LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082625\
Data File : BP025587.D
Acq On : 26 Aug 2025 15:55
Operator : CG/JU
Sample : Q2940-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

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-BP081425.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP025587.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.031	11	16	31	rVB	3731965	5599473	46.81%	7.532%
2	4.943	334	341	352	rBV2	227903	340518	2.85%	0.458%
3	5.408	411	420	429	rBV	1676036	2551046	21.33%	3.431%
4	6.972	676	686	698	rBV	943768	1573928	13.16%	2.117%
5	7.331	741	747	751	rBV	82576	120236	1.01%	0.162%
6	7.566	782	787	795	rVB	75215	123595	1.03%	0.166%
7	7.778	815	823	831	rBV	770365	1219799	10.20%	1.641%
8	8.919	1009	1017	1026	rBV	2318661	3911722	32.70%	5.262%
9	10.549	1287	1294	1307	rBV	929961	1728021	14.45%	2.324%
10	10.831	1335	1342	1350	rVB	142284	223111	1.87%	0.300%
11	13.007	1704	1712	1726	rBV	5493884	8756766	73.21%	11.779%
12	14.395	1941	1948	1955	rBV	1480488	2212468	18.50%	2.976%
13	14.619	1974	1986	1991	rBV3	160115	519340	4.34%	0.699%
14	15.772	2177	2182	2190	rBV	699723	1038710	8.68%	1.397%
15	15.907	2198	2205	2218	rBV	4964168	7422365	62.06%	9.984%
16	16.683	2333	2337	2342	rBV	83815	131456	1.10%	0.177%
17	17.195	2418	2424	2431	rBV	1810740	2636213	22.04%	3.546%
18	18.136	2579	2584	2592	rBV	1080283	1714441	14.33%	2.306%
19	19.236	2762	2771	2779	rBV	4259915	8289660	69.31%	11.150%
20	19.301	2779	2782	2798	rVB	355726	853777	7.14%	1.148%
21	19.454	2804	2808	2818	rBV	369880	589357	4.93%	0.793%
22	19.683	2843	2847	2853	rVB4	132692	236140	1.97%	0.318%
23	19.919	2881	2887	2894	rBV	9006667	11960923	100.00%	16.089%
24	19.989	2896	2899	2901	rBV4	96765	134014	1.12%	0.180%
25	20.124	2918	2922	2928	rVB	461251	769464	6.43%	1.035%
26	20.501	2984	2986	2990	rVB3	338504	318450	2.66%	0.428%
27	20.677	3013	3016	3020	rVB	261404	300272	2.51%	0.404%
28	20.771	3030	3032	3035	rVB2	236952	243938	2.04%	0.328%
29	21.213	3102	3107	3113	rVB	721843	1398105	11.69%	1.881%
30	21.295	3118	3121	3131	rVB	683961	961516	8.04%	1.293%
31	21.630	3173	3178	3184	rVB	2018046	3081978	25.77%	4.146%
32	23.289	3458	3460	3462	rBV2	182809	168742	1.41%	0.227%
33	25.007	3745	3752	3761	rBV	1229851	3214322	26.87%	4.324%

Sum of corrected areas: 74343866

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082625\
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Acq On : 26 Aug 2025 15:55
Operator : CG/JU
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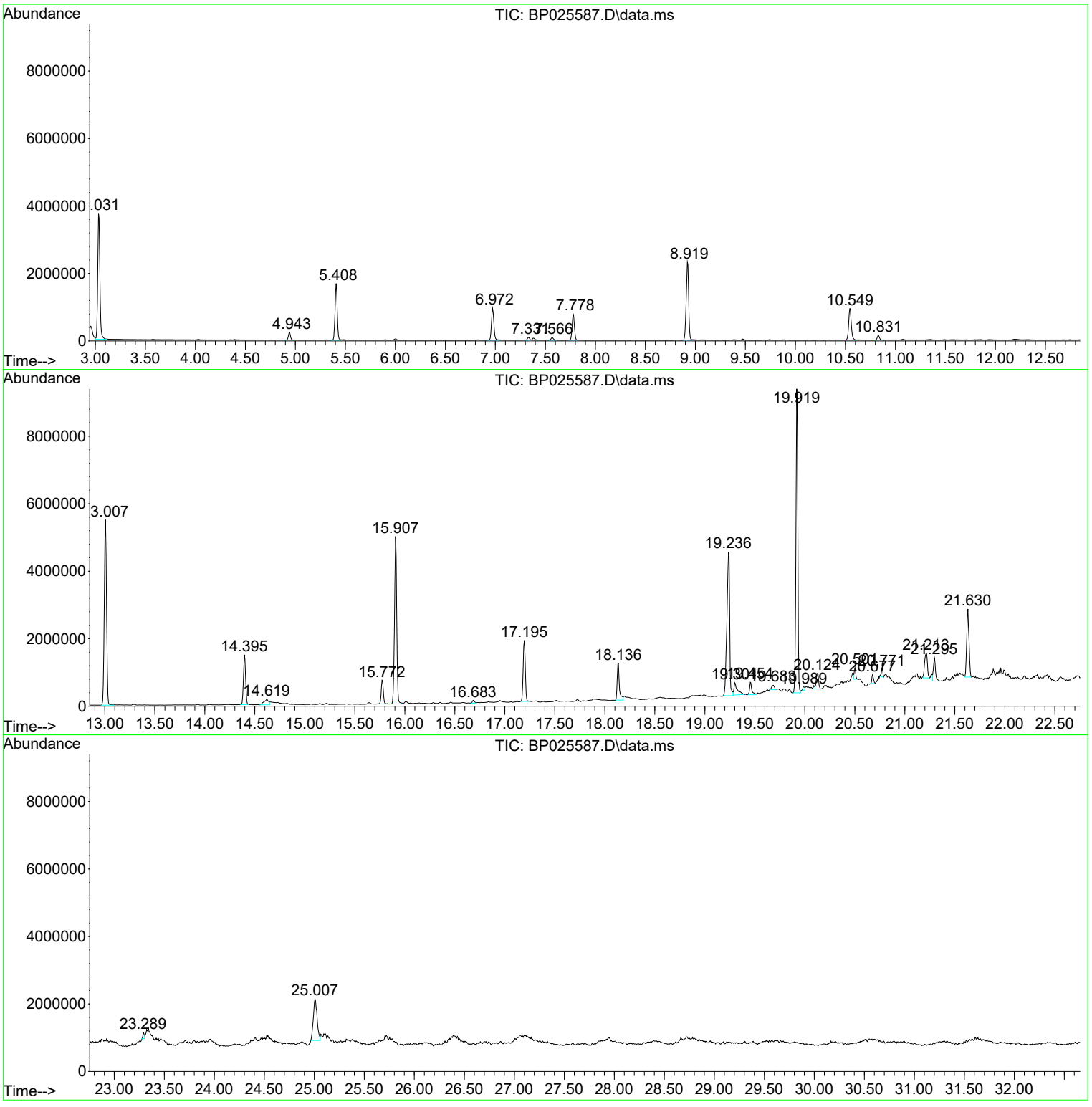
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.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\BP082625\
Data File : BP025587.D
Acq On : 26 Aug 2025 15:55
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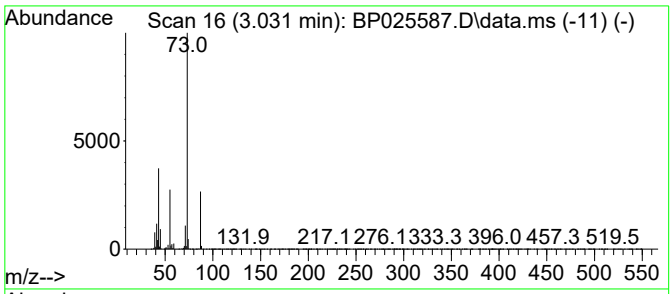
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

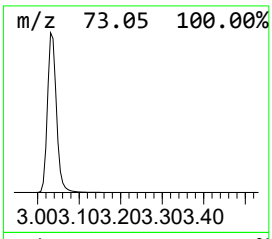
****

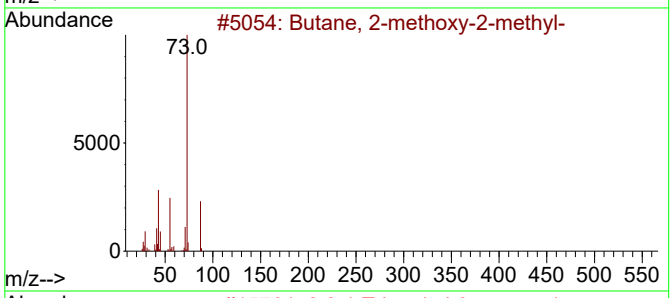
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

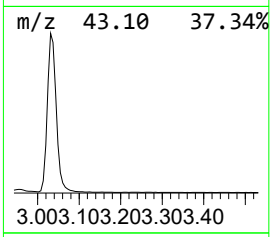
R.T.	EstConc	Area	Relative to ISTD	R.T.
3.031	91.81 ng	5599470	1,4-Dichlorobenzene-d4	7.778

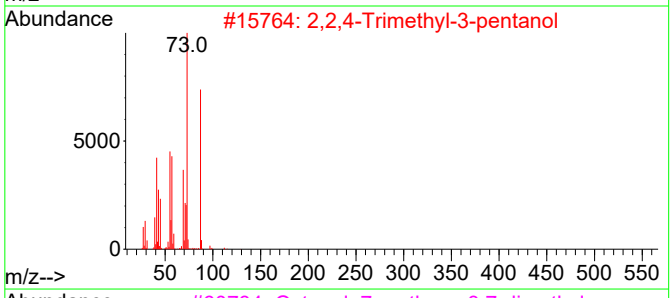
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3	Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	23
4	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17

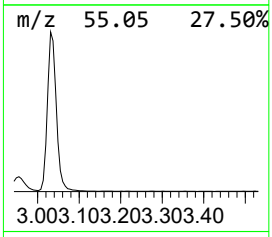


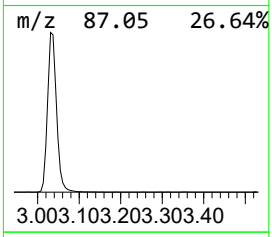


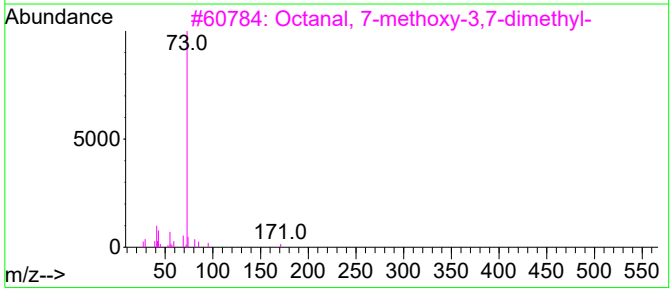


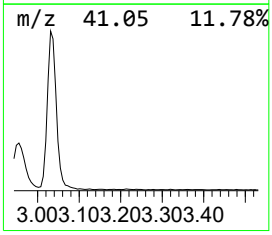












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082625\
Data File : BP025587.D
Acq On : 26 Aug 2025 15:55
Operator : CG/JU
Sample : Q2940-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

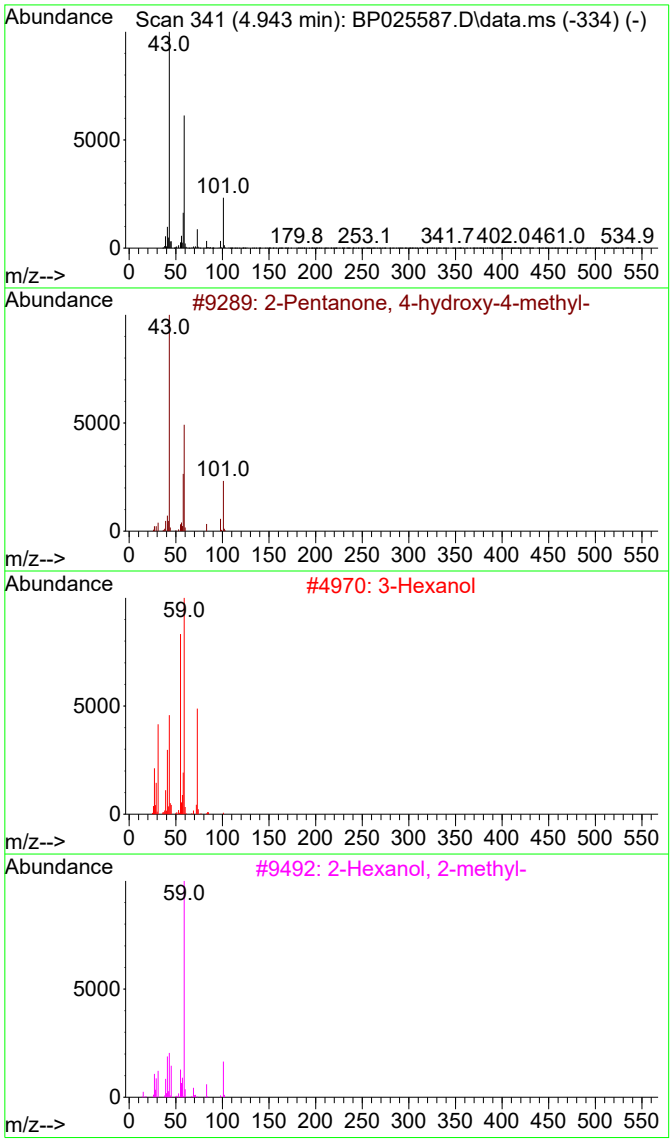
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

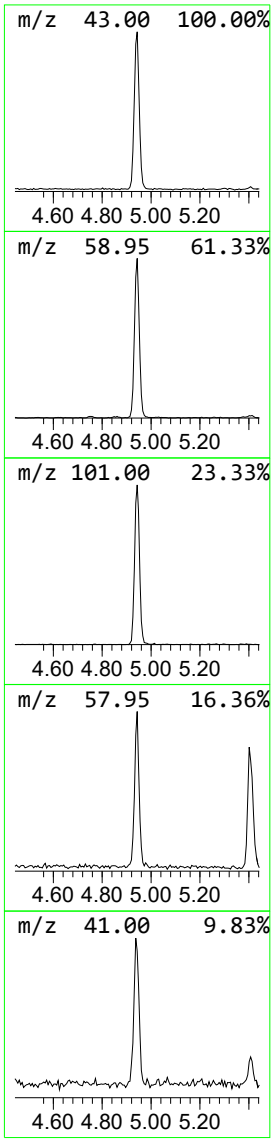
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.943	5.58 ng	340518	1,4-Dichlorobenzene-d4	7.778

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	3-Hexanol	102	C6H14O	000623-37-0	25
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25
4	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	23
5	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	23





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082625\
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Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

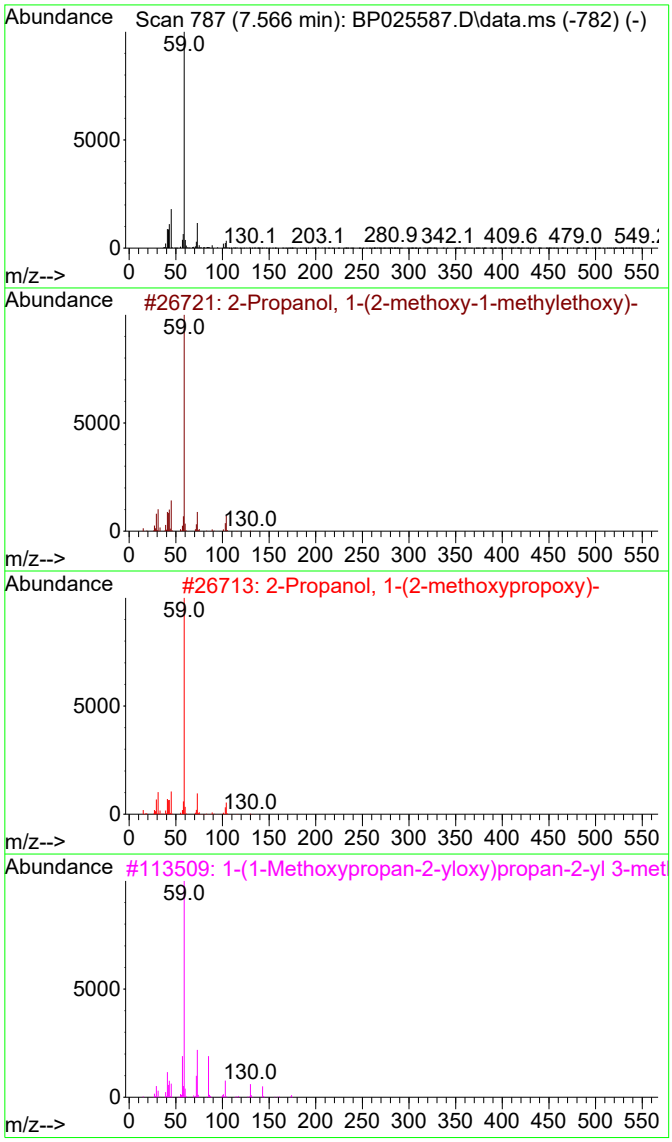
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

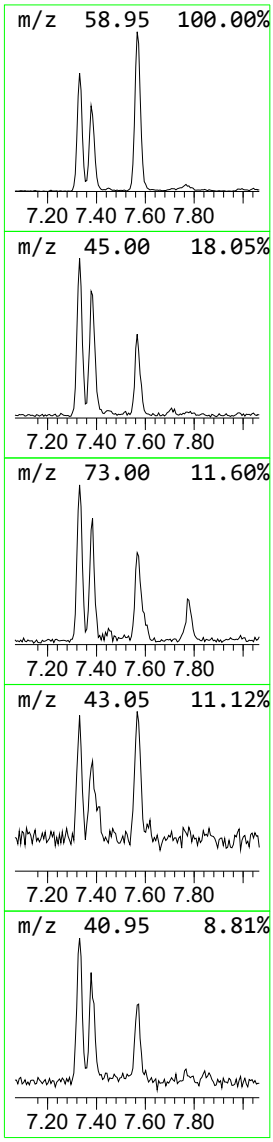
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.566	2.03 ng	123595	1,4-Dichlorobenzene-d4	7.778

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	78
2	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	78
3	1-(1-Methoxypropan-2-yloxy)propa...	232	C12H24O4	1000367-13-2	72
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64
5	Dipropylene glycol (isomer 3)	134	C6H14O3	1000506-25-5	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082625\
Data File : BP025587.D
Acq On : 26 Aug 2025 15:55
Operator : CG/JU
Sample : Q2940-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

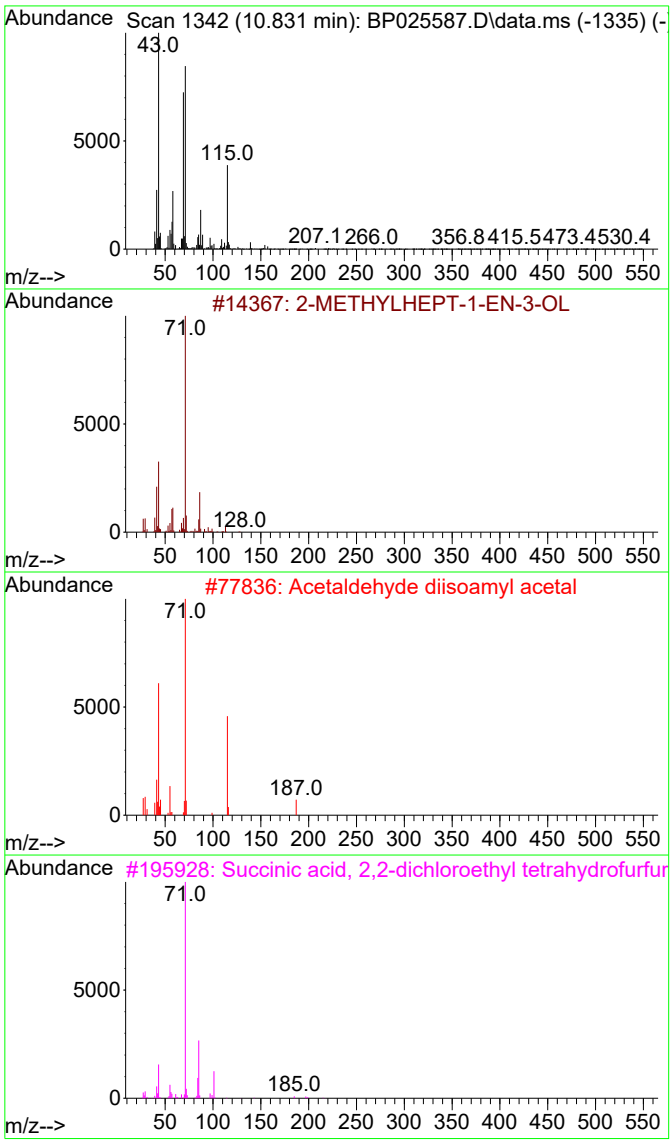
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

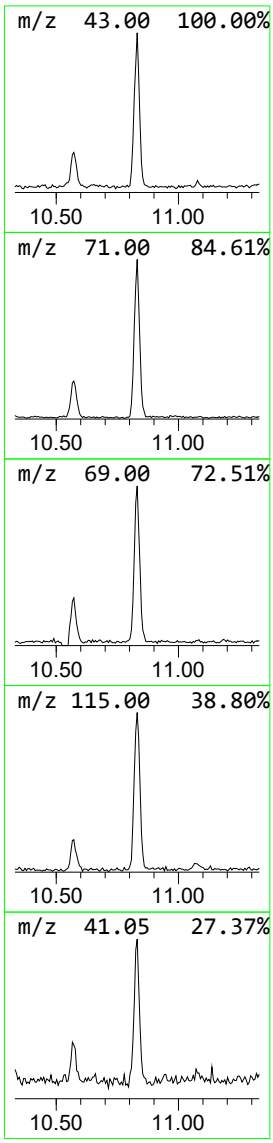
****

Peak Number 4 unknown10.831 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.831	2.58 ng	223111	Naphthalene-d8	10.543

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-METHYLHEPT-1-EN-3-OL	128	C8H16O	013019-19-7	43
2	Acetaldehyde diisoamyl acetal	202	C12H26O2	1000430-80-1	40
3	Succinic acid, 2,2-dichloroethyl...	298	C11H16Cl2O5	1000390-71-6	38
4	1,3-Dimethylbutyl isobutyrate	172	C10H20O2	1000429-31-6	38
5	Glutaric acid, octyl tetrahydrof...	328	C18H32O5	1000359-66-4	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082625\
Data File : BP025587.D
Acq On : 26 Aug 2025 15:55
Operator : CG/JU
Sample : Q2940-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

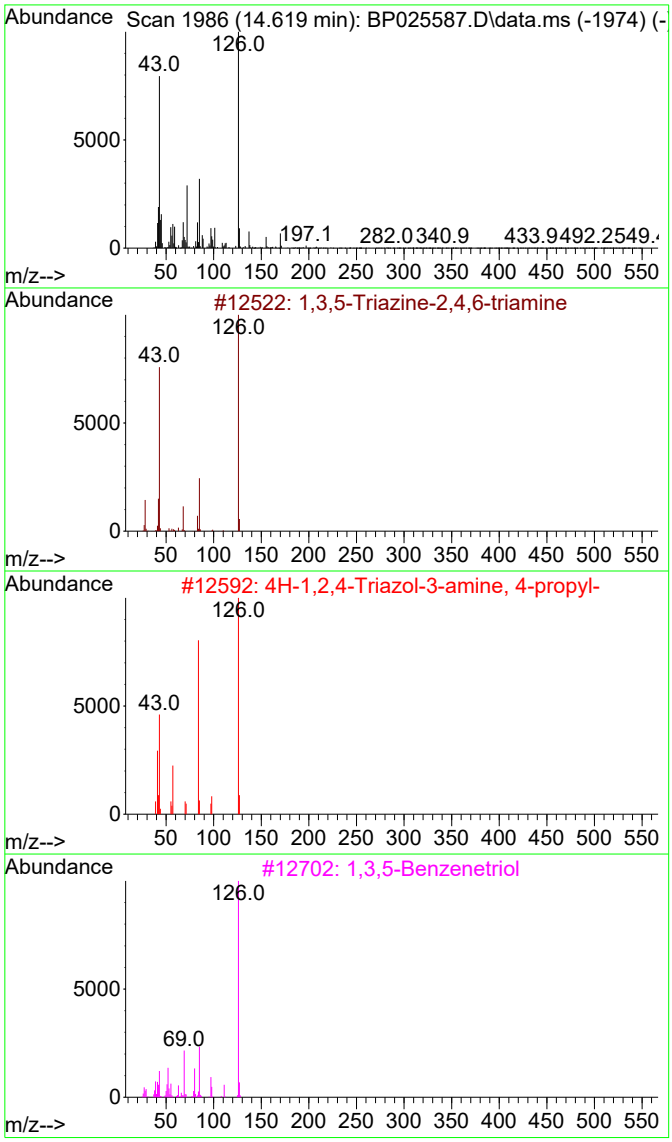
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

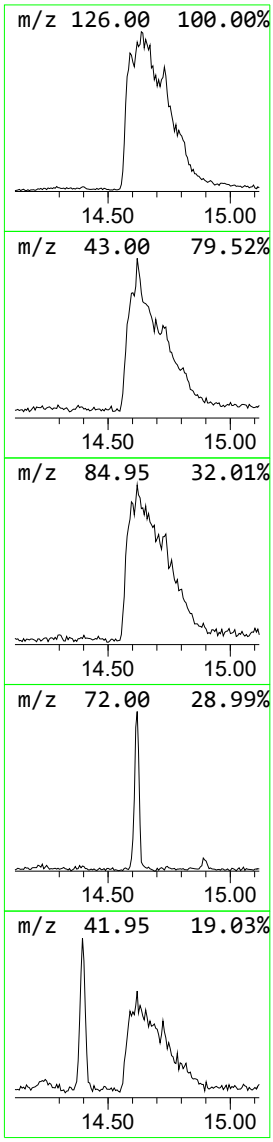
****

Peak Number 5 1,3,5-Triazine-2,4,6-triamine Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.619	4.69 ng	519340	Acenaphthene-d10	14.395

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,5-Triazine-2,4,6-triamine	126	C3H6N6	000108-78-1	52
2	4H-1,2,4-Triazol-3-amine, 4-propyl-	126	C5H10N4	058661-97-5	50
3	1,3,5-Benzenetriol	126	C6H6O3	000108-73-6	49
4	4-Methyl-2-pyrimidinethiol	126	C5H6N2S	035071-17-1	47
5	3H-Pyrazol-3-one, 1,2-dihydro-1,...	126	C6H10N2O	003201-26-1	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082625\
Data File : BP025587.D
Acq On : 26 Aug 2025 15:55
Operator : CG/JU
Sample : Q2940-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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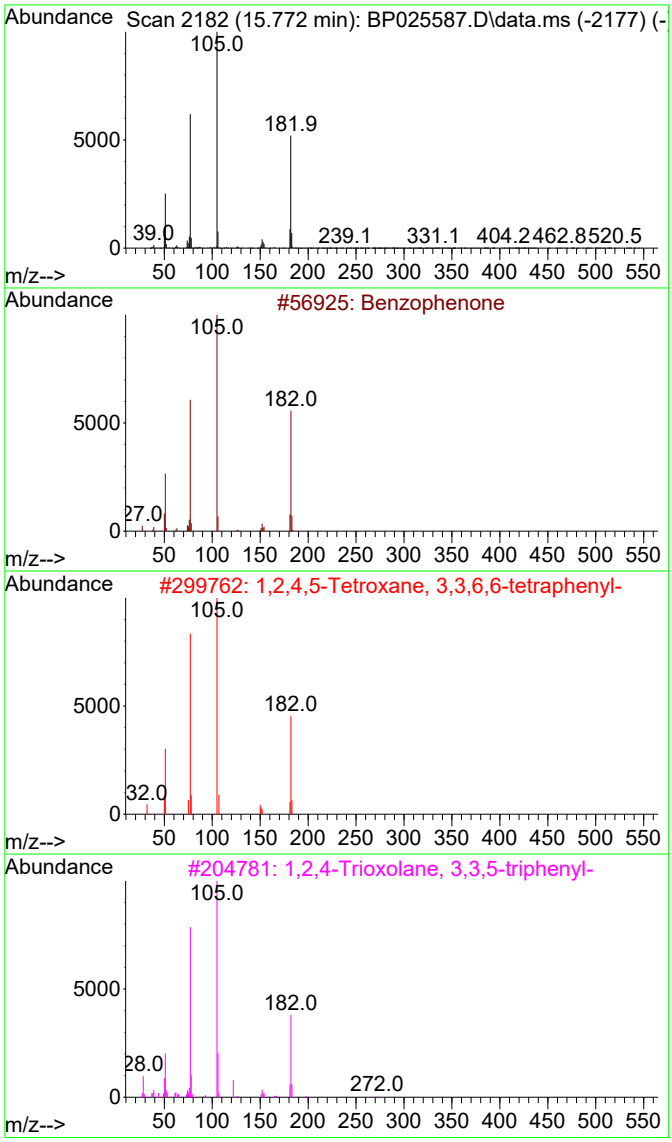
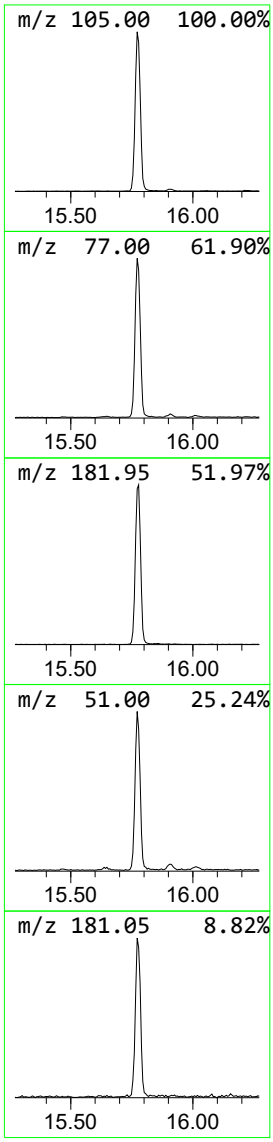
****

Peak Number 6 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.772	9.39 ng	1038710	Acenaphthene-d10	14.395

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	64
3	1,2,4-Trioxolane, 3,3,5-triphenyl-	304	C20H16O3	023246-12-0	45
4	Benzenamine, N-(3-pyridinylmethy...	182	C12H10N2	029722-97-2	45
5	Azobenzene	182	C12H10N2	000103-33-3	42





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082625\
Data File : BP025587.D
Acq On : 26 Aug 2025 15:55
Operator : CG/JU
Sample : Q2940-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

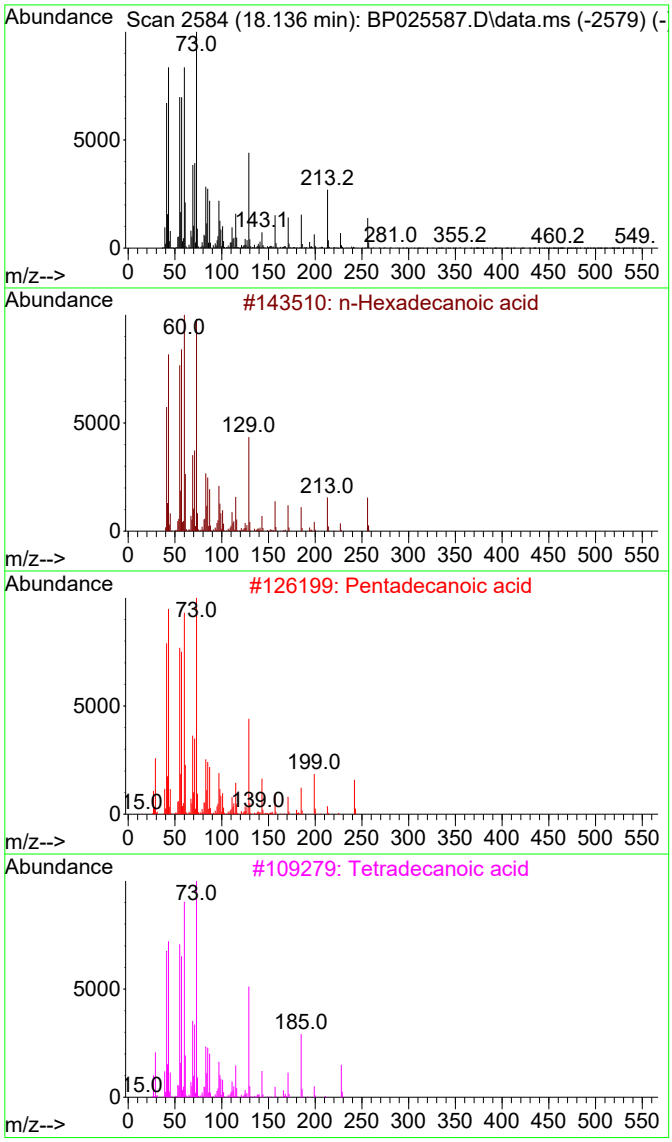
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

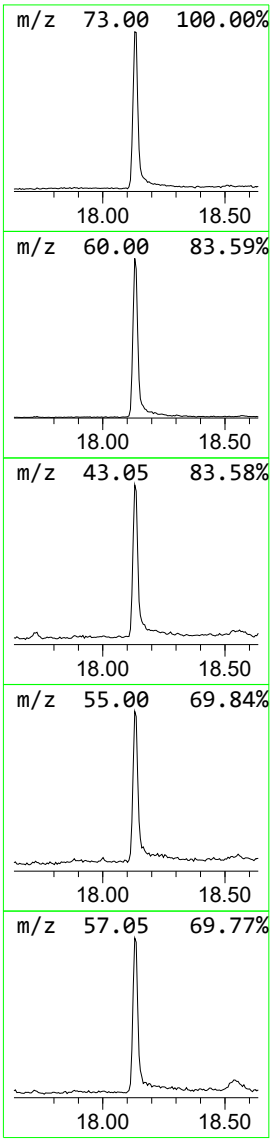
****

Peak Number 7 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.136	13.01 ng	1714440	Phenanthrene-d10	17.195

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	93
4	Octadecanoic acid	284	C18H36O2	000057-11-4	93
5	Tridecanoic acid	214	C13H26O2	000638-53-9	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082625\
Data File : BP025587.D
Acq On : 26 Aug 2025 15:55
Operator : CG/JU
Sample : Q2940-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

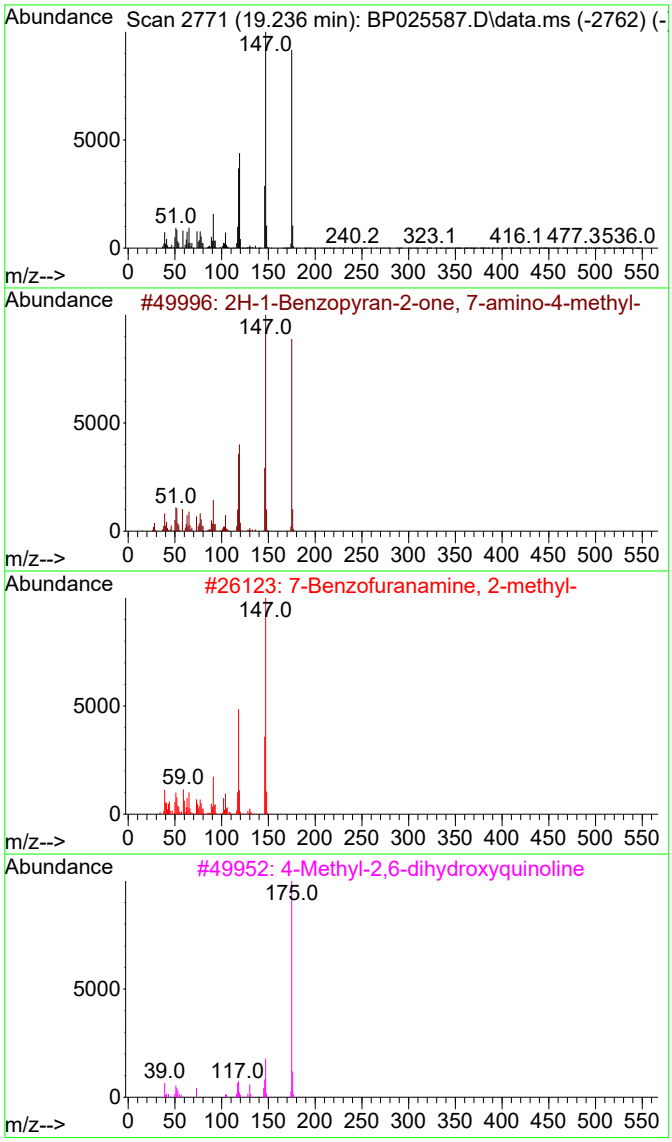
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

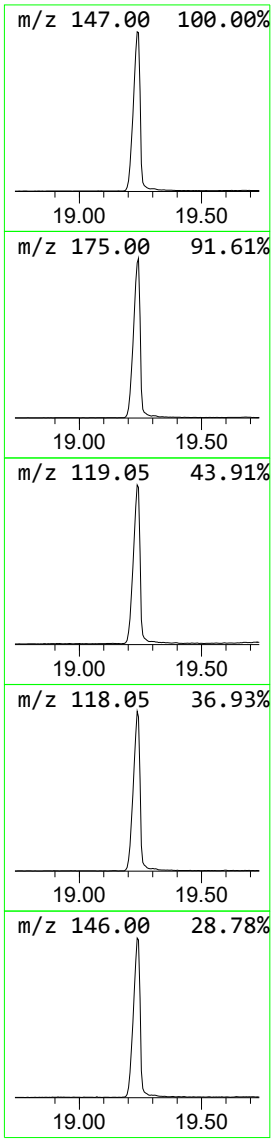
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 2H-1-Benzopyran-2-one, 7-am... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.236	62.89 ng	8289660	Phenanthrene-d10	17.195

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-1-Benzopyran-2-one, 7-amino-4...	175	C10H9NO2	026093-31-2	98
2	7-Benzofuranamine, 2-methyl-	147	C9H9NO	1000327-46-4	64
3	4-Methyl-2,6-dihydroxyquinoline	175	C10H9NO2	034982-01-9	52
4	1H-Inden-1-one, 5-amino-2,3-dihy...	147	C9H9NO	003470-54-0	49
5	2,6-Dimethylphenyl isocyanate	147	C9H9NO	028556-81-2	49





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Data File : BP025587.D
Acq On : 26 Aug 2025 15:55
Operator : CG/JU
Sample : Q2940-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.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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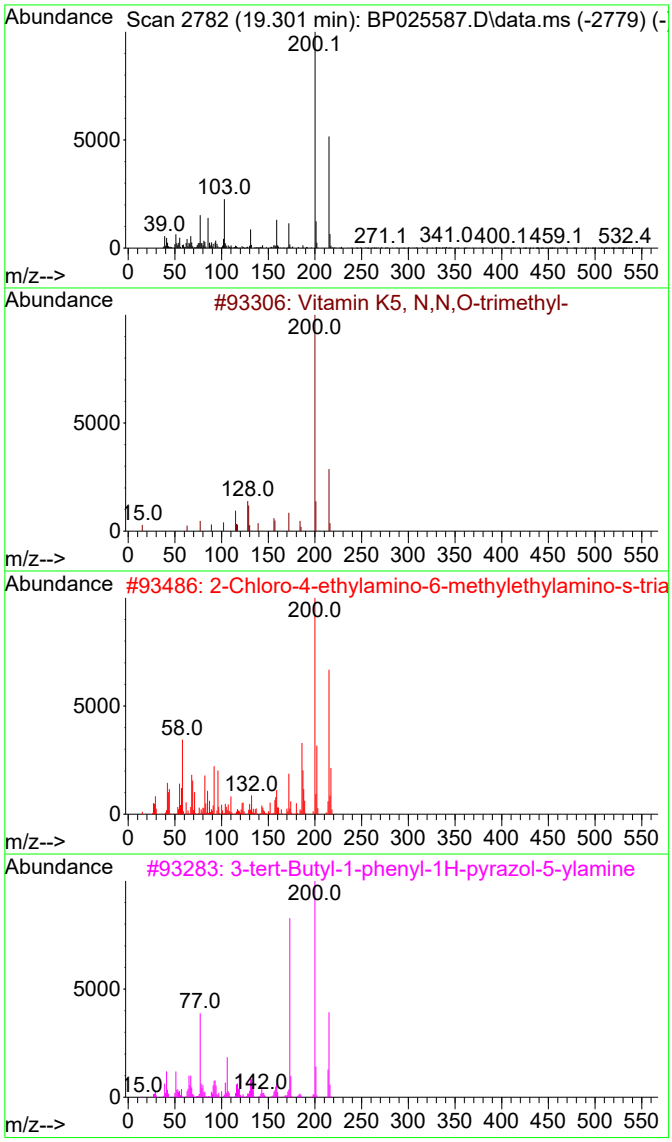
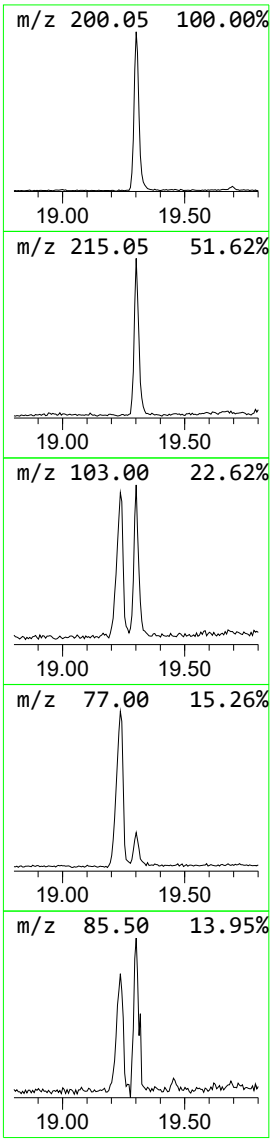
Peak Number 9 Vitamin K5, N,N,O-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.301	6.48 ng	853777	Phenanthrene-d10	17.195

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Vitamin K5, N,N,O-trimethyl-	215	C14H17NO	1000453-07-0	58
2	2-Chloro-4-ethylamino-6-methylet...	215	C8H14ClN5	310901-77-0	47
3	3-tert-Butyl-1-phenyl-1H-pyrazol...	215	C13H17N3	126208-61-5	45
4	Atrazine	215	C8H14ClN5	001912-24-9	45
5	Phenoxathiin	200	C12H8OS	000262-20-4	43





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Sample : Q2940-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.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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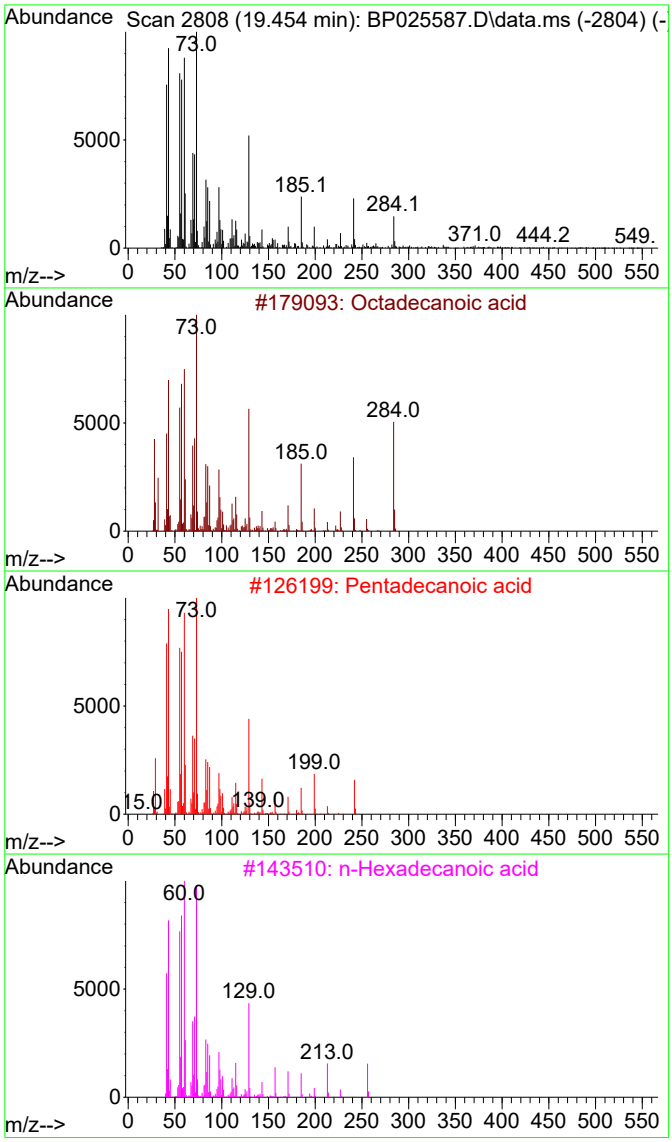
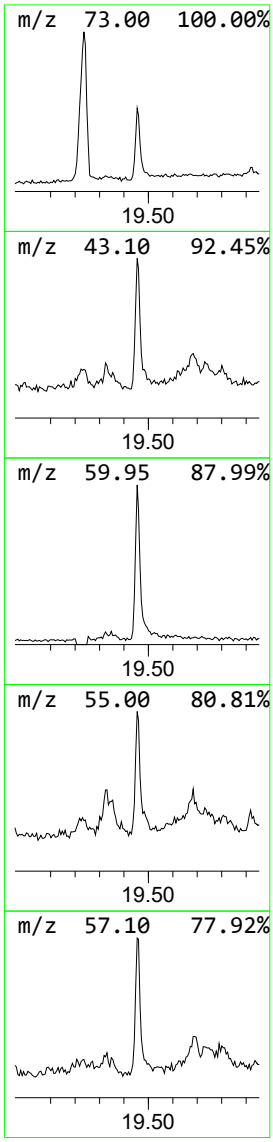
Peak Number 10 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.454	3.82 ng	589357	Chrysene-d12	21.630

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	93
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	74
4	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	72
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	72





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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.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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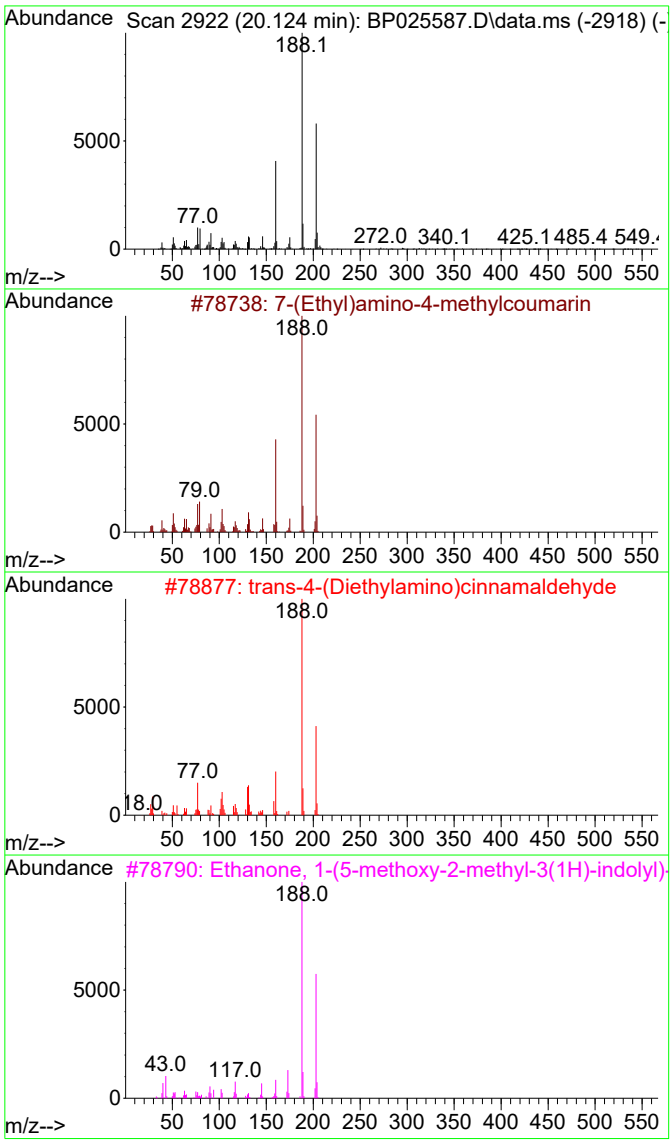
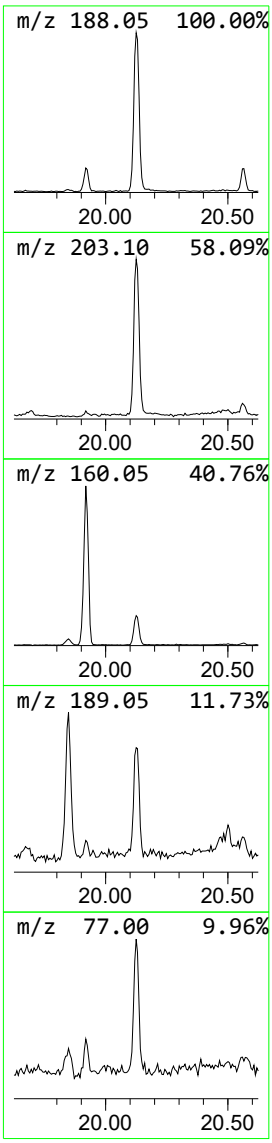
Peak Number 11 7-(Ethyl)amino-4-methylcoum... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.125	4.99 ng	769464	Chrysene-d12	21.630

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-(Ethyl)amino-4-methylcoumarin	203	C12H13NO2	1000395-85-1	97
2	trans-4-(Diethylamino)cinnamalde...	203	C13H17NO	022411-59-2	80
3	Ethanone, 1-(5-methoxy-2-methyl-...	203	C12H13NO2	055895-80-2	62
4	5H-Pyrazolo[3,4-b]quinolin-3-ami...	188	C10H12N4	1000362-40-8	43
5	1-naphthalenamine, 4,8-dimethoxy-	203	C12H13NO2	1000396-05-9	43





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Acq On : 26 Aug 2025 15:55
Operator : CG/JU
Sample : Q2940-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.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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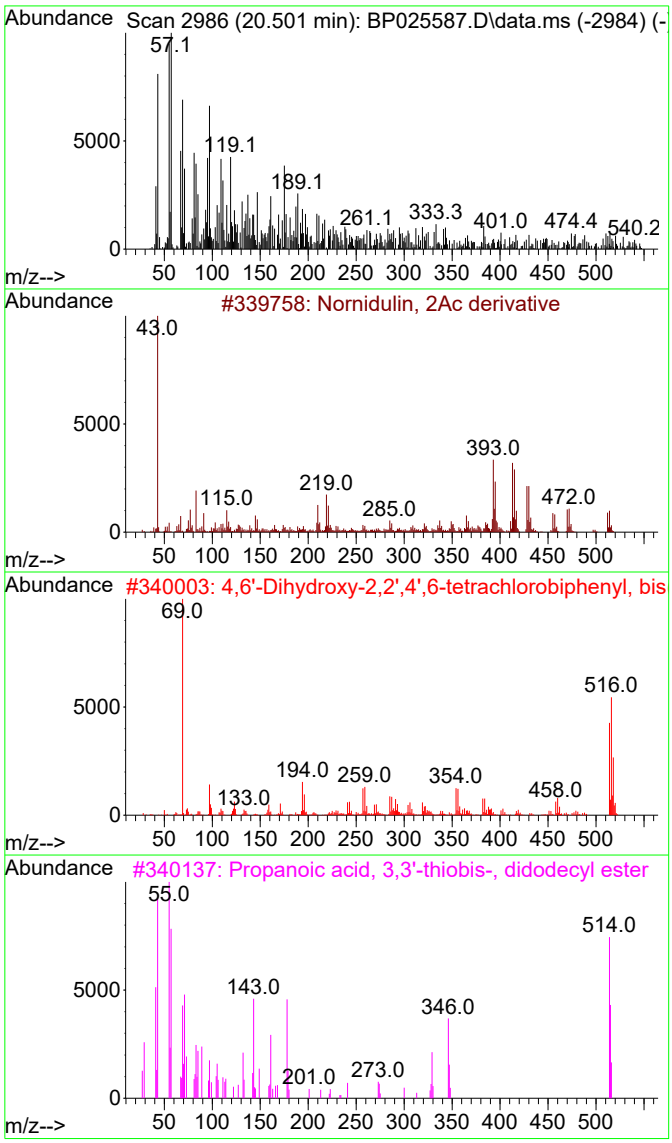
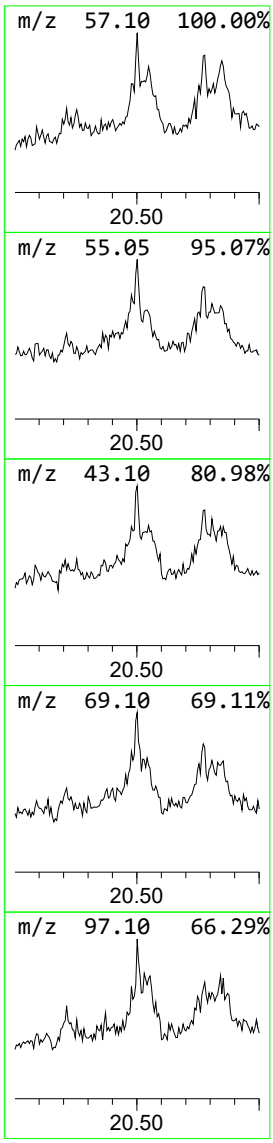
Peak Number 12 unknown20.501 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.501	2.07 ng	318450	Chrysene-d12	21.630

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nornidulin, 2Ac derivative	512	C23H19Cl3O7	1000480-54-4	20
2	4,6'-Dihydroxy-2,2',4',6-tetrach...	514	C16H4Cl4F6O4	1010448-04-3	9
3	Propanoic acid, 3,3'-thiobis-, d...	514	C30H58O4S	000123-28-4	9
4	Lanost-9(11)-en-18-oic acid, 23-...	514	C32H50O5	036872-76-1	9
5	4,4'-Dihydroxy-2,2',6,6'-tetrach...	514	C16H4Cl4F6O4	1000448-04-2	6





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Acq On : 26 Aug 2025 15:55
Operator : CG/JU
Sample : Q2940-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.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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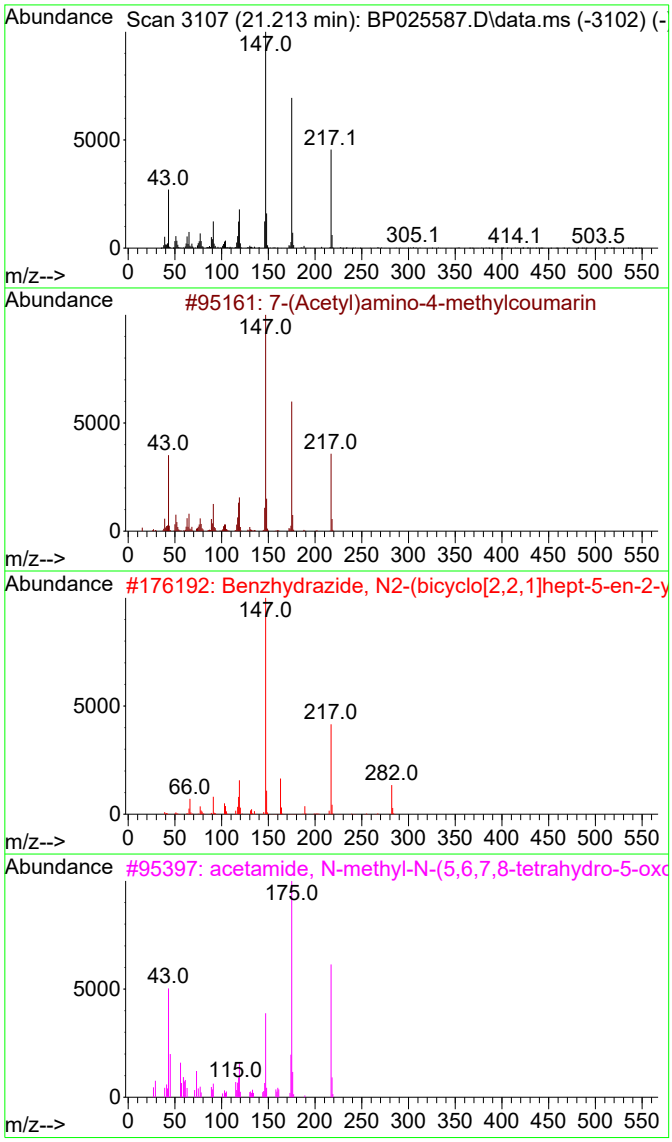
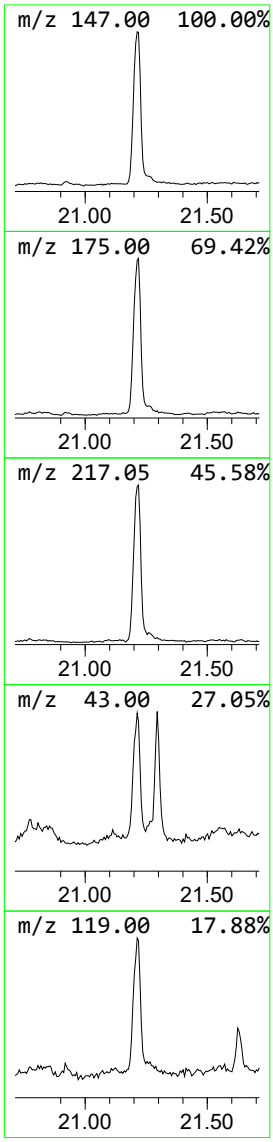
Peak Number 13 7-(Acetyl)amino-4-methylcou... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.213	9.07 ng	1398110	Chrysene-d12	21.630

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-(Acetyl)amino-4-methylcoumarin	217	C12H11NO3	1000395-85-5	95
2	Benzhydrazide, N2-(bicyclo[2,2,1...	282	C18H22N2O	1000273-03-1	53
3	acetamide, N-methyl-N-(5,6,7,8-t...	217	C13H15NO2	1000400-74-0	47
4	2H-1-Benzopyran-2-one, 7-amino-4...	175	C10H9NO2	026093-31-2	46
5	2,6-Dimethylphenyl isocyanate	147	C9H9NO	028556-81-2	43





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Acq On : 26 Aug 2025 15:55
Operator : CG/JU
Sample : Q2940-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.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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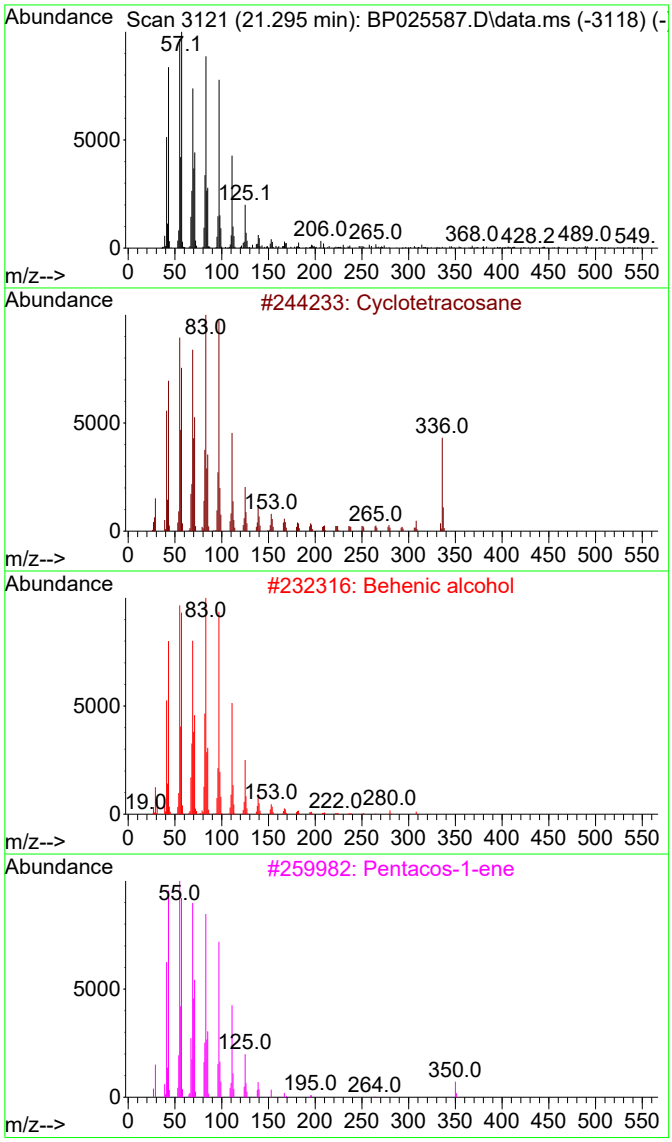
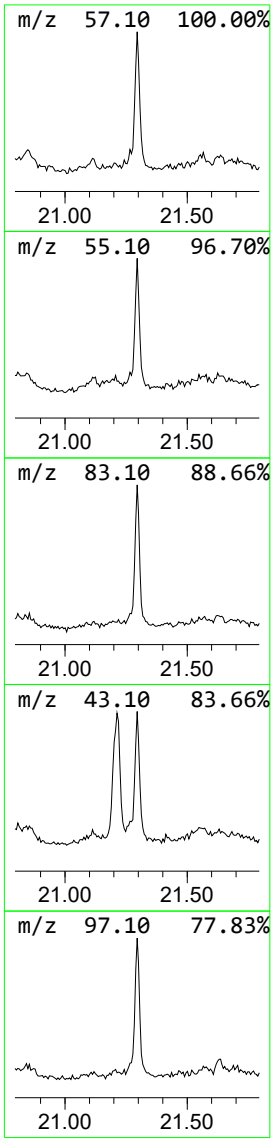
Peak Number 14 Cyclotetracosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.295	6.24 ng	961516	Chrysene-d12	21.630

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetracosane	336	C24H48	000297-03-0	97
2	Behenic alcohol	326	C22H46O	000661-19-8	95
3	Pentacos-1-ene	350	C25H50	016980-85-1	94
4	1-Heneicosanol	312	C21H44O	015594-90-8	94
5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	93





Tentatively Identified Compound (LSC) summary

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Data File : BP025587.D
Acq On : 26 Aug 2025 15:55
Operator : CG/JU
Sample : Q2940-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP(1-6)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.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...	3.031	91.8	ng	5599470	1	7.778	1219800	20.0
2-Pentanone, 4-...	4.943	5.6	ng	340518	1	7.778	1219800	20.0
2-Propanol, 1-(...	7.566	2.0	ng	123595	1	7.778	1219800	20.0
unknown10.831	10.831	2.6	ng	223111	2	10.543	1728020	20.0
1,3,5-Triazine-...	14.619	4.7	ng	519340	3	14.395	2212470	20.0
Benzophenone	15.772	9.4	ng	1038710	3	14.395	2212470	20.0
n-Hexadecanoic ...	18.136	13.0	ng	1714440	4	17.195	2636210	20.0
2H-1-Benzopyran...	19.236	62.9	ng	8289660	4	17.195	2636210	20.0
Vitamin K5, N,N...	19.301	6.5	ng	853777	4	17.195	2636210	20.0
Octadecanoic acid	19.454	3.8	ng	589357	5	21.630	3081980	20.0
7-(Ethyl)amino-...	20.125	5.0	ng	769464	5	21.630	3081980	20.0
unknown20.501	20.501	2.1	ng	318450	5	21.630	3081980	20.0
7-(Acetyl)amino...	21.213	9.1	ng	1398110	5	21.630	3081980	20.0
Cyclotetracosane	21.295	6.2	ng	961516	5	21.630	3081980	20.0