LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007740.D
Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGB8

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\SFAM-EPA-BP102521.M
Title : SVOA CALIBRATION

Signal : TIC: BP007740.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.799	85	87	98	rBV	46811	77123	1.16%	0.139%
2	3.958	109	114	121	rBV2	41068	102677	1.55%	0.184%
3	4.022	121	125	131	rVB	64529	103487	1.56%	0.186%
4	4.211	154	157	163	rVB3	23101	34182	0.52%	0.061%
5	4.487	200	204	211	rVB6	15622	26417	0.40%	0.047%
6	5.034	292	297	303	rVB	35875	53469	0.81%	0.096%
7	6.316	510	515	520	rBV	24577	39208	0.59%	0.070%
8	6.622	562	567	574	rVB	45051	70637	1.07%	0.127%
9	6.934	613	620	625	rBV7	12918	32855	0.50%	0.059%
10	7.093	639	647	658	rVB	624521	1004043	15.16%	1.804%
11	7.257	668	675	684	rBV	538276	859626	12.98%	1.544%
12	7.381	689	696	704	rBV3	53038	119292	1.80%	0.214%
13	7.463	704	710	722	rVB	670883	1058212	15.98%	1.901%
14	7.840	768	774	782	rVB	256382	386002	5.83%	0.693%
15	7.934	782	790	797	rBV	1069120	1746333	26.36%	3.137%
16	8.010	797	803	810	rVV2	37611	91118	1.38%	0.164%
17	8.093	811	817	825	rVV9	16306	60466	0.91%	0.109%
18	8.163	825	829	836	rVB3	36993	65386	0.99%	0.117%
19	8.640	899	910	924	rBV	430858	751351	11.34%	1.350%
20	8.752	924	929	935	rVB	28192	48021	0.72%	0.086%
21	9.087	978	986	999	rBV	608881	1007348	15.21%	1.810%
22	9.810	1100	1109	1117	rBV	481864	798931	12.06%	1.435%
23	10.351	1194	1201	1211	rBV	762655	1259502	19.01%	2.263%
24	10.516	1220	1229	1241	rBV2	109347	208459	3.15%	0.374%
25	10.734	1257	1266	1277	rBV	1441525	2474506	37.36%	4.445%
26	10.869	1283	1289	1298	rBV	255136	462205	6.98%	0.830%
27	11.157	1334	1338	1345	rBV7	18988	41566	0.63%	0.075%
28	11.369	1370	1374	1377	rBV4	19426	25068	0.38%	0.045%
29	11.446	1381	1387	1393	rVB8	38373	92772	1.40%	0.167%
30	11.516	1393	1399	1406	rBV8	35107	101844	1.54%	0.183%
31	11.657	1418	1423	1426	rBV6	20952	41215	0.62%	0.074%
32	11.781	1440	1444	1451	rBV6	92299	225992	3.41%	0.406%
33	12.010	1478	1483	1484	rBV3	30526	47623	0.72%	0.086%
34	12.198	1511	1515	1519	rBV5	46757	93250	1.41%	0.168%
35	12.481	1558	1563	1569	rBV5	131619	301669	4.55%	0.542%
36	12.540	1569	1573	1576	rBV4	63635	110030	1.66%	0.198%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007740.D
Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGB8

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\SFAM-EPA-BP102521.M
Title : SVOA CALIBRATION

37	12.604	1581	1584	1587	rBV5	66877	101416	1.53%	0.182%
38	12.651	1587	1592	1597	rVV3	217694	370097	5.59%	0.665%
39	12.728	1603	1605	1610	rVB2	108141	120617	1.82%	0.217%
40	12.934	1637	1640	1641	rBV3	106328	110763	1.67%	0.199%
41	13.016	1652	1654	1659	rVB2	214154	276346	4.17%	0.496%
42	13.081	1660	1665	1668	rBV6	105475	246730	3.72%	0.443%
43	13.122	1669	1672	1676	rVB	335992	452217	6.83%	0.812%
44	13.275	1695	1698	1704	rVB4	136824	202337	3.05%	0.363%
45	13.363	1709	1713	1716	rBV3	261104	419129	6.33%	0.753%
46	13.516	1735	1739	1743	rVB3	403960	630775	9.52%	1.133%
47	13.569	1745	1748	1750	rBV2	208349	293346	4.43%	0.527%
48	13.851	1792	1796	1801	rBV6	408254	720086	10.87%	1.294%
49	13.981	1814	1818	1820	rBV	1050433	1568473	23.68%	2.818%
50	14.075	1830	1834	1841	rBV3	844348	1739705	26.26%	3.125%
51	14.234	1857	1861	1862	rBV3	584863	722158	10.90%	1.297%
52	14.263	1862	1866	1871	rVB	1354170	2076679	31.35%	3.731%
53	14.322	1871	1876	1879	rBV5	595715	1149757	17.36%	2.065%
54	14.410	1887	1891	1898	rVB3	1355420	2540509	38.35%	4.564%
55	14.481	1899	1903	1908	rBV4	572495	912644	13.78%	1.640%
56	14.569	1909	1918	1923	rBV4	2673262	6623997	100.00%	11.900%
57	14.904	1971	1975	1978	rBV	657503	1003673	15.15%	1.803%
58	15.104	2007	2009	2018	rVB6	754851	1282569	19.36%	2.304%
59	15.369	2049	2054	2059	rBV2	1757698	2975624	44.92%	5.346%
60	15.563	2084	2087	2092	rVB	1638989	2280014	34.42%	4.096%
61	16.333	2214	2218	2224	rVB	3051883	4348798	65.65%	7.812%
62	17.163	2356	2359	2366	rVB2	1903344	2899502	43.77%	5.209%
63	17.328	2384	2387	2396	rVB	1928699	3007450	45.40%	5.403%
64	17.428	2401	2404	2409	rVB	1935906	2567537	38.76%	4.612%

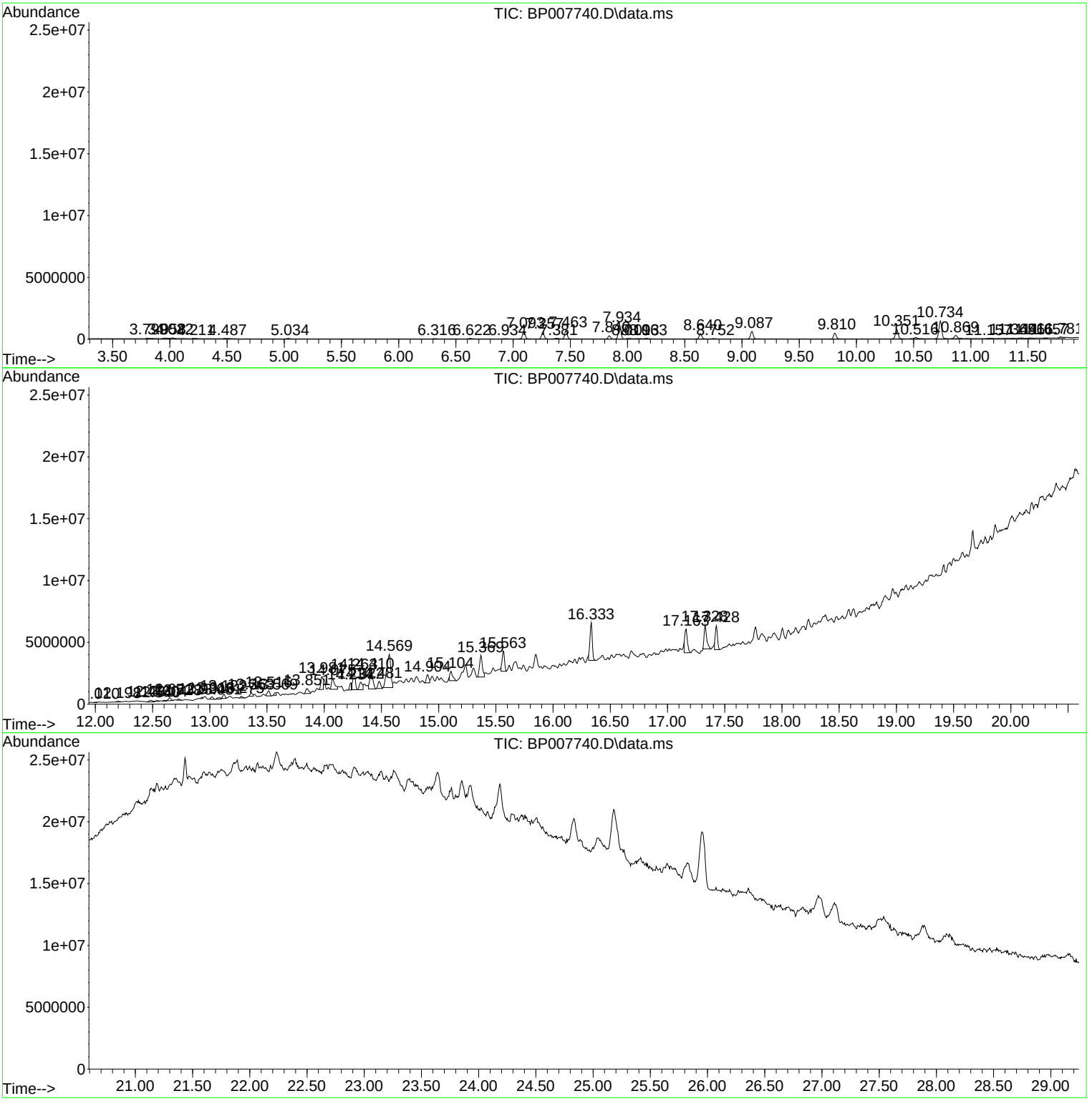
Sum of corrected areas: 55664833

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007740.D
Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007740.D
Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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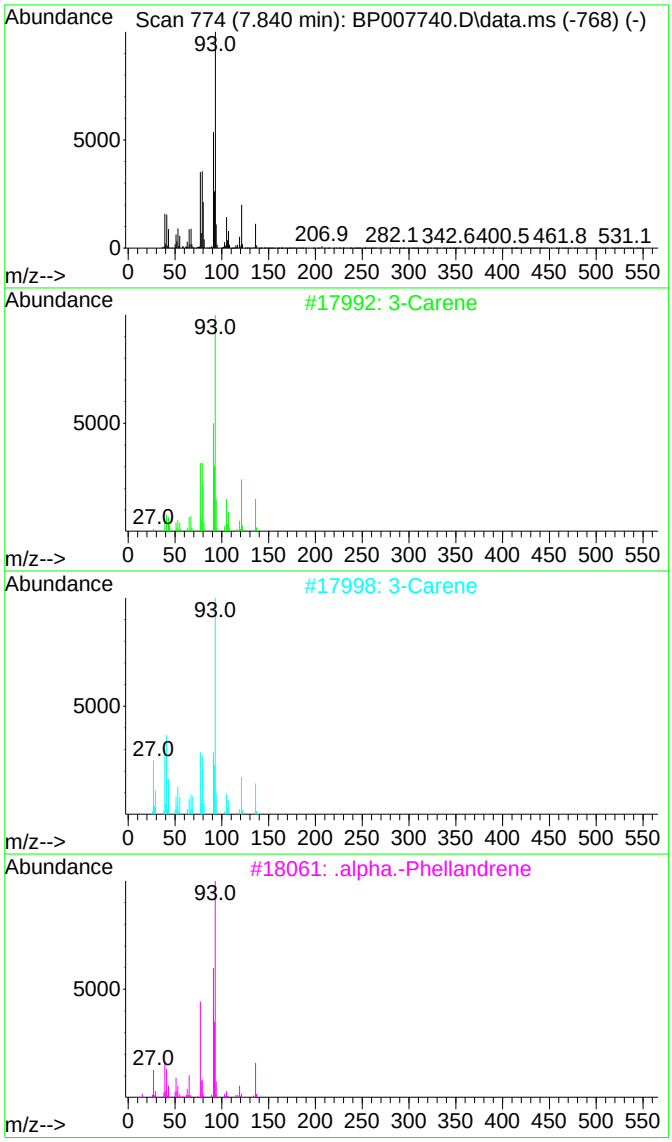
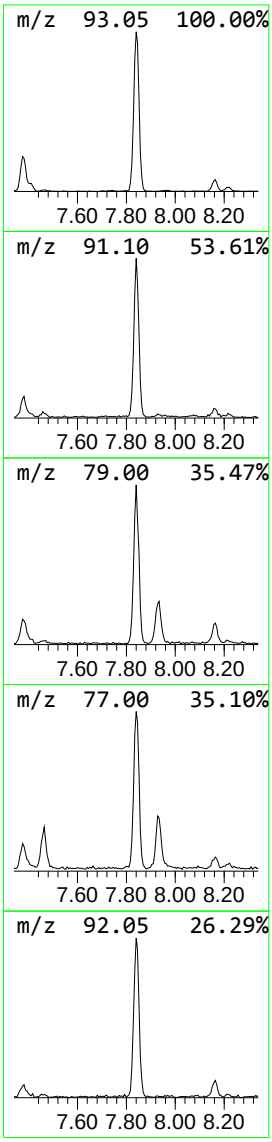
Peak Number 1 3-Carene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.840	4.42 ng/ul	386002	1,4-Dichlorobenzene-d4	7.934

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Carene	136	C10H16	013466-78-9	97
2	3-Carene	136	C10H16	013466-78-9	94
3	.alpha.-Phellandrene	136	C10H16	000099-83-2	93
4	(+)-4-Carene	136	C10H16	029050-33-7	93
5	.gamma.-Terpinene	136	C10H16	000099-85-4	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007740.D
Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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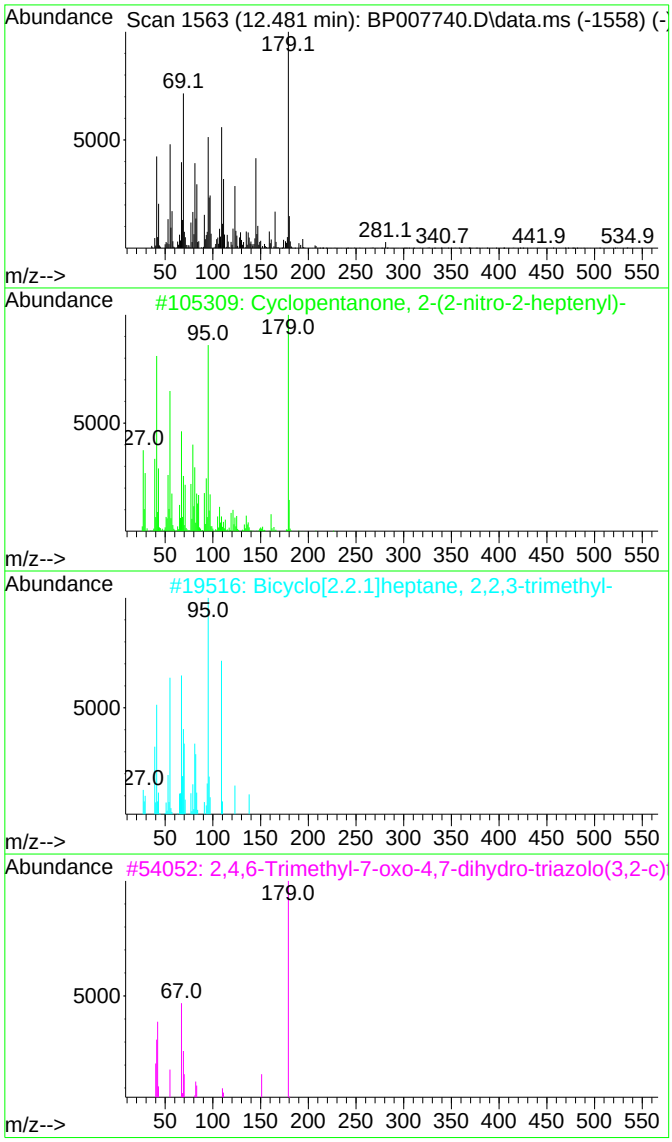
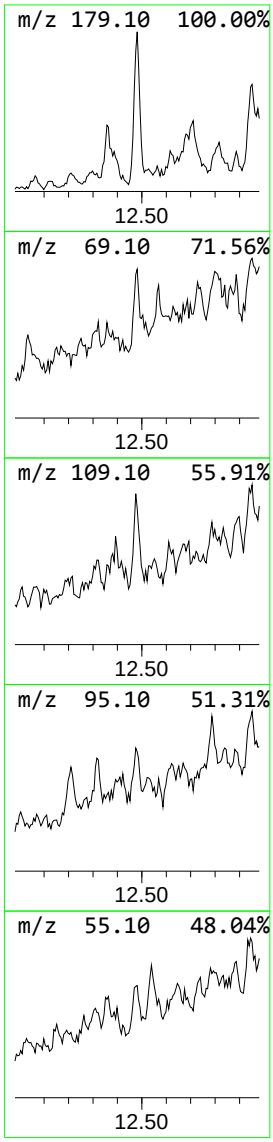
Peak Number 2 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.481	2.44 ng/ul	301669	Naphthalene-d8	10.734

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentanone, 2-(2-nitro-2-hep...	225	C12H19NO3	104313-57-7	25
2	Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	000473-19-8	22
3	2,4,6-Trimethyl-7-oxo-4,7-dihydr...	179	C7H9N5O	025623-78-3	22
4	Tricyclo[4.3.1.1(3,8)]undecane, ...	180	C12H20O	021898-92-0	15
5	.alpha.,.alpha.-Diisopropyl-o-me...	222	C14H22O2	010277-90-4	15





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007740.D
Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGB8

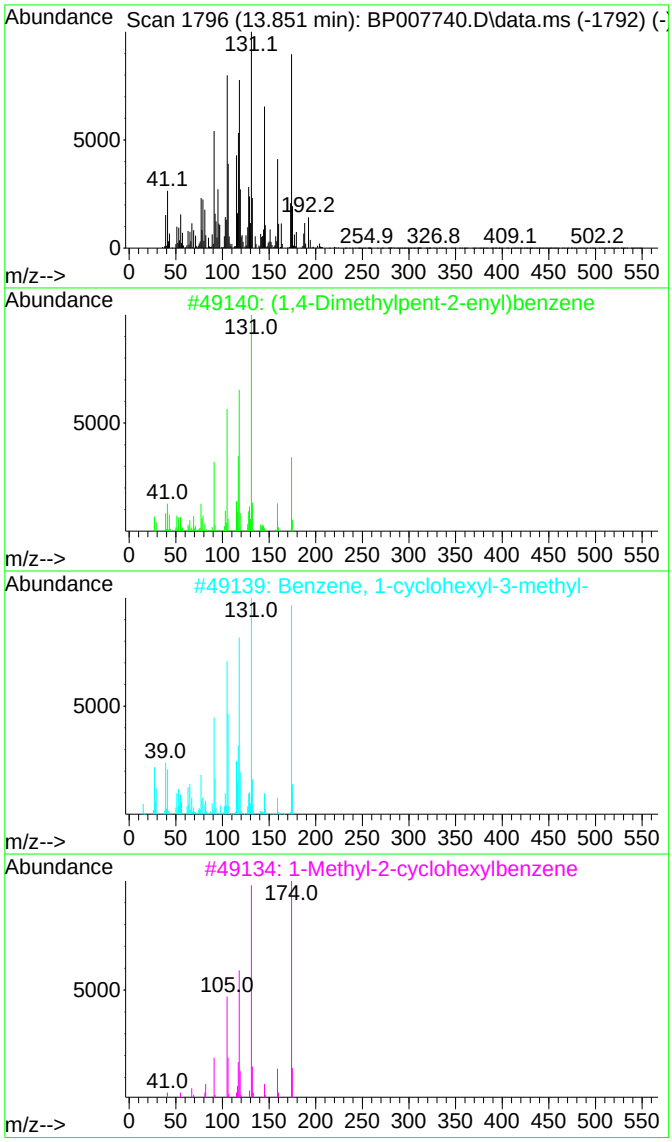
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

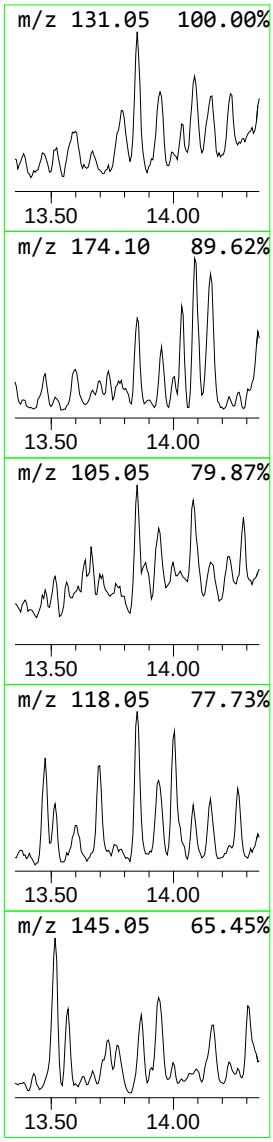
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (1,4-Dimethylpent-2-enyl)be... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.851	2.17 ng/ul	720086	Acenaphthene-d10	14.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1,4-Dimethylpent-2-enyl)benzene	174	C13H18	1000184-97-9	70
2	Benzene, 1-cyclohexyl-3-methyl-	174	C13H18	004575-46-6	70
3	1-Methyl-2-cyclohexylbenzene	174	C13H18	004501-35-3	53
4	1H-1,3,2-Benzodiazaborole, 2-but...	174	C10H15BN2	031748-14-8	46
5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007740.D
Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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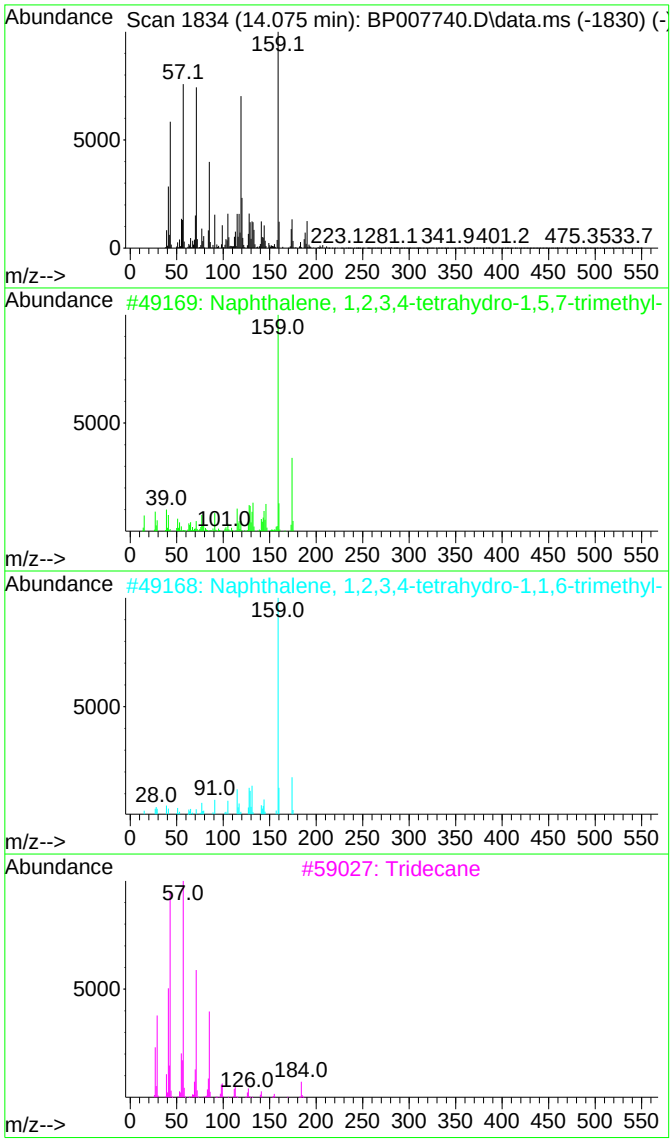
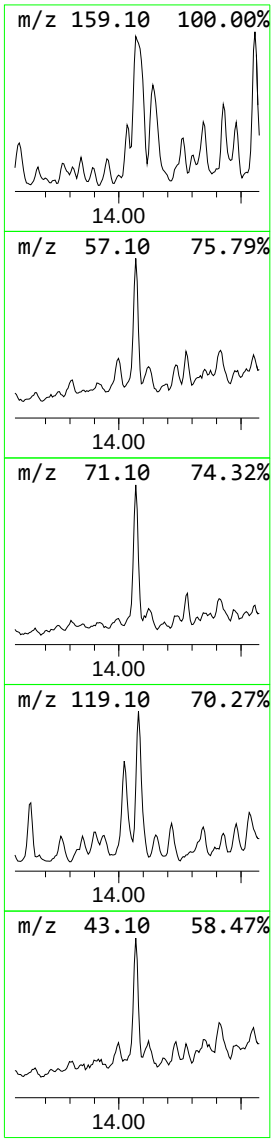
Peak Number 4 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.075	5.25 ng/ul	1739710	Acenaphthene-d10	14.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	021693-55-0	47
2	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	000475-03-6	47
3	Tridecane	184	C13H28	000629-50-5	40
4	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	000475-03-6	38
5	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	030316-36-0	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007740.D
Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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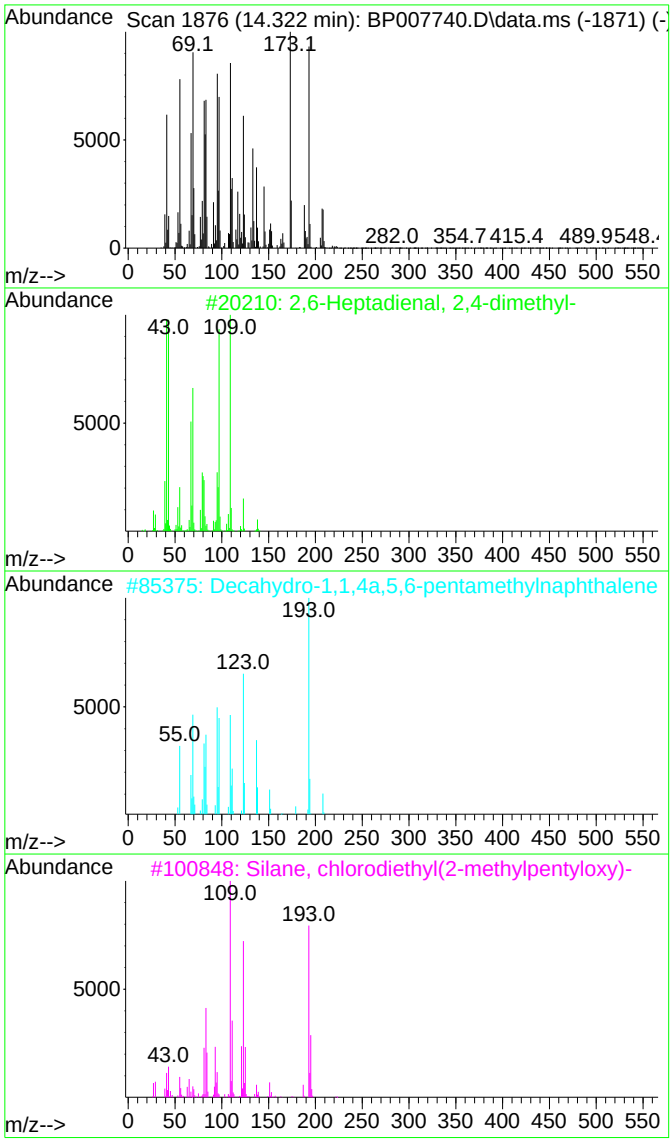
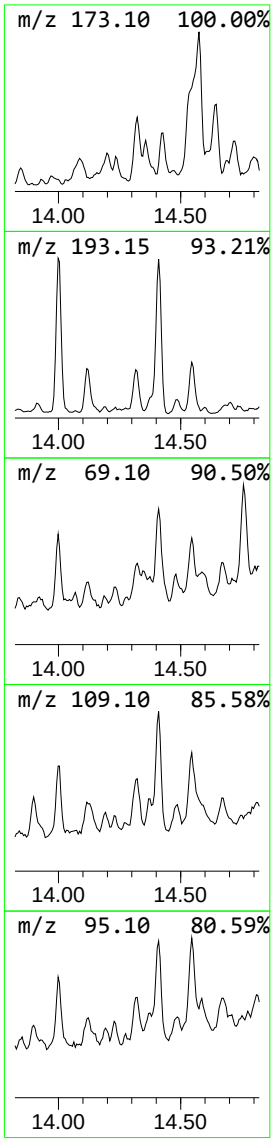
Peak Number 5 unknown-03 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.322	3.47 ng/ul	1149760	Acenaphthene-d10	14.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6-Heptadienal, 2,4-dimethyl-	138	C9H14O	085136-08-9	42
2	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	30
3	Silane, chlorodiethyl(2-methylpe...	222	C10H23ClOSi	1000363-12-7	25
4	11,13-Dimethyl-12-tetradecen-1-o...	282	C18H34O2	1000130-81-0	25
5	Cyclohexane, 1,3-dimethyl-2-meth...	124	C9H16	020348-74-7	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007740.D
Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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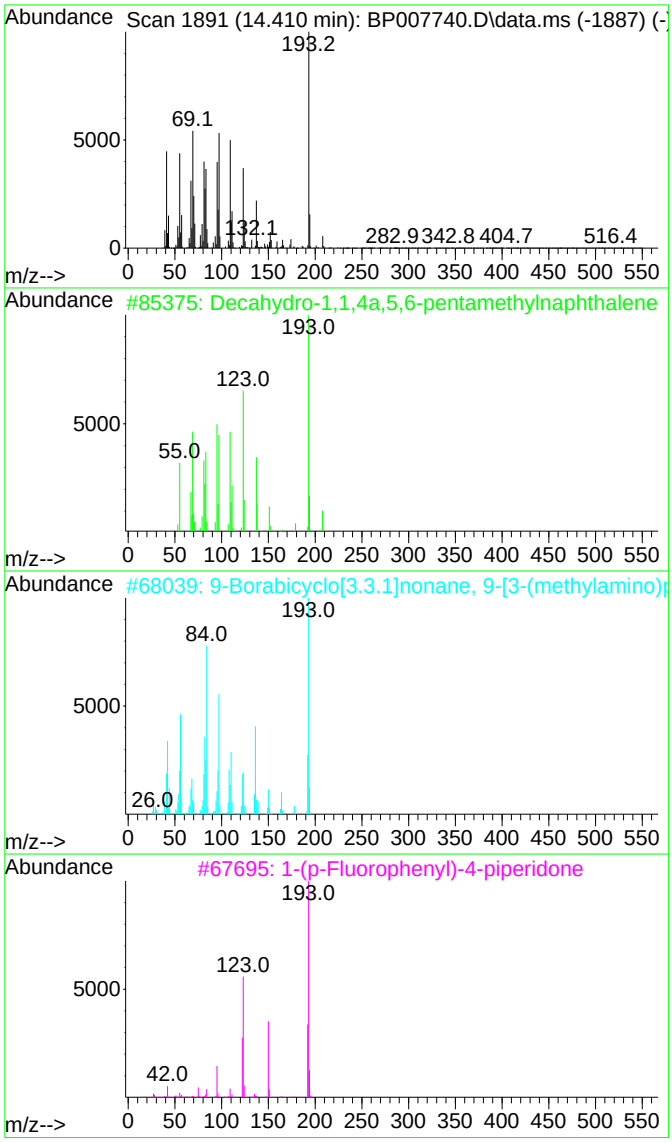
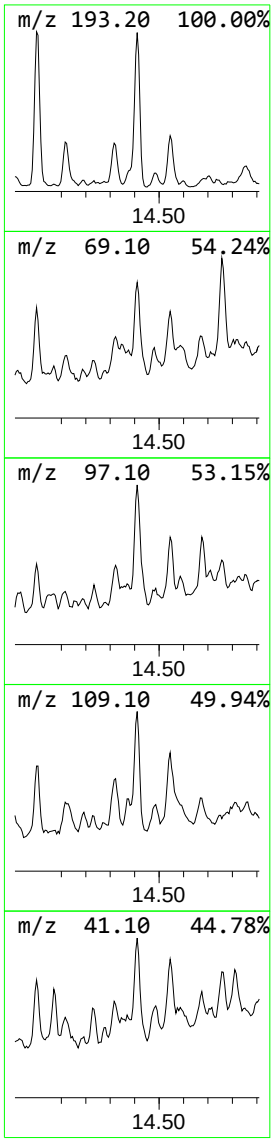
Peak Number 6 unknown-04 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.410	7.67 ng/ul	2540510	Acenaphthene-d10	14.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	47
2	9-Borabicyclo[3.3.1]nonane, 9-[3...	193	C12H24BN	1010162-56-0	43
3	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	38
4	Silane, chlorodiethylhexyloxy-	222	C10H23ClOSi	1000363-74-4	38
5	1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007740.D
Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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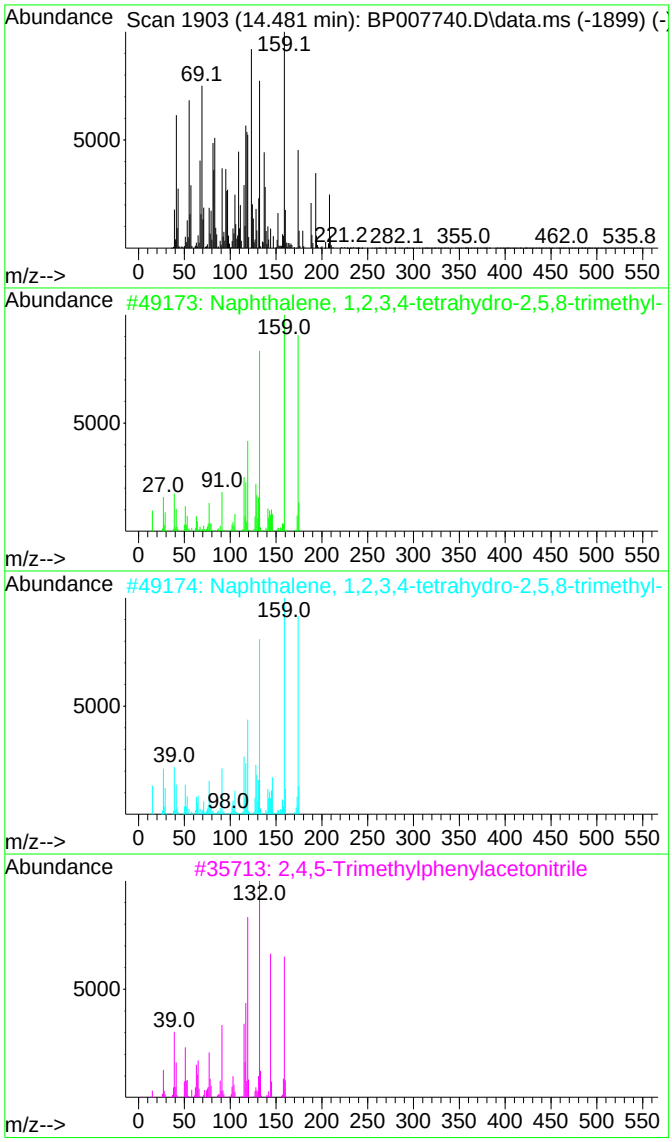
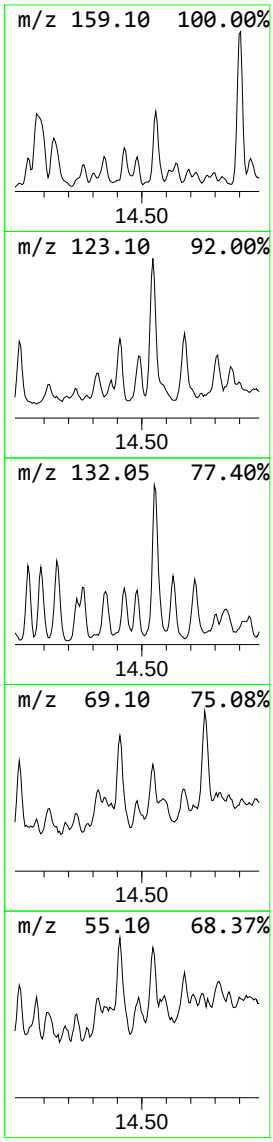
Peak Number 7 unknown-05 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.481	2.76 ng/ul	912644	Acenaphthene-d10	14.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	030316-17-7	38
2	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	030316-17-7	35
3	2,4,5-Trimethylphenylacetonitrile	159	C11H13N	075279-58-2	35
4	Ethanone, 1-[4-(1-methyl-2-prope...	174	C12H14O	109586-49-4	25
5	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	000475-03-6	15





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Data File : BP007740.D
Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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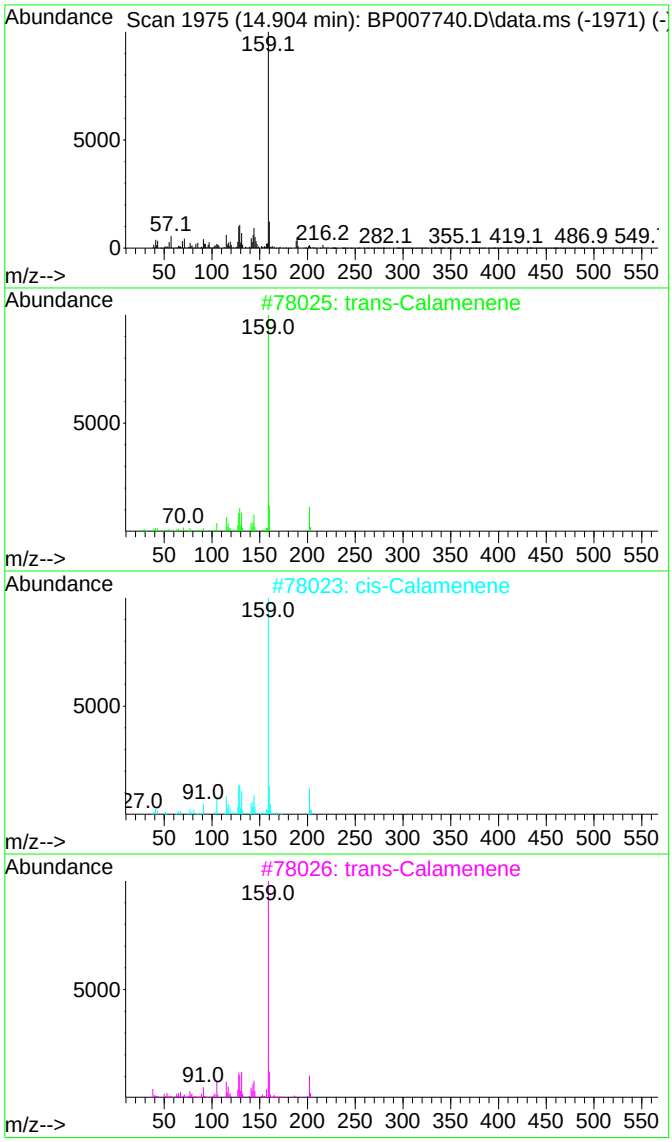
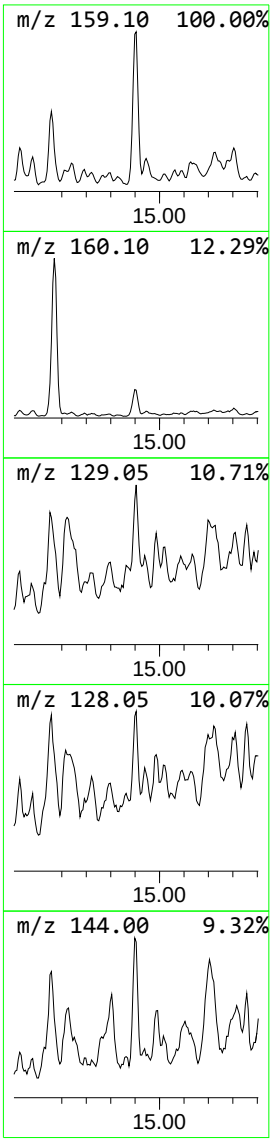
Peak Number 8 trans-Calamenene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.904	3.03 ng/ul	1003670	Acenaphthene-d10	14.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	trans-Calamenene	202	C15H22	073209-42-4	95
2	cis-Calamenene	202	C15H22	072937-55-4	93
3	trans-Calamenene	202	C15H22	073209-42-4	87
4	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	030316-36-0	86
5	4-isopropyl-1,6-dimethyl-1,2,3,4...	202	C15H22	1000378-99-6	86





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Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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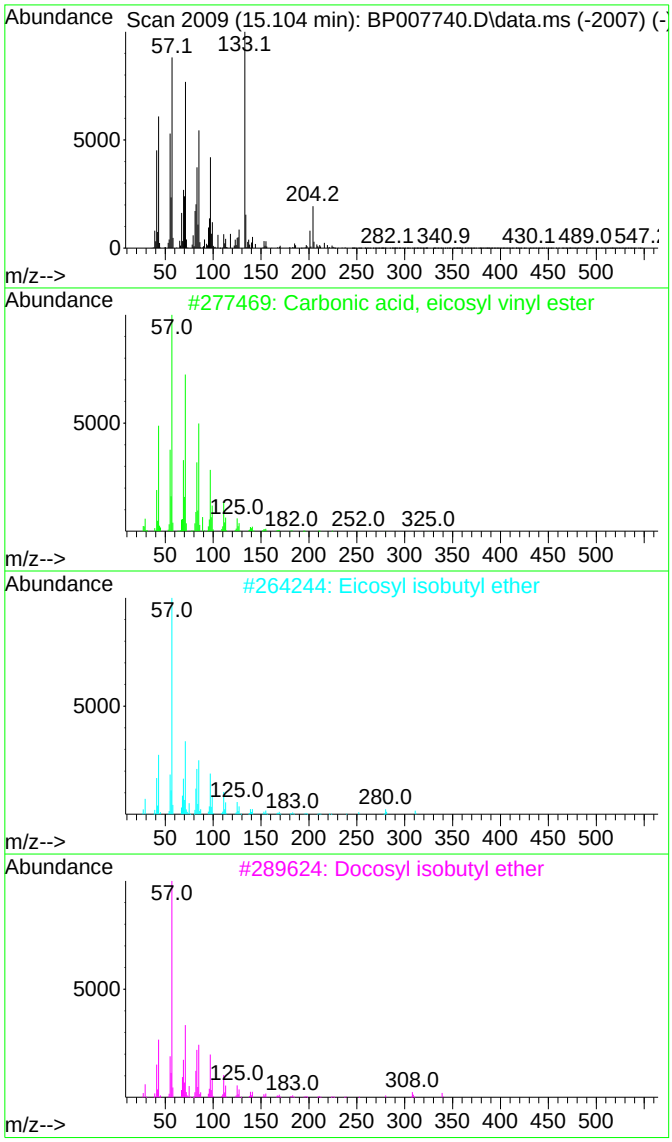
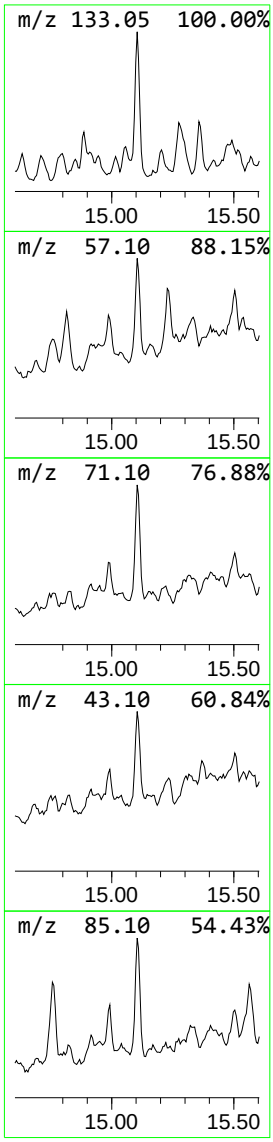
Peak Number 9 unknown-06 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.104	3.87 ng/ul	1282570	Acenaphthene-d10	14.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	47
2	Eicosyl isobutyl ether	354	C24H50O	1000406-33-0	46
3	Docosyl isobutyl ether	382	C26H54O	1000406-33-1	43
4	2-Methyltetracosane	352	C25H52	001560-78-7	43
5	Hexadecane	226	C16H34	000544-76-3	41





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Data File : BP007740.D
Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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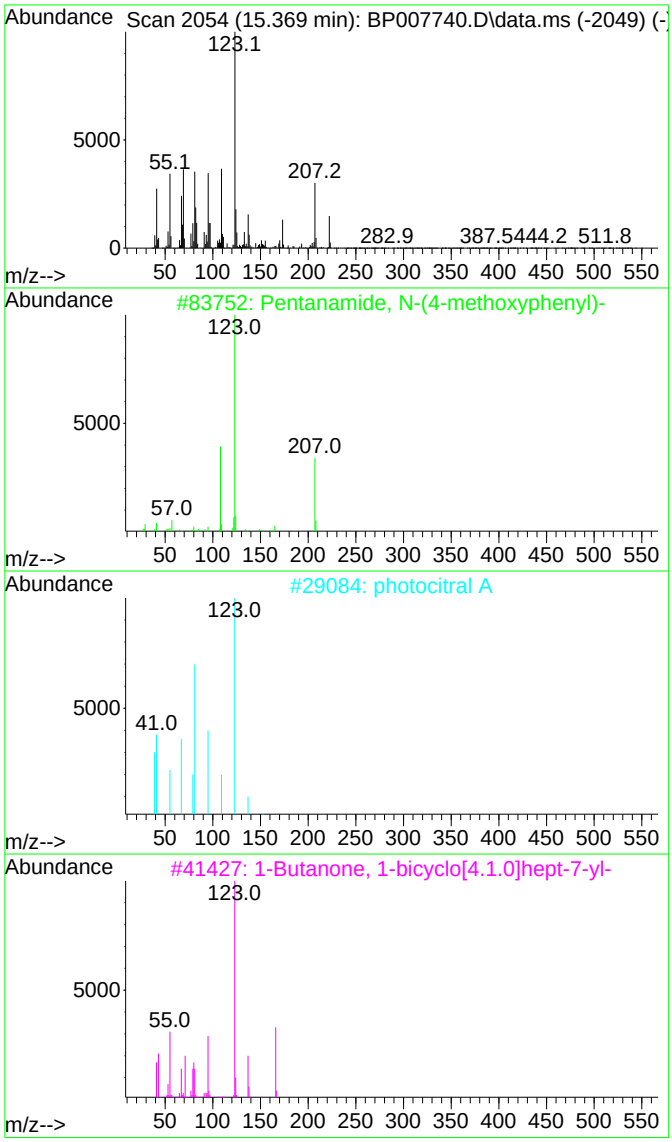
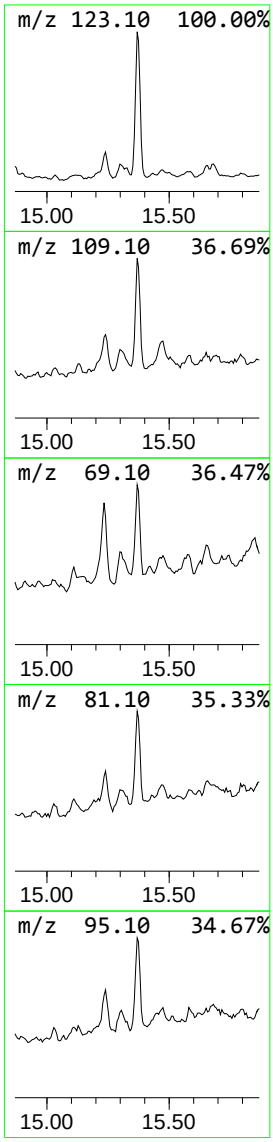
Peak Number 10 unknown-07 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.369	8.98 ng/ul	2975620	Acenaphthene-d10	14.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentanamide, N-(4-methoxyphenyl)-	207	C12H17NO2	1000306-89-3	49
2	photocitral A	152	C10H16O	055253-28-6	47
3	1-Butanone, 1-bicyclo[4.1.0]hept...	166	C11H18O	054764-62-4	47
4	Benzamide, 3-fluoro-N-(2-phenyle...	299	C19H22FNO	1000451-30-2	47
5	Ethanone, 1-[3-[2-methyl-2-(5-me...	222	C13H18O3	080114-28-9	46





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Data File : BP007740.D
Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGB8

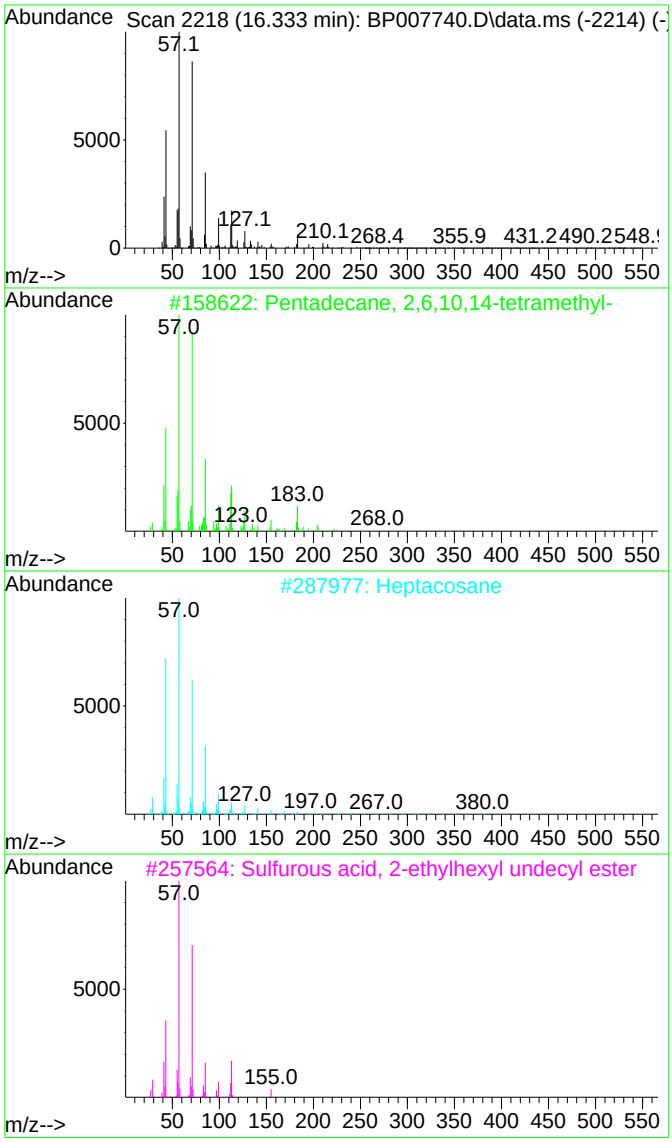
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

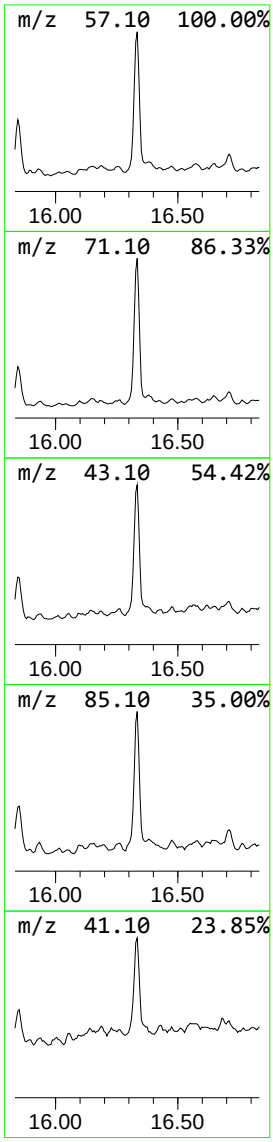
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.334	28.92 ng/ul	4348800	Phenanthrene-d10	17.328

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	90
2	Heptacosane	380	C27H56	000593-49-7	72
3	Sulfurous acid, 2-ethylhexyl und...	348	C19H40O3S	1000309-19-4	64
4	Dodecane	170	C12H26	000112-40-3	59
5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007740.D
Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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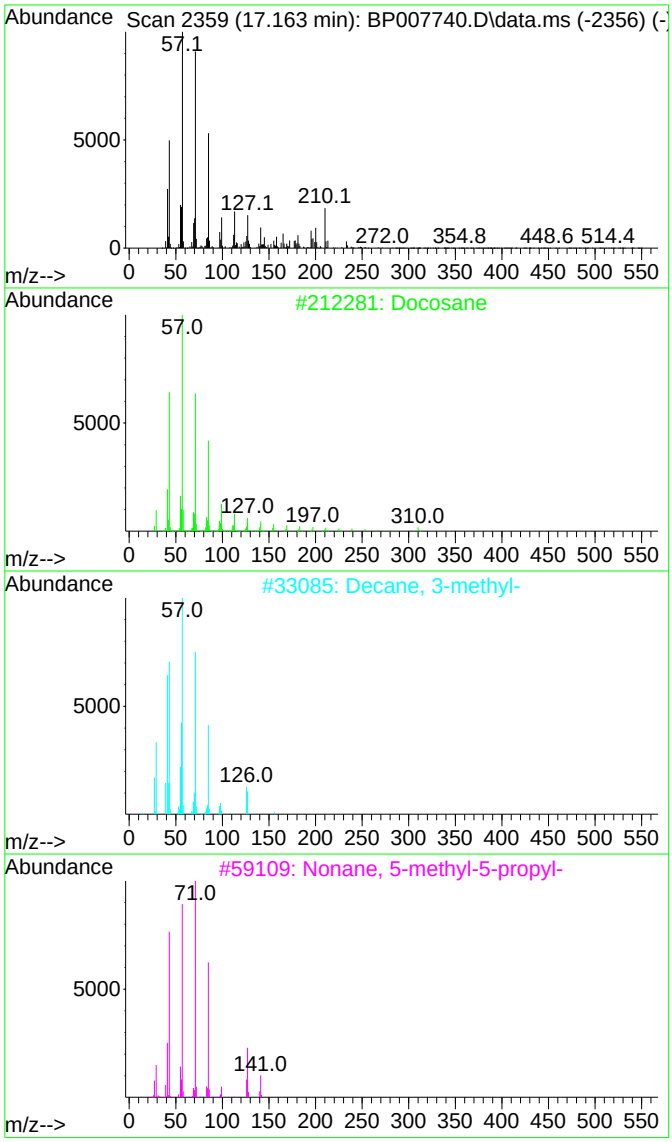
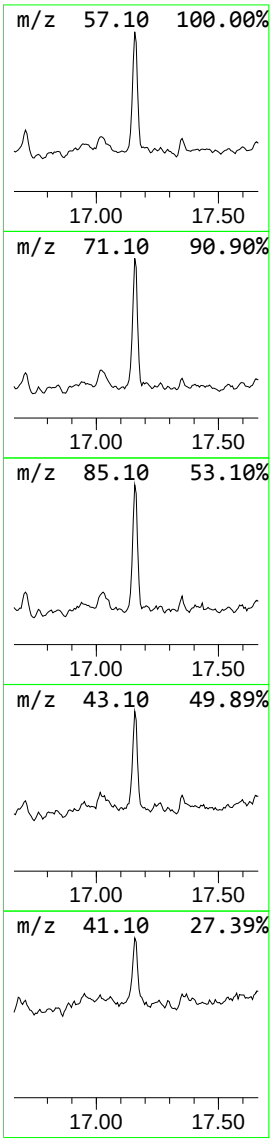
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.163	19.28 ng/ul	2899500	Phenanthrene-d10	17.328

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane	310	C22H46	000629-97-0	72
2	Decane, 3-methyl-	156	C11H24	013151-34-3	70
3	Nonane, 5-methyl-5-propyl-	184	C13H28	017312-75-3	64
4	Hexadecane	226	C16H34	000544-76-3	64
5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	60





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007740.D
Acq On : 28 Oct 2021 03:27
Operator : CG/JU
Sample : M4215-02
Misc :
ALS Vial : 97 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JDGB8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
3-Carene	7.840	4.4	ng/ul	386002	1	7.934	1746330	20.0
unknown-01	12.481	2.4	ng/ul	301669	2	10.734	2474510	20.0
(1,4-Dimethylpe...	13.851	2.2	ng/ul	720086	3	14.569	6624000	20.0
unknown-02	14.075	5.3	ng/ul	1739710	3	14.569	6624000	20.0
unknown-03	14.322	3.5	ng/ul	1149760	3	14.569	6624000	20.0
unknown-04	14.410	7.7	ng/ul	2540510	3	14.569	6624000	20.0
unknown-05	14.481	2.8	ng/ul	912644	3	14.569	6624000	20.0
trans-Calamenene	14.904	3.0	ng/ul	1003670	3	14.569	6624000	20.0
unknown-06	15.104	3.9	ng/ul	1282570	3	14.569	6624000	20.0
unknown-07	15.369	9.0	ng/ul	2975620	3	14.569	6624000	20.0
(DEL) Alkane: S...	16.334	28.9	ng/ul	4348800	4	17.328	3007450	20.0
(DEL) Alkane: S...	17.163	19.3	ng/ul	2899500	4	17.328	3007450	20.0