

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
 Data File : BN031107.D
 Acq On : 21 Apr 2024 13:49
 Operator : MA/JU
 Sample : P2161-03
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CORD4

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_N\Methods\SFAM-EPA-BN041924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN031107.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.146	24	27	32	rVB	16882	20798	0.29%	0.038%
2	3.558	96	97	114	rBV	81393	147427	2.09%	0.271%
3	3.864	147	149	156	rVB3	7568	10527	0.15%	0.019%
4	4.975	331	338	352	rBV	1073029	1765713	24.98%	3.249%
5	6.616	613	617	623	rBV6	6210	11117	0.16%	0.020%
6	6.805	644	649	659	rVB	96803	163643	2.31%	0.301%
7	6.975	672	678	694	rVB	532094	832783	11.78%	1.532%
8	7.164	701	710	722	rBV	460064	735500	10.40%	1.353%
9	7.516	764	770	779	rBV	143234	232416	3.29%	0.428%
10	7.634	783	790	793	rBV	500443	875026	12.38%	1.610%
11	7.663	793	795	805	rVB	532825	781432	11.05%	1.438%
12	7.981	842	849	856	rBV	101706	166844	2.36%	0.307%
13	8.334	902	909	920	rBV	339935	553183	7.83%	1.018%
14	8.787	978	986	995	rBV	644187	1038027	14.68%	1.910%
15	8.952	1009	1014	1020	rVB5	7660	12966	0.18%	0.024%
16	9.181	1049	1053	1058	rVB	10330	16313	0.23%	0.030%
17	9.505	1101	1108	1124	rBV	472915	774497	10.96%	1.425%
18	9.910	1173	1177	1181	rBV4	4982	8913	0.13%	0.016%
19	10.034	1191	1198	1214	rBV	757815	1262049	17.85%	2.322%
20	10.275	1232	1239	1244	rBV	188648	309094	4.37%	0.569%
21	10.316	1244	1246	1252	rVV	16341	22630	0.32%	0.042%
22	10.410	1255	1262	1270	rBV	747428	1244468	17.60%	2.290%
23	10.557	1279	1287	1304	rBV	460834	801587	11.34%	1.475%
24	10.793	1322	1327	1332	rBV2	23604	35393	0.50%	0.065%
25	12.004	1530	1533	1539	rVB2	15567	22545	0.32%	0.041%
26	12.446	1604	1608	1615	rVB6	6170	11212	0.16%	0.021%
27	12.757	1656	1661	1668	rBV6	3901	8513	0.12%	0.016%
28	13.063	1708	1713	1717	rBV4	6096	8825	0.12%	0.016%
29	13.175	1724	1732	1738	rBV	126115	195920	2.77%	0.361%
30	13.693	1814	1820	1833	rBV	1633839	2237647	31.65%	4.118%
31	13.969	1859	1867	1880	rBV2	1893705	2764644	39.11%	5.087%
32	14.275	1911	1919	1930	rBV	1206450	1779514	25.17%	3.275%
33	14.498	1951	1957	1969	rBV3	55678	123680	1.75%	0.228%
34	14.598	1969	1974	1977	rVV7	4963	8468	0.12%	0.016%
35	14.693	1985	1990	1997	rBV8	4775	11012	0.16%	0.020%
36	14.845	2010	2016	2027	rVB3	11267	30162	0.43%	0.056%

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 ClientSampleId :
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Integration Parameters: LSCINT.P

Integrator: RTE

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Filtering: 5

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Stop Thrs : 0

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Peak separation: 5

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37	15.275	2082	2089	2101	rBV	2538548	3557341	50.32%	6.546%
38	15.398	2104	2110	2121	rBV	298506	412514	5.84%	0.759%
39	15.516	2126	2130	2139	rVB3	11911	20461	0.29%	0.038%
40	16.063	2219	2223	2231	rVB6	8751	14622	0.21%	0.027%
41	16.434	2282	2286	2293	rBV5	21461	33189	0.47%	0.061%
42	17.028	2381	2387	2394	rBV	1682728	2311719	32.70%	4.254%
43	17.128	2398	2404	2415	rBV	2982134	4109938	58.14%	7.563%
44	17.704	2498	2502	2508	rVB7	12213	16657	0.24%	0.031%
45	17.939	2535	2542	2567	rBV	929227	1576325	22.30%	2.901%
46	18.604	2652	2655	2659	rBV5	9548	13900	0.20%	0.026%
47	18.945	2710	2713	2717	rBV6	13374	19374	0.27%	0.036%
48	18.992	2717	2721	2725	rBV2	41081	56012	0.79%	0.103%
49	19.063	2729	2733	2736	rBV	39344	51581	0.73%	0.095%
50	19.098	2736	2739	2741	rBV3	21057	25707	0.36%	0.047%
51	19.139	2741	2746	2756	rVV	322767	518378	7.33%	0.954%
52	19.222	2756	2760	2775	rVB	429759	631193	8.93%	1.162%
53	19.428	2789	2795	2803	rBV	3758410	5297658	74.94%	9.749%
54	20.216	2927	2929	2936	rBV7	17794	21802	0.31%	0.040%
55	20.322	2943	2947	2952	rBV	108668	130940	1.85%	0.241%
56	20.957	3051	3055	3066	rBV4	99630	203412	2.88%	0.374%
57	21.263	3101	3107	3117	rBV2	2308822	3204591	45.33%	5.897%
58	23.968	3557	3567	3582	rBV2	2931924	7069422	100.00%	13.009%
59	24.163	3591	3600	3613	rBV	1667336	3926326	55.54%	7.225%
60	25.710	3856	3863	3886	rVB5	207218	774884	10.96%	1.426%
61	26.168	3934	3941	3953	rVB5	436585	1320024	18.67%	2.429%

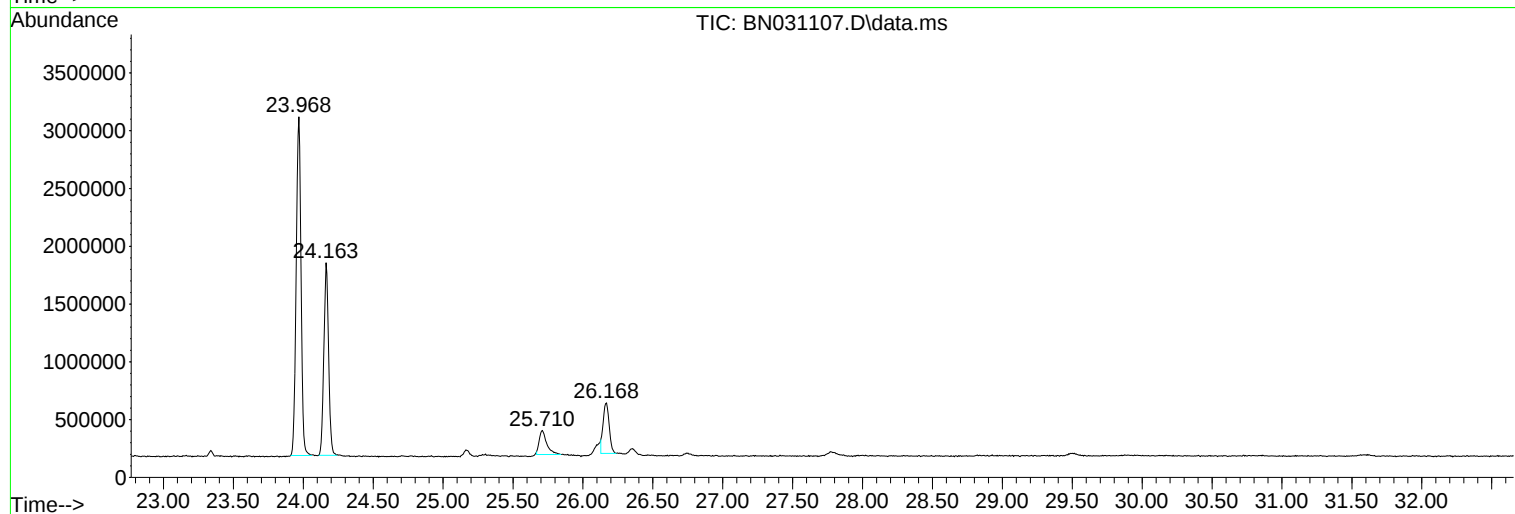
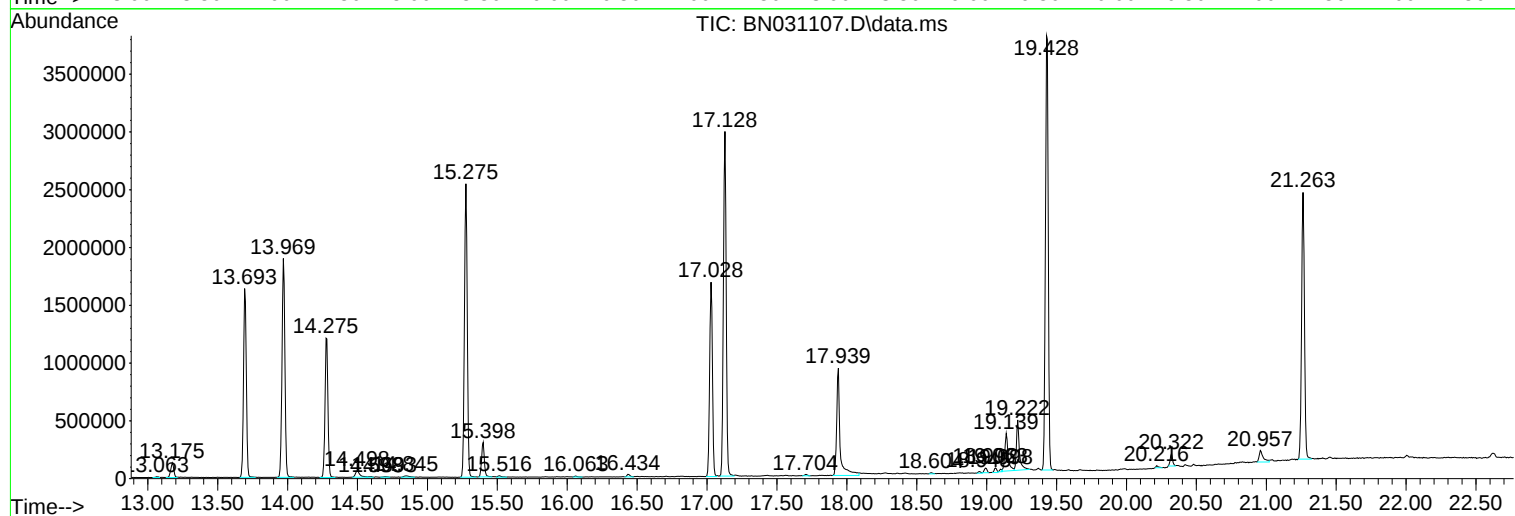
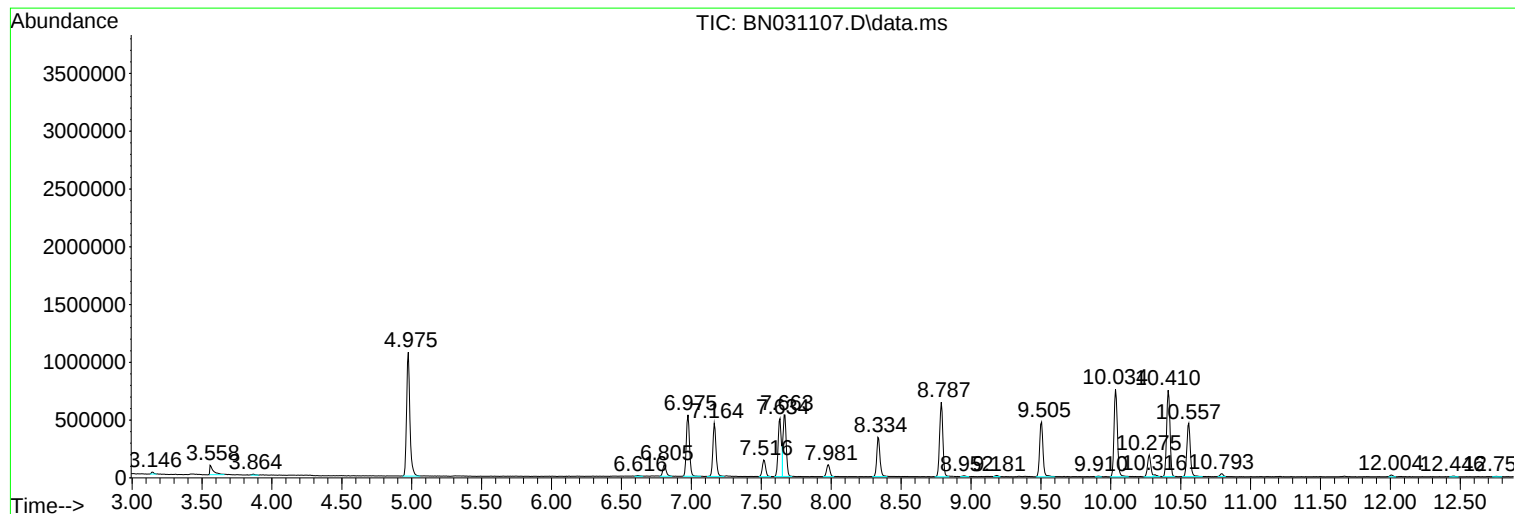
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ClientSampleId :
C0RD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
BNA_N
ClientSampleId :
CORD4

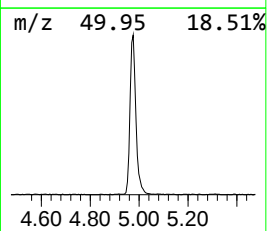
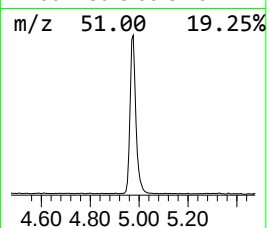
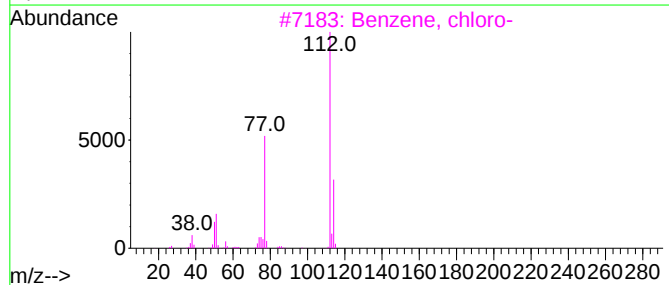
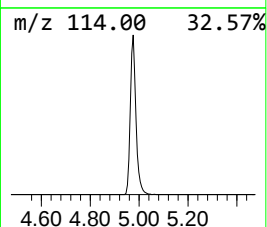
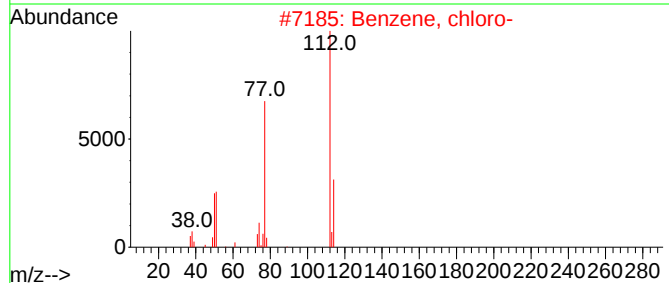
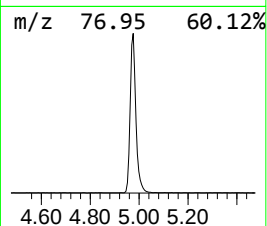
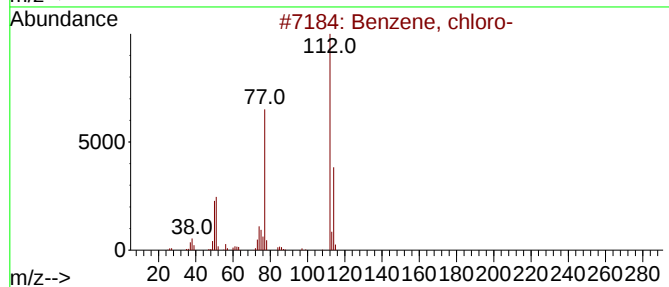
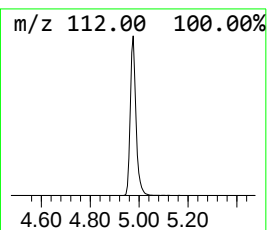
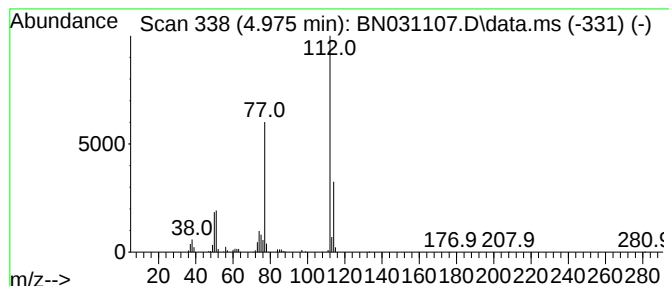
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.975	40.36 ng/ul	1765710	1,4-Dichlorobenzene-d4	7.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	96
2			Benzene, chloro-	112	C6H5Cl	000108-90-7	96
3			Benzene, chloro-	112	C6H5Cl	000108-90-7	95
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	94



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ALS Vial : 58 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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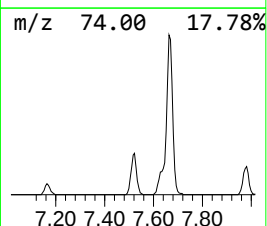
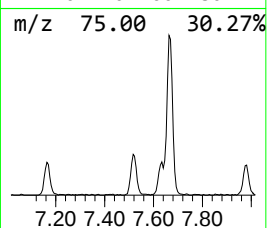
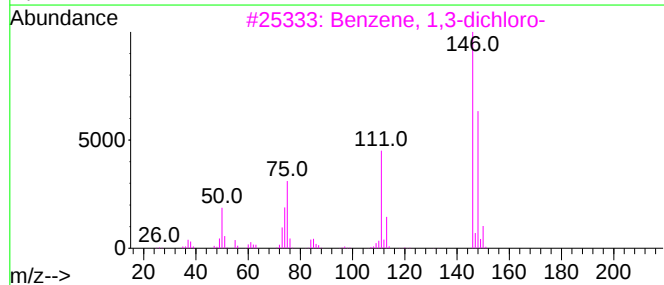
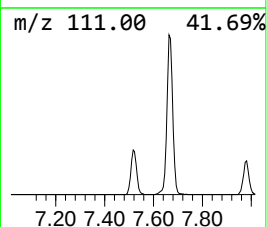
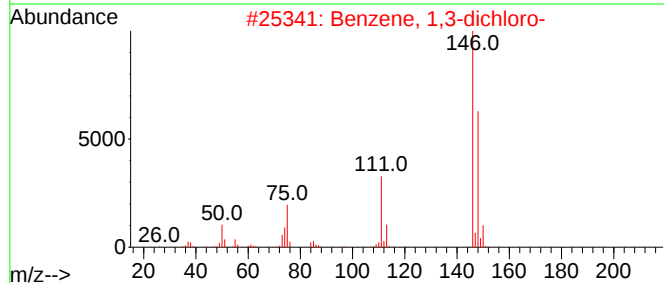
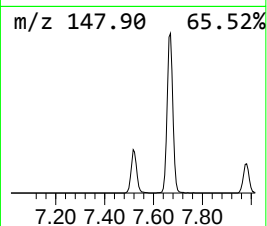
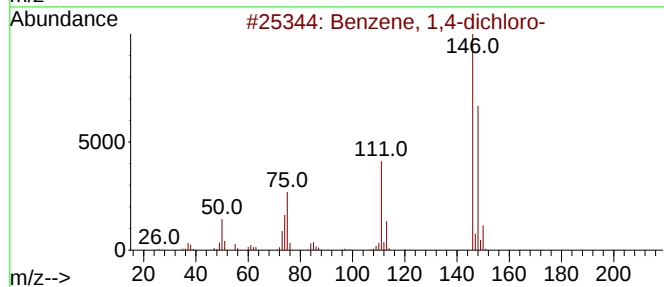
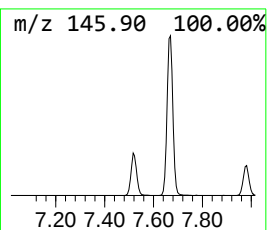
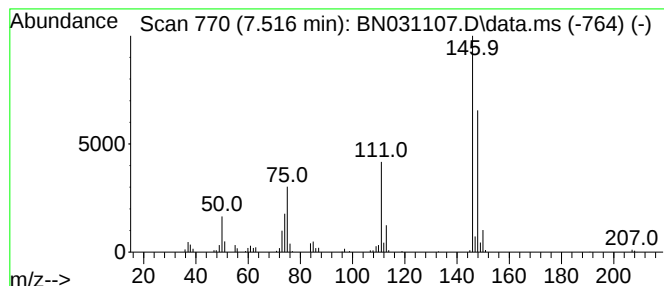
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,4-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.516	5.31 ng/ul	232416	1,4-Dichlorobenzene-d4	7.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
2			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96



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ALS Vial : 58 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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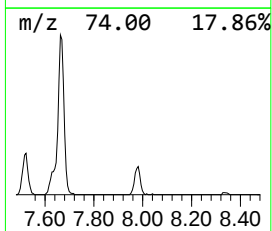
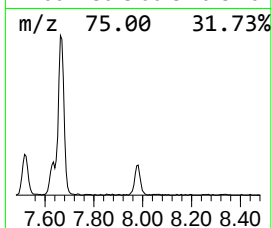
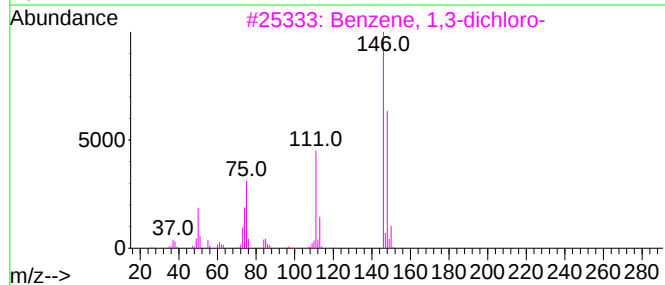
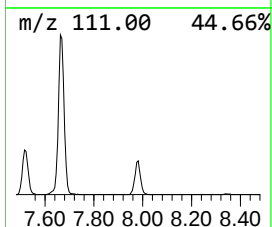
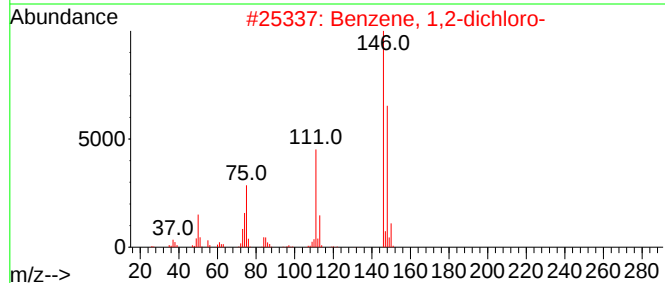
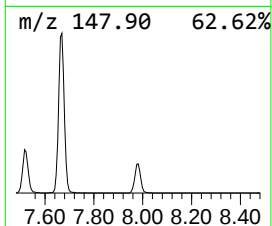
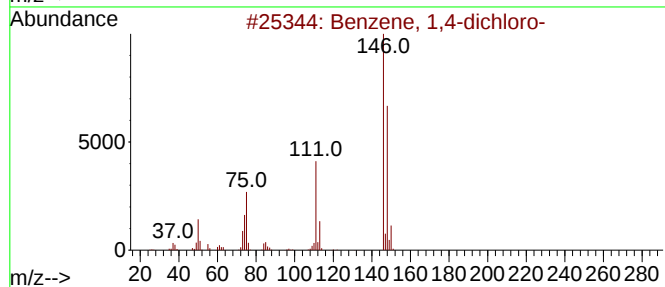
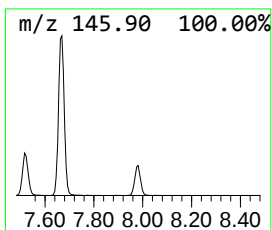
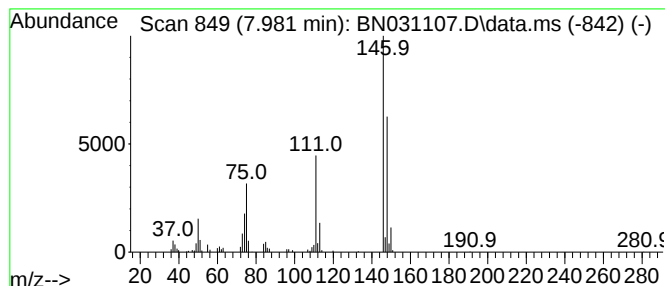
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,2-dichloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.981	3.81 ng/ul	166844	1,4-Dichlorobenzene-d4	7.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
2			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96



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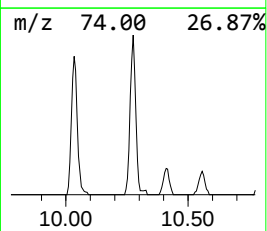
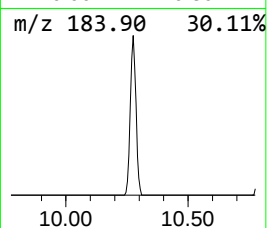
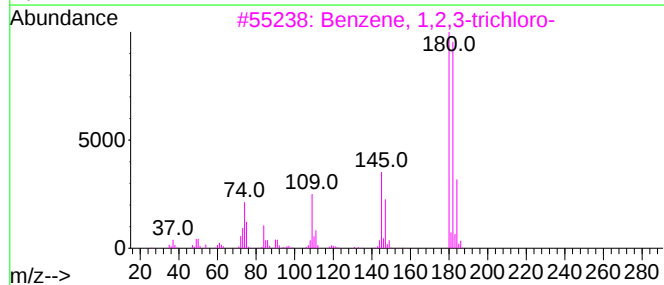
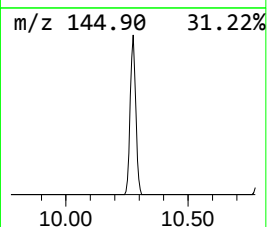
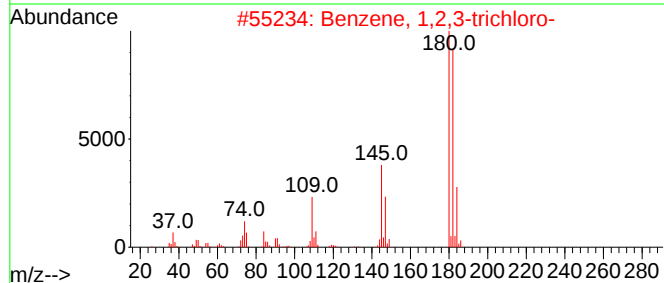
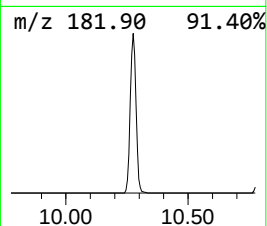
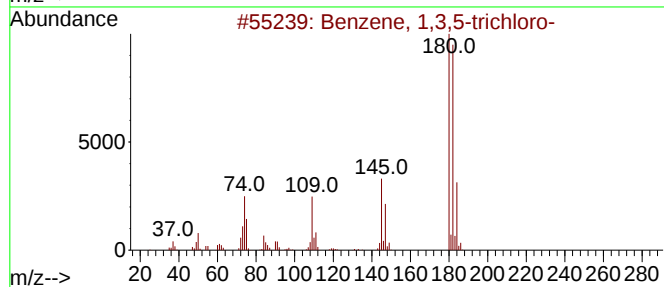
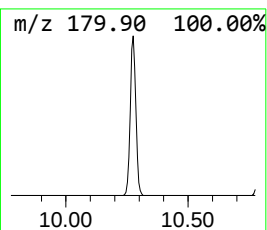
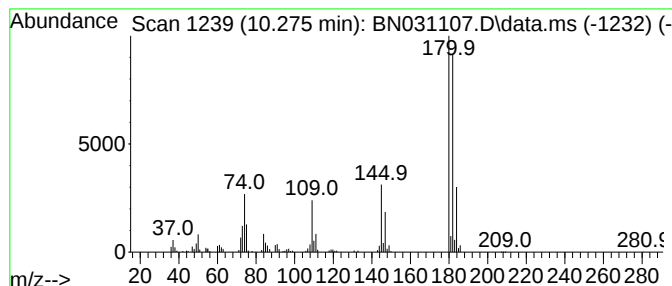
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1,3,5-trichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.275	4.97 ng/ul	309094	Naphthalene-d8	10.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
2			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
3			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
4			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	97
5			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	96



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ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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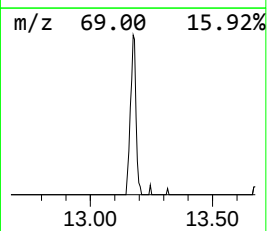
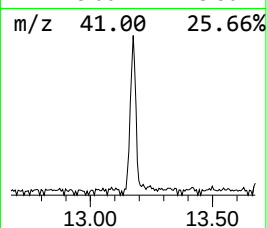
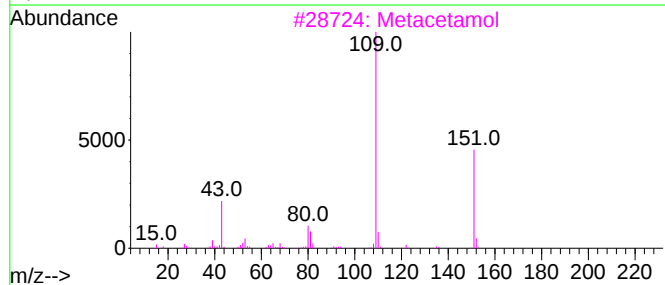
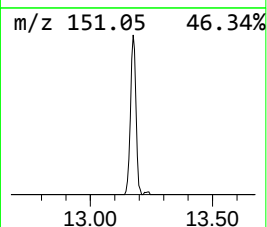
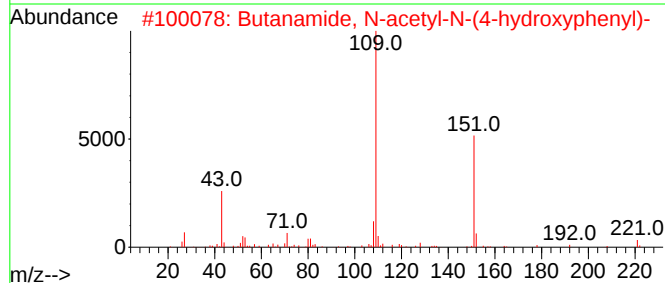
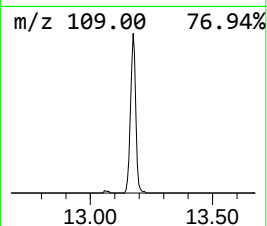
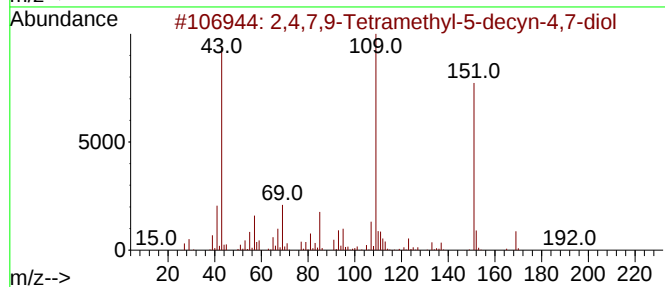
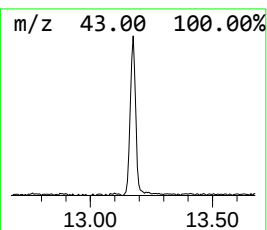
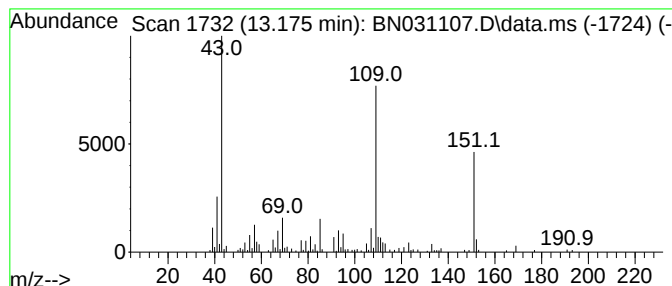
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.175	2.20 ng/ul	195920	Acenaphthene-d10	14.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91		
2	Butanamide, N-acetyl-N-(4-hydrox...	221	C12H15NO3	1000107-08-7	59		
3	Metacetamol	151	C8H9NO2	000621-42-1	58		
4	Diacetamate	193	C10H11NO3	002623-33-8	53		
5	Metacetamol	151	C8H9NO2	000621-42-1	53		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
Data File : BN031107.D
Acq On : 21 Apr 2024 13:49
Operator : MA/JU
Sample : P2161-03
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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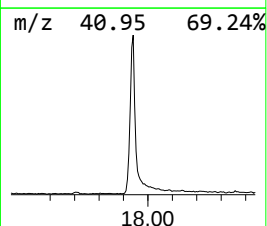
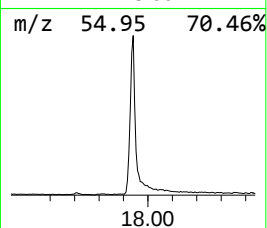
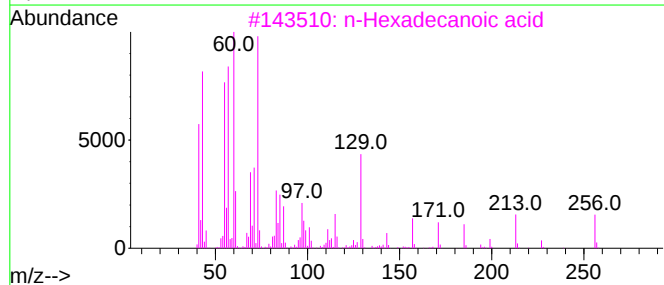
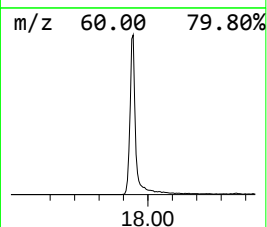
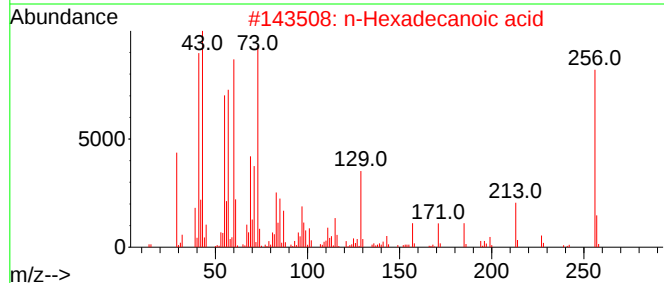
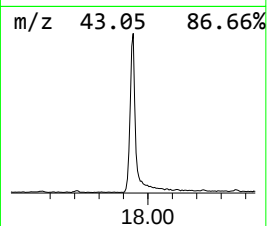
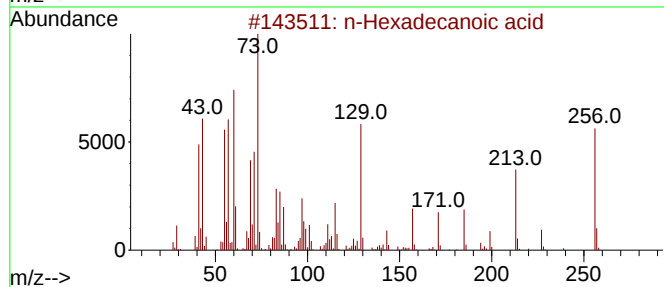
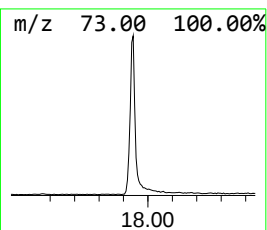
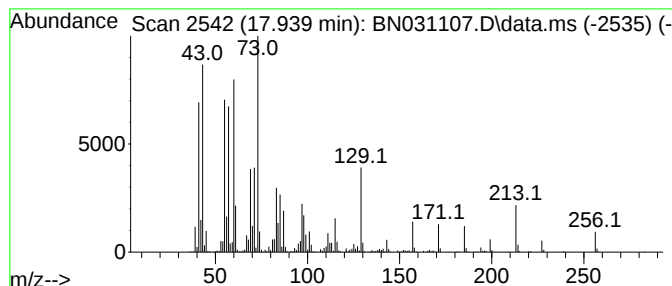
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.939	13.64 ng/ul	1576330	Phenanthrene-d10	17.028

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	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Tridecanoic acid	214	C13H26O2	000638-53-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
Data File : BN031107.D
Acq On : 21 Apr 2024 13:49
Operator : MA/JU
Sample : P2161-03
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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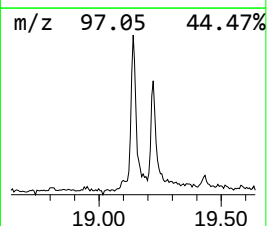
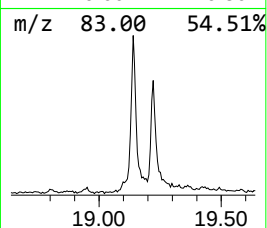
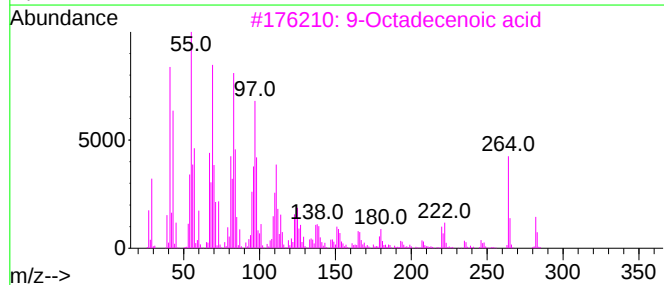
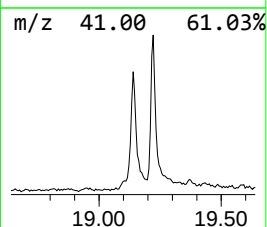
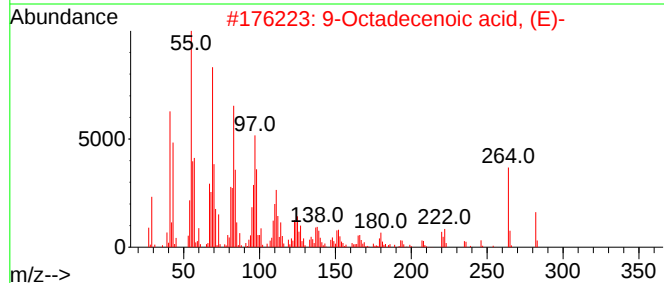
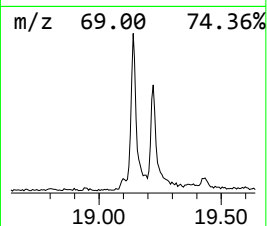
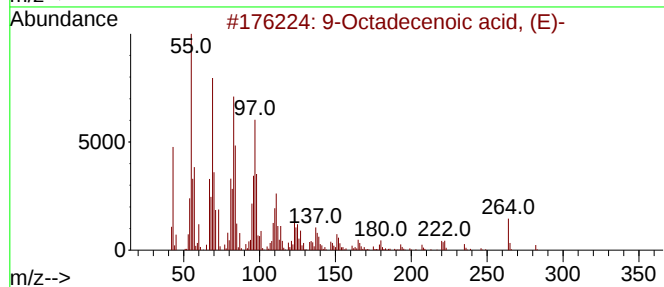
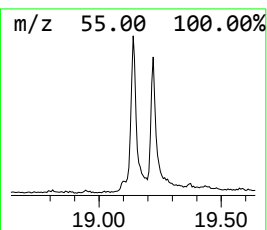
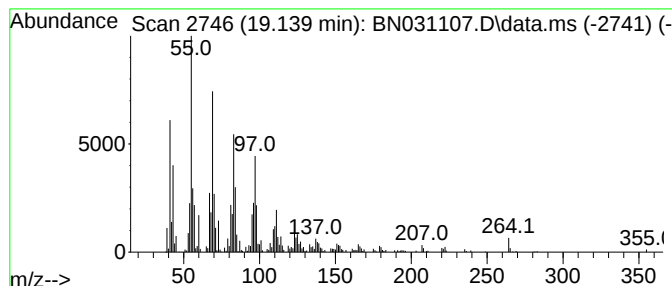
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9-Octadecenoic acid, (E)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.139	4.48 ng/ul	518378	Phenanthrene-d10	17.028

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	94
2			9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	91
3			9-Octadecenoic acid	282	C18H34O2	002027-47-6	91
4			cis-Vaccenic acid	282	C18H34O2	000506-17-2	91
5			n-Propyl 11-octadecenoate	324	C21H40O2	1000336-71-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
Data File : BN031107.D
Acq On : 21 Apr 2024 13:49
Operator : MA/JU
Sample : P2161-03
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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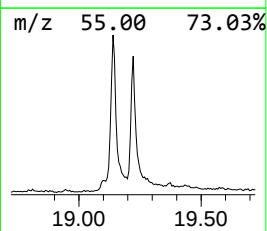
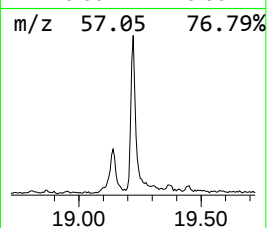
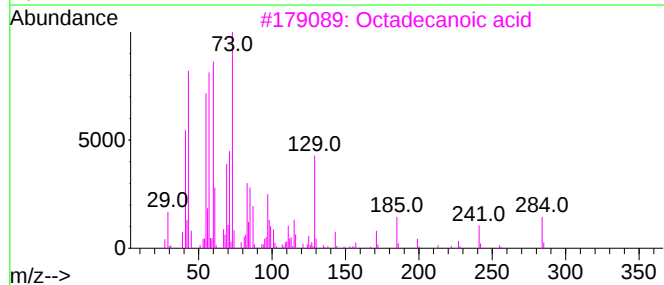
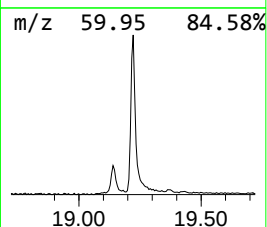
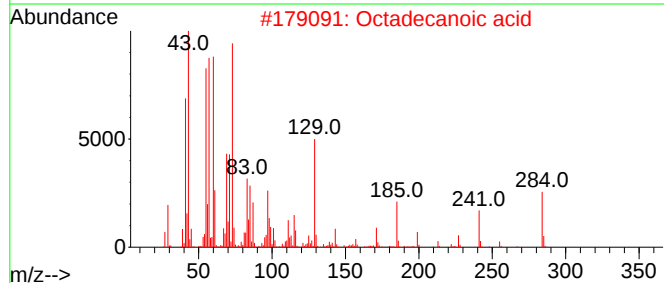
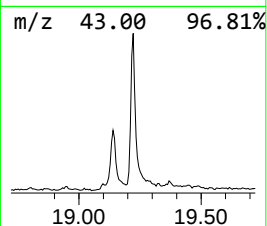
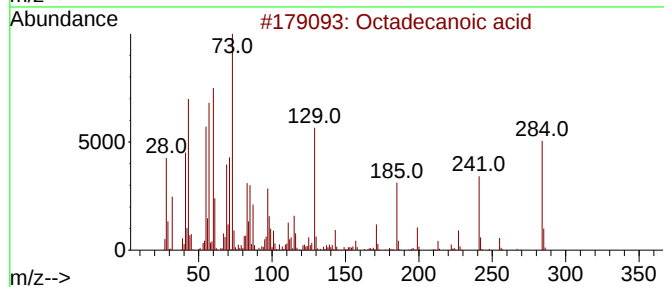
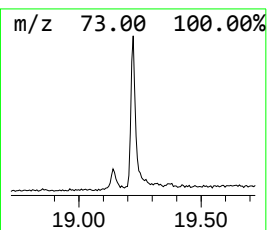
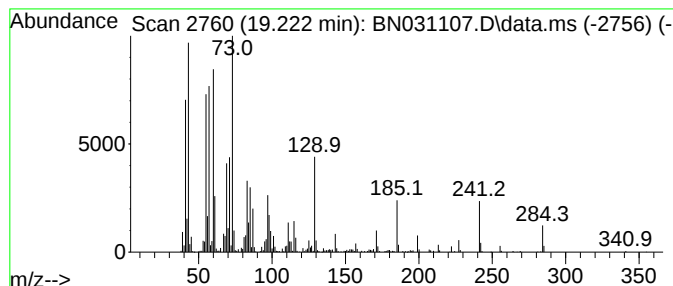
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.222	3.94 ng/ul	631193	Chrysene-d12	21.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	98
5			Octadecanoic acid	284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
Data File : BN031107.D
Acq On : 21 Apr 2024 13:49
Operator : MA/JU
Sample : P2161-03
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CORD4

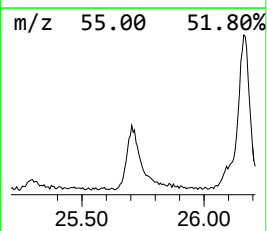
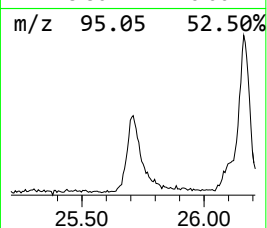
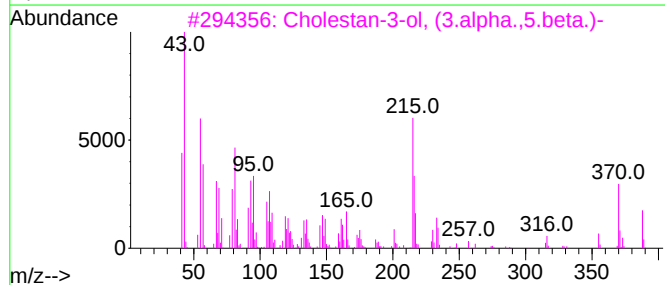
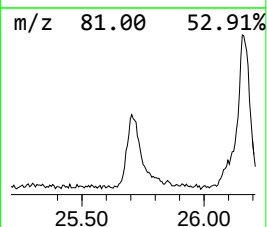
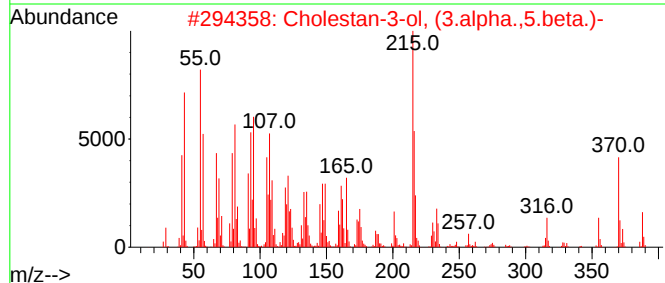
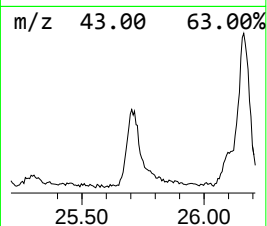
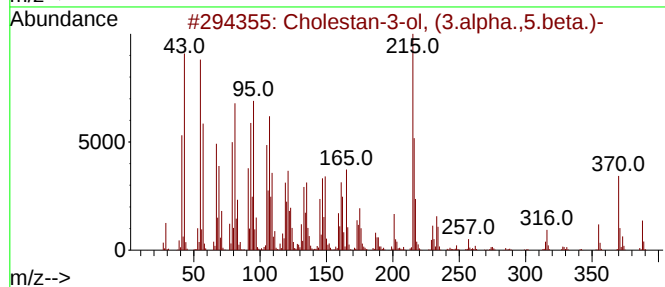
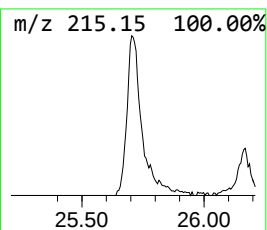
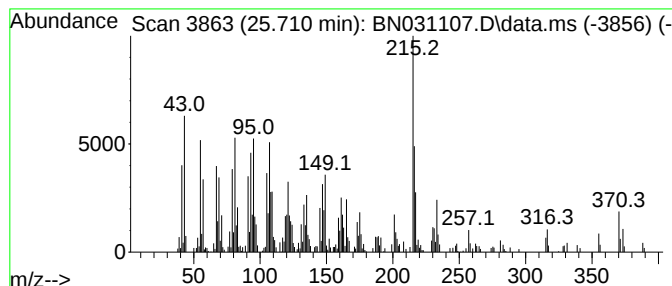
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Cholestan-3-ol, (3.alpha.,5... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.710	3.95 ng/ul	774884	Perylene-d12	24.163

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cholestan-3-ol, (3.alpha.,5.beta.)-	388	C27H48O	000516-92-7	95
2			Cholestan-3-ol, (3.alpha.,5.beta.)-	388	C27H48O	000516-92-7	95
3			Cholestan-3-ol, (3.alpha.,5.beta.)-	388	C27H48O	000516-92-7	93
4			Cholestan-3-ol, (3.alpha.,5.beta.)-	388	C27H48O	000516-92-7	93
5			5.beta.-Cholestan-3.alpha.-ol, 2...	458	C31H54O2	1000514-95-7	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
Data File : BN031107.D
Acq On : 21 Apr 2024 13:49
Operator : MA/JU
Sample : P2161-03
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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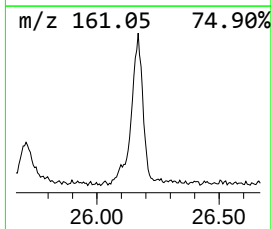
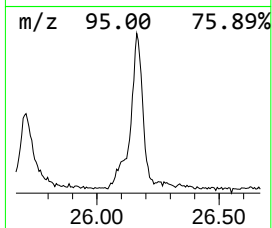
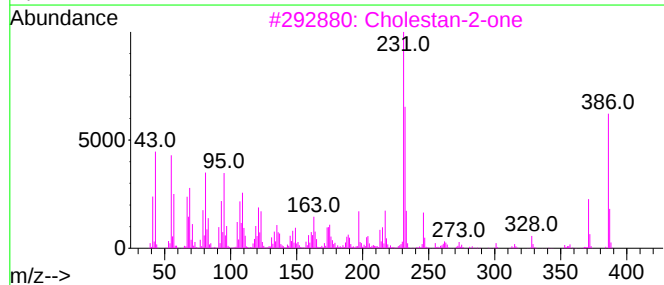
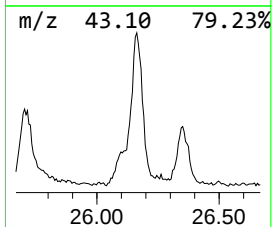
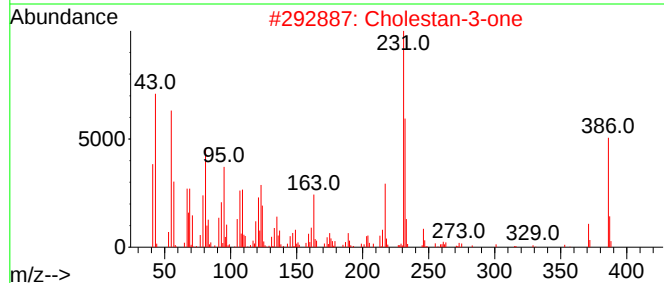
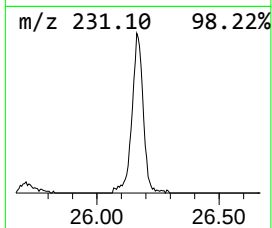
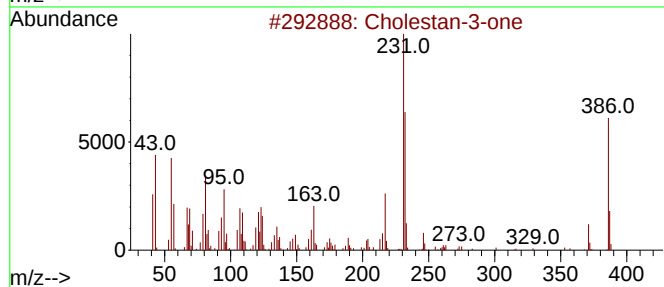
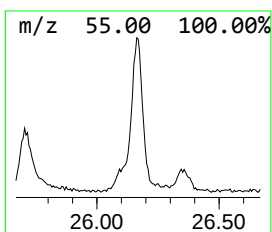
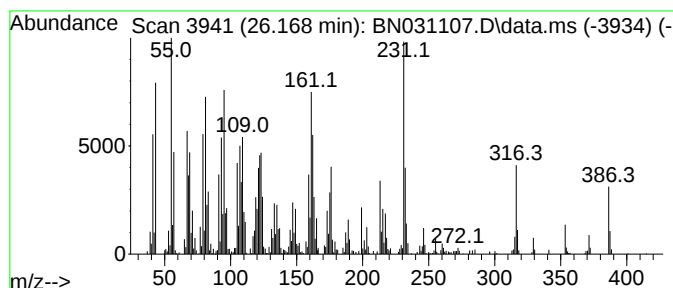
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Cholestan-3-one Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.168	6.72 ng/ul	1320020	Perylene-d12	24.163

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cholestan-3-one	386	C27H46O	015600-08-5	96
2			Cholestan-3-one	386	C27H46O	015600-08-5	72
3			Cholestan-2-one	386	C27H46O	1010210-43-4	70
4			Cholestane, 4,5-epoxy-, (4.alpha...	386	C27H46O	006079-19-2	55
5			Cholestan-3-one	386	C27H46O	015600-08-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
Data File : BN031107.D
Acq On : 21 Apr 2024 13:49
Operator : MA/JU
Sample : P2161-03
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
CORD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.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
Benzene, chloro-	4.975	40.4	ng/ul	1765710	1	7.634	875026	20.0
Benzene, 1,4-di...	7.516	5.3	ng/ul	232416	1	7.634	875026	20.0
Benzene, 1,2-di...	7.981	3.8	ng/ul	166844	1	7.634	875026	20.0
Benzene, 1,3,5-...	10.275	5.0	ng/ul	309094	2	10.410	1244470	20.0
2,4,7,9-Tetrame...	13.175	2.2	ng/ul	195920	3	14.275	1779510	20.0
n-Hexadecanoic ...	17.939	13.6	ng/ul	1576330	4	17.028	2311720	20.0
9-Octadecenoic ...	19.139	4.5	ng/ul	518378	4	17.028	2311720	20.0
Octadecanoic acid	19.222	3.9	ng/ul	631193	5	21.263	3204590	20.0
Cholestan-3-ol,...	25.710	4.0	ng/ul	774884	6	24.163	3926330	20.0
Cholestan-3-one	26.168	6.7	ng/ul	1320020	6	24.163	3926330	20.0