LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
Data File : BP025536.D
Acq On : 21 Aug 2025 12:12
Operator : CG/JU
Sample : Q2819-12
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
17M-S

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: BP025536.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.037	12	17	30	rVB	1029946	1322365	15.37%	2.093%
2	4.943	335	341	348	rVB	248472	347210	4.04%	0.550%
3	5.402	412	419	432	rBV	2532189	3747610	43.56%	5.932%
4	6.961	676	684	698	rBV	2487590	3973034	46.18%	6.289%
5	7.784	815	824	832	rBV	794349	1293405	15.03%	2.047%
6	8.919	1009	1017	1029	rBV	1582082	2712377	31.53%	4.294%
7	10.549	1285	1294	1307	rBV	1052335	1818367	21.14%	2.878%
8	13.013	1704	1713	1724	rBV	4070516	6150966	71.50%	9.737%
9	14.401	1938	1949	1957	rBV	1522507	2414246	28.06%	3.822%
10	15.784	2177	2184	2191	rBV	693812	998003	11.60%	1.580%
11	15.919	2196	2207	2216	rBV	3290064	4871857	56.63%	7.712%
12	16.331	2271	2277	2279	rBV3	61259	100497	1.17%	0.159%
13	16.601	2317	2323	2330	rBV5	58522	107299	1.25%	0.170%
14	16.672	2330	2335	2343	rBV2	59639	154551	1.80%	0.245%
15	17.201	2418	2425	2430	rBV	1926538	2832100	32.92%	4.483%
16	17.242	2430	2432	2438	rVB	196091	258782	3.01%	0.410%
17	18.101	2573	2578	2579	rBV2	162272	187587	2.18%	0.297%
18	18.136	2580	2584	2593	rVB	866144	1242907	14.45%	1.967%
19	18.525	2643	2650	2658	rBV3	1926437	3454903	40.16%	5.469%
20	18.595	2658	2662	2666	rVV4	108476	164609	1.91%	0.261%
21	18.654	2668	2672	2677	rVB2	138190	175439	2.04%	0.278%
22	19.019	2729	2734	2738	rBV6	51294	87754	1.02%	0.139%
23	19.178	2755	2761	2767	rBV6	163642	291668	3.39%	0.462%
24	19.266	2771	2776	2782	rVB3	182509	308053	3.58%	0.488%
25	19.325	2782	2786	2791	rBV	454946	626882	7.29%	0.992%
26	19.466	2804	2810	2811	rBV	449585	589963	6.86%	0.934%
27	19.648	2835	2841	2846	rBV2	933501	1286808	14.96%	2.037%
28	19.701	2846	2850	2859	rVV2	390169	682419	7.93%	1.080%
29	19.777	2860	2863	2868	rVV4	123466	176551	2.05%	0.279%
30	19.919	2881	2887	2894	rBV	6349769	8602709	100.00%	13.618%
31	20.225	2936	2939	2944	rBV	102331	134715	1.57%	0.213%
32	20.666	3010	3014	3019	rVB3	290494	348289	4.05%	0.551%
33	21.313	3119	3124	3137	rBV2	990793	1603931	18.64%	2.539%
34	21.648	3173	3181	3187	rBV	1804863	3310497	38.48%	5.240%
35	21.701	3187	3190	3197	rVB	200154	302397	3.52%	0.479%
36	22.577	3333	3339	3354	rVB	616932	1365912	15.88%	2.162%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
Data File : BP025536.D
Acq On : 21 Aug 2025 12:12
Operator : CG/JU
Sample : Q2819-12
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
17M-S

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

37	23.666	3520	3524	3536	rVB2	296056	620764	7.22%	0.983%
38	24.230	3615	3620	3633	rVB2	486071	1309119	15.22%	2.072%
39	25.018	3745	3754	3765	rBV	1056648	3077571	35.77%	4.872%
40	26.530	4008	4011	4012	rBV2	112311	118187	1.37%	0.187%

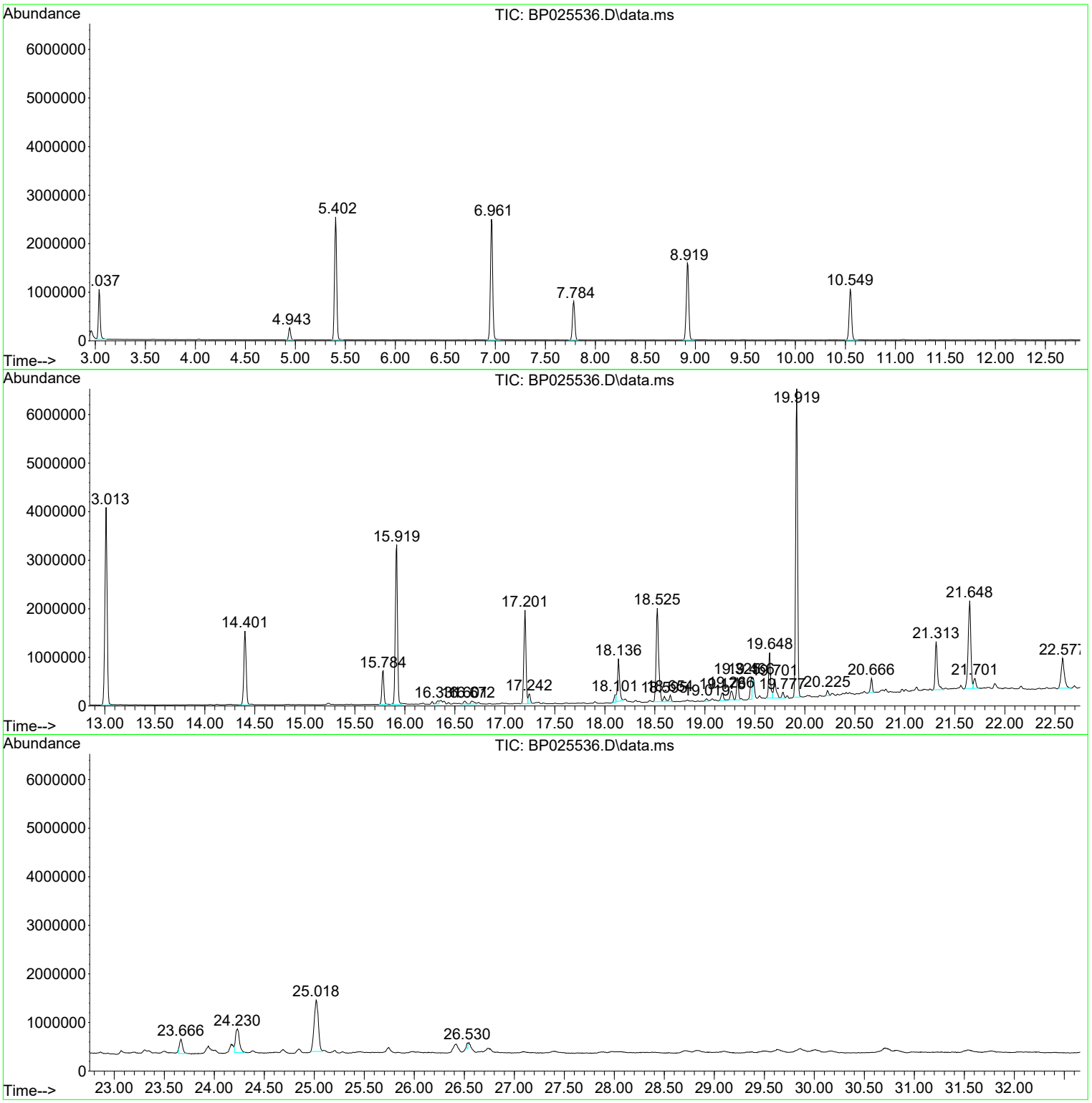
Sum of corrected areas: 63172303

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
Data File : BP025536.D
Acq On : 21 Aug 2025 12:12
Operator : CG/JU
Sample : Q2819-12
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
17M-S

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\BP082025\
Data File : BP025536.D
Acq On : 21 Aug 2025 12:12
Operator : CG/JU
Sample : Q2819-12
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
17M-S

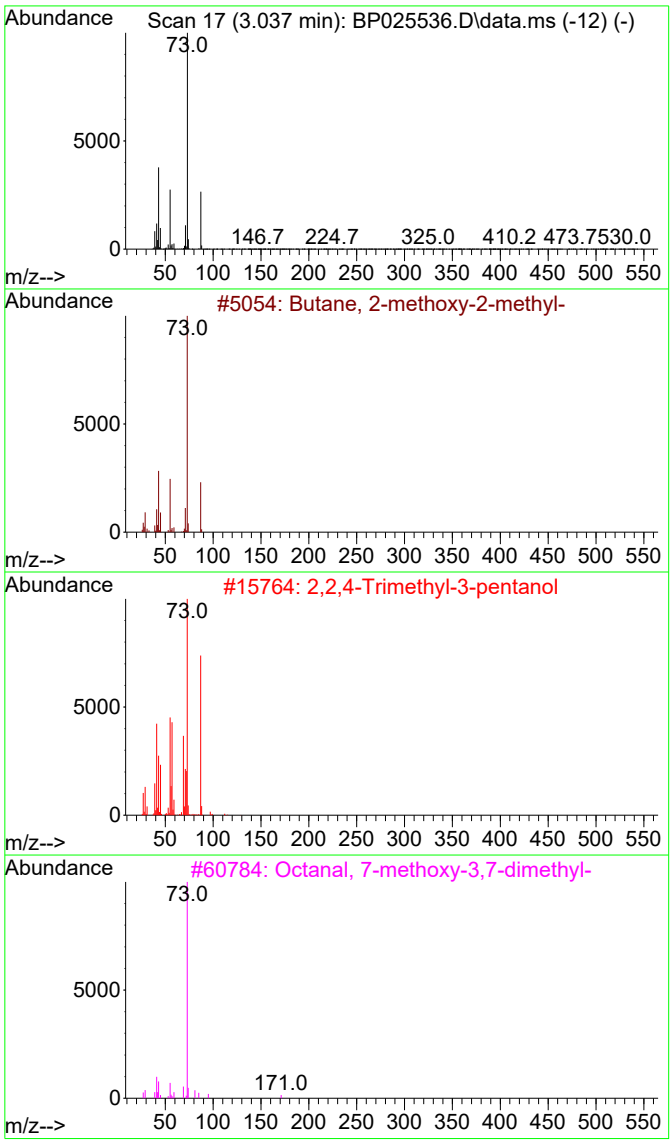
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

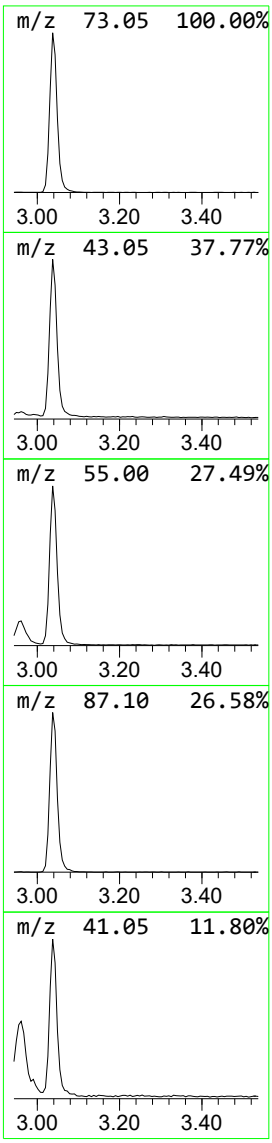
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.037	20.45 ng	1322370	1,4-Dichlorobenzene-d4	7.784

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	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
Data File : BP025536.D
Acq On : 21 Aug 2025 12:12
Operator : CG/JU
Sample : Q2819-12
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
17M-S

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

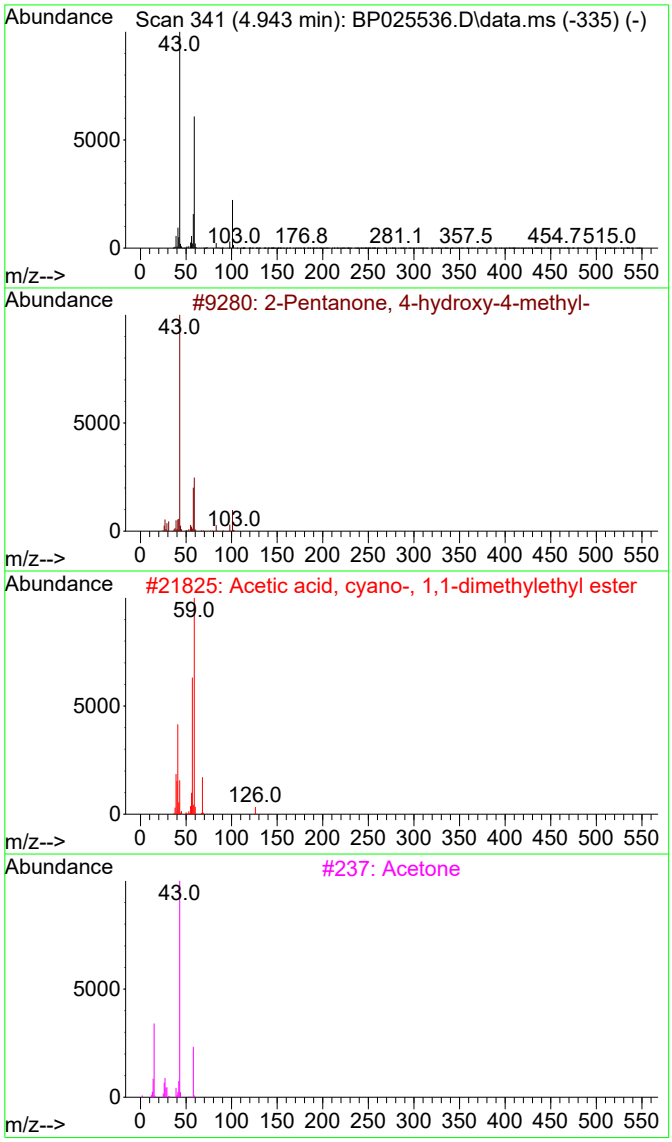
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

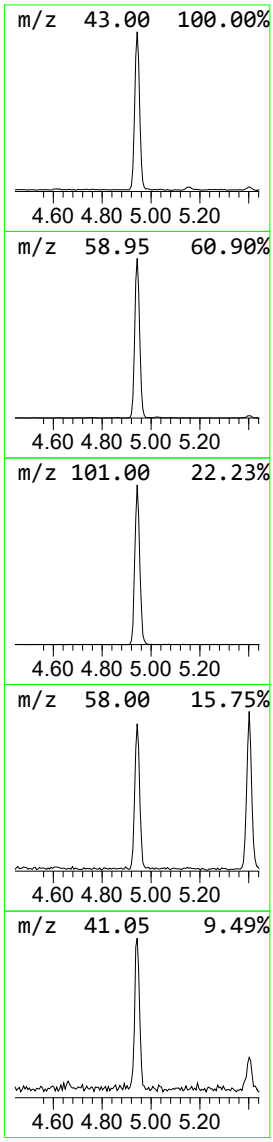
****

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.943	5.37 ng	347210	1,4-Dichlorobenzene-d4	7.784

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
3	Acetone	58	C3H6O	000067-64-1	10
4	(+-)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
Data File : BP025536.D
Acq On : 21 Aug 2025 12:12
Operator : CG/JU
Sample : Q2819-12
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
17M-S

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

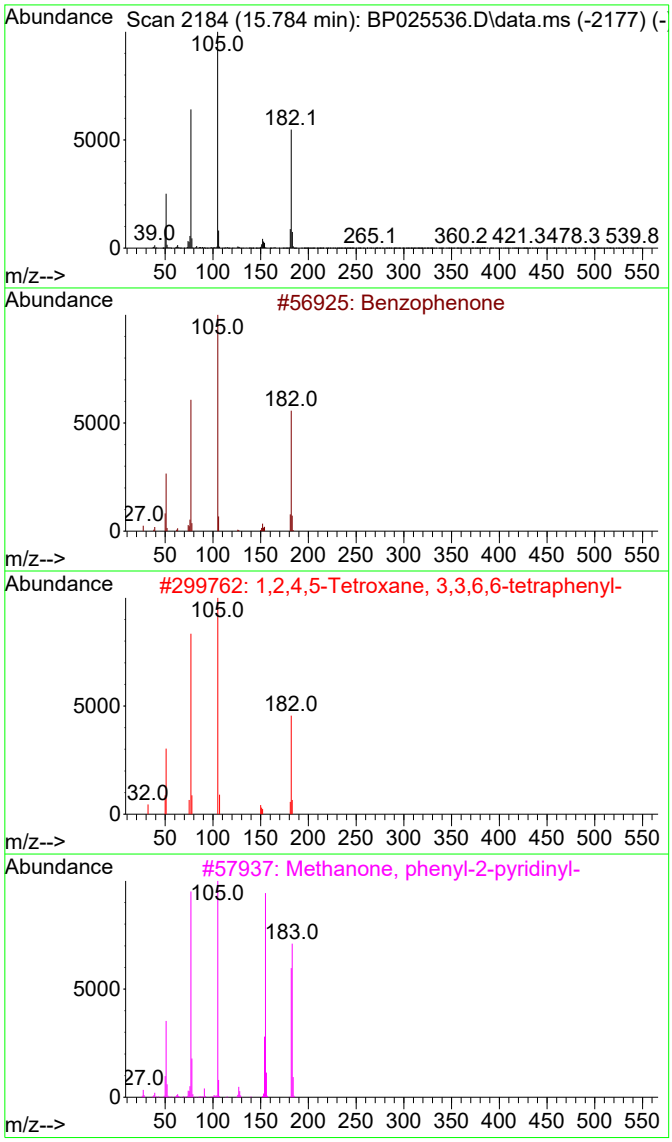
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

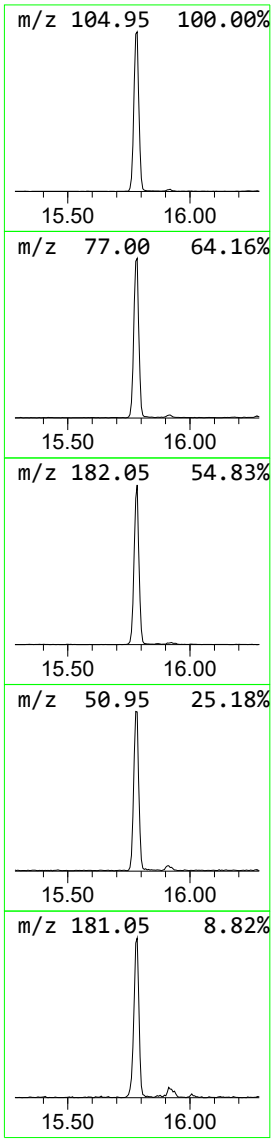
****

Peak Number 3 Benzophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.784	8.27 ng	998003	Acenaphthene-d10	14.401

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	Methanone, phenyl-2-pyridinyl-	183	C12H9NO	000091-02-1	56
4	Azobenzene	182	C12H10N2	000103-33-3	45
5	Benzenamine, N-(3-pyridinylmethy...	182	C12H10N2	029722-97-2	45





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
Data File : BP025536.D
Acq On : 21 Aug 2025 12:12
Operator : CG/JU
Sample : Q2819-12
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
17M-S

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

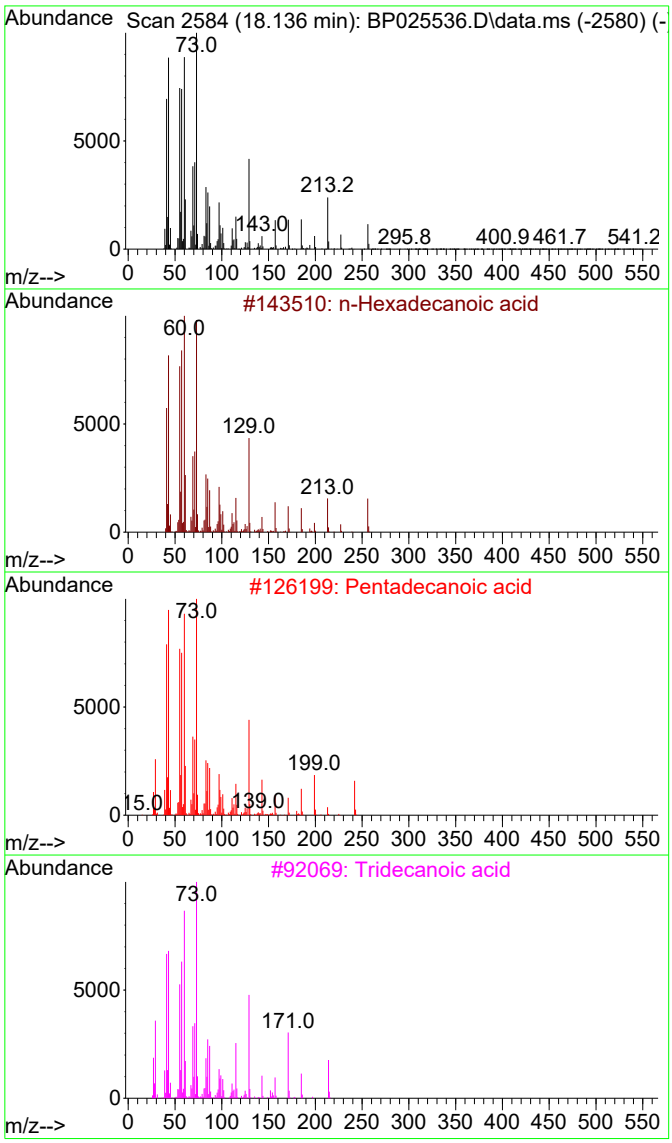
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

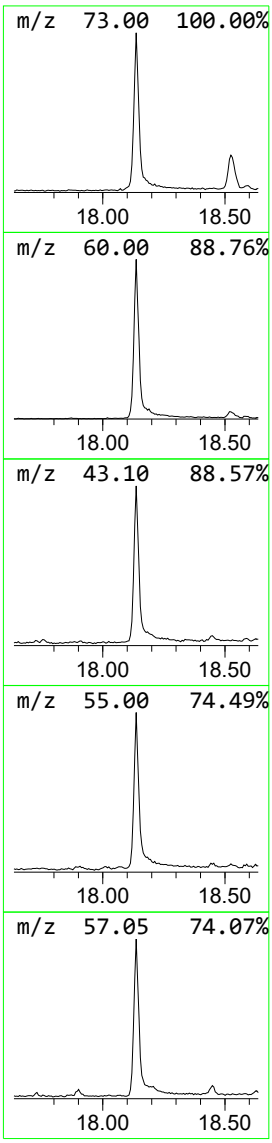
****

Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.136	8.78 ng	1242910	Phenanthrene-d10	17.201

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	Tridecanoic acid	214	C13H26O2	000638-53-9	92
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5	n-Decanoic acid	172	C10H20O2	000334-48-5	76





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
Data File : BP025536.D
Acq On : 21 Aug 2025 12:12
Operator : CG/JU
Sample : Q2819-12
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
17M-S

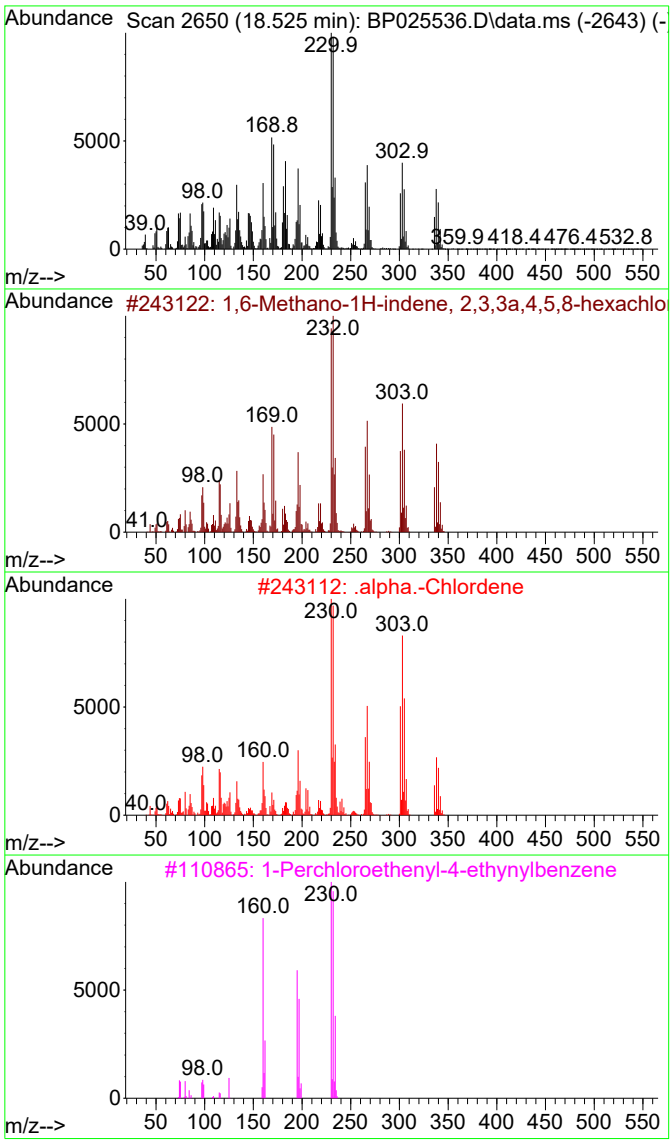
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

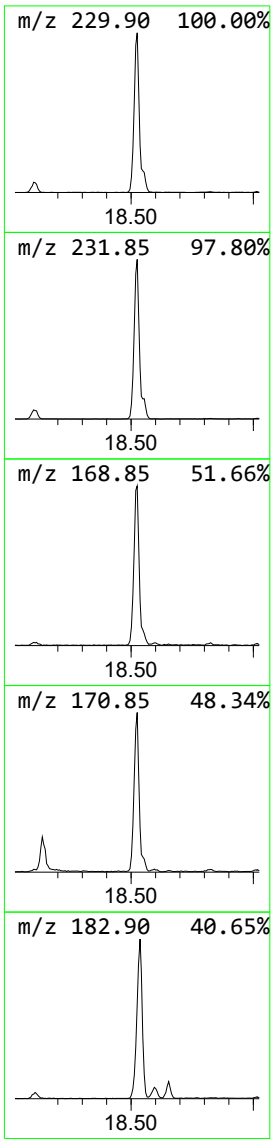
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1,6-Methano-1H-indene, 2,3,... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.525	24.40 ng	3454900	Phenanthrene-d10	17.201

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,6-Methano-1H-indene, 2,3,3a,4,...	336	C10H6Cl6	056641-38-4	98
2	.alpha.-Chlordene	336	C10H6Cl6	056534-02-2	91
3	1-Perchloroethenyl-4-ethynylbenzene	230	C10H5Cl3	1000283-76-9	90
4	1-(5-Hydroxy-3-pyridin-4-yl-5-tr...	301	C13H14F3N3O2	1000311-57-2	59
5	Naphthalene, 2,3,6-trichloro-	230	C10H5Cl3	055720-40-6	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
Data File : BP025536.D
Acq On : 21 Aug 2025 12:12
Operator : CG/JU
Sample : Q2819-12
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ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
17M-S

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

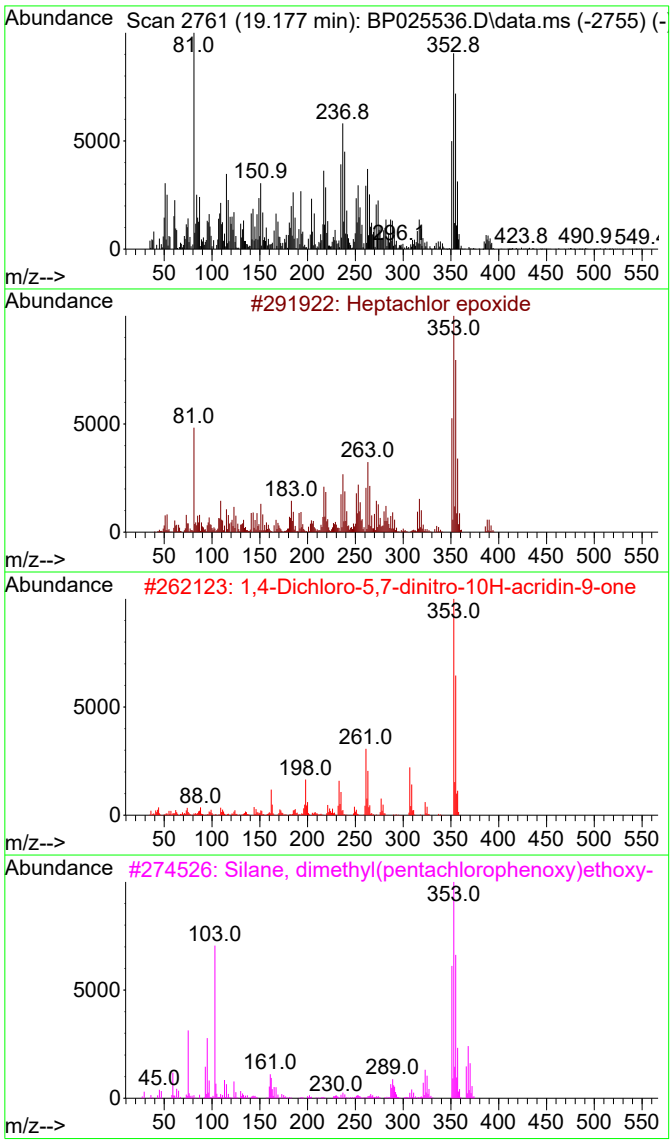
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

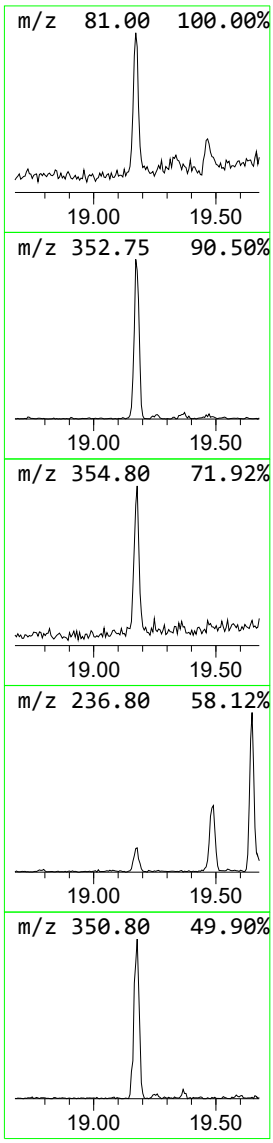
****

Peak Number 6 Heptachlor epoxide Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.177	2.06 ng	291668	Phenanthrene-d10	17.201

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptachlor epoxide	386	C10H5Cl7O	001024-57-3	95
2	1,4-Dichloro-5,7-dinitro-10H-acr...	353	C13H5Cl2N3O5	148877-13-8	38
3	Silane, dimethyl(pentachlorophen...	366	C10H11Cl5O2Si	1010347-47-6	32
4	Titanium, chlorobis[(1,2,3,4,5-....	353	C20H30ClTi	073348-99-9	27
5	7-Chloro-3,4-dihydro-3-[4-methox...	353	C20H16ClNO3	144155-15-7	20





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Data File : BP025536.D
Acq On : 21 Aug 2025 12:12
Operator : CG/JU
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Misc :
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BNA_P
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17M-S

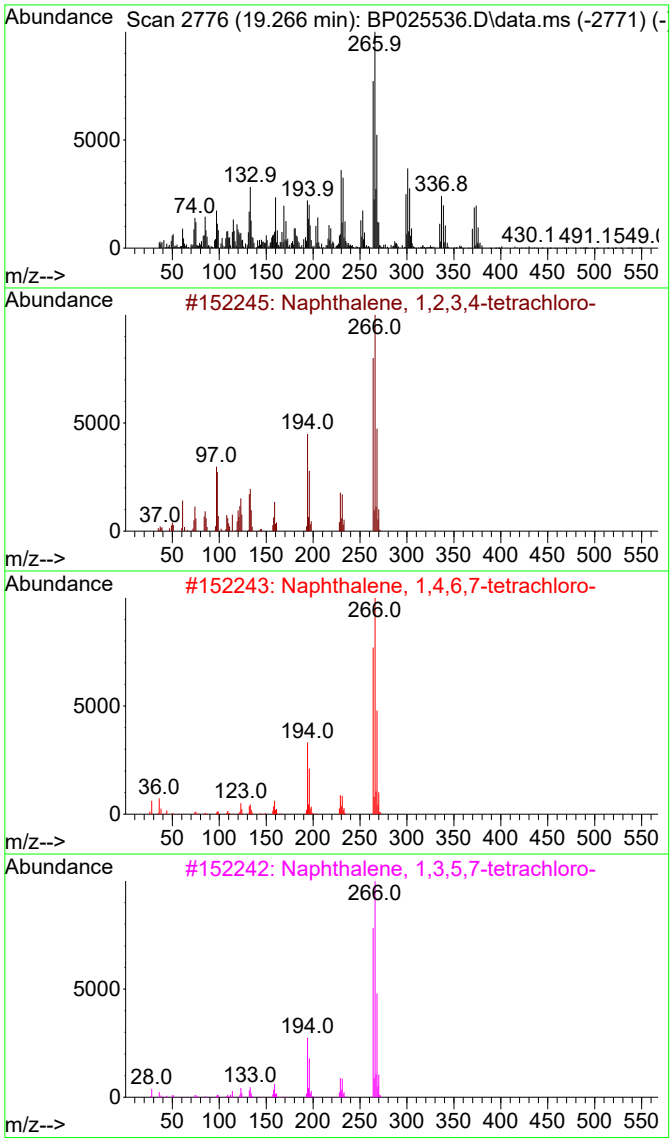
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

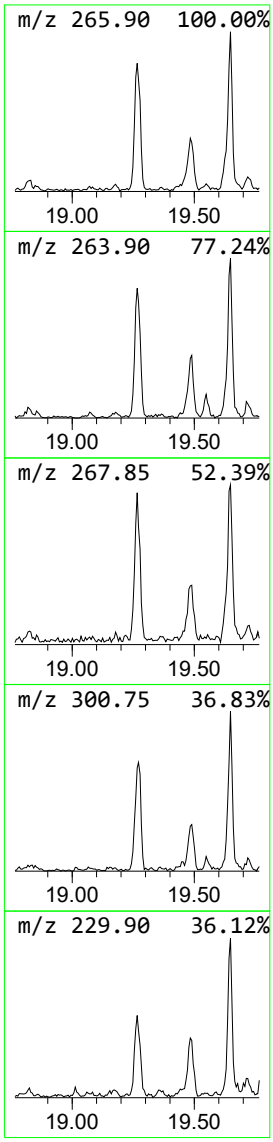
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Naphthalene, 1,2,3,4-tetrac... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.266	2.18 ng	308053	Phenanthrene-d10	17.201

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrachloro-	264	C10H4Cl4	020020-02-4	78
2	Naphthalene, 1,4,6,7-tetrachloro-	264	C10H4Cl4	055720-43-9	78
3	Naphthalene, 1,3,5,7-tetrachloro-	264	C10H4Cl4	053555-64-9	78
4	Tetrachloroisophthalonitrile	264	C8Cl4N2	001897-45-6	53
5	Tetrachloroterephthalonitrile	264	C8Cl4N2	001897-41-2	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
Data File : BP025536.D
Acq On : 21 Aug 2025 12:12
Operator : CG/JU
Sample : Q2819-12
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
17M-S

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

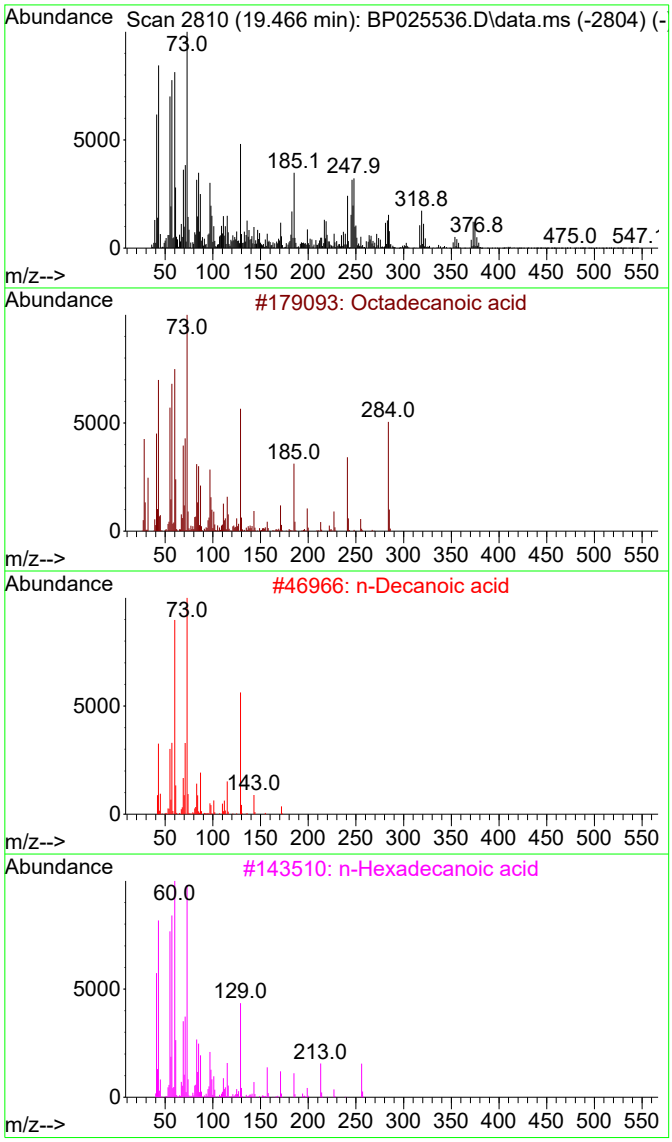
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

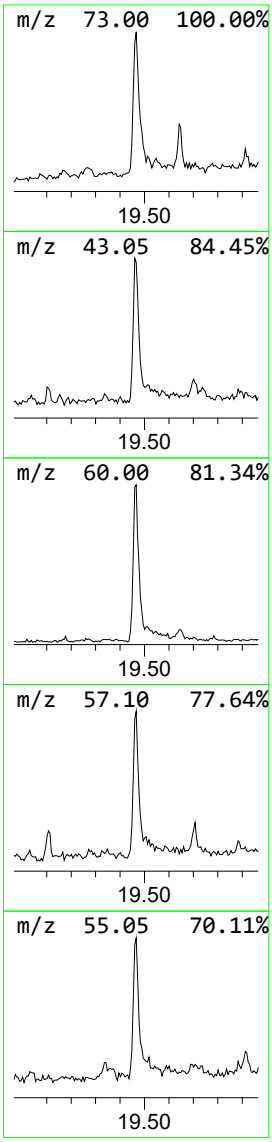
****

Peak Number 8 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.466	3.56 ng	589963	Chrysene-d12	21.648

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	97
2	n-Decanoic acid	172	C10H20O2	000334-48-5	89
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	89
4	Tricosanoic acid	354	C23H46O2	002433-96-7	72
5	Tridecanoic acid	214	C13H26O2	000638-53-9	64





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Data File : BP025536.D
Acq On : 21 Aug 2025 12:12
Operator : CG/JU
Sample : Q2819-12
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
17M-S

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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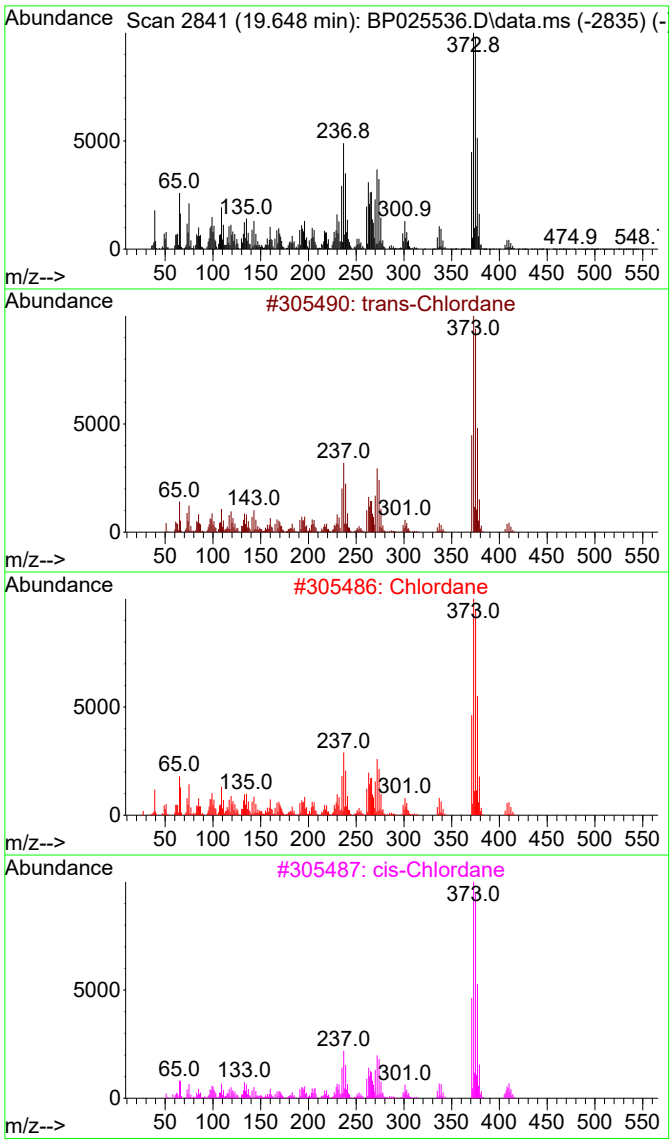
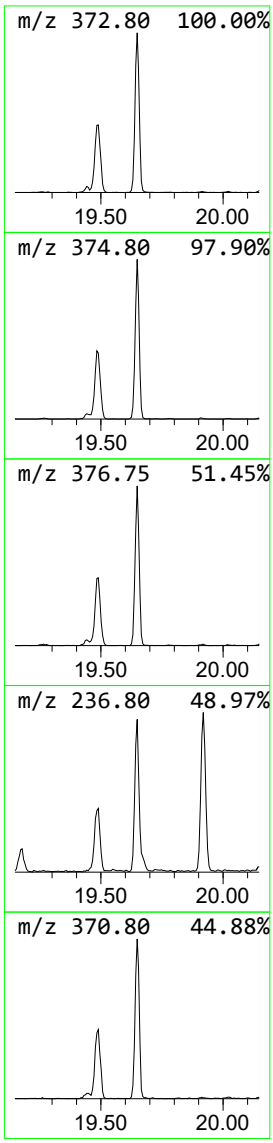
Peak Number 9 trans-Chlordane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.648	7.77 ng	1286810	Chrysene-d12	21.648

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	trans-Chlordane	406	C10H6Cl8	005103-74-2	99
2	Chlordane	406	C10H6Cl8	000057-74-9	99
3	cis-Chlordane	406	C10H6Cl8	005103-71-9	96
4	4,7-Methanoindene, 1,2,3,5,6,7,8...	406	C10H6Cl8	1000017-10-9	91
5	6,8-Dichloro-2-(1-adamantanyl)-4...	375	C20H19Cl2NO2	049647-91-8	12





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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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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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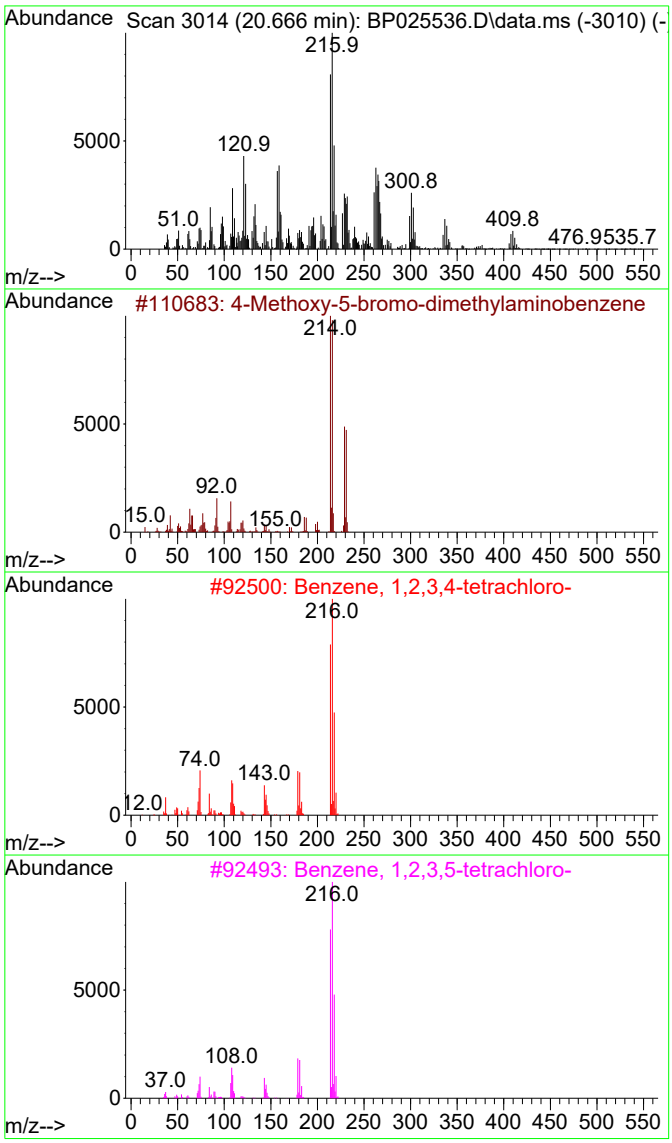
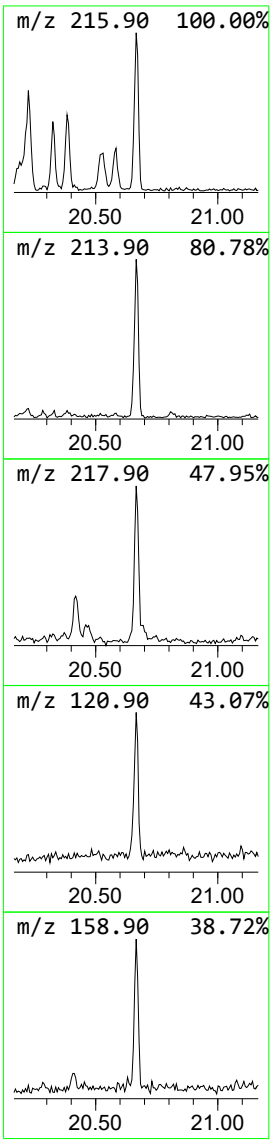
Peak Number 10 4-Methoxy-5-bromo-dimethyla... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.666	2.10 ng	348289	Chrysene-d12	21.648

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methoxy-5-bromo-dimethylaminob...	229	C9H12BrNO	1010374-81-1	50
2	Benzene, 1,2,3,4-tetrachloro-	214	C6H2Cl4	000634-66-2	45
3	Benzene, 1,2,3,5-tetrachloro-	214	C6H2Cl4	000634-90-2	45
4	Benzene, 1,2,4,5-tetrachloro-	214	C6H2Cl4	000095-94-3	45
5	2,4-Dichloro-6-n-butoxy-1,5-naph...	270	C12H12Cl2N2O	1010213-83-1	38





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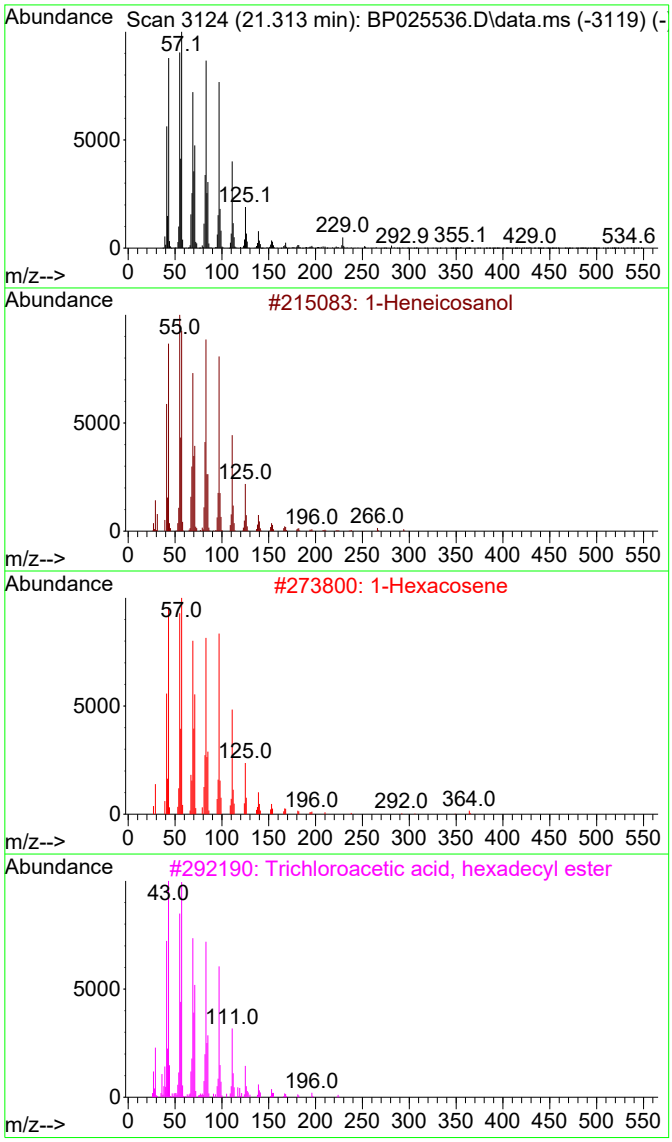
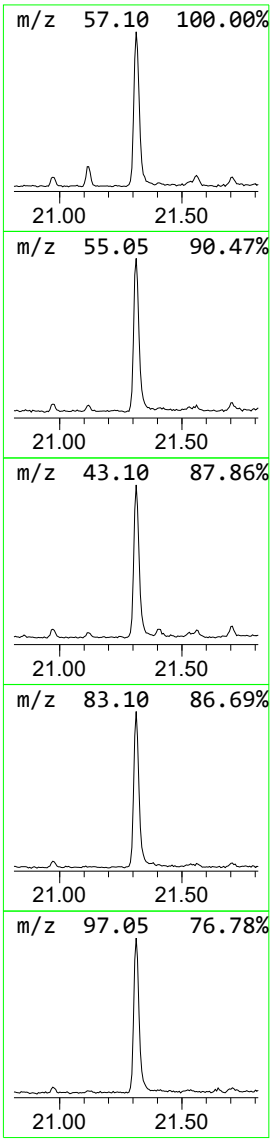
Peak Number 11 1-Heneicosanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.313	9.69 ng	1603930	Chrysene-d12	21.648

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	1-Hexacosene	364	C26H52	018835-33-1	95
3	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
4	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
5	Pentacos-1-ene	350	C25H50	016980-85-1	94





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Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.M
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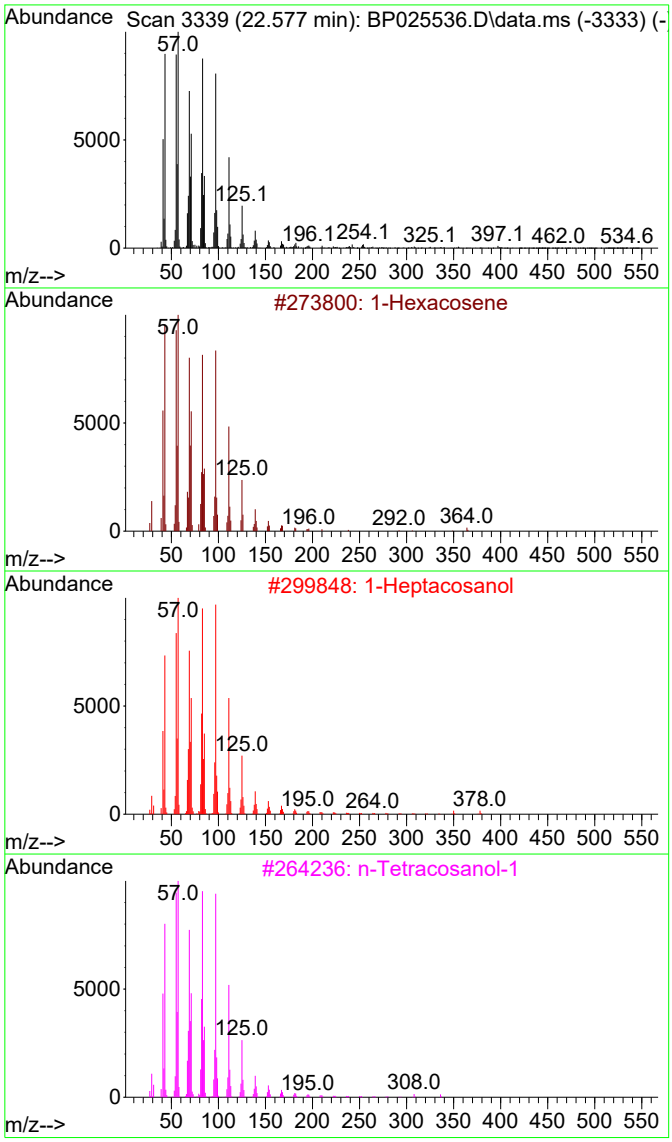
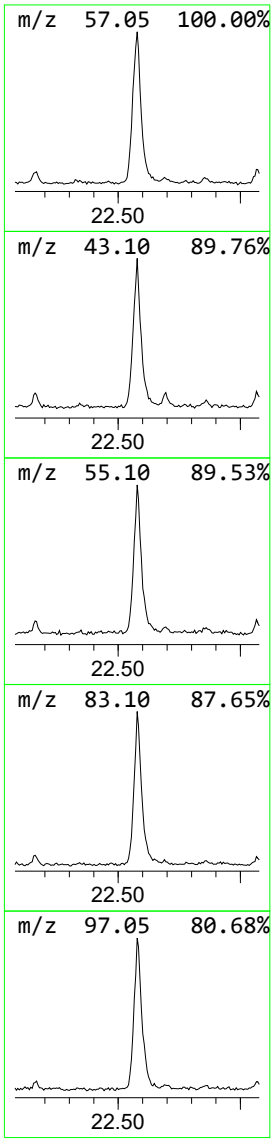
Peak Number 12 1-Hexacosene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.577	8.25 ng	1365910	Chrysene-d12	21.648

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene	364	C26H52	018835-33-1	95
2	1-Heptacosanol	396	C27H56O	002004-39-9	95
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4	Octacosanol	410	C28H58O	000557-61-9	94
5	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94





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Operator : CG/JU
Sample : Q2819-12
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
17M-S

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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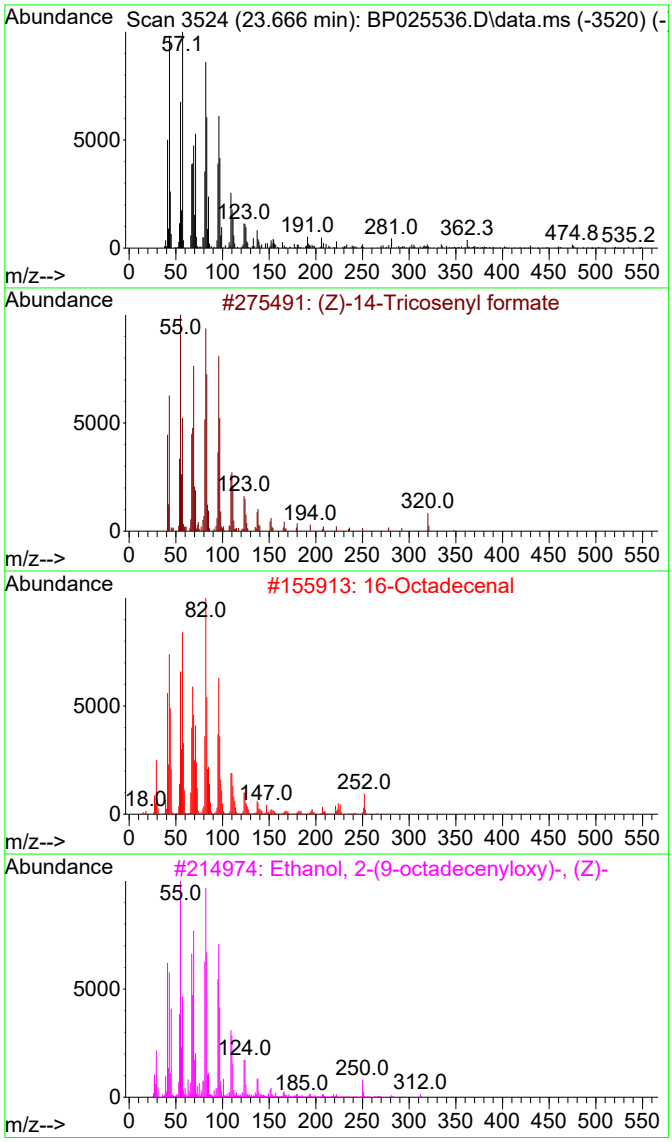
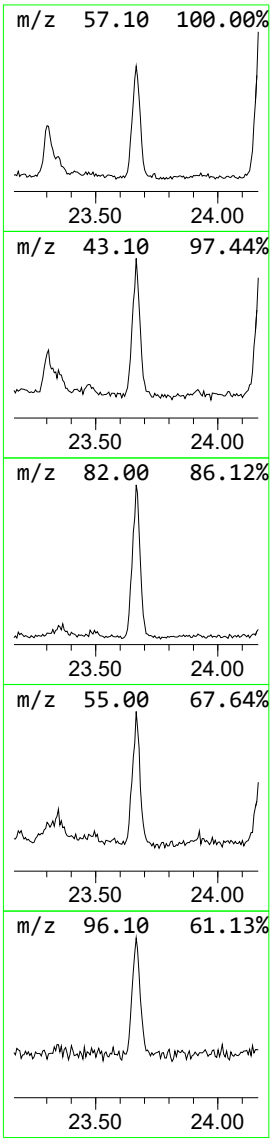
Peak Number 13 (Z)-14-Tricosenyl formate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.666	4.03 ng	620764	Perylene-d12	25.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	95
2	16-Octadecenal	266	C18H34O	056554-87-1	94
3	Ethanol, 2-(9-octadecenyloxy)-, ...	312	C20H40O2	005353-25-3	93
4	Eicosen-1-ol, cis-9-	296	C20H40O	112248-30-3	92
5	Dotriacontanal	464	C32H64O	057878-00-9	91





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TIC Library : C:\Database\NIST20.L
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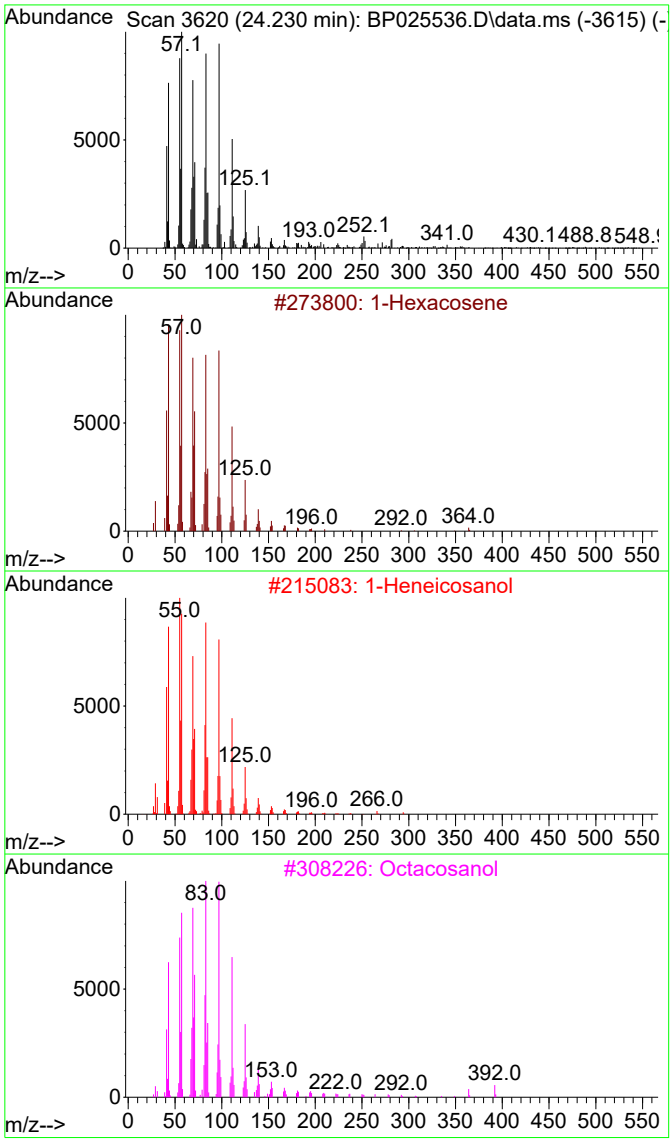
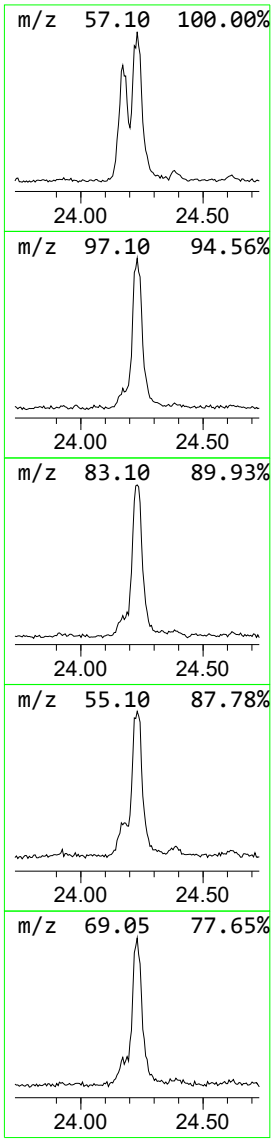
Peak Number 14 Octacosanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.230	8.51 ng	1309120	Perylene-d12	25.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene	364	C26H52	018835-33-1	96
2	1-Heneicosanol	312	C21H44O	015594-90-8	95
3	Octacosanol	410	C28H58O	000557-61-9	94
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5	n-Nonadecanol-1	284	C19H40O	001454-84-8	94





Tentatively Identified Compound (LSC) summary

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ClientSampleId :
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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...	3.037	20.4	ng	1322370	1	7.784	1293410	20.0
2-Pentanone, 4-...	4.943	5.4	ng	347210	1	7.784	1293410	20.0
Benzophenone	15.784	8.3	ng	998003	3	14.401	2414250	20.0
n-Hexadecanoic ...	18.136	8.8	ng	1242910	4	17.201	2832100	20.0
1,6-Methano-1H-...	18.525	24.4	ng	3454900	4	17.201	2832100	20.0
Heptachlor epoxide	19.177	2.1	ng	291668	4	17.201	2832100	20.0
Naphthalene, 1,...	19.266	2.2	ng	308053	4	17.201	2832100	20.0
Octadecanoic acid	19.466	3.6	ng	589963	5	21.648	3310500	20.0
trans-Chlordane	19.648	7.8	ng	1286810	5	21.648	3310500	20.0
4-Methoxy-5-bro...	20.666	2.1	ng	348289	5	21.648	3310500	20.0
1-Heneicosanol	21.313	9.7	ng	1603930	5	21.648	3310500	20.0
1-Hexacosene	22.577	8.3	ng	1365910	5	21.648	3310500	20.0
(Z)-14-Tricosen...	23.666	4.0	ng	620764	6	25.018	3077570	20.0
Octacosanol	24.230	8.5	ng	1309120	6	25.018	3077570	20.0