LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052225\
Data File : BP024763.D
Acq On : 22 May 2025 17:43
Operator : RC/JU
Sample : Q2097-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
RT3419

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_P\Methods\8270E-BP051325.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP024763.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.822	288	295	307	rBV	101790	155437	2.95%	0.390%
2	5.287	367	374	397	rBV	1114123	1767901	33.51%	4.430%
3	6.857	633	641	665	rBV	955357	1808181	34.27%	4.531%
4	7.651	768	776	791	rBV	458944	721144	13.67%	1.807%
5	8.804	964	972	991	rBV	637442	1198862	22.72%	3.004%
6	10.422	1237	1247	1263	rBV	546929	1014191	19.22%	2.541%
7	11.510	1417	1432	1440	rBV	33480	152625	2.89%	0.382%
8	12.157	1537	1542	1548	rBV3	55891	109839	2.08%	0.275%
9	12.616	1617	1620	1627	rBV7	66604	156445	2.96%	0.392%
10	12.704	1630	1635	1641	rVB3	149279	249242	4.72%	0.625%
11	12.769	1642	1646	1649	rBV5	67424	123373	2.34%	0.309%
12	12.892	1657	1667	1676	rBV	2166598	3379204	64.04%	8.468%
13	13.057	1687	1695	1702	rBV3	280917	717257	13.59%	1.797%
14	13.698	1800	1804	1808	rBV	942390	1289913	24.45%	3.232%
15	13.769	1813	1816	1819	rBV2	265210	321928	6.10%	0.807%
16	14.022	1854	1859	1862	rBV5	630280	1093565	20.73%	2.740%
17	14.110	1870	1874	1882	rVB	1792219	2866130	54.32%	7.182%
18	14.186	1882	1887	1891	rBV3	429463	904225	17.14%	2.266%
19	14.251	1893	1898	1900	rBV	827655	1363177	25.84%	3.416%
20	14.286	1900	1904	1913	rVB	1215223	2093172	39.67%	5.245%
21	14.822	1991	1995	2000	rBV2	1005495	1519098	28.79%	3.807%
22	15.086	2035	2040	2044	rBV2	1984663	3261361	61.81%	8.173%
23	15.351	2080	2085	2086	rBV2	704585	1142685	21.66%	2.863%
24	15.804	2160	2162	2168	rVB2	1630056	2163980	41.01%	5.423%
25	16.416	2264	2266	2269	rBV2	986740	1186433	22.49%	2.973%
26	19.804	2838	2842	2849	rVB	3662884	5276447	100.00%	13.222%
27	21.510	3128	3132	3137	rVB2	1100509	1755454	33.27%	4.399%
28	24.780	3681	3688	3708	rVB	748192	2114450	40.07%	5.299%

Sum of corrected areas: 39905719

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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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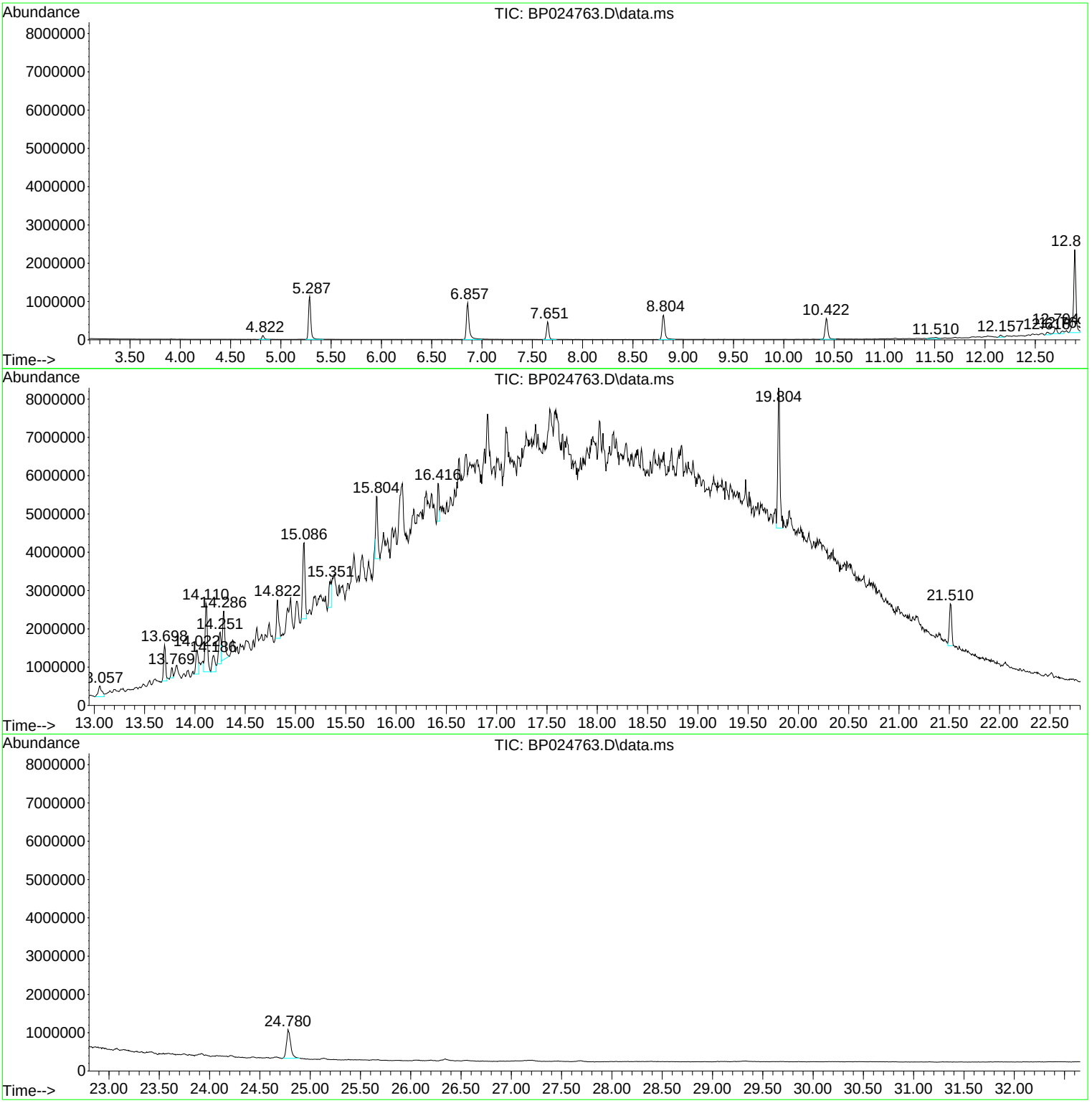
ClientSampleId :
RT3419

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.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_P\Data\BP052225\
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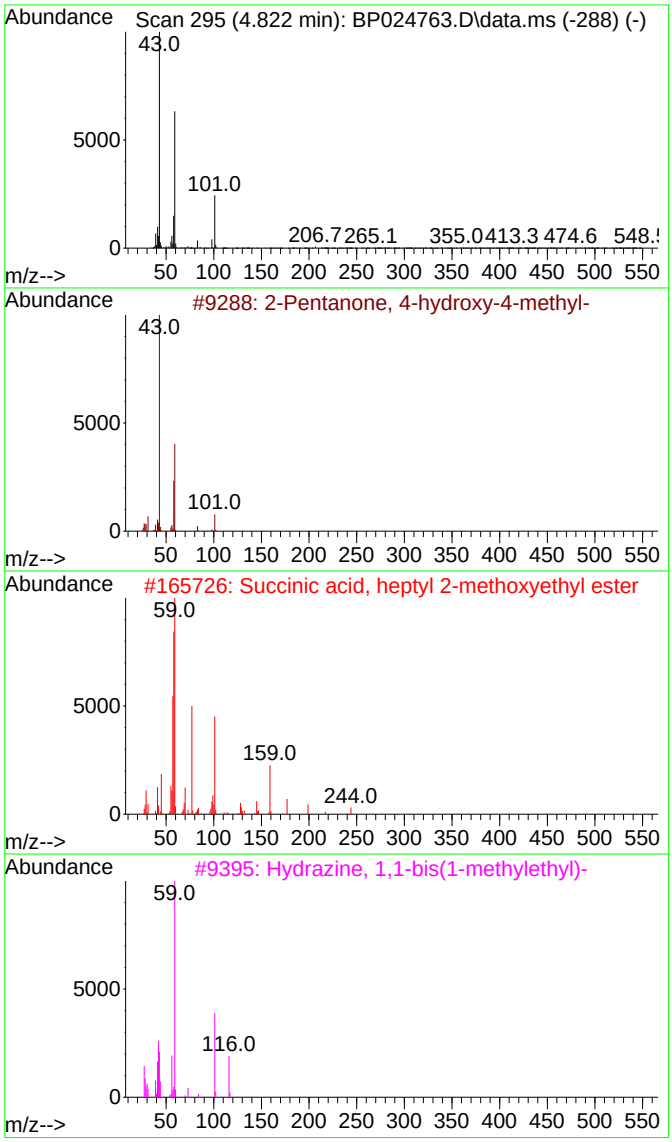
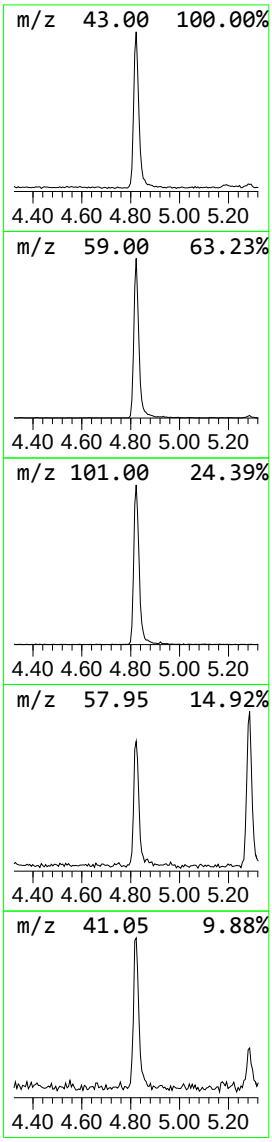
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.822	4.31 ng	155437	1,4-Dichlorobenzene-d4	7.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	Succinic acid, heptyl 2-methoxye...	274	C14H26O5	1000325-80-7	33
3	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	28
4	n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	28
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25





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Data File : BP024763.D
Acq On : 22 May 2025 17:43
Operator : RC/JU
Sample : Q2097-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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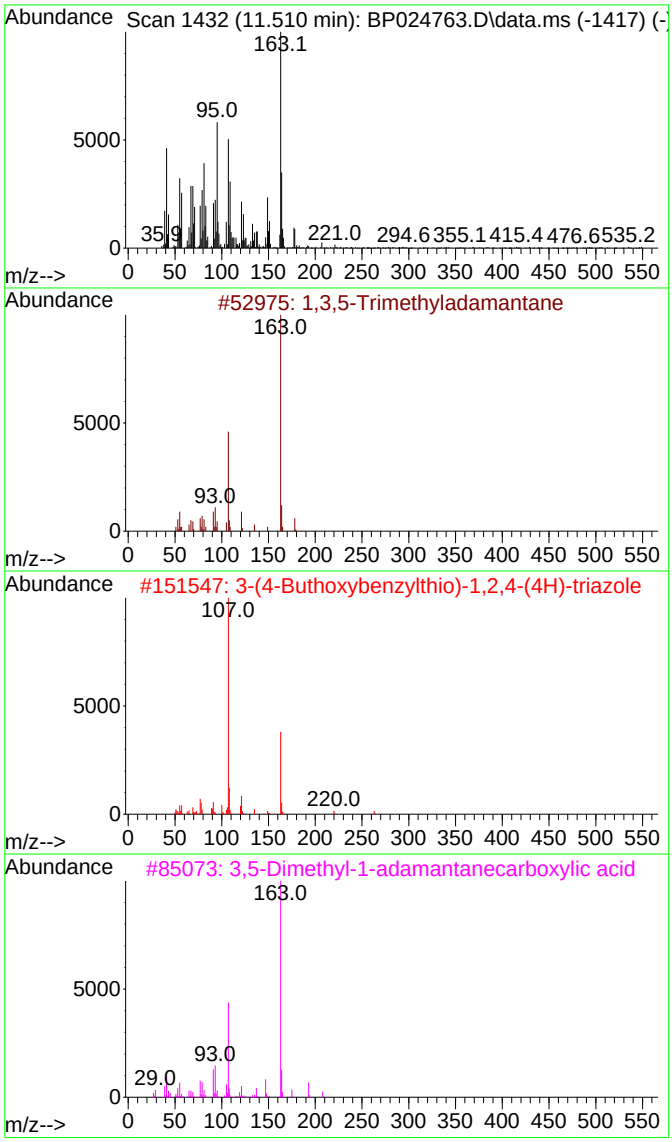
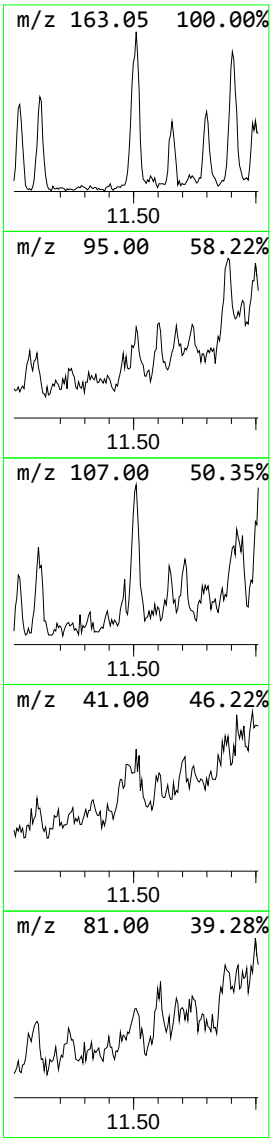
****

Peak Number 2 unknown11.510 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.510	3.01 ng	152625	Naphthalene-d8	10.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,5-Trimethyladamantane	178	C13H22	000707-35-7	47
2	3-(4-Buthoxybenzylthio)-1,2,4-(4...	263	C13H17N3OS	057736-49-9	43
3	3,5-Dimethyl-1-adamantanecarboxy...	208	C13H20O2	014670-94-1	37
4	1,5-Dimethylbicyclo(3.2.2)non-3-...	164	C11H16O	1000426-97-9	30
5	Pyrido[2,3-d]pyridazine-5(6H)-th...	163	C7H5N3S	015370-85-1	27





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Operator : RC/JU
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BNA_P
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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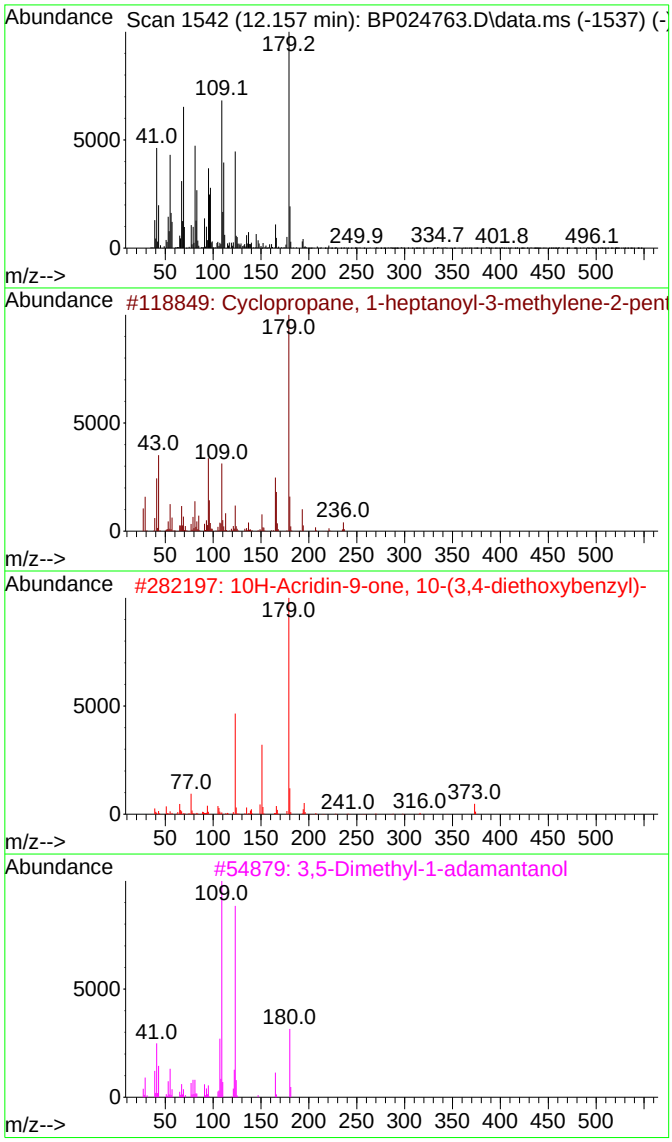
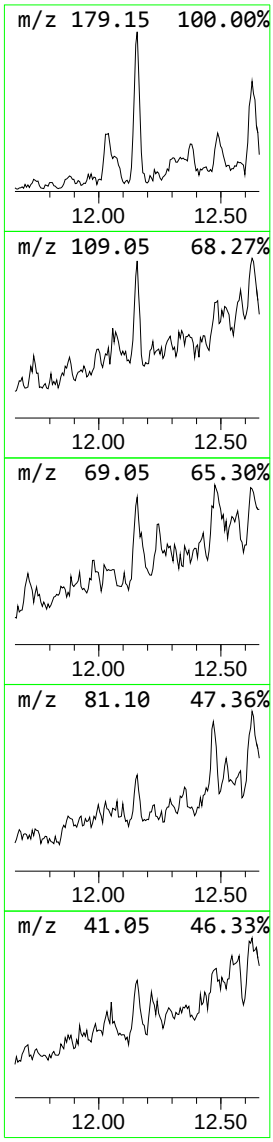
Peak Number 3 unknown12.157 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.157	2.17 ng	109839	Naphthalene-d8	10.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1-heptanoyl-3-meth...	236	C16H28O	1000157-26-0	38
2	10H-Acridin-9-one, 10-(3,4-dieth...	373	C24H23NO3	1000302-73-5	35
3	3,5-Dimethyl-1-adamantanol	180	C12H20O	000707-37-9	25
4	pyridine, 4-(2-phenylethynyl)-	179	C13H9N	1000404-81-1	14
5	4,7(1H,8H)-Pteridinedione, 2-amino-	179	C6H5N5O2	000529-69-1	14





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Operator : RC/JU
Sample : Q2097-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
RT3419

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

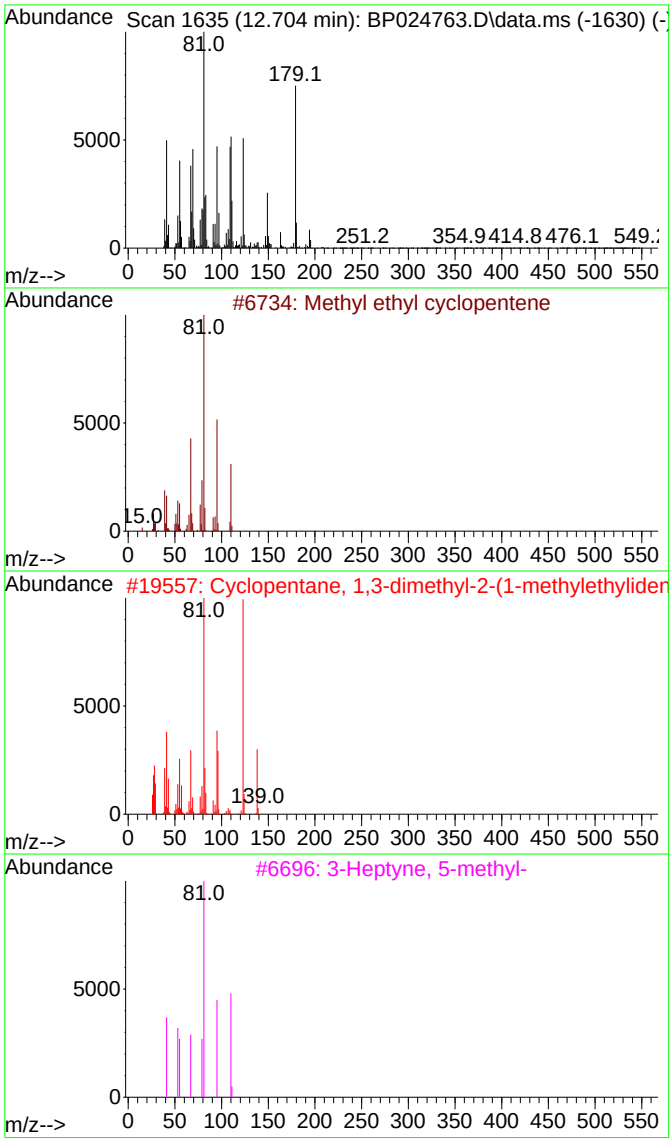
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

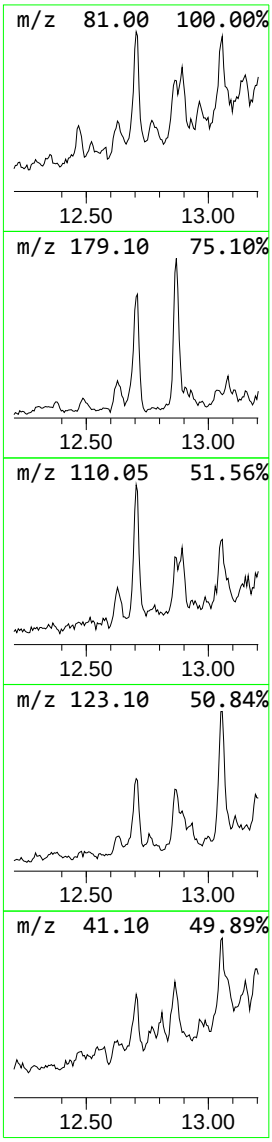
****

Peak Number 4 unknown12.704 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.704	2.38 ng	249242	Acenaphthene-d10	14.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl ethyl cyclopentene	110	C8H14	019780-56-4	45
2	Cyclopentane, 1,3-dimethyl-2-(1-...	138	C10H18	061142-30-1	43
3	3-Heptyne, 5-methyl-	110	C8H14	061228-09-9	38
4	Cyclohexane, ethylidene-	110	C8H14	001003-64-1	35
5	Trichloroacetic acid, 2-ethylcyc...	272	C10H15Cl3O2	1000282-57-3	27





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Operator : RC/JU
Sample : Q2097-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
RT3419

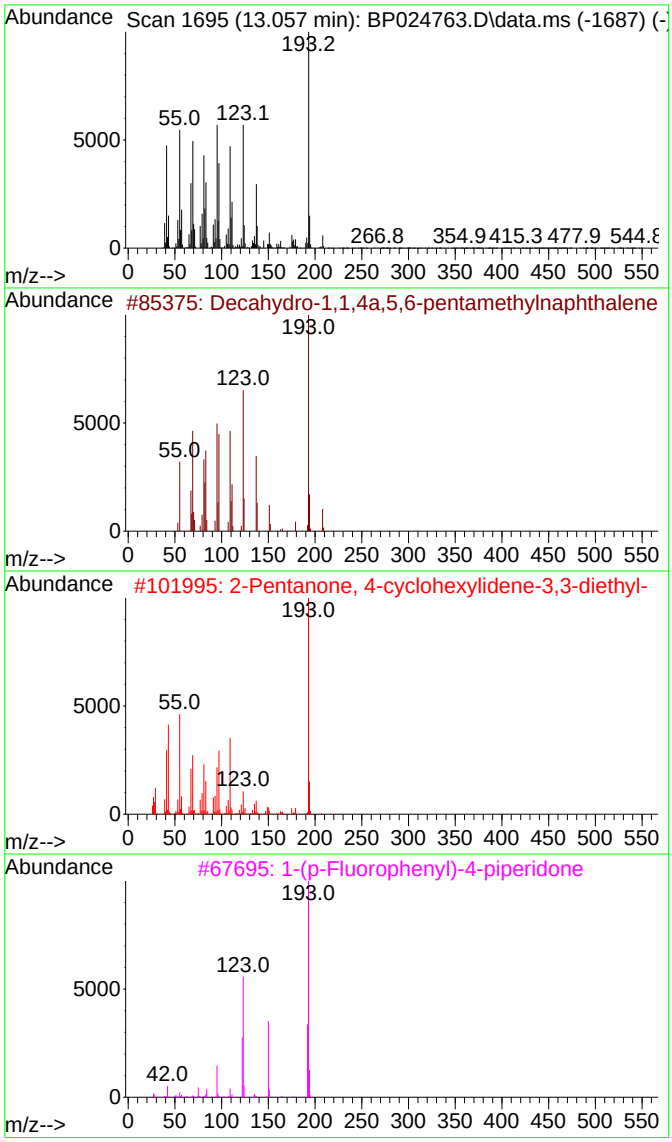
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

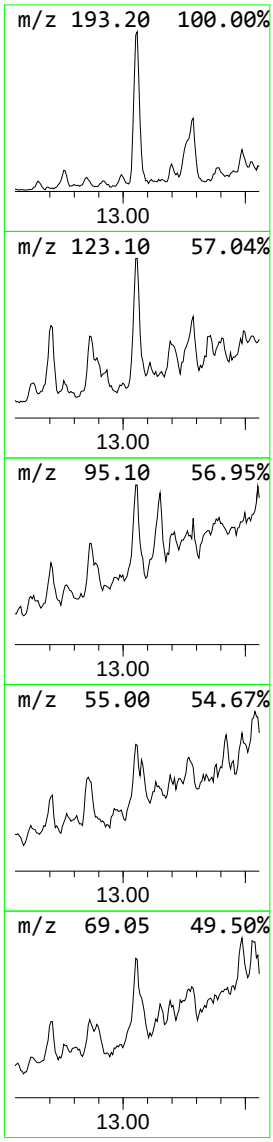
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Decahydro-1,1,4a,5,6-pentam... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.057	6.85 ng	717257	Acenaphthene-d10	14.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	55
2	2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	52
3	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	38
4	N-(4-Bromo-2-chlorophenyl)-2-[4-...	425	C17H17BrClN3O3	1000482-48-9	35
5	N-(2,5-Dichlorophenyl)-2-[4-(fur...	381	C17H17Cl2N3O3	1010482-48-3	35





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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
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TIC Library : C:\Database\NIST20.L
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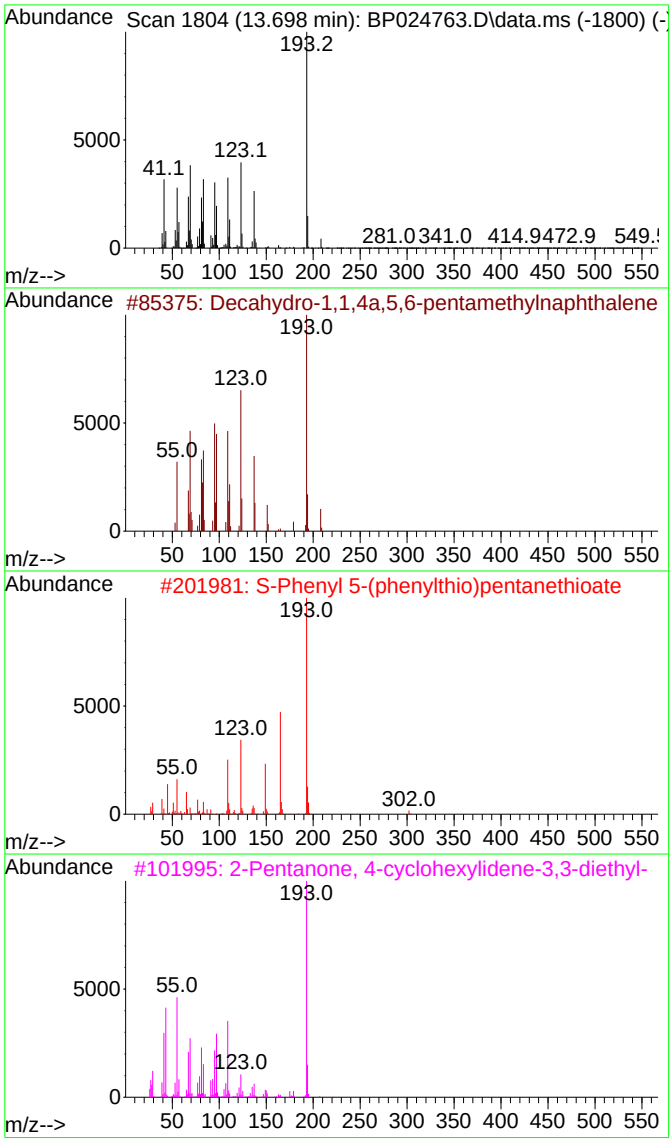
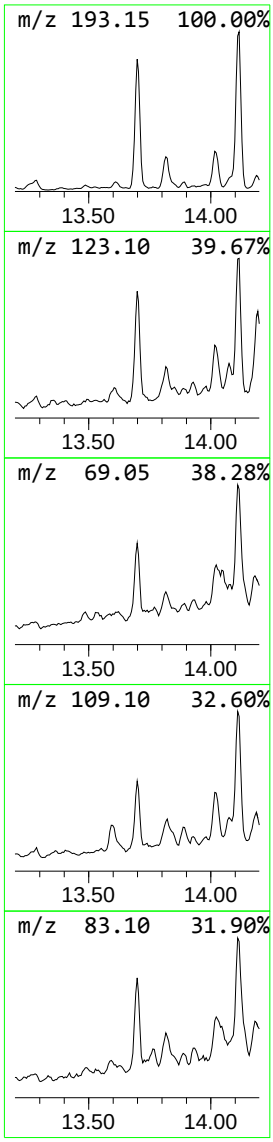
Peak Number 6 unknown13.698 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.698	12.32 ng	1289910	Acenaphthene-d10	14.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	64
2	S-Phenyl 5-(phenylthio)pentaneth...	302	C17H18OS2	1000234-40-7	38
3	2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	28
4	Iminostilbene	193	C14H11N	000256-96-2	27
5	9-Phenanthrenamine	193	C14H11N	000947-73-9	27





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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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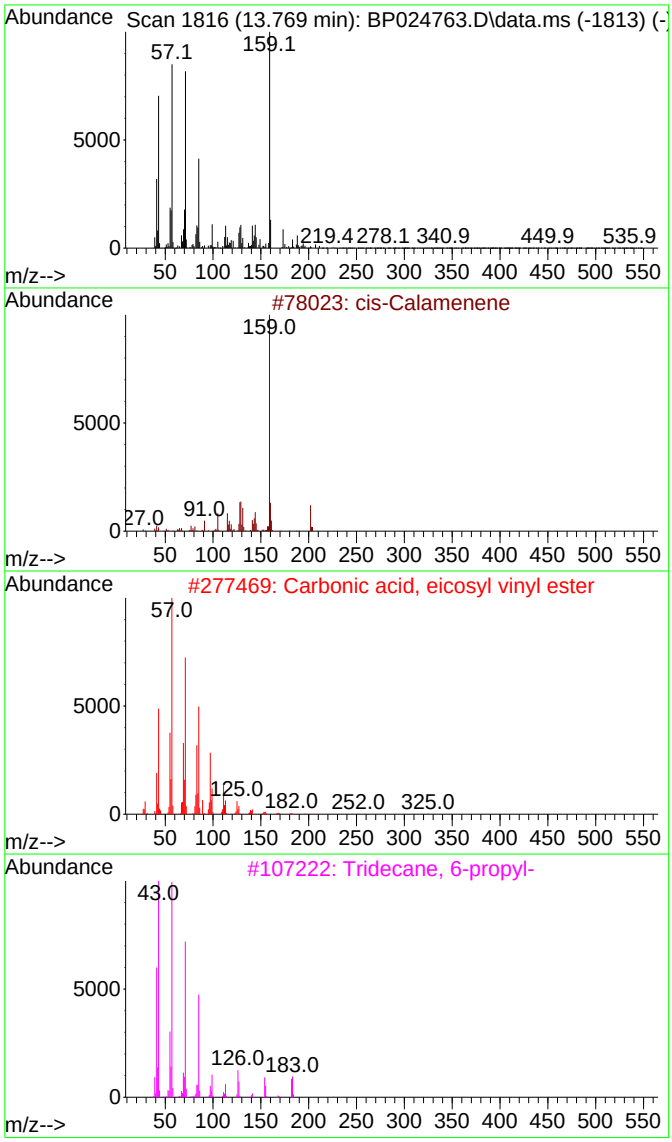
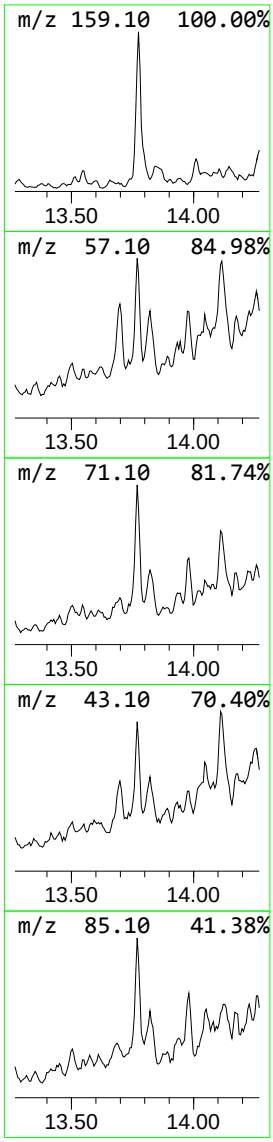
****

Peak Number 7 unknown13.769 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.769	3.08 ng	321928	Acenaphthene-d10	14.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	cis-Calamenene	202	C15H22	072937-55-4	46
2	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	43
3	Tridecane, 6-propyl-	226	C16H34	055045-10-8	43
4	Dodecane, 4-methyl-	184	C13H28	006117-97-1	43
5	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	43





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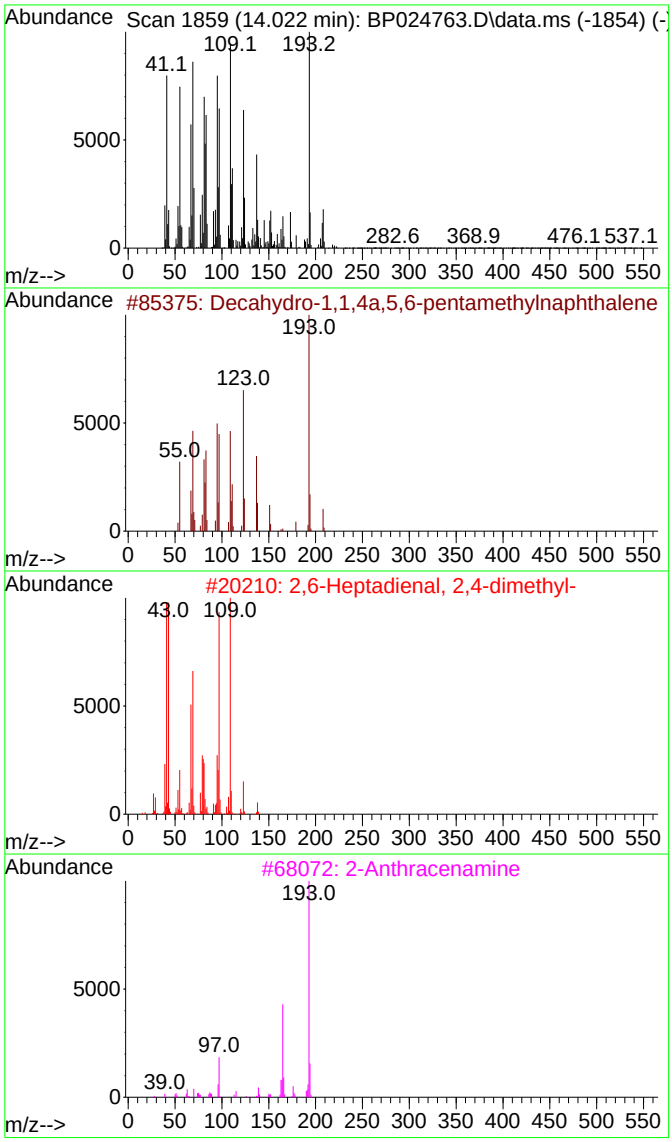
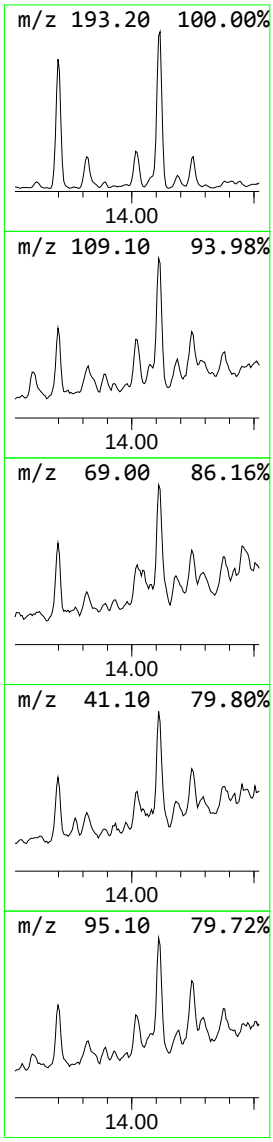
****

Peak Number 8 unknown14.022 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.022	10.45 ng	1093570	Acenaphthene-d10	14.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	62
2	2,6-Heptadienal, 2,4-dimethyl-	138	C9H14O	085136-08-9	46
3	2-Anthracenamine	193	C14H11N	000613-13-8	43
4	Silane, chlorodiethyl(2-methylpe...	222	C10H23ClOSi	1000363-79-1	38
5	Silane, chlorodiethyl(2-methylpe...	222	C10H23ClOSi	1000363-12-7	38





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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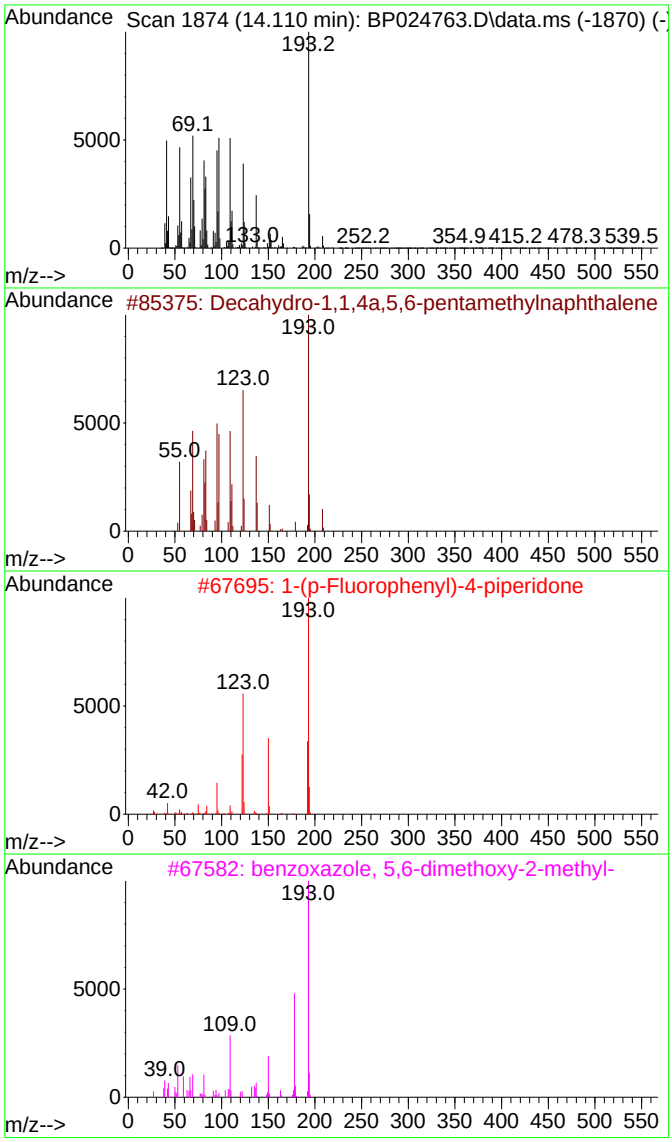
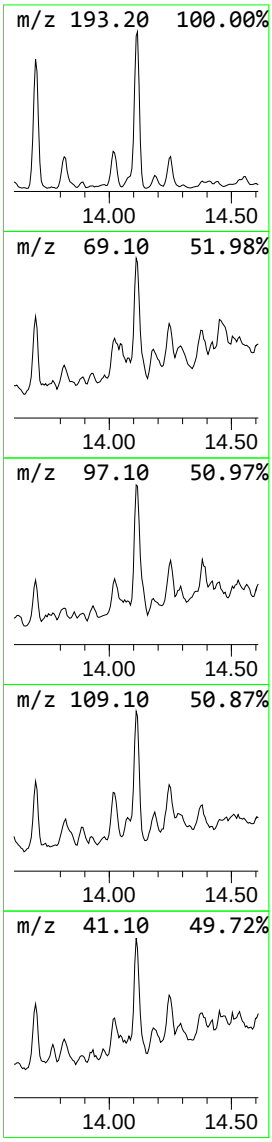
Peak Number 9 unknown14.110 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.110	27.39 ng	2866130	Acenaphthene-d10	14.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	50
2	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	43
3	benzoxazole, 5,6-dimethoxy-2-met...	193	C10H11NO3	1000395-98-9	22
4	Butan-2-one, 1-[3-(3,3-dimethyl-...	278	C16H26N2O2	1000303-34-2	22
5	3-Cyclohexene-1-carboxaldehyde, ...	152	C10H16O	040702-26-9	22





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Data File : BP024763.D
Acq On : 22 May 2025 17:43
Operator : RC/JU
Sample : Q2097-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
RT3419

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

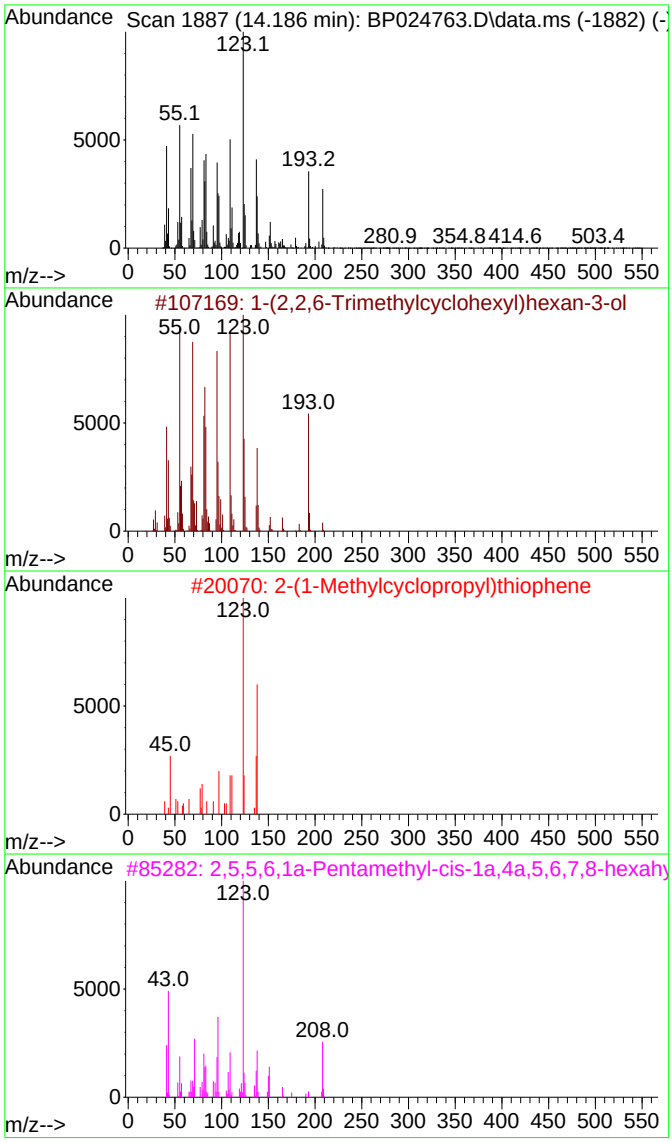
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

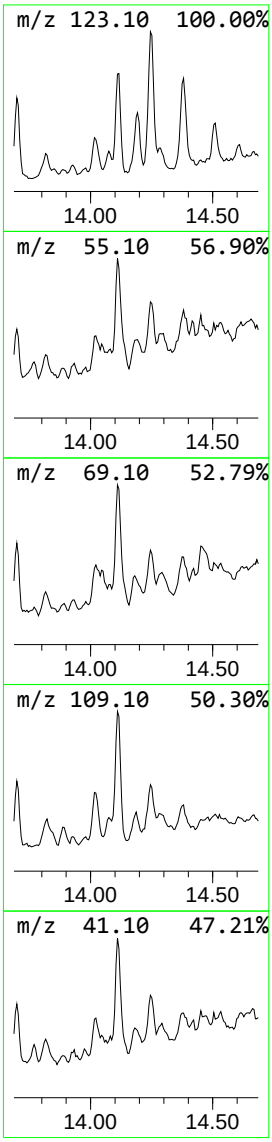
****

Peak Number 10 unknown14.186 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.186	8.64 ng	904225	Acenaphthene-d10	14.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(2,2,6-Trimethylcyclohexyl)hex...	226	C15H30O	070788-30-6	49
2	2-(1-Methylcyclopropyl)thiophene	138	C8H10S	1000100-02-7	38
3	2,5,5,6,1a-Pentamethyl-cis-1a,4a...	208	C14H24O	1010215-77-9	38
4	Cyclopentene, 1,3-dimethyl-2-(1-...	138	C10H18	061142-32-3	38
5	1,3-Benzenediamine, 4-methoxy-	138	C7H10N2O	000615-05-4	38





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Data File : BP024763.D
Acq On : 22 May 2025 17:43
Operator : RC/JU
Sample : Q2097-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
RT3419

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

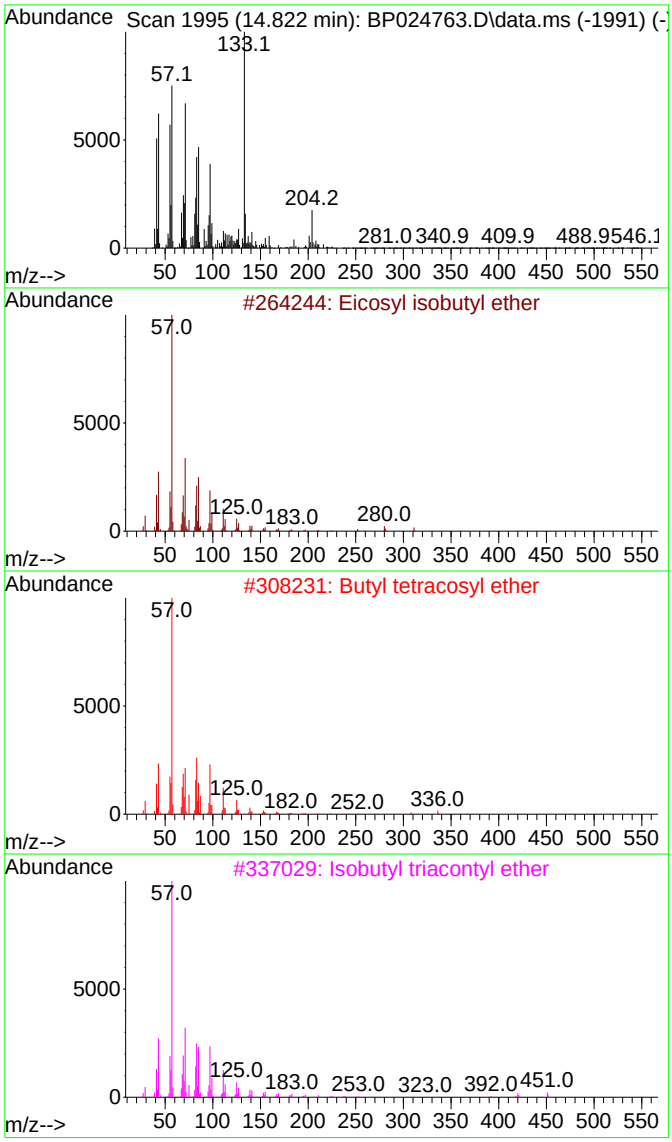
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

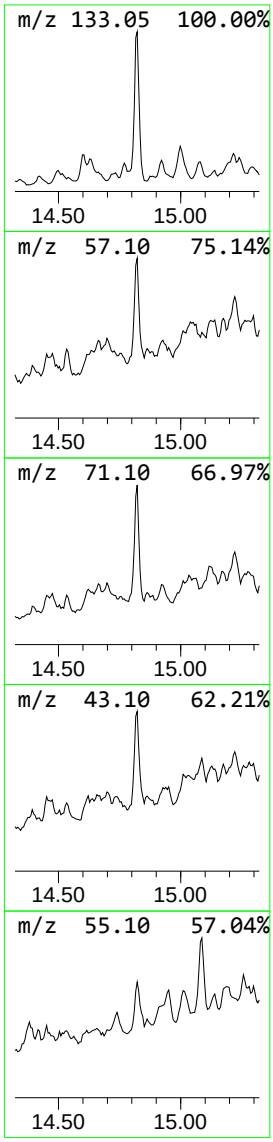
****

Peak Number 11 unknown14.822 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.822	14.51 ng	1519100	Acenaphthene-d10	14.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosyl isobutyl ether	354	C24H50O	1000406-33-0	46
2	Butyl tetracosyl ether	410	C28H58O	1000406-40-9	35
3	Isobutyl triacontyl ether	495	C34H70O	1000406-33-5	35
4	Isobutyl octacosyl ether	467	C32H66O	1000406-33-4	35
5	Tritetracontane	605	C43H88	007098-21-7	30





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Data File : BP024763.D
Acq On : 22 May 2025 17:43
Operator : RC/JU
Sample : Q2097-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
RT3419

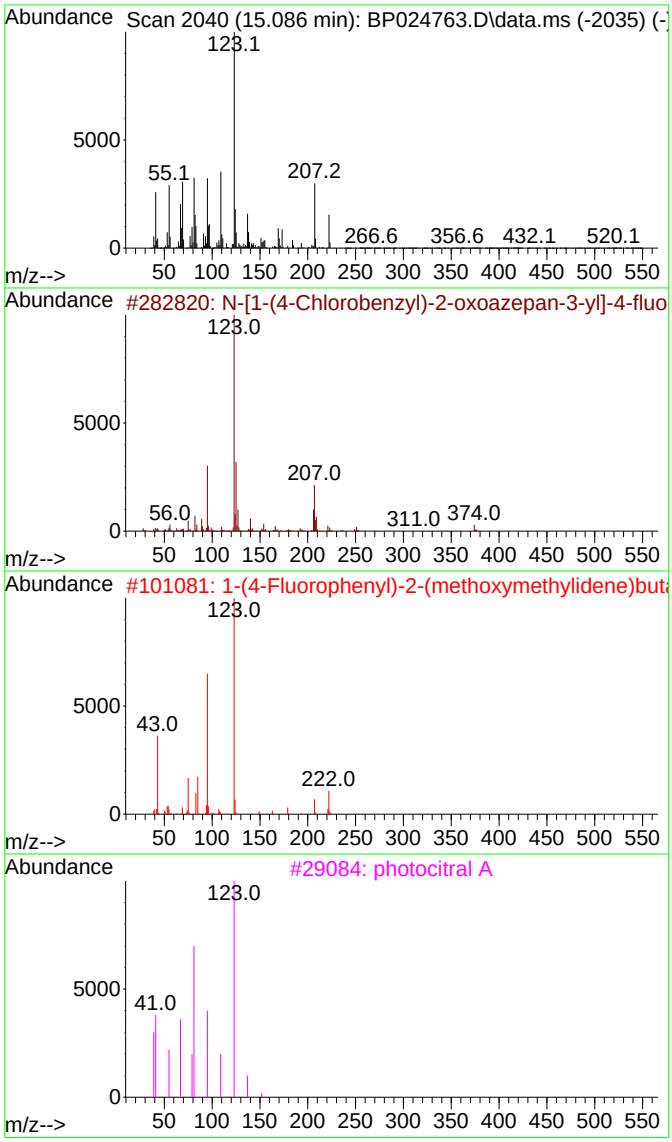
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

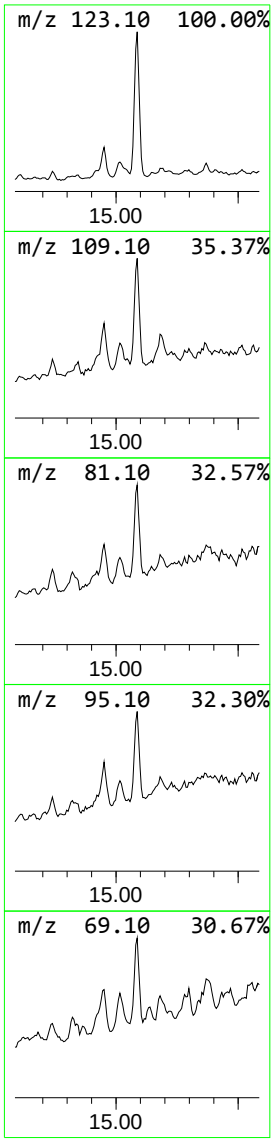
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 N-[1-(4-Chlorobenzyl)-2-oxo... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.086	31.16 ng	3261360	Acenaphthene-d10	14.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N-[1-(4-Chlorobenzyl)-2-oxoazepa...	374	C20H20ClFN2O2	1000481-70-9	50
2	1-(4-Fluorophenyl)-2-(methoxymet...	222	C12H11FO3	1000388-14-4	50
3	photocitral A	152	C10H16O	055253-28-6	50
4	Pentanamide, N-(4-methoxyphenyl)-	207	C12H17NO2	1000306-89-3	49
5	1-Butanone, 1-bicyclo[4.1.0]hept...	166	C11H18O	054764-62-4	47





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Data File : BP024763.D
Acq On : 22 May 2025 17:43
Operator : RC/JU
Sample : Q2097-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
RT3419

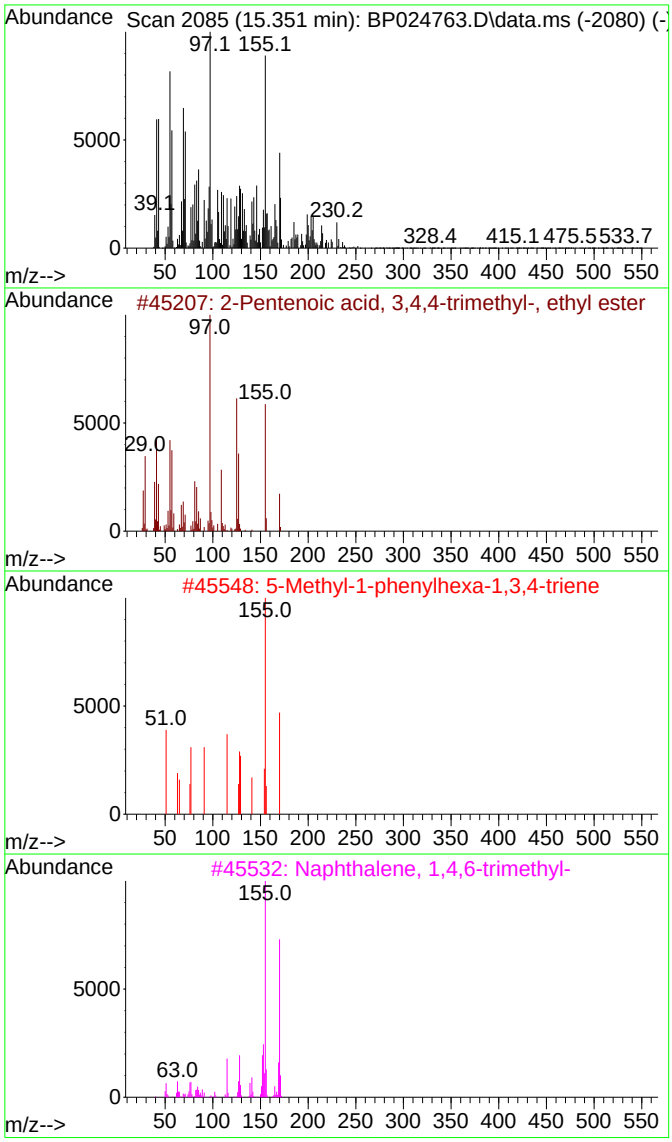
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

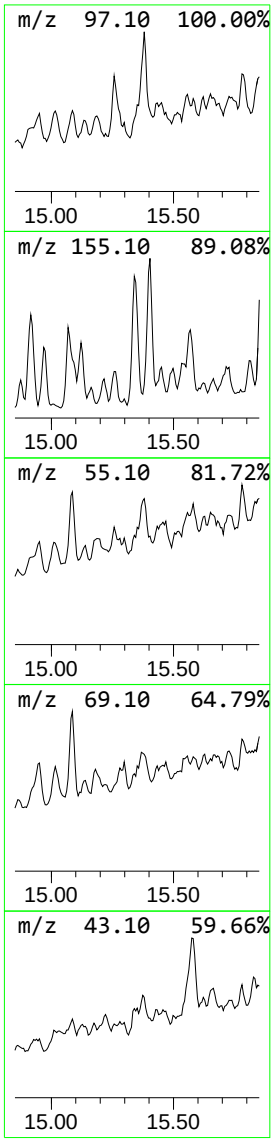
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 2-Pentenoic acid, 3,4,4-tri... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.351	10.92 ng	1142690	Acenaphthene-d10	14.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentenoic acid, 3,4,4-trimethy...	170	C10H18O2	091140-23-7	52
2	5-Methyl-1-phenylhexa-1,3,4-triene	170	C13H14	103240-79-5	38
3	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	38
4	4-(2-Thienyl)butyric acid	170	C8H10O2S	004653-11-6	30
5	2,5-Dimethoxythiophenol	170	C8H10O2S	001483-27-8	30





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052225\
Data File : BP024763.D
Acq On : 22 May 2025 17:43
Operator : RC/JU
Sample : Q2097-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
RT3419

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.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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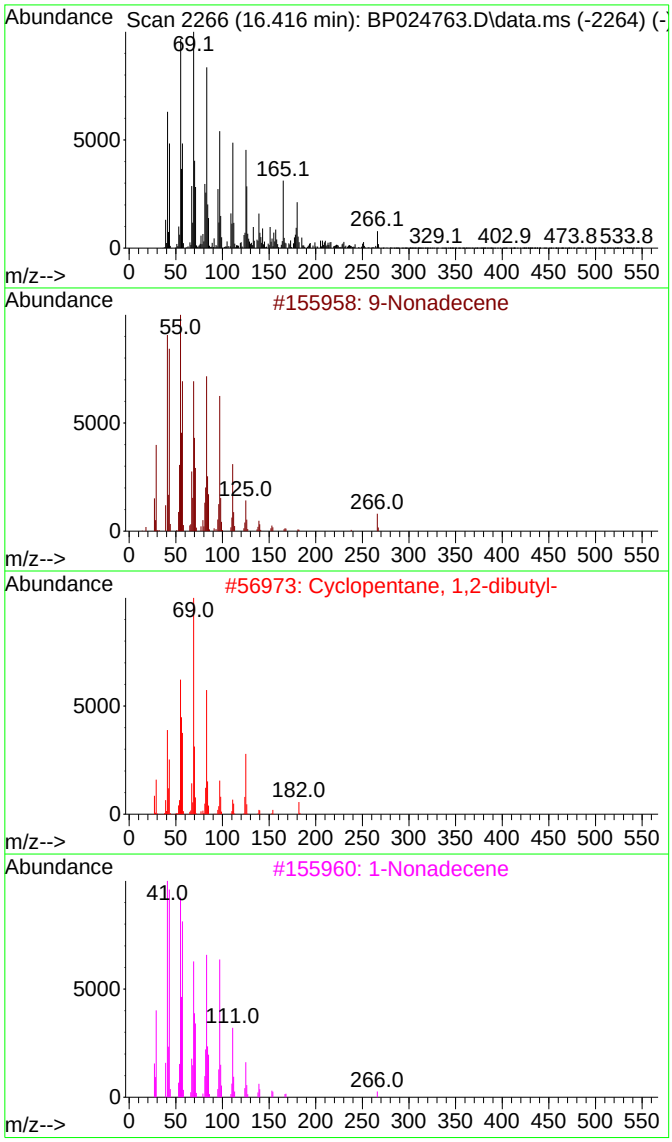
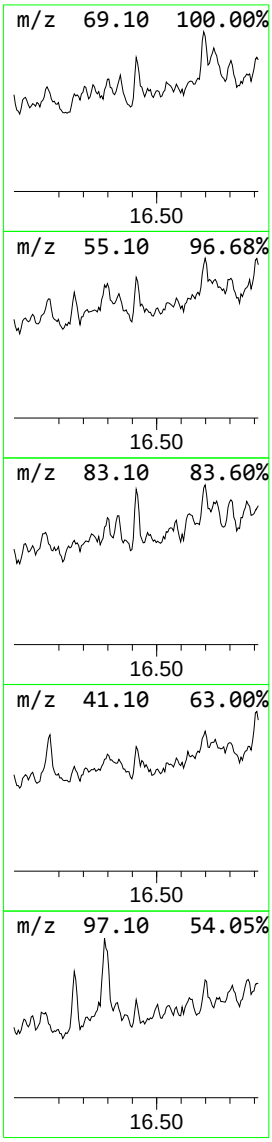
Peak Number 14 9-Nonadecene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.416	20.00 ng	1186430	Phenanthrene-d10	17.092

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Nonadecene	266	C19H38	031035-07-1	84
2	Cyclopentane, 1,2-dibutyl-	182	C13H26	062199-52-4	64
3	1-Nonadecene	266	C19H38	018435-45-5	48
4	Z-5-Nonadecene	266	C19H38	1000131-11-8	47
5	Cyclohexane, 1,1'-propylidenebis-	208	C15H28	054934-91-7	46





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052225\
Data File : BP024763.D
Acq On : 22 May 2025 17:43
Operator : RC/JU
Sample : Q2097-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
RT3419

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.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-...	4.822	4.3	ng	155437	1	7.651	721144	20.0
unknown11.510	11.510	3.0	ng	152625	2	10.422	1014190	20.0
unknown12.157	12.157	2.2	ng	109839	2	10.422	1014190	20.0
unknown12.704	12.704	2.4	ng	249242	3	14.286	2093170	20.0
Decahydro-1,1,4...	13.057	6.8	ng	717257	3	14.286	2093170	20.0
unknown13.698	13.698	12.3	ng	1289910	3	14.286	2093170	20.0
unknown13.769	13.769	3.1	ng	321928	3	14.286	2093170	20.0
unknown14.022	14.022	10.4	ng	1093570	3	14.286	2093170	20.0
unknown14.110	14.110	27.4	ng	2866130	3	14.286	2093170	20.0
unknown14.186	14.186	8.6	ng	904225	3	14.286	2093170	20.0
unknown14.822	14.822	14.5	ng	1519100	3	14.286	2093170	20.0
N-[1-(4-Chlorob...	15.086	31.2	ng	3261360	3	14.286	2093170	20.0
2-Pentenoic aci...	15.351	10.9	ng	1142690	3	14.286	2093170	20.0
9-Nonadecene	16.416	20.0	ng	1186430	4	17.092	1186430	20.0