LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051920\
Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

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-BM051320MA.M
Title : SVOA CALIBRATION

Signal : TIC: BM026035.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.081	29	33	42	rVB	60515	94081	0.20%	0.081%
2	6.693	636	647	660	rVB	250227	432975	0.92%	0.374%
3	6.851	668	674	680	rBV	820591	1235291	2.63%	1.068%
4	6.904	680	683	688	rVB	52172	75551	0.16%	0.065%
5	7.040	699	706	714	rBV	873276	1327882	2.82%	1.148%
6	7.498	778	784	791	rBV	811024	1255852	2.67%	1.086%
7	7.622	801	805	816	rVB2	59018	108712	0.23%	0.094%
8	7.869	842	847	854	rVB	69139	108508	0.23%	0.094%
9	8.210	898	905	916	rVV	648755	1068131	2.27%	0.923%
10	8.310	916	922	931	rVV2	38265	90106	0.19%	0.078%
11	8.645	973	979	996	rVB2	1027804	1976867	4.20%	1.709%
12	8.839	1006	1012	1020	rVB3	136531	249619	0.53%	0.216%
13	8.939	1020	1029	1034	rBV3	74734	153061	0.33%	0.132%
14	9.292	1083	1089	1094	rBV	61085	93564	0.20%	0.081%
15	9.357	1094	1100	1108	rVV	881572	1412413	3.00%	1.221%
16	9.481	1116	1121	1129	rVB4	100295	194968	0.41%	0.169%
17	9.563	1131	1135	1141	rVB2	63982	96860	0.21%	0.084%
18	9.675	1149	1154	1164	rBV3	146748	281286	0.60%	0.243%
19	9.892	1185	1191	1200	rVB	1203854	1971496	4.19%	1.704%
20	9.981	1202	1206	1213	rVB2	51793	96912	0.21%	0.084%
21	10.163	1226	1237	1241	rBV5	65041	179178	0.38%	0.155%
22	10.263	1247	1254	1260	rBV	1177105	1878681	4.00%	1.624%
23	10.328	1260	1265	1271	rVV4	108277	223767	0.48%	0.193%
24	10.392	1271	1276	1290	rVB2	103876	256721	0.55%	0.222%
25	10.839	1345	1352	1355	rBV2	128067	233496	0.50%	0.202%
26	10.875	1355	1358	1365	rVB	183056	291924	0.62%	0.252%
27	10.986	1371	1377	1382	rBV2	64101	109059	0.23%	0.094%
28	11.110	1391	1398	1403	rBV6	45672	95020	0.20%	0.082%
29	11.163	1403	1407	1415	rVB5	42929	100550	0.21%	0.087%
30	11.304	1426	1431	1435	rBV	75262	117318	0.25%	0.101%
31	11.375	1436	1443	1450	rVB6	101886	231885	0.49%	0.200%
32	11.492	1457	1463	1470	rVB4	133609	286733	0.61%	0.248%
33	11.563	1470	1475	1477	rBV4	48780	87001	0.19%	0.075%
34	11.592	1477	1480	1485	rVV2	75147	161402	0.34%	0.140%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051920\
Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

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-BM051320MA.M
Title : SVOA CALIBRATION

35	11.651	1485	1490	1497	rVB3	390988	694676	1.48%	0.601%
36	11.769	1503	1510	1516	rBV5	50892	155200	0.33%	0.134%
37	11.904	1528	1533	1537	rBV4	103804	194596	0.41%	0.168%
38	11.945	1537	1540	1545	rVB3	103098	146727	0.31%	0.127%
39	11.998	1545	1549	1553	rVB	63670	90067	0.19%	0.078%
40	12.104	1562	1567	1572	rBV	147972	249293	0.53%	0.216%
41	12.163	1572	1577	1582	rVV3	81596	195320	0.42%	0.169%
42	12.233	1582	1589	1591	rVV2	112194	233100	0.50%	0.202%
43	12.269	1591	1595	1601	rVB2	236748	381421	0.81%	0.330%
44	12.339	1601	1607	1613	rBV3	195796	413140	0.88%	0.357%
45	12.404	1613	1618	1623	rBV2	79804	131929	0.28%	0.114%
46	12.463	1623	1628	1632	rBV3	145257	220661	0.47%	0.191%
47	12.627	1651	1656	1660	rVV3	146954	197271	0.42%	0.171%
48	12.674	1660	1664	1672	rVB	193134	289734	0.62%	0.250%
49	12.857	1691	1695	1699	rBV	83666	130530	0.28%	0.113%
50	12.898	1699	1702	1707	rVB3	70058	85257	0.18%	0.074%
51	12.963	1708	1713	1719	rVB8	48183	95980	0.20%	0.083%
52	13.063	1726	1730	1732	rBV	46429	69455	0.15%	0.060%
53	13.104	1732	1737	1742	rVB4	165741	270650	0.58%	0.234%
54	13.186	1746	1751	1758	rVB3	77217	129764	0.28%	0.112%
55	13.274	1760	1766	1769	rBV3	169106	305869	0.65%	0.264%
56	13.427	1786	1792	1797	rVB7	63511	132180	0.28%	0.114%
57	13.480	1797	1801	1805	rVB2	55027	72199	0.15%	0.062%
58	13.569	1810	1816	1820	rVV	2054672	2798191	5.95%	2.419%
59	13.616	1820	1824	1835	rVB	507733	767318	1.63%	0.663%
60	13.833	1854	1861	1867	rVB	2338255	3459486	7.36%	2.991%
61	14.145	1908	1914	1925	rVB	1550943	2351414	5.00%	2.033%
62	14.363	1940	1951	1958	rBV3	221889	500354	1.06%	0.433%
63	14.480	1965	1971	1981	rVB5	79098	221415	0.47%	0.191%
64	14.698	2003	2008	2012	rVV	351567	517292	1.10%	0.447%
65	14.739	2012	2015	2017	rVV3	85433	119861	0.25%	0.104%
66	14.780	2017	2022	2033	rVB	1535467	2164951	4.60%	1.872%
67	14.910	2041	2044	2049	rVV2	61062	90141	0.19%	0.078%
68	14.986	2052	2057	2062	rVB4	92982	158834	0.34%	0.137%
69	15.145	2076	2084	2090	rVB	3042409	4158099	8.84%	3.595%
70	15.204	2090	2094	2100	rBV7	37608	79869	0.17%	0.069%
71	15.274	2100	2106	2113	rBV2	1115004	2092929	4.45%	1.809%

LSC Area Percent Report

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Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

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-BM051320MA.M
Title : SVOA CALIBRATION

72	15.363	2118	2121	2128	rVB	88704	125754	0.27%	0.109%
73	15.598	2157	2161	2169	rVB3	81787	173618	0.37%	0.150%
74	15.686	2169	2176	2184	rBV5	118834	265376	0.56%	0.229%
75	15.868	2203	2207	2215	rBV2	410216	623927	1.33%	0.539%
76	16.292	2276	2279	2286	rVB6	77650	106269	0.23%	0.092%
77	16.410	2296	2299	2303	rBV	61923	72590	0.15%	0.063%
78	16.451	2303	2306	2315	rVB5	89710	219437	0.47%	0.190%
79	16.574	2316	2327	2341	rBV	27723460	47014663	100.00%	40.644%
80	16.680	2342	2345	2349	rVB6	71232	96001	0.20%	0.083%
81	16.768	2357	2360	2365	rVB	150353	196767	0.42%	0.170%
82	16.898	2377	2382	2387	rVB	1898832	2518007	5.36%	2.177%
83	16.992	2393	2398	2406	rBV2	3081805	4389646	9.34%	3.795%
84	17.098	2411	2416	2426	rVB6	79060	220718	0.47%	0.191%
85	17.398	2464	2467	2471	rVB2	66869	86148	0.18%	0.074%
86	17.521	2484	2488	2491	rBV2	74943	104352	0.22%	0.090%
87	17.821	2535	2539	2542	rBV2	151645	223279	0.47%	0.193%
88	17.886	2547	2550	2554	rVB	78956	94071	0.20%	0.081%
89	19.292	2783	2789	2797	rBV	3780596	4827771	10.27%	4.174%
90	20.839	3048	3052	3057	rBV	282444	367707	0.78%	0.318%
91	21.092	3090	3095	3101	rVB	2242940	2786303	5.93%	2.409%
92	21.280	3125	3127	3132	rVB	102360	101045	0.21%	0.087%
93	21.703	3195	3199	3203	rBV	188953	239348	0.51%	0.207%
94	22.139	3270	3273	3278	rVB	271107	326152	0.69%	0.282%
95	22.615	3351	3354	3364	rVB	341092	503618	1.07%	0.435%
96	23.080	3427	3433	3440	rBV	2848337	4759360	10.12%	4.114%
97	23.150	3441	3445	3449	rVV	380284	570875	1.21%	0.494%
98	23.215	3450	3456	3466	rVB	1612385	2842006	6.04%	2.457%
99	23.750	3543	3547	3554	rVB	295620	495024	1.05%	0.428%
100	24.444	3661	3665	3672	rVB	203353	383222	0.82%	0.331%

Sum of corrected areas: 115675097

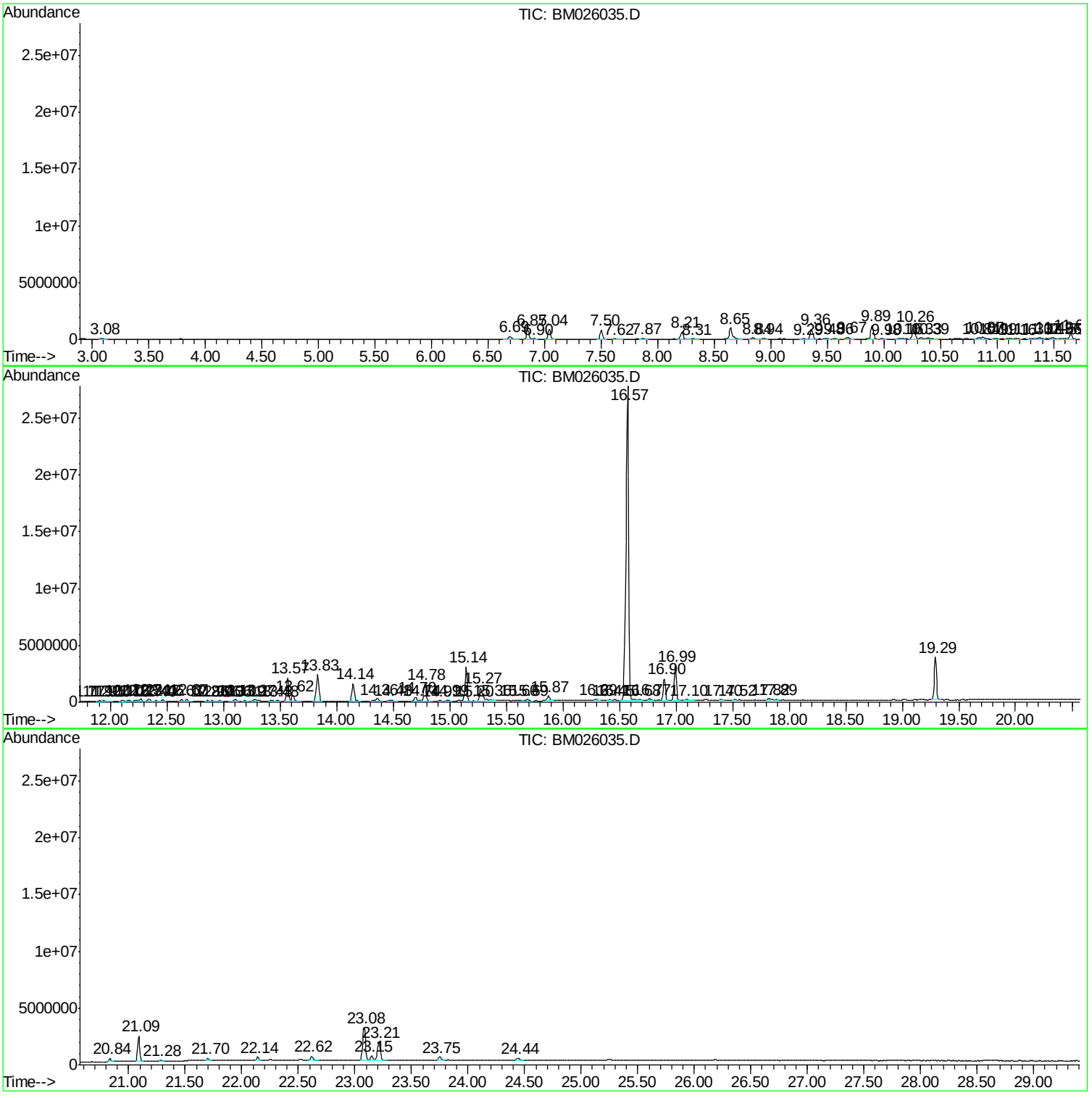
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051920\
Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.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\BM051920\
Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

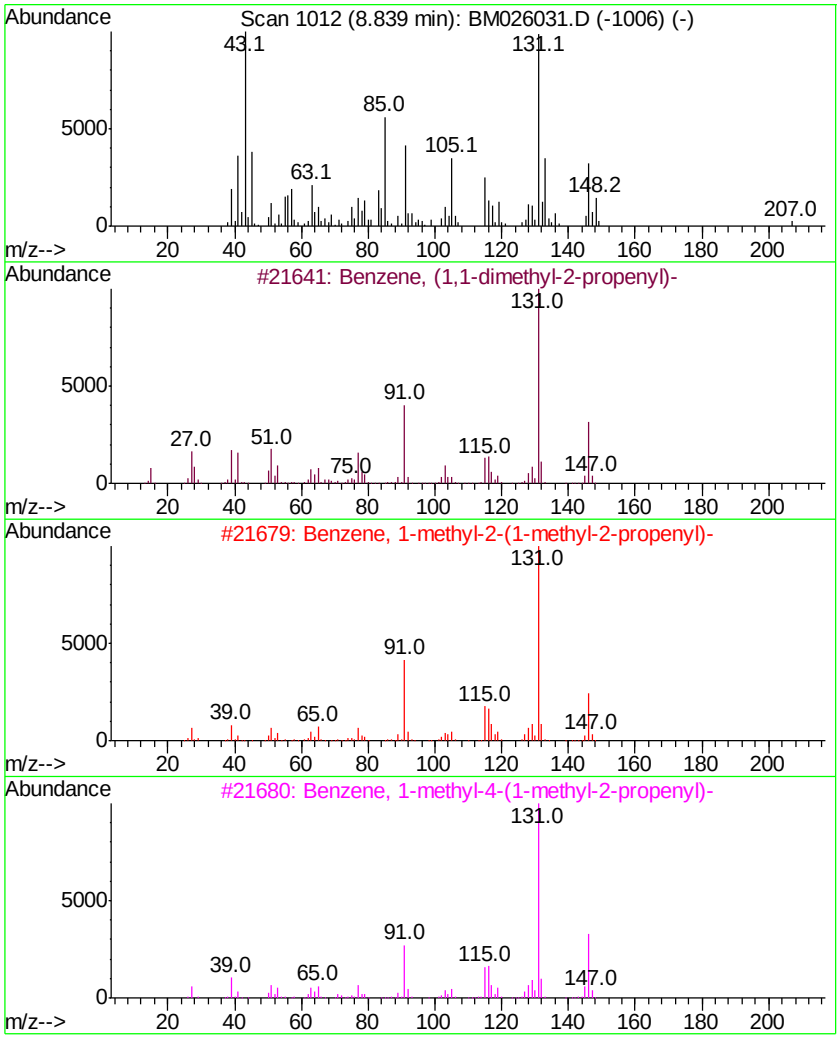
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

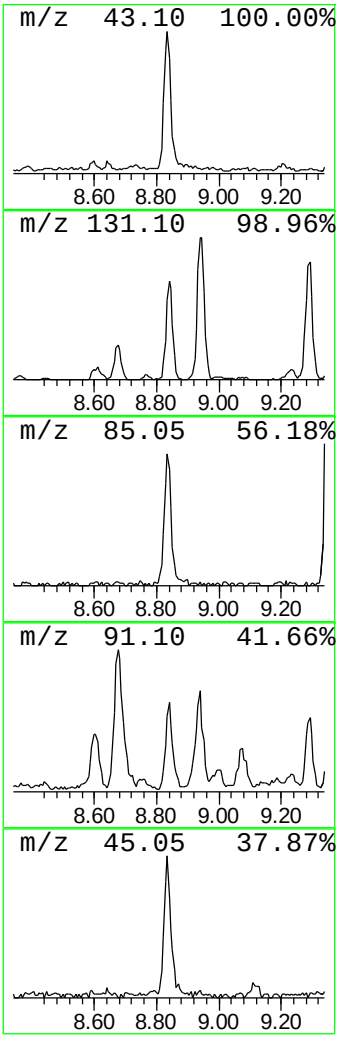
****

Peak Number 1 Benzene, (1,1-dimethyl-2-pr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.84	3.98 ng/ul	249619	1,4-Dichlorobenzene-d4	7.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1,1-dimethyl-2-propenyl)-	146	C11H14	018321-36-3	91
2	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	64
3	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	64
4	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	62
5	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	60





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051920\
Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

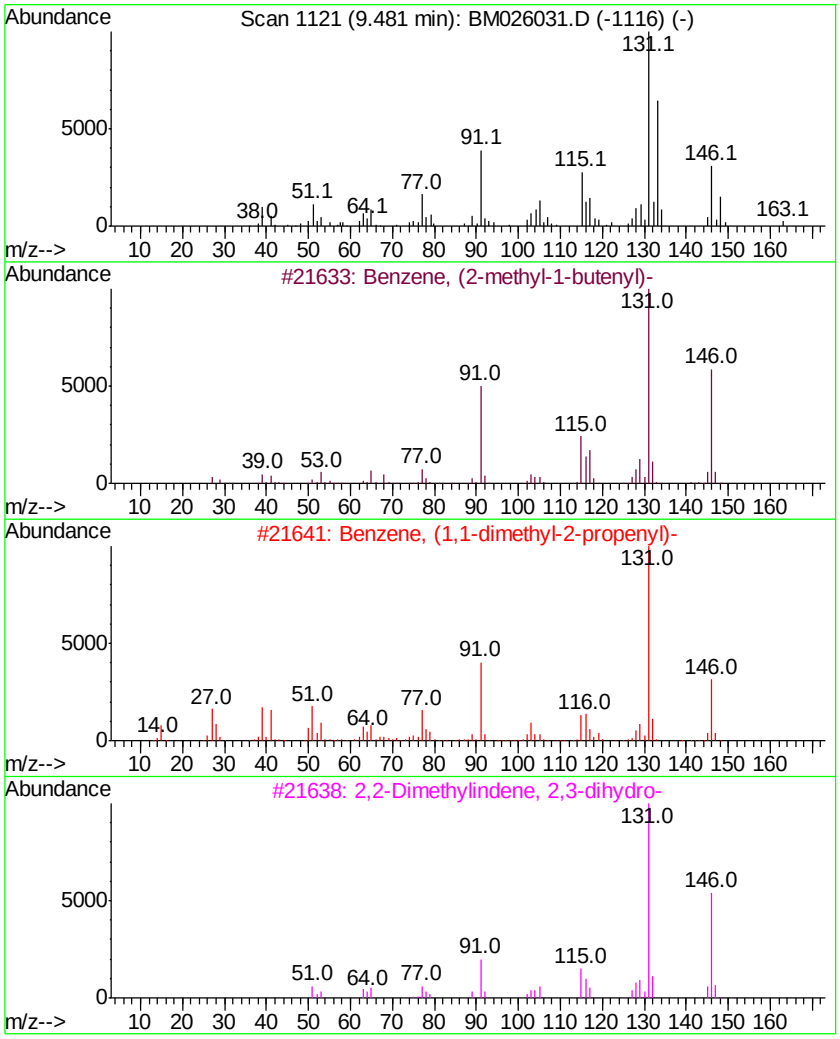
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

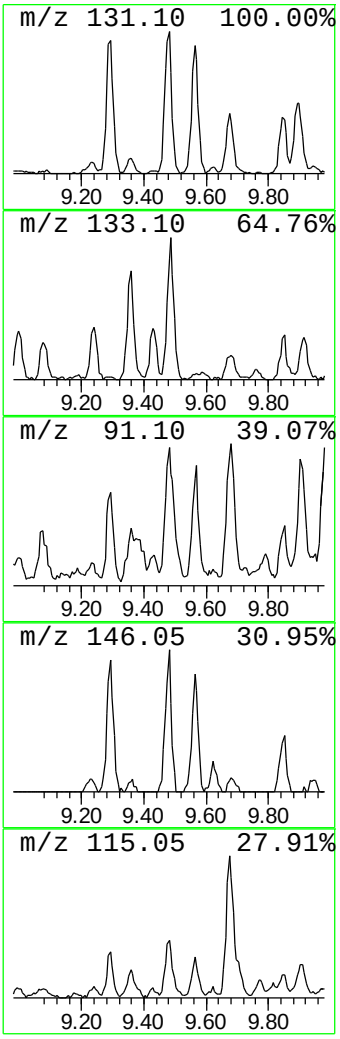
****

Peak Number 2 Benzene, (2-methyl-1-butenyl)- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.48	2.08 ng/ul	194968	Naphthalene-d8	10.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	96
2	Benzene, (1,1-dimethyl-2-propenyl)-	146	C11H14	018321-36-3	91
3	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	89
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	86
5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	86





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051920\
Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

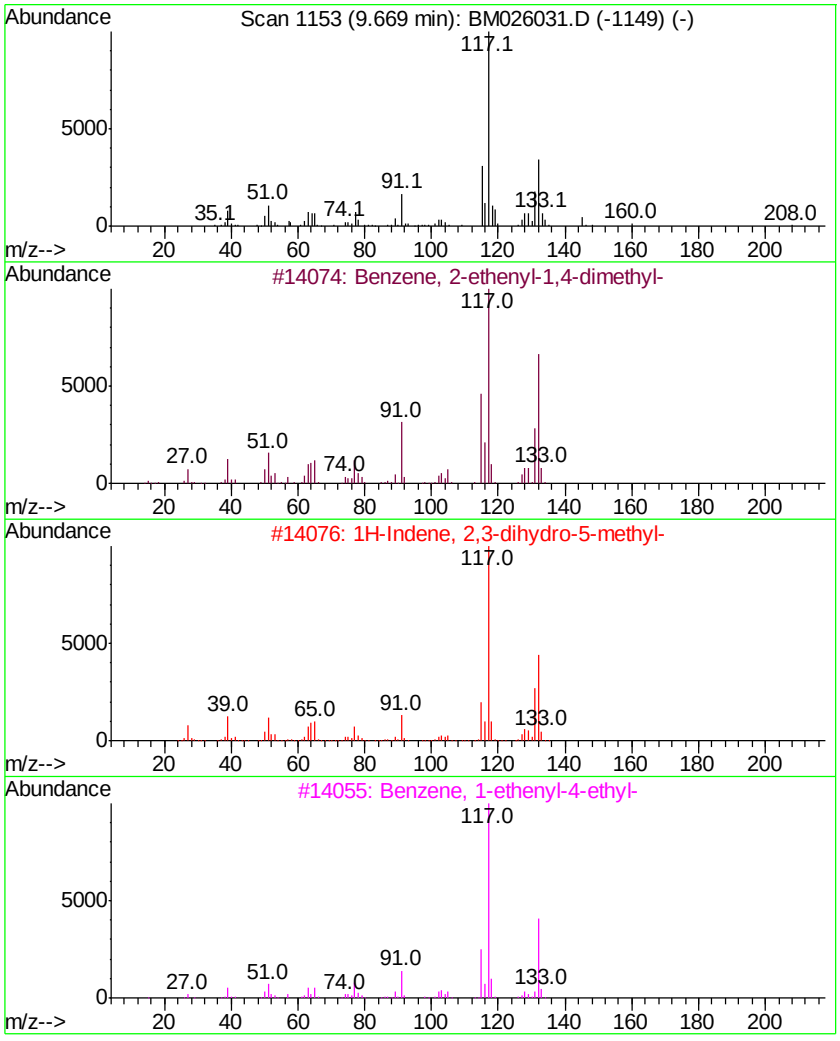
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

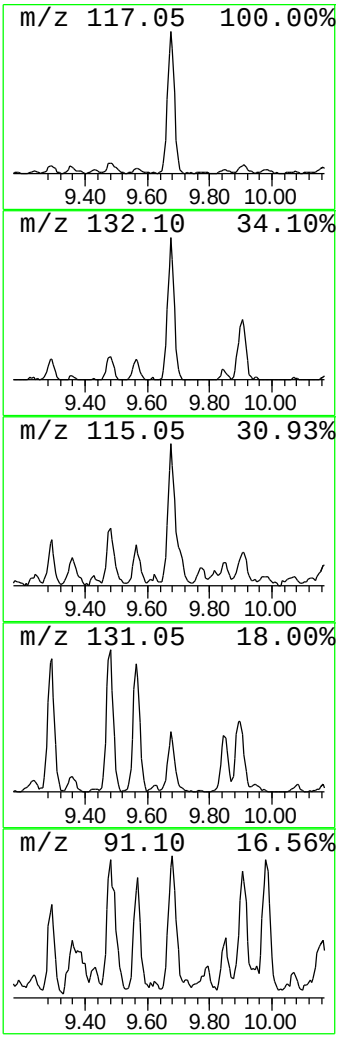
****

Peak Number 3 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.67	2.99 ng/ul	281286	Naphthalene-d8	10.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	94
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	87





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051920\
Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

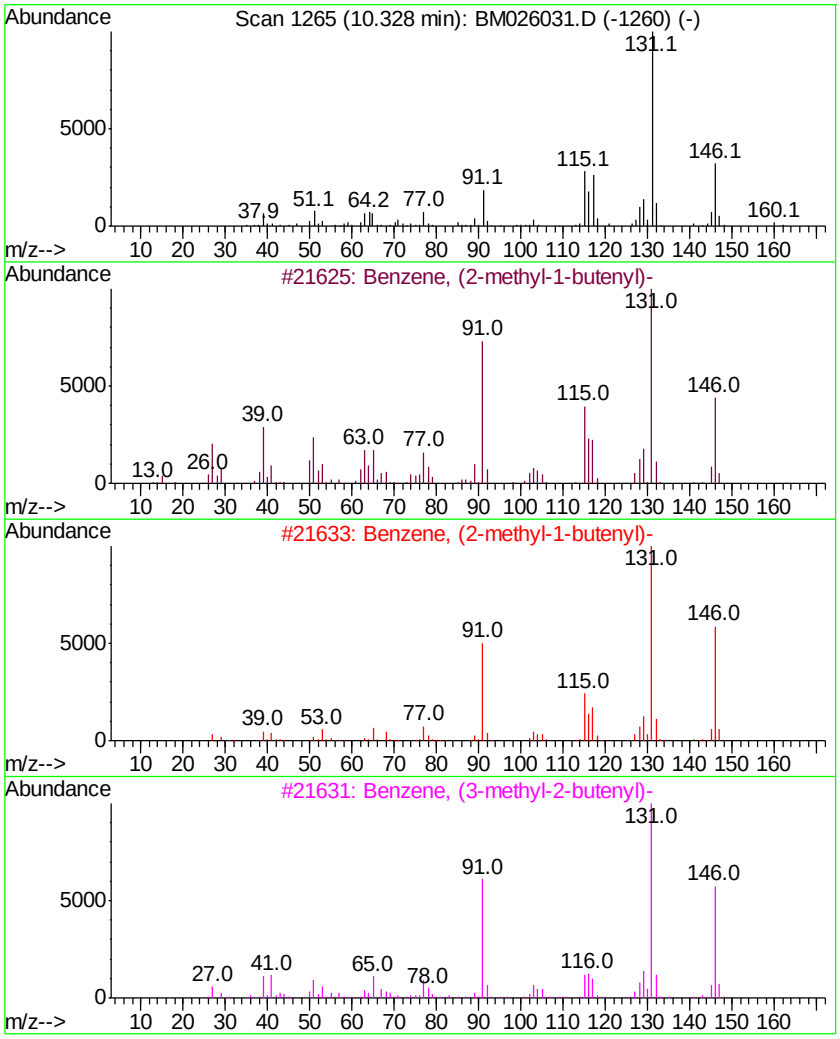
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

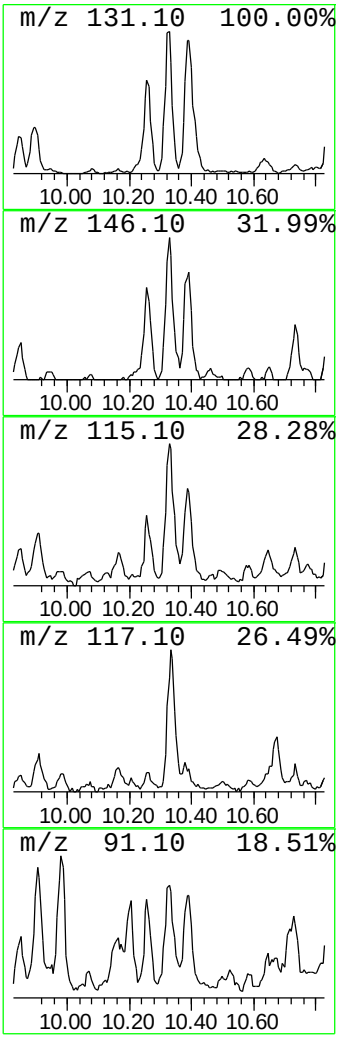
****

Peak Number 4 Benzene, (3-methyl-2-butenyl)- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.33	2.38 ng/ul	223767	Naphthalene-d8	10.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	94
2	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	93
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	87
4	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	87
5	Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	87





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051920\
Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

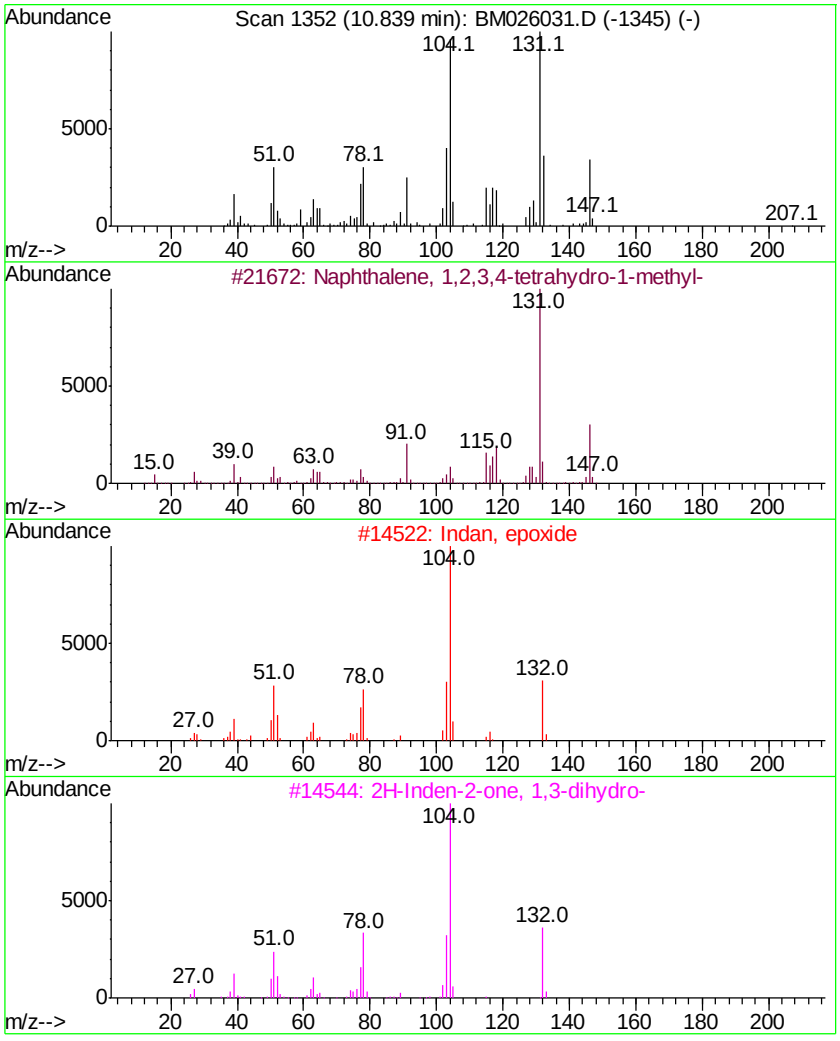
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

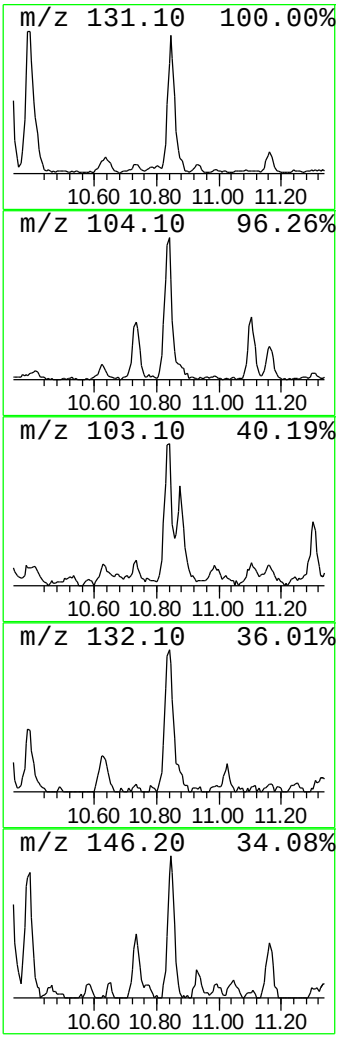
****

Peak Number 5 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.84	2.49 ng/ul	233496	Naphthalene-d8	10.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	90
2	Indan, epoxide	132	C9H8O	000768-22-9	62
3	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	60
4	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	60
5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	60





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Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

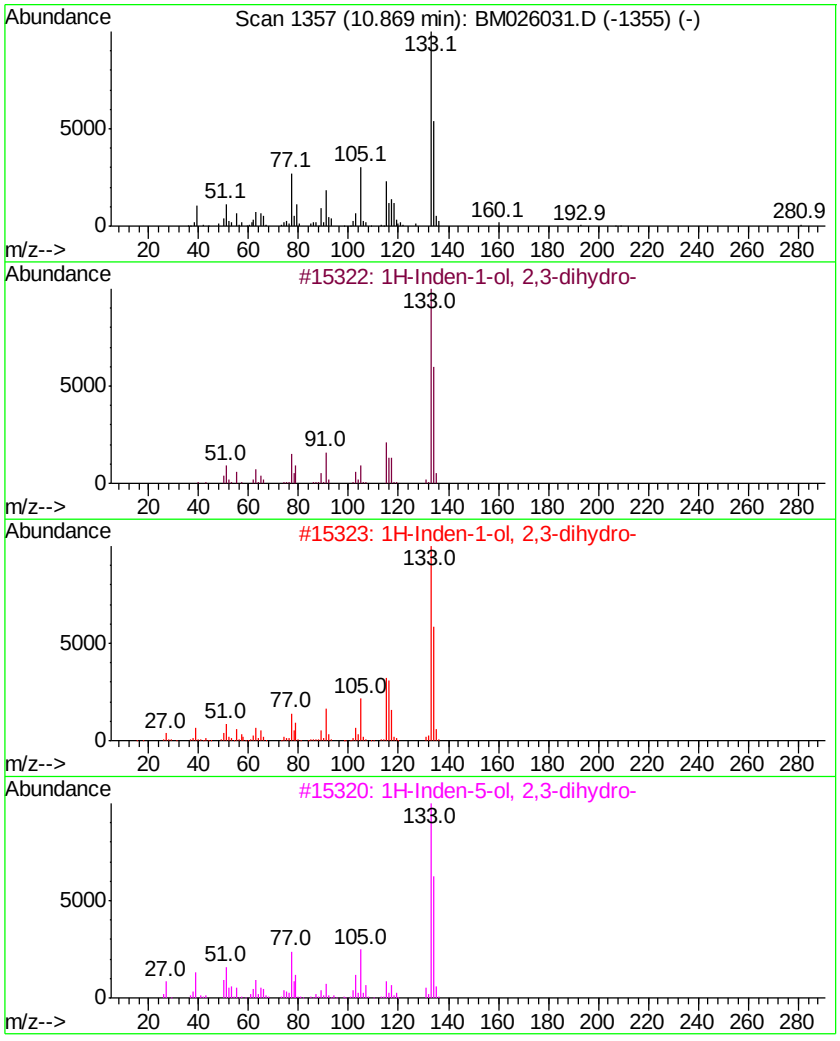
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

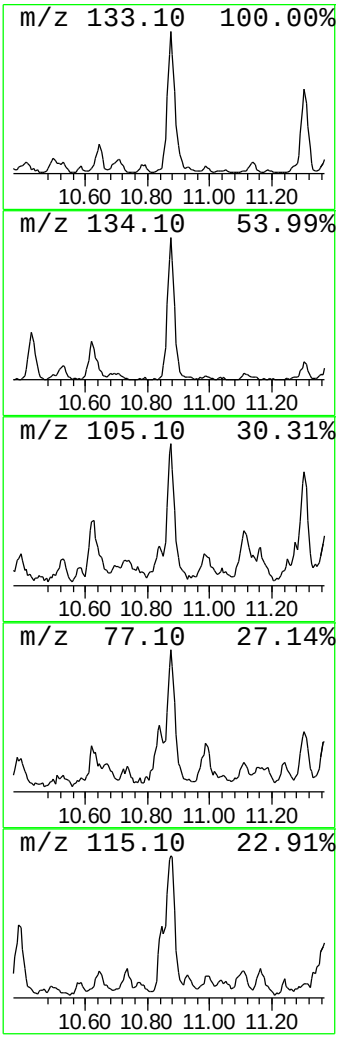
****

Peak Number 6 1H-Inden-1-ol, 2,3-dihydro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.87	3.11 ng/ul	291924	Naphthalene-d8	10.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Inden-1-ol, 2,3-dihydro-	134	C9H10O	006351-10-6	93
2	1H-Inden-1-ol, 2,3-dihydro-	134	C9H10O	006351-10-6	80
3	1H-Inden-5-ol, 2,3-dihydro-	134	C9H10O	001470-94-6	74
4	1H-Inden-1-ol, 2,3-dihydro-	134	C9H10O	006351-10-6	72
5	2,3-Dihydro-2,2,4-trimethyl-1,4-...	189	C12H15NO	026030-81-9	47





Library Search Compound Report

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Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

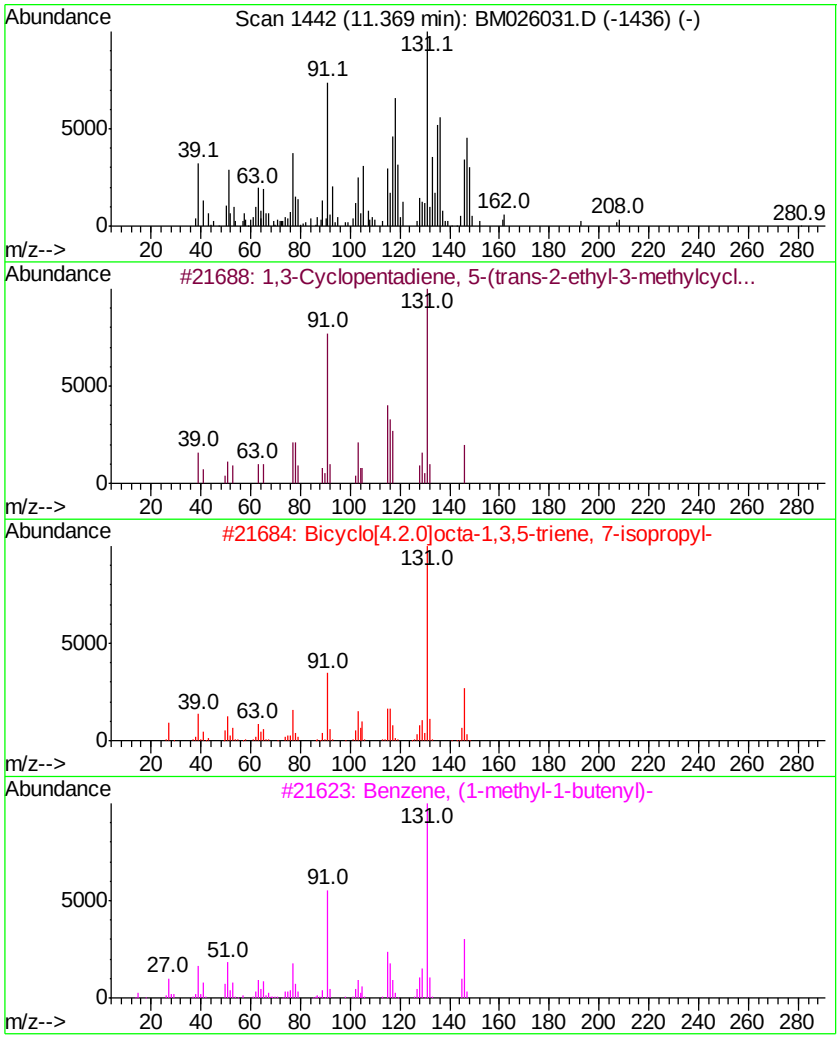
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

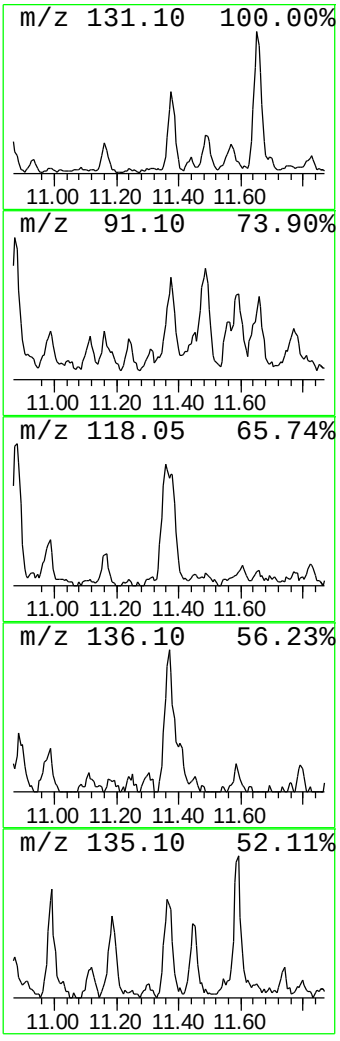
****

Peak Number 7 1,3-Cyclopentadiene, 5-(tra... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.37	2.47 ng/ul	231885	Naphthalene-d8	10.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Cyclopentadiene, 5-(trans-2-...	146	C11H14	079209-36-2	80
2	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	70
3	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	51
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	46
5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	43





Library Search Compound Report

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Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

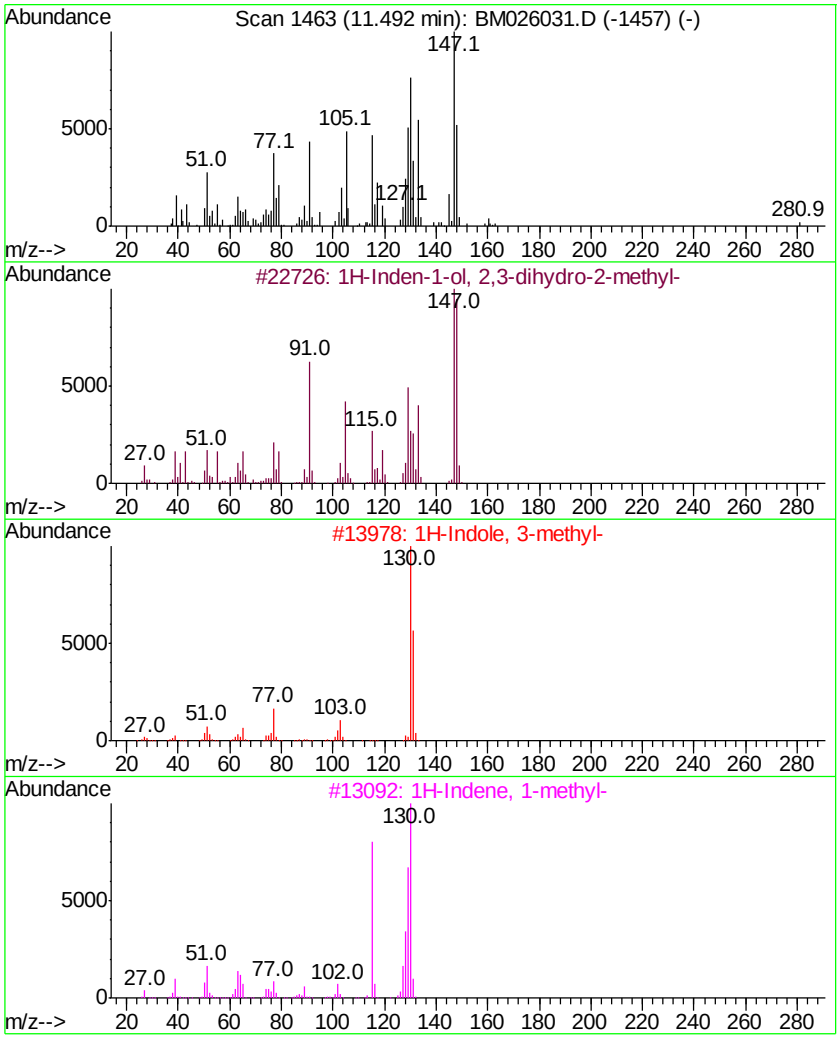
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

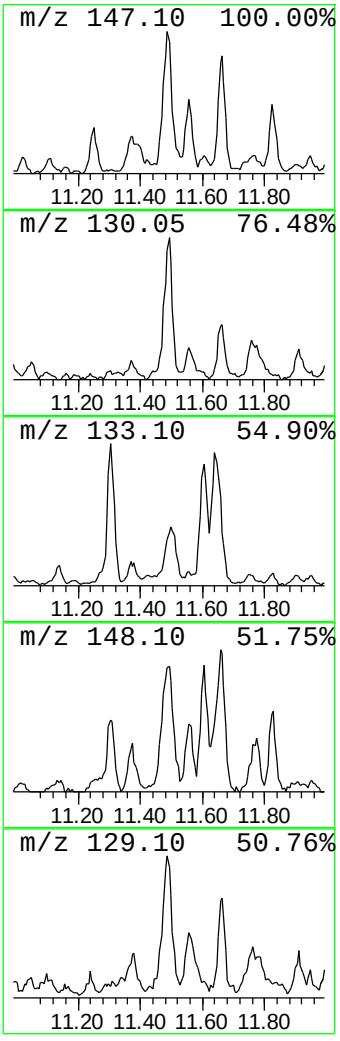
****

Peak Number 8 1H-Inden-1-ol, 2,3-dihydro-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.49	3.05 ng/ul	286733	Naphthalene-d8	10.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Inden-1-ol, 2,3-dihydro-2-met...	148	C10H12O	017496-18-3	52
2	1H-Indole, 3-methyl-	131	C9H9N	000083-34-1	48
3	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	42
4	Benzene, 1,3-butadienyl-	130	C10H10	001515-78-2	42
5	Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	42





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Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

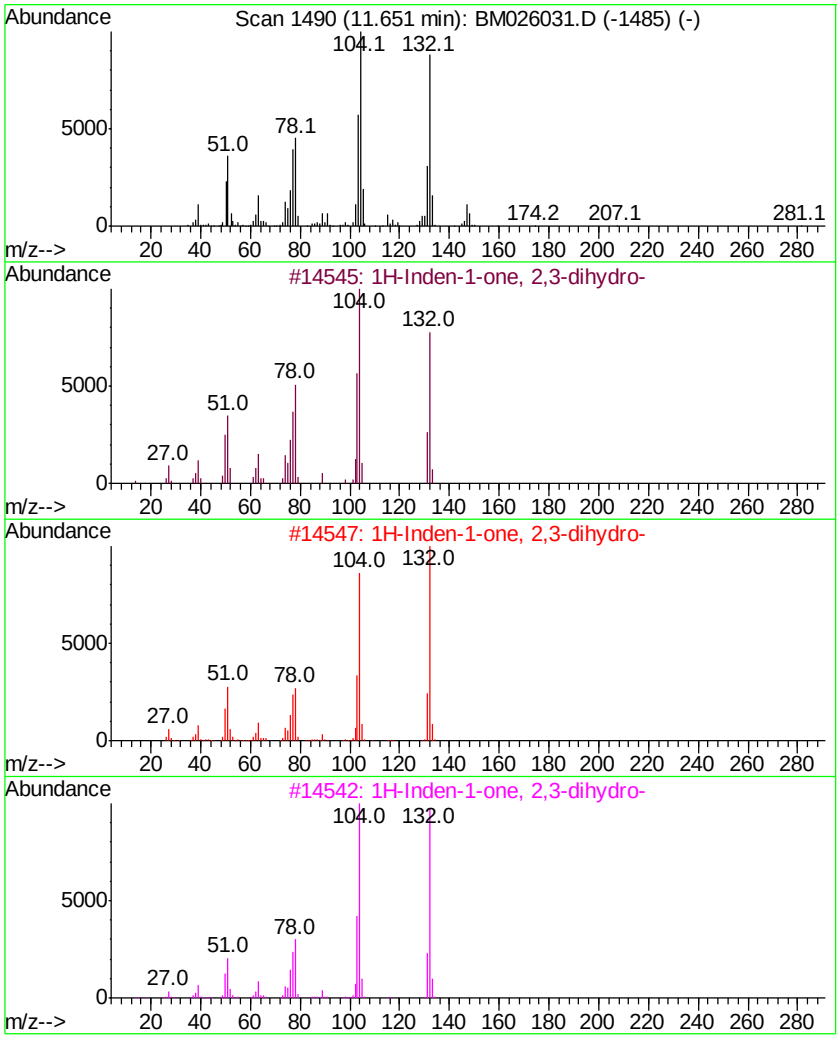
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

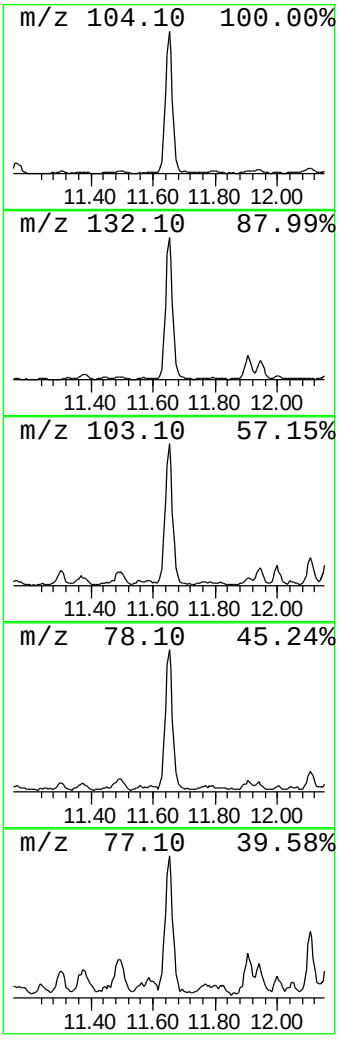
****

Peak Number 9 1H-Inden-1-one, 2,3-dihydro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	7.40 ng/ul	694676	Naphthalene-d8	10.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	93
2	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	93
3	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	81
4	Isoquinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000091-21-4	64
5	Styrene	104	C8H8	000100-42-5	60





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Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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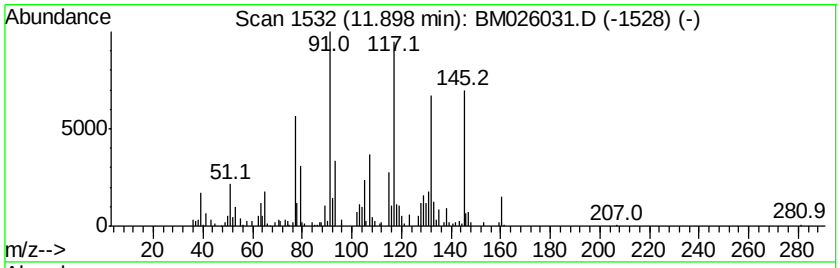
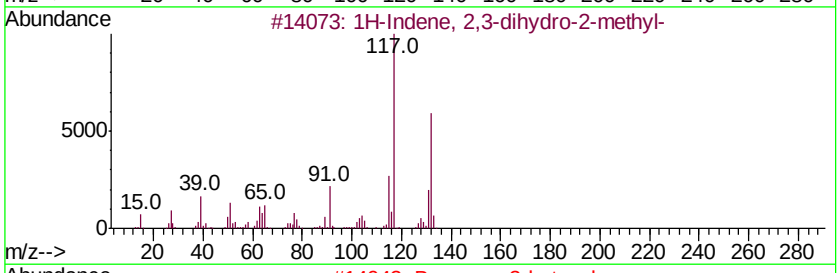
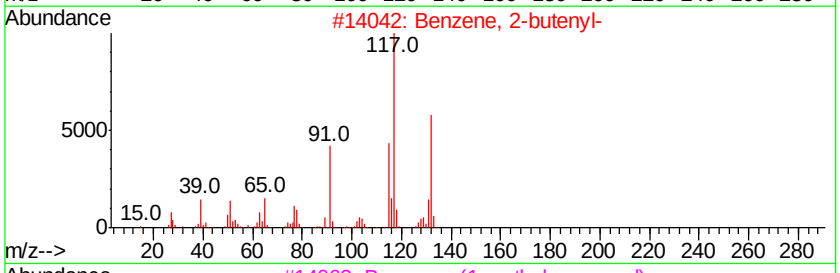
Peak Number 10 1H-Indene, 2,3-dihydro-2-me... Concentration Rank 26

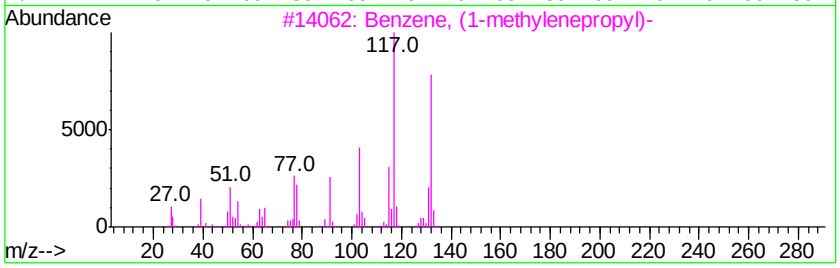
R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	2.07 ng/ul	194596	Naphthalene-d8	10.26

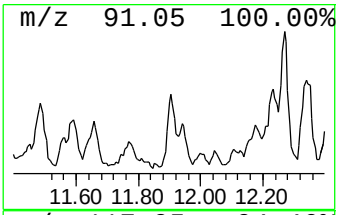
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	86
2	Benzene, 2-butenyl-	132	C10H12	001560-06-1	55
3	Benzene, (1-methylenepropyl)-	132	C10H12	002039-93-2	55
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	55
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	55

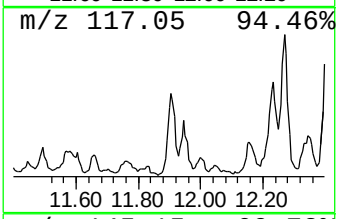


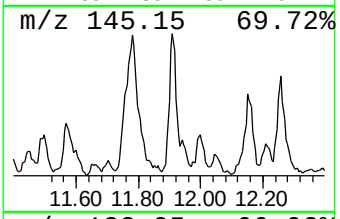


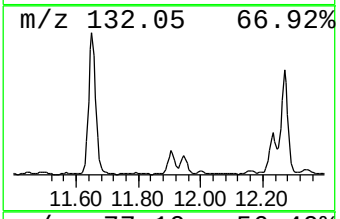


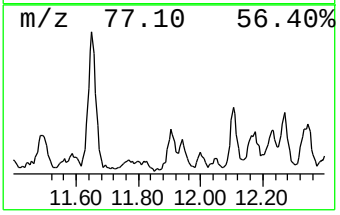












Library Search Compound Report

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Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

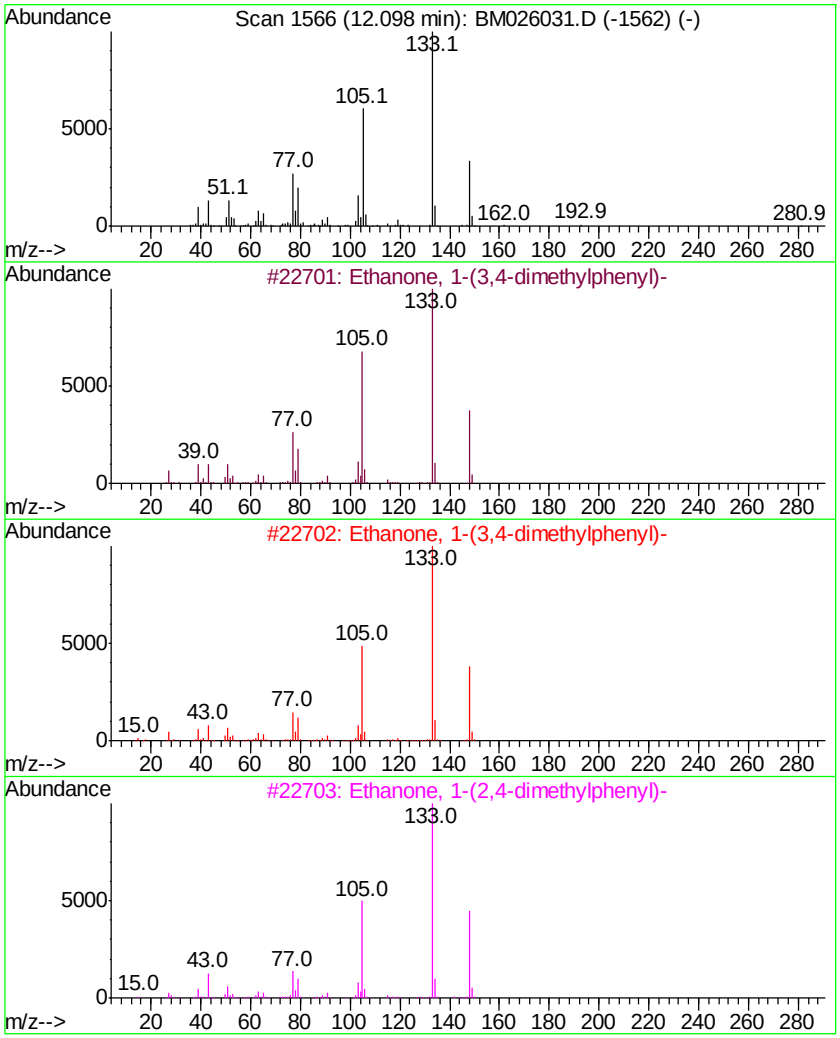
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

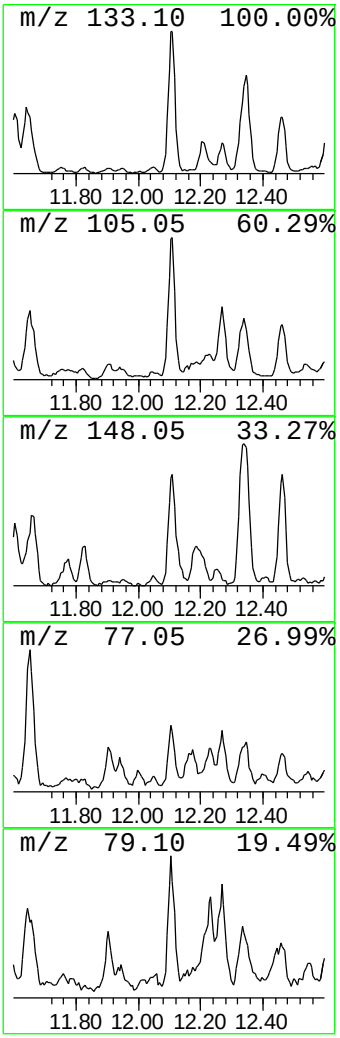
****

Peak Number 11 Ethanone, 1-(3,4-dimethylph... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.10	2.65 ng/ul	249293	Naphthalene-d8	10.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 1-(3,4-dimethylphenyl)-	148	C10H12O	003637-01-2	97
2	Ethanone, 1-(3,4-dimethylphenyl)-	148	C10H12O	003637-01-2	95
3	Ethanone, 1-(2,4-dimethylphenyl)-	148	C10H12O	000089-74-7	95
4	Ethanone, 1-(2,4-dimethylphenyl)-	148	C10H12O	000089-74-7	94
5	Ethanone, 1-(2,5-dimethylphenyl)-	148	C10H12O	002142-73-6	94





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Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

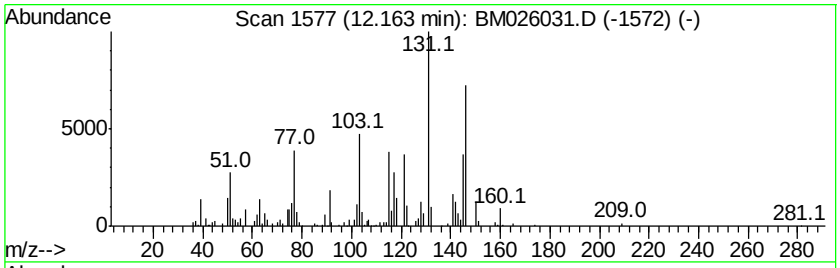
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

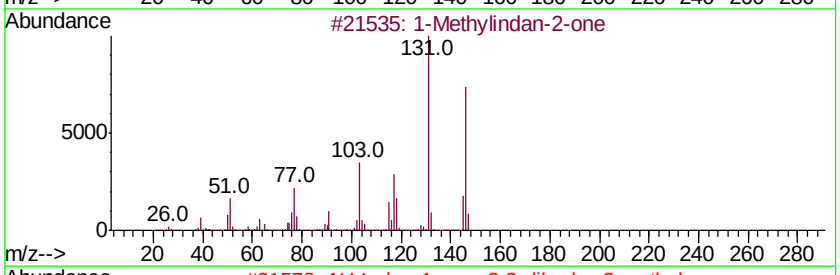
****

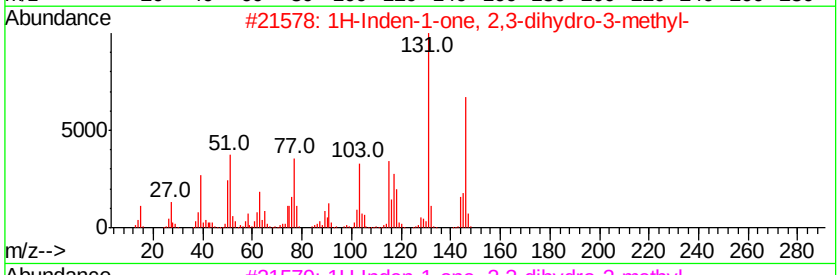
Peak Number 12 1-Methylindan-2-one Concentration Rank 24

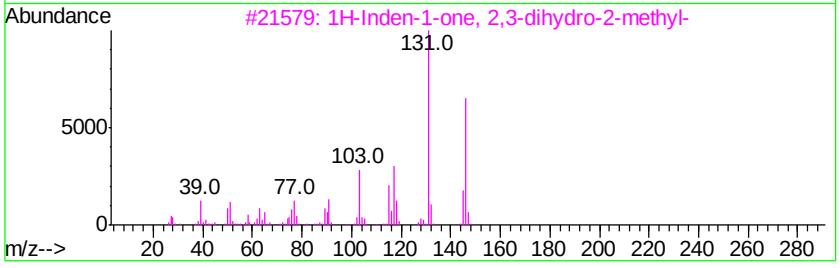
R.T.	EstConc	Area	Relative to ISTD	R.T.
12.16	2.08 ng/ul	195320	Naphthalene-d8	10.26

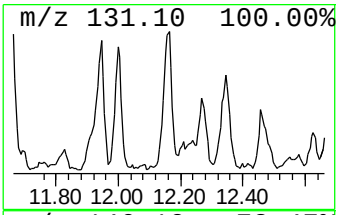
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Methylindan-2-one	146	C10H10O	035587-60-1	76
2	1H-Inden-1-one, 2,3-dihydro-3-me...	146	C10H10O	006072-57-7	74
3	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	46
4	2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	46
5	3-Buten-2-one, 4-phenyl-	146	C10H10O	000122-57-6	45

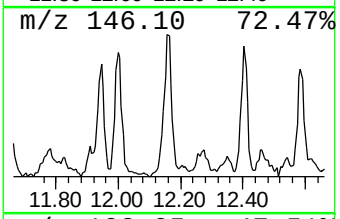


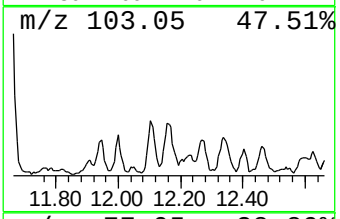


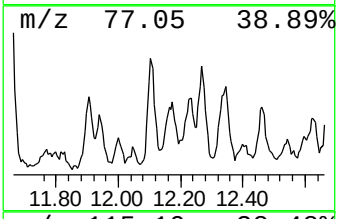


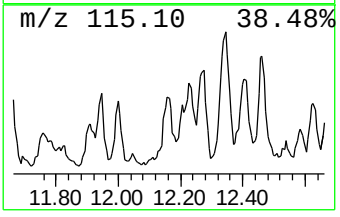












Library Search Compound Report

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Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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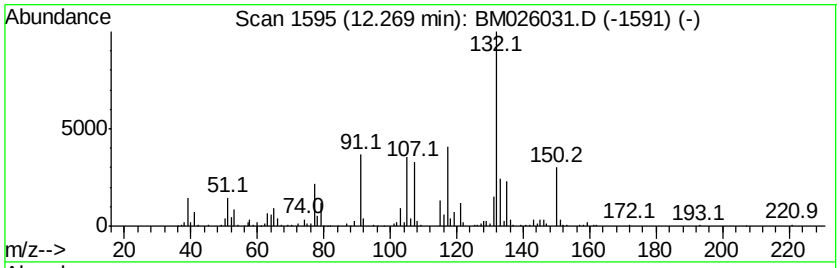
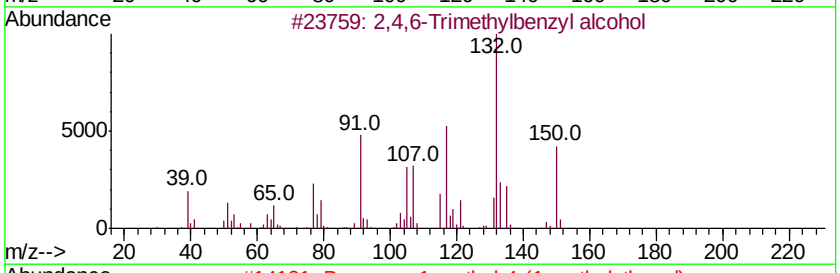
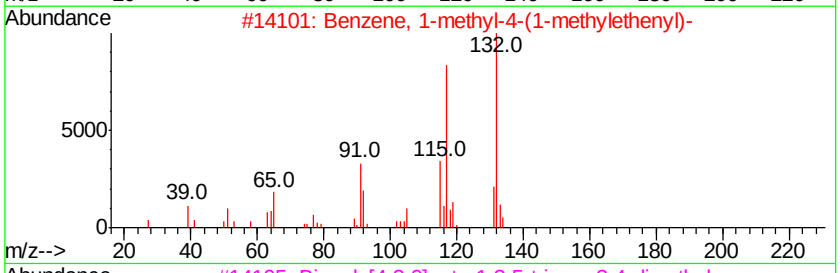
Peak Number 13 2,4,6-Trimethylbenzyl alcohol Concentration Rank 9

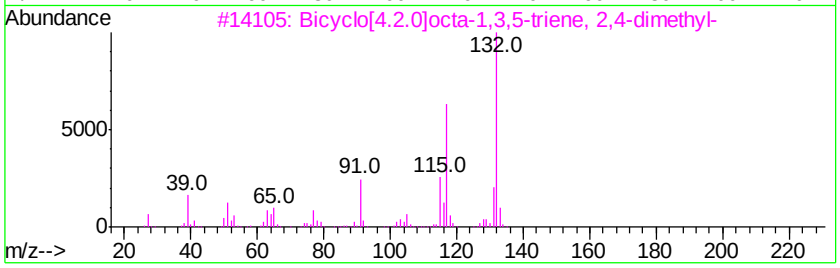
R.T.	EstConc	Area	Relative to ISTD	R.T.
12.27	3.24 ng/ul	381421	Acenaphthene-d10	14.15

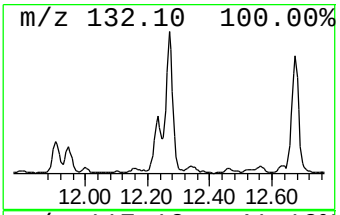
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,6-Trimethylbenzyl alcohol	150	C10H14O	1000342-65-8	94
2	Benzene, 1-methyl-4-(1-methyleth...	132	C10H12	001195-32-0	55
3	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	50
4	Benzene, 4-ethenyl-1,2-dimethyl-	132	C10H12	027831-13-6	46
5	2-Cyanomethyl-1,2,3,4-tetrahydro...	172	C11H12N2	005100-75-4	38

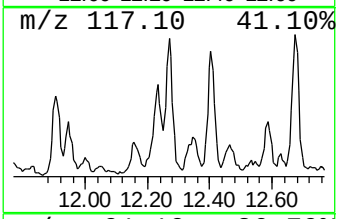


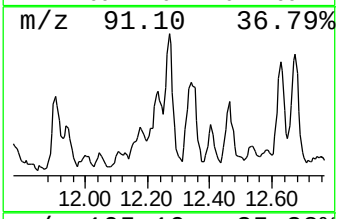


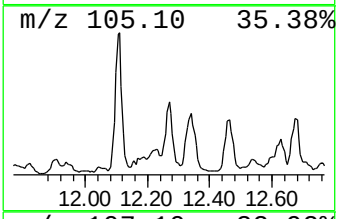


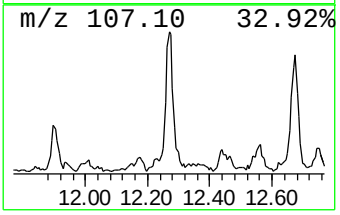












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Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

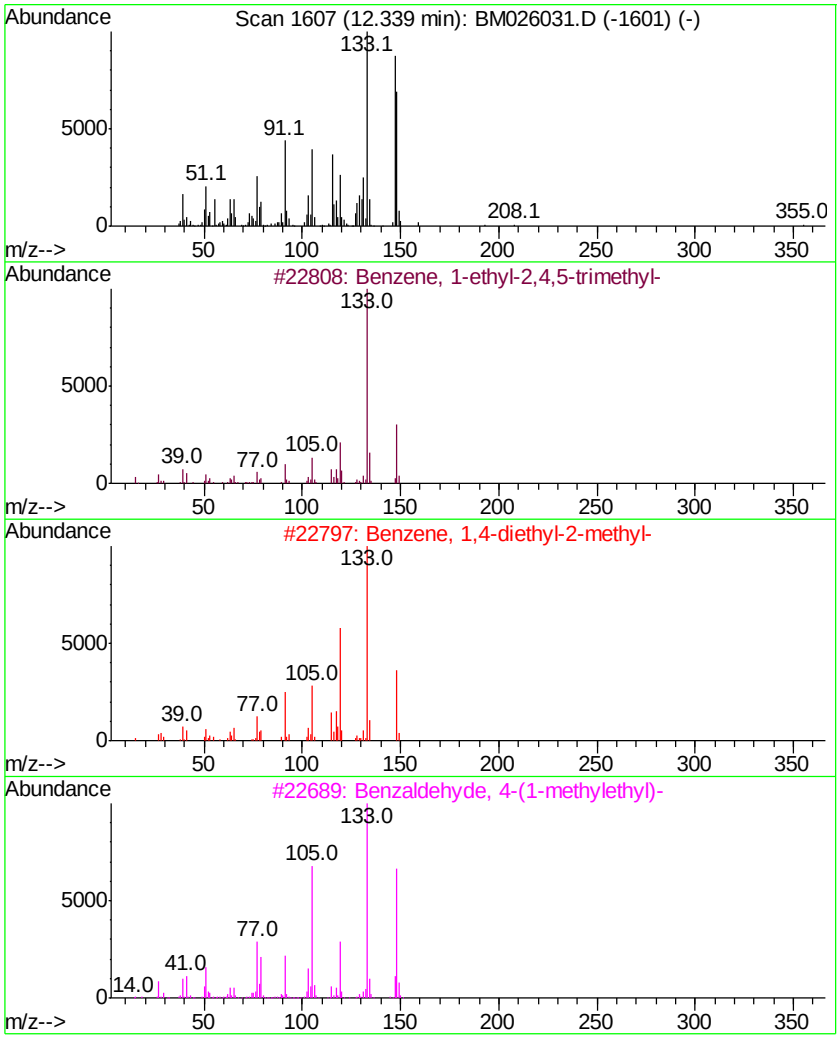
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

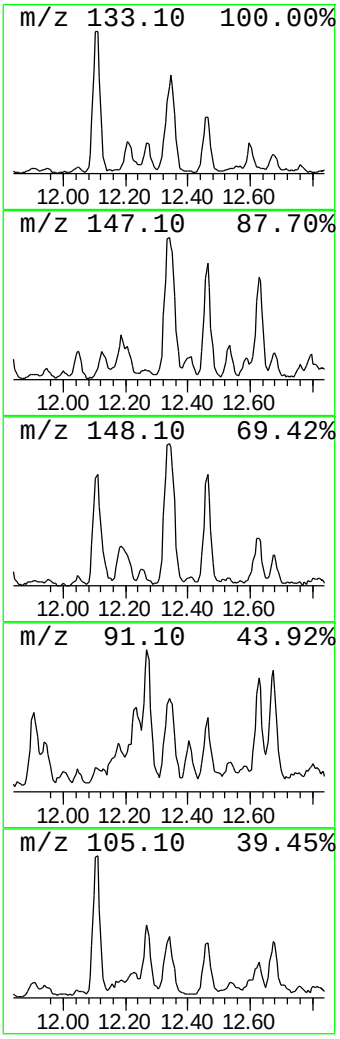
****

Peak Number 14 Benzene, 1-ethyl-2,4,5-trim... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.34	3.51 ng/ul	413140	Acenaphthene-d10	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4,5-trimethyl-	148	C11H16	017851-27-3	55
2	Benzene, 1,4-diethyl-2-methyl-	148	C11H16	013632-94-5	55
3	Benzaldehyde, 4-(1-methylethyl)-	148	C10H12O	000122-03-2	55
4	1H-Inden-1-ol, 2,3-dihydro-2-met...	148	C10H12O	017496-18-3	53
5	Pyrrole-3-carbonitrile, 5-formyl...	148	C8H8N2O	032487-71-1	50





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Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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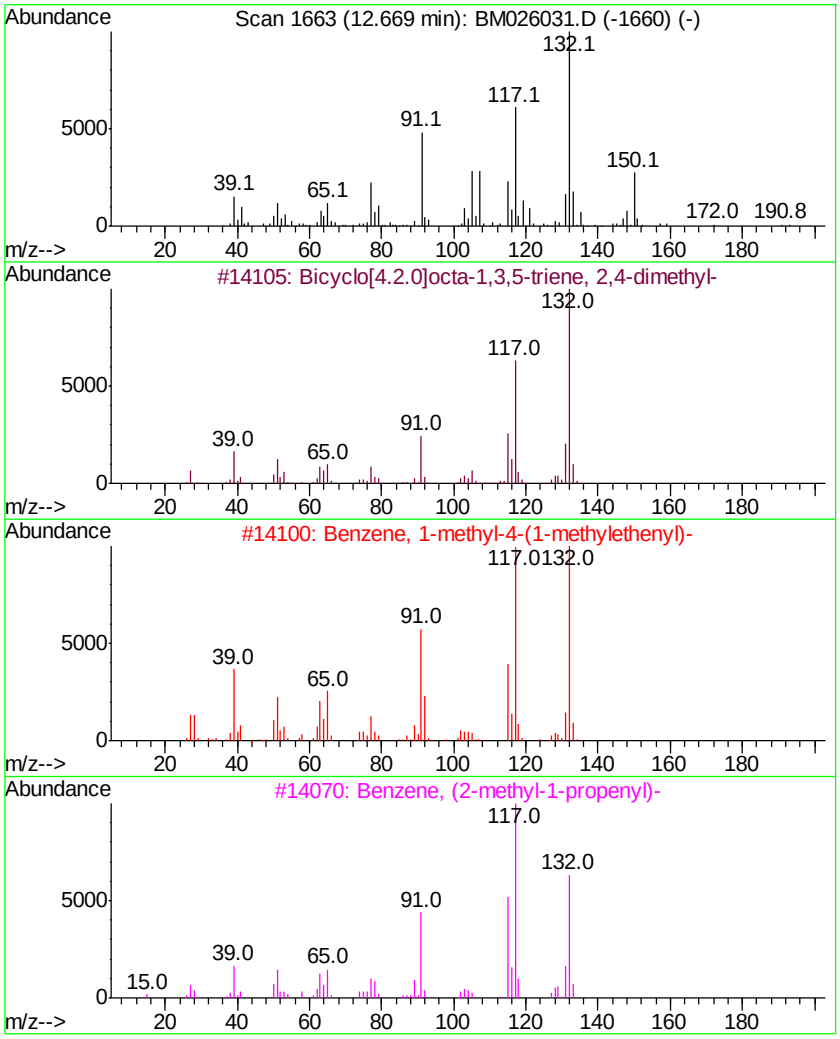
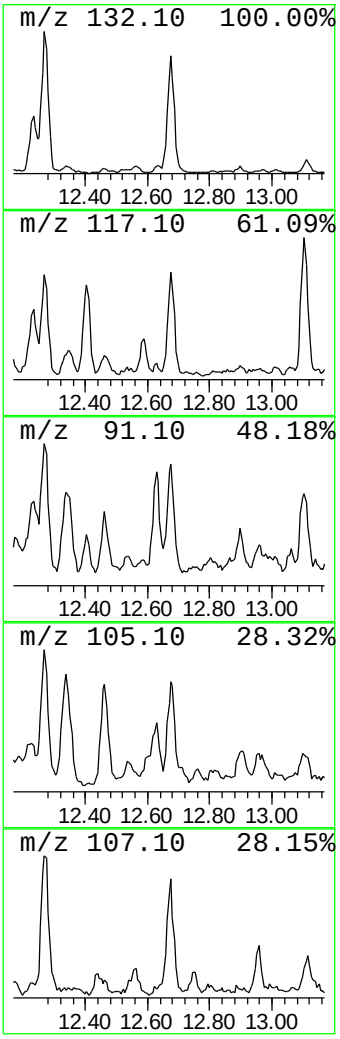
Peak Number 15 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	2.46 ng/ul	289734	Acenaphthene-d10	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	64
2	Benzene, 1-methyl-4-(1-methyleth...	132	C10H12	001195-32-0	64
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	64
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	64
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	58





Library Search Compound Report

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Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

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Quant Title : SVOA CALIBRATION

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

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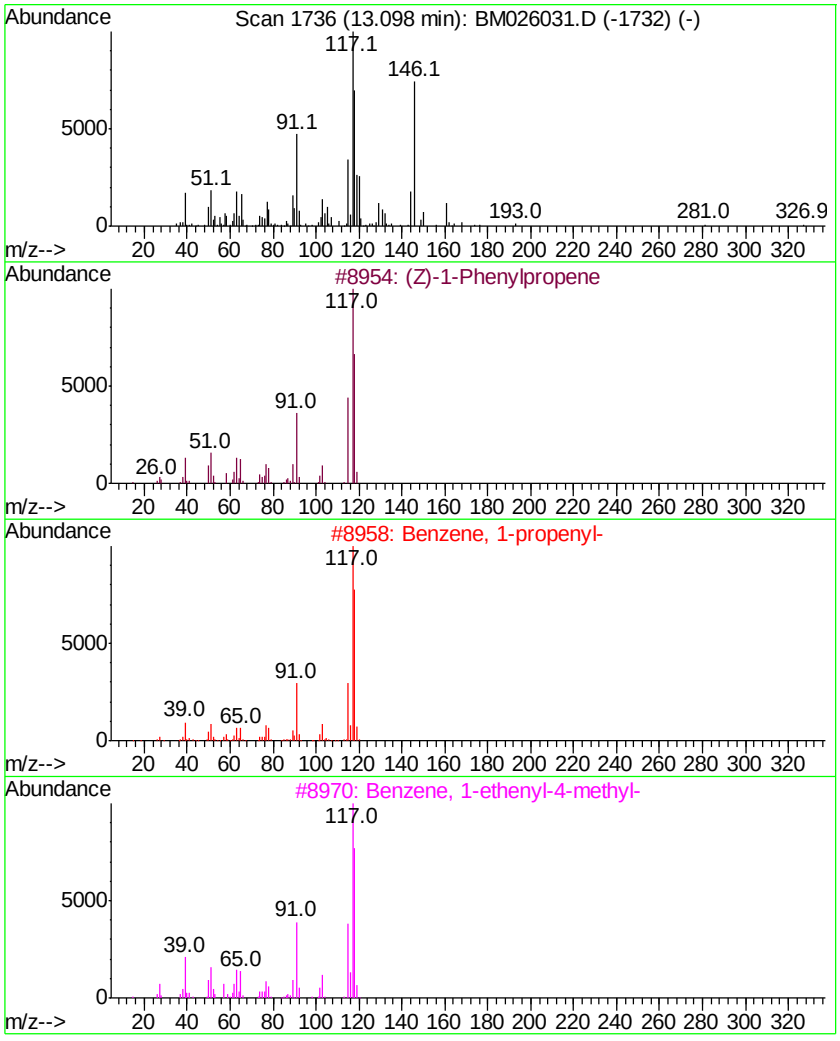
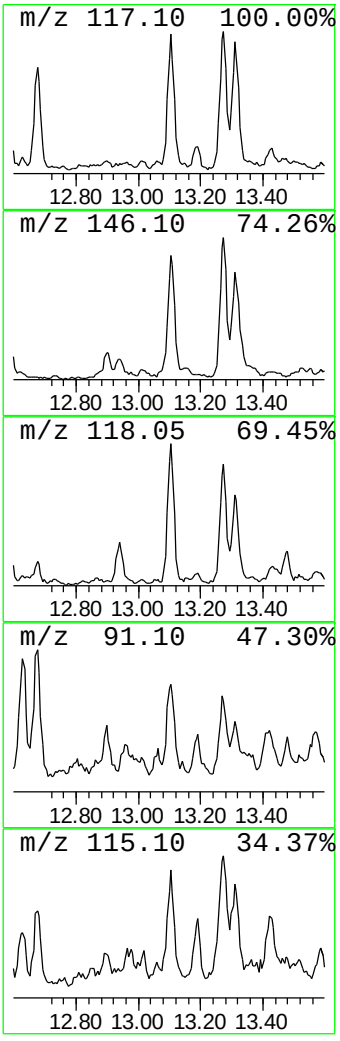
Peak Number 16 (Z)-1-Phenylpropene Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.10	2.30 ng/ul	270650	Acenaphthene-d10	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	90
2	Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
3	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	64
4	Benzene, cyclopentyl-	146	C11H14	000700-88-9	64
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64





Library Search Compound Report

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Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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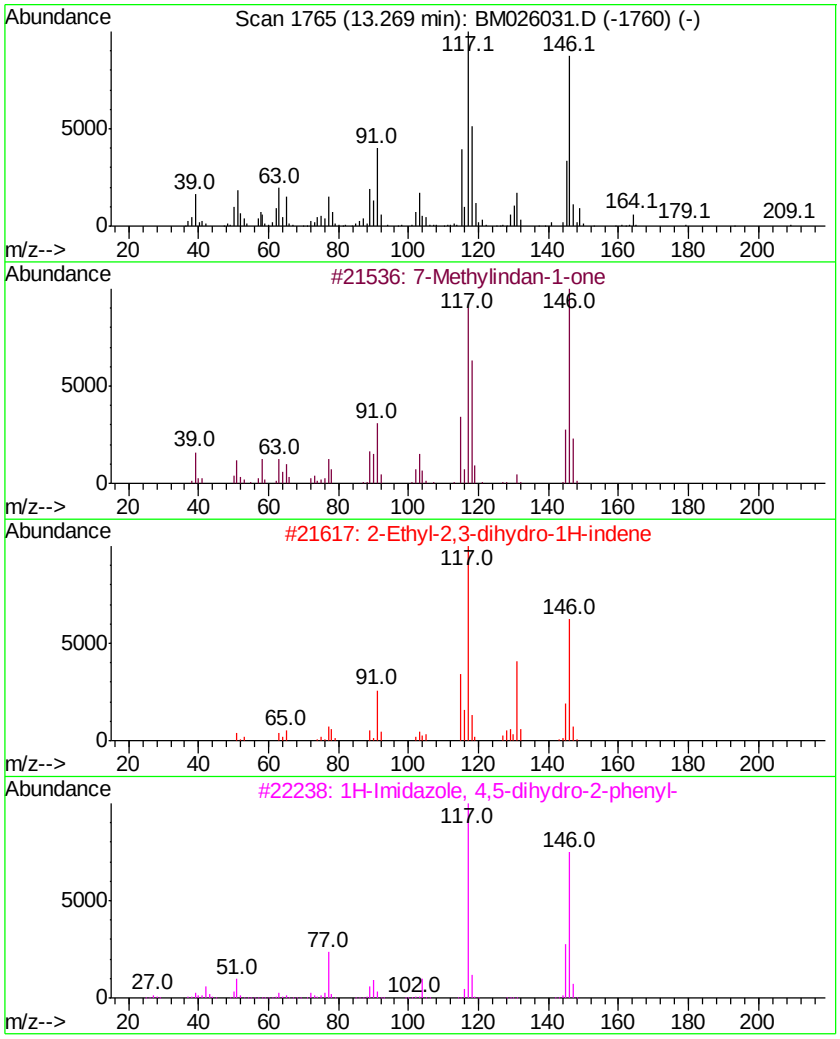
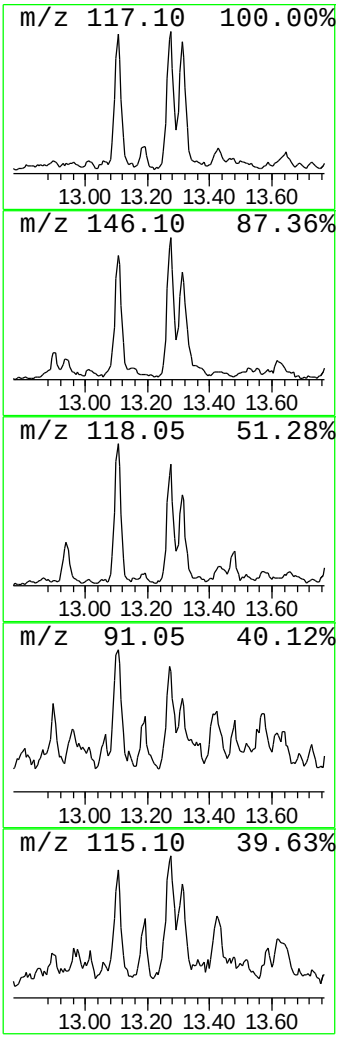
Peak Number 17 7-Methylindan-1-one Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.27	2.60 ng/ul	305869	Acenaphthene-d10	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Methylindan-1-one	146	C10H10O	039627-61-7	80
2	2-Ethyl-2,3-dihydro-1H-indene	146	C11H14	056147-63-8	68
3	1H-Imidazole, 4,5-dihydro-2-phenyl-	146	C9H10N2	000936-49-2	60
4	Bicyclo[4.2.1]nona-2,4,7-triene,...	146	C11H14	1000164-42-6	53
5	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	42





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Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

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Quant Title : SVOA CALIBRATION

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

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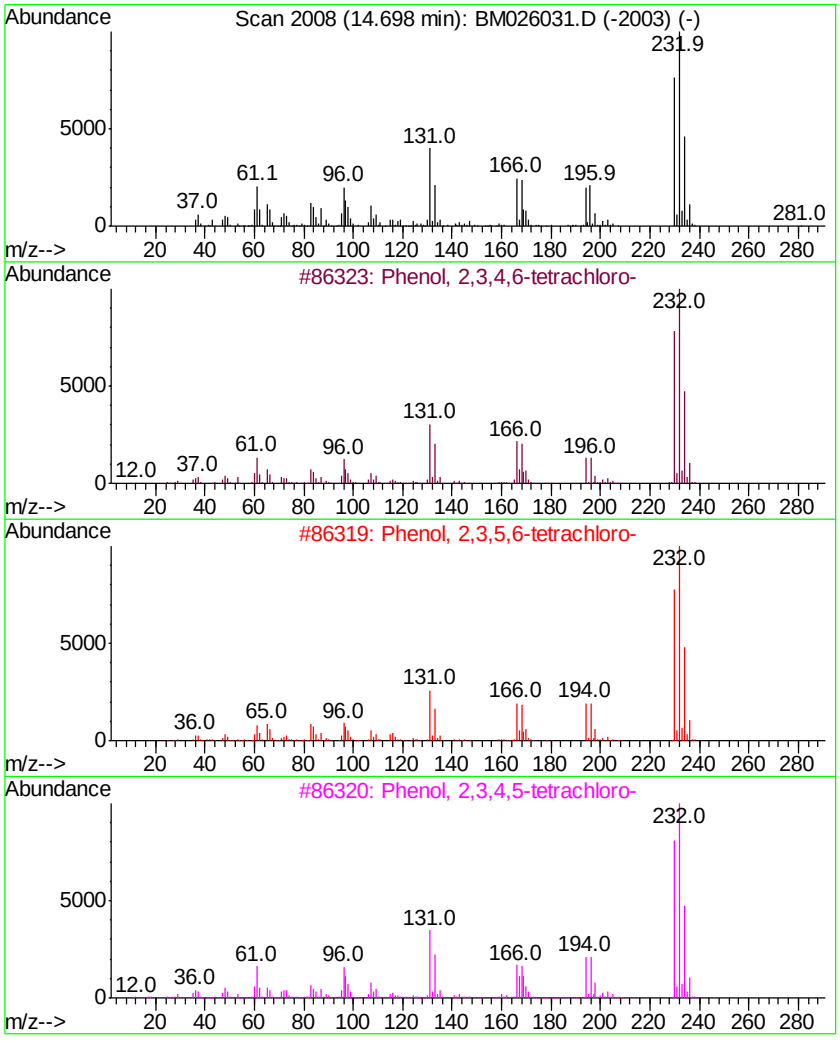
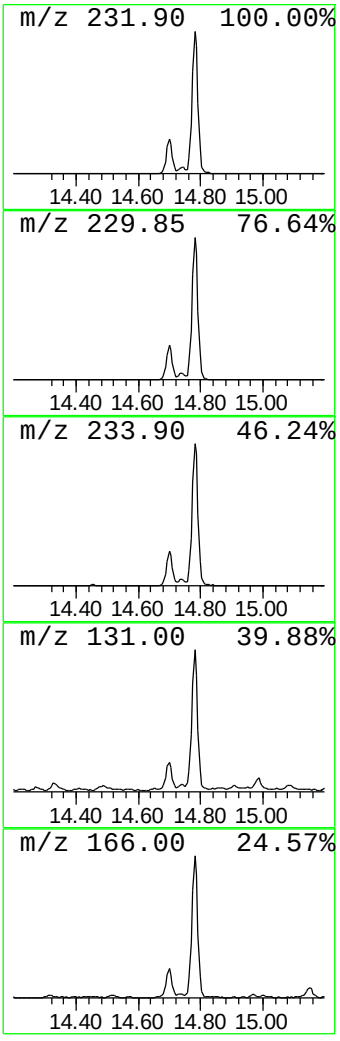
Peak Number 18 Phenol, 2,3,4,6-tetrachloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.70	4.40 ng/ul	517292	Acenaphthene-d10	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
2	Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	96
3	Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	96
4	Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	96
5	Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	95





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Sample : L2681-07
Misc :
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Instrument : BNA_M
ClientSampleId : C5CA9

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Quant Title : SVOA CALIBRATION

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

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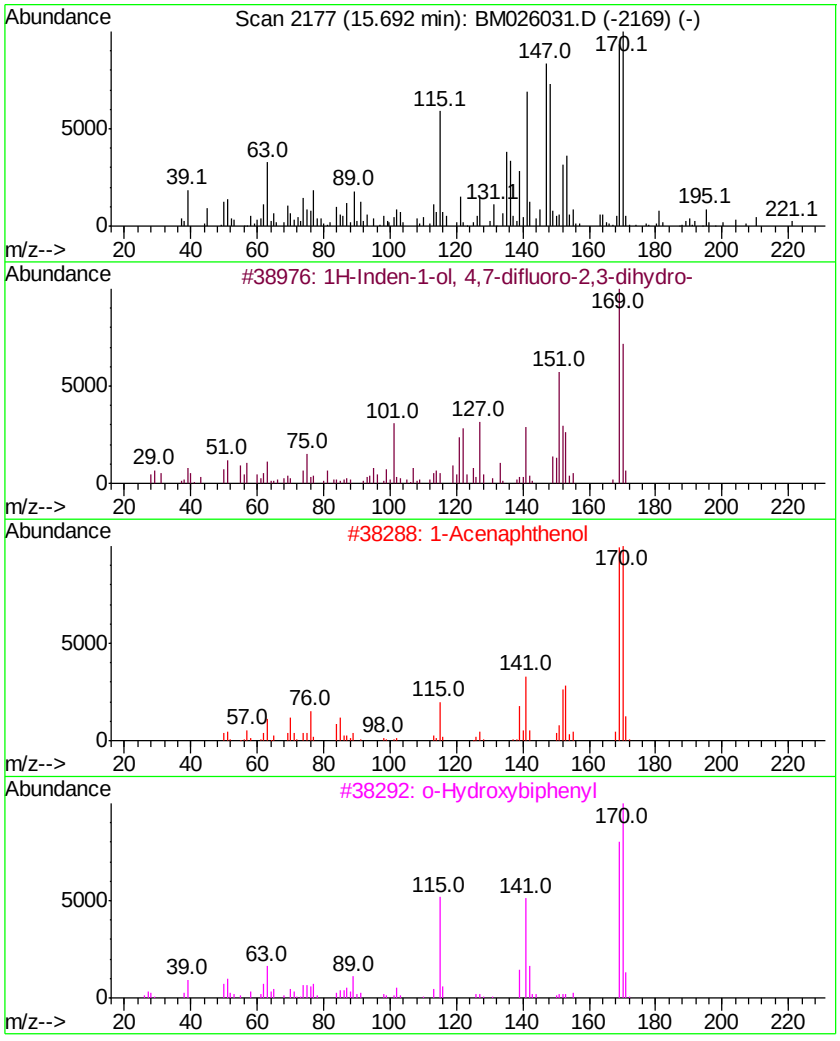
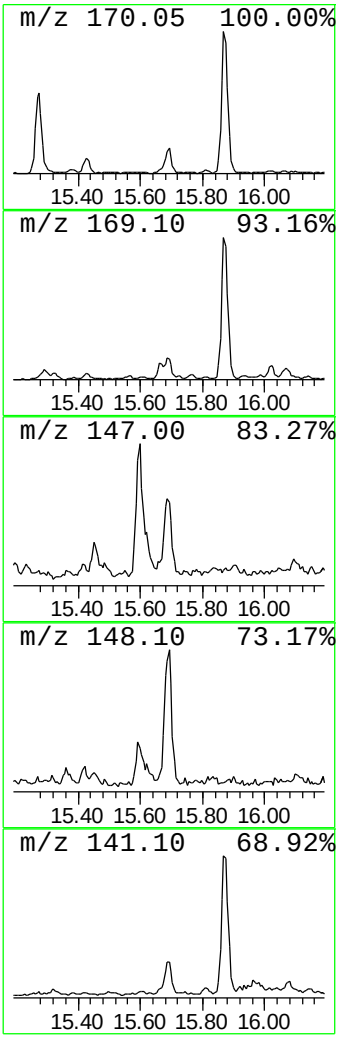
Peak Number 19 unknown-01 Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.69	2.11 ng/ul	265376	Phenanthrene-d10	16.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Inden-1-ol, 4,7-difluoro-2,3-...	170	C9H8F2O	130408-17-2	43
2	1-Acenaphthenol	170	C12H10O	006306-07-6	38
3	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	38
4	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	30
5	1-Naphthalenecarboxaldehyde, 4-m...	170	C12H10O	033738-48-6	25





Library Search Compound Report

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Sample : L2681-07
Misc :
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Instrument : BNA_M
ClientSampleId : C5CA9

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Quant Title : SVOA CALIBRATION

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

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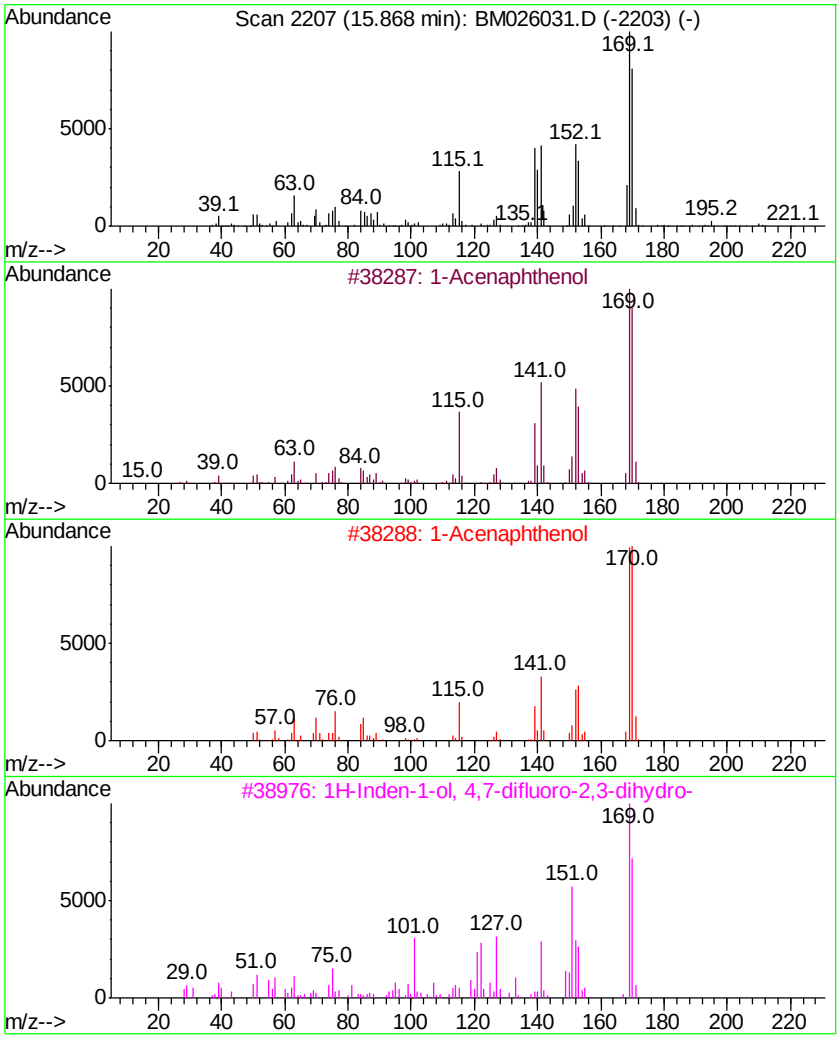
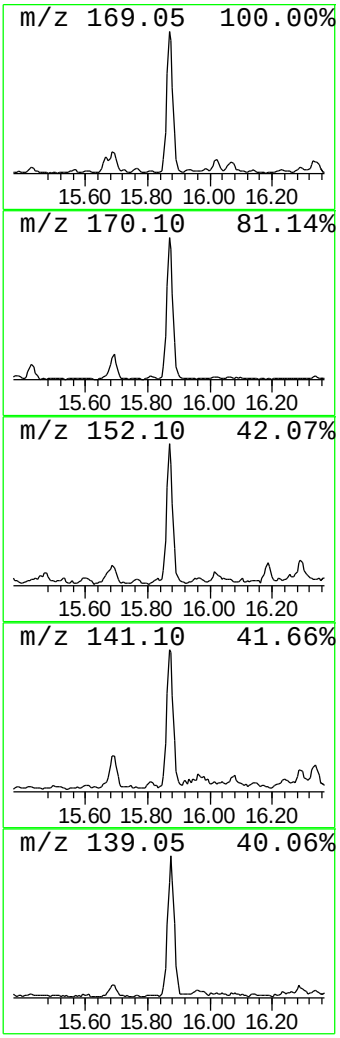
Peak Number 20 1-Acenaphthenol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.87	4.96 ng/ul	623927	Phenanthrene-d10	16.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Acenaphthenol	170	C12H10O	006306-07-6	96
2	1-Acenaphthenol	170	C12H10O	006306-07-6	95
3	1H-Inden-1-ol, 4,7-difluoro-2,3-...	170	C9H8F2O	130408-17-2	62
4	1-Acenaphthenol	170	C12H10O	006306-07-6	60
5	5-Chloro-2-methoxy-2,4,6-cyclohe...	170	C8H7ClO2	013187-38-7	38





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Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

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Quant Title : SVOA CALIBRATION

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

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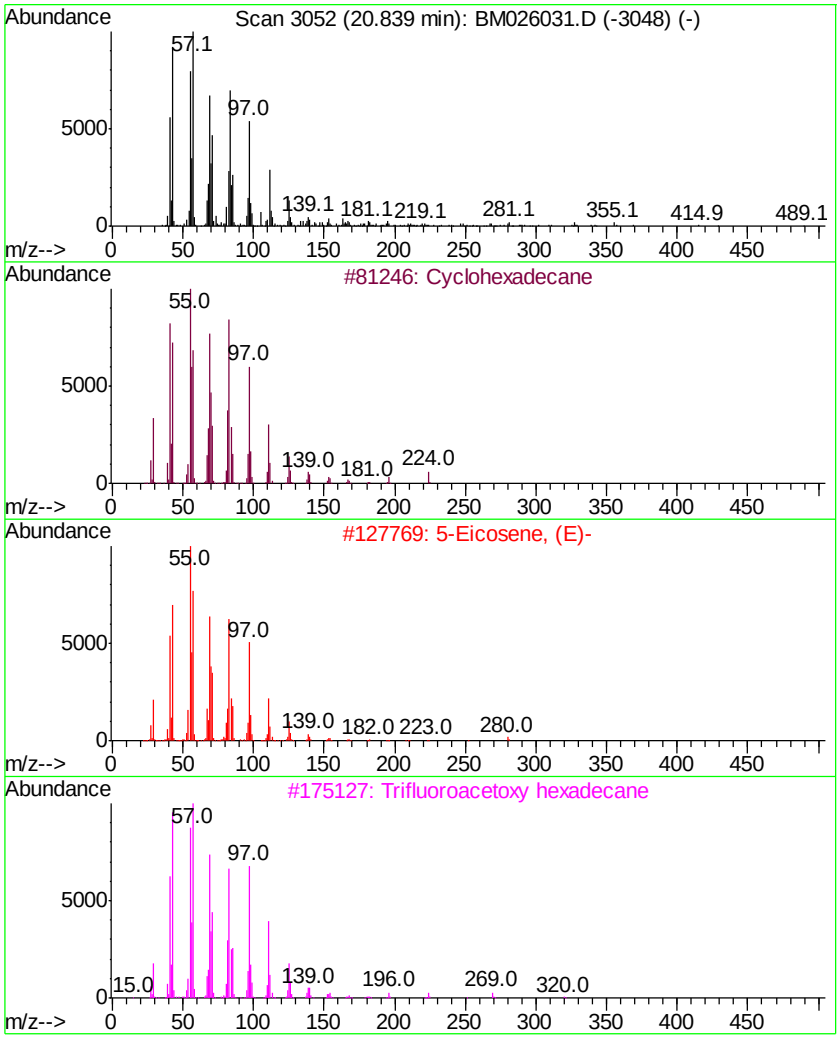
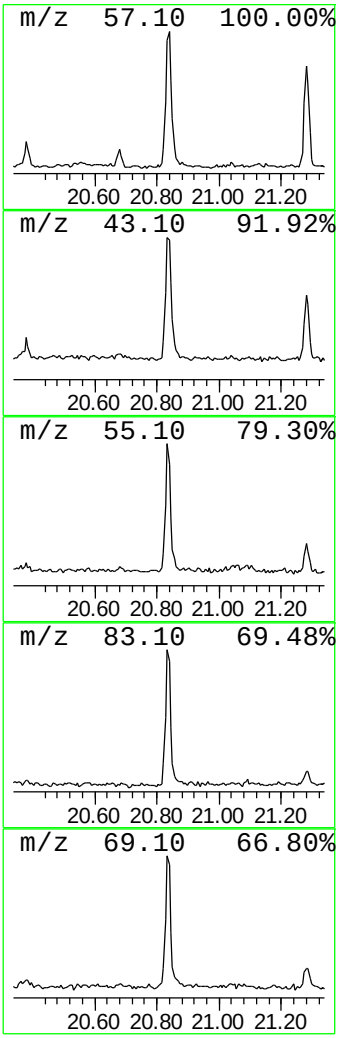
Peak Number 21 (DEL) Alkane: Cyclic20.84 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.84	2.64 ng/ul	367707	Chrysene-d12	21.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	98
2	5-Eicosene, (E)-	280	C20H40	074685-30-6	97
3	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
4	Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	94
5	Cyclotetracosane	336	C24H48	000297-03-0	94





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ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

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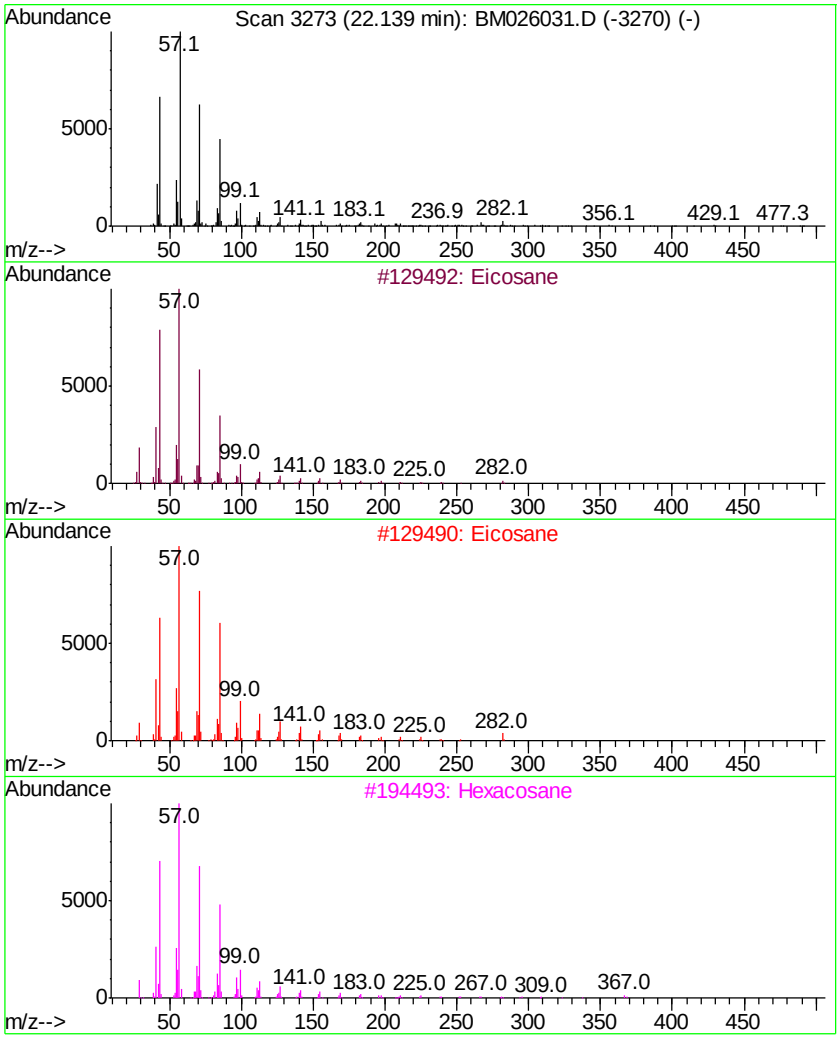
Quant Title : SVOA CALIBRATION

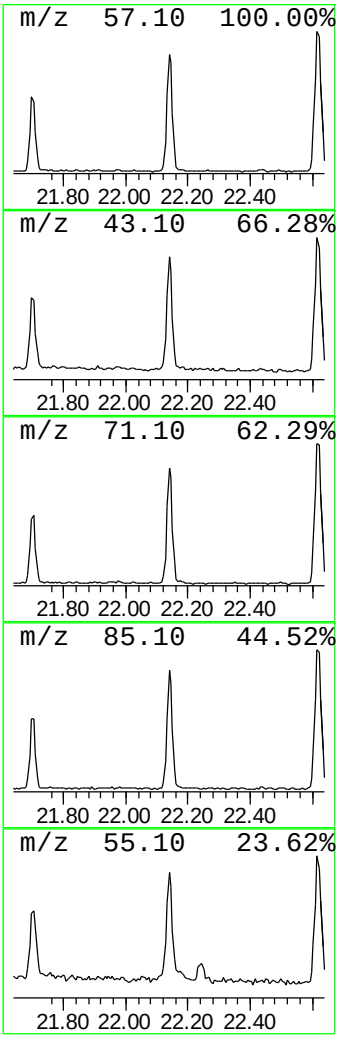
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.14	2.34 ng/ul	326152	Chrysene-d12	21.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	97
2	Eicosane	282	C20H42	000112-95-8	95
3	Hexacosane	366	C26H54	000630-01-3	90
4	Heptacosane	380	C27H56	000593-49-7	90
5	Tetracosane	338	C24H50	000646-31-1	90





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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

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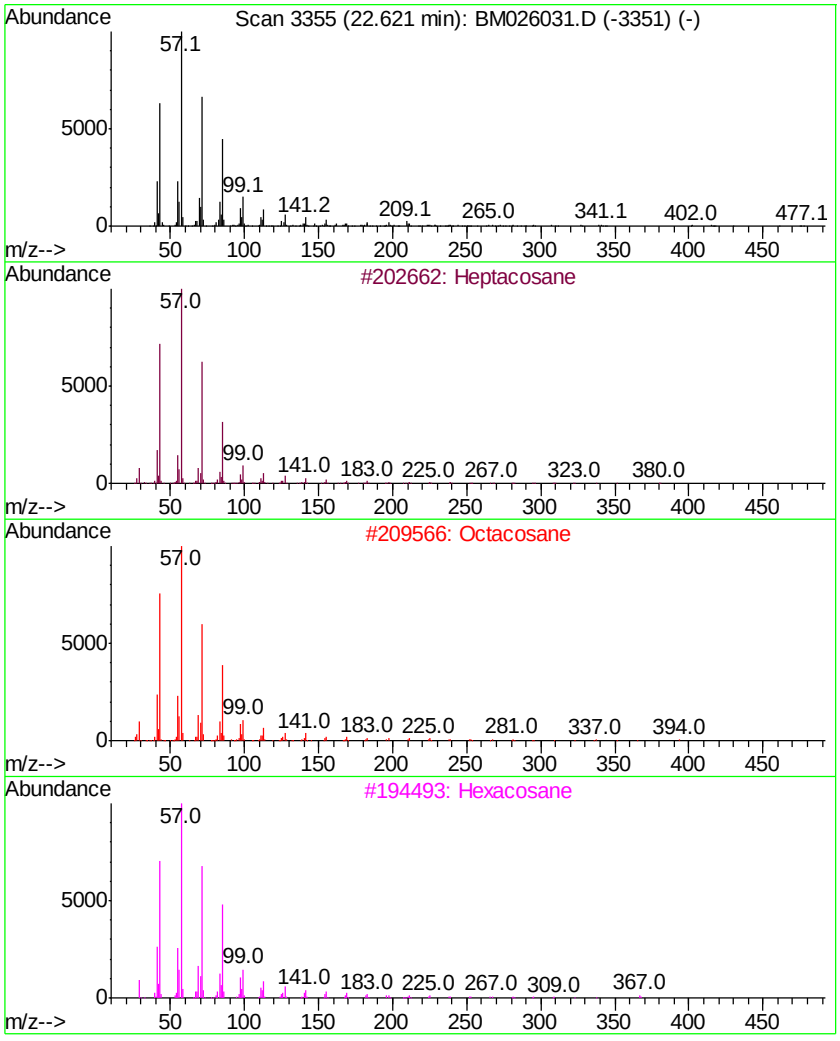
Quant Title : SVOA CALIBRATION

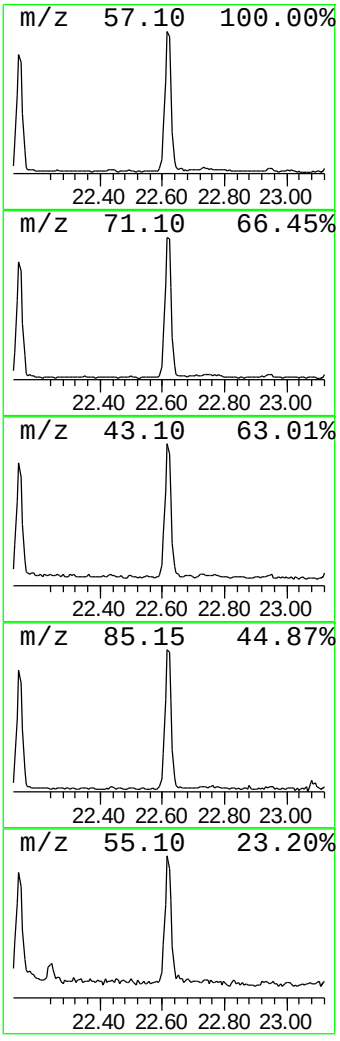
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.62	3.54 ng/ul	503618	Perylene-d12	23.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	91
2	Octacosane	394	C28H58	000630-02-4	91
3	Hexacosane	366	C26H54	000630-01-3	91
4	Tetracosane	338	C24H50	000646-31-1	90
5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	90





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051920\
Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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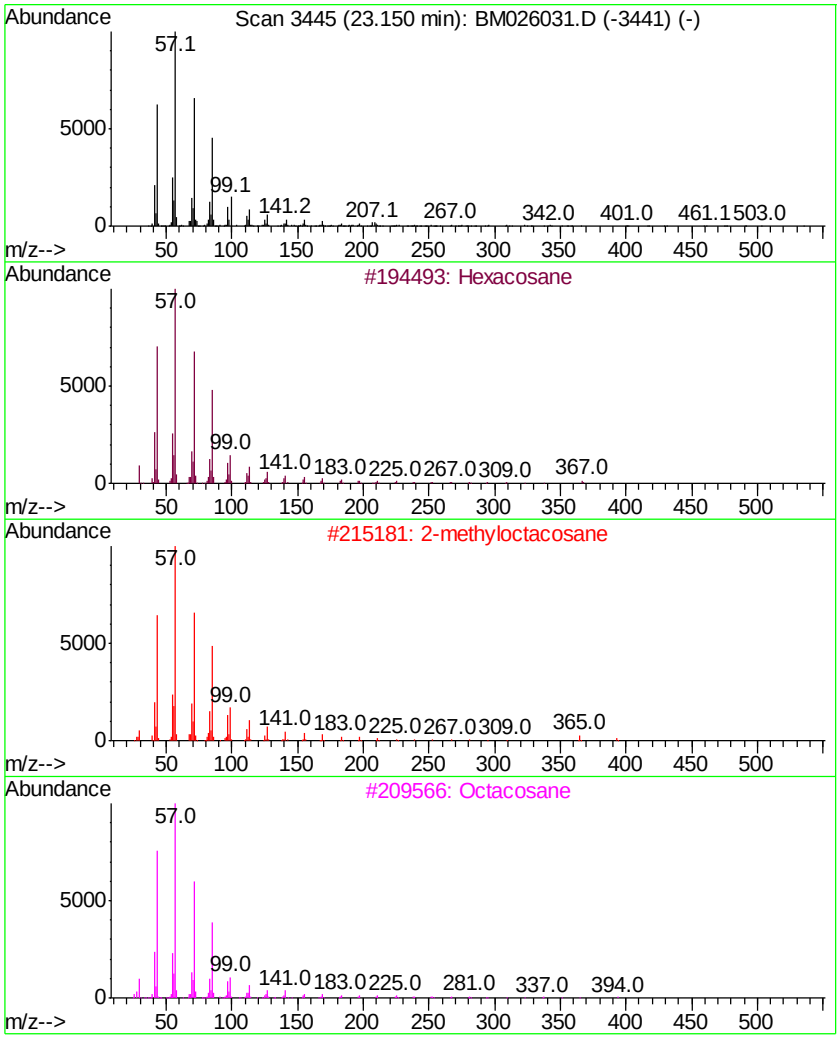
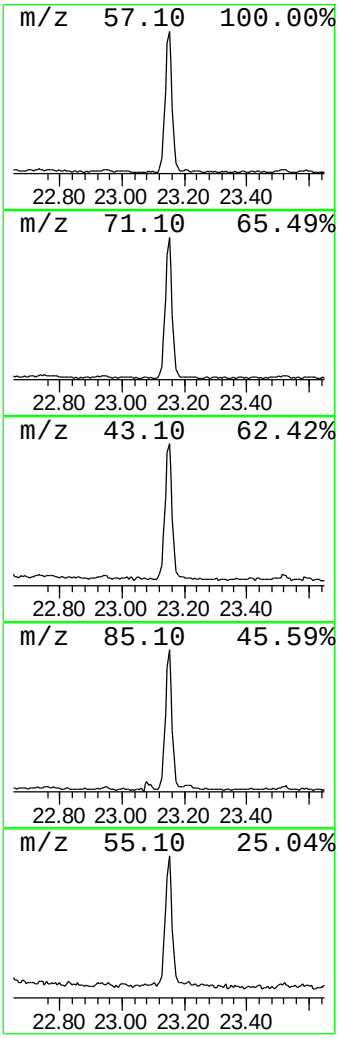
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.15	4.02 ng/ul	570875	Perylene-d12	23.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	90
2	2-methyloctacosane	408	C29H60	1000376-72-8	90
3	Octacosane	394	C28H58	000630-02-4	90
4	Heneicosane	296	C21H44	000629-94-7	87
5	Heptacosane	380	C27H56	000593-49-7	87





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051920\
Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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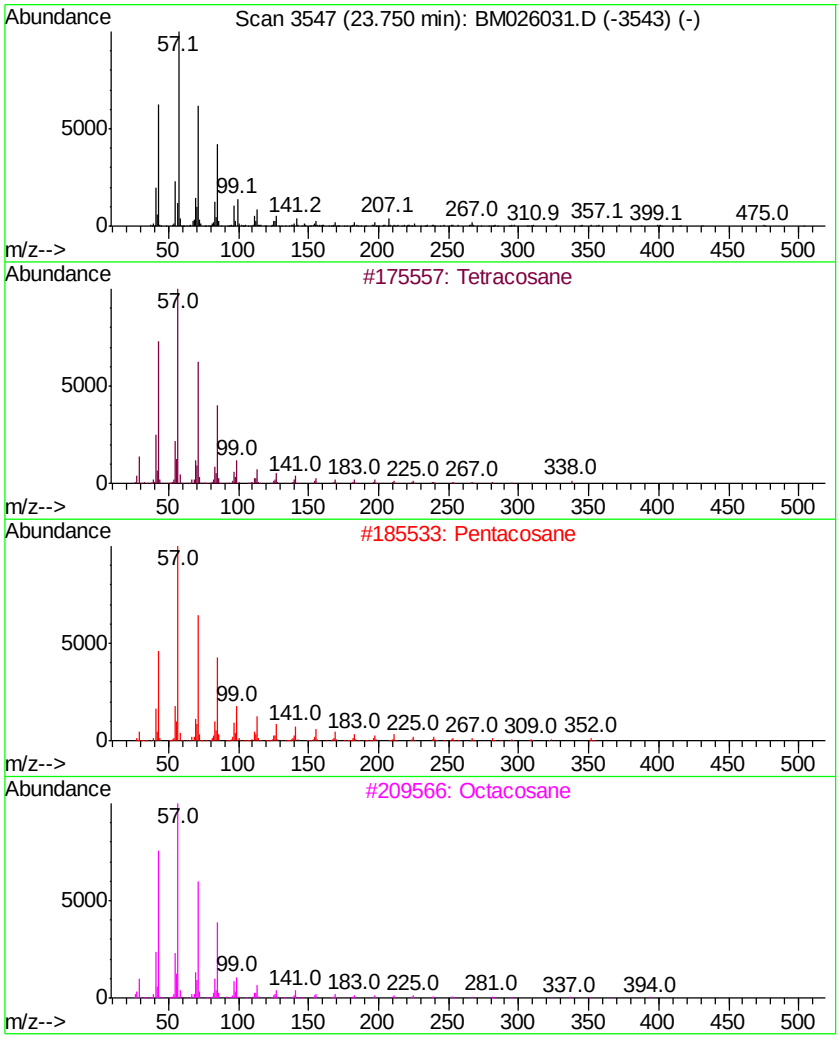
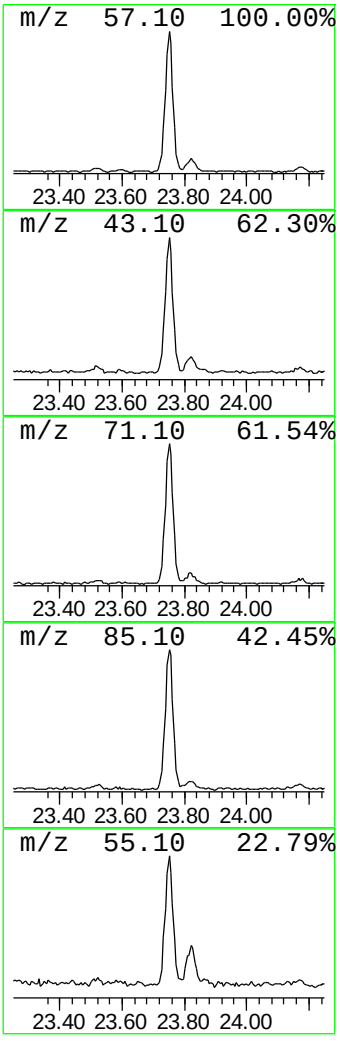
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.75	3.48 ng/ul	495024	Perylene-d12	23.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	90
2	Pentacosane	352	C25H52	000629-99-2	90
3	Octacosane	394	C28H58	000630-02-4	90
4	Heneicosane	296	C21H44	000629-94-7	90
5	Tetratetracontane	619	C44H90	007098-22-8	90





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051920\
Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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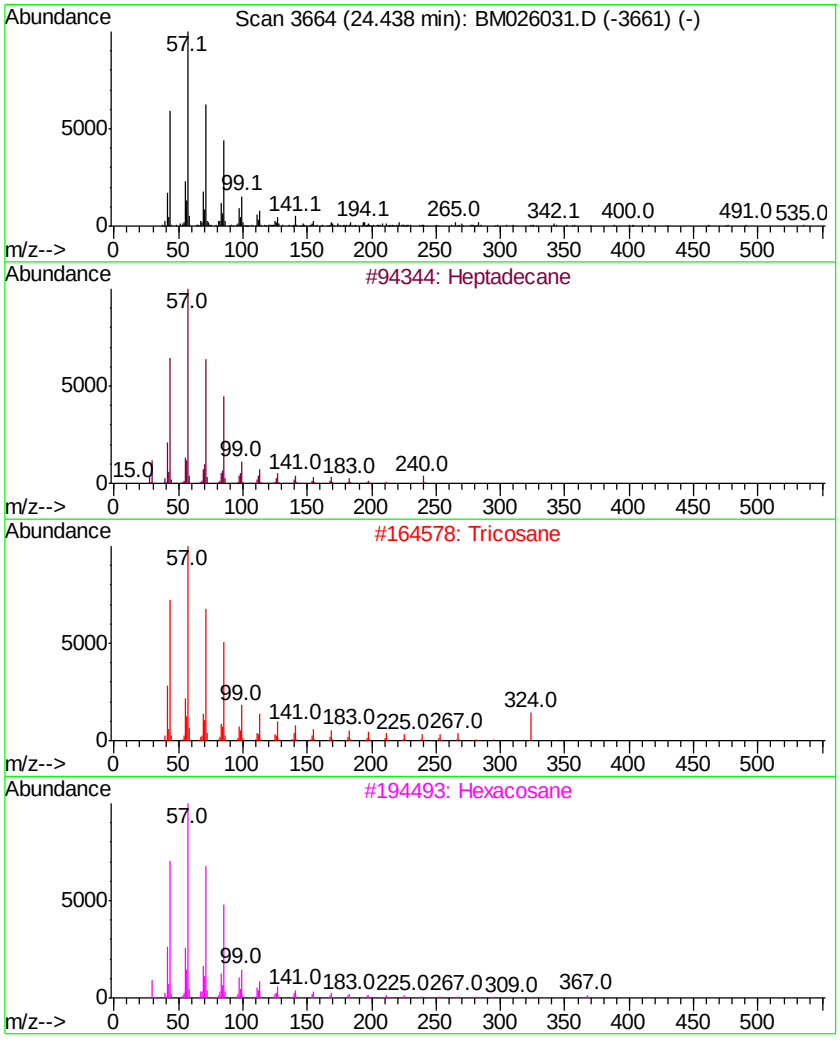
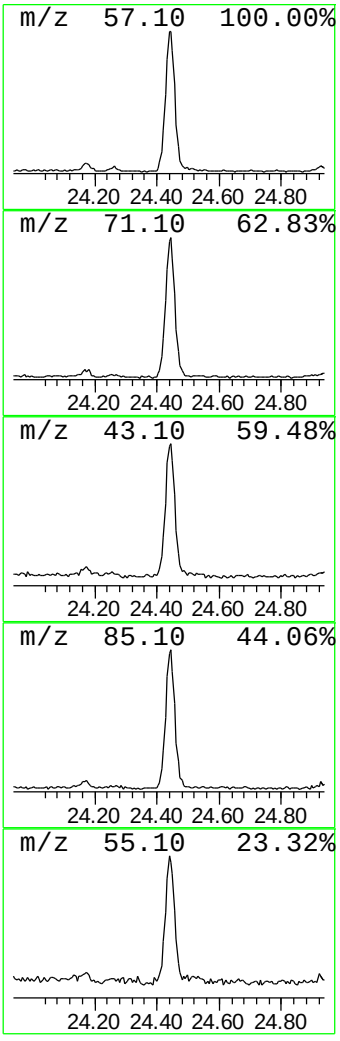
Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.44	2.70 ng/ul	383222	Perylene-d12	23.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	93
2	Tricosane	324	C23H48	000638-67-5	89
3	Hexacosane	366	C26H54	000630-01-3	87
4	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	83
5	Tetratetracontane	619	C44H90	007098-22-8	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM051920\
Data File : BM026031.D
Acq On : 19 May 2020 23:52
Operator : MA/SJ
Sample : L2681-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C5CA9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.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
Benzene, (1,1-dim...	8.84	4.0	ng/ul	249619	1	7.50	1255850	20.0
Benzene, (2-methy...	9.48	2.1	ng/ul	194968	2	10.26	1878680	20.0
Benzene, 2-etheny...	9.67	3.0	ng/ul	281286	2	10.26	1878680	20.0
Benzene, (3-methy...	10.33	2.4	ng/ul	223767	2	10.26	1878680	20.0
Naphthalene, 1,2,...	10.84	2.5	ng/ul	233496	2	10.26	1878680	20.0
1H-Inden-1-ol, 2,...	10.87	3.1	ng/ul	291924	2	10.26	1878680	20.0
1,3-Cyclopentadie...	11.37	2.5	ng/ul	231885	2	10.26	1878680	20.0
1H-Inden-1-ol, 2,...	11.49	3.0	ng/ul	286733	2	10.26	1878680	20.0
1H-Inden-1-one, 2...	11.65	7.4	ng/ul	694676	2	10.26	1878680	20.0
1H-Indene, 2,3-di...	11.90	2.1	ng/ul	194596	2	10.26	1878680	20.0
Ethanone, 1-(3,4-...	12.10	2.6	ng/ul	249293	2	10.26	1878680	20.0
1-Methylindan-2-one	12.16	2.1	ng/ul	195320	2	10.26	1878680	20.0
2,4,6-Trimethylbe...	12.27	3.2	ng/ul	381421	3	14.15	2351410	20.0
Benzene, 1-ethyl-...	12.34	3.5	ng/ul	413140	3	14.15	2351410	20.0
Bicyclo[4.2.0]oct...	12.67	2.5	ng/ul	289734	3	14.15	2351410	20.0
(Z)-1-Phenylpropene	13.10	2.3	ng/ul	270650	3	14.15	2351410	20.0
7-Methylindan-1-one	13.27	2.6	ng/ul	305869	3	14.15	2351410	20.0
Phenol, 2,3,4,6-t...	14.70	4.4	ng/ul	517292	3	14.15	2351410	20.0
unknown-01	15.69	2.1	ng/ul	265376	4	16.90	2518010	20.0
1-Acenaphthenol	15.87	5.0	ng/ul	623927	4	16.90	2518010	20.0
(DEL) Alkane: Cyc...	20.84	2.6	ng/ul	367707	5	21.09	2786300	20.0
(DEL) Alkane: Str...	22.14	2.3	ng/ul	326152	5	21.09	2786300	20.0
(DEL) Alkane: Str...	22.62	3.5	ng/ul	503618	6	23.22	2842010	20.0
(DEL) Alkane: Str...	23.15	4.0	ng/ul	570875	6	23.22	2842010	20.0
(DEL) Alkane: Str...	23.75	3.5	ng/ul	495024	6	23.22	2842010	20.0
(DEL) Alkane: Str...	24.44	2.7	ng/ul	383222	6	23.22	2842010	20.0