LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031622\
Data File : BP009429.D
Acq On : 16 Mar 2022 18:53
Operator : CG/JU
Sample : N1951-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DB-8-(15-20)-3-12-22

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-BP030522.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP009429.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.058	8	12	33	rVB	447591	823089	8.71%	1.065%
2	5.416	406	413	435	rBV	3692295	6091812	64.50%	7.879%
3	7.005	670	683	697	rBV	3330633	6105085	64.64%	7.896%
4	7.805	812	819	828	rBV	1360838	2418369	25.60%	3.128%
5	8.963	1009	1016	1029	rVB	2173591	3923648	41.54%	5.075%
6	10.604	1287	1295	1307	rBV	1608994	3238358	34.29%	4.188%
7	11.304	1405	1414	1421	rVB9	246630	644629	6.82%	0.834%
8	11.651	1468	1473	1477	rBV3	259308	555713	5.88%	0.719%
9	12.351	1587	1592	1598	rBV5	377838	843935	8.94%	1.091%
10	12.816	1667	1671	1678	rBV6	287822	615315	6.51%	0.796%
11	12.892	1680	1684	1690	rVB3	736037	1225292	12.97%	1.585%
12	12.957	1692	1695	1699	rVV4	283262	461787	4.89%	0.597%
13	13.069	1707	1714	1722	rBV	5340629	9445204	100.00%	12.216%
14	13.240	1739	1743	1753	rVB	1135195	2056108	21.77%	2.659%
15	13.881	1847	1852	1857	rBV	3027526	4840554	51.25%	6.260%
16	14.198	1902	1906	1911	rBV4	1516519	2896508	30.67%	3.746%
17	14.292	1918	1922	1929	rVB	4705084	7700009	81.52%	9.959%
18	14.363	1930	1934	1939	rBV4	1051966	1968145	20.84%	2.545%
19	14.428	1940	1945	1946	rBV	2402595	3320254	35.15%	4.294%
20	14.992	2038	2041	2048	rBV4	1479723	2648433	28.04%	3.425%
21	15.257	2081	2086	2093	rBV2	4118463	6889708	72.94%	8.911%
22	15.957	2200	2205	2211	rBV3	4481227	8607591	91.13%	11.132%

Sum of corrected areas: 77319546

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031622\
Data File : BP009429.D
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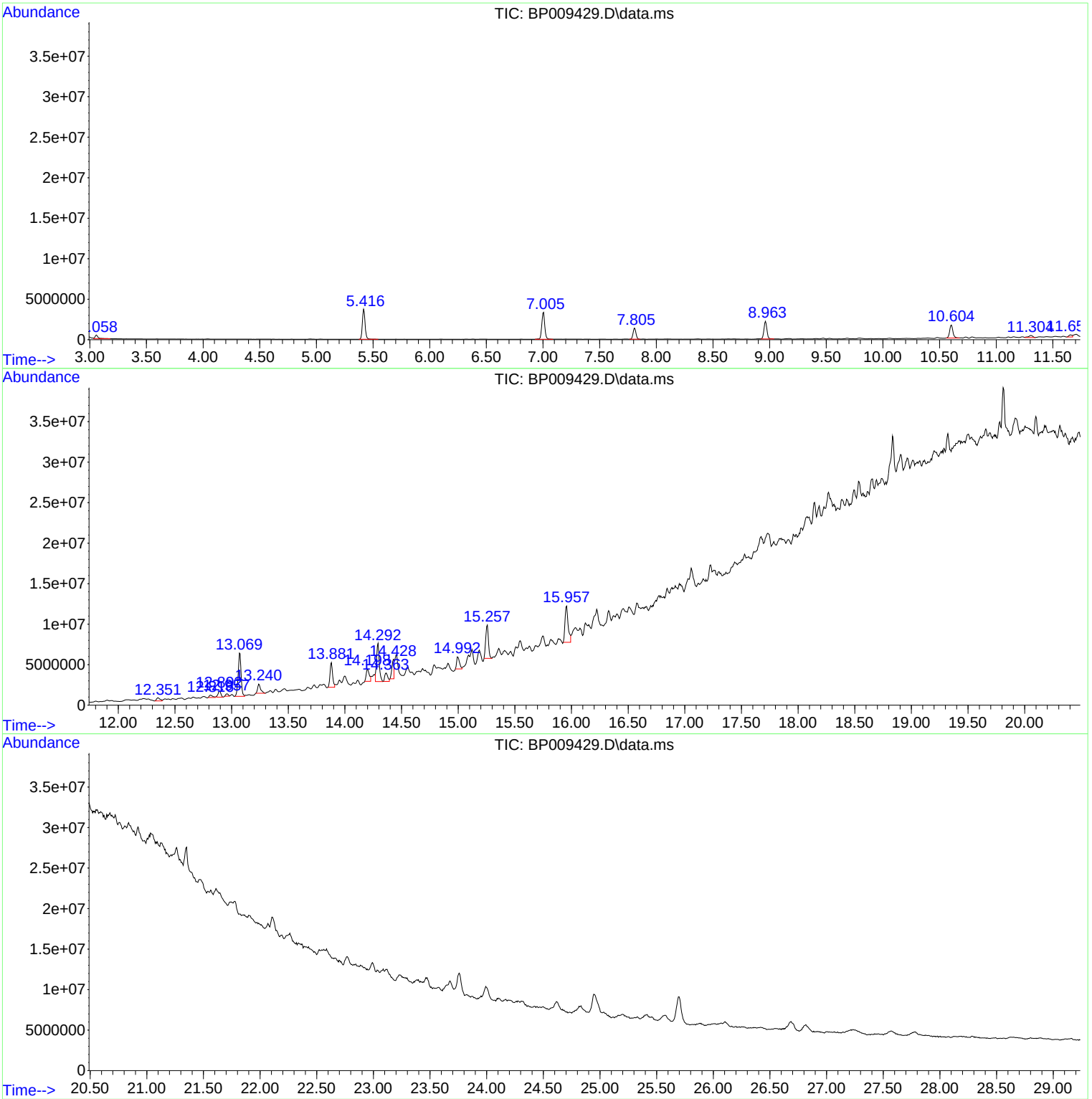
Sample : N1951-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DB-8-(15-20)-3-12-22

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.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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Data File : BP009429.D
Acq On : 16 Mar 2022 18:53
Operator : CG/JU
Sample : N1951-09
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

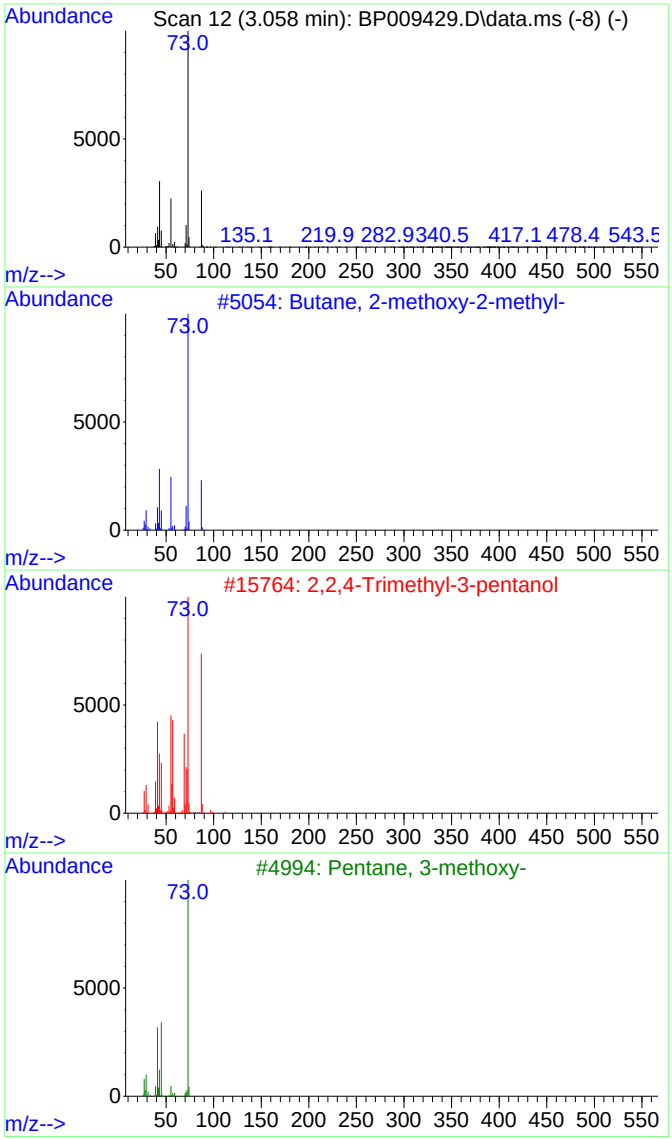
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

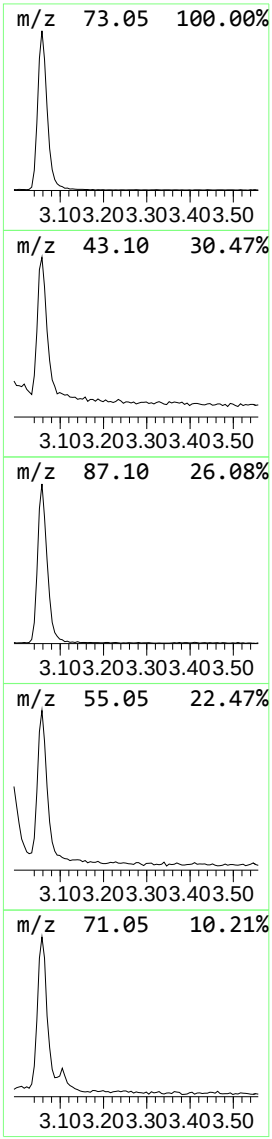
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.058	6.81 ng	823089	1,4-Dichlorobenzene-d4	7.810

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	90
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	12
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9





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Data File : BP009429.D
Acq On : 16 Mar 2022 18:53
Operator : CG/JU
Sample : N1951-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DB-8-(15-20)-3-12-22

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.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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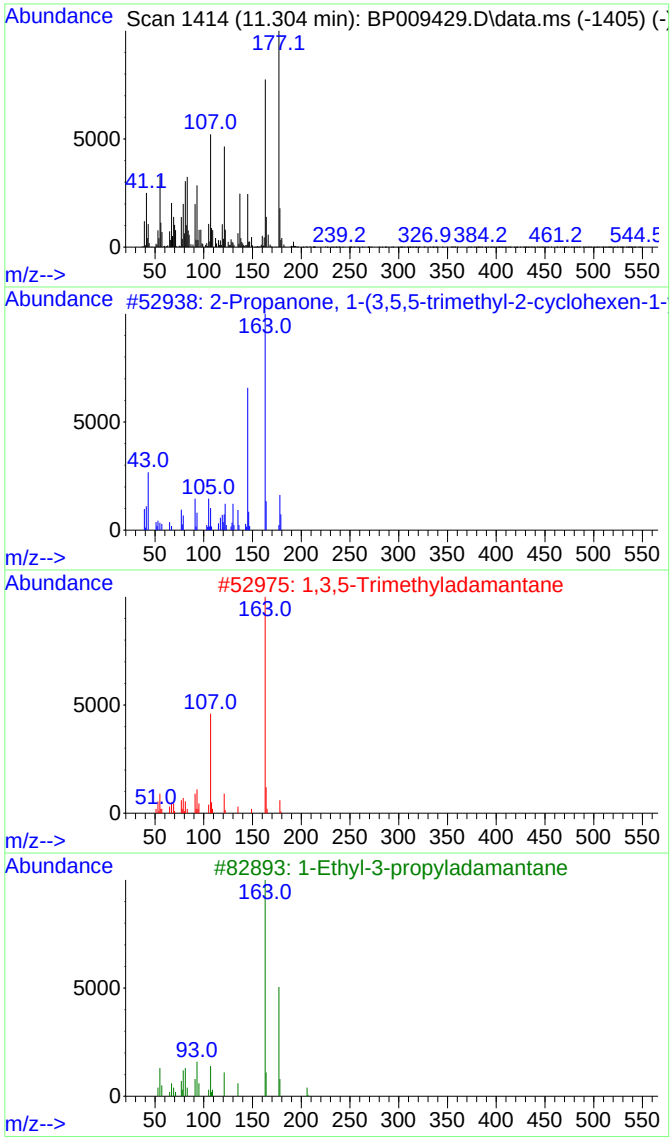
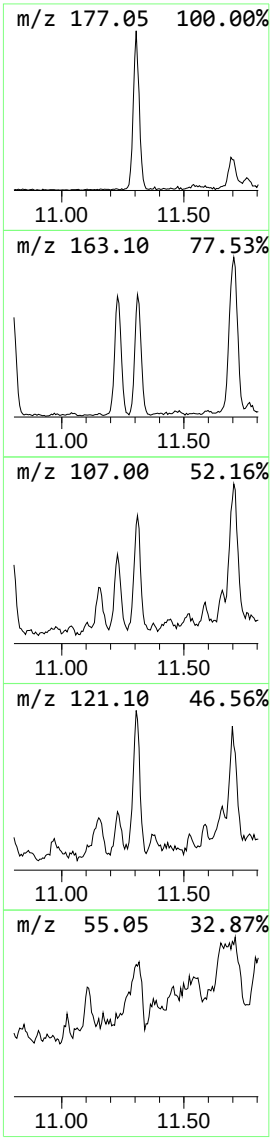
Peak Number 2 unknown11.304 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.304	3.98 ng	644629	Naphthalene-d8	10.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanone, 1-(3,5,5-trimethyl-...	178	C12H18O	016695-72-0	46
2	1,3,5-Trimethyladamantane	178	C13H22	000707-35-7	38
3	1-Ethyl-3-propyladamantane	206	C15H26	019385-94-5	38
4	3-(4-Buthoxybenzylthio)-1,2,4-(4...	263	C13H17N3OS	057736-49-9	35
5	1,2-Dimethyl-5-nitroadamantane	209	C12H19NO2	006588-68-7	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031622\
Data File : BP009429.D
Acq On : 16 Mar 2022 18:53
Operator : CG/JU
Sample : N1951-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DB-8-(15-20)-3-12-22

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.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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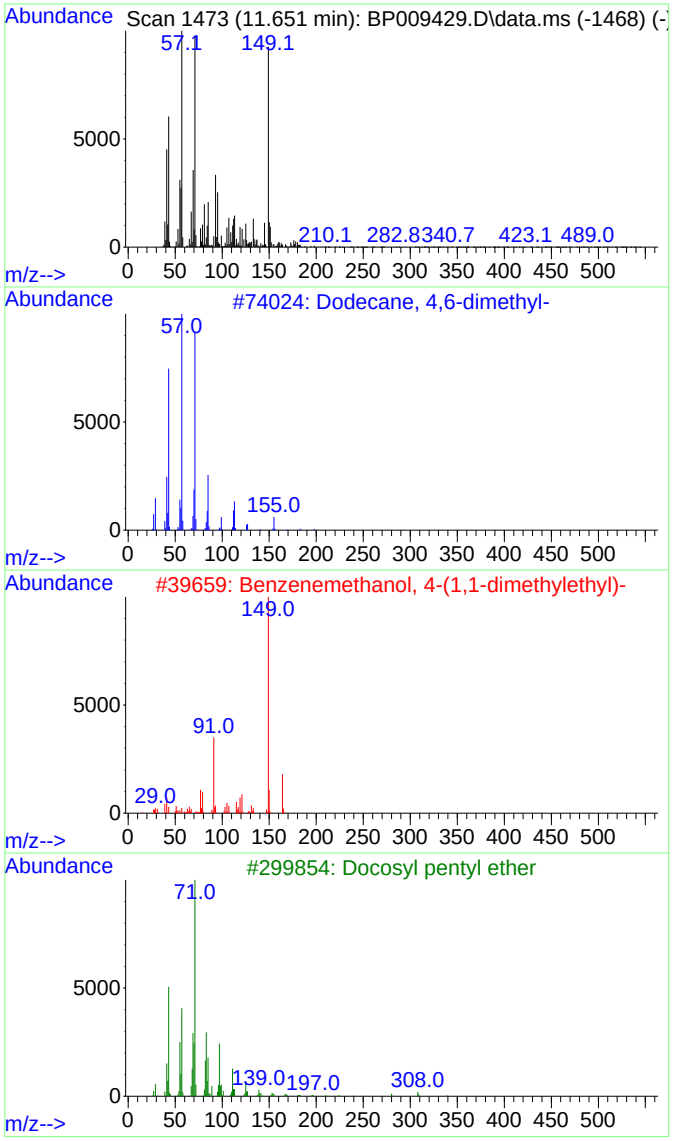
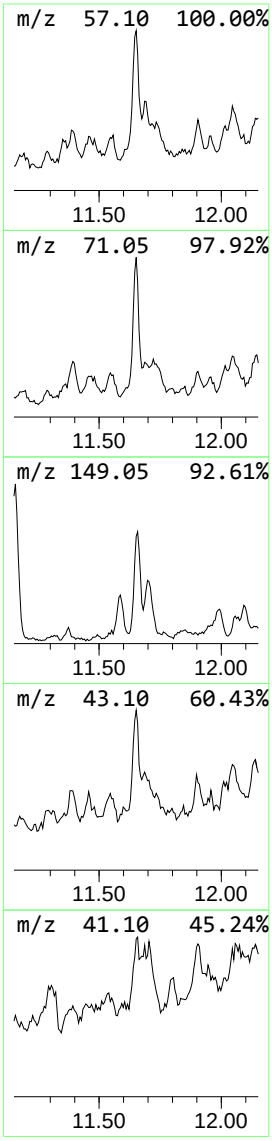
Peak Number 3 Dodecane, 4,6-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.651	3.43 ng	555713	Naphthalene-d8	10.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	51
2	Benzenemethanol, 4-(1,1-dimethyl...	164	C11H16O	000877-65-6	27
3	Docosyl pentyl ether	396	C27H56O	1000406-42-1	27
4	Bicyclo[2.2.2]oct-5-en-2-yldimet...	151	C10H17N	033162-94-6	25
5	7-Methylthieno[3,2-b]pyridine	149	C8H7NS	013362-83-9	22





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031622\
Data File : BP009429.D
Acq On : 16 Mar 2022 18:53
Operator : CG/JU
Sample : N1951-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DB-8-(15-20)-3-12-22

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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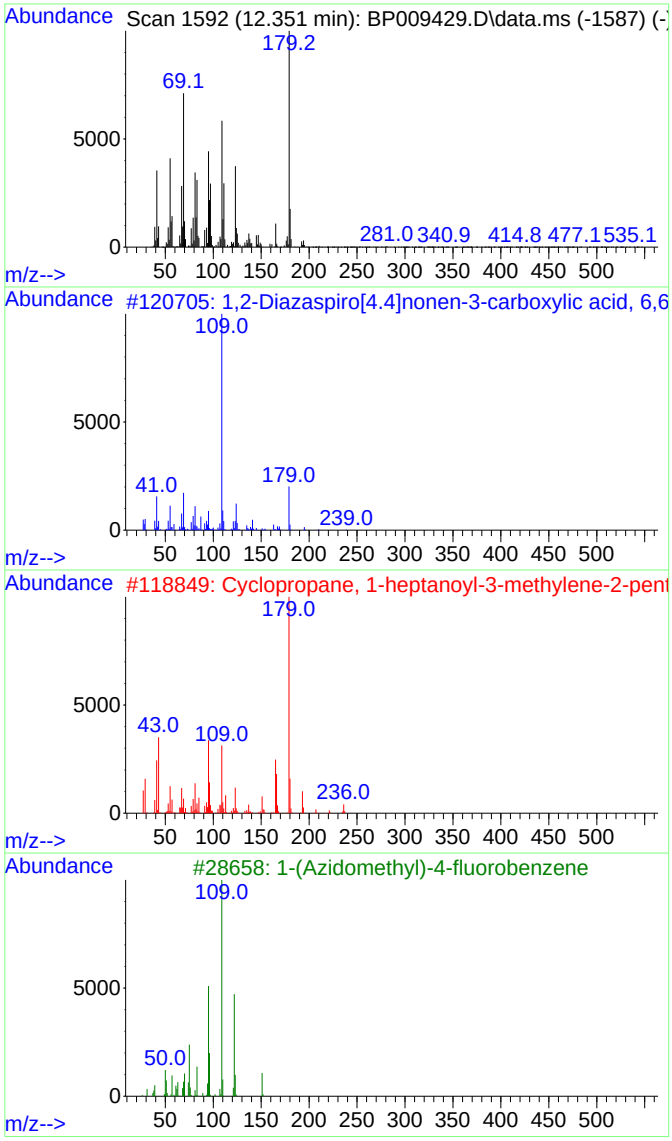
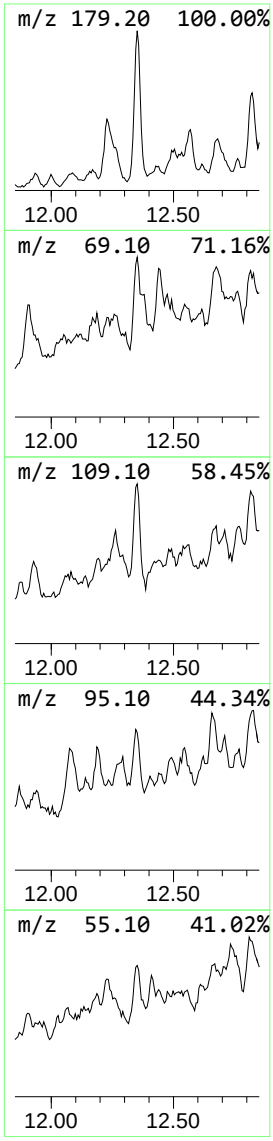
Peak Number 4 unknown12.351 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.351	5.21 ng	843935	Naphthalene-d8	10.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Diazaspiro[4.4]nonen-3-carbo...	238	C13H22N2O2	1000164-25-8	43
2	Cyclopropane, 1-heptanoyl-3-meth...	236	C16H28O	1000157-26-0	37
3	1-(Azidomethyl)-4-fluorobenzene	151	C7H6FN3	159979-96-1	27
4	Tricyclo[4.3.1.1(3,8)]undecane, ...	180	C12H20O	021898-92-0	18
5	4,2-Cresotic acid, 6-methoxy-, b...	538	C29H30O10	019314-74-0	16





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031622\
Data File : BP009429.D
Acq On : 16 Mar 2022 18:53
Operator : CG/JU
Sample : N1951-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DB-8-(15-20)-3-12-22

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

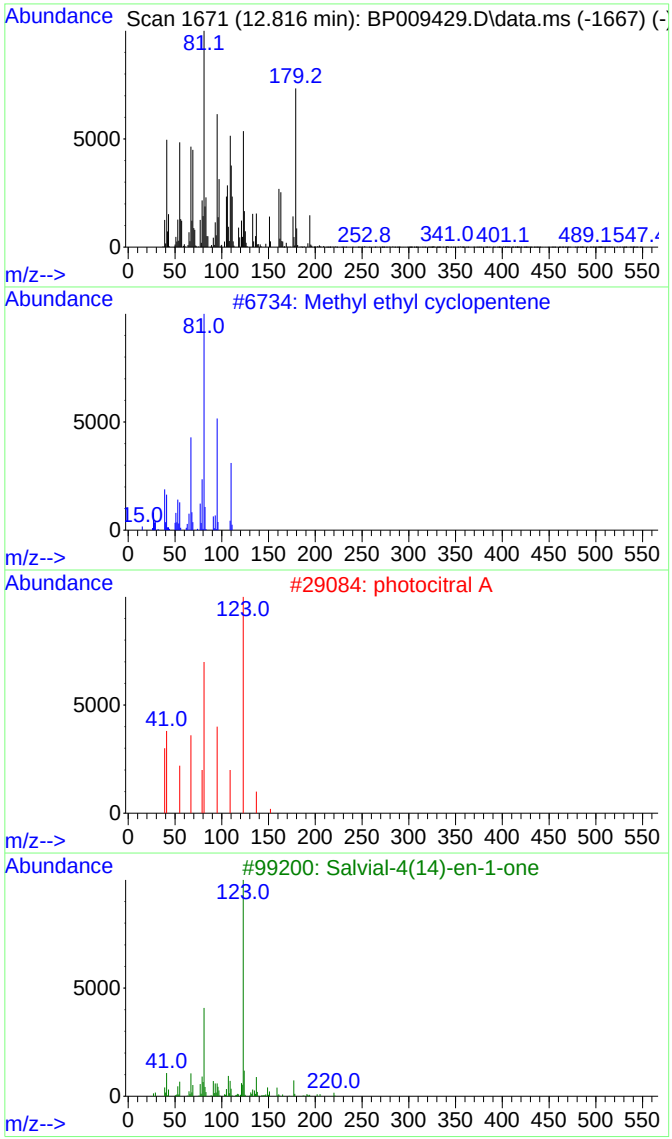
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

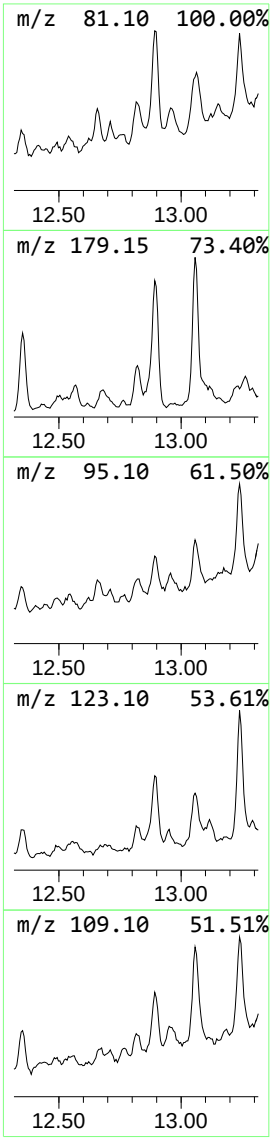
****

Peak Number 5 unknown12.816 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.816	3.71 ng	615315	Acenaphthene-d10	14.451

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl ethyl cyclopentene	110	C8H14	019780-56-4	42
2	photocitral A	152	C10H16O	055253-28-6	25
3	Salvial-4(14)-en-1-one	220	C15H24O	073809-82-2	22
4	Cyclohexene, 1-ethyl-	110	C8H14	001453-24-3	18
5	Cyclohexane, 1,4-dibromo-	240	C6H10Br2	035076-92-7	14





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031622\
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Operator : CG/JU
Sample : N1951-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DB-8-(15-20)-3-12-22

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

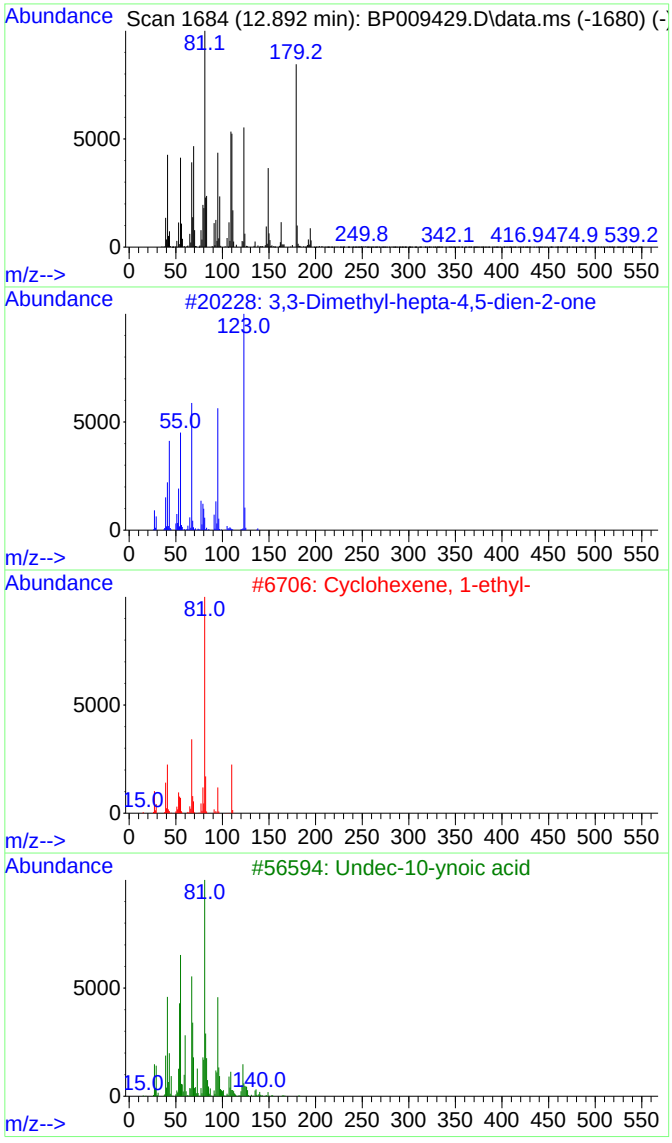
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

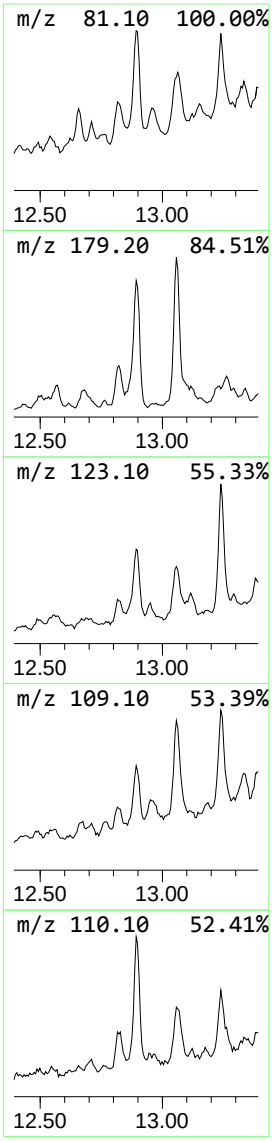
****

Peak Number 6 unknown12.893 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.893	7.38 ng	1225290	Acenaphthene-d10	14.451

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3-Dimethyl-hepta-4,5-dien-2-one	138	C9H14O	1000190-54-1	35
2	Cyclohexene, 1-ethyl-	110	C8H14	001453-24-3	30
3	Undec-10-ynoic acid	182	C11H18O2	002777-65-3	27
4	Furan, 2,3,5-trimethyl-	110	C7H10O	010504-04-8	27
5	trans-(2-Ethylcyclopentyl)methyl...	170	C10H18O2	1000099-13-9	27





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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.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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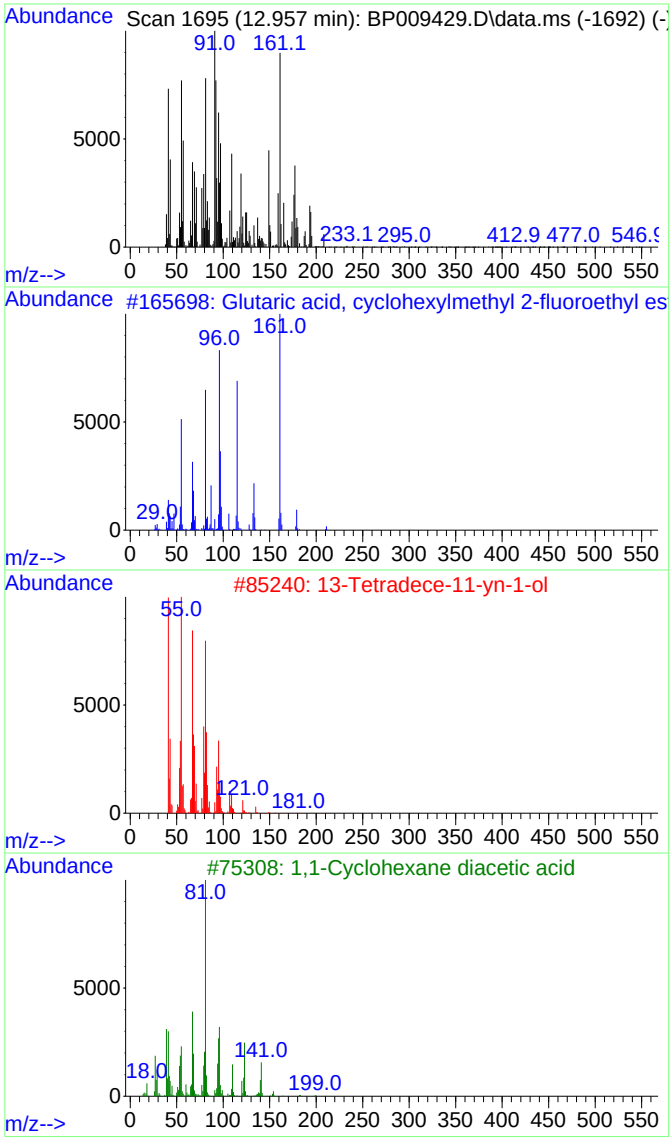
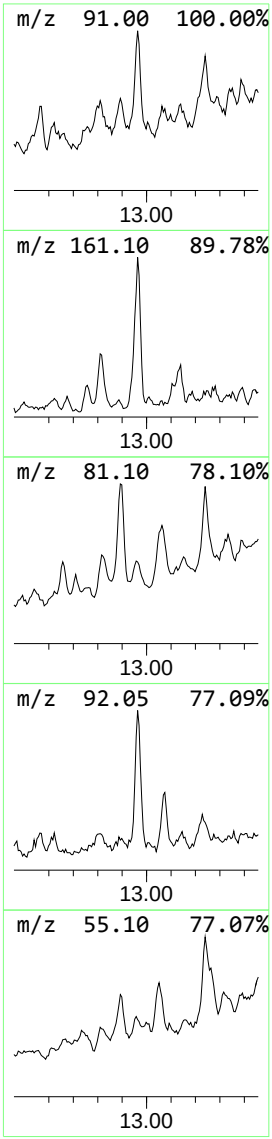
Peak Number 7 unknown12.957 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.957	2.78 ng	461787	Acenaphthene-d10	14.451

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Glutaric acid, cyclohexylmethyl ...	274	C14H23FO4	1000393-71-3	25
2	13-Tetradece-11-yn-1-ol	208	C14H24O	1000131-00-4	14
3	1,1-Cyclohexane diacetic acid	200	C10H16O4	009355-11-7	11
4	1-Ethyl-5-methylcyclopentene	110	C8H14	097797-57-4	10
5	4-Amino-2-chloro-3-fluorophenol	161	C6H5ClFNO	1003710-18-6	10





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TIC Library : C:\Database\NIST20.L
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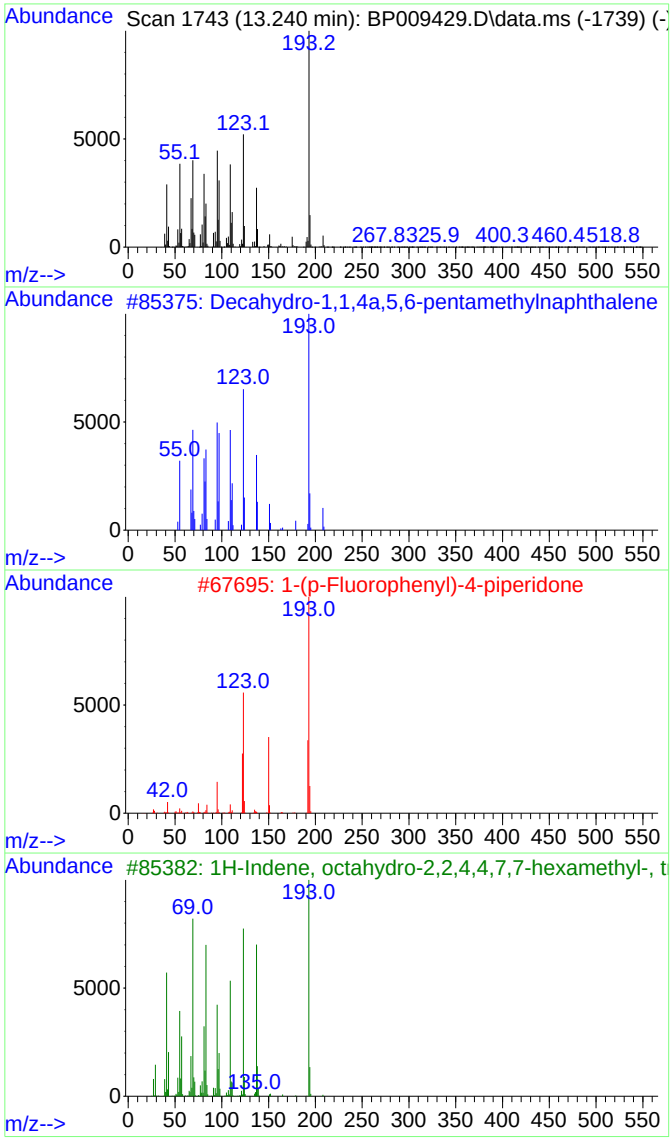
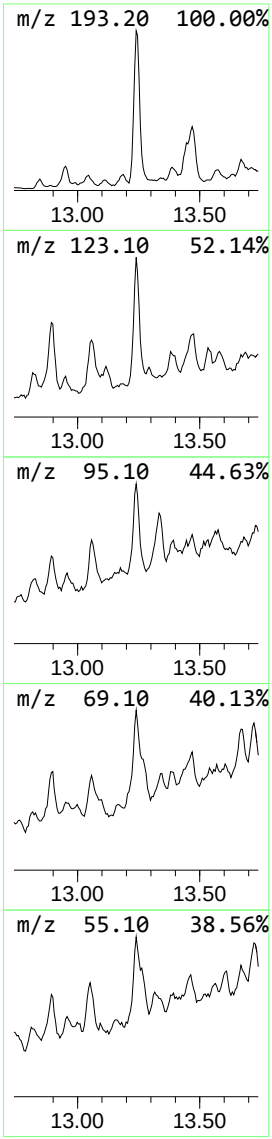
****
Peak Number 8 Decahydro-1,1,4a,5,6-pentam... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.240	12.39 ng	2056110	Acenaphthene-d10	14.451

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	60
2	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	43
3	1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	38
4	2,4,6-Trimethyl-2-(4-methyl-pent...	206	C14H22O	1000193-58-6	35
5	benzoxazole, 5,6-dimethoxy-2-met...	193	C10H11NO3	1000395-98-9	27





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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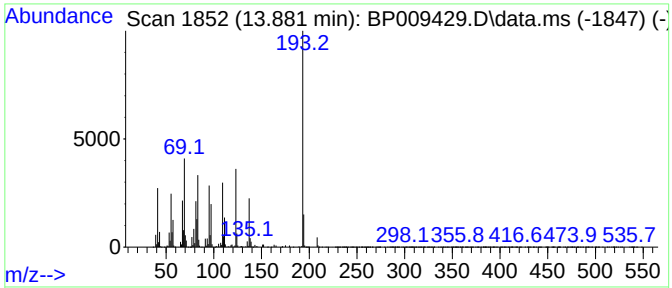
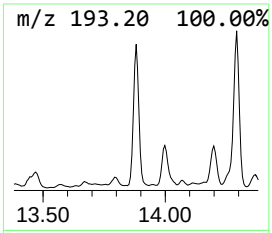
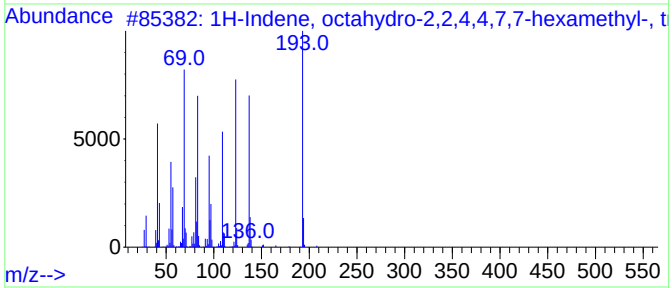
Peak Number 9 unknown13.881 Concentration Rank 2

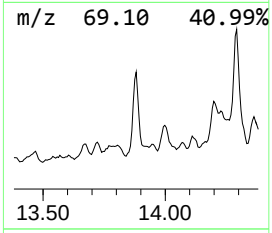
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.881	29.16 ng	4840550	Acenaphthene-d10	14.451

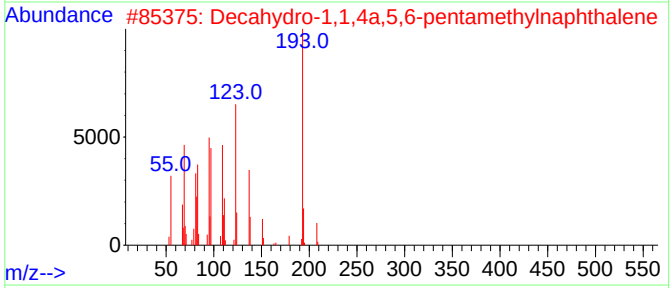
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	45
2	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	43
3	Borinic acid, diethyl-, 3,3,5-tr...	208	C13H25BO	057387-76-5	38
4	2-Anthracenamine	193	C14H11N	000613-13-8	38
5	9-Phenanthrenamine	193	C14H11N	000947-73-9	35

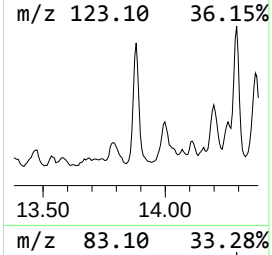


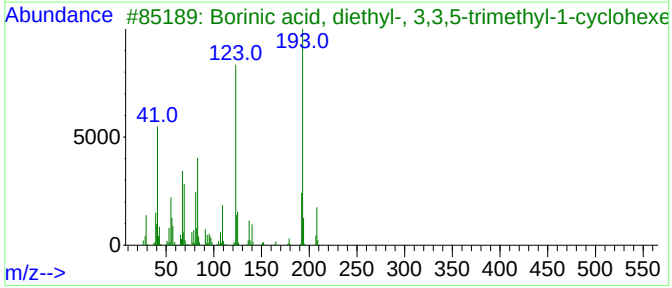


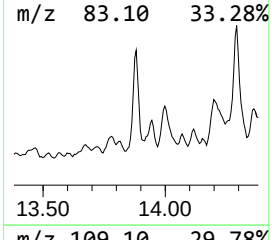


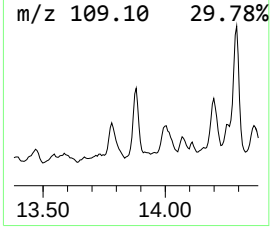












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031622\
Data File : BP009429.D
Acq On : 16 Mar 2022 18:53
Operator : CG/JU
Sample : N1951-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DB-8-(15-20)-3-12-22

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.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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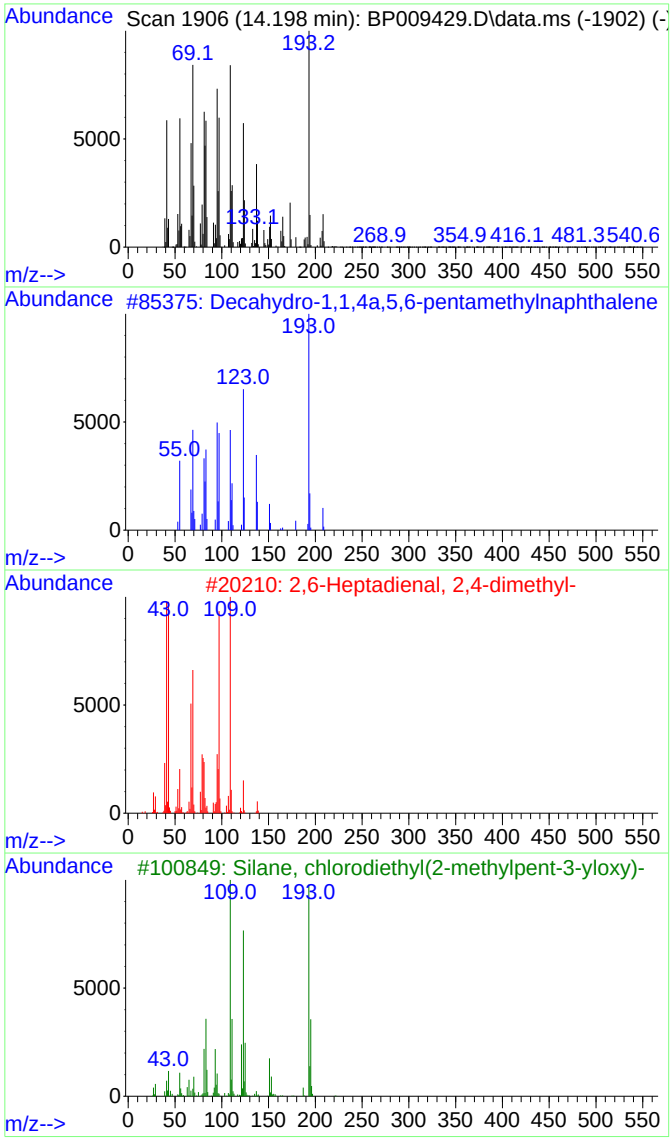
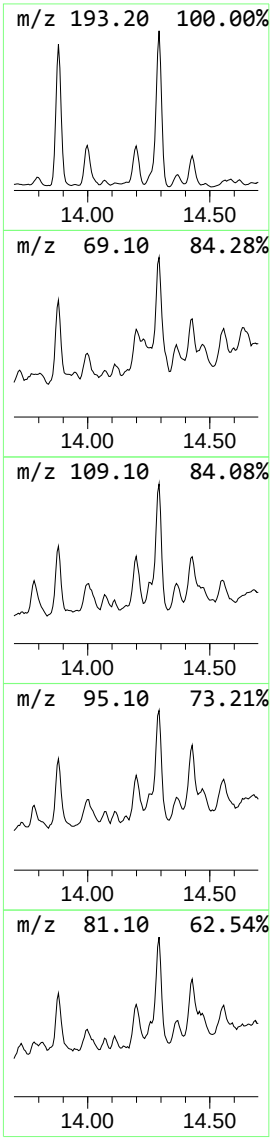
Peak Number 10 unknown14.198 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.198	17.45 ng	2896510	Acenaphthene-d10	14.451

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	49
2	2,6-Heptadienal, 2,4-dimethyl-	138	C9H14O	085136-08-9	46
3	Silane, chlorodiethyl(2-methylpe...	222	C10H23ClOSi	1000363-79-1	35
4	1-(2-Furyl)-4,4,4-trifluoro-1,3-...	206	C8H5F3O3	000326-90-9	27
5	Iminostilbene	193	C14H11N	000256-96-2	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031622\
Data File : BP009429.D
Acq On : 16 Mar 2022 18:53
Operator : CG/JU
Sample : N1951-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DB-8-(15-20)-3-12-22

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.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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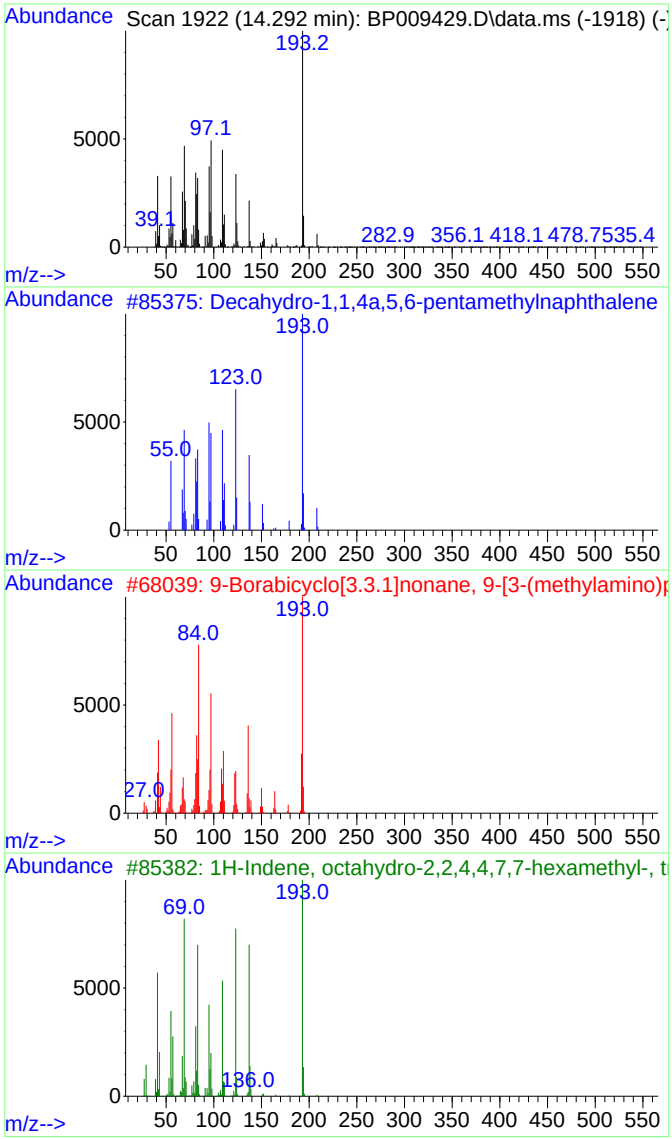
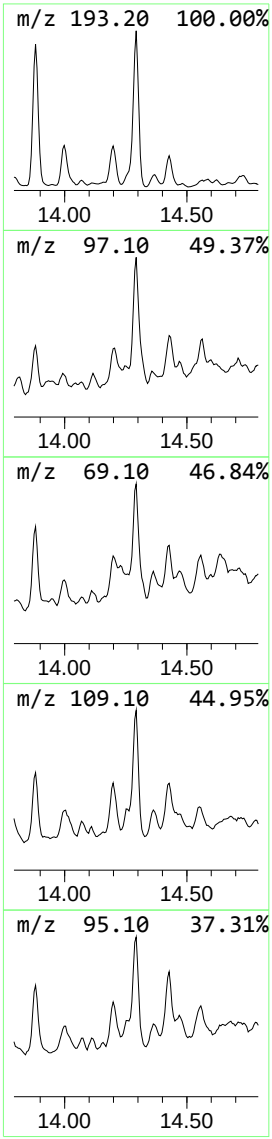
Peak Number 11 unknown14.292 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.292	46.38 ng	7700010	Acenaphthene-d10	14.451

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	52
2	9-Borabicyclo[3.3.1]nonane, 9-[3...	193	C12H24BN	1010162-56-0	47
3	1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	38
4	2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	36
5	Borinic acid, diethyl-, 3,3,5-tr...	208	C13H25BO	057387-76-5	32





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Data File : BP009429.D
Acq On : 16 Mar 2022 18:53
Operator : CG/JU
Sample : N1951-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DB-8-(15-20)-3-12-22

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.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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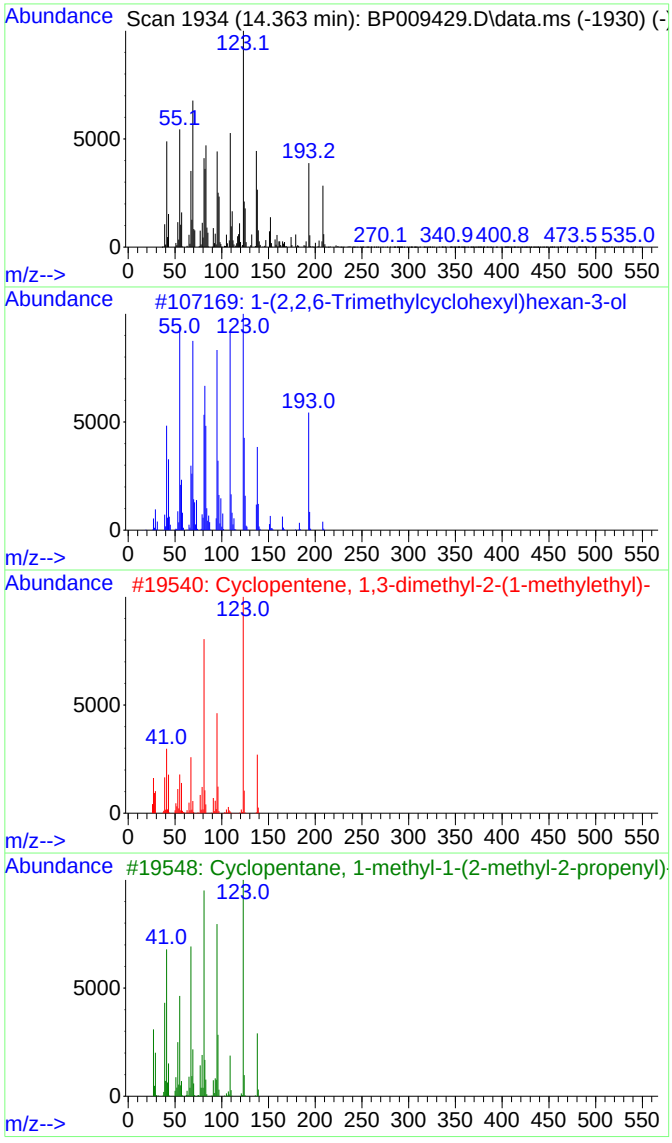
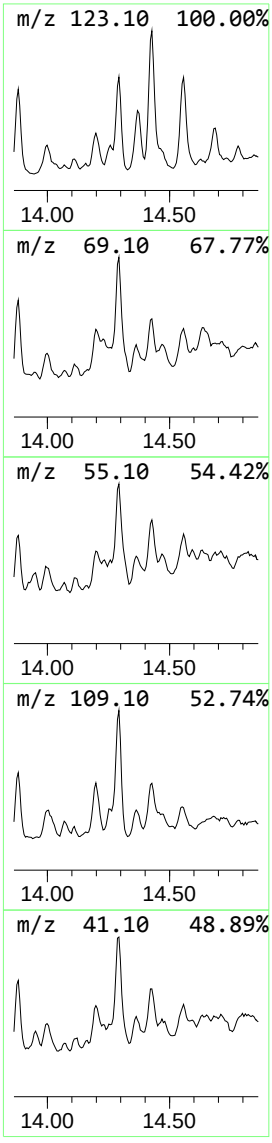
Peak Number 12 1-(2,2,6-Trimethylcyclohexy... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.363	11.86 ng	1968150	Acenaphthene-d10	14.451

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(2,2,6-Trimethylcyclohexyl)hex...	226	C15H30O	070788-30-6	52
2	Cyclopentene, 1,3-dimethyl-2-(1-...	138	C10H18	061142-32-3	38
3	Cyclopentane, 1-methyl-1-(2-meth...	138	C10H18	074764-47-9	35
4	1,3-Benzenediamine, 4-methoxy-	138	C7H10N2O	000615-05-4	35
5	Diethyl 1-(carbethoxy)ethylphosp...	238	C9H19O5P	1000306-25-7	30





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031622\
Data File : BP009429.D
Acq On : 16 Mar 2022 18:53
Operator : CG/JU
Sample : N1951-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DB-8-(15-20)-3-12-22

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.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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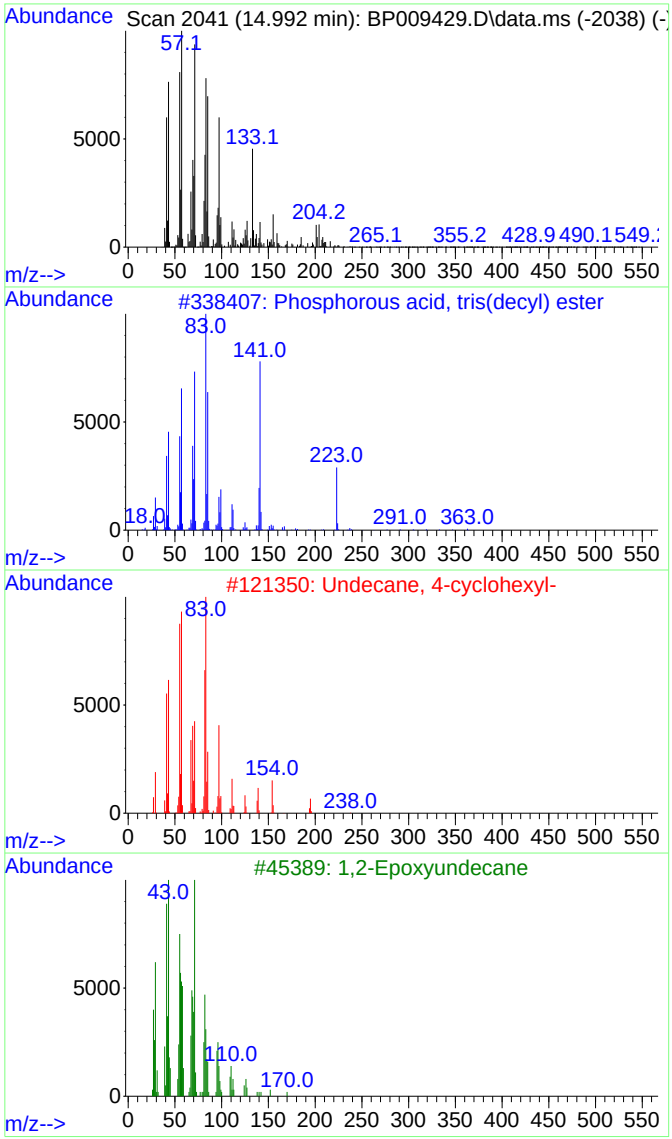
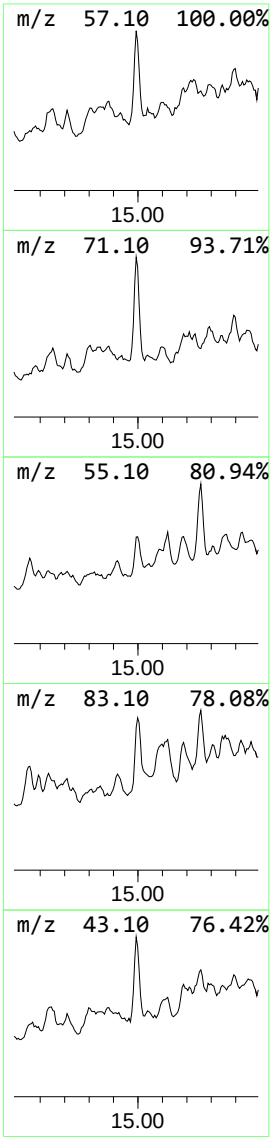
Peak Number 13 unknown14.992 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.992	15.95 ng	2648430	Acenaphthene-d10	14.451

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phosphorous acid, tris(decyl) ester	502	C30H63O3P	002929-86-4	49
2	Undecane, 4-cyclohexyl-	238	C17H34	013151-79-6	38
3	1,2-Epoxyundecane	170	C11H22O	017322-97-3	38
4	Tridecane, 7-propyl-	226	C16H34	055045-09-5	30
5	Dodecane, 4-methyl-	184	C13H28	006117-97-1	30





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031622\
Data File : BP009429.D
Acq On : 16 Mar 2022 18:53
Operator : CG/JU
Sample : N1951-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DB-8-(15-20)-3-12-22

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.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
Butane, 2-metho...	3.058	6.8	ng	823089	1	7.810	2418370	20.0
unknown11.304	11.304	4.0	ng	644629	2	10.604	3238360	20.0
Dodecane, 4,6-d...	11.651	3.4	ng	555713	2	10.604	3238360	20.0
unknown12.351	12.351	5.2	ng	843935	2	10.604	3238360	20.0
unknown12.816	12.816	3.7	ng	615315	3	14.451	3320250	20.0
unknown12.893	12.893	7.4	ng	1225290	3	14.451	3320250	20.0
unknown12.957	12.957	2.8	ng	461787	3	14.451	3320250	20.0
Decahydro-1,1,4...	13.240	12.4	ng	2056110	3	14.451	3320250	20.0
unknown13.881	13.881	29.2	ng	4840550	3	14.451	3320250	20.0
unknown14.198	14.198	17.4	ng	2896510	3	14.451	3320250	20.0
unknown14.292	14.292	46.4	ng	7700010	3	14.451	3320250	20.0
1-(2,2,6-Trimet...	14.363	11.9	ng	1968150	3	14.451	3320250	20.0
unknown14.992	14.992	15.9	ng	2648430	3	14.451	3320250	20.0