

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
 Data File : BP010380.D
 Acq On : 17 May 2022 04:48
 Operator : CG/JU
 Sample : N2831-06
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YB2K3

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-BP051322.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010380.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.311	34	38	41	rBV3	24514	33309	0.32%	0.048%
2	3.340	41	43	48	rVB	28292	38187	0.37%	0.055%
3	3.628	89	92	102	rVB	18164	33219	0.32%	0.048%
4	3.770	113	116	133	rBV	145891	316315	3.04%	0.457%
5	4.087	167	170	172	rBV3	8962	10801	0.10%	0.016%
6	4.423	223	227	232	rBV	12431	18671	0.18%	0.027%
7	4.540	236	247	255	rVV	59473	197294	1.89%	0.285%
8	4.617	255	260	270	rVB2	60447	116015	1.11%	0.168%
9	5.011	322	327	333	rVB6	5661	11913	0.11%	0.017%
10	5.946	480	486	491	rBV	19887	31878	0.31%	0.046%
11	7.064	670	676	689	rBV2	145432	284416	2.73%	0.411%
12	7.228	695	704	713	rBV	591338	982455	9.43%	1.419%
13	7.299	713	716	724	rVB	25560	45070	0.43%	0.065%
14	7.428	731	738	751	rBV	603479	1029988	9.88%	1.488%
15	7.528	751	755	761	rVB7	6690	10786	0.10%	0.016%
16	7.899	809	818	825	rBV	548147	930191	8.93%	1.344%
17	7.964	825	829	838	rVB	48030	83986	0.81%	0.121%
18	8.605	925	938	951	rBV	406311	869473	8.34%	1.256%
19	8.722	956	958	965	rVB8	7947	13186	0.13%	0.019%
20	9.058	1001	1015	1028	rBV	785369	1402873	13.46%	2.026%
21	9.381	1062	1070	1076	rBV5	12072	30995	0.30%	0.045%
22	9.499	1085	1090	1096	rVB5	8528	13744	0.13%	0.020%
23	9.781	1126	1138	1156	rBV	642333	1108749	10.64%	1.601%
24	10.158	1197	1202	1207	rBV4	7189	13018	0.12%	0.019%
25	10.228	1209	1214	1221	rVV5	16715	34376	0.33%	0.050%
26	10.352	1224	1235	1253	rVV	576166	1589817	15.26%	2.296%
27	10.699	1285	1294	1305	rBV	811520	1367565	13.12%	1.975%
28	10.834	1308	1317	1328	rBV	740007	1362715	13.08%	1.968%
29	10.928	1331	1333	1340	rVB8	9555	17613	0.17%	0.025%
30	11.552	1433	1439	1444	rBV	67011	95232	0.91%	0.138%
31	11.675	1444	1460	1478	rVV	2145565	6653723	63.85%	9.611%
32	11.957	1504	1508	1512	rBV3	12236	20627	0.20%	0.030%
33	12.193	1543	1548	1552	rVB8	9425	16213	0.16%	0.023%
34	12.287	1559	1564	1567	rBV3	14505	25573	0.25%	0.037%
35	12.334	1567	1572	1583	rVB	228474	368937	3.54%	0.533%
36	12.834	1653	1657	1663	rVB8	6888	12641	0.12%	0.018%

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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-BP051322.M
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37	12.899	1663	1668	1672	rBV6	8586	17286	0.17%	0.025%
38	13.363	1743	1747	1754	rVB3	13101	23548	0.23%	0.034%
39	13.881	1829	1835	1839	rBV2	46142	71789	0.69%	0.104%
40	13.940	1839	1845	1857	rVB	1837357	2550617	24.48%	3.684%
41	14.222	1881	1893	1905	rBV	2028076	3106076	29.81%	4.486%
42	14.316	1906	1909	1916	rVB7	14233	23618	0.23%	0.034%
43	14.393	1919	1922	1926	rVV6	6577	11032	0.11%	0.016%
44	14.434	1926	1929	1933	rVV4	10500	16915	0.16%	0.024%
45	14.528	1936	1945	1953	rBV	1362941	2003367	19.23%	2.894%
46	14.610	1957	1959	1963	rVV4	13389	18365	0.18%	0.027%
47	14.657	1963	1967	1971	rVV3	28609	36815	0.35%	0.053%
48	14.734	1975	1980	1992	rBV2	85559	197032	1.89%	0.285%
49	14.875	2001	2004	2009	rBV6	8015	14674	0.14%	0.021%
50	14.951	2014	2017	2022	rVB6	13905	21581	0.21%	0.031%
51	15.122	2034	2046	2047	rBV3	74486	230491	2.21%	0.333%
52	15.522	2108	2114	2127	rVB	2703083	3823216	36.69%	5.522%
53	15.640	2129	2134	2137	rVV	399796	659857	6.33%	0.953%
54	15.663	2137	2138	2149	rVV	365503	409955	3.93%	0.592%
55	15.763	2151	2155	2162	rVB2	34679	53790	0.52%	0.078%
56	15.863	2169	2172	2176	rBV6	13955	20751	0.20%	0.030%
57	16.034	2197	2201	2206	rVB2	21649	30239	0.29%	0.044%
58	16.310	2244	2248	2253	rVB2	27834	39663	0.38%	0.057%
59	16.422	2264	2267	2271	rBV5	7050	11555	0.11%	0.017%
60	16.493	2275	2279	2283	rVB5	13271	18738	0.18%	0.027%
61	17.057	2369	2375	2376	rBV3	90749	144198	1.38%	0.208%
62	17.081	2376	2379	2386	rVB	146509	222824	2.14%	0.322%
63	17.281	2407	2413	2422	rVB	1681633	2407819	23.11%	3.478%
64	17.381	2424	2430	2440	rVV	2927843	4236710	40.66%	6.120%
65	17.510	2449	2452	2455	rBV	32866	40191	0.39%	0.058%
66	17.951	2522	2527	2533	rBV2	280909	374306	3.59%	0.541%
67	18.187	2560	2567	2574	rBV2	427403	996061	9.56%	1.439%
68	18.398	2595	2603	2623	rBV	6158914	10420109	100.00%	15.051%
69	19.398	2770	2773	2783	rVB3	82089	127232	1.22%	0.184%
70	19.616	2804	2810	2822	rVB	4031321	5194604	49.85%	7.503%
71	20.051	2882	2884	2888	rVB4	29439	32180	0.31%	0.046%
72	20.139	2896	2899	2903	rVB2	121073	128090	1.23%	0.185%
73	20.622	2978	2981	2985	rVB	100953	110371	1.06%	0.159%
74	21.010	3045	3047	3053	rBV6	52890	67473	0.65%	0.097%
75	21.075	3054	3058	3068	rBV	646847	882003	8.46%	1.274%
76	21.363	3102	3107	3121	rBV	2088663	2839357	27.25%	4.101%

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Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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77	21.516	3129	3133	3138	rVB	203346	246927	2.37%	0.357%
78	21.980	3209	3212	3220	rVB2	226729	330157	3.17%	0.477%
79	22.486	3294	3298	3304	rVB	207866	300931	2.89%	0.435%
80	23.045	3390	3393	3400	rVB2	187190	287917	2.76%	0.416%
81	23.639	3486	3494	3500	rBV2	2244583	4650166	44.63%	6.717%
82	23.792	3513	3520	3530	rVB2	1256820	2590769	24.86%	3.742%

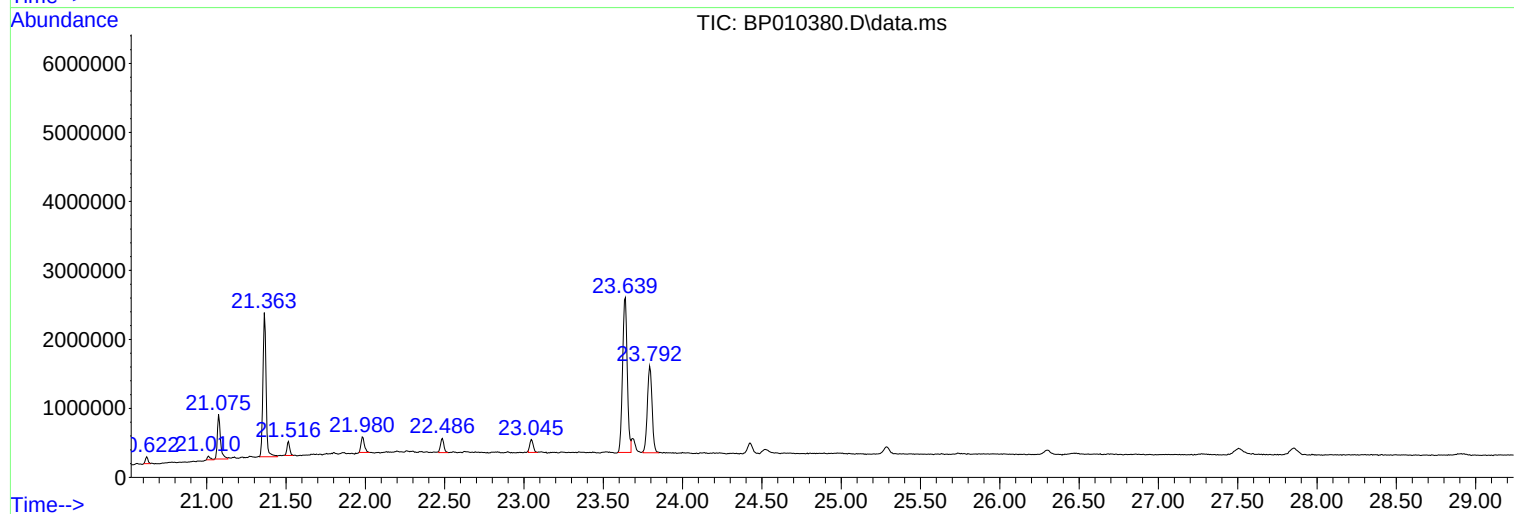
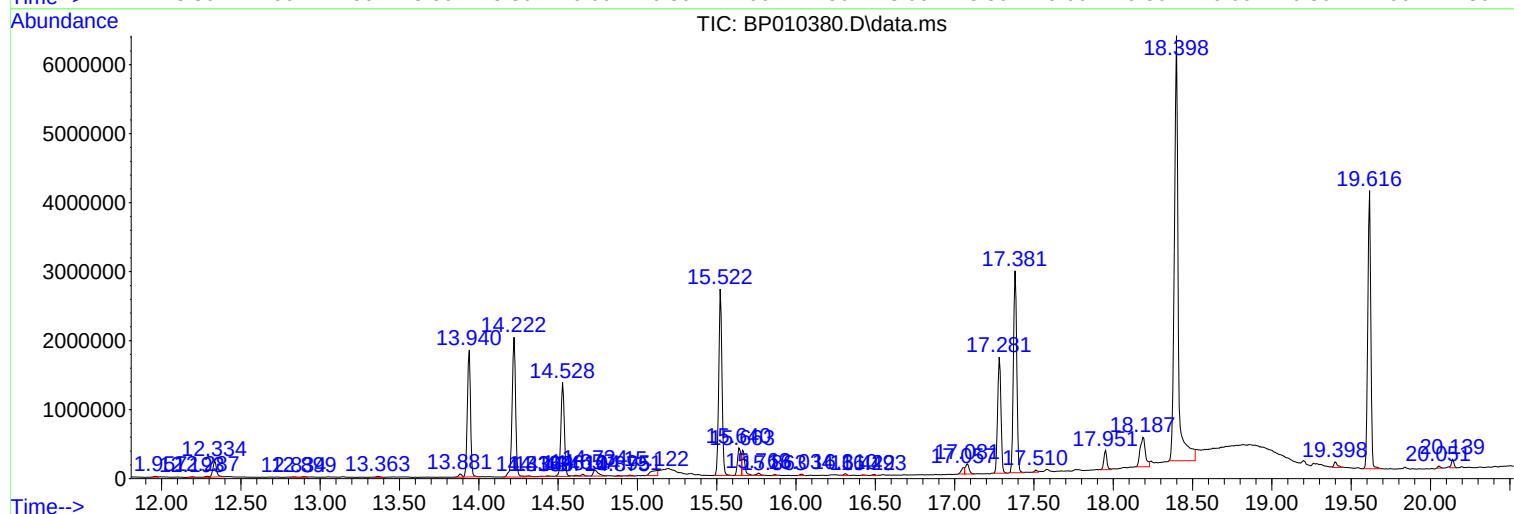
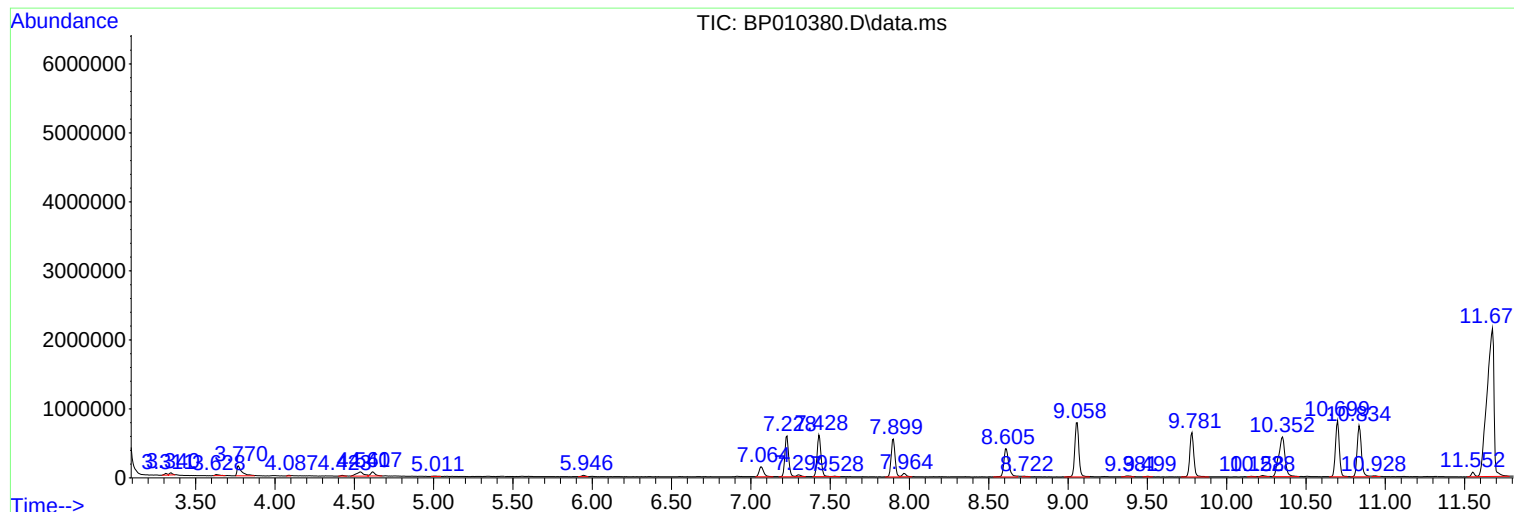
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 28 Sample Multiplier: 1

Instrument :
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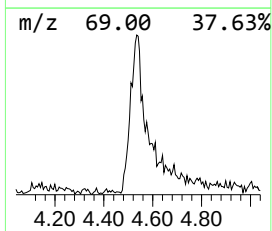
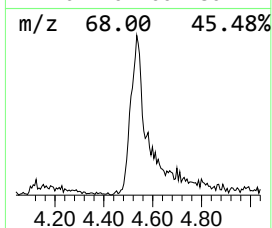
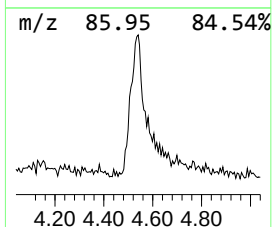
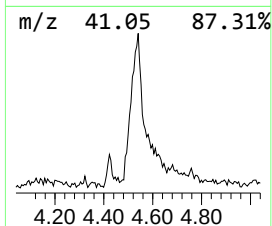
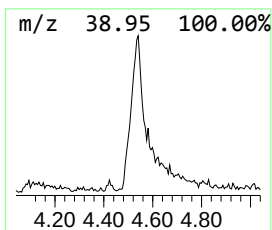
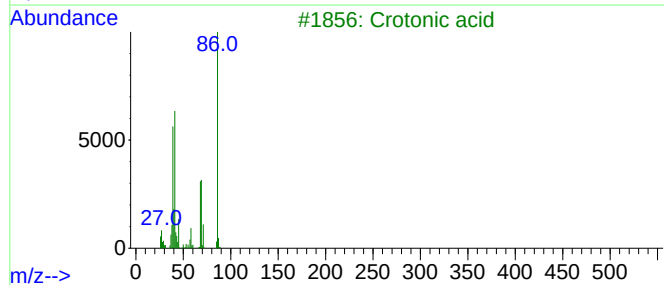
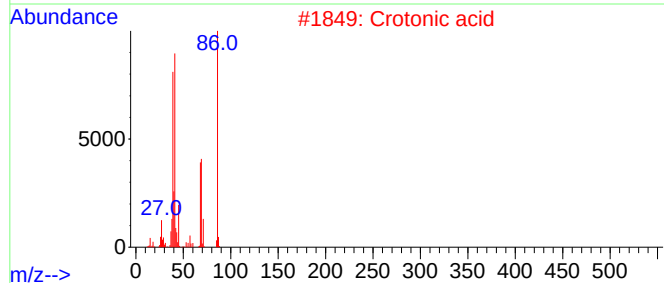
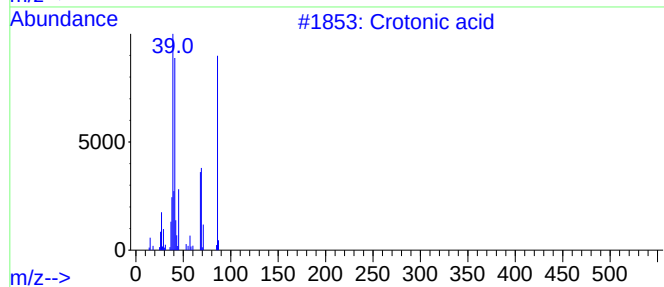
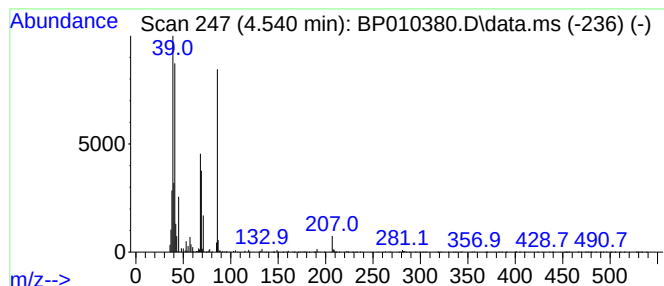
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Crotonic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.540	4.24 ng/ul	197294	1,4-Dichlorobenzene-d4	7.899

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Crotonic acid	86	C4H6O2	003724-65-0	96
2			Crotonic acid	86	C4H6O2	003724-65-0	95
3			Crotonic acid	86	C4H6O2	003724-65-0	91
4			2-Butenoic acid, (E)-	86	C4H6O2	000107-93-7	90
5			Crotonic acid	86	C4H6O2	003724-65-0	90



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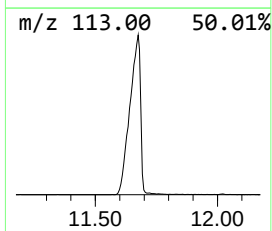
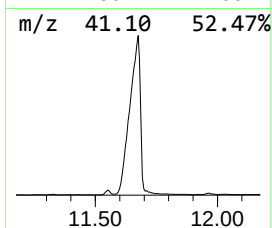
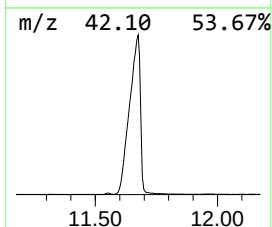
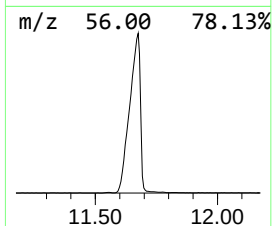
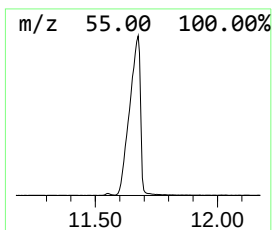
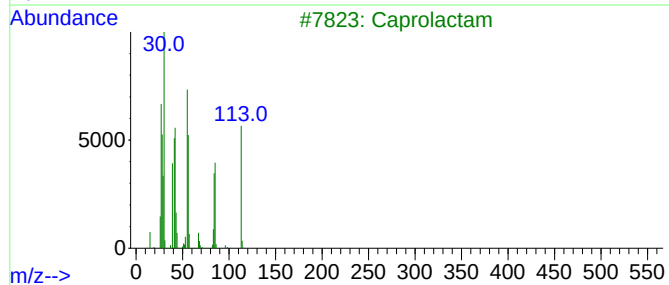
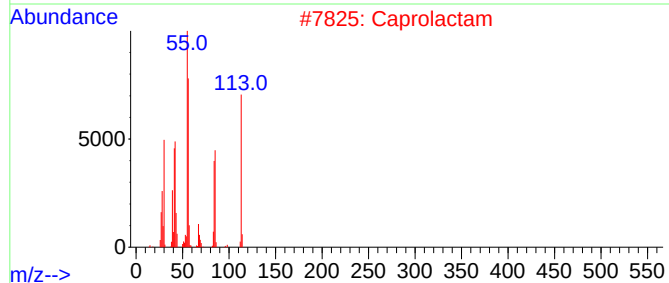
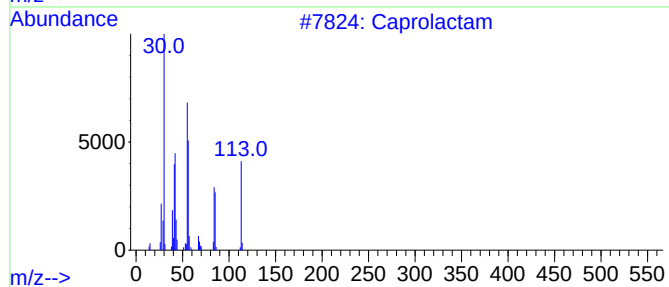
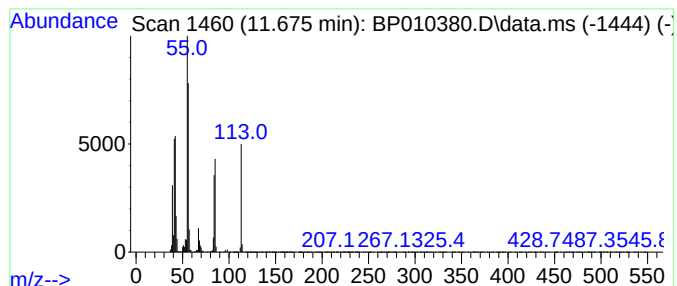
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Capro lactam Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.675	97.31 ng/ul	6653720	Naphthalene-d8	10.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Caprolactam	113	C6H11NO	000105-60-2	91
2			Caprolactam	113	C6H11NO	000105-60-2	90
3			Caprolactam	113	C6H11NO	000105-60-2	64
4			1-Hexene	84	C6H12	000592-41-6	55
5			Caprolactam	113	C6H11NO	000105-60-2	52



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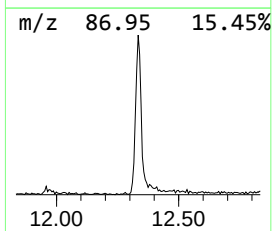
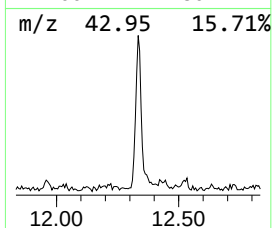
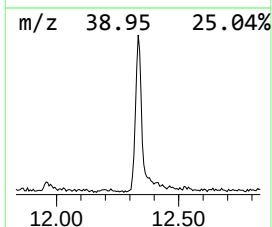
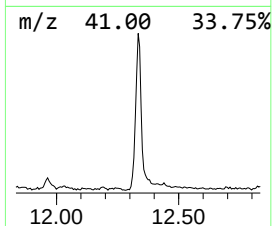
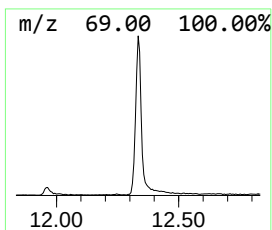
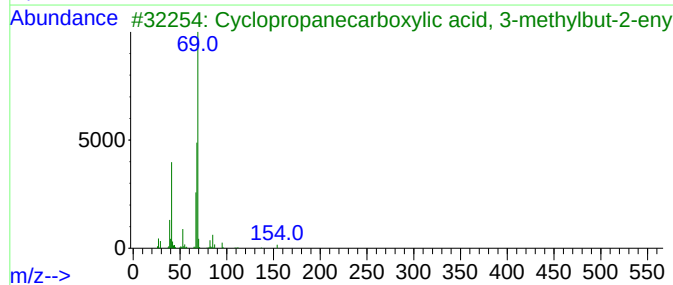
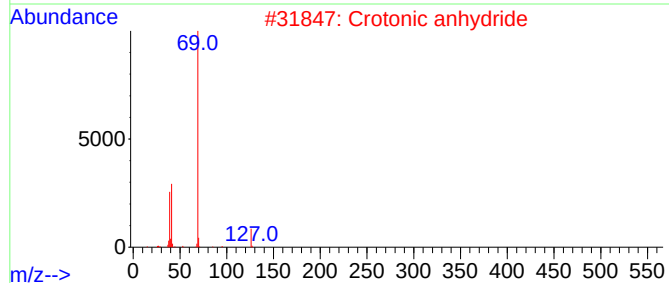
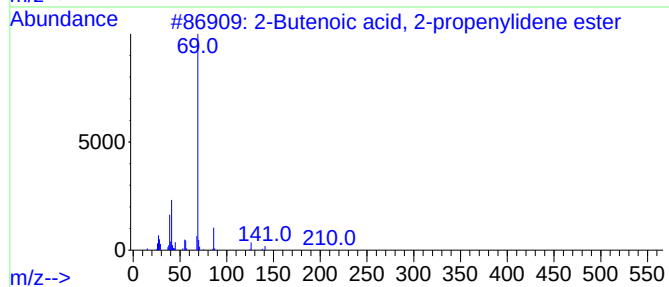
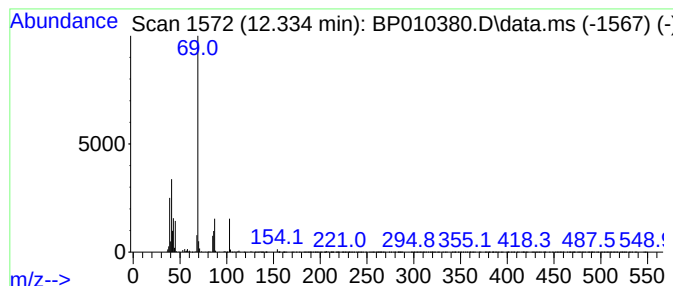
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 2-Butenoic acid, 2-propenyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.334	5.40 ng/ul	368937	Naphthalene-d8	10.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Butenoic acid, 2-propenylidene...	210	C11H14O4	055030-70-1	50
2			Crotonic anhydride	154	C8H10O3	000623-68-7	47
3			Cyclopropanecarboxylic acid, 3-m...	154	C9H14O2	1000299-37-4	40
4			Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	39
5			Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	39



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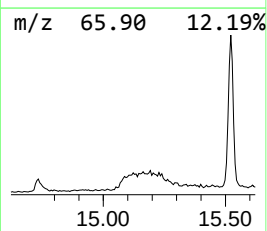
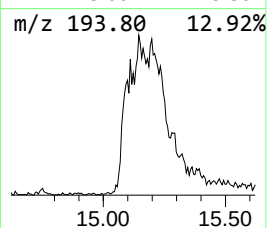
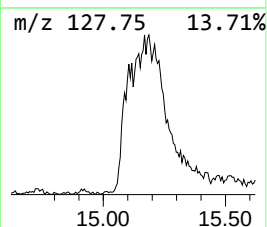
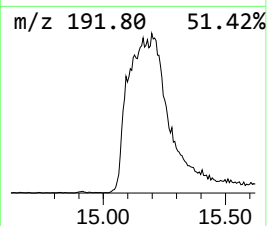
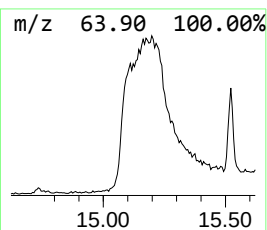
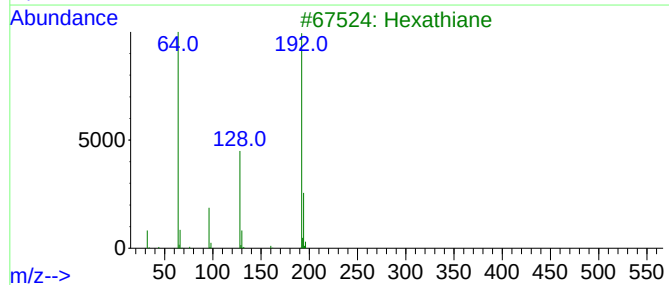
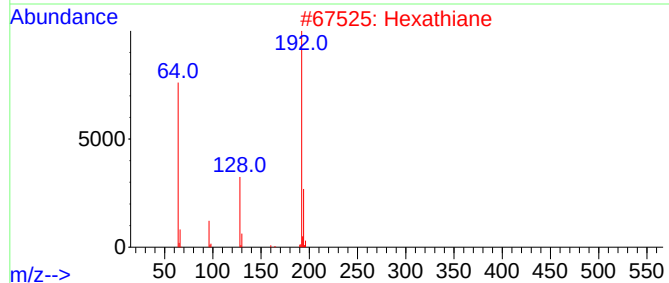
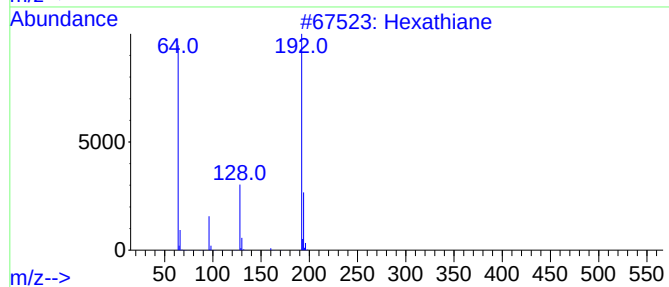
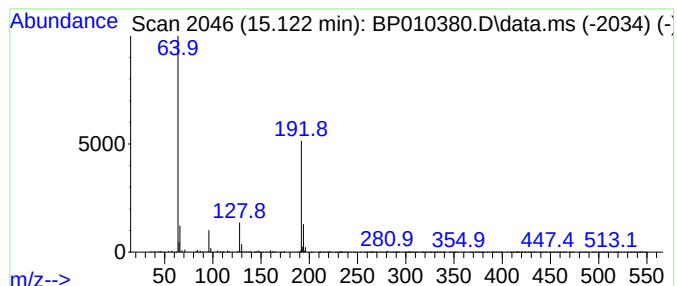
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Hexathiane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.122	2.30 ng/ul	230491	Acenaphthene-d10	14.528

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexathiane	192	S6	013798-23-7	94
2			Hexathiane	192	S6	013798-23-7	91
3			Hexathiane	192	S6	013798-23-7	91
4			Cyclic octaatomic sulfur	256	S8	010544-50-0	56
5			Di(tert-butyl) trisulfide, perfl...	534	C8F18S3	016005-64-4	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
Data File : BP010380.D
Acq On : 17 May 2022 04:48
Operator : CG/JU
Sample : N2831-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YB2K3

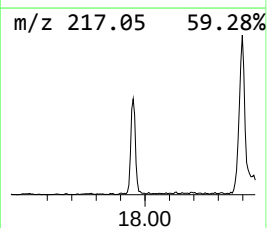
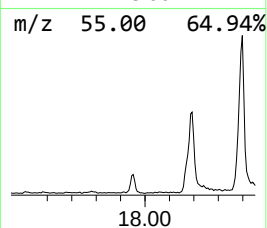
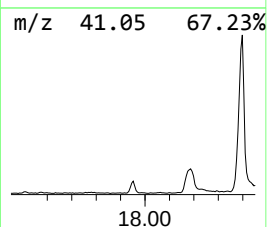
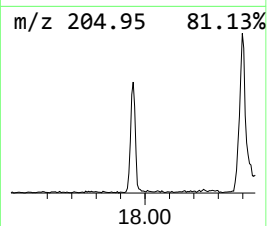
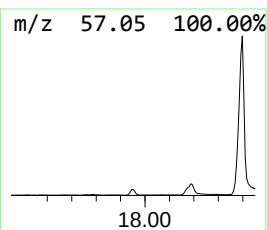
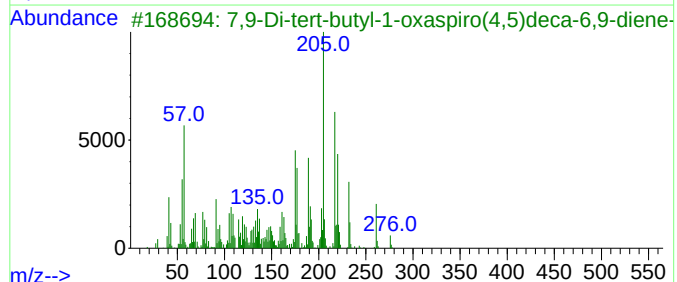
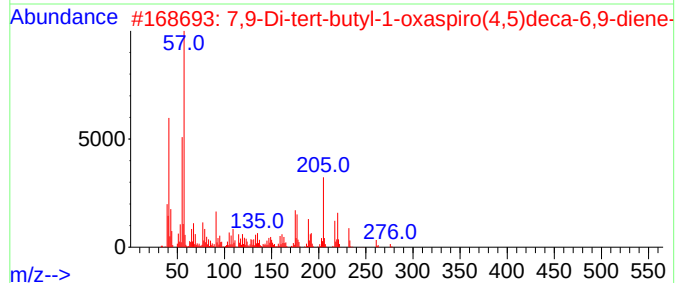
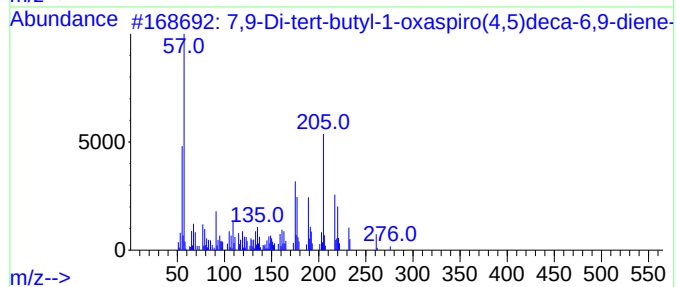
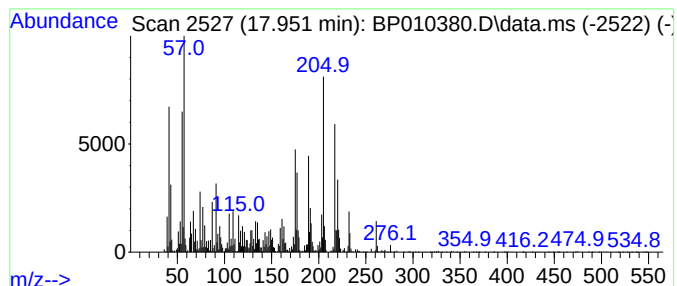
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.951	3.11 ng/ul	374306	Phenanthrene-d10	17.281

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	99		
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	98		
3	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	94		
4	(Z)-((2-Benzylideneheptyl)oxy)tr...	276	C17H28OSi	1000495-65-9	51		
5	1-(3-Amino-4,6-dimethylthieno[2,...	220	C11H12N2OS	052505-42-7	42		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
Data File : BP010380.D
Acq On : 17 May 2022 04:48
Operator : CG/JU
Sample : N2831-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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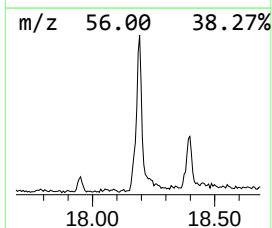
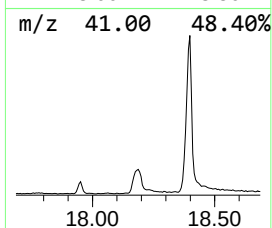
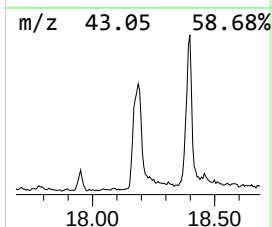
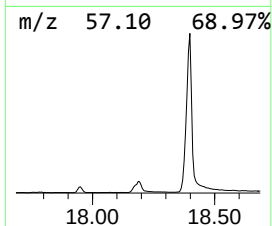
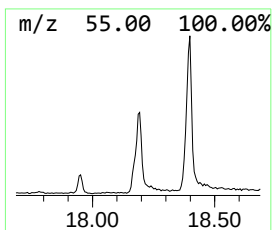
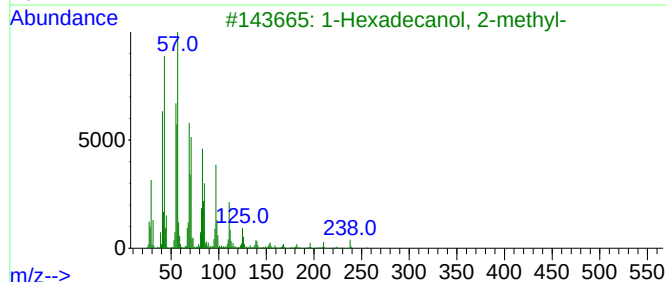
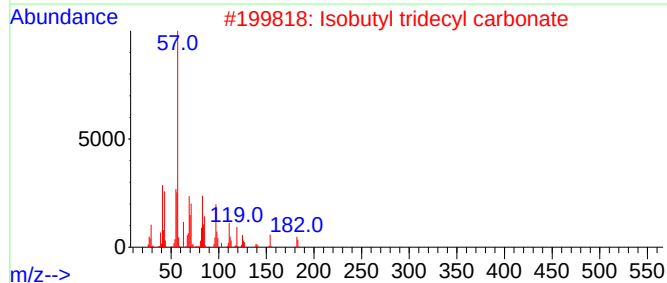
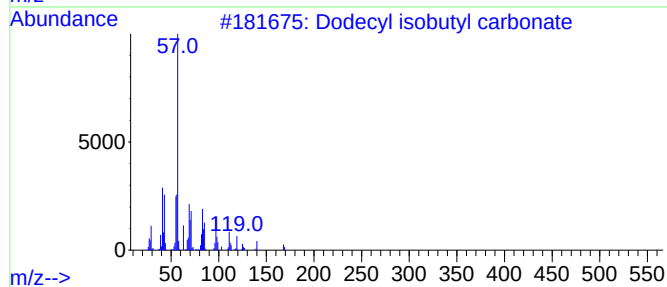
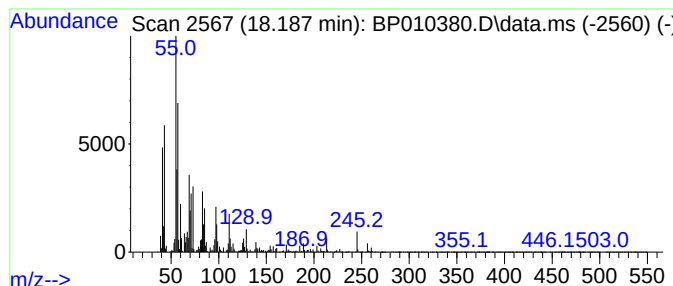
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Dodecyl isobutyl carbonate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.186	8.27 ng/ul	996061	Phenanthrene-d10	17.281

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecyl isobutyl carbonate	286	C17H34O3	959067-22-2	52
2			Isobutyl tridecyl carbonate	300	C18H36O3	959275-44-6	52
3			1-Hexadecanol, 2-methyl-	256	C17H36O	002490-48-4	52
4			2- Bromopropionic acid, pentadec...	362	C18H35BrO2	1000292-44-2	46
5			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
Data File : BP010380.D
Acq On : 17 May 2022 04:48
Operator : CG/JU
Sample : N2831-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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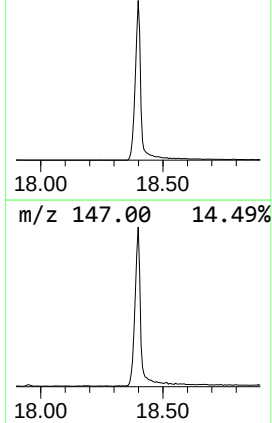
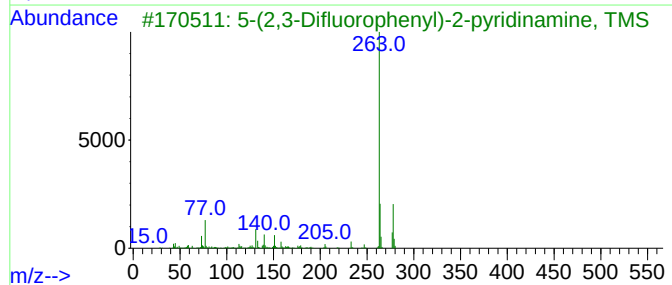
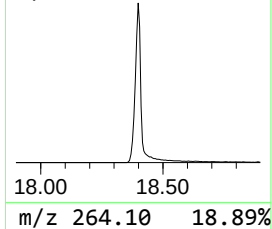
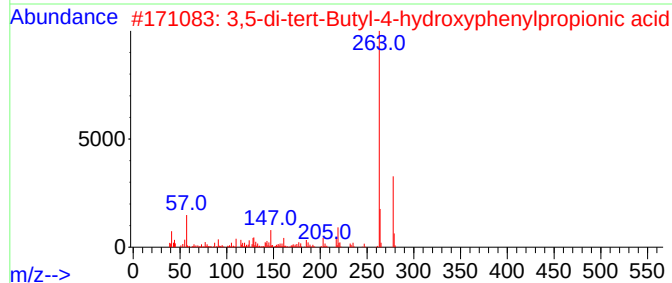
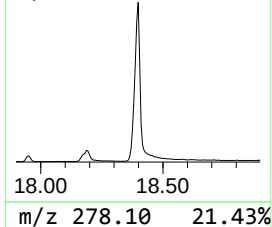
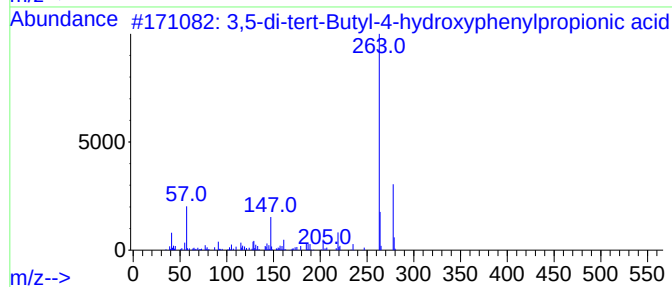
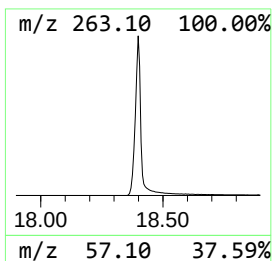
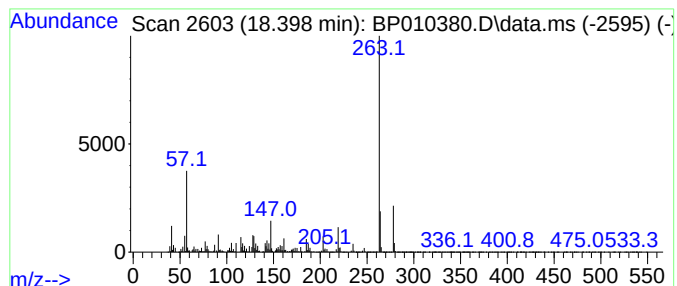
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.398	86.55 ng/ul	10420100	Phenanthrene-d10	17.281

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	97
2			3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	94
3			5-(2,3-Difluorophenyl)-2-pyridin...	278	C14H16F2N2Si	1000504-49-4	74
4			6,7-Dimethoxy-4(1H)-quinazolinon...	278	C13H18N2O3Si	1000479-65-5	72
5			Cyanic acid, (1-methylethylidene...	278	C17H14N2O2	001156-51-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
Data File : BP010380.D
Acq On : 17 May 2022 04:48
Operator : CG/JU
Sample : N2831-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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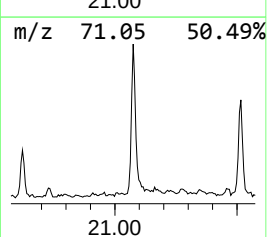
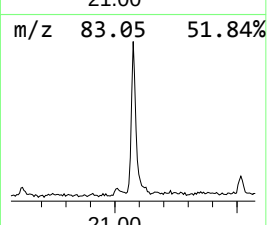
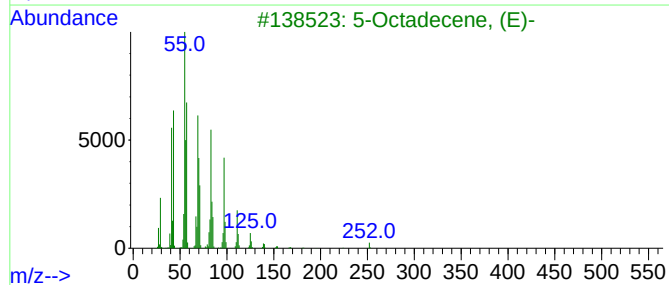
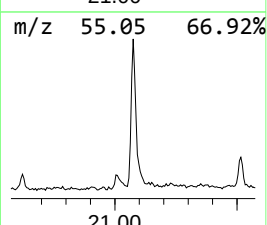
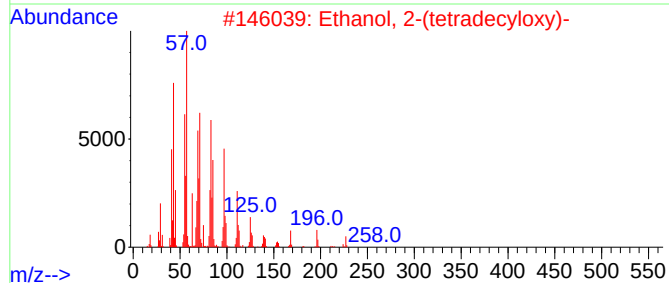
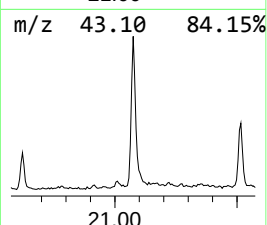
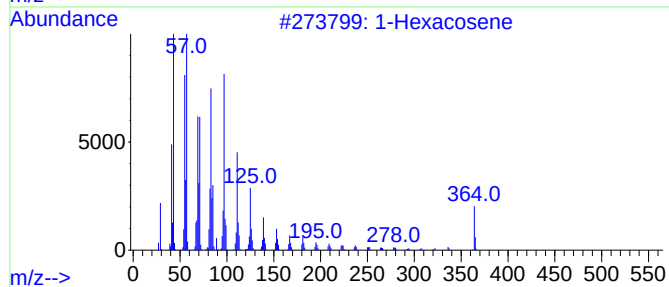
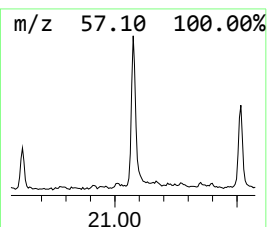
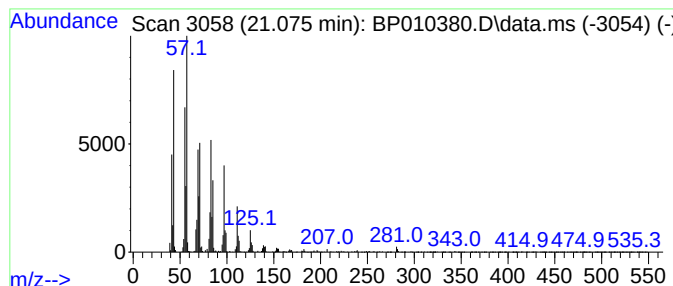
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.075	6.21 ng/ul	882003	Chrysene-d12	21.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene			364	C26H52	018835-33-1	98
2	Ethanol, 2-(tetradecyloxy)-			258	C16H34O2	002136-70-1	95
3	5-Octadecene, (E)-			252	C18H36	007206-21-5	93
4	1-Triacontanol			438	C30H62O	000593-50-0	91
5	17-Pentatriacontene			491	C35H70	006971-40-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
Data File : BP010380.D
Acq On : 17 May 2022 04:48
Operator : CG/JU
Sample : N2831-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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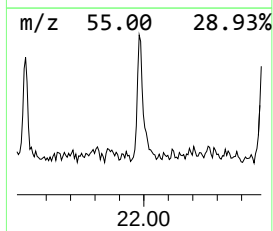
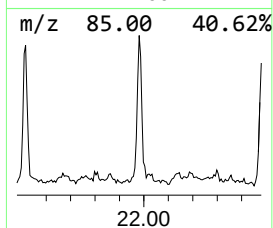
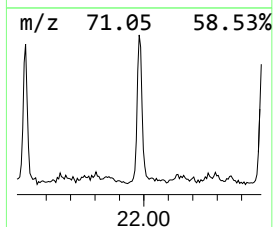
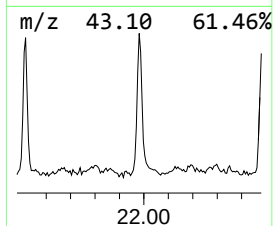
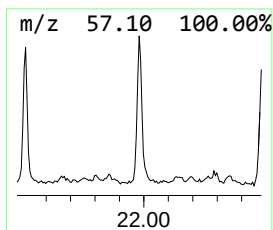
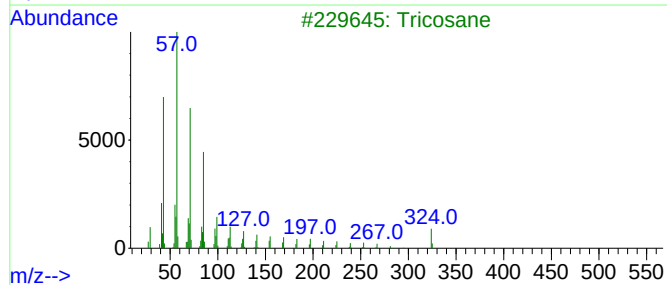
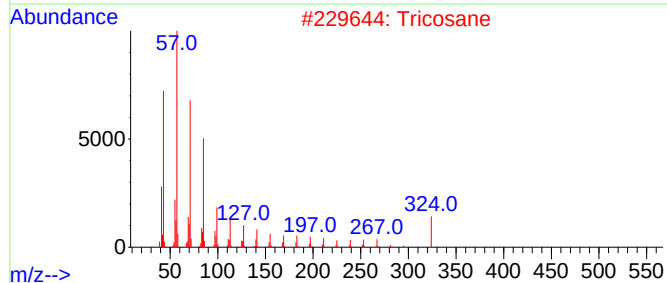
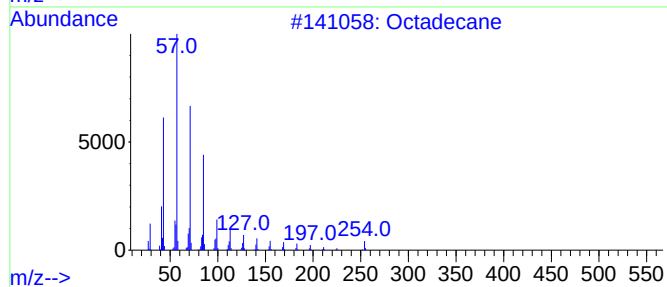
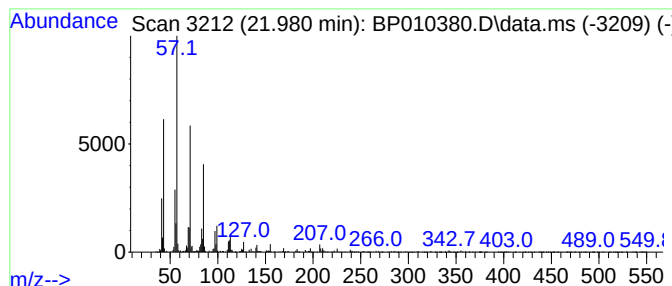
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.980	2.33 ng/ul	330157	Chrysene-d12	21.363

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane	254	C18H38	000593-45-3	95
2		Tricosane	324	C23H48	000638-67-5	94
3		Tricosane	324	C23H48	000638-67-5	93
4		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90
5		Eicosane, 10-methyl-	296	C21H44	054833-23-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
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Misc :
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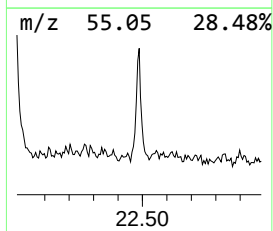
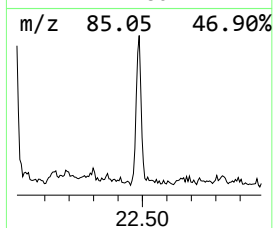
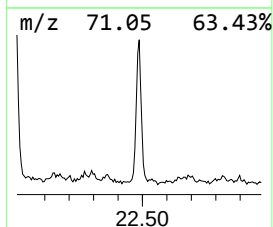
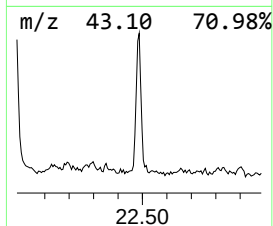
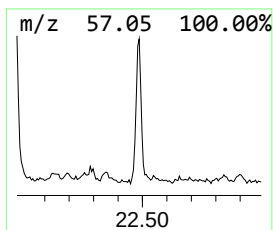
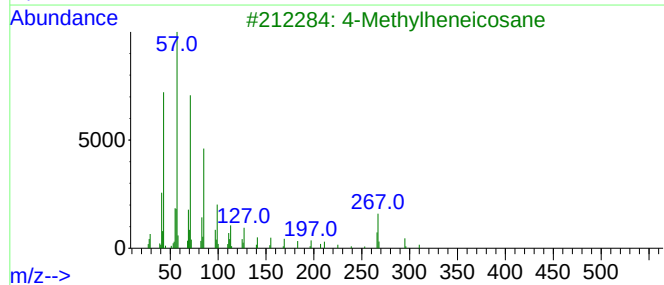
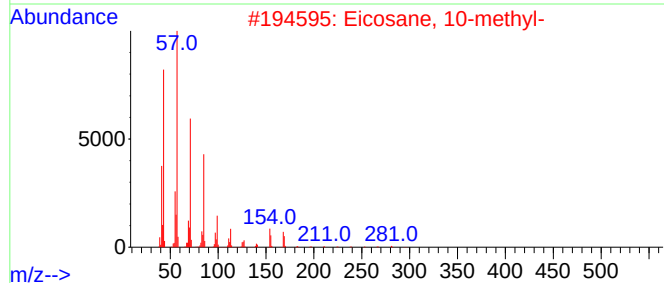
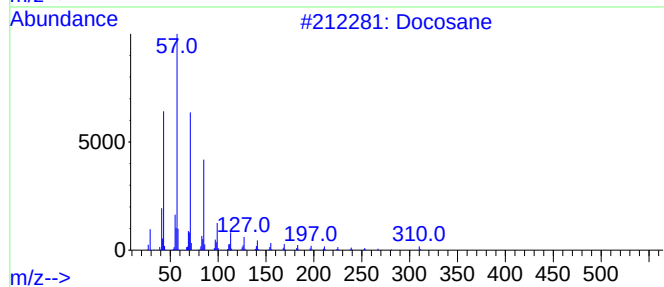
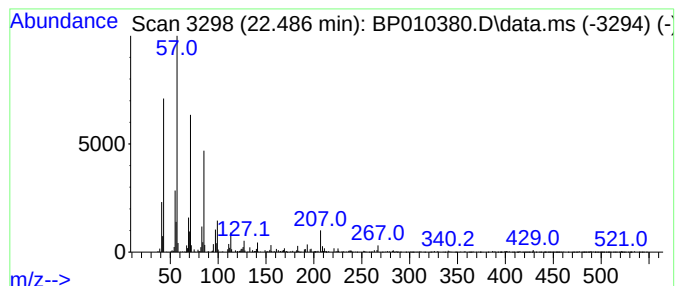
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
22.486	2.12 ng/ul	300931	Chrysene-d12		21.363	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Docosane		310	C22H46	000629-97-0	93
2	Eicosane, 10-methyl-		296	C21H44	054833-23-7	91
3	4-Methylheneicosane		310	C22H46	025117-29-7	91
4	Heptadecane, 9-hexyl-		324	C23H48	055124-79-3	91
5	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
Data File : BP010380.D
Acq On : 17 May 2022 04:48
Operator : CG/JU
Sample : N2831-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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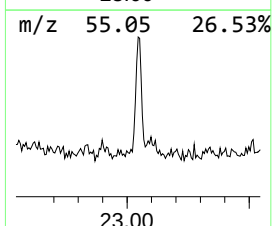
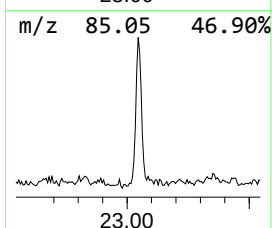
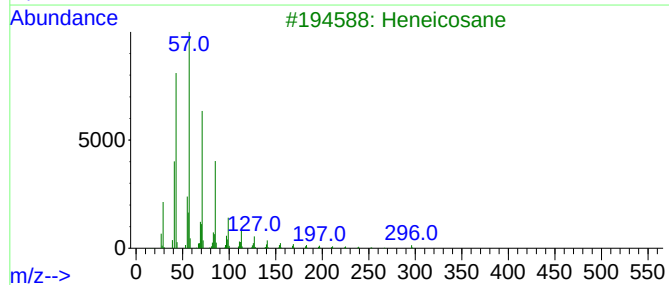
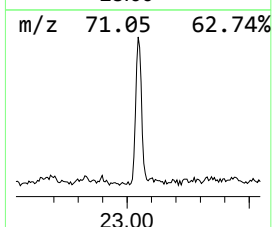
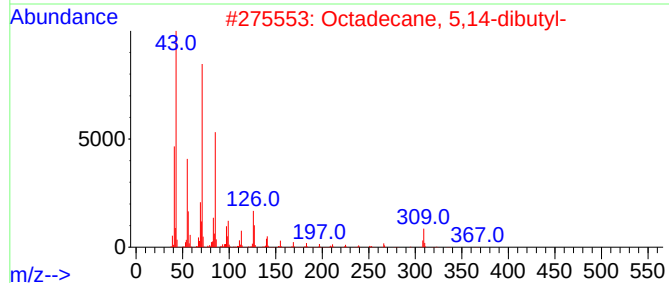
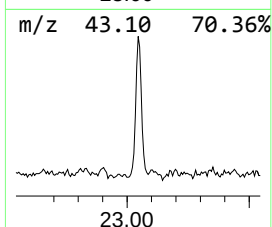
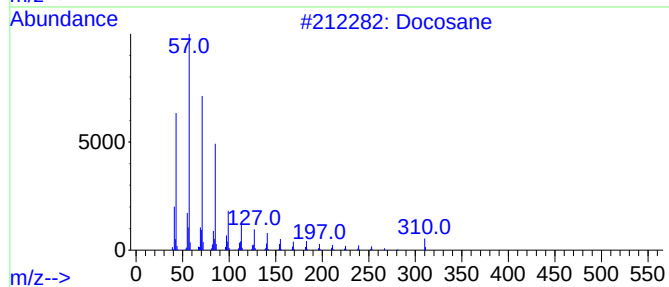
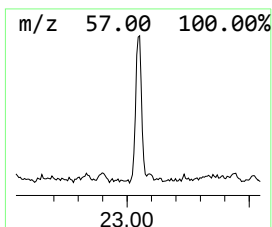
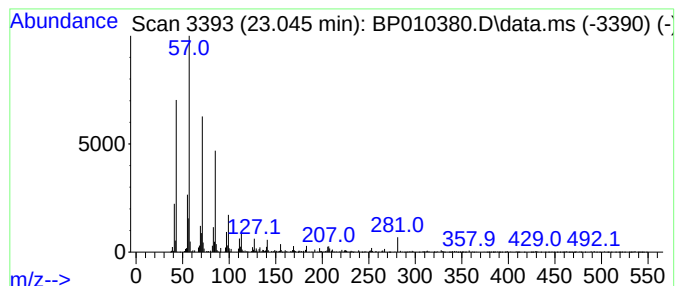
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.045	2.22 ng/ul	287917	Perylene-d12	23.792

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Docosane	310	C22H46	000629-97-0	98
2		Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	93
3		Heneicosane	296	C21H44	000629-94-7	91
4		Tricosane, 2-methyl-	338	C24H50	001928-30-9	91
5		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
Data File : BP010380.D
Acq On : 17 May 2022 04:48
Operator : CG/JU
Sample : N2831-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YB2K3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.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
Crotonic acid	4.540	4.2	ng/ul	197294	1	7.899	930191	20.0
Capro lactam	11.675	97.3	ng/ul	6653720	2	10.699	1367570	20.0
2-Butenoic acid...	12.334	5.4	ng/ul	368937	2	10.699	1367570	20.0
Hexathiane	15.122	2.3	ng/ul	230491	3	14.528	2003370	20.0
7,9-Di-tert-but...	17.951	3.1	ng/ul	374306	4	17.281	2407820	20.0
Dodecyl isobuty...	18.186	8.3	ng/ul	996061	4	17.281	2407820	20.0
3,5-di-tert-But...	18.398	86.5	ng/ul	10420100	4	17.281	2407820	20.0
(DEL) Alkane: S...	21.075	6.2	ng/ul	882003	5	21.363	2839360	20.0
(DEL) Alkane: S...	21.980	2.3	ng/ul	330157	5	21.363	2839360	20.0
(DEL) Alkane: S...	22.486	2.1	ng/ul	300931	5	21.363	2839360	20.0
(DEL) Alkane: S...	23.045	2.2	ng/ul	287917	6	23.792	2590770	20.0