LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081721\
Data File : BF125082.D
Acq On : 17 Aug 2021 22:46
Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

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_F\METHODS\8270-BF081721.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF125082.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.210	15	21	27	rBV	1492240	1871832	46.39%	5.079%
2	3.434	223	229	236	rBV	25530	47851	1.19%	0.130%
3	5.145	509	520	526	rBV	159647	204045	5.06%	0.554%
4	5.534	580	586	589	rBV	3908009	3883988	96.26%	10.540%
5	6.534	751	756	759	rBV	3489601	4034912	100.00%	10.949%
6	6.916	818	821	825	rBV	1234469	1070324	26.53%	2.904%
7	6.975	828	831	834	rBV	60276	47289	1.17%	0.128%
8	7.481	912	917	920	rBV	2838927	2549789	63.19%	6.919%
9	7.816	971	974	978	rVB	58579	44501	1.10%	0.121%
10	8.098	1018	1022	1025	rBV	963351	705727	17.49%	1.915%
11	8.204	1036	1040	1043	rBV	1634633	1431444	35.48%	3.884%
12	9.281	1218	1223	1226	rBV	4469282	4018954	99.60%	10.906%
13	9.957	1334	1338	1341	rBV	1837157	1476830	36.60%	4.008%
14	10.745	1468	1472	1475	rBV	3047042	2704295	67.02%	7.338%
15	11.445	1587	1591	1593	rBV	1807872	1505411	37.31%	4.085%
16	11.463	1593	1594	1597	rVB	110840	86585	2.15%	0.235%
17	11.833	1654	1657	1660	rBV	74187	55154	1.37%	0.150%
18	11.963	1674	1679	1685	rBV	653929	681332	16.89%	1.849%
19	12.657	1794	1797	1799	rBV	219246	199775	4.95%	0.542%
20	12.751	1809	1813	1821	rBV	403372	400875	9.94%	1.088%
21	12.857	1827	1831	1832	rBV3	54728	52723	1.31%	0.143%
22	12.886	1833	1836	1838	rVV	136833	117341	2.91%	0.318%
23	13.033	1856	1861	1864	rVB	3922708	3592383	89.03%	9.748%
24	13.210	1884	1891	1893	rBV3	31739	58642	1.45%	0.159%
25	13.833	1986	1997	1999	rBV5	43489	110616	2.74%	0.300%
26	13.921	2009	2012	2019	rBV	528977	488219	12.10%	1.325%
27	14.063	2032	2036	2038	rBV	1823454	1759509	43.61%	4.775%
28	14.080	2038	2039	2042	rVV	1060584	808315	20.03%	2.193%
29	14.104	2042	2043	2046	rVB	82364	66123	1.64%	0.179%
30	14.557	2117	2120	2123	rBV	169249	150266	3.72%	0.408%
31	14.768	2153	2156	2161	rBV6	35927	55766	1.38%	0.151%
32	14.951	2184	2187	2193	rVB6	37889	53143	1.32%	0.144%
33	15.021	2196	2199	2202	rVB2	49912	52102	1.29%	0.141%
34	15.133	2214	2218	2221	rBV	137725	194038	4.81%	0.527%
35	15.221	2230	2233	2234	rBV	60149	60906	1.51%	0.165%
36	15.245	2234	2237	2241	rVB	145459	166756	4.13%	0.453%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081721\
Data File : BF125082.D
Acq On : 17 Aug 2021 22:46
Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

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_F\METHODS\8270-BF081721.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	15.445	2268	2271	2278	rVB	92486	114120	2.83%	0.310%
38	15.510	2278	2282	2288	rBV	88852	118243	2.93%	0.321%
39	15.580	2289	2294	2298	rBV2	786773	972676	24.11%	2.639%
40	16.068	2373	2377	2381	rBV	76636	117770	2.92%	0.320%
41	16.115	2381	2385	2396	rVB	103695	195116	4.84%	0.529%
42	16.898	2514	2518	2524	rVB9	38102	77818	1.93%	0.211%
43	17.074	2546	2548	2560	rVB2	36864	92699	2.30%	0.252%
44	17.292	2578	2585	2594	rBV2	46686	162790	4.03%	0.442%
45	17.462	2610	2614	2620	rVB9	23911	47454	1.18%	0.129%
46	17.703	2651	2655	2665	rVB8	69259	144933	3.59%	0.393%

Sum of corrected areas: 36851380

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081721\
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Acq On : 17 Aug 2021 22:46
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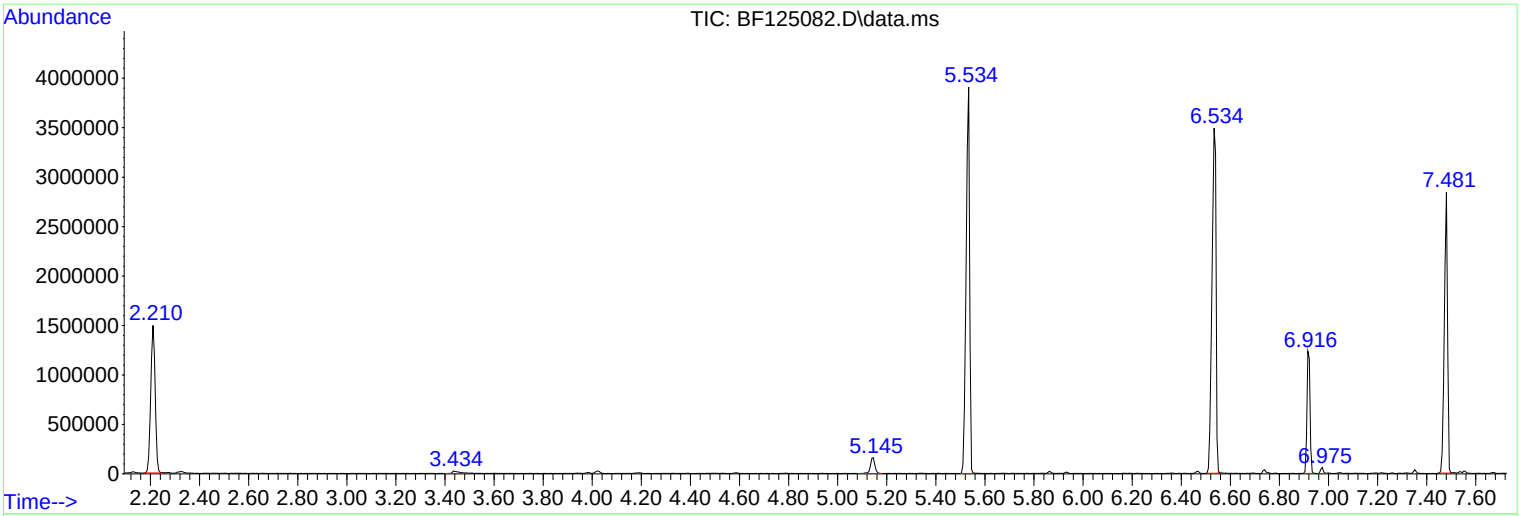
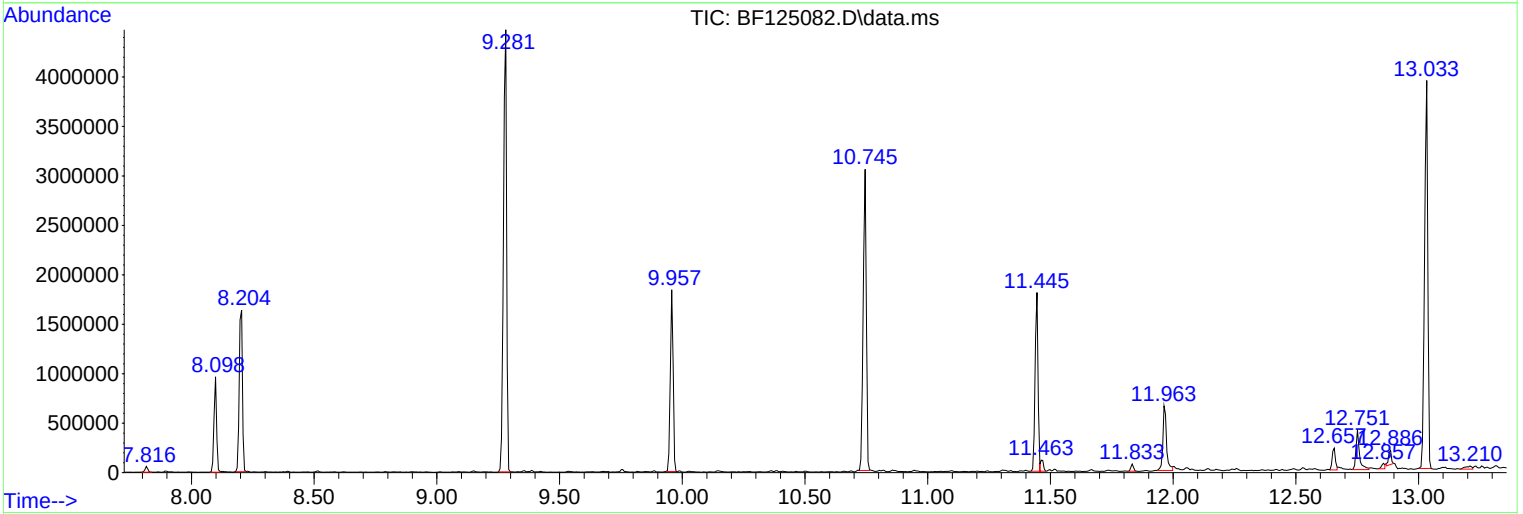
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

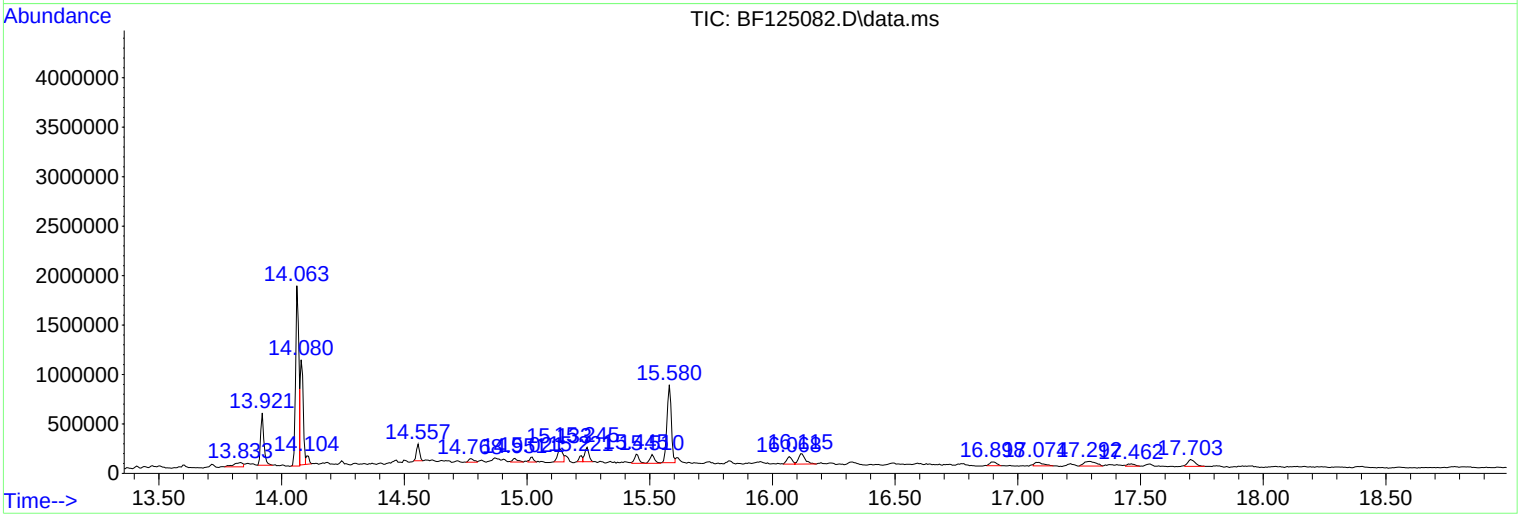
Instrument :
BNA_F
ClientSampleId :
340

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.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_F\Data\BF081721\
Data File : BF125082.D
Acq On : 17 Aug 2021 22:46
Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

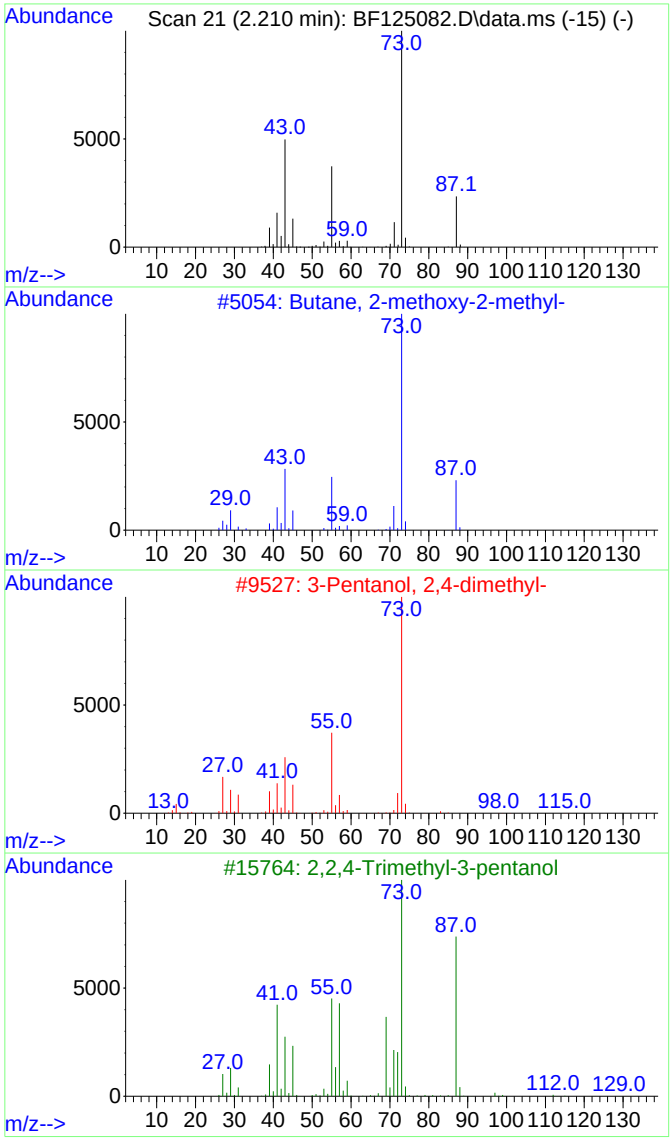
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

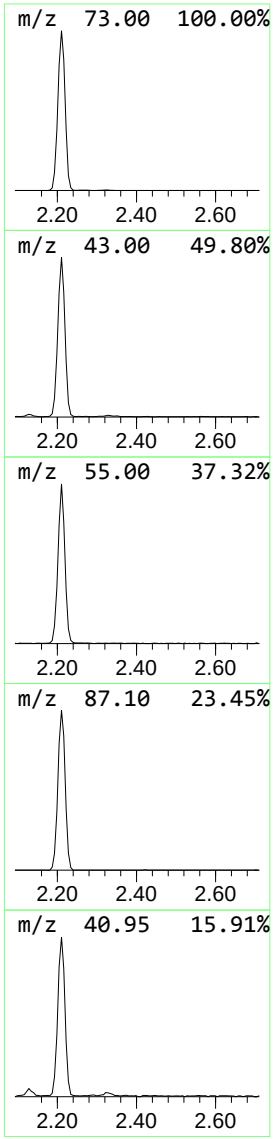
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.210	34.98 ng	1871830	1,4-Dichlorobenzene-d4	6.922

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	43
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081721\
Data File : BF125082.D
Acq On : 17 Aug 2021 22:46
Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

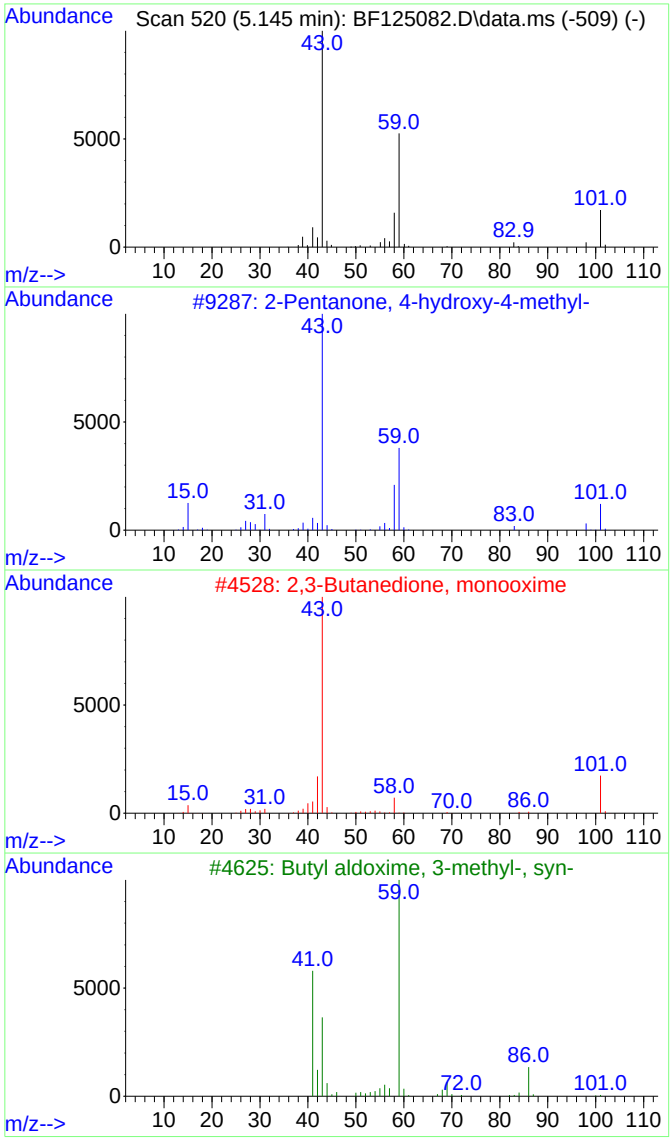
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

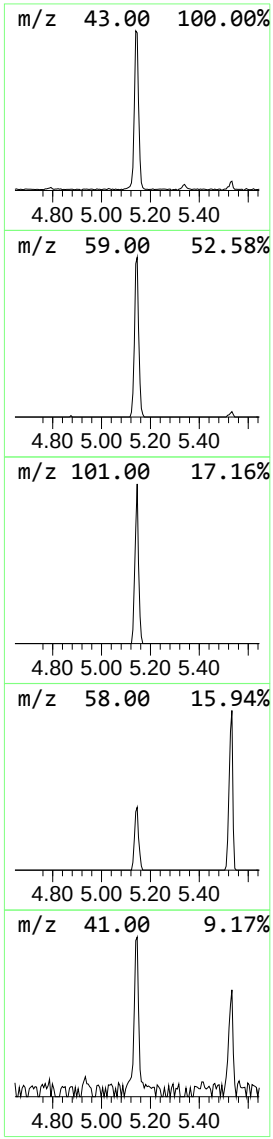
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.145	3.81 ng	204045	1,4-Dichlorobenzene-d4	6.922

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
3	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081721\
Data File : BF125082.D
Acq On : 17 Aug 2021 22:46
Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

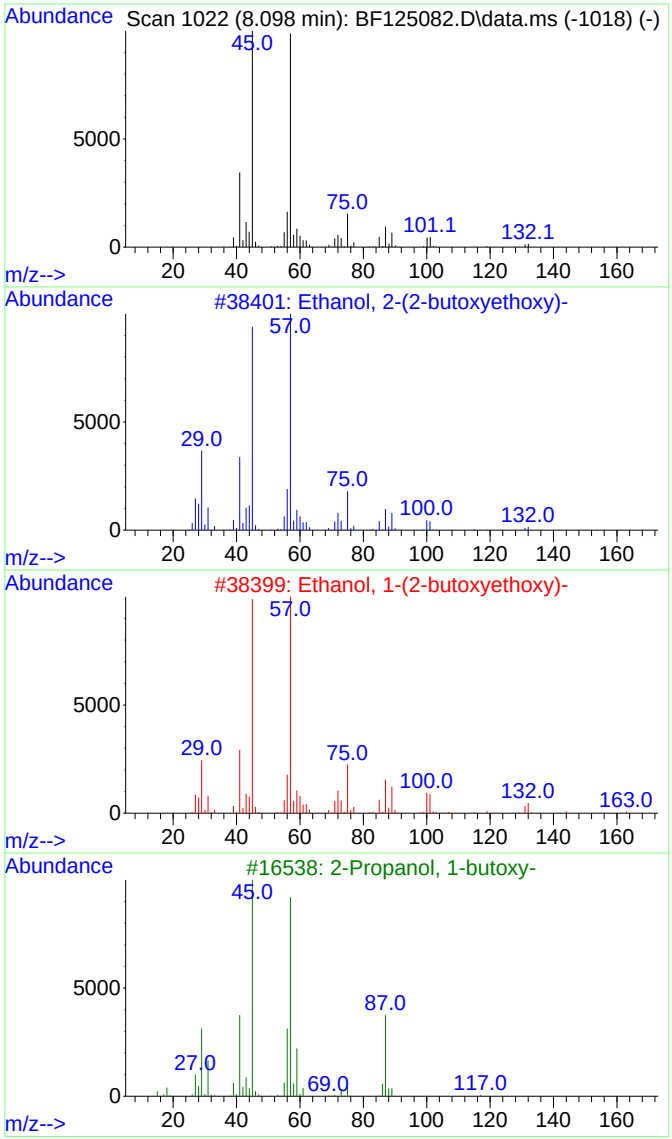
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

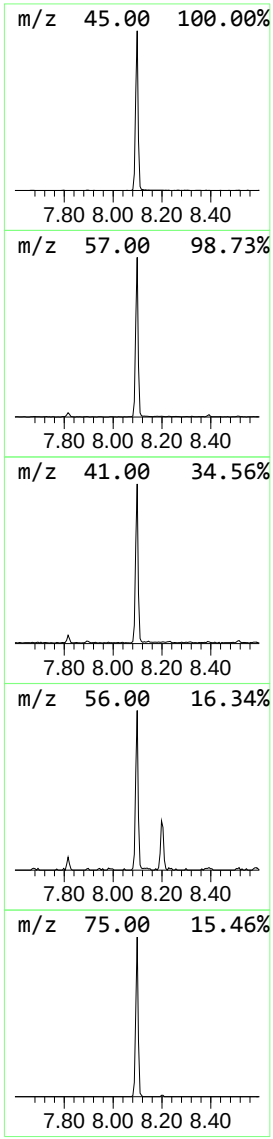
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.098	9.86 ng	705727	Naphthalene-d8	8.204

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3	2-Propanol, 1-butoxy-	132	C7H16O2	005131-66-8	59
4	3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	53
5	Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081721\
Data File : BF125082.D
Acq On : 17 Aug 2021 22:46
Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

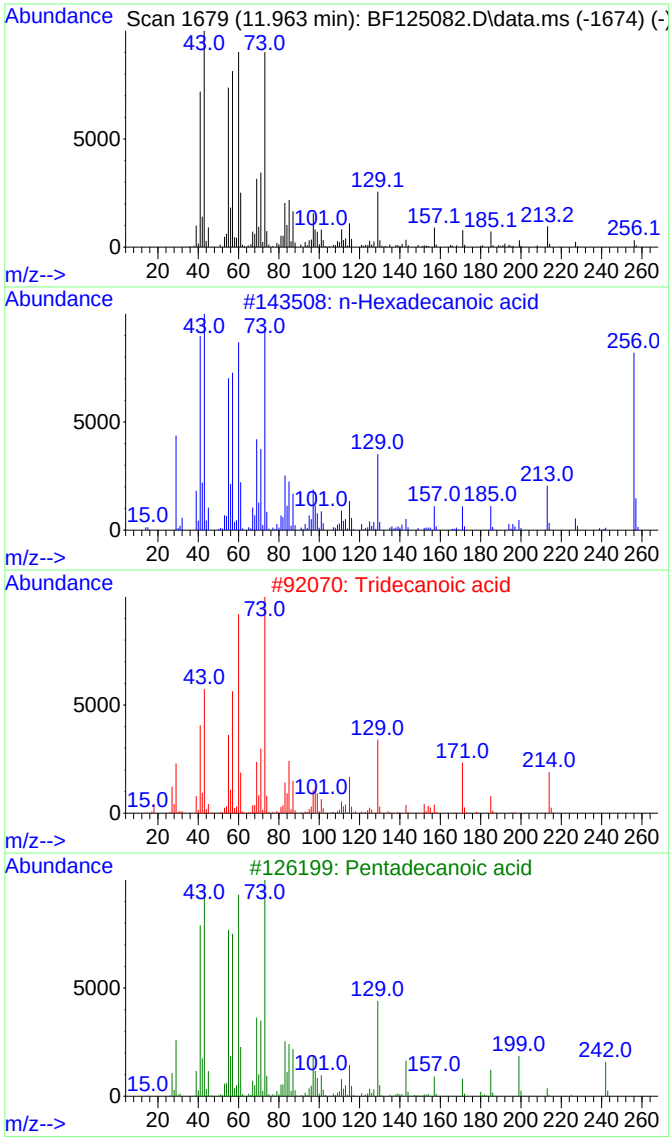
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

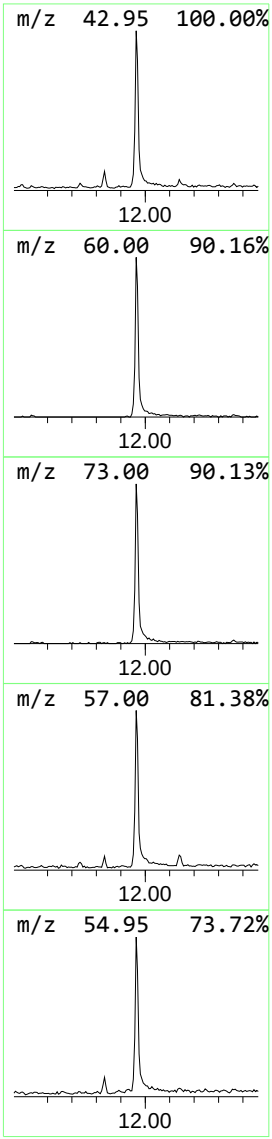
****

Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.963	9.05 ng	681332	Phenanthrene-d10	11.445

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tridecanoic acid	214	C13H26O2	000638-53-9	93
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4	8-Methylnonanoic acid	172	C10H20O2	005963-14-4	64
5	Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	15





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081721\
Data File : BF125082.D
Acq On : 17 Aug 2021 22:46
Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

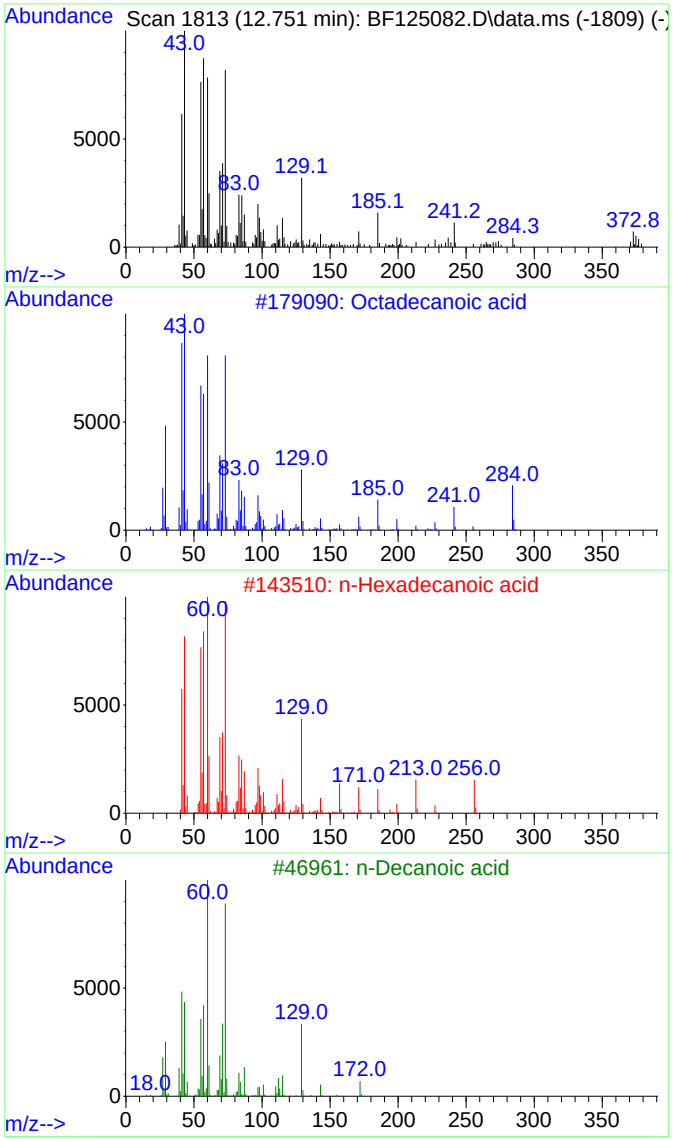
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

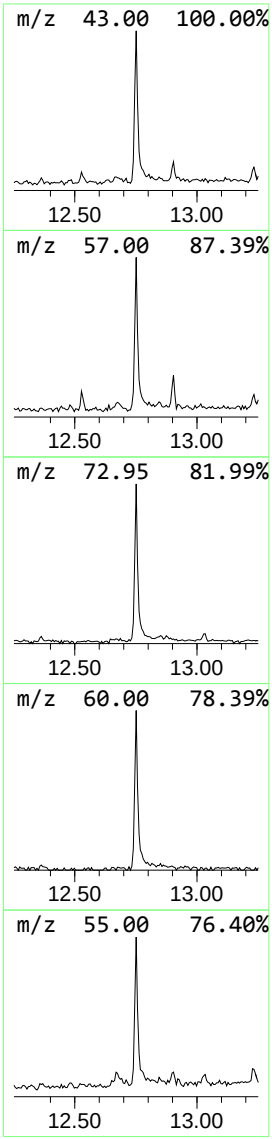
****

Peak Number 5 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.751	5.33 ng	400875	Phenanthrene-d10	11.445

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	97
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
3	n-Decanoic acid	172	C10H20O2	000334-48-5	68
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
5	Undecanoic acid	186	C11H22O2	000112-37-8	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081721\
Data File : BF125082.D
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Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

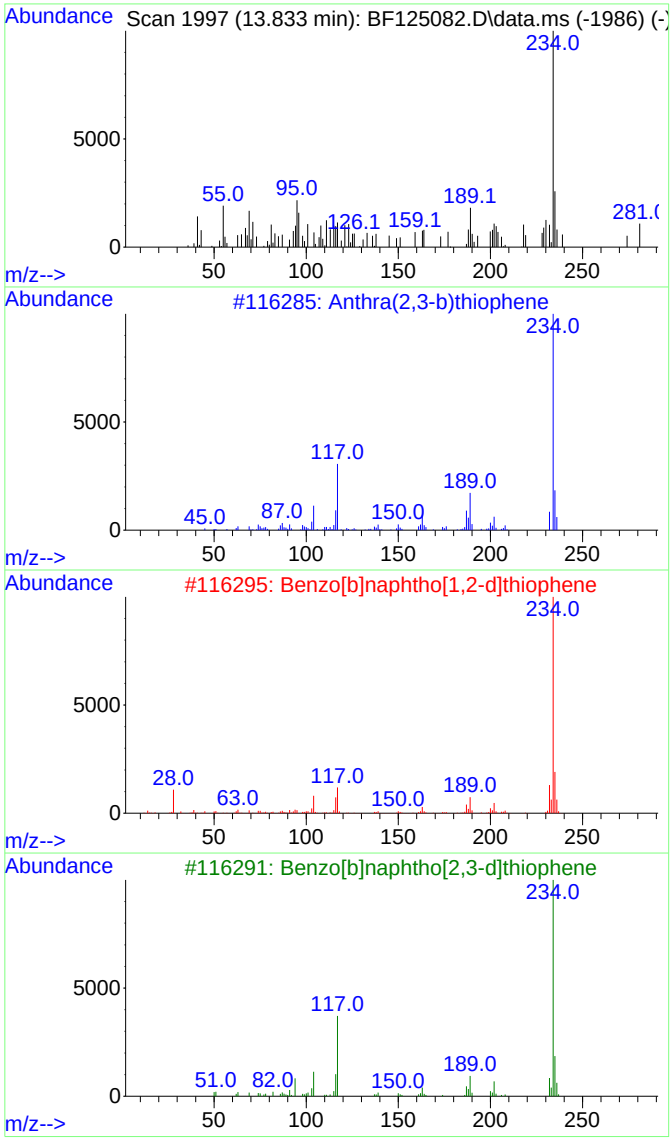
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

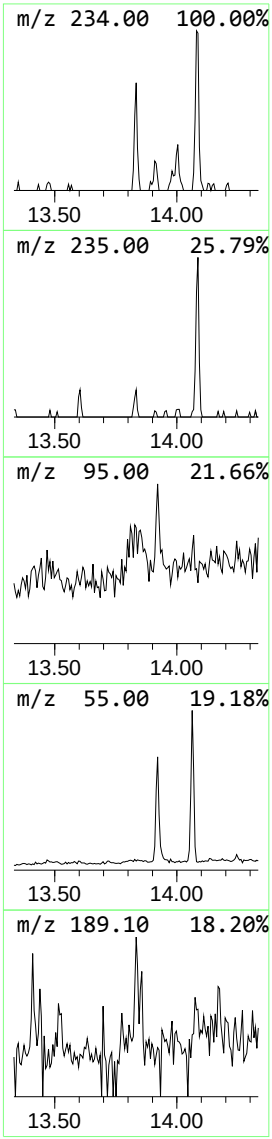
****

Peak Number 6 Anthra(2,3-b)thiophene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.833	2.74 ng	110616	Chrysene-d12	14.080

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthra(2,3-b)thiophene	234	C16H10S	022108-55-0	60
2	Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	55
3	Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	53
4	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	52
5	Quinoxalino[2,3-b]quinoxaline, 5...	234	C14H10N4	000531-46-4	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081721\
Data File : BF125082.D
Acq On : 17 Aug 2021 22:46
Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

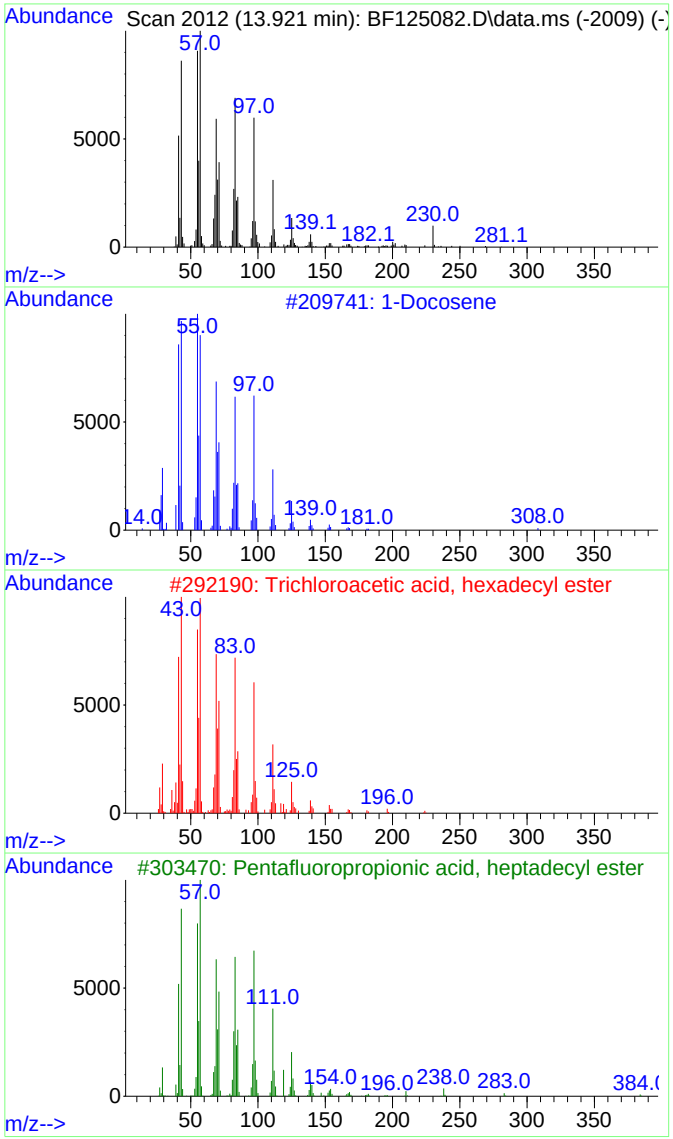
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

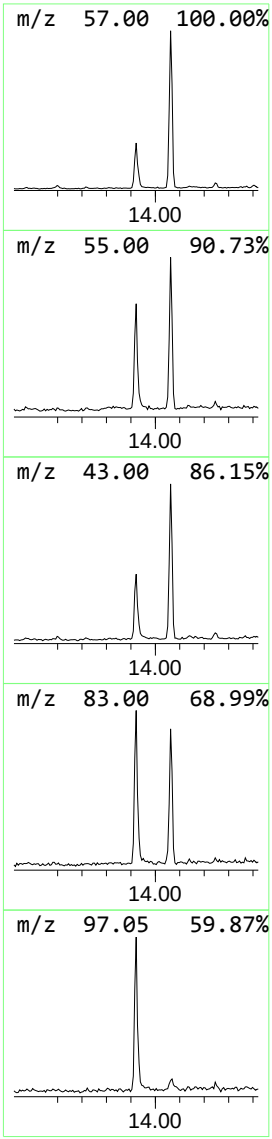
****

Peak Number 7 1-Docosene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.921	12.08 ng	488219	Chrysene-d12	14.080

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	94
2	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	91
4	1-Hexacosene	364	C26H52	018835-33-1	91
5	Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081721\
Data File : BF125082.D
Acq On : 17 Aug 2021 22:46
Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

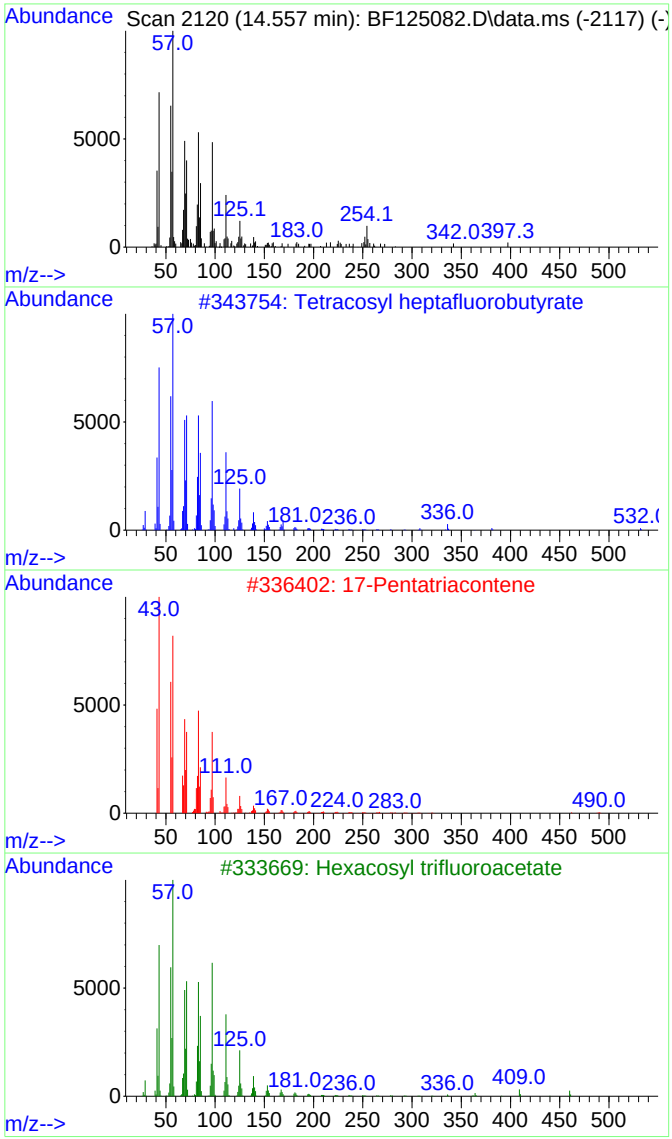
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

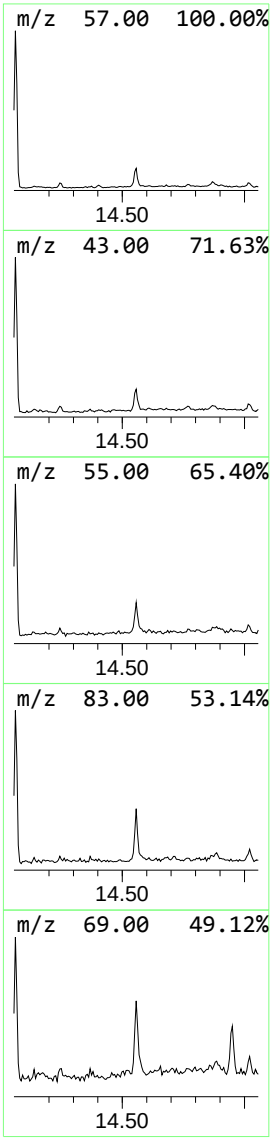
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Tetracosyl heptafluorobutyrate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.557	3.72 ng	150266	Chrysene-d12	14.080

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosyl heptafluorobutyrate	550	C28H49F7O2	1000351-83-7	87
2	17-Pentatriacontene	491	C35H70	006971-40-0	87
3	Hexacosyl trifluoroacetate	478	C28H53F3O2	1000351-75-0	87
4	1-Hexacosene	364	C26H52	018835-33-1	87
5	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	86





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Acq On : 17 Aug 2021 22:46
Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.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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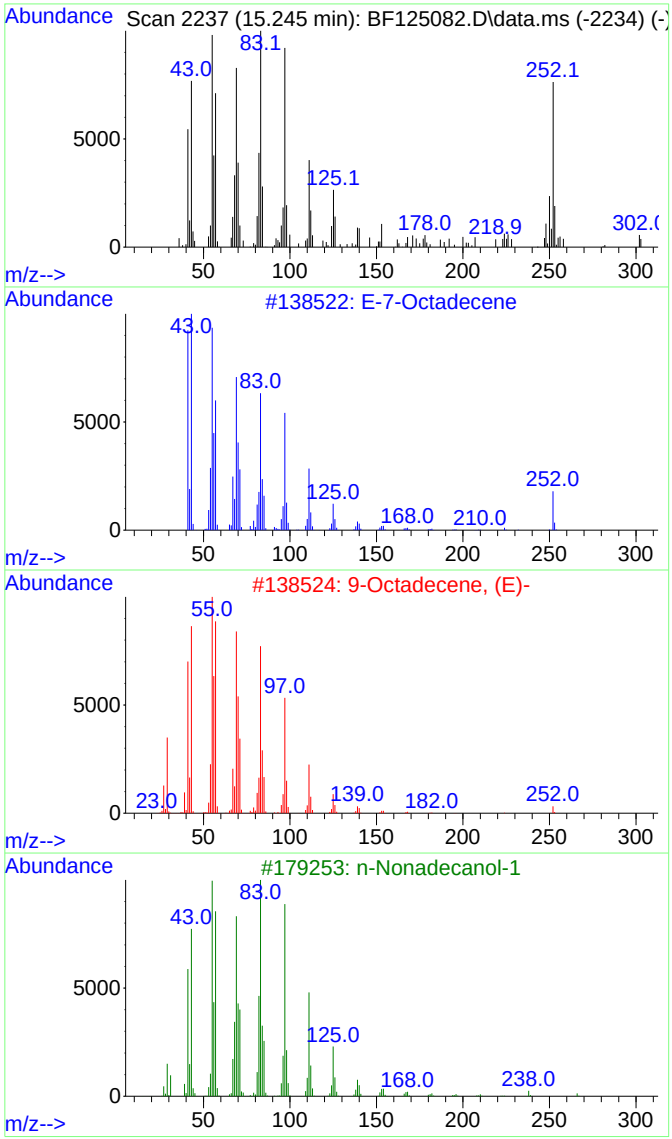
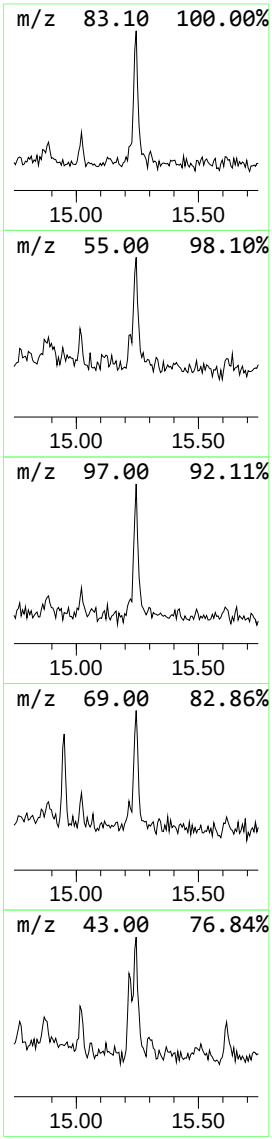
Peak Number 9 E-7-Octadecene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.245	3.43 ng	166756	Perylene-d12	15.580

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	E-7-Octadecene	252	C18H36	1000130-92-0	94
2	9-Octadecene, (E)-	252	C18H36	007206-25-9	93
3	n-Nonadecanol-1	284	C19H40O	001454-84-8	89
4	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	76
5	Behenic alcohol	326	C22H46O	000661-19-8	76





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Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.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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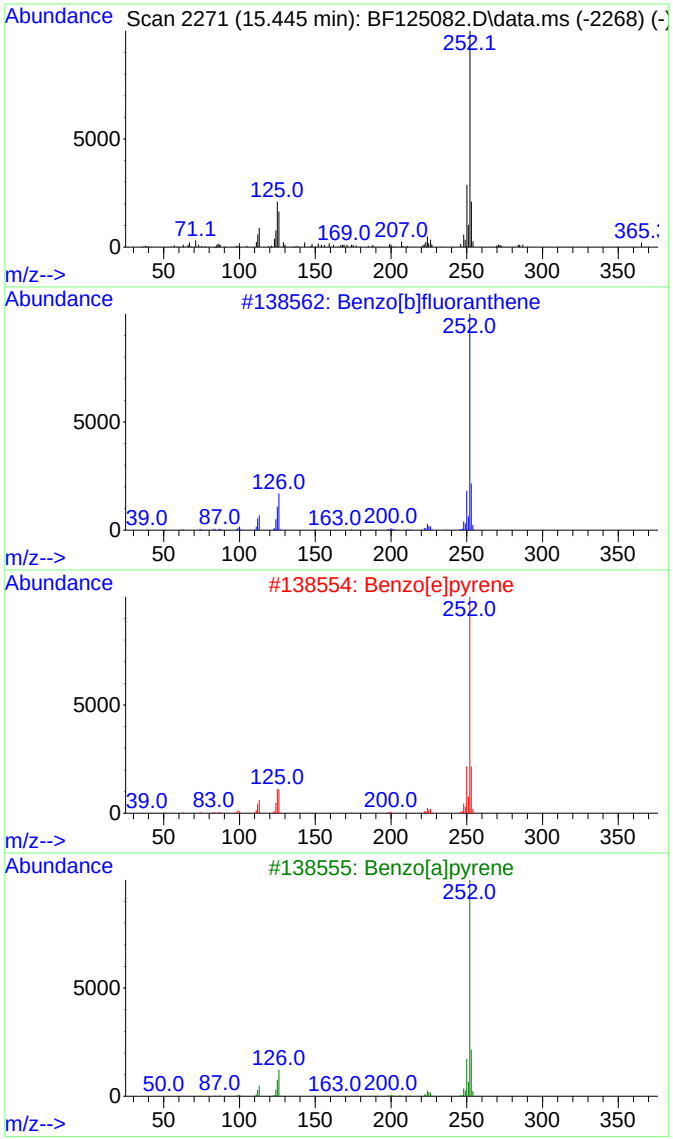
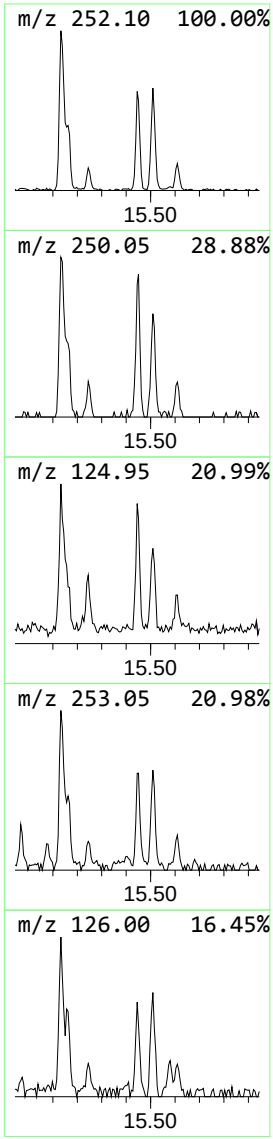
Peak Number 10 Benzo[e]pyrene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.445	2.35 ng	114120	Perylene-d12	15.580

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]fluoranthene	252	C20H12	000205-99-2	96
2	Benzo[e]pyrene	252	C20H12	000192-97-2	95
3	Benzo[a]pyrene	252	C20H12	000050-32-8	94
4	Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
5	Benzo[j]fluoranthene	252	C20H12	000205-82-3	90





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Acq On : 17 Aug 2021 22:46
Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.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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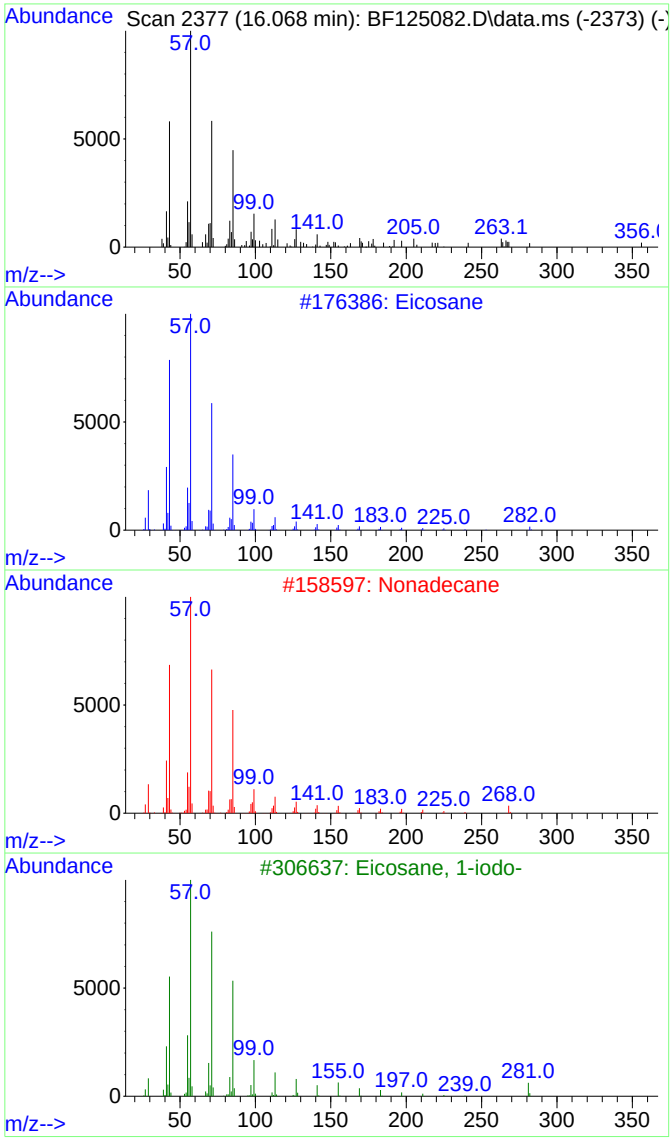
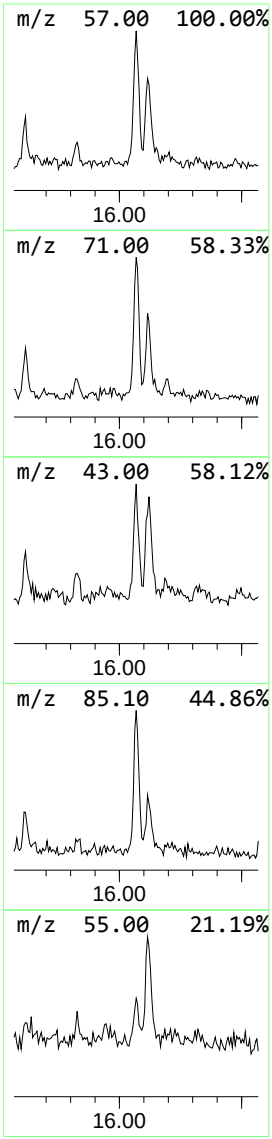
Peak Number 12 Eicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.068	2.42 ng	117770	Perylene-d12	15.580

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	94
2	Nonadecane	268	C19H40	000629-92-5	91
3	Eicosane, 1-iodo-	408	C20H41I	1000406-31-8	87
4	4-Methylheneicosane	310	C22H46	025117-29-7	87
5	Docosane, 1-iodo-	436	C22H45I	1000406-31-9	83





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Acq On : 17 Aug 2021 22:46
Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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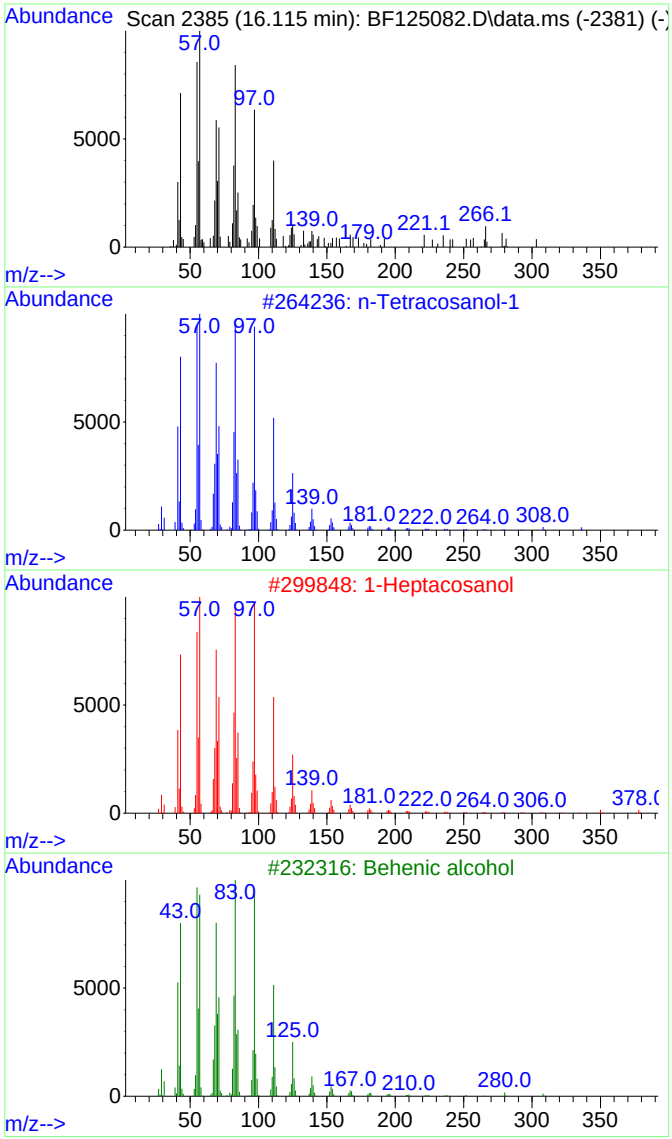
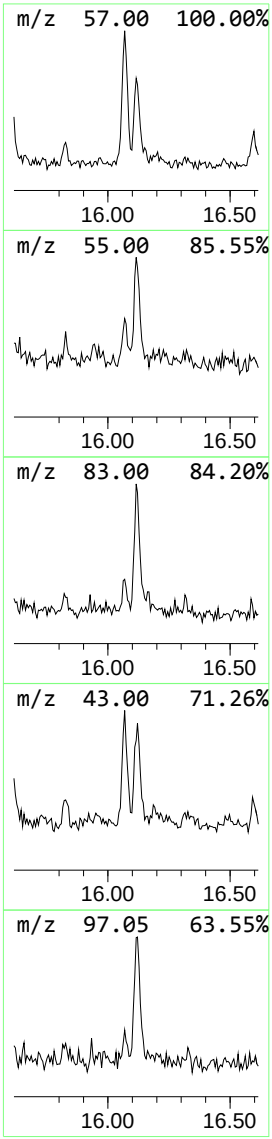
Peak Number 13 n-Tetracosanol-1 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.115	4.01 ng	195116	Perylene-d12	15.580

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Tetracosanol-1	354	C24H50O	000506-51-4	81
2	1-Heptacosanol	396	C27H56O	002004-39-9	81
3	Behenic alcohol	326	C22H46O	000661-19-8	76
4	17-Pentatriacontene	491	C35H70	006971-40-0	74
5	1-Tetracosene	336	C24H48	010192-32-2	62





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Acq On : 17 Aug 2021 22:46
Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.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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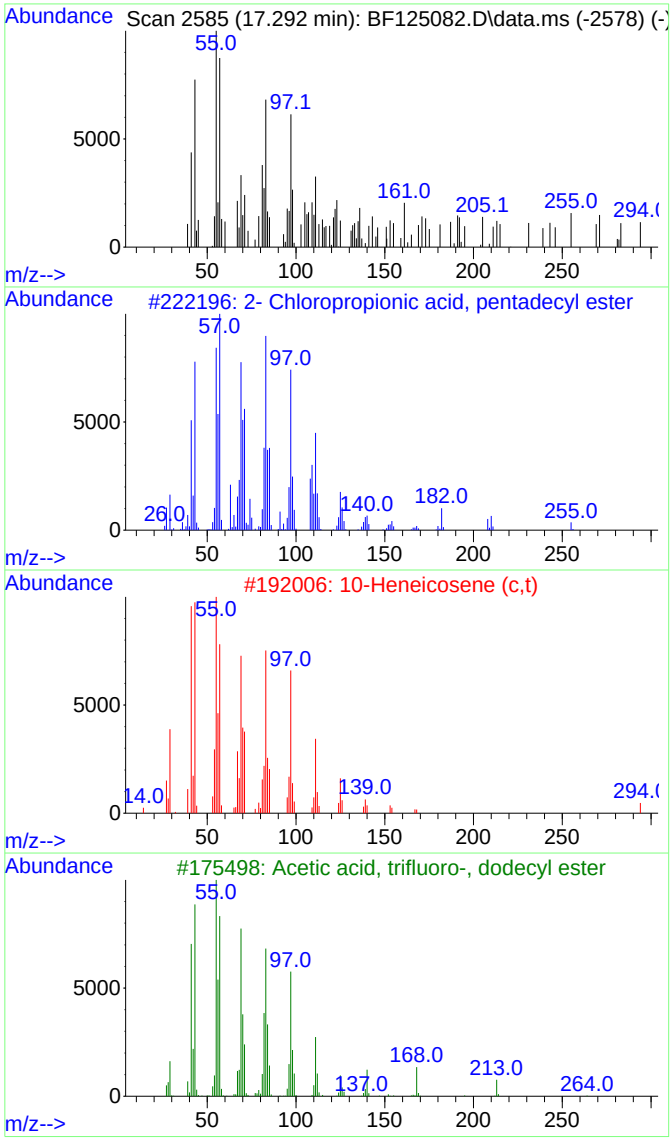
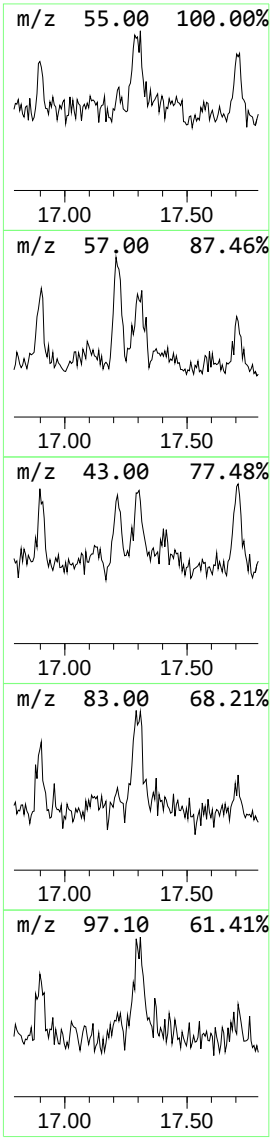
Peak Number 14 unknown17.292 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.292	3.35 ng	162790	Perylene-d12	15.580

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2- Chloropropionic acid, pentade...	318	C18H35ClO2	1000292-44-1	49
2	10-Heneicosene (c,t)	294	C21H42	095008-11-0	46
3	Acetic acid, trifluoro-, dodecyl...	282	C14H25F3O2	006222-00-0	46
4	Cis-1,3-di(n-hexyl)cyclohexane	252	C18H36	086701-17-9	27
5	Cyclohexane, methyl-	98	C7H14	000108-87-2	18





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Acq On : 17 Aug 2021 22:46
Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.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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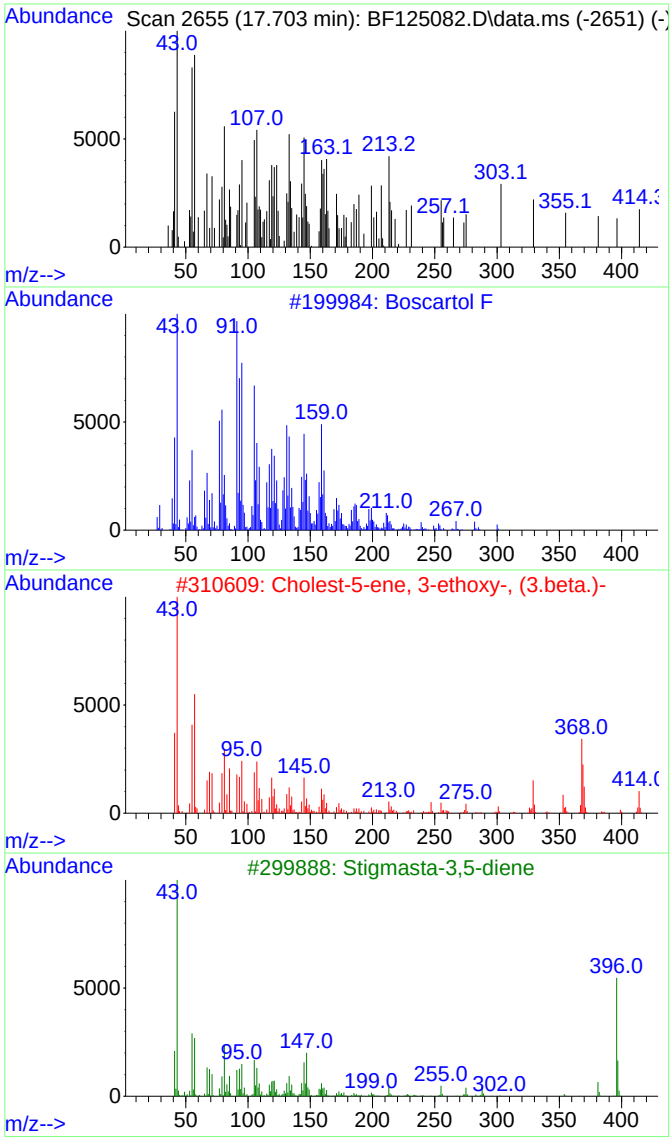
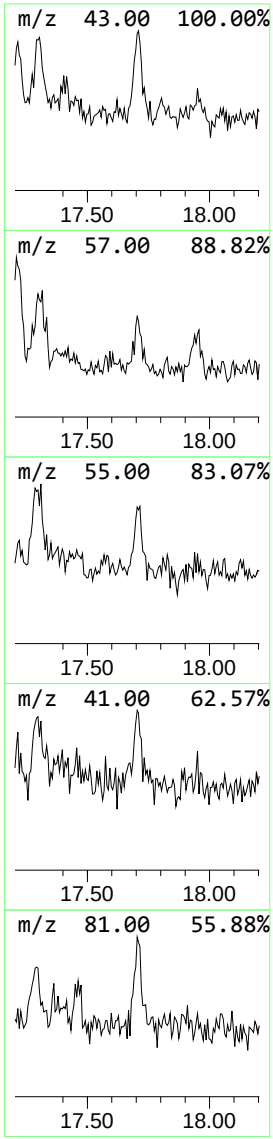
Peak Number 15 unknown17.703 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.703	2.98 ng	144933	Perylene-d12	15.580

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Boscartol F	300	C20H28O2	1486443-17-7	25
2	Cholest-5-ene, 3-ethoxy-, (3.bet...	414	C29H50O	000986-19-6	14
3	Stigmasta-3,5-diene	396	C29H48	004970-37-0	14
4	Ledene oxide-(II)	220	C15H24O	1000159-36-7	12
5	4-Tolylselenourea	214	C8H10N2Se	016518-82-4	10





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081721\
Data File : BF125082.D
Acq On : 17 Aug 2021 22:46
Operator : JU/CG
Sample : M3421-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 340

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.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...	2.210	35.0	ng	1871830	1	6.922	1070320	20.0
2-Pentanone, 4-...	5.145	3.8	ng	204045	1	6.922	1070320	20.0
Ethanol, 2-(2-b...	8.098	9.9	ng	705727	2	8.204	1431440	20.0
n-Hexadecanoic ...	11.963	9.1	ng	681332	4	11.445	1505410	20.0
Octadecanoic acid	12.751	5.3	ng	400875	4	11.445	1505410	20.0
Anthra(2,3-b)th...	13.833	2.7	ng	110616	5	14.080	808315	20.0
1-Docosene	13.921	12.1	ng	488219	5	14.080	808315	20.0
Tetracosyl hept...	14.557	3.7	ng	150266	5	14.080	808315	20.0
E-7-Octadecene	15.245	3.4	ng	166756	6	15.580	972676	20.0
Benzo[e]pyrene	15.445	2.4	ng	114120	6	15.580	972676	20.0
Eicosane	16.068	2.4	ng	117770	6	15.580	972676	20.0
n-Tetracosanol-1	16.115	4.0	ng	195116	6	15.580	972676	20.0
unknown17.292	17.292	3.4	ng	162790	6	15.580	972676	20.0
unknown17.703	17.703	3.0	ng	144933	6	15.580	972676	20.0