29%	1.282%
18	8.834	990	994	1000	rVB7	7080	13607	0.34%	0.033%
19	9.058	1024	1032	1037	rBV	14630	24623	0.62%	0.060%
20	9.240	1059	1063	1070	rBV2	22060	38933	0.98%	0.095%
21	9.387	1080	1088	1102	rBV	230525	447637	11.31%	1.091%
22	9.646	1124	1132	1133	rBV2	19056	33874	0.86%	0.083%
23	9.763	1146	1152	1159	rBV8	5521	16150	0.41%	0.039%
24	9.863	1165	1169	1175	rVB8	5171	10354	0.26%	0.025%
25	9.940	1175	1182	1206	rBV	340185	739088	18.67%	1.801%
26	10.275	1225	1239	1244	rBV	625376	1119539	28.28%	2.727%
27	10.322	1244	1247	1253	rVV	73511	129275	3.27%	0.315%
28	10.387	1253	1258	1262	rVV5	17001	39414	1.00%	0.096%
29	10.452	1262	1269	1285	rVV2	85288	233307	5.89%	0.568%
30	10.663	1299	1305	1308	rBV4	5779	10613	0.27%	0.026%
31	10.840	1324	1335	1346	rBV4	11977	41313	1.04%	0.101%
32	11.116	1374	1382	1386	rBV4	7411	21532	0.54%	0.052%
33	11.163	1388	1390	1399	rVB10	7123	17044	0.43%	0.042%
34	11.275	1405	1409	1413	rVV4	11072	17566	0.44%	0.043%
35	11.640	1464	1471	1474	rBV5	9358	18241	0.46%	0.044%
36	11.681	1474	1478	1484	rVB2	26120	42879	1.08%	0.104%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

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

37	11.869	1506	1510	1516	rVV5	6423	9938	0.25%	0.024%
38	11.946	1516	1523	1530	rVV	49339	78171	1.97%	0.190%
39	12.163	1553	1560	1567	rVB2	18011	37160	0.94%	0.091%
40	12.328	1581	1588	1595	rBV2	7524	23257	0.59%	0.057%
41	12.651	1640	1643	1650	rVV3	12250	19301	0.49%	0.047%
42	12.875	1676	1681	1686	rBV6	10294	20765	0.52%	0.051%
43	12.951	1689	1694	1706	rVB3	40618	98547	2.49%	0.240%
44	13.163	1725	1730	1738	rVB2	9845	22811	0.58%	0.056%
45	13.316	1748	1756	1761	rBV4	22777	66191	1.67%	0.161%
46	13.451	1775	1779	1785	rBV2	23671	40540	1.02%	0.099%
47	13.510	1785	1789	1790	rVV	18718	23831	0.60%	0.058%
48	13.546	1790	1795	1803	rVV	605928	980629	24.77%	2.389%
49	13.610	1803	1806	1811	rVB3	18238	28988	0.73%	0.071%
50	13.698	1818	1821	1824	rBV4	8650	12477	0.32%	0.030%
51	13.834	1837	1844	1861	rVB	787764	1193904	30.16%	2.909%
52	13.969	1863	1867	1873	rVB9	8576	14365	0.36%	0.035%
53	14.040	1874	1879	1884	rBV	54013	65503	1.65%	0.160%
54	14.140	1889	1896	1902	rVV	1086035	1541360	38.94%	3.755%
55	14.193	1902	1905	1909	rVB5	19453	30175	0.76%	0.074%
56	14.369	1931	1935	1943	rBV5	9432	19638	0.50%	0.048%
57	14.451	1944	1949	1962	rBV2	120219	329028	8.31%	0.802%
58	14.551	1963	1966	1969	rVV2	21051	26153	0.66%	0.064%
59	14.775	1996	2004	2010	rBV7	17010	38575	0.97%	0.094%
60	14.828	2010	2013	2017	rVV	16662	21013	0.53%	0.051%
61	14.940	2025	2032	2036	rBV9	10091	24338	0.61%	0.059%
62	15.010	2040	2044	2049	rVB	39282	58500	1.48%	0.143%
63	15.157	2063	2069	2075	rBV	1097990	1514607	38.26%	3.690%
64	15.322	2092	2097	2108	rBV	143200	280096	7.08%	0.682%
65	15.428	2111	2115	2119	rVB4	22802	35805	0.90%	0.087%
66	15.528	2126	2132	2145	rVB	609861	807896	20.41%	1.968%
67	15.857	2184	2188	2190	rBV2	15599	18077	0.46%	0.044%
68	15.892	2190	2194	2195	rBV	55283	65738	1.66%	0.160%
69	16.057	2217	2222	2226	rBV8	15556	33544	0.85%	0.082%
70	16.104	2226	2230	2237	rVB	56801	87538	2.21%	0.213%
71	16.357	2268	2273	2278	rBV9	20365	40940	1.03%	0.100%
72	16.481	2292	2294	2297	rBV4	11909	14456	0.37%	0.035%
73	16.604	2310	2315	2322	rVB	79128	127169	3.21%	0.310%
74	16.704	2329	2332	2337	rVB2	30126	40646	1.03%	0.099%
75	16.763	2339	2342	2350	rVB4	30496	56222	1.42%	0.137%
76	16.857	2354	2358	2365	rBV2	82539	114045	2.88%	0.278%

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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

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77	16.945	2365	2373	2377	rBV	1417246	1888075	47.70%	4.600%
78	16.987	2377	2380	2385	rVB	199595	287010	7.25%	0.699%
79	17.051	2385	2391	2402	rVB2	930880	1590592	40.18%	3.875%
80	17.145	2404	2407	2414	rBV8	38509	87500	2.21%	0.213%
81	17.469	2459	2462	2466	rVB2	45164	55630	1.41%	0.136%
82	17.769	2510	2513	2518	rVB5	68466	100413	2.54%	0.245%
83	17.828	2518	2523	2526	rBV	308581	414057	10.46%	1.009%
84	17.886	2528	2533	2545	rVB2	1168696	1828512	46.20%	4.455%
85	18.045	2557	2560	2565	rBV2	51955	68901	1.74%	0.168%
86	18.192	2580	2585	2591	rBV6	105264	243996	6.16%	0.594%
87	18.681	2666	2668	2673	rVB3	51840	53300	1.35%	0.130%
88	18.851	2694	2697	2703	rVB3	60944	92401	2.33%	0.225%
89	18.910	2705	2707	2710	rBV3	46873	47219	1.19%	0.115%
90	19.092	2732	2738	2742	rBV	443034	803158	20.29%	1.957%
91	19.222	2756	2760	2766	rVB2	377049	560546	14.16%	1.366%
92	19.439	2792	2797	2801	rBV	1284682	1818728	45.95%	4.431%
93	19.510	2805	2809	2813	rVB	178684	282158	7.13%	0.687%
94	21.057	3067	3072	3080	rBV2	801414	1653272	41.77%	4.028%
95	21.380	3121	3127	3131	rBV3	1426667	2477200	62.58%	6.035%
96	22.192	3261	3265	3270	rBV2	354056	579603	14.64%	1.412%
97	23.680	3512	3518	3525	rVB	1179673	2379784	60.12%	5.798%
98	23.769	3527	3533	3545	rBV3	1014647	2848888	71.97%	6.941%
99	24.551	3660	3666	3674	rVB	762402	2022813	51.10%	4.928%
100	25.874	3883	3891	3908	rVB3	957812	3958240	100.00%	9.643%

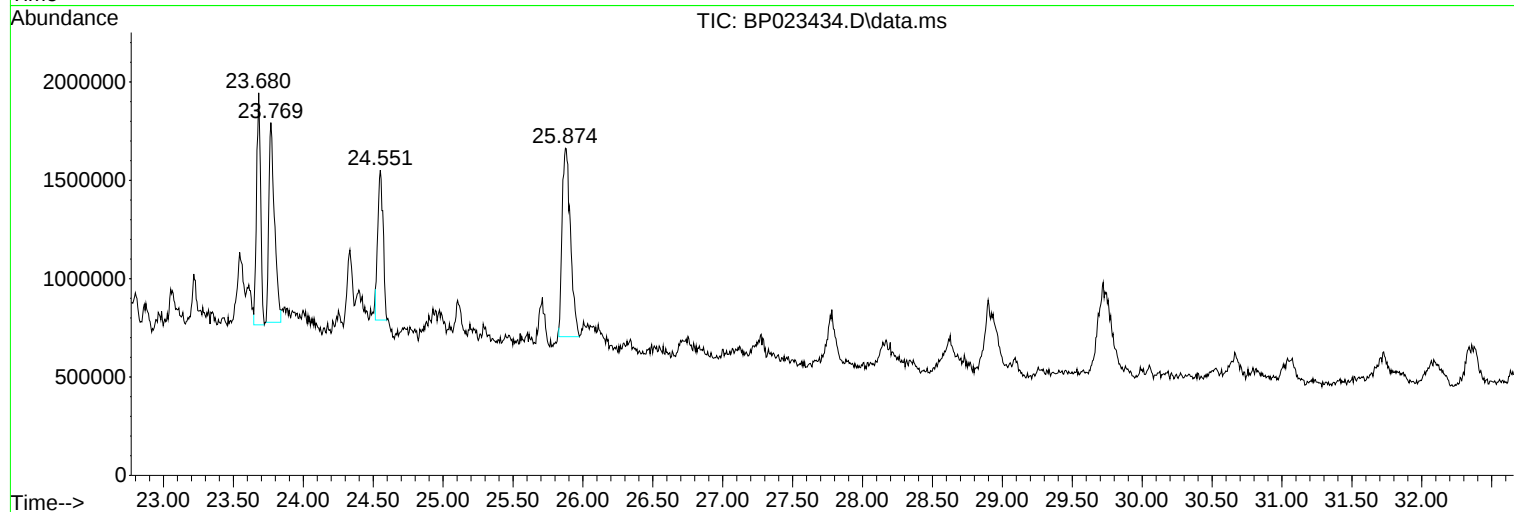
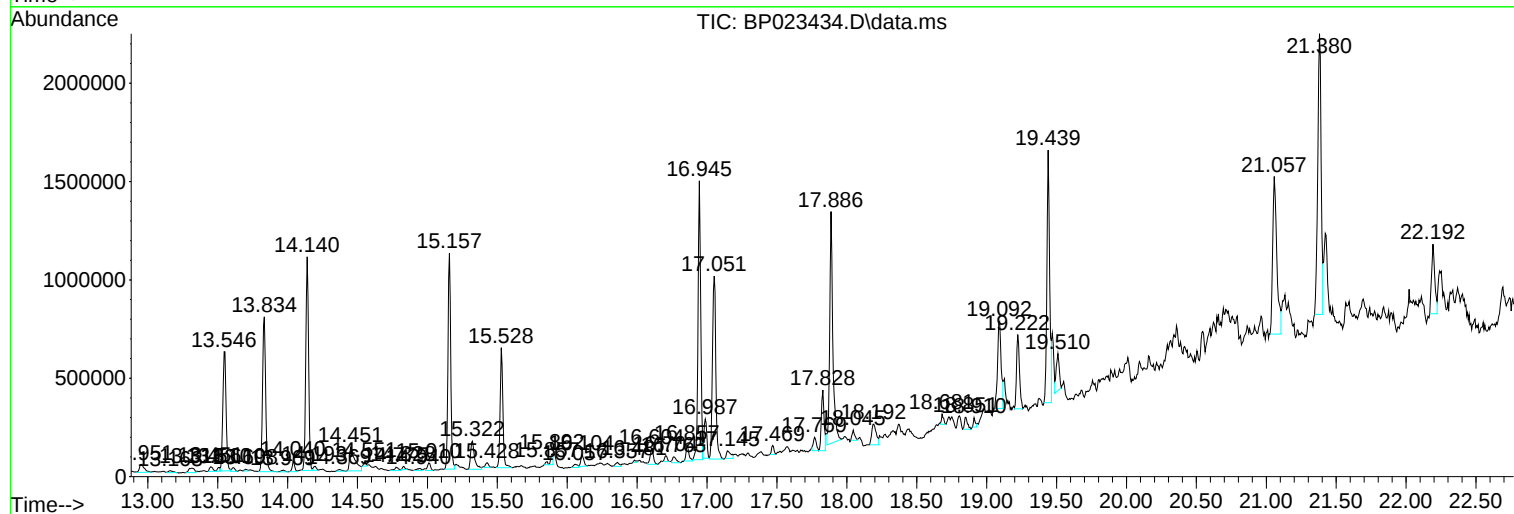
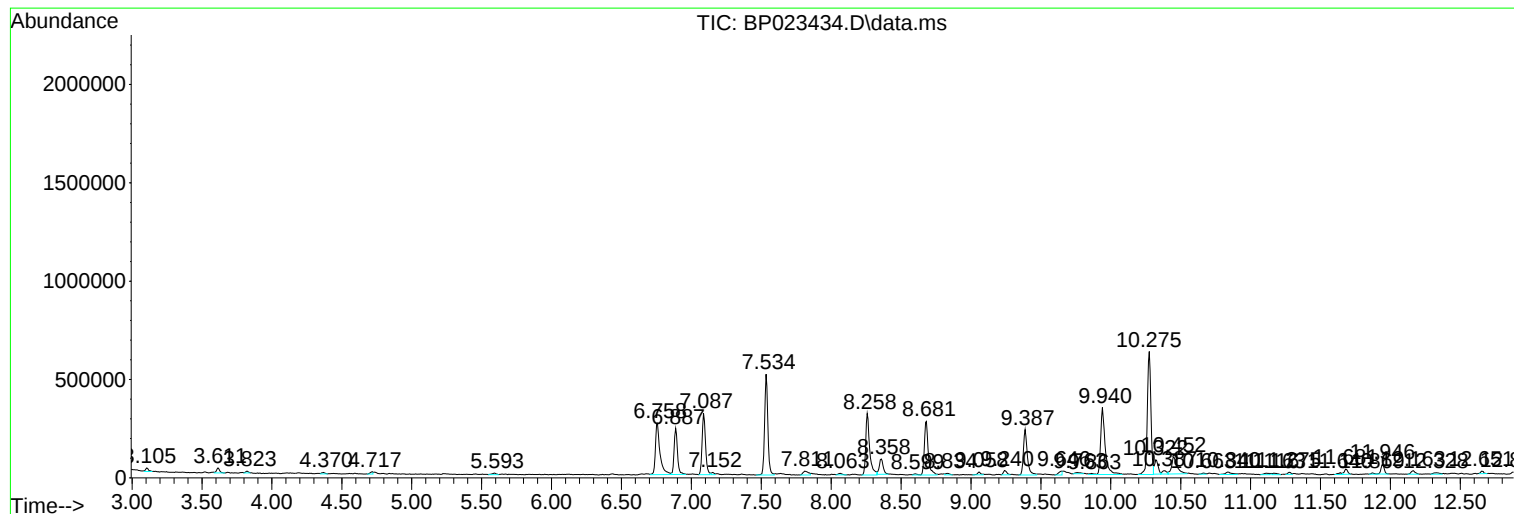
Sum of corrected areas: 41046733

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Instrument :
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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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
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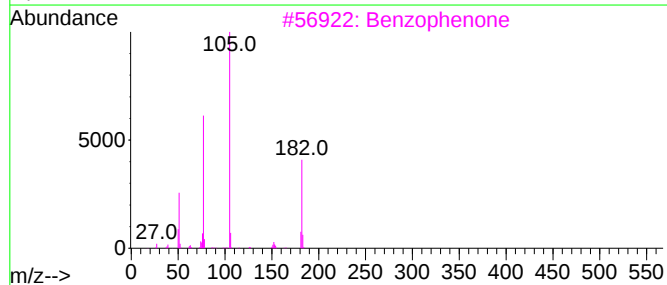
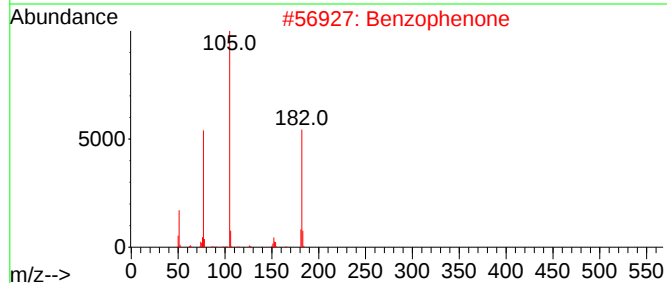
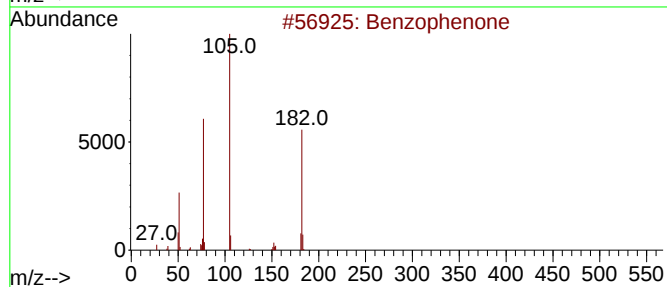
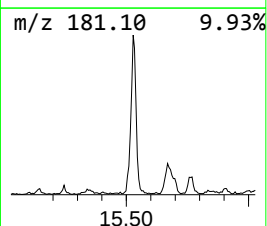
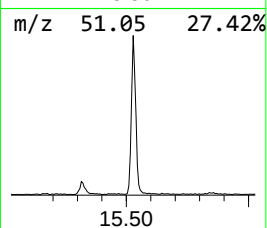
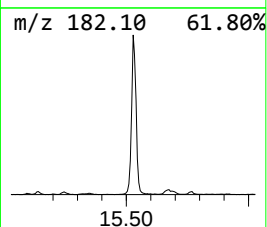
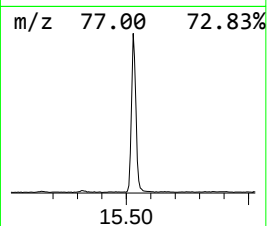
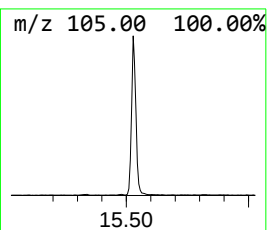
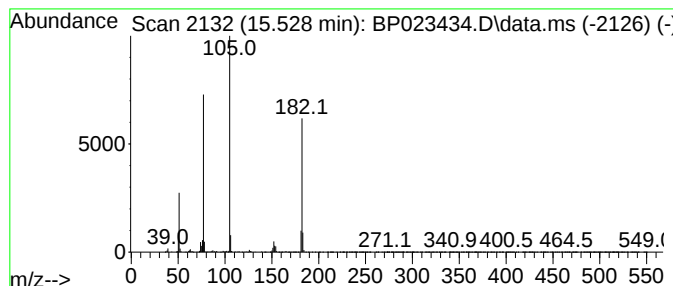
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.528	10.48 ng/ul	807896	Acenaphthene-d10	14.140

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	91
5			Benzophenone	182	C13H10O	000119-61-9	91



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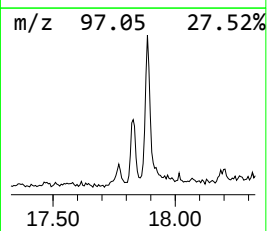
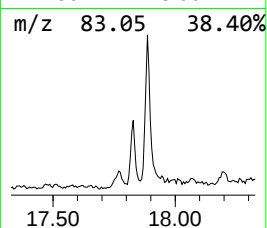
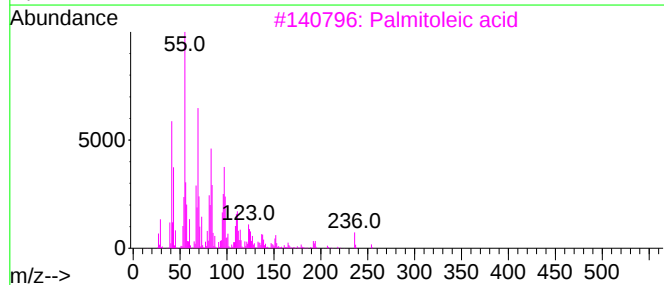
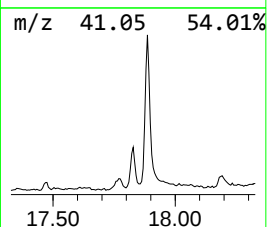
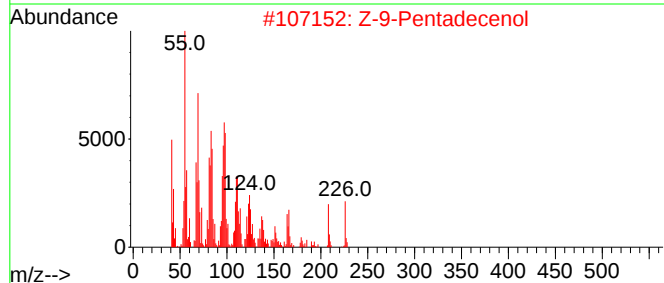
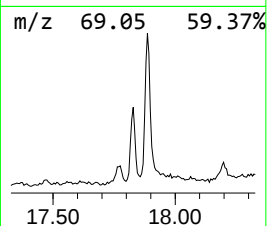
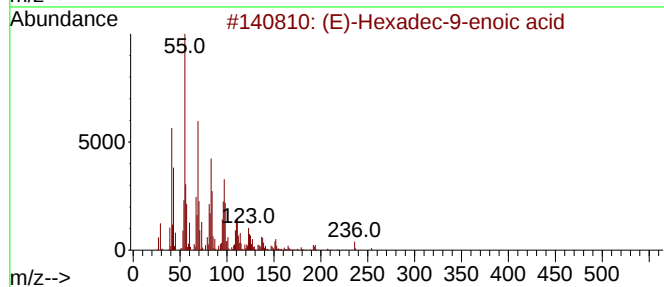
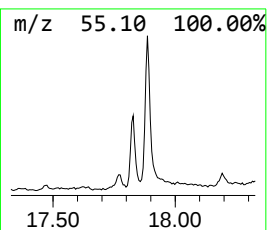
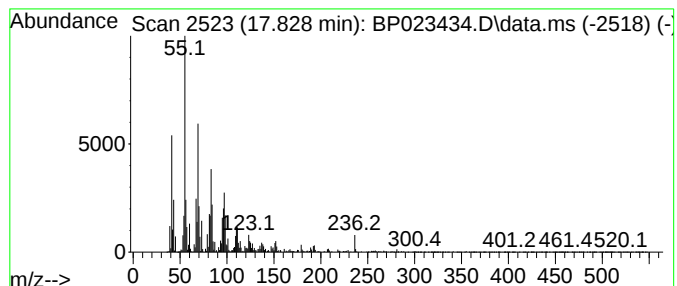
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (E)-Hexadec-9-enoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.828	4.39 ng/ul	414057	Phenanthrene-d10	16.945

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	98
2			Z-9-Pentadecenol	226	C15H30O	1000130-76-9	93
3			Palmitoleic acid	254	C16H30O2	000373-49-9	91
4			Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	91
5			trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	91



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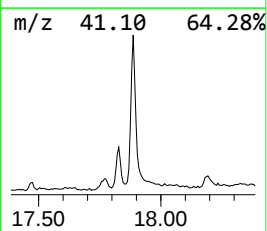
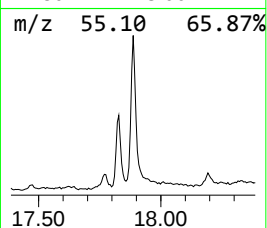
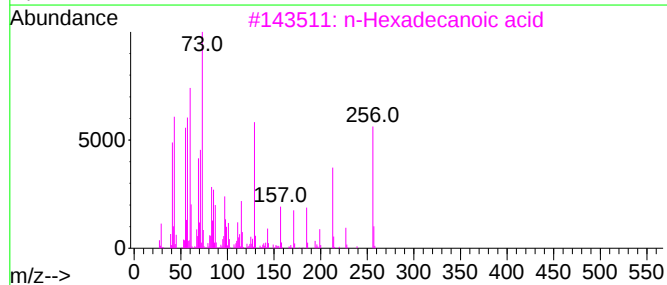
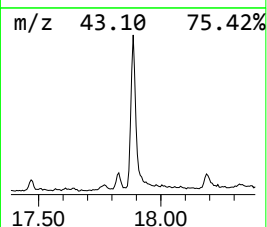
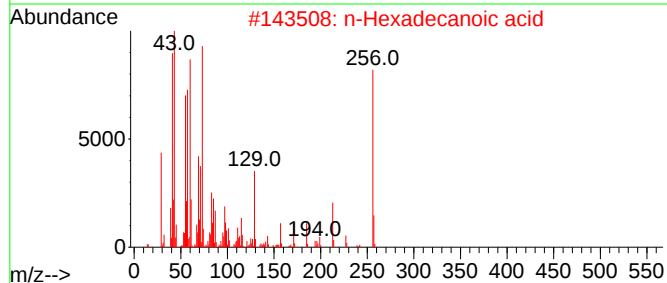
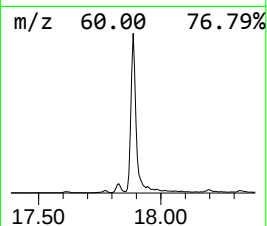
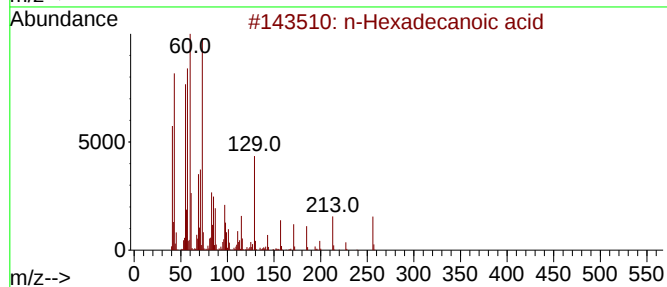
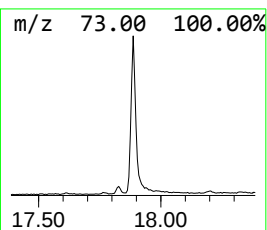
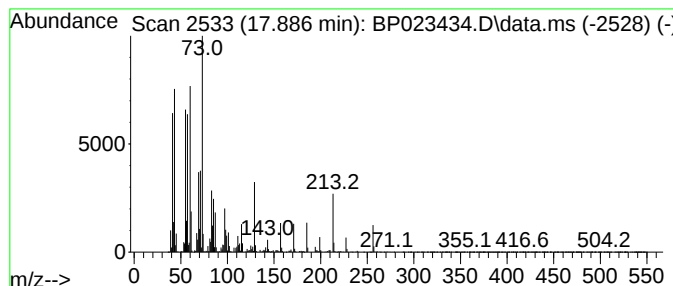
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.887	19.37 ng/ul	1828510	Phenanthrene-d10	16.945

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	99
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	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023434.D
Acq On : 16 Dec 2024 14:27
Operator : RC/JU
Sample : P5153-21
Misc :
ALS Vial : 8 Sample Multiplier: 1

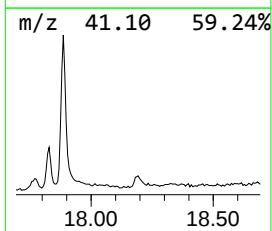
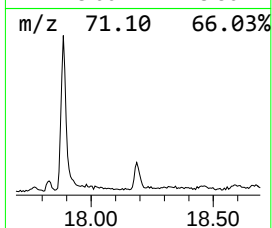
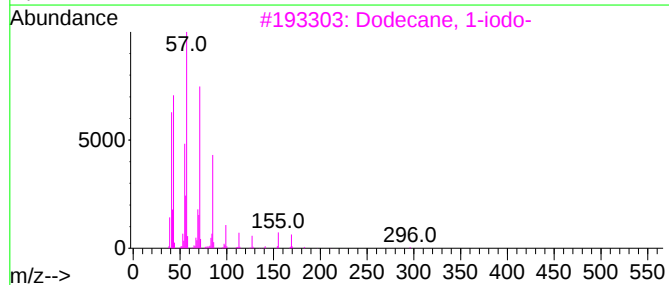
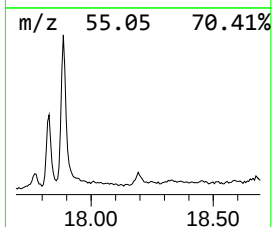
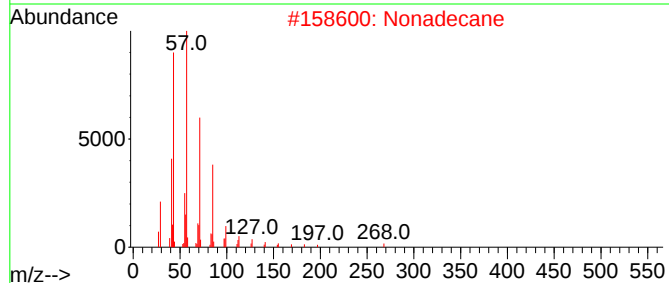
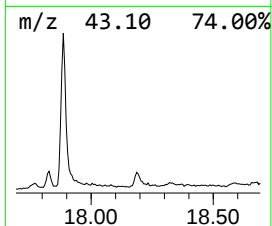
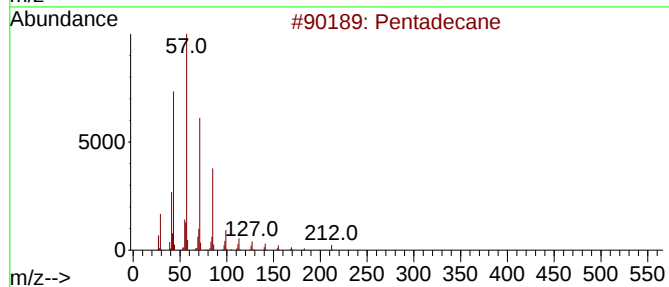
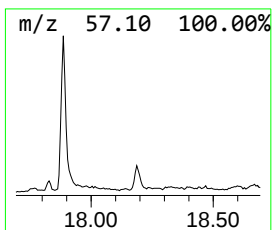
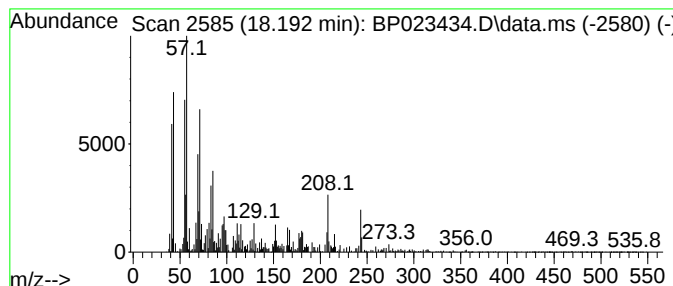
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.192	2.58 ng/ul	243996	Phenanthrene-d10		16.945	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecane		212	C15H32	000629-62-9	93
2	Nonadecane		268	C19H40	000629-92-5	55
3	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	51
4	Hentriacontane, 3-methyl-		451	C32H66	004981-99-1	51
5	Docosane		310	C22H46	000629-97-0	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023434.D
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Sample : P5153-21
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Instrument :
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ClientSampleId :
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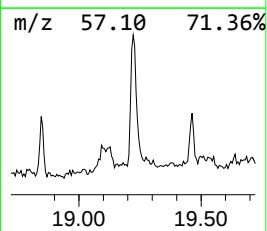
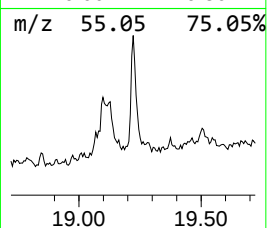
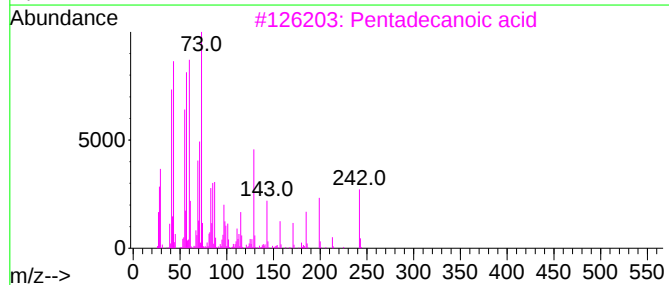
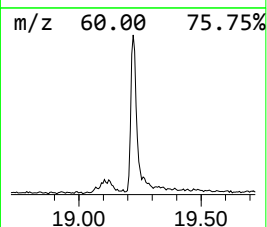
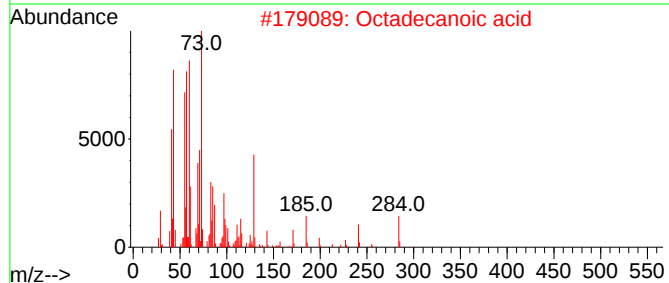
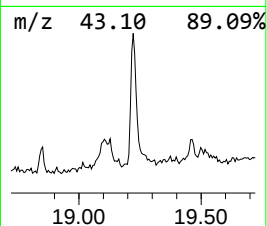
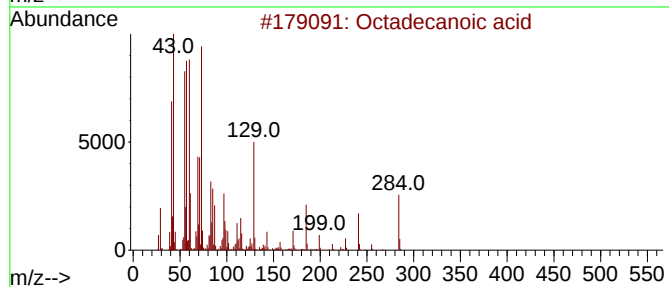
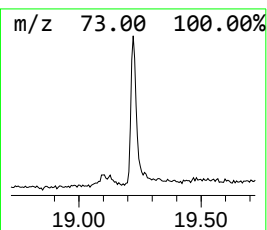
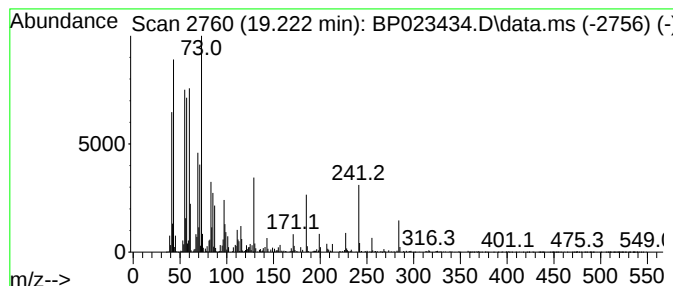
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.222	4.53 ng/ul	560546	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	98
2			Octadecanoic acid	284	C18H36O2	000057-11-4	95
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	83
5			Octadecanoic acid	284	C18H36O2	000057-11-4	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023434.D
Acq On : 16 Dec 2024 14:27
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Sample : P5153-21
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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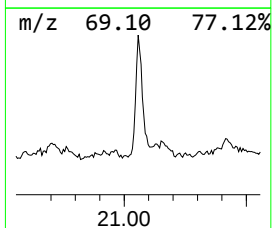
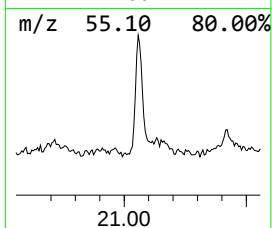
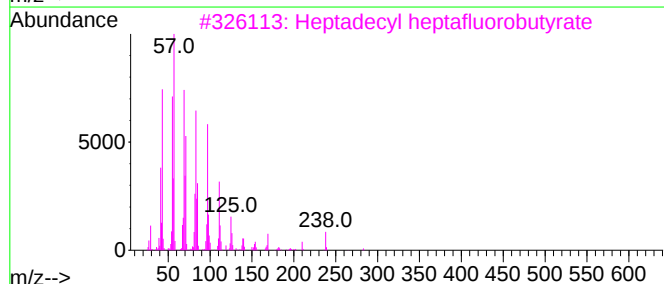
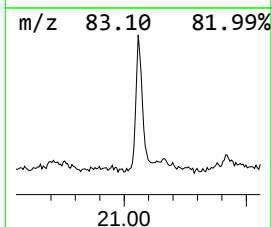
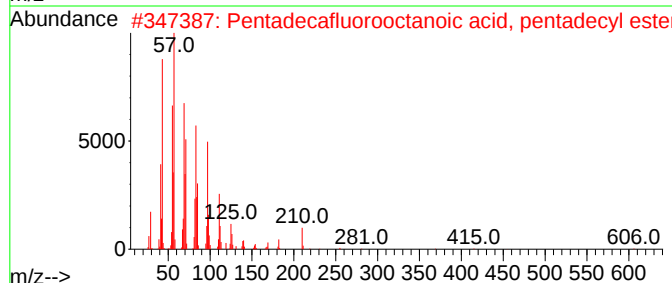
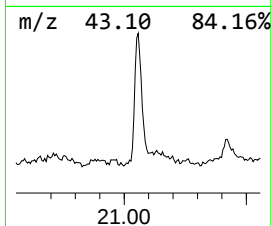
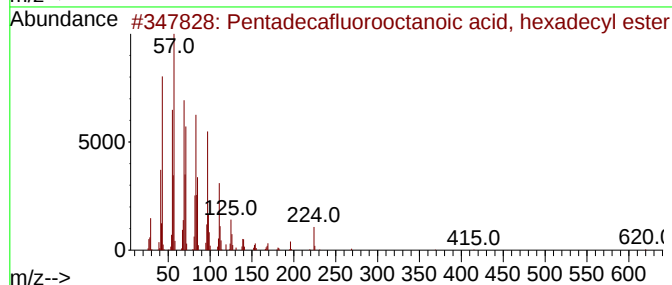
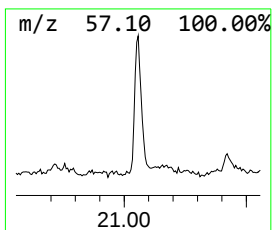
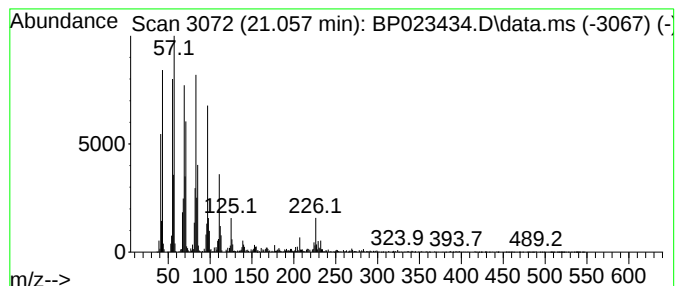
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Pentadecafluorooctanoic aci... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.057	13.35 ng/ul	1653270	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	94
2			Pentadecafluorooctanoic acid, pe...	624	C23H31F15O2	1000406-04-5	94
3			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
4			Cyclohexadecane	224	C16H32	000295-65-8	93
5			Nonacos-1-ene	406	C29H58	018835-35-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
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Sample : P5153-21
Misc :
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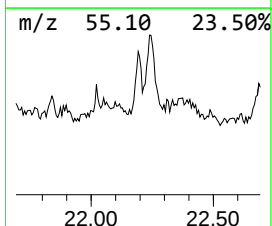
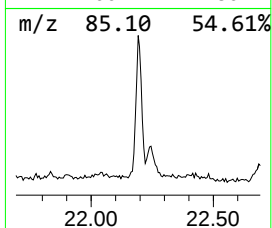
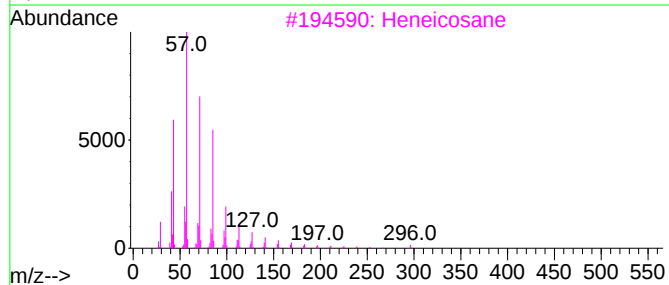
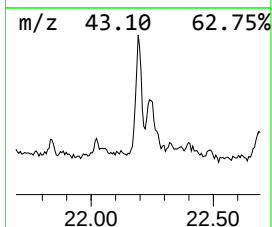
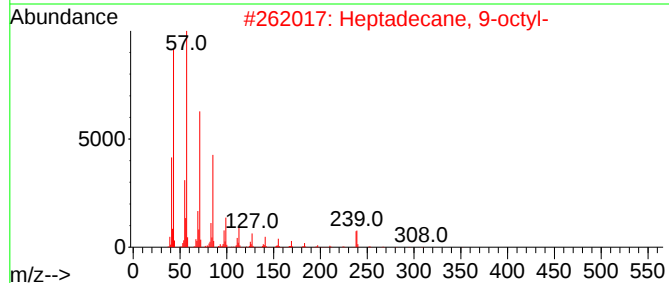
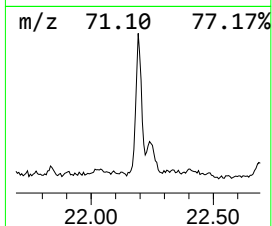
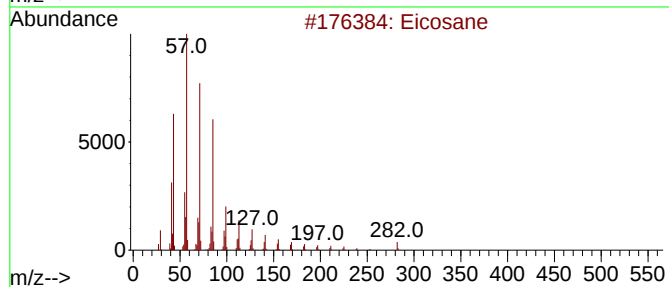
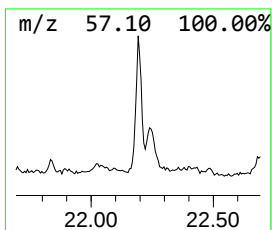
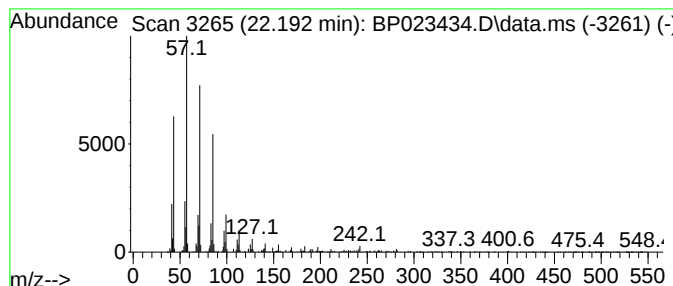
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.192	4.68 ng/ul	579603	Chrysene-d12	21.380	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	97
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3	Heneicosane	296	C21H44	000629-94-7	91
4	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	87
5	Nonadecane	268	C19H40	000629-92-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023434.D
Acq On : 16 Dec 2024 14:27
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Instrument :
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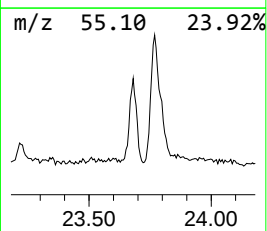
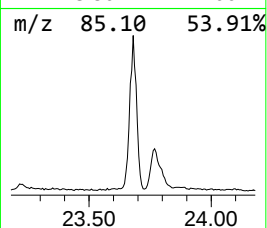
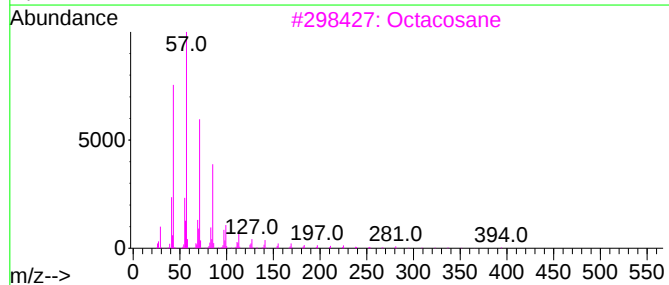
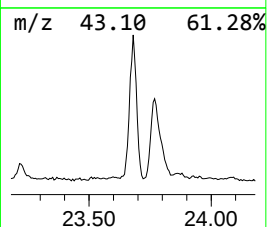
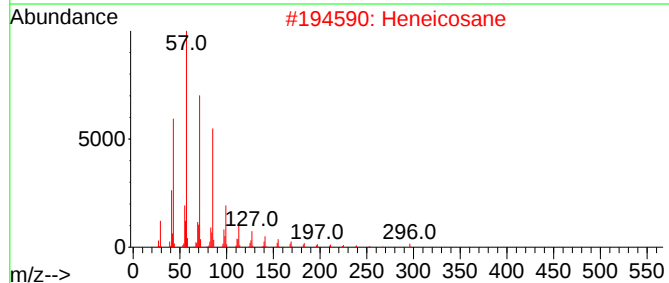
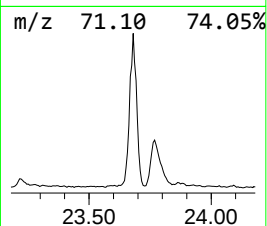
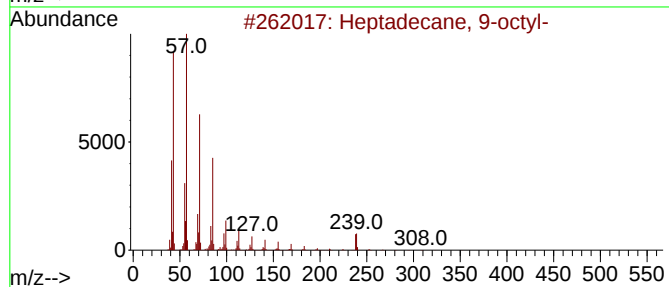
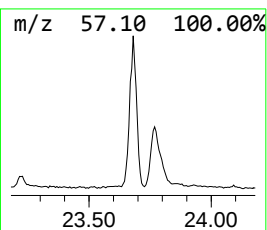
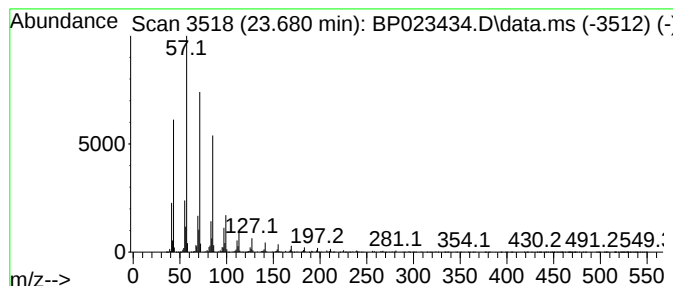
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Quant Title : SVOA CALIBRATION

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.
23.680	23.53 ng/ul	2379780	Perylene-d12	24.551

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2		Heneicosane	296	C21H44	000629-94-7	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
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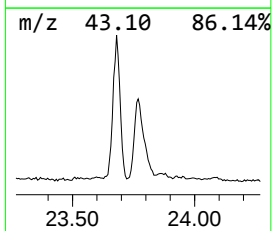
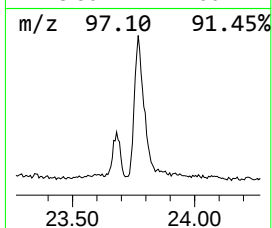
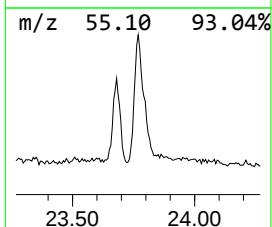
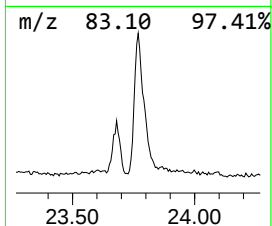
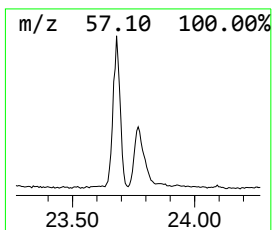
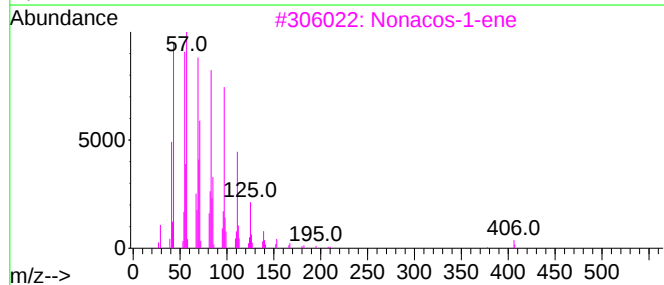
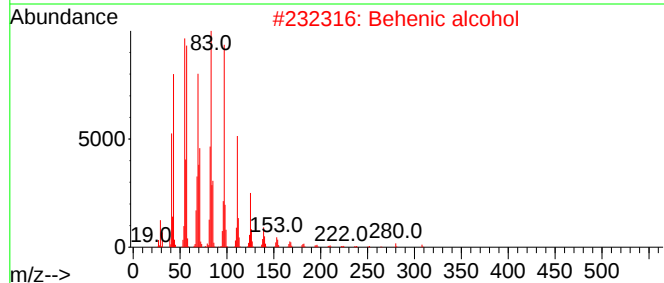
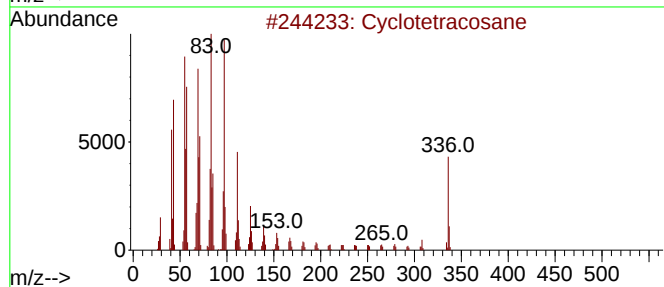
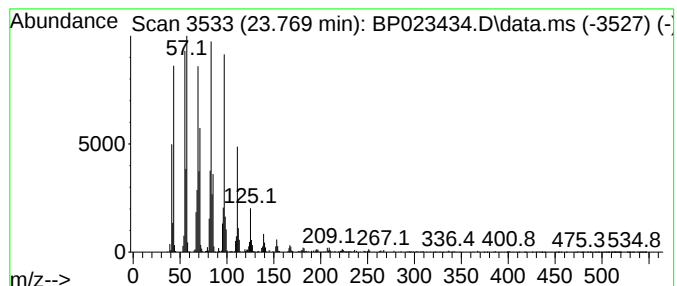
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Cyclic23.769 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.769	28.17 ng/ul	2848890	Perylene-d12	24.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	97
2			Behenic alcohol	326	C22H46O	000661-19-8	94
3			Nonacos-1-ene	406	C29H58	018835-35-3	94
4			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5			Heptacos-1-ene	378	C27H54	015306-27-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023434.D
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Sample : P5153-21
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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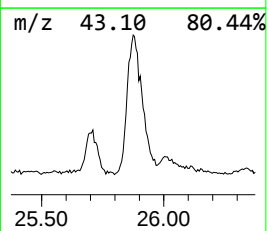
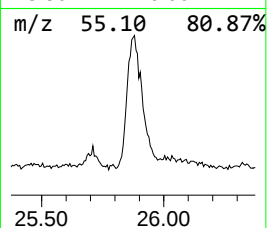
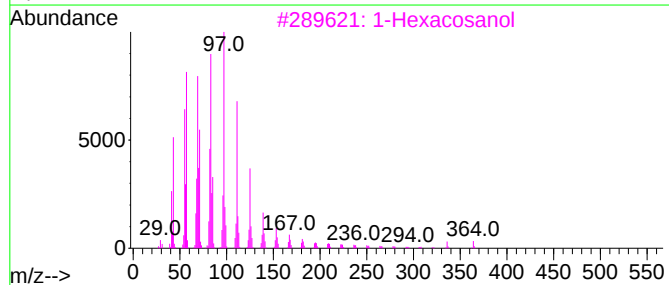
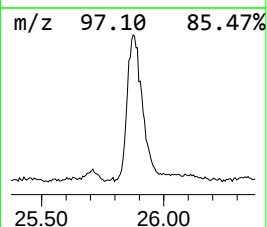
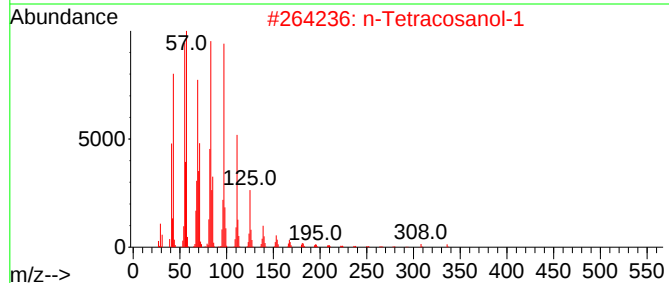
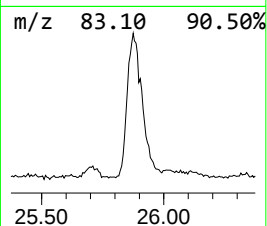
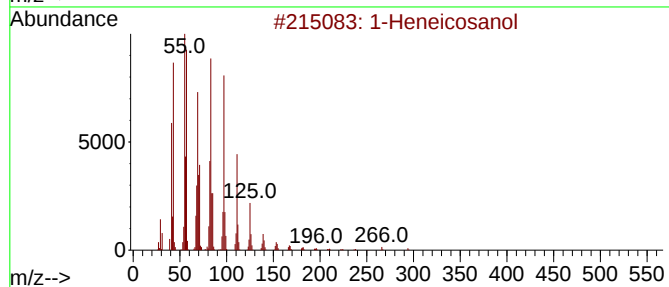
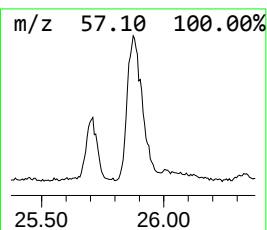
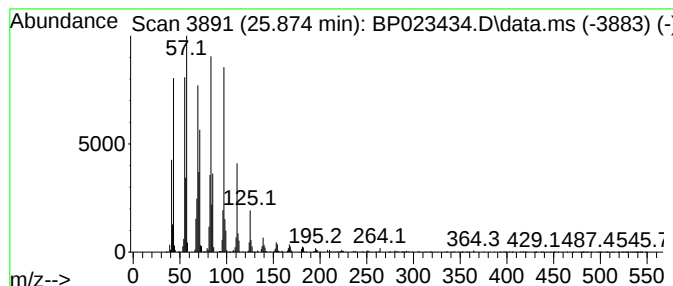
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 1-Heneicosanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.874	39.14 ng/ul	3958240	Perylene-d12	24.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	93
2			n-Tetracosanol-1	354	C24H50O	000506-51-4	93
3			1-Hexacosanol	382	C26H54O	000506-52-5	91
4			Triacontan-15-ol	438	C30H62O	031849-26-0	91
5			Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023434.D
Acq On : 16 Dec 2024 14:27
Operator : RC/JU
Sample : P5153-21
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E28R4

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	15.528	10.5	ng/ul	807896	3	14.140	1541360	20.0
(E)-Hexadec-9-e...	17.828	4.4	ng/ul	414057	4	16.945	1888080	20.0
n-Hexadecanoic ...	17.887	19.4	ng/ul	1828510	4	16.945	1888080	20.0
(DEL) Alkane: S...	18.192	2.6	ng/ul	243996	4	16.945	1888080	20.0
Octadecanoic acid	19.222	4.5	ng/ul	560546	5	21.380	2477200	20.0
Pentadecafluoro...	21.057	13.4	ng/ul	1653270	5	21.380	2477200	20.0
(DEL) Alkane: S...	22.192	4.7	ng/ul	579603	5	21.380	2477200	20.0
(DEL) Alkane: S...	23.680	23.5	ng/ul	2379780	6	24.551	2022810	20.0
(DEL) Alkane: C...	23.769	28.2	ng/ul	2848890	6	24.551	2022810	20.0
1-Heneicosanol	25.874	39.1	ng/ul	3958240	6	24.551	2022810	20.0