

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013164.D
 Acq On : 20 Dec 2022 13:38
 Operator : CG/JU
 Sample : N6003-14
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXQ5

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

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

Signal : TIC: BP013164.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.334	22	25	35	rVB	204700	265226	1.67%	0.138%
2	3.628	72	75	82	rVB	45899	63667	0.40%	0.033%
3	3.764	95	98	113	rBV	1364786	2032126	12.82%	1.056%
4	3.870	113	116	122	rVB	69417	91967	0.58%	0.048%
5	3.999	133	138	144	rVB	65408	103654	0.65%	0.054%
6	4.093	151	154	158	rVB	40552	52854	0.33%	0.027%
7	4.476	215	219	222	rVB5	16860	25415	0.16%	0.013%
8	4.523	225	227	235	rVB5	17626	23144	0.15%	0.012%
9	5.040	309	315	323	rBV2	393629	576714	3.64%	0.300%
10	5.134	327	331	334	rBV4	13348	20855	0.13%	0.011%
11	6.546	568	571	576	rBV6	11559	21176	0.13%	0.011%
12	6.628	579	585	591	rVB	129812	205240	1.29%	0.107%
13	6.893	625	630	633	rBV6	10477	18490	0.12%	0.010%
14	6.934	633	637	642	rBV4	26049	45500	0.29%	0.024%
15	7.111	659	667	680	rBV	3038469	4921426	31.04%	2.558%
16	7.234	682	688	690	rVV	50864	77549	0.49%	0.040%
17	7.281	690	696	705	rVB	2606252	3970552	25.04%	2.064%
18	7.387	710	714	719	rVB3	37820	57388	0.36%	0.030%
19	7.481	723	730	740	rVV	3175060	5026188	31.70%	2.612%
20	7.569	740	745	754	rVB5	24276	47301	0.30%	0.025%
21	7.775	775	780	786	rVB5	10376	20694	0.13%	0.011%
22	7.952	802	810	818	rBV	2052878	3283768	20.71%	1.707%
23	8.016	818	821	826	rVV	18454	24952	0.16%	0.013%
24	8.664	921	931	943	rBV	3664871	5765767	36.36%	2.997%
25	8.781	946	951	957	rVB	58362	102332	0.65%	0.053%
26	9.122	1002	1009	1020	rBV	3021161	4924739	31.06%	2.560%
27	9.281	1032	1036	1042	rVB9	9156	18335	0.12%	0.010%
28	9.528	1070	1078	1084	rVB2	64596	122157	0.77%	0.063%
29	9.846	1123	1132	1144	rBV	2568513	4278758	26.99%	2.224%
30	9.987	1153	1156	1162	rBV8	8839	17151	0.11%	0.009%
31	10.063	1164	1169	1177	rVB4	73159	141974	0.90%	0.074%
32	10.387	1216	1224	1233	rBV	3937809	6459397	40.74%	3.357%
33	10.546	1246	1251	1258	rVB	48105	81882	0.52%	0.043%
34	10.769	1279	1289	1296	rBV	2852604	4750367	29.96%	2.469%
35	10.905	1301	1312	1329	rVV	2764538	4871876	30.73%	2.532%
36	11.299	1374	1379	1382	rBV6	16135	28268	0.18%	0.015%

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

37	11.434	1394	1402	1409	rBV6	8067	16149	0.10%	0.008%
38	11.552	1414	1422	1427	rBV7	17101	38704	0.24%	0.020%
39	12.081	1507	1512	1516	rVV2	14318	21902	0.14%	0.011%
40	12.175	1524	1528	1533	rVB	40639	57601	0.36%	0.030%

41	12.257	1538	1542	1545	rBV2	15748	24829	0.16%	0.013%
42	12.293	1545	1548	1552	rVV5	11647	21444	0.14%	0.011%
43	12.352	1552	1558	1564	rVB2	69051	115883	0.73%	0.060%
44	12.963	1657	1662	1664	rBV6	11450	17695	0.11%	0.009%
45	13.122	1684	1689	1692	rBV	15333	20919	0.13%	0.011%

46	13.410	1729	1738	1746	rVB2	85040	134258	0.85%	0.070%
47	13.493	1746	1752	1753	rVV5	17745	33874	0.21%	0.018%
48	13.516	1753	1756	1760	rVV3	27502	45343	0.29%	0.024%
49	13.581	1761	1767	1771	rVB6	19363	45127	0.28%	0.023%
50	13.916	1818	1824	1830	rBV4	63437	95713	0.60%	0.050%

51	14.004	1830	1839	1848	rBV	6401883	8861125	55.89%	4.606%
52	14.122	1853	1859	1864	rVB5	36650	58736	0.37%	0.031%
53	14.181	1865	1869	1873	rBV	50338	63404	0.40%	0.033%
54	14.228	1874	1877	1881	rBV3	13470	18301	0.12%	0.010%
55	14.293	1881	1888	1900	rVV	7541005	10979913	69.25%	5.707%

56	14.451	1911	1915	1917	rBV5	20765	30040	0.19%	0.016%
57	14.481	1917	1920	1925	rVB	46423	59141	0.37%	0.031%
58	14.599	1933	1940	1947	rVB	4096023	5988243	37.77%	3.112%
59	14.793	1966	1973	1994	rBV	2260688	3555333	22.42%	1.848%
60	14.946	1995	1999	2004	rVV2	48269	81665	0.52%	0.042%

61	15.010	2004	2010	2015	rVB5	72544	139060	0.88%	0.072%
62	15.375	2068	2072	2078	rBV8	15811	38964	0.25%	0.020%
63	15.434	2080	2082	2087	rVB	23850	23856	0.15%	0.012%
64	15.593	2102	2109	2116	rBV	9133298	13108596	82.67%	6.813%
65	15.710	2122	2129	2142	rBV	2657186	3681053	23.22%	1.913%

66	15.828	2146	2149	2156	rVB2	43879	77133	0.49%	0.040%
67	15.957	2165	2171	2182	rVB	3708865	4969084	31.34%	2.583%
68	16.163	2203	2206	2210	rVB6	19173	24540	0.15%	0.013%
69	16.222	2212	2216	2222	rVB8	23940	33035	0.21%	0.017%
70	16.298	2226	2229	2232	rBV5	19324	24183	0.15%	0.013%

71	16.522	2262	2267	2273	rVB	185210	251115	1.58%	0.131%
72	16.734	2299	2303	2308	rBV3	39796	56368	0.36%	0.029%
73	16.881	2326	2328	2333	rVB5	14059	20642	0.13%	0.011%
74	17.234	2384	2388	2391	rBV2	61140	88212	0.56%	0.046%
75	17.298	2397	2399	2402	rBV4	20709	20075	0.13%	0.010%

76	17.351	2402	2408	2419	rVV	4467969	6464048	40.77%	3.360%
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 Title : SVOA CALIBRATION

77	17.451	2419	2425	2434	rVV	8382983	12165550	76.73%	6.323%
78	17.540	2437	2440	2443	rVB4	44944	61190	0.39%	0.032%
79	17.634	2452	2456	2465	rVB5	27679	45373	0.29%	0.024%
80	18.010	2517	2520	2524	rVB3	30521	34545	0.22%	0.018%
81	18.110	2534	2537	2541	rVB4	44693	54543	0.34%	0.028%
82	18.163	2543	2546	2551	rBV5	69508	98099	0.62%	0.051%
83	18.222	2552	2556	2565	rBV	1068074	1557598	9.82%	0.810%
84	18.557	2609	2613	2619	rBV4	56696	94389	0.60%	0.049%
85	18.634	2622	2626	2630	rBV	272824	303977	1.92%	0.158%
86	19.116	2703	2708	2712	rBV5	74311	121562	0.77%	0.063%
87	19.204	2720	2723	2727	rBV2	64282	82370	0.52%	0.043%
88	19.263	2729	2733	2738	rBV	116145	156093	0.98%	0.081%
89	19.328	2739	2744	2746	rBV2	169204	309624	1.95%	0.161%
90	19.357	2747	2749	2753	rVB	174787	206615	1.30%	0.107%
91	19.451	2762	2765	2778	rVB	266737	421089	2.66%	0.219%
92	19.686	2799	2805	2816	rBV	9808839	13376647	84.36%	6.953%
93	21.128	3046	3050	3055	rBV2	1815960	2642854	16.67%	1.374%
94	21.439	3098	3103	3108	rBV	4375085	5848384	36.88%	3.040%
95	22.045	3204	3206	3209	rBV	293206	366606	2.31%	0.191%
96	23.133	3387	3391	3397	rVB	1062516	1684197	10.62%	0.875%
97	23.763	3491	3498	3512	rVB2	6822063	14274436	90.03%	7.419%
98	23.922	3519	3525	3538	rVB2	3116965	6596900	41.61%	3.429%
99	24.539	3624	3630	3640	rVB	1882821	4095767	25.83%	2.129%
100	25.480	3785	3790	3814	rVB3	3749013	15855744	100.00%	8.241%

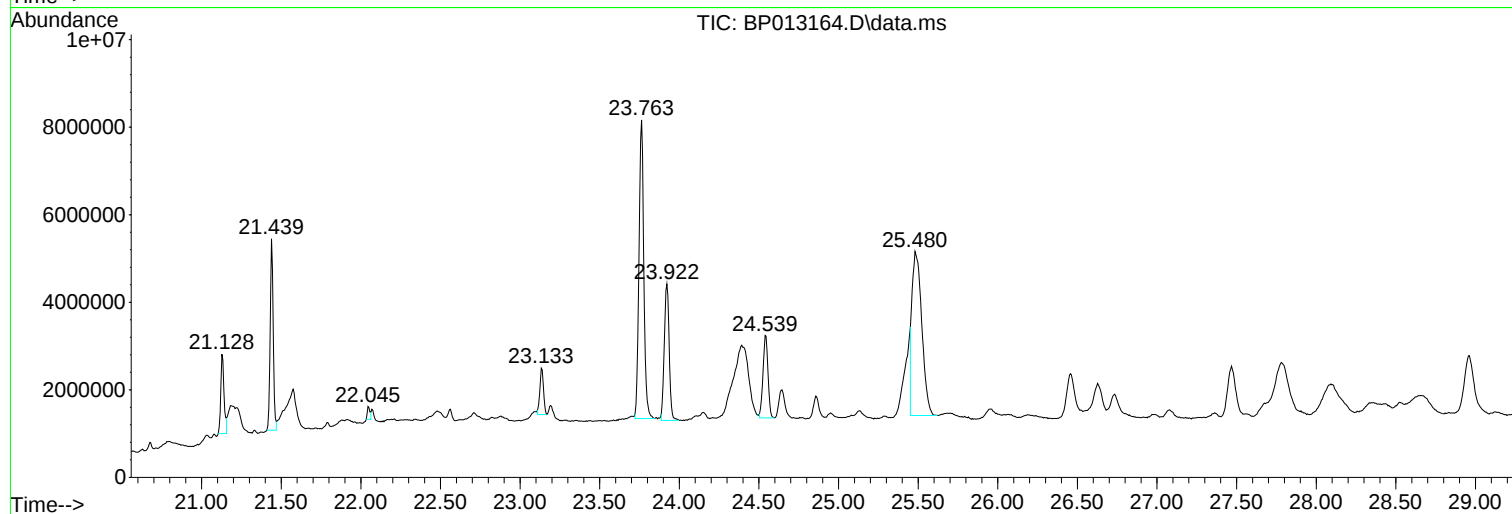
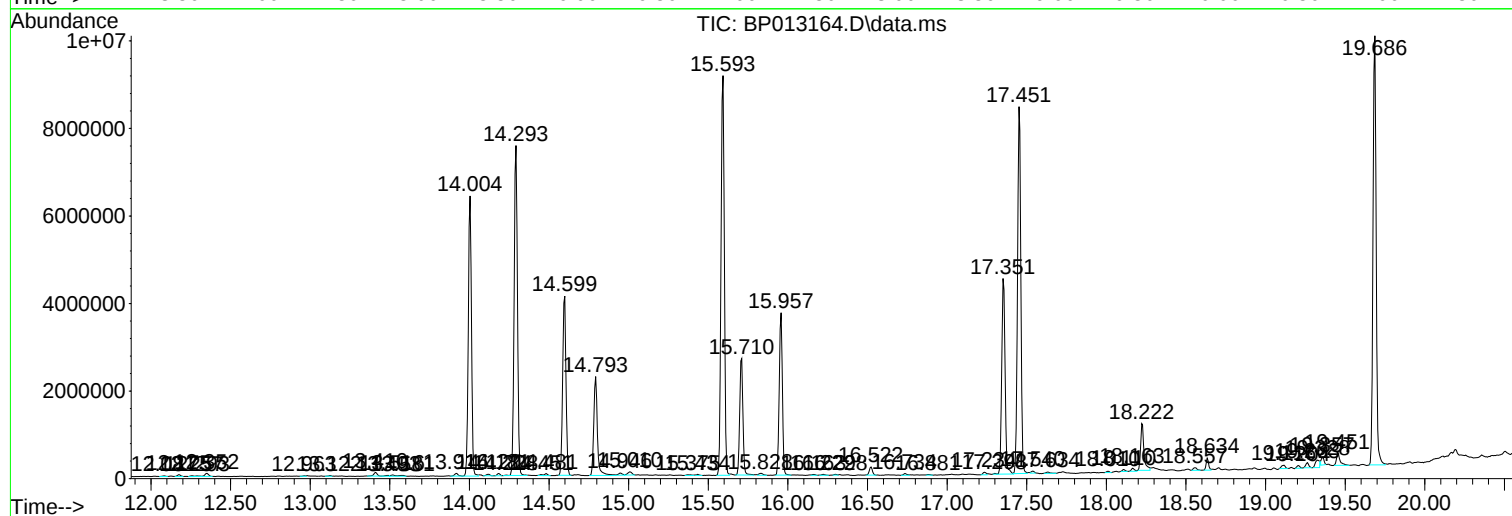
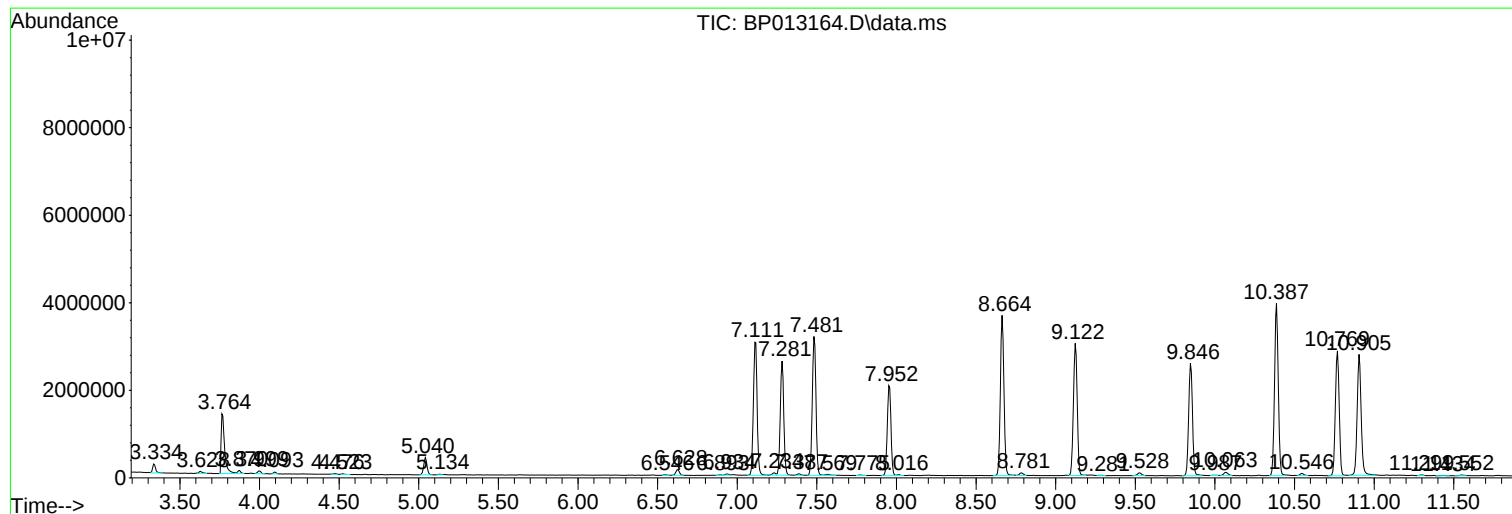
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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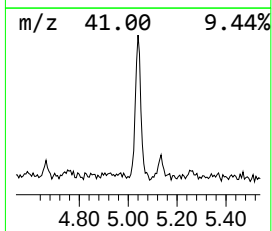
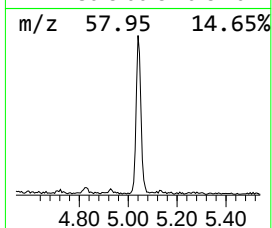
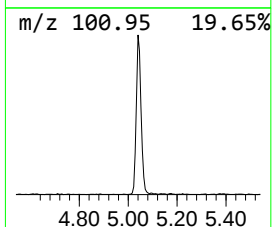
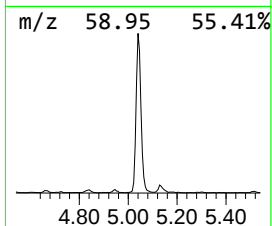
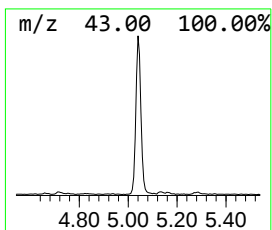
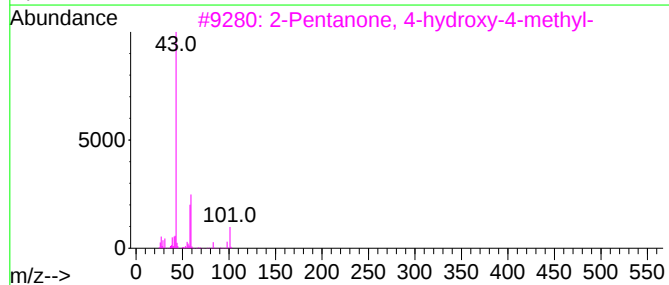
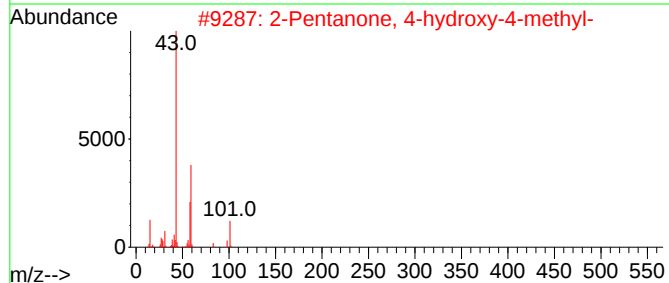
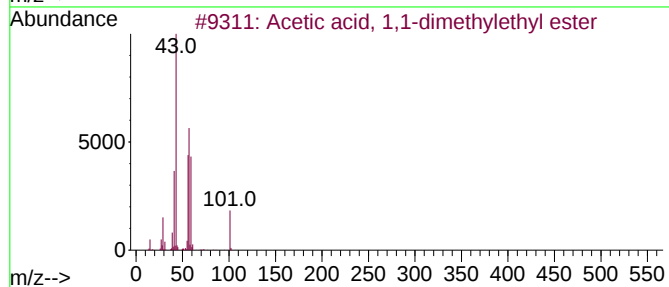
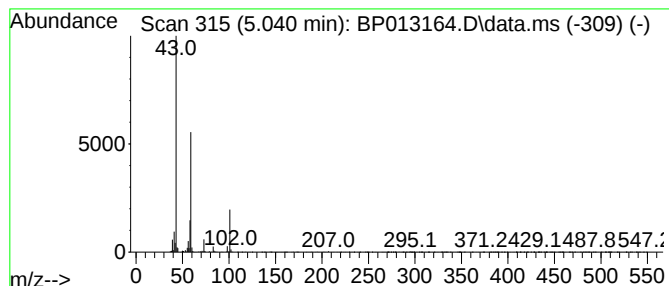
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Acetic acid, 1,1-dimethylet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.040	3.51 ng/ul	576714	1,4-Dichlorobenzene-d4	7.958

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	50
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
4			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
5			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	33



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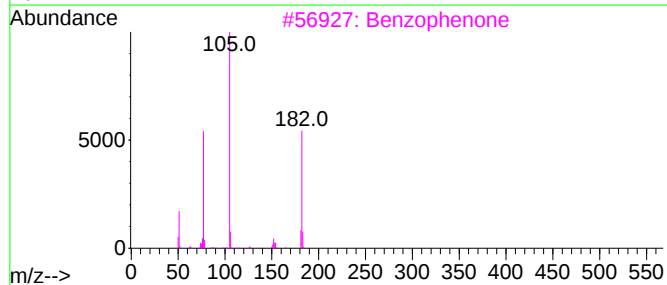
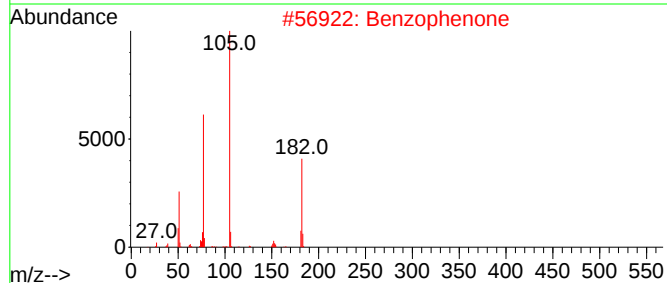
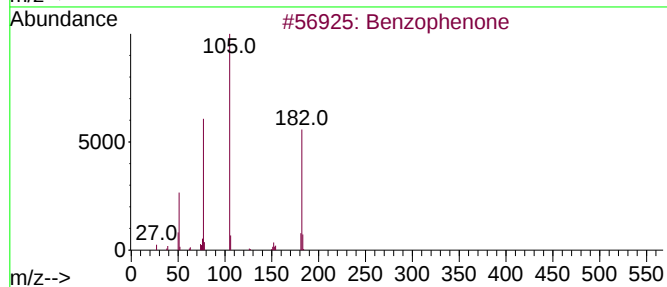
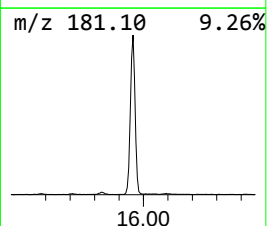
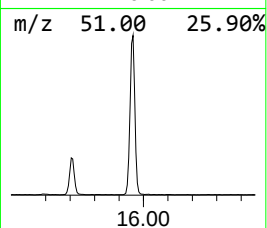
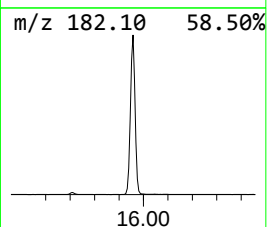
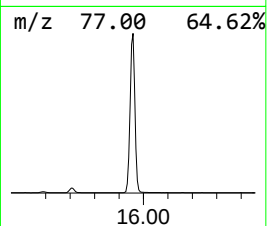
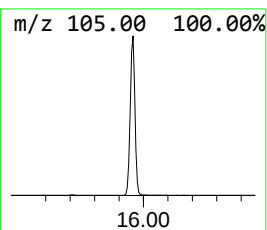
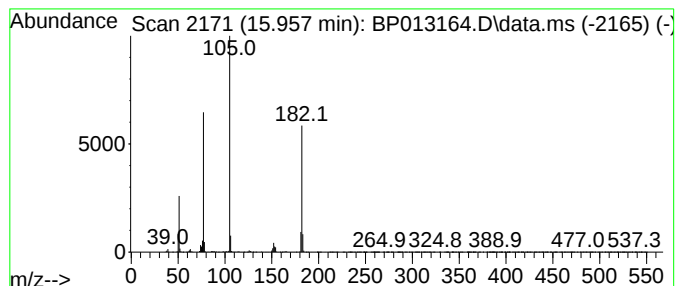
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.957	16.60 ng/ul	4969080	Acenaphthene-d10	14.599

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	93
5			Benzophenone	182	C13H10O	000119-61-9	91



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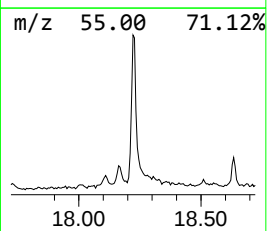
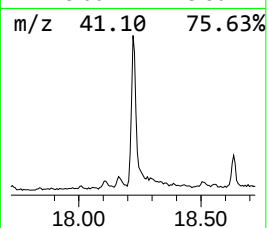
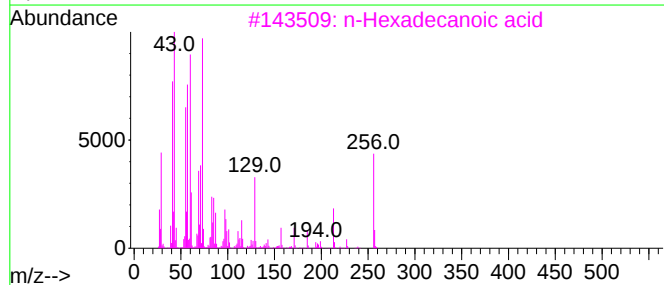
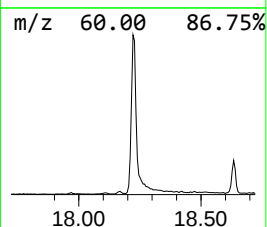
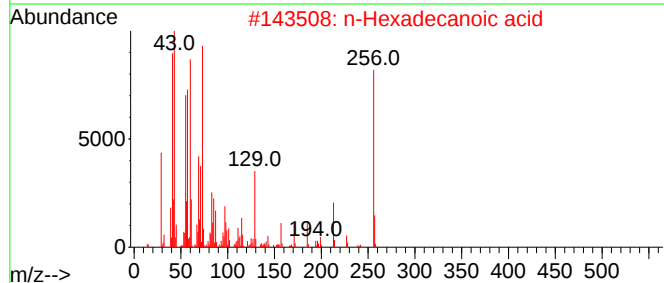
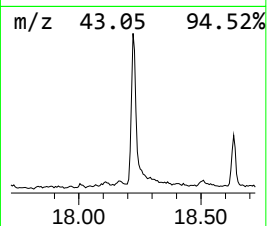
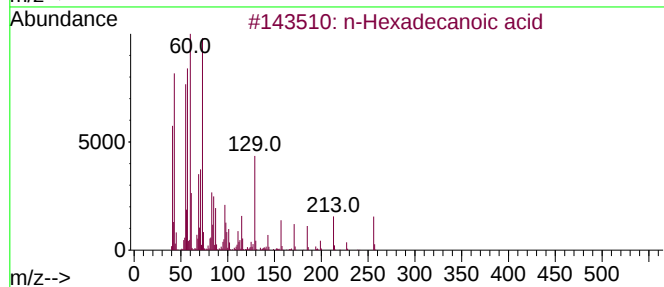
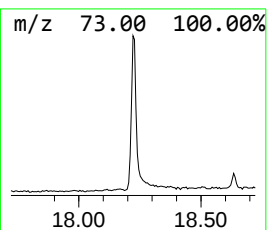
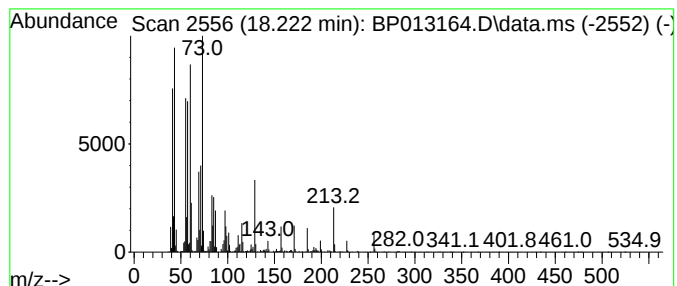
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.222	4.82 ng/ul	1557600	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5			Tridecanoic acid	214	C13H26O2	000638-53-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013164.D
Acq On : 20 Dec 2022 13:38
Operator : CG/JU
Sample : N6003-14
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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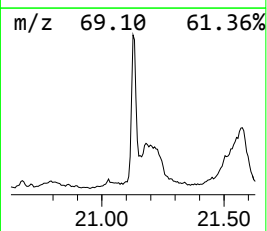
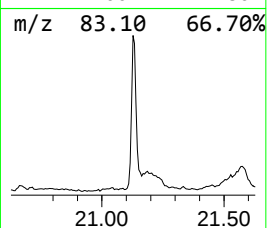
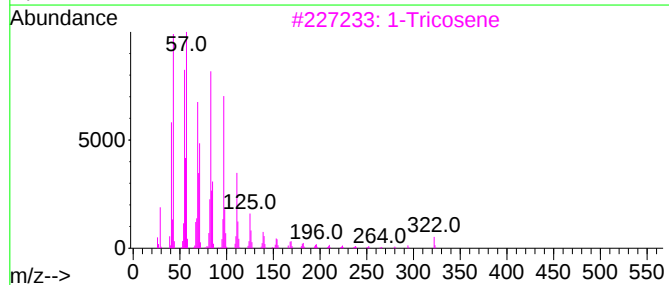
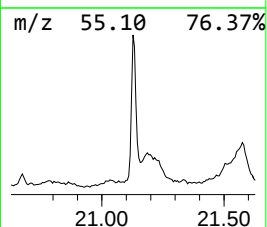
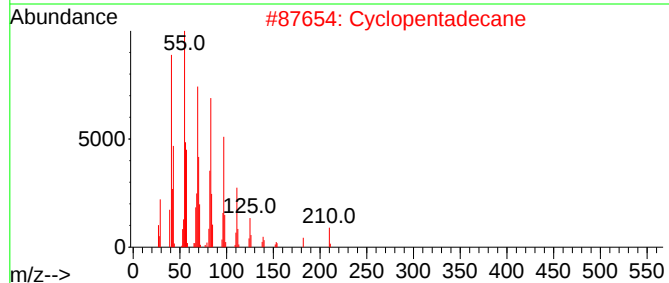
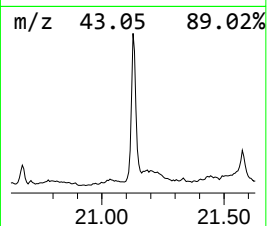
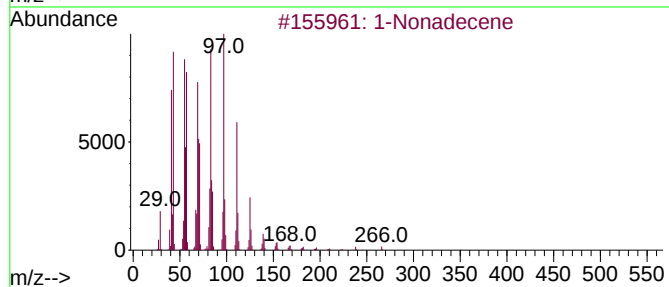
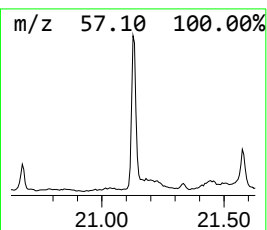
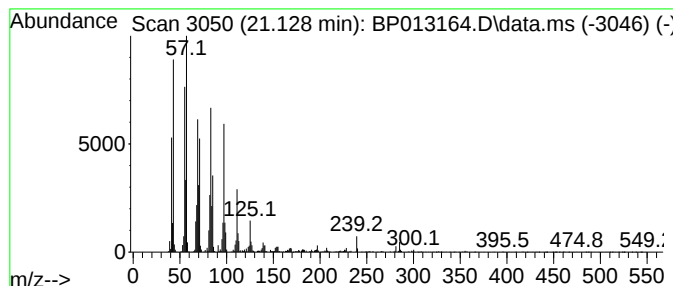
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.128	9.04 ng/ul	2642850	Chrysene-d12	21.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene			266	C19H38	018435-45-5	97
2	Cyclopentadecane			210	C15H30	000295-48-7	97
3	1-Tricosene			322	C23H46	018835-32-0	95
4	Pentadecafluorooctanoic acid, oc...			666	C26H37F15O2	1000406-04-8	94
5	Pentafluoropropionic acid, hepta...			402	C20H35F5O2	959218-78-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013164.D
Acq On : 20 Dec 2022 13:38
Operator : CG/JU
Sample : N6003-14
Misc :
ALS Vial : 45 Sample Multiplier: 1

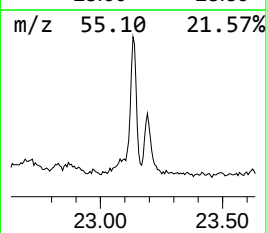
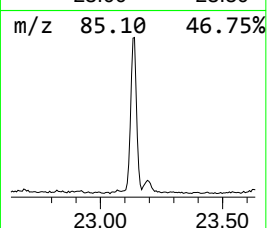
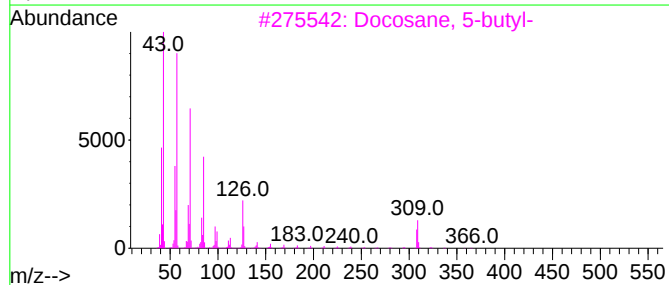
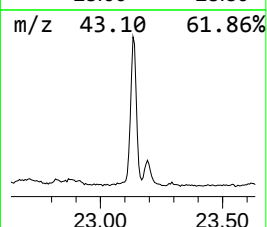
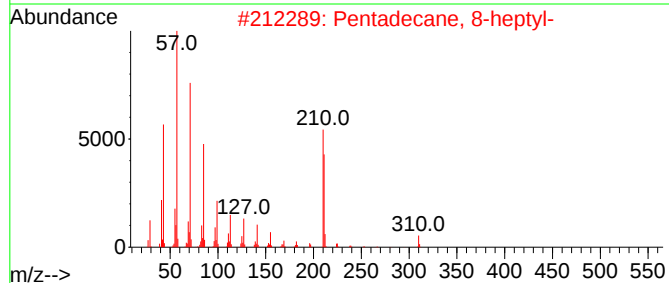
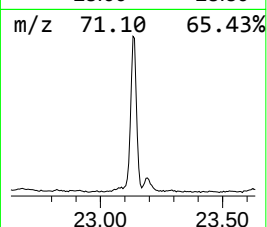
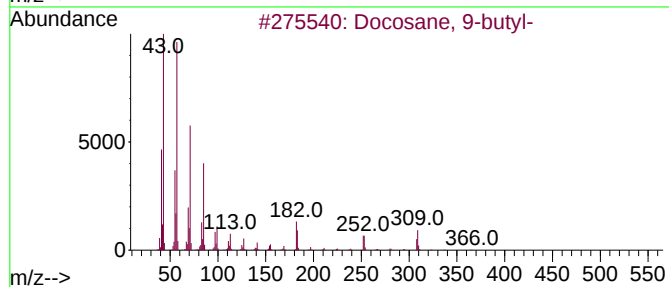
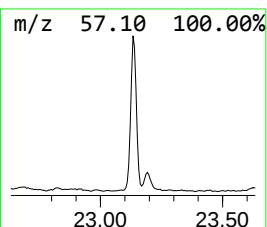
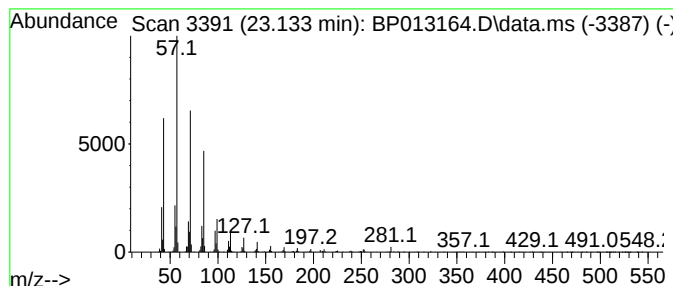
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.133	5.11 ng/ul	1684200	Perylene-d12	23.922	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 9-butyl-	366	C26H54	055282-14-9	94
2	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	94
3	Docosane, 5-butyl-	366	C26H54	055282-16-1	93
4	Heptacosane	380	C27H56	000593-49-7	91
5	Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013164.D
Acq On : 20 Dec 2022 13:38
Operator : CG/JU
Sample : N6003-14
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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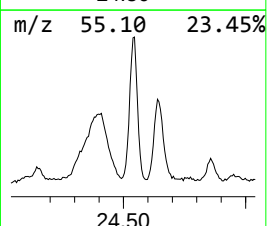
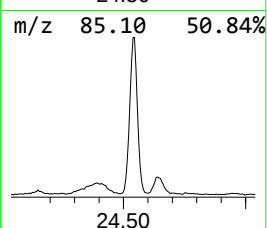
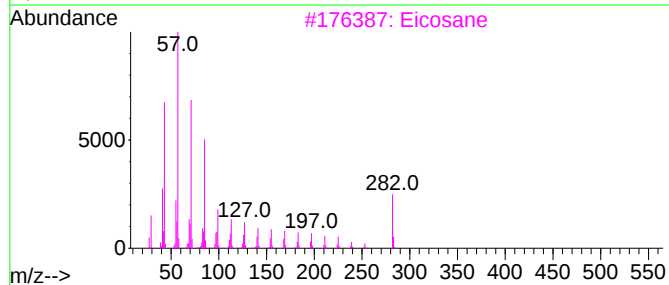
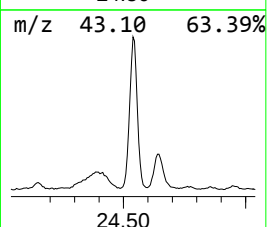
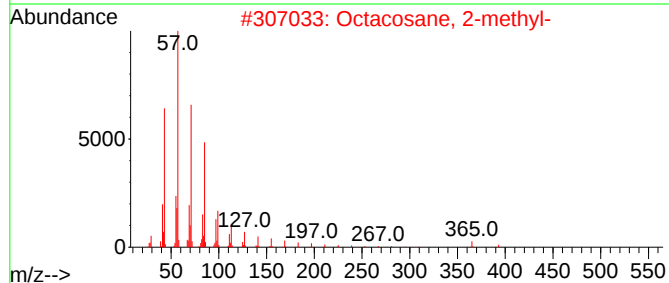
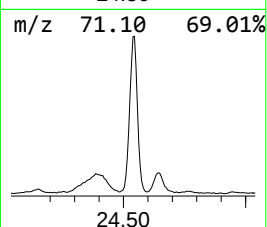
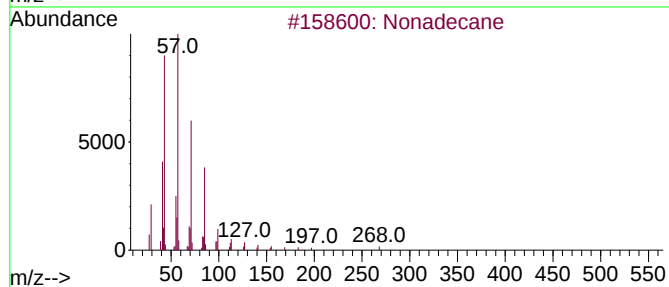
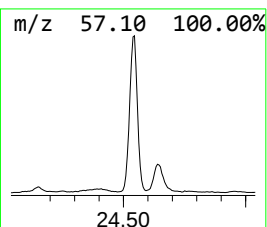
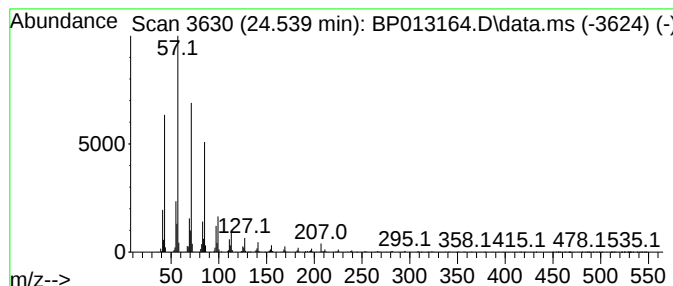
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.539	12.42 ng/ul	4095770	Perylene-d12	23.922

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	95
2		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
3		Eicosane	282	C20H42	000112-95-8	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Hexatriacontane	507	C36H74	000630-06-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013164.D
Acq On : 20 Dec 2022 13:38
Operator : CG/JU
Sample : N6003-14
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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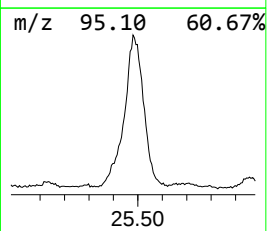
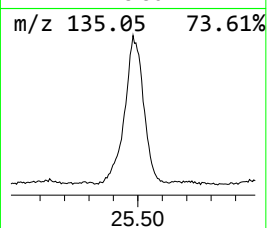
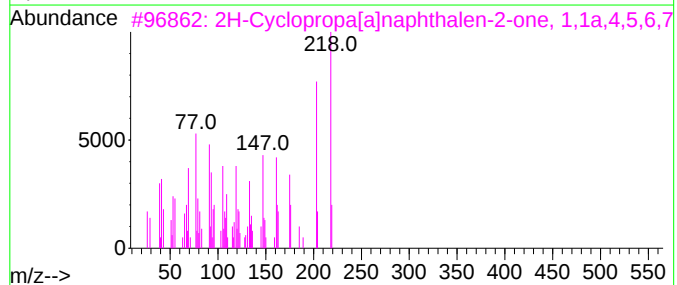
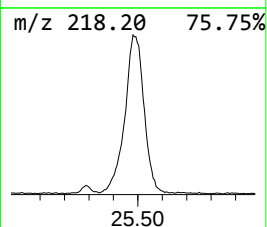
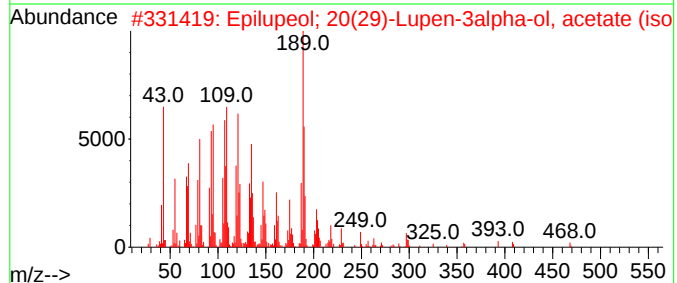
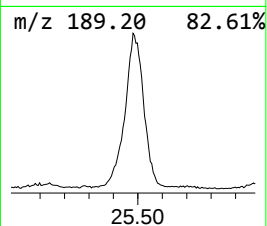
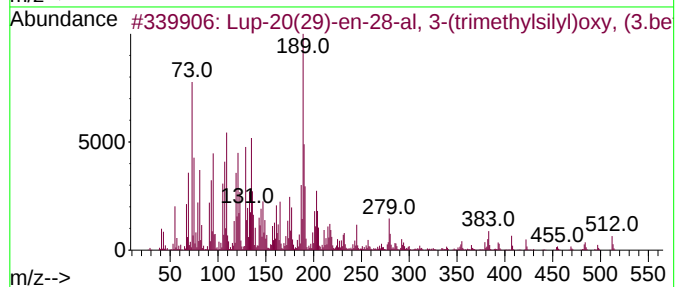
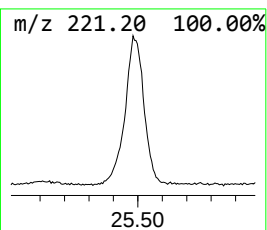
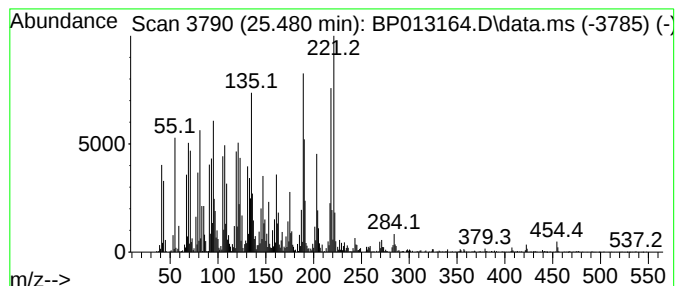
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Lup-20(29)-en-28-al, 3-(tri... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.480	48.07 ng/ul	15855700	Perylene-d12	23.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Lup-20(29)-en-28-al, 3-(trimethy...	512	C33H56O2Si	1000414-39-2	50
2			Epilupeol; 20(29)-Lupen-3alpha-o...	468	C32H52O2	1010513-01-8	41
3			2H-Cyclopropa[a]naphthalen-2-one...	218	C15H22O	006831-17-0	41
4			Methyl 6-(methylsulfonyl)-1H-ind...	221	C11H11NO2S	202584-20-1	35
5			1-Hydroxypyrene	218	C16H10O	005315-79-7	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013164.D
Acq On : 20 Dec 2022 13:38
Operator : CG/JU
Sample : N6003-14
Misc :
ALS Vial : 45 Sample Multiplier: 1

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Acetic acid, 1,...	5.040	3.5	ng/ul	576714	1	7.958	3283770	20.0
Benzophenone	15.957	16.6	ng/ul	4969080	3	14.599	5988240	20.0
n-Hexadecanoic ...	18.222	4.8	ng/ul	1557600	4	17.351	6464050	20.0
(DEL) Alkane: S...	21.128	9.0	ng/ul	2642850	5	21.439	5848380	20.0
(DEL) Alkane: S...	23.133	5.1	ng/ul	1684200	6	23.922	6596900	20.0
(DEL) Alkane: S...	24.539	12.4	ng/ul	4095770	6	23.922	6596900	20.0
Lup-20(29)-en-2...	25.480	48.1	ng/ul	15855700	6	23.922	6596900	20.0