LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121620\
Data File : BP004319.D
Acq On : 16 Dec 2020 12:08
Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.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	5.893	455	460	462	rBV	171015	261456	3.03%	0.336%
2	5.922	462	465	474	rVB	178038	265951	3.08%	0.342%
3	6.487	557	561	565	rVB3	36564	51976	0.60%	0.067%
4	6.834	617	620	625	rVB3	19993	28250	0.33%	0.036%
5	6.999	643	648	651	rBV	200334	344497	3.99%	0.443%
6	7.169	671	677	686	rBV	739433	1214124	14.07%	1.561%
7	7.369	705	711	718	rBV	827762	1344329	15.58%	1.729%
8	7.434	718	722	731	rVB	39290	63771	0.74%	0.082%
9	7.611	750	752	760	rVB4	11866	22519	0.26%	0.029%
10	7.840	784	791	796	rBV	840469	1367037	15.84%	1.758%
11	7.899	796	801	809	rVB2	93335	177377	2.06%	0.228%
12	7.987	813	816	820	rVB4	14042	19051	0.22%	0.025%
13	8.075	825	831	836	rBV	78562	139969	1.62%	0.180%
14	8.446	890	894	899	rBV2	34748	54839	0.64%	0.071%
15	8.540	903	910	920	rVV	617369	1031740	11.96%	1.327%
16	8.652	920	929	935	rVB5	43128	131245	1.52%	0.169%
17	8.928	970	976	981	rBV	347867	565770	6.56%	0.728%
18	8.993	981	987	998	rVV	975826	1645644	19.07%	2.116%
19	9.087	998	1003	1004	rVV4	19806	29872	0.35%	0.038%
20	9.110	1004	1007	1010	rVV	43370	65527	0.76%	0.084%
21	9.158	1010	1015	1025	rVB	429158	682728	7.91%	0.878%
22	9.381	1049	1053	1056	rBV2	7946	13208	0.15%	0.017%
23	9.422	1057	1060	1065	rVB4	15534	24775	0.29%	0.032%
24	9.552	1077	1082	1090	rVB3	41905	70216	0.81%	0.090%
25	9.716	1102	1110	1124	rBV	810992	1391534	16.13%	1.790%
26	9.887	1130	1139	1145	rVV3	137603	337510	3.91%	0.434%
27	9.946	1145	1149	1160	rVB9	24724	56277	0.65%	0.072%
28	10.052	1162	1167	1171	rBV8	7764	14480	0.17%	0.019%
29	10.110	1171	1177	1187	rVB2	71069	180122	2.09%	0.232%
30	10.252	1194	1201	1222	rBV	1189044	2071346	24.01%	2.664%
31	10.416	1223	1229	1237	rVV3	41112	78938	0.91%	0.102%
32	10.493	1237	1242	1245	rVV6	11443	21975	0.25%	0.028%
33	10.540	1245	1250	1258	rVB2	55096	101815	1.18%	0.131%
34	10.634	1258	1266	1271	rBV	1185536	2114216	24.51%	2.719%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121620\
Data File : BP004319.D
Acq On : 16 Dec 2020 12:08
Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Title : SVOA CALIBRATION

35	10.669	1271	1272	1281	rVV	279005	354328	4.11%	0.456%
36	10.769	1282	1289	1298	rVB	88421	161168	1.87%	0.207%
37	10.928	1308	1316	1323	rBV4	44209	93626	1.09%	0.120%
38	10.993	1323	1327	1331	rBV4	8914	18255	0.21%	0.023%
39	11.169	1350	1357	1363	rBV3	22025	37954	0.44%	0.049%
40	11.228	1363	1367	1372	rVV2	10762	18117	0.21%	0.023%
41	11.281	1372	1376	1385	rVB	25325	50123	0.58%	0.064%
42	11.381	1388	1393	1395	rBV5	7302	13214	0.15%	0.017%
43	11.422	1396	1400	1410	rVV7	16761	35185	0.41%	0.045%
44	11.804	1455	1465	1470	rBV3	26159	50353	0.58%	0.065%
45	11.863	1470	1475	1482	rVV3	16204	31269	0.36%	0.040%
46	11.946	1484	1489	1499	rVB5	11908	35519	0.41%	0.046%
47	12.122	1515	1519	1525	rVV4	9536	15051	0.17%	0.019%
48	12.228	1528	1537	1542	rBV	25250	45125	0.52%	0.058%
49	12.451	1569	1575	1584	rVB2	38450	69812	0.81%	0.090%
50	12.699	1614	1617	1620	rVV	20787	29108	0.34%	0.037%
51	12.740	1620	1624	1633	rVV	65953	105971	1.23%	0.136%
52	12.840	1635	1641	1652	rVV	297705	513116	5.95%	0.660%
53	12.934	1652	1657	1661	rVB	26908	41106	0.48%	0.053%
54	13.087	1676	1683	1695	rBV	441426	656114	7.60%	0.844%
55	13.316	1718	1722	1726	rBV3	15915	21382	0.25%	0.027%
56	13.381	1726	1733	1742	rBV3	1180501	2046402	23.72%	2.632%
57	13.581	1761	1767	1782	rVV	108030	209858	2.43%	0.270%
58	13.710	1784	1789	1791	rVV5	11138	17339	0.20%	0.022%
59	13.746	1791	1795	1800	rVV	37686	56033	0.65%	0.072%
60	13.893	1813	1820	1824	rVV	2337535	3360710	38.95%	4.322%
61	13.934	1824	1827	1837	rVV	207805	292746	3.39%	0.376%
62	14.016	1837	1841	1848	rVV2	52485	74455	0.86%	0.096%
63	14.175	1854	1868	1882	rBV	2821069	4303055	49.88%	5.534%
64	14.340	1889	1896	1901	rBV	58806	92835	1.08%	0.119%
65	14.487	1914	1921	1930	rVV2	1790226	2736540	31.72%	3.519%
66	14.569	1930	1935	1939	rVB6	13502	20382	0.24%	0.026%
67	14.681	1949	1954	1961	rBV	231682	362513	4.20%	0.466%
68	14.740	1961	1964	1971	rVB	19938	29797	0.35%	0.038%
69	14.904	1985	1992	1997	rBV	9245	19092	0.22%	0.025%
70	15.081	2018	2022	2025	rBV3	15760	20076	0.23%	0.026%
71	15.187	2032	2040	2043	rBV6	11661	21930	0.25%	0.028%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121620\
Data File : BP004319.D
Acq On : 16 Dec 2020 12:08
Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Title : SVOA CALIBRATION

72	15.228	2043	2047	2053	rVB	26481	36920	0.43%	0.047%
73	15.304	2053	2060	2065	rBV2	123592	193742	2.25%	0.249%
74	15.410	2072	2078	2081	rBV4	9945	14889	0.17%	0.019%
75	15.481	2083	2090	2101	rBV	3566064	5342344	61.92%	6.871%
76	15.610	2106	2112	2126	rVV2	1699702	2620189	30.37%	3.370%
77	15.722	2126	2131	2138	rVV2	49829	93775	1.09%	0.121%
78	15.851	2147	2153	2160	rVB	40750	65000	0.75%	0.084%
79	16.063	2181	2189	2194	rBV10	16594	32249	0.37%	0.041%
80	16.181	2202	2209	2213	rBV4	15774	29131	0.34%	0.037%
81	16.269	2218	2224	2231	rVB6	18493	35418	0.41%	0.046%
82	17.034	2345	2354	2361	rBV5	12854	33364	0.39%	0.043%
83	17.245	2383	2390	2397	rBV	2194600	3245196	37.61%	4.174%
84	17.345	2400	2407	2419	rBV2	4462459	6557426	76.01%	8.433%
85	17.481	2426	2430	2436	rBV9	12284	26575	0.31%	0.034%
86	17.557	2440	2443	2450	rVV	79479	131723	1.53%	0.169%
87	17.645	2454	2458	2467	rVB10	16478	34697	0.40%	0.045%
88	17.916	2498	2504	2518	rBV	1229136	1526944	17.70%	1.964%
89	18.139	2538	2542	2547	rBV4	26914	42065	0.49%	0.054%
90	18.204	2549	2553	2558	rVB	23512	32782	0.38%	0.042%
91	18.339	2572	2576	2581	rBV4	14807	27450	0.32%	0.035%
92	18.975	2679	2684	2692	rBV	171174	302826	3.51%	0.389%
93	19.233	2724	2728	2731	rBV	61122	80546	0.93%	0.104%
94	19.281	2732	2736	2741	rVB	268412	303018	3.51%	0.390%
95	19.475	2766	2769	2779	rBV	51900	94968	1.10%	0.122%
96	19.586	2782	2788	2798	rBV	5950709	8273480	95.90%	10.640%
97	21.057	3035	3038	3045	rBV4	113568	173045	2.01%	0.223%
98	21.339	3081	3086	3093	rBV	2960953	3811622	44.18%	4.902%
99	23.557	3456	3463	3479	rVB	4135440	8627565	100.00%	11.096%
100	23.710	3483	3489	3498	rVB	1988679	3921925	45.46%	5.044%

Sum of corrected areas: 77754617

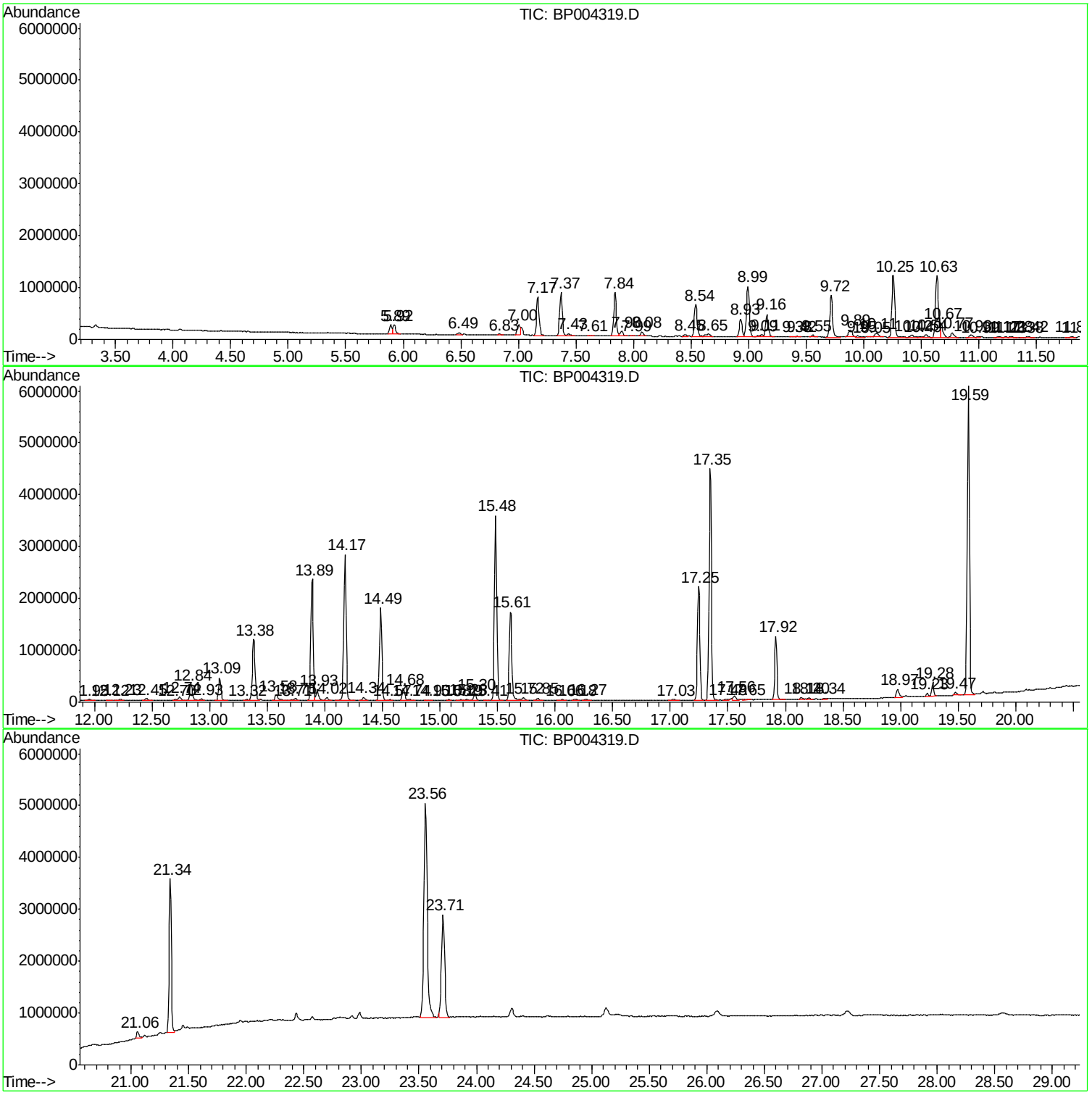
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121620\
Data File : BP004319.D
Acq On : 16 Dec 2020 12:08
Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.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 P\DATA\BP121620\
Data File : BP004319.D
Acq On : 16 Dec 2020 12:08
Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

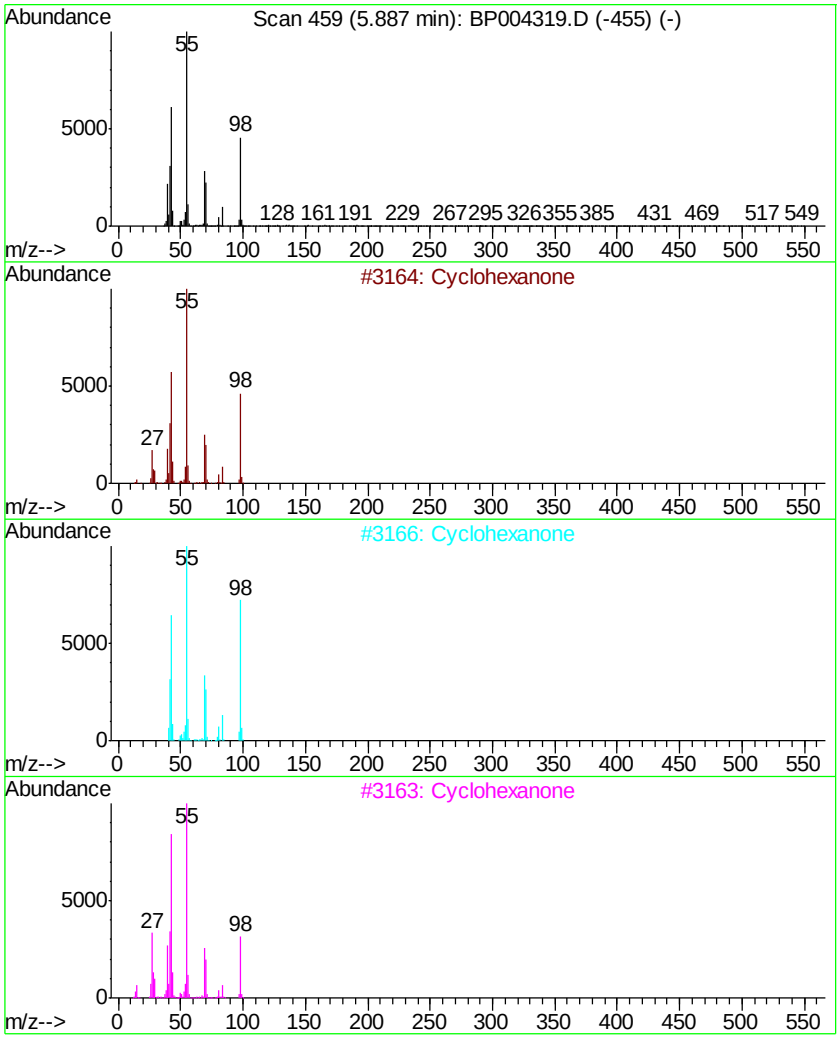
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

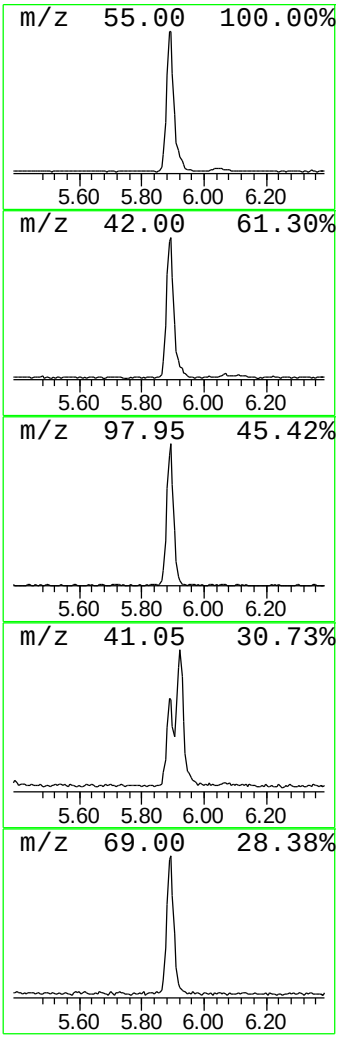
****

Peak Number 1 Cyclohexanone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.89	3.83 ng/ul	261456	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone	98	C6H10O	000108-94-1	94
2	Cyclohexanone	98	C6H10O	000108-94-1	93
3	Cyclohexanone	98	C6H10O	000108-94-1	91
4	Cyclohexanone	98	C6H10O	000108-94-1	83
5	Cyclopentanone, 2-methyl-	98	C6H10O	001120-72-5	50





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121620\
Data File : BP004319.D
Acq On : 16 Dec 2020 12:08
Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

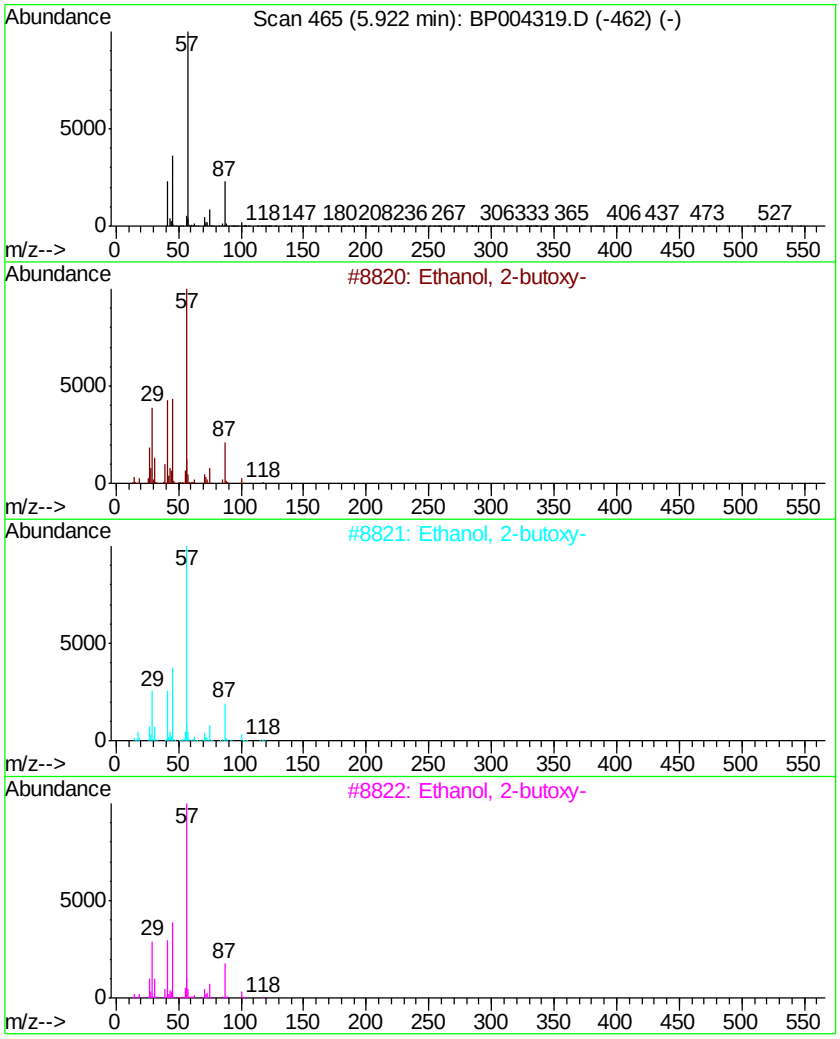
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

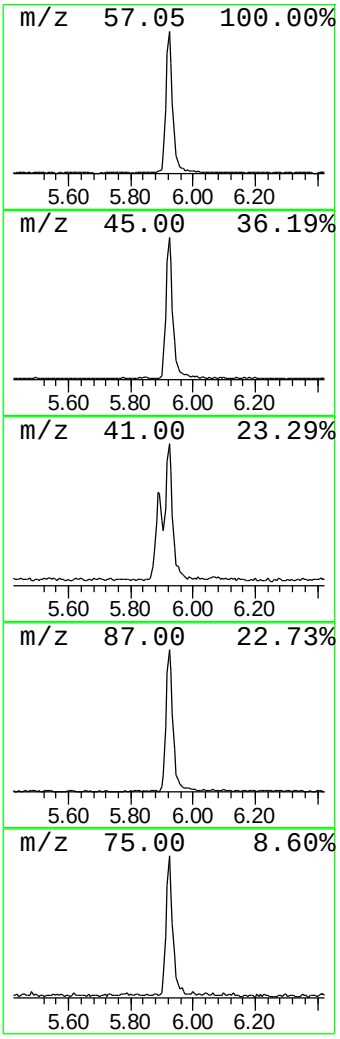
****

Peak Number 2 Ethanol, 2-butoxy- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.92	3.89 ng/ul	265951	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	72
2	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	43
3	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	43
4	3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	33
5	1,3-Propanediol, 2-methyl-, dipr...	202	C10H18O4	054932-83-1	12





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121620\
Data File : BP004319.D
Acq On : 16 Dec 2020 12:08
Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

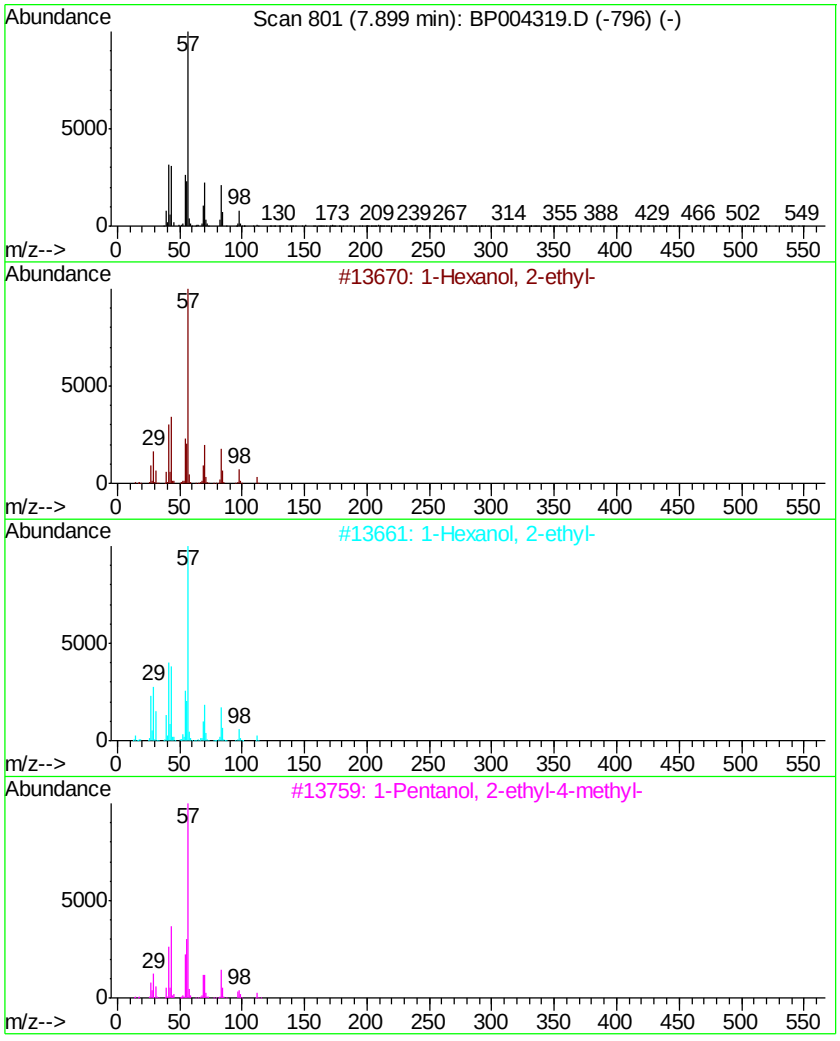
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

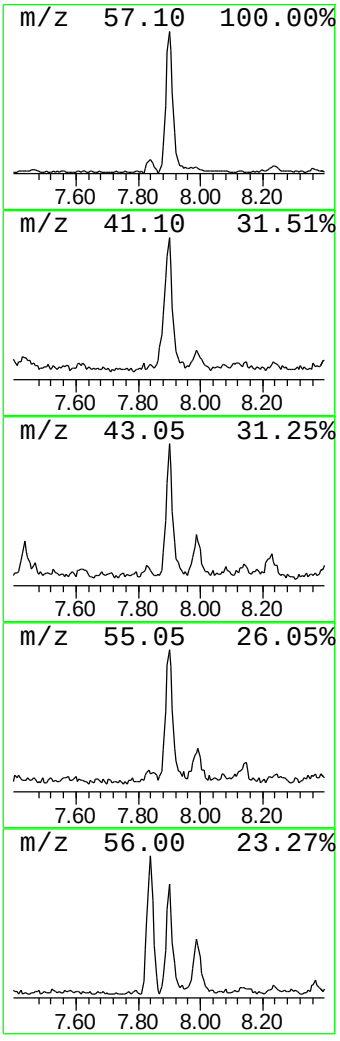
****

Peak Number 3 1-Hexanol, 2-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.90	2.60 ng/ul	177377	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
2	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
3	1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	64
4	Carbonic acid, isobutyl 2-ethylh...	230	C13H26O3	1000357-82-9	53
5	Heptyl isobutyl carbonate	216	C12H24O3	959068-08-7	53





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121620\
Data File : BP004319.D
Acq On : 16 Dec 2020 12:08
Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

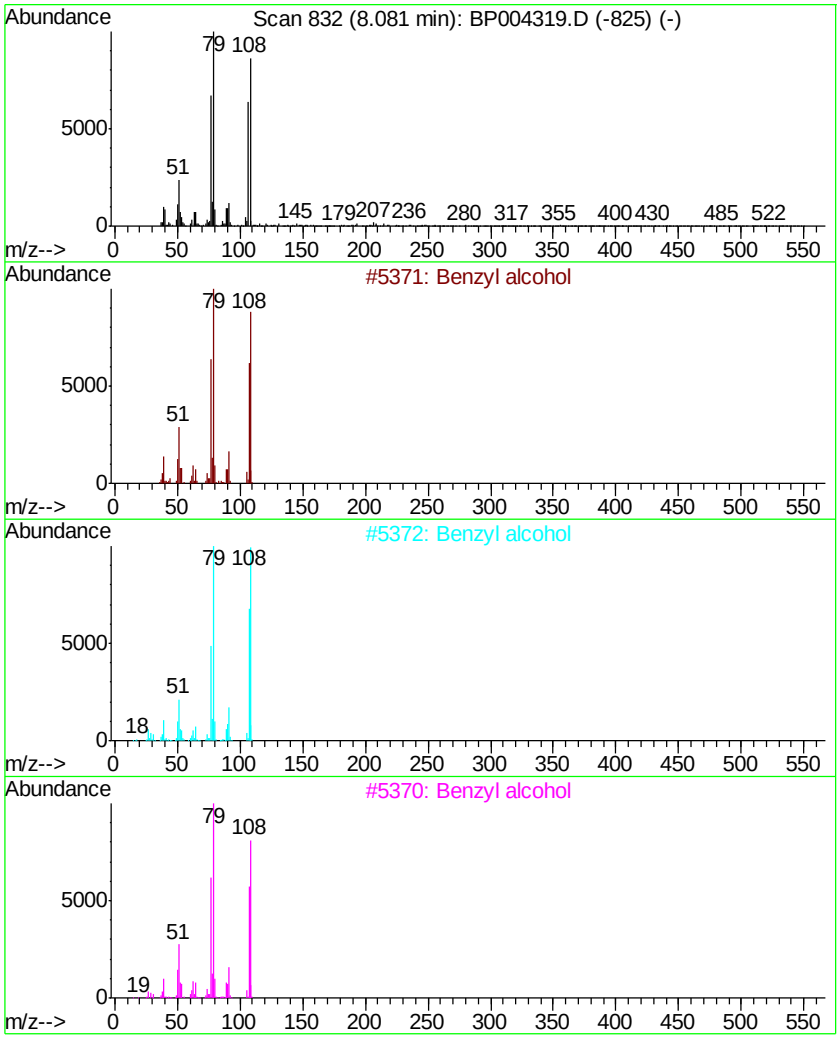
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

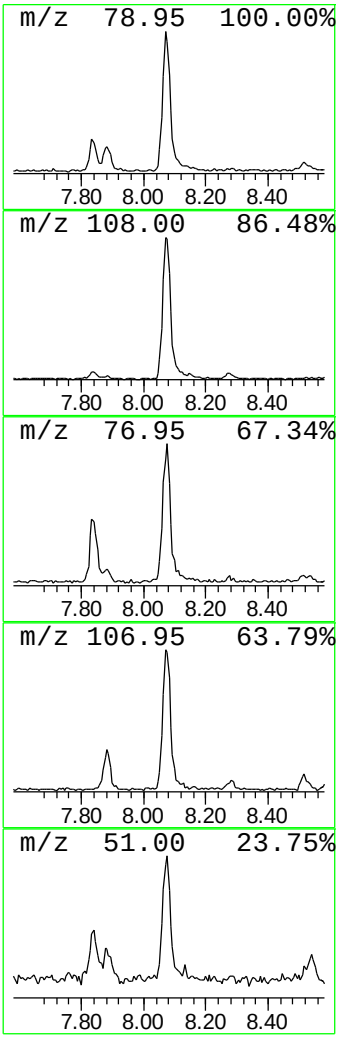
****

Peak Number 4 Benzyl alcohol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.08	2.05 ng/ul	139969	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzyl alcohol	108	C7H8O	000100-51-6	94
2	Benzyl alcohol	108	C7H8O	000100-51-6	94
3	Benzyl alcohol	108	C7H8O	000100-51-6	94
4	Benzyl alcohol	108	C7H8O	000100-51-6	94
5	N-Cbz-glycylglycine p-nitropheny...	387	C18H17N3O7	013574-81-7	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121620\
Data File : BP004319.D
Acq On : 16 Dec 2020 12:08
Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

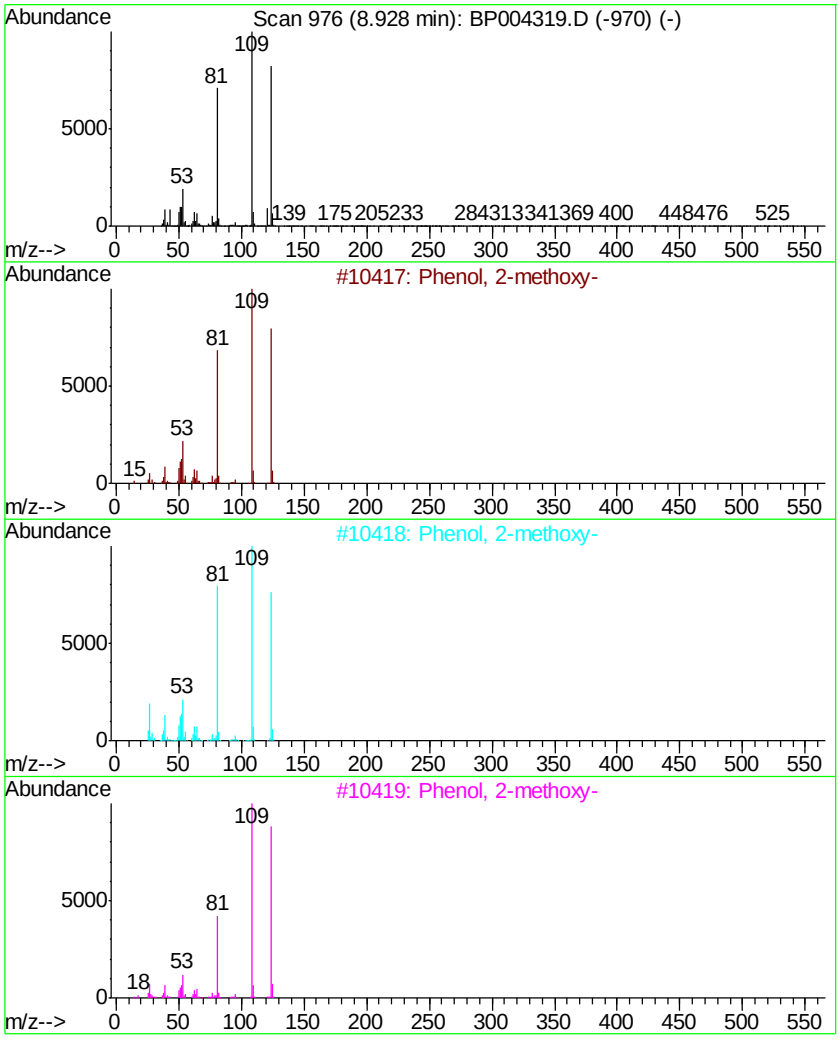
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

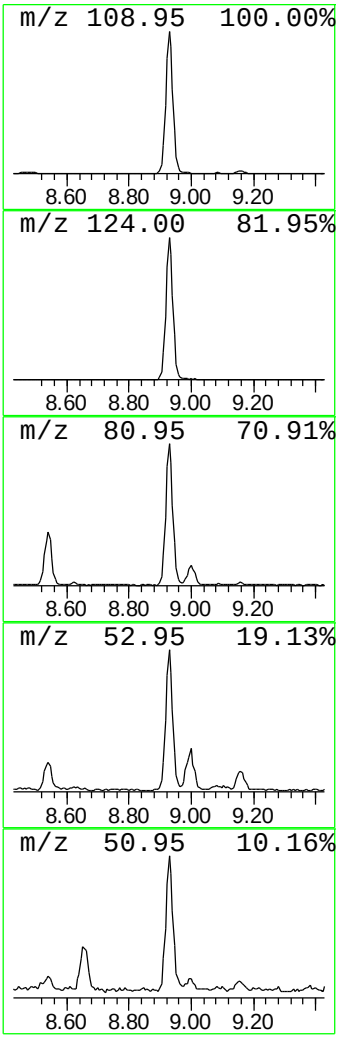
****

Peak Number 5 Phenol, 2-methoxy- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.93	8.28 ng/ul	565770	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	96
2	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	90
3	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	90
4	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	87
5	2-Acetyl-5-methylfuran	124	C7H8O2	001193-79-9	86





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Data File : BP004319.D
Acq On : 16 Dec 2020 12:08
Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

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

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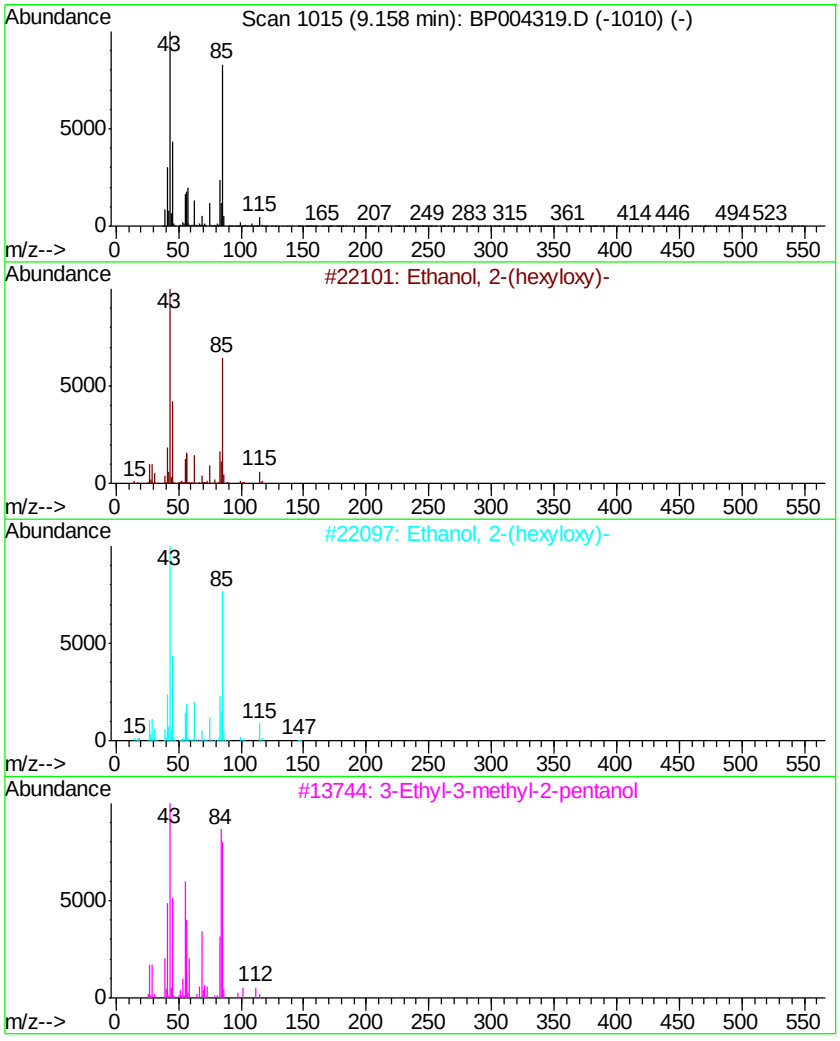
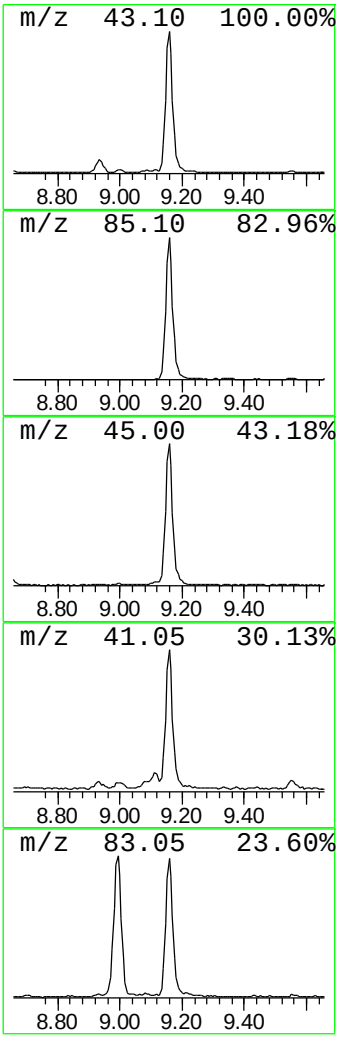
Peak Number 6 Ethanol, 2-(hexyloxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.16	9.99 ng/ul	682728	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	86
2	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	83
3	3-Ethyl-3-methyl-2-pentanol	130	C8H18O	066576-22-5	56
4	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	53
5	Hexane, 1-(hexyloxy)-2-methyl-	200	C13H28O	074421-17-3	50





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Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

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

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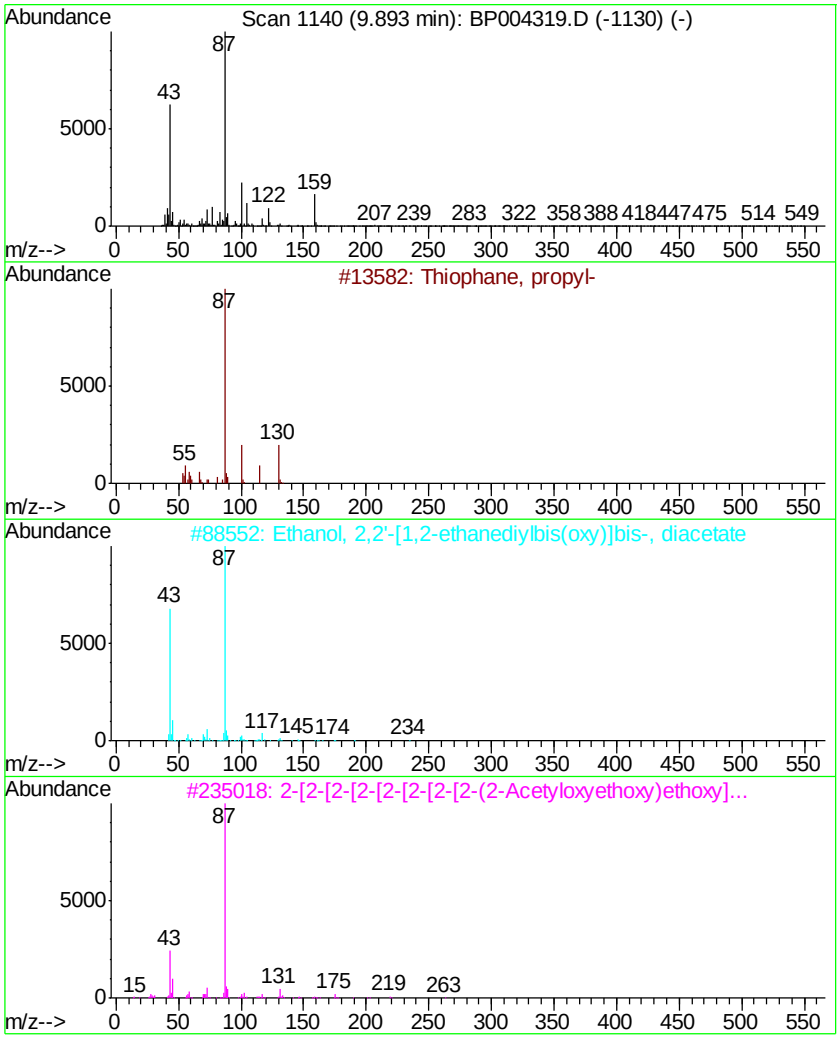
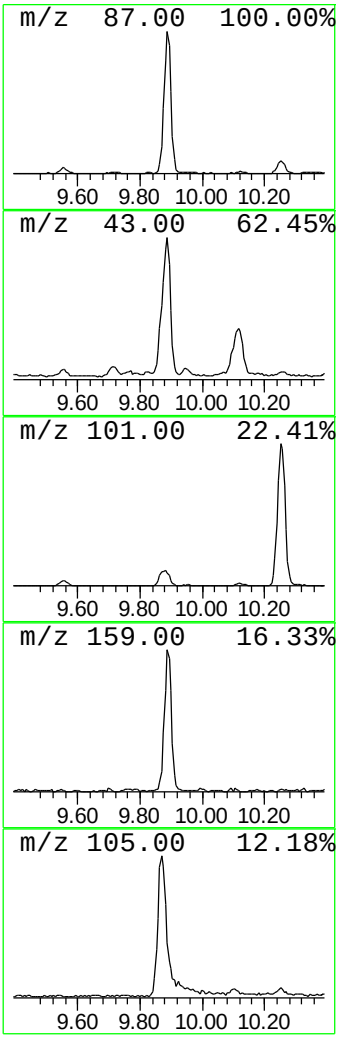
Peak Number 7 Thiophane, propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.89	3.19 ng/ul	337510	Naphthalene-d8	10.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Thiophane, propyl-	130	C7H14S	001551-34-4	50
2	Ethanol, 2,2'-[1,2-ethanediylbis...	234	C10H18O6	000111-21-7	47
3	2-[2-[2-[2-[2-[2-[2-(2-Acetyl...	498	C22H42O12	1000351-90-5	47
4	2-[2-[2-[2-[2-[2-(2-Acetyloxy...	454	C20H38O11	1000351-90-6	47
5	Ethyl acetoacetate ethylene acetal	174	C8H14O4	006413-10-1	47





Library Search Compound Report

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Data File : BP004319.D
Acq On : 16 Dec 2020 12:08
Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

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

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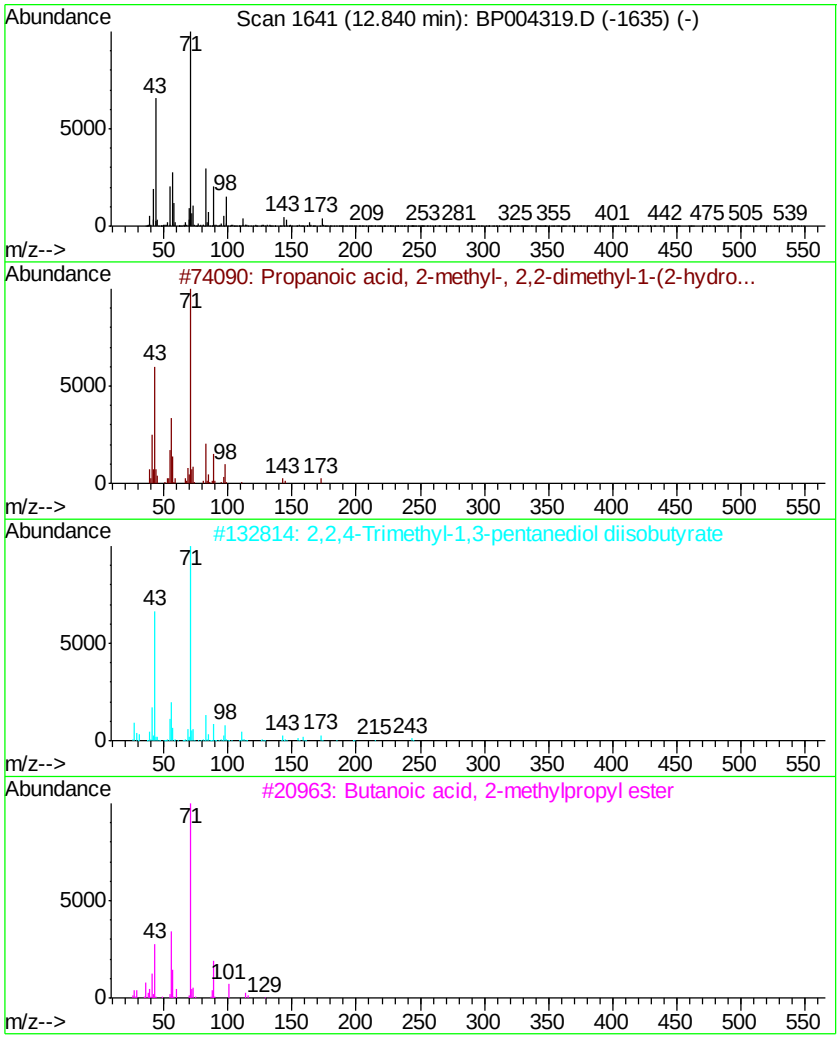
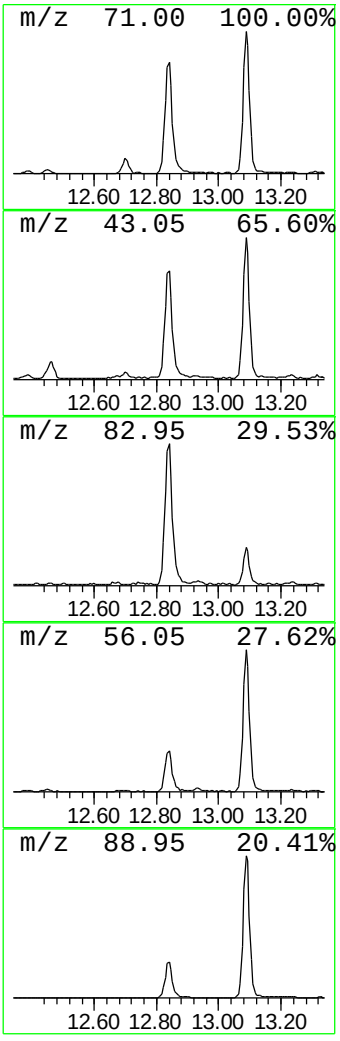
Peak Number 8 Propanoic acid, 2-methyl-, ... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.84	3.75 ng/ul	513116	Acenaphthene-d10	14.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-methyl-, 2,2-d...	216	C12H24O3	074367-33-2	78
2	2,2,4-Trimethyl-1,3-pentanediol ...	286	C16H30O4	006846-50-0	59
3	Butanoic acid, 2-methylpropyl ester	144	C8H16O2	000539-90-2	42
4	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	38
5	Isophytol	296	C20H40O	000505-32-8	38





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Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

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

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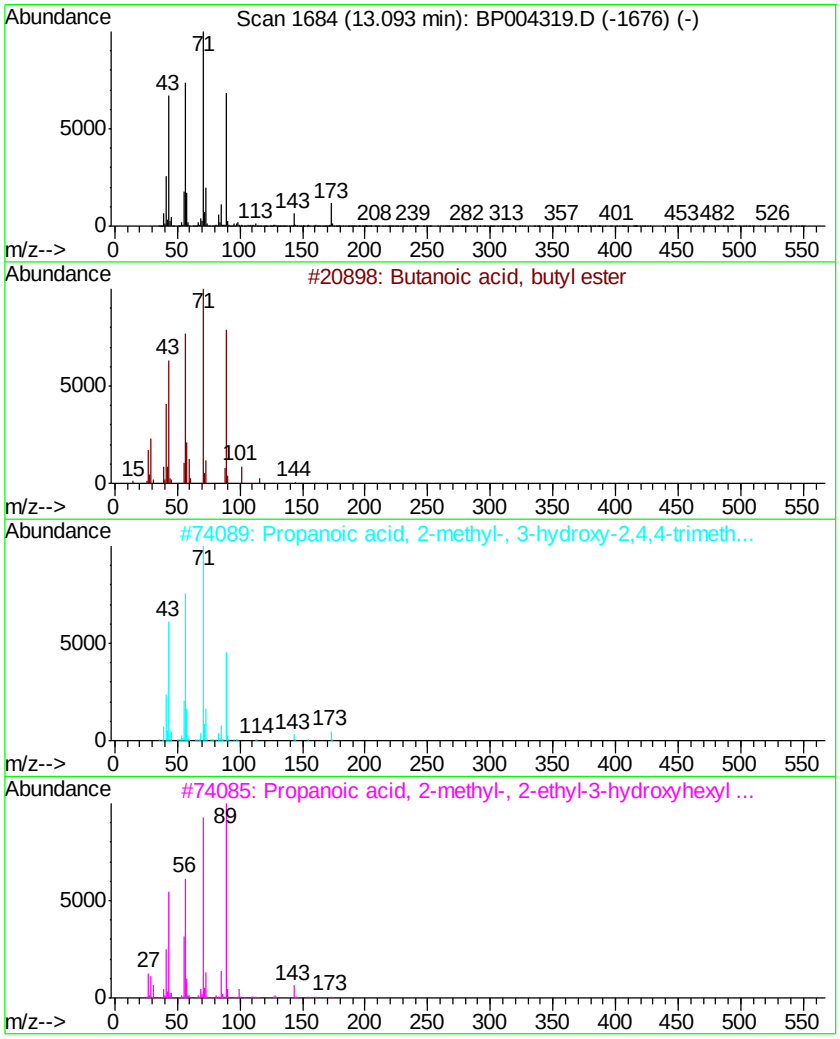
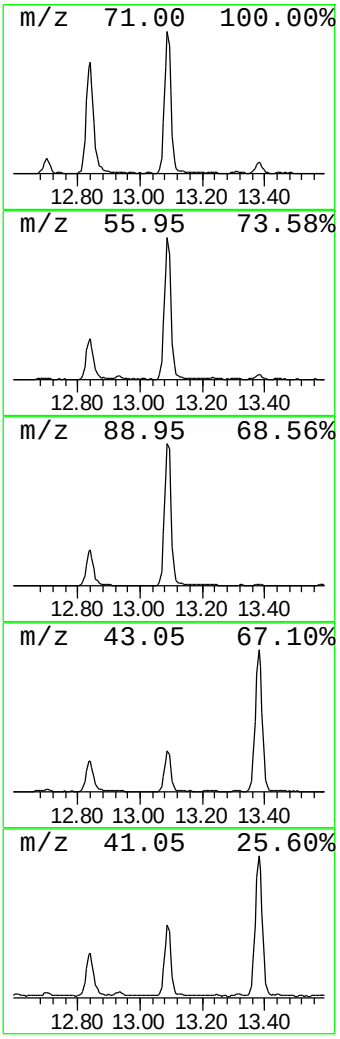
Peak Number 9 Butanoic acid, butyl ester Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.09	4.80 ng/ul	656114	Acenaphthene-d10	14.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	87
2	Propanoic acid, 2-methyl-, 3-hyd...	216	C12H24O3	074367-34-3	80
3	Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	64
4	Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	64
5	Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	53





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Acq On : 16 Dec 2020 12:08
Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

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

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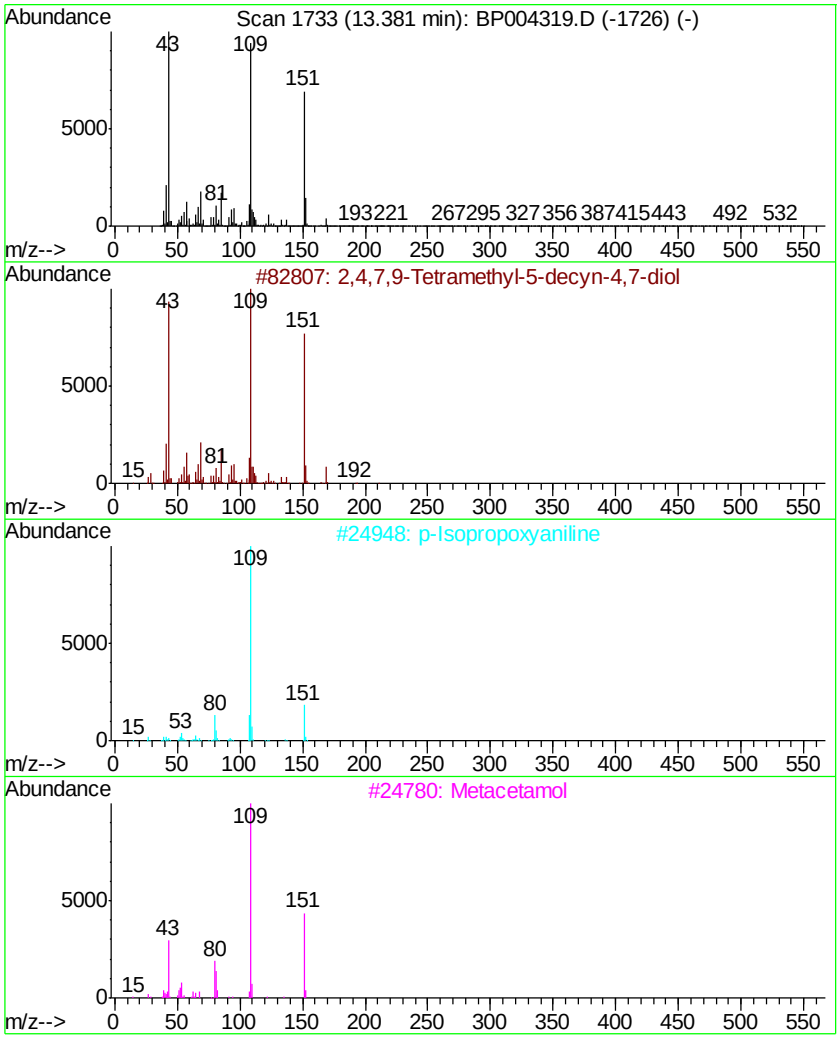
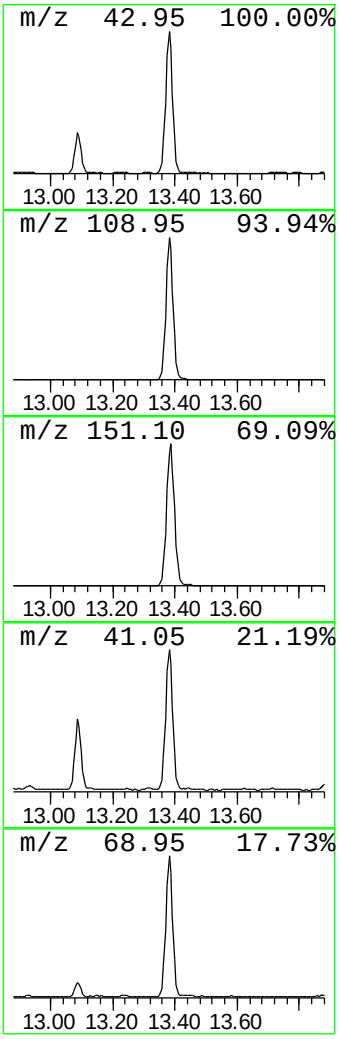
Peak Number 10 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.38	14.96 ng/ul	2046400	Acenaphthene-d10	14.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91
2	p-Isopropoxyaniline	151	C9H13NO	007664-66-6	53
3	Metacetamol	151	C8H9NO2	000621-42-1	53
4	Metacetamol	151	C8H9NO2	000621-42-1	53
5	3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	53





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Acq On : 16 Dec 2020 12:08
Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

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

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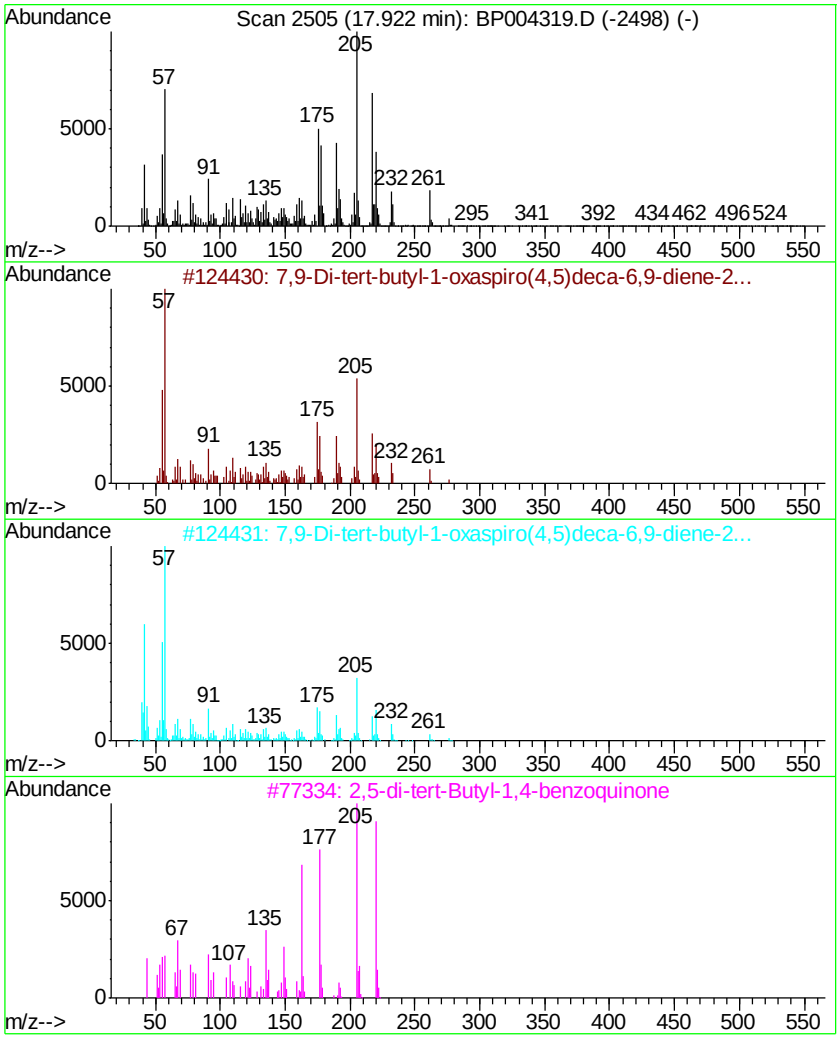
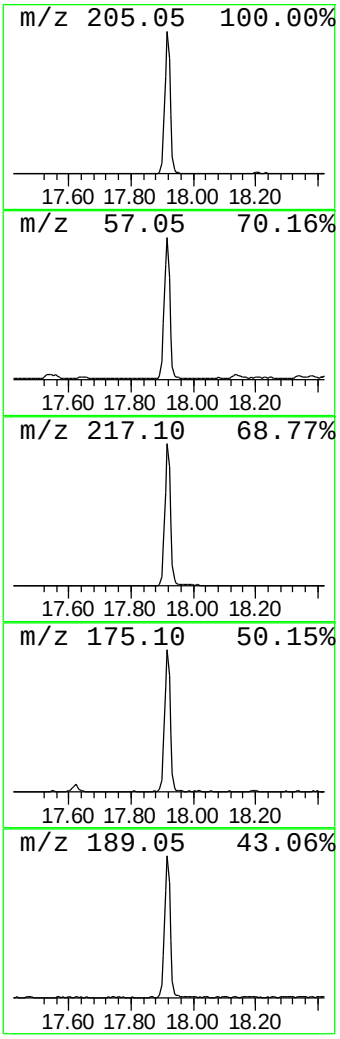
Peak Number 11 7,9-Di-tert-butyl-1-oxaspir... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.92	9.41 ng/ul	1526940	Phenanthrene-d10	17.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	96
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	89
3	2,5-di-tert-Butyl-1,4-benzoquinone	220	C14H20O2	002460-77-7	45
4	2,5-Cyclohexadien-1-one, 2,6-bis...	232	C16H24O	006738-27-8	42
5	Ethanone, 1-(5,6,7,8-tetrahydro-...	220	C14H20O2	071596-88-8	35





Tentatively Identified Compound (LSC) summary

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Data File : BP004319.D
Acq On : 16 Dec 2020 12:08
Operator : CG/JU
Sample : L5071-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A54K2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.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
Cyclohexanone	5.89	3.8	ng/ul	261456	1	7.84	1367040	20.0
Ethanol, 2-butoxy-	5.92	3.9	ng/ul	265951	1	7.84	1367040	20.0
1-Hexanol, 2-ethyl-	7.90	2.6	ng/ul	177377	1	7.84	1367040	20.0
Benzyl alcohol	8.08	2.0	ng/ul	139969	1	7.84	1367040	20.0
Phenol, 2-methoxy-	8.93	8.3	ng/ul	565770	1	7.84	1367040	20.0
Ethanol, 2-(hexyl...	9.16	10.0	ng/ul	682728	1	7.84	1367040	20.0
Thiophane, propyl-	9.89	3.2	ng/ul	337510	2	10.63	2114220	20.0
Propanoic acid, 2...	12.84	3.8	ng/ul	513116	3	14.49	2736540	20.0
Butanoic acid, bu...	13.09	4.8	ng/ul	656114	3	14.49	2736540	20.0
2,4,7,9-Tetrameth...	13.38	15.0	ng/ul	2046400	3	14.49	2736540	20.0
7,9-Di-tert-butyl...	17.92	9.4	ng/ul	1526940	4	17.25	3245200	20.0