LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121219\
Data File : BM024098.D
Acq On : 13 Dec 2019 19:00
Operator : JU
Sample : K6167-20
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

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-BM112719MA.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	3.028	21	24	27	rBV	94740	118294	0.21%	0.078%
2	3.063	27	30	43	rVB	91709	165394	0.29%	0.110%
3	4.028	190	194	203	rBV	62224	104334	0.18%	0.069%
4	4.822	325	329	334	rBV	51513	76222	0.13%	0.050%
5	5.769	485	490	500	rBV3	34801	86125	0.15%	0.057%
6	6.669	634	643	654	rBV2	305398	657969	1.16%	0.436%
7	6.822	663	669	680	rBV	1001761	1668200	2.95%	1.105%
8	7.010	694	701	714	rBV	1121019	1882106	3.33%	1.246%
9	7.469	772	779	789	rBV	1239130	1968998	3.49%	1.304%
10	7.551	789	793	809	rVB	86072	166676	0.30%	0.110%
11	8.192	896	902	912	rBV	809851	1375316	2.43%	0.911%
12	8.622	969	975	986	rVB	1200019	2022060	3.58%	1.339%
13	9.057	1039	1049	1052	rBV	86539	153130	0.27%	0.101%
14	9.098	1052	1056	1063	rVB	191655	312358	0.55%	0.207%
15	9.339	1089	1097	1106	rBV	1013301	1702384	3.01%	1.127%
16	9.875	1180	1188	1207	rBV	1509577	2802745	4.96%	1.856%
17	10.039	1210	1216	1220	rBV3	35155	56748	0.10%	0.038%
18	10.239	1243	1250	1257	rBV	1610818	2692385	4.77%	1.783%
19	10.410	1273	1279	1287	rVB2	44331	90717	0.16%	0.060%
20	10.651	1314	1320	1332	rBV	626406	1084329	1.92%	0.718%
21	10.998	1373	1379	1387	rBV5	31778	65934	0.12%	0.044%
22	11.098	1392	1396	1401	rVV	34454	69649	0.12%	0.046%
23	11.157	1401	1406	1412	rVB3	40288	69917	0.12%	0.046%
24	11.551	1465	1473	1477	rBV3	29365	63387	0.11%	0.042%
25	12.069	1555	1561	1566	rBV	71452	153679	0.27%	0.102%
26	12.127	1566	1571	1582	rVB	342371	541836	0.96%	0.359%
27	12.610	1645	1653	1659	rBV4	44248	82162	0.15%	0.054%
28	13.039	1721	1726	1732	rBV2	98823	178055	0.32%	0.118%
29	13.551	1807	1813	1818	rBV	2903045	3930251	6.96%	2.603%
30	13.598	1818	1821	1829	rVB	185804	254901	0.45%	0.169%
31	13.692	1832	1837	1838	rBV	98130	132046	0.23%	0.087%
32	13.815	1852	1858	1871	rVB	3450396	4975184	8.81%	3.295%
33	13.963	1878	1883	1884	rBV2	112329	158435	0.28%	0.105%
34	14.127	1905	1911	1919	rBV	2377009	3462146	6.13%	2.293%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121219\
Data File : BM024098.D
Acq On : 13 Dec 2019 19:00
Operator : JU
Sample : K6167-20
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

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-BM112719MA.M
Title : SVOA CALIBRATION

35	14.245	1923	1931	1951	rBV	4403315	8624797	15.27%	5.712%
36	14.392	1951	1956	1965	rVB2	267401	522407	0.92%	0.346%
37	14.910	2039	2044	2049	rVB2	88074	148722	0.26%	0.098%
38	15.133	2075	2082	2091	rBV	4249417	6126762	10.85%	4.058%
39	15.280	2102	2107	2117	rBV	1339353	1970105	3.49%	1.305%
40	15.551	2143	2153	2157	rBV	33754296	56485519	100.00%	37.409%
41	15.774	2187	2191	2198	rVV2	48904	89142	0.16%	0.059%
42	16.674	2340	2344	2346	rBV4	52116	68431	0.12%	0.045%
43	16.886	2373	2380	2389	rBV	2707700	3922100	6.94%	2.598%
44	16.980	2390	2396	2406	rVV	4508064	6425151	11.37%	4.255%
45	17.598	2496	2501	2507	rVB	174972	244556	0.43%	0.162%
46	17.809	2532	2537	2551	rBV	768145	1297928	2.30%	0.860%
47	19.103	2753	2757	2764	rBV	416446	636431	1.13%	0.421%
48	19.286	2783	2788	2794	rBV2	5270671	7222780	12.79%	4.784%
49	20.833	3047	3051	3057	rBV2	998461	1252317	2.22%	0.829%
50	21.091	3090	3095	3101	rBV	3450132	4390283	7.77%	2.908%
51	21.274	3122	3126	3130	rBV	390416	547043	0.97%	0.362%
52	21.691	3194	3197	3201	rVB	504906	520573	0.92%	0.345%
53	22.133	3268	3272	3277	rVB	777224	921191	1.63%	0.610%
54	22.615	3350	3354	3360	rVB	824584	1147288	2.03%	0.760%
55	23.097	3430	3436	3441	rBV	4502959	7915496	14.01%	5.242%
56	23.144	3441	3444	3450	rVB	844495	1248851	2.21%	0.827%
57	23.227	3452	3458	3467	rVB	2639492	4816093	8.53%	3.190%
58	23.750	3543	3547	3555	rVB	638084	1125400	1.99%	0.745%

Sum of corrected areas: 150993438

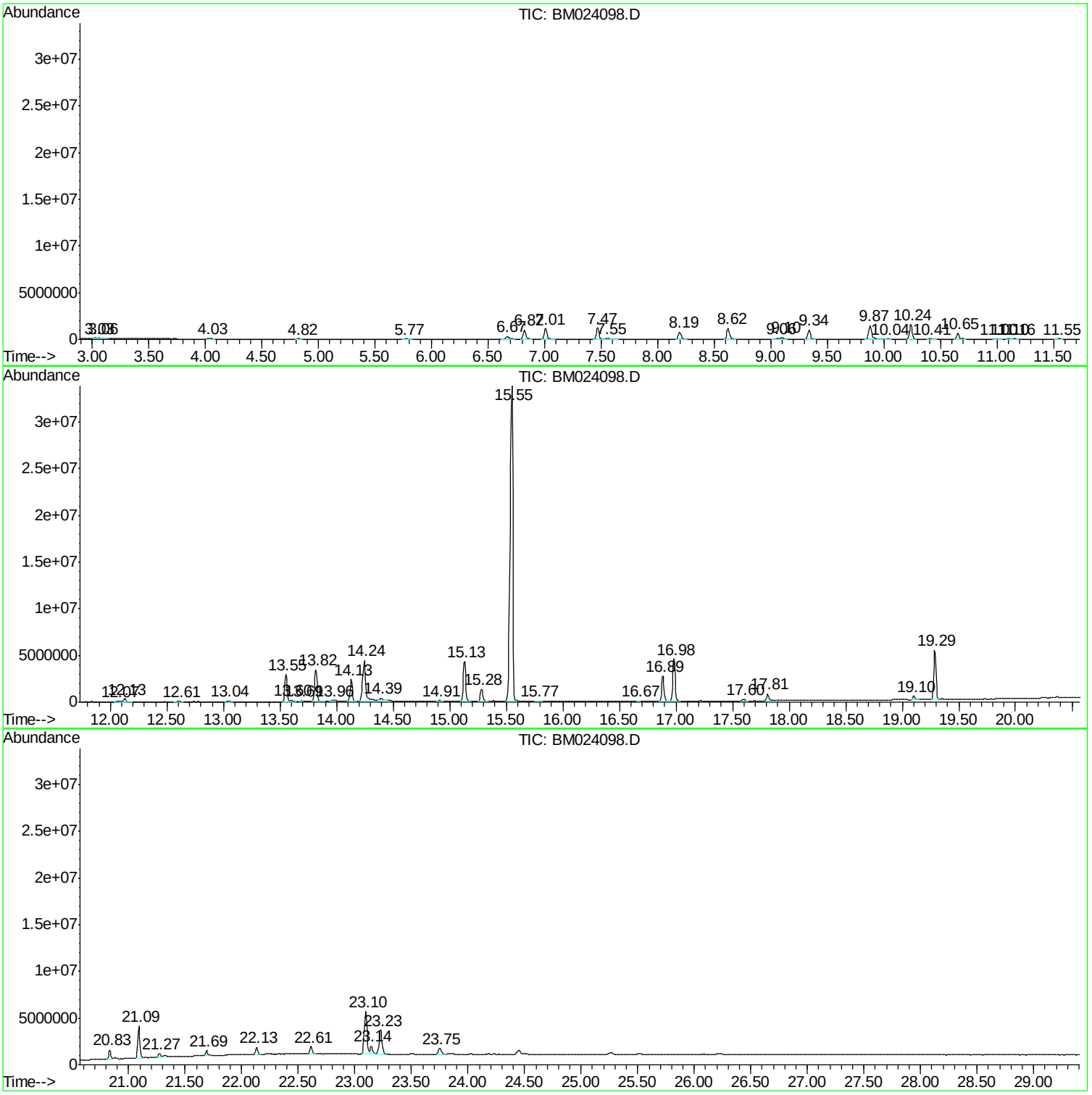
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121219\
Data File : BM024098.D
Acq On : 13 Dec 2019 19:00
Operator : JU
Sample : K6167-20
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.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\BM121219\
Data File : BM024098.D
Acq On : 13 Dec 2019 19:00
Operator : JU
Sample : K6167-20
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

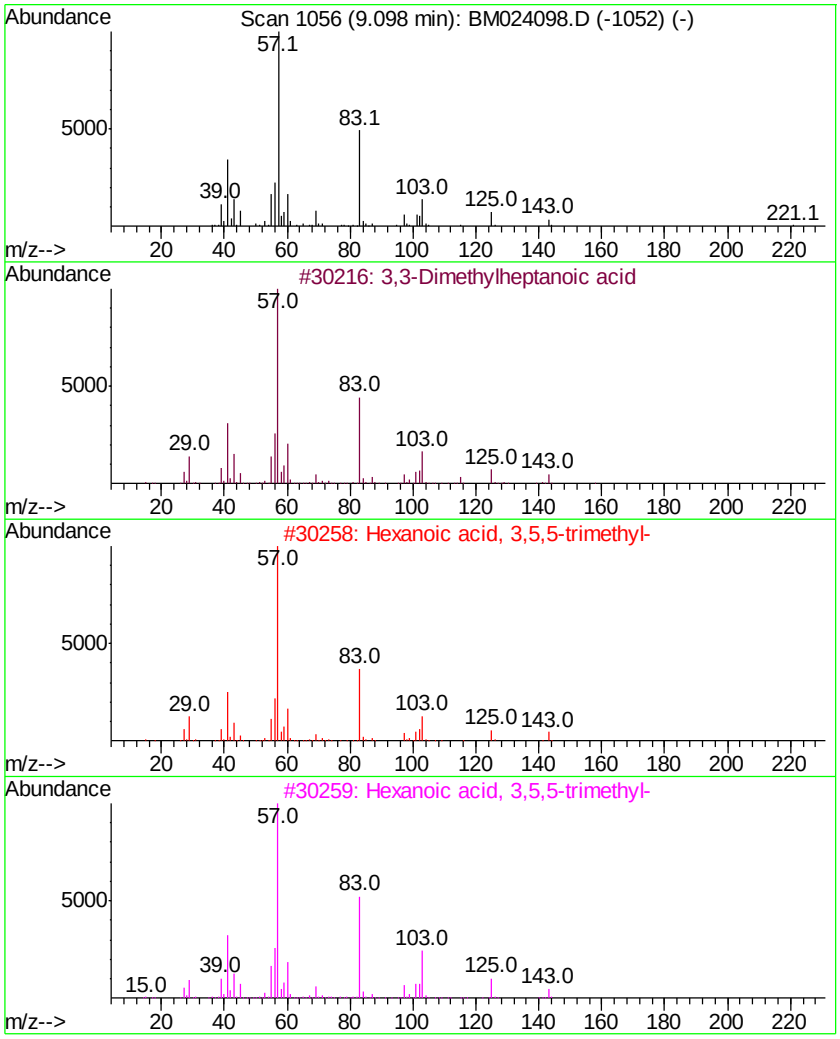
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

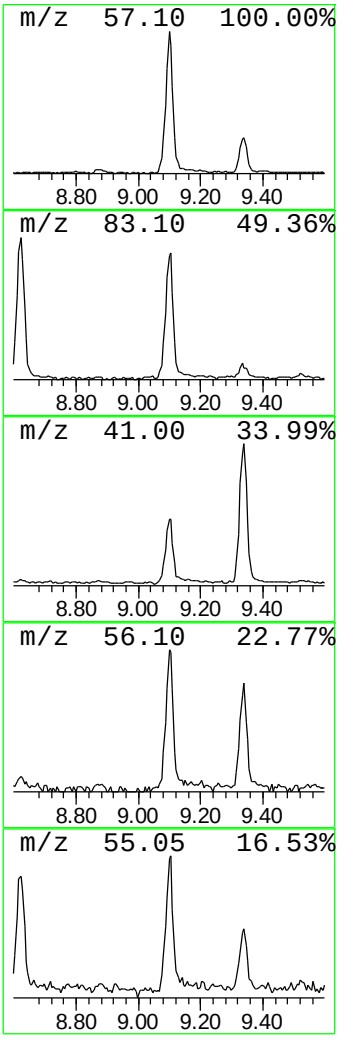
****

Peak Number 1 3,3-Dimethylheptanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.10	2.32 ng/ul	312358	Naphthalene-d8	10.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3-Dimethylheptanoic acid	158	C9H18O2	067061-30-7	83
2	Hexanoic acid, 3,5,5-trimethyl-	158	C9H18O2	003302-10-1	78
3	Hexanoic acid, 3,5,5-trimethyl-	158	C9H18O2	003302-10-1	78
4	Hexanoic acid, 3,5,5-trimethyl-	158	C9H18O2	003302-10-1	72
5	Hexanoyl chloride, 3,5,5-trimethyl-	176	C9H17ClO	036727-29-4	53





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121219\
Data File : BM024098.D
Acq On : 13 Dec 2019 19:00
Operator : JU
Sample : K6167-20
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

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

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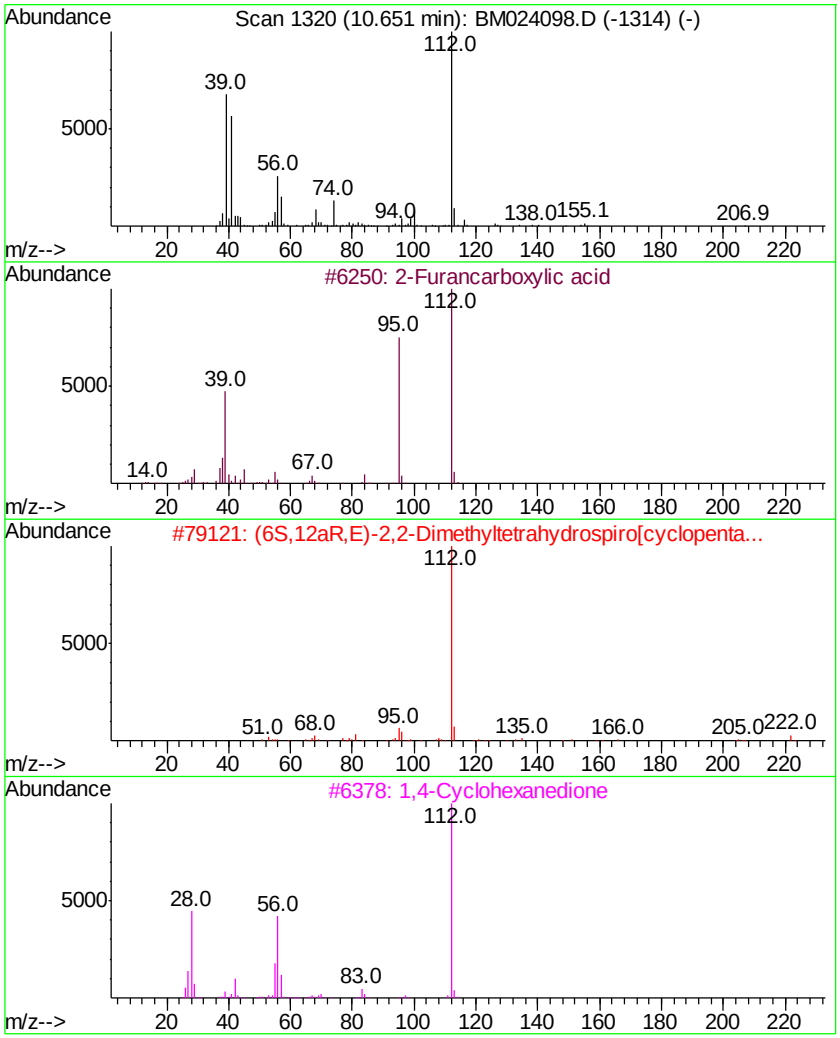
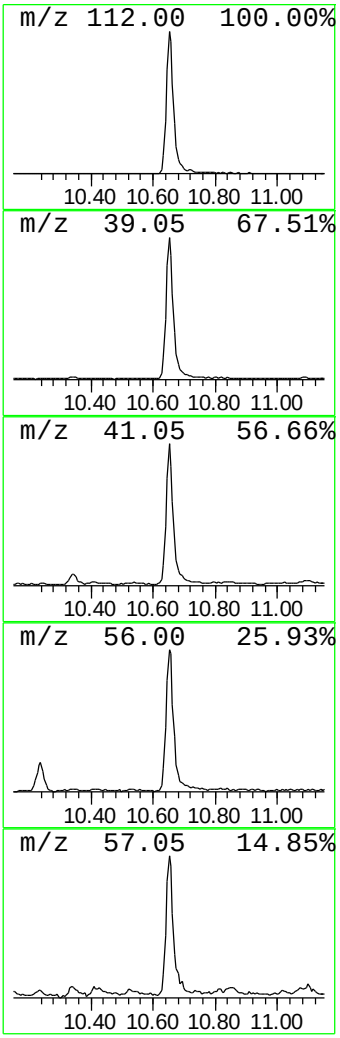
Peak Number 2 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	8.05 ng/ul	1084330	Naphthalene-d8	10.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Furancarboxylic acid	112	C5H4O3	000088-14-2	42
2	(6S,12aR,E)-2,2-Dimethyltetrahyd...	222	C13H22N2O	109175-49-7	40
3	1,4-Cyclohexanedione	112	C6H8O2	000637-88-7	38
4	3-Nitropyrrole	112	C4H4N2O2	005930-94-9	38
5	Aminomaleimide	112	C4H4N2O2	037770-94-8	38





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121219\
Data File : BM024098.D
Acq On : 13 Dec 2019 19:00
Operator : JU
Sample : K6167-20
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

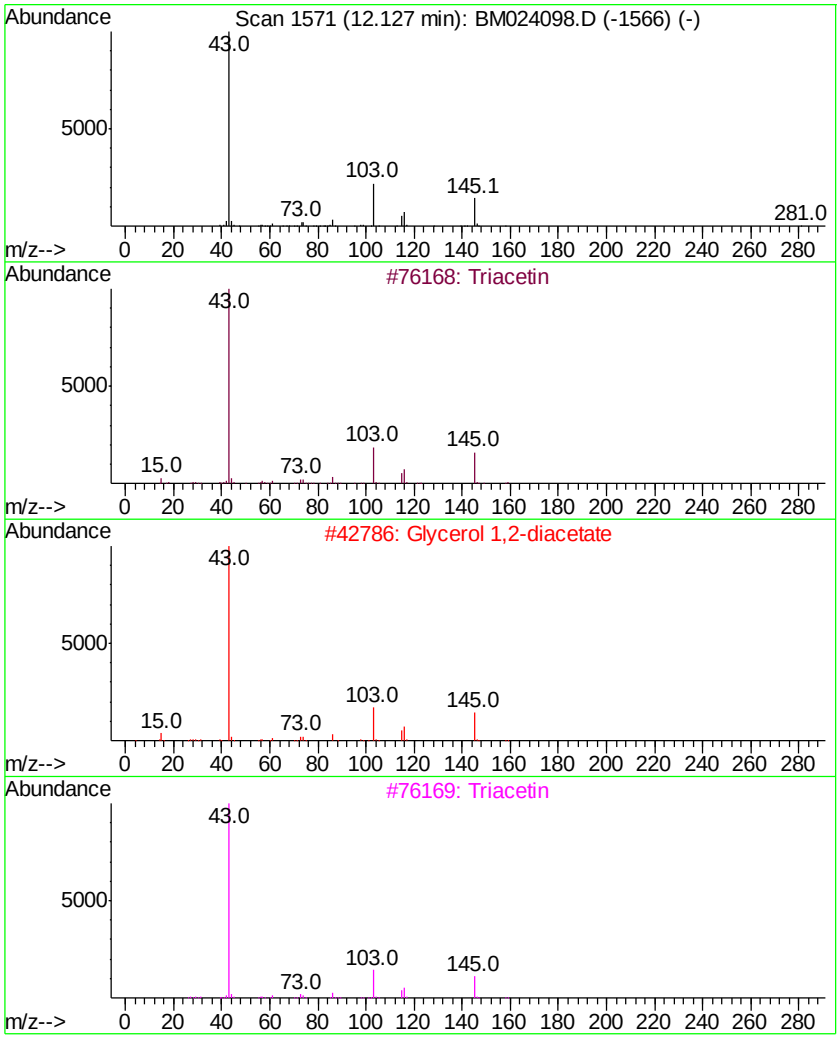
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

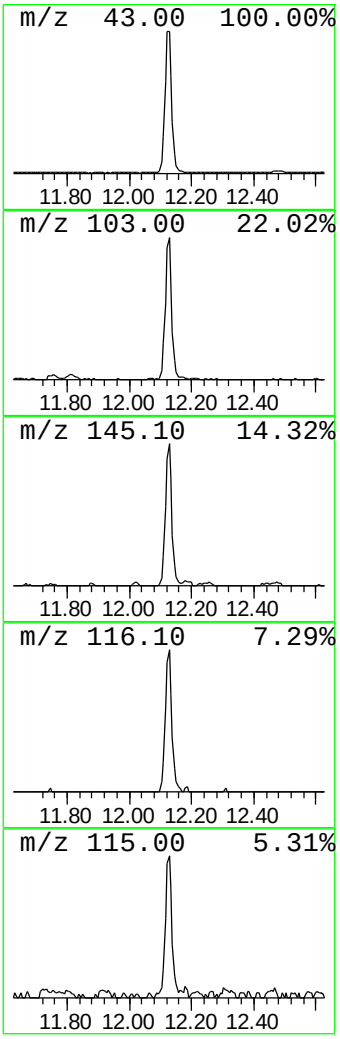
****

Peak Number 3 Triacetin Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	4.02 ng/ul	541836	Naphthalene-d8	10.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triacetin	218	C9H14O6	000102-76-1	90
2	Glycerol 1,2-diacetate	176	C7H12O5	000102-62-5	83
3	Triacetin	218	C9H14O6	000102-76-1	64
4	Triacetin	218	C9H14O6	000102-76-1	53
5	1,2,3,4-Butanetetrol, tetraaceta...	290	C12H18O8	007208-40-4	42





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121219\
Data File : BM024098.D
Acq On : 13 Dec 2019 19:00
Operator : JU
Sample : K6167-20
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

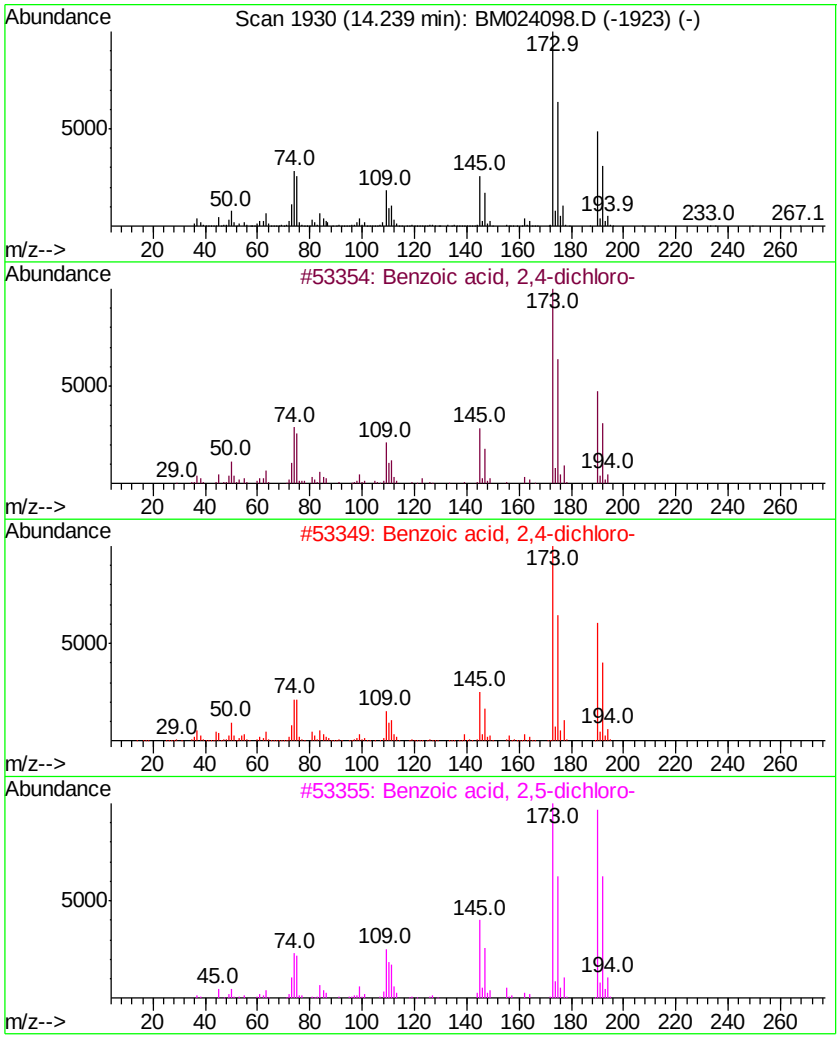
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

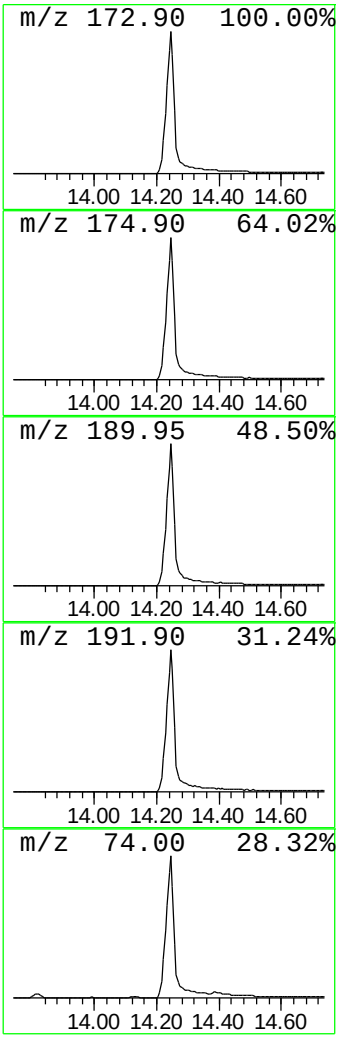
****

Peak Number 4 Benzoic acid, 2,4-dichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.24	49.82 ng/ul	8624800	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 2,4-dichloro-	190	C7H4Cl2O2	000050-84-0	99
2	Benzoic acid, 2,4-dichloro-	190	C7H4Cl2O2	000050-84-0	98
3	Benzoic acid, 2,5-dichloro-	190	C7H4Cl2O2	000050-79-3	96
4	Benzoic acid, 2,6-dichloro-	190	C7H4Cl2O2	000050-30-6	96
5	Benzoic acid, 2,5-dichloro-	190	C7H4Cl2O2	000050-79-3	96





Library Search Compound Report

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Data File : BM024098.D
Acq On : 13 Dec 2019 19:00
Operator : JU
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Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

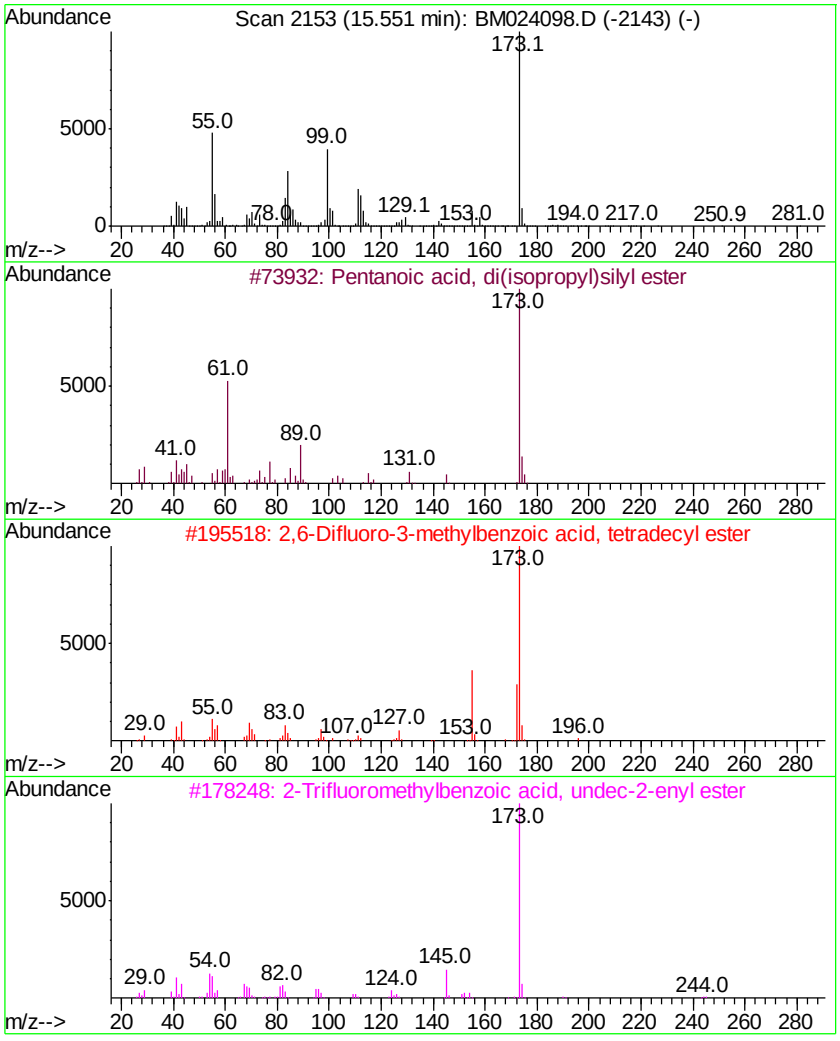
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

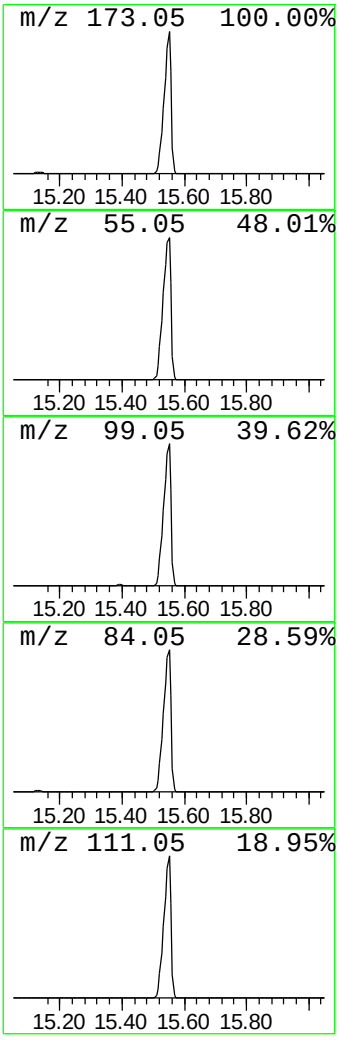
****

Peak Number 5 unknown-02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	288.04 ng/ul	56485500	Phenanthrene-d10	16.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentanoic acid, di(isopropyl)sil...	216	C11H24O2Si	1000279-48-0	35
2	2,6-Difluoro-3-methylbenzoic aci...	368	C22H34F2O2	1000338-85-1	32
3	2-Trifluoromethylbenzoic acid, u...	342	C19H25F3O2	1000299-40-8	25
4	2,6-Difluoro-3-methylbenzoic aci...	396	C24H38F2O2	1000338-85-3	23
5	2,6-Difluoro-3-methylbenzoic aci...	354	C21H32F2O2	1000338-85-0	23





Library Search Compound Report

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Data File : BM024098.D
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Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

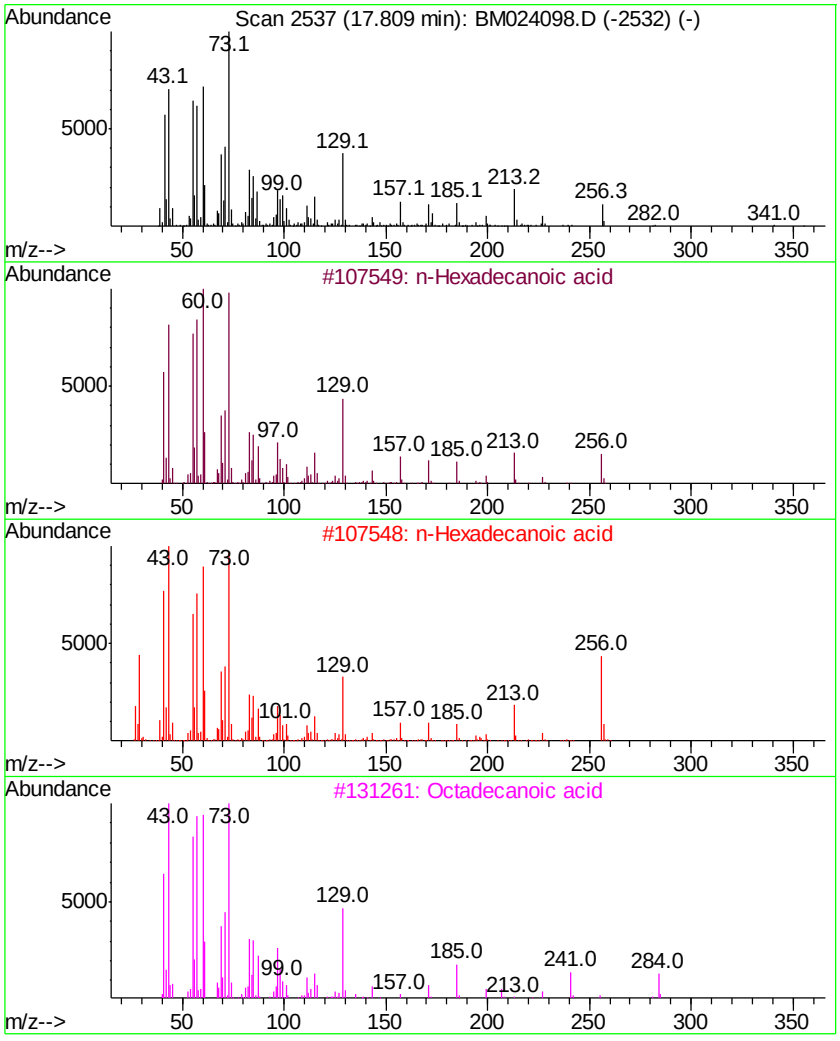
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

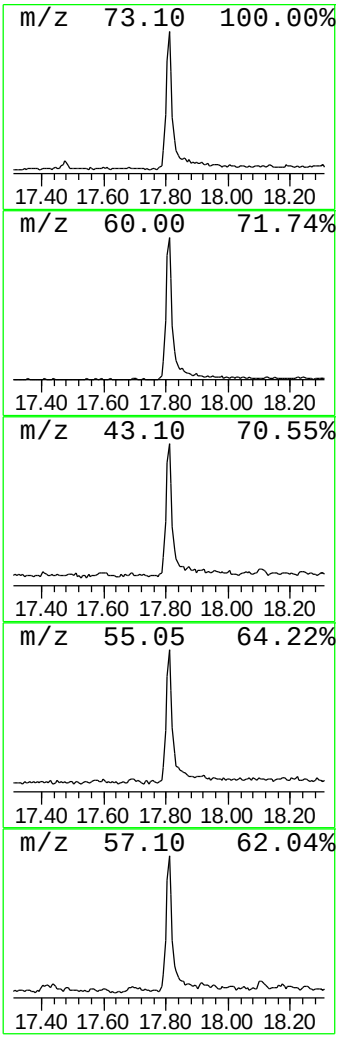
****

Peak Number 6 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.81	6.62 ng/ul	1297930	Phenanthrene-d10	16.89

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	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	81
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	76
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	72





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121219\
Data File : BM024098.D
Acq On : 13 Dec 2019 19:00
Operator : JU
Sample : K6167-20
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

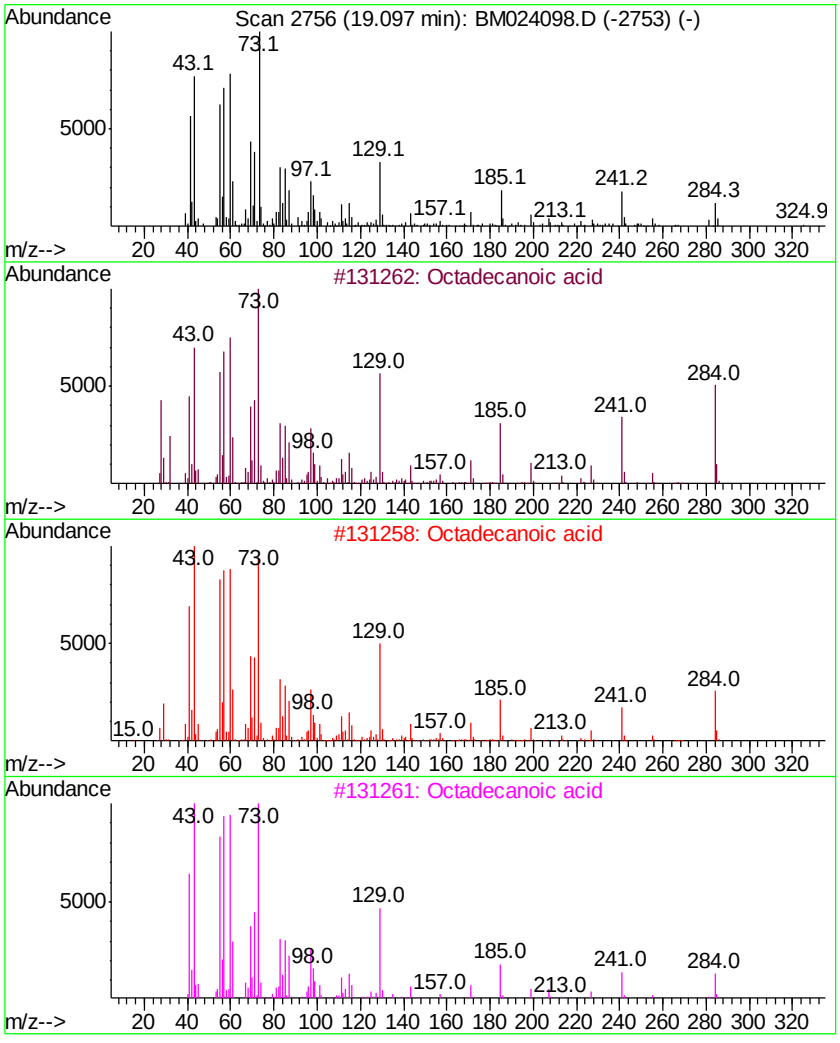
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

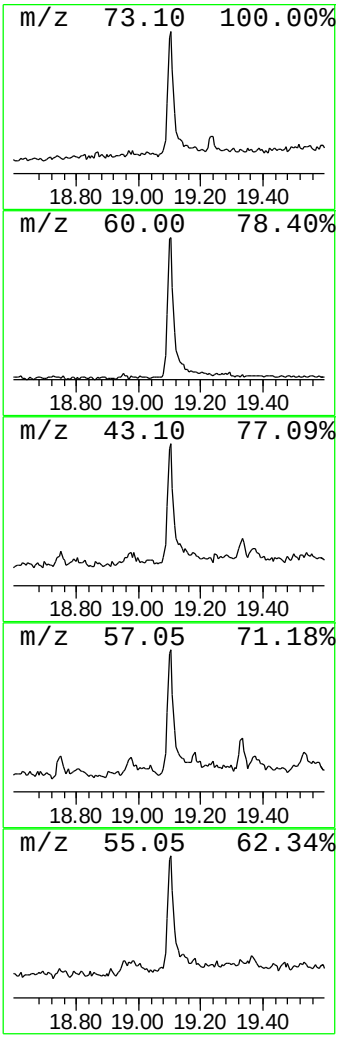
****

Peak Number 7 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.10	2.90 ng/ul	636431	Chrysene-d12	21.09

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	94
5	Octadecanoic acid	284	C18H36O2	000057-11-4	90





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121219\
Data File : BM024098.D
Acq On : 13 Dec 2019 19:00
Operator : JU
Sample : K6167-20
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

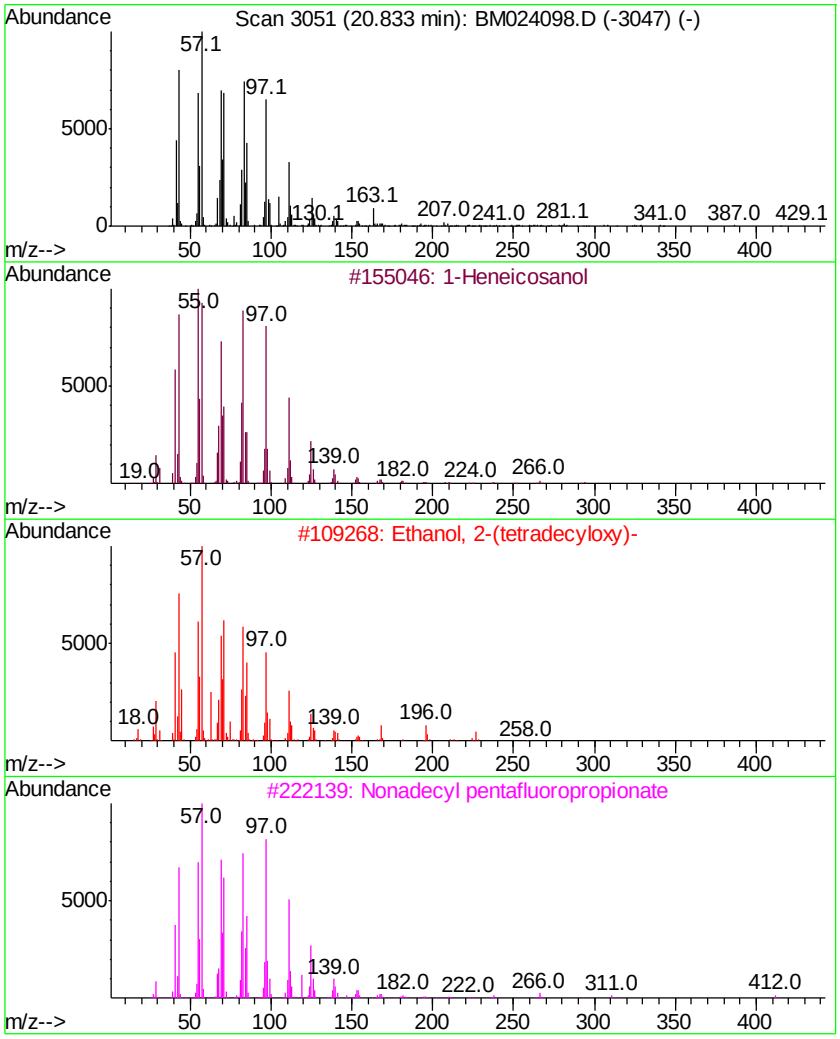
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

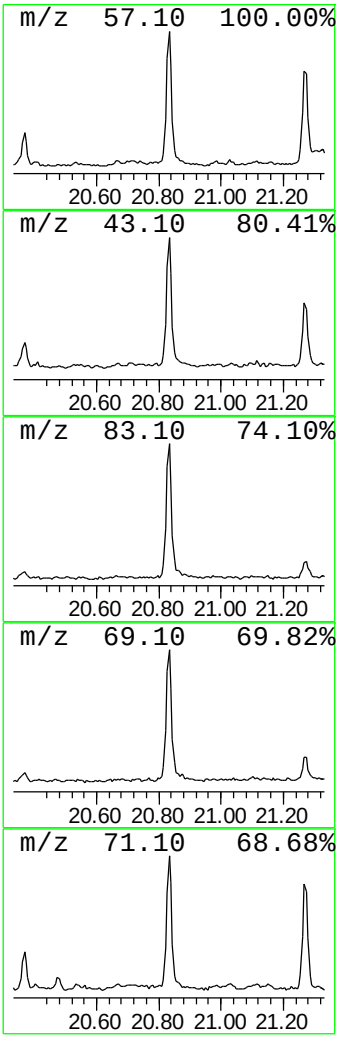
****

Peak Number 8 1-Heneicosanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.83	5.70 ng/ul	1252320	Chrysene-d12	21.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	93
2	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
3	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
4	Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	90
5	3-Eicosene, (E)-	280	C20H40	074685-33-9	90





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Data File : BM024098.D
Acq On : 13 Dec 2019 19:00
Operator : JU
Sample : K6167-20
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

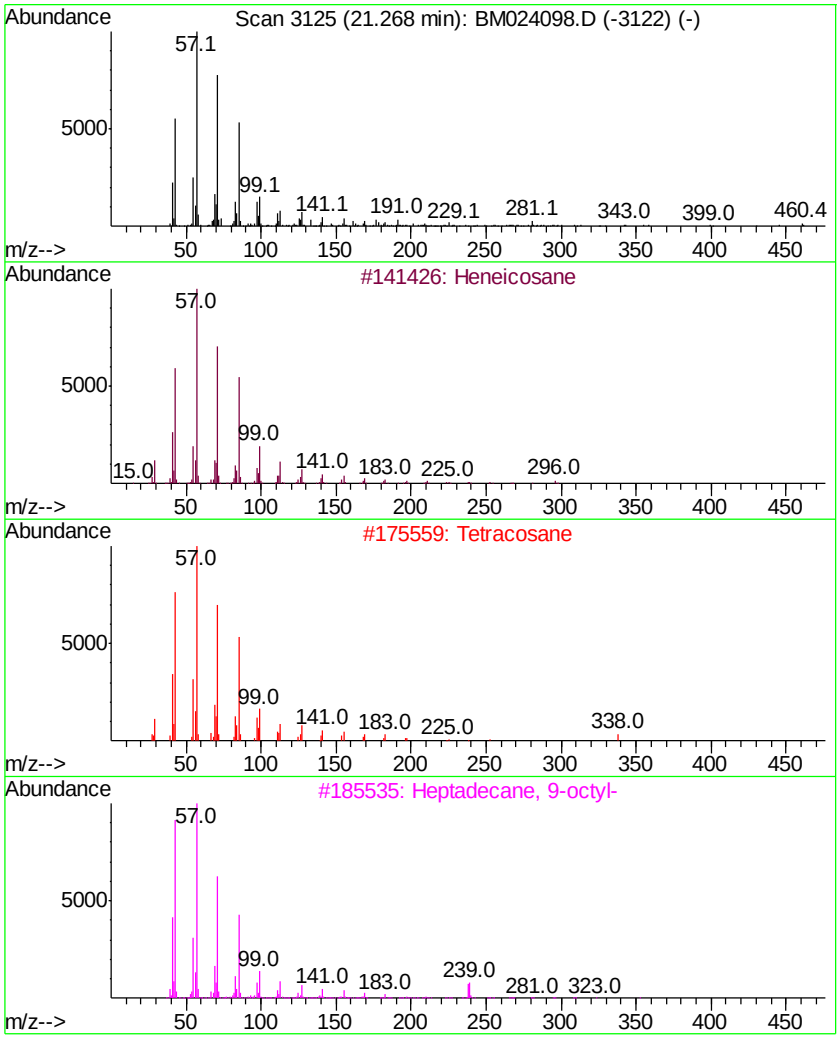
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

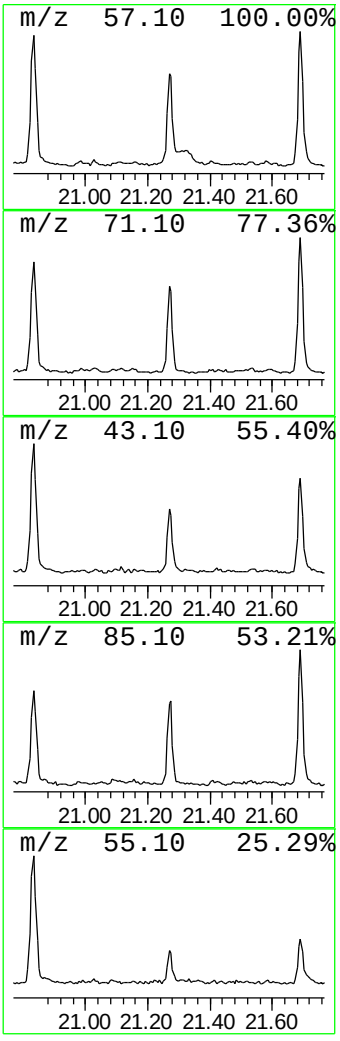
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.27	2.49 ng/ul	547043	Chrysene-d12	21.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	90
2	Tetracosane	338	C24H50	000646-31-1	87
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
4	Eicosane	282	C20H42	000112-95-8	86
5	Tetracosane	338	C24H50	000646-31-1	86





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121219\
Data File : BM024098.D
Acq On : 13 Dec 2019 19:00
Operator : JU
Sample : K6167-20
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

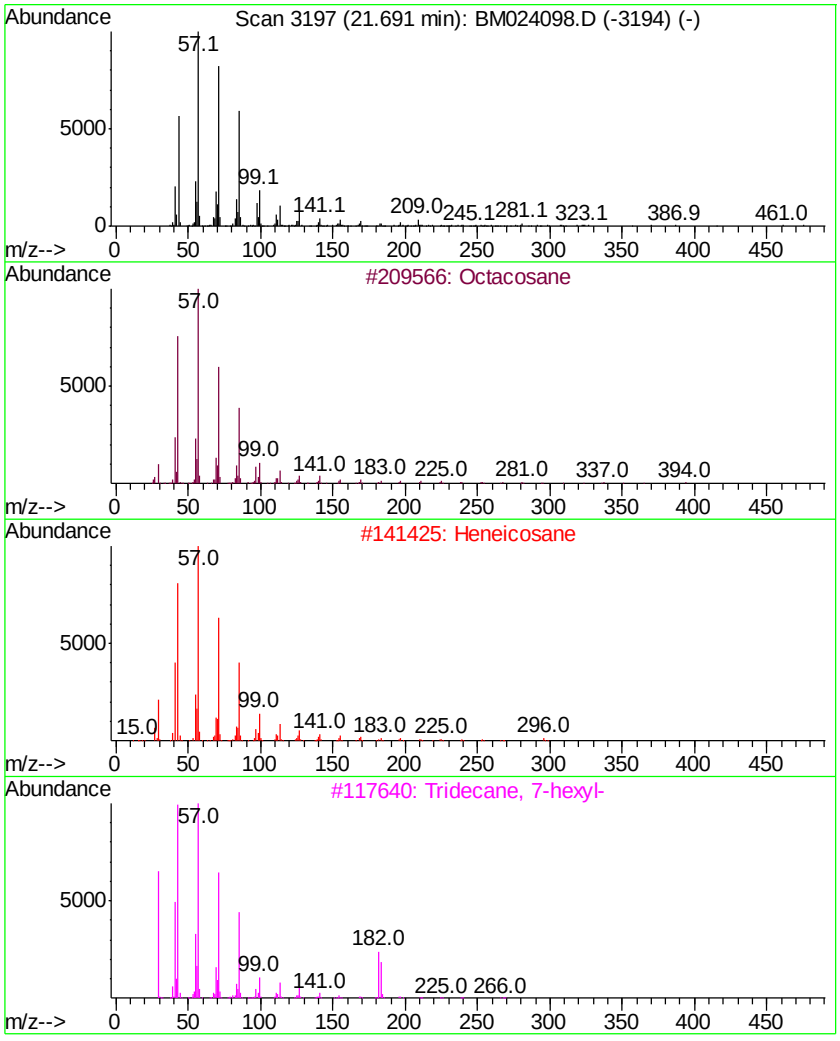
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

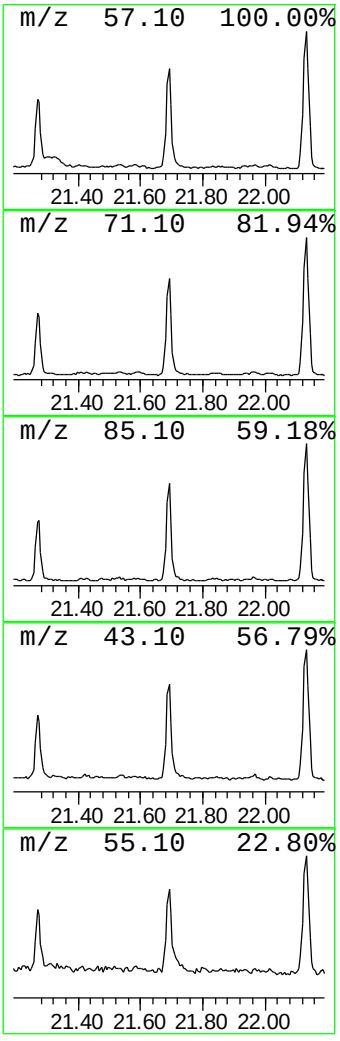
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.69	2.37 ng/ul	520573	Chrysene-d12	21.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	87
4	Tridecane, 6-propyl-	226	C16H34	055045-10-8	87
5	Tridecane, 7-propyl-	226	C16H34	055045-09-5	83





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Acq On : 13 Dec 2019 19:00
Operator : JU
Sample : K6167-20
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

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

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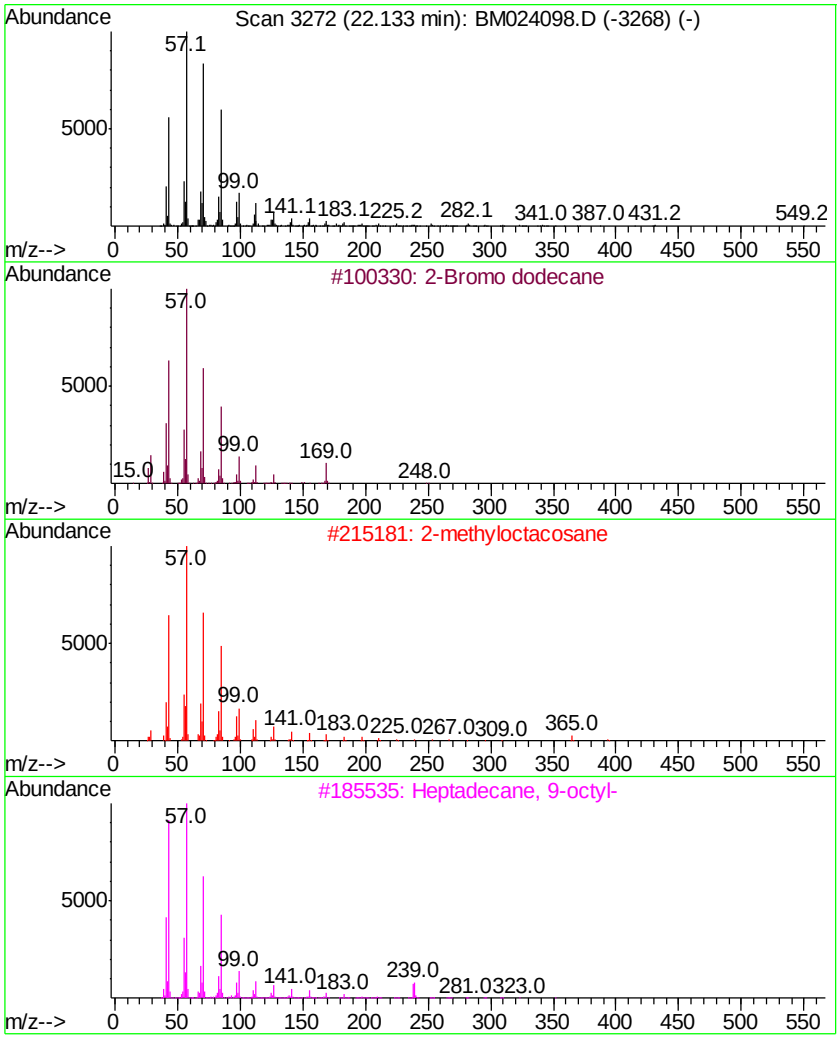
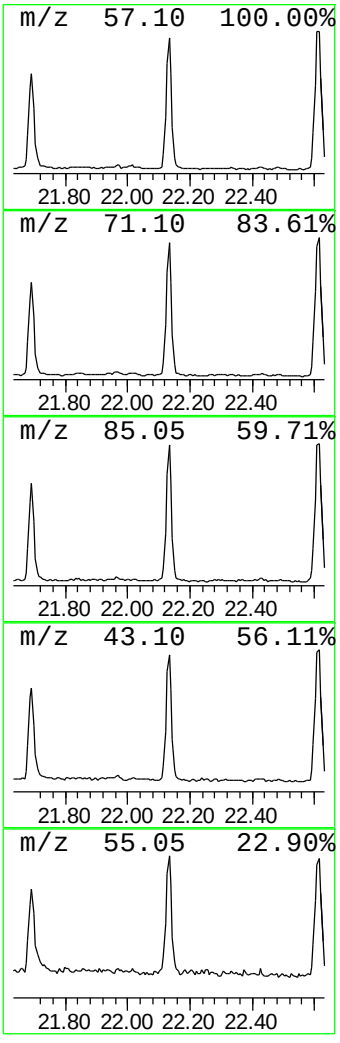
Peak Number 11 2-Bromo dodecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.13	4.20 ng/ul	921191	Chrysene-d12	21.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromo dodecane	248	C12H25Br	013187-99-0	93
2	2-methyloctacosane	408	C29H60	1000376-72-8	91
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4	Hexacosane	366	C26H54	000630-01-3	90
5	Heptacosane	380	C27H56	000593-49-7	87





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Data File : BM024098.D
Acq On : 13 Dec 2019 19:00
Operator : JU
Sample : K6167-20
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

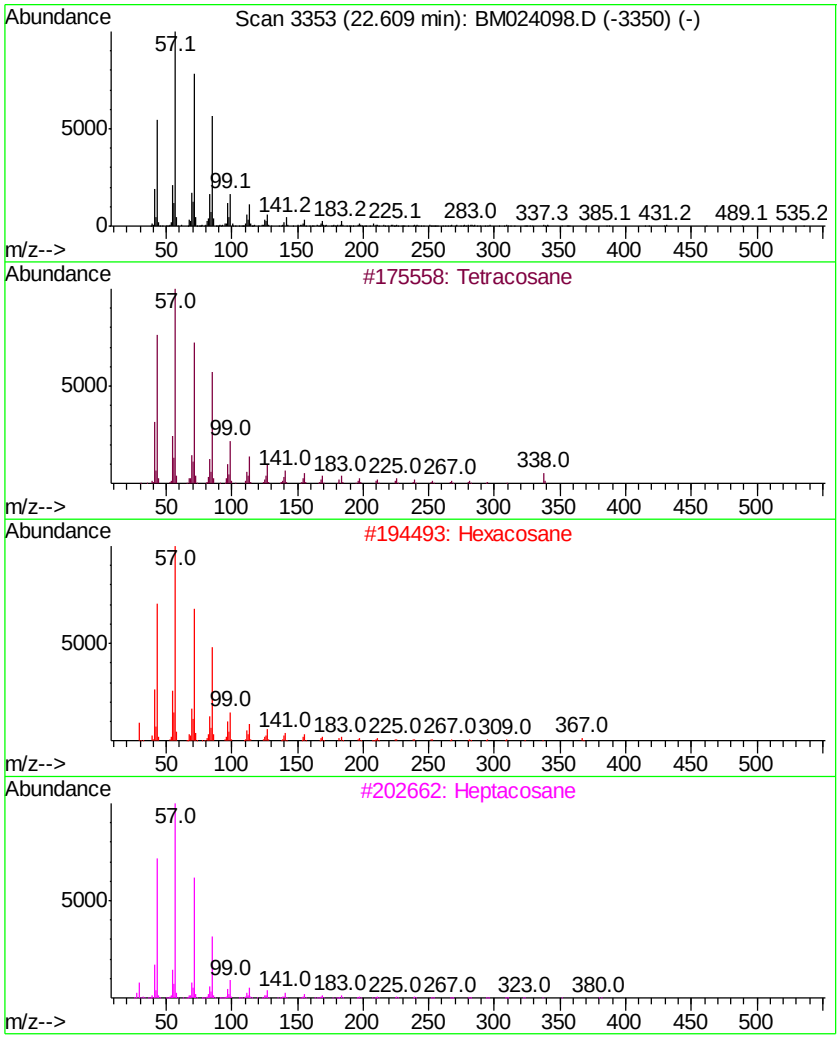
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

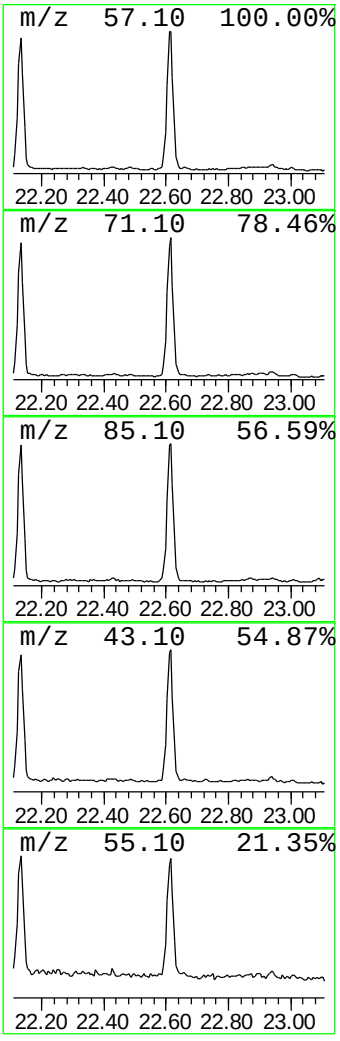
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.61	4.76 ng/ul	1147290	Perylene-d12	23.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	97
2	Hexacosane	366	C26H54	000630-01-3	94
3	Heptacosane	380	C27H56	000593-49-7	91
4	Octacosane	394	C28H58	000630-02-4	91
5	2-methyloctacosane	408	C29H60	1000376-72-8	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121219\
Data File : BM024098.D
Acq On : 13 Dec 2019 19:00
Operator : JU
Sample : K6167-20
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

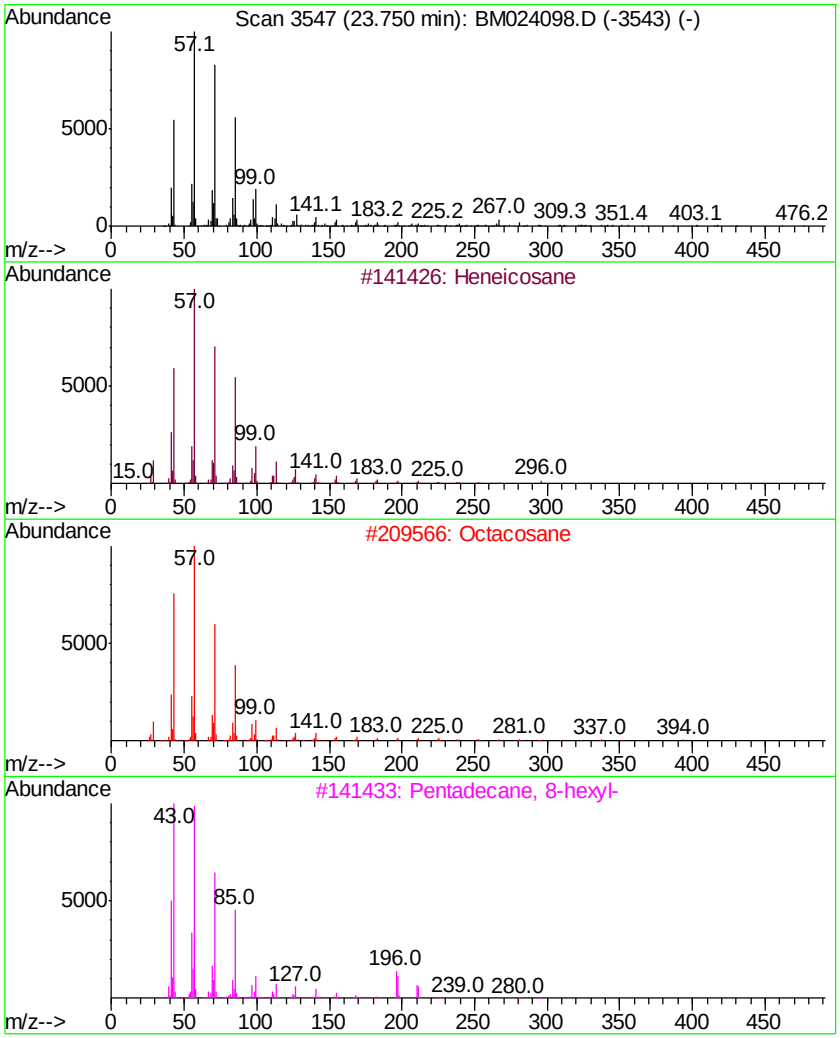
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

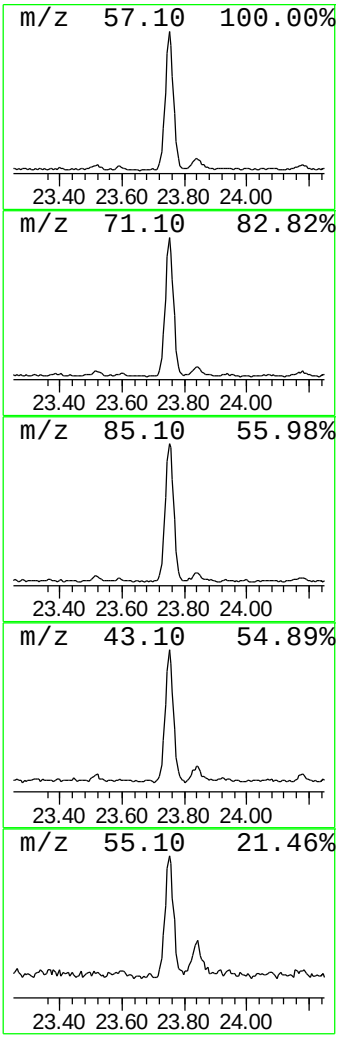
****

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.75	4.67 ng/ul	1125400	Perylene-d12	23.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Octacosane	394	C28H58	000630-02-4	91
3	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
4	Docosane, 11-butyl-	366	C26H54	013475-76-8	91
5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM121219\
Data File : BM024098.D
Acq On : 13 Dec 2019 19:00
Operator : JU
Sample : K6167-20
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFR92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.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
3,3-Dimethylhepta...	9.10	2.3	ng/ul	312358	2	10.24	2692390	20.0
unknown-01	10.65	8.1	ng/ul	1084330	2	10.24	2692390	20.0
Triacetin	12.13	4.0	ng/ul	541836	2	10.24	2692390	20.0
Benzoic acid, 2,4...	14.24	49.8	ng/ul	8624800	3	14.13	3462150	20.0
unknown-02	15.55	288.0	ng/ul	56485500	4	16.89	3922100	20.0
n-Hexadecanoic acid	17.81	6.6	ng/ul	1297930	4	16.89	3922100	20.0
Octadecanoic acid	19.10	2.9	ng/ul	636431	5	21.09	4390280	20.0
1-Heneicosanol	20.83	5.7	ng/ul	1252320	5	21.09	4390280	20.0
(DEL) Alkane: Str...	21.27	2.5	ng/ul	547043	5	21.09	4390280	20.0
(DEL) Alkane: Str...	21.69	2.4	ng/ul	520573	5	21.09	4390280	20.0
2-Bromo dodecane	22.13	4.2	ng/ul	921191	5	21.09	4390280	20.0
(DEL) Alkane: Str...	22.61	4.8	ng/ul	1147290	6	23.23	4816090	20.0
(DEL) Alkane: Str...	23.75	4.7	ng/ul	1125400	6	23.23	4816090	20.0