LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN070518\
Data File : BN001968.D
Acq On : 05 Jul 2018 15:34
Operator : JU/SJ
Sample : J3776-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAZQ9

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-BN061918.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	3.023	3	6	13	rVB	1056279	1403762	8.71%	0.792%
2	3.081	13	16	26	rVB2	272113	390009	2.42%	0.220%
3	3.170	27	31	33	rBV	203900	249171	1.55%	0.141%
4	3.270	44	48	54	rBV	138664	185286	1.15%	0.105%
5	3.752	126	130	141	rVB	71063	107691	0.67%	0.061%
6	4.069	179	184	188	rBV	64006	96339	0.60%	0.054%
7	4.117	188	192	196	rVV	237238	309255	1.92%	0.175%
8	4.170	196	201	213	rVB	810930	1138729	7.07%	0.643%
9	4.334	225	229	236	rVB	81301	121792	0.76%	0.069%
10	4.875	318	321	328	rVV	54320	80212	0.50%	0.045%
11	5.199	368	376	382	rBV	127669	227040	1.41%	0.128%
12	5.334	394	399	403	rVV	47367	72546	0.45%	0.041%
13	5.405	405	411	415	rVV5	34809	73637	0.46%	0.042%
14	5.452	415	419	426	rVB	60457	94949	0.59%	0.054%
15	5.534	426	433	438	rBV	1227678	1710176	10.61%	0.965%
16	5.587	438	442	451	rVB	484232	689777	4.28%	0.389%
17	5.793	473	477	483	rVB4	47638	82451	0.51%	0.047%
18	5.981	506	509	515	rVB	48031	72805	0.45%	0.041%
19	6.399	575	580	582	rVV	62324	87699	0.54%	0.049%
20	6.422	582	584	590	rVB2	53022	77378	0.48%	0.044%
21	6.599	605	614	624	rBV4	41882	104921	0.65%	0.059%
22	6.728	624	636	647	rBV	160919	393113	2.44%	0.222%
23	6.887	647	663	675	rBV2	777699	1525906	9.47%	0.861%
24	6.981	675	679	684	rBV	58698	87071	0.54%	0.049%
25	7.052	684	691	696	rBV	2235732	3477312	21.58%	1.963%
26	7.105	696	700	706	rVB	602667	861289	5.34%	0.486%
27	7.175	706	712	717	rVB	344280	514337	3.19%	0.290%
28	7.246	717	724	733	rBV	2362313	3844742	23.86%	2.170%
29	7.469	758	762	770	rVB8	29261	81835	0.51%	0.046%
30	7.710	796	803	811	rBV	2272300	3719246	23.08%	2.099%
31	7.899	828	835	843	rVB2	94580	155847	0.97%	0.088%
32	8.010	850	854	864	rVB6	40743	94972	0.59%	0.054%
33	8.287	895	901	911	rVB2	211139	391524	2.43%	0.221%
34	8.410	915	922	929	rVV	1965761	3220053	19.98%	1.817%

LSC Area Percent Report

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Operator : JU/SJ
Sample : J3776-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAZQ9

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-BN061918.M
Title : SVOA CALIBRATION

35	8.522	932	941	947	rVB2	510618	1213794	7.53%	0.685%
36	8.710	967	973	978	rVB	89562	137833	0.86%	0.078%
37	8.787	978	986	990	rBV2	94892	180472	1.12%	0.102%
38	8.857	990	998	1005	rBV	2757736	4652599	28.87%	2.626%
39	8.981	1013	1019	1024	rBV	111135	177666	1.10%	0.100%
40	9.093	1033	1038	1044	rVB2	48894	87423	0.54%	0.049%
41	9.246	1060	1064	1069	rBV3	59381	104828	0.65%	0.059%
42	9.569	1111	1119	1126	rBV	2339373	3942563	24.47%	2.225%
43	9.704	1134	1142	1148	rBV3	49113	119070	0.74%	0.067%
44	9.822	1154	1162	1172	rVV2	324266	633189	3.93%	0.357%
45	10.004	1187	1193	1196	rBV5	35669	74310	0.46%	0.042%
46	10.046	1196	1200	1203	rVV3	87382	150925	0.94%	0.085%
47	10.104	1203	1210	1220	rVB	3578830	6043156	37.50%	3.411%
48	10.346	1247	1251	1258	rVB	82829	122774	0.76%	0.069%
49	10.481	1264	1274	1287	rVB	3307003	5962516	37.00%	3.365%
50	11.122	1377	1383	1389	rBV2	63979	119827	0.74%	0.068%
51	11.293	1405	1412	1423	rVV2	411917	791608	4.91%	0.447%
52	11.446	1432	1438	1446	rVB6	66228	150086	0.93%	0.085%
53	11.834	1493	1504	1517	rVB5	93695	243618	1.51%	0.138%
54	11.957	1520	1525	1532	rBV3	102929	191859	1.19%	0.108%
55	12.216	1563	1569	1575	rBV	95410	158204	0.98%	0.089%
56	12.516	1616	1620	1627	rVB5	52317	106996	0.66%	0.060%
57	12.622	1632	1638	1648	rBV	350232	622233	3.86%	0.351%
58	12.746	1655	1659	1663	rBV3	66224	97231	0.60%	0.055%
59	12.787	1663	1666	1673	rVB3	59205	93776	0.58%	0.053%
60	12.981	1690	1699	1706	rVB	257487	429428	2.66%	0.242%
61	13.122	1718	1723	1732	rVB5	79284	183325	1.14%	0.103%
62	13.304	1750	1754	1767	rVV4	150964	320182	1.99%	0.181%
63	13.687	1814	1819	1824	rVV	79717	134105	0.83%	0.076%
64	13.751	1824	1830	1841	rVV	6751605	9929504	61.62%	5.605%
65	14.028	1868	1877	1886	rBV	7094337	10654631	66.12%	6.014%
66	14.098	1886	1889	1893	rVB	67687	80627	0.50%	0.046%
67	14.151	1893	1898	1903	rBV	104363	175343	1.09%	0.099%
68	14.340	1918	1930	1937	rBV3	5070130	8427130	52.30%	4.757%
69	14.540	1958	1964	1974	rBV	642737	1094755	6.79%	0.618%
70	14.769	1999	2003	2008	rVV2	89203	119317	0.74%	0.067%
71	14.822	2008	2012	2016	rVB	105736	137311	0.85%	0.078%

LSC Area Percent Report

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Operator : JU/SJ
Sample : J3776-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAZQ9

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-BN061918.M
Title : SVOA CALIBRATION

72	15.087	2053	2057	2064	rVV	153740	264861	1.64%	0.149%
73	15.169	2064	2071	2079	rVB2	480126	863576	5.36%	0.487%
74	15.334	2092	2099	2106	rBV	9876744	14298376	88.73%	8.070%
75	15.451	2112	2119	2130	rBV	3386237	4703045	29.19%	2.655%
76	15.569	2135	2139	2150	rVB	154558	260721	1.62%	0.147%
77	15.992	2207	2211	2219	rBV4	40751	82425	0.51%	0.047%
78	16.869	2356	2360	2369	rVB3	49441	90566	0.56%	0.051%
79	17.081	2390	2396	2406	rVB2	5977507	8664310	53.77%	4.890%
80	17.181	2406	2413	2424	rBV2	9990421	14855302	92.19%	8.385%
81	17.381	2442	2447	2453	rBV	432940	544882	3.38%	0.308%
82	17.992	2546	2551	2559	rBV	650128	1008614	6.26%	0.569%
83	18.063	2559	2563	2570	rVB	118495	190523	1.18%	0.108%
84	18.739	2675	2678	2685	rVB	155771	214206	1.33%	0.121%
85	19.116	2737	2742	2747	rBV	137741	210753	1.31%	0.119%
86	19.280	2765	2770	2780	rBV	1051884	1648067	10.23%	0.930%
87	19.369	2781	2785	2792	rVV	153156	308314	1.91%	0.174%
88	19.428	2793	2795	2798	rVV3	65535	83178	0.52%	0.047%
89	19.480	2798	2804	2813	rVB	11102632	16114106	100.00%	9.095%
90	20.427	2961	2965	2969	rVB	81544	102033	0.63%	0.058%
91	21.280	3104	3110	3116	rBV	6285525	7854530	48.74%	4.433%
92	21.451	3137	3139	3144	rVB2	66072	74011	0.46%	0.042%
93	21.892	3211	3214	3217	rBV	99068	100833	0.63%	0.057%
94	22.357	3289	3293	3299	rVB2	184276	266273	1.65%	0.150%
95	22.863	3375	3379	3385	rVB	260932	388549	2.41%	0.219%
96	23.374	3458	3466	3472	rBV	6371878	12019596	74.59%	6.784%
97	23.427	3472	3475	3481	rVV	310589	476097	2.95%	0.269%
98	23.510	3482	3489	3500	rVB	3572307	6832102	42.40%	3.856%
99	24.074	3580	3585	3590	rVB	288032	500181	3.10%	0.282%
100	24.815	3707	3711	3722	rVB2	253143	542785	3.37%	0.306%

Sum of corrected areas: 177169901

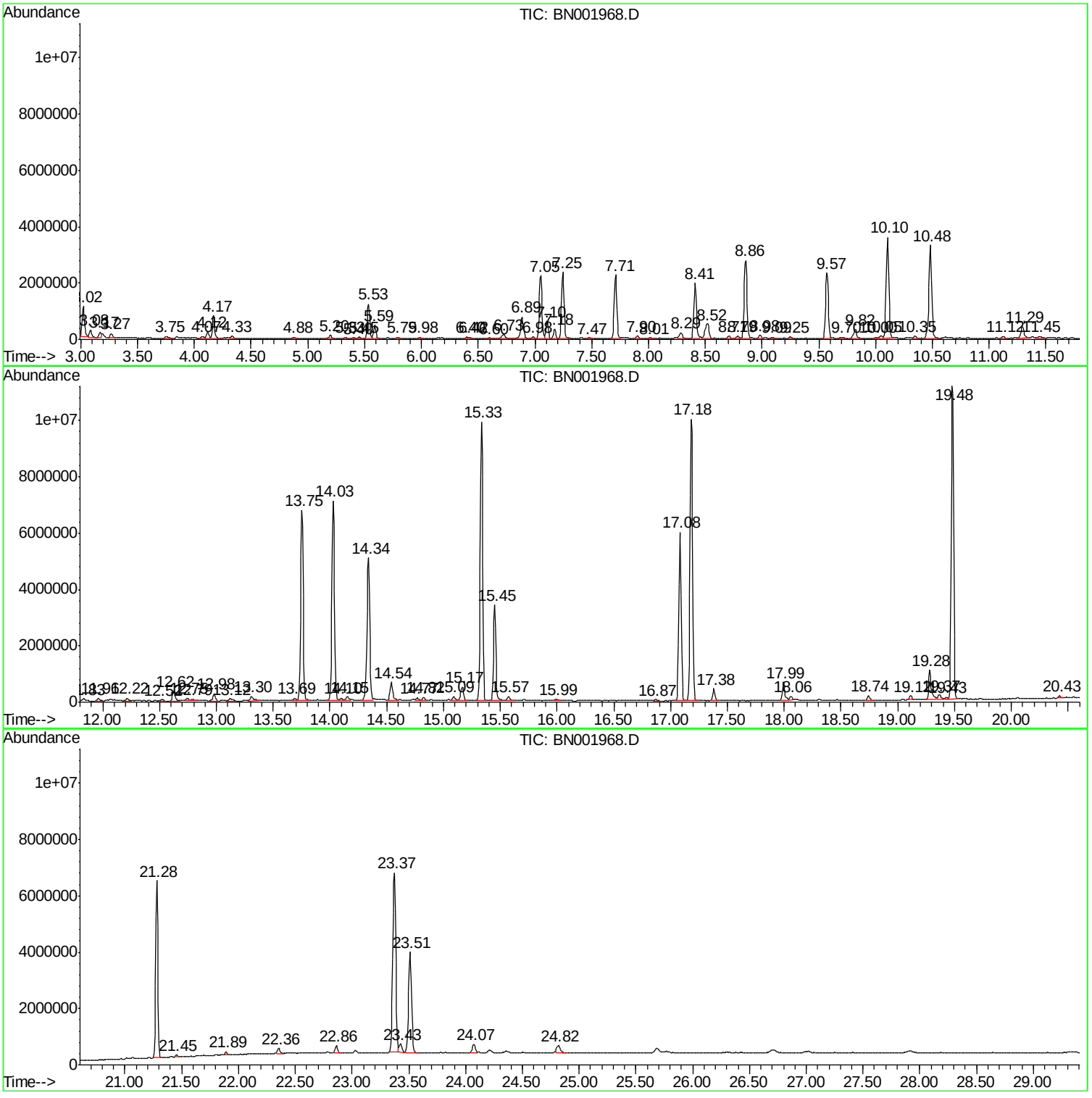
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN070518\
Data File : BN001968.D
Acq On : 05 Jul 2018 15:34
Operator : JU/SJ
Sample : J3776-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DAZQ9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN070518\
Data File : BN001968.D
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Operator : JU/SJ
Sample : J3776-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAZQ9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

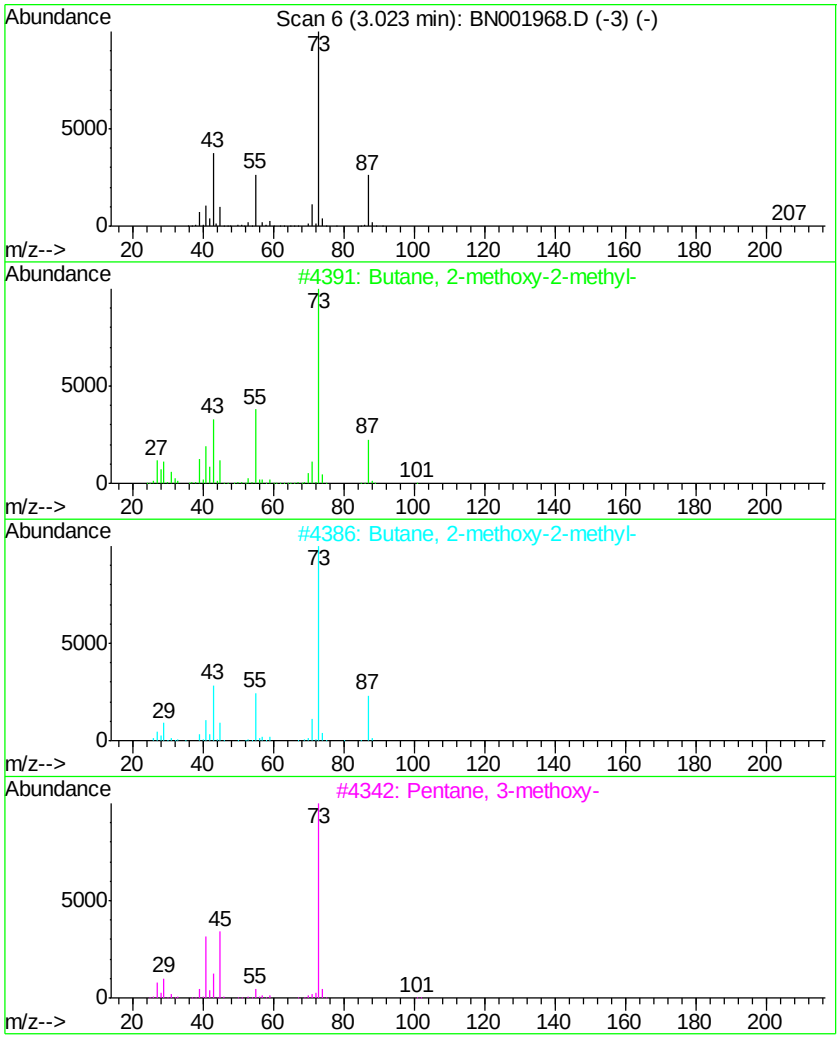
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

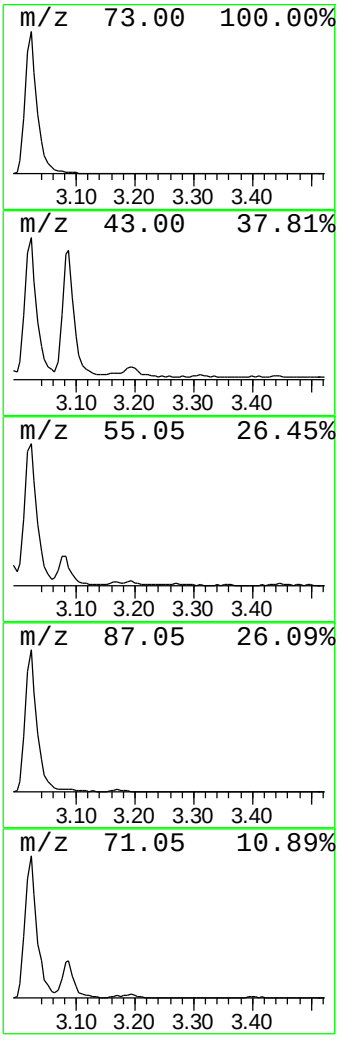
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.02	7.55 ng/ul	1403760	1,4-Dichlorobenzene-d4	7.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4	N-Ethylformamide	73	C3H7NO	000627-45-2	9
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9





Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN070518\
Data File : BN001968.D
Acq On : 05 Jul 2018 15:34
Operator : JU/SJ
Sample : J3776-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAZQ9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

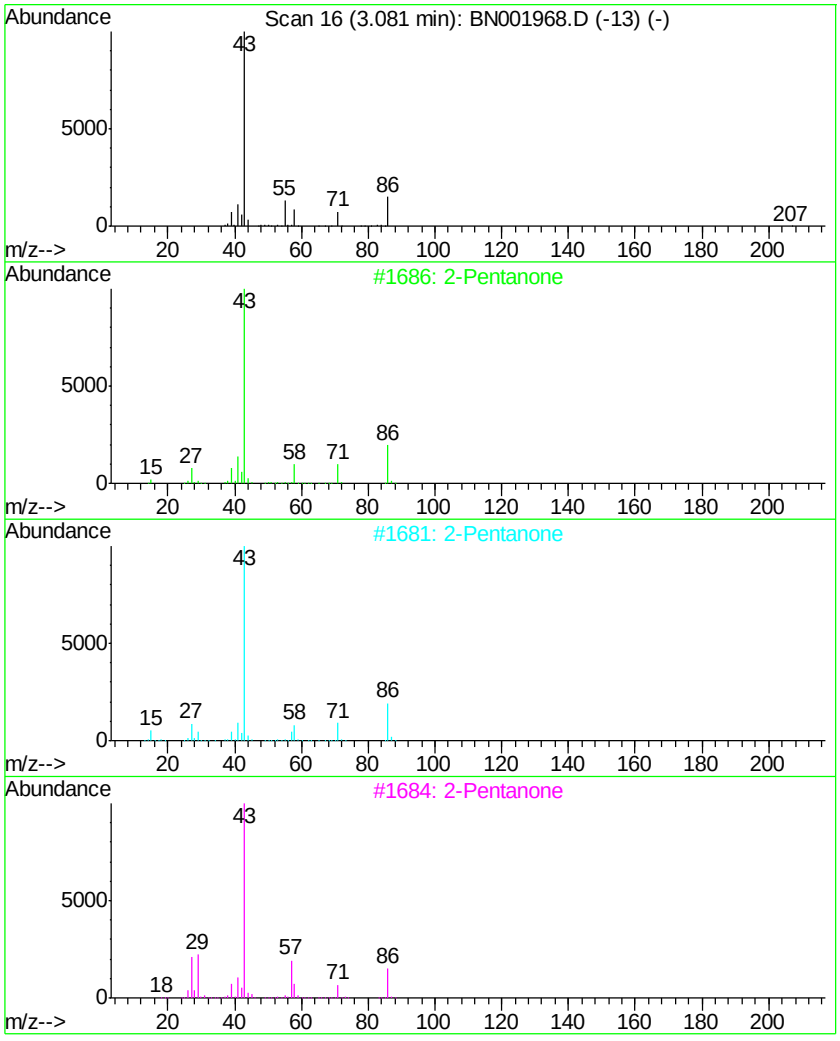
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

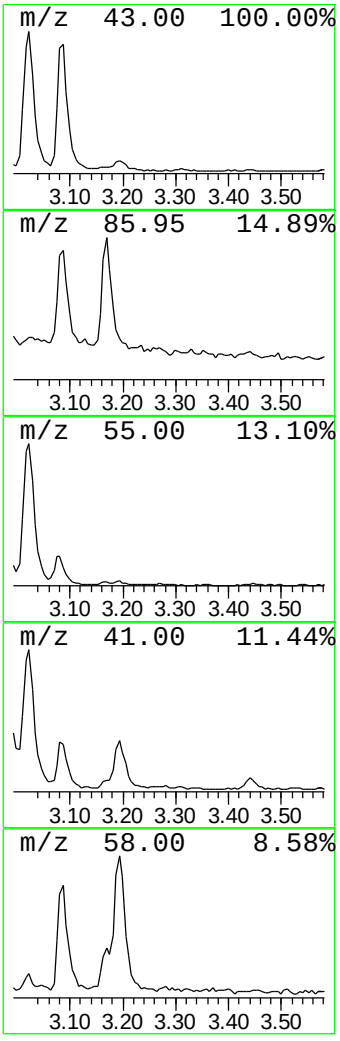
****

Peak Number 2 2-Pentanone Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.08	2.10 ng/ul	390009	1,4-Dichlorobenzene-d4	7.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone	86	C5H10O	000107-87-9	72
2	2-Pentanone	86	C5H10O	000107-87-9	64
3	2-Pentanone	86	C5H10O	000107-87-9	64
4	2,3-Butanedione	86	C4H6O2	000431-03-8	52
5	2,3-Butanedione	86	C4H6O2	000431-03-8	47





Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN070518\
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Sample : J3776-03
Misc :
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Instrument : BNA_N
ClientSampleId : DAZQ9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

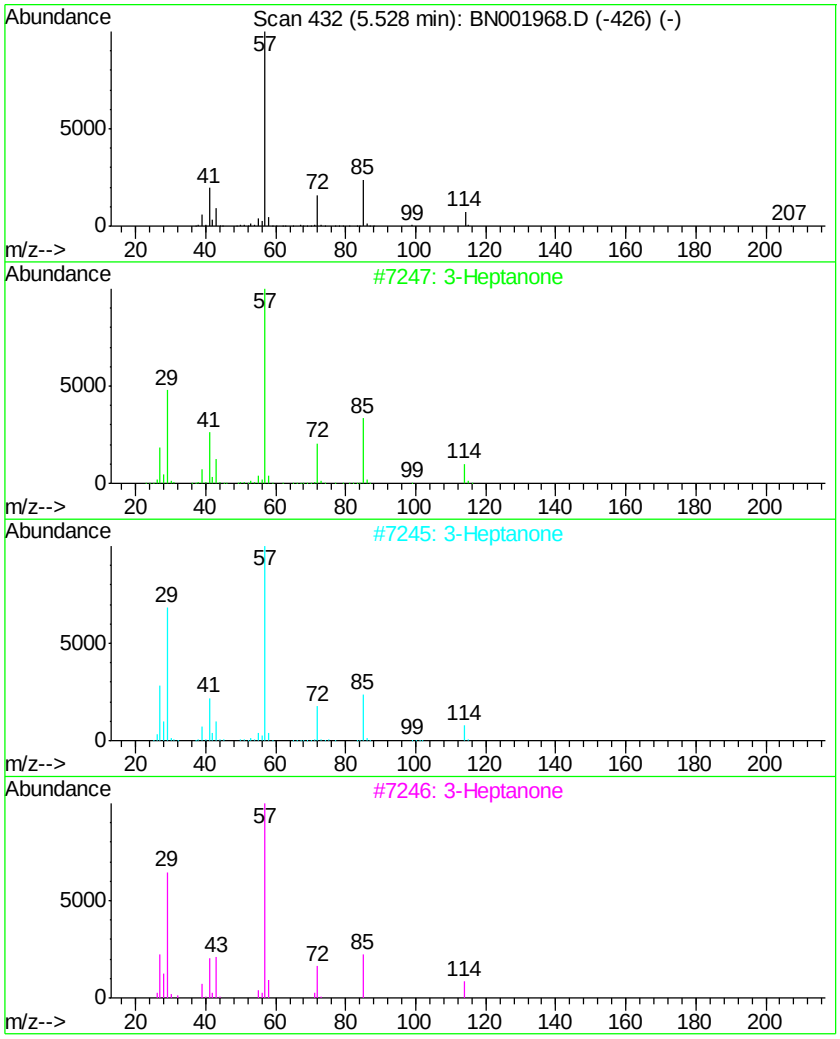
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

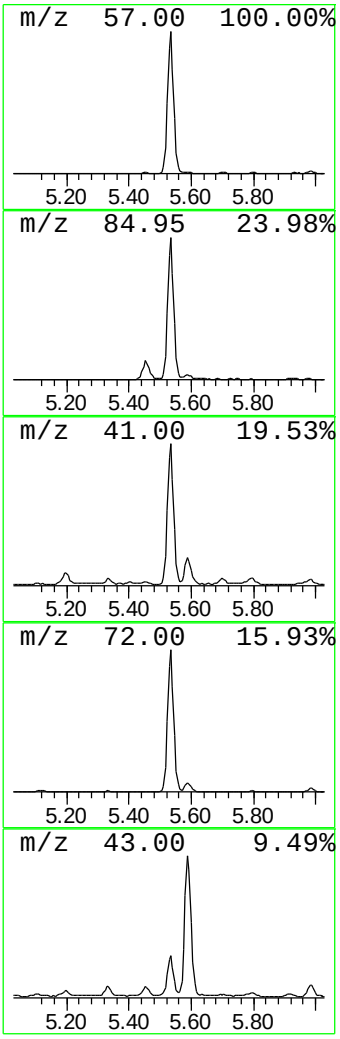
****

Peak Number 4 3-Heptanone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.53	9.20 ng/ul	1710180	1,4-Dichlorobenzene-d4	7.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Heptanone	114	C7H14O	000106-35-4	90
2	3-Heptanone	114	C7H14O	000106-35-4	87
3	3-Heptanone	114	C7H14O	000106-35-4	72
4	3-Hexanone, 5-methyl-	114	C7H14O	000623-56-3	64
5	3-Heptanone	114	C7H14O	000106-35-4	50





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Sample : J3776-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAZQ9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

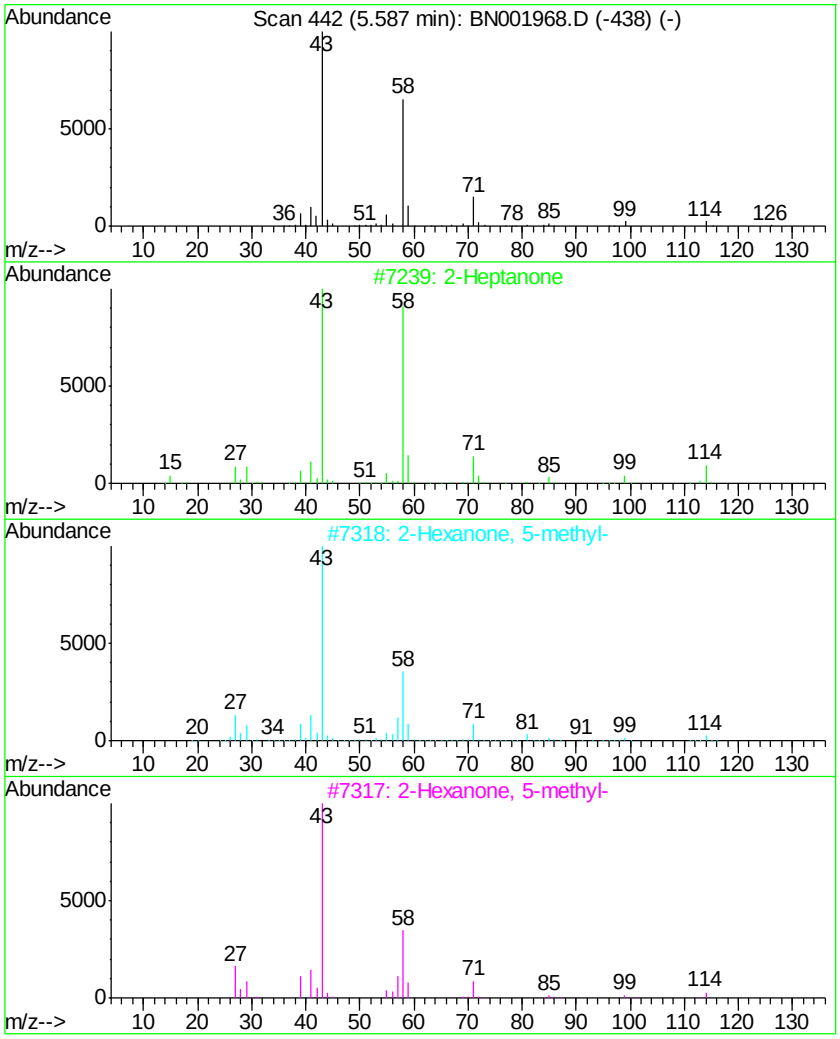
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

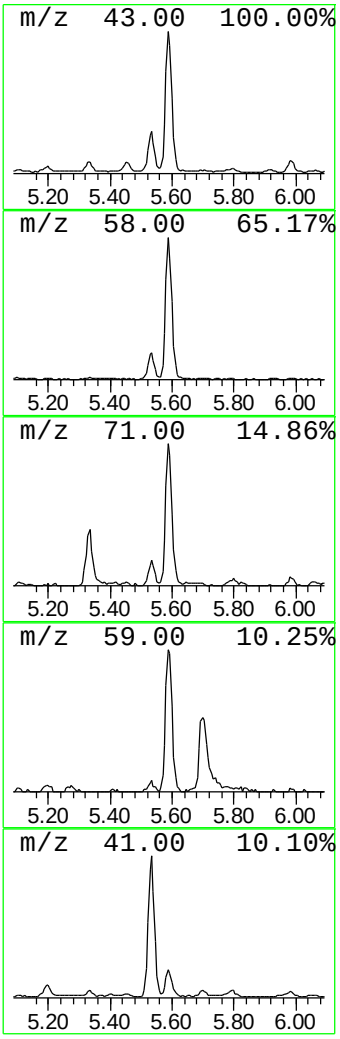
****

Peak Number 5 2-Heptanone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.59	3.71 ng/ul	689777	1,4-Dichlorobenzene-d4	7.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone	114	C7H14O	000110-43-0	78
2	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	64
3	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	64
4	2-Heptanone	114	C7H14O	000110-43-0	64
5	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	58





Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN070518\
Data File : BN001968.D
Acq On : 05 Jul 2018 15:34
Operator : JU/SJ
Sample : J3776-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAZQ9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

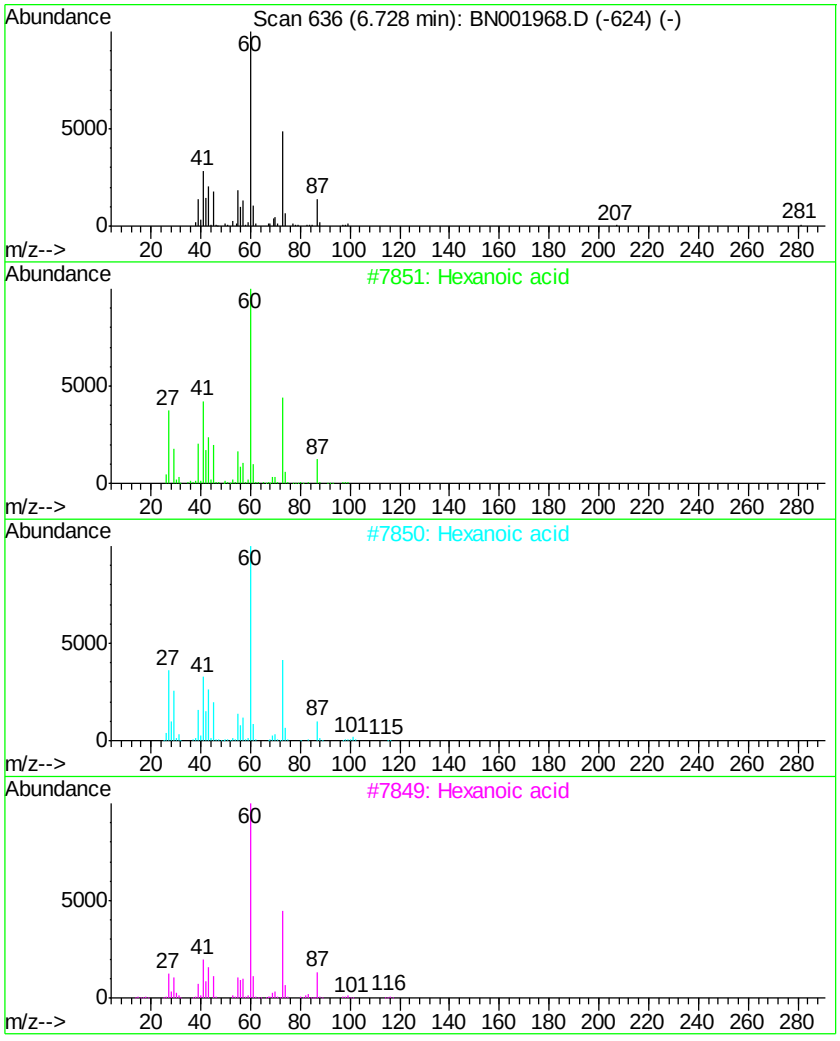
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

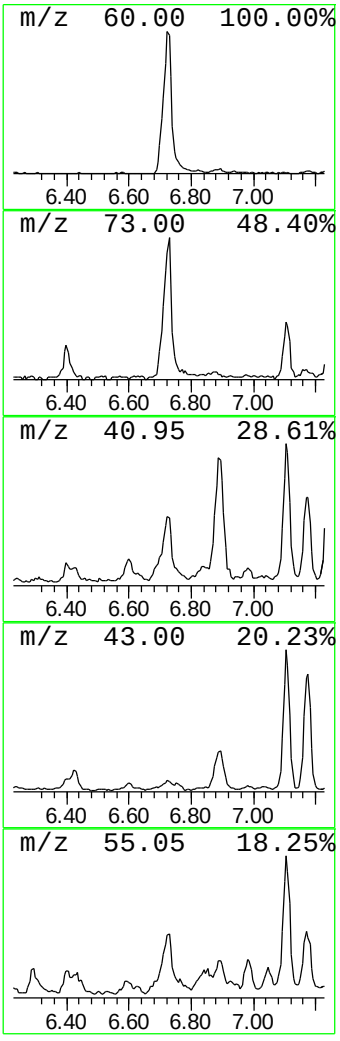
****

Peak Number 6 Hexanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	2.11 ng/ul	393113	1,4-Dichlorobenzene-d4	7.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid	116	C6H12O2	000142-62-1	90
2	Hexanoic acid	116	C6H12O2	000142-62-1	83
3	Hexanoic acid	116	C6H12O2	000142-62-1	83
4	Pentanoic acid	102	C5H10O2	000109-52-4	78
5	Octanoic Acid	144	C8H16O2	000124-07-2	64





Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN070518\
Data File : BN001968.D
Acq On : 05 Jul 2018 15:34
Operator : JU/SJ
Sample : J3776-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAZQ9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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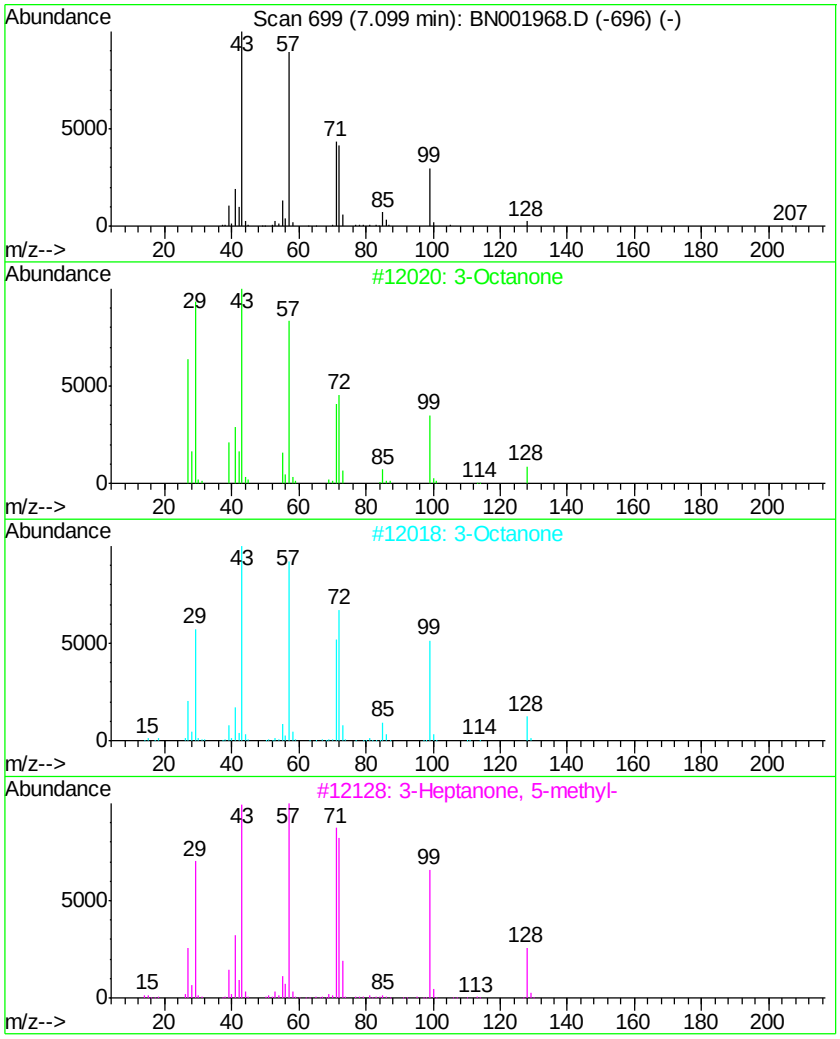
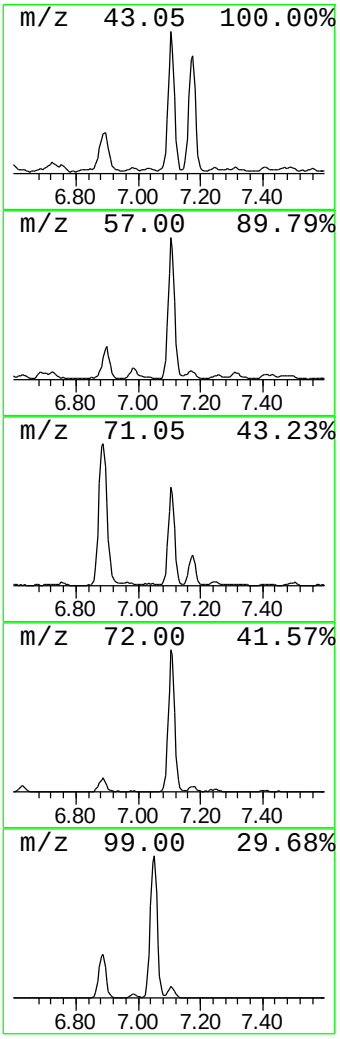
Peak Number 7 3-Octanone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.10	4.63 ng/ul	861289	1,4-Dichlorobenzene-d4	7.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Octanone	128	C8H16O	000106-68-3	78
2	3-Octanone	128	C8H16O	000106-68-3	74
3	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	56
4	3-Octanone	128	C8H16O	000106-68-3	50
5	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	50





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Sample : J3776-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAZQ9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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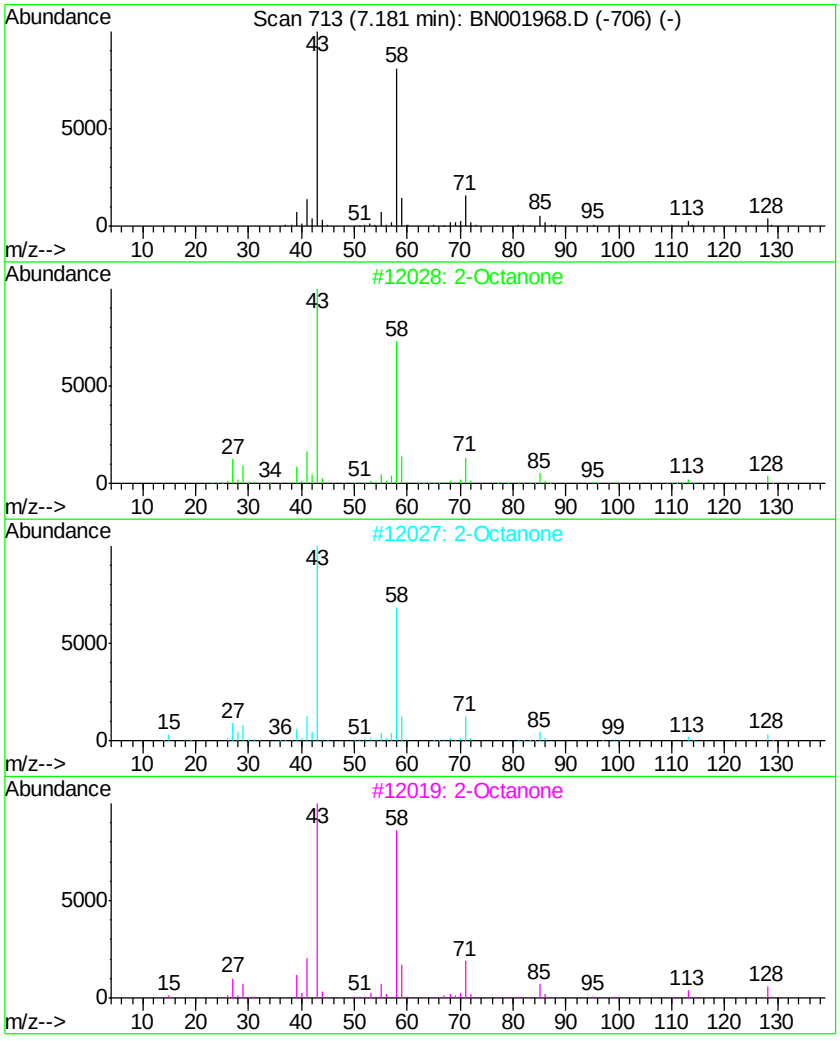
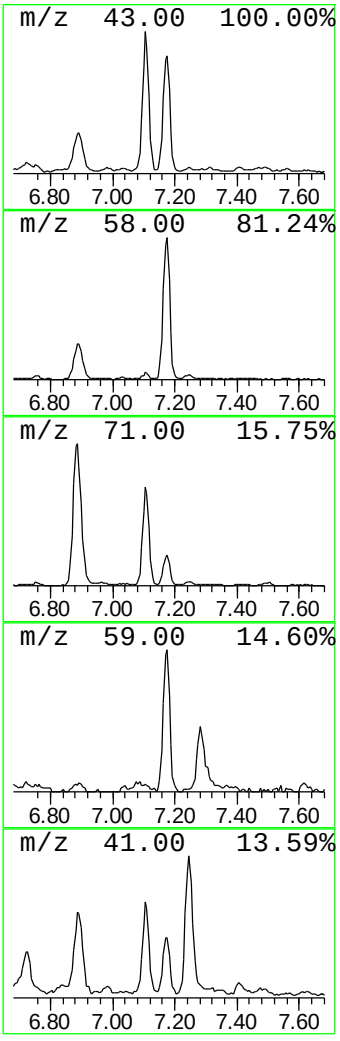
Peak Number 8 2-Octanone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.18	2.77 ng/ul	514337	1,4-Dichlorobenzene-d4	7.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Octanone	128	C8H16O	000111-13-7	94
2	2-Octanone	128	C8H16O	000111-13-7	91
3	2-Octanone	128	C8H16O	000111-13-7	91
4	2-Octanone	128	C8H16O	000111-13-7	90
5	2-Octanone	128	C8H16O	000111-13-7	78





Library Search Compound Report

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Sample : J3776-03
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Instrument : BNA_N
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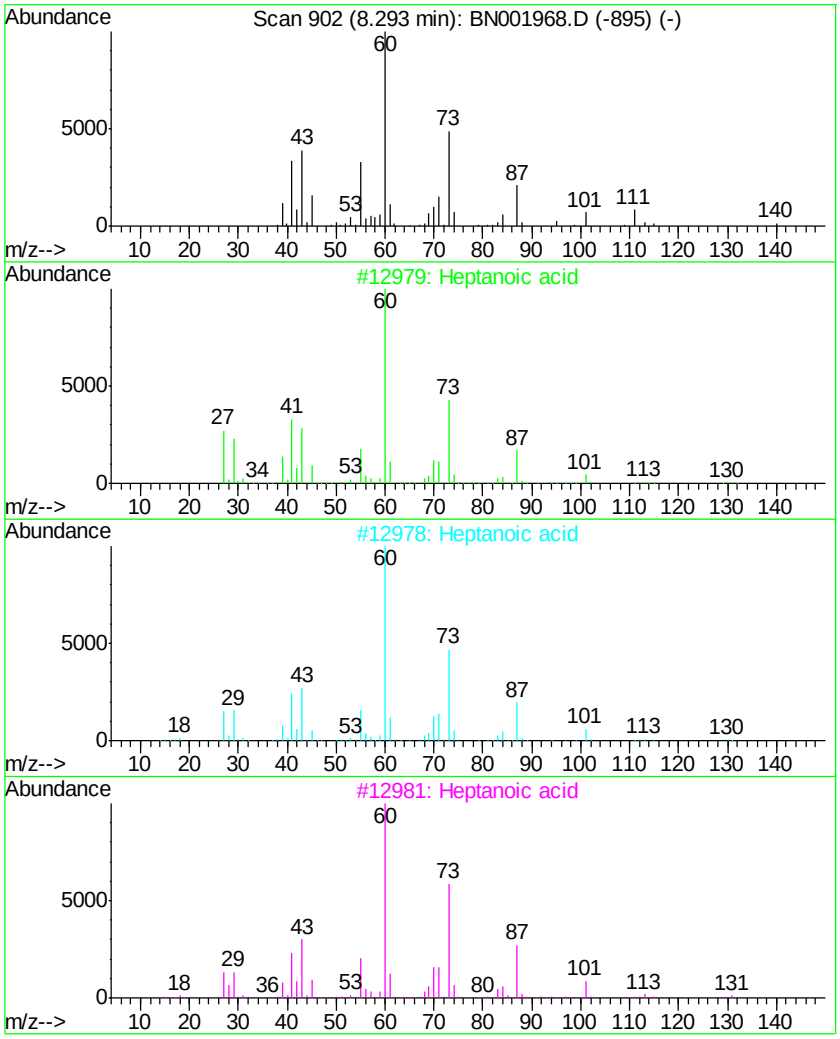
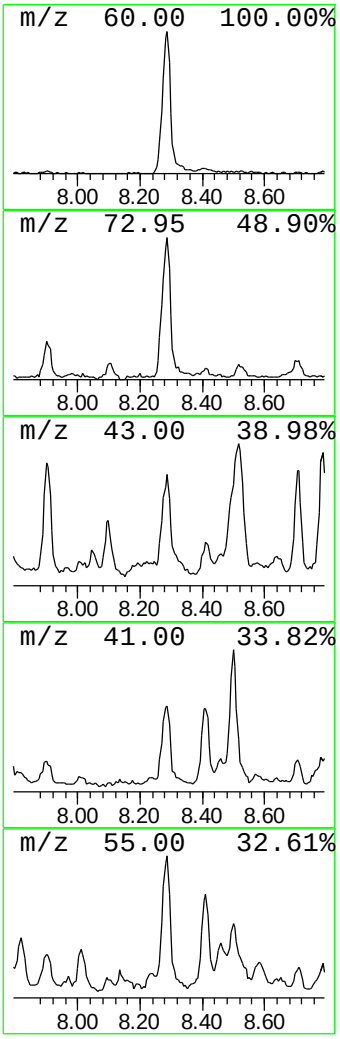
Peak Number 9 Heptanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.29	2.11 ng/ul	391524	1,4-Dichlorobenzene-d4	7.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptanoic acid	130	C7H14O2	000111-14-8	90
2	Heptanoic acid	130	C7H14O2	000111-14-8	90
3	Heptanoic acid	130	C7H14O2	000111-14-8	86
4	Hexanoic acid	116	C6H12O2	000142-62-1	59
5	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	53





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Misc :
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Instrument : BNA_N
ClientSampleId : DAZQ9

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TIC Library : C:\DATABASE\NIST02.L
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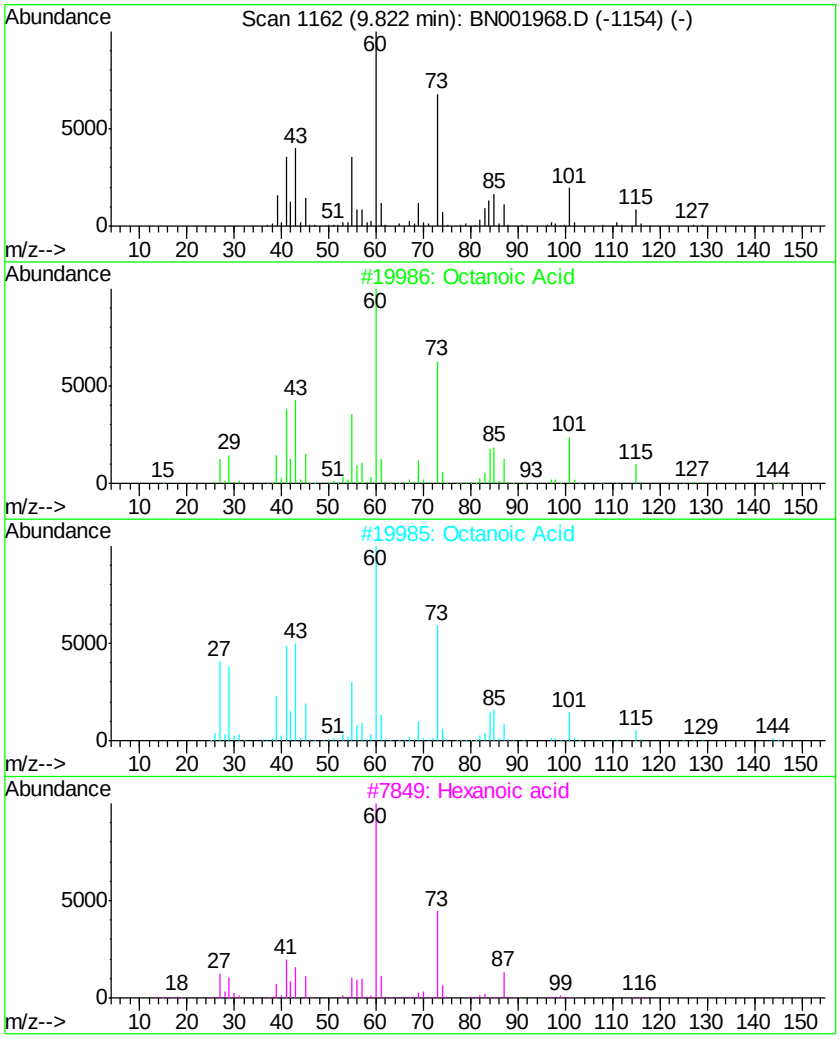
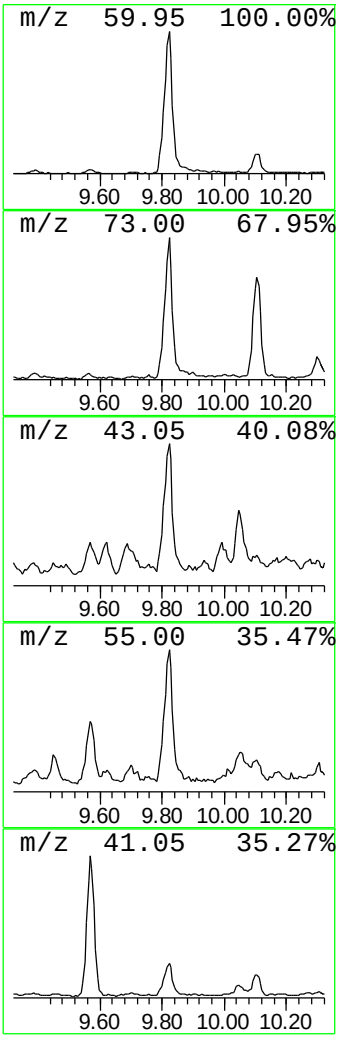
Peak Number 10 Octanoic Acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	2.12 ng/ul	633189	Naphthalene-d8	10.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanoic Acid	144	C8H16O2	000124-07-2	91
2	Octanoic Acid	144	C8H16O2	000124-07-2	90
3	Hexanoic acid	116	C6H12O2	000142-62-1	64
4	Hexanoic acid	116	C6H12O2	000142-62-1	64
5	Octanoic Acid	144	C8H16O2	000124-07-2	64





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Operator : JU/SJ
Sample : J3776-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAZQ9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
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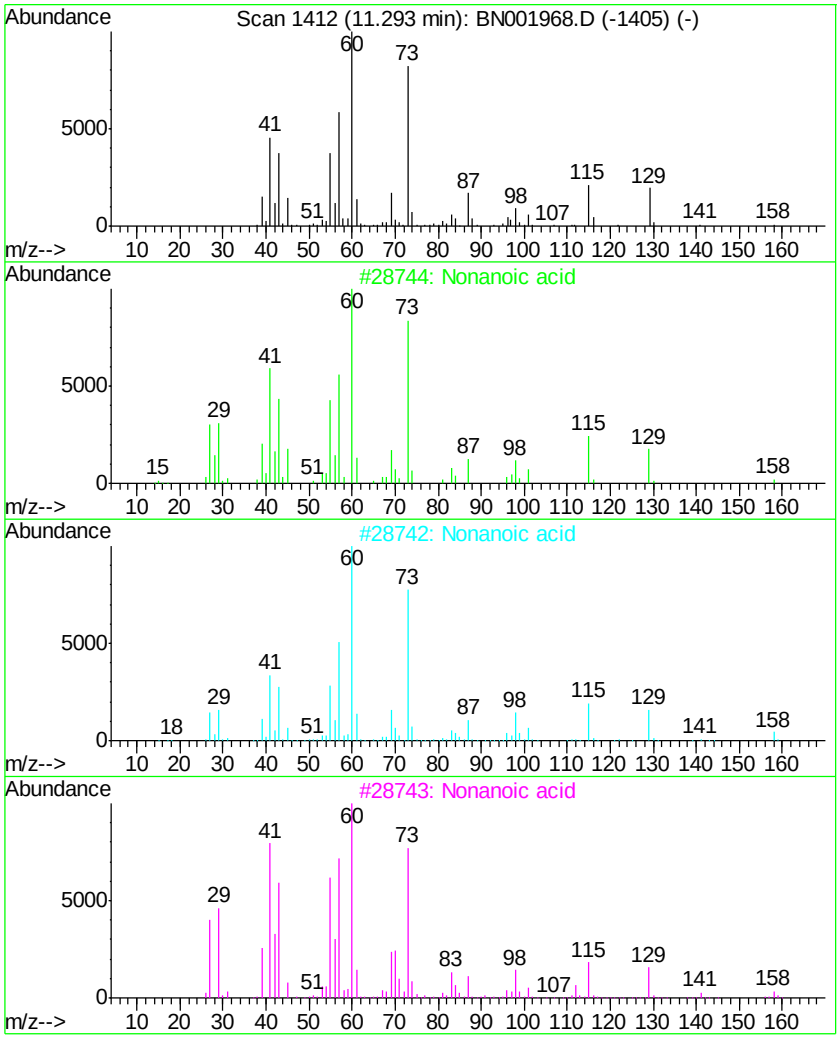
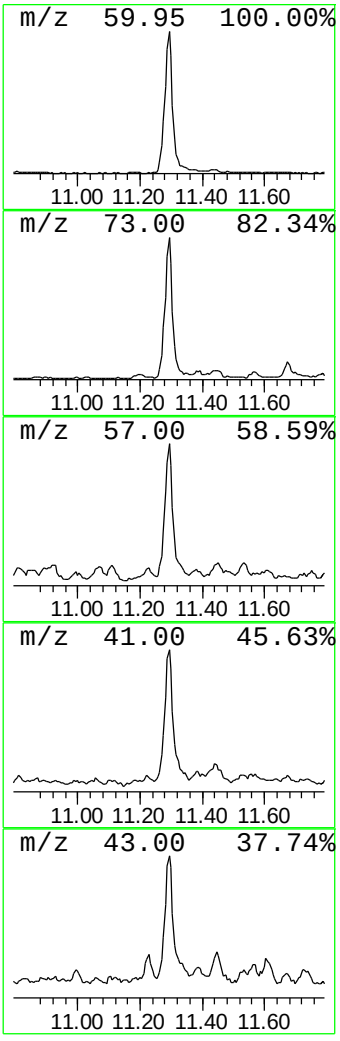
Peak Number 11 Nonanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.29	2.66 ng/ul	791608	Naphthalene-d8	10.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanoic acid	158	C9H18O2	000112-05-0	86
2	Nonanoic acid	158	C9H18O2	000112-05-0	78
3	Nonanoic acid	158	C9H18O2	000112-05-0	64
4	Hexanoic acid	116	C6H12O2	000142-62-1	47
5	Hexanoic acid	116	C6H12O2	000142-62-1	47





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Quant Title : SVOA CALIBRATION

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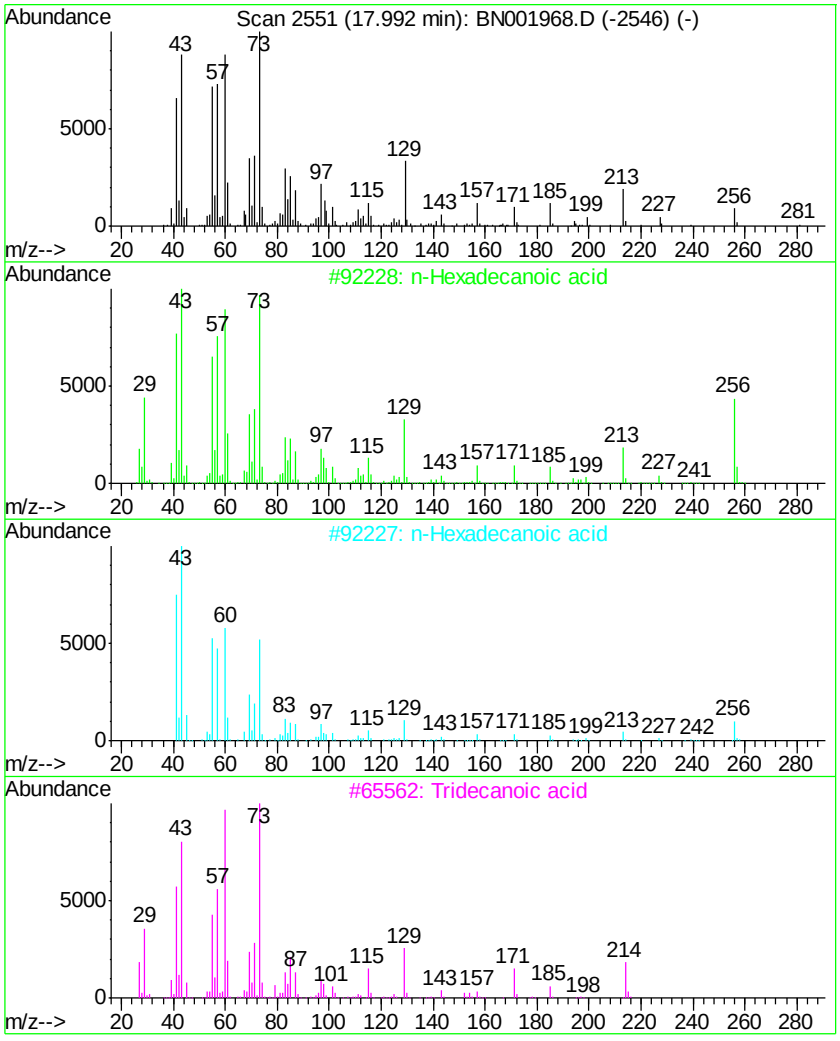
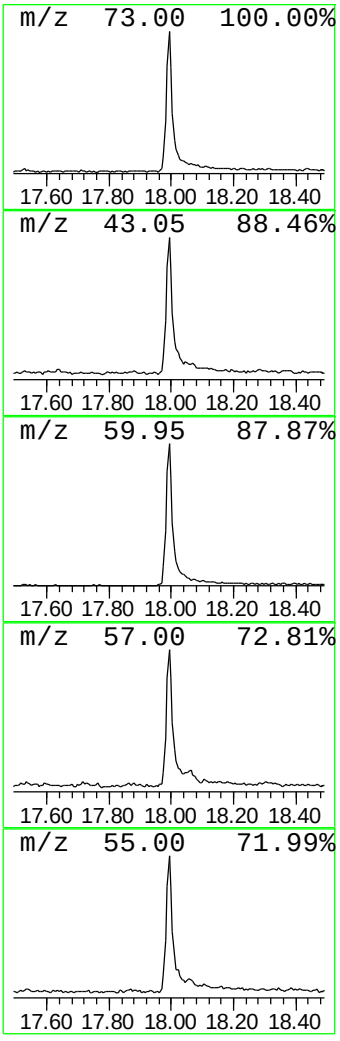
Peak Number 12 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.99	2.33 ng/ul	1008610	Phenanthrene-d10	17.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
3	Tridecanoic acid	214	C13H26O2	000638-53-9	76
4	Tridecanoic acid	214	C13H26O2	000638-53-9	76
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	74





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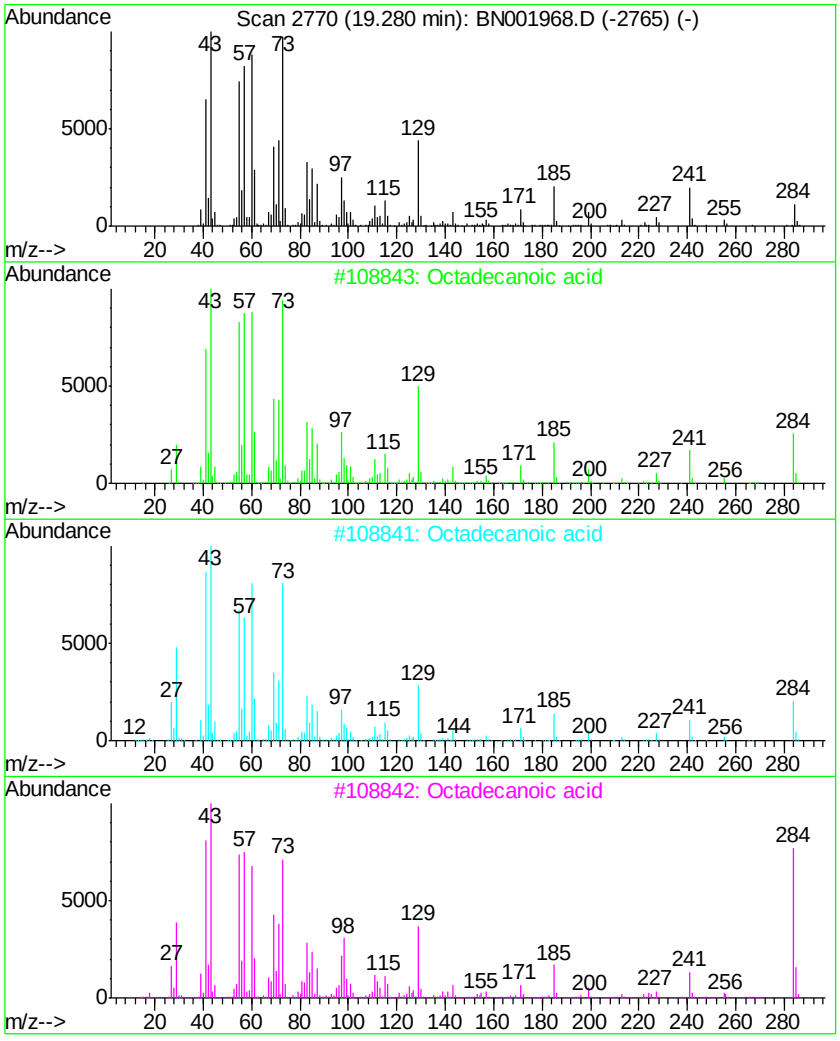
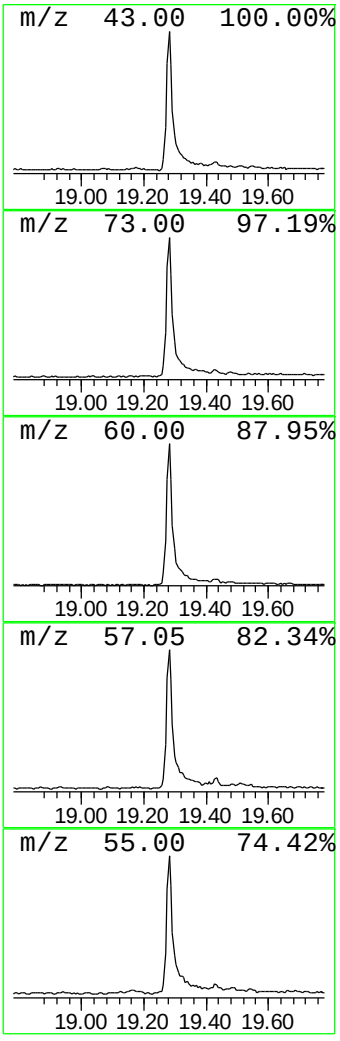
Peak Number 13 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.28	4.20 ng/ul	1648070	Chrysene-d12	21.28

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	98
3	Octadecanoic acid	284	C18H36O2	000057-11-4	95
4	Octadecanoic acid	284	C18H36O2	000057-11-4	94
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	90





Tentatively Identified Compound (LSC) summary

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methoxy...	3.02	7.5	ng/ul	1403760	1	7.71	3719250	20.0
2-Pentanone	3.08	2.1	ng/ul	390009	1	7.71	3719250	20.0
3-Heptanone	5.53	9.2	ng/ul	1710180	1	7.71	3719250	20.0
2-Heptanone	5.59	3.7	ng/ul	689777	1	7.71	3719250	20.0
Hexanoic acid	6.73	2.1	ng/ul	393113	1	7.71	3719250	20.0
3-Octanone	7.10	4.6	ng/ul	861289	1	7.71	3719250	20.0
2-Octanone	7.18	2.8	ng/ul	514337	1	7.71	3719250	20.0
Heptanoic acid	8.29	2.1	ng/ul	391524	1	7.71	3719250	20.0
Octanoic Acid	9.82	2.1	ng/ul	633189	2	10.48	5962520	20.0
Nonanoic acid	11.29	2.7	ng/ul	791608	2	10.48	5962520	20.0
n-Hexadecanoic acid	17.99	2.3	ng/ul	1008610	4	17.08	8664310	20.0
Octadecanoic acid	19.28	4.2	ng/ul	1648070	5	21.28	7854530	20.0