LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112621\
Data File : BM033273.D
Acq On : 26 Nov 2021 20:05
Operator : CG/JU
Sample : M4817-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A6879

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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\8270-BM112421.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM033273.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.055	287	292	299	rBV	54246	85133	1.25%	0.181%
2	5.507	363	369	380	rVB	1122620	1695774	24.94%	3.615%
3	7.096	632	639	653	rVB	690449	1180384	17.36%	2.516%
4	7.931	773	781	787	rBV	503935	827496	12.17%	1.764%
5	7.990	787	791	795	rVB	45266	68232	1.00%	0.145%
6	8.760	914	922	929	rVB	43515	80981	1.19%	0.173%
7	9.095	972	979	988	rBV	1610128	2741173	40.32%	5.843%
8	9.972	1122	1128	1136	rVB5	28995	70625	1.04%	0.151%
9	10.125	1149	1154	1163	rBV4	57347	117836	1.73%	0.251%
10	10.501	1211	1218	1227	rBV	153405	274357	4.04%	0.585%
11	10.731	1245	1257	1261	rBV	665380	1171539	17.23%	2.497%
12	10.778	1261	1265	1272	rVB	473117	815409	11.99%	1.738%
13	11.466	1376	1382	1390	rBV	48276	99945	1.47%	0.213%
14	11.913	1453	1458	1464	rBV2	43730	77580	1.14%	0.165%
15	11.984	1464	1470	1480	rBV2	41363	150556	2.21%	0.321%
16	12.107	1487	1491	1500	rVB6	29033	70001	1.03%	0.149%
17	12.389	1525	1539	1557	rBV2	1806357	5065885	74.52%	10.798%
18	12.601	1569	1575	1586	rVB	296243	536471	7.89%	1.144%
19	13.178	1666	1673	1682	rBV	3728375	5623002	82.71%	11.986%
20	13.354	1697	1703	1715	rBV2	273422	658390	9.68%	1.403%
21	13.519	1723	1731	1739	rBV	381598	709606	10.44%	1.513%
22	13.583	1739	1742	1755	rVB	69661	171474	2.52%	0.366%
23	13.719	1758	1765	1766	rBV2	78810	126016	1.85%	0.269%
24	13.878	1785	1792	1797	rBV	151663	299944	4.41%	0.639%
25	13.930	1797	1801	1812	rVB2	108992	215655	3.17%	0.460%
26	14.113	1828	1832	1835	rBV	55392	70427	1.04%	0.150%
27	14.278	1855	1860	1864	rBV	51173	74981	1.10%	0.160%
28	14.366	1871	1875	1884	rBV	64912	122210	1.80%	0.261%
29	14.554	1897	1907	1913	rBV	975980	1568156	23.07%	3.343%
30	14.619	1913	1918	1925	rVB3	60924	123608	1.82%	0.263%
31	14.766	1938	1943	1951	rVB4	28496	74631	1.10%	0.159%
32	14.948	1969	1974	1981	rVB3	50270	95851	1.41%	0.204%
33	15.166	2006	2011	2016	rBV2	45407	86198	1.27%	0.184%
34	15.595	2080	2084	2088	rVB2	46840	71624	1.05%	0.153%
35	16.036	2153	2159	2169	rVB	2457613	3625378	53.33%	7.728%
36	16.272	2192	2199	2205	rVB4	52993	126739	1.86%	0.270%

LSC Area Percent Report

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Operator : CG/JU
Sample : M4817-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
A6879

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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\8270-BM112421.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	16.571	2246	2250	2252	rBV2	60881	84782	1.25%	0.181%
38	16.795	2281	2288	2305	rBV	2112540	3298864	48.52%	7.032%
39	17.289	2367	2372	2376	rBV	1132789	1692533	24.90%	3.608%
40	17.407	2388	2392	2403	rVB	200806	309519	4.55%	0.660%
41	17.501	2404	2408	2416	rVB	209575	317806	4.67%	0.677%
42	17.948	2481	2484	2489	rVB3	54578	68385	1.01%	0.146%
43	18.171	2517	2522	2526	rBV2	438172	625425	9.20%	1.333%
44	19.442	2735	2738	2748	rBV	198895	284195	4.18%	0.606%
45	19.671	2774	2777	2780	rBV	106327	119700	1.76%	0.255%
46	19.901	2810	2816	2825	rVB	5421104	6798398	100.00%	14.491%
47	21.154	3025	3029	3035	rVB	447055	546620	8.04%	1.165%
48	21.454	3075	3080	3085	rVB	1369914	1819194	26.76%	3.878%
49	23.783	3470	3476	3484	rVB2	982427	1974792	29.05%	4.209%

Sum of corrected areas: 46913480

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Instrument :
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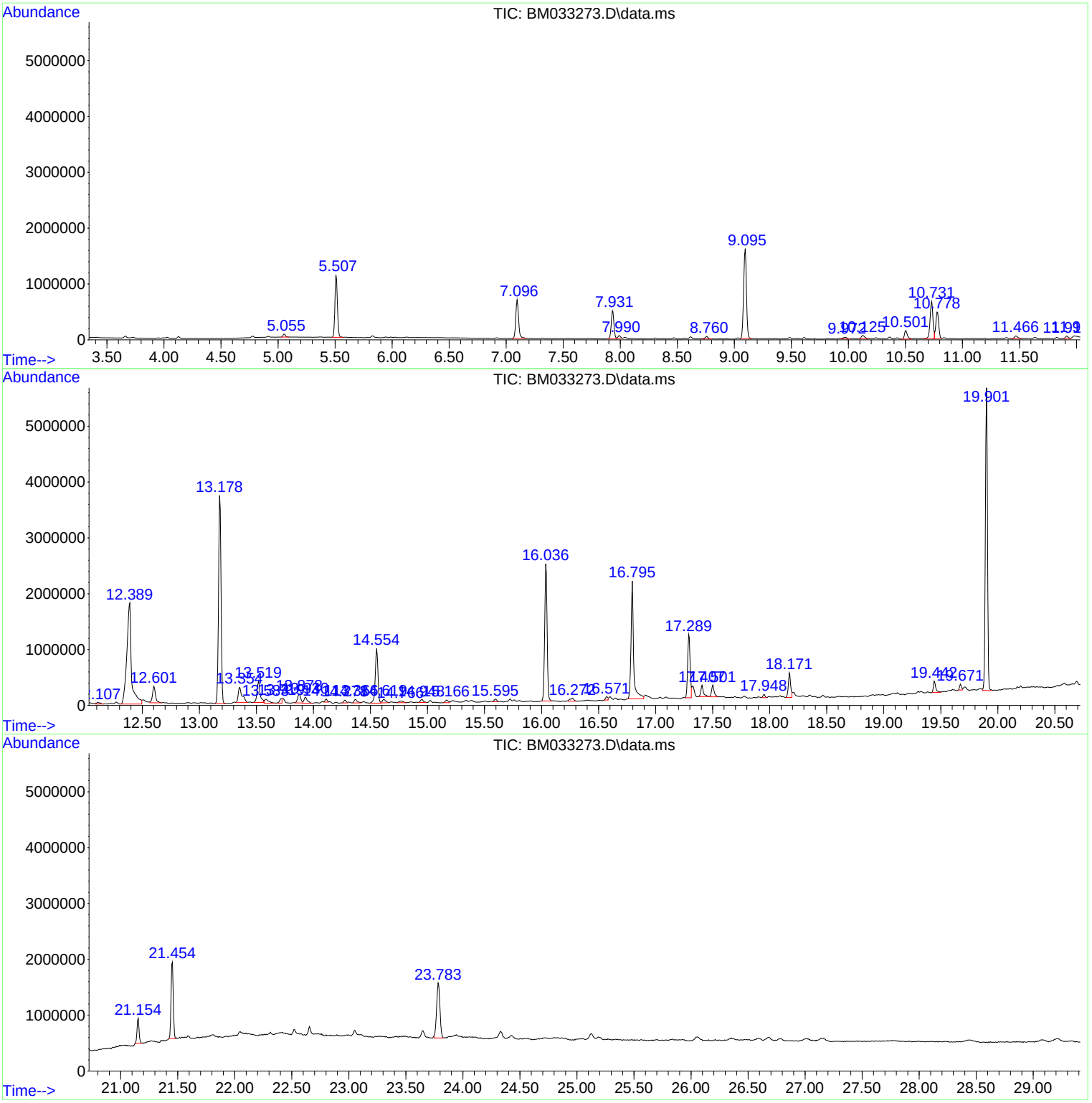
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112621\
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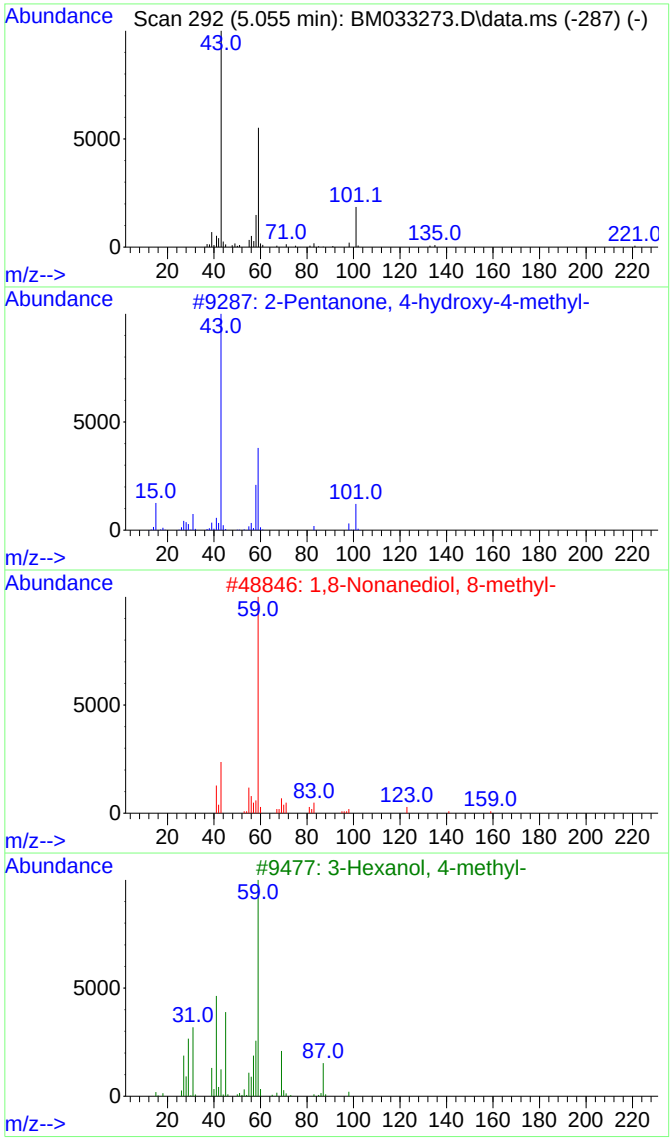
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

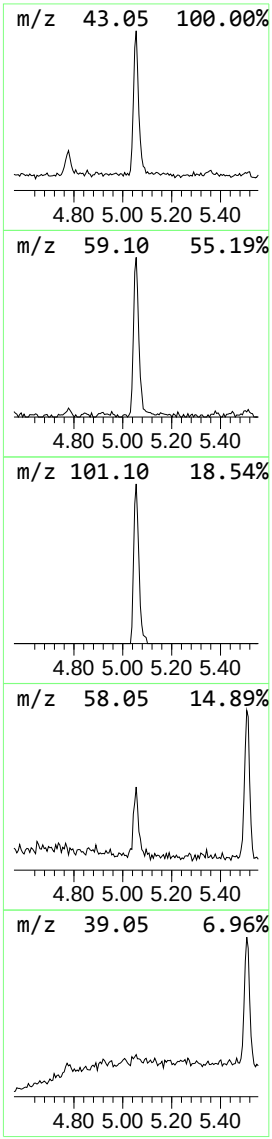
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.055	2.06 ng	85133	1,4-Dichlorobenzene-d4	7.931

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	32
3	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
5	Threo-2-methyl-3,4-dibromo-2-but...	244	C5H10Br2O	1000288-19-2	25





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Operator : CG/JU
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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

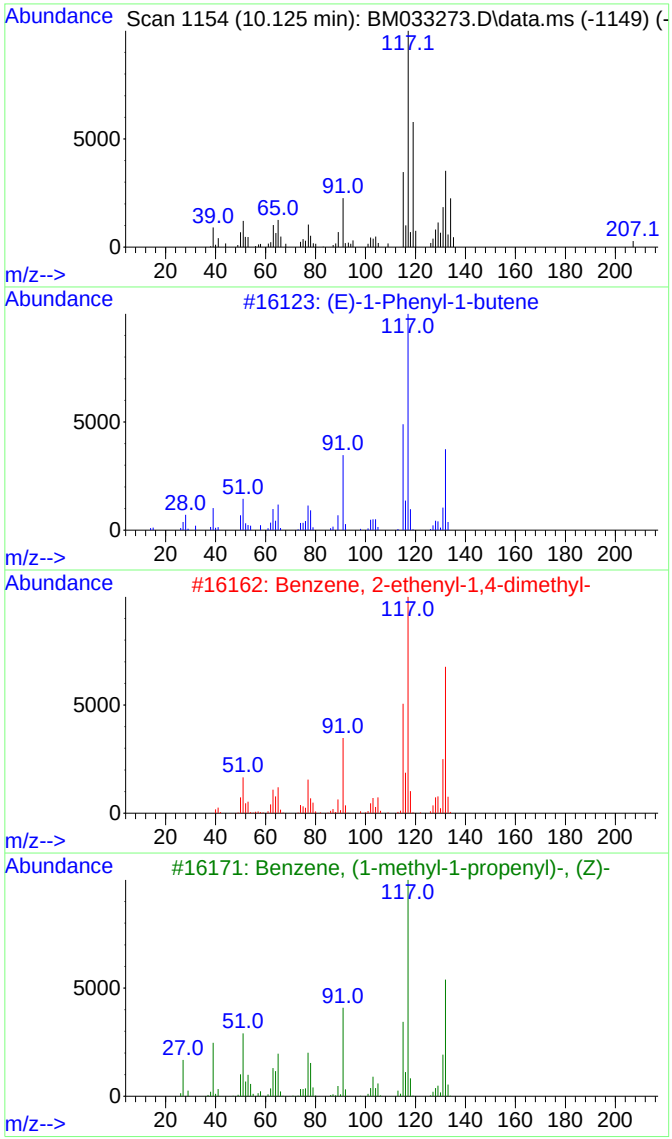
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

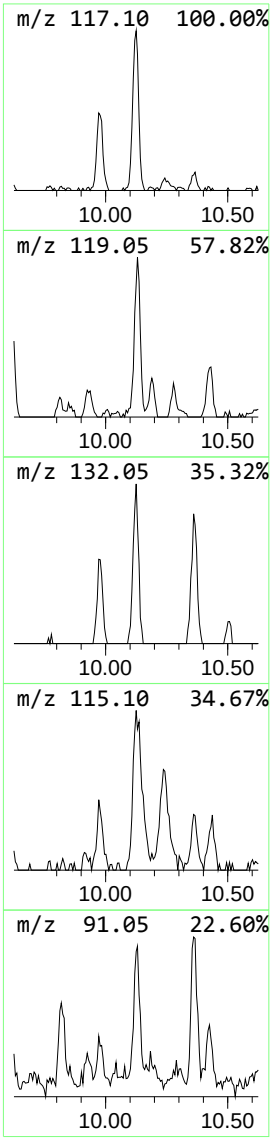
****

Peak Number 2 (E)-1-Phenyl-1-butene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.125	2.01 ng	117836	Naphthalene-d8	10.731

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	91
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	86
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	64
4	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	60
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	60





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Data File : BM033273.D
Acq On : 26 Nov 2021 20:05
Operator : CG/JU
Sample : M4817-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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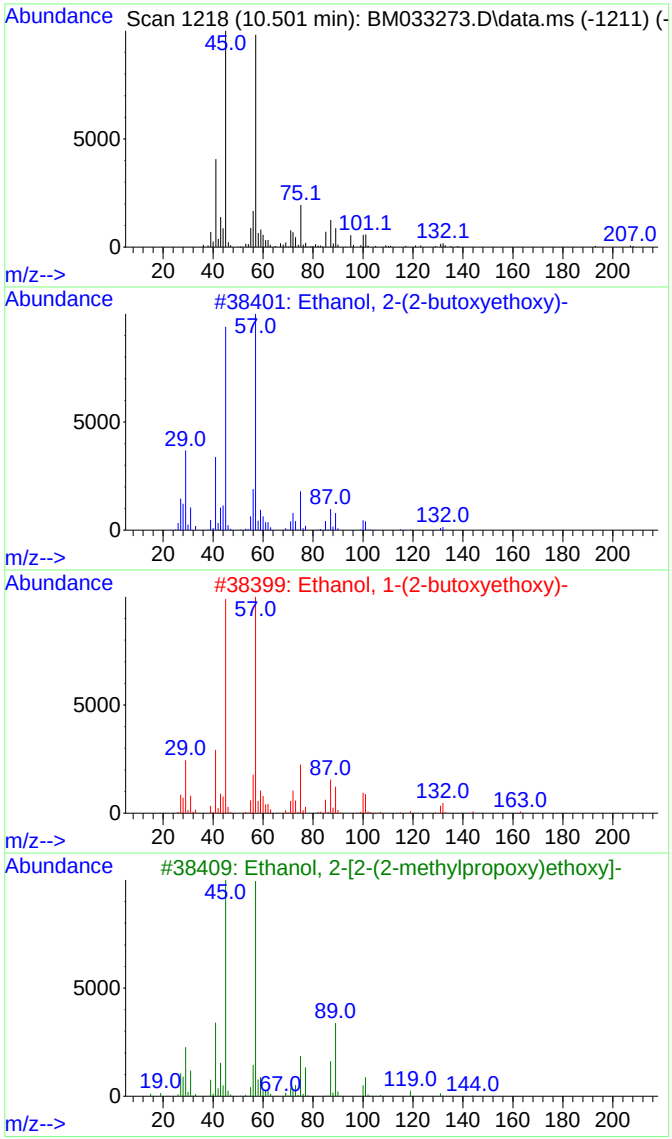
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

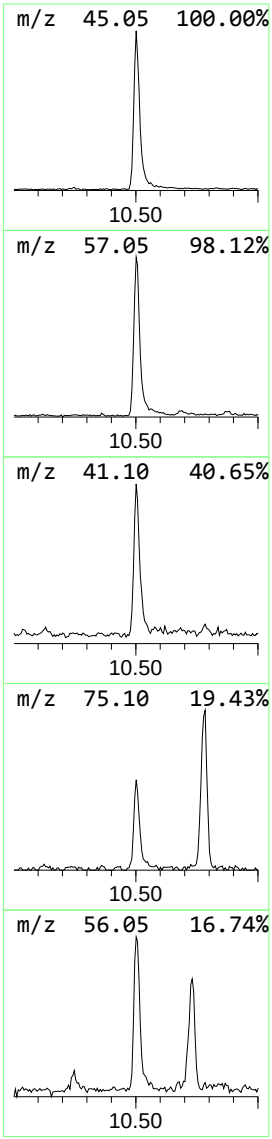
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.501	4.68 ng	274357	Naphthalene-d8	10.731

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
3	Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	56
4	Sulfurous acid, bis(1-methylprop...	194	C8H18O3S	024769-51-5	50
5	Di-sec-Butyl ether	130	C8H18O	006863-58-7	50





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Sample : M4817-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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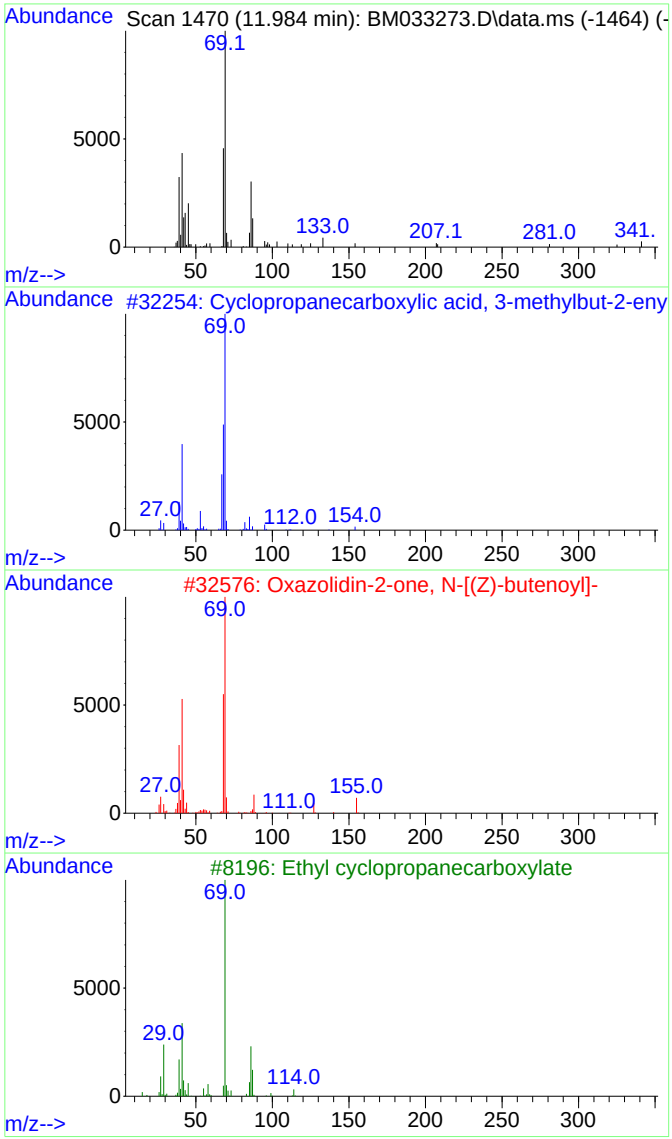
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

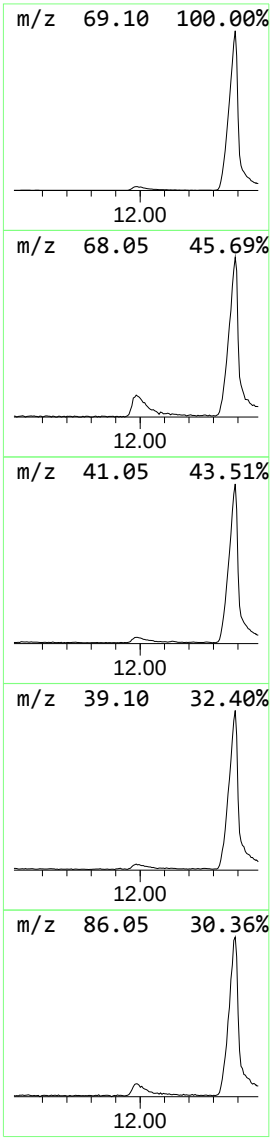
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Cyclopropanecarboxylic acid... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.983	2.57 ng	150556	Naphthalene-d8	10.731

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropanecarboxylic acid, 3-m...	154	C9H14O2	1000299-37-4	52
2	Oxazolidin-2-one, N-[(Z)-butenoyl]-	155	C7H9NO3	1000156-45-4	50
3	Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	50
4	2-Methyl-1,4-butanediol, bis(pen...	396	C11H10F10O4	1000376-24-0	50
5	Glutaric acid, 3-methylbut-2-en-...	268	C15H24O4	1000404-95-9	42





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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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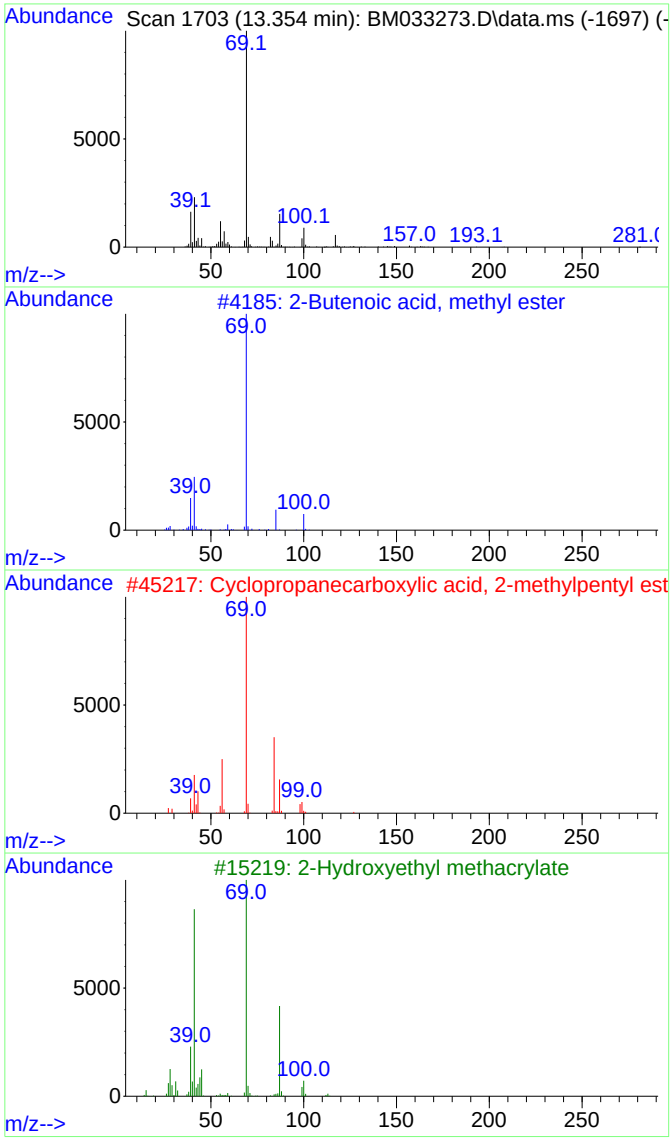
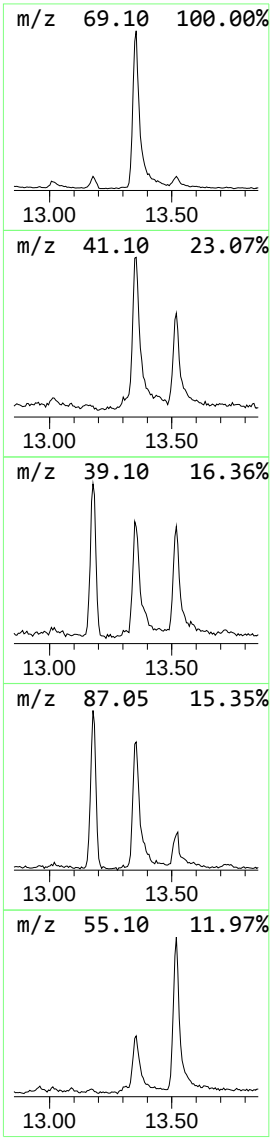
****

Peak Number 5 unknown13.354 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.354	8.40 ng	658390	Acenaphthene-d10	14.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butenoic acid, methyl ester	100	C5H8O2	018707-60-3	42
2	Cyclopropanecarboxylic acid, 2-m...	170	C10H18O2	1000354-65-8	39
3	2-Hydroxyethyl methacrylate	130	C6H10O3	000868-77-9	38
4	2-Butenoic acid, 2-propenyl ester	126	C7H10O2	020474-93-5	36
5	4-Hexen-3-one	98	C6H10O	002497-21-4	36





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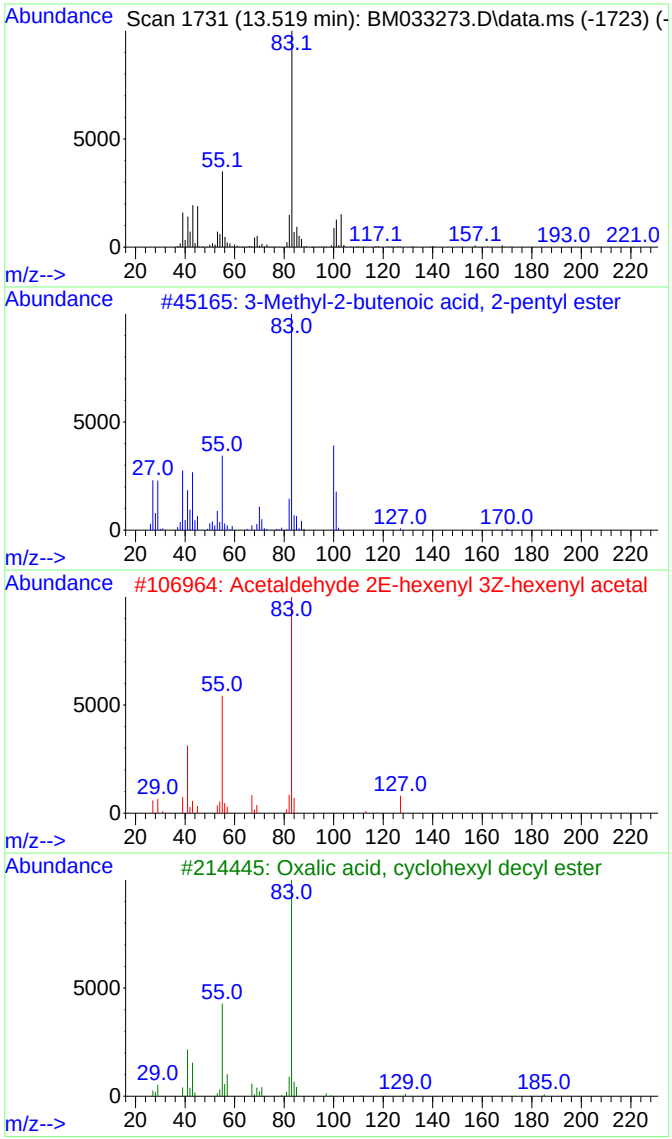
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

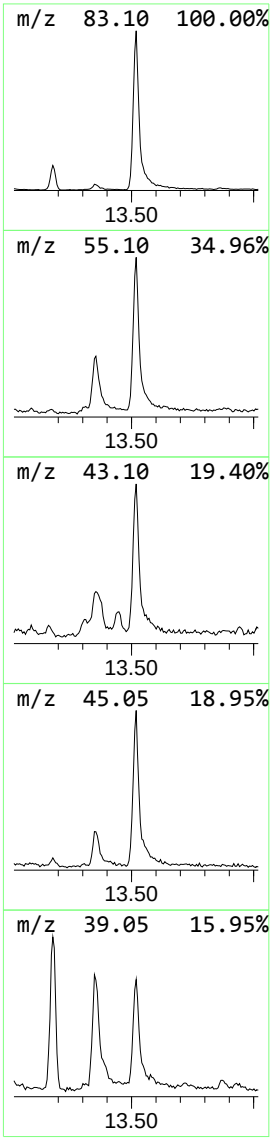
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 3-Methyl-2-butenoic acid, 2... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.519	9.05 ng	709606	Acenaphthene-d10	14.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methyl-2-butenoic acid, 2-pent...	170	C10H18O2	150462-84-3	64
2	Acetaldehyde 2E-hexenyl 3Z-hexen...	226	C14H26O2	1000431-45-5	50
3	Oxalic acid, cyclohexyl decyl ester	312	C18H32O4	1000309-31-2	50
4	Oxalic acid, cyclohexyl undecyl ...	326	C19H34O4	1000309-31-3	50
5	3-Methyl-2-butenoic acid, cyclop...	168	C10H16O2	1000279-23-3	45





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

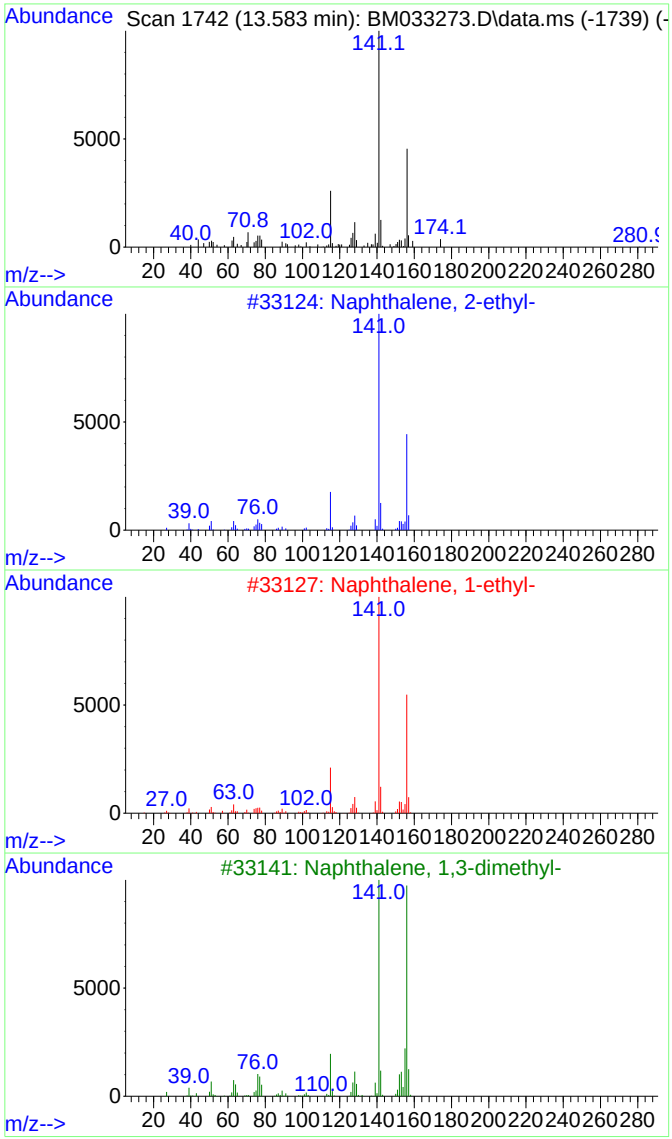
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

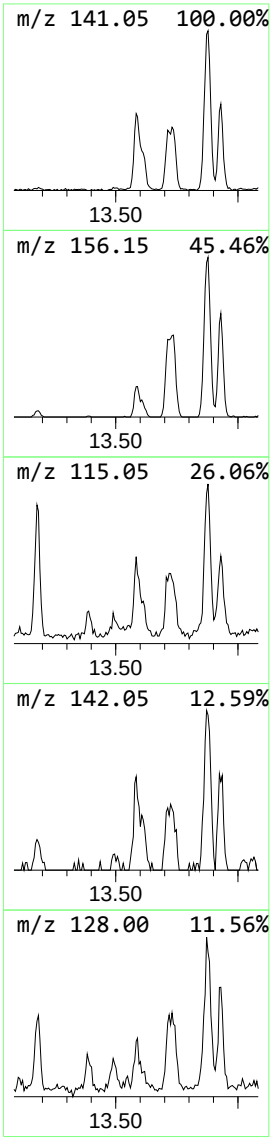
****

Peak Number 7 Naphthalene, 2-ethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.583	2.19 ng	171474	Acenaphthene-d10	14.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	94
2	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	94
3	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	91
4	2,4-Difluoromesitylene	156	C9H10F2	000392-61-0	64
5	2-Thiophenecarboxylic acid, 5-et...	156	C7H8O2S	023229-72-3	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112621\
Data File : BM033273.D
Acq On : 26 Nov 2021 20:05
Operator : CG/JU
Sample : M4817-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A6879

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

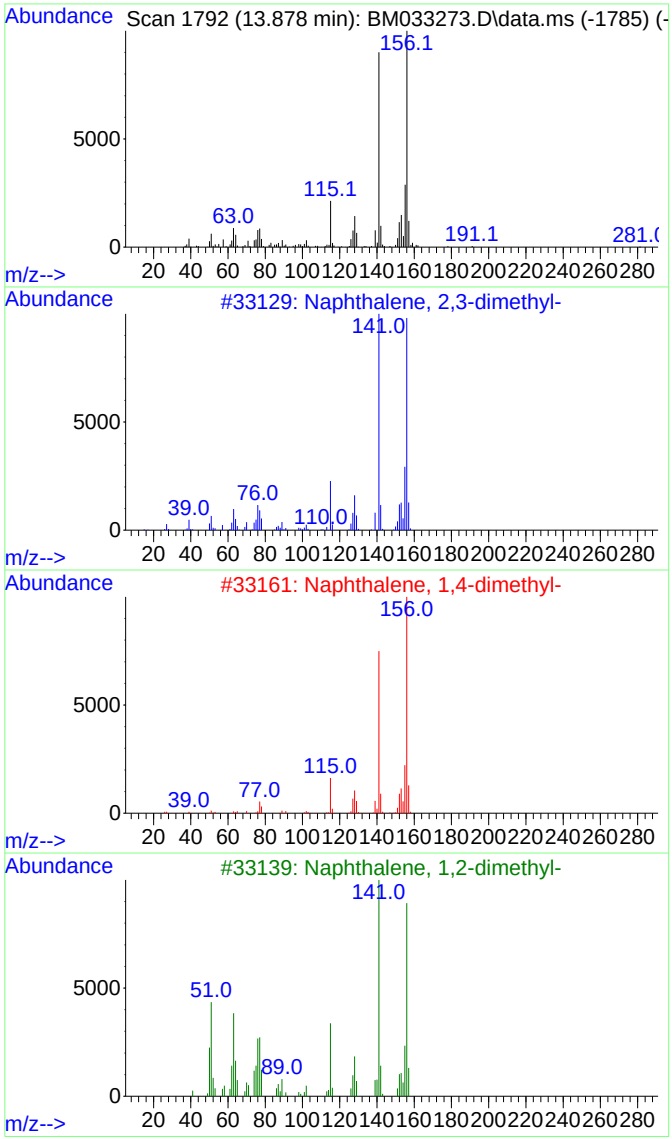
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

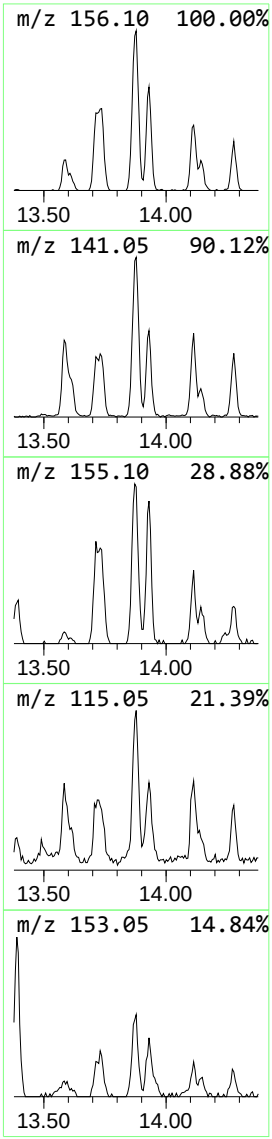
****

Peak Number 8 Naphthalene, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.878	3.83 ng	299944	Acenaphthene-d10	14.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
3	Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	97
4	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96





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Data File : BM033273.D
Acq On : 26 Nov 2021 20:05
Operator : CG/JU
Sample : M4817-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A6879

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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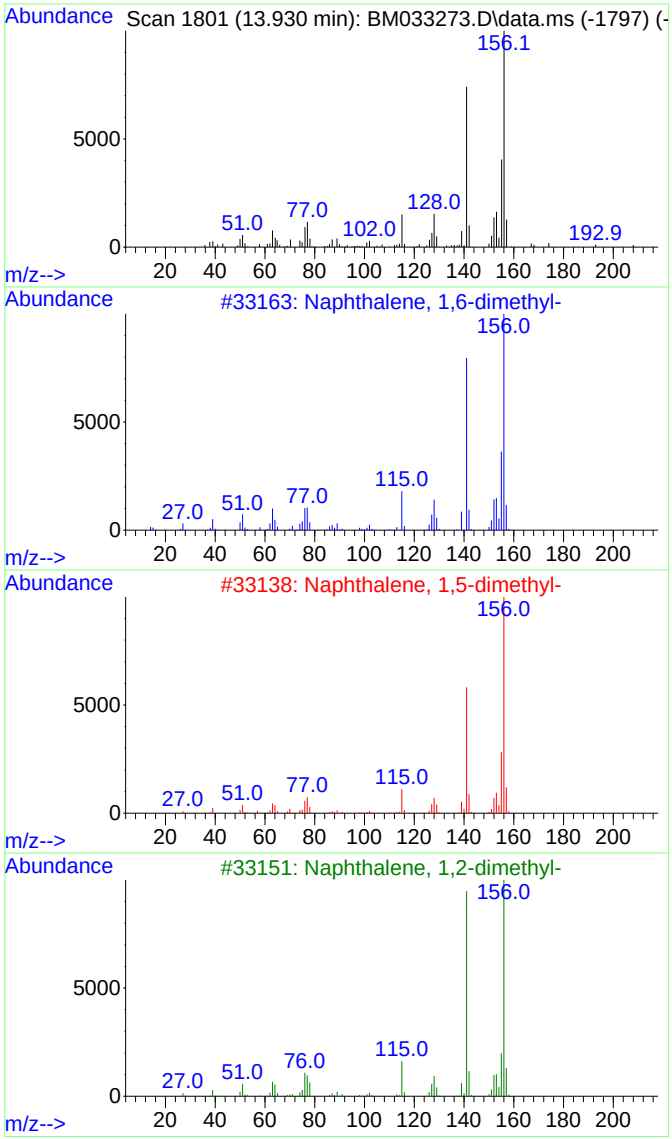
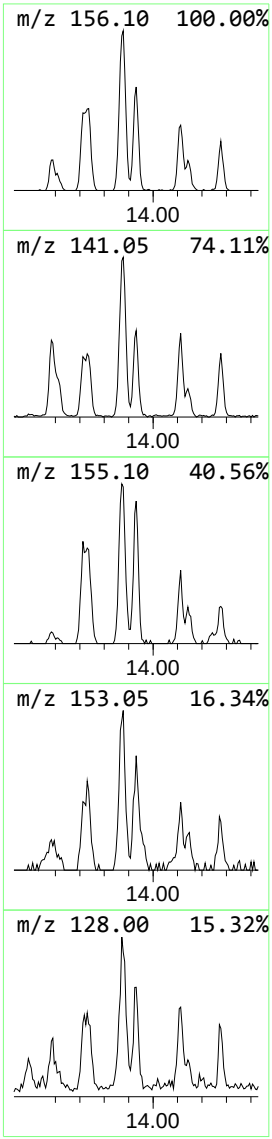
Peak Number 9 Naphthalene, 1,6-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.931	2.75 ng	215655	Acenaphthene-d10	14.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
2	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
3	Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96
4	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96





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Acq On : 26 Nov 2021 20:05
Operator : CG/JU
Sample : M4817-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A6879

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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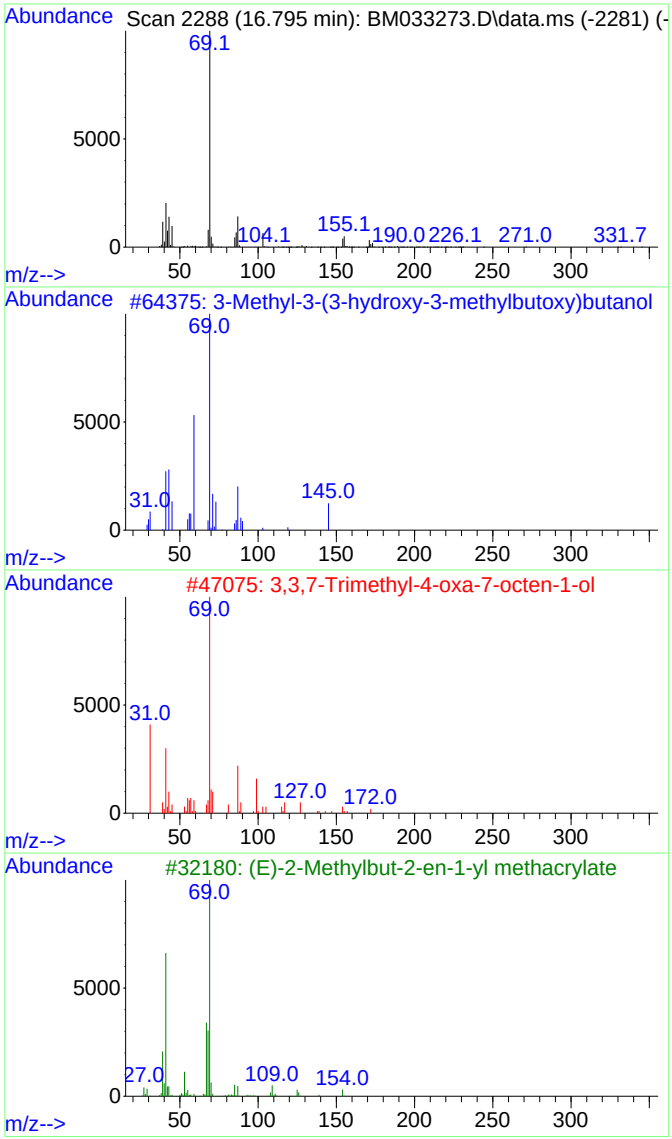
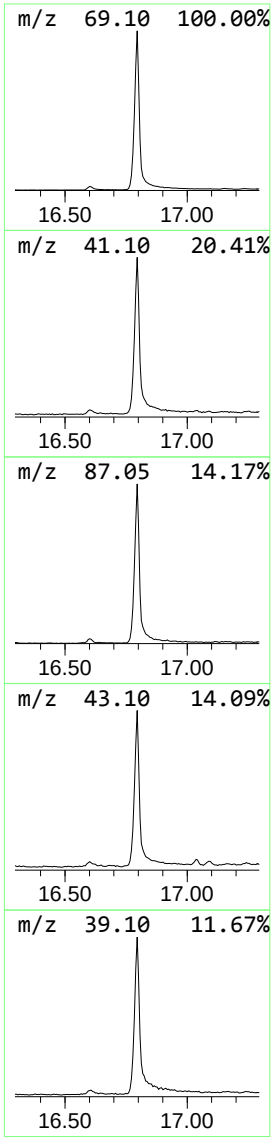
Peak Number 10 unknown16.795 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.795	38.98 ng	3298860	Phenanthrene-d10	17.289

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methyl-3-(3-hydroxy-3-methylbu...	190	C10H22O3	1000432-16-8	40
2	3,3,7-Trimethyl-4-oxa-7-octen-1-ol	172	C10H20O2	069161-47-3	9
3	(E)-2-Methylbut-2-en-1-yl methac...	154	C9H14O2	088142-95-4	9
4	Cyclopropanecarboxylic acid,isob...	142	C8H14O2	064969-79-5	9
5	Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	9





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Operator : CG/JU
Sample : M4817-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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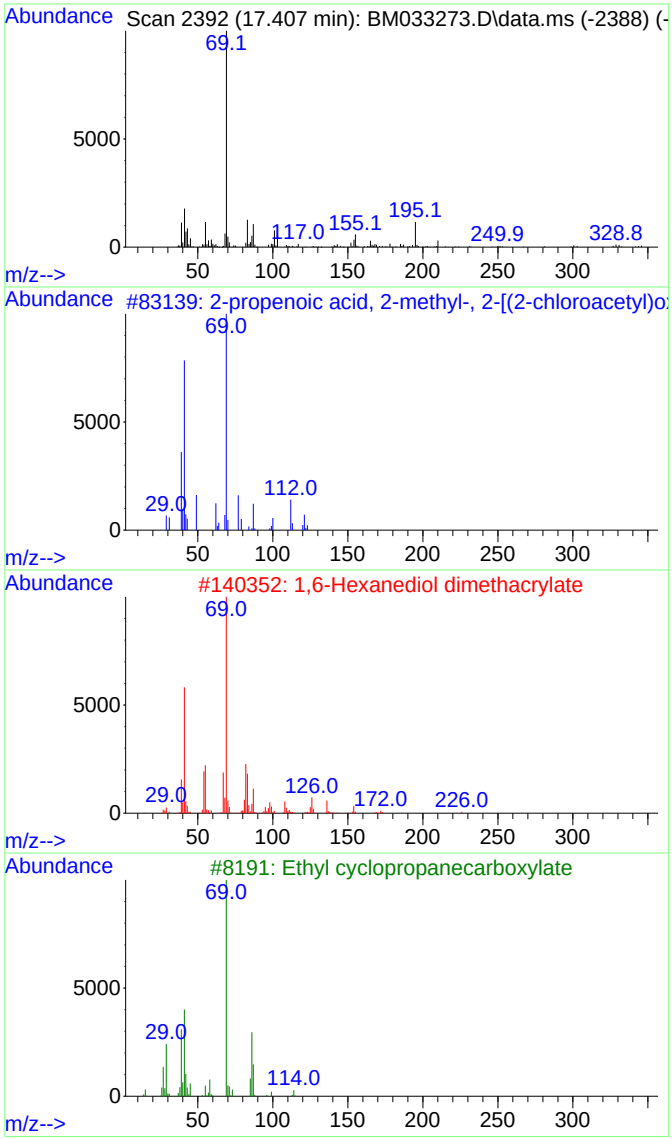
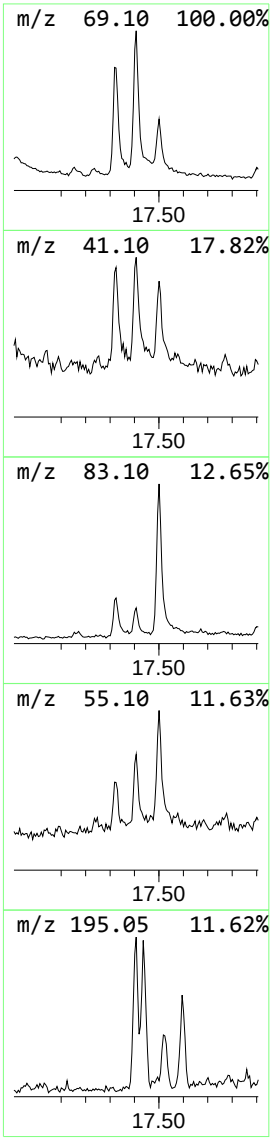
Peak Number 11 unknown17.407 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.407	3.66 ng	309519	Phenanthrene-d10	17.289

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-propenoic acid, 2-methyl-, 2-[...	206	C8H11ClO4	1000400-53-0	42
2	1,6-Hexanediol dimethacrylate	254	C14H22O4	006606-59-3	38
3	Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	36
4	2-Propenoic acid, 2-methyl-, 2-e...	338	C18H26O6	003290-92-4	33
5	5-Thia-10-azaspiro[3.6]decane-7-...	402	C15H16F6N2O2S	1010259-93-7	28





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Acq On : 26 Nov 2021 20:05
Operator : CG/JU
Sample : M4817-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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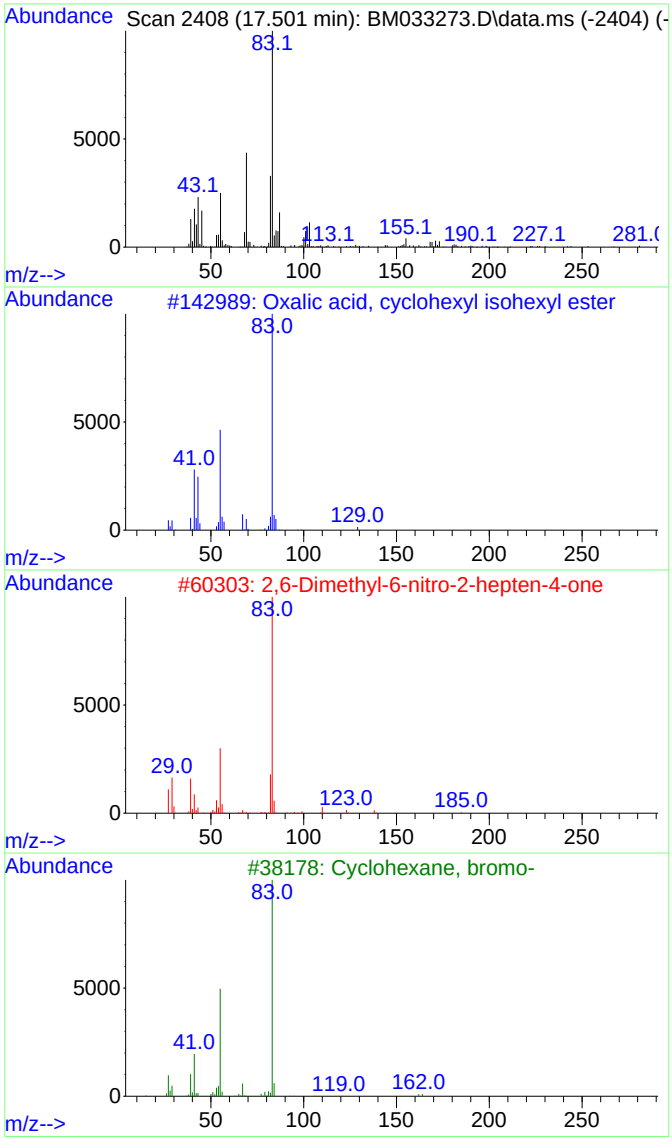
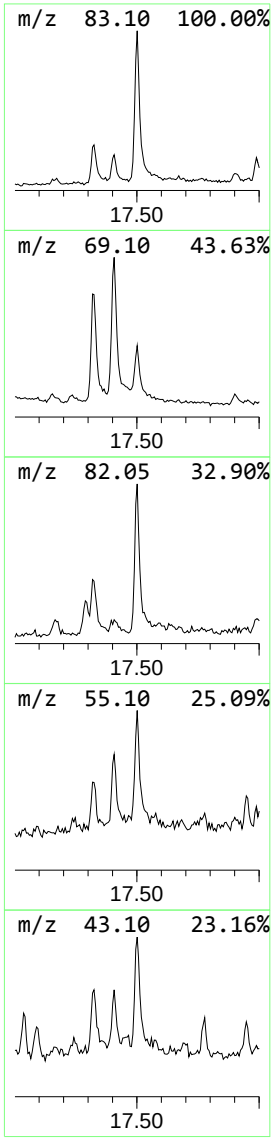
Peak Number 12 unknown17.501 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.501	3.76 ng	317806	Phenanthrene-d10	17.289

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxalic acid, cyclohexyl isohexyl...	256	C14H24O4	1010309-30-7	47
2	2,6-Dimethyl-6-nitro-2-hepten-4-one	185	C9H15NO3	073583-56-9	47
3	Cyclohexane, bromo-	162	C6H11Br	000108-85-0	43
4	2-Butenoic acid, 3-methyl-, ethy...	128	C7H12O2	000638-10-8	40
5	Neopentyl (E)-2-methylbut-2-enoate	170	C10H18O2	1000373-71-0	40





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Acq On : 26 Nov 2021 20:05
Operator : CG/JU
Sample : M4817-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A6879

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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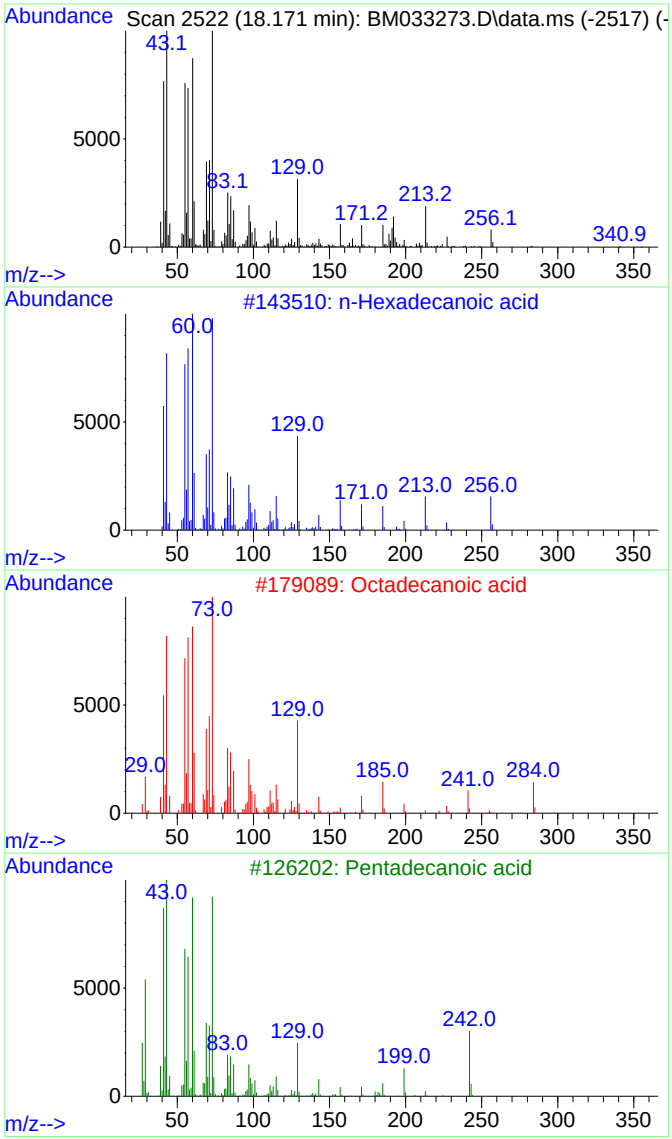
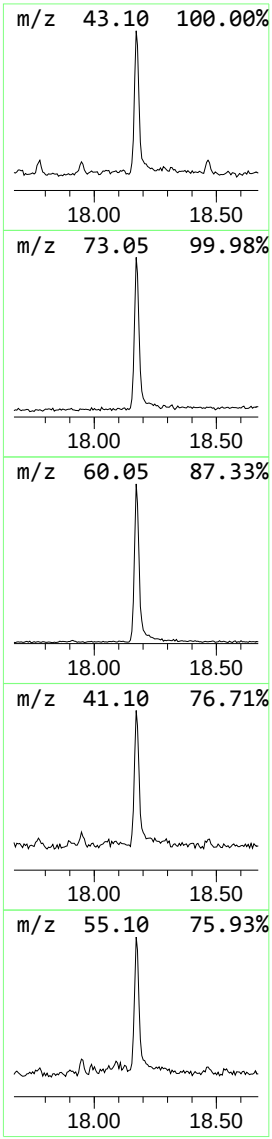
Peak Number 13 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.171	7.39 ng	625425	Phenanthrene-d10	17.289

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	90
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4	Tridecanoic acid	214	C13H26O2	000638-53-9	80
5	Octanoic acid	144	C8H16O2	000124-07-2	45





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112621\
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Acq On : 26 Nov 2021 20:05
Operator : CG/JU
Sample : M4817-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A6879

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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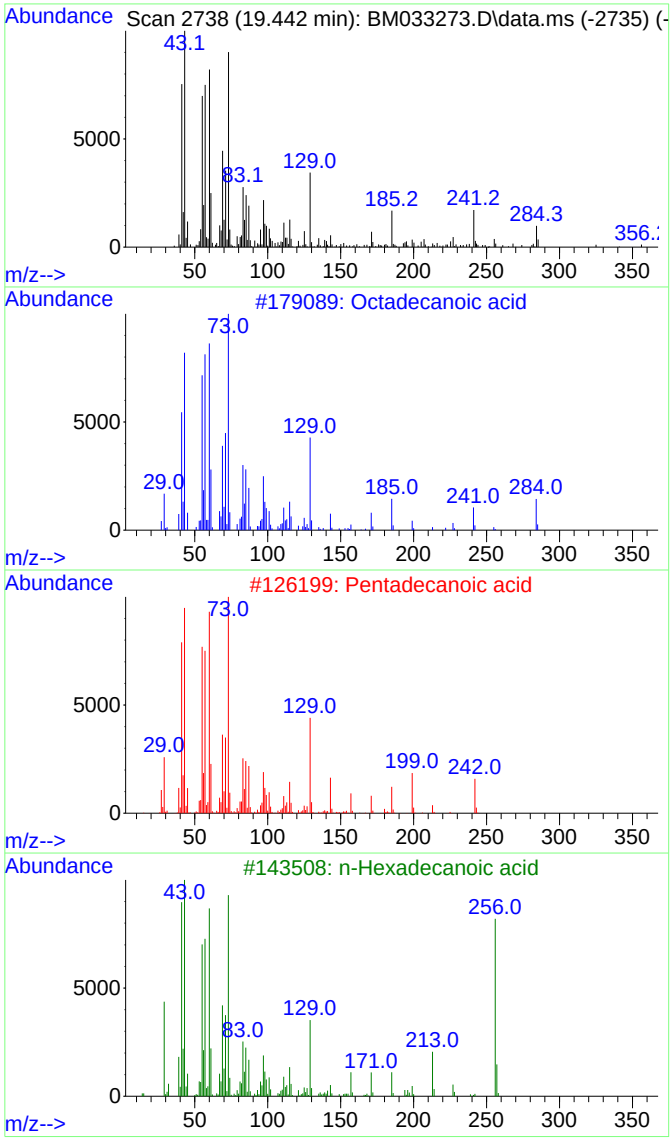
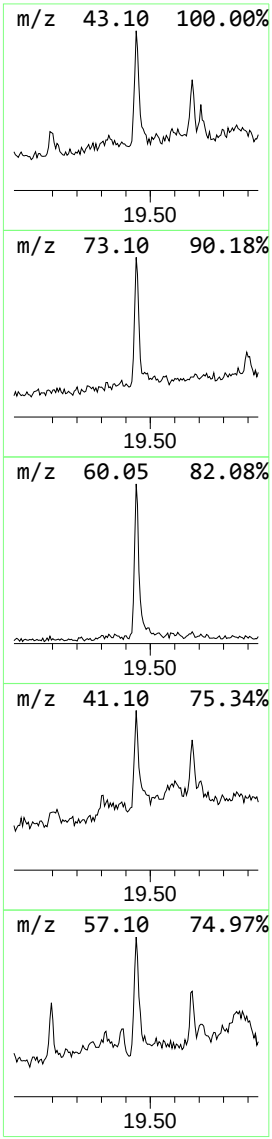
Peak Number 14 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.442	3.12 ng	284195	Chrysene-d12	21.453

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
4	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	91
5	Tridecanoic acid	214	C13H26O2	000638-53-9	90





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Data File : BM033273.D
Acq On : 26 Nov 2021 20:05
Operator : CG/JU
Sample : M4817-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A6879

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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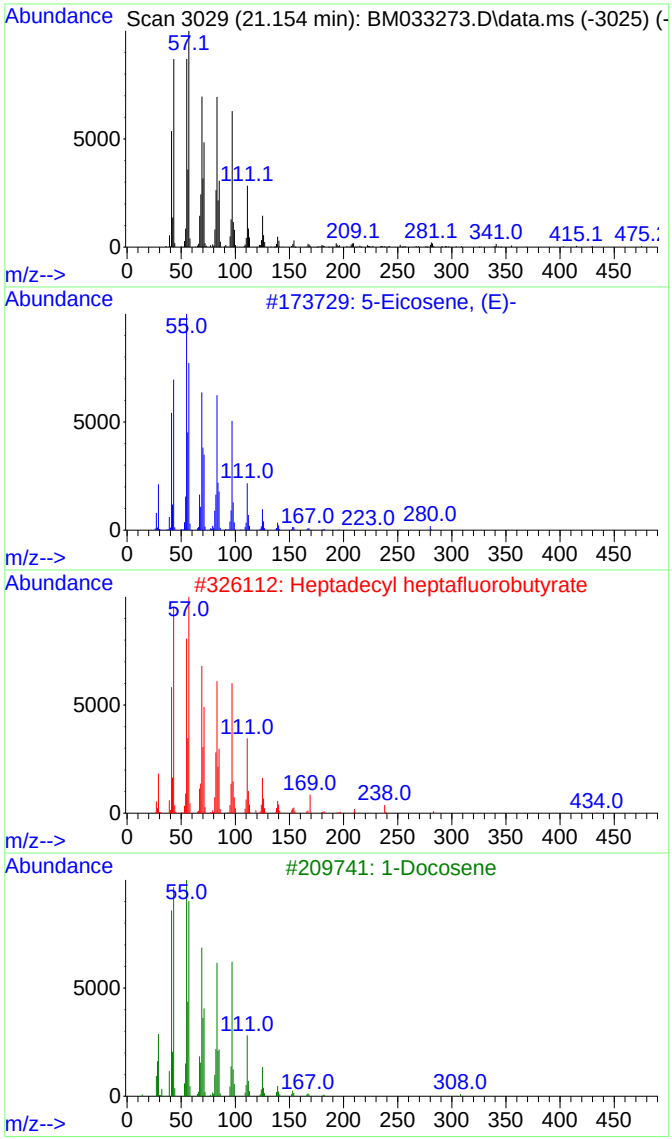
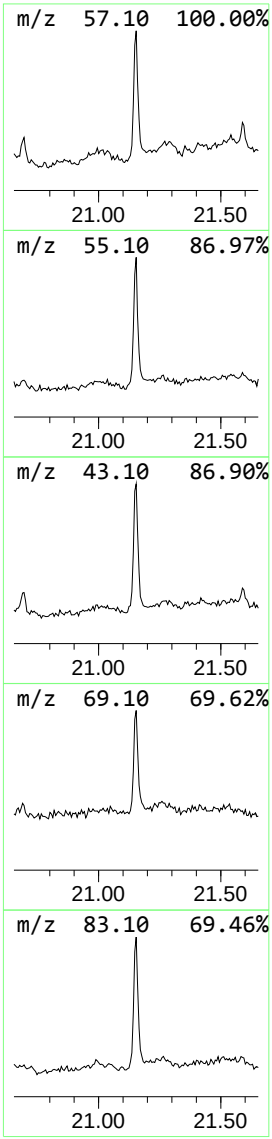
Peak Number 15 5-Eicosene, (E)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.154	6.01 ng	546620	Chrysene-d12	21.453

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	96
2	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3	1-Docosene	308	C22H44	001599-67-3	94
4	Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	94
5	Cyclotridecane	182	C13H26	000295-02-3	93





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112621\
Data File : BM033273.D
Acq On : 26 Nov 2021 20:05
Operator : CG/JU
Sample : M4817-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
2-Pentanone, 4-...	5.055	2.1	ng	85133	1	7.931	827496	20.0
(E)-1-Phenyl-1-...	10.125	2.0	ng	117836	2	10.731	1171540	20.0
Ethanol, 2-(2-b...	10.501	4.7	ng	274357	2	10.731	1171540	20.0
Cyclopropanecar...	11.983	2.6	ng	150556	2	10.731	1171540	20.0
unknown13.354	13.354	8.4	ng	658390	3	14.554	1568160	20.0
3-Methyl-2-bute...	13.519	9.1	ng	709606	3	14.554	1568160	20.0
Naphthalene, 2-...	13.583	2.2	ng	171474	3	14.554	1568160	20.0
Naphthalene, 2,...	13.878	3.8	ng	299944	3	14.554	1568160	20.0
Naphthalene, 1,...	13.931	2.8	ng	215655	3	14.554	1568160	20.0
unknown16.795	16.795	39.0	ng	3298860	4	17.289	1692530	20.0
unknown17.407	17.407	3.7	ng	309519	4	17.289	1692530	20.0
unknown17.501	17.501	3.8	ng	317806	4	17.289	1692530	20.0
n-Hexadecanoic ...	18.171	7.4	ng	625425	4	17.289	1692530	20.0
Octadecanoic acid	19.442	3.1	ng	284195	5	21.453	1819190	20.0
5-Eicosene, (E)-	21.154	6.0	ng	546620	5	21.453	1819190	20.0