LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023483.D
Acq On : 30 Oct 2019 15:51
Operator : JU
Sample : K5434-10
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : JLNH1

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_M\METHODS\SOM-EPA-BM102319MA.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.799	642	648	660	rVB2	1147980	2040477	7.86%	1.229%
2	6.957	670	675	687	rVB	862945	1383429	5.33%	0.833%
3	7.151	702	708	717	rBV	1172320	1845271	7.11%	1.111%
4	7.616	780	787	797	rBV	1583211	2512089	9.67%	1.513%
5	8.328	901	908	916	rBV	1130805	1816738	7.00%	1.094%
6	8.398	916	920	931	rVB4	51475	121626	0.47%	0.073%
7	8.763	975	982	990	rBV	1074077	1789299	6.89%	1.077%
8	9.234	1058	1062	1071	rVB2	24386	43802	0.17%	0.026%
9	9.487	1098	1105	1112	rBV	839092	1407230	5.42%	0.847%
10	10.022	1188	1196	1208	rBV	1417745	2342028	9.02%	1.410%
11	10.392	1251	1259	1270	rBV	2148513	3580602	13.79%	2.156%
12	10.534	1276	1283	1299	rBV	1049294	1907024	7.34%	1.148%
13	11.992	1525	1531	1536	rBV3	23894	39101	0.15%	0.024%
14	13.680	1812	1818	1823	rBV	2462080	3560563	13.71%	2.144%
15	13.727	1823	1826	1833	rVB	1013353	1358877	5.23%	0.818%
16	13.957	1858	1865	1878	rBV	2897600	4305242	16.58%	2.592%
17	14.263	1910	1917	1928	rBV	3171670	4763885	18.35%	2.868%
18	14.480	1948	1954	1974	rBV	674509	1315196	5.07%	0.792%
19	14.633	1976	1980	1986	rVB8	22122	38725	0.15%	0.023%
20	14.698	1986	1991	1995	rBV5	74914	112795	0.43%	0.068%
21	15.098	2050	2059	2061	rBV7	15467	30893	0.12%	0.019%
22	15.263	2080	2087	2098	rBV	3782915	5384655	20.74%	3.242%
23	15.386	2102	2108	2123	rVV	749722	1135001	4.37%	0.683%
24	15.504	2123	2128	2135	rVB2	43021	67405	0.26%	0.041%
25	16.568	2305	2309	2315	rVB	36699	49258	0.19%	0.030%
26	16.804	2343	2349	2352	rBV4	17836	31917	0.12%	0.019%
27	17.010	2378	2384	2395	rBV2	3788107	5522902	21.27%	3.325%
28	17.110	2395	2401	2413	rVV	4313238	5923489	22.81%	3.566%
29	17.933	2537	2541	2546	rBV2	77654	114172	0.44%	0.069%
30	18.233	2590	2592	2597	rVB6	25254	27347	0.11%	0.016%
31	18.321	2602	2607	2613	rVV6	62528	102404	0.39%	0.062%
32	18.398	2617	2620	2626	rVB7	21172	34553	0.13%	0.021%
33	19.045	2727	2730	2733	rBV	41490	46759	0.18%	0.028%
34	19.298	2770	2773	2779	rVB	42235	55271	0.21%	0.033%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023483.D
Acq On : 30 Oct 2019 15:51
Operator : JU
Sample : K5434-10
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : JLNH1

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_M\METHODS\SOM-EPA-BM102319MA.M
Title : SVOA CALIBRATION

35	19.409	2786	2792	2799	rBV	5202216	7177411	27.64%	4.321%
36	19.998	2888	2892	2896	rVB	200935	220789	0.85%	0.133%
37	20.521	2978	2981	2982	rBV2	200268	237397	0.91%	0.143%
38	20.545	2982	2985	2990	rVB	842137	1009717	3.89%	0.608%
39	20.827	3030	3033	3036	rBV2	169296	217878	0.84%	0.131%
40	20.868	3036	3040	3048	rVB	1370914	1957644	7.54%	1.179%
41	20.956	3051	3055	3060	rVV	1341176	1528839	5.89%	0.921%
42	21.074	3071	3075	3081	rBV3	359632	699001	2.69%	0.421%
43	21.209	3093	3098	3104	rBV	5024844	6389856	24.61%	3.847%
44	21.392	3127	3129	3133	rVB2	204493	197051	0.76%	0.119%
45	21.821	3198	3202	3211	rBV	2198119	2844770	10.96%	1.713%
46	22.280	3277	3280	3283	rBV	290492	345478	1.33%	0.208%
47	22.774	3361	3364	3368	rBV	680320	884193	3.41%	0.532%
48	23.262	3441	3447	3454	rBV	4126338	7482179	28.82%	4.505%
49	23.397	3464	3470	3481	rVB	3981128	7587902	29.22%	4.569%
50	26.791	4036	4047	4055	rBV2	8172924	25965145	100.00%	15.633%
51	26.874	4056	4061	4071	rVB2	1712931	4537608	17.48%	2.732%
52	27.144	4100	4107	4119	rVV5	1466158	4974920	19.16%	2.995%
53	27.297	4123	4133	4153	rVB3	6393504	23026515	88.68%	13.864%
54	29.238	4453	4463	4485	rVB2	3333853	13993059	53.89%	8.425%

Sum of corrected areas: 166087377

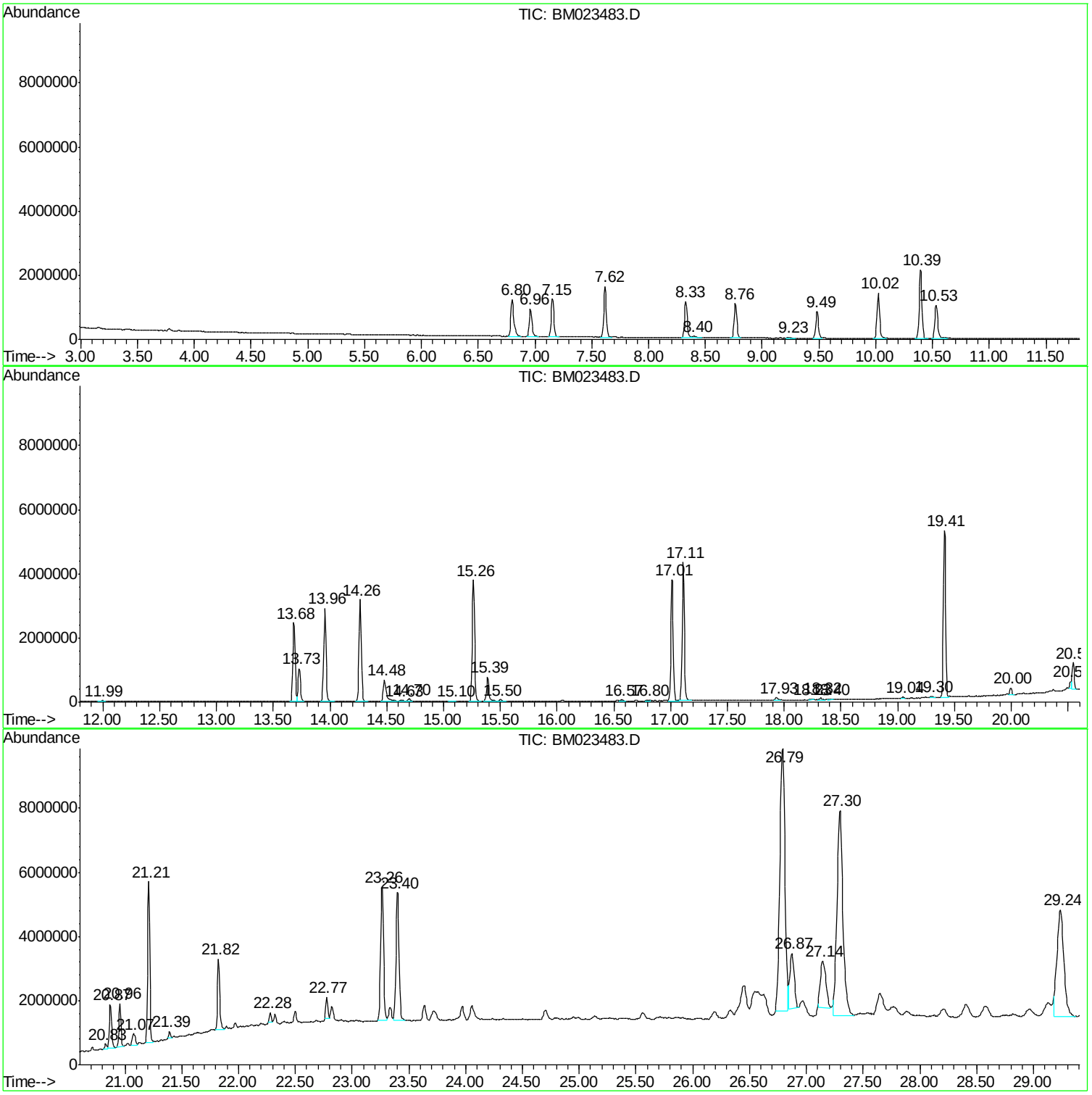
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023483.D
Acq On : 30 Oct 2019 15:51
Operator : JU
Sample : K5434-10
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : JLNH1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023483.D
Acq On : 30 Oct 2019 15:51
Operator : JU
Sample : K5434-10
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : JLNH1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

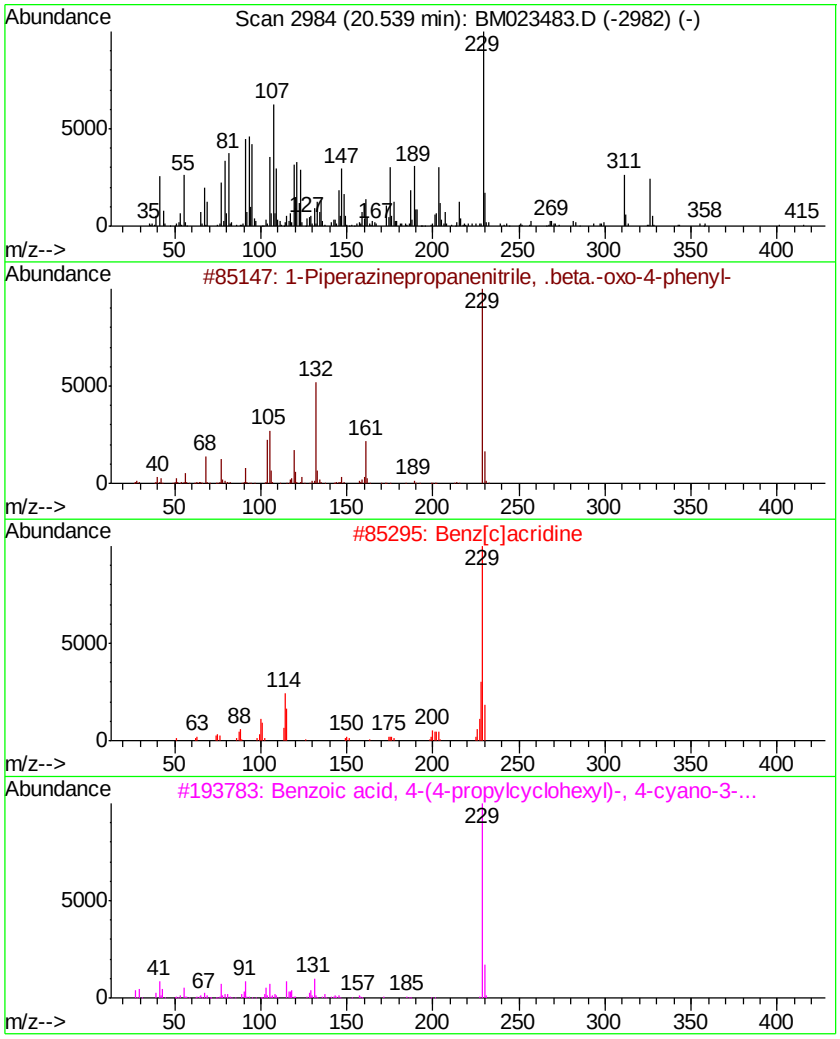
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

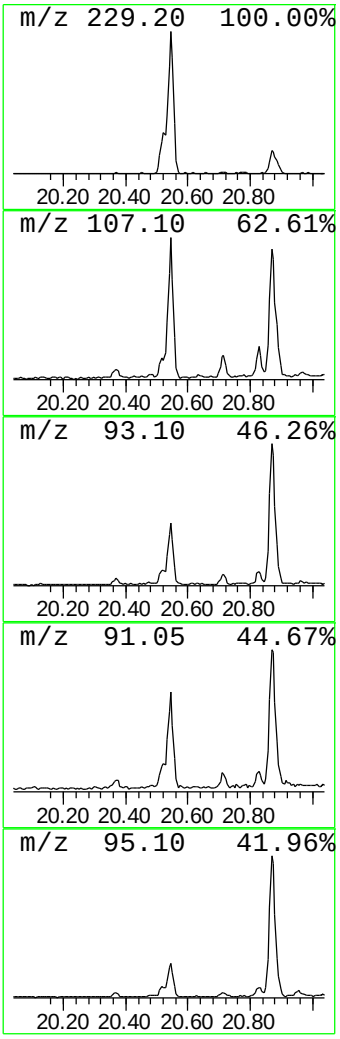
****

Peak Number 1 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.54	3.16 ng/ul	1009720	Chrysene-d12	21.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Piperazinepropanenitrile, .bet...	229	C13H15N3O	014761-40-1	30
2	Benz[c]acridine	229	C17H11N	000225-51-4	22
3	Benzoic acid, 4-(4-propylcyclohe...	365	C23H24FNO2	092118-82-6	22
4	Benzene, 1,3-bis(1-isocyanato-1-...	244	C14H16N2O2	002778-42-9	22
5	Picric acid	229	C6H3N3O7	000088-89-1	14





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023483.D
Acq On : 30 Oct 2019 15:51
Operator : JU
Sample : K5434-10
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : JLNH1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

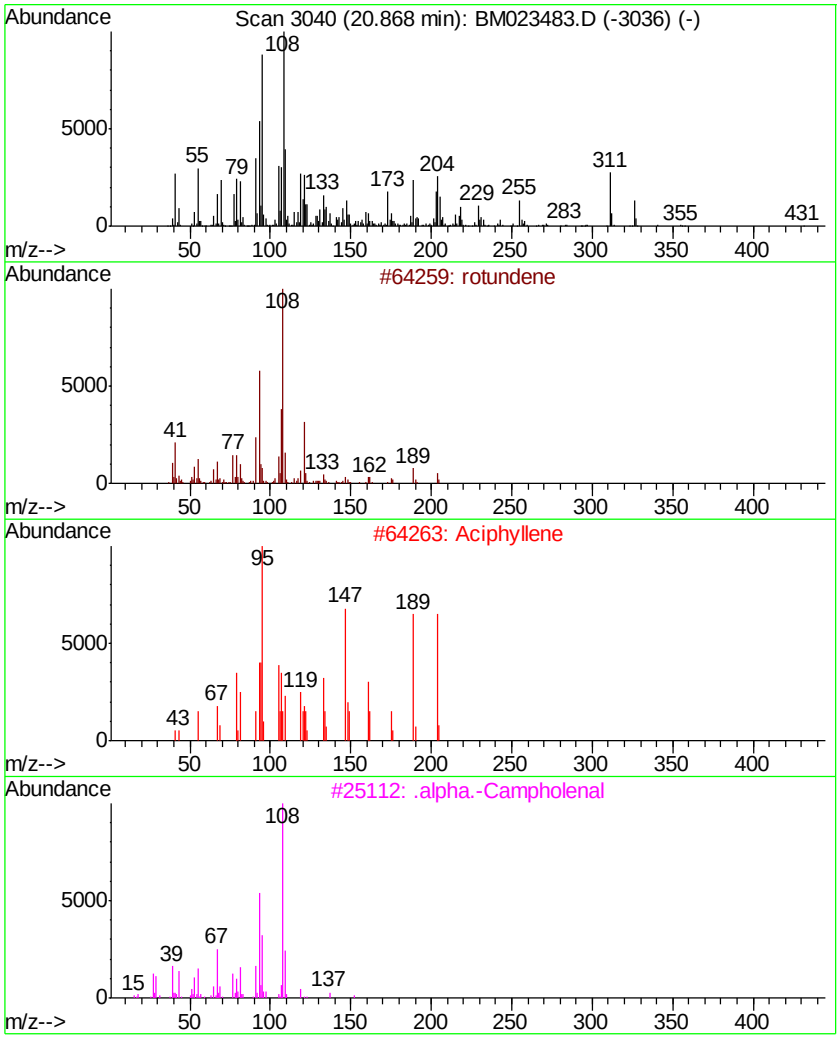
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

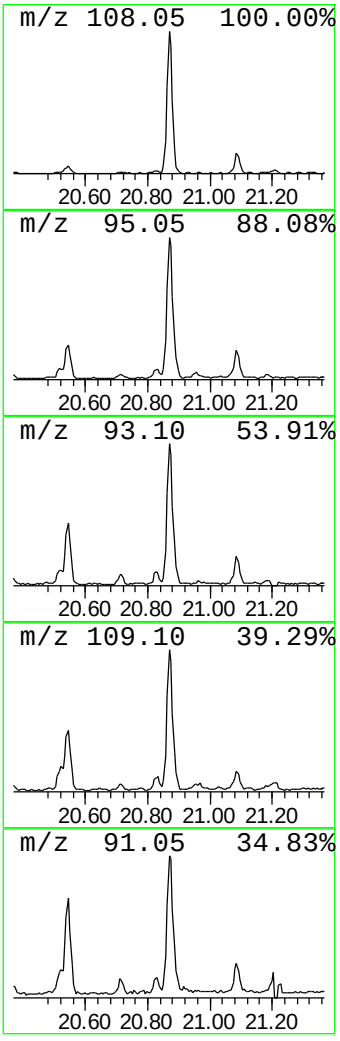
****

Peak Number 2 rotundene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.87	6.13 ng/ul	1957640	Chrysene-d12	21.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	rotundene	204	C15H24	1000374-17-0	55
2	Aciphyllene	204	C15H24	087745-31-1	45
3	.alpha.-Campholenal	152	C10H16O	004501-58-0	38
4	1H-Benzocycloheptene, 2,4a,5,6,7...	204	C15H24	003853-83-6	30
5	(3Ar)-(+)-8,8-dimethyl-4,5,6,7-t...	213	C10H15NO2S	107869-45-4	30





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023483.D
Acq On : 30 Oct 2019 15:51
Operator : JU
Sample : K5434-10
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : JLNH1

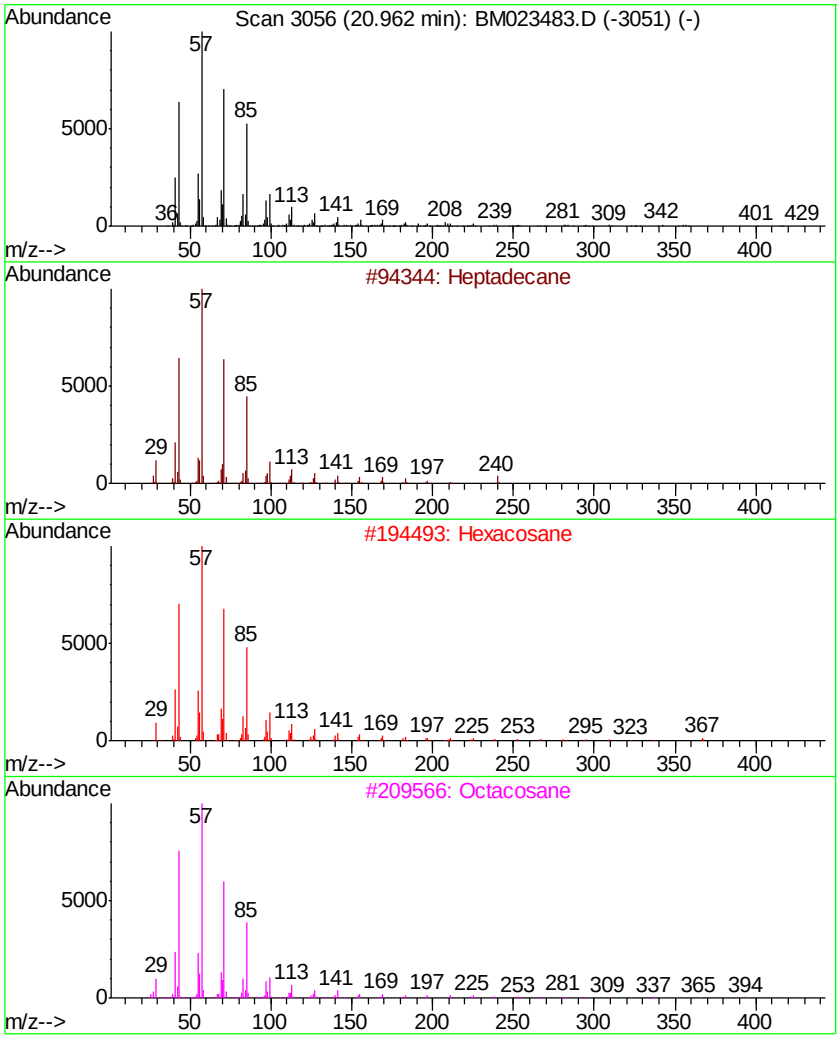
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

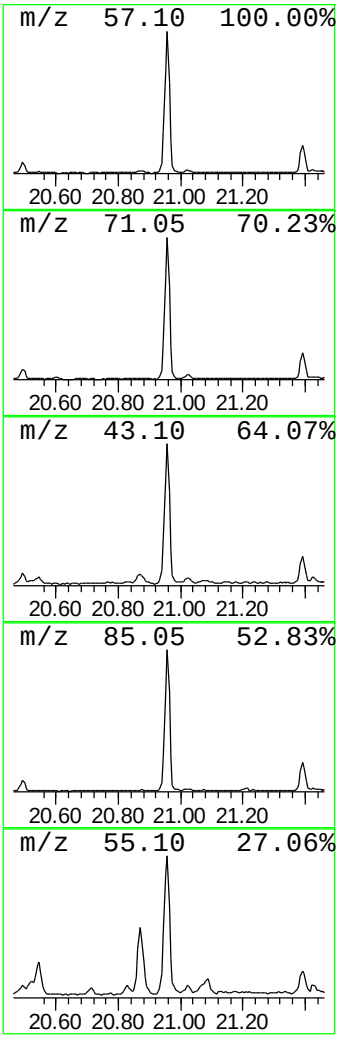
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.96	4.79 ng/ul	1528840	Chrysene-d12	21.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	93
2	Hexacosane	366	C26H54	000630-01-3	91
3	Octacosane	394	C28H58	000630-02-4	91
4	Tetracosane	338	C24H50	000646-31-1	91
5	Octacosane	394	C28H58	000630-02-4	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023483.D
Acq On : 30 Oct 2019 15:51
Operator : JU
Sample : K5434-10
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : JLNH1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

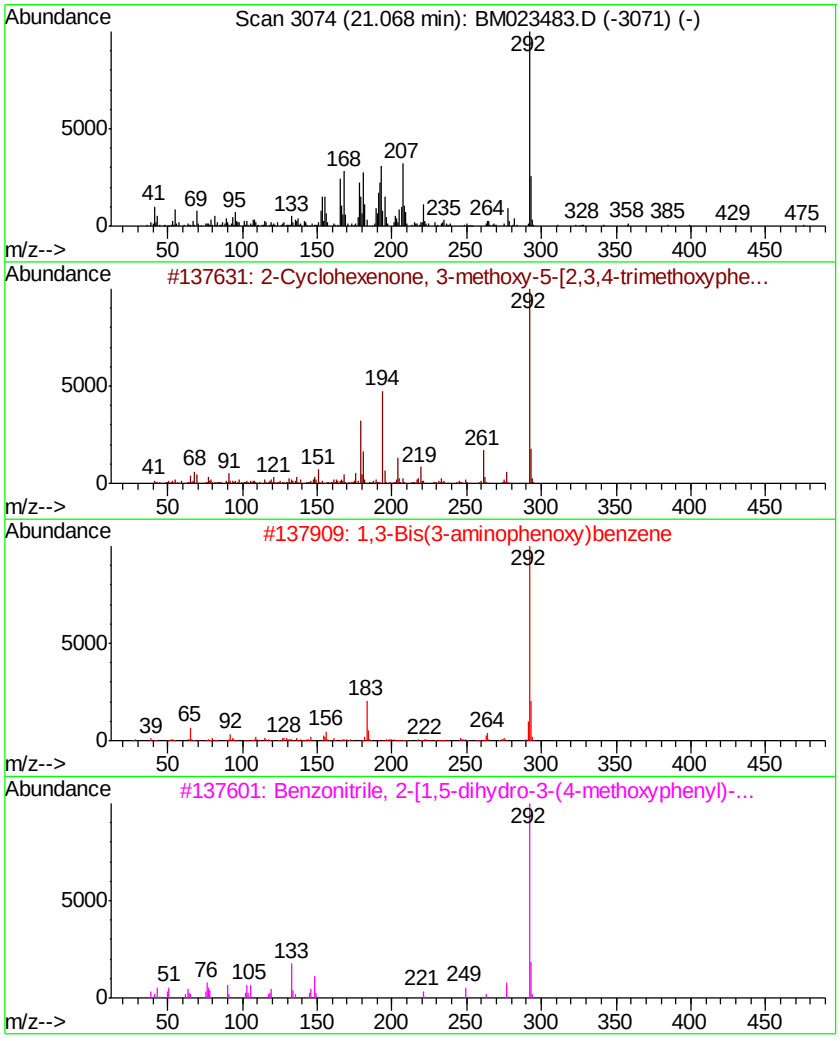
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

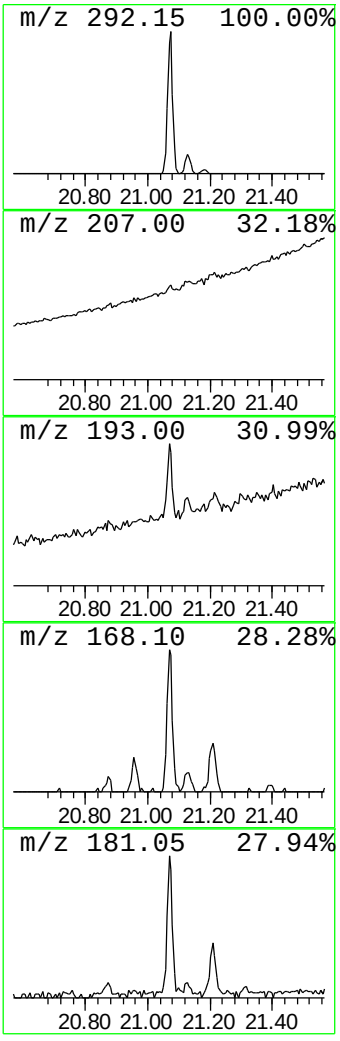
****

Peak Number 4 unknown-02 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.07	2.19 ng/ul	699001	Chrysene-d12	21.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Cyclohexenone, 3-methoxy-5-[2,...	292	C16H20O5	119101-25-6	49
2	1,3-Bis(3-aminophenoxy)benzene	292	C18H16N2O2	010526-07-5	38
3	Benzonitrile, 2-[1,5-dihydro-3-(...	292	C16H12N4O2	1000349-81-5	35
4	Cyclohexano[a]pyrene, 7,8-difluo...	292	C20H14F2	1000116-80-4	30
5	2,4-Diamino-5-[3,4,5-trimethoxyp...	292	C13H16N4O4	034594-35-9	30





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023483.D
Acq On : 30 Oct 2019 15:51
Operator : JU
Sample : K5434-10
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : JLNH1

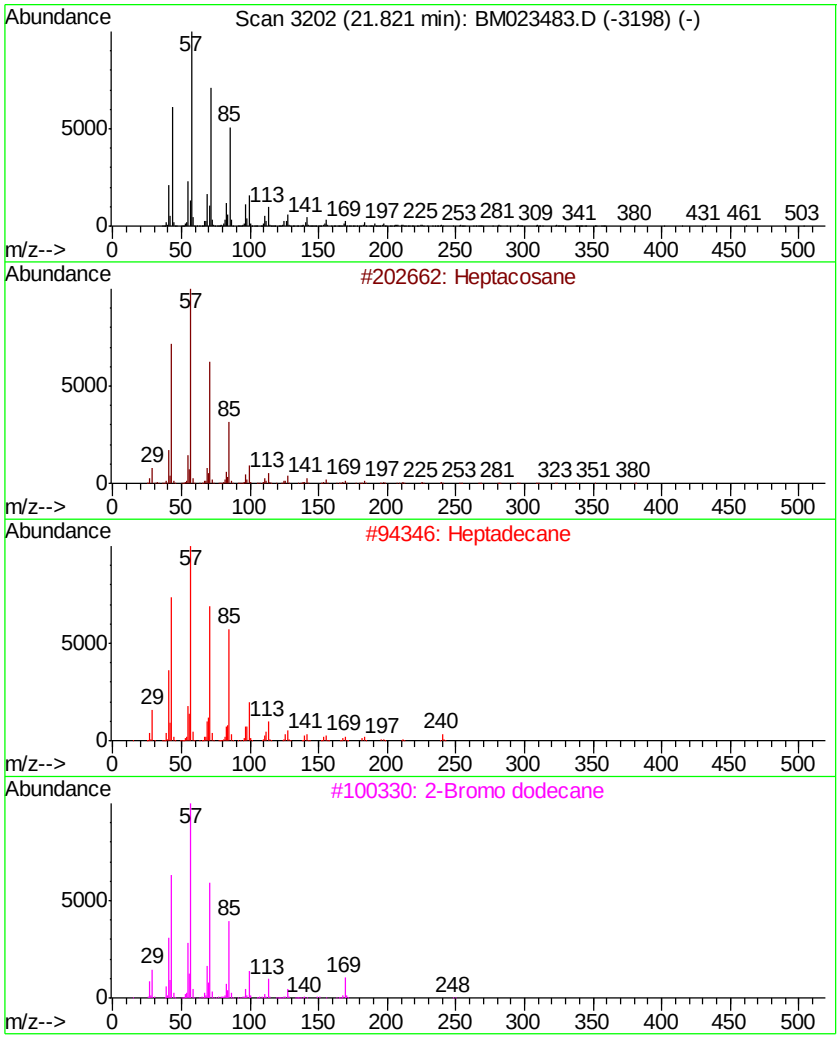
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

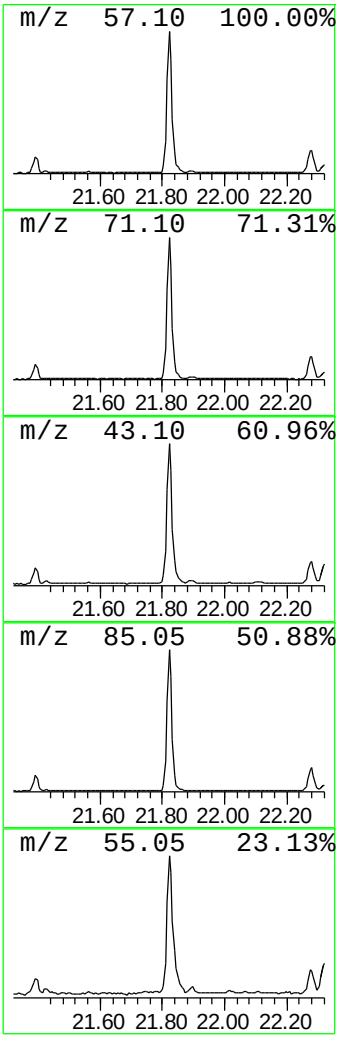
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.82	8.90 ng/ul	2844770	Chrysene-d12	21.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	98
2	Heptadecane	240	C17H36	000629-78-7	93
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	93
4	Heneicosane	296	C21H44	000629-94-7	91
5	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023483.D
Acq On : 30 Oct 2019 15:51
Operator : JU
Sample : K5434-10
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : JLNH1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

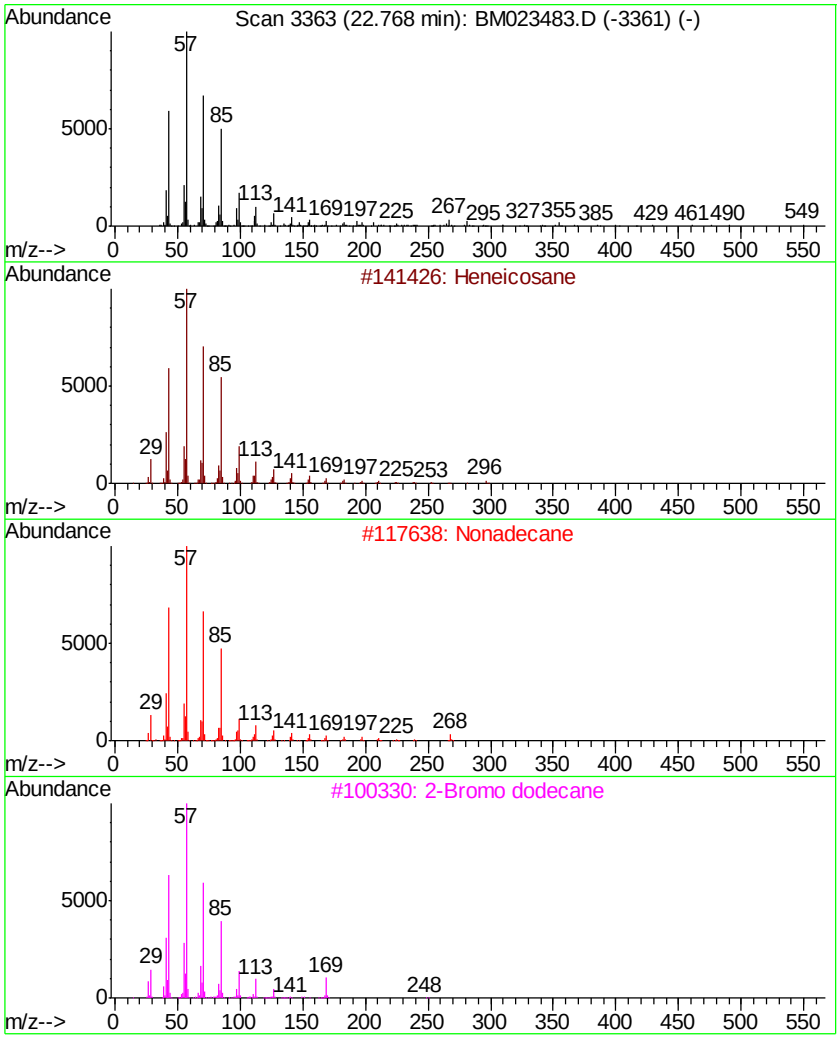
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

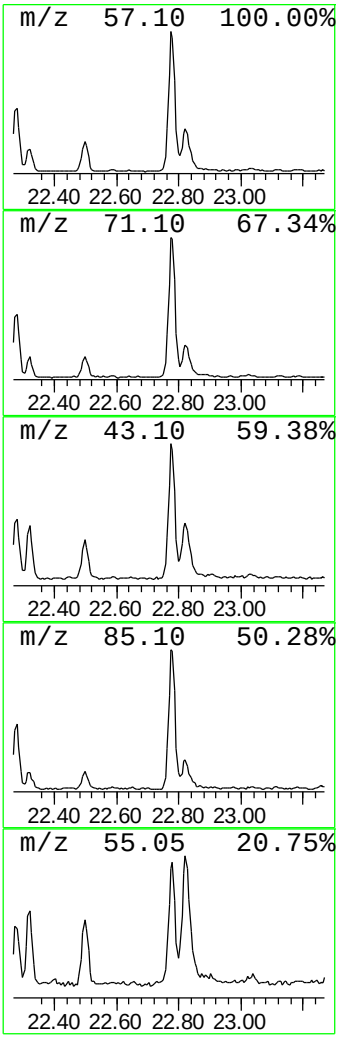
****

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.77	2.33 ng/ul	884193	Perylene-d12	23.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	97
2	Nonadecane	268	C19H40	000629-92-5	93
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	93
4	Hexacosane	366	C26H54	000630-01-3	90
5	Tetracosane	338	C24H50	000646-31-1	90





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023483.D
Acq On : 30 Oct 2019 15:51
Operator : JU
Sample : K5434-10
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : JLNH1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

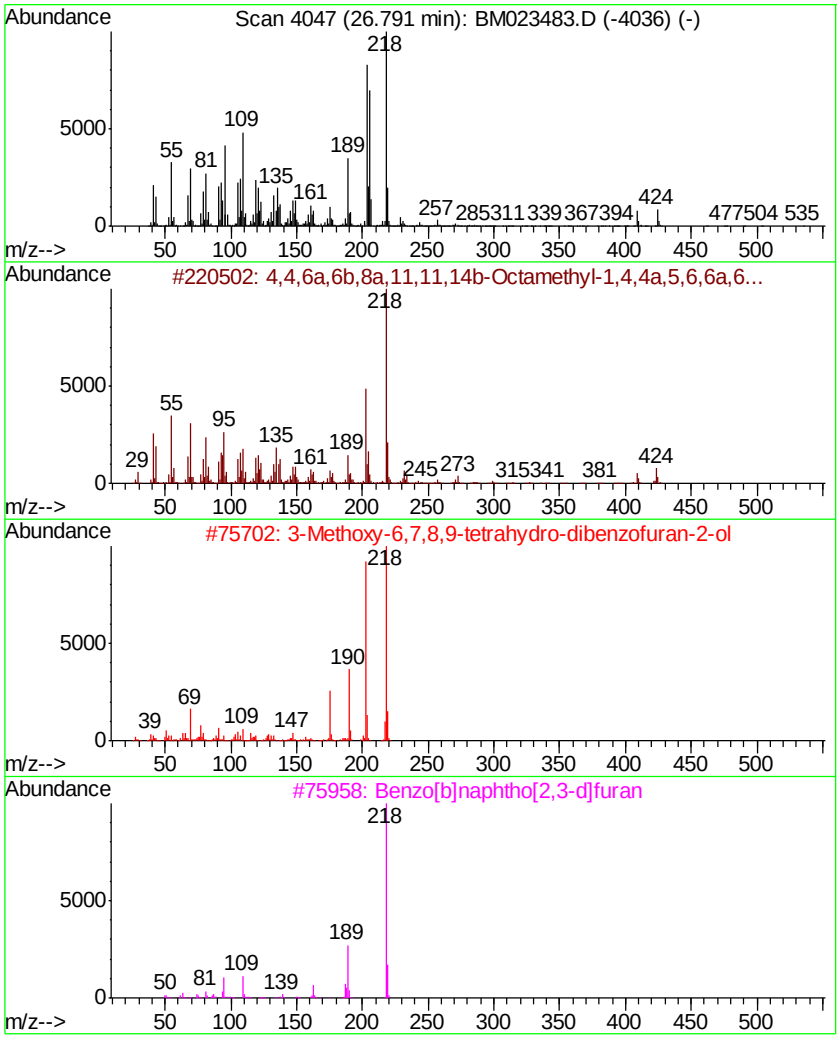
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

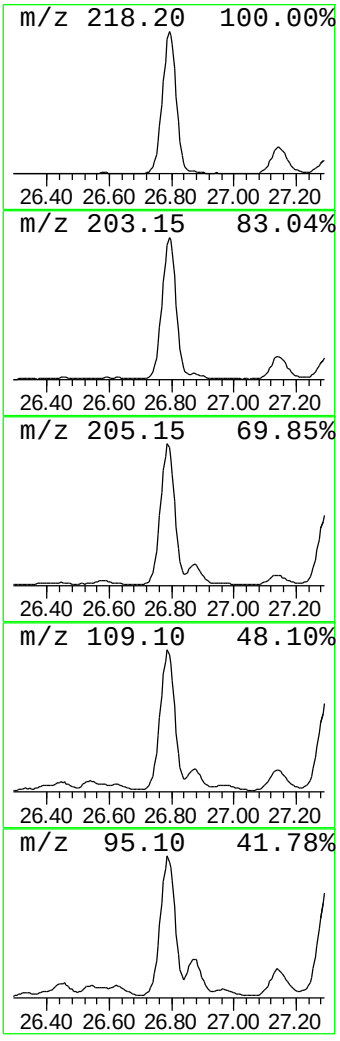
****

Peak Number 7 4,4,6a,6b,8a,11,11,14b-Octa... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.79	68.44 ng/ul	25965100	Perylene-d12	23.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,4,6a,6b,8a,11,11,14b-Octamethy...	424	C30H48O	1000194-62-4	55
2	3-Methoxy-6,7,8,9-tetrahydro-dib...	218	C13H14O3	1000186-57-9	46
3	Benzo[b]naphtho[2,3-d]furan	218	C16H10O	000243-42-5	45
4	4-Allyl-5-pyridin-3-yl-2,4-dihyd...	218	C10H10N4S	1000300-40-5	43
5	Pyrene, hexadecahydro-	218	C16H26	002435-85-0	38





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023483.D
Acq On : 30 Oct 2019 15:51
Operator : JU
Sample : K5434-10
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : JLNH1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

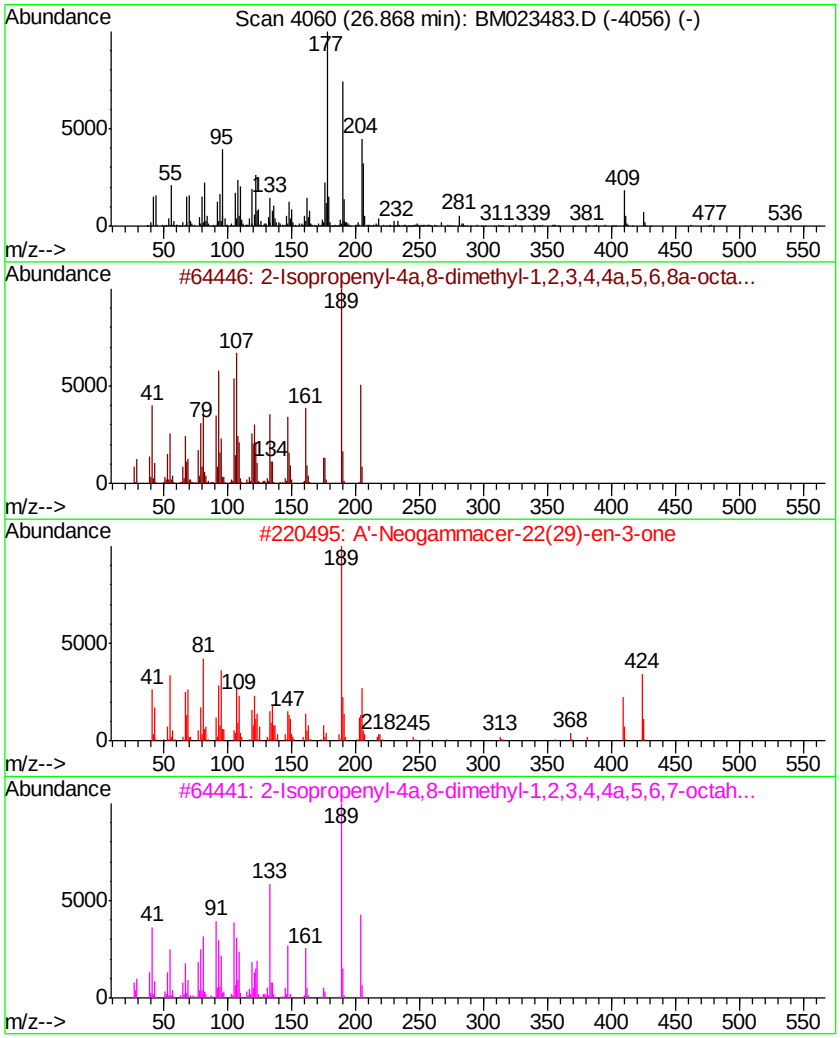
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

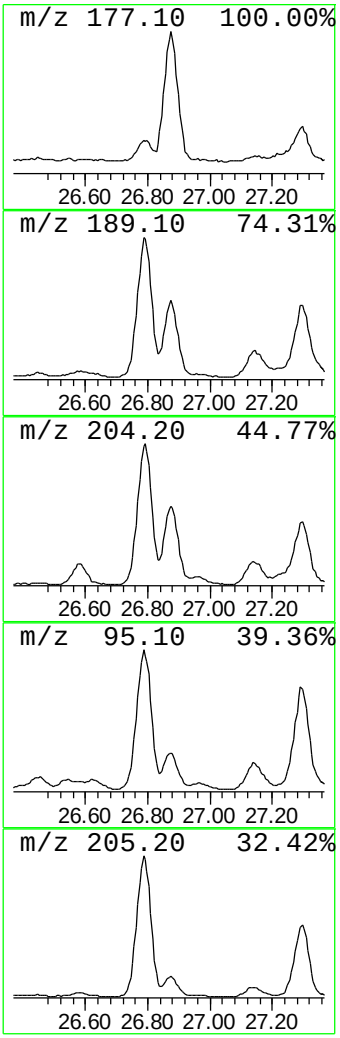
****

Peak Number 8 unknown-03 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.87	11.96 ng/ul	4537610	Perylene-d12	23.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Isopropenyl-4a,8-dimethyl-1,2,...	204	C15H24	1000193-57-0	45
2	A'-Neogammacer-22(29)-en-3-one	424	C30H48O	025615-11-6	38
3	2-Isopropenyl-4a,8-dimethyl-1,2,...	204	C15H24	1000192-43-5	35
4	3,4-Dihydrocoumarin, 4,4,7,8-tet...	204	C13H16O2	040614-36-6	30
5	1-(4,5-Dimethyl-6H-thiopyran-2-y...	204	C9H10F2OS	1000318-25-0	30





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023483.D
Acq On : 30 Oct 2019 15:51
Operator : JU
Sample : K5434-10
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : JLNH1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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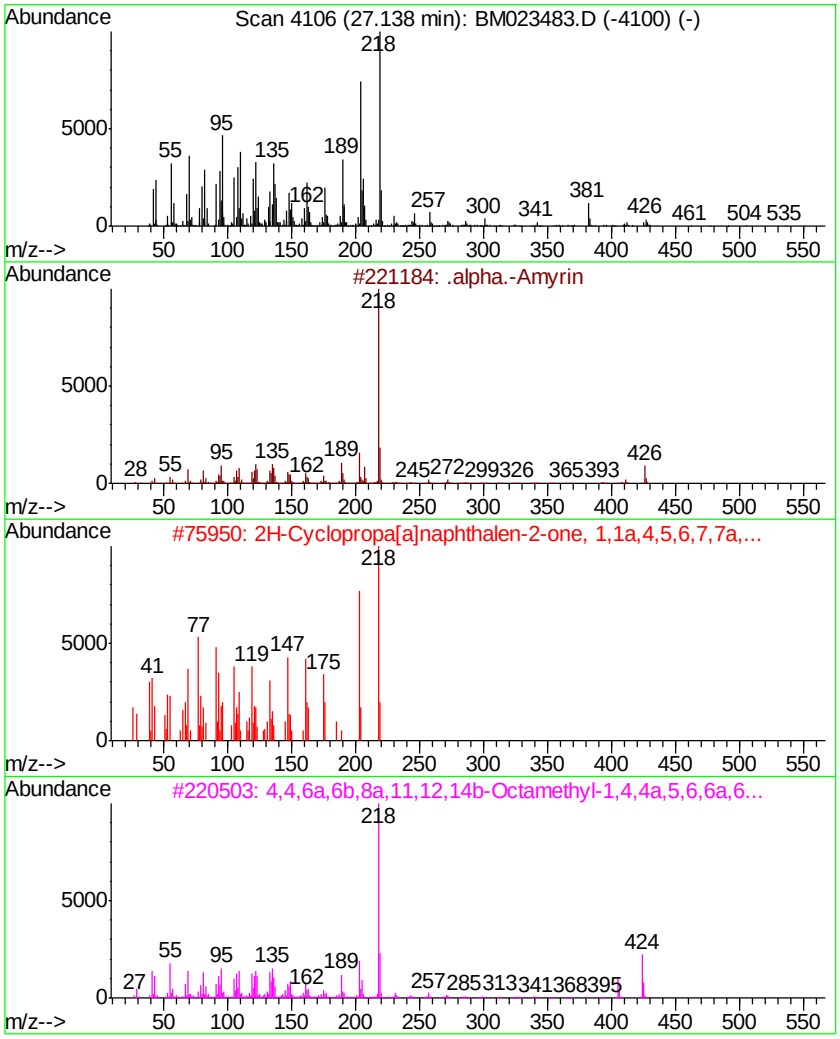
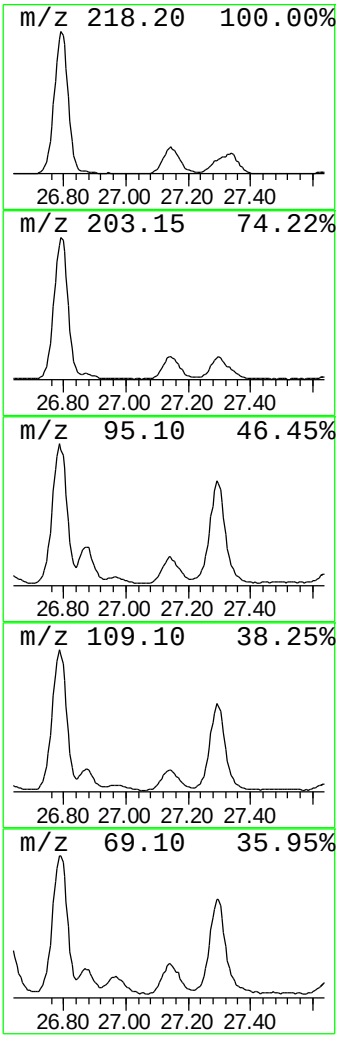
Peak Number 9 .alpha.-Amyrin Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.14	13.11 ng/ul	4974920	Perylene-d12	23.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Amyrin	426	C30H50O	000638-95-9	83
2	2H-Cyclopropa[a]naphthalen-2-one...	218	C15H22O	006831-17-0	70
3	4,4,6a,6b,8a,11,12,14b-Octamethy...	424	C30H48O	1000194-64-2	64
4	.alpha.-Amyrin	426	C30H50O	000638-95-9	49
5	2(1H)Naphthalenone, 3,5,6,7,8,8a...	218	C15H22O	1000188-66-5	45





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023483.D
Acq On : 30 Oct 2019 15:51
Operator : JU
Sample : K5434-10
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : JLNH1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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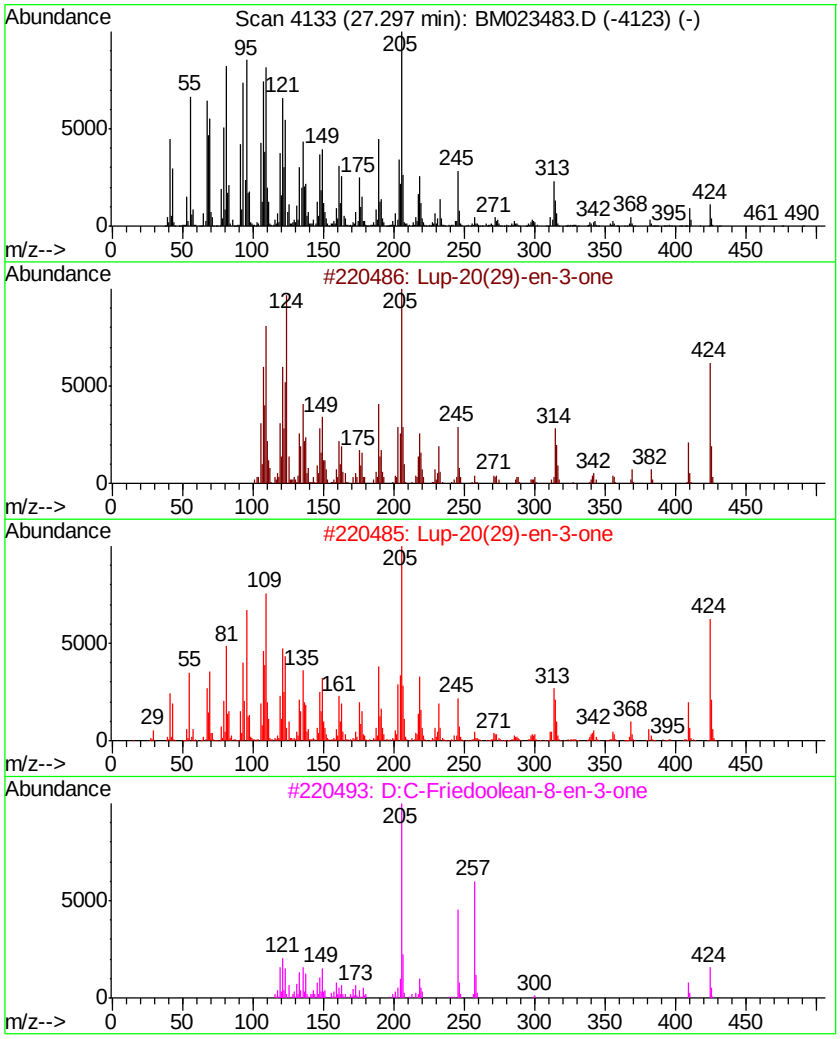
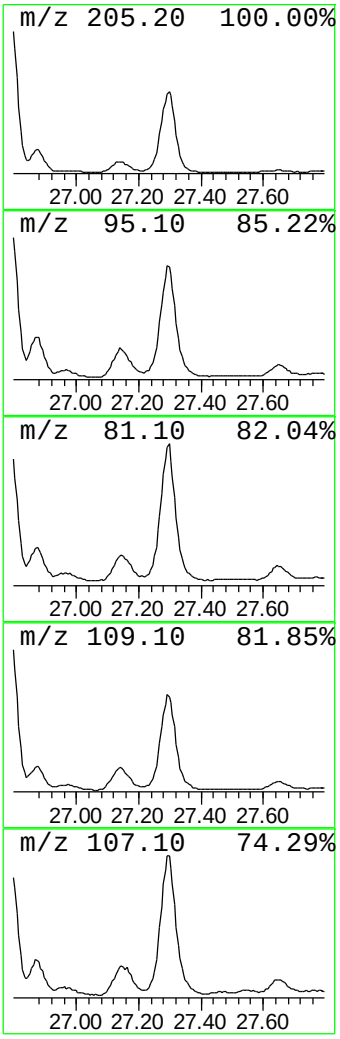
Peak Number 10 Lup-20(29)-en-3-one Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.30	60.69 ng/ul	23026500	Perylene-d12	23.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Lup-20(29)-en-3-one	424	C30H48O	001617-70-5	86
2	Lup-20(29)-en-3-one	424	C30H48O	001617-70-5	64
3	D:C-Friedoolean-8-en-3-one	424	C30H48O	022611-26-3	62
4	Guaia-9,11-diene	204	C15H24	1000374-19-8	52
5	6,10-Dimethyl-3-(1-methylethylid...	206	C15H26	069239-71-0	51





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023483.D
Acq On : 30 Oct 2019 15:51
Operator : JU
Sample : K5434-10
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : JLNH1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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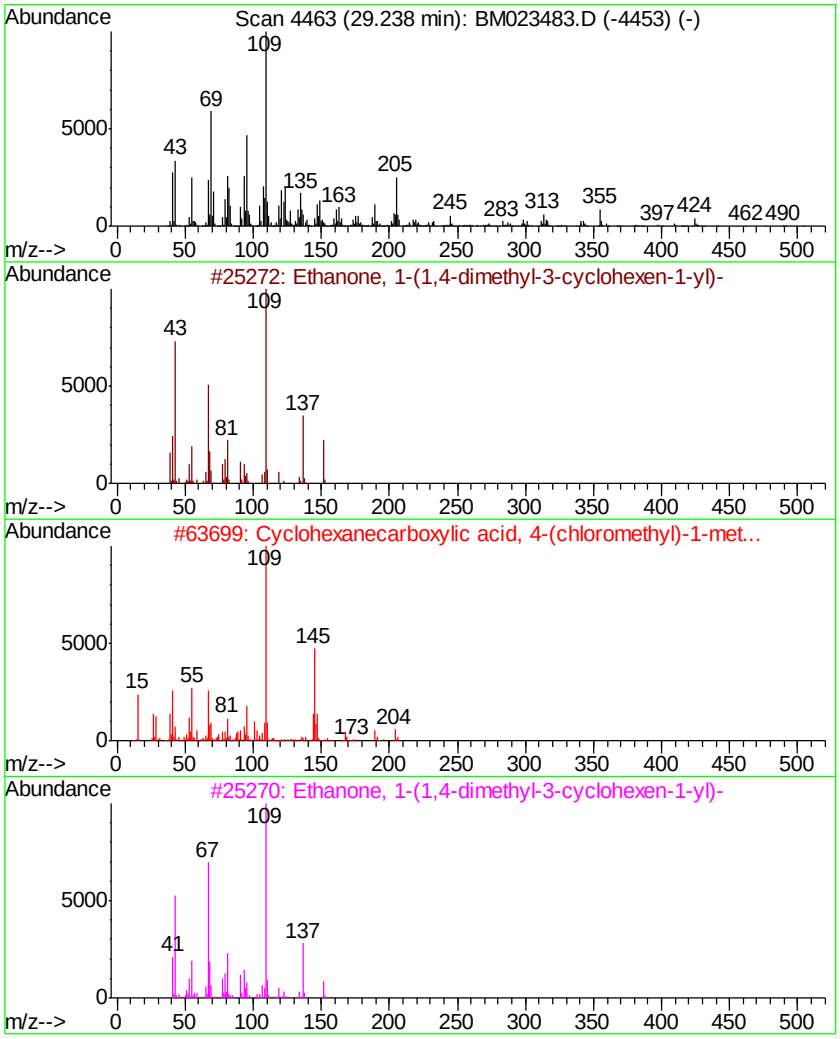
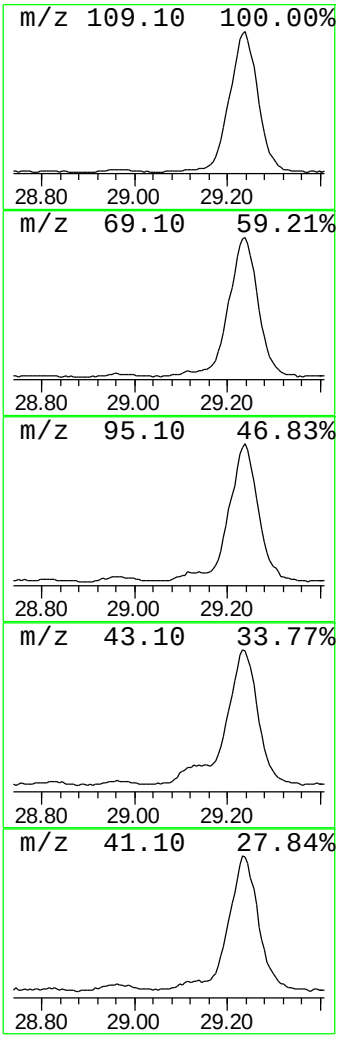
Peak Number 11 unknown-04 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.24	36.88 ng/ul	13993100	Perylene-d12	23.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 1-(1,4-dimethyl-3-cycl...	152	C10H16O	043219-68-7	38
2	Cyclohexanecarboxylic acid, 4-(c...	204	C10H17ClO2	055590-77-7	35
3	Ethanone, 1-(1,4-dimethyl-3-cycl...	152	C10H16O	043219-68-7	35
4	p-Isopropoxyaniline	151	C9H13NO	007664-66-6	30
5	(+)-(Z)-Longipinane	206	C15H26	1000156-12-3	30





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM102919\
Data File : BM023483.D
Acq On : 30 Oct 2019 15:51
Operator : JU
Sample : K5434-10
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : JLNH1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard		
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown-01	20.54	3.2	ng/ul	1009720	5	21.21	6389860	20.0
rotundene	20.87	6.1	ng/ul	1957640	5	21.21	6389860	20.0
(DEL) Alkane: Str...	20.96	4.8	ng/ul	1528840	5	21.21	6389860	20.0
unknown-02	21.07	2.2	ng/ul	699001	5	21.21	6389860	20.0
(DEL) Alkane: Str...	21.82	8.9	ng/ul	2844770	5	21.21	6389860	20.0
(DEL) Alkane: Str...	22.77	2.3	ng/ul	884193	6	23.40	7587900	20.0
4,4,6a,6b,8a,11,1...	26.79	68.4	ng/ul	25965100	6	23.40	7587900	20.0
unknown-03	26.87	12.0	ng/ul	4537610	6	23.40	7587900	20.0
.alpha.-Amyrin	27.14	13.1	ng/ul	4974920	6	23.40	7587900	20.0
Lup-20(29)-en-3-one	27.30	60.7	ng/ul	23026500	6	23.40	7587900	20.0
unknown-04	29.24	36.9	ng/ul	13993100	6	23.40	7587900	20.0