LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022120\
Data File : BP001841.D
Acq On : 22 Feb 2020 01:48
Operator : CG/JU
Sample : L1506-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBCZ2

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\SOM-EPA-BP021820MA.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	2.928	3	7	24	rVB	882171	1337276	8.53%	0.779%
2	3.175	45	49	59	rVB	143376	197698	1.26%	0.115%
3	3.446	92	95	102	rBV2	23750	41296	0.26%	0.024%
4	3.817	154	158	167	rVB2	11083	18508	0.12%	0.011%
5	3.905	169	173	178	rVB	14925	21907	0.14%	0.013%
6	4.393	251	256	261	rVB4	12346	22667	0.14%	0.013%
7	4.799	320	325	330	rBV	29826	44327	0.28%	0.026%
8	6.369	582	592	598	rVB7	9984	24003	0.15%	0.014%
9	6.834	663	671	689	rVB	2474895	4035832	25.75%	2.350%
10	6.993	689	698	712	rBV	2057122	3159690	20.16%	1.840%
11	7.193	723	732	743	rBV	2693117	4303777	27.46%	2.506%
12	7.458	773	777	783	rBV	16296	29072	0.19%	0.017%
13	7.652	800	810	820	rBV	1555656	2464450	15.72%	1.435%
14	7.734	820	824	829	rVB	19833	27679	0.18%	0.016%
15	7.887	844	850	856	rBV	75496	119455	0.76%	0.070%
16	8.357	918	930	942	rBV	3086697	5075280	32.38%	2.956%
17	8.799	997	1005	1018	rBV	2400571	3959163	25.26%	2.306%
18	8.981	1028	1036	1042	rVB6	9471	18777	0.12%	0.011%
19	9.287	1083	1088	1095	rVB	82621	132059	0.84%	0.077%
20	9.516	1118	1127	1136	rBV	2078929	3382542	21.58%	1.970%
21	10.052	1209	1218	1236	rBV	3411173	5636168	35.96%	3.282%
22	10.428	1273	1282	1290	rBV	2199260	3640105	23.22%	2.120%
23	10.557	1297	1304	1320	rBV	3121756	5380182	34.32%	3.133%
24	10.840	1350	1352	1360	rVB7	10131	18510	0.12%	0.011%
25	12.022	1546	1553	1559	rVB2	61138	102850	0.66%	0.060%
26	12.669	1655	1663	1674	rVB7	21560	46012	0.29%	0.027%
27	13.187	1747	1751	1752	rBV	50832	56671	0.36%	0.033%
28	13.704	1832	1839	1844	rBV	5536395	8057469	51.40%	4.692%
29	13.751	1844	1847	1855	rVB	777944	994621	6.35%	0.579%
30	13.981	1878	1886	1895	rBV	6703997	10113739	64.52%	5.890%
31	14.292	1931	1939	1946	rVB	3321734	5081648	32.42%	2.959%
32	14.381	1950	1954	1961	rVB10	12267	21862	0.14%	0.013%
33	14.492	1963	1973	1984	rBV	2718313	4028005	25.70%	2.346%
34	14.569	1984	1986	1994	rVB8	16605	32579	0.21%	0.019%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022120\
Data File : BP001841.D
Acq On : 22 Feb 2020 01:48
Operator : CG/JU
Sample : L1506-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBCZ2

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\SOM-EPA-BP021820MA.M
Title : SVOA CALIBRATION

35	14.716	2006	2011	2013	rBV3	18502	26271	0.17%	0.015%
36	15.128	2077	2081	2085	rVB2	17201	22852	0.15%	0.013%
37	15.187	2085	2091	2099	rBV3	17751	34786	0.22%	0.020%
38	15.287	2101	2108	2117	rBV	8254761	12460913	79.50%	7.257%
39	15.404	2122	2128	2144	rBV	2524109	3500163	22.33%	2.038%
40	15.528	2144	2149	2156	rVB2	125030	187840	1.20%	0.109%
41	15.675	2169	2174	2181	rVB	700494	894612	5.71%	0.521%
42	15.863	2201	2206	2209	rBV5	11136	23794	0.15%	0.014%
43	15.957	2216	2222	2225	rBV7	10099	20612	0.13%	0.012%
44	16.075	2238	2242	2248	rVB2	36619	53614	0.34%	0.031%
45	16.339	2285	2287	2295	rVV8	9991	19548	0.12%	0.011%
46	16.475	2305	2310	2319	rBV2	80271	135897	0.87%	0.079%
47	16.722	2349	2352	2357	rVB3	24061	31881	0.20%	0.019%
48	16.828	2365	2370	2380	rBV2	24991	73624	0.47%	0.043%
49	17.039	2400	2406	2414	rVV	4464648	6323837	40.34%	3.683%
50	17.139	2416	2423	2434	rVV2	9146850	13004219	82.96%	7.573%
51	17.245	2437	2441	2450	rBV	84109	122517	0.78%	0.071%
52	17.404	2465	2468	2473	rBV3	10830	17249	0.11%	0.010%
53	17.839	2537	2542	2549	rBV	214393	361881	2.31%	0.211%
54	17.904	2550	2553	2558	rVV2	71852	122351	0.78%	0.071%
55	17.969	2559	2564	2586	rVV	1849323	2888540	18.43%	1.682%
56	18.251	2608	2612	2618	rBV	53913	84137	0.54%	0.049%
57	18.798	2702	2705	2711	rVB5	49424	61131	0.39%	0.036%
58	18.922	2722	2726	2730	rBV2	109097	134376	0.86%	0.078%
59	19.028	2739	2744	2748	rBV	145119	208900	1.33%	0.122%
60	19.086	2750	2754	2766	rVV2	108519	262673	1.68%	0.153%
61	19.204	2770	2774	2781	rVV	223125	317212	2.02%	0.185%
62	19.275	2782	2786	2795	rVB2	129234	262115	1.67%	0.153%
63	19.386	2799	2805	2812	rVB	11815384	15674613	100.00%	9.128%
64	19.910	2891	2894	2898	rBV	77568	105051	0.67%	0.061%
65	20.010	2908	2911	2915	rVB4	43642	49612	0.32%	0.029%
66	20.086	2921	2924	2936	rBV7	53028	132292	0.84%	0.077%
67	20.557	3001	3004	3011	rBV6	78793	122614	0.78%	0.071%
68	20.874	3053	3058	3073	rBV	2562942	3333456	21.27%	1.941%
69	21.133	3096	3102	3111	rBV	5781739	7360483	46.96%	4.287%
70	21.245	3118	3121	3129	rBV	272550	385650	2.46%	0.225%
71	21.310	3129	3132	3141	rVB	179139	232615	1.48%	0.135%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022120\
Data File : BP001841.D
Acq On : 22 Feb 2020 01:48
Operator : CG/JU
Sample : L1506-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBCZ2

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\SOM-EPA-BP021820MA.M
Title : SVOA CALIBRATION

72	21.751	3202	3207	3224	rBV2	325988	596297	3.80%	0.347%
73	22.192	3277	3282	3289	rBV3	471762	1153198	7.36%	0.672%
74	22.333	3301	3306	3312	rVB	2743974	3792085	24.19%	2.208%
75	22.716	3367	3371	3385	rVB	493722	852010	5.44%	0.496%
76	23.210	3447	3455	3463	rBV	8332066	15487620	98.81%	9.020%
77	23.286	3464	3468	3472	rVV	420697	694290	4.43%	0.404%
78	23.345	3472	3478	3490	rVB	3695971	6950797	44.34%	4.048%
79	23.939	3574	3579	3586	rVV	404960	746400	4.76%	0.435%
80	24.010	3586	3591	3605	rVB	348888	758887	4.84%	0.442%

Sum of corrected areas: 171711394

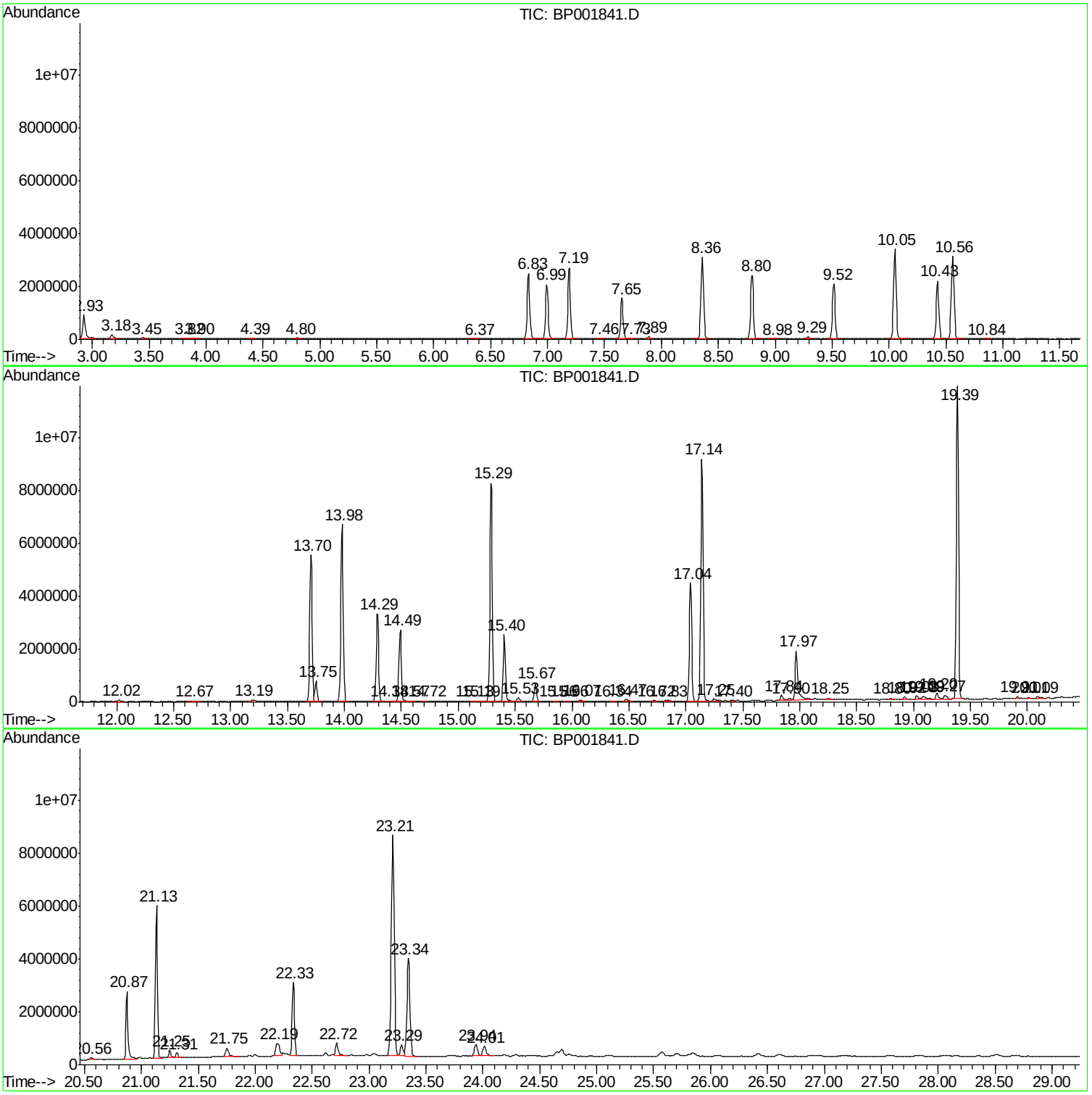
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022120\
Data File : BP001841.D
Acq On : 22 Feb 2020 01:48
Operator : CG/JU
Sample : L1506-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBCZ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.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 P\DATA\BP022120\
Data File : BP001841.D
Acq On : 22 Feb 2020 01:48
Operator : CG/JU
Sample : L1506-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBCZ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

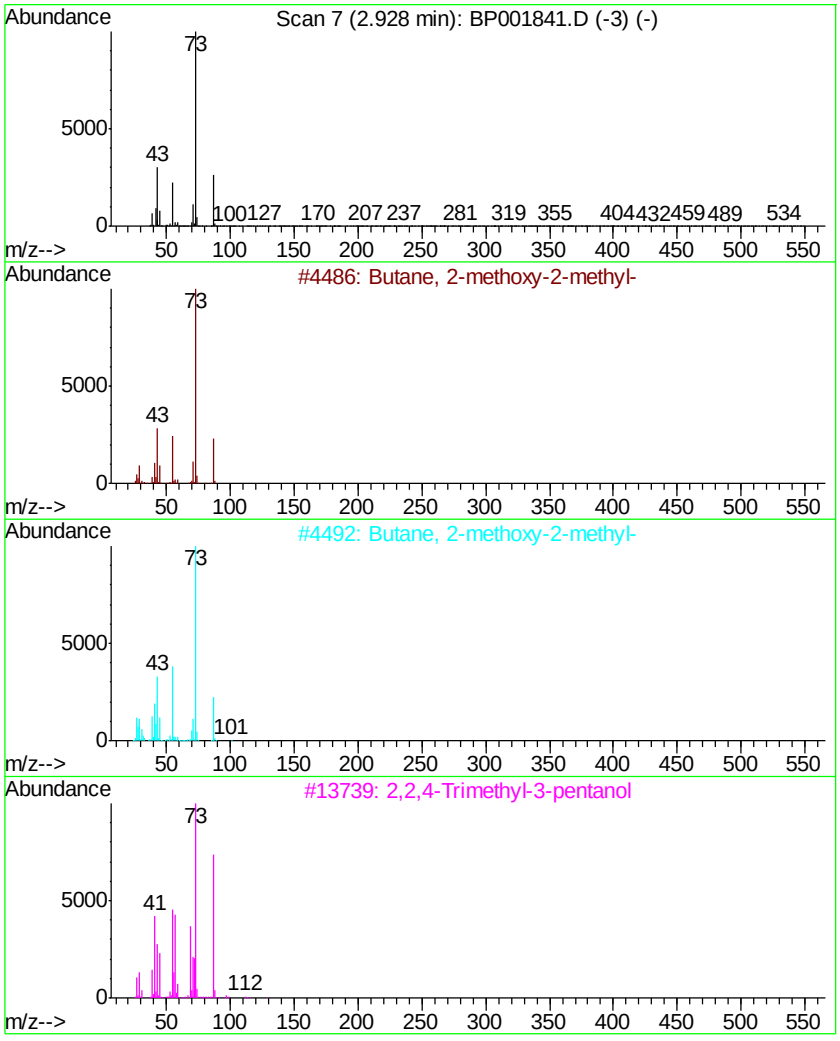
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

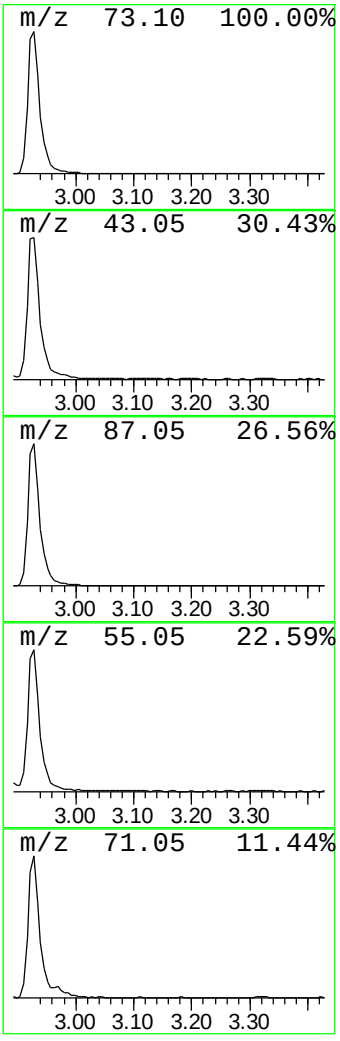
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.93	10.85 ng/ul	1337280	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	90
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022120\
Data File : BP001841.D
Acq On : 22 Feb 2020 01:48
Operator : CG/JU
Sample : L1506-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBCZ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

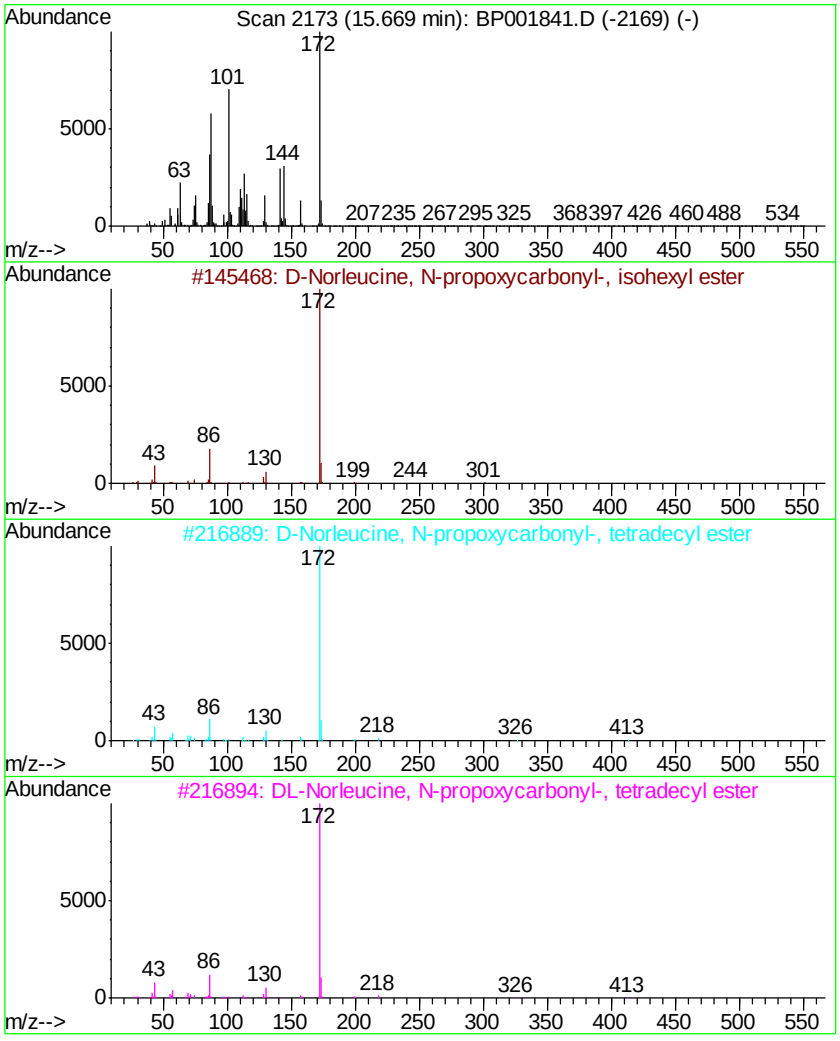
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

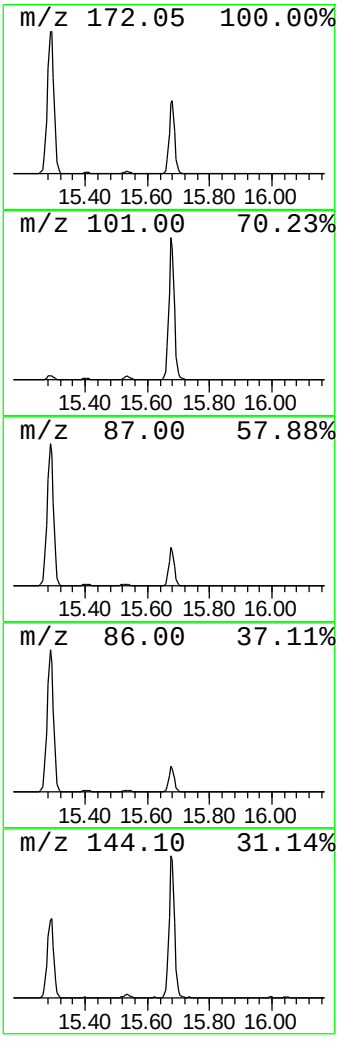
****

Peak Number 2 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	2.83 ng/ul	894612	Phenanthrene-d10	17.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	D-Norleucine, N-propoxycarbonyl-...	301	C16H31NO4	1000338-42-9	35
2	D-Norleucine, N-propoxycarbonyl-...	413	C24H47NO4	1000338-43-4	30
3	DL-Norleucine, N-propoxycarbonyl...	413	C24H47NO4	1000339-03-6	30
4	D-Norleucine, N-propoxycarbonyl-...	427	C25H49NO4	1000338-43-5	27
5	2,5-Dimethyl-3-phenylfuran	172	C12H12O	019842-57-0	27





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022120\
Data File : BP001841.D
Acq On : 22 Feb 2020 01:48
Operator : CG/JU
Sample : L1506-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBCZ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

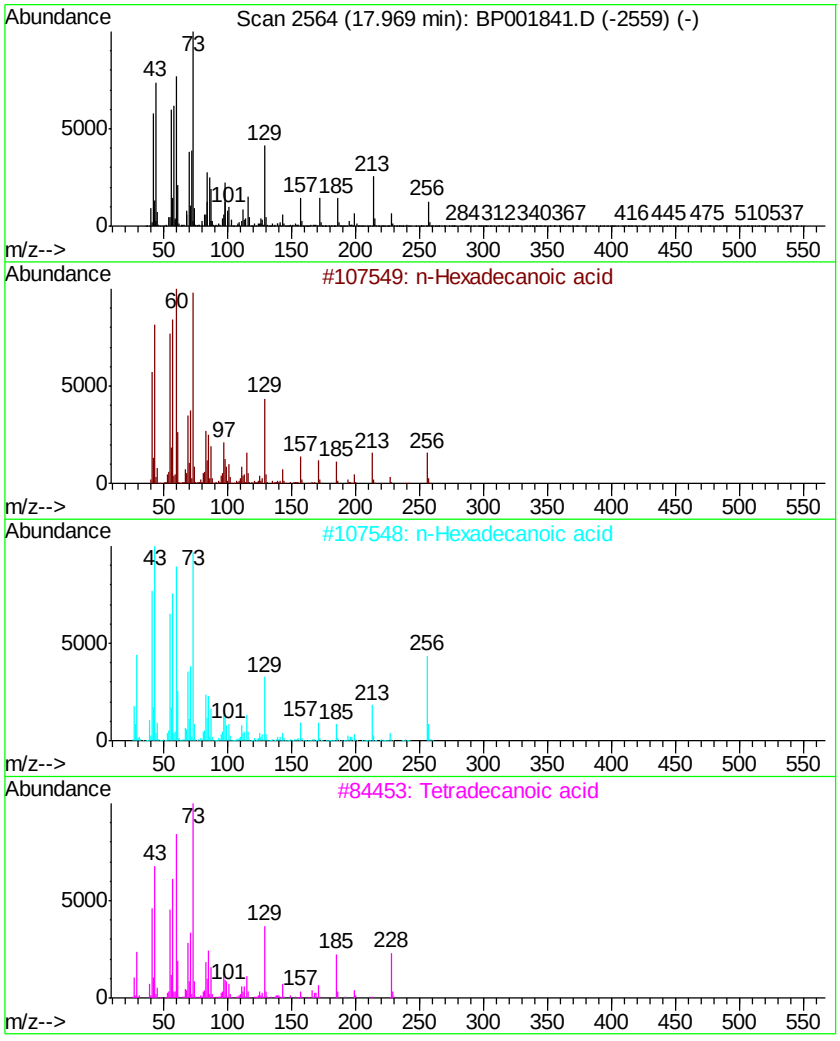
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

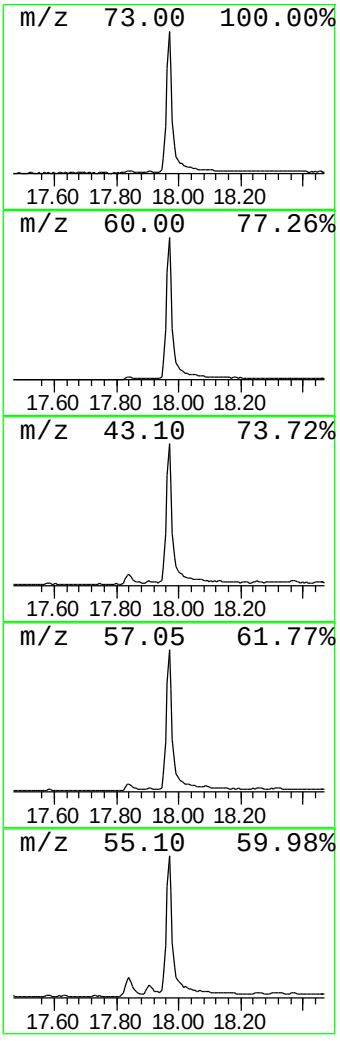
****

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.97	9.14 ng/ul	2888540	Phenanthrene-d10	17.04

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	97
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	93
4	Tridecanoic acid	214	C13H26O2	000638-53-9	91
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	81





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022120\
Data File : BP001841.D
Acq On : 22 Feb 2020 01:48
Operator : CG/JU
Sample : L1506-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBCZ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

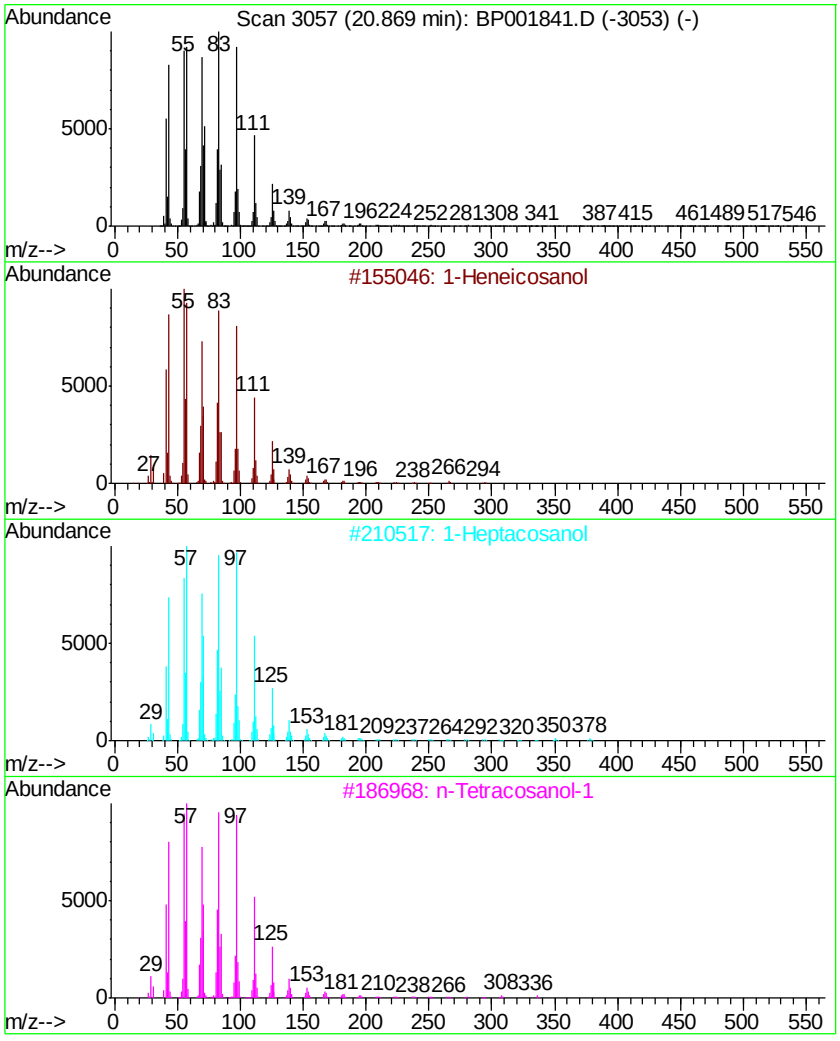
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

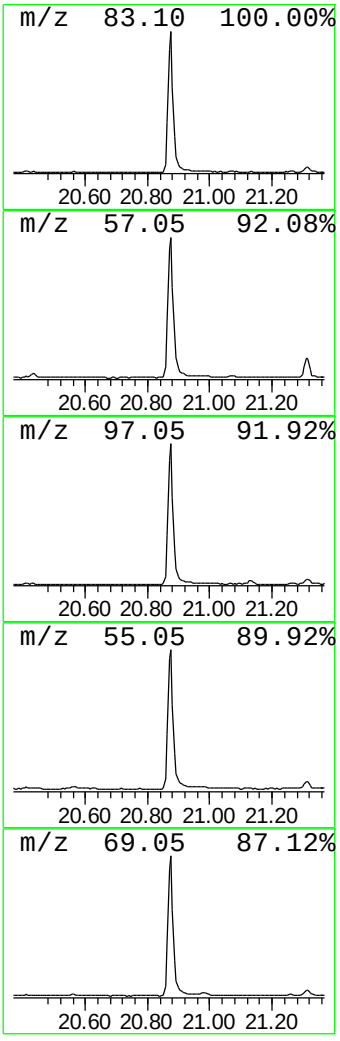
****

Peak Number 4 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.87	9.06 ng/ul	3333460	Chrysene-d12	21.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	1-Heptacosanol	396	C27H56O	002004-39-9	95
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4	Behenic alcohol	326	C22H46O	000661-19-8	94
5	1-Nonadecene	266	C19H38	018435-45-5	94





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022120\
Data File : BP001841.D
Acq On : 22 Feb 2020 01:48
Operator : CG/JU
Sample : L1506-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBCZ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

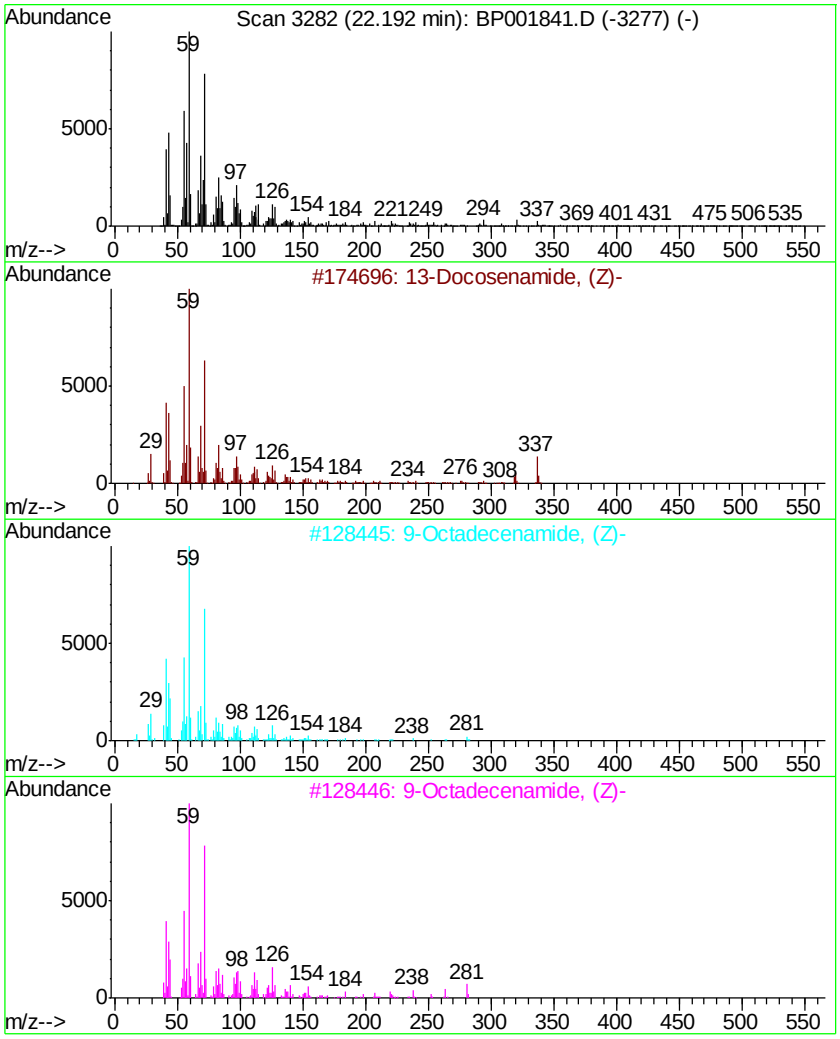
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

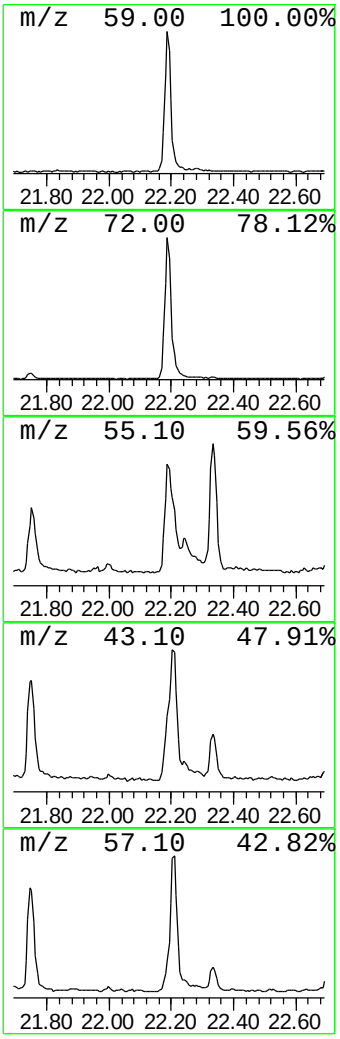
****

Peak Number 5 13-Docosenamide, (Z)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.19	3.13 ng/ul	1153200	Chrysene-d12	21.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	90
2	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	87
3	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	80
4	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	80
5	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	74





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022120\
Data File : BP001841.D
Acq On : 22 Feb 2020 01:48
Operator : CG/JU
Sample : L1506-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBCZ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

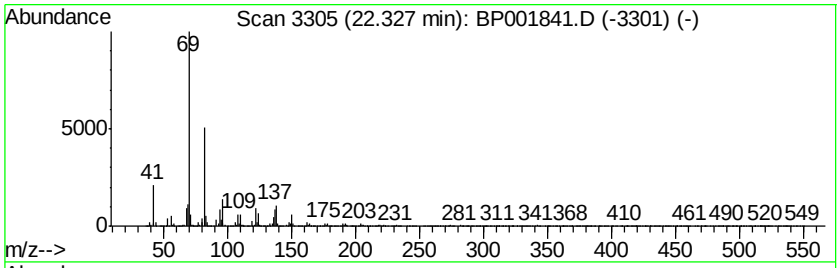
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

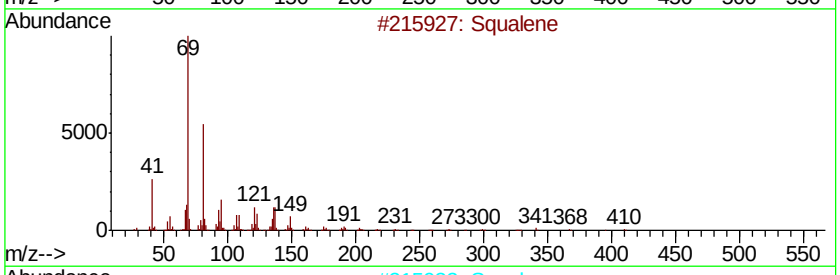
****

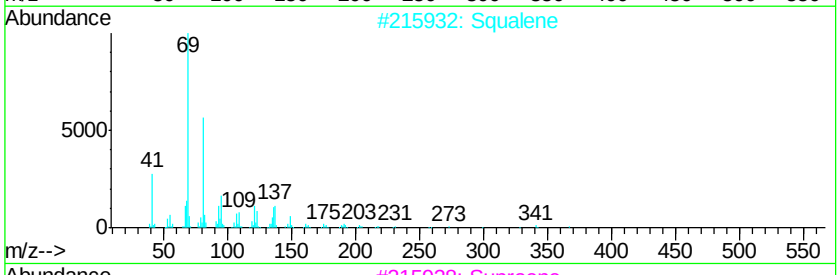
Peak Number 6 Squalene Concentration Rank 1

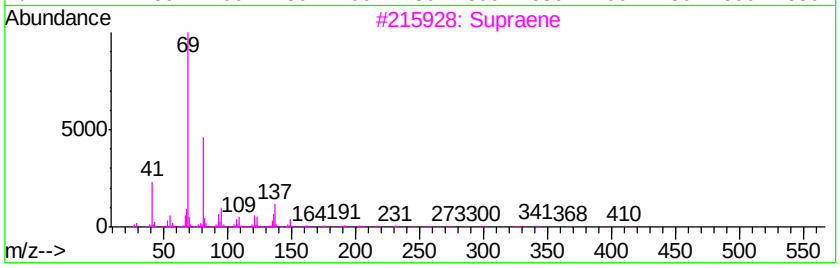
R.T.	EstConc	Area	Relative to ISTD	R.T.
22.33	10.91 ng/ul	3792090	Perylene-d12	23.34

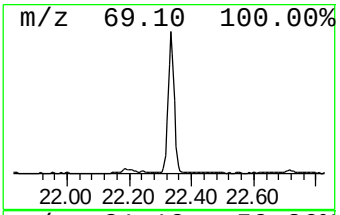
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Squalene	410	C30H50	000111-02-4	95
2	Squalene	410	C30H50	000111-02-4	94
3	Supraene	410	C30H50	007683-64-9	91
4	Squalene	410	C30H50	000111-02-4	91
5	Squalene	410	C30H50	000111-02-4	91

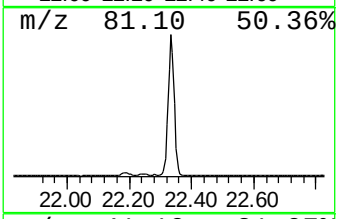


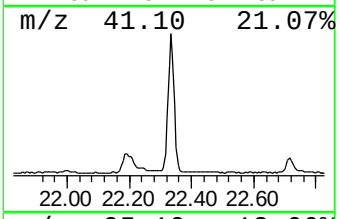


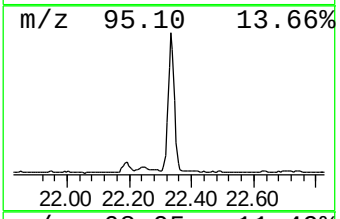


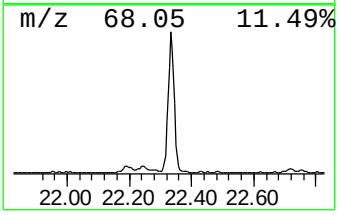












Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022120\
Data File : BP001841.D
Acq On : 22 Feb 2020 01:48
Operator : CG/JU
Sample : L1506-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBCZ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

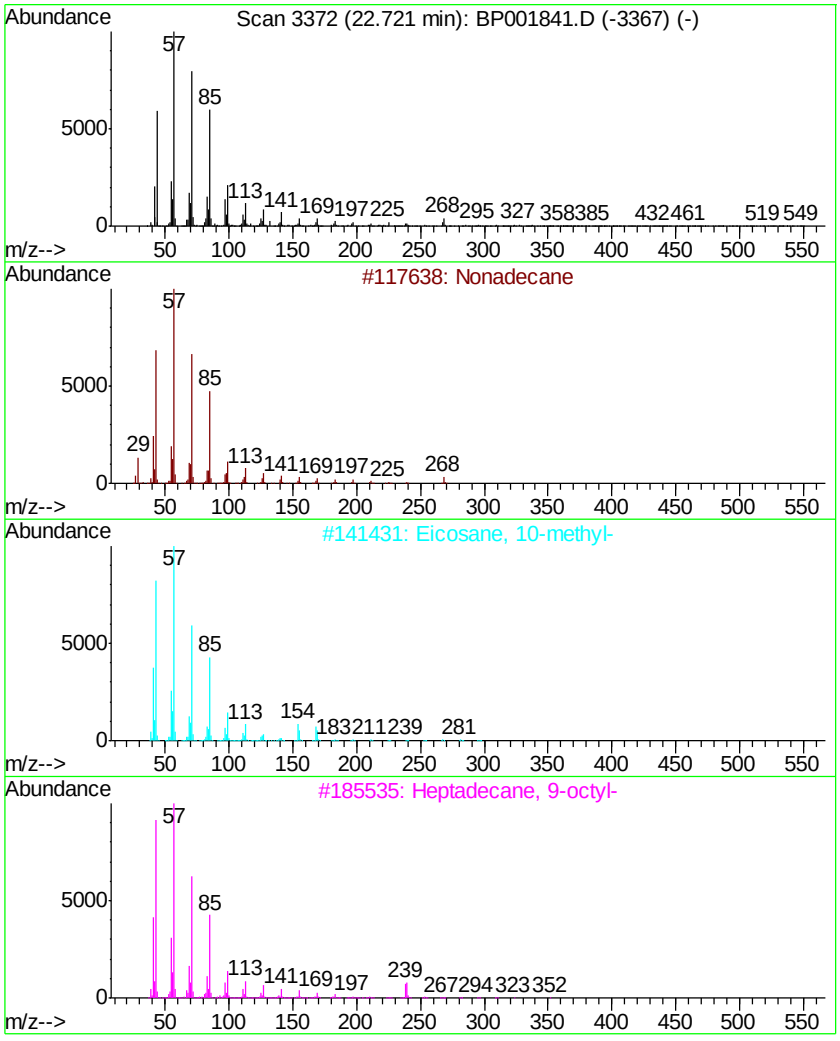
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

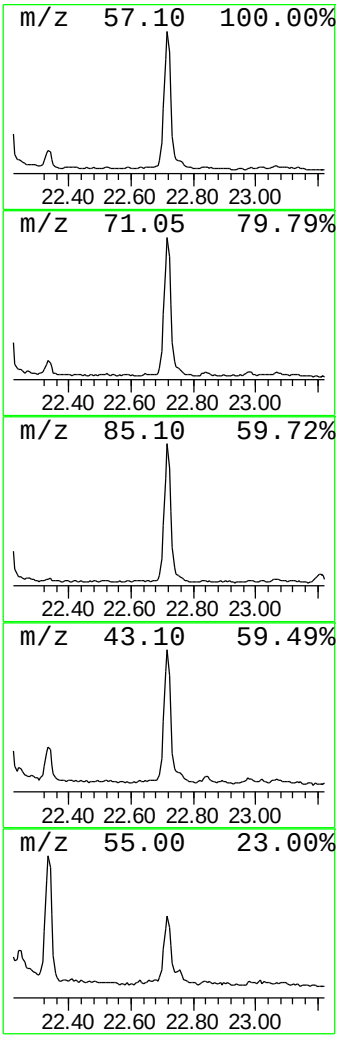
****

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.72	2.45 ng/ul	852010	Perylene-d12	23.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	96
2	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4	Docosane	310	C22H46	000629-97-0	91
5	Heptacosane	380	C27H56	000593-49-7	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022120\
Data File : BP001841.D
Acq On : 22 Feb 2020 01:48
Operator : CG/JU
Sample : L1506-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBCZ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

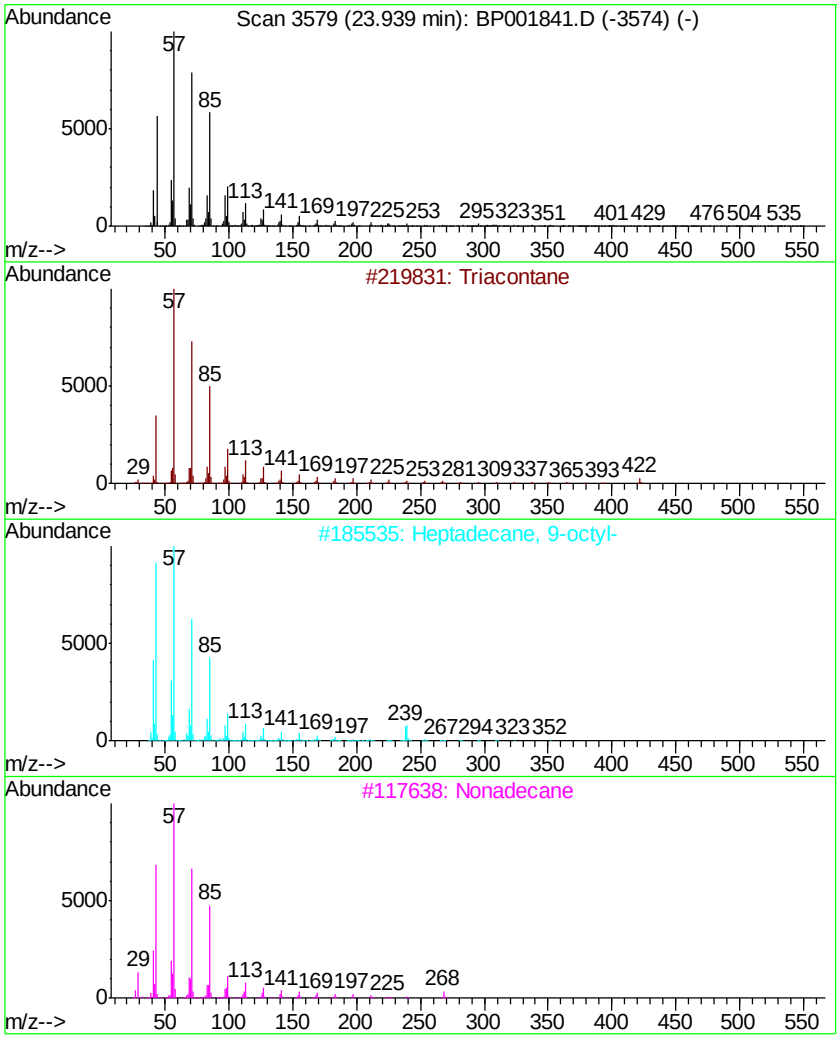
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

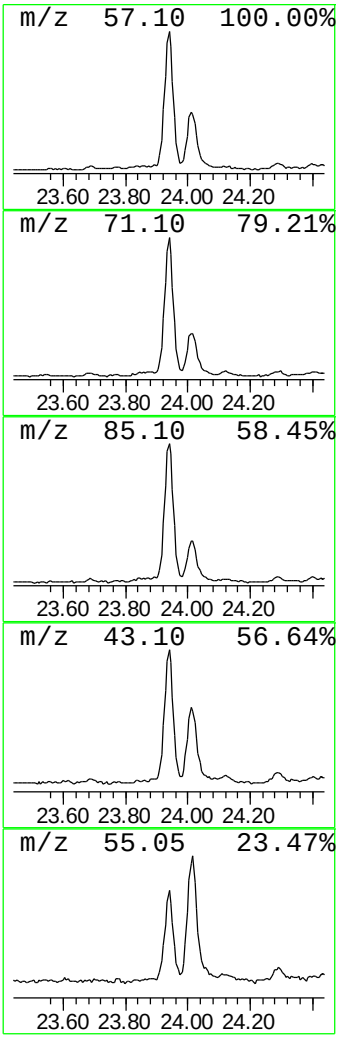
****

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.94	2.15 ng/ul	746400	Perylene-d12	23.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triacontane	422	C30H62	000638-68-6	95
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
3	Nonadecane	268	C19H40	000629-92-5	93
4	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93
5	Docosane, 11-butyl-	366	C26H54	013475-76-8	93





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022120\
Data File : BP001841.D
Acq On : 22 Feb 2020 01:48
Operator : CG/JU
Sample : L1506-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBCZ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

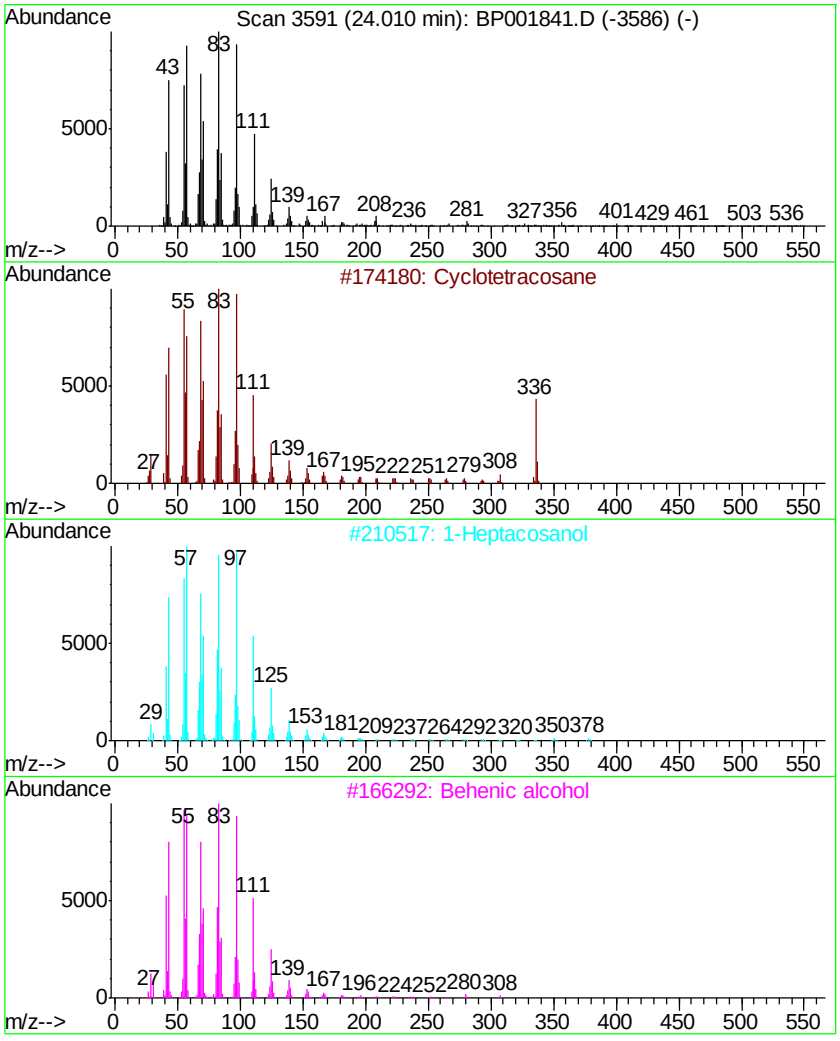
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

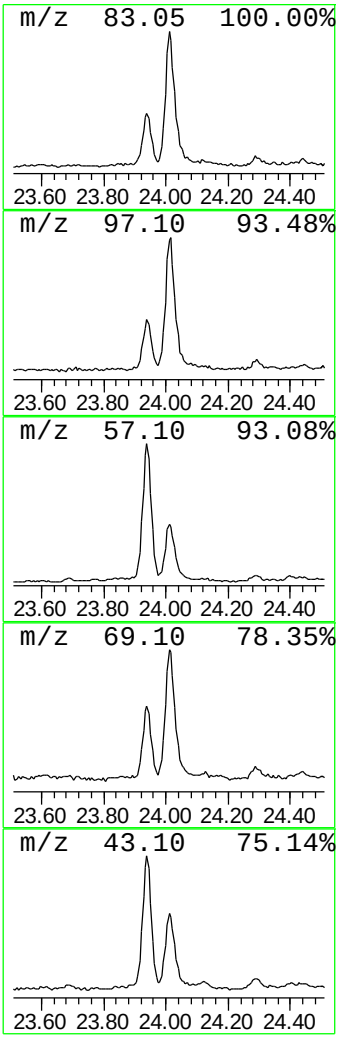
****

Peak Number 9 (DEL) Alkane: Cyclic24.01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.01	2.18 ng/ul	758887	Perylene-d12	23.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetracosane	336	C24H48	000297-03-0	97
2	1-Heptacosanol	396	C27H56O	002004-39-9	95
3	Behenic alcohol	326	C22H46O	000661-19-8	95
4	1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	95
5	Octacosyl acetate	452	C30H60O2	018206-97-8	94





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP022120\
Data File : BP001841.D
Acq On : 22 Feb 2020 01:48
Operator : CG/JU
Sample : L1506-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBCZ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.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
Butane, 2-methoxy...	2.93	10.9	ng/ul	1337280	1	7.65	2464450	20.0
unknown-01	15.67	2.8	ng/ul	894612	4	17.04	6323840	20.0
n-Hexadecanoic acid	17.97	9.1	ng/ul	2888540	4	17.04	6323840	20.0
1-Heneicosanol	20.87	9.1	ng/ul	3333460	5	21.13	7360480	20.0
13-Docosenamide, ...	22.19	3.1	ng/ul	1153200	5	21.13	7360480	20.0
Squalene	22.33	10.9	ng/ul	3792090	6	23.34	6950800	20.0
(DEL) Alkane: Str...	22.72	2.5	ng/ul	852010	6	23.34	6950800	20.0
(DEL) Alkane: Str...	23.94	2.1	ng/ul	746400	6	23.34	6950800	20.0
(DEL) Alkane: Cyc...	24.01	2.2	ng/ul	758887	6	23.34	6950800	20.0