LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011322\
Data File : BP008795.D
Acq On : 13 Jan 2022 17:57
Operator : CG/JU
Sample : N1119-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-NORTH-SIDEWALL

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-BP010622.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP008795.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.064	8	13	25	rVB	2833103	3342740	12.07%	3.147%
2	5.375	401	406	411	rBV	210050	300496	1.09%	0.283%
3	5.446	411	418	425	rVV	4294787	6453250	23.30%	6.075%
4	5.528	425	432	447	rVB2	104552	347929	1.26%	0.328%
5	6.846	649	656	664	rBV2	202798	429618	1.55%	0.404%
6	7.040	679	689	705	rBV	3721480	6460940	23.33%	6.083%
7	7.858	821	828	837	rBV	1213053	2046937	7.39%	1.927%
8	8.234	886	892	898	rVB	206172	345066	1.25%	0.325%
9	8.358	908	913	918	rBV	180072	288561	1.04%	0.272%
10	8.969	1008	1017	1020	rBV	308446	520937	1.88%	0.490%
11	9.022	1020	1026	1038	rVB	2319040	4417327	15.95%	4.159%
12	9.916	1172	1178	1182	rBV	178448	284879	1.03%	0.268%
13	10.063	1193	1203	1212	rBV2	368651	695328	2.51%	0.655%
14	10.663	1294	1305	1311	rBV	1510119	2769867	10.00%	2.608%
15	13.122	1716	1723	1747	rBV	6291983	9785907	35.34%	9.213%
16	14.498	1949	1957	1965	rBV	2002398	3194614	11.54%	3.008%
17	15.992	2204	2211	2233	rBV	4287675	6785877	24.51%	6.388%
18	17.251	2419	2425	2440	rBV	2275363	3586839	12.95%	3.377%
19	18.151	2573	2578	2586	rBV	618216	1037242	3.75%	0.976%
20	19.380	2783	2787	2799	rBV2	516917	1055390	3.81%	0.994%
21	19.816	2856	2861	2870	rBV	10155401	12101657	43.70%	11.393%
22	20.492	2971	2976	2982	rBV2	1235540	2109624	7.62%	1.986%
23	21.010	3061	3064	3068	rBV2	455551	511553	1.85%	0.482%
24	21.057	3069	3072	3078	rVV2	851963	1093796	3.95%	1.030%
25	21.345	3116	3121	3130	rVB	2945861	3874848	13.99%	3.648%
26	21.469	3136	3142	3162	rBV	18040384	27690622	100.00%	26.069%
27	23.774	3527	3534	3543	rVB2	2151432	4688760	16.93%	4.414%

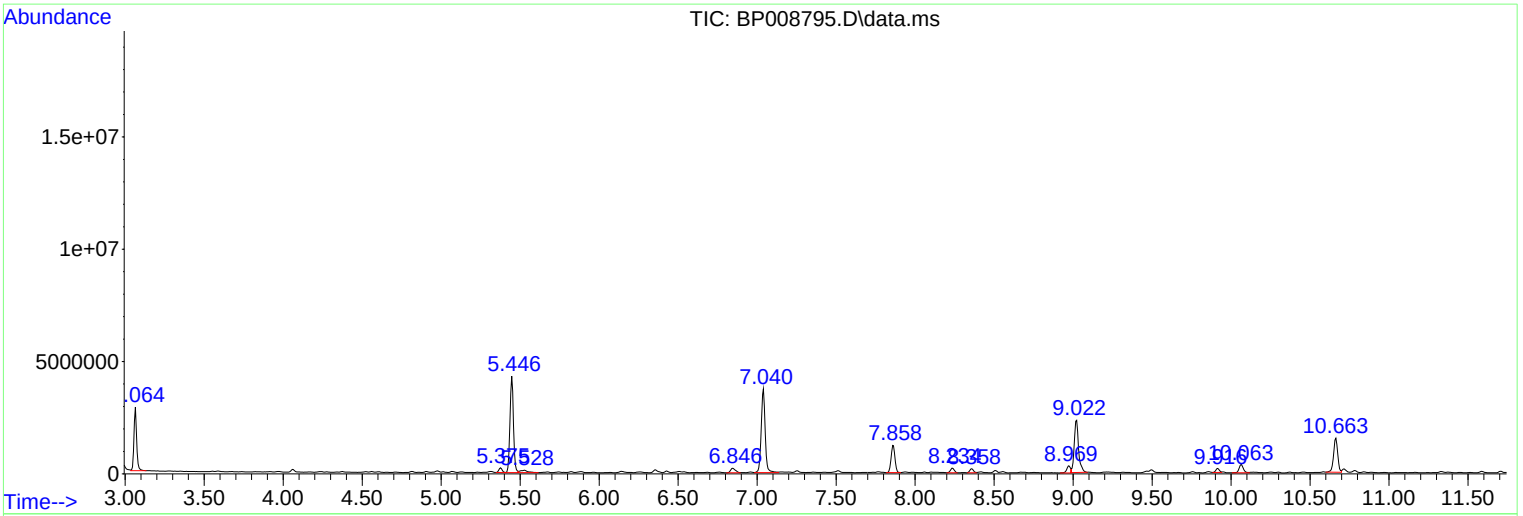
Sum of corrected areas: 106220604

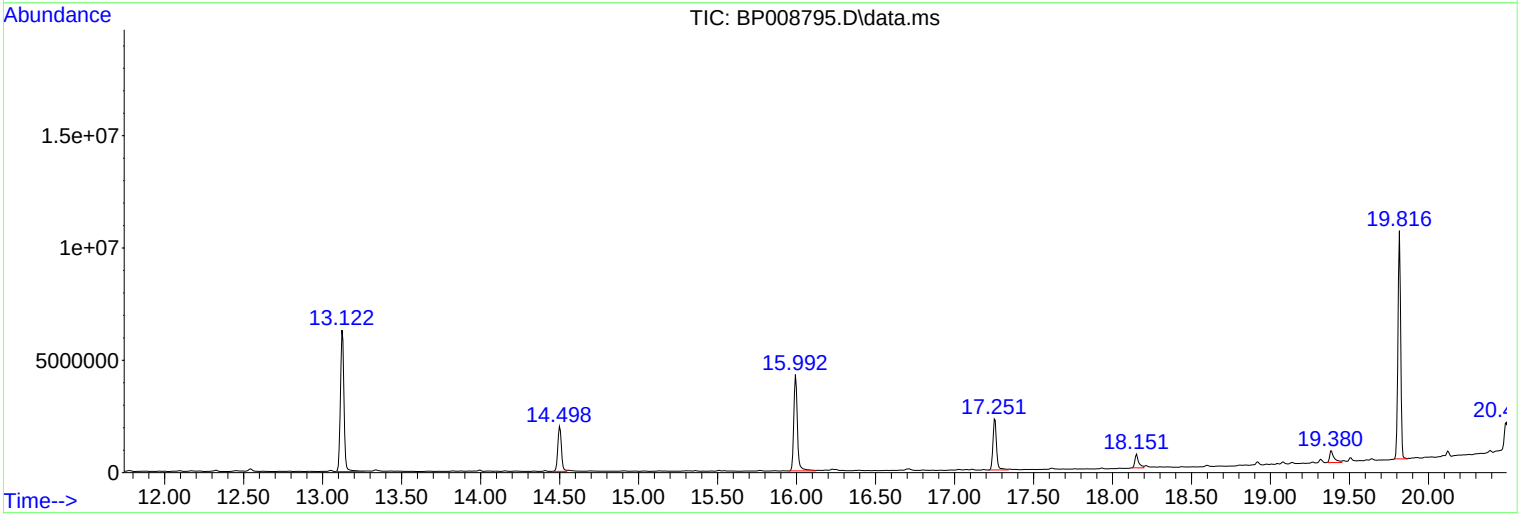
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011322\
Data File : BP008795.D
Acq On : 13 Jan 2022 17:57
Operator : CG/JU
Sample : N1119-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

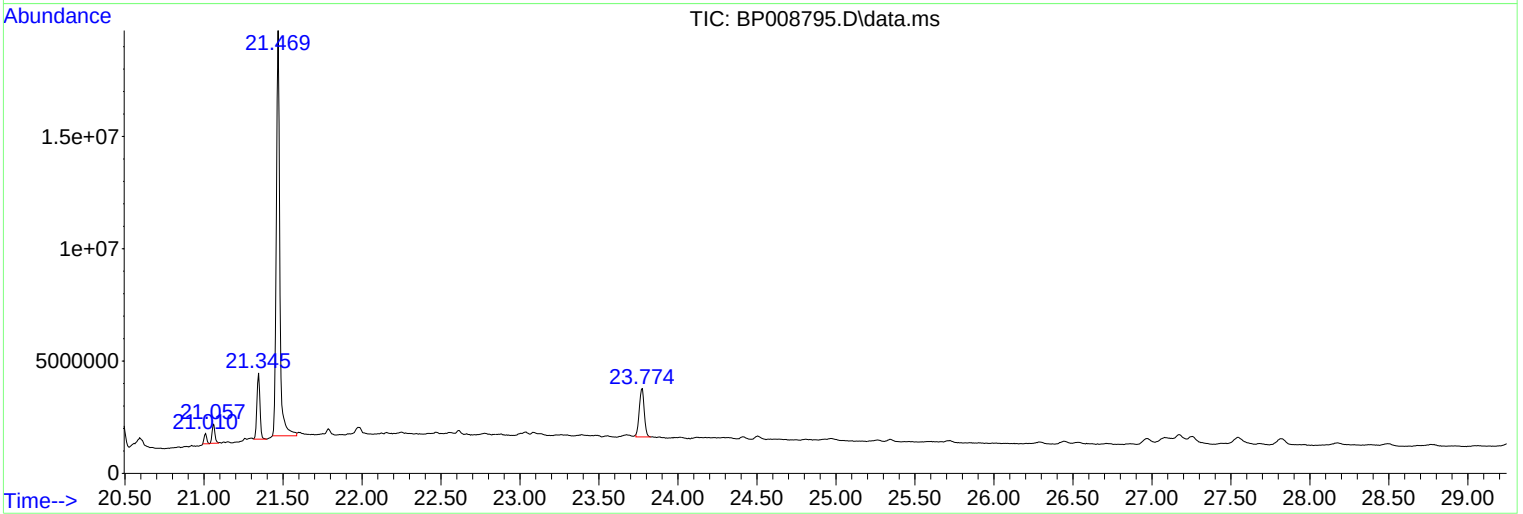
Instrument :
BNA_P
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-NORTH-SIDEWALL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.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 : BP008795.D
Acq On : 13 Jan 2022 17:57
Operator : CG/JU
Sample : N1119-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
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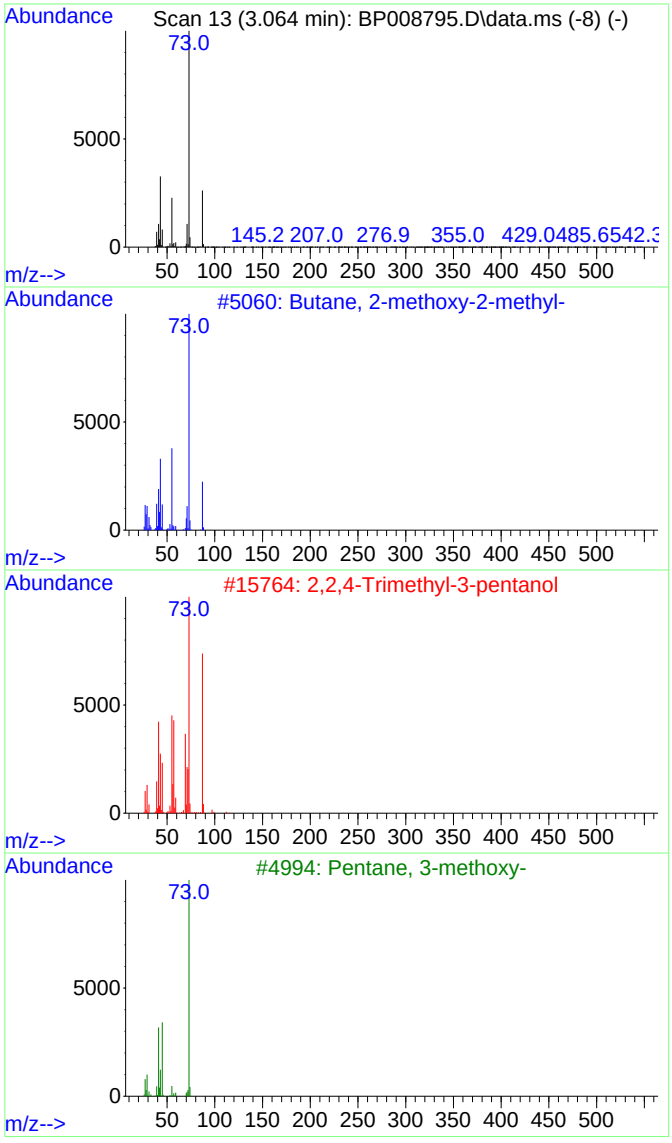
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

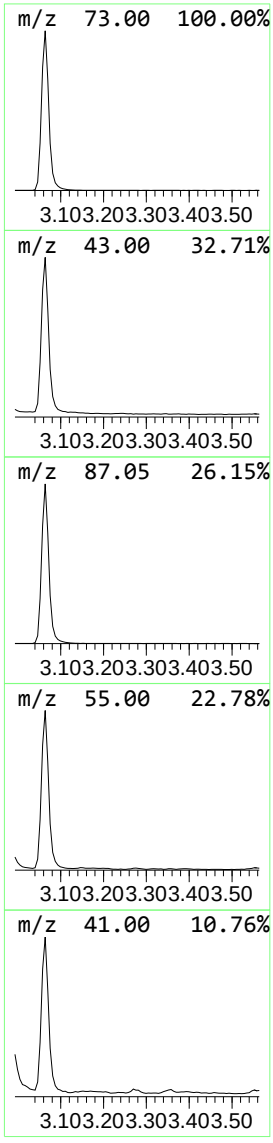
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.064	32.66 ng	3342740	1,4-Dichlorobenzene-d4	7.863

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011322\
Data File : BP008795.D
Acq On : 13 Jan 2022 17:57
Operator : CG/JU
Sample : N1119-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-NORTH-SIDEWALL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

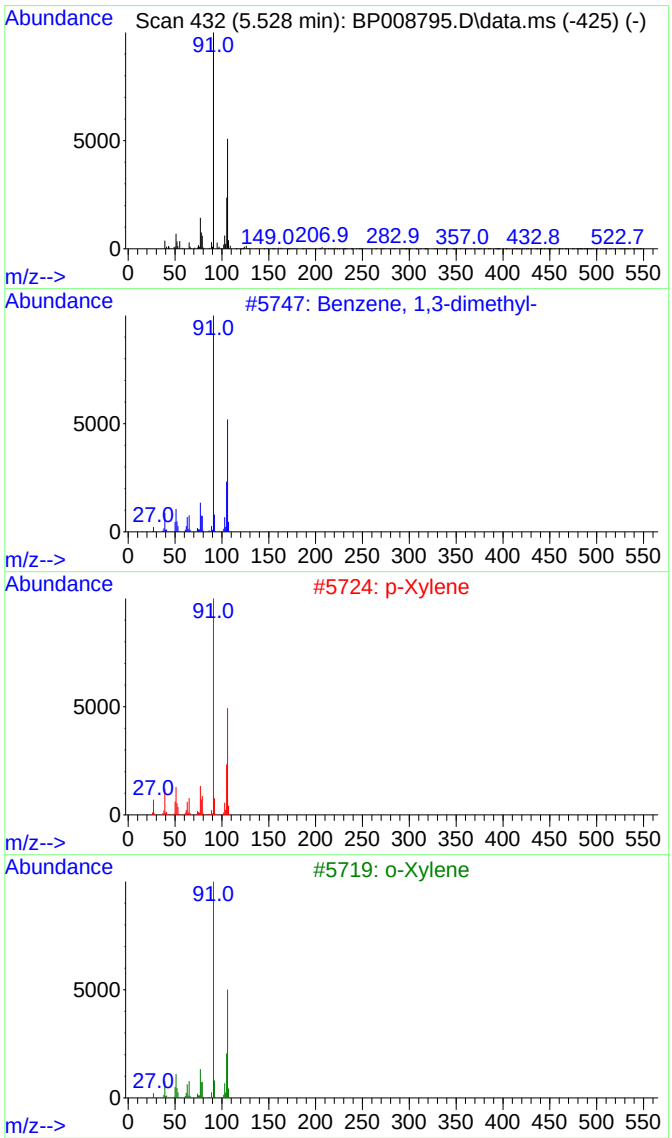
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

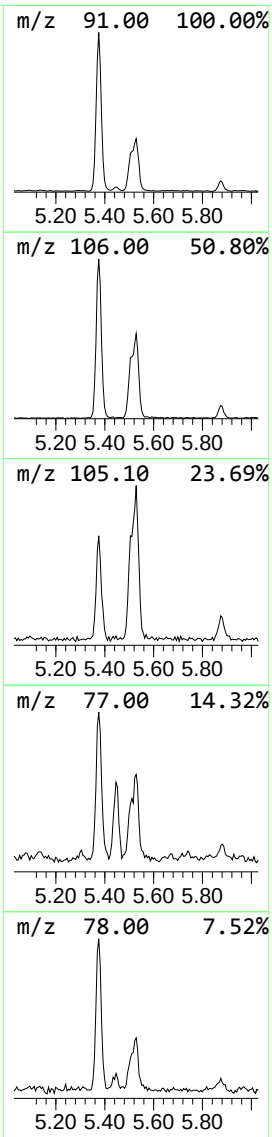
****

Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.528	3.40 ng	347929	1,4-Dichlorobenzene-d4	7.863

Hit# of 4	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	90
2	p-Xylene	106	C8H10	000106-42-3	90
3	o-Xylene	106	C8H10	000095-47-6	90
4	Ethylbenzene	106	C8H10	000100-41-4	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011322\
Data File : BP008795.D
Acq On : 13 Jan 2022 17:57
Operator : CG/JU
Sample : N1119-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-NORTH-SIDEWALL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

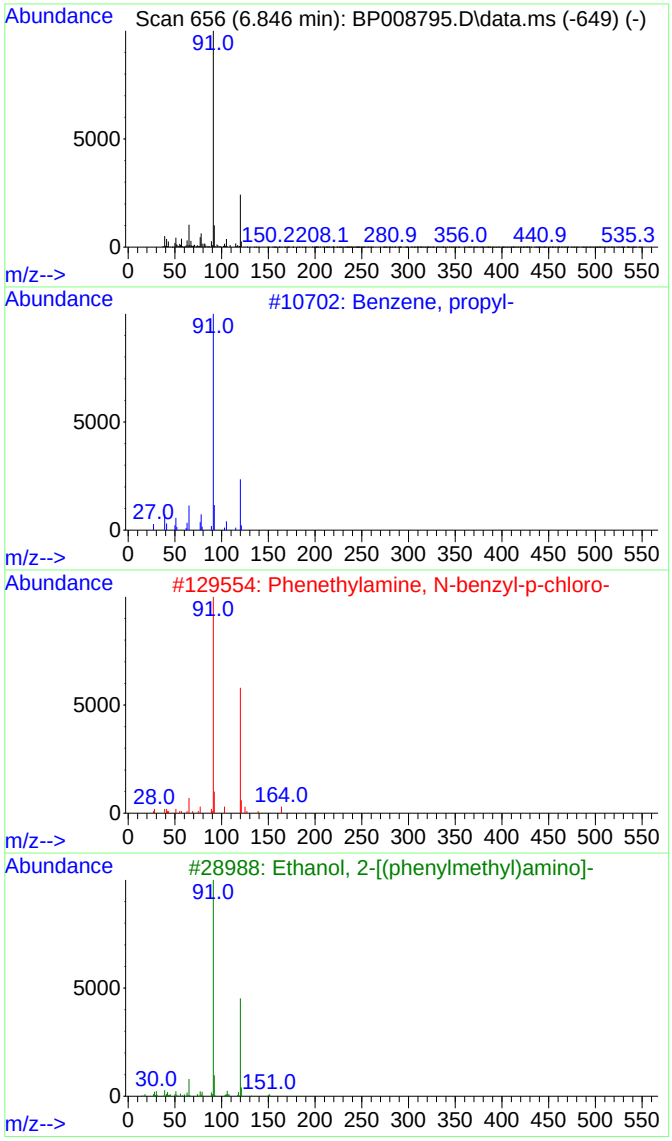
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

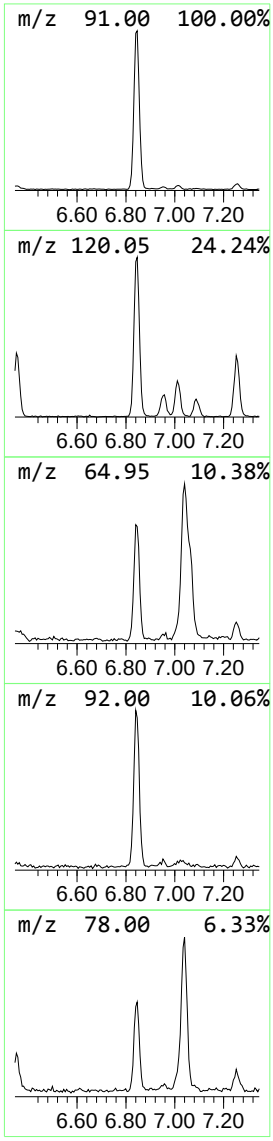
****

Peak Number 4 Benzene, propyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.846	4.20 ng	429618	1,4-Dichlorobenzene-d4	7.863

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	93
2	Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	64
3	Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	64
4	1-Benzylamino-2-benzyloxyethane	241	C16H19NO	038336-06-0	64
5	Benzeneacetaldehyde	120	C8H8O	000122-78-1	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011322\
Data File : BP008795.D
Acq On : 13 Jan 2022 17:57
Operator : CG/JU
Sample : N1119-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-NORTH-SIDEWALL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

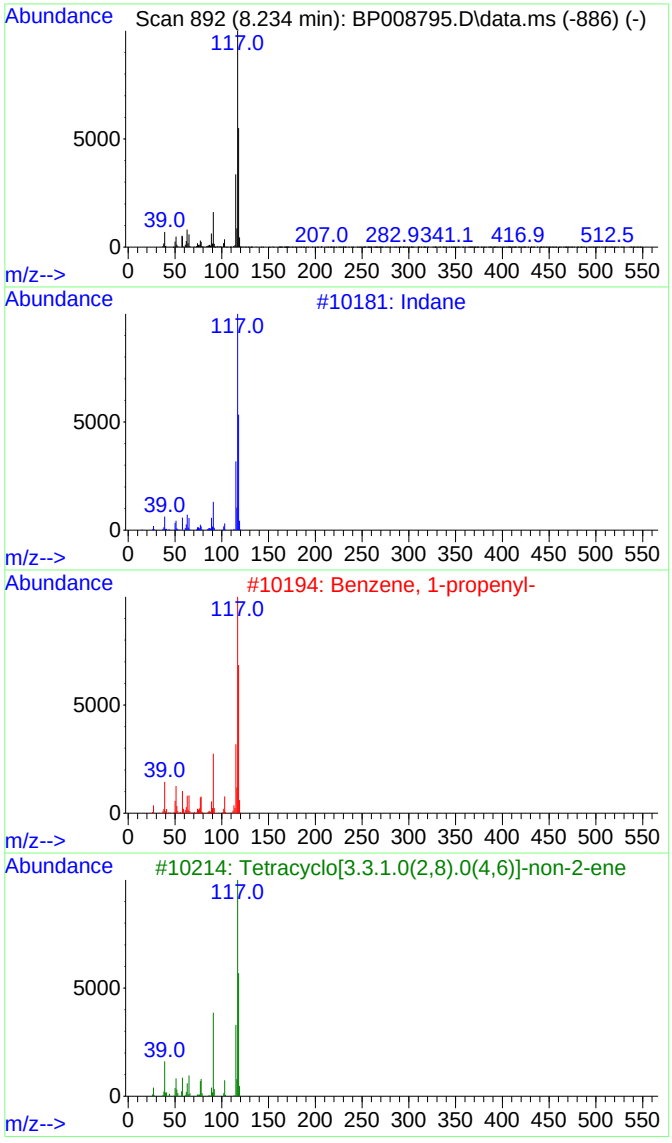
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

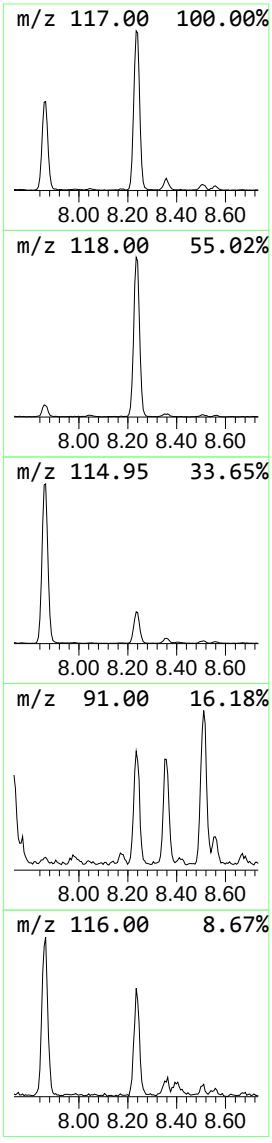
****

Peak Number 5 Indane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.234	3.37 ng	345066	1,4-Dichlorobenzene-d4	7.863

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	95
2	Benzene, 1-propenyl-	118	C9H10	000637-50-3	83
3	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
4	Deltacyclene	118	C9H10	007785-10-6	72
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011322\
Data File : BP008795.D
Acq On : 13 Jan 2022 17:57
Operator : CG/JU
Sample : N1119-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-NORTH-SIDEWALL

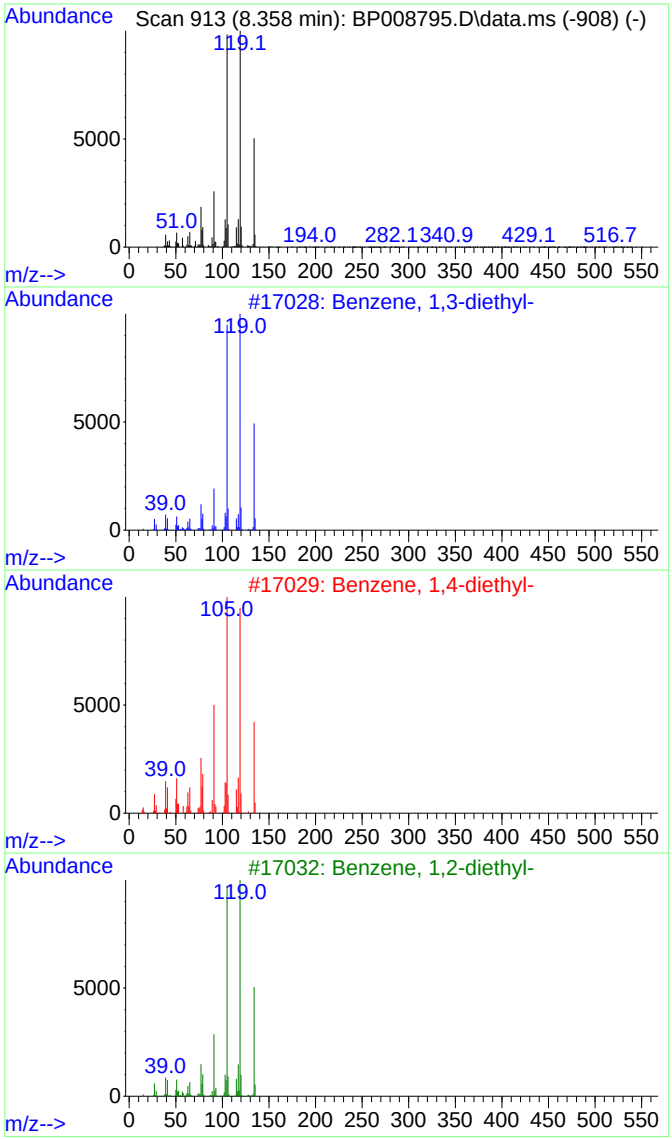
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

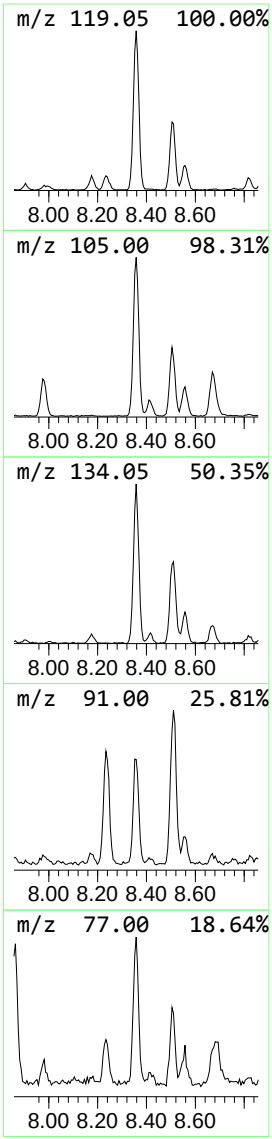
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,3-diethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.358	2.82 ng	288561	1,4-Dichlorobenzene-d4	7.863

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
3	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	94
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5	o-Cymene	134	C10H14	000527-84-4	81





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011322\
Data File : BP008795.D
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ALS Vial : 11 Sample Multiplier: 1

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BNA_P
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NORTH-EAST-CORNER-EXCAVATION-NORTH-SIDEWALL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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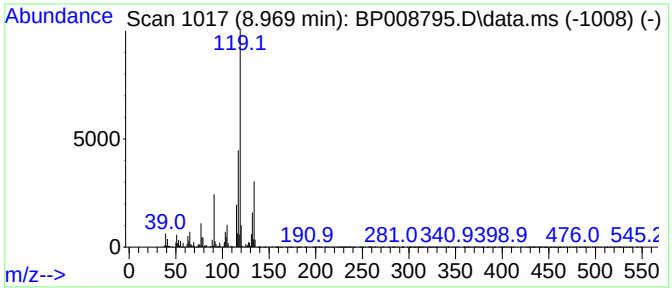
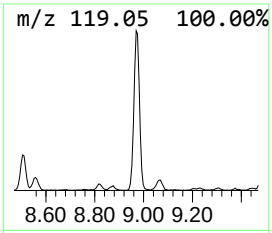
****

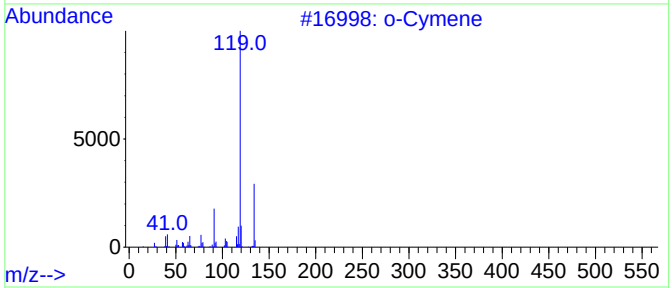
Peak Number 7 o-Cymene Concentration Rank 6

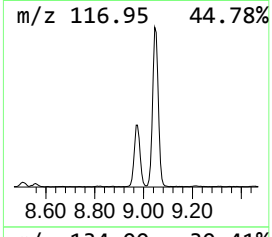
R.T.	EstConc	Area	Relative to ISTD	R.T.
8.969	5.09 ng	520937	1,4-Dichlorobenzene-d4	7.863

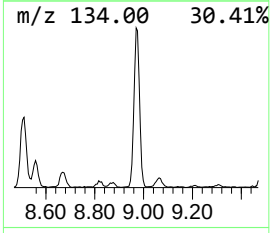
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	64
2	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	64
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	64
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	64
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	64

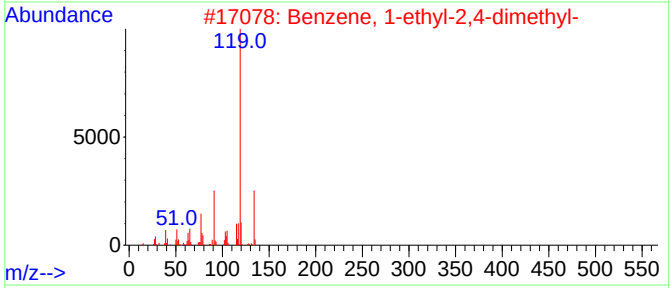


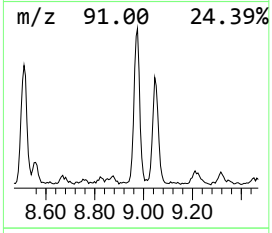


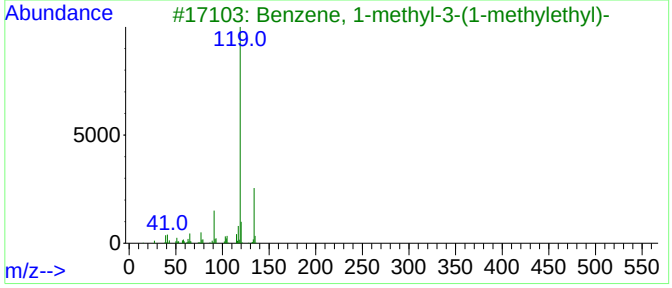


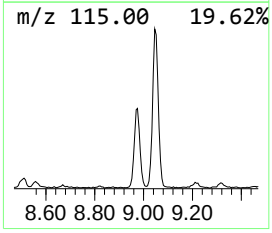












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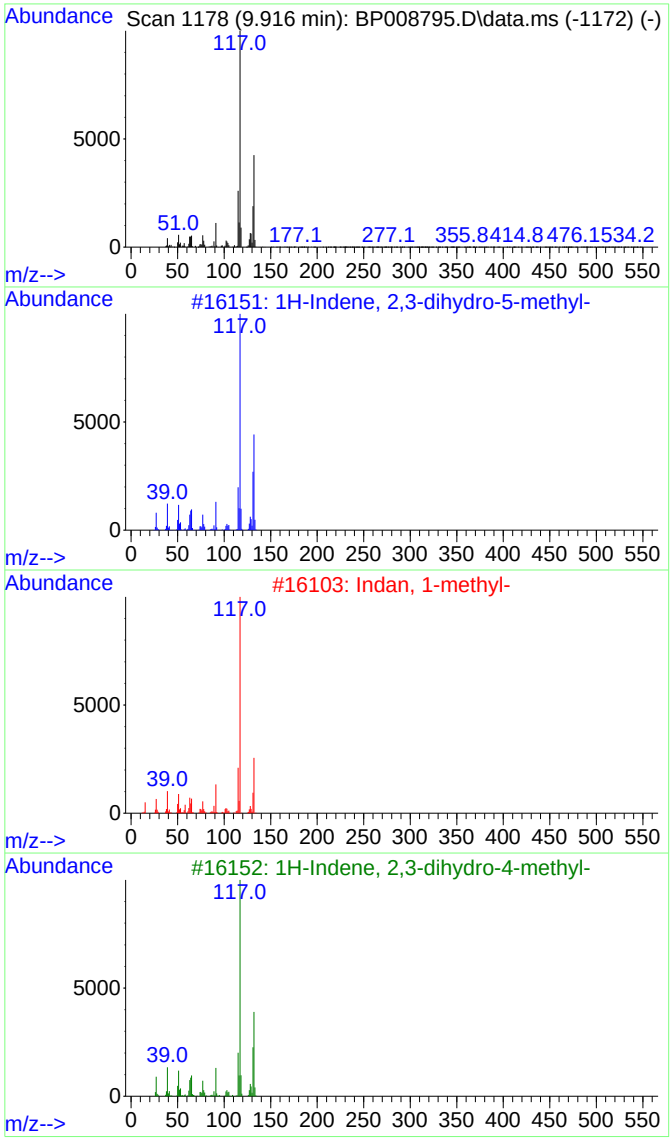
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

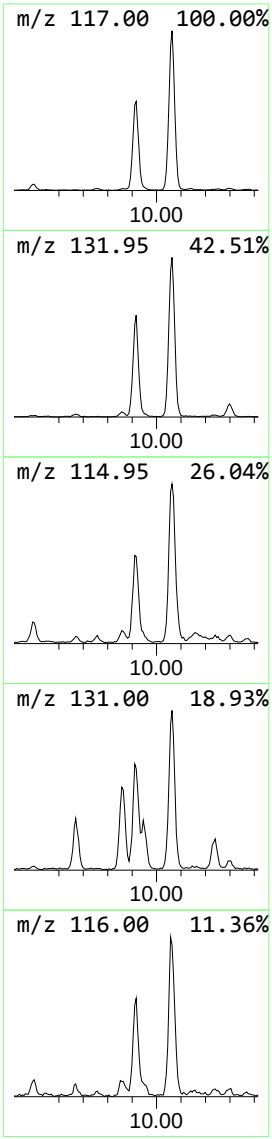
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.916	2.06 ng	284879	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2	Indan, 1-methyl-	132	C10H12	000767-58-8	90
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011322\
Data File : BP008795.D
Acq On : 13 Jan 2022 17:57
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ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-NORTH-SIDEWALL

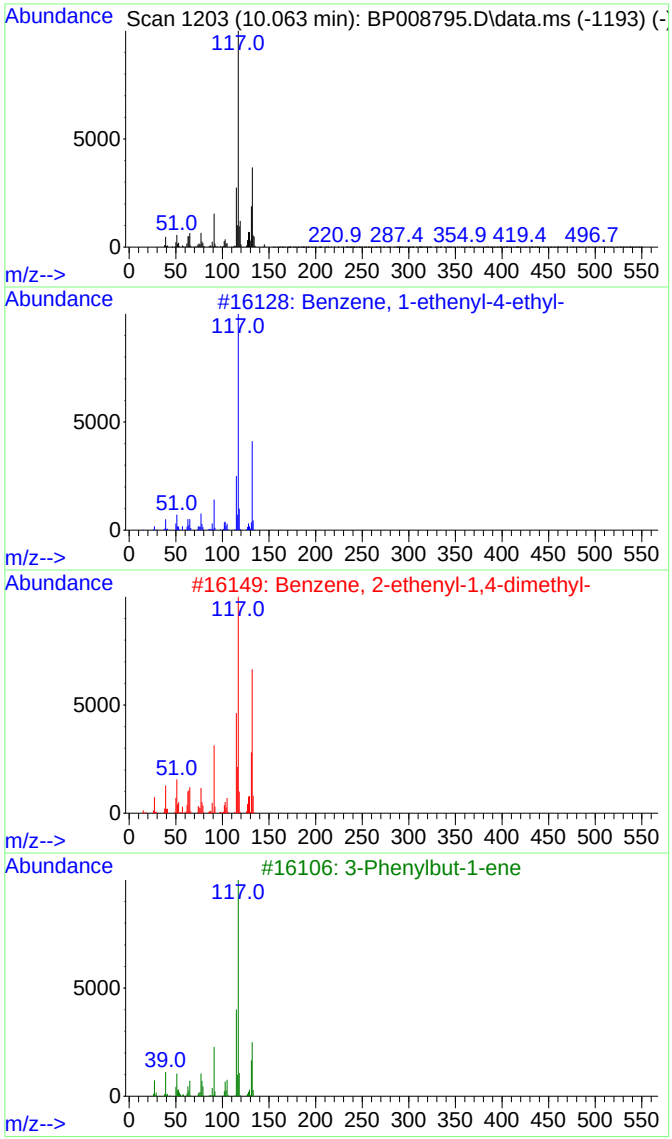
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

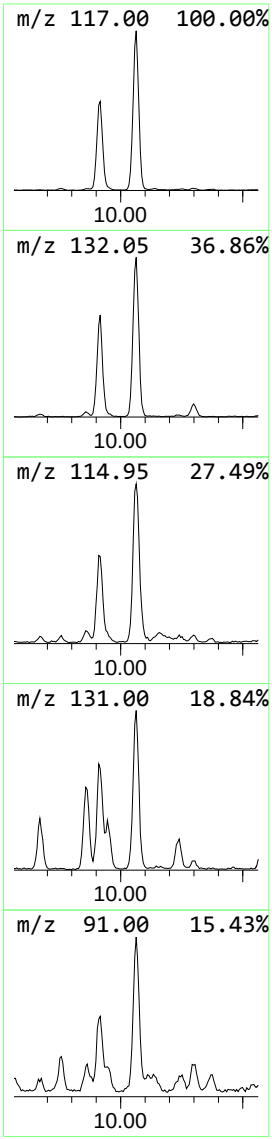
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.063	5.02 ng	695328	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	90
5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87





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Data File : BP008795.D
Acq On : 13 Jan 2022 17:57
Operator : CG/JU
Sample : N1119-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-NORTH-SIDEWALL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.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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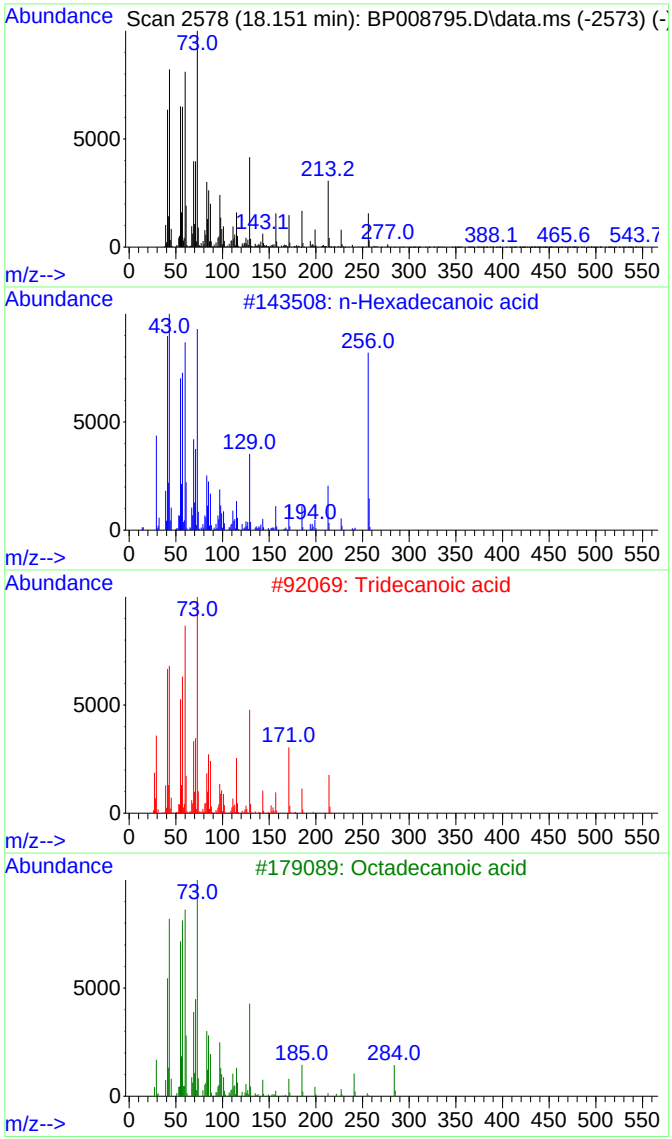
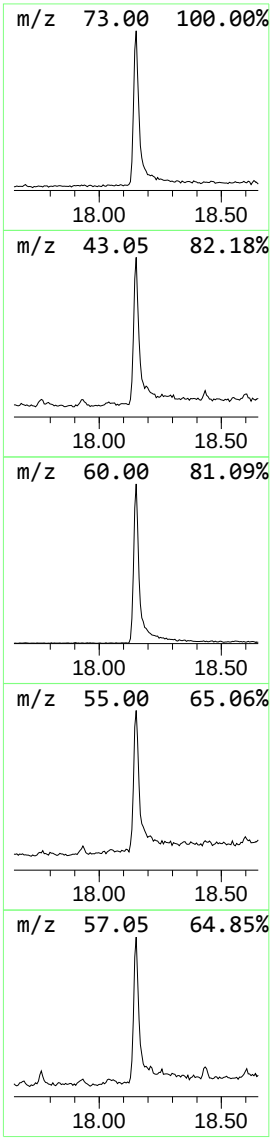
Peak Number 10 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.151	5.78 ng	1037240	Phenanthrene-d10	17.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tridecanoic acid	214	C13H26O2	000638-53-9	93
3	Octadecanoic acid	284	C18H36O2	000057-11-4	90
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	86





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Data File : BP008795.D
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Operator : CG/JU
Sample : N1119-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-NORTH-SIDEWALL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.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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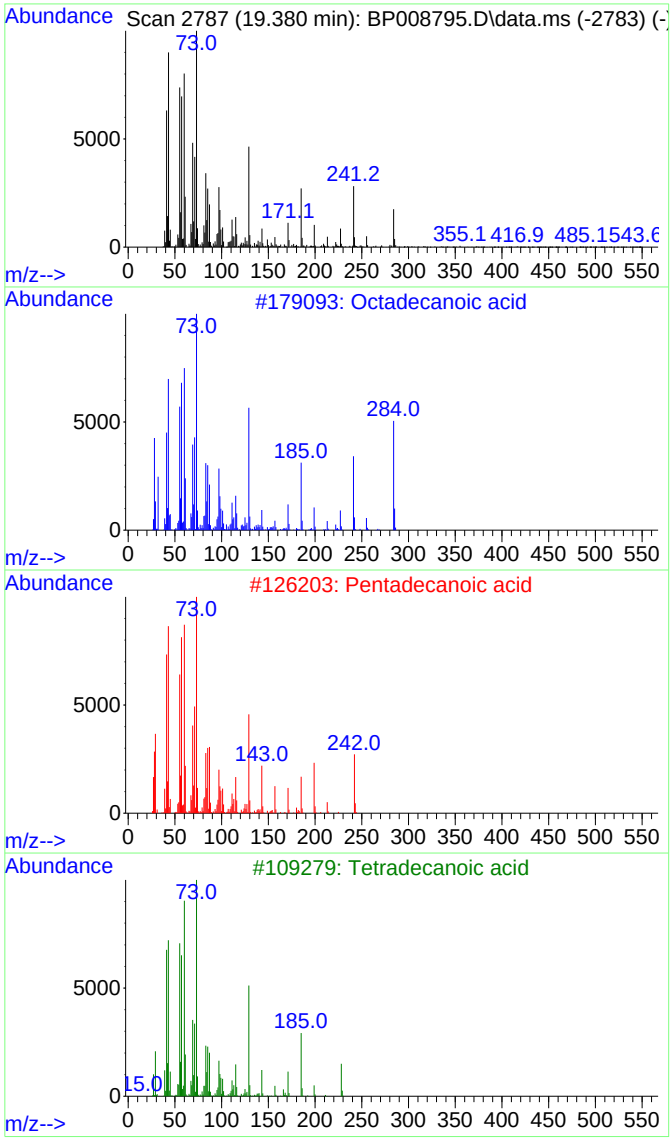
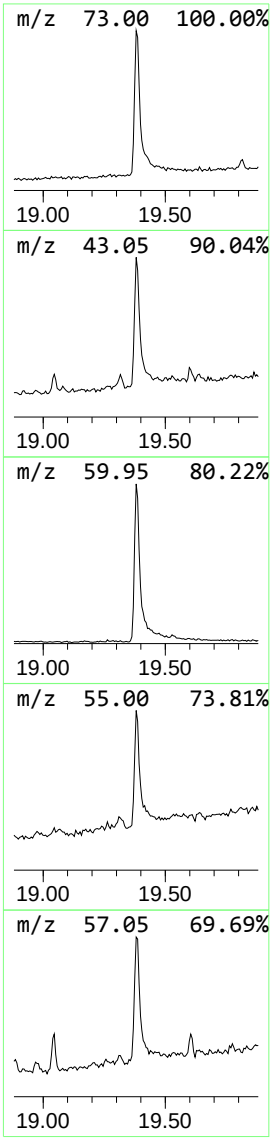
Peak Number 11 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.381	5.45 ng	1055390	Chrysene-d12	21.345

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	89
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
5	n-Decanoic acid	172	C10H20O2	000334-48-5	70





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Data File : BP008795.D
Acq On : 13 Jan 2022 17:57
Operator : CG/JU
Sample : N1119-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-NORTH-SIDEWALL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.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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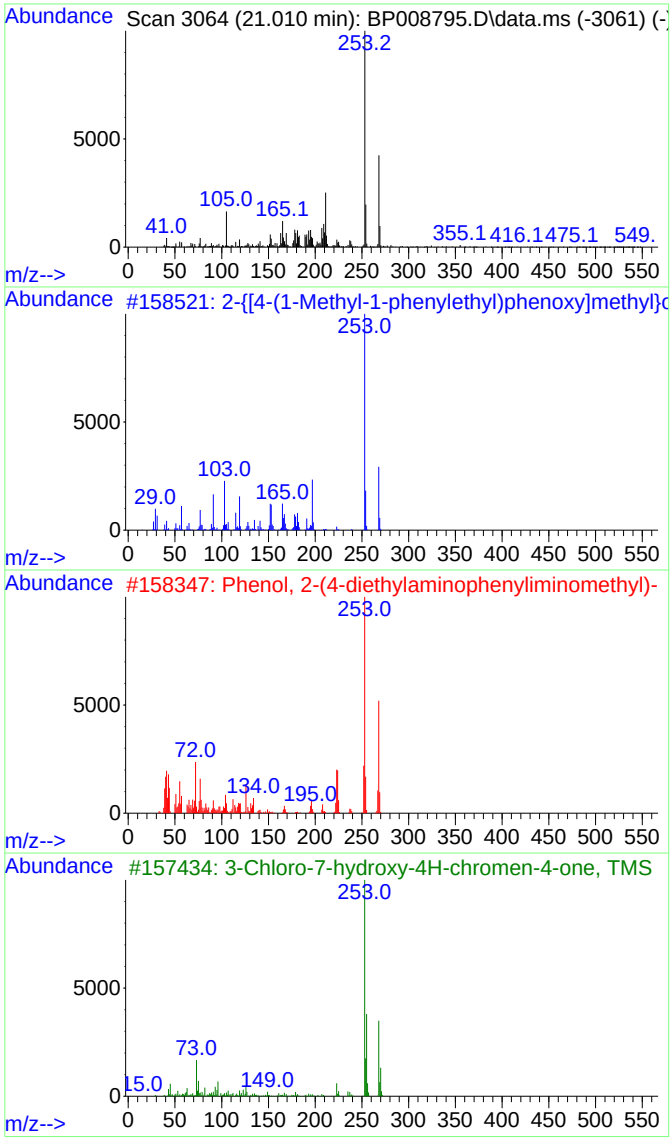
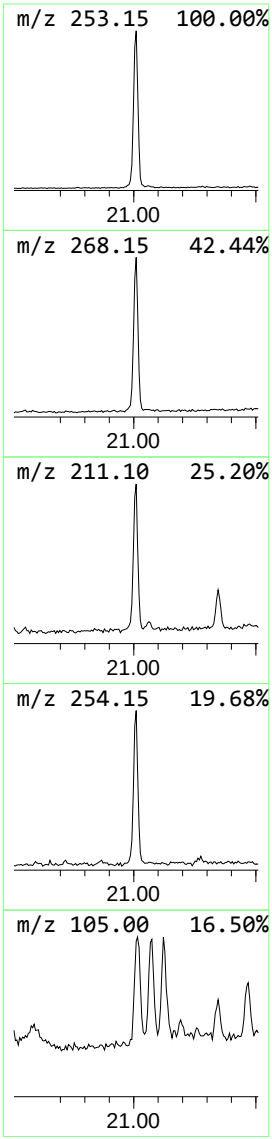
Peak Number 12 2-{[4-(1-Methyl-1-phenyleth... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.010	2.64 ng	511553	Chrysene-d12	21.345

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-{[4-(1-Methyl-1-phenylethyl)ph...	268	C18H20O2	061578-04-9	76
2	Phenol, 2-(4-diethylaminophenyli...	268	C17H20N2O	004938-45-8	74
3	3-Chloro-7-hydroxy-4H-chromen-4-...	268	C12H13ClO3Si	1000514-35-2	64
4	ethanone, 1,1'-dibenzo[b,d]thien...	268	C16H12O2S	1000401-69-1	64
5	1H-Inden-5-ol, 2,3-dihydro-3-(4-...	268	C18H20O2	010527-11-4	64





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Acq On : 13 Jan 2022 17:57
Operator : CG/JU
Sample : N1119-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-NORTH-SIDEWALL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.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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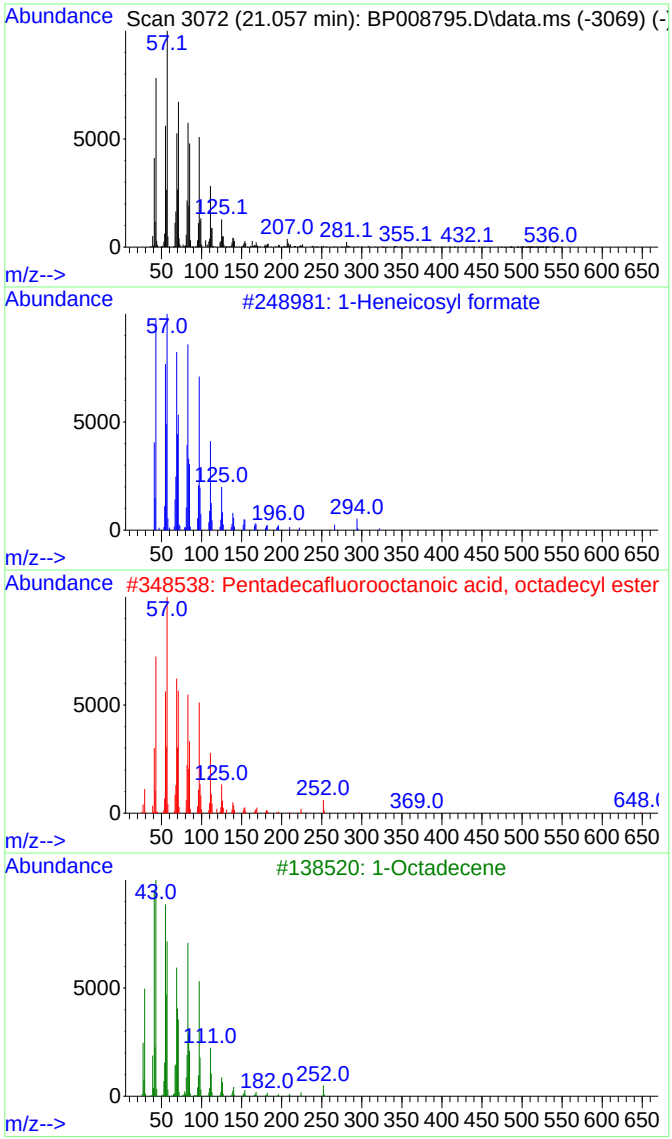
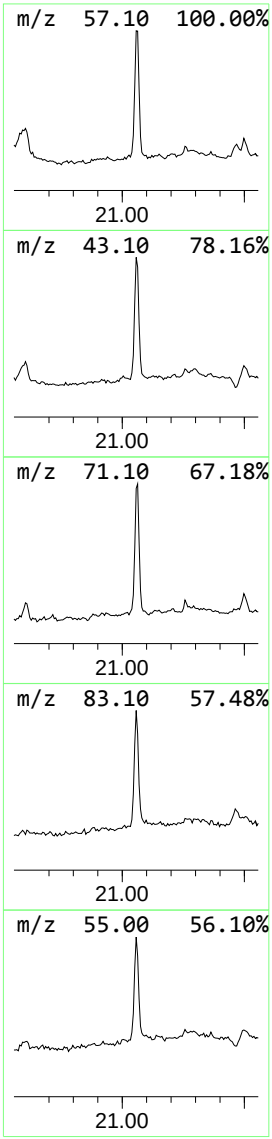
Peak Number 13 1-Heneicosyl formate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.057	5.65 ng	1093800	Chrysene-d12	21.345

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate	340	C22H44O2	077899-03-7	93
2	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
3	1-Octadecene	252	C18H36	000112-88-9	91
4	Methoxyacetic acid, heptadecyl e...	328	C20H40O3	1000282-99-1	91
5	1-Docosene	308	C22H44	001599-67-3	91





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Data File : BP008795.D
Acq On : 13 Jan 2022 17:57
Operator : CG/JU
Sample : N1119-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-NORTH-SIDEWALL

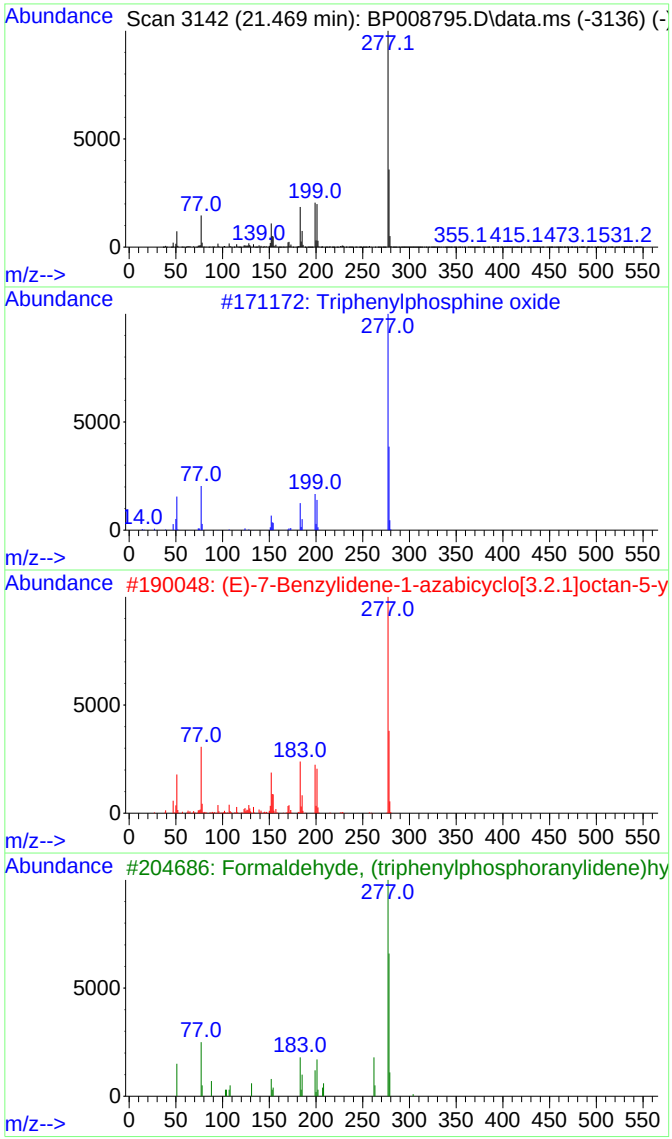
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

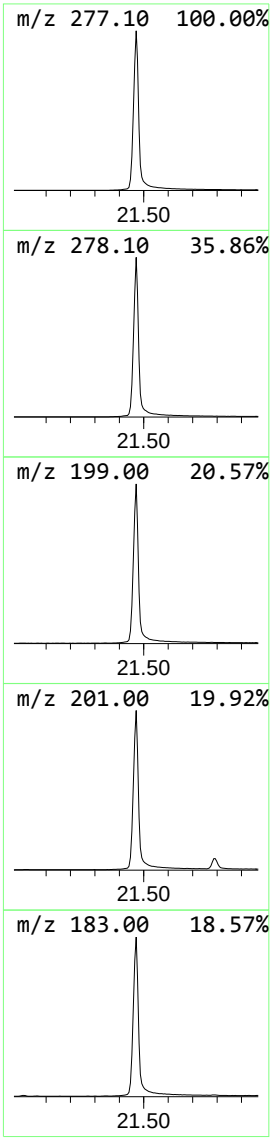
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Triphenylphosphine oxide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.469	142.93 ng	27690600	Chrysene-d12	21.345

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	96
2	(E)-7-Benzylidene-1-azabicyclo[3...	293	C15H19NO3S	1000483-83-4	80
3	Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	45
4	5-Methyl-8-phenylfuro[3',2':5,6]...	277	C16H11N3O2	1000460-07-9	25
5	Vanadium, cycloheptatrienyl-(pen...	277	C17H22V	1000155-20-5	9





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011322\
Data File : BP008795.D
Acq On : 13 Jan 2022 17:57
Operator : CG/JU
Sample : N1119-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
NORTH-EAST-CORNER-EXCAVATION-NORTH-SIDEWALL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.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...	3.064	32.7	ng	3342740	1	7.863	2046940	20.0
Benzene, 1,3-di...	5.528	3.4	ng	347929	1	7.863	2046940	20.0
Benzene, propyl-	6.846	4.2	ng	429618	1	7.863	2046940	20.0
Indane	8.234	3.4	ng	345066	1	7.863	2046940	20.0
Benzene, 1,3-di...	8.358	2.8	ng	288561	1	7.863	2046940	20.0
o-Cymene	8.969	5.1	ng	520937	1	7.863	2046940	20.0
1H-Indene, 2,3-...	9.916	2.1	ng	284879	2	10.663	2769870	20.0
Benzene, 1-ethe...	10.063	5.0	ng	695328	2	10.663	2769870	20.0
n-Hexadecanoic ...	18.151	5.8	ng	1037240	4	17.251	3586840	20.0
Octadecanoic acid	19.381	5.5	ng	1055390	5	21.345	3874850	20.0
2-{[4-(1-Methyl...	21.010	2.6	ng	511553	5	21.345	3874850	20.0
1-Heneicosyl fo...	21.057	5.7	ng	1093800	5	21.345	3874850	20.0
Triphenylphosph...	21.469	142.9	ng	27690600	5	21.345	3874850	20.0