LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062119\
Data File : BN006470.D
Acq On : 21 Jun 2019 17:30
Operator : HP/JU
Sample : K3388-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : A4214

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\SOM-EPA-BN061919MA.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	3.134	22	25	37	rVB	51529	80110	1.76%	0.119%
2	3.552	92	96	102	rBV	42227	58553	1.29%	0.087%
3	3.646	107	112	121	rVV	296577	393493	8.64%	0.584%
4	3.716	121	124	128	rVV	28769	39963	0.88%	0.059%
5	3.764	128	132	138	rVB	34712	57130	1.25%	0.085%
6	3.858	145	148	154	rVB	24502	31262	0.69%	0.046%
7	4.405	236	241	252	rVB	43065	71052	1.56%	0.105%
8	4.558	263	267	270	rBV4	8285	12864	0.28%	0.019%
9	4.658	278	284	288	rBV7	8113	14857	0.33%	0.022%
10	4.752	295	300	308	rVB	62138	98332	2.16%	0.146%
11	4.852	312	317	325	rBV	49264	84560	1.86%	0.126%
12	5.158	364	369	377	rVB	35175	56559	1.24%	0.084%
13	5.287	386	391	399	rBV3	9549	19356	0.42%	0.029%
14	5.628	444	449	458	rBV2	35932	85286	1.87%	0.127%
15	5.728	463	466	475	rVB5	6820	12759	0.28%	0.019%
16	6.122	527	533	536	rBV	13886	21293	0.47%	0.032%
17	6.234	547	552	557	rBV4	9216	17427	0.38%	0.026%
18	6.499	592	597	602	rBV3	9555	16280	0.36%	0.024%
19	6.610	610	616	620	rBV5	10397	17569	0.39%	0.026%
20	6.805	639	649	664	rBV	796461	1453638	31.91%	2.158%
21	6.963	666	676	683	rBV	707577	1083008	23.77%	1.608%
22	7.157	703	709	720	rBV	850183	1360267	29.86%	2.020%
23	7.240	720	723	728	rBV	12672	21359	0.47%	0.032%
24	7.622	780	788	796	rBV	1197793	1900405	41.72%	2.822%
25	7.840	820	825	829	rBV3	8604	11835	0.26%	0.018%
26	8.152	872	878	884	rVB7	7223	14473	0.32%	0.021%
27	8.334	903	909	923	rBV	861446	1436028	31.52%	2.132%
28	8.446	923	928	938	rVB2	34336	66650	1.46%	0.099%
29	8.775	977	984	992	rVV	787735	1334750	29.30%	1.982%
30	8.840	992	995	1004	rVV	26285	55549	1.22%	0.082%
31	8.940	1008	1012	1019	rVB6	13098	27503	0.60%	0.041%
32	9.187	1048	1054	1060	rVB	19368	30098	0.66%	0.045%
33	9.499	1100	1107	1121	rBV	640289	1111254	24.39%	1.650%
34	10.034	1192	1198	1215	rBV	991825	1705806	37.45%	2.533%

LSC Area Percent Report

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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : A4214

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\SOM-EPA-BN061919MA.M
Title : SVOA CALIBRATION

35	10.404	1252	1261	1268	rBV	1760189	2911674	63.92%	4.323%
36	10.551	1280	1286	1305	rBV	268482	610879	13.41%	0.907%
37	11.375	1420	1426	1430	rBV7	5588	11720	0.26%	0.017%
38	11.834	1498	1504	1512	rVB4	15664	25365	0.56%	0.038%
39	11.916	1512	1518	1519	rBV3	9383	13722	0.30%	0.020%
40	12.004	1526	1533	1539	rVB2	14311	25785	0.57%	0.038%
41	12.081	1539	1546	1552	rBV	14852	27527	0.60%	0.041%
42	12.298	1579	1583	1588	rVB2	7900	12933	0.28%	0.019%
43	12.469	1607	1612	1619	rVB6	10330	18398	0.40%	0.027%
44	12.869	1675	1680	1686	rBV6	7410	13203	0.29%	0.020%
45	13.034	1700	1708	1712	rBV4	20878	33193	0.73%	0.049%
46	13.098	1712	1719	1723	rVV3	20429	34679	0.76%	0.051%
47	13.145	1723	1727	1730	rVV4	29783	41464	0.91%	0.062%
48	13.210	1733	1738	1741	rVV5	9624	20182	0.44%	0.030%
49	13.240	1741	1743	1750	rVB6	10283	15816	0.35%	0.023%
50	13.434	1771	1776	1783	rBV8	9588	25349	0.56%	0.038%
51	13.598	1797	1804	1809	rBV7	10520	22708	0.50%	0.034%
52	13.692	1813	1820	1824	rVV	1862844	2614200	57.39%	3.882%
53	13.740	1824	1828	1835	rVB	661023	903068	19.82%	1.341%
54	13.969	1855	1867	1881	rBV	2146718	3301936	72.48%	4.903%
55	14.098	1886	1889	1896	rVB6	9111	18625	0.41%	0.028%
56	14.181	1899	1903	1909	rVB	27334	39627	0.87%	0.059%
57	14.281	1909	1920	1927	rBV2	2543250	3828102	84.04%	5.684%
58	14.339	1927	1930	1935	rVB3	35928	48193	1.06%	0.072%
59	14.398	1937	1940	1947	rVB5	21027	31345	0.69%	0.047%
60	14.510	1951	1959	1971	rBV	365697	753850	16.55%	1.119%
61	14.651	1979	1983	1991	rBV3	39529	81541	1.79%	0.121%
62	14.716	1991	1994	1998	rVB4	14490	18406	0.40%	0.027%
63	14.804	2007	2009	2018	rVB10	7823	15148	0.33%	0.022%
64	14.957	2032	2035	2041	rVB7	10622	15081	0.33%	0.022%
65	15.086	2051	2057	2063	rBV	46798	75495	1.66%	0.112%
66	15.139	2063	2066	2072	rVB2	29024	40016	0.88%	0.059%
67	15.281	2083	2090	2096	rBV	2795957	3975711	87.28%	5.903%
68	15.339	2097	2100	2105	rVV5	22477	35873	0.79%	0.053%
69	15.422	2109	2114	2123	rBV	290096	444799	9.76%	0.660%
70	15.522	2126	2131	2142	rVV4	43911	105162	2.31%	0.156%
71	15.669	2153	2156	2165	rVB	76085	115040	2.53%	0.171%

LSC Area Percent Report

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Sample : K3388-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : A4214

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

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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Title : SVOA CALIBRATION

72	15.839	2180	2185	2188	rBV4	24096	36893	0.81%	0.055%
73	15.910	2193	2197	2200	rBV3	40918	54911	1.21%	0.082%
74	16.010	2209	2214	2215	rBV3	65129	90352	1.98%	0.134%
75	16.222	2246	2250	2257	rVB	42166	67732	1.49%	0.101%
76	16.316	2262	2266	2269	rBV4	13233	20391	0.45%	0.030%
77	16.381	2275	2277	2285	rVB3	19471	29861	0.66%	0.044%
78	16.639	2317	2321	2327	rVB	163409	208769	4.58%	0.310%
79	16.798	2345	2348	2351	rBV	60213	73540	1.61%	0.109%
80	16.851	2354	2357	2363	rVB5	35776	54546	1.20%	0.081%
81	17.039	2382	2389	2394	rBV	3103088	4498759	98.76%	6.680%
82	17.075	2394	2395	2399	rVV	97211	105066	2.31%	0.156%
83	17.133	2399	2405	2417	rVB2	2932560	4197723	92.15%	6.233%
84	17.945	2538	2543	2551	rBV	727409	1128719	24.78%	1.676%
85	19.110	2738	2741	2745	rVB	228958	262239	5.76%	0.389%
86	19.239	2759	2763	2771	rBV2	313408	640668	14.06%	0.951%
87	19.445	2793	2798	2806	rBV	3054078	4555364	100.00%	6.764%
88	20.504	2976	2978	2981	rVB	201243	187288	4.11%	0.278%
89	20.951	3051	3054	3056	rBV	574799	681898	14.97%	1.012%
90	20.974	3056	3058	3067	rVB	827285	1181403	25.93%	1.754%
91	21.174	3089	3092	3097	rVB	380810	413934	9.09%	0.615%
92	21.263	3102	3107	3111	rVV	2643897	3499566	76.82%	5.196%
93	21.427	3130	3135	3140	rBV3	519900	892531	19.59%	1.325%
94	21.480	3141	3144	3148	rBV	679632	777688	17.07%	1.155%
95	21.851	3204	3207	3210	rBV	738155	782673	17.18%	1.162%
96	22.310	3282	3285	3293	rVB	639508	819983	18.00%	1.218%
97	22.815	3367	3371	3377	rVB2	1086467	1584141	34.78%	2.352%
98	23.357	3458	3463	3469	rBV2	1748516	3537377	77.65%	5.252%
99	23.498	3482	3487	3494	rVB	1591982	2708799	59.46%	4.022%
100	24.010	3571	3574	3586	rVB	610361	1129690	24.80%	1.677%

Sum of corrected areas: 67349647

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062119\
Data File : BN006470.D
Acq On : 21 Jun 2019 17:30
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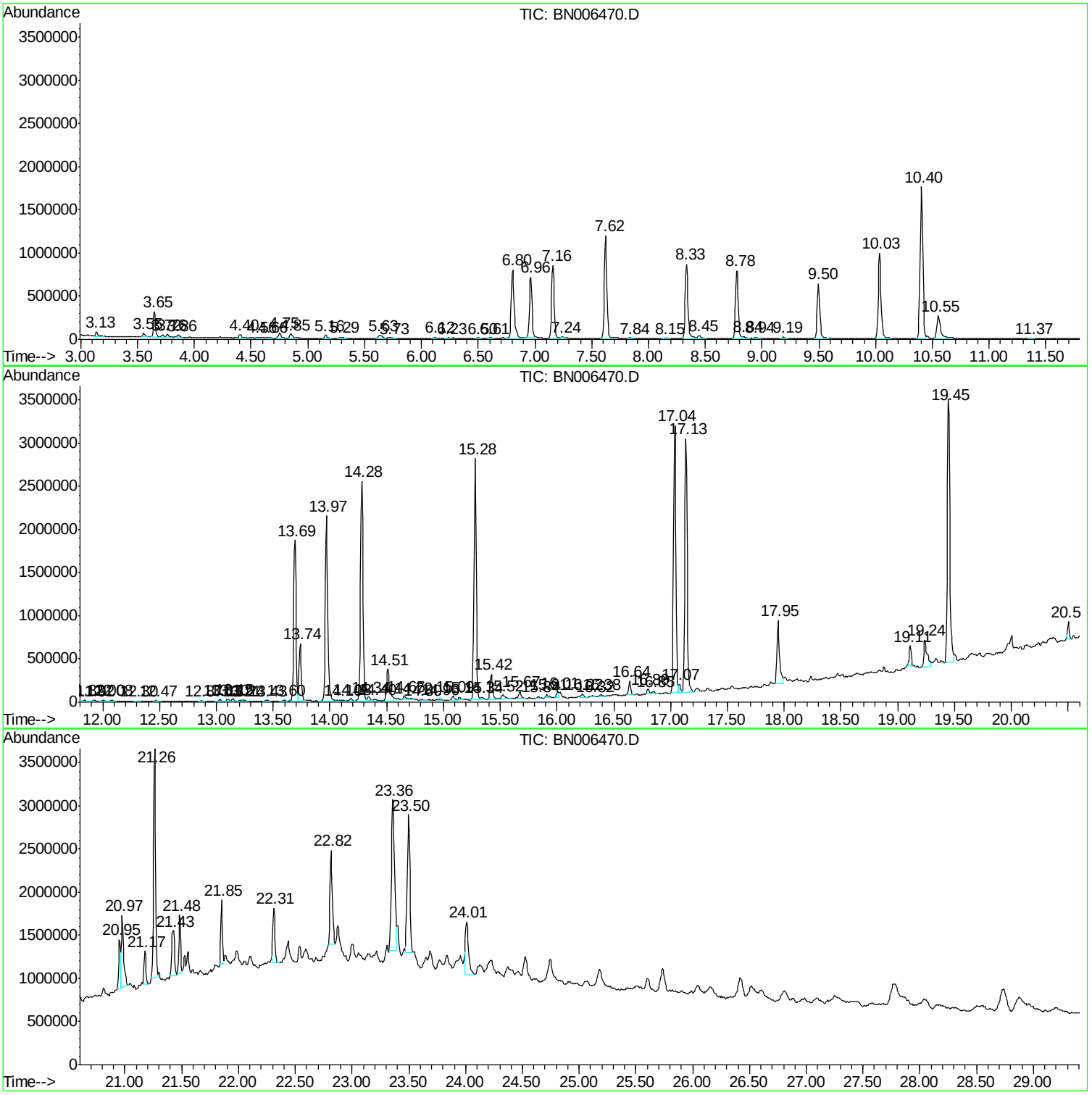
Sample : K3388-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A4214

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.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_N\DATA\BN062119\
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Instrument : BNA_N
ClientSampleId : A4214

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Quant Title : SVOA CALIBRATION

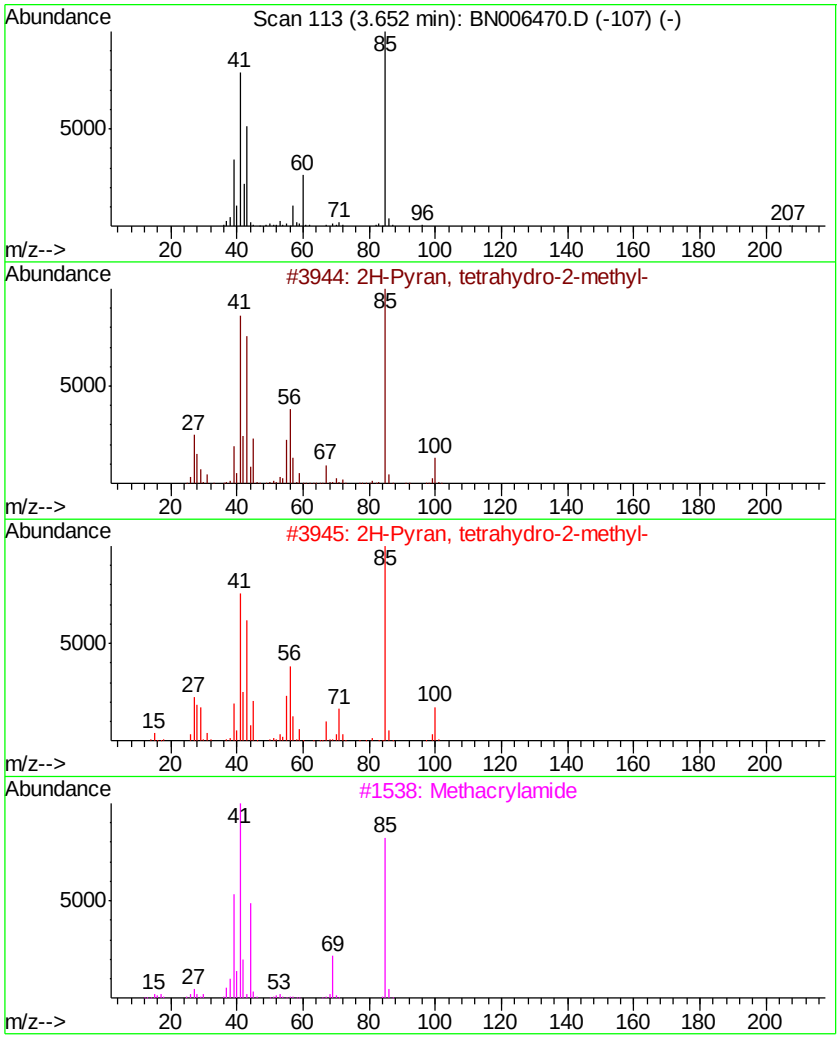
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

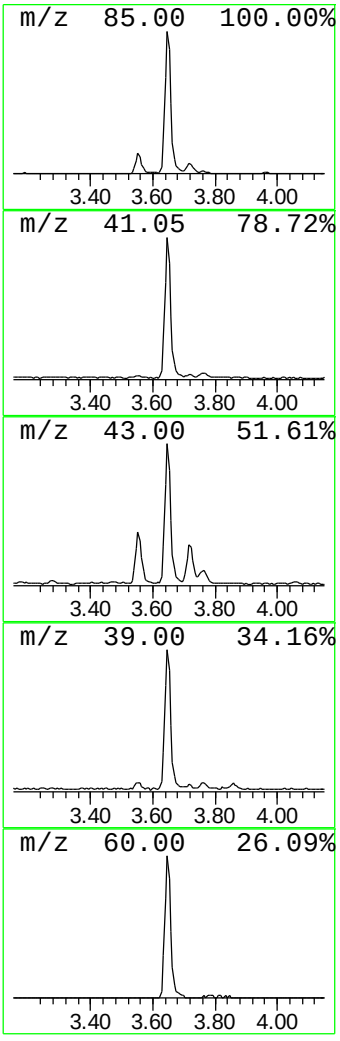
****

Peak Number 1 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.65	4.14 ng/ul	393493	1,4-Dichlorobenzene-d4	7.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	64
2	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	38
3	Methacrylamide	85	C4H7NO	000079-39-0	36
4	2-Pyrrolidinone	85	C4H7NO	000616-45-5	28
5	2(3H)-Furanone, 5-acetyldihydro-	128	C6H8O3	029393-32-6	28





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062119\
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Sample : K3388-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : A4214

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Quant Title : SVOA CALIBRATION

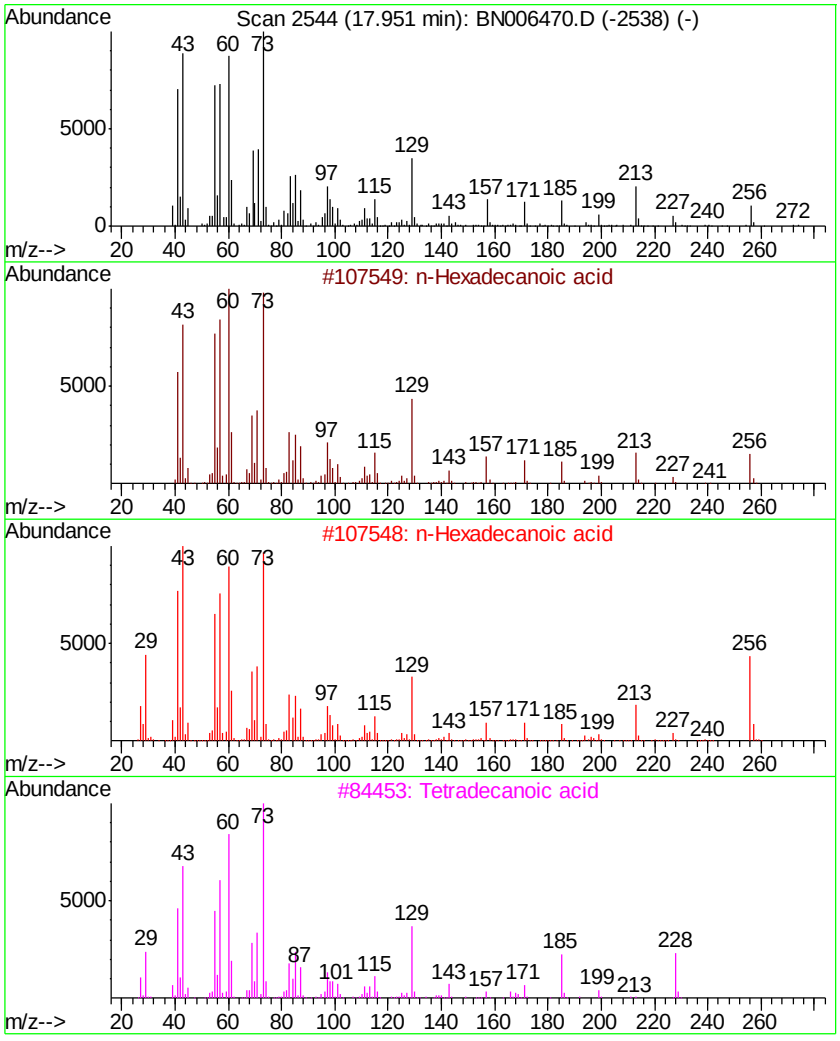
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

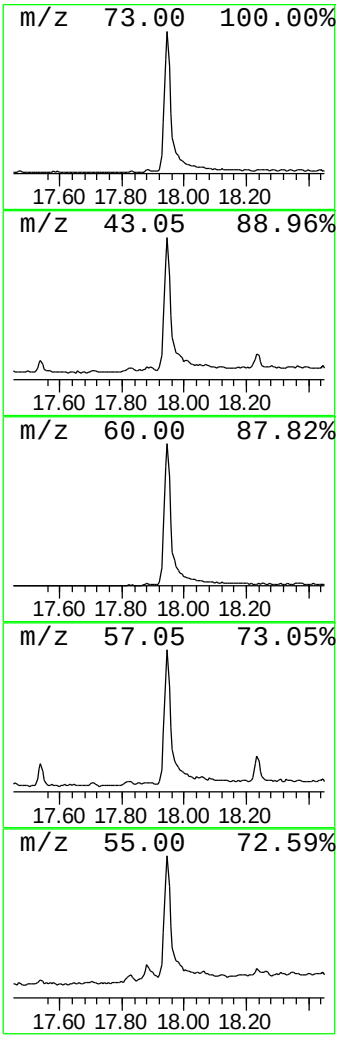
****

Peak Number 2 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.95	5.02 ng/ul	1128720	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	98
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	91
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
5	Tridecanoic acid	214	C13H26O2	000638-53-9	76





Library Search Compound Report

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Instrument : BNA_N
ClientSampleId : A4214

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Quant Title : SVOA CALIBRATION

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

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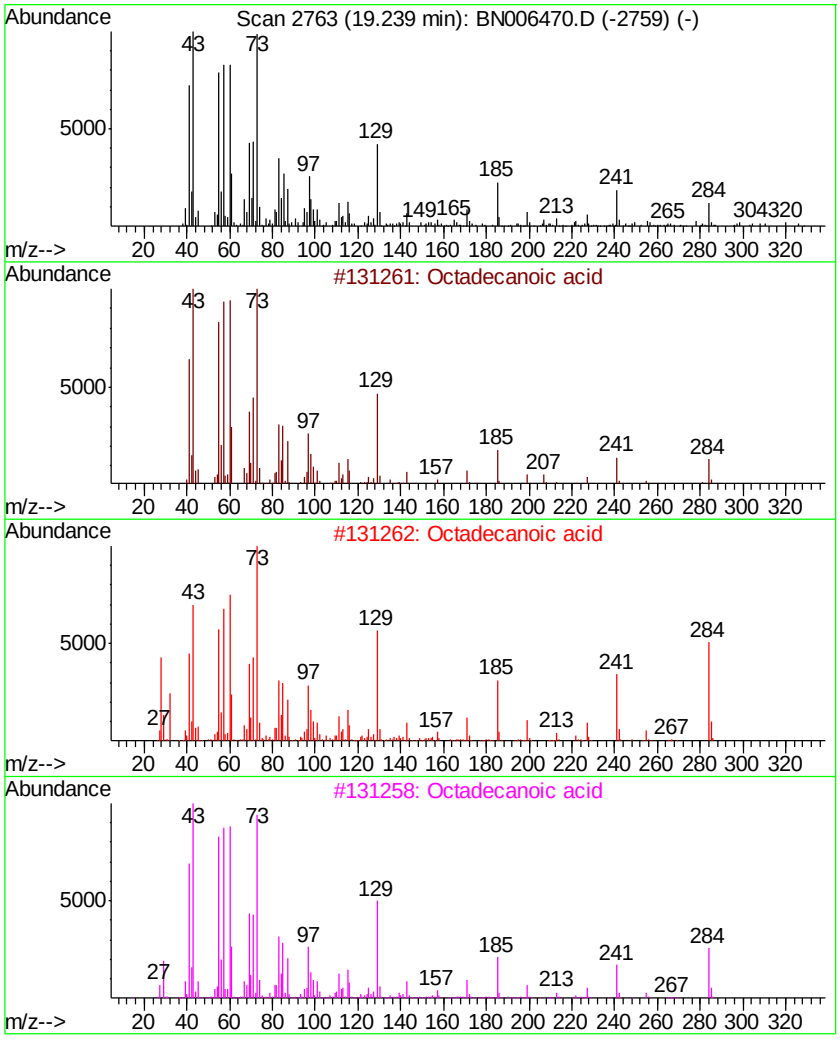
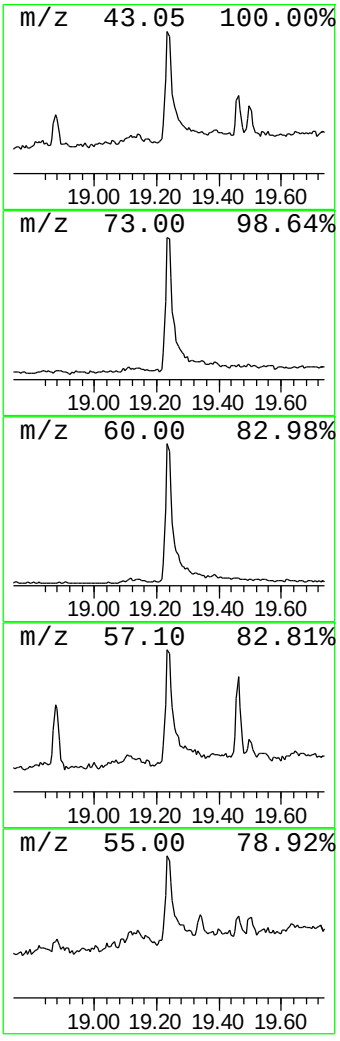
Peak Number 3 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.24	3.66 ng/ul	640668	Chrysene-d12	21.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	99
4	Octadecanoic acid	284	C18H36O2	000057-11-4	98
5	Octadecanoic acid	284	C18H36O2	000057-11-4	96





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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : A4214

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Quant Title : SVOA CALIBRATION

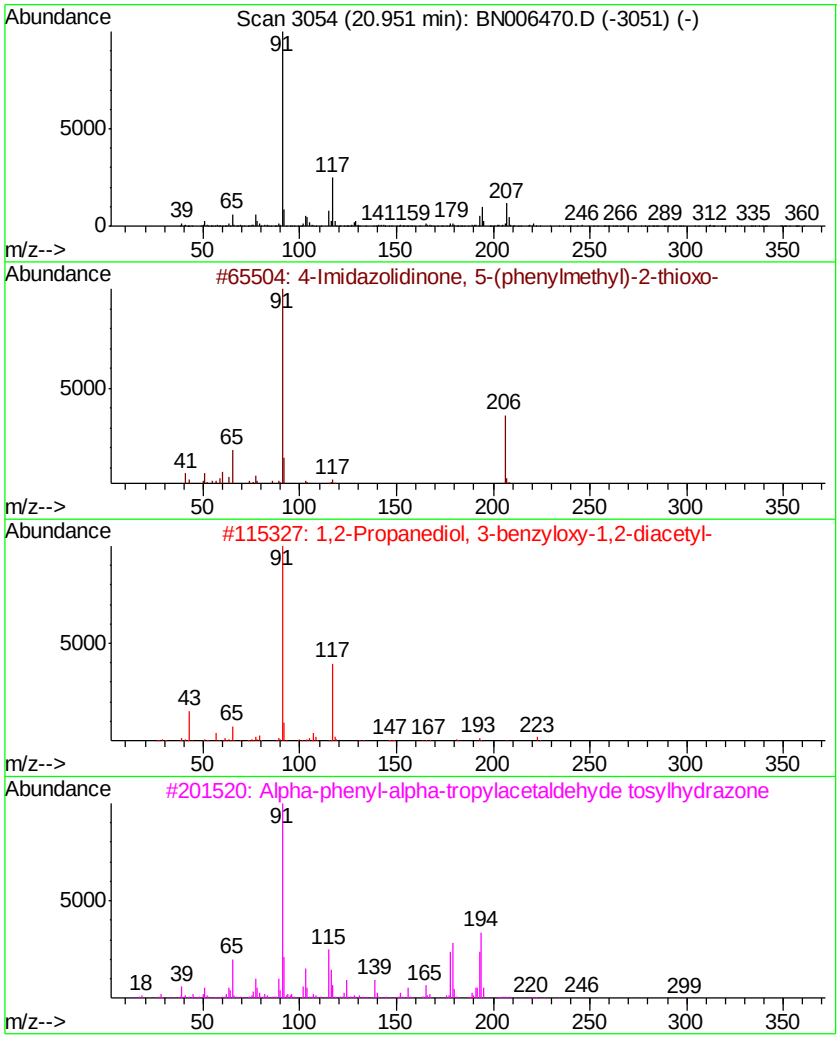
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

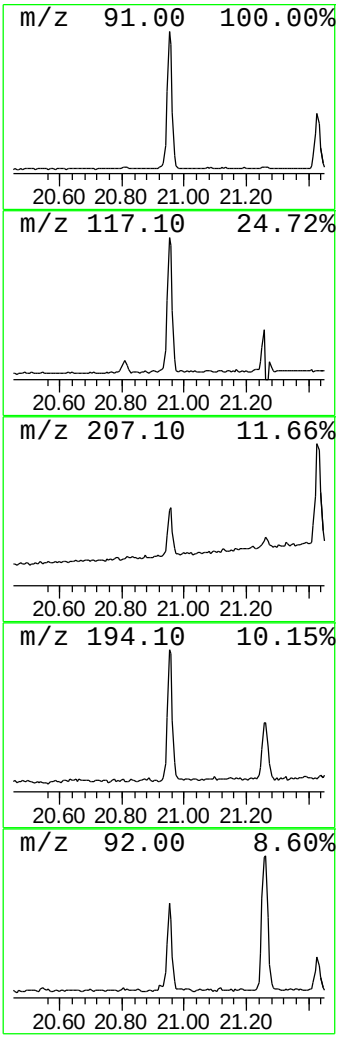
****

Peak Number 4 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.95	3.90 ng/ul	681898	Chrysene-d12	21.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Imidazolidinone, 5-(phenylmeth...	206	C10H10N2OS	006330-09-2	23
2	1,2-Propanediol, 3-benzyloxy-1,2...	266	C14H18O5	013754-10-4	12
3	Alpha-phenyl-alpha-tropylacetald...	378	C22H22N2O2S	022532-16-7	12
4	Pentanoic acid, 2-benzenmethanim...	262	C13H18N4O2	005411-76-7	10
5	2,5-Diphenyltetrazole	222	C13H10N4	018039-45-7	9





Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062119\
Data File : BN006470.D
Acq On : 21 Jun 2019 17:30
Operator : HP/JU
Sample : K3388-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : A4214

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Quant Title : SVOA CALIBRATION

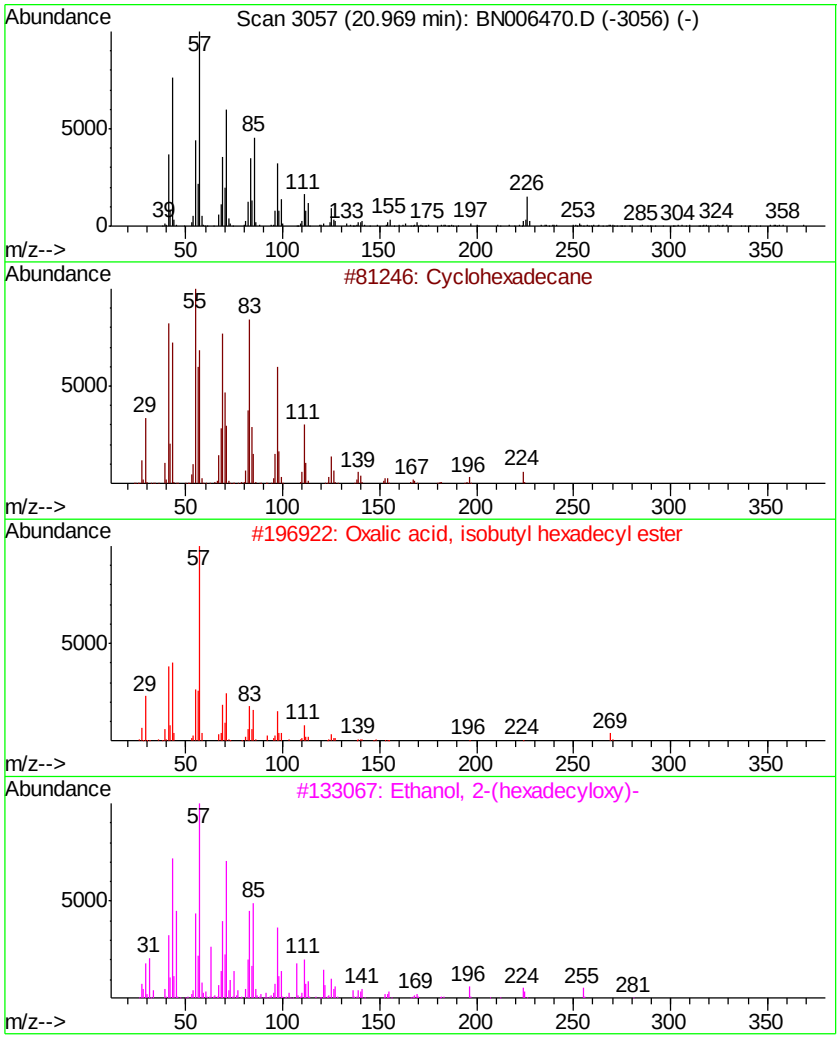
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

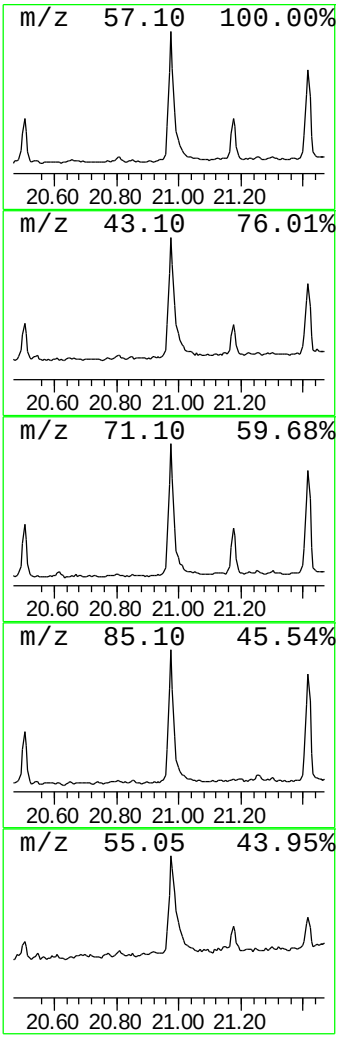
****

Peak Number 5 (DEL) Alkane: Cyclic20.97 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.97	6.75 ng/ul	1181400	Chrysene-d12	21.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	94
2	Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	83
3	Ethanol, 2-(hexadecyloxy)-	286	C18H38O2	002136-71-2	83
4	Hexadecane	226	C16H34	000544-76-3	64
5	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	53





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Data File : BN006470.D
Acq On : 21 Jun 2019 17:30
Operator : HP/JU
Sample : K3388-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : A4214

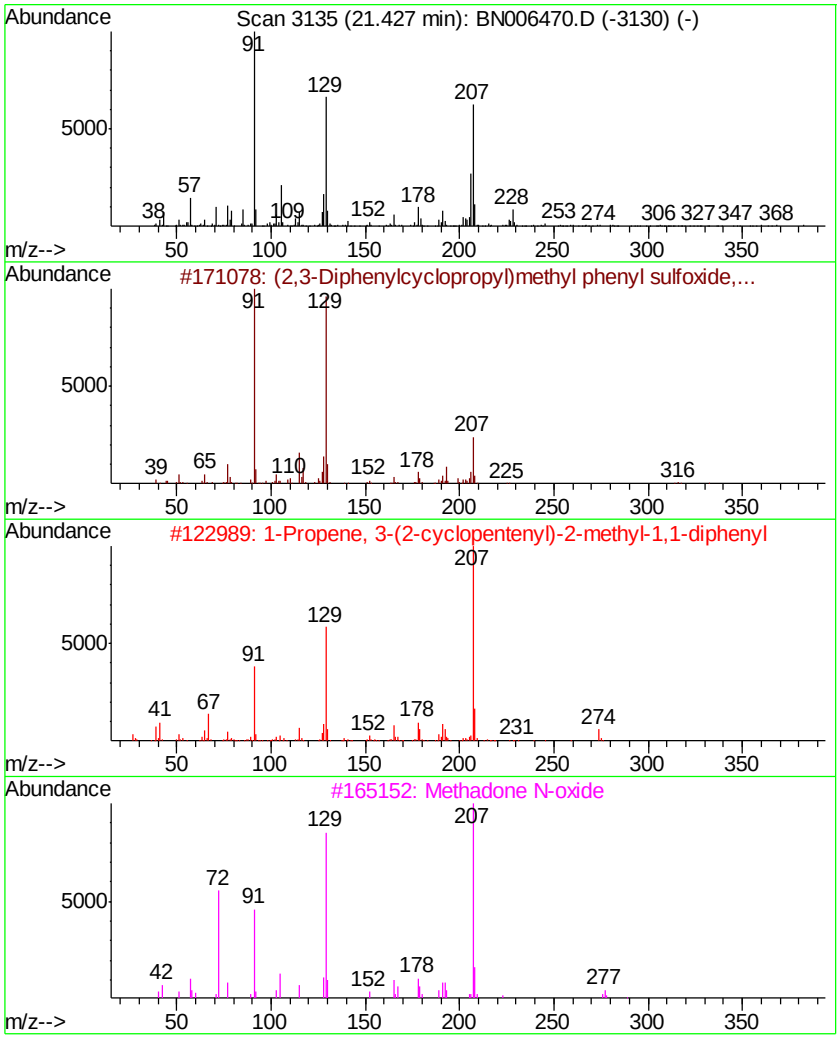
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Quant Title : SVOA CALIBRATION

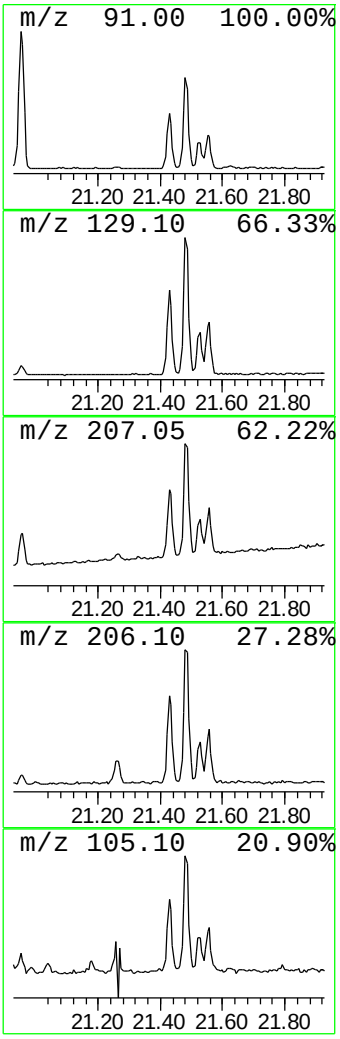
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (2,3-Diphenylcyclopropyl)me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.43	5.10 ng/ul	892531	Chrysene-d12	21.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	53
2	1-Propene, 3-(2-cyclopentenyl)-2...	274	C21H22	1000154-23-3	50
3	Methadone N-oxide	325	C21H27NO2	1000120-80-7	49
4	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	46
5	1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	30





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Acq On : 21 Jun 2019 17:30
Operator : HP/JU
Sample : K3388-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : A4214

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Quant Title : SVOA CALIBRATION

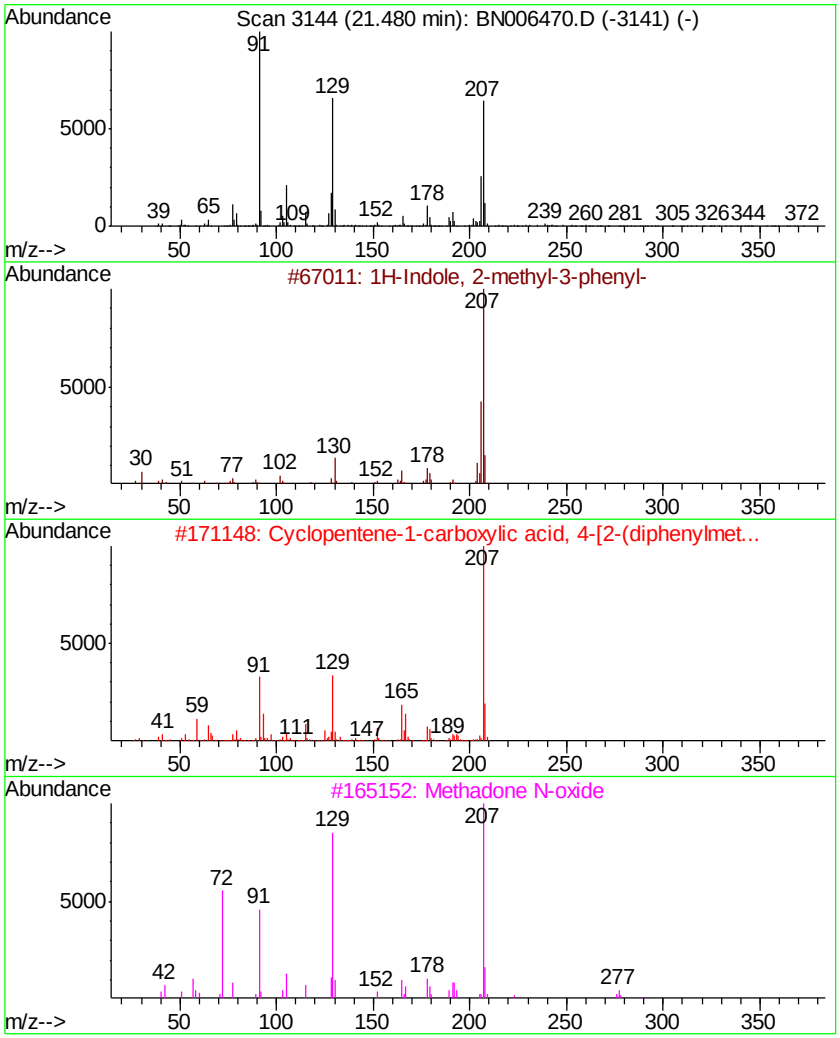
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

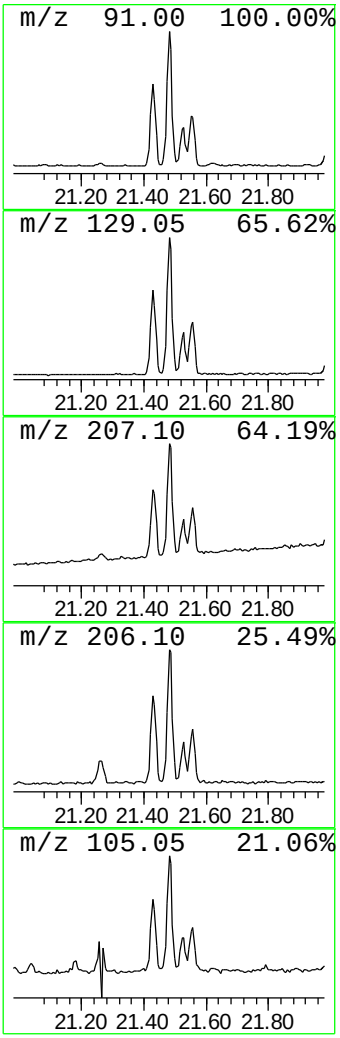
****

Peak Number 7 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.48	4.44 ng/ul	777688	Chrysene-d12	21.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	41
2	Cyclopentene-1-carboxylic acid, ...	332	C23H24O2	1000159-40-6	38
3	Methadone N-oxide	325	C21H27NO2	1000120-80-7	38
4	5-Methyl-2-phenylindolizine	207	C15H13N	036944-99-7	38
5	1-benzylindole	207	C15H13N	1000334-31-3	30





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Operator : HP/JU
Sample : K3388-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : A4214

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Quant Title : SVOA CALIBRATION

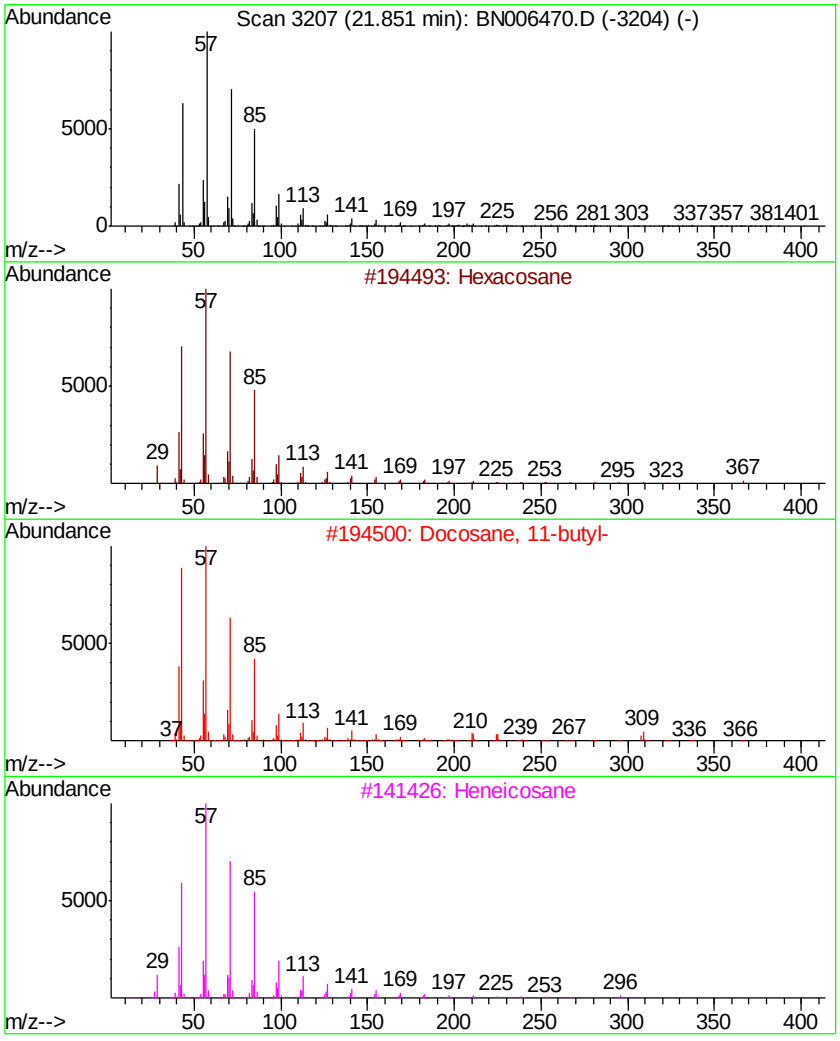
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

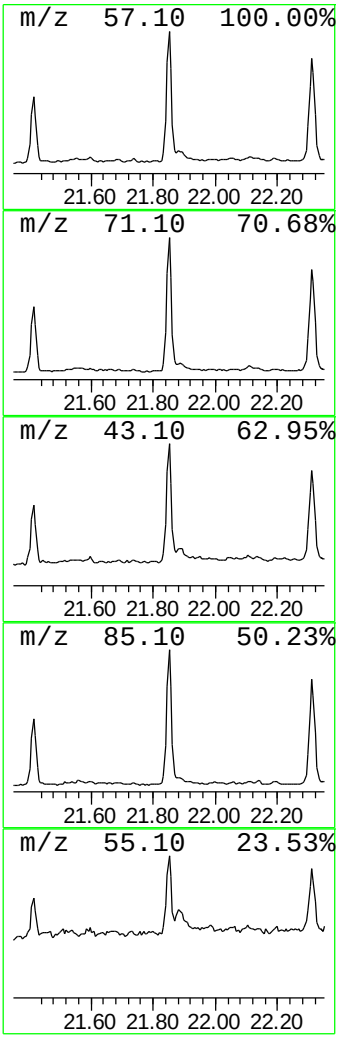
****

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.85	4.47 ng/ul	782673	Chrysene-d12	21.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	91
2	Docosane, 11-butyl-	366	C26H54	013475-76-8	91
3	Heneicosane	296	C21H44	000629-94-7	91
4	Triacontane	422	C30H62	000638-68-6	91
5	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91





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Data File : BN006470.D
Acq On : 21 Jun 2019 17:30
Operator : HP/JU
Sample : K3388-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : A4214

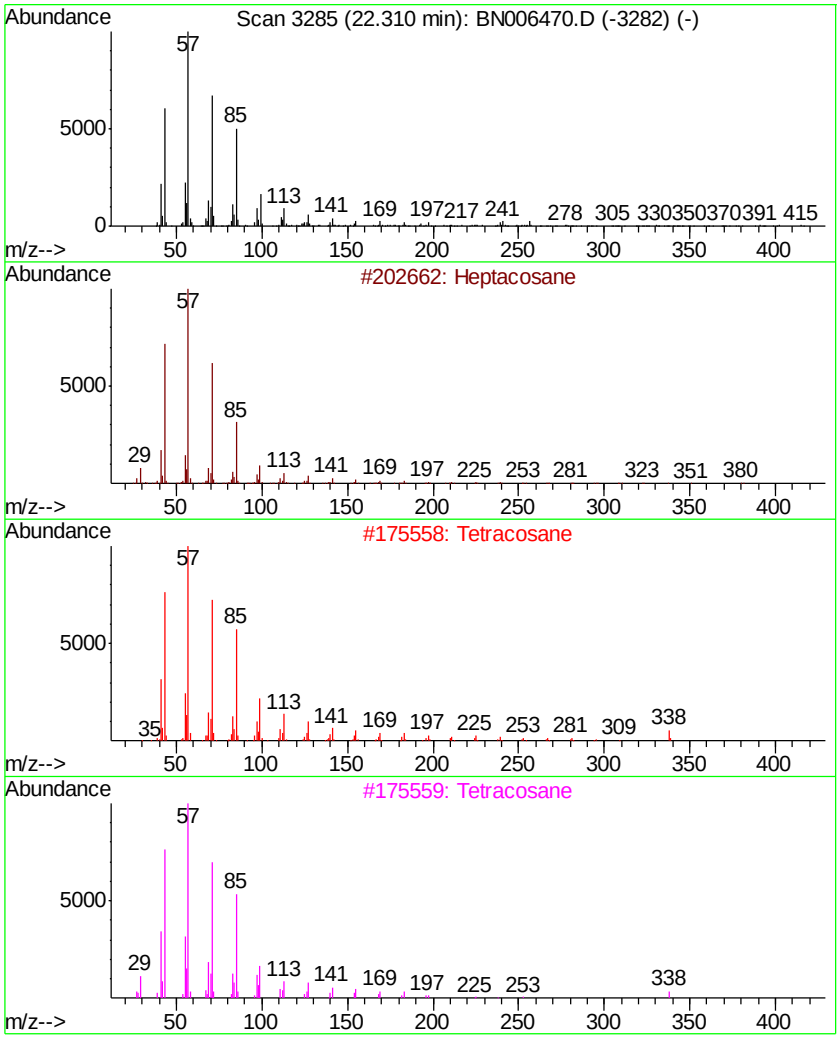
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Quant Title : SVOA CALIBRATION

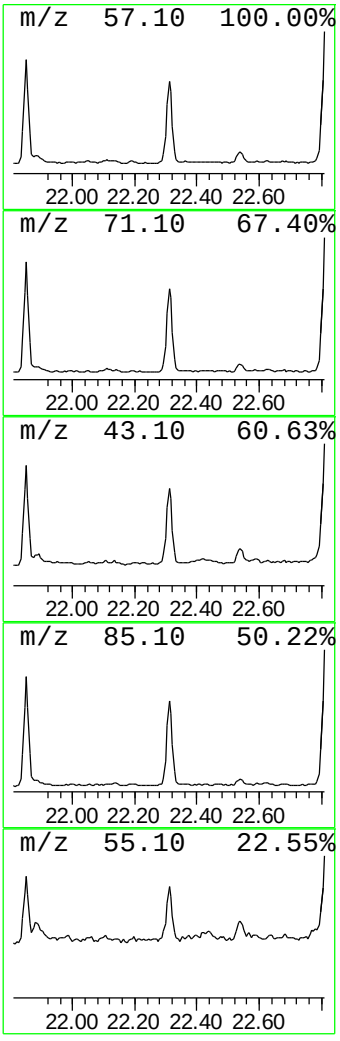
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.31	4.69 ng/ul	819983	Chrysene-d12	21.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	98
2	Tetracosane	338	C24H50	000646-31-1	96
3	Tetracosane	338	C24H50	000646-31-1	95
4	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	93
5	Heneicosane	296	C21H44	000629-94-7	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062119\
Data File : BN006470.D
Acq On : 21 Jun 2019 17:30
Operator : HP/JU
Sample : K3388-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : A4214

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Quant Title : SVOA CALIBRATION

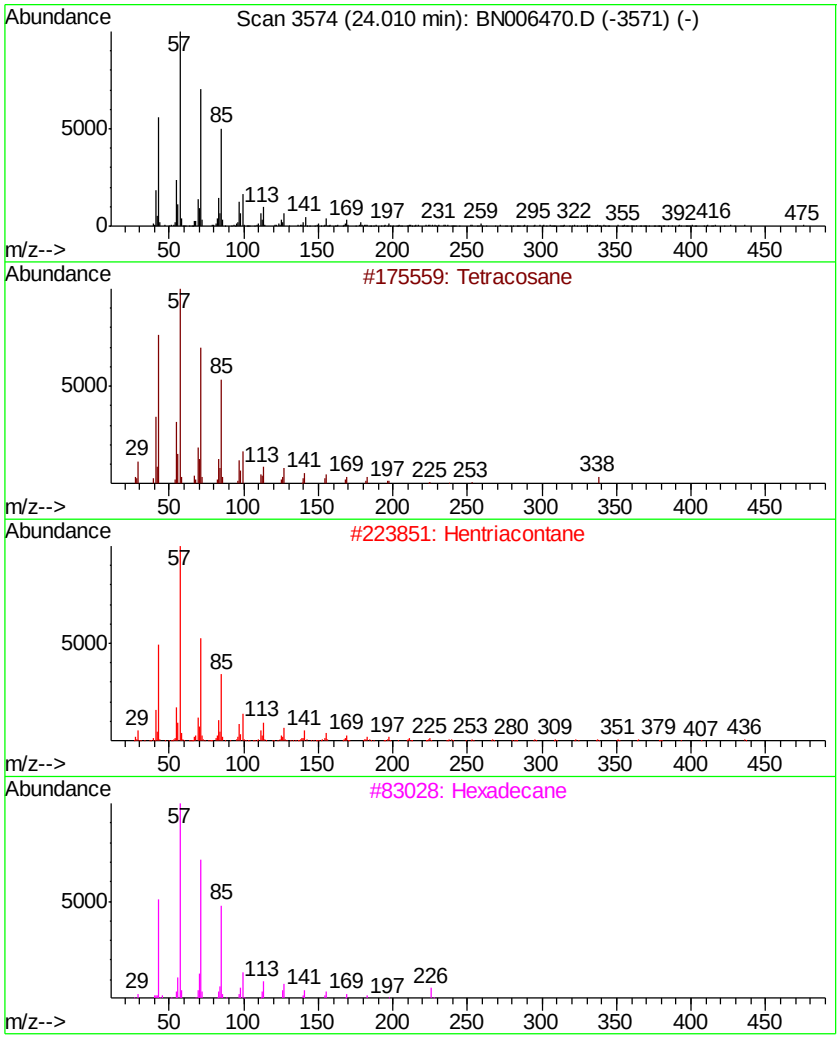
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

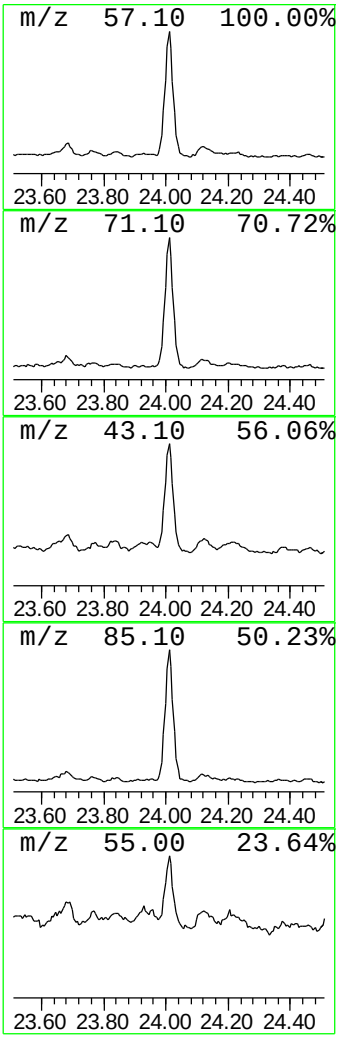
****

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.01	8.34 ng/ul	1129690	Perylene-d12	23.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	93
2	Hentriacontane	437	C31H64	000630-04-6	93
3	Hexadecane	226	C16H34	000544-76-3	91
4	Heneicosane	296	C21H44	000629-94-7	90
5	Hexacosane	366	C26H54	000630-01-3	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062119\
Data File : BN006470.D
Acq On : 21 Jun 2019 17:30
Operator : HP/JU
Sample : K3388-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : A4214

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.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
2H-Pyran, tetrahy...	3.65	4.1	ng/ul	393493	1	7.62	1900410	20.0
n-Hexadecanoic acid	17.95	5.0	ng/ul	1128720	4	17.04	4498760	20.0
Octadecanoic acid	19.24	3.7	ng/ul	640668	5	21.26	3499570	20.0
unknown-01	20.95	3.9	ng/ul	681898	5	21.26	3499570	20.0
(DEL) Alkane: Cyc...	20.97	6.8	ng/ul	1181400	5	21.26	3499570	20.0
(2,3-Diphenylcycl...	21.43	5.1	ng/ul	892531	5	21.26	3499570	20.0
unknown-02	21.48	4.4	ng/ul	777688	5	21.26	3499570	20.0
(DEL) Alkane: Str...	21.85	4.5	ng/ul	782673	5	21.26	3499570	20.0
(DEL) Alkane: Str...	22.31	4.7	ng/ul	819983	5	21.26	3499570	20.0
(DEL) Alkane: Str...	24.01	8.3	ng/ul	1129690	6	23.50	2708800	20.0