LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090819\
Data File : BM022608.D
Acq On : 09 Sep 2019 17:55
Operator : HP/JU
Sample : K4706-07DL 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF573DL

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-BM090519MA.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.046	6	10	29	rVB	366856	587334	2.38%	0.785%
2	5.375	399	406	412	rBV	34006	53485	0.22%	0.072%
3	5.504	423	428	439	rBV	65247	136114	0.55%	0.182%
4	5.875	481	491	497	rVB3	46908	101039	0.41%	0.135%
5	7.010	679	684	693	rVV	47134	76659	0.31%	0.103%
6	7.087	693	697	703	rVB	20736	30888	0.13%	0.041%
7	7.181	708	713	721	rBV	90499	152412	0.62%	0.204%
8	7.387	741	748	754	rBV	106126	171536	0.70%	0.229%
9	7.510	762	769	775	rBV2	78912	124174	0.50%	0.166%
10	7.575	775	780	788	rVB	61053	101325	0.41%	0.136%
11	7.851	820	827	837	rBV	1045743	1653334	6.71%	2.211%
12	7.975	842	848	854	rBV	29754	47325	0.19%	0.063%
13	8.234	884	892	900	rBV	545848	898906	3.65%	1.202%
14	8.392	906	919	927	rBV	892475	1414521	5.74%	1.892%
15	8.551	940	946	953	rVB	91423	148625	0.60%	0.199%
16	9.004	1018	1023	1028	rBV	108180	167301	0.68%	0.224%
17	9.210	1051	1058	1064	rBV	70882	111554	0.45%	0.149%
18	9.334	1068	1079	1085	rBV	179581	386475	1.57%	0.517%
19	9.392	1085	1089	1097	rVV	57448	92579	0.38%	0.124%
20	9.728	1137	1146	1153	rBV	66413	114363	0.46%	0.153%
21	9.904	1171	1176	1183	rBV2	31730	60248	0.24%	0.081%
22	10.057	1196	1202	1212	rBV3	73588	218664	0.89%	0.292%
23	10.151	1212	1218	1222	rVV	64523	124758	0.51%	0.167%
24	10.186	1222	1224	1231	rVV3	29254	41025	0.17%	0.055%
25	10.263	1231	1237	1247	rVB	147342	243285	0.99%	0.325%
26	10.651	1295	1303	1306	rVV	1431534	2574824	10.45%	3.444%
27	10.704	1306	1312	1320	rVV	14376287	24642429	100.00%	32.956%
28	10.816	1323	1331	1341	rVB	1309641	2233365	9.06%	2.987%
29	11.063	1368	1373	1381	rVB3	17428	36078	0.15%	0.048%
30	12.022	1530	1536	1543	rBV3	17615	33258	0.13%	0.044%
31	12.127	1549	1554	1561	rBV5	15690	29573	0.12%	0.040%
32	12.204	1561	1567	1575	rVB	38581	64748	0.26%	0.087%
33	12.310	1577	1585	1592	rBV	1017048	1617096	6.56%	2.163%
34	12.427	1598	1605	1611	rBV2	44650	88855	0.36%	0.119%

LSC Area Percent Report

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Acq On : 09 Sep 2019 17:55
Operator : HP/JU
Sample : K4706-07DL 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF573DL

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-BM090519MA.M
Title : SVOA CALIBRATION

35	12.527	1611	1622	1631	rVB	692293	1100612	4.47%	1.472%
36	12.945	1688	1693	1701	rVB4	21497	38649	0.16%	0.052%
37	13.316	1749	1756	1762	rVB	192224	284122	1.15%	0.380%
38	13.516	1785	1790	1802	rBV2	34715	66324	0.27%	0.089%
39	13.657	1808	1814	1822	rVB4	50154	128594	0.52%	0.172%
40	13.804	1832	1839	1845	rBV	98924	172169	0.70%	0.230%
41	13.886	1845	1853	1859	rVV	298906	490243	1.99%	0.656%
42	14.045	1874	1880	1883	rBV	32885	49540	0.20%	0.066%
43	14.174	1895	1902	1912	rBV2	319977	705721	2.86%	0.944%
44	14.486	1945	1955	1960	rVV	2380615	3642995	14.78%	4.872%
45	14.545	1960	1965	1971	rVV	812822	1162122	4.72%	1.554%
46	14.604	1971	1975	1980	rVV	210960	303519	1.23%	0.406%
47	14.657	1980	1984	1993	rVB	85441	139350	0.57%	0.186%
48	14.757	1995	2001	2005	rBV	72536	107866	0.44%	0.144%
49	14.880	2016	2022	2028	rVB2	598352	888395	3.61%	1.188%
50	14.951	2028	2034	2045	rVB	730641	1027099	4.17%	1.374%
51	15.068	2049	2054	2057	rBV4	18573	28027	0.11%	0.037%
52	15.186	2071	2074	2083	rVB	33891	51543	0.21%	0.069%
53	15.268	2083	2088	2091	rBV5	13826	28423	0.12%	0.038%
54	15.474	2117	2123	2127	rBV	449589	637731	2.59%	0.853%
55	15.527	2127	2132	2137	rVB	572078	761091	3.09%	1.018%
56	15.574	2137	2140	2145	rVB	56210	75650	0.31%	0.101%
57	15.651	2145	2153	2159	rBV	100950	203116	0.82%	0.272%
58	15.745	2164	2169	2173	rVV3	50864	104646	0.42%	0.140%
59	15.821	2177	2182	2186	rVB2	38259	57094	0.23%	0.076%
60	15.898	2191	2195	2198	rBV2	18235	27491	0.11%	0.037%
61	15.968	2203	2207	2208	rBV2	21636	26372	0.11%	0.035%
62	16.027	2213	2217	2226	rVB	94111	141149	0.57%	0.189%
63	16.192	2237	2245	2254	rBV	197884	309778	1.26%	0.414%
64	16.321	2261	2267	2272	rBV7	18251	40819	0.17%	0.055%
65	16.404	2276	2281	2288	rVV2	137053	240299	0.98%	0.321%
66	16.474	2288	2293	2297	rVV	116419	162767	0.66%	0.218%
67	16.527	2297	2302	2306	rVV3	24838	52981	0.21%	0.071%
68	16.574	2307	2310	2316	rVB7	20166	33387	0.14%	0.045%
69	16.686	2324	2329	2334	rBV5	22876	46176	0.19%	0.062%
70	16.998	2377	2382	2386	rBV2	102310	149390	0.61%	0.200%
71	17.039	2386	2389	2396	rVB	49262	68935	0.28%	0.092%

LSC Area Percent Report

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Sample : K4706-07DL 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF573DL

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

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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Title : SVOA CALIBRATION

72	17.115	2397	2402	2403	rBV2	19680	26490	0.11%	0.035%
73	17.168	2407	2411	2414	rBV2	27091	35243	0.14%	0.047%
74	17.221	2414	2420	2424	rBV2	3059657	4327439	17.56%	5.787%
75	17.262	2424	2427	2432	rVV	566933	750908	3.05%	1.004%
76	17.315	2432	2436	2445	rVB	579379	879494	3.57%	1.176%
77	17.392	2445	2449	2450	rBV	34730	38289	0.16%	0.051%
78	17.515	2463	2470	2476	rBV5	34566	81733	0.33%	0.109%
79	17.574	2476	2480	2482	rVV	32566	49547	0.20%	0.066%
80	17.615	2482	2487	2495	rVV	853694	1161234	4.71%	1.553%
81	17.709	2497	2503	2509	rBV2	113801	236095	0.96%	0.316%
82	17.768	2509	2513	2518	rVB	98663	145327	0.59%	0.194%
83	17.956	2541	2545	2549	rBV2	18680	26735	0.11%	0.036%
84	18.009	2549	2554	2557	rBV3	15488	29688	0.12%	0.040%
85	18.045	2557	2560	2565	rVB	48784	67458	0.27%	0.090%
86	18.133	2570	2575	2583	rVB2	182404	292261	1.19%	0.391%
87	18.204	2583	2587	2594	rVB	89487	128738	0.52%	0.172%
88	18.292	2594	2602	2605	rBV	65367	111469	0.45%	0.149%
89	18.327	2605	2608	2611	rVB2	21739	27842	0.11%	0.037%
90	18.427	2623	2625	2629	rVB	28706	34880	0.14%	0.047%
91	18.527	2637	2642	2648	rBV3	37361	60768	0.25%	0.081%
92	19.274	2764	2769	2774	rVB	132952	184649	0.75%	0.247%
93	19.603	2820	2825	2841	rBV3	645634	991420	4.02%	1.326%
94	19.986	2886	2890	2898	rBV2	101481	198415	0.81%	0.265%
95	20.068	2898	2904	2913	rBV	563745	839371	3.41%	1.123%
96	20.580	2987	2991	2995	rBV	52786	71353	0.29%	0.095%
97	20.703	3009	3012	3018	rVB3	69719	110179	0.45%	0.147%
98	21.392	3124	3129	3140	rBV	4453499	5338288	21.66%	7.139%
99	23.533	3488	3493	3499	rVB	420593	754031	3.06%	1.008%
100	23.680	3511	3518	3530	rVB	3025531	5641524	22.89%	7.545%

Sum of corrected areas: 74773208

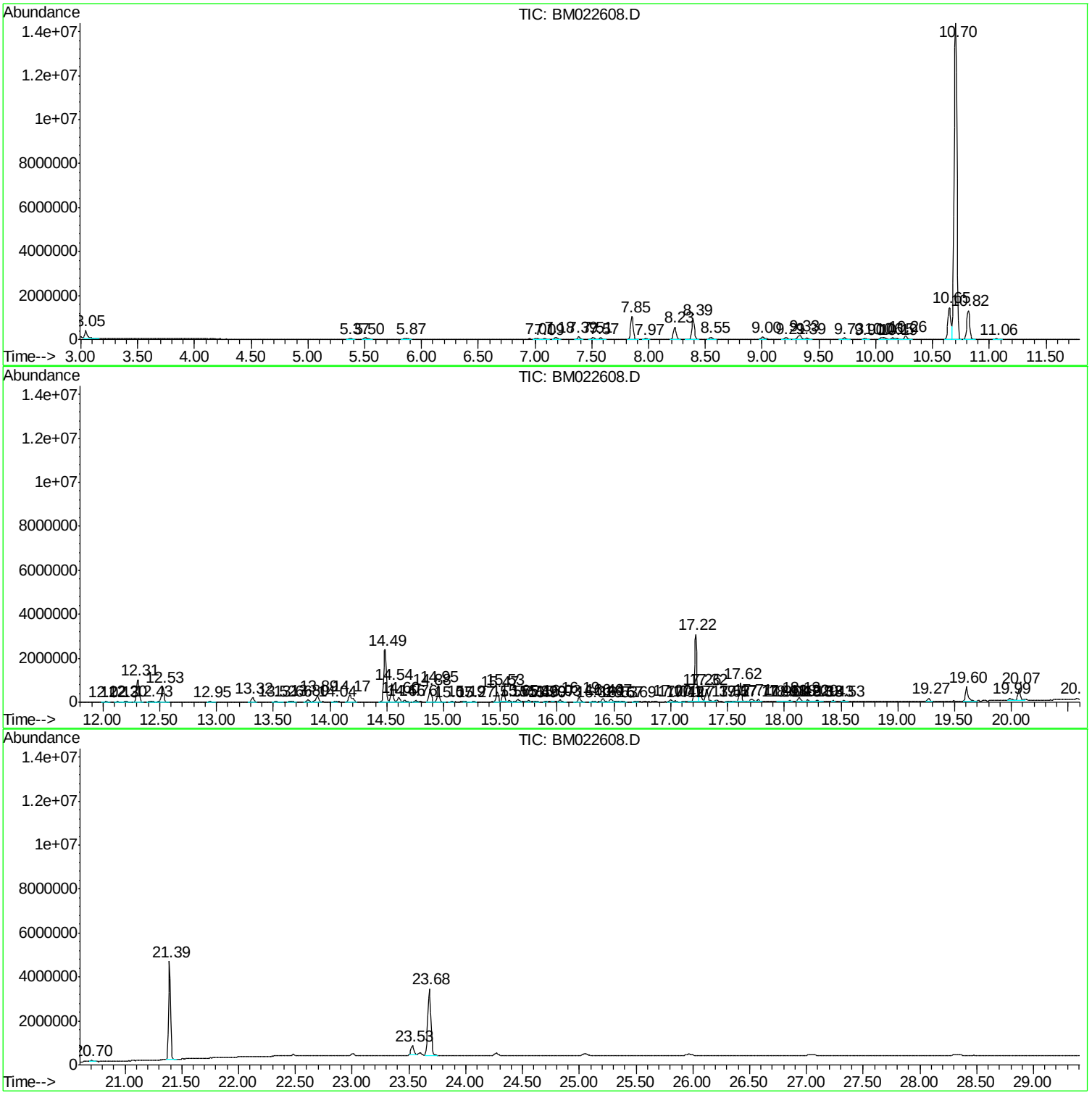
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090819\
Data File : BM022608.D
Acq On : 09 Sep 2019 17:55
Operator : HP/JU
Sample : K4706-07DL 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF573DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090819\
Data File : BM022608.D
Acq On : 09 Sep 2019 17:55
Operator : HP/JU
Sample : K4706-07DL 10X
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ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF573DL

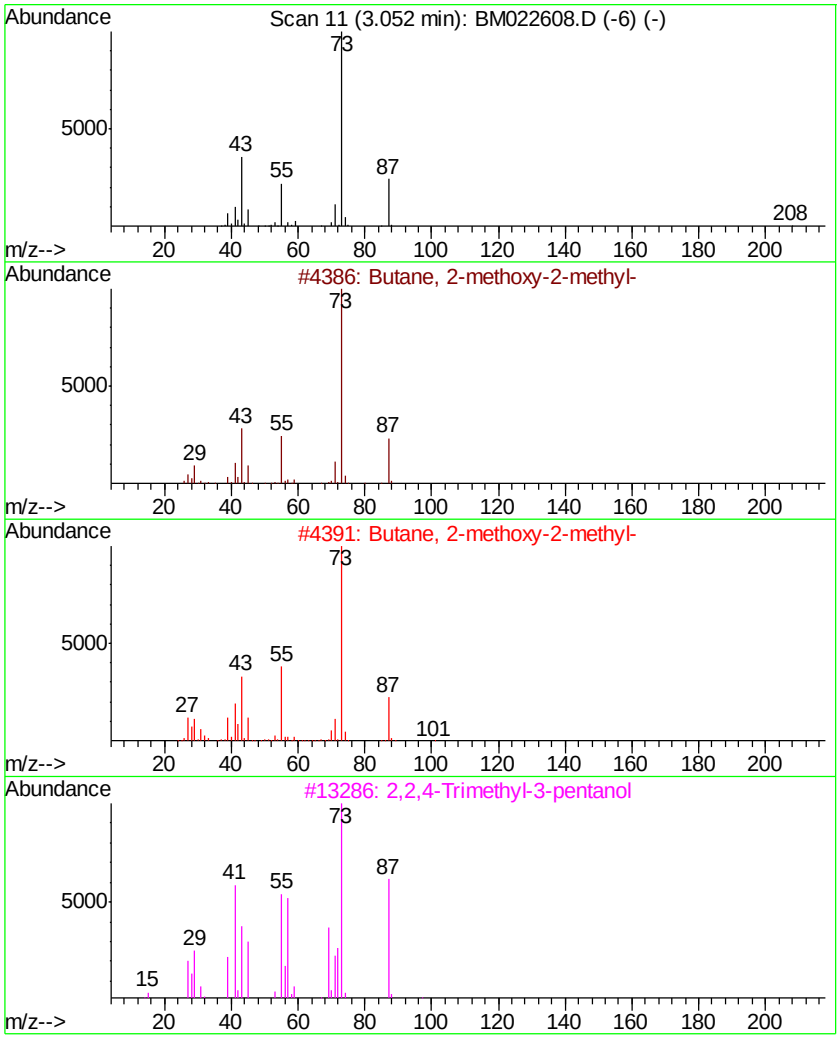
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

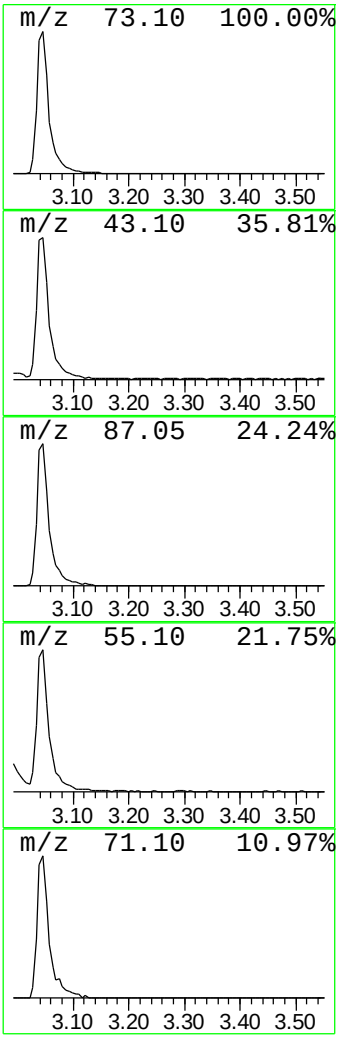
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.05	7.10 ng/ul	587334	1,4-Dichlorobenzene-d4	7.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090819\
Data File : BM022608.D
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Sample : K4706-07DL 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF573DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

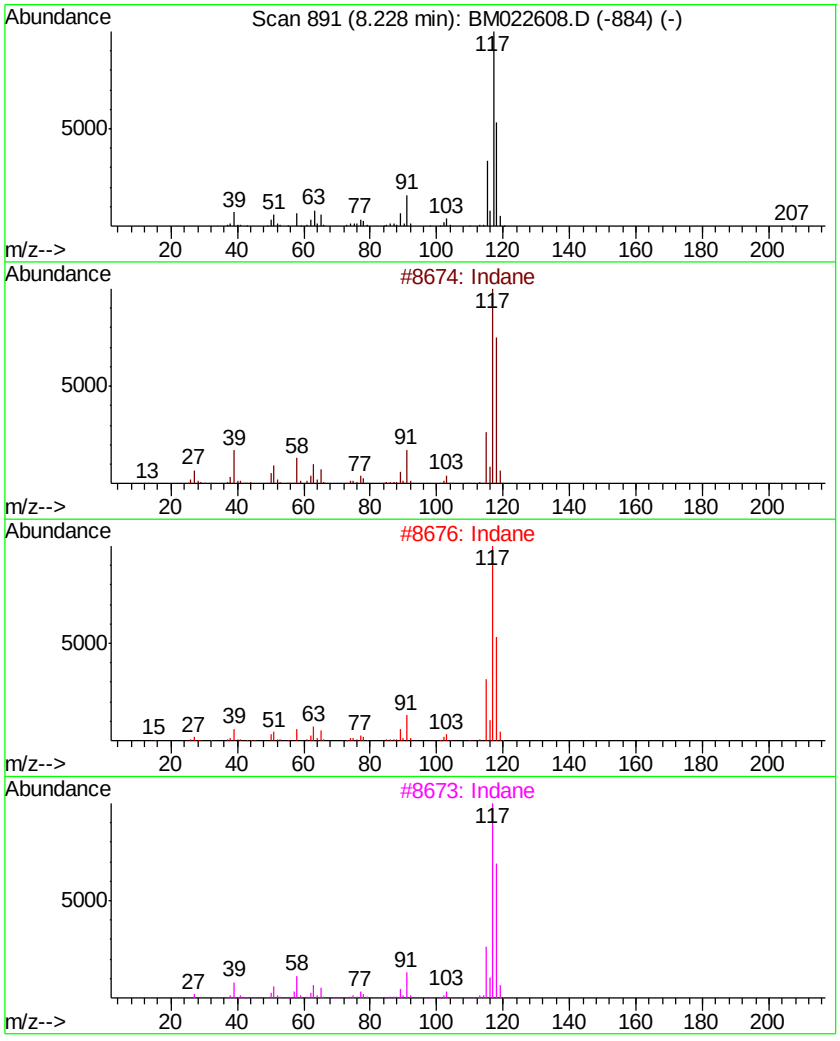
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

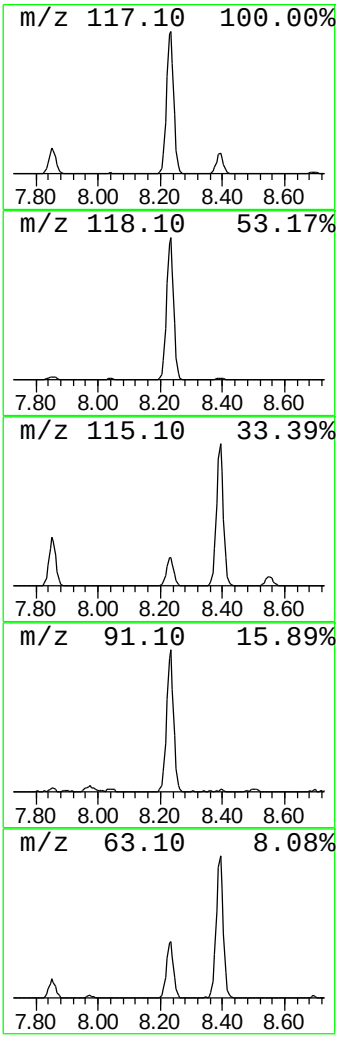
****

Peak Number 2 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.23	10.87 ng/ul	898906	1,4-Dichlorobenzene-d4	7.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Indane	118	C9H10	000496-11-7	87
3	Indane	118	C9H10	000496-11-7	83
4	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	60





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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

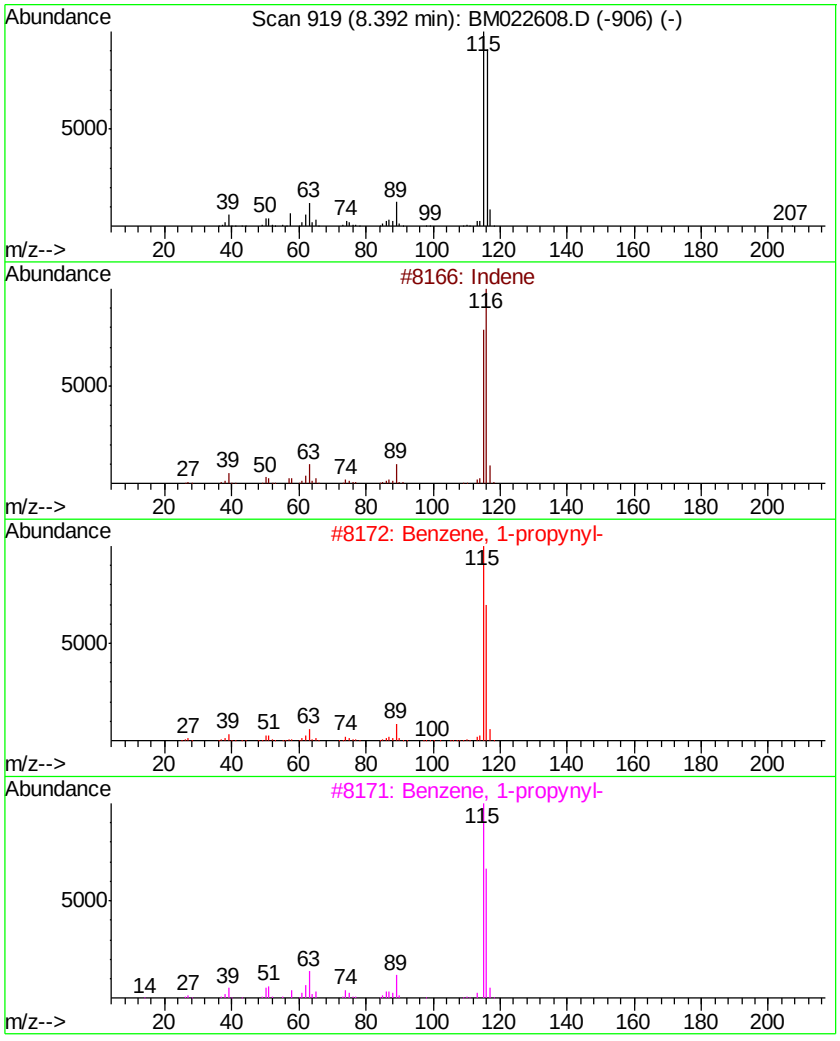
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

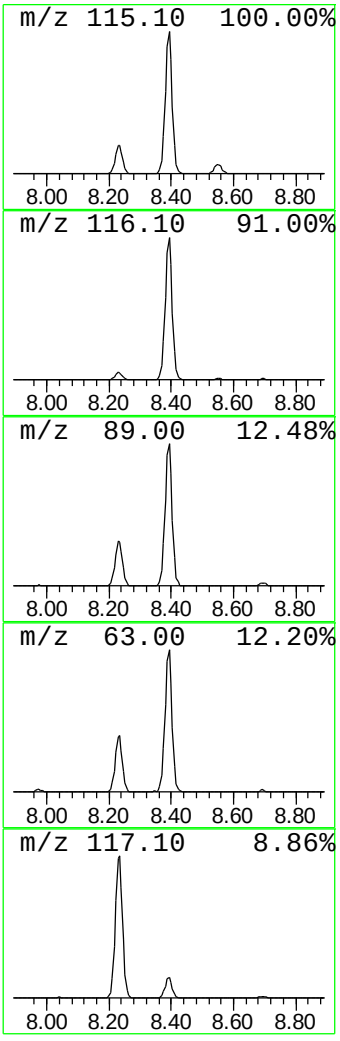
****

Peak Number 3 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.39	17.11 ng/ul	1414520	1,4-Dichlorobenzene-d4	7.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	97
2	Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
3	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
4	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91
5	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090819\
Data File : BM022608.D
Acq On : 09 Sep 2019 17:55
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Sample : K4706-07DL 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF573DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

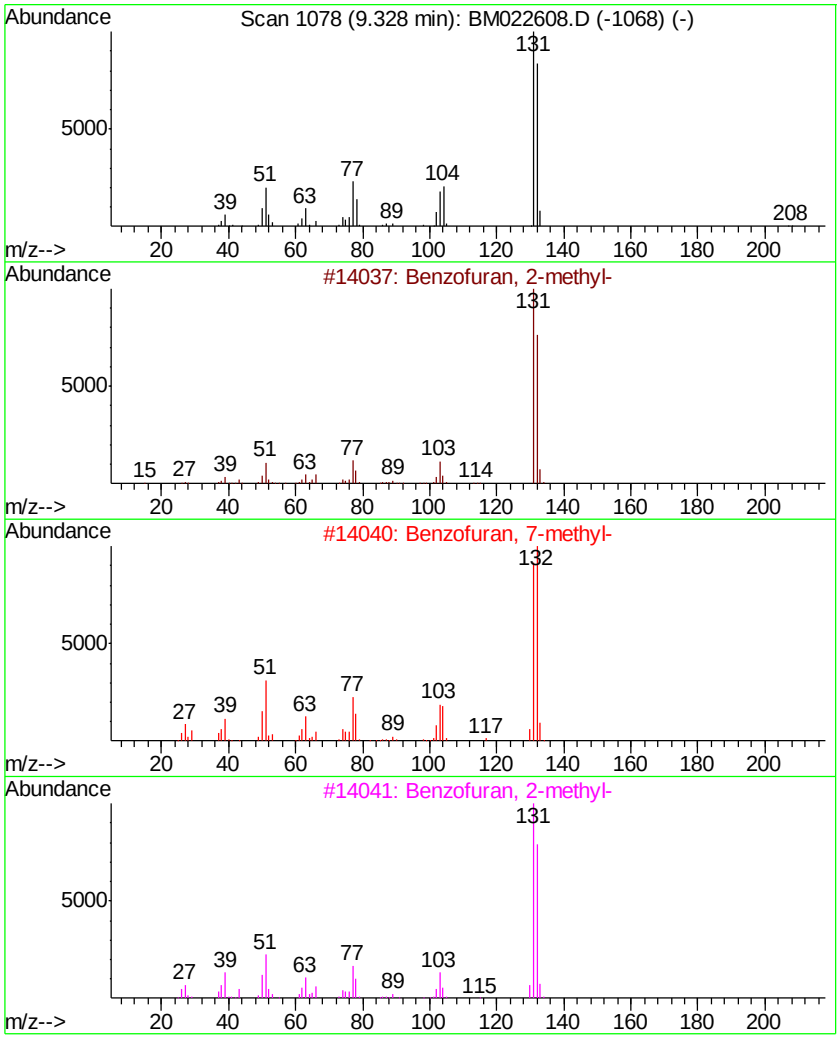
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

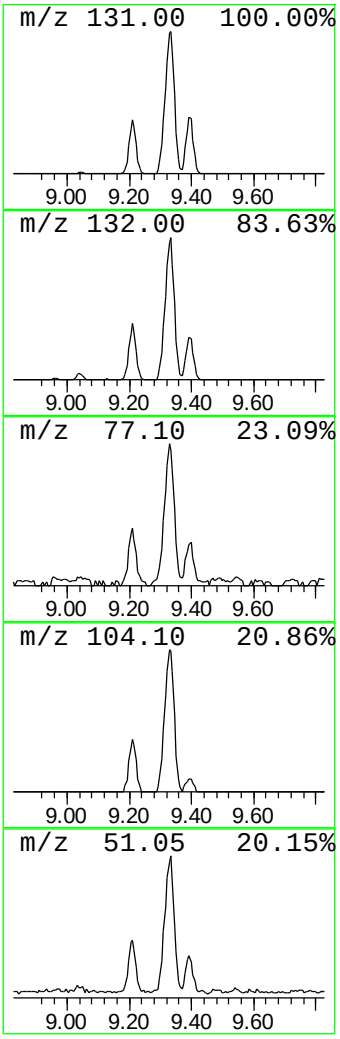
****

Peak Number 4 Benzofuran, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.33	3.00 ng/ul	386475	Naphthalene-d8	10.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	91
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
5	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	90





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090819\
Data File : BM022608.D
Acq On : 09 Sep 2019 17:55
Operator : HP/JU
Sample : K4706-07DL 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF573DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

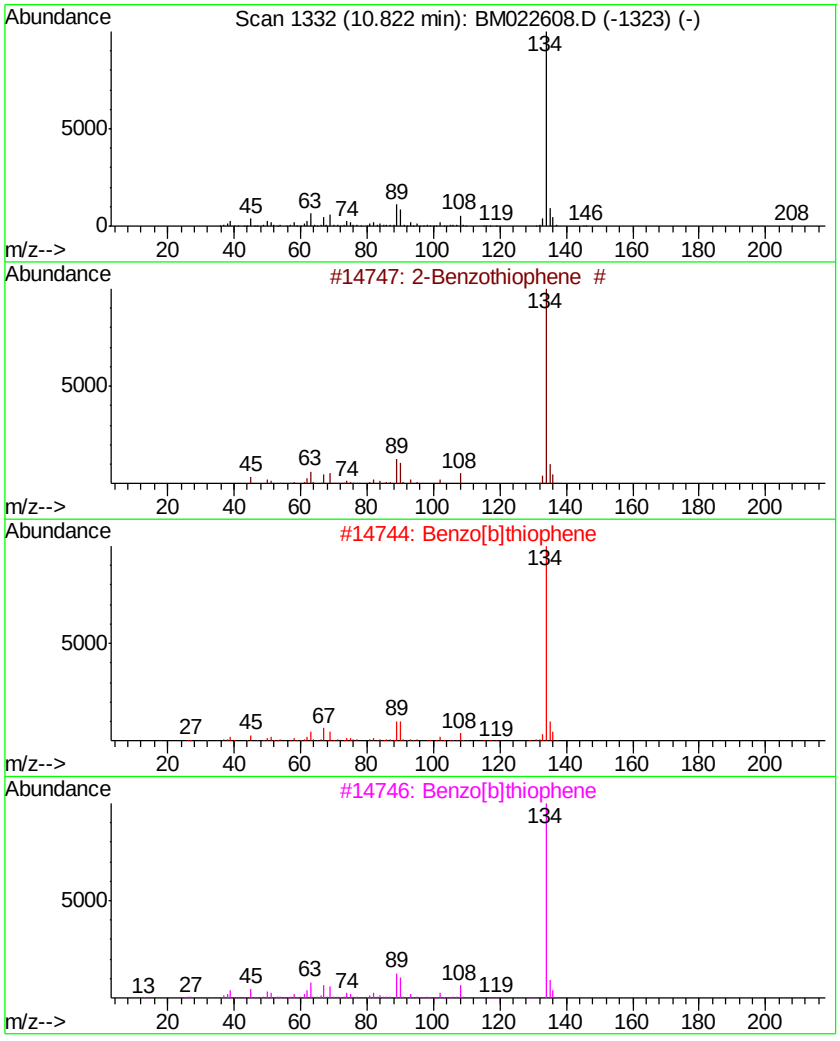
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

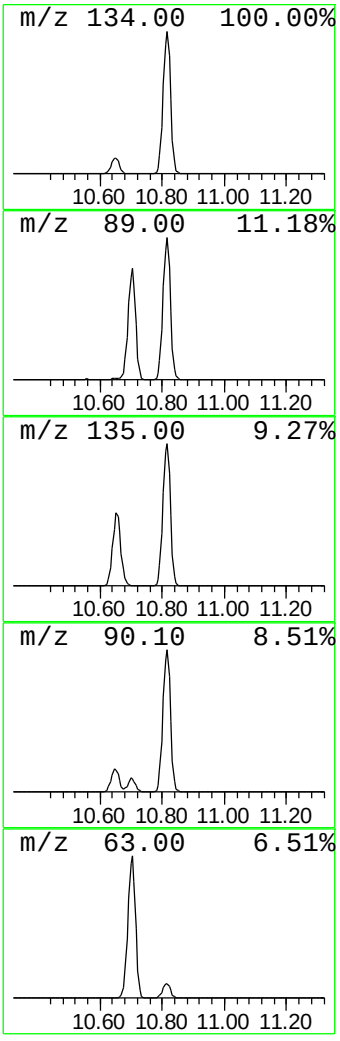
****

Peak Number 5 2-Benzothiophene # Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.82	17.35 ng/ul	2233370	Naphthalene-d8	10.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Benzothiophene #	134	C8H6S	000270-82-6	97
2	Benzo[b]thiophene	134	C8H6S	000095-15-8	94
3	Benzo[b]thiophene	134	C8H6S	000095-15-8	94
4	Benzo[b]thiophene	134	C8H6S	000095-15-8	94
5	Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	94





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Data File : BM022608.D
Acq On : 09 Sep 2019 17:55
Operator : HP/JU
Sample : K4706-07DL 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF573DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

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

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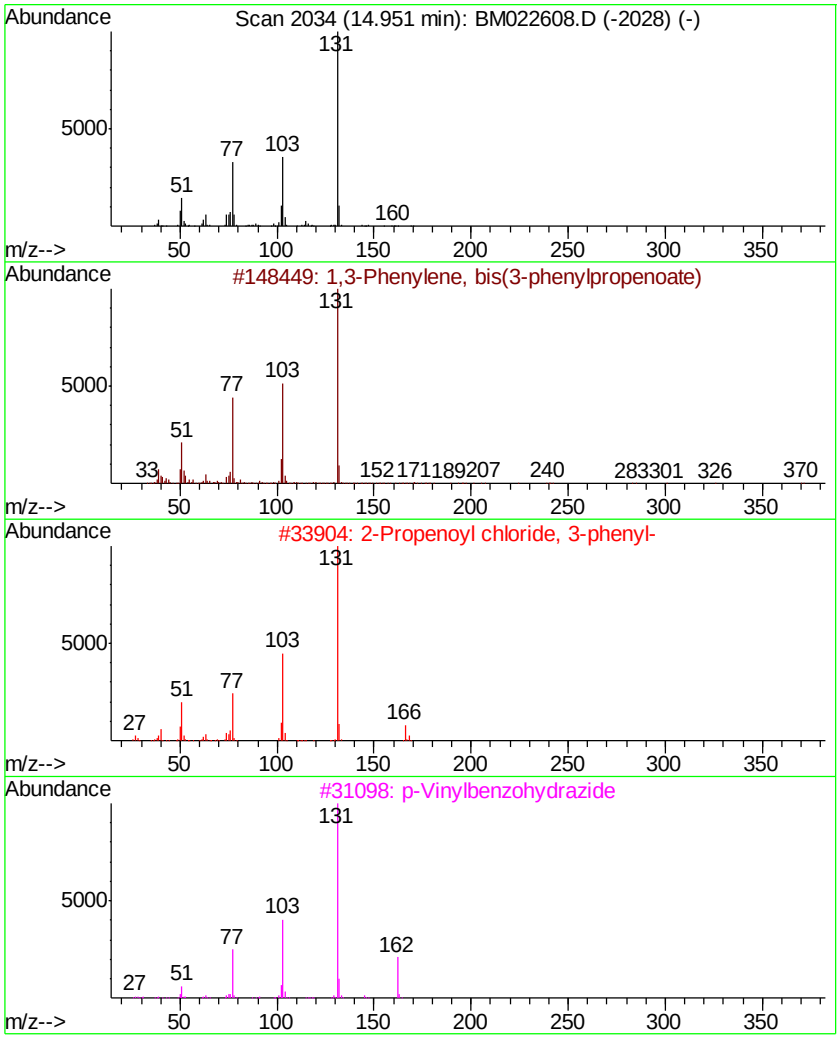
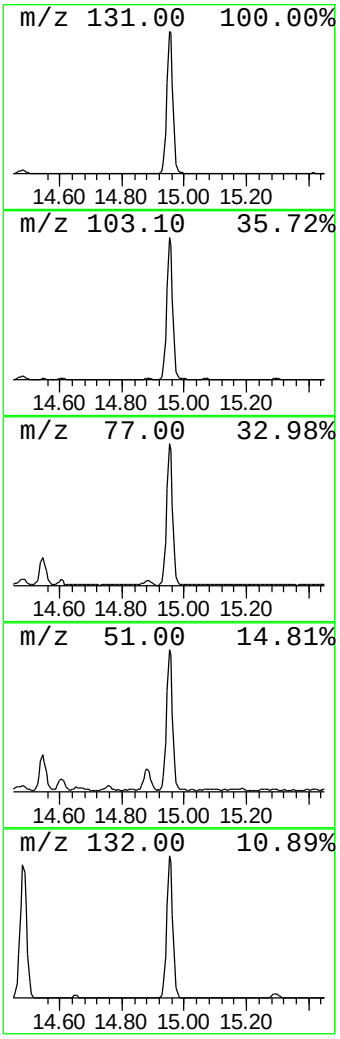
Peak Number 6 1,3-Phenylene, bis(3-phenyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.95	5.64 ng/ul	1027100	Acenaphthene-d10	14.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Phenylene, bis(3-phenylprope...	370	C24H18O4	1000259-62-4	83
2	2-Propenoyl chloride, 3-phenyl-	166	C9H7ClO	000102-92-1	78
3	p-Vinylbenzohydrazide	162	C9H10N2O	1000244-74-9	78
4	Propanedioic acid, nitrile, hydr...	229	C12H11N3O2	294878-35-6	78
5	Vinyl trans-cinnamate	174	C11H10O2	017719-70-9	78





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Data File : BM022608.D
Acq On : 09 Sep 2019 17:55
Operator : HP/JU
Sample : K4706-07DL 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF573DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

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

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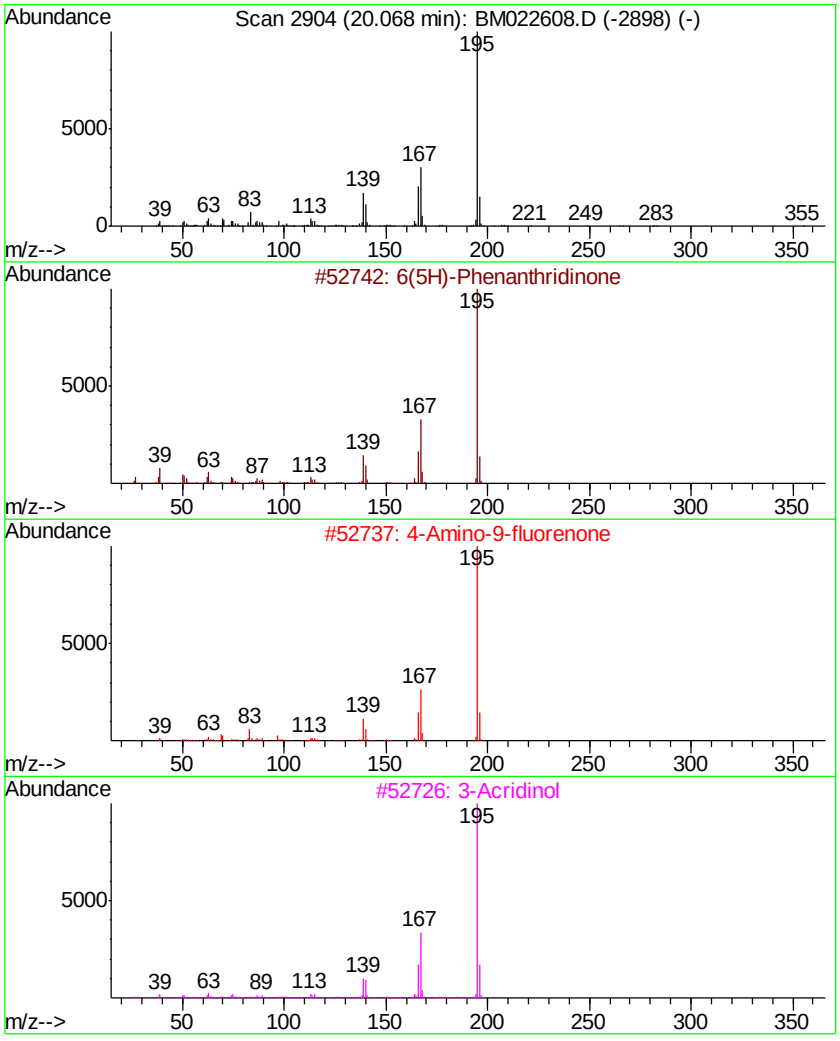
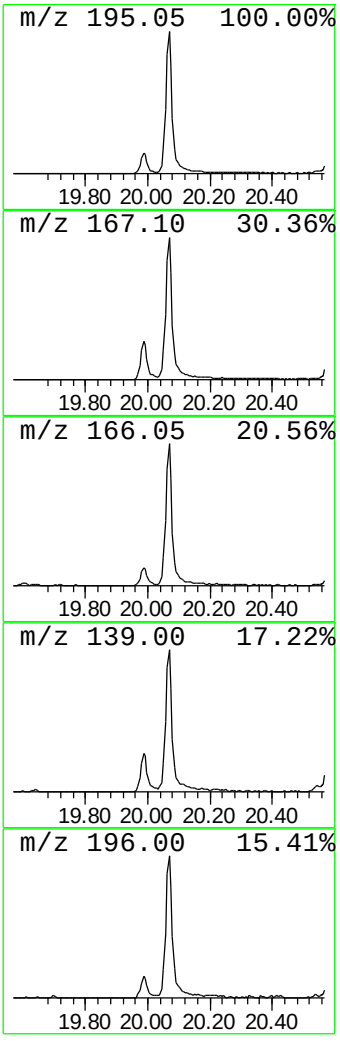
Peak Number 7 6(5H)-Phenanthridinone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.07	3.14 ng/ul	839371	Chrysene-d12	21.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6(5H)-Phenanthridinone	195	C13H9NO	001015-89-0	96
2	4-Amino-9-fluorenone	195	C13H9NO	004269-15-2	96
3	3-Acridinol	195	C13H9NO	007132-70-9	94
4	4-Amino-9-fluorenone	195	C13H9NO	004269-15-2	94
5	9(10H)-Acridinone	195	C13H9NO	000578-95-0	94





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090819\
Data File : BM022608.D
Acq On : 09 Sep 2019 17:55
Operator : HP/JU
Sample : K4706-07DL 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF573DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

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

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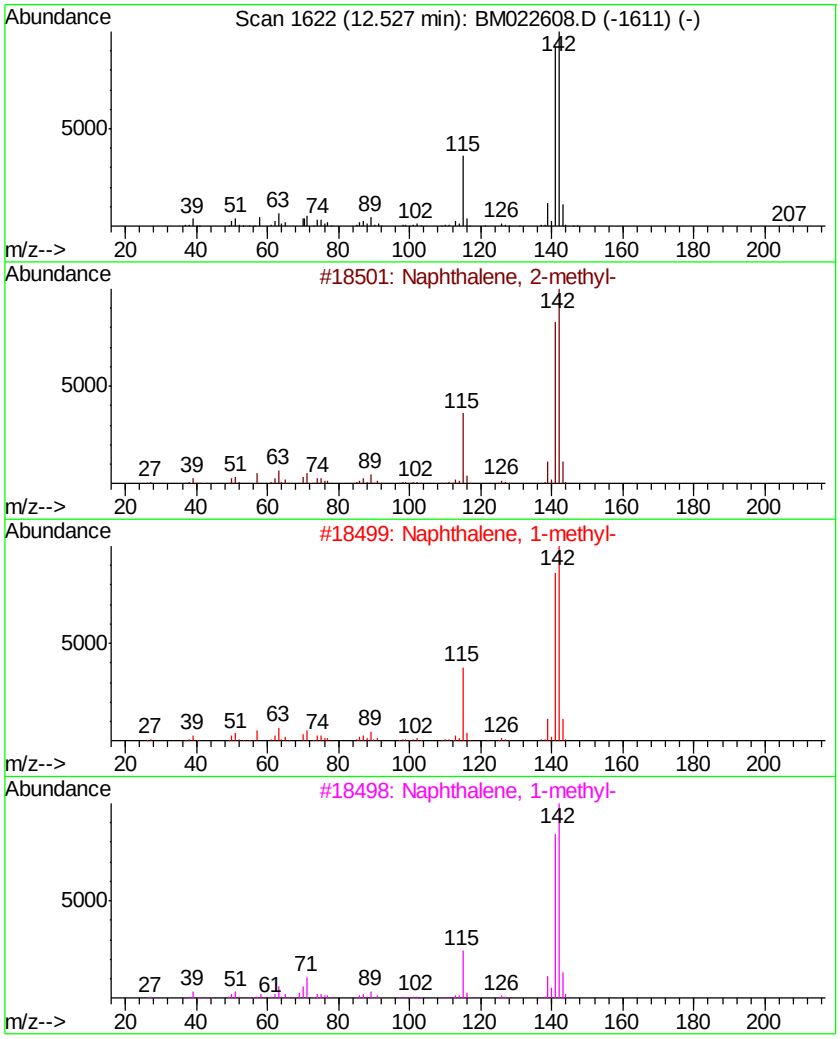
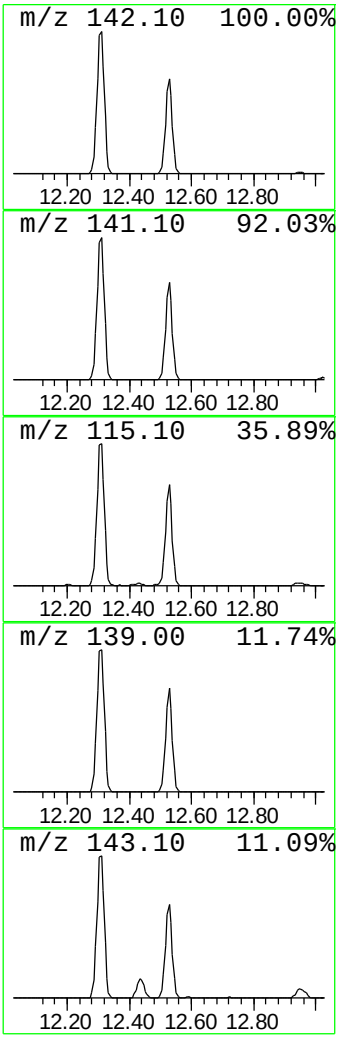
Peak Number 8 Naphthalene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.53	8.55 ng/ul	1100610	Naphthalene-d8	10.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
4	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
5	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM090819\
Data File : BM022608.D
Acq On : 09 Sep 2019 17:55
Operator : HP/JU
Sample : K4706-07DL 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF573DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methoxy...	3.05	7.1	ng/ul	587334	1	7.85	1653330	20.0
Indane	8.23	10.9	ng/ul	898906	1	7.85	1653330	20.0
Indene	8.39	17.1	ng/ul	1414520	1	7.85	1653330	20.0
Benzofuran, 2-met...	9.33	3.0	ng/ul	386475	2	10.65	2574820	20.0
2-Benzothiophene #	10.82	17.4	ng/ul	2233370	2	10.65	2574820	20.0
1,3-Phenylene, bi...	14.95	5.6	ng/ul	1027100	3	14.49	3643000	20.0
6(5H)-Phenanthrid...	20.07	3.1	ng/ul	839371	5	21.39	5338290	20.0
Naphthalene, 1-me...	12.53	8.6	ng/ul	1100610	2	10.65	2574820	20.0