LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031522\
Data File : BP009414.D
Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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: BP009414.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	22	rVB	127051	213821	2.07%	0.193%
2	5.340	393	400	407	rBV	232763	377280	3.65%	0.340%
3	5.422	407	414	432	rVB	1269602	2225754	21.56%	2.008%
4	5.816	472	481	493	rBV	339964	605660	5.87%	0.547%
5	6.316	559	566	573	rBV	143773	252487	2.45%	0.228%
6	6.799	641	648	654	rBV	78464	133295	1.29%	0.120%
7	7.010	671	684	698	rBV	1147280	2265659	21.94%	2.044%
8	7.810	811	820	833	rBV	1122417	1983245	19.21%	1.790%
9	8.969	1010	1017	1028	rBV	775975	1459951	14.14%	1.317%
10	10.604	1287	1295	1303	rBV	1459124	2738421	26.52%	2.471%
11	13.075	1708	1715	1732	rBV	2354187	3764060	36.45%	3.397%
12	13.286	1745	1751	1762	rBV2	139631	264264	2.56%	0.238%
13	13.504	1783	1788	1798	rVB	218006	341710	3.31%	0.308%
14	14.175	1896	1902	1917	rVB	369757	673279	6.52%	0.608%
15	14.369	1931	1935	1939	rVB	102842	149616	1.45%	0.135%
16	14.457	1942	1950	1953	rVV	2293131	3767836	36.49%	3.400%
17	14.492	1953	1956	1971	rVB	837559	1238005	11.99%	1.117%
18	15.322	2093	2097	2102	rVB2	119210	153066	1.48%	0.138%
19	15.733	2164	2167	2172	rVB	93833	134662	1.30%	0.122%
20	15.951	2197	2204	2212	rBV	1978839	3164630	30.65%	2.856%
21	16.192	2240	2245	2247	rBV	198095	296072	2.87%	0.267%
22	16.398	2276	2280	2289	rVB	320980	464592	4.50%	0.419%
23	16.563	2305	2308	2312	rVB	104114	129749	1.26%	0.117%
24	16.822	2346	2352	2357	rVB	1303737	1805760	17.49%	1.629%
25	16.992	2377	2381	2384	rBV2	96374	136922	1.33%	0.124%
26	17.039	2385	2389	2398	rVB2	163861	298951	2.90%	0.270%
27	17.216	2413	2419	2423	rBV	2984745	4520866	43.78%	4.079%
28	17.257	2423	2426	2432	rVB	570961	833363	8.07%	0.752%
29	17.351	2437	2442	2448	rVB	431325	697891	6.76%	0.630%
30	17.627	2485	2489	2493	rBV	168875	236105	2.29%	0.213%
31	17.733	2504	2507	2514	rVB2	104574	139806	1.35%	0.126%
32	18.121	2569	2573	2581	rBV2	944794	1423329	13.78%	1.284%
33	18.610	2654	2656	2661	rVB	142925	167264	1.62%	0.151%
34	19.116	2739	2742	2749	rVB	241338	348209	3.37%	0.314%
35	19.239	2757	2763	2771	rBV	2941378	4390373	42.52%	3.962%
36	19.363	2780	2784	2792	rBV2	969449	1840258	17.82%	1.661%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031522\
Data File : BP009414.D
Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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

37	19.592	2814	2823	2831	rBV	2317577	3936448	38.12%	3.552%
38	19.792	2852	2857	2861	rBV	3891550	5072900	49.13%	4.578%
39	21.027	3063	3067	3076	rBV	2168116	3850668	37.29%	3.475%
40	21.245	3100	3104	3109	rVB	8854097	10325357	100.00%	9.317%
41	21.321	3111	3117	3121	rBV2	4493881	7902035	76.53%	7.130%
42	21.445	3134	3138	3143	rBV	2431900	3639497	35.25%	3.284%
43	21.504	3143	3148	3152	rVV	4265809	6144822	59.51%	5.545%
44	21.557	3153	3157	3162	rVV	5480377	8011407	77.59%	7.229%
45	21.604	3162	3165	3168	rVV	2010670	2765545	26.78%	2.496%
46	21.633	3168	3170	3179	rVB3	2086672	2970494	28.77%	2.680%
47	23.009	3399	3404	3419	rVB	1587128	5132681	49.71%	4.632%
48	23.633	3506	3510	3519	rVB	1193922	2484314	24.06%	2.242%
49	23.739	3522	3528	3534	rBV2	2291616	4948568	47.93%	4.465%

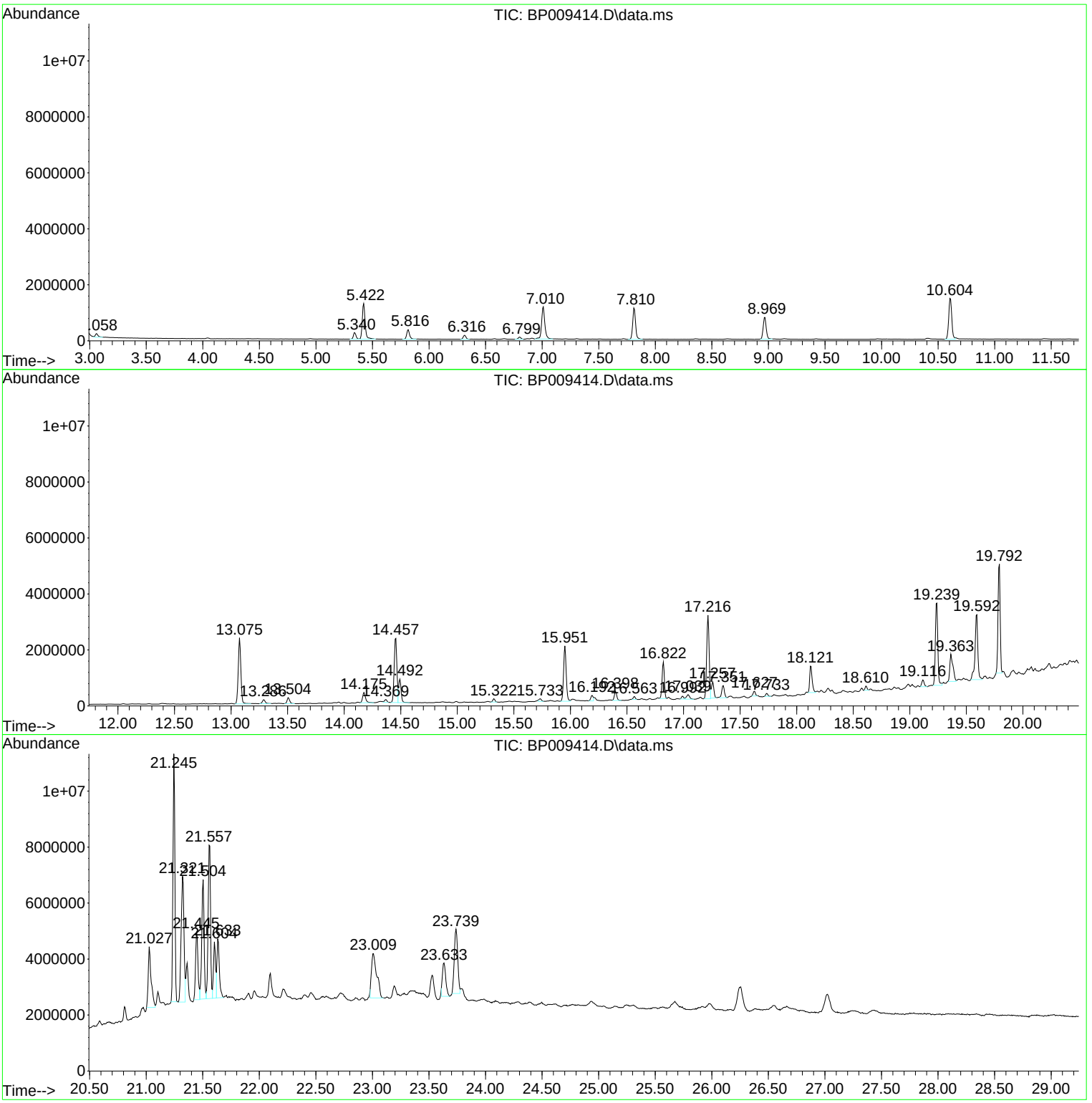
Sum of corrected areas: 110820947

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031522\
Data File : BP009414.D
Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031522\
Data File : BP009414.D
Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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

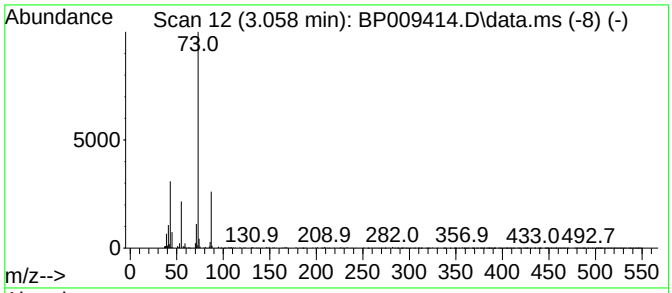
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

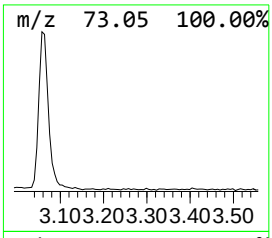
****

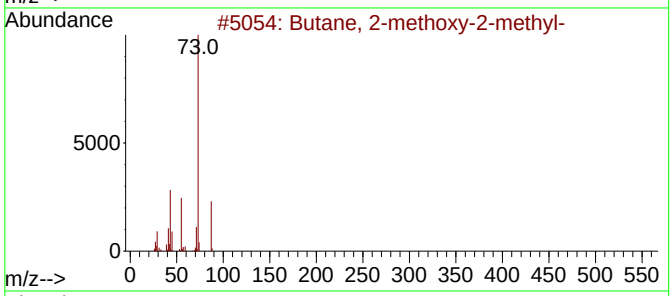
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 13

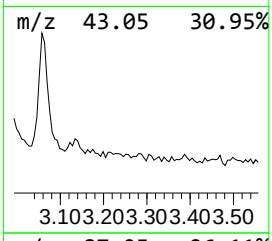
R.T.	EstConc	Area	Relative to ISTD	R.T.
3.058	2.16 ng	213821	1,4-Dichlorobenzene-d4	7.810

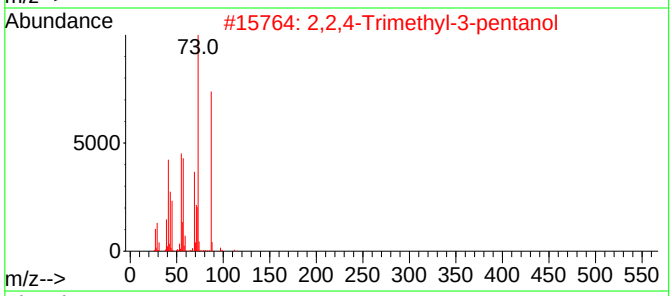
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	38
3	3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	28
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5	1,3-Dioxolane, 2-(6-heptynyl)-	168	C10H16O2	025595-82-8	9

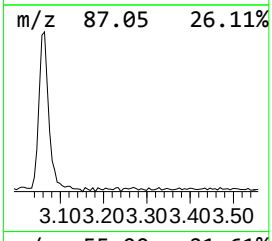


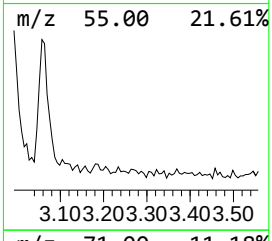


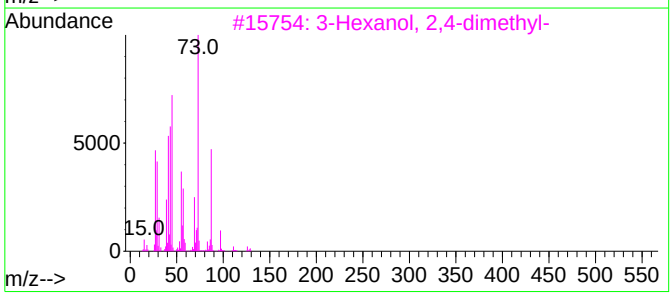


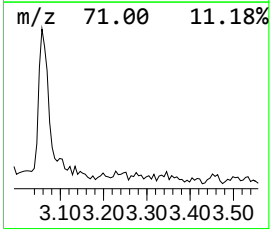












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031522\
Data File : BP009414.D
Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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

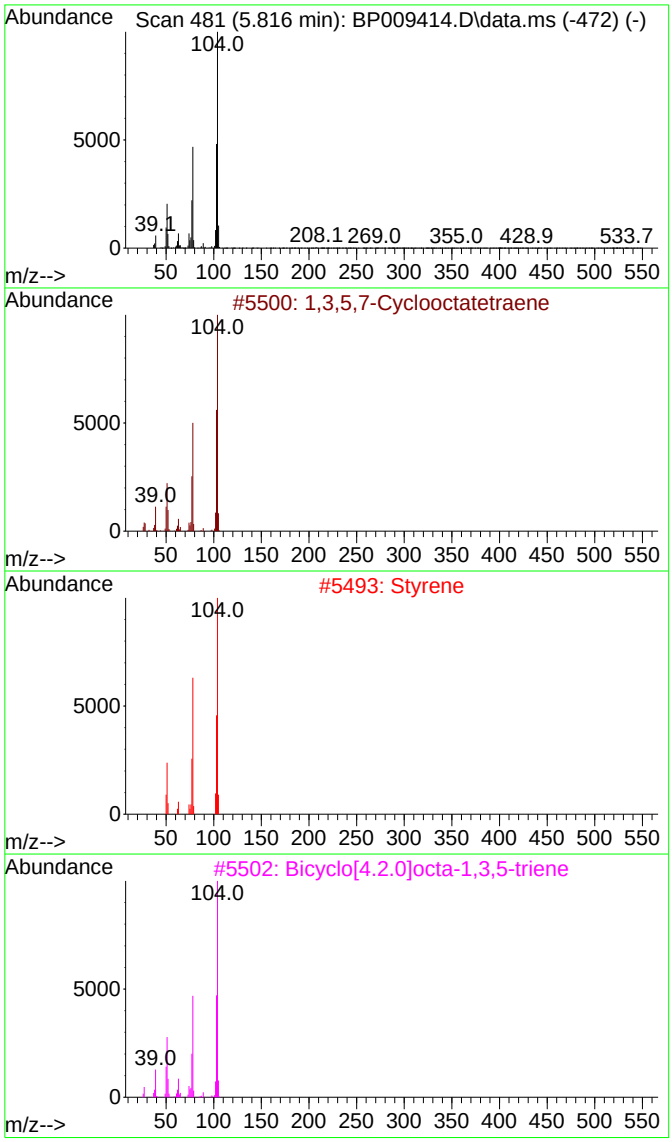
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

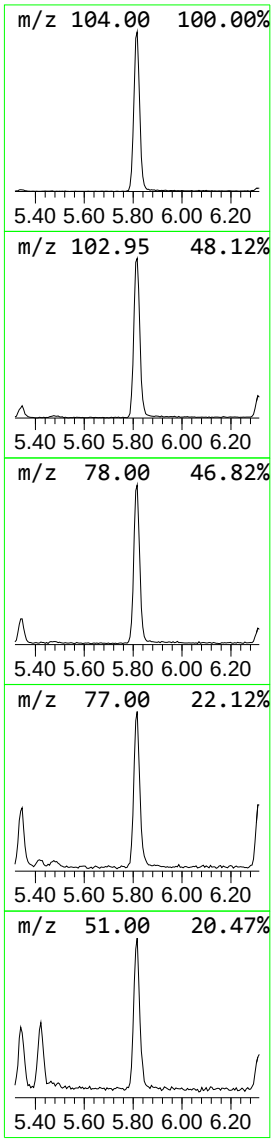
****

Peak Number 3 1,3,5,7-Cyclooctatetraene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.816	6.11 ng	605660	1,4-Dichlorobenzene-d4	7.810

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,5,7-Cyclooctatetraene	104	C8H8	000629-20-9	97
2	Styrene	104	C8H8	000100-42-5	97
3	Bicyclo[4.2.0]octa-1,3,5-triene	104	C8H8	000694-87-1	91
4	Butanedioic acid, phenyl-	194	C10H10O4	000635-51-8	72
5	2-Propenoic acid, 3-phenyl-, 2-p...	252	C17H16O2	000103-53-7	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031522\
Data File : BP009414.D
Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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

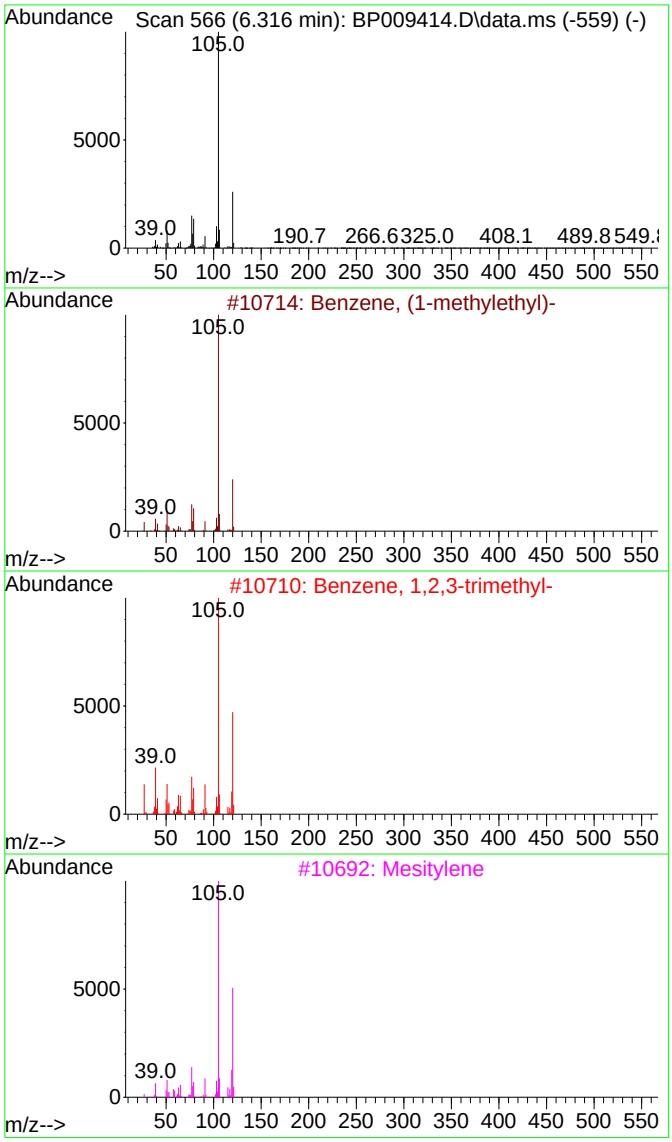
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

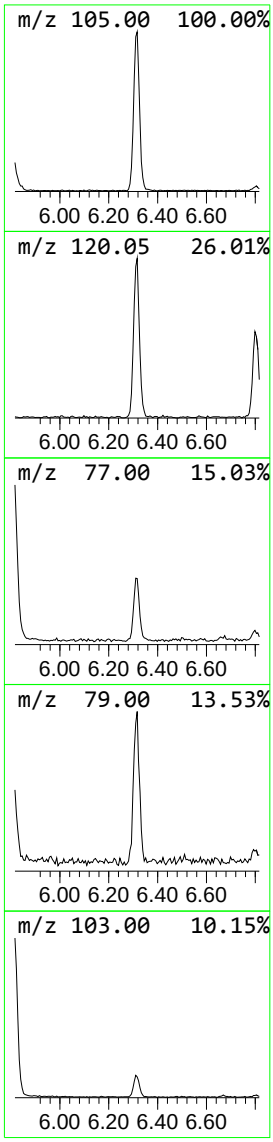
****

Peak Number 4 Benzene, (1-methylethyl)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.316	2.55 ng	252487	1,4-Dichlorobenzene-d4	7.810

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3	Mesitylene	120	C9H12	000108-67-8	90
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	87
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031522\
Data File : BP009414.D
Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

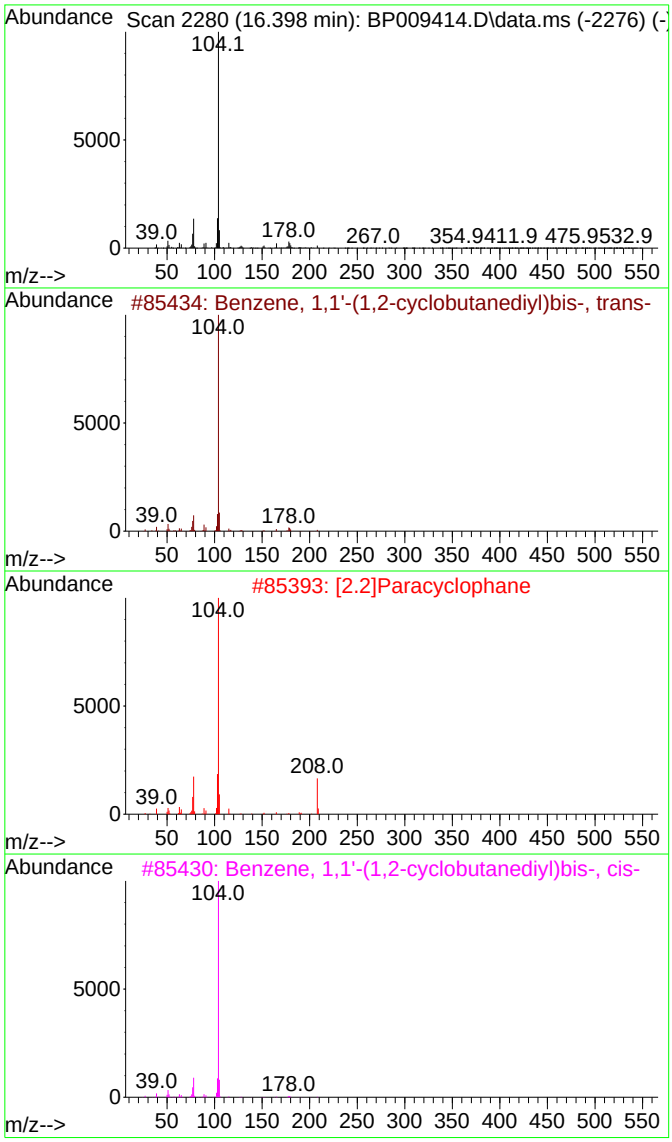
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

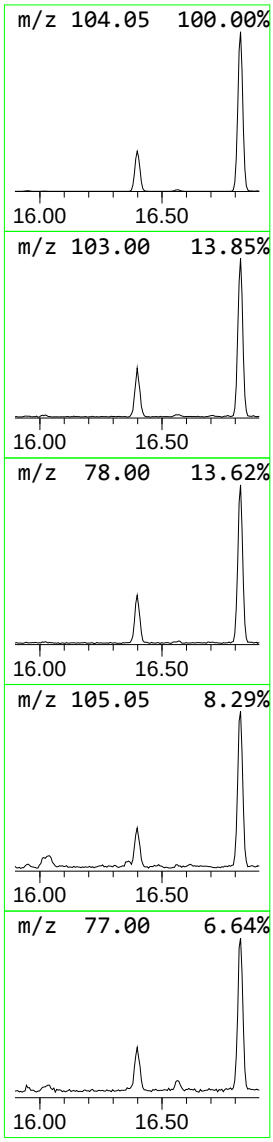
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1,1'-(1,2-cyclobut... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.398	2.06 ng	464592	Phenanthrene-d10	17.216

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	90
2	[2.2]Paracyclophane	208	C16H16	001633-22-3	87
3	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	007694-30-6	86
4	Cyclobutane, 1,2-diphenyl-	208	C16H16	003018-21-1	78
5	Styrene	104	C8H8	000100-42-5	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031522\
Data File : BP009414.D
Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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

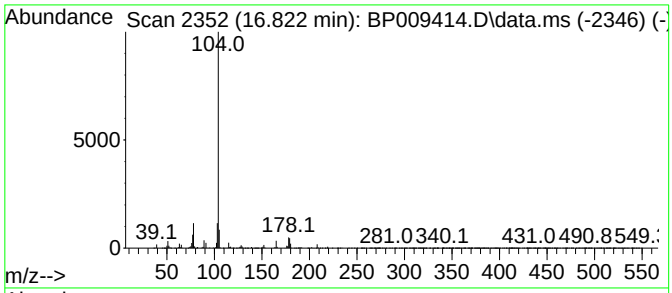
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

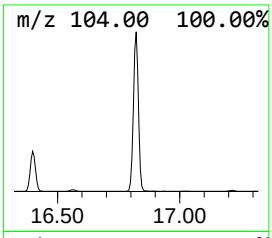
****

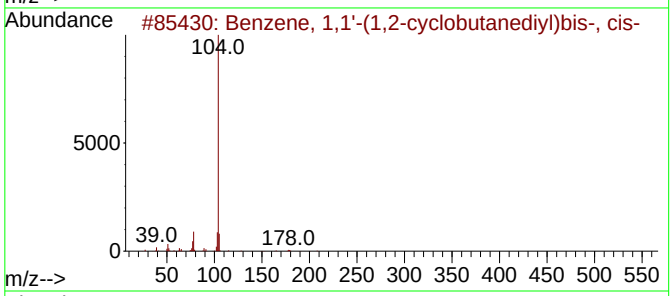
Peak Number 6 [2.2]Paracyclophane Concentration Rank 5

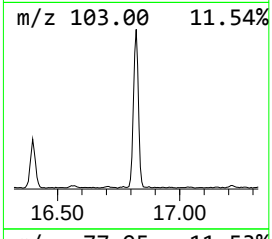
R.T.	EstConc	Area	Relative to ISTD	R.T.
16.822	7.99 ng	1805760	Phenanthrene-d10	17.216

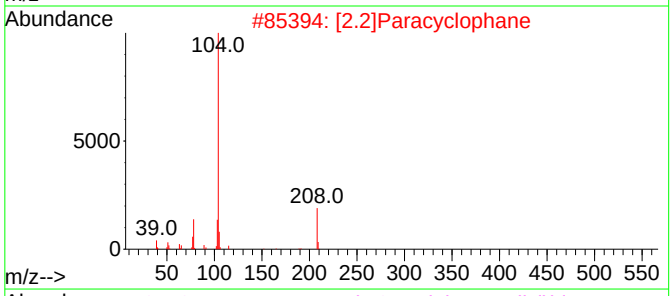
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	007694-30-6	86
2	[2.2]Paracyclophane	208	C16H16	001633-22-3	74
3	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	72
4	5,5-Dimethyl-4-phenyldihydrofura...	190	C12H14O2	013133-96-5	59
5	Carphedon, N-(trifluoroacetyl)	314	C14H13F3N2O3	1000445-42-6	50

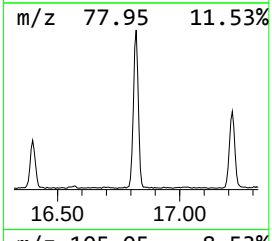


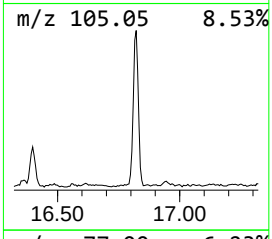


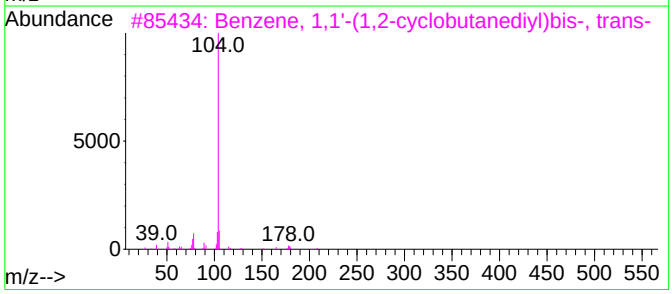


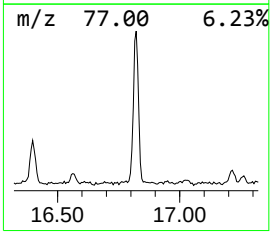












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031522\
Data File : BP009414.D
Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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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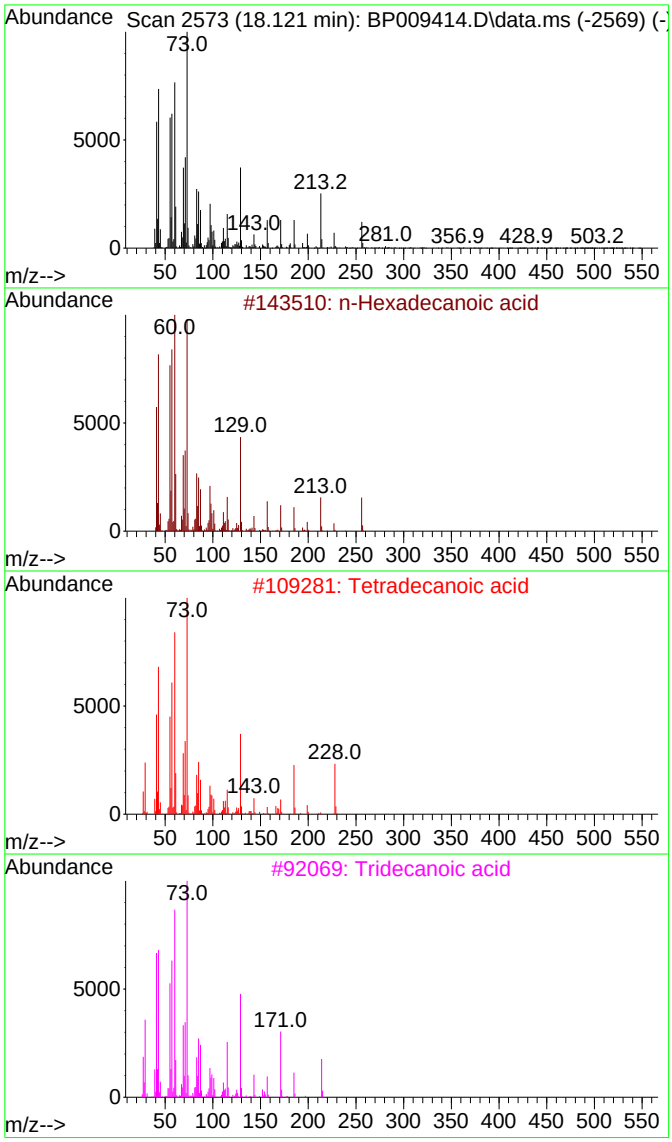
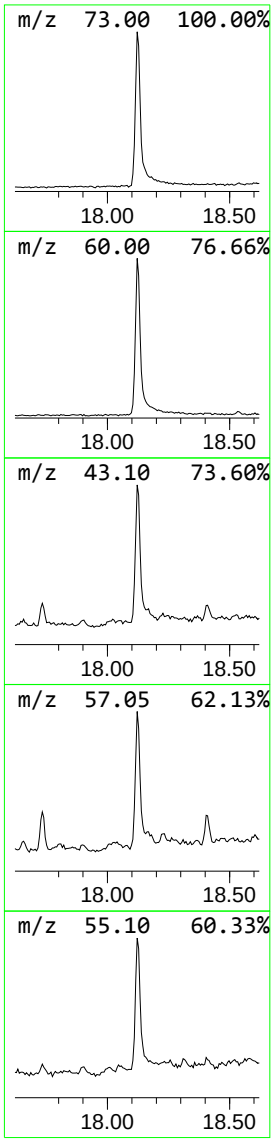
Peak Number 7 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.121	6.30 ng	1423330	Phenanthrene-d10	17.216

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3	Tridecanoic acid	214	C13H26O2	000638-53-9	93
4	Octadecanoic acid	284	C18H36O2	000057-11-4	87
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	76





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Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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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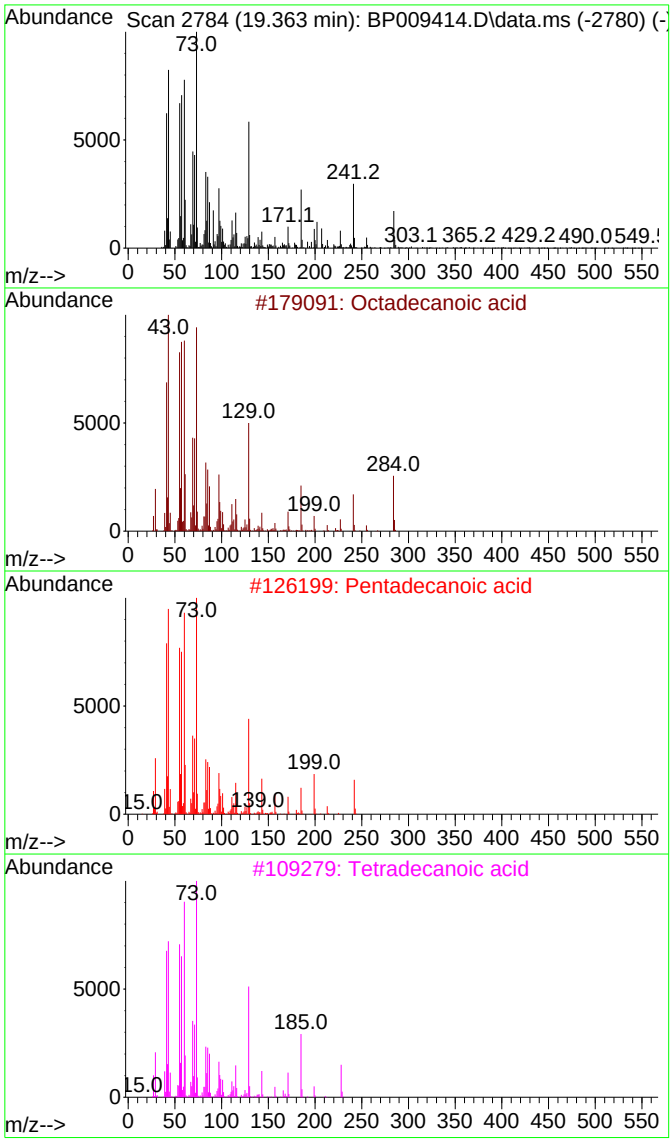
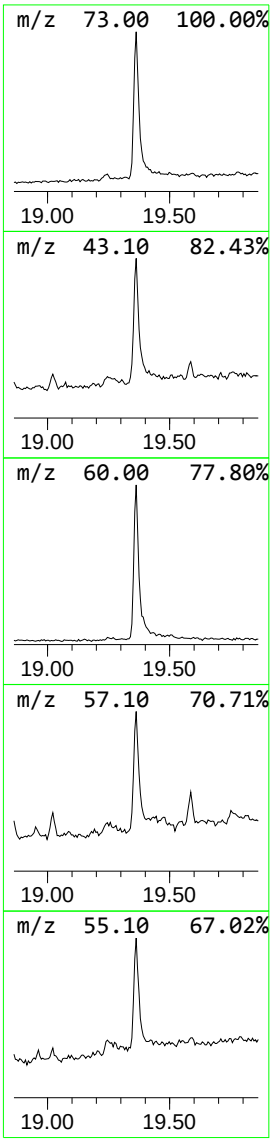
Peak Number 8 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.363	4.66 ng	1840260	Chrysene-d12	21.327

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	91
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	72
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64





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Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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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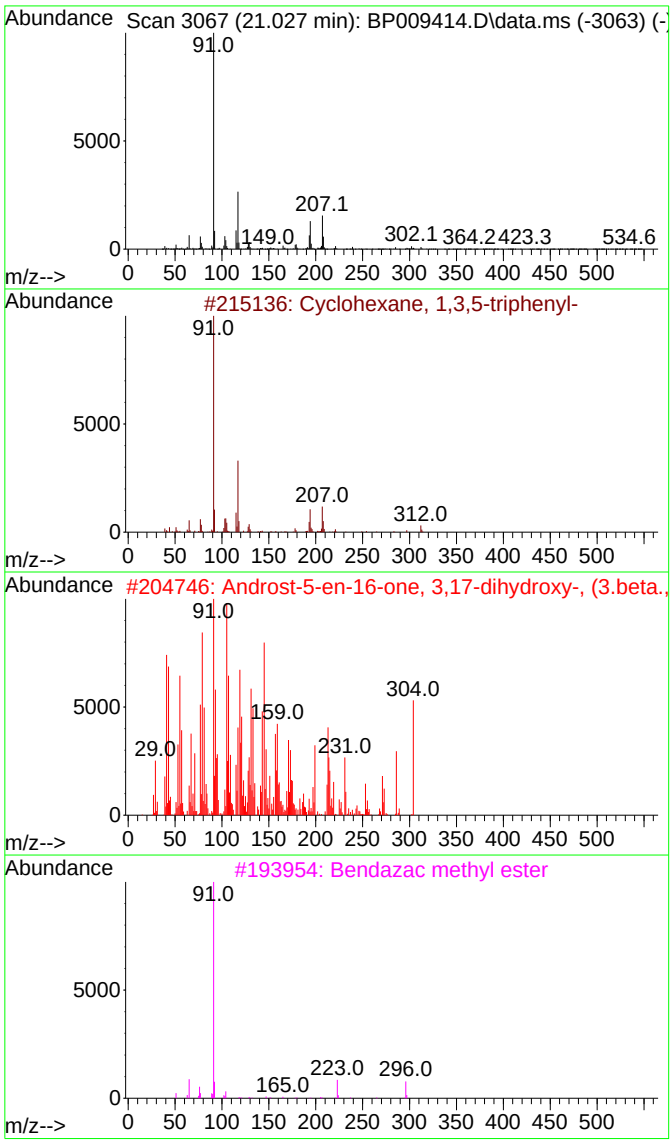
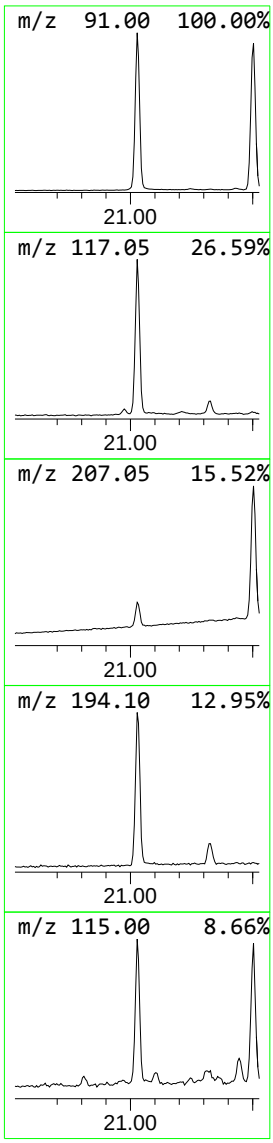
Peak Number 9 Cyclohexane, 1,3,5-triphenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.027	9.75 ng	3850670	Chrysene-d12	21.327

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,3,5-triphenyl-	312	C24H24	028336-57-4	53
2	Androst-5-en-16-one, 3,17-dihydr...	304	C19H28O3	001159-66-6	22
3	Bendazac methyl ester	296	C17H16N2O3	1010119-77-6	14
4	Salicylic acid, trimethylsilyl e...	300	C17H20O3Si	1000463-82-7	14
5	Nialamide, 1TMS derivative	370	C19H26N4O2Si	1000485-79-3	12





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Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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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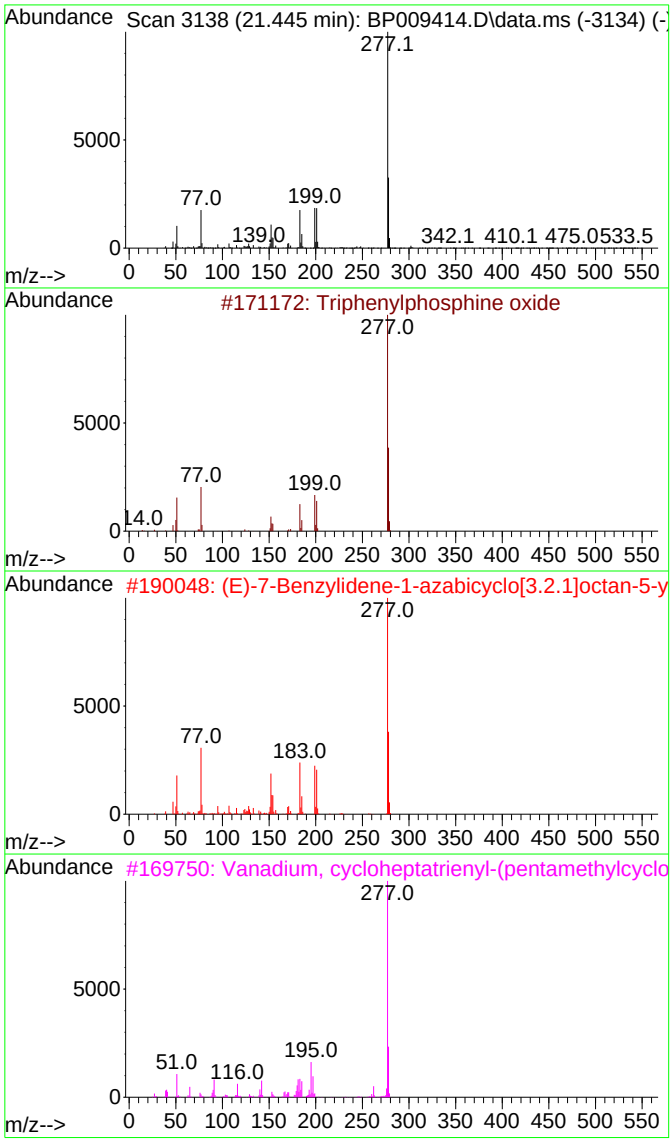
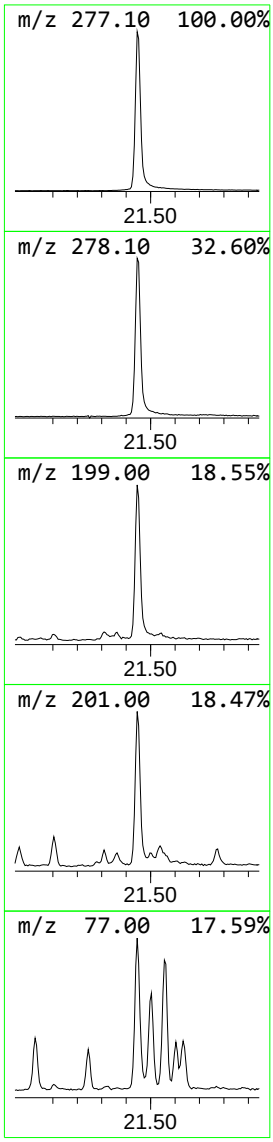
Peak Number 10 Triphenylphosphine oxide Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.445	9.21 ng	3639500	Chrysene-d12	21.327

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	94
2	(E)-7-Benzylidene-1-azabicyclo[3...	293	C15H19NO3S	1000483-83-4	52
3	Vanadium, cycloheptatrienyl-(pen...	277	C17H22V	1000155-20-5	40
4	(Carbethoxymethyl)-triphenylphos...	428	C22H22BrO2P	001530-45-6	40
5	Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	38





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Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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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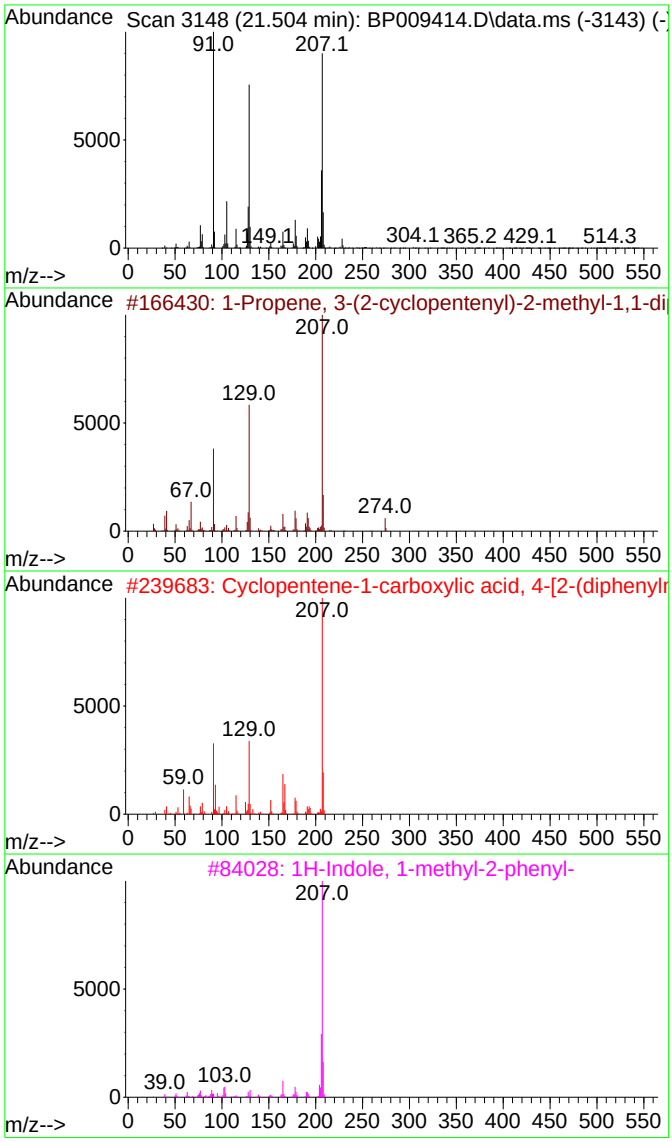
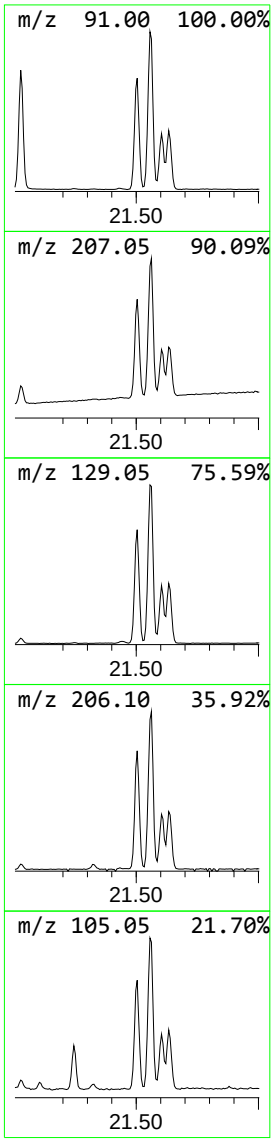
Peak Number 11 1-Propene, 3-(2-cyclopenten... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.504	15.55 ng	6144820	Chrysene-d12	21.327

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propene, 3-(2-cyclopentenyl)-2...	274	C21H22	1010154-23-3	59
2	Cyclopentene-1-carboxylic acid, ...	332	C23H24O2	1000159-40-6	45
3	1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	42
4	2-(4-Methylphenyl)-1H-indole	207	C15H13N	055577-25-8	41
5	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	38





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Data File : BP009414.D
Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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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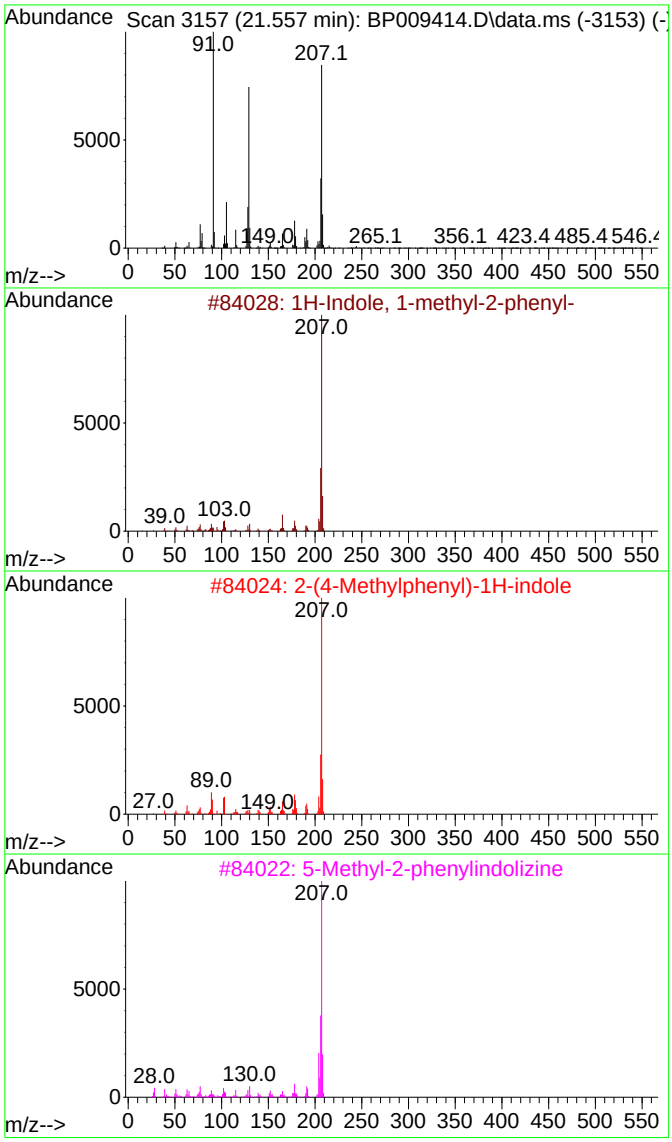
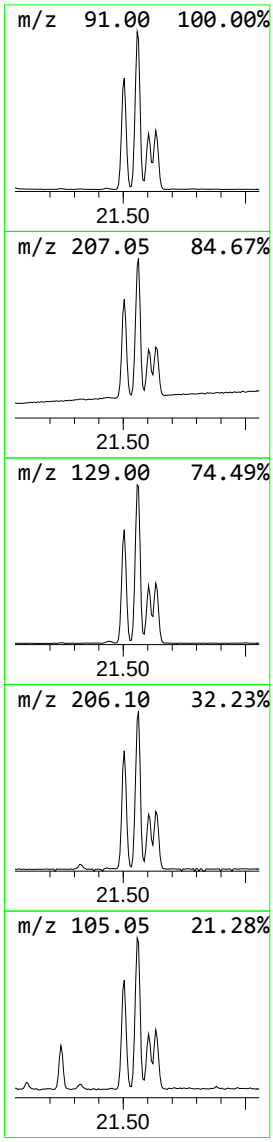
Peak Number 12 unknown21.557 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.557	20.28 ng	8011410	Chrysene-d12	21.327

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	46
2	2-(4-Methylphenyl)-1H-indole	207	C15H13N	055577-25-8	46
3	5-Methyl-2-phenylindolizine	207	C15H13N	036944-99-7	45
4	Cyclopentene-1-carboxylic acid, ...	332	C23H24O2	1000159-40-6	45
5	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	43





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Data File : BP009414.D
Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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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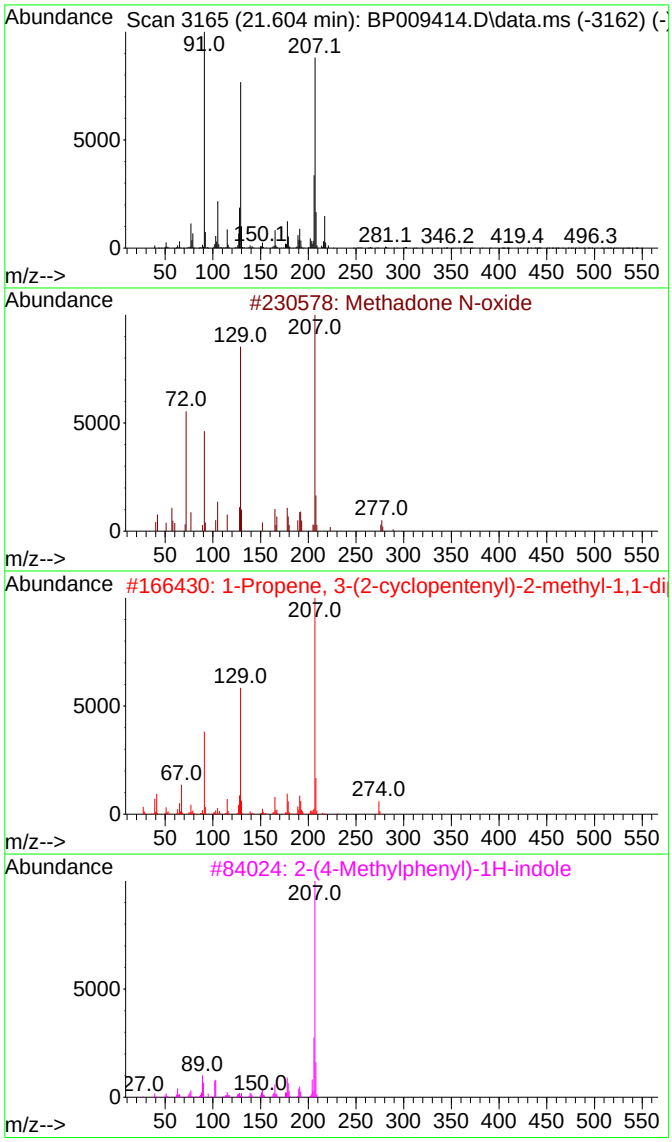
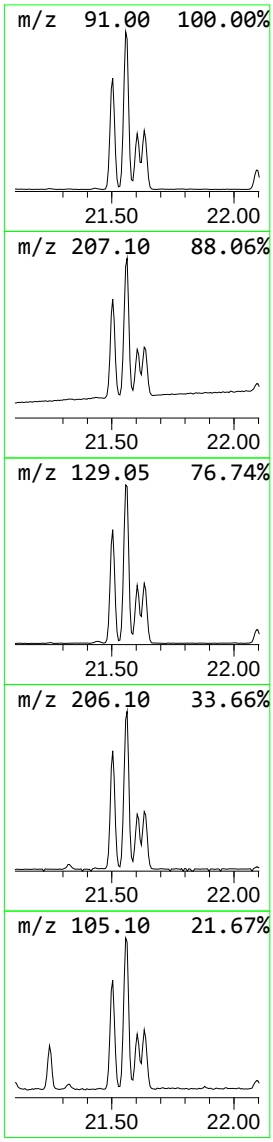
Peak Number 13 Methadone N-oxide Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.604	7.00 ng	2765550	Chrysene-d12	21.327

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methadone N-oxide	325	C21H27NO2	1000120-80-7	64
2	1-Propene, 3-(2-cyclopentenyl)-2...	274	C21H22	1010154-23-3	59
3	2-(4-Methylphenyl)-1H-indole	207	C15H13N	055577-25-8	50
4	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	45
5	1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	41





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Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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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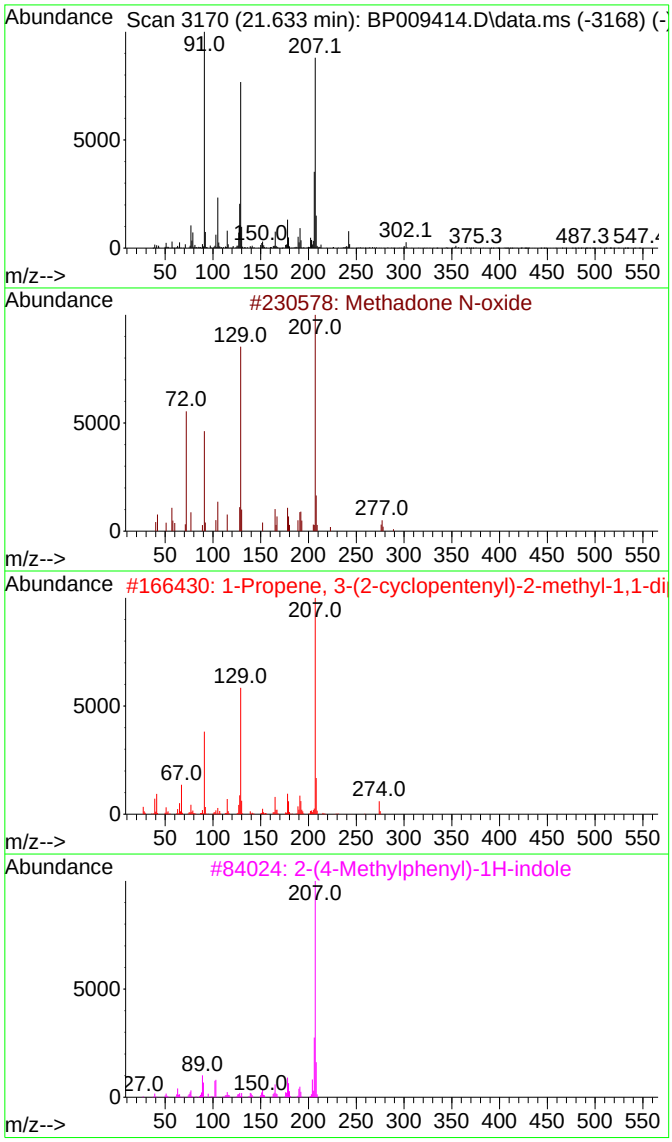
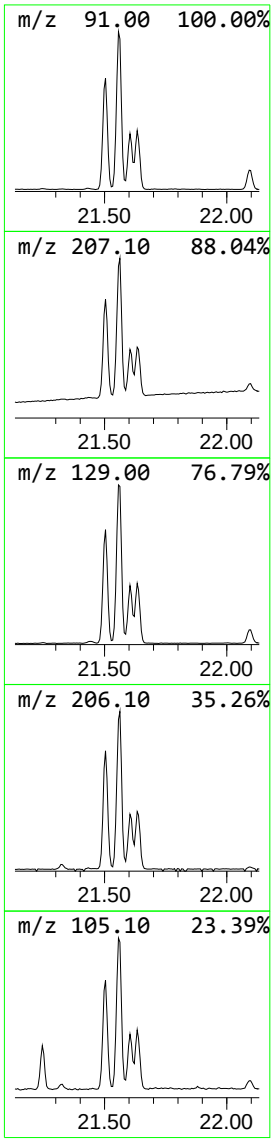
Peak Number 14 1-Propene, 3-(2-cyclopenten... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.633	7.52 ng	2970490	Chrysene-d12	21.327

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methadone N-oxide	325	C21H27NO2	1000120-80-7	72
2	1-Propene, 3-(2-cyclopentenyl)-2...	274	C21H22	1010154-23-3	50
3	2-(4-Methylphenyl)-1H-indole	207	C15H13N	055577-25-8	46
4	7-Methyl -2-phenyl-1H-indole	207	C15H13N	059541-82-1	43
5	5-Methyl-2-phenylindolizine	207	C15H13N	036944-99-7	43





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031522\
Data File : BP009414.D
Acq On : 16 Mar 2022 00:08
Operator : CG/JU
Sample : N1894-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP3(0-10)

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	2.2	ng	213821	1	7.810	1983250	20.0
1,3,5,7-Cyclooc...	5.816	6.1	ng	605660	1	7.810	1983250	20.0
Benzene, (1-met...	6.316	2.5	ng	252487	1	7.810	1983250	20.0
Benzene, 1,1'-(...	16.398	2.1	ng	464592	4	17.216	4520870	20.0
[2.2]Paracyclop...	16.822	8.0	ng	1805760	4	17.216	4520870	20.0
n-Hexadecanoic ...	18.121	6.3	ng	1423330	4	17.216	4520870	20.0
Octadecanoic acid	19.363	4.7	ng	1840260	5	21.327	7902040	20.0
Cyclohexane, 1,...	21.027	9.8	ng	3850670	5	21.327	7902040	20.0
Triphenylphosph...	21.445	9.2	ng	3639500	5	21.327	7902040	20.0
1-Propene, 3-(2...	21.504	15.6	ng	6144820	5	21.327	7902040	20.0
unknown21.557	21.557	20.3	ng	8011410	5	21.327	7902040	20.0
Methadone N-oxide	21.604	7.0	ng	2765550	5	21.327	7902040	20.0
1-Propene, 3-(2...	21.633	7.5	ng	2970490	5	21.327	7902040	20.0