

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
 Data File : BN022744.D
 Acq On : 18 Nov 2022 06:31
 Operator : CG/JU
 Sample : N5574-01
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 B0AA0

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-BN111522.M
 Title : SVOA CALIBRATION

Signal : TIC: BN022744.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.064	9	13	28	rVB	288217	470106	21.09%	1.979%
2	3.181	30	33	46	rVB	16784	27731	1.24%	0.117%
3	3.281	46	50	56	rBV2	6751	10746	0.48%	0.045%
4	3.475	80	83	87	rVB3	3903	5455	0.24%	0.023%
5	3.622	106	108	118	rBV	58210	104258	4.68%	0.439%
6	3.852	141	147	153	rVB	7879	13701	0.61%	0.058%
7	3.946	159	163	171	rVB	11528	18506	0.83%	0.078%
8	4.481	243	254	269	rBV	925308	1512283	67.83%	6.365%
9	4.669	284	286	293	rBV4	1901	2717	0.12%	0.011%
10	4.740	293	298	299	rVV	6077	7875	0.35%	0.033%
11	4.775	299	304	314	rVV	100682	165444	7.42%	0.696%
12	4.875	316	321	323	rVV4	1466	2457	0.11%	0.010%
13	4.922	324	329	337	rVV	55214	84659	3.80%	0.356%
14	4.999	337	342	352	rVB	43506	68639	3.08%	0.289%
15	6.087	523	527	532	rBV3	5835	9153	0.41%	0.039%
16	7.022	681	686	695	rVB	42642	69236	3.11%	0.291%
17	7.187	707	714	722	rBV	299060	474915	21.30%	1.999%
18	7.258	722	726	734	rVB2	9770	17107	0.77%	0.072%
19	7.381	739	747	756	rBV	344958	567310	25.45%	2.388%
20	7.558	772	777	784	rVB2	5595	9258	0.42%	0.039%
21	7.852	820	827	834	rBV	233475	371909	16.68%	1.565%
22	7.916	834	838	846	rVB	11129	19231	0.86%	0.081%
23	8.581	945	951	965	rBV	103245	179578	8.06%	0.756%
24	8.728	971	976	982	rBV3	9175	14960	0.67%	0.063%
25	9.034	1017	1028	1037	rBV	371228	650015	29.16%	2.736%
26	9.134	1039	1045	1058	rVB	188368	311070	13.95%	1.309%
27	9.405	1088	1091	1099	rVB4	2210	4304	0.19%	0.018%
28	9.757	1143	1151	1163	rBV	306189	510516	22.90%	2.149%
29	9.934	1176	1181	1186	rBV3	2549	6111	0.27%	0.026%
30	10.299	1236	1243	1254	rBV	389151	682156	30.60%	2.871%
31	10.452	1263	1269	1280	rVB2	45339	83757	3.76%	0.353%
32	10.569	1282	1289	1299	rBV2	54648	113574	5.09%	0.478%
33	10.669	1299	1306	1317	rVB	366749	607519	27.25%	2.557%
34	10.834	1330	1334	1341	rVB3	2162	3734	0.17%	0.016%
35	11.481	1439	1444	1454	rBV3	10340	21885	0.98%	0.092%
36	11.940	1518	1522	1530	rVB2	10312	18736	0.84%	0.079%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
 Title : SVOA CALIBRATION

37	12.146	1553	1557	1563	rVB2	3258	5160	0.23%	0.022%
38	12.269	1573	1578	1582	rBV2	5054	7418	0.33%	0.031%
39	12.869	1676	1680	1686	rVB5	1371	3044	0.14%	0.013%
40	12.946	1688	1693	1697	rVV	4621	6308	0.28%	0.027%
41	13.716	1817	1824	1832	rBV3	4856	11642	0.52%	0.049%
42	13.940	1849	1862	1874	rVB	997279	1384824	62.12%	5.829%
43	14.216	1901	1909	1922	rBV	506540	716606	32.14%	3.016%
44	14.522	1955	1961	1972	rVB	598343	876591	39.32%	3.690%
45	14.751	1995	2000	2018	rBV	79118	150097	6.73%	0.632%
46	15.116	2056	2062	2063	rBV3	1235	2270	0.10%	0.010%
47	15.169	2063	2071	2079	rVB2	8024	18032	0.81%	0.076%
48	15.316	2089	2096	2099	rVB5	1538	2808	0.13%	0.012%
49	15.522	2124	2131	2140	rBV	1321400	1883479	84.48%	7.928%
50	15.651	2147	2153	2166	rBV	525682	715748	32.11%	3.013%
51	15.763	2166	2172	2176	rBV2	2629	3898	0.17%	0.016%
52	16.216	2242	2249	2260	rBV2	70596	110689	4.97%	0.466%
53	16.934	2362	2371	2377	rBV	59281	81315	3.65%	0.342%
54	17.151	2404	2408	2413	rVB2	5182	7881	0.35%	0.033%
55	17.281	2423	2430	2436	rBV	758724	1062434	47.66%	4.472%
56	17.381	2440	2447	2457	rVV	1033720	1492294	66.94%	6.281%
57	17.469	2457	2462	2466	rVV	53289	80517	3.61%	0.339%
58	17.522	2466	2471	2491	rVV	237006	439131	19.70%	1.848%
59	17.681	2493	2498	2503	rVB2	8728	13132	0.59%	0.055%
60	17.945	2539	2543	2547	rBV2	10528	13278	0.60%	0.056%
61	18.104	2565	2570	2576	rBV	50884	72312	3.24%	0.304%
62	18.169	2576	2581	2588	rVV	26454	37042	1.66%	0.156%
63	18.239	2588	2593	2608	rVB	30166	64746	2.90%	0.273%
64	18.522	2637	2641	2647	rBV6	6935	10823	0.49%	0.046%
65	18.681	2663	2668	2674	rBV2	25815	41426	1.86%	0.174%
66	18.945	2709	2713	2717	rBV5	1801	3539	0.16%	0.015%
67	19.186	2751	2754	2755	rBV2	2254	2479	0.11%	0.010%
68	19.245	2760	2764	2771	rVB3	16555	28212	1.27%	0.119%
69	19.316	2772	2776	2782	rBV2	26305	38715	1.74%	0.163%
70	19.457	2796	2800	2805	rBV3	20142	26504	1.19%	0.112%
71	19.604	2823	2825	2827	rBV3	4207	3990	0.18%	0.017%
72	19.680	2832	2838	2847	rBV2	1635870	2200376	98.70%	9.261%
73	20.869	3037	3040	3041	rBV3	12018	12794	0.57%	0.054%
74	20.904	3041	3046	3053	rVB	139837	193405	8.68%	0.814%
75	21.480	3139	3144	3156	rBV	965124	1232190	55.27%	5.186%
76	23.727	3518	3526	3540	rVB2	1092037	2229372	100.00%	9.383%

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ClientSampleId :
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Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
Title : SVOA CALIBRATION

77	23.880	3545	3552	3565	rvB2	612905	1231571	55.24%	5.184%
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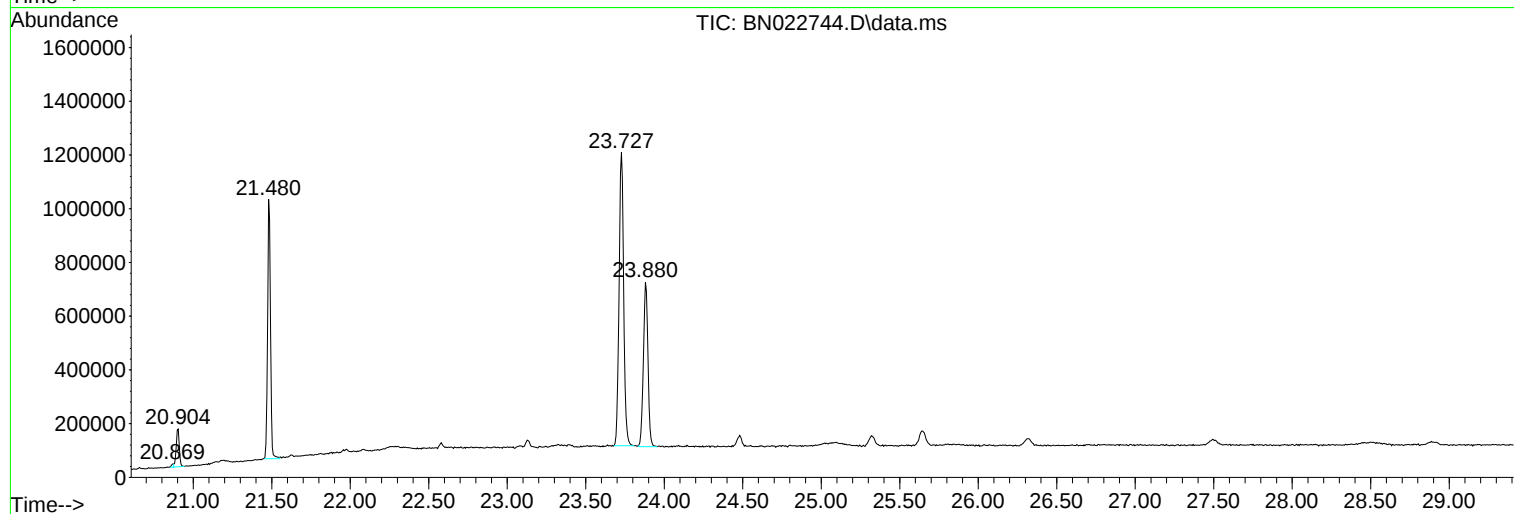
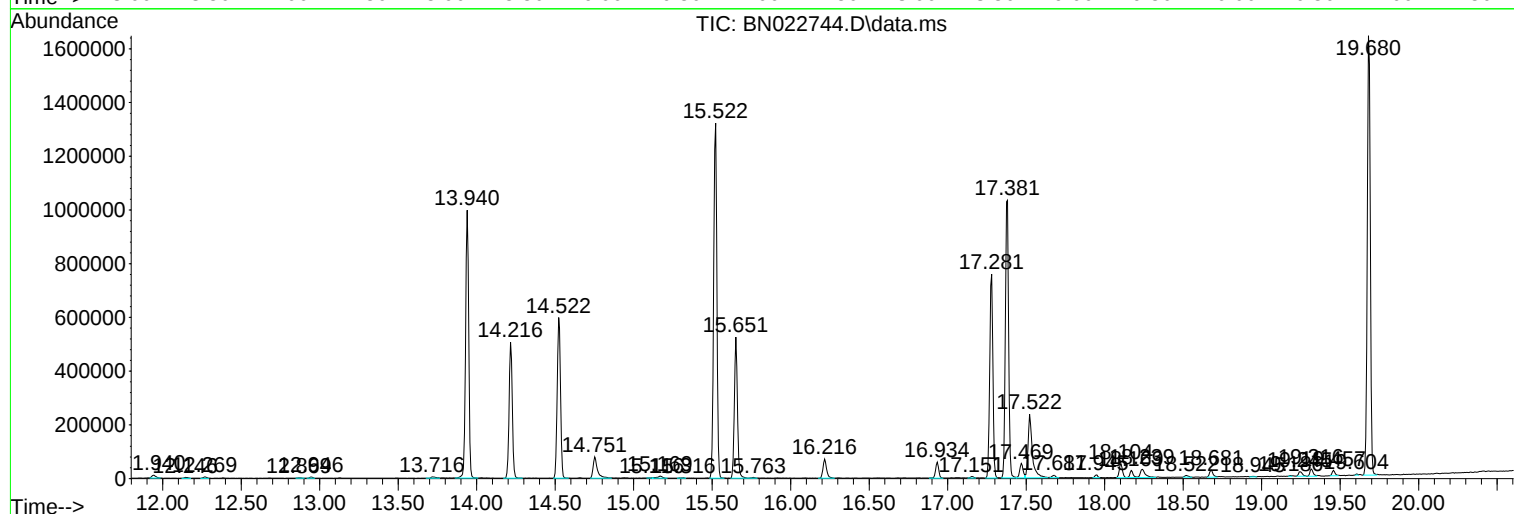
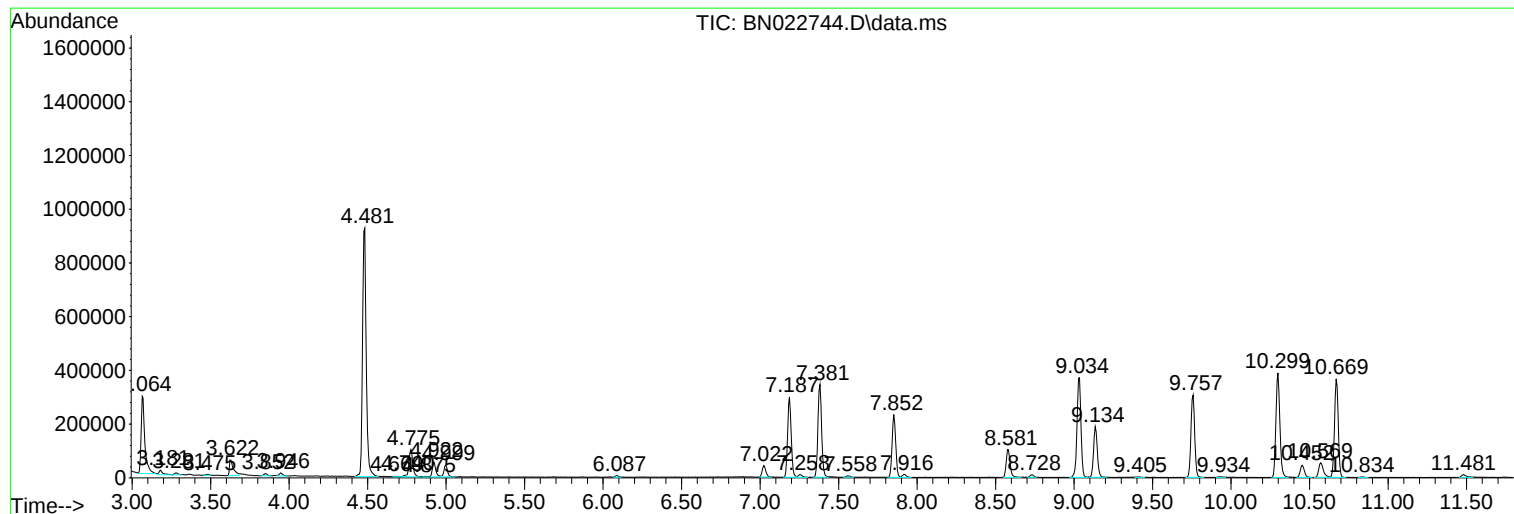
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
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Operator : CG/JU
Sample : N5574-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
B0AA0

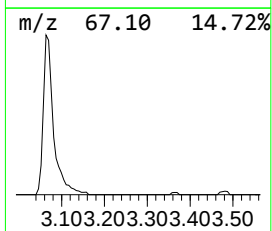
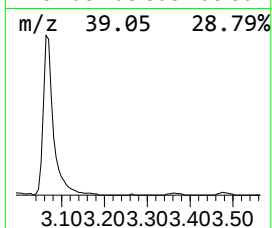
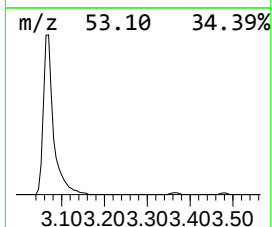
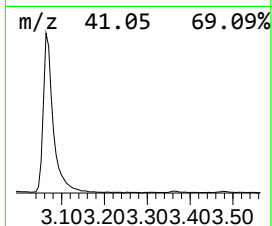
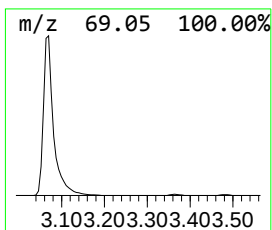
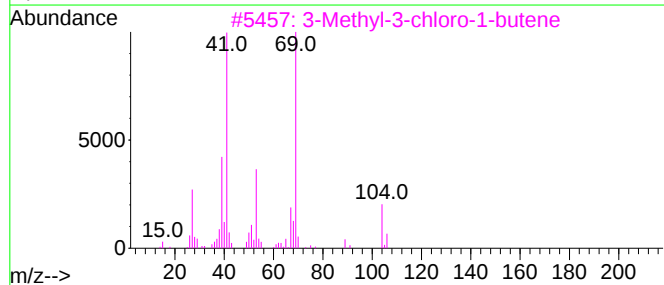
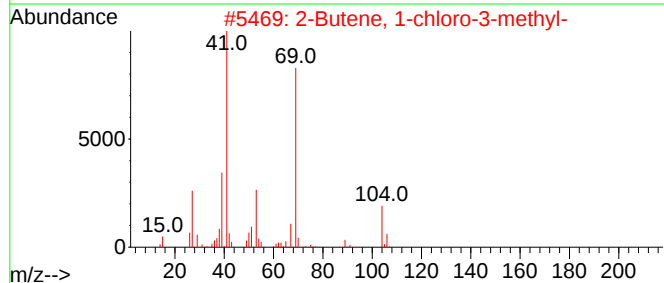
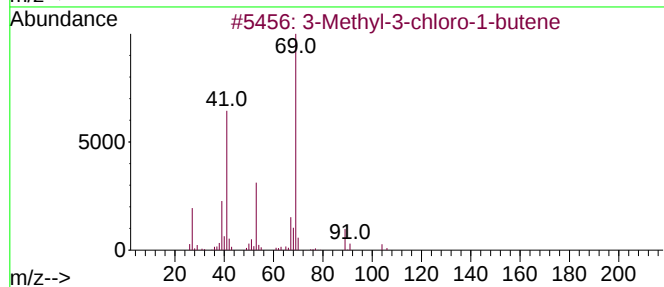
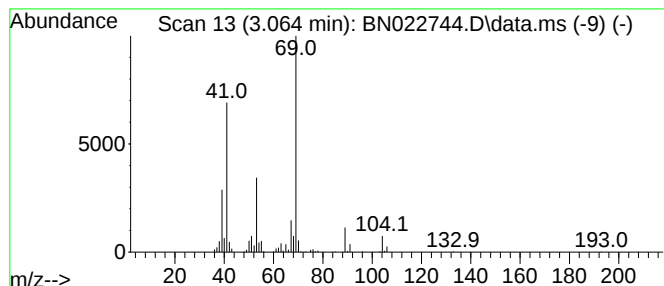
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 3-Methyl-3-chloro-1-butene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.064	25.28 ng/ul	470106	1,4-Dichlorobenzene-d4	7.852

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Methyl-3-chloro-1-butene	104	C5H9Cl	002190-48-9	78
2			2-Butene, 1-chloro-3-methyl-	104	C5H9Cl	000503-60-6	72
3			3-Methyl-3-chloro-1-butene	104	C5H9Cl	002190-48-9	72
4			2-Butene, 1-chloro-3-methyl-	104	C5H9Cl	000503-60-6	64
5			2-Butene, 1-chloro-3-methyl-	104	C5H9Cl	000503-60-6	64



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ALS Vial : 32 Sample Multiplier: 1

Instrument :
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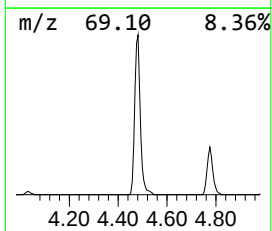
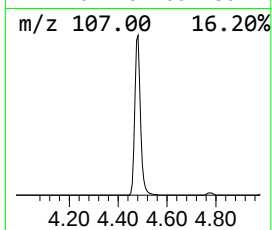
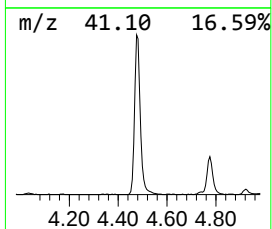
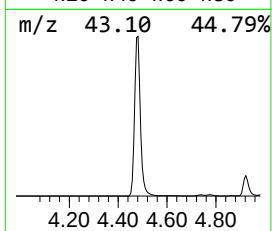
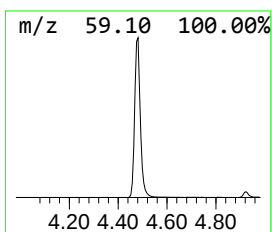
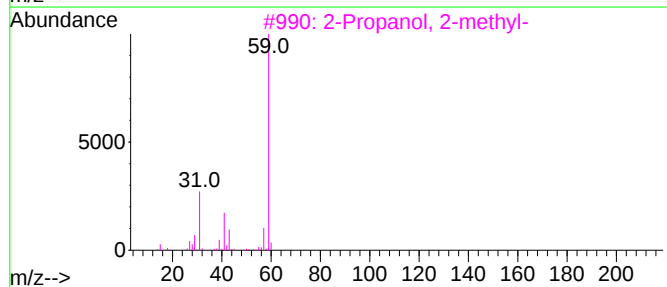
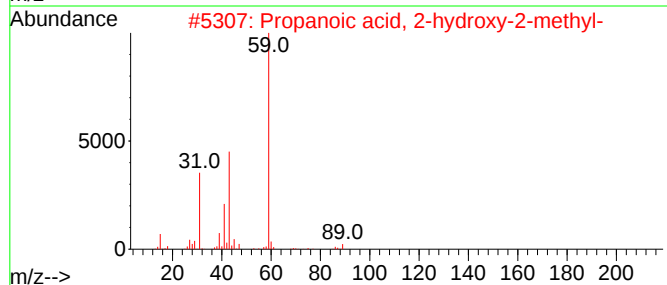
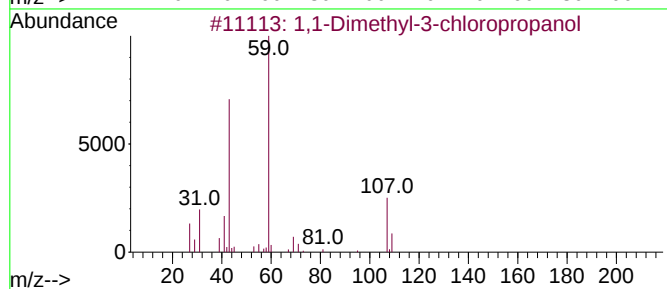
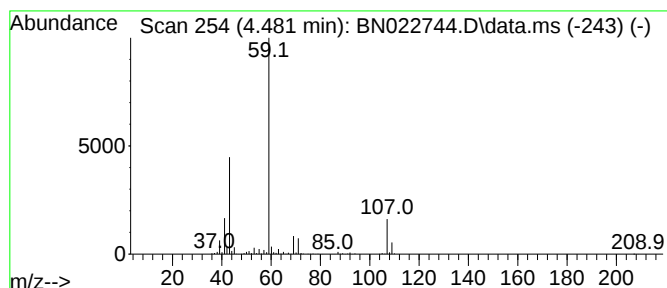
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1,1-Dimethyl-3-chloropropanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.481	81.33 ng/ul	1512280	1,4-Dichlorobenzene-d4	7.852

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1-Dimethyl-3-chloropropanol	122	C5H11ClO	001985-88-2	74
2			Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	59
3			2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	40
4			4-Methyl-2-pentanol, methyl ether	116	C7H16O	1000333-93-5	38
5			Propanoic acid, 2-hydroxy-2-meth...	118	C5H10O3	002110-78-3	38



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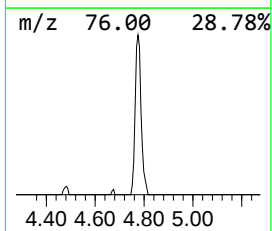
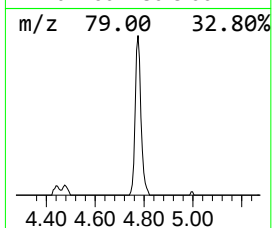
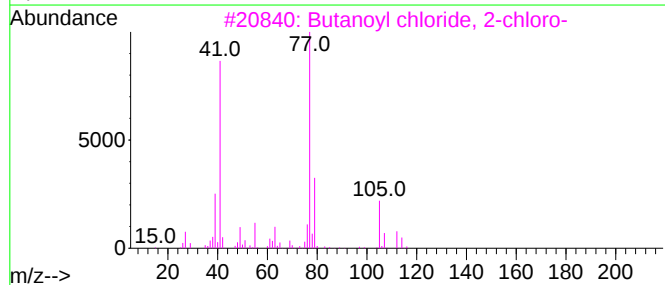
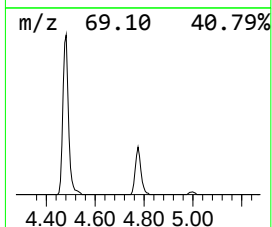
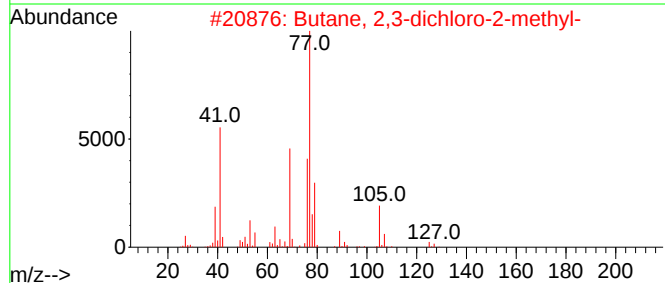
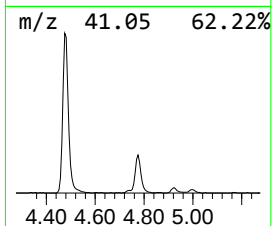
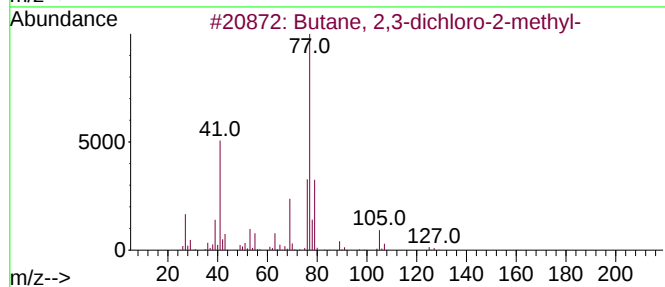
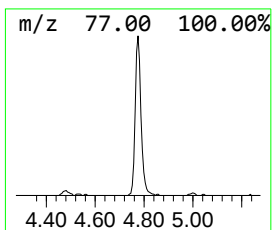
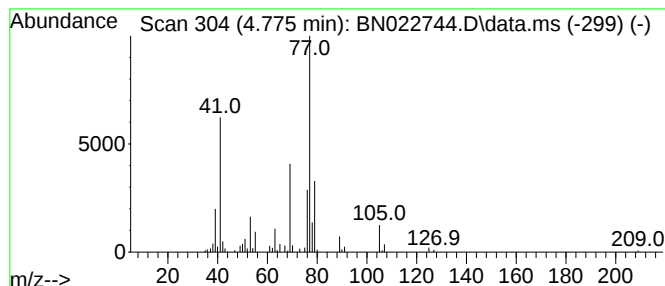
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Butane, 2,3-dichloro-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.775	8.90 ng/ul	165444	1,4-Dichlorobenzene-d4	7.852

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dichloro-2-methyl-	140	C5H10Cl2	000507-45-9	90
2			Butane, 2,3-dichloro-2-methyl-	140	C5H10Cl2	000507-45-9	50
3			Butanoyl chloride, 2-chloro-	140	C4H6Cl2O	007623-11-2	45
4			Propane, 1,1,2-trichloro-2-methyl-	160	C4H7Cl3	029559-52-2	40
5			Propane, 2-bromo-1-chloro-	156	C3H6BrCl	003017-95-6	38



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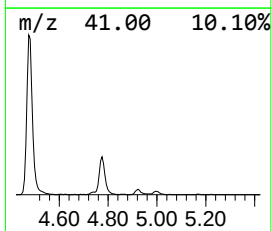
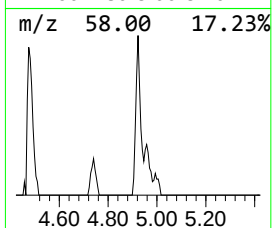
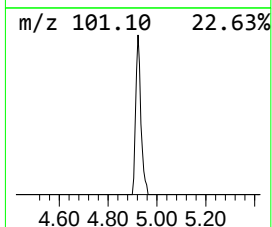
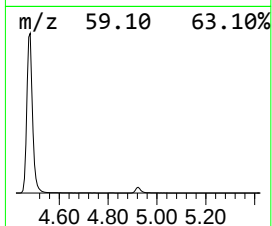
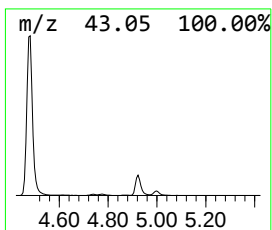
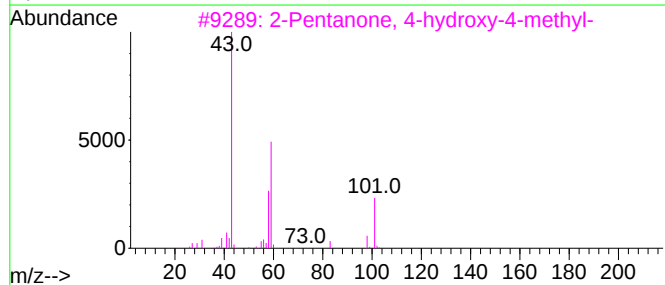
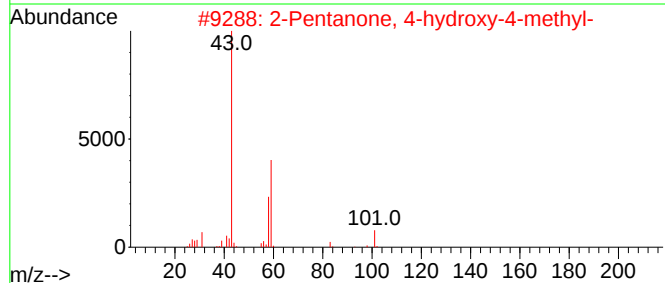
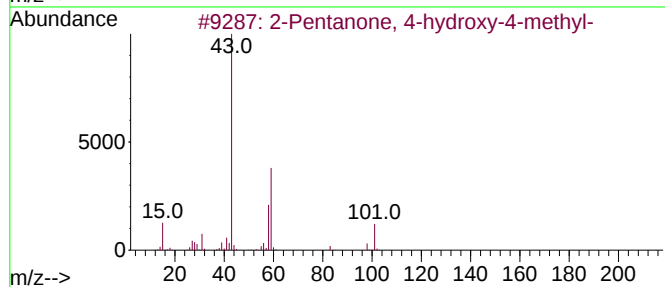
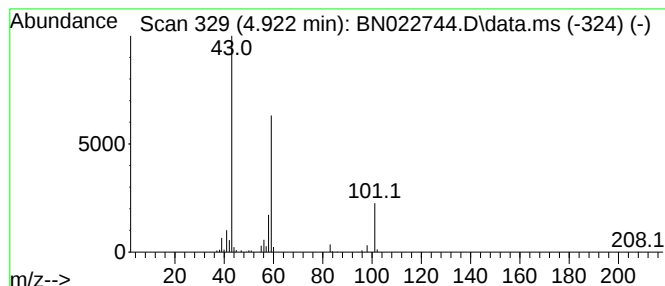
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.922	4.55 ng/ul	84659	1,4-Dichlorobenzene-d4	7.852

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
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	38		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
Data File : BN022744.D
Acq On : 18 Nov 2022 06:31
Operator : CG/JU
Sample : N5574-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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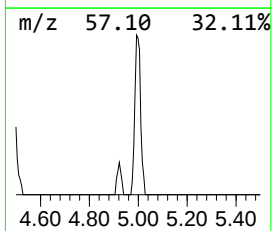
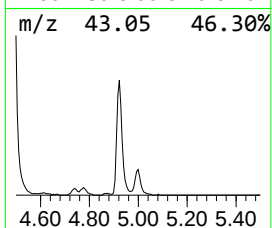
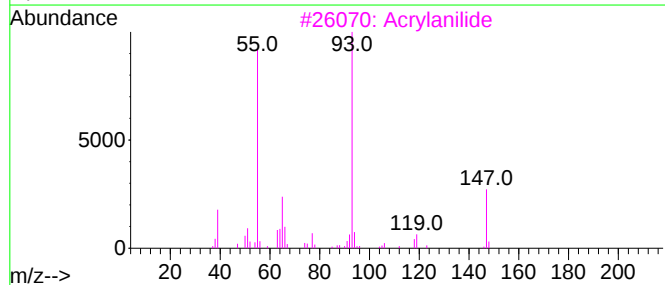
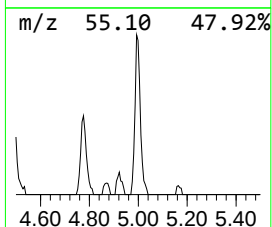
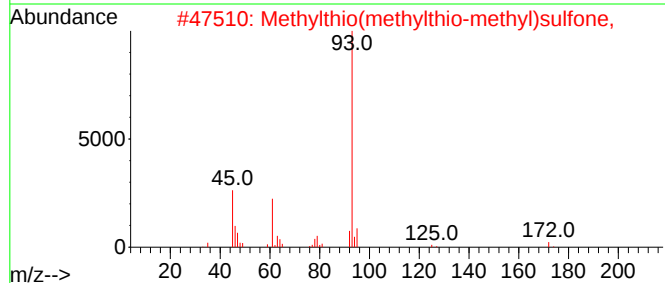
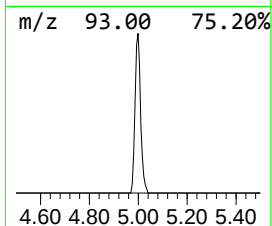
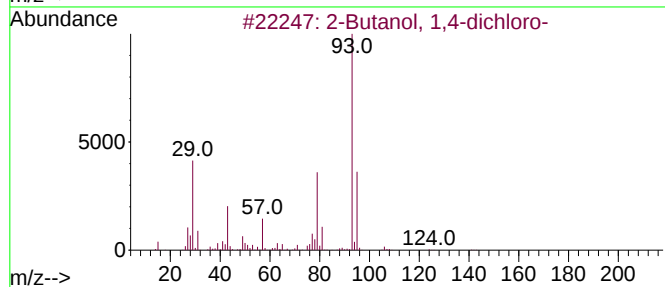
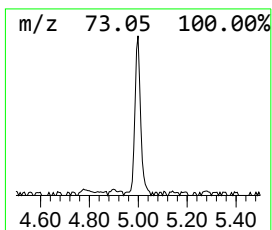
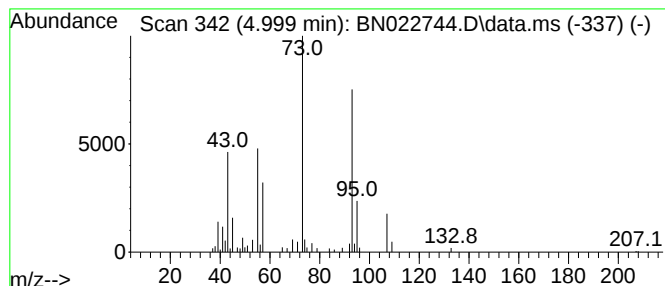
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.999	3.69 ng/ul	68639	1,4-Dichlorobenzene-d4	7.852

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Butanol, 1,4-dichloro-	142	C4H8Cl2O	002419-74-1	33
2			Methylthio(methylthio-methyl)sul...	172	C3H8O2S3	058000-45-6	28
3			Acrylanilide	147	C9H9NO	002210-24-4	28
4			1,3-Dioxolane, 2-(chloromethyl)-	122	C4H7ClO2	002568-30-1	9
5			Propanamide, N-phenyl-	149	C9H11NO	000620-71-3	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
Data File : BN022744.D
Acq On : 18 Nov 2022 06:31
Operator : CG/JU
Sample : N5574-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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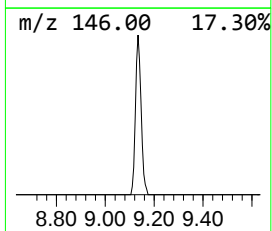
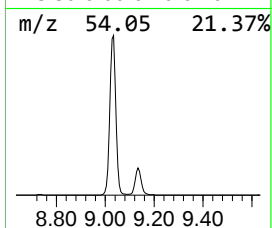
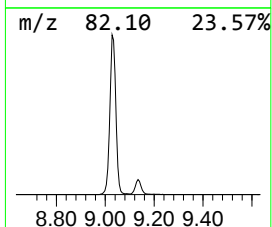
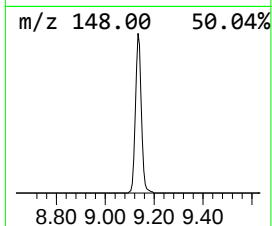
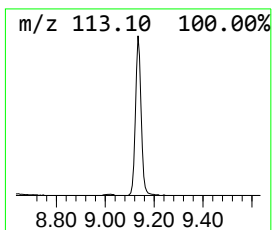
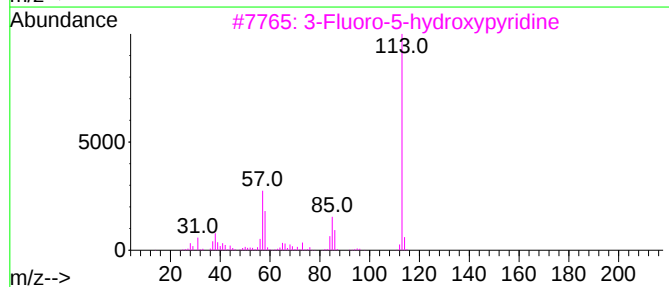
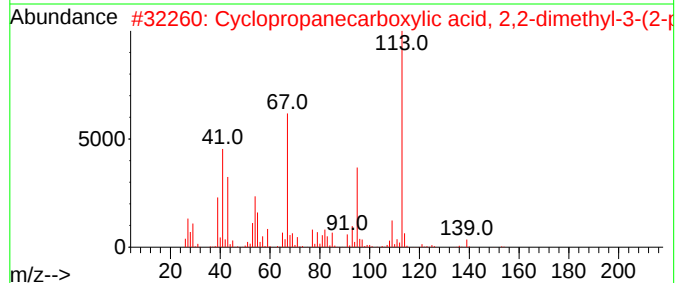
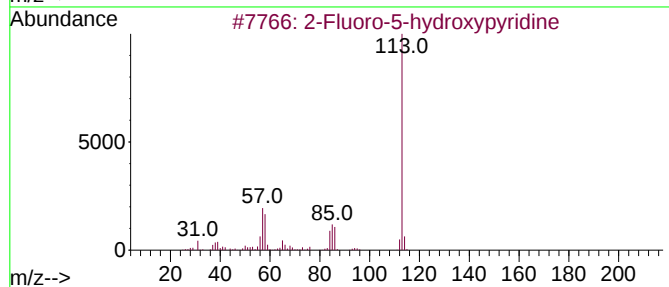
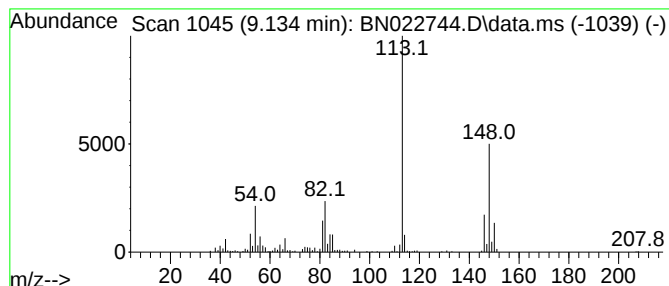
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.134	16.73 ng/ul	311070	1,4-Dichlorobenzene-d4	7.852

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Fluoro-5-hydroxypyridine	113	C5H4FNO	055758-32-2	38
2			Cyclopropanecarboxylic acid, 2,2...	154	C9H14O2	074685-76-0	37
3			3-Fluoro-5-hydroxypyridine	113	C5H4FNO	209328-55-2	35
4			3-Fluoro-4-pyridinol	113	C5H4FNO	022282-73-1	35
5			4-Fluoro-2-pyridone	113	C5H4FNO	096530-75-5	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
Data File : BN022744.D
Acq On : 18 Nov 2022 06:31
Operator : CG/JU
Sample : N5574-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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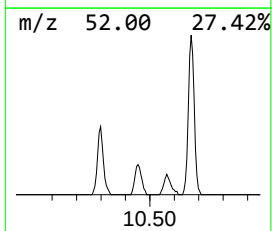
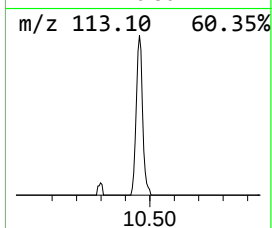
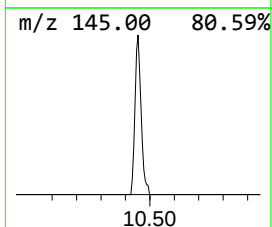
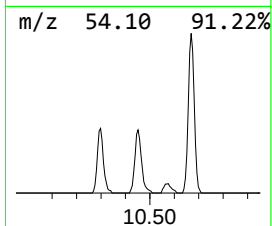
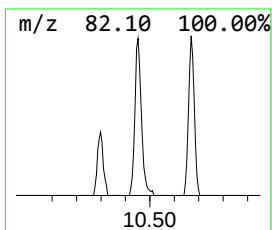
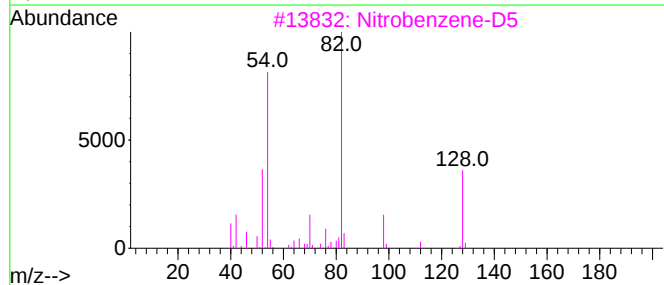
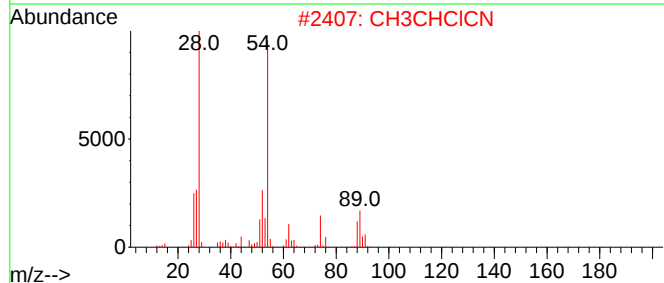
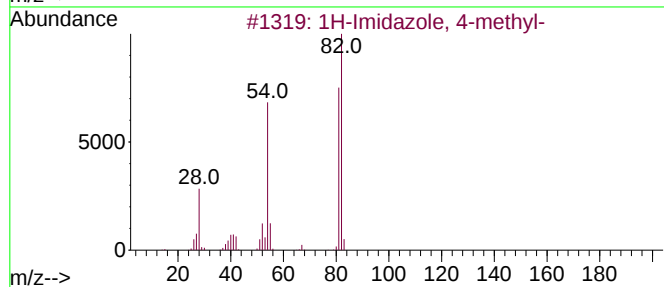
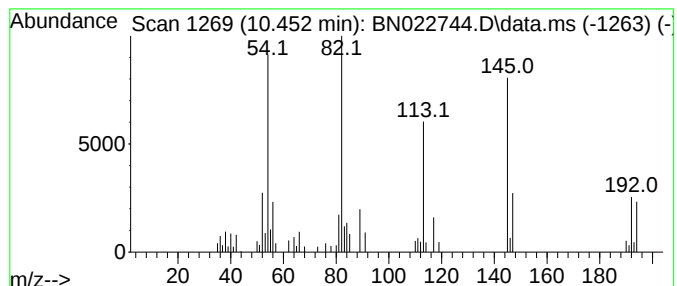
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-03 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.451	2.76 ng/ul	83757	Naphthalene-d8	10.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	47
2			CH3CHClCN	89	C3H4ClN	001617-17-0	38
3			Nitrobenzene-D5	128	C6D5NO2	004165-60-0	36
4			1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	28
5			CH3CHClCN	89	C3H4ClN	001617-17-0	28



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
Data File : BN022744.D
Acq On : 18 Nov 2022 06:31
Operator : CG/JU
Sample : N5574-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

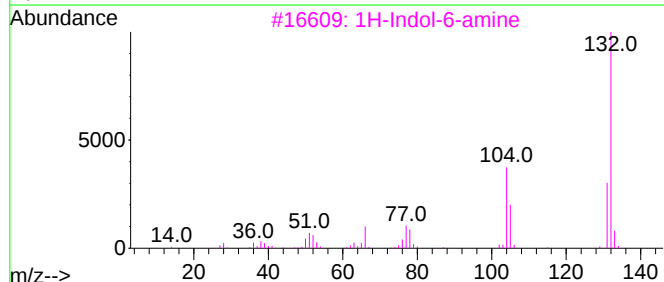
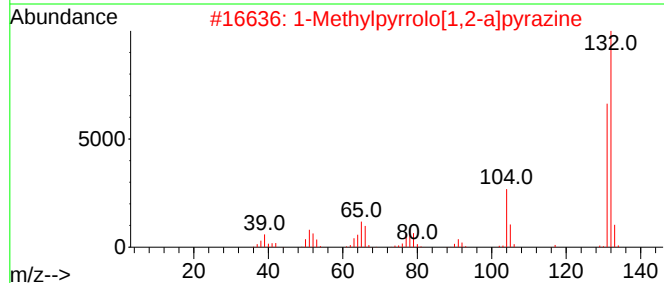
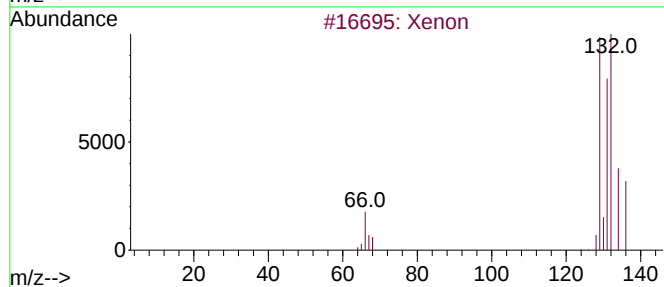
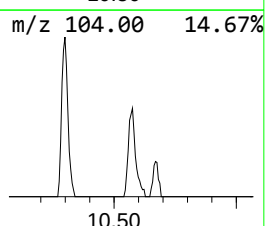
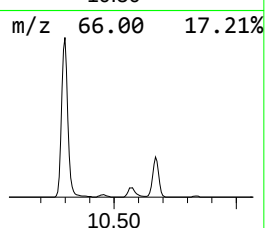
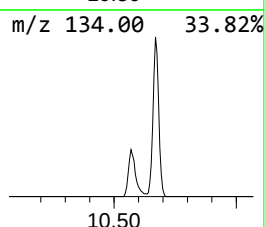
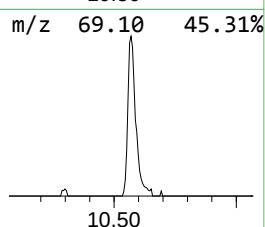
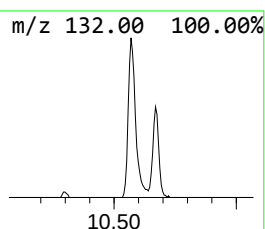
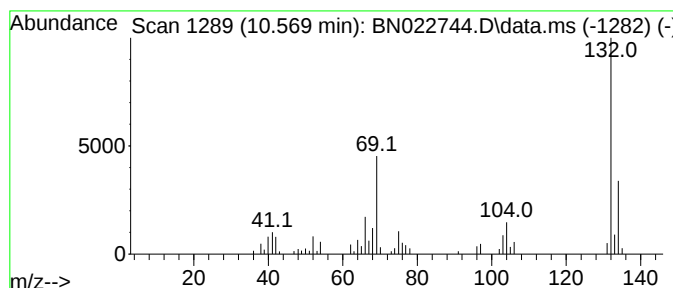
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Xenon Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.569	3.74 ng/ul	113574	Naphthalene-d8	10.669	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Xenon	132	Xe	007440-63-3	59
2	1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	58
3	1H-Indol-6-amine	132	C8H8N2	005318-27-4	53
4	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	53
5	5-Methylimidazo[1,2-a]pyridine	132	C8H8N2	000933-69-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
Data File : BN022744.D
Acq On : 18 Nov 2022 06:31
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Sample : N5574-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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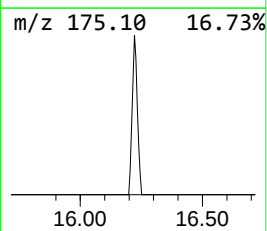
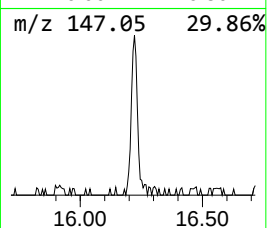
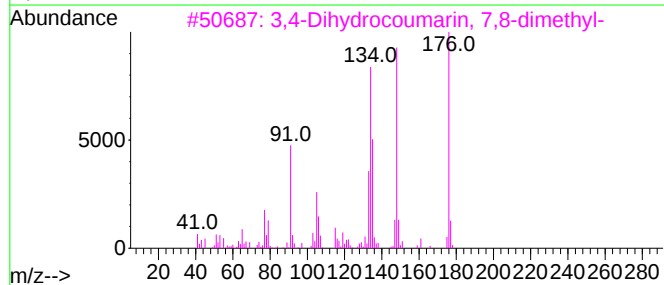
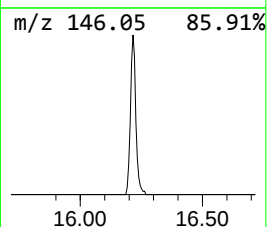
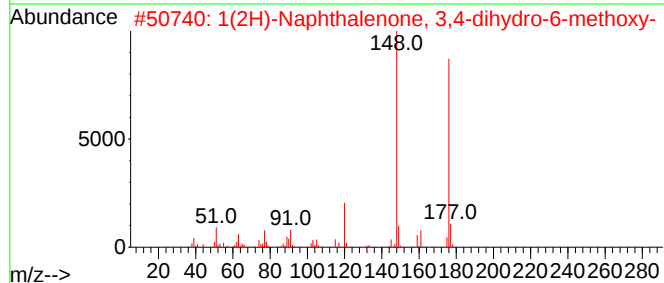
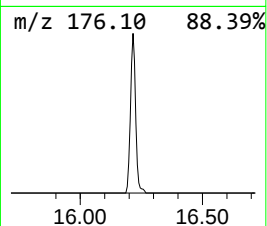
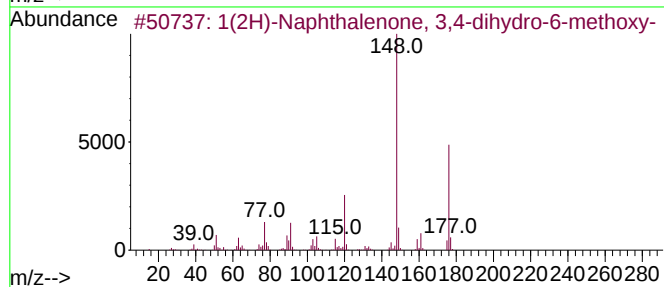
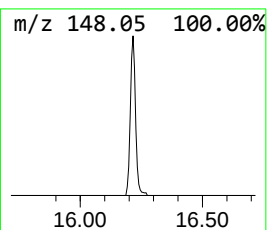
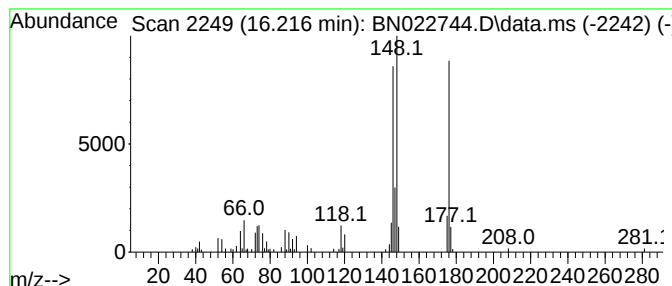
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-04 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.216	2.08 ng/ul	110689	Phenanthrene-d10	17.281

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Naphthalenone, 3,4-dihydro...	176	C11H12O2	001078-19-9	47		
2	1(2H)-Naphthalenone, 3,4-dihydro...	176	C11H12O2	001078-19-9	47		
3	3,4-Dihydrocoumarin, 7,8-dimethyl-	176	C11H12O2	040614-32-2	47		
4	Pyridine-3-carbonitrile, 1-ethyl...	176	C10H12N2O	1010316-81-3	43		
5	Furan, 3-bromo-	146	C4H3BrO	022037-28-1	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
Data File : BN022744.D
Acq On : 18 Nov 2022 06:31
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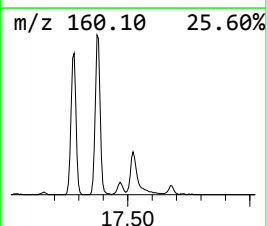
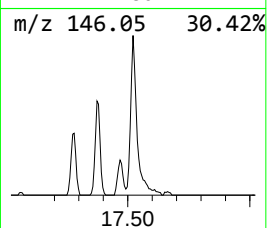
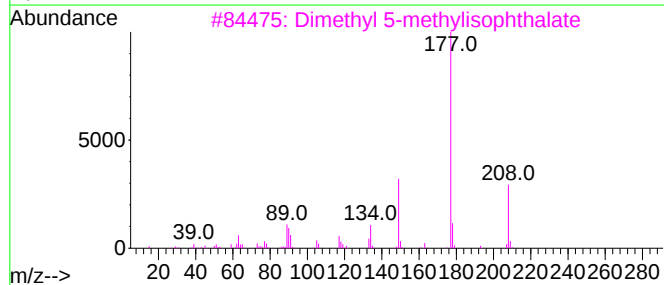
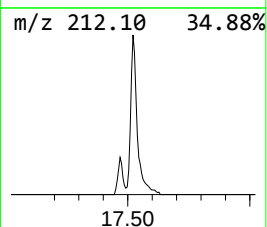
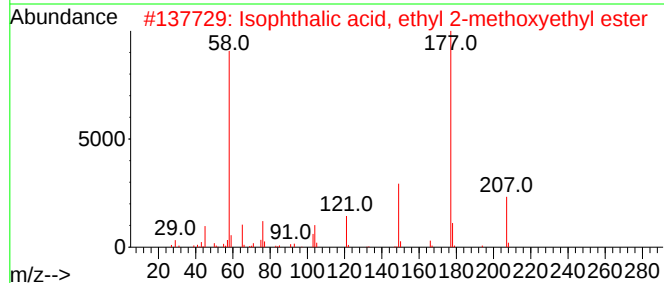
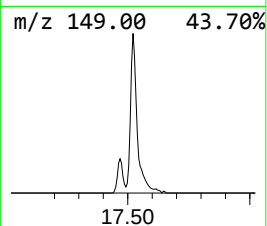
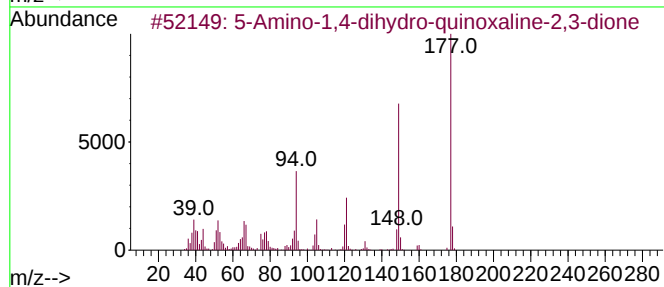
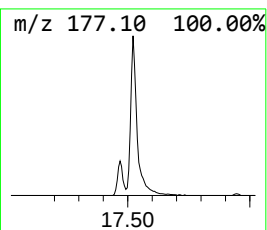
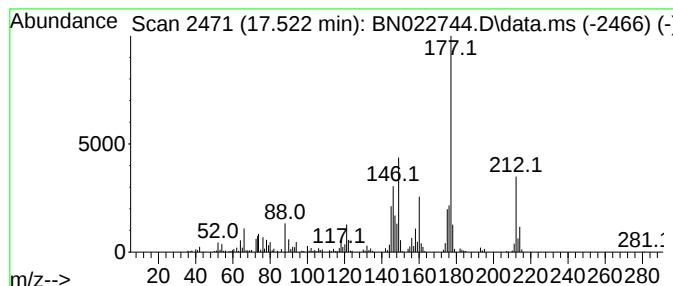
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-05 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.522	8.27 ng/ul	439131	Phenanthrene-d10	17.281

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		5-Amino-1,4-dihydro-quinoxaline-...	177	C8H7N3O2	076097-87-5	38
2		Isophthalic acid, ethyl 2-methox...	252	C13H16O5	1000345-85-7	38
3		Dimethyl 5-methylisophthalate	208	C11H12O4	017649-58-0	38
4		Isophthalic acid, ethyl 3-methyl...	262	C15H18O4	1000343-93-2	38
5		4-Chloro-2,5,6-trimethylthieno[2...	212	C9H9ClN2S	083548-58-7	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
Data File : BN022744.D
Acq On : 18 Nov 2022 06:31
Operator : CG/JU
Sample : N5574-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
B0AA0

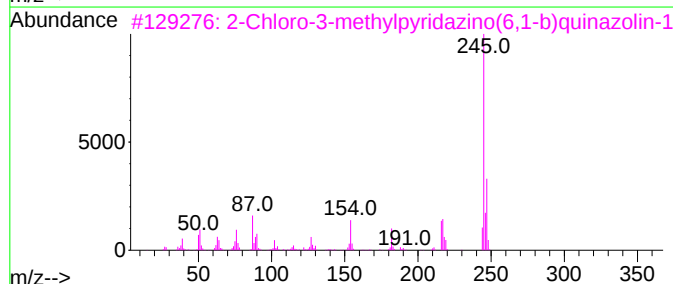
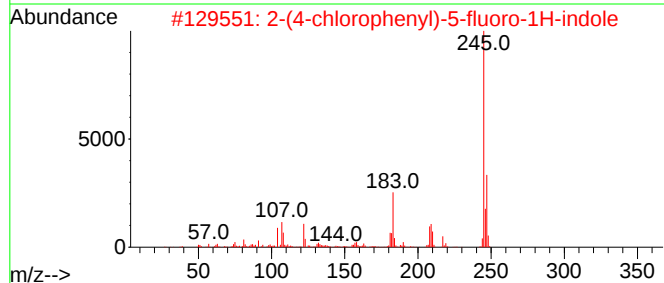
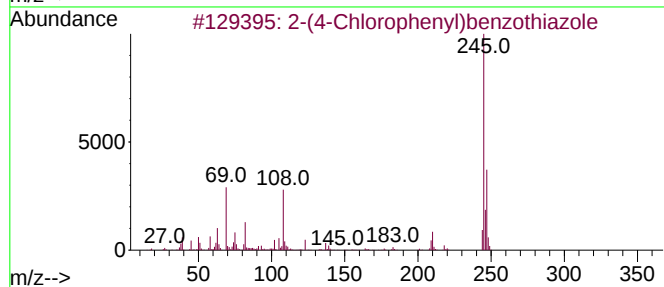
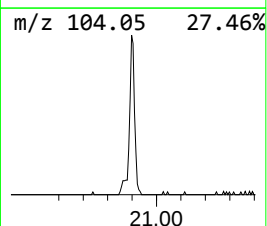
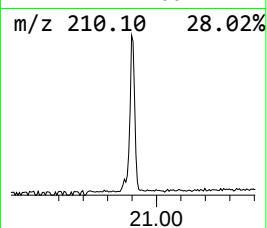
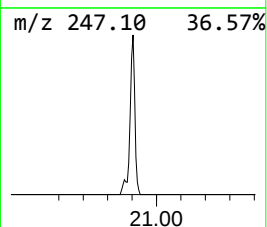
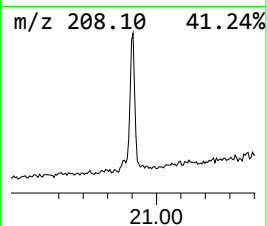
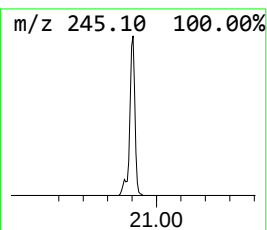
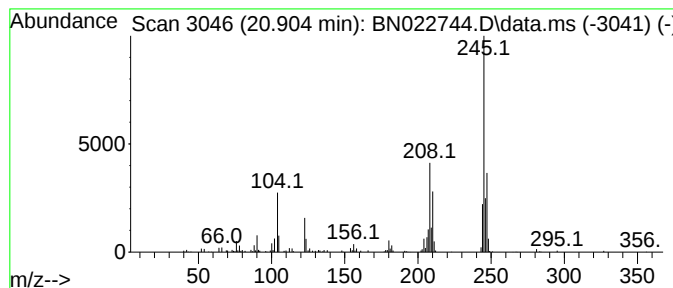
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-06 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.904	3.14 ng/ul	193405	Chrysene-d12	21.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(4-Chlorophenyl)benzothiazole	245	C13H8ClNS	006265-91-4	49		
2	2-(4-chlorophenyl)-5-fluoro-1H-i...	245	C14H9ClFN	881040-32-0	49		
3	2-Chloro-3-methylpyridazino(6,1-...	245	C12H8ClN3O	001703-03-3	46		
4	2-Chloro-4-methylpyridazino(6,1-...	245	C12H8ClN3O	001703-04-4	43		
5	4-Chloro-2-methyl-6-trifluoromet...	245	C11H7ClF3N	867167-05-3	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
Data File : BN022744.D
Acq On : 18 Nov 2022 06:31
Operator : CG/JU
Sample : N5574-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
B0AA0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.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
3-Methyl-3-chlo...	3.064	25.3	ng/u1	470106	1	7.852	371909	20.0
1,1-Dimethyl-3-...	4.481	81.3	ng/u1	1512280	1	7.852	371909	20.0
Butane, 2,3-dic...	4.775	8.9	ng/u1	165444	1	7.852	371909	20.0
2-Pentanone, 4-...	4.922	4.5	ng/u1	84659	1	7.852	371909	20.0
unknown-01	4.999	3.7	ng/u1	68639	1	7.852	371909	20.0
unknown-02	9.134	16.7	ng/u1	311070	1	7.852	371909	20.0
unknown-03	10.451	2.8	ng/u1	83757	2	10.669	607519	20.0
Xenon	10.569	3.7	ng/u1	113574	2	10.669	607519	20.0
unknown-04	16.216	2.1	ng/u1	110689	4	17.281	1062430	20.0
unknown-05	17.522	8.3	ng/u1	439131	4	17.281	1062430	20.0
unknown-06	20.904	3.1	ng/u1	193405	5	21.480	1232190	20.0