LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022324\
Data File : BP019852.D
Acq On : 23 Feb 2024 20:03
Operator : MA/JU
Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
2

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-BP013124.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP019852.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.128	3	7	19	rVB	686394	835923	18.70%	2.265%
2	4.075	163	168	178	rVB3	43560	77986	1.74%	0.211%
3	4.646	260	265	270	rBV	84368	123888	2.77%	0.336%
4	4.699	270	274	284	rVB	48078	75658	1.69%	0.205%
5	5.022	320	329	340	rBV	247003	370708	8.29%	1.004%
6	5.481	399	407	416	rBV	1070520	1639754	36.69%	4.443%
7	7.081	671	679	693	rBV	984575	1680384	37.60%	4.553%
8	7.905	813	819	826	rVB	299334	465083	10.41%	1.260%
9	9.075	1010	1018	1028	rBV	642542	1126944	25.21%	3.053%
10	10.704	1287	1295	1303	rBV	364153	640087	14.32%	1.734%
11	13.169	1707	1714	1729	rBV	1803194	2761742	61.79%	7.483%
12	14.269	1895	1901	1908	rBV	78108	131211	2.94%	0.356%
13	14.545	1941	1948	1955	rBV	589869	904264	20.23%	2.450%
14	14.604	1955	1958	1967	rVB2	28960	49784	1.11%	0.135%
15	15.598	2122	2127	2140	rVB2	45429	91792	2.05%	0.249%
16	16.045	2196	2203	2212	rBV	1711560	2541047	56.85%	6.885%
17	16.281	2238	2243	2251	rVB10	22897	52179	1.17%	0.141%
18	17.128	2383	2387	2394	rVB	37044	67936	1.52%	0.184%
19	17.310	2412	2418	2422	rBV	776069	1173241	26.25%	3.179%
20	17.351	2422	2425	2432	rVV	633502	879279	19.67%	2.382%
21	17.445	2436	2441	2446	rVB	192267	286034	6.40%	0.775%
22	18.180	2561	2566	2567	rBV	124329	152784	3.42%	0.414%
23	18.210	2567	2571	2580	rVV2	349209	709077	15.86%	1.921%
24	18.304	2583	2587	2592	rVV	74938	111907	2.50%	0.303%
25	18.369	2592	2598	2601	rVV3	205659	363238	8.13%	0.984%
26	18.404	2601	2604	2609	rVB	94735	136053	3.04%	0.369%
27	18.669	2645	2649	2652	rBV	87301	112403	2.51%	0.305%
28	18.710	2652	2656	2664	rVB	60192	96079	2.15%	0.260%
29	19.069	2713	2717	2721	rBV2	92834	147197	3.29%	0.399%
30	19.210	2737	2741	2747	rBV4	60836	107606	2.41%	0.292%
31	19.327	2755	2761	2767	rBV	1771077	2361212	52.83%	6.398%
32	19.445	2778	2781	2783	rBV2	74609	85419	1.91%	0.231%
33	19.469	2783	2785	2790	rVB	87141	99730	2.23%	0.270%
34	19.651	2811	2816	2817	rBV	256329	349771	7.83%	0.948%
35	19.680	2817	2821	2829	rVV	1483726	2022427	45.25%	5.480%
36	19.751	2830	2833	2839	rVB2	85926	133183	2.98%	0.361%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022324\
Data File : BP019852.D
Acq On : 23 Feb 2024 20:03
Operator : MA/JU
Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
2

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-BP013124.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	19.880	2848	2855	2861	rBV	3481126	4469698	100.00%	12.110%
38	19.998	2870	2875	2883	rBV2	113293	289226	6.47%	0.784%
39	20.163	2900	2903	2910	rVB	256260	378524	8.47%	1.026%
40	20.263	2917	2920	2924	rBV	173530	203773	4.56%	0.552%
41	20.316	2927	2929	2934	rVB	138329	180719	4.04%	0.490%
42	21.133	3065	3068	3074	rVB3	344110	456766	10.22%	1.238%
43	21.410	3109	3115	3120	rBV2	1312902	2752765	61.59%	7.459%
44	21.457	3120	3123	3132	rVB	787139	1076814	24.09%	2.918%
45	23.116	3400	3405	3411	rBV	715836	1744068	39.02%	4.725%
46	23.751	3508	3513	3520	rVB	635117	1311110	29.33%	3.552%
47	23.857	3527	3531	3537	rVB	588150	1081268	24.19%	2.930%

Sum of corrected areas: 36907741

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022324\
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ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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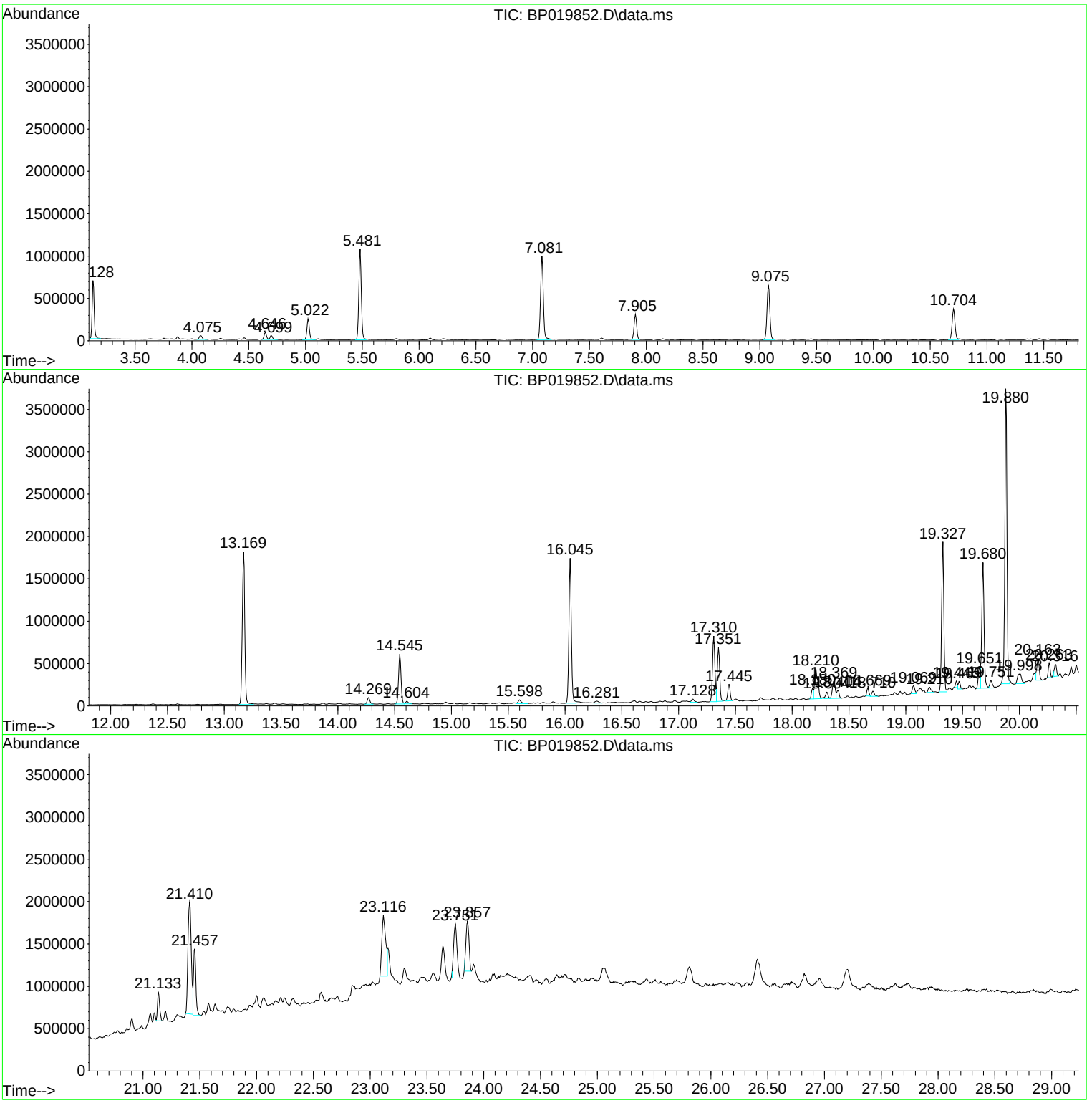
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022324\
Data File : BP019852.D
Acq On : 23 Feb 2024 20:03
Operator : MA/JU
Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

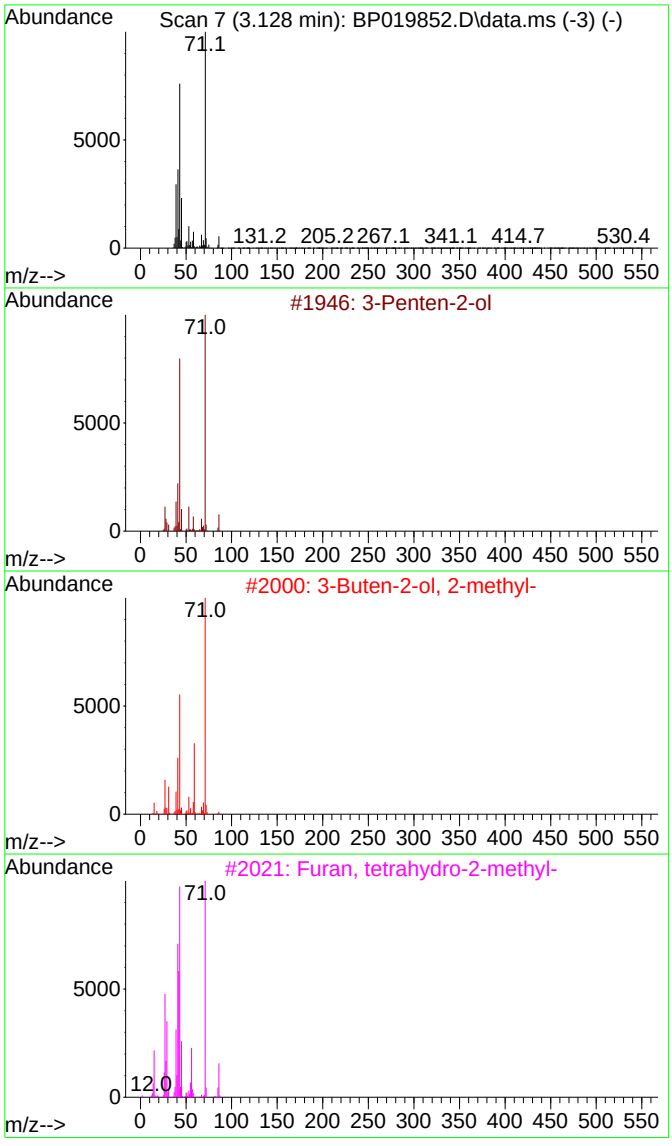
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

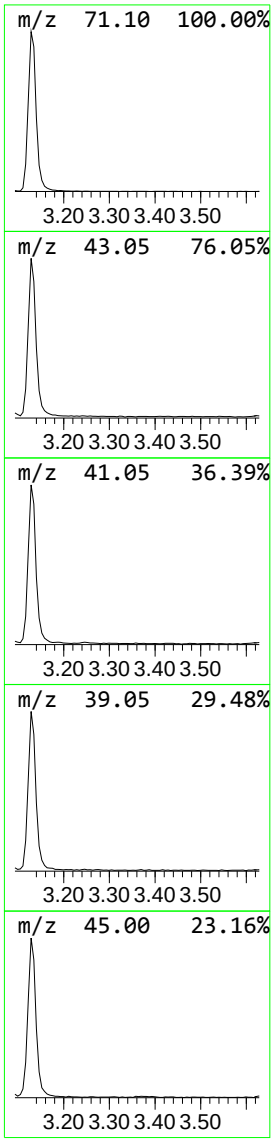
****

Peak Number 1 3-Penten-2-ol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.128	35.95 ng	835923	1,4-Dichlorobenzene-d4	7.905

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Penten-2-ol	86	C5H10O	001569-50-2	80
2	3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	64
3	Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	50
4	Butanoic acid, anhydride	158	C8H14O3	000106-31-0	45
5	Furan, tetrahydro-2-(methoxymeth...	116	C6H12O2	019354-27-9	45





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022324\
Data File : BP019852.D
Acq On : 23 Feb 2024 20:03
Operator : MA/JU
Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

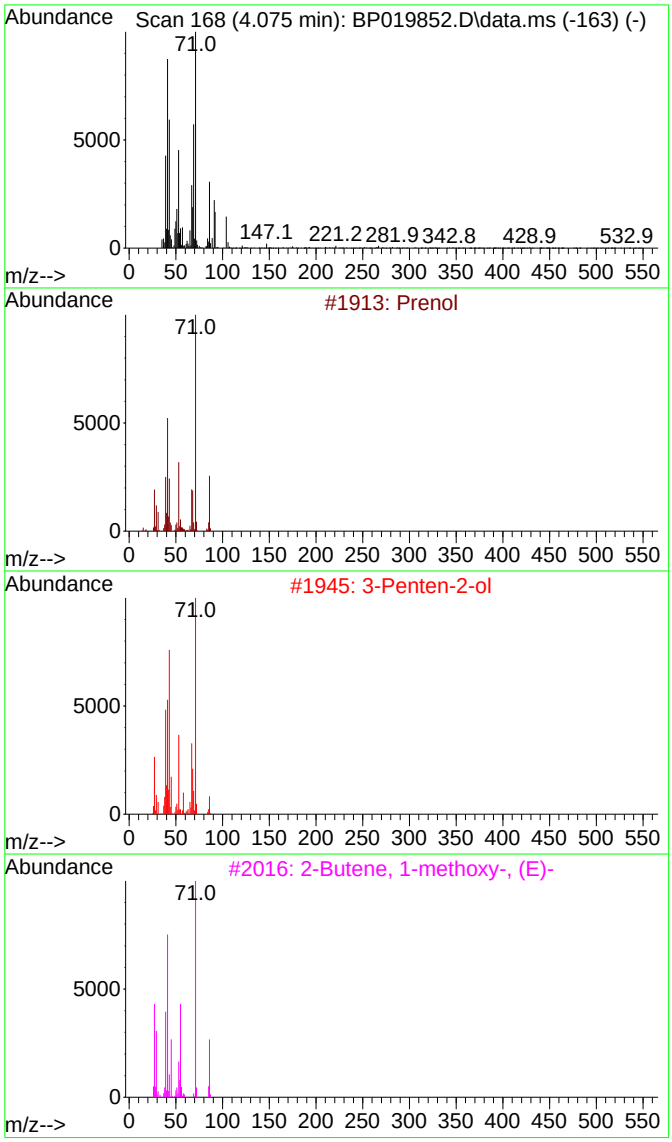
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

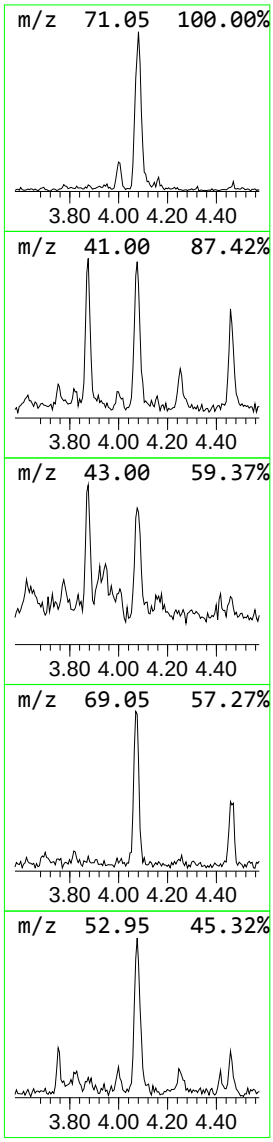
****

Peak Number 2 Prenol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.075	3.35 ng	77986	1,4-Dichlorobenzene-d4	7.905

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Prenol	86	C5H10O	000556-82-1	64
2	3-Penten-2-ol	86	C5H10O	001569-50-2	35
3	2-Butene, 1-methoxy-, (E)-	86	C5H10O	010034-14-7	27
4	2-Butyne, 1-methoxy-	84	C5H8O	002768-41-4	25
5	1-Butene, 1-chloro-3-methyl-	104	C5H9Cl	023010-00-6	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022324\
Data File : BP019852.D
Acq On : 23 Feb 2024 20:03
Operator : MA/JU
Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

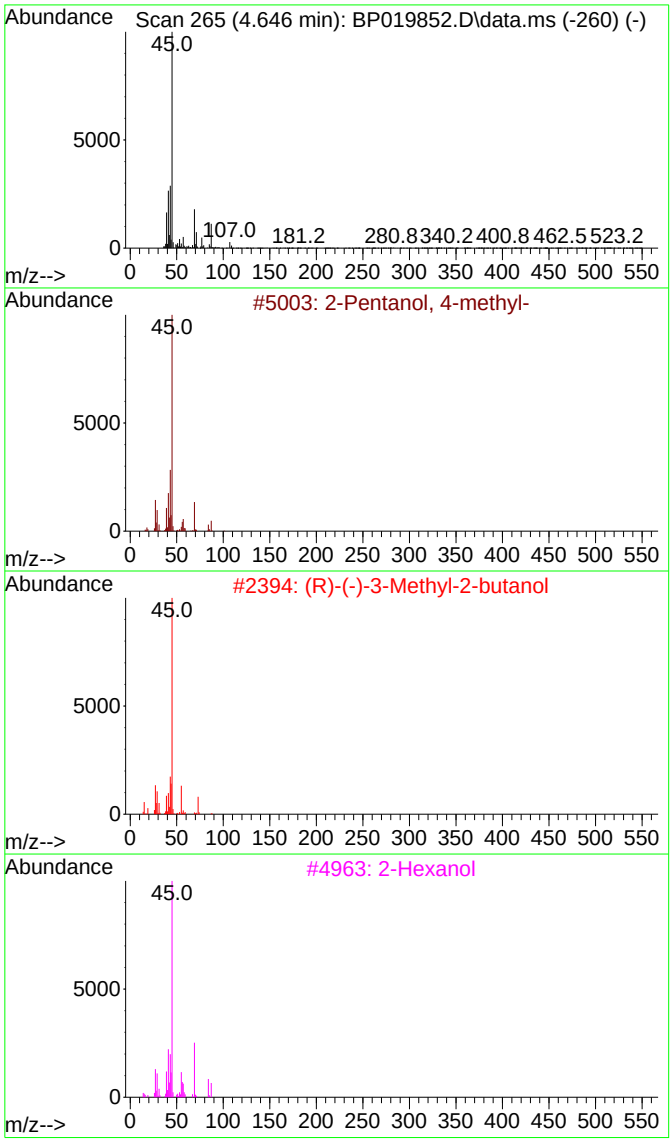
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

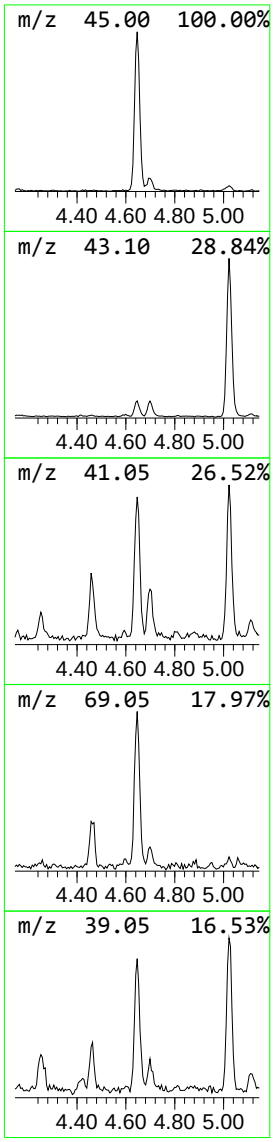
****

Peak Number 3 2-Pentanol, 4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.646	5.33 ng	123888	1,4-Dichlorobenzene-d4	7.905

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	64
2	(R)-(-)-3-Methyl-2-butanol	88	C5H12O	001572-93-6	40
3	2-Hexanol	102	C6H14O	000626-93-7	39
4	4-Penten-2-ol	86	C5H10O	000625-31-0	36
5	Formic acid, 1-methylethyl ester	88	C4H8O2	000625-55-8	33





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022324\
Data File : BP019852.D
Acq On : 23 Feb 2024 20:03
Operator : MA/JU
Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
2

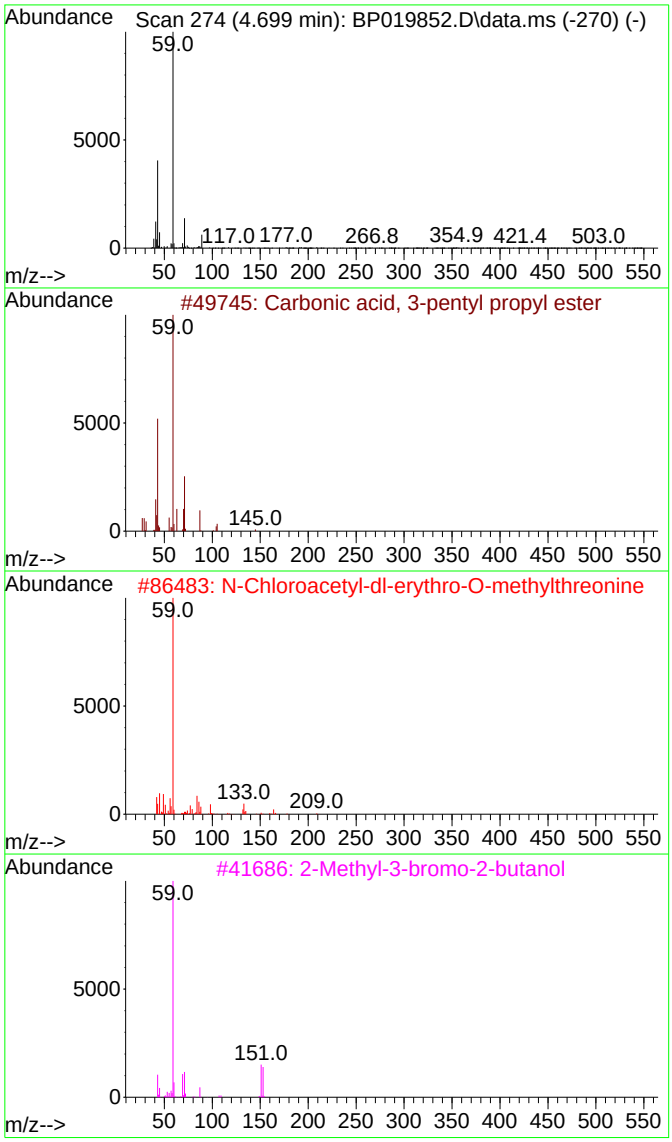
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

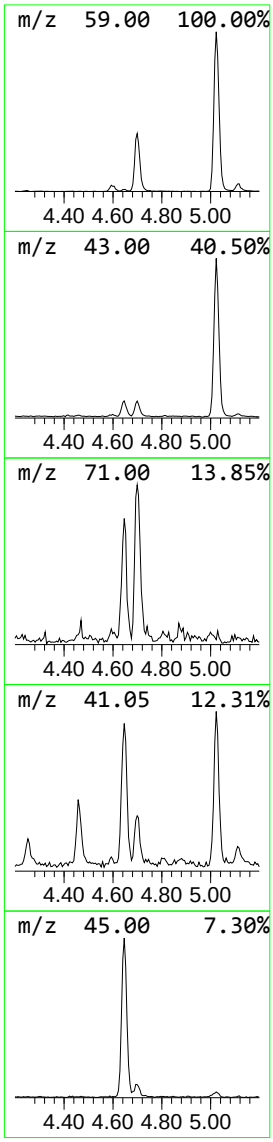
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Carbonic acid, 3-pentyl pro... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.699	3.25 ng	75658	1,4-Dichlorobenzene-d4	7.905

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, 3-pentyl propyl e...	174	C9H18O3	1000325-64-0	56
2	N-Chloroacetyl-dl-erythro-O-meth...	209	C7H12ClNO4	1010214-48-1	9
3	2-Methyl-3-bromo-2-butanol	166	C5H11BrO	002588-77-4	9
4	Hydrazine, 1-methyl-1-(2-methylp...	102	C5H14N2	020240-63-5	9
5	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022324\
Data File : BP019852.D
Acq On : 23 Feb 2024 20:03
Operator : MA/JU
Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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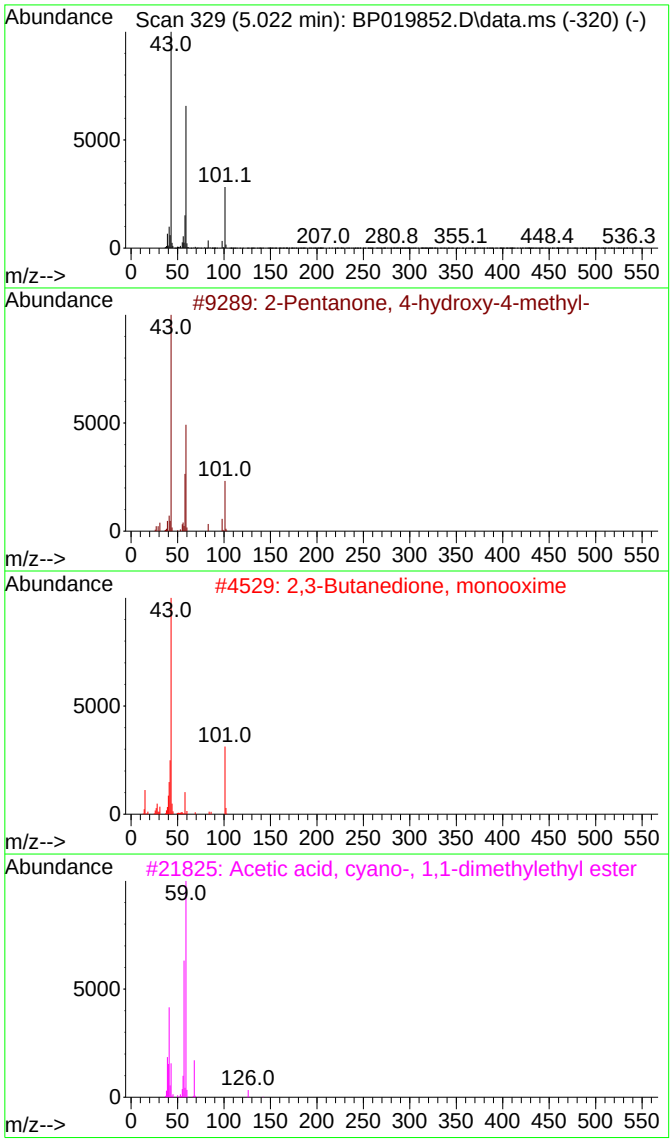
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

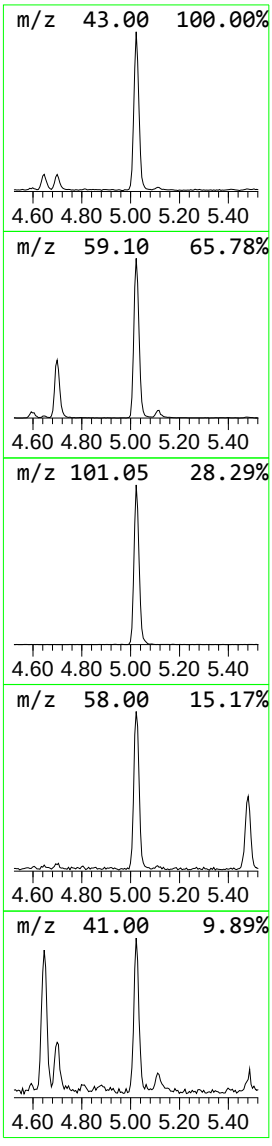
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.022	15.94 ng	370708	1,4-Dichlorobenzene-d4	7.905

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022324\
Data File : BP019852.D
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Operator : MA/JU
Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

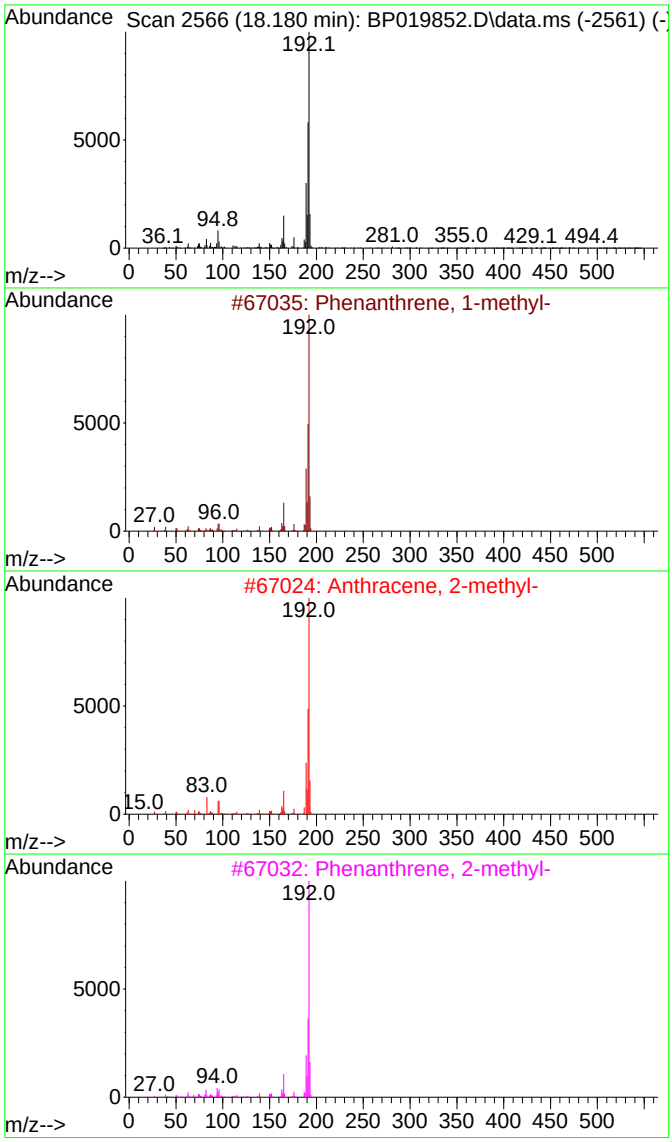
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

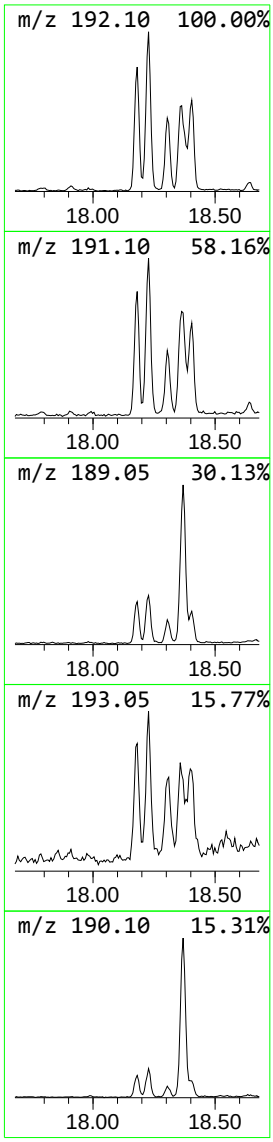
****

Peak Number 7 Phenanthrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.180	2.60 ng	152784	Phenanthrene-d10	17.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	93
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022324\
Data File : BP019852.D
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Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

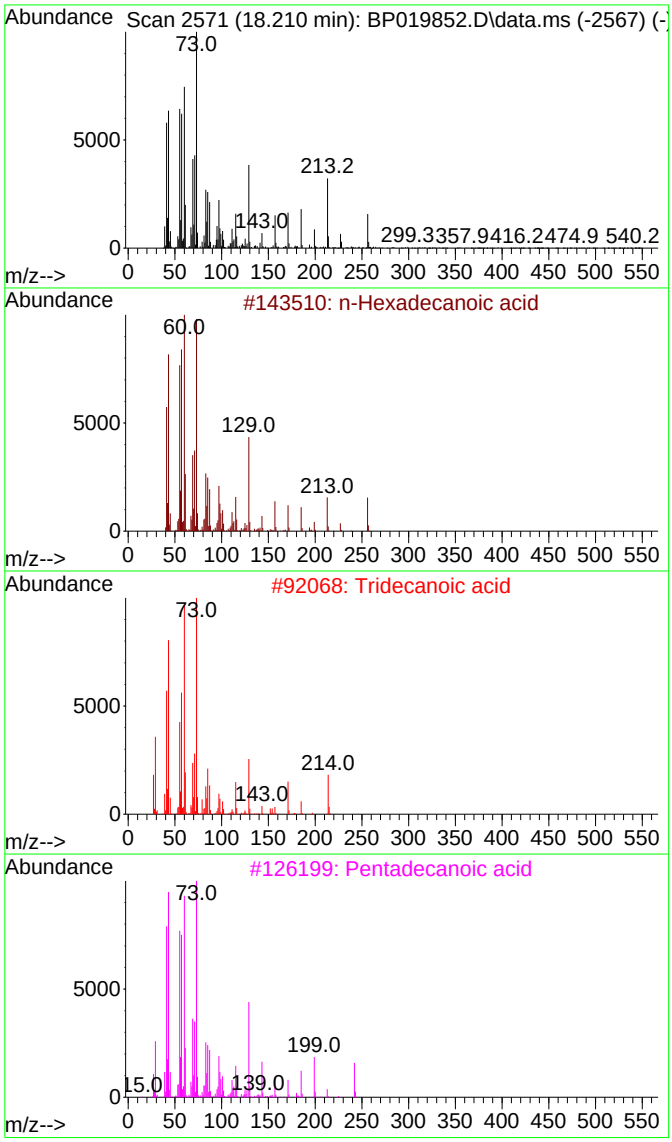
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

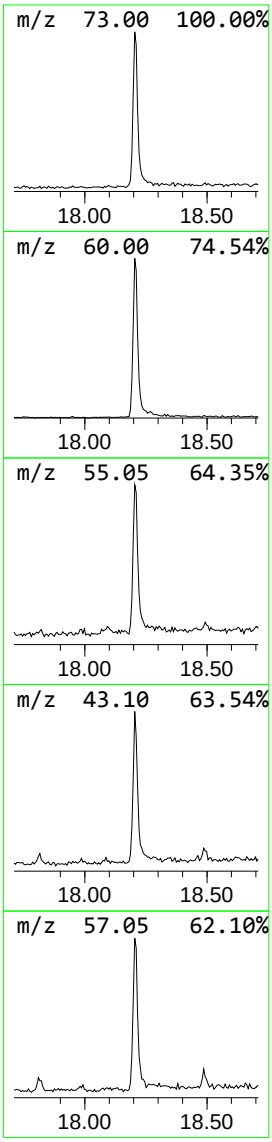
****

Peak Number 8 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.210	12.09 ng	709077	Phenanthrene-d10	17.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tridecanoic acid	214	C13H26O2	000638-53-9	90
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	81
4	n-Decanoic acid	172	C10H20O2	000334-48-5	76
5	Octadecanoic acid	284	C18H36O2	000057-11-4	74





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022324\
Data File : BP019852.D
Acq On : 23 Feb 2024 20:03
Operator : MA/JU
Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
2

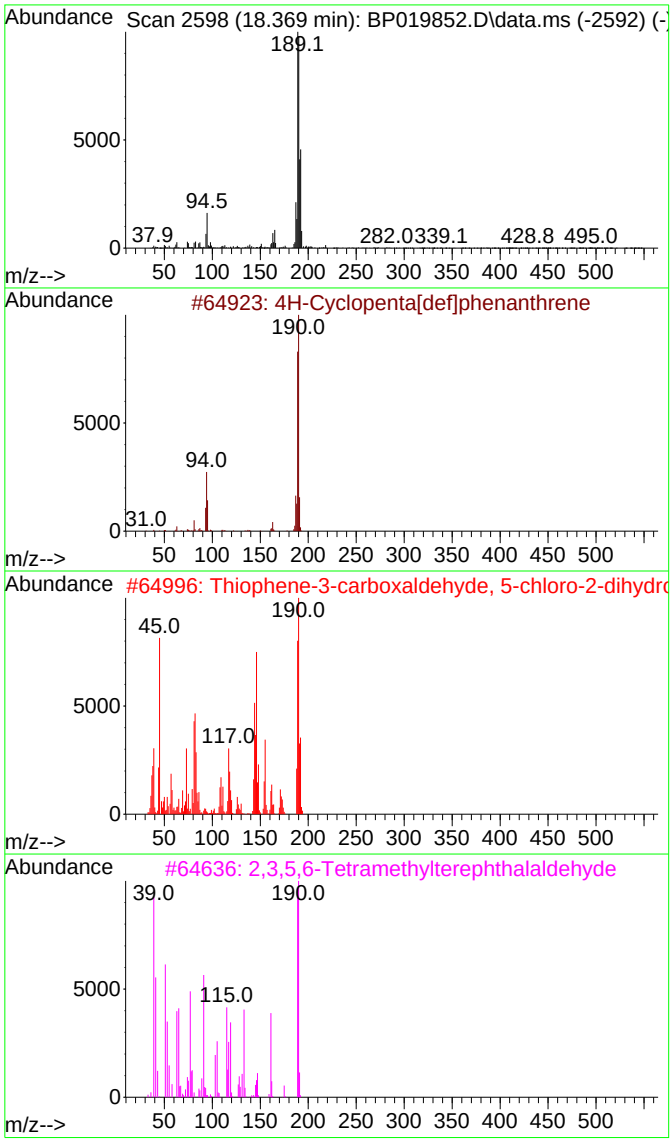
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

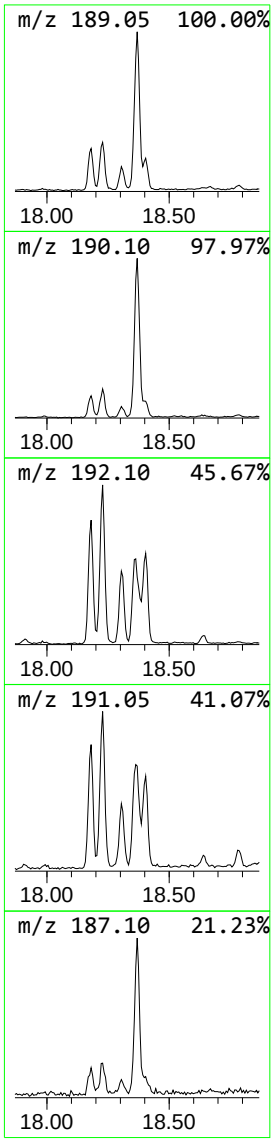
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.369	6.19 ng	363238	Phenanthrene-d10	17.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
2	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	53
3	2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	43
4	Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	42
5	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	38





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Acq On : 23 Feb 2024 20:03
Operator : MA/JU
Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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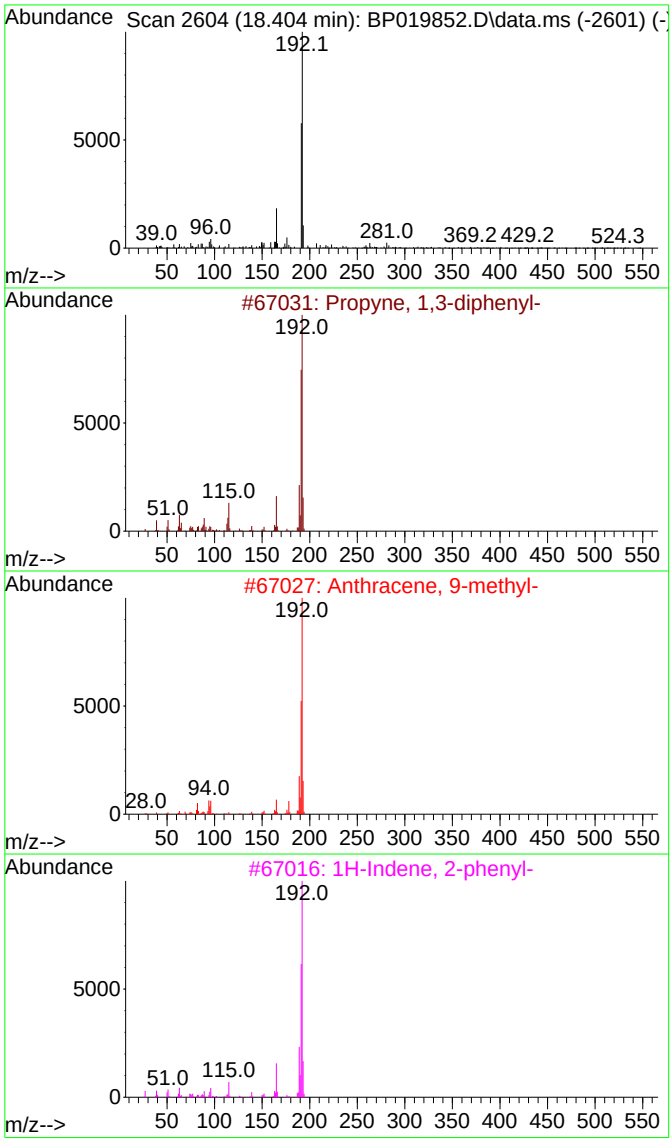
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

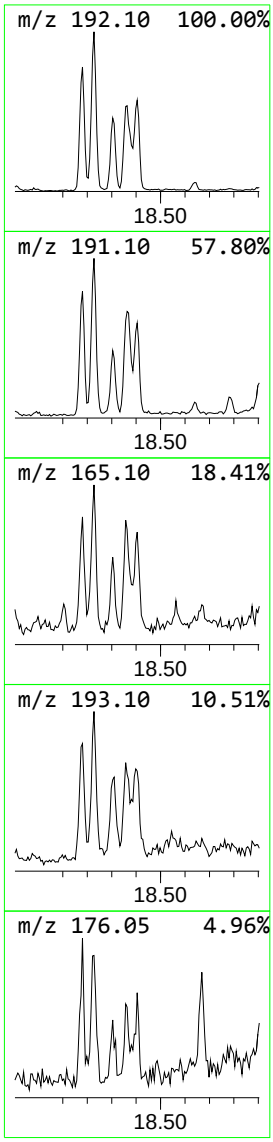
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Propyne, 1,3-diphenyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.404	2.32 ng	136053	Phenanthrene-d10	17.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propyne, 1,3-diphenyl-	192	C15H12	004980-70-5	86
2	Anthracene, 9-methyl-	192	C15H12	000779-02-2	86
3	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	74
4	1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	72
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	72





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Acq On : 23 Feb 2024 20:03
Operator : MA/JU
Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

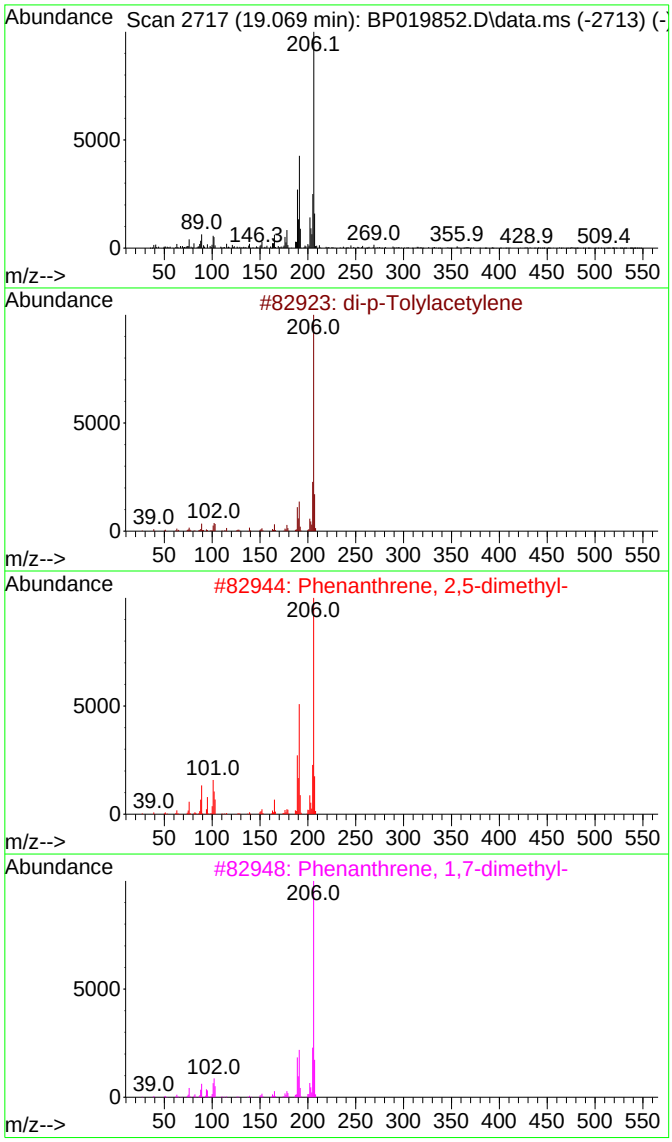
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

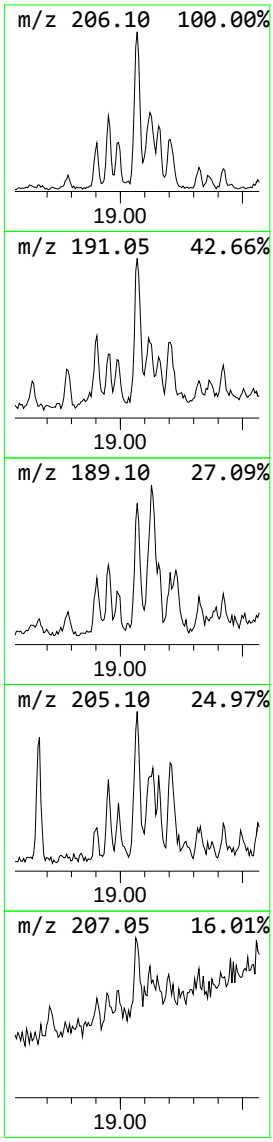
****

Peak Number 11 di-p-Tolylacetylene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.069	2.51 ng	147197	Phenanthrene-d10	17.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	di-p-Tolylacetylene	206	C16H14	002789-88-0	94
2	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
3	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	93
4	9,10-Dimethylanthracene	206	C16H14	000781-43-1	91
5	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	90





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Acq On : 23 Feb 2024 20:03
Operator : MA/JU
Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.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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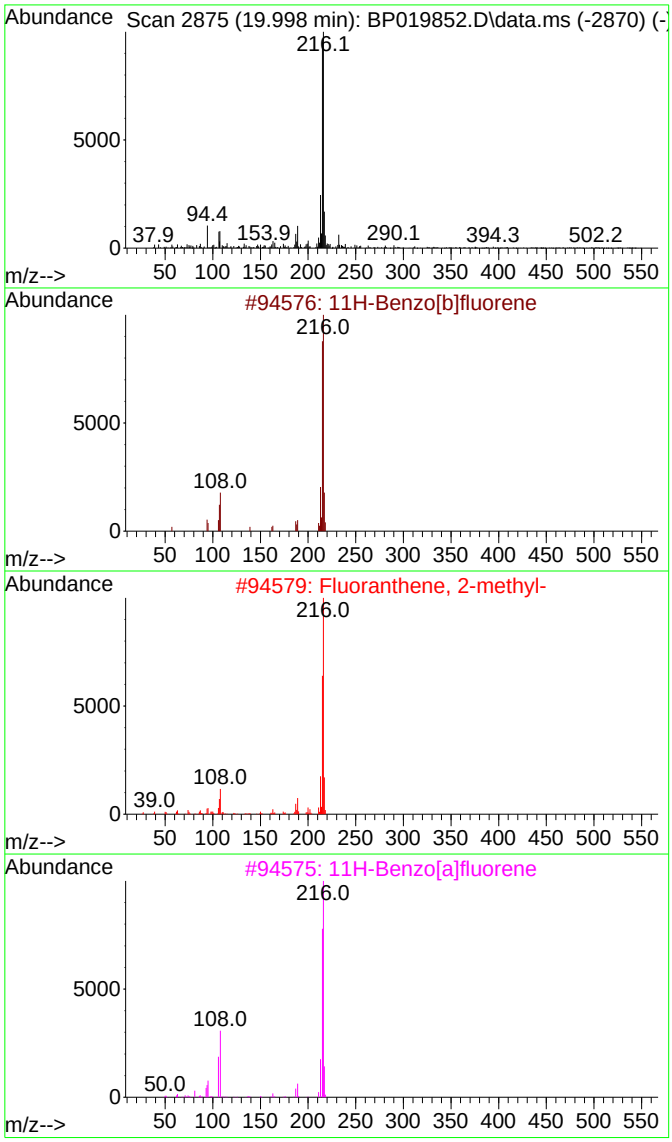
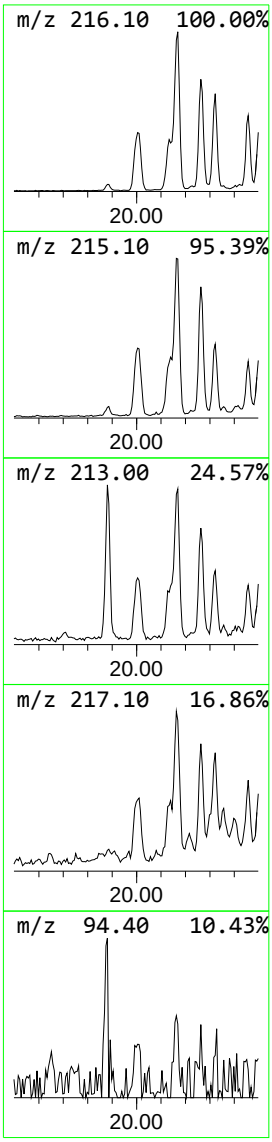
Peak Number 12 11H-Benzo[b]fluorene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.998	2.10 ng	289226	Chrysene-d12	21.416

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	96
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
4	Pyrene, 1-methyl-	216	C17H12	002381-21-7	81
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74





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Acq On : 23 Feb 2024 20:03
Operator : MA/JU
Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.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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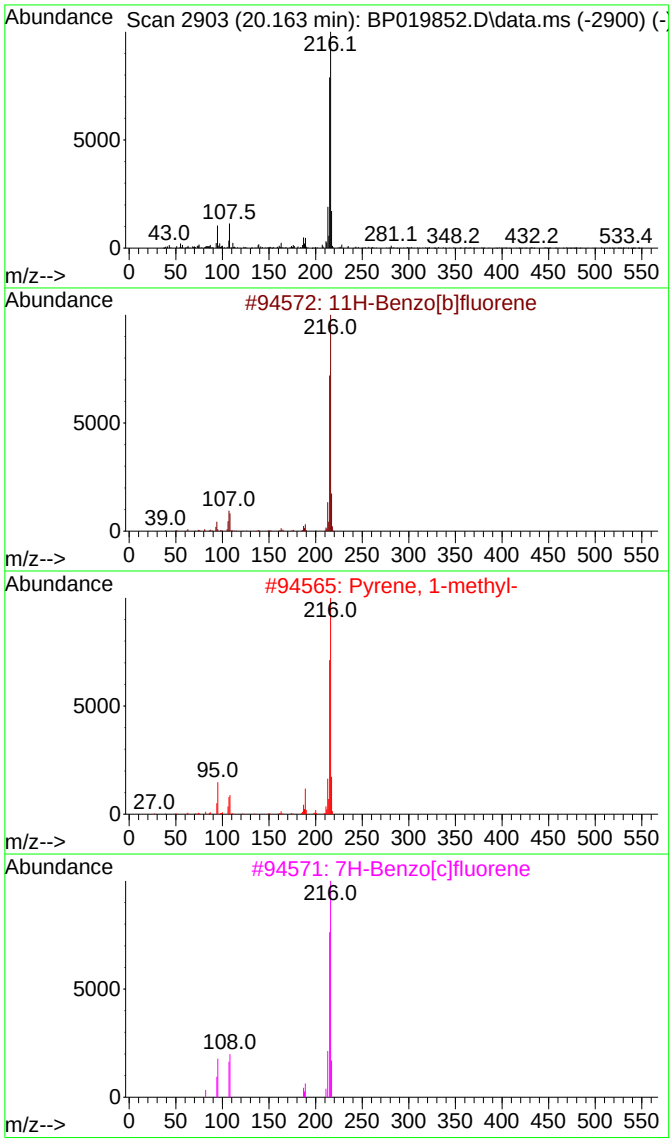
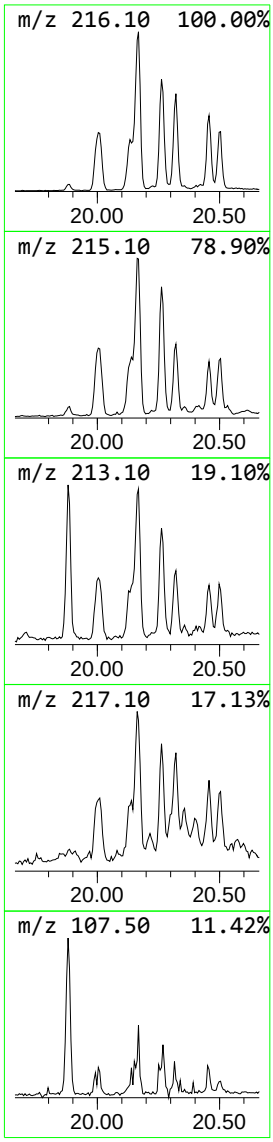
Peak Number 13 Pyrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.163	2.75 ng	378524	Chrysene-d12	21.416

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
3	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	87
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	83





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Acq On : 23 Feb 2024 20:03
Operator : MA/JU
Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.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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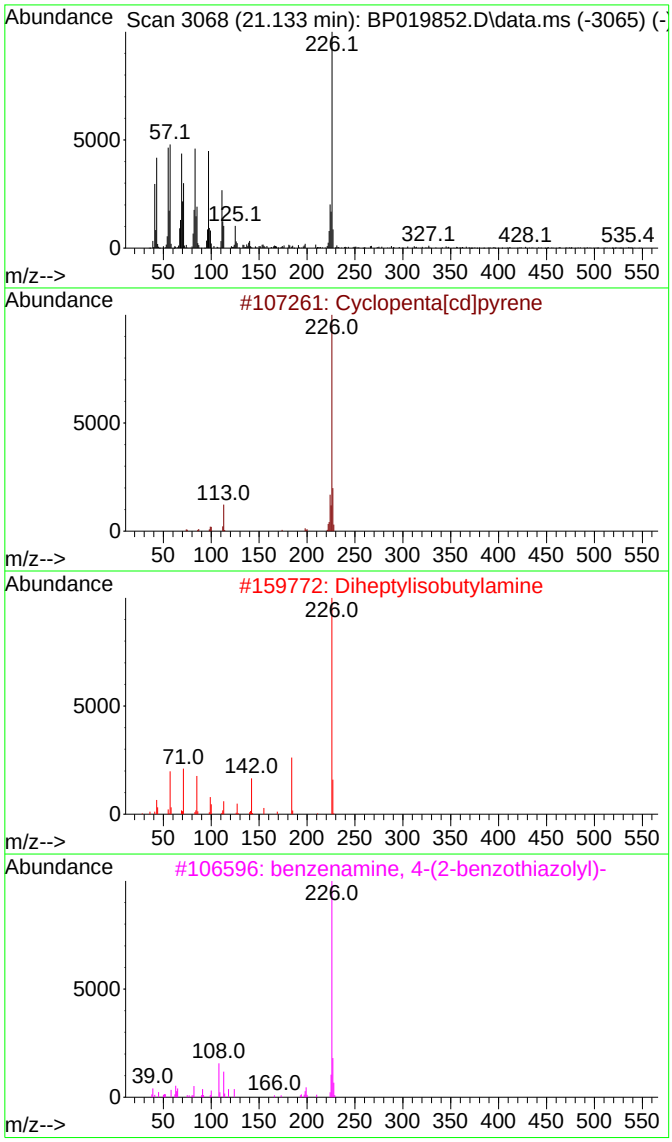
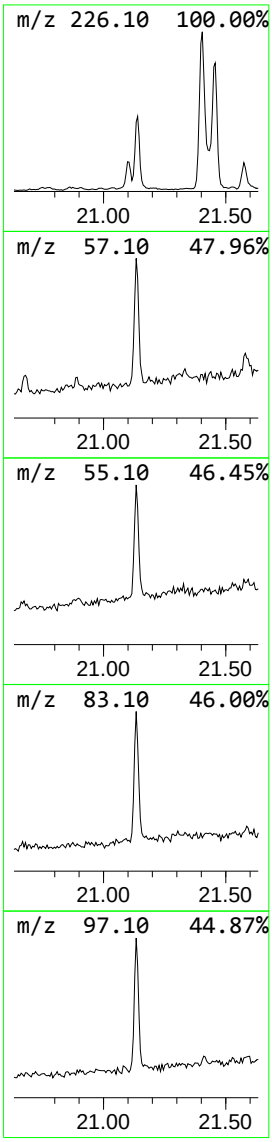
Peak Number 14 unknown21.133 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.133	3.32 ng	456766	Chrysene-d12	21.416

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	22
2	Diheptylisobutylamine	269	C18H39N	1000310-32-0	12
3	benzenamine, 4-(2-benzothiazolyl)-	226	C13H10N2S	1000400-93-4	10
4	9H-Carbazole, 9-methyl-3-nitro-	226	C13H10N2O2	061166-05-0	10
5	L-Proline, N-(3-chloro-2-fluorob...	425	C23H33ClFNO3	1000345-94-3	10





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022324\
Data File : BP019852.D
Acq On : 23 Feb 2024 20:03
Operator : MA/JU
Sample : P1537-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.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
3-Penten-2-ol	3.128	36.0	ng	835923	1	7.905	465083	20.0
Prenol	4.075	3.4	ng	77986	1	7.905	465083	20.0
2-Pentanol, 4-m...	4.646	5.3	ng	123888	1	7.905	465083	20.0
Carbonic acid, ...	4.699	3.3	ng	75658	1	7.905	465083	20.0
2-Pentanone, 4-...	5.022	15.9	ng	370708	1	7.905	465083	20.0
Phenanthrene, 1...	18.180	2.6	ng	152784	4	17.310	1173240	20.0
n-Hexadecanoic ...	18.210	12.1	ng	709077	4	17.310	1173240	20.0
4H-Cyclopenta[d...	18.369	6.2	ng	363238	4	17.310	1173240	20.0
Propyne, 1,3-di...	18.404	2.3	ng	136053	4	17.310	1173240	20.0
di-p-Tolylacety...	19.069	2.5	ng	147197	4	17.310	1173240	20.0
11H-Benzo[b]flu...	19.998	2.1	ng	289226	5	21.416	2752770	20.0
Pyrene, 1-methyl-	20.163	2.8	ng	378524	5	21.416	2752770	20.0
unknown21.133	21.133	3.3	ng	456766	5	21.416	2752770	20.0