LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040820\
Data File : BN010434.D
Acq On : 08 Apr 2020 15:44
Operator : CG/JU
Sample : L2155-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBFA1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.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	2.834	3	8	14	rBV4	179973	516763	1.20%	0.096%
2	2.893	14	18	26	rVV	394355	1018020	2.36%	0.188%
3	2.964	26	30	56	rVB	1482394	2673033	6.19%	0.494%
4	5.181	402	407	414	rVB	279268	462019	1.07%	0.085%
5	5.311	423	429	438	rBV	2438180	4359824	10.10%	0.806%
6	5.669	484	490	500	rVB	2171407	3825269	8.86%	0.707%
7	5.934	527	535	545	rBV5	227745	747963	1.73%	0.138%
8	6.134	563	569	578	rVB	864907	1546716	3.58%	0.286%
9	6.287	587	595	598	rBV4	283795	571450	1.32%	0.106%
10	6.322	598	601	607	rVB	376651	561898	1.30%	0.104%
11	6.610	644	650	653	rBV3	237225	437053	1.01%	0.081%
12	6.722	660	669	674	rBV2	511818	1007861	2.33%	0.186%
13	6.787	674	680	687	rVV	2043161	3809502	8.82%	0.704%
14	6.852	687	691	697	rVB	1293318	1943328	4.50%	0.359%
15	6.952	698	708	714	rBV	1480119	2491310	5.77%	0.461%
16	7.016	714	719	723	rVB	423723	600952	1.39%	0.111%
17	7.146	734	741	747	rVB	1886900	3028897	7.02%	0.560%
18	7.269	756	762	768	rBV2	1930827	3090594	7.16%	0.572%
19	7.522	800	805	813	rVV	457055	927200	2.15%	0.171%
20	7.604	813	819	825	rVV2	1994966	3352829	7.77%	0.620%
21	7.722	835	839	848	rVV2	625105	1236334	2.86%	0.229%
22	7.904	863	870	878	rVV7	246549	666540	1.54%	0.123%
23	8.157	908	913	920	rVB	580345	1012102	2.34%	0.187%
24	8.246	924	928	932	rVV	871176	1304109	3.02%	0.241%
25	8.304	932	938	945	rVB	2417577	3842008	8.90%	0.710%
26	8.410	948	956	964	rBV3	350454	1005631	2.33%	0.186%
27	8.557	976	981	985	rBV	363718	493118	1.14%	0.091%
28	8.604	985	989	995	rVB	360806	578544	1.34%	0.107%
29	8.704	998	1006	1008	rBV	639515	1092115	2.53%	0.202%
30	8.740	1008	1012	1017	rVB	1637879	2506299	5.81%	0.463%
31	8.846	1026	1030	1039	rVB	761882	1203044	2.79%	0.222%
32	8.957	1039	1049	1055	rBV6	524271	1314078	3.04%	0.243%
33	9.034	1055	1062	1067	rBV3	260144	575730	1.33%	0.106%
34	9.222	1090	1094	1098	rVV2	274593	455142	1.05%	0.084%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040820\
Data File : BN010434.D
Acq On : 08 Apr 2020 15:44
Operator : CG/JU
Sample : L2155-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBFA1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Title : SVOA CALIBRATION

35	9.281	1101	1104	1114	rVB2	380017	752331	1.74%	0.139%
36	9.457	1128	1134	1142	rBV2	1500791	2618820	6.07%	0.484%
37	9.598	1150	1158	1167	rVB5	298158	814649	1.89%	0.151%
38	9.787	1186	1190	1194	rVB4	365018	603110	1.40%	0.112%
39	9.857	1199	1202	1208	rVB2	341303	506525	1.17%	0.094%
40	9.993	1219	1225	1235	rVB	2674662	4718208	10.93%	0.873%
41	10.363	1281	1288	1294	rBV	2766744	4640543	10.75%	0.858%
42	10.416	1294	1297	1302	rVV2	320412	599327	1.39%	0.111%
43	10.498	1305	1311	1318	rVB	1577307	2662027	6.17%	0.492%
44	11.193	1420	1429	1439	rBV	333287	719816	1.67%	0.133%
45	13.645	1840	1846	1851	rVV	4795225	6448521	14.94%	1.192%
46	13.692	1851	1854	1860	rVB	1256742	1612055	3.73%	0.298%
47	13.922	1886	1893	1902	rBV	5306854	8031153	18.60%	1.485%
48	14.010	1902	1908	1914	rVB	2812740	3364549	7.79%	0.622%
49	14.234	1939	1946	1952	rVV	3998587	6018515	13.94%	1.113%
50	14.428	1973	1979	1993	rBV	1396969	2156830	5.00%	0.399%
51	15.228	2108	2115	2121	rBV	7224462	10283726	23.82%	1.902%
52	15.345	2130	2135	2143	rBV	1273405	1764253	4.09%	0.326%
53	16.975	2406	2412	2416	rBV	5255460	7586649	17.57%	1.403%
54	17.016	2416	2419	2424	rVV	2729850	3779084	8.75%	0.699%
55	17.075	2424	2429	2443	rVB	8051300	11773614	27.27%	2.177%
56	17.380	2477	2481	2486	rVB	292487	433822	1.00%	0.080%
57	17.851	2557	2561	2565	rBV	305703	408845	0.95%	0.076%
58	17.904	2565	2570	2578	rBV2	1048246	1653033	3.83%	0.306%
59	18.492	2665	2670	2674	rVB	474344	565903	1.31%	0.105%
60	18.563	2674	2682	2688	rBV	13877004	23301695	53.97%	4.309%
61	18.704	2702	2706	2713	rVB	633266	832518	1.93%	0.154%
62	19.045	2758	2764	2771	rBV	7864607	10653797	24.68%	1.970%
63	19.380	2815	2821	2824	rBV	10793810	17090283	39.58%	3.160%
64	19.410	2824	2826	2834	rVV	7436089	7613687	17.64%	1.408%
65	19.480	2835	2838	2847	rVB2	345651	537262	1.24%	0.099%
66	19.592	2853	2857	2861	rBV	409541	529782	1.23%	0.098%
67	19.739	2876	2882	2887	rBV5	574041	1279206	2.96%	0.237%
68	19.839	2895	2899	2902	rBV3	331893	457499	1.06%	0.085%
69	19.904	2902	2910	2917	rVB	1456552	3081698	7.14%	0.570%
70	20.010	2924	2928	2932	rVB	956091	1116638	2.59%	0.206%
71	20.069	2933	2938	2942	rVV2	1041850	1564778	3.62%	0.289%

LSC Area Percent Report

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Acq On : 08 Apr 2020 15:44
Operator : CG/JU
Sample : L2155-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBFA1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Title : SVOA CALIBRATION

72	20.104	2942	2944	2948	rVB2	398693	461548	1.07%	0.085%
73	20.210	2958	2962	2965	rBV	721666	889811	2.06%	0.165%
74	20.251	2966	2969	2973	rVV	653458	906149	2.10%	0.168%
75	20.292	2973	2976	2979	rVB2	402446	428214	0.99%	0.079%
76	20.657	3035	3038	3045	rVB	832147	1319384	3.06%	0.244%
77	20.827	3062	3067	3070	rBV	2371286	3056608	7.08%	0.565%
78	20.869	3070	3074	3077	rVV	1169180	1436790	3.33%	0.266%
79	20.921	3077	3083	3098	rVB5	3197829	6324227	14.65%	1.169%
80	21.033	3099	3102	3104	rBV	503418	548937	1.27%	0.102%
81	21.063	3105	3107	3116	rVB	746395	977547	2.26%	0.181%
82	21.174	3120	3126	3131	rVV3	13415592	29129972	67.47%	5.387%
83	21.221	3131	3134	3144	rVB	11833639	15129276	35.04%	2.798%
84	21.345	3150	3155	3158	rBV2	1434837	2053622	4.76%	0.380%
85	21.492	3176	3180	3185	rVB2	2070710	3075126	7.12%	0.569%
86	21.780	3226	3229	3235	rVB2	1409717	2021484	4.68%	0.374%
87	21.892	3243	3248	3255	rBV	20060993	25902714	60.00%	4.790%
88	22.221	3298	3304	3308	rBV	28878248	37881195	87.74%	7.005%
89	22.727	3382	3390	3394	rBV	18482599	43173673	100.00%	7.984%
90	22.880	3412	3416	3422	rVB	4229024	6112110	14.16%	1.130%
91	23.010	3434	3438	3446	rBV	1225197	2462318	5.70%	0.455%
92	23.110	3451	3455	3460	rVV2	3205881	5538632	12.83%	1.024%
93	23.180	3461	3467	3472	rVV	12424658	24738929	57.30%	4.575%
94	23.227	3472	3475	3479	rVV	7552838	13692968	31.72%	2.532%
95	23.274	3479	3483	3490	rVB	13407572	23128530	53.57%	4.277%
96	23.362	3492	3498	3502	rVV	5055059	10079305	23.35%	1.864%
97	23.410	3503	3506	3514	rVB2	3724600	7038514	16.30%	1.302%
98	25.145	3794	3801	3809	rBV	8667967	19617404	45.44%	3.628%
99	25.521	3856	3865	3874	rVB	10244770	27944087	64.72%	5.168%
100	26.174	3968	3976	3992	rVB2	6000176	18862635	43.69%	3.488%

Sum of corrected areas: 540763735

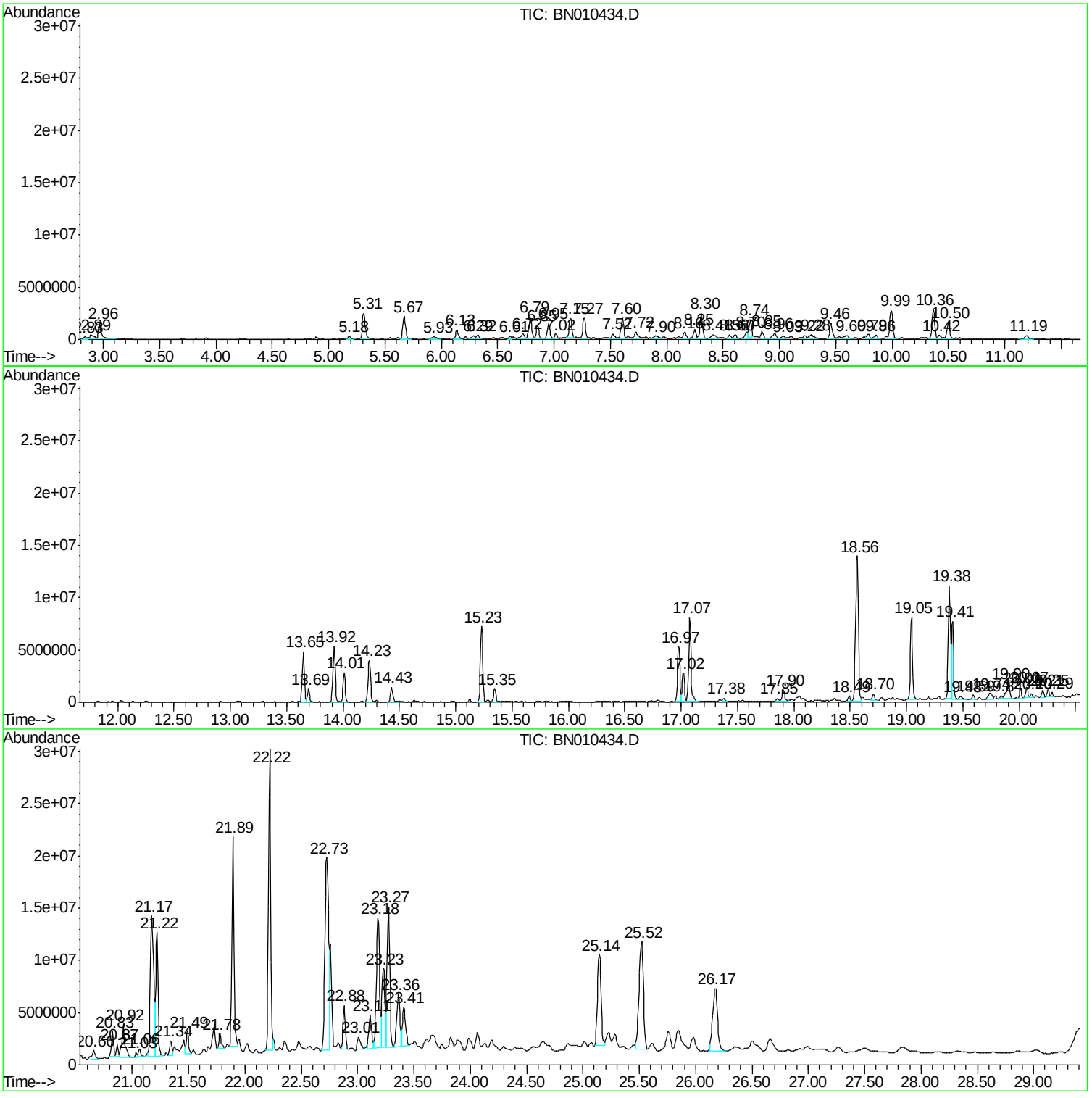
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040820\
Data File : BN010434.D
Acq On : 08 Apr 2020 15:44
Operator : CG/JU
Sample : L2155-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBFA1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040820\
Data File : BN010434.D
Acq On : 08 Apr 2020 15:44
Operator : CG/JU
Sample : L2155-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBFA1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

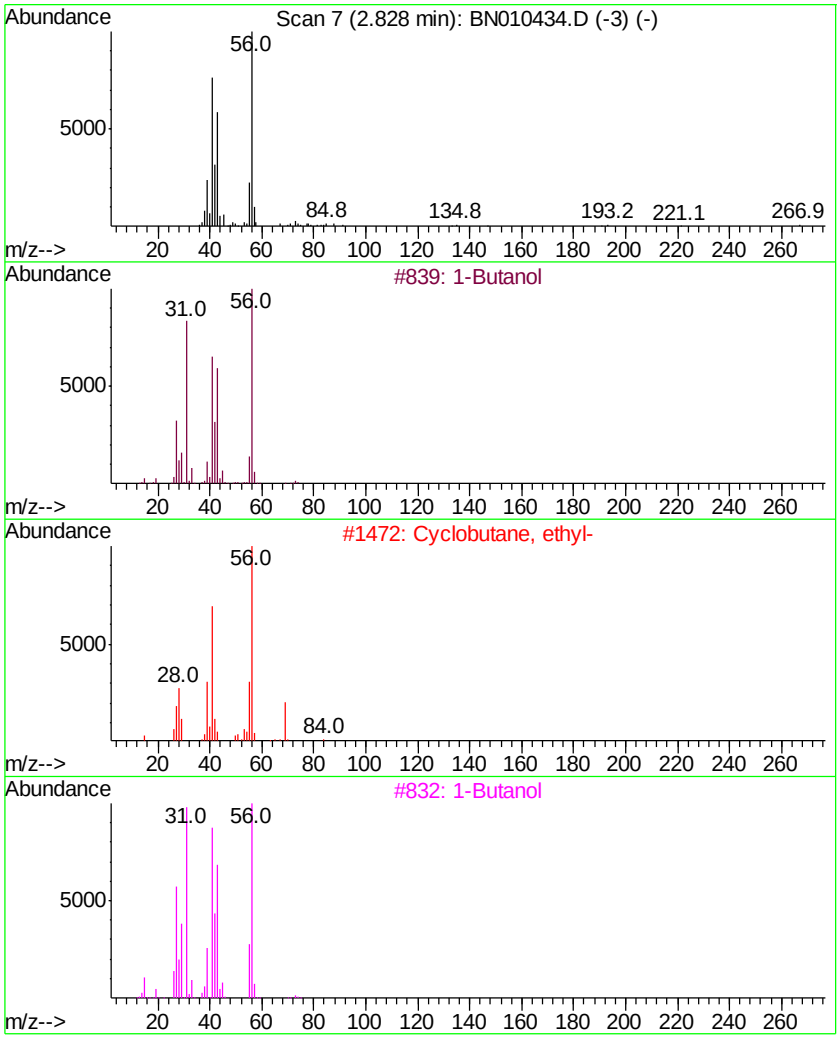
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

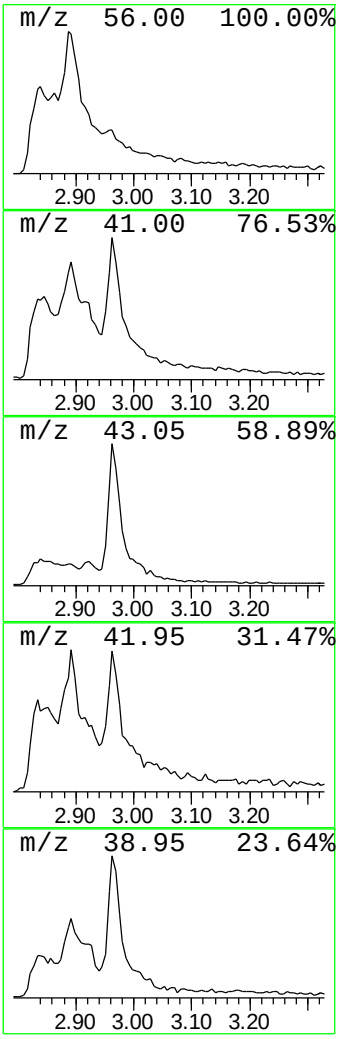
****

Peak Number 1 1-Butanol Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.83	3.08 ng/ul	516763	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol	74	C4H10O	000071-36-3	80
2	Cyclobutane, ethyl-	84	C6H12	004806-61-5	59
3	1-Butanol	74	C4H10O	000071-36-3	46
4	Succinic anhydride	100	C4H4O3	000108-30-5	45
5	1-Butanol	74	C4H10O	000071-36-3	43





Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040820\
Data File : BN010434.D
Acq On : 08 Apr 2020 15:44
Operator : CG/JU
Sample : L2155-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBFA1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

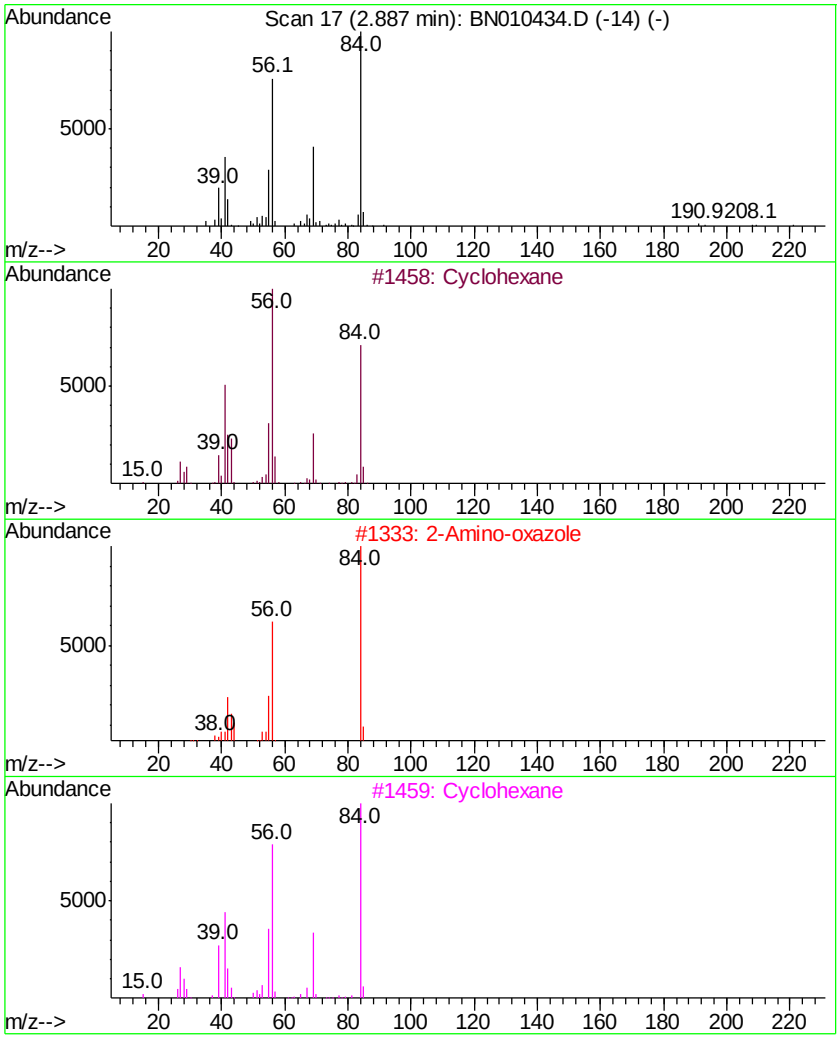
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

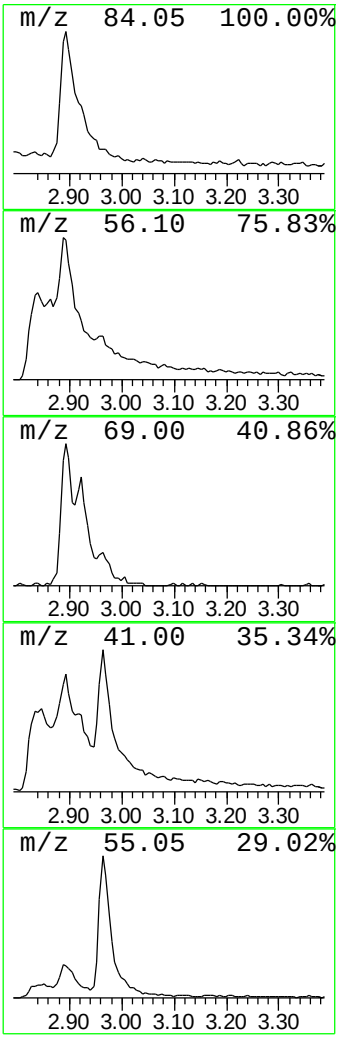
****

Peak Number 2 (DEL) Alkane: Cyclic2.89 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.89	6.07 ng/ul	1018020	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane	84	C6H12	000110-82-7	72
2	2-Amino-oxazole	84	C3H4N2O	004570-45-0	72
3	Cyclohexane	84	C6H12	000110-82-7	72
4	Cyclohexane	84	C6H12	000110-82-7	72
5	Cyclohexane	84	C6H12	000110-82-7	64





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040820\
Data File : BN010434.D
Acq On : 08 Apr 2020 15:44
Operator : CG/JU
Sample : L2155-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBFA1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

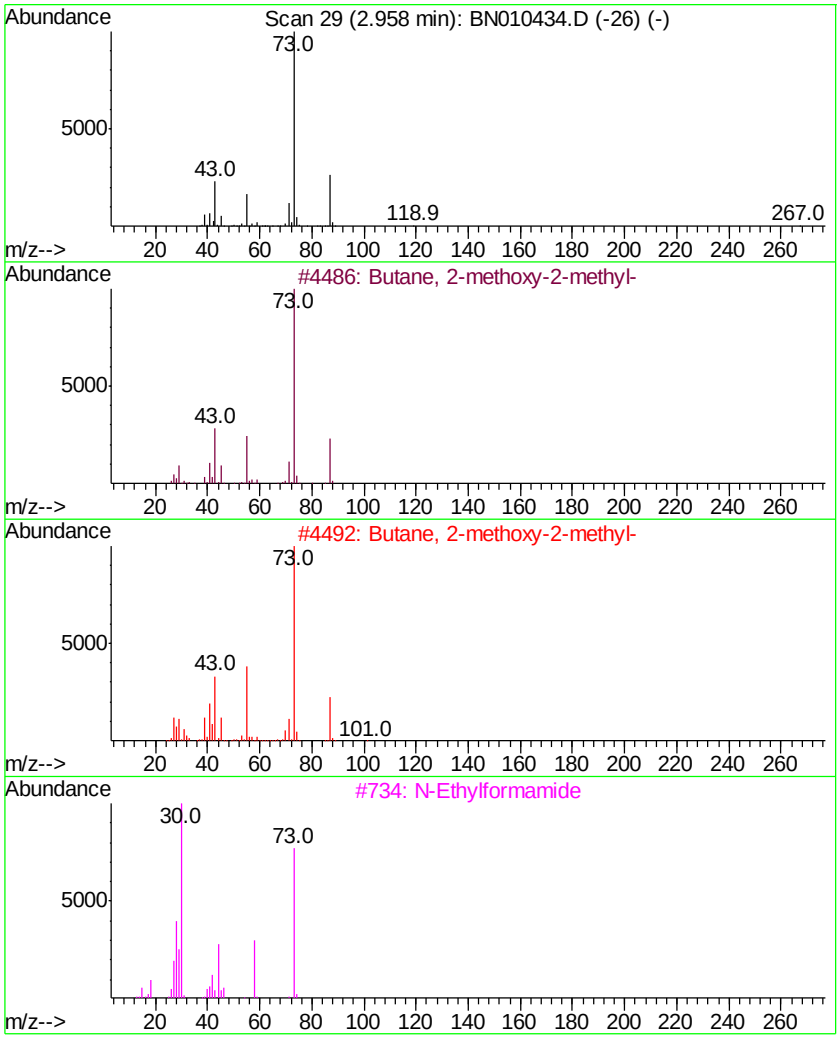
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

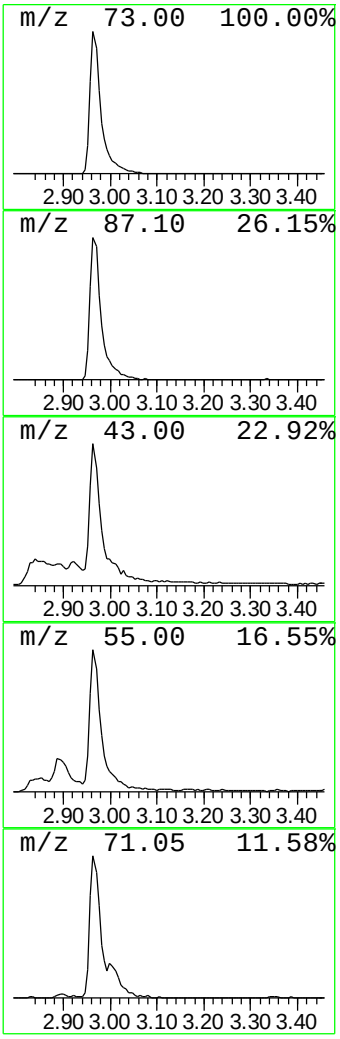
****

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.96	15.94 ng/ul	2673030	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3	N-Ethylformamide	73	C3H7NO	000627-45-2	9
4	Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	1000186-02-3	9
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040820\
Data File : BN010434.D
Acq On : 08 Apr 2020 15:44
Operator : CG/JU
Sample : L2155-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBFA1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

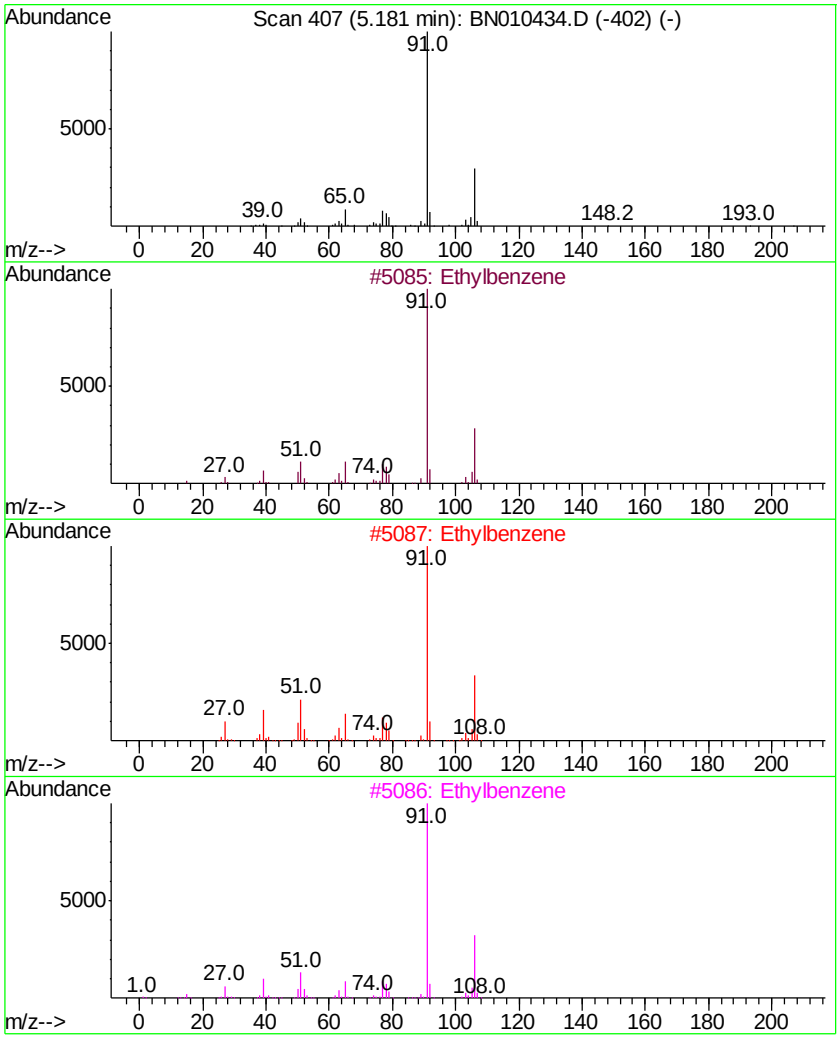
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

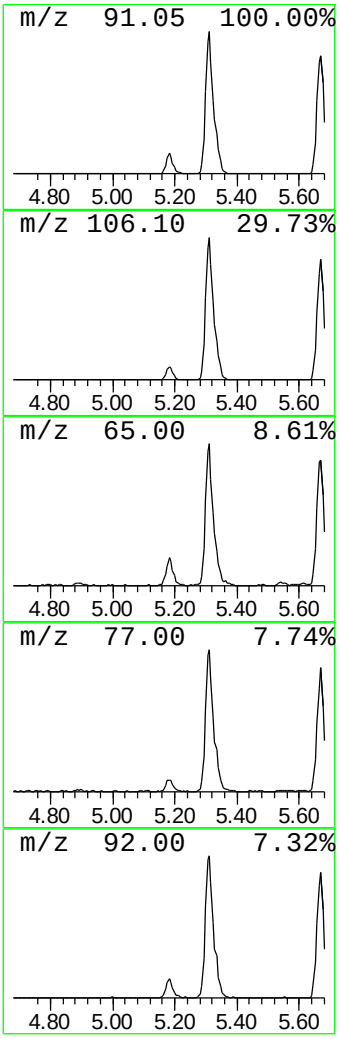
****

Peak Number 4 Ethylbenzene Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	2.76 ng/ul	462019	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethylbenzene	106	C8H10	000100-41-4	93
2	Ethylbenzene	106	C8H10	000100-41-4	90
3	Ethylbenzene	106	C8H10	000100-41-4	90
4	Ethylbenzene	106	C8H10	000100-41-4	90
5	p-Xylene	106	C8H10	000106-42-3	68





Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040820\
Data File : BN010434.D
Acq On : 08 Apr 2020 15:44
Operator : CG/JU
Sample : L2155-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBFA1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

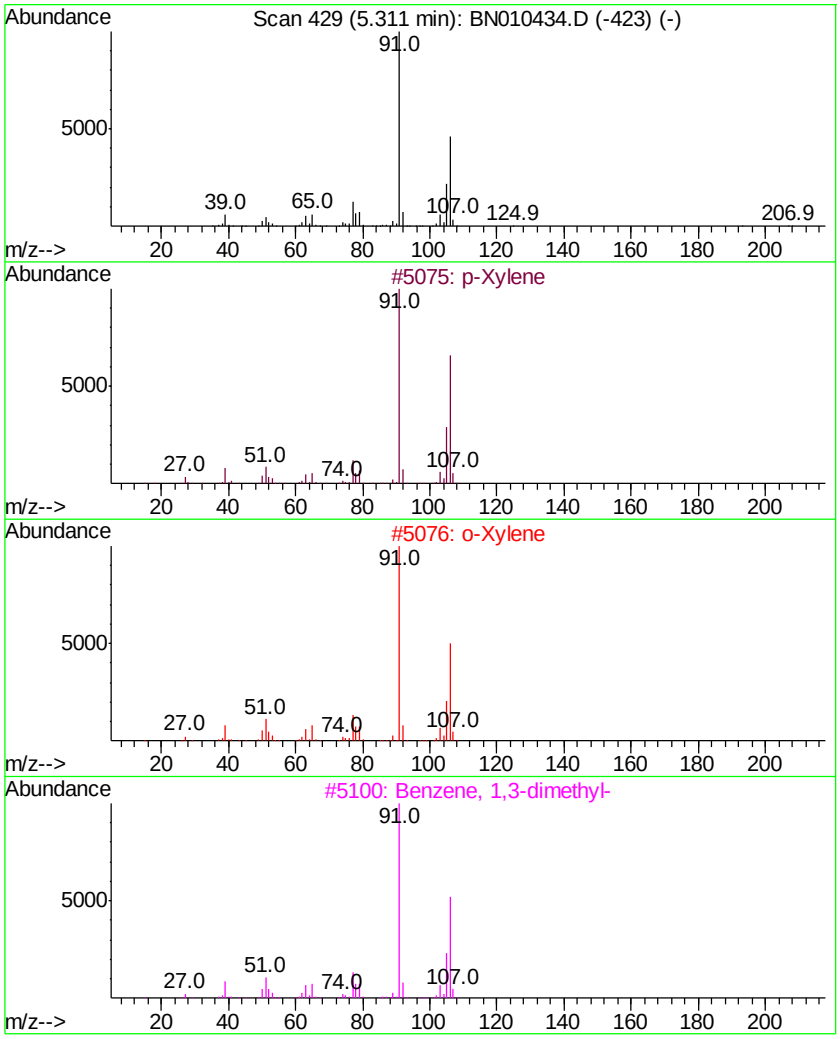
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

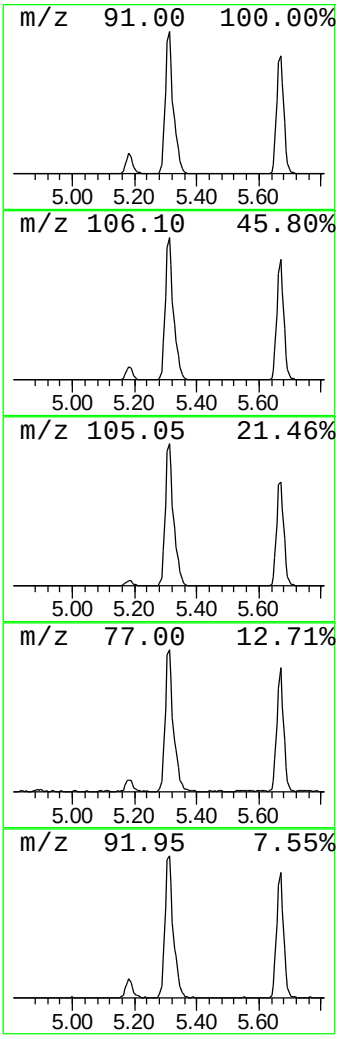
****

Peak Number 5 p-Xylene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.31	26.01 ng/ul	4359820	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Xylene	106	C8H10	000106-42-3	95
2	o-Xylene	106	C8H10	000095-47-6	95
3	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
4	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
5	p-Xylene	106	C8H10	000106-42-3	94





Library Search Compound Report

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Acq On : 08 Apr 2020 15:44
Operator : CG/JU
Sample : L2155-02
Misc :
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Instrument : BNA_N
ClientSampleId : DBFA1

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Quant Title : SVOA CALIBRATION

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

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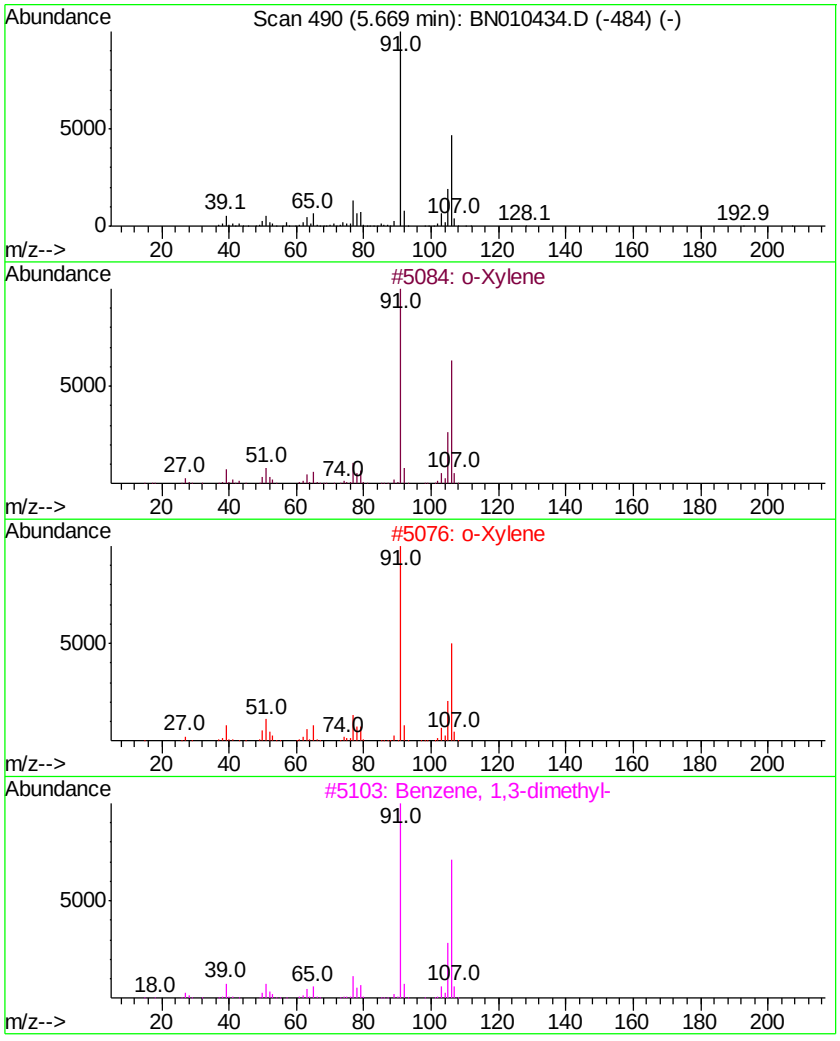
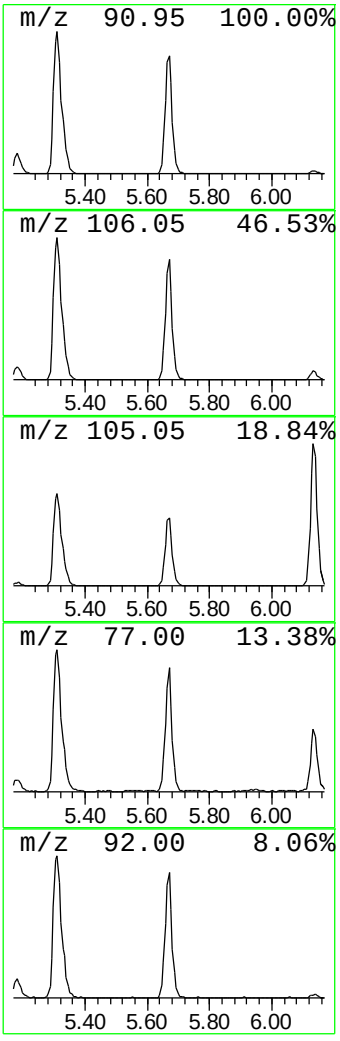
Peak Number 6 o-Xylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.67	22.82 ng/ul	3825270	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Xylene	106	C8H10	000095-47-6	95
2	o-Xylene	106	C8H10	000095-47-6	95
3	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
4	p-Xylene	106	C8H10	000106-42-3	95
5	p-Xylene	106	C8H10	000106-42-3	95





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Quant Title : SVOA CALIBRATION

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

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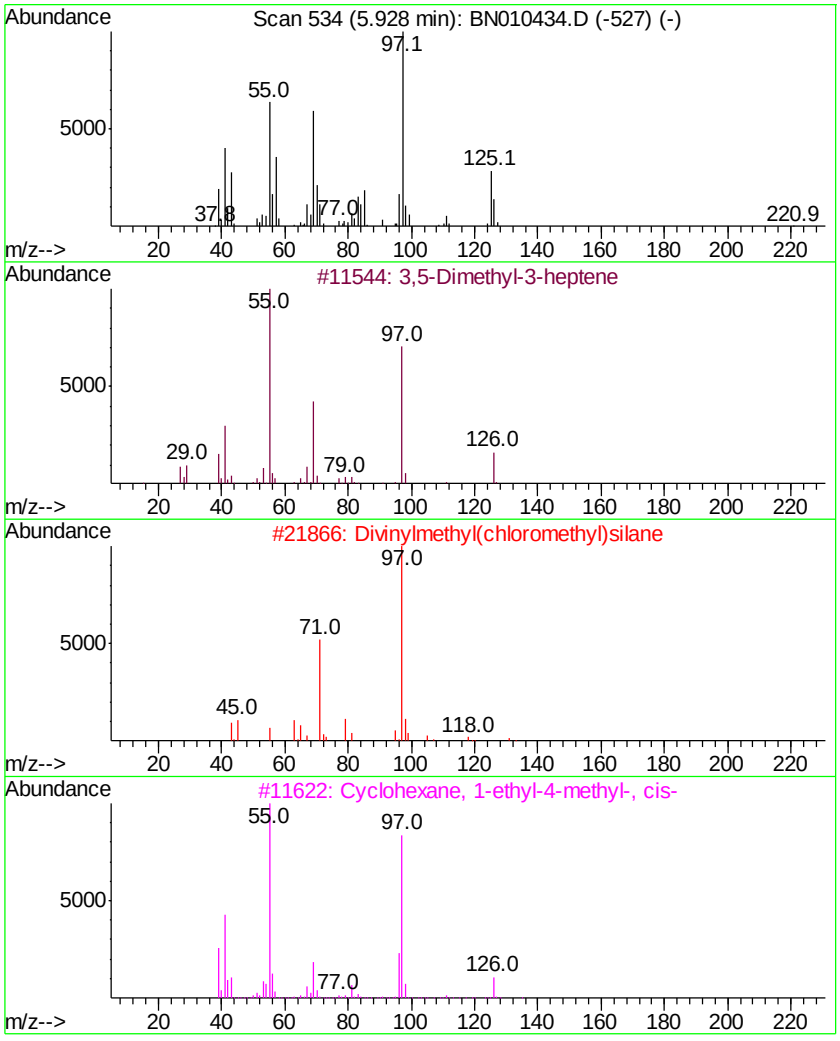
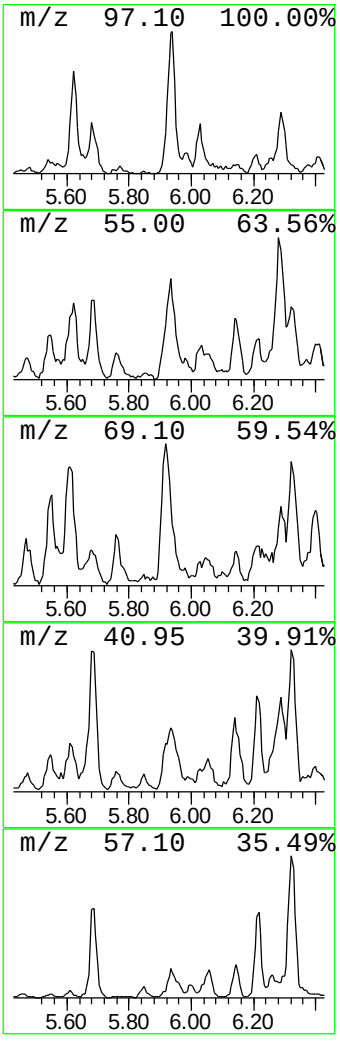
Peak Number 7 3,5-Dimethyl-3-heptene Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.93	4.46 ng/ul	747963	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	53
2	Divinylmethyl(chloromethyl)silane	146	C6H11ClSi	025202-02-2	50
3	Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	49
4	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	49
5	3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	49





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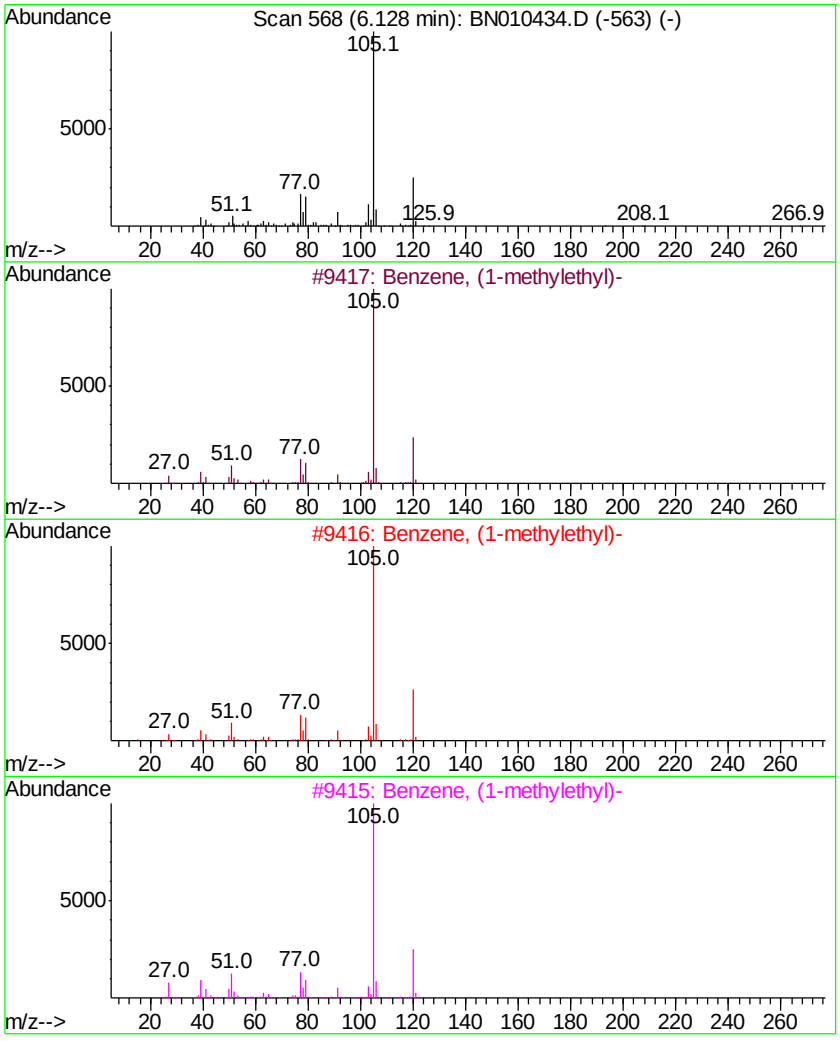
Quant Title : SVOA CALIBRATION

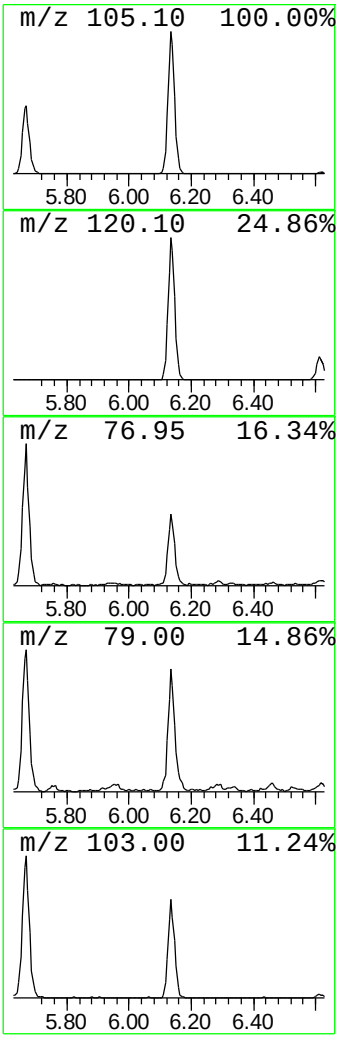
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, (1-methylethyl)- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.13	9.23 ng/ul	1546720	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
2	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
3	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90





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ClientSampleId : DBFA1

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Quant Title : SVOA CALIBRATION

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

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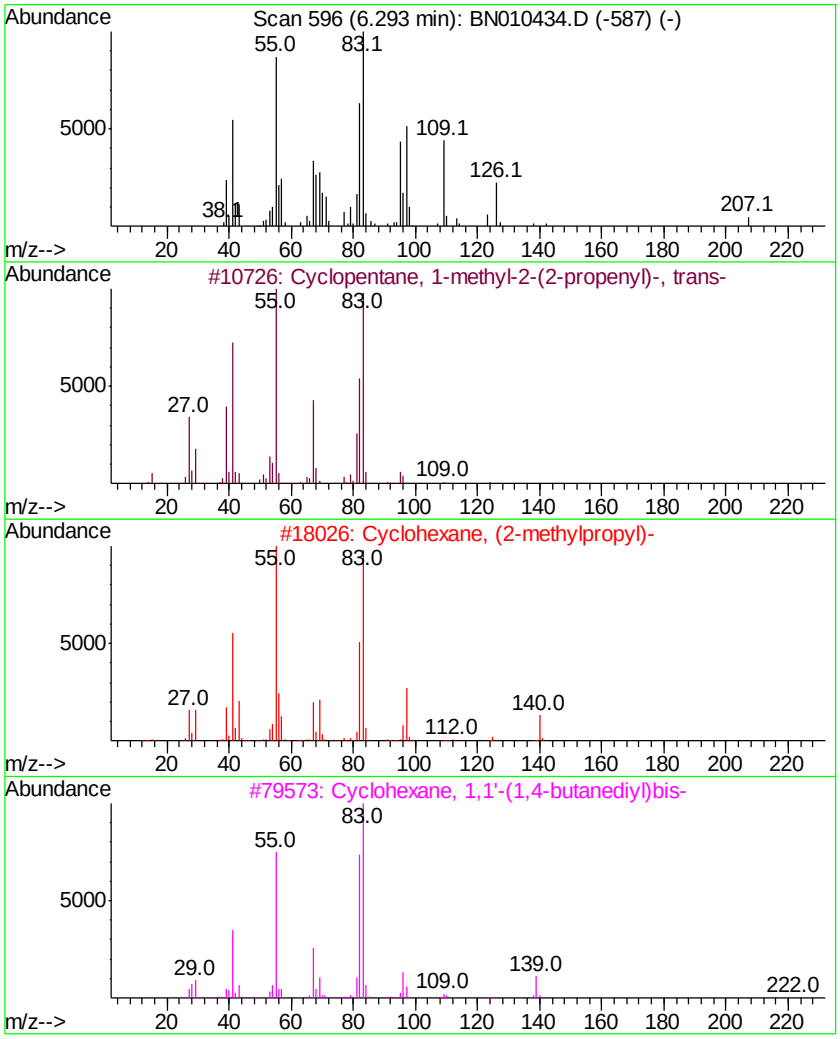
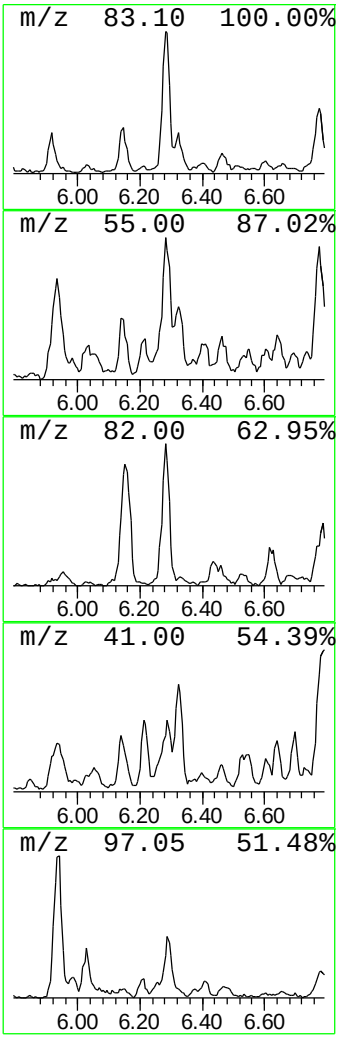
Peak Number 9 Cyclopentane, 1-methyl-2-(2... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.29	3.41 ng/ul	571450	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1-methyl-2-(2-prop...	124	C9H16	050746-53-7	58
2	Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	53
3	Cyclohexane, 1,1'-(1,4-butanediy...	222	C16H30	006165-44-2	53
4	Cyclohexane, propyl-	126	C9H18	001678-92-8	52
5	Cyclohexane, propyl-	126	C9H18	001678-92-8	52





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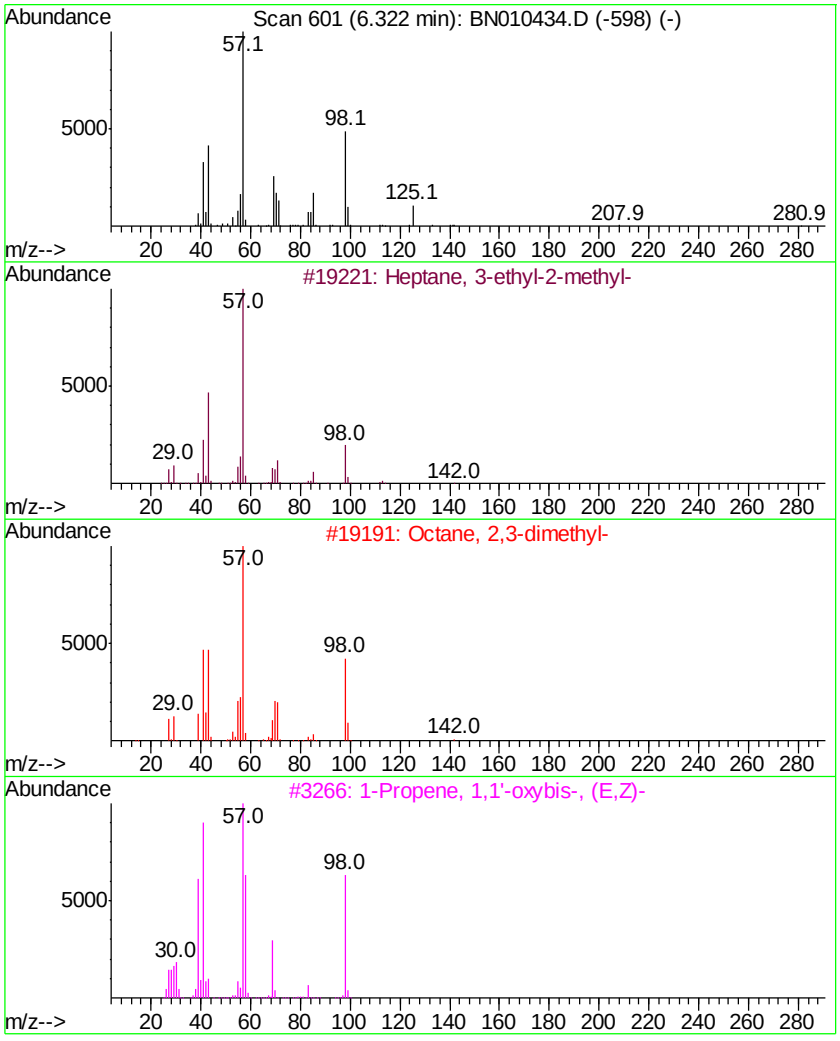
Quant Title : SVOA CALIBRATION

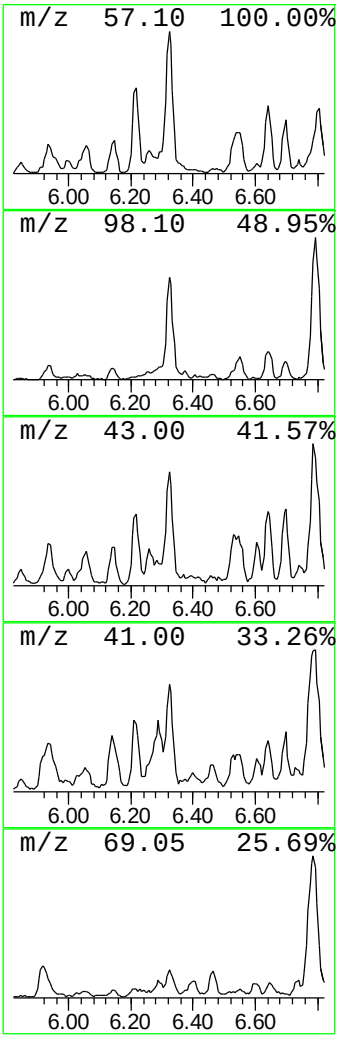
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.32	3.35 ng/ul	561898	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	56
2	Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	52
3	1-Propene, 1,1'-oxybis-, (E,Z)-	98	C6H10O	004696-29-1	46
4	Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	43
5	Hexane, 3-ethyl-2,5-dimethyl-	142	C10H22	052897-04-8	43





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Instrument : BNA_N
ClientSampleId : DBFA1

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Quant Title : SVOA CALIBRATION

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

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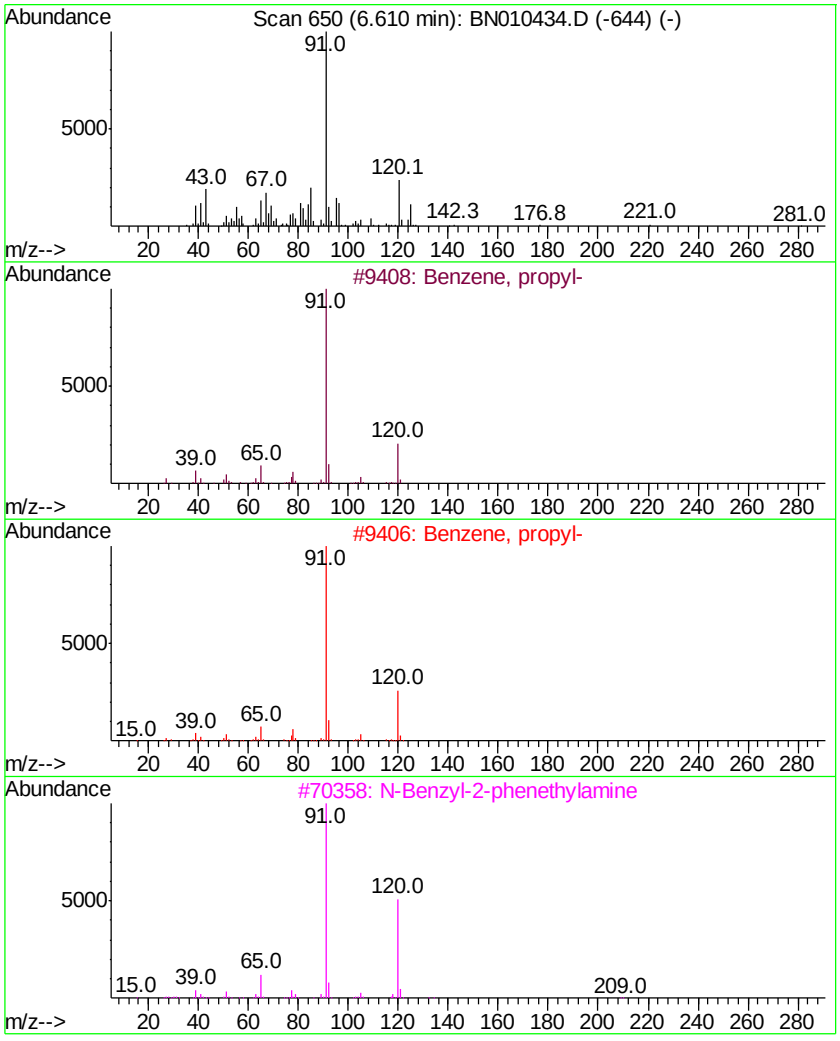
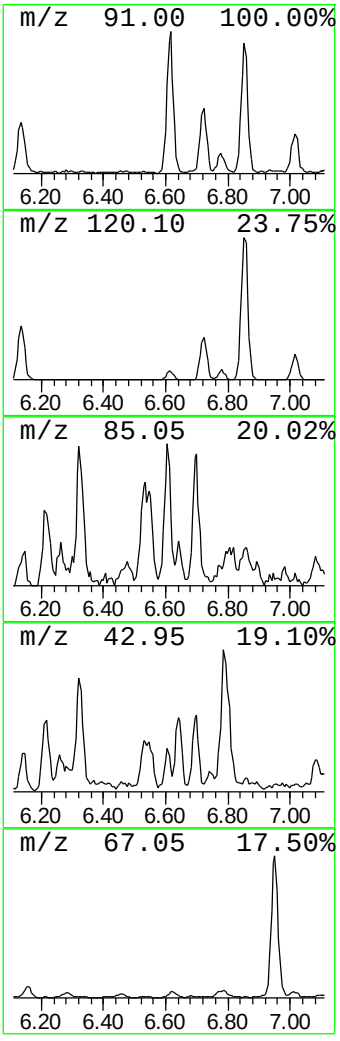
Peak Number 11 unknown-01 Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	2.61 ng/ul	437053	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	42
2	Benzene, propyl-	120	C9H12	000103-65-1	30
3	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	27
4	Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	27
5	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	27





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Instrument : BNA_N
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Quant Title : SVOA CALIBRATION

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

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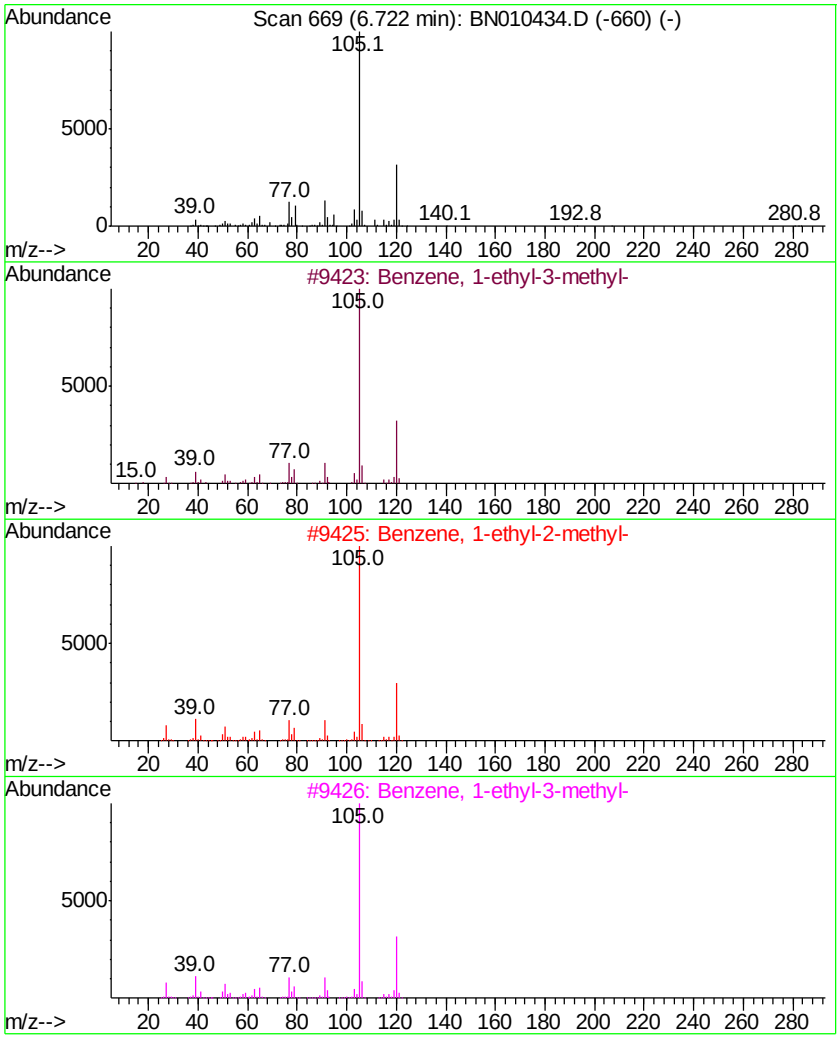
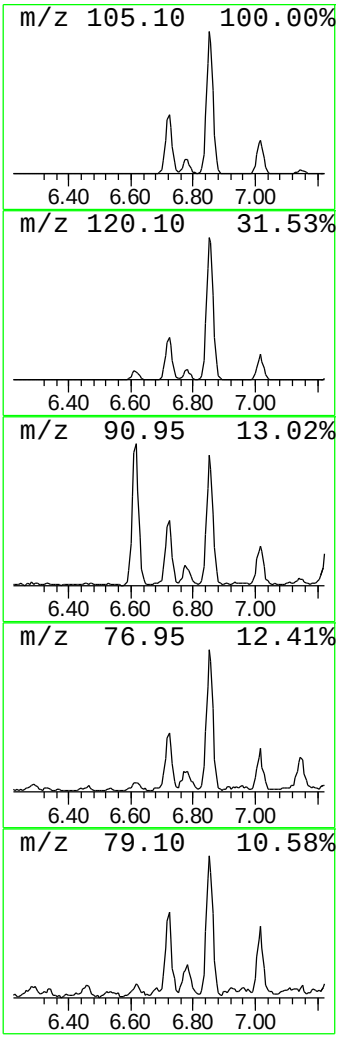
Peak Number 12 Benzene, 1-ethyl-3-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.72	6.01 ng/ul	1007860	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	93
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93





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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 13 Benzene, 1,2,3-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.85	11.59 ng/ul	1943330	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Mesitylene	120	C9H12	000108-67-8	94
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
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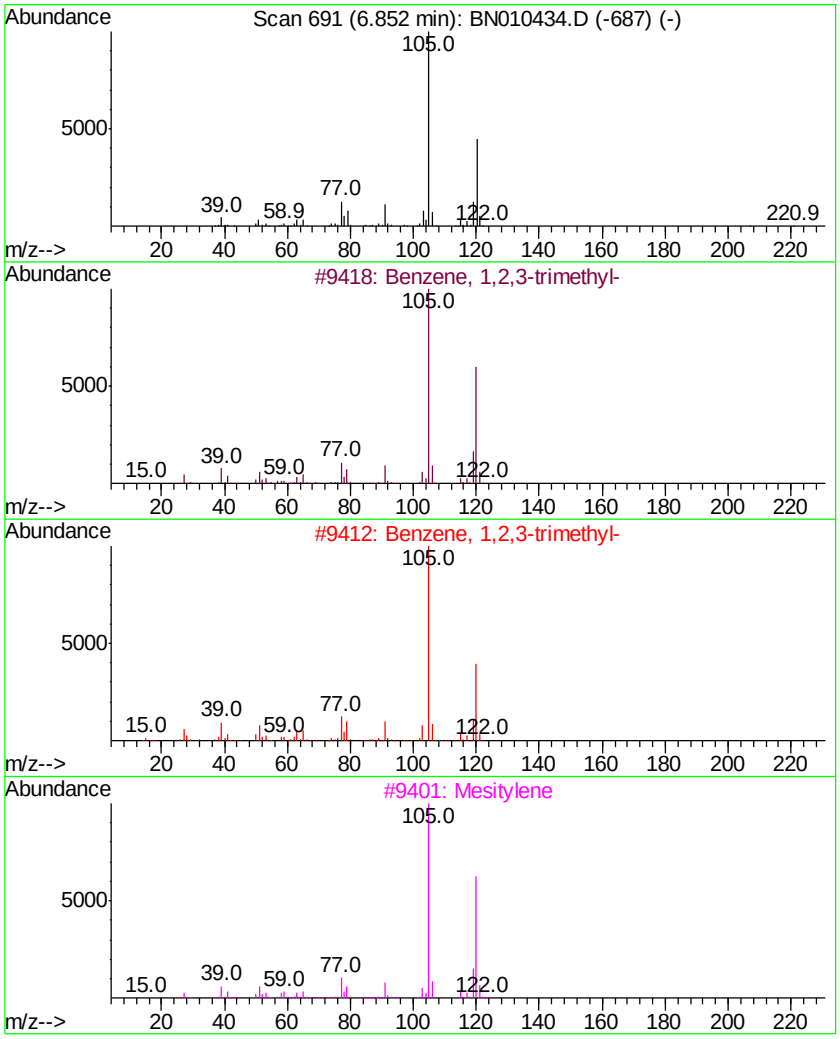

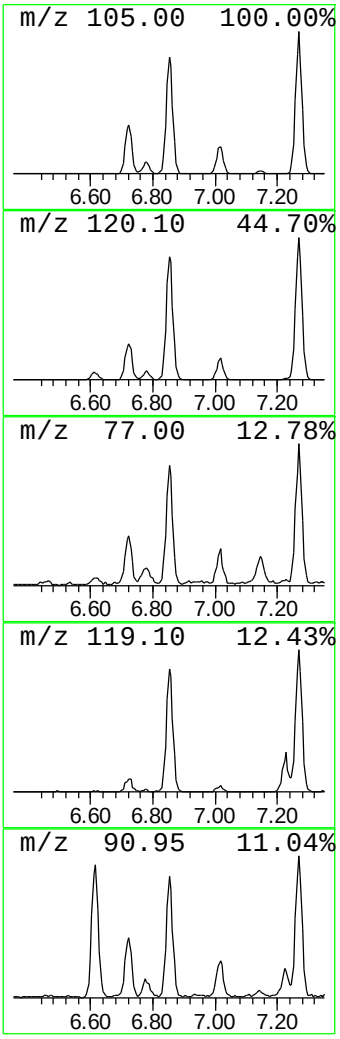

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Quant Title : SVOA CALIBRATION

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

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Peak Number 14 Benzene, 1-ethyl-2-methyl- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.02	3.58 ng/ul	600952	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	76
2	Mesitylene	120	C9H12	000108-67-8	76
3	Mesitylene	120	C9H12	000108-67-8	76
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	76
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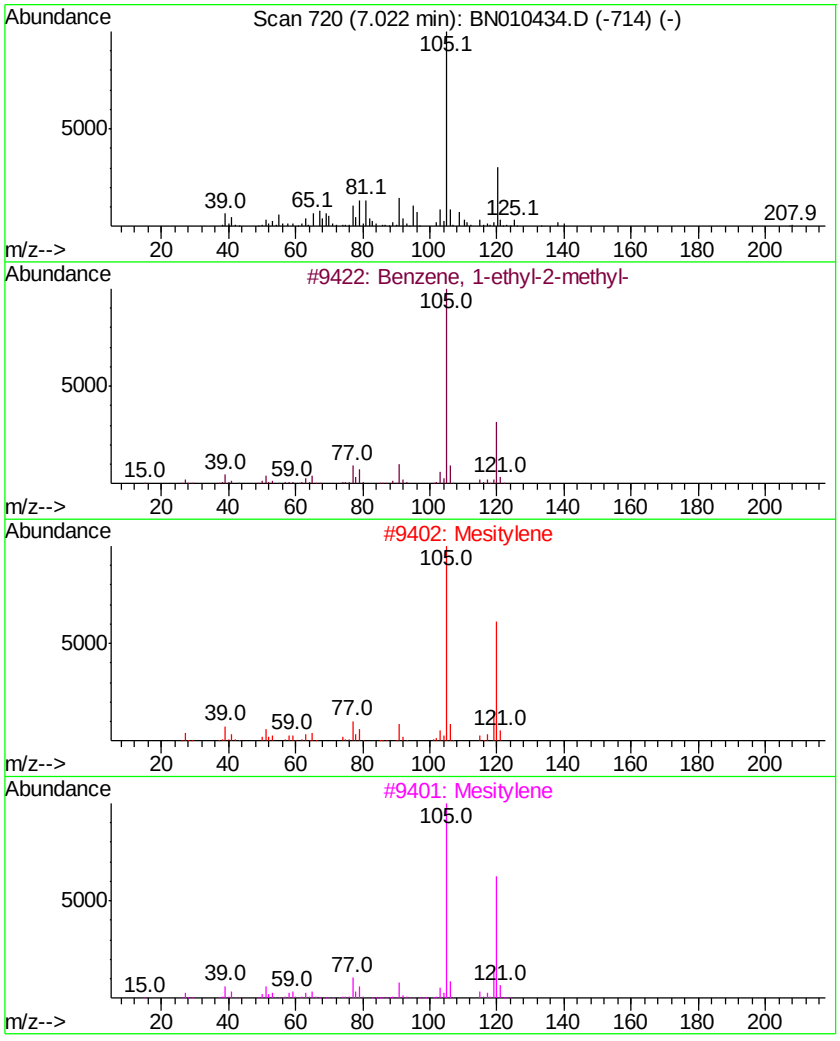

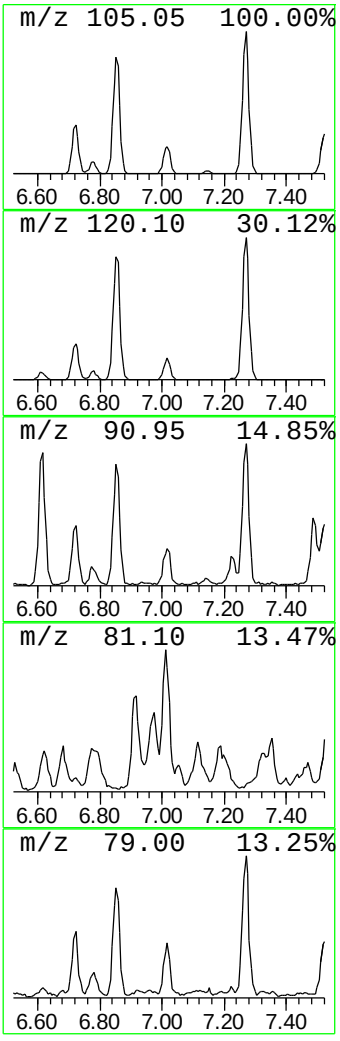

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Quant Title : SVOA CALIBRATION

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

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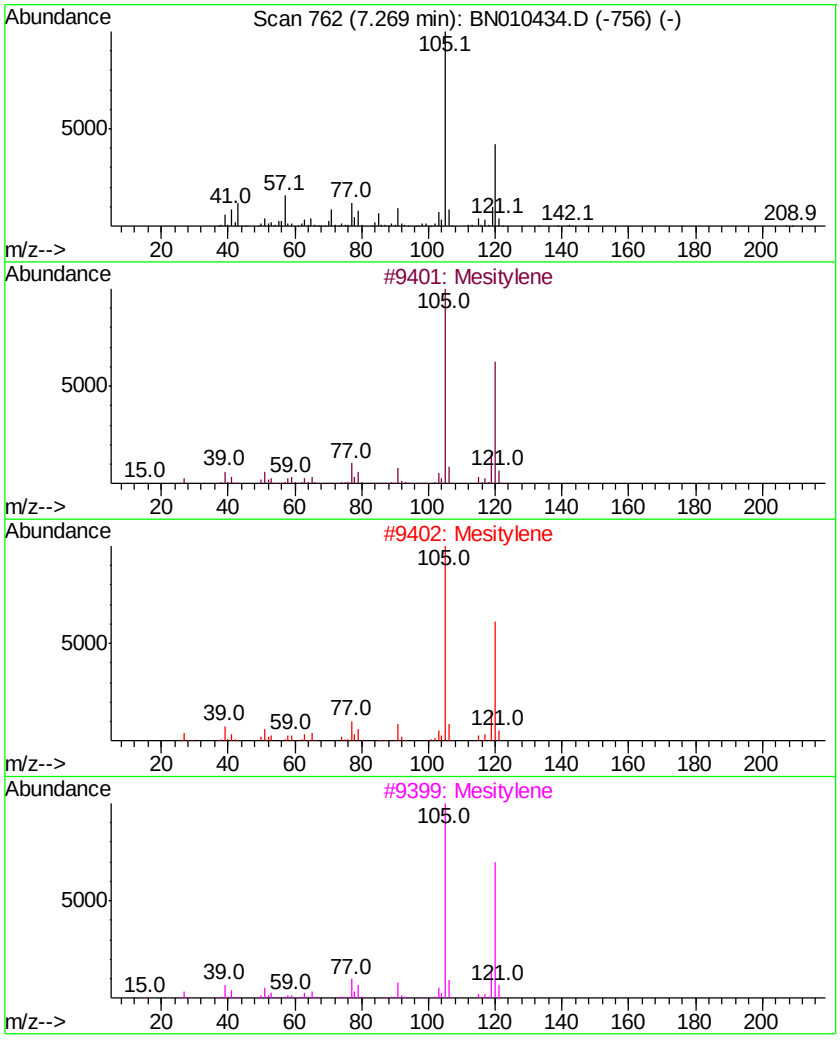
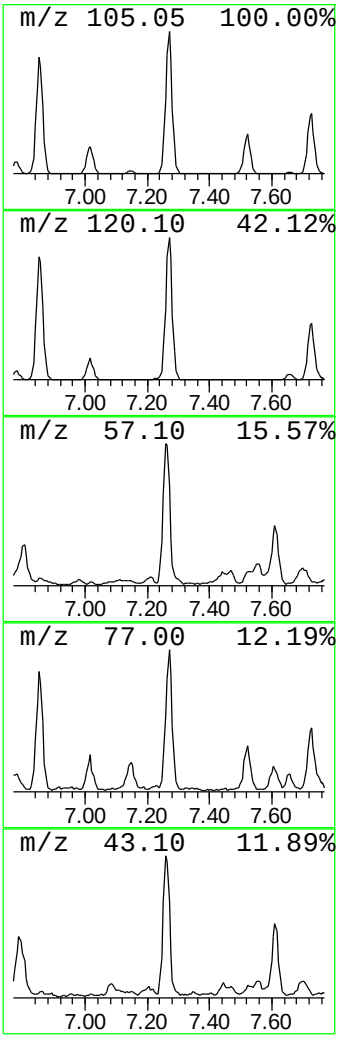
Peak Number 15 Mesitylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.27	18.44 ng/ul	3090590	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	91
2	Mesitylene	120	C9H12	000108-67-8	91
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Mesitylene	120	C9H12	000108-67-8	91





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Quant Title : SVOA CALIBRATION

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

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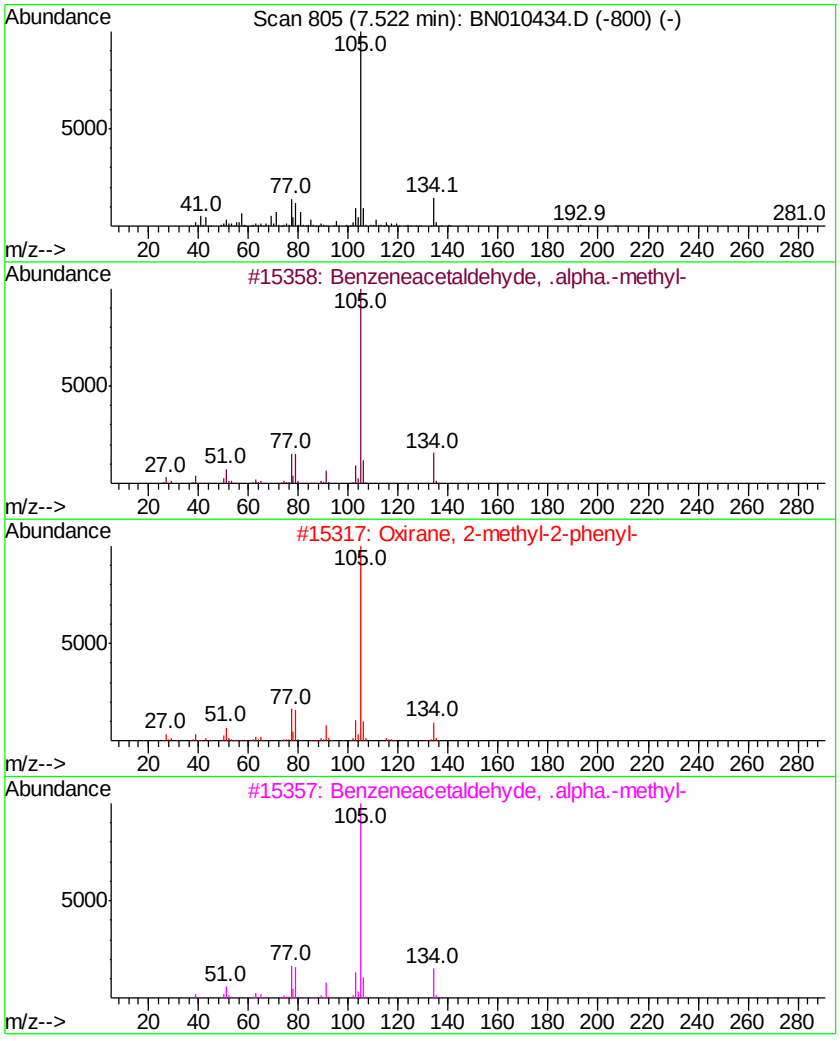
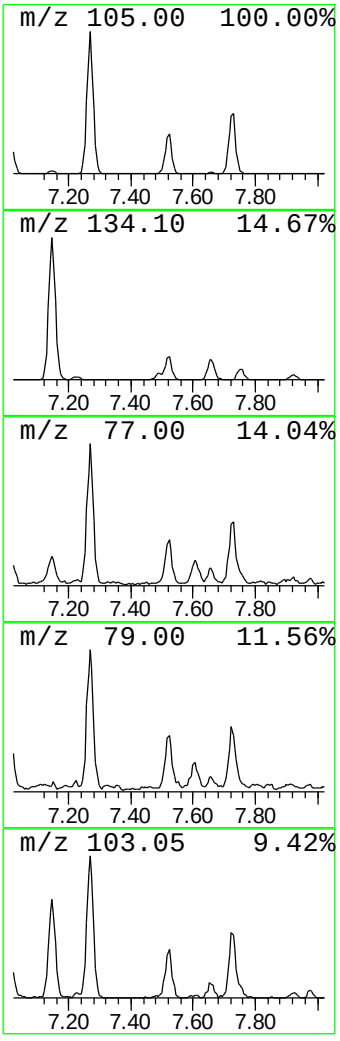
Peak Number 16 Benzeneacetaldehyde, .alpha... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.52	5.53 ng/ul	927200	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	87
2	Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	83
3	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	83
4	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	83
5	2-Tolyloxirane	134	C9H10O	002783-26-8	80





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Quant Title : SVOA CALIBRATION

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

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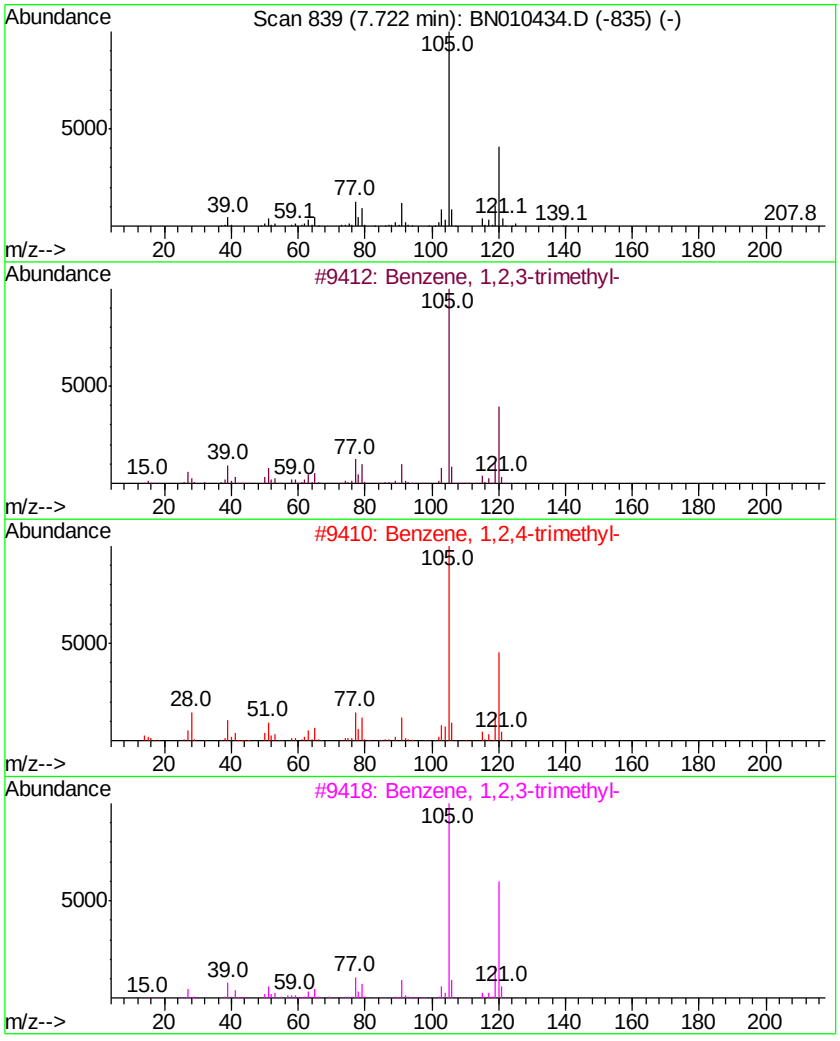
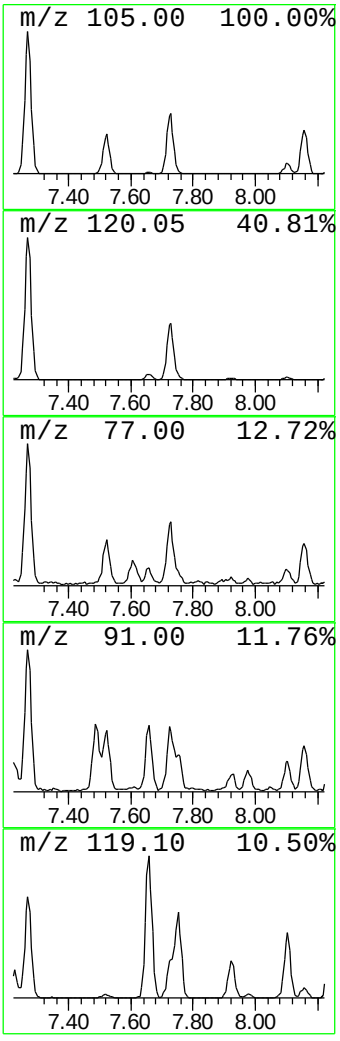
Peak Number 17 Benzene, 1,2,4-trimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.72	7.37 ng/ul	1236330	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
5	Mesitylene	120	C9H12	000108-67-8	91





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Sample : L2155-02
Misc :
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Instrument : BNA_N
ClientSampleId : DBFA1

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Quant Title : SVOA CALIBRATION

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

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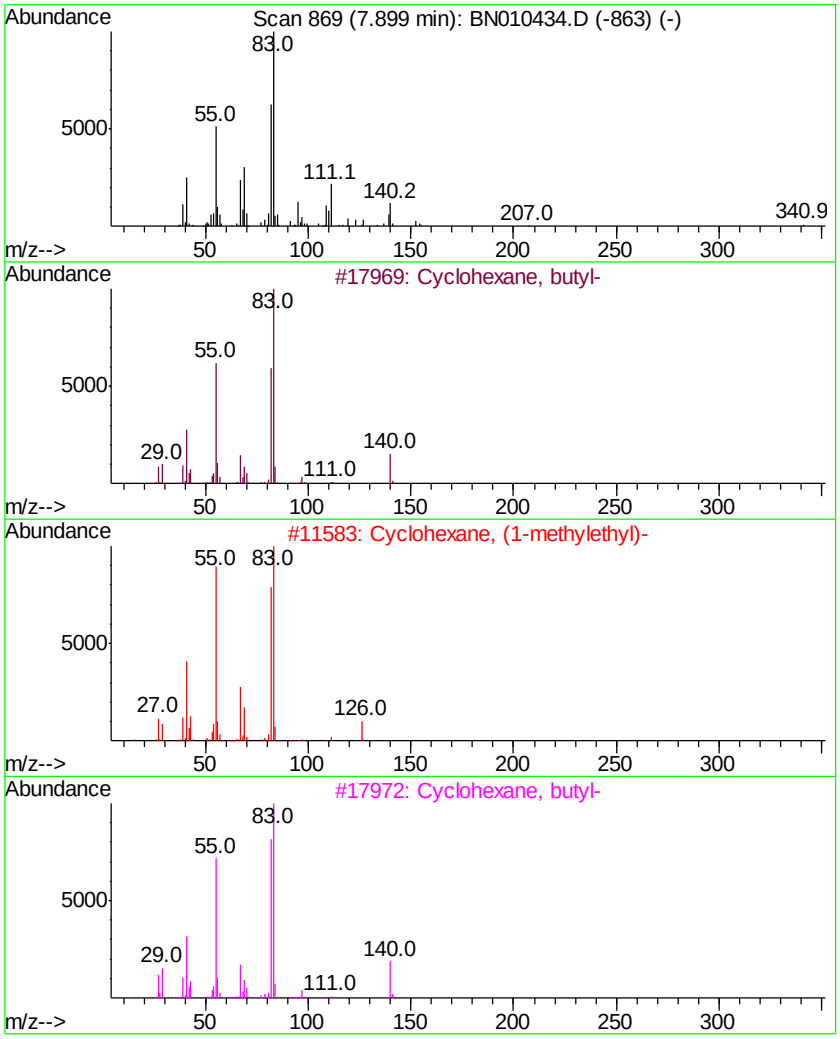
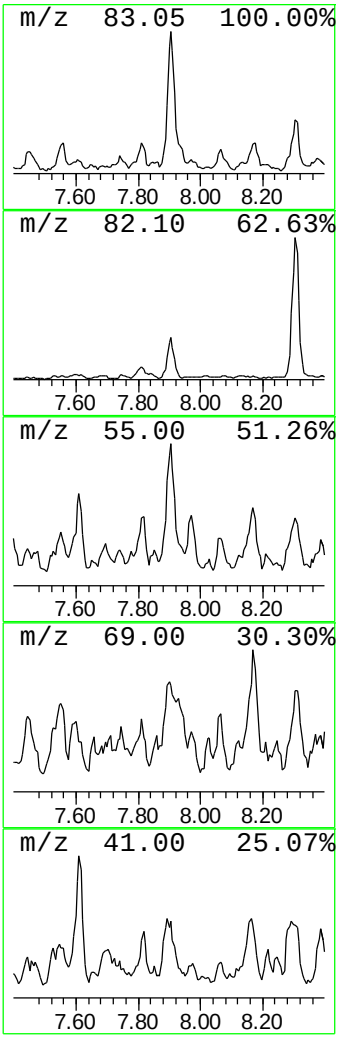
Peak Number 18 (DEL) Alkane: Cyclic7.90 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.90	3.98 ng/ul	666540	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, butyl-	140	C10H20	001678-93-9	62
2	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	59
3	Cyclohexane, butyl-	140	C10H20	001678-93-9	58
4	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	53
5	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	53





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ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBFA1

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Quant Title : SVOA CALIBRATION

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

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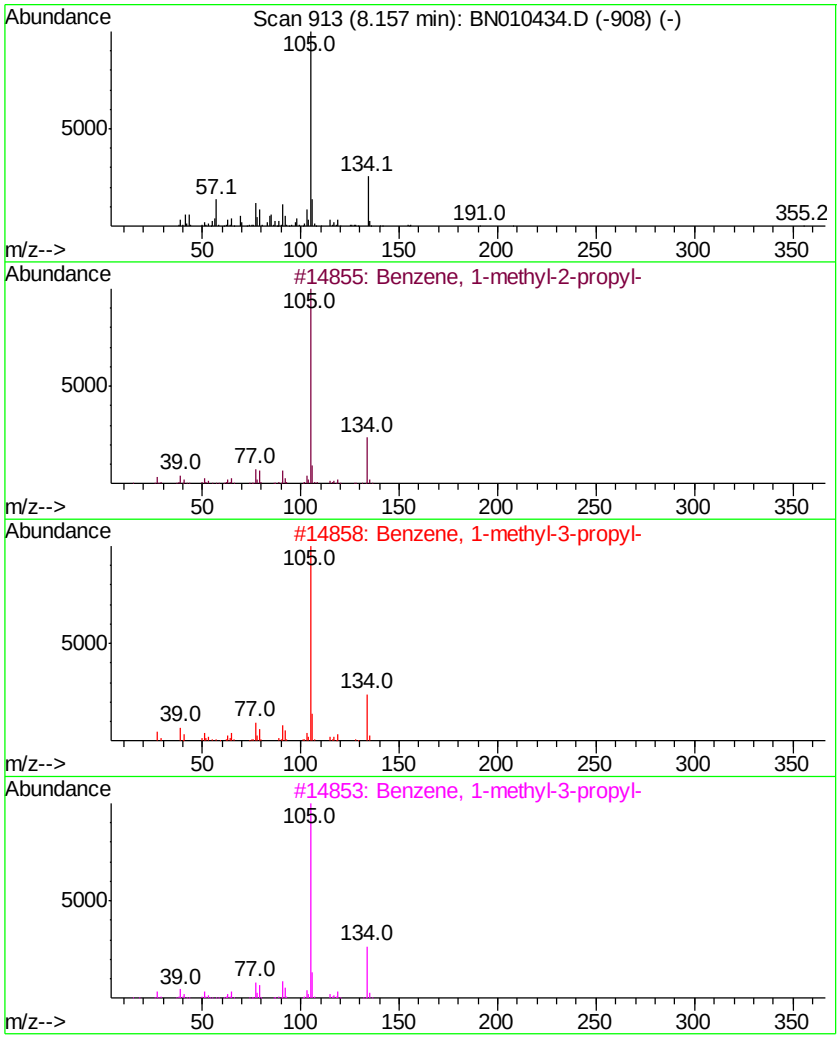
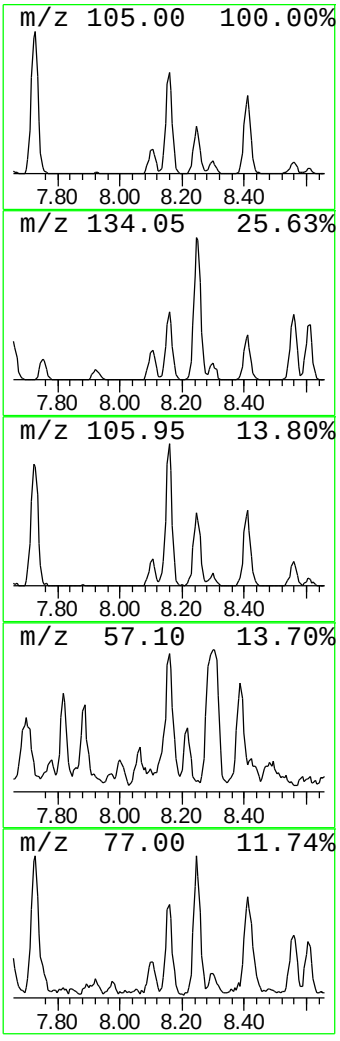
Peak Number 19 Benzene, 1-methyl-2-propyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.16	6.04 ng/ul	1012100	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	93
2	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	87
3	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	87
4	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	81
5	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	68





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Quant Title : SVOA CALIBRATION

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

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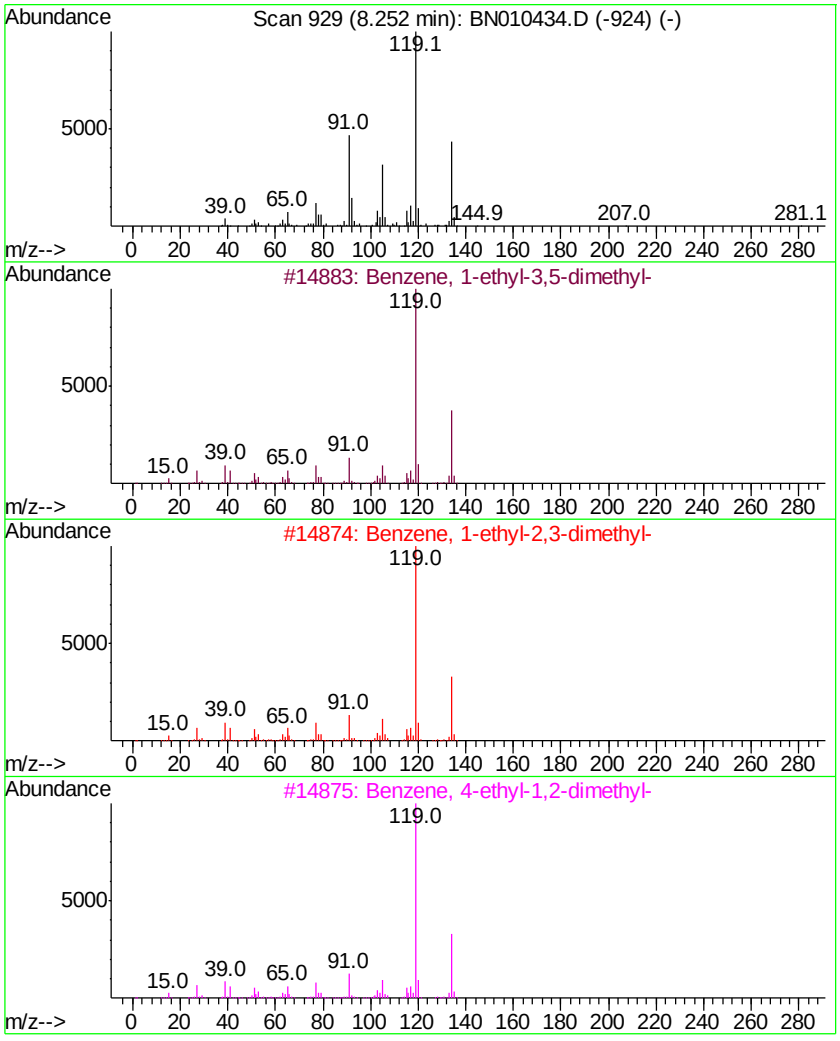
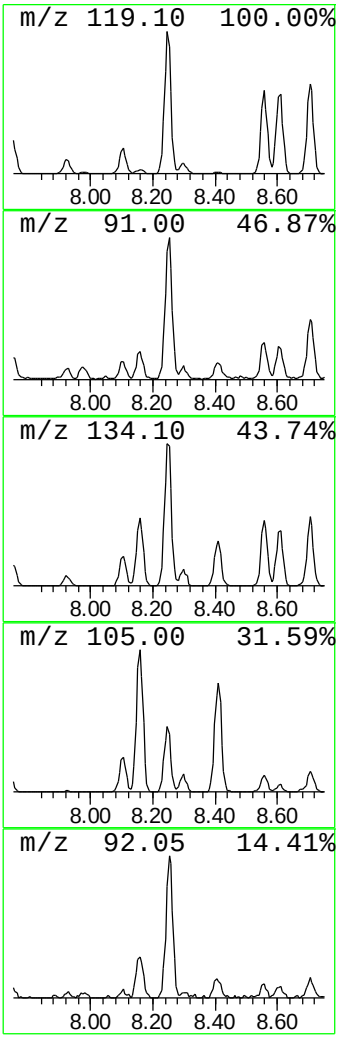
Peak Number 20 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.25	7.78 ng/ul	1304110	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	95
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	93
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91





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ClientSampleId : DBFA1

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Quant Title : SVOA CALIBRATION

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

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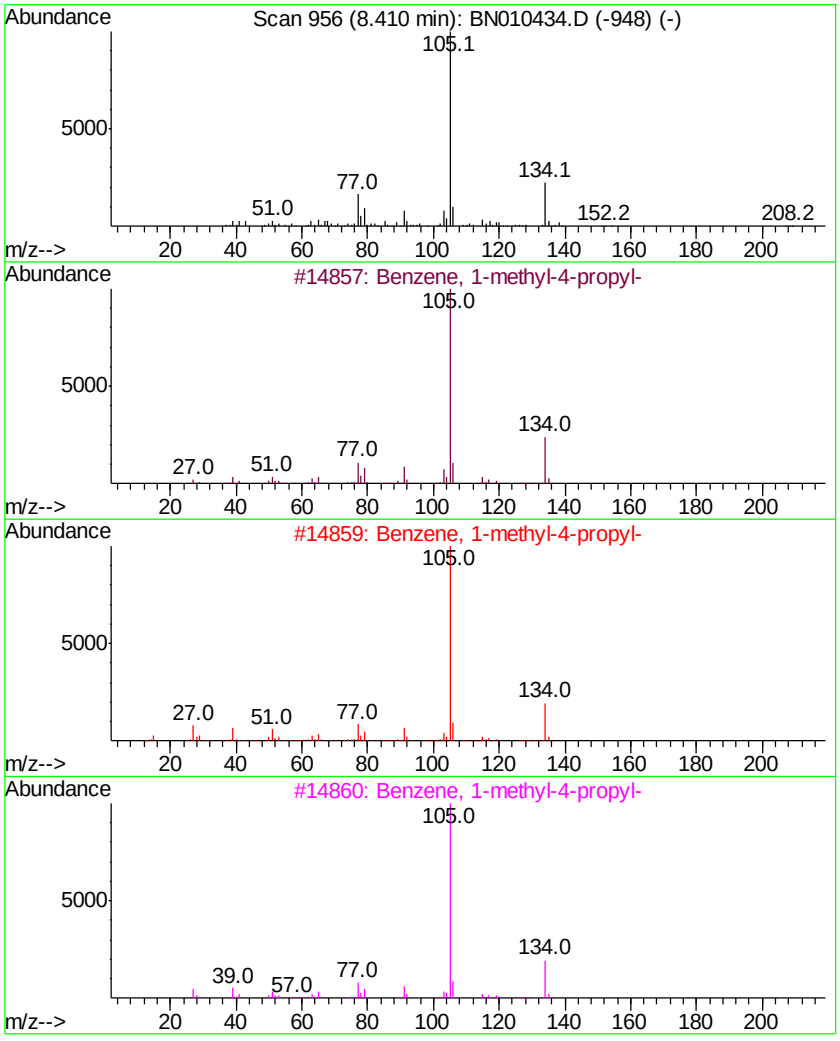
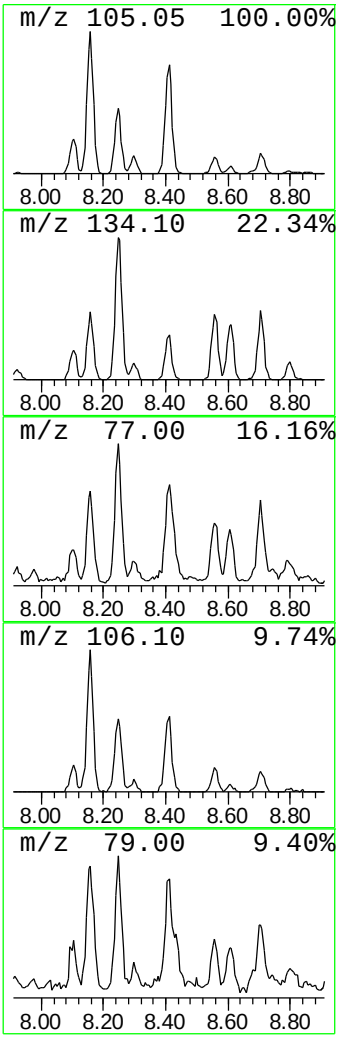
Peak Number 21 Benzene, 1-methyl-4-propyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.41	6.00 ng/ul	1005630	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	94
2	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
3	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
4	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	90
5	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	90





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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 22 p-Cymene Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.56	2.94 ng/ul	493118	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3	p-Cymene	134	C10H14	000099-87-6	95
4	o-Cymene	134	C10H14	000527-84-4	95
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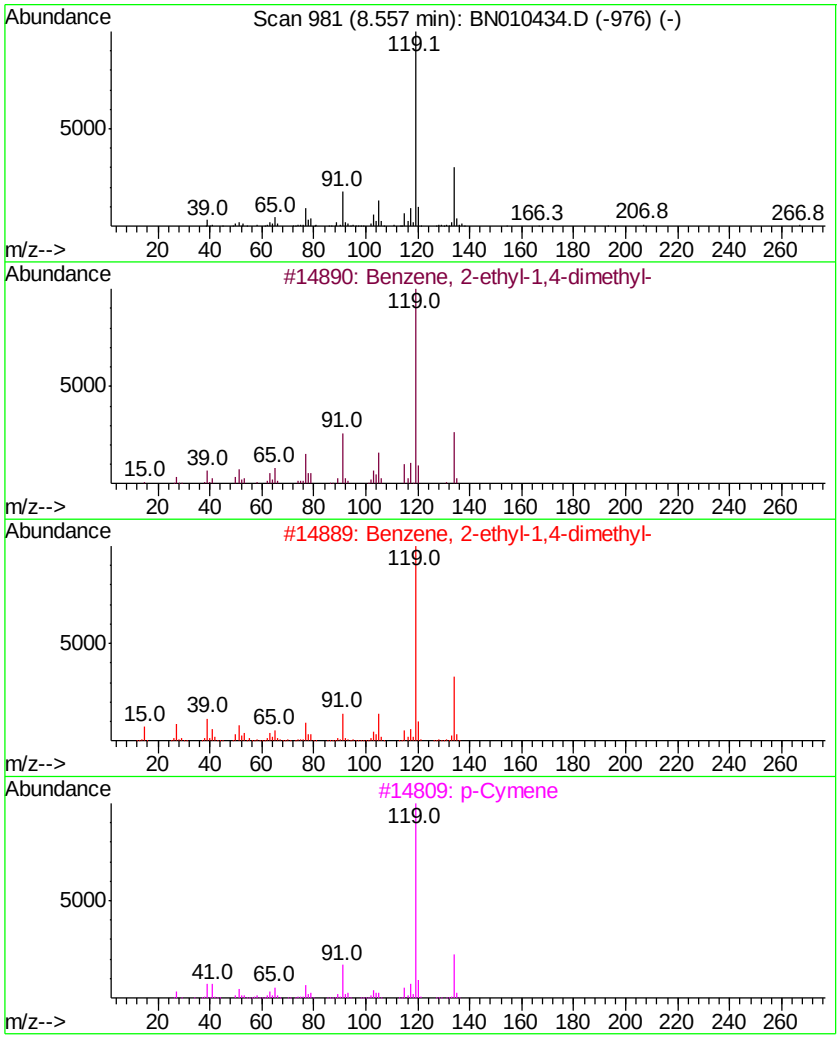

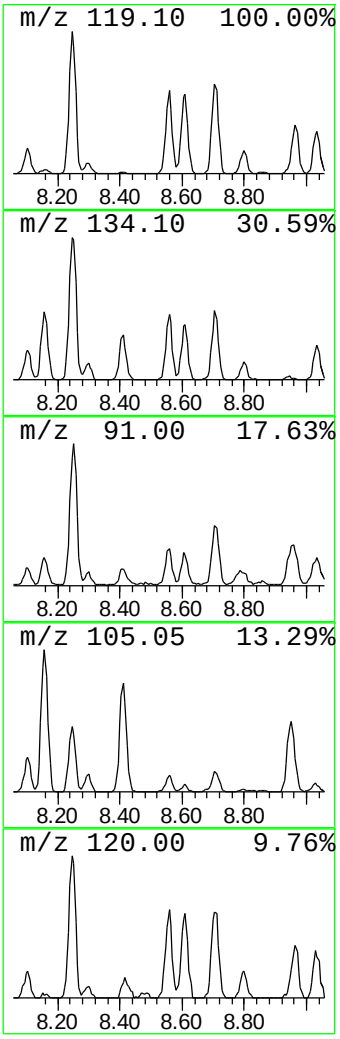

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Quant Title : SVOA CALIBRATION

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

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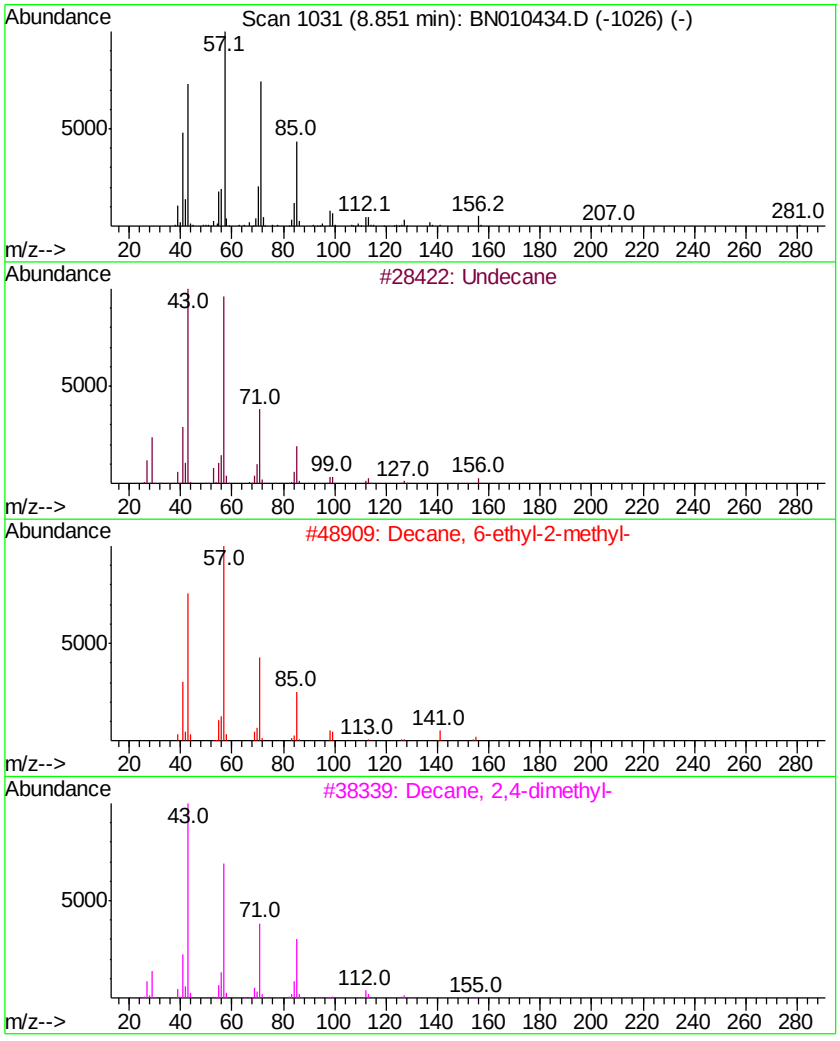
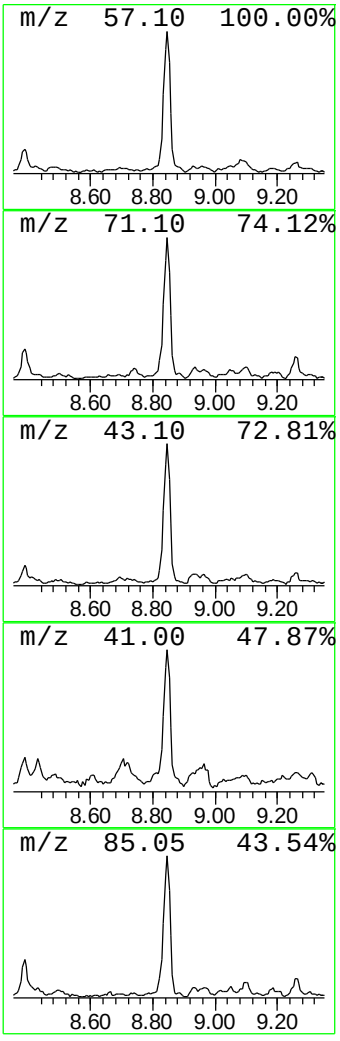
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.85	7.18 ng/ul	1203040	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane	156	C11H24	001120-21-4	87
2	Decane, 6-ethyl-2-methyl-	184	C13H28	062108-21-8	64
3	Decane, 2,4-dimethyl-	170	C12H26	002801-84-5	64
4	Decane	142	C10H22	000124-18-5	59
5	Undecane	156	C11H24	001120-21-4	58





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Quant Title : SVOA CALIBRATION

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

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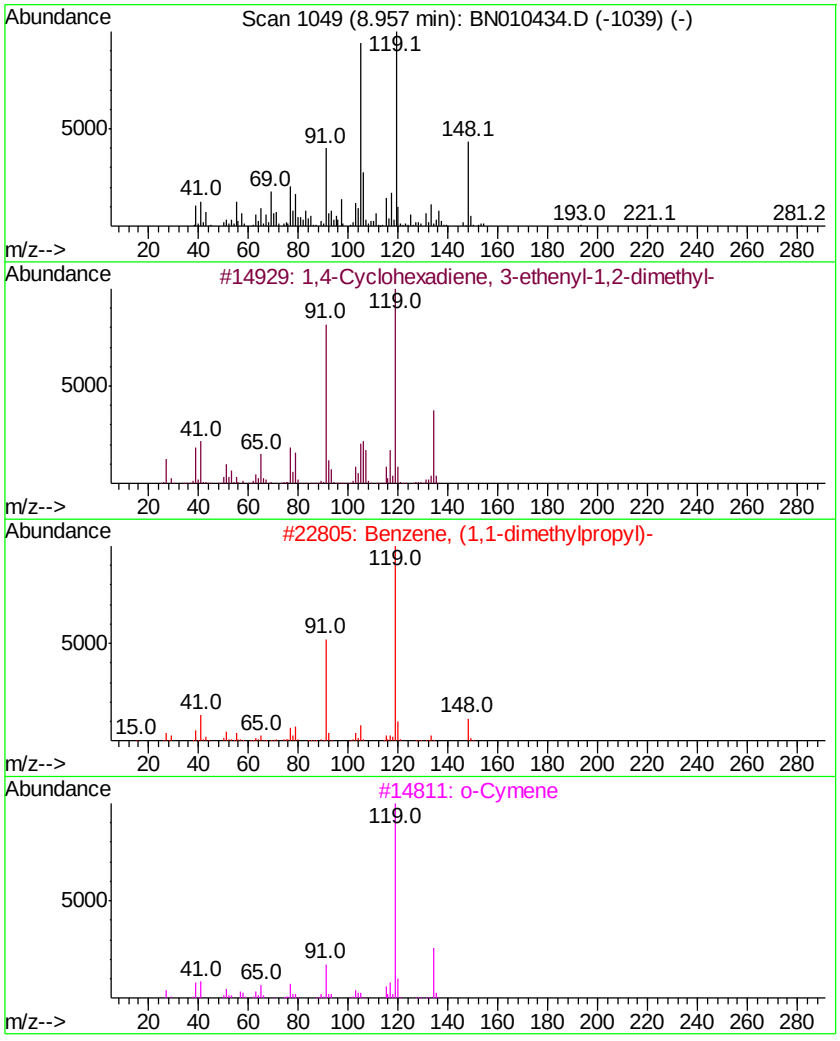
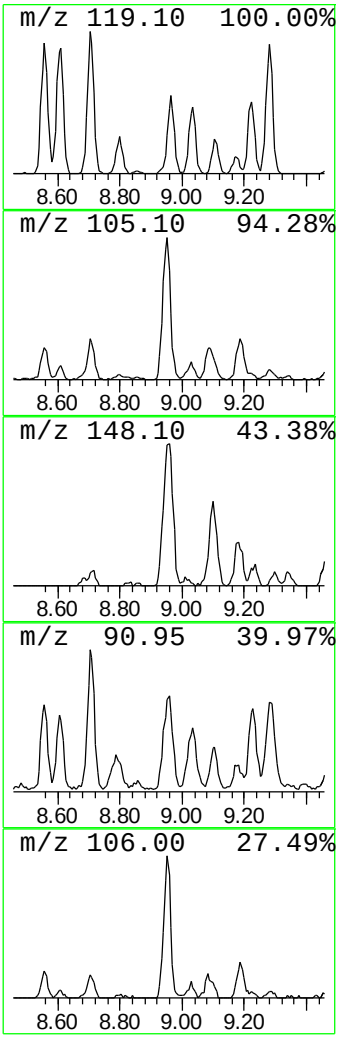
Peak Number 25 1,4-Cyclohexadiene, 3-ethen... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.96	7.84 ng/ul	1314080	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Cyclohexadiene, 3-ethenyl-1,...	134	C10H14	062338-57-2	53
2	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	52
3	o-Cymene	134	C10H14	000527-84-4	49
4	p-Cymene	134	C10H14	000099-87-6	46
5	Benzene, (1-methylbutyl)-	148	C11H16	002719-52-0	42





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Quant Title : SVOA CALIBRATION

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

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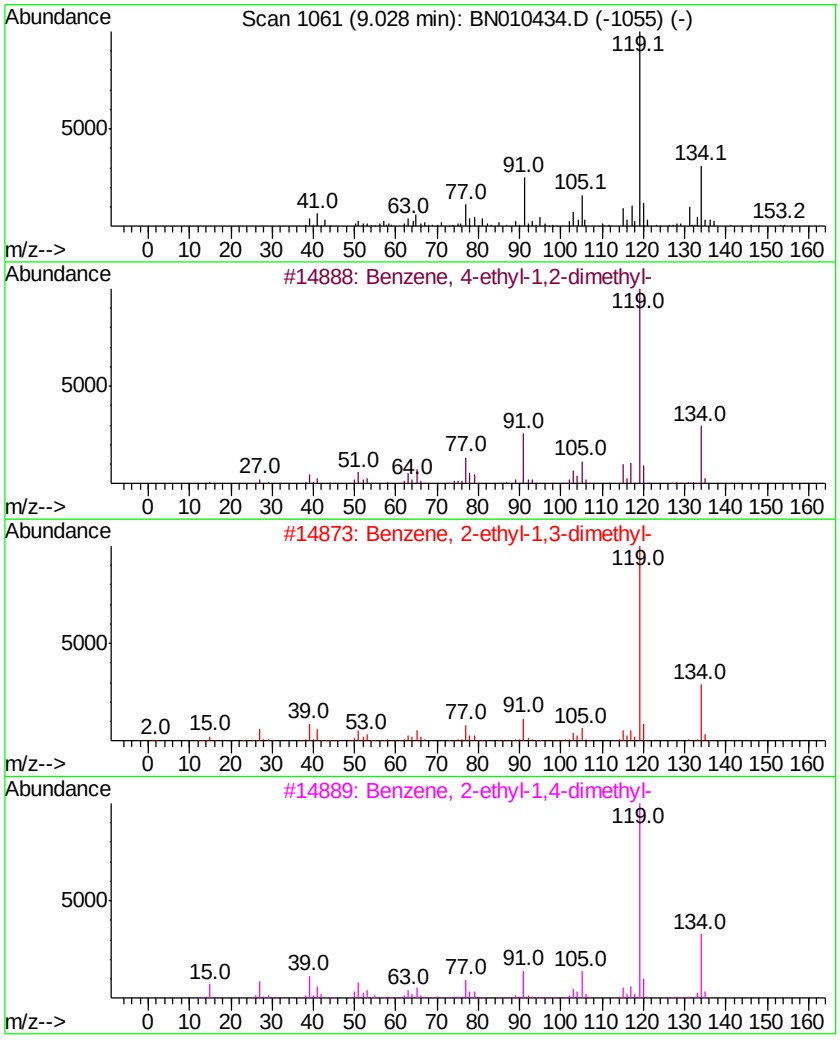
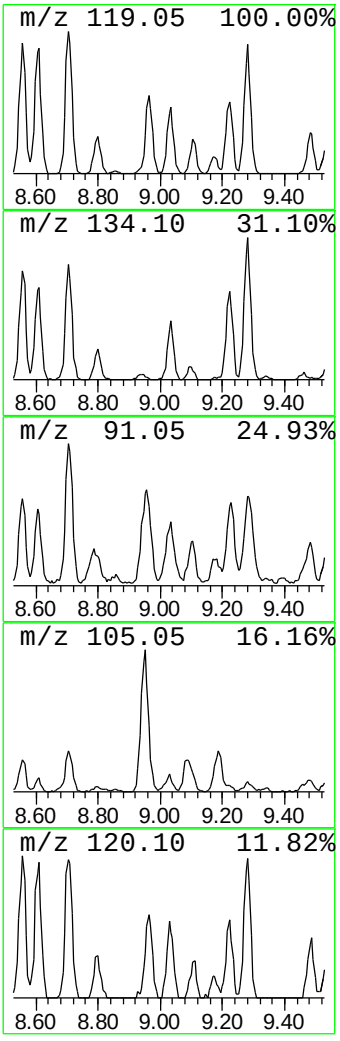
Peak Number 26 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.03	2.48 ng/ul	575730	Naphthalene-d8	10.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93





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Quant Title : SVOA CALIBRATION

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

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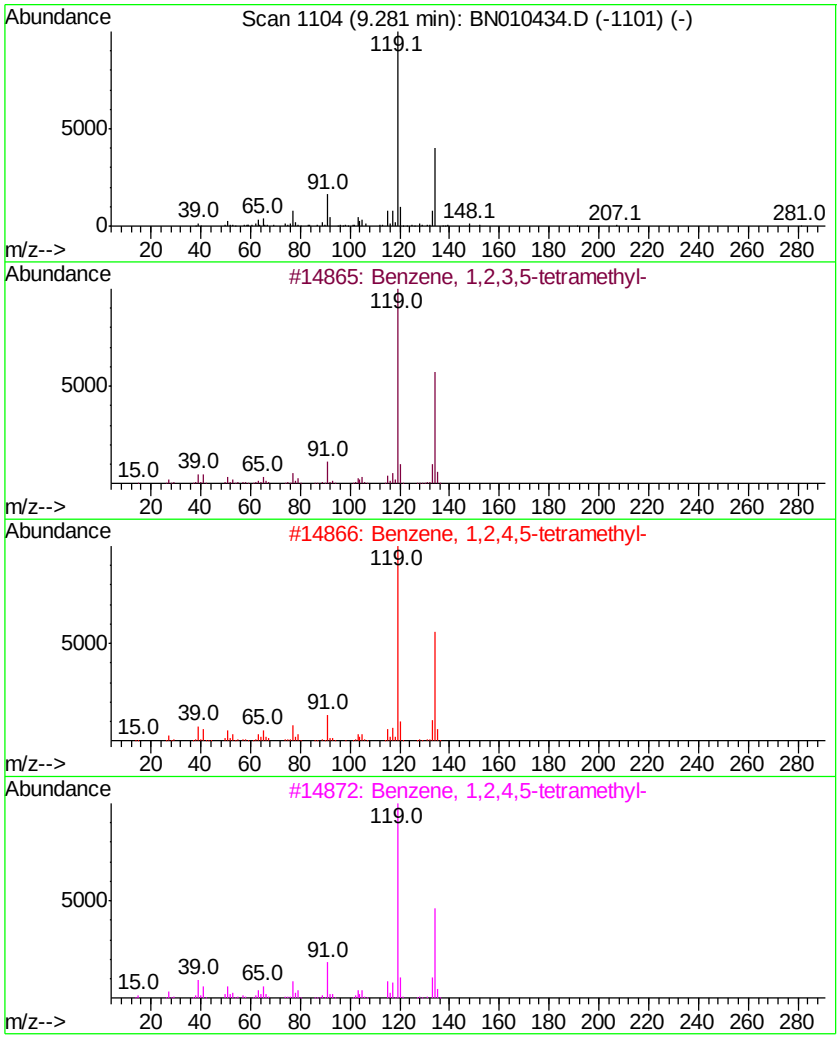
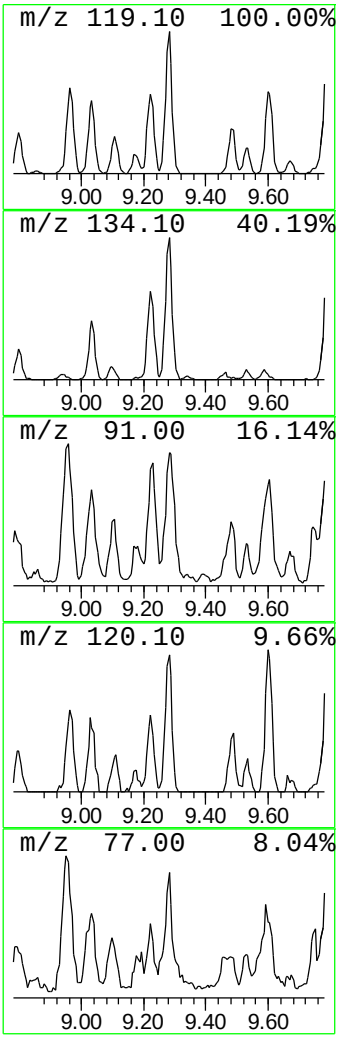
Peak Number 27 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.28	3.24 ng/ul	752331	Naphthalene-d8	10.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91





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TIC Integration Parameters: LSCINT.P

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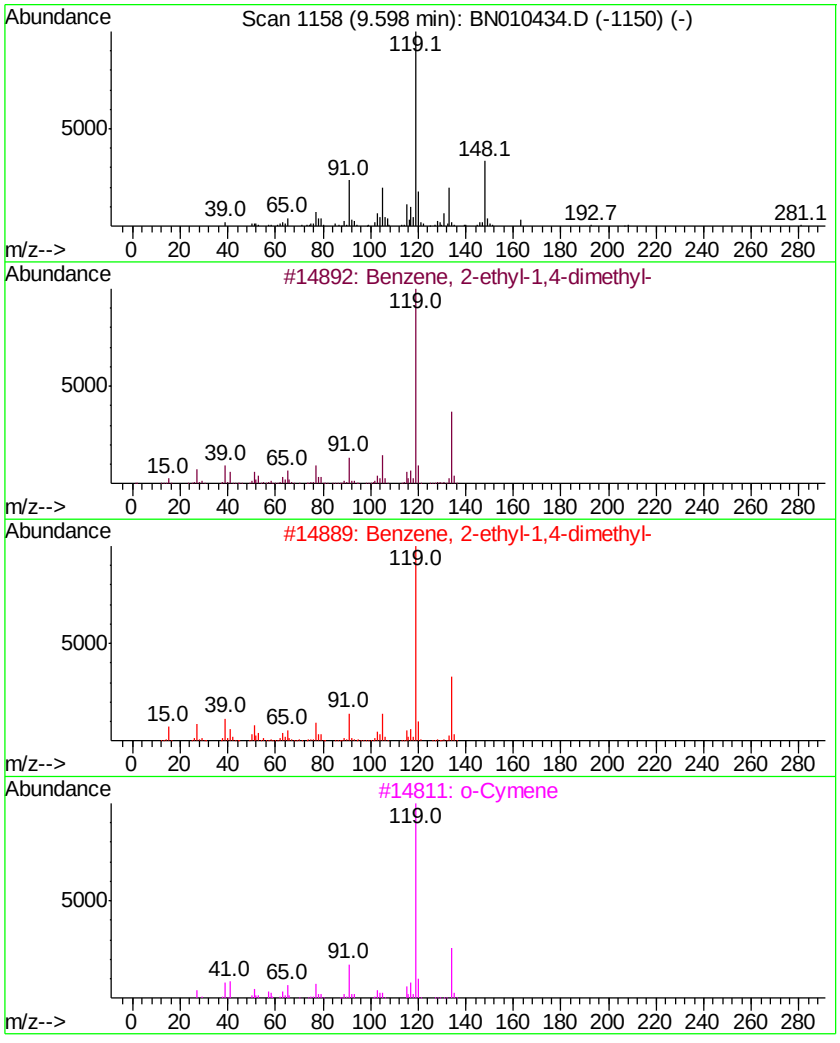
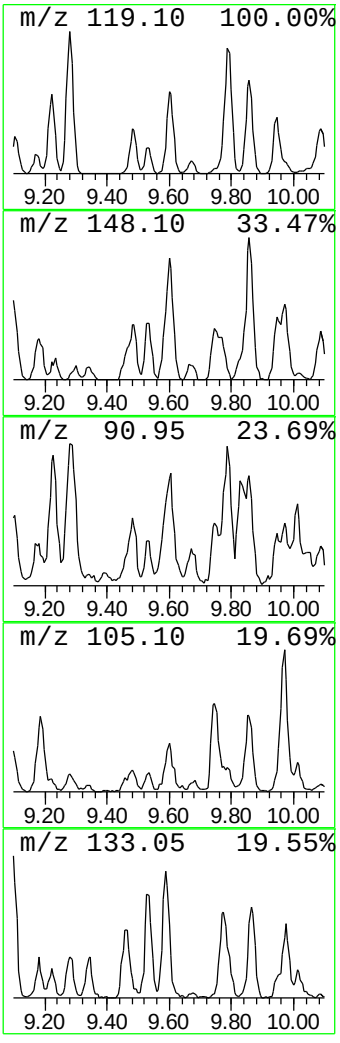
Peak Number 28 o-Cymene Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.60	3.51 ng/ul	814649	Naphthalene-d8	10.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	70
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	70
3	o-Cymene	134	C10H14	000527-84-4	55
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	55
5	o-Cymene	134	C10H14	000527-84-4	55





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TIC Integration Parameters: LSCINT.P

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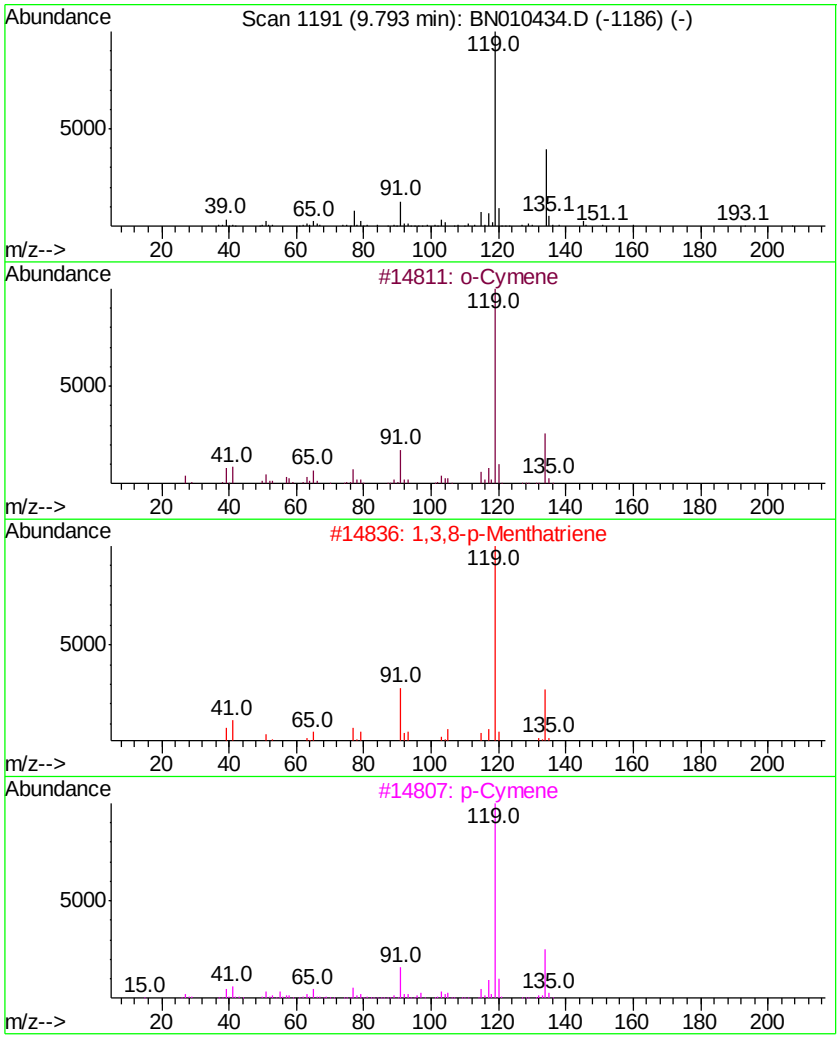
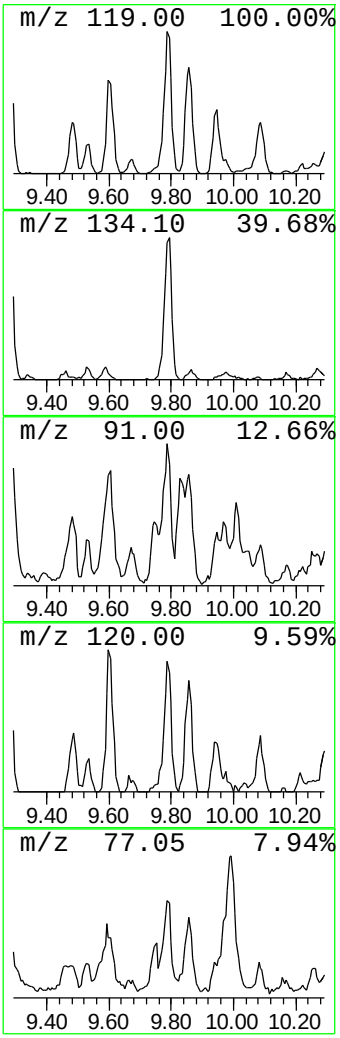
Peak Number 29 1,3,8-p-Menthatriene Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.79	2.60 ng/ul	603110	Naphthalene-d8	10.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	90
2	1,3,8-p-Menthatriene	134	C10H14	018368-95-1	90
3	p-Cymene	134	C10H14	000099-87-6	87
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	87
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	87





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Data File : BN010434.D
Acq On : 08 Apr 2020 15:44
Operator : CG/JU
Sample : L2155-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBFA1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

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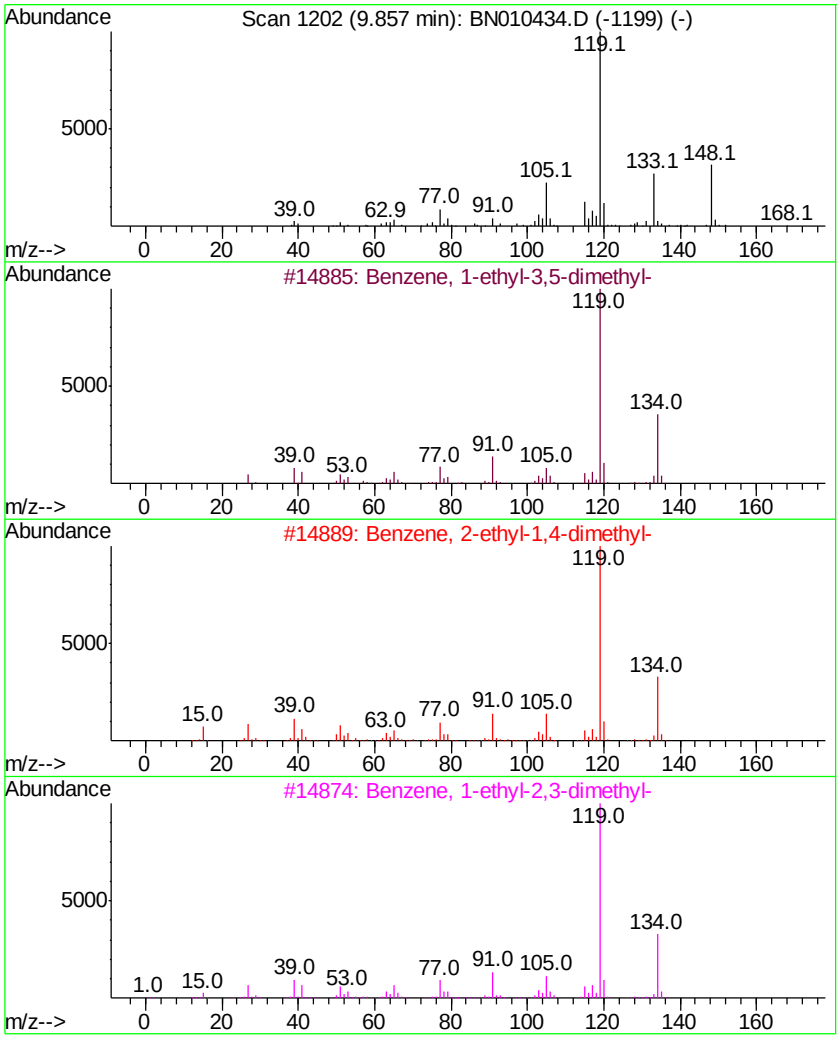
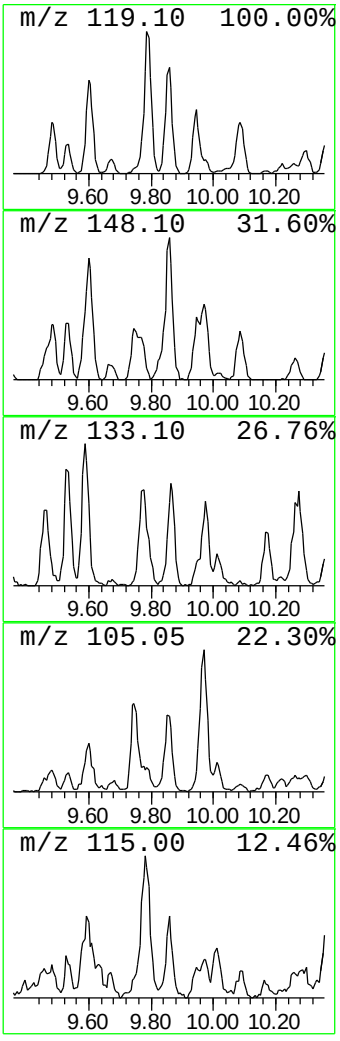
Peak Number 30 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.86	2.18 ng/ul	506525	Naphthalene-d8	10.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	58
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	58
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	58
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	53
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	52





Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040820\
Data File : BN010434.D
Acq On : 08 Apr 2020 15:44
Operator : CG/JU
Sample : L2155-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBFA1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

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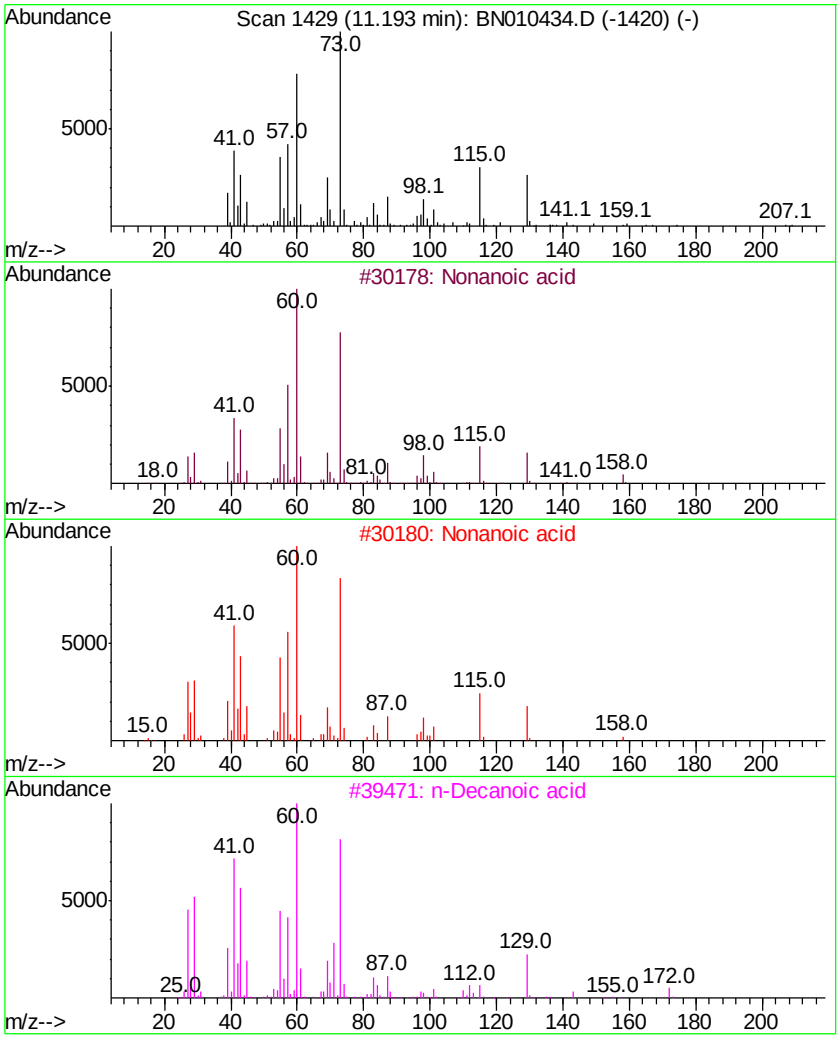
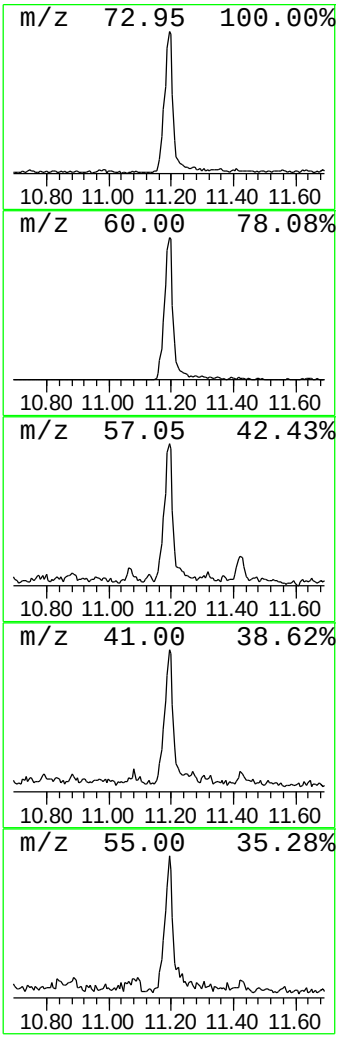
Peak Number 31 Nonanoic acid Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.19	3.10 ng/ul	719816	Naphthalene-d8	10.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanoic acid	158	C9H18O2	000112-05-0	64
2	Nonanoic acid	158	C9H18O2	000112-05-0	59
3	n-Decanoic acid	172	C10H20O2	000334-48-5	59
4	Hexanoic acid	116	C6H12O2	000142-62-1	50
5	n-Decanoic acid	172	C10H20O2	000334-48-5	50





Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040820\
Data File : BN010434.D
Acq On : 08 Apr 2020 15:44
Operator : CG/JU
Sample : L2155-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBFA1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

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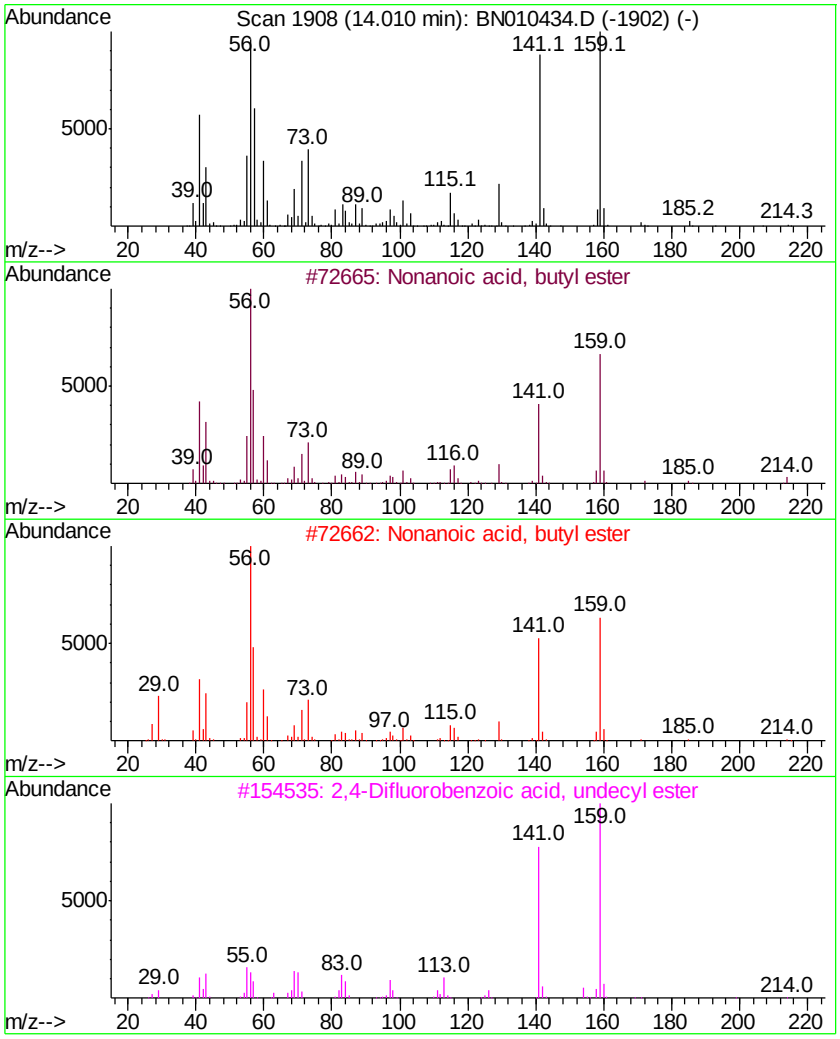
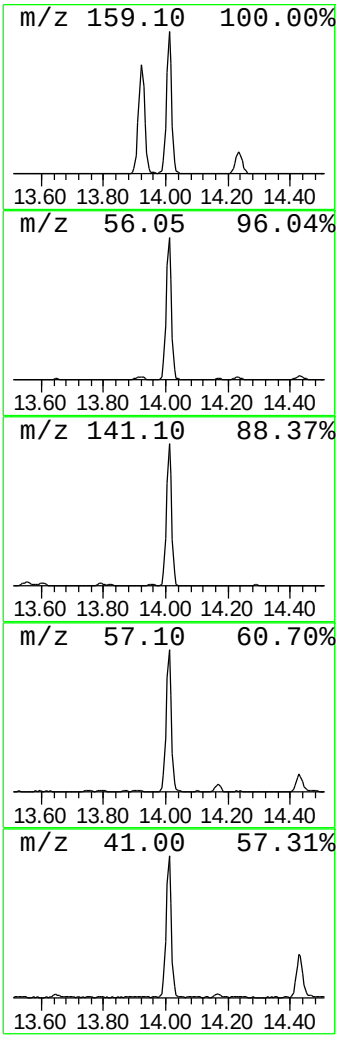
Peak Number 32 Nonanoic acid, butyl ester Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.01	11.18 ng/ul	3364550	Acenaphthene-d10	14.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanoic acid, butyl ester	214	C13H26O2	050623-57-9	90
2	Nonanoic acid, butyl ester	214	C13H26O2	050623-57-9	70
3	2,4-Difluorobenzoic acid, undecy...	312	C18H26F2O2	1000338-78-9	46
4	2,4-Difluorobenzoic acid, 8-chlo...	304	C15H19ClF2O2	1000360-56-2	43
5	2,4-Difluorobenzoic acid, nonyl ...	284	C16H22F2O2	1000338-78-7	43





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040820\
Data File : BN010434.D
Acq On : 08 Apr 2020 15:44
Operator : CG/JU
Sample : L2155-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBFA1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

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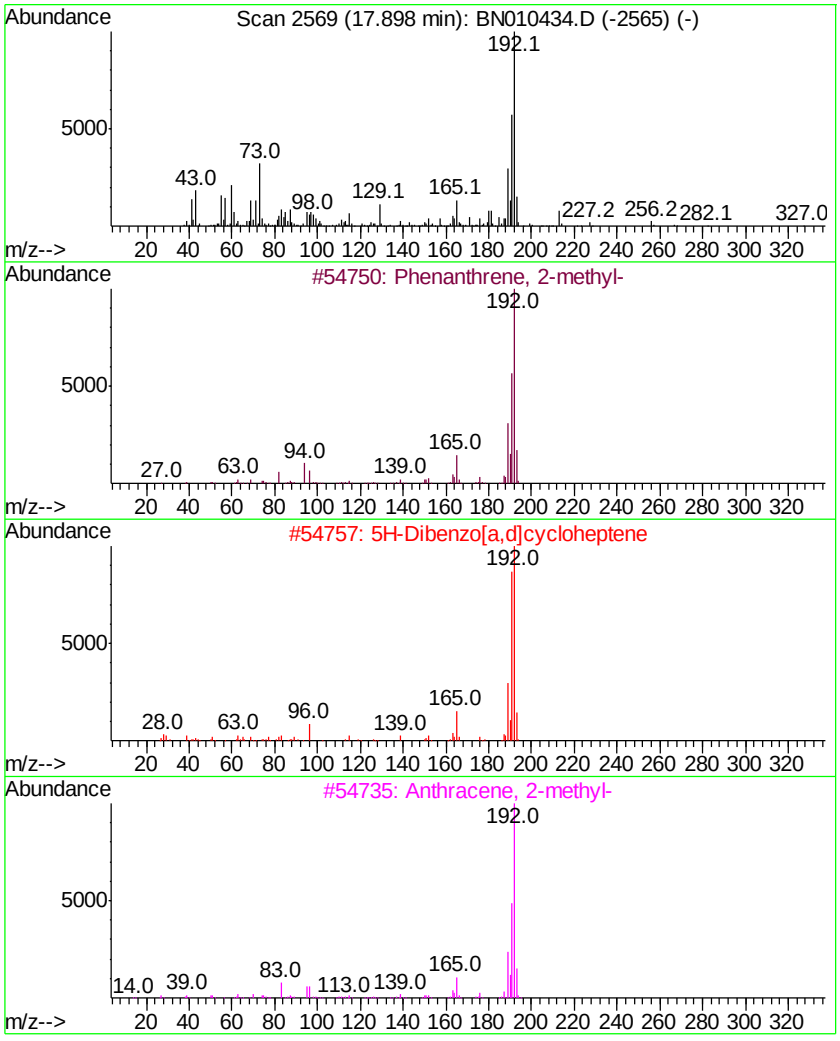
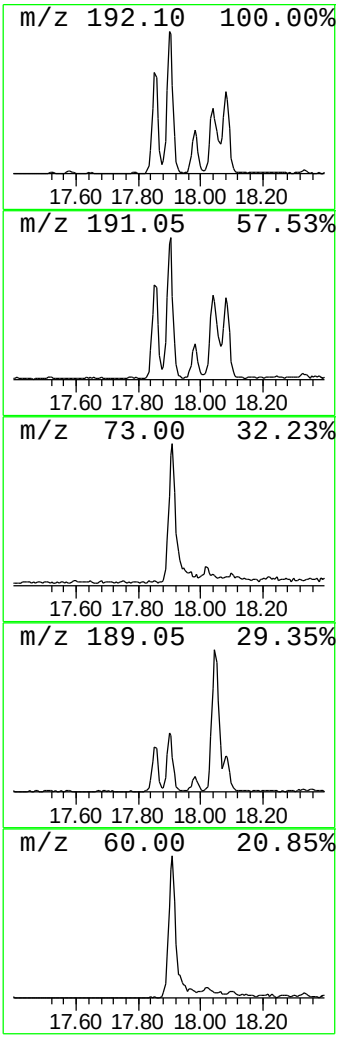
Peak Number 33 5H-Dibenzo[a,d]cycloheptene Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.90	4.36 ng/ul	1653030	Phenanthrene-d10	16.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	96
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	93





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN040820\
Data File : BN010434.D
Acq On : 08 Apr 2020 15:44
Operator : CG/JU
Sample : L2155-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBFA1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.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
1-Butanol	2.83	3.1	ng/ul	516763	1	7.60	3352830	20.0
(DEL) Alkane: Cyc...	2.89	6.1	ng/ul	1018020	1	7.60	3352830	20.0
Butane, 2-methoxy...	2.96	15.9	ng/ul	2673030	1	7.60	3352830	20.0
Ethylbenzene	5.18	2.8	ng/ul	462019	1	7.60	3352830	20.0
p-Xylene	5.31	26.0	ng/ul	4359820	1	7.60	3352830	20.0
o-Xylene	5.67	22.8	ng/ul	3825270	1	7.60	3352830	20.0
3,5-Dimethyl-3-he...	5.93	4.5	ng/ul	747963	1	7.60	3352830	20.0
Benzene, (1-methy...	6.13	9.2	ng/ul	1546720	1	7.60	3352830	20.0
Cyclopentane, 1-m...	6.29	3.4	ng/ul	571450	1	7.60	3352830	20.0
(DEL) Alkane: Str...	6.32	3.4	ng/ul	561898	1	7.60	3352830	20.0
unknown-01	6.61	2.6	ng/ul	437053	1	7.60	3352830	20.0
Benzene, 1-ethyl-...	6.72	6.0	ng/ul	1007860	1	7.60	3352830	20.0
Benzene, 1,2,3-tr...	6.85	11.6	ng/ul	1943330	1	7.60	3352830	20.0
Benzene, 1-ethyl-...	7.02	3.6	ng/ul	600952	1	7.60	3352830	20.0
Mesitylene	7.27	18.4	ng/ul	3090590	1	7.60	3352830	20.0
Benzeneacetaldehy...	7.52	5.5	ng/ul	927200	1	7.60	3352830	20.0
Benzene, 1,2,4-tr...	7.72	7.4	ng/ul	1236330	1	7.60	3352830	20.0
(DEL) Alkane: Cyc...	7.90	4.0	ng/ul	666540	1	7.60	3352830	20.0
Benzene, 1-methyl...	8.16	6.0	ng/ul	1012100	1	7.60	3352830	20.0
Benzene, 1-ethyl-...	8.25	7.8	ng/ul	1304110	1	7.60	3352830	20.0
Benzene, 1-methyl...	8.41	6.0	ng/ul	1005630	1	7.60	3352830	20.0
p-Cymene	8.56	2.9	ng/ul	493118	1	7.60	3352830	20.0
(DEL) Alkane: Str...	8.85	7.2	ng/ul	1203040	1	7.60	3352830	20.0
1,4-Cyclohexadien...	8.96	7.8	ng/ul	1314080	1	7.60	3352830	20.0
Benzene, 4-ethyl-...	9.03	2.5	ng/ul	575730	2	10.36	4640540	20.0
Benzene, 1,2,3,5-...	9.28	3.2	ng/ul	752331	2	10.36	4640540	20.0
o-Cymene	9.60	3.5	ng/ul	814649	2	10.36	4640540	20.0
1,3,8-p-Menthatriene	9.79	2.6	ng/ul	603110	2	10.36	4640540	20.0
Benzene, 2-ethyl-...	9.86	2.2	ng/ul	506525	2	10.36	4640540	20.0
Nonanoic acid	11.19	3.1	ng/ul	719816	2	10.36	4640540	20.0
Nonanoic acid, bu...	14.01	11.2	ng/ul	3364550	3	14.23	6018520	20.0
5H-Dibenzo[a,d]cy...	17.90	4.4	ng/ul	1653030	4	16.97	7586650	20.0