LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101919\
Data File : BP000919.D
Acq On : 19 Oct 2019 17:03
Operator : JU
Sample : K5354-15
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFBB8

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-BP100919MA.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	2.346	41	44	54	rBV	21426	31834	0.80%	0.094%
2	3.704	270	275	278	rBV2	3415	5336	0.13%	0.016%
3	3.740	278	281	286	rVB	5635	7829	0.20%	0.023%
4	4.363	383	387	392	rBV3	4340	5829	0.15%	0.017%
5	5.604	595	598	604	rVB	6850	6732	0.17%	0.020%
6	5.693	611	613	622	rVB	3666	5022	0.13%	0.015%
7	6.298	713	716	718	rBV	4551	5154	0.13%	0.015%
8	6.387	727	731	734	rBV	134975	150894	3.79%	0.443%
9	6.422	734	737	749	rVV	665858	551743	13.84%	1.621%
10	6.510	749	752	762	rVV	711215	605202	15.18%	1.778%
11	6.740	787	791	795	rBV	791302	641137	16.08%	1.883%
12	6.804	800	802	809	rVV	47258	45892	1.15%	0.135%
13	6.892	815	817	825	rVB3	5595	7504	0.19%	0.022%
14	7.128	854	857	871	rVV	510992	479390	12.03%	1.408%
15	7.275	879	882	883	rBV	41899	33247	0.83%	0.098%
16	7.298	883	886	893	rVB	815947	648501	16.27%	1.905%
17	7.387	898	901	906	rVB2	27025	24521	0.62%	0.072%
18	7.516	919	923	926	rVB	4450	4528	0.11%	0.013%
19	7.557	926	930	934	rBV2	3311	4069	0.10%	0.012%
20	7.622	938	941	950	rVV	627623	540645	13.56%	1.588%
21	7.875	981	984	992	rBV	961568	846434	21.23%	2.486%
22	7.987	1001	1003	1005	rBV	16072	12092	0.30%	0.036%
23	8.022	1005	1009	1012	rVV	1136530	859414	21.56%	2.524%
24	8.045	1012	1013	1017	rVV3	11234	8661	0.22%	0.025%
25	8.087	1017	1020	1028	rVV	57813	55249	1.39%	0.162%
26	8.151	1028	1031	1036	rVB2	10836	10473	0.26%	0.031%
27	8.304	1054	1057	1061	rBV	9435	9249	0.23%	0.027%
28	8.522	1091	1094	1095	rBV2	3926	4031	0.10%	0.012%
29	8.586	1102	1105	1110	rBV3	6475	7343	0.18%	0.022%
30	8.675	1114	1120	1123	rBV2	11041	12609	0.32%	0.037%
31	8.704	1123	1125	1129	rVV	17062	14785	0.37%	0.043%
32	8.910	1156	1160	1163	rBV4	3183	4031	0.10%	0.012%
33	8.981	1169	1172	1175	rBV2	5296	5514	0.14%	0.016%
34	9.022	1175	1179	1183	rVV2	3161	4520	0.11%	0.013%

LSC Area Percent Report

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Operator : JU
Sample : K5354-15
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFBB8

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-BP100919MA.M
Title : SVOA CALIBRATION

35	9.092	1187	1191	1196	rVB3	17076	17806	0.45%	0.052%
36	9.228	1204	1214	1217	rBV	744544	670011	16.81%	1.968%
37	9.481	1253	1257	1259	rBV	1445779	1217541	30.54%	3.576%
38	9.498	1259	1260	1265	rVB	118702	83419	2.09%	0.245%
39	9.628	1279	1282	1286	rBV	1774482	1587292	39.82%	4.662%
40	9.722	1293	1298	1301	rBV3	4977	6923	0.17%	0.020%
41	9.786	1305	1309	1312	rBV	1346855	1102083	27.65%	3.237%
42	9.916	1328	1331	1342	rBV	53595	80543	2.02%	0.237%
43	10.010	1342	1347	1351	rVB4	7301	14082	0.35%	0.041%
44	10.169	1371	1374	1377	rVV	25831	20693	0.52%	0.061%
45	10.204	1377	1380	1383	rVV2	11899	11240	0.28%	0.033%
46	10.310	1394	1398	1401	rBV	2215578	1891872	47.46%	5.557%
47	10.386	1407	1411	1417	rBV	392324	363672	9.12%	1.068%
48	10.433	1417	1419	1428	rVV	33352	41883	1.05%	0.123%
49	10.504	1428	1431	1439	rVB2	21024	22548	0.57%	0.066%
50	10.733	1467	1470	1473	rVB2	9923	9329	0.23%	0.027%
51	10.828	1483	1486	1489	rVB3	4959	5407	0.14%	0.016%
52	10.998	1512	1515	1522	rBV	67706	64728	1.62%	0.190%
53	11.057	1522	1525	1528	rVV	8098	7230	0.18%	0.021%
54	11.098	1529	1532	1535	rVB2	6068	5268	0.13%	0.015%
55	11.157	1540	1542	1549	rVB2	6331	7862	0.20%	0.023%
56	11.280	1560	1563	1567	rVB	1408188	1217038	30.53%	3.575%
57	11.339	1569	1573	1577	rBV	2498426	2090770	52.45%	6.141%
58	11.404	1581	1584	1587	rVB2	5402	4447	0.11%	0.013%
59	11.451	1590	1592	1597	rVB	10973	9558	0.24%	0.028%
60	11.675	1627	1630	1634	rBV	242084	195142	4.90%	0.573%
61	11.827	1653	1656	1658	rBV	8244	9271	0.23%	0.027%
62	11.857	1658	1661	1665	rVV	18561	20521	0.51%	0.060%
63	11.898	1665	1668	1676	rVV	15309	15104	0.38%	0.044%
64	12.492	1766	1769	1774	rBV	28231	26297	0.66%	0.077%
65	12.651	1792	1796	1799	rVB	22789	21298	0.53%	0.063%
66	12.722	1803	1808	1812	rBV	2568674	2526279	63.37%	7.420%
67	12.886	1833	1836	1840	rBV	7600	6361	0.16%	0.019%
68	13.127	1874	1877	1881	rBV	6716	7079	0.18%	0.021%
69	13.386	1914	1921	1930	rBV	3500514	3986425	100.00%	11.709%
70	13.457	1930	1933	1935	rVV	58268	64301	1.61%	0.189%
71	13.474	1935	1936	1943	rVB	39348	37292	0.94%	0.110%

LSC Area Percent Report

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Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFBB8

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_P\METHODS\SOM-EPA-BP100919MA.M
Title : SVOA CALIBRATION

72	13.751	1979	1983	1987	rBV	17080	18987	0.48%	0.056%
73	13.792	1987	1990	1991	rBV	23884	23902	0.60%	0.070%
74	13.816	1991	1994	2003	rVB2	136592	203353	5.10%	0.597%
75	13.957	2014	2018	2022	rBV	1535898	1323906	33.21%	3.889%
76	14.074	2031	2038	2044	rBV4	11923	20259	0.51%	0.060%
77	14.133	2044	2048	2056	rVV	383749	328588	8.24%	0.965%
78	14.445	2097	2101	2111	rBV	628552	624982	15.68%	1.836%
79	14.633	2130	2133	2136	rBV2	6987	6691	0.17%	0.020%
80	14.751	2149	2153	2163	rVB	934190	910466	22.84%	2.674%
81	14.957	2184	2188	2192	rBV	43858	46969	1.18%	0.138%
82	15.086	2206	2210	2228	rBV	892174	1014327	25.44%	2.979%
83	15.345	2246	2254	2263	rBV	2197512	2507911	62.91%	7.366%
84	15.433	2264	2269	2272	rVV	1170999	1396756	35.04%	4.103%
85	15.468	2272	2275	2290	rVB	747776	963696	24.17%	2.831%
86	15.733	2317	2320	2325	rVB	18763	26130	0.66%	0.077%
87	15.786	2326	2329	2335	rVB	12300	18539	0.47%	0.054%
88	15.904	2343	2349	2360	rBV	499697	774689	19.43%	2.275%
89	16.210	2398	2401	2407	rVB2	15057	22787	0.57%	0.067%
90	16.410	2428	2435	2450	rBV	252025	467492	11.73%	1.373%
91	16.998	2529	2535	2545	rVB	77464	158314	3.97%	0.465%
92	17.698	2648	2654	2662	rVB2	24676	58589	1.47%	0.172%

Sum of corrected areas: 34045362

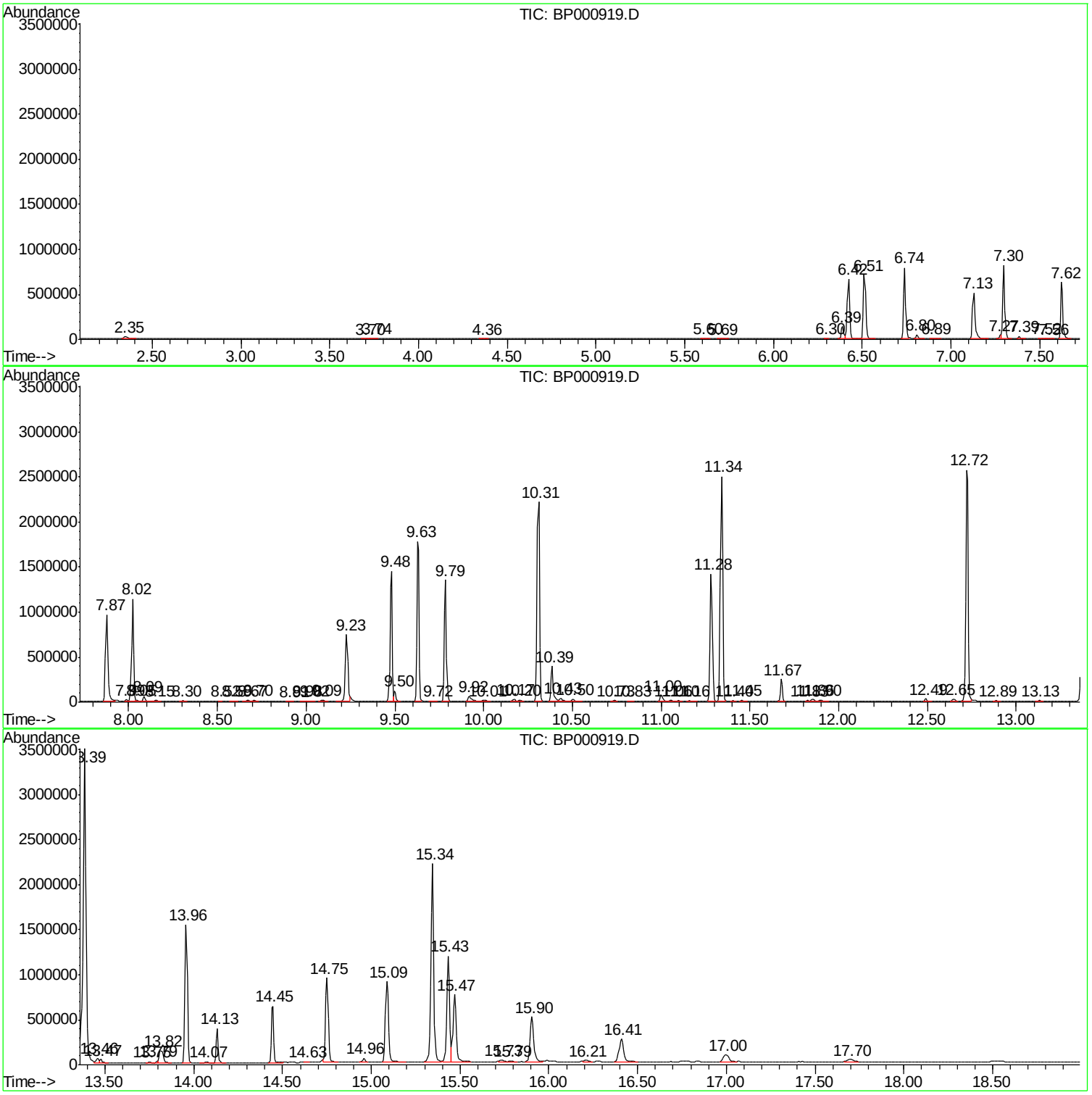
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101919\
Data File : BP000919.D
Acq On : 19 Oct 2019 17:03
Operator : JU
Sample : K5354-15
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFBB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101919\
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Quant Title : SVOA CALIBRATION

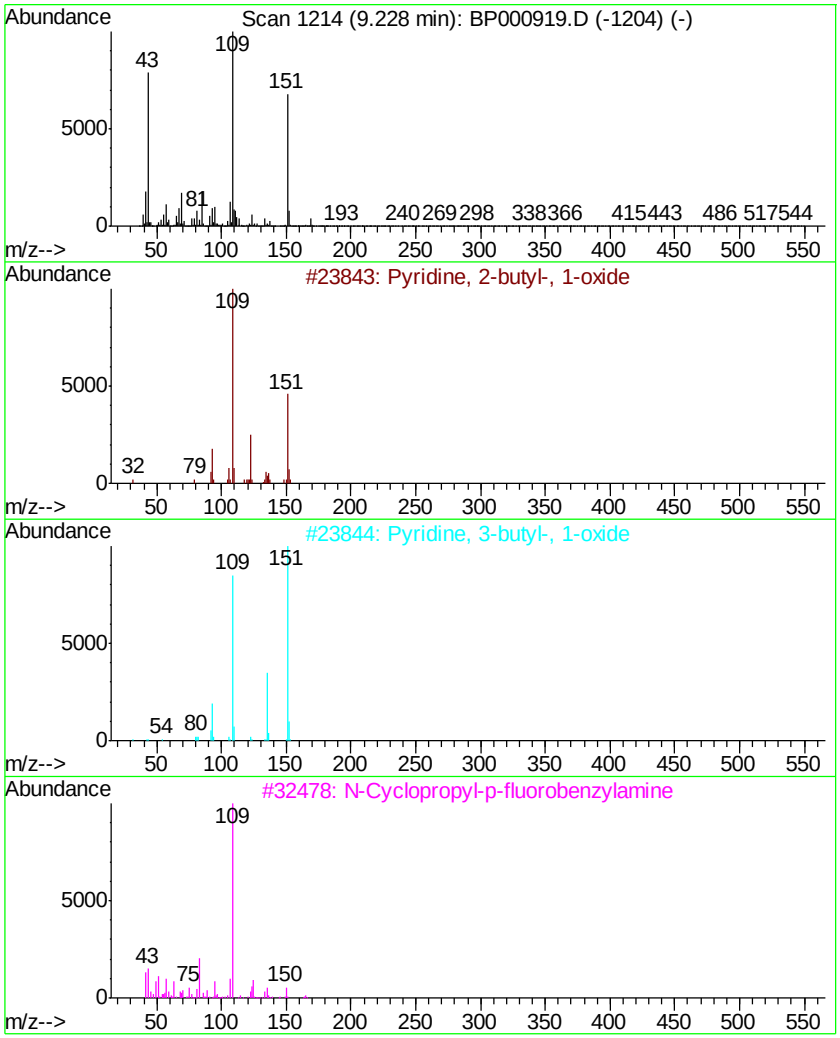
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

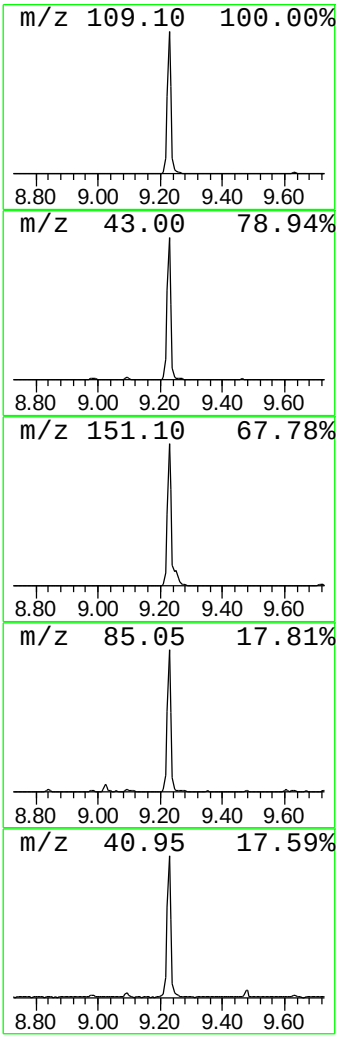
****

Peak Number 1 Pyridine, 2-butyl-, 1-oxide Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.23	12.16 ng/ul	670011	Acenaphthene-d10	9.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	59
2	Pyridine, 3-butyl-, 1-oxide	151	C9H13NO	031396-33-5	53
3	N-Cyclopropyl-p-fluorobenzylamine	165	C10H12FN	1000250-88-9	47
4	3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	42
5	Ethanone, 1-(1,3-dimethyl-3-cycl...	152	C10H16O	051733-68-7	38





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Data File : BP000919.D
Acq On : 19 Oct 2019 17:03
Operator : JU
Sample : K5354-15
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFBB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
Quant Title : SVOA CALIBRATION

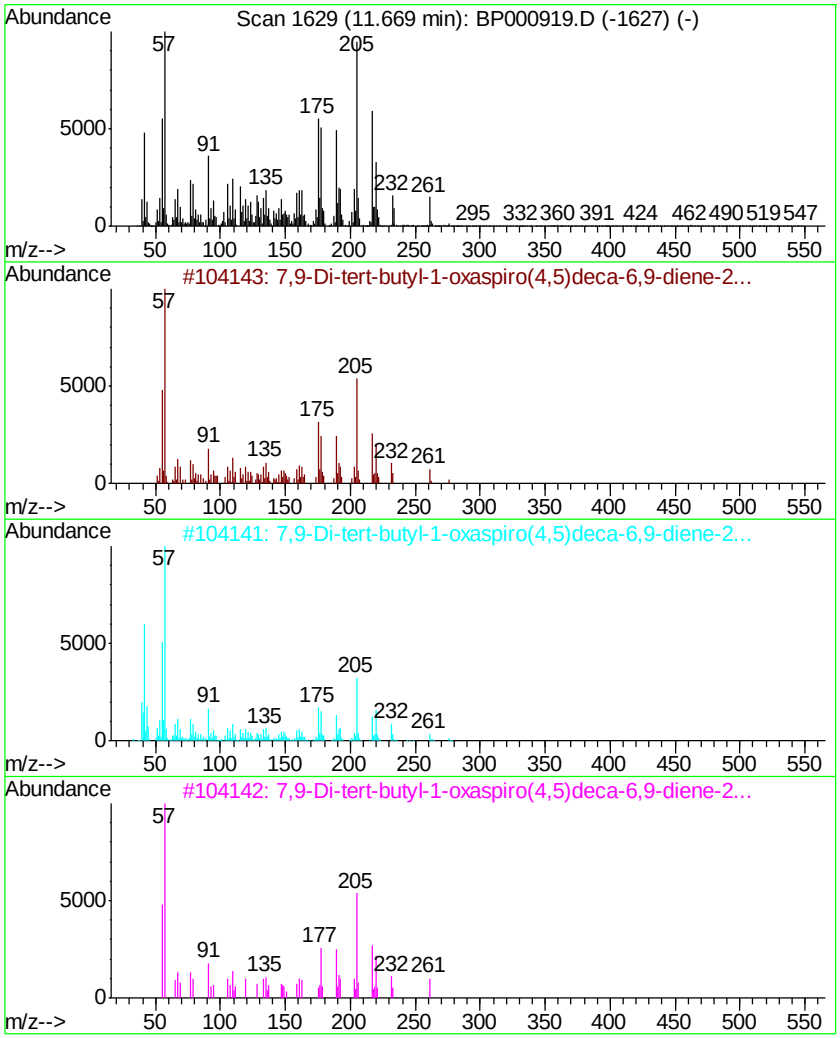
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

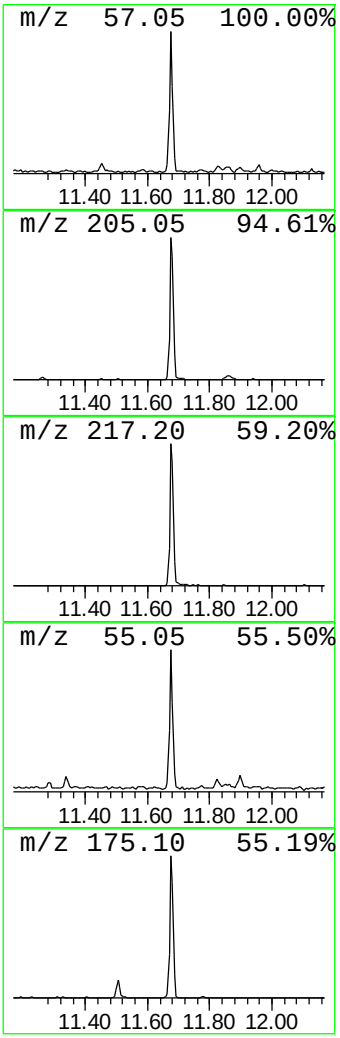
****

Peak Number 2 7,9-Di-tert-butyl-1-oxaspir... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.67	3.21 ng/ul	195142	Phenanthrene-d10	11.28

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	86
4	2,5-Cyclohexadien-1-one, 2,6-bis...	232	C16H24O	006738-27-8	83
5	Ethanone, 1-(5,6,7,8-tetrahydro-...	220	C14H20O2	071596-88-8	46





Library Search Compound Report

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
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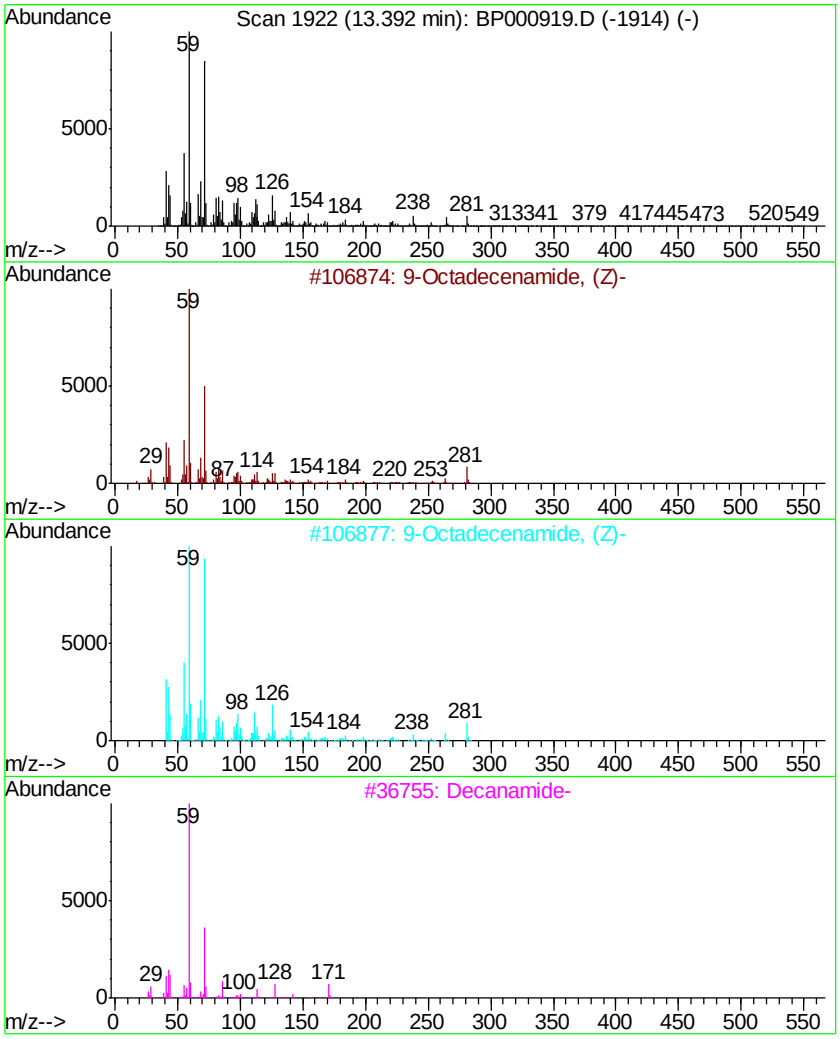
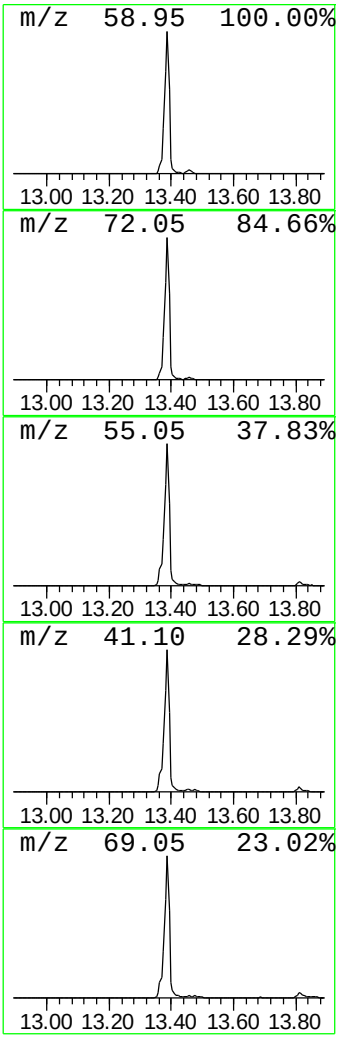
Peak Number 3 9-Octadecenamide, (Z)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.39	60.22 ng/ul	3986430	Chrysene-d12	13.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	93
2	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	70
3	Decanamide-	171	C10H21NO	002319-29-1	59
4	Pentanamide	101	C5H11NO	000626-97-1	53
5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	49





Library Search Compound Report

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Instrument : BNA_P
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
Quant Title : SVOA CALIBRATION

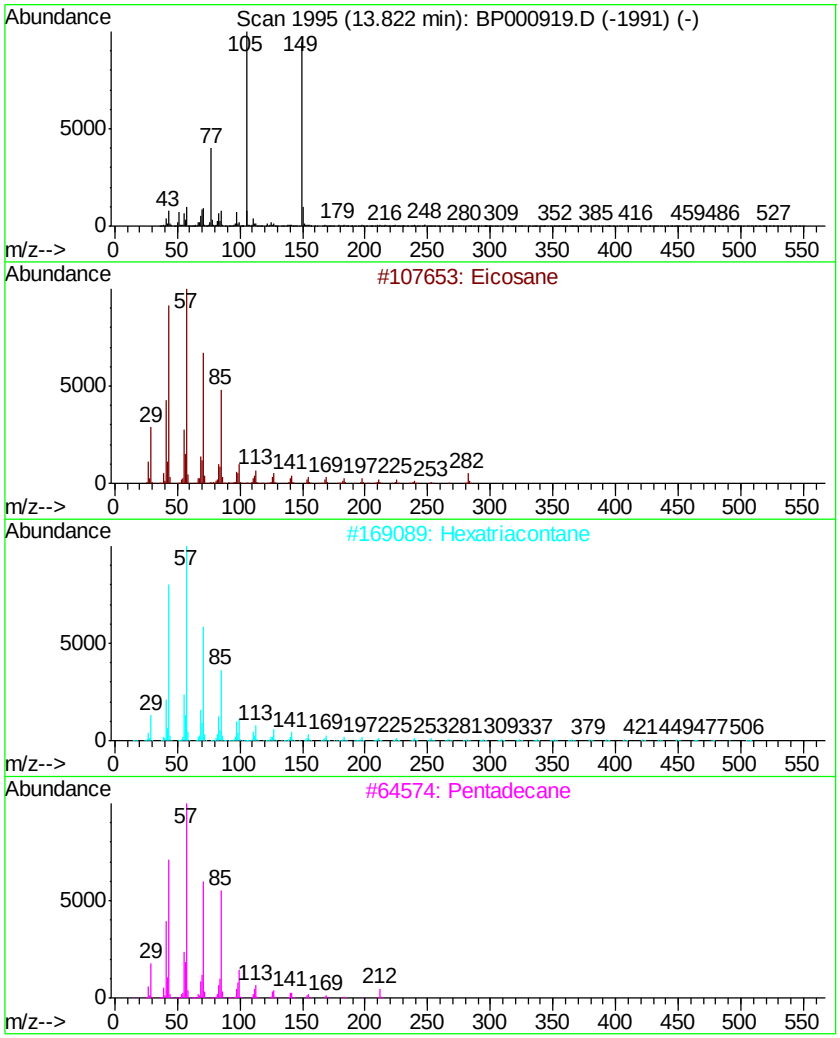
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

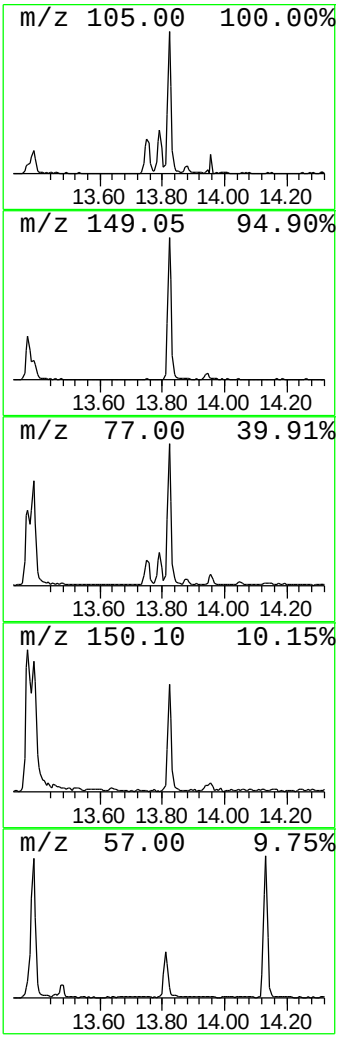
****

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	3.07 ng/ul	203353	Chrysene-d12	13.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	90
2	Hexatriacontane	507	C36H74	000630-06-8	83
3	Pentadecane	212	C15H32	000629-62-9	78
4	Tetracosane, 1-bromo-	416	C24H49Br	006946-24-3	64
5	Tetratetracontane	619	C44H90	007098-22-8	62





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Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFBB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
Quant Title : SVOA CALIBRATION

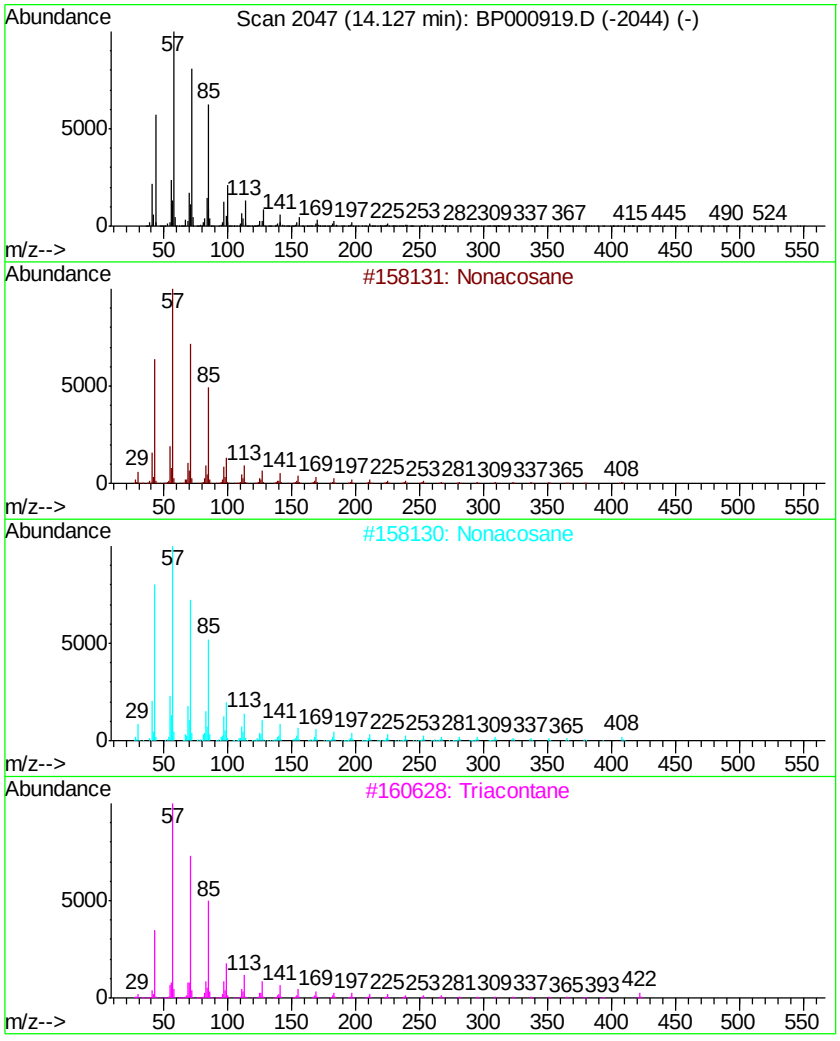
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

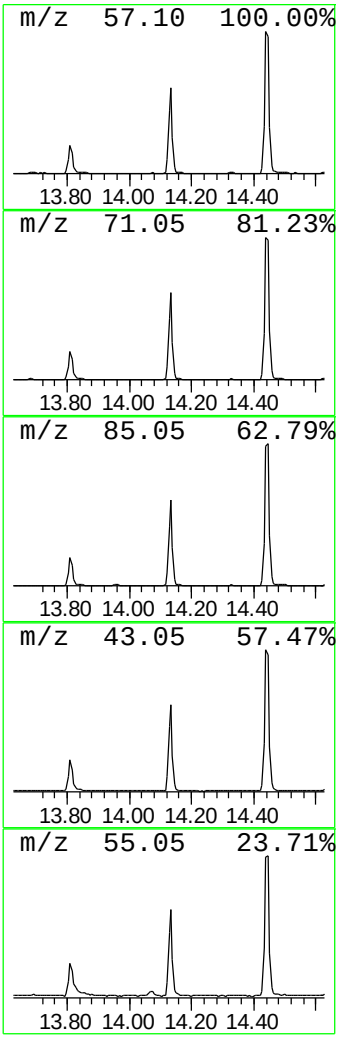
****

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.13	4.96 ng/ul	328588	Chrysene-d12	13.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonacosane	408	C29H60	000630-03-5	96
2	Nonacosane	408	C29H60	000630-03-5	94
3	Triacontane	422	C30H62	000638-68-6	94
4	Octacosane	394	C28H58	000630-02-4	93
5	Heneicosane	296	C21H44	000629-94-7	91





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ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFBB8

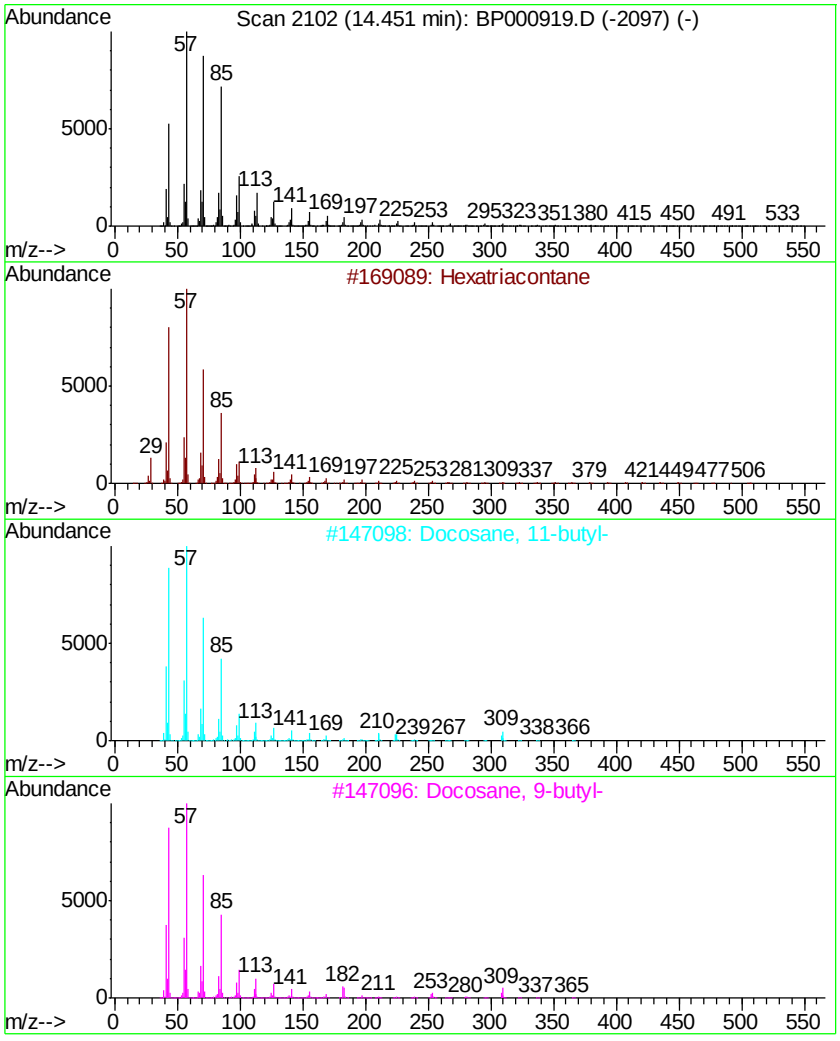
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
Quant Title : SVOA CALIBRATION

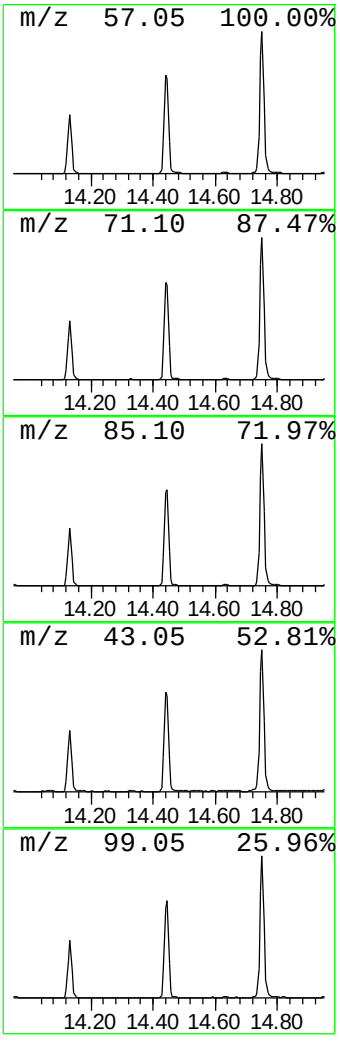
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.45	9.44 ng/ul	624982	Chrysene-d12	13.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexatriacontane	507	C36H74	000630-06-8	96
2	Docosane, 11-butyl-	366	C26H54	013475-76-8	95
3	Docosane, 9-butyl-	366	C26H54	055282-14-9	95
4	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	94
5	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101919\
Data File : BP000919.D
Acq On : 19 Oct 2019 17:03
Operator : JU
Sample : K5354-15
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFBB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
Quant Title : SVOA CALIBRATION

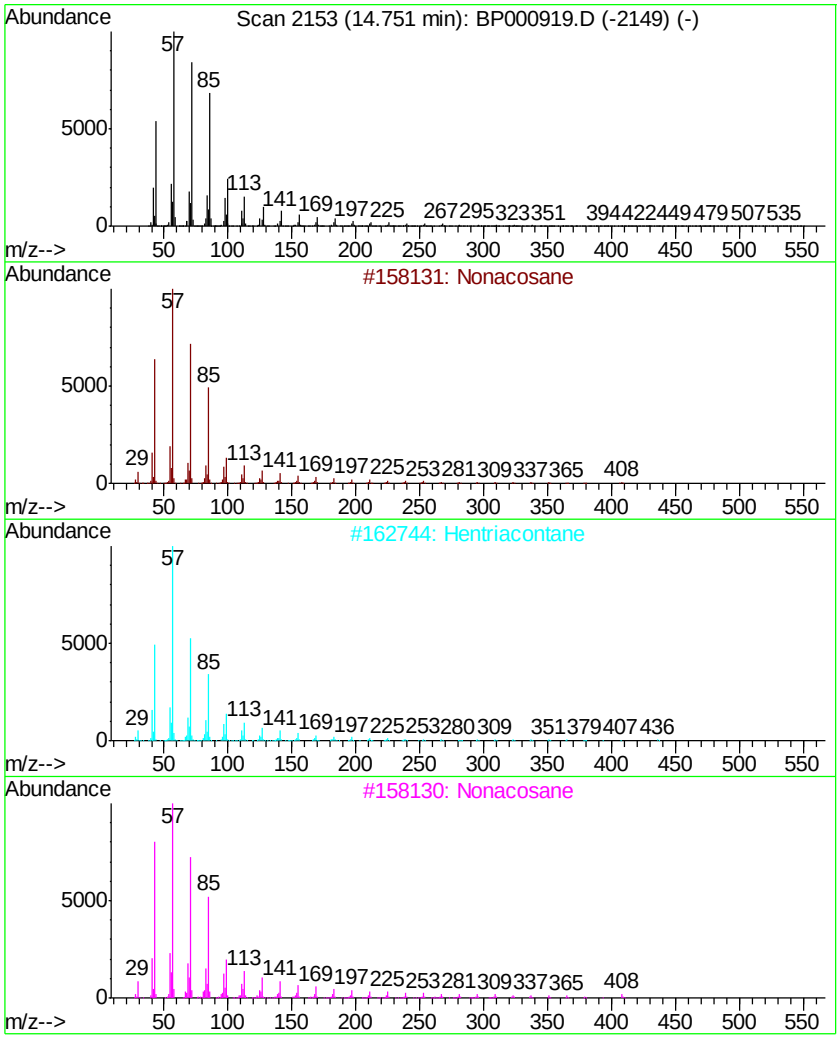
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

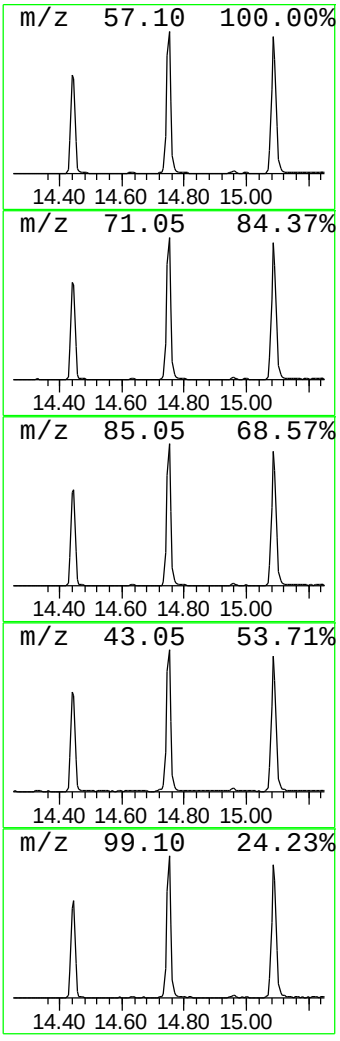
****

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.75	13.04 ng/ul	910466	Perylene-d12	15.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonacosane	408	C29H60	000630-03-5	98
2	Hentriacontane	437	C31H64	000630-04-6	96
3	Nonacosane	408	C29H60	000630-03-5	91
4	Heptacosane	380	C27H56	000593-49-7	91
5	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91





Library Search Compound Report

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Data File : BP000919.D
Acq On : 19 Oct 2019 17:03
Operator : JU
Sample : K5354-15
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFBB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
Quant Title : SVOA CALIBRATION

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

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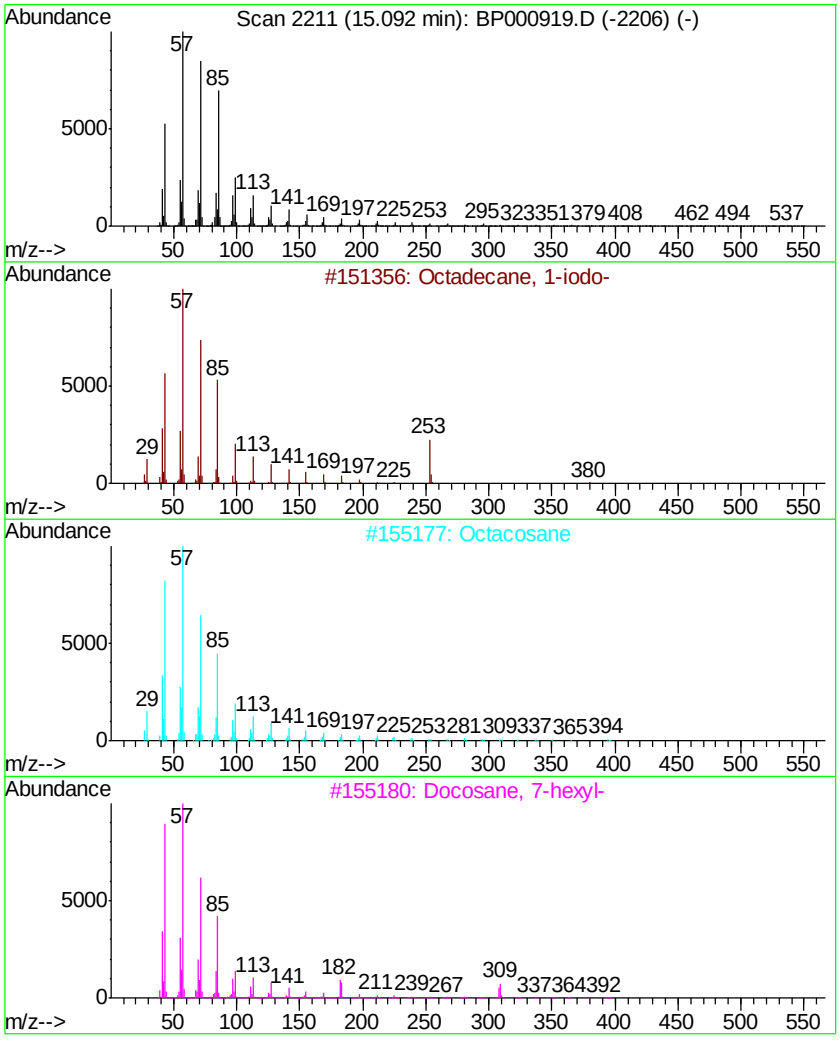
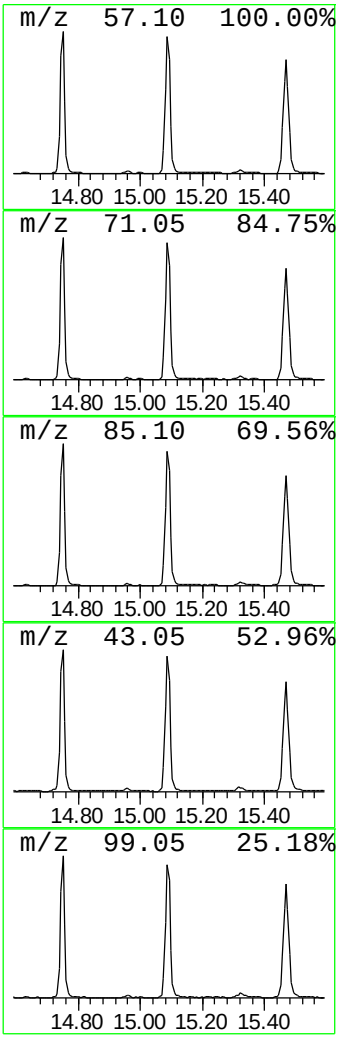
Peak Number 8 Octadecane, 1-iodo- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.09	14.52 ng/ul	1014330	Perylene-d12	15.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	97
2	Octacosane	394	C28H58	000630-02-4	96
3	Docosane, 7-hexyl-	394	C28H58	055373-86-9	94
4	Heptadecane	240	C17H36	000629-78-7	94
5	Heptadecane	240	C17H36	000629-78-7	93





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Data File : BP000919.D
Acq On : 19 Oct 2019 17:03
Operator : JU
Sample : K5354-15
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFBB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
Quant Title : SVOA CALIBRATION

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

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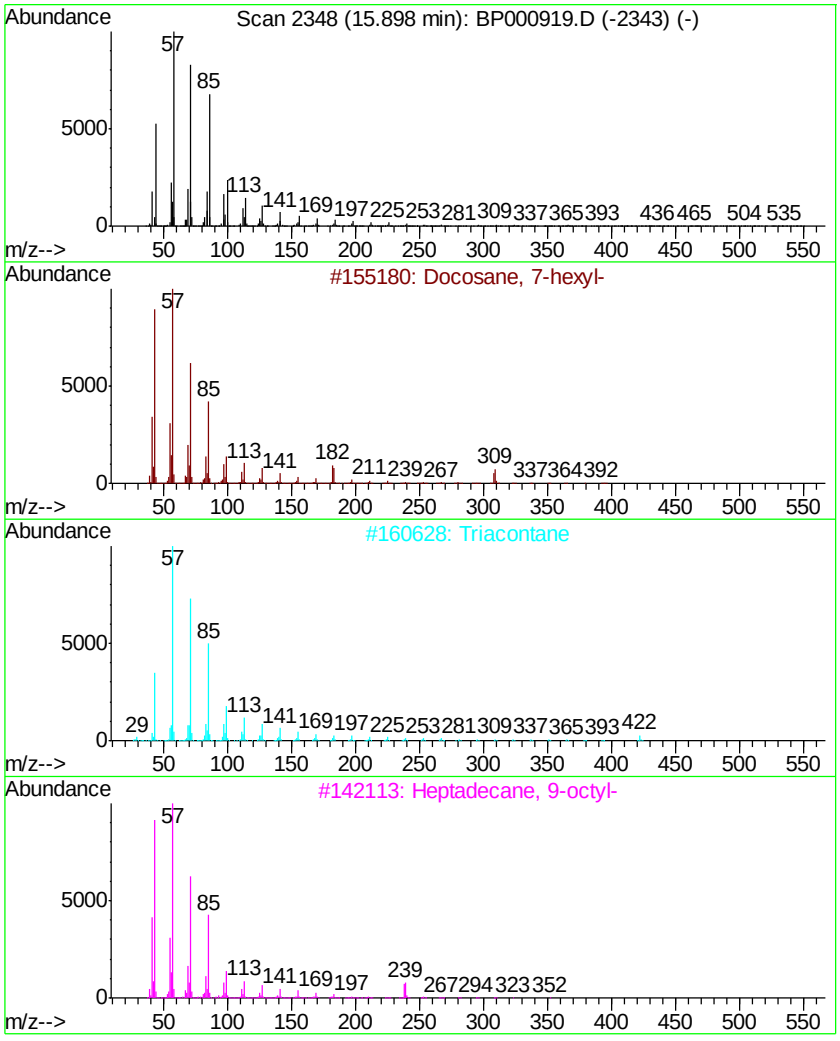
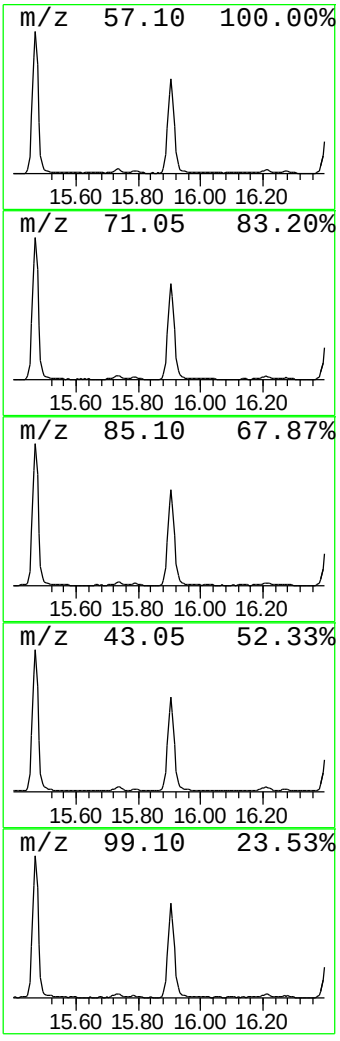
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.90	11.09 ng/ul	774689	Perylene-d12	15.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
2	Triacontane	422	C30H62	000638-68-6	91
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4	Hentriacontane	437	C31H64	000630-04-6	90
5	Tetratetracontane	619	C44H90	007098-22-8	90





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101919\
Data File : BP000919.D
Acq On : 19 Oct 2019 17:03
Operator : JU
Sample : K5354-15
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFBB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
Quant Title : SVOA CALIBRATION

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

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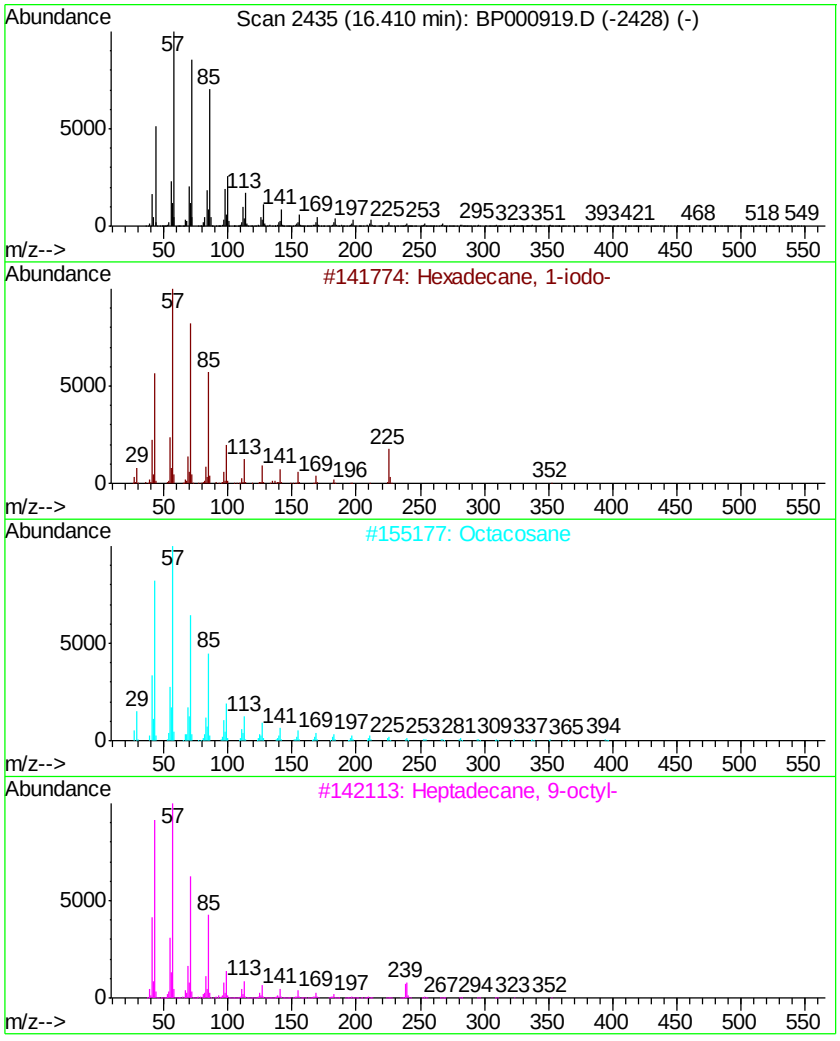
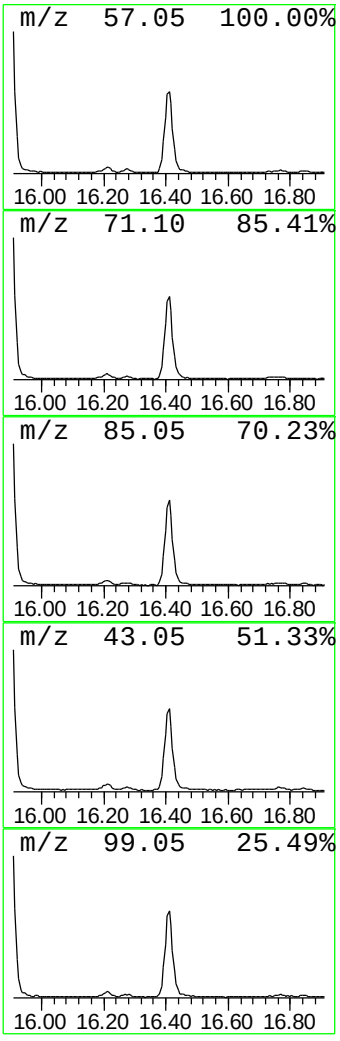
Peak Number 10 Hexadecane, 1-iodo- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.41	6.69 ng/ul	467492	Perylene-d12	15.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	91
2	Octacosane	394	C28H58	000630-02-4	91
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
4	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	90
5	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101919\
Data File : BP000919.D
Acq On : 19 Oct 2019 17:03
Operator : JU
Sample : K5354-15
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFBB8

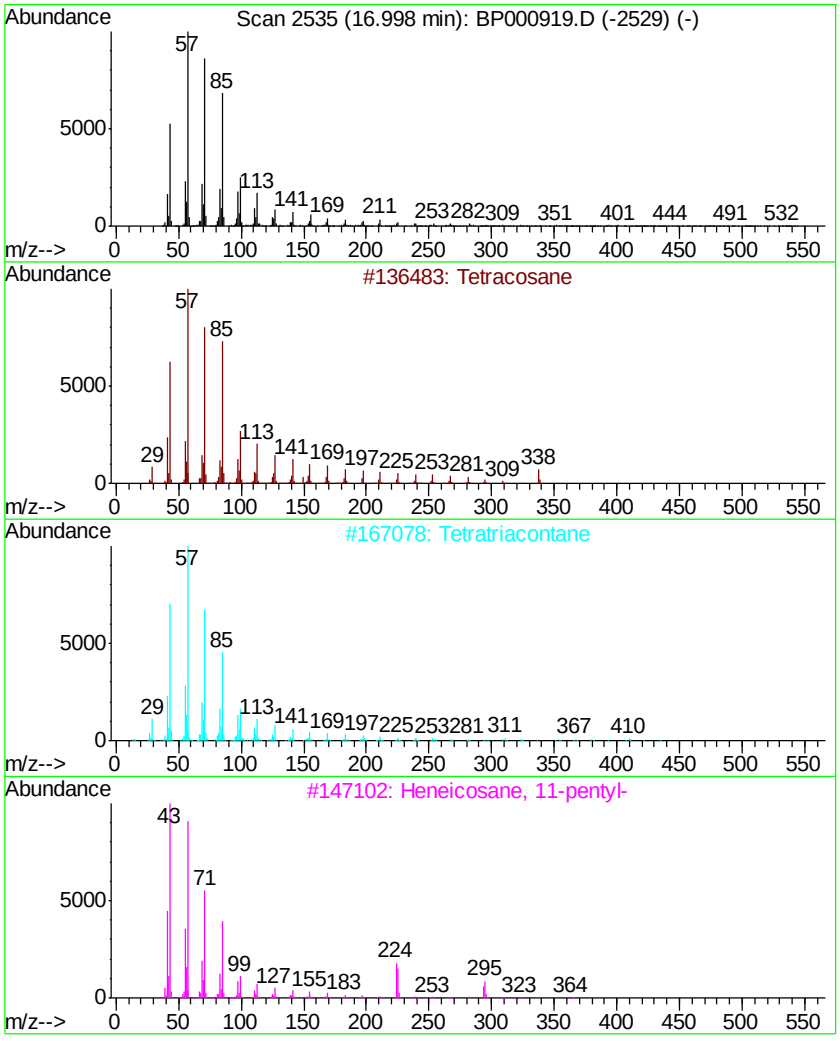
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
Quant Title : SVOA CALIBRATION

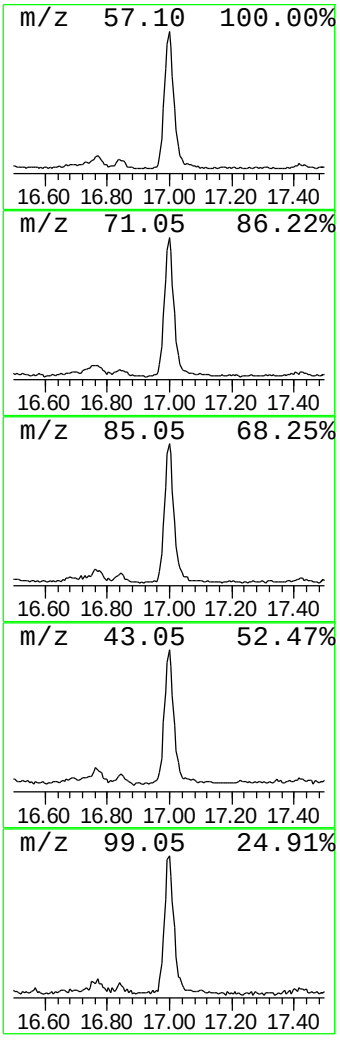
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.00	2.27 ng/ul	158314	Perylene-d12	15.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	93
2	Tetratriacontane	479	C34H70	014167-59-0	91
3	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	91
4	Nonadecane	268	C19H40	000629-92-5	89
5	Hentriacontane	437	C31H64	000630-04-6	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101919\
Data File : BP000919.D
Acq On : 19 Oct 2019 17:03
Operator : JU
Sample : K5354-15
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFBB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
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
Pyridine, 2-butyl...	9.23	12.2	ng/ul	670011	3	9.79	1102080	20.0
7,9-Di-tert-butyl...	11.67	3.2	ng/ul	195142	4	11.28	1217040	20.0
9-Octadecenamide,...	13.39	60.2	ng/ul	3986430	5	13.96	1323910	20.0
(DEL) Alkane: Str...	13.82	3.1	ng/ul	203353	5	13.96	1323910	20.0
(DEL) Alkane: Str...	14.13	5.0	ng/ul	328588	5	13.96	1323910	20.0
(DEL) Alkane: Str...	14.45	9.4	ng/ul	624982	5	13.96	1323910	20.0
(DEL) Alkane: Str...	14.75	13.0	ng/ul	910466	6	15.43	1396760	20.0
Octadecane, 1-iodo-	15.09	14.5	ng/ul	1014330	6	15.43	1396760	20.0
(DEL) Alkane: Str...	15.90	11.1	ng/ul	774689	6	15.43	1396760	20.0
Hexadecane, 1-iodo-	16.41	6.7	ng/ul	467492	6	15.43	1396760	20.0
(DEL) Alkane: Str...	17.00	2.3	ng/ul	158314	6	15.43	1396760	20.0