LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013160.D
Acq On : 20 Dec 2022 11:22
Operator : CG/JU
Sample : N6003-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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\SFAM-EPA-BP120722.M
Title : SVOA CALIBRATION

Signal : TIC: BP013160.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.334	22	25	32	rVB	194799	249417	0.73%	0.127%
2	3.764	95	98	112	rBV	1549862	2223374	6.49%	1.136%
3	3.870	113	116	122	rVV	119378	170357	0.50%	0.087%
4	3.999	133	138	145	rVB2	62351	103104	0.30%	0.053%
5	4.093	151	154	160	rVB	39047	53265	0.16%	0.027%
6	5.040	308	315	324	rBV2	238865	364810	1.06%	0.186%
7	7.111	660	667	678	rBV	2874694	4561872	13.31%	2.331%
8	7.281	689	696	707	rBV	2425797	3709259	10.82%	1.896%
9	7.481	723	730	738	rBV	2987516	4668609	13.62%	2.386%
10	7.952	803	810	817	rBV	1716451	2718284	7.93%	1.389%
11	8.663	924	931	942	rBV	2952051	4699627	13.71%	2.402%
12	8.787	947	952	960	rVB	82926	143063	0.42%	0.073%
13	9.122	1001	1009	1017	rBV	2744351	4446855	12.97%	2.273%
14	9.528	1072	1078	1084	rVB2	57811	95187	0.28%	0.049%
15	9.846	1122	1132	1145	rBV	2245657	3820777	11.15%	1.953%
16	10.069	1163	1170	1181	rVB6	35069	71757	0.21%	0.037%
17	10.387	1217	1224	1234	rBV	3400820	5653141	16.49%	2.889%
18	10.769	1280	1289	1301	rBV	2228683	3758552	10.96%	1.921%
19	10.904	1304	1312	1326	rBV	2152333	3786857	11.05%	1.935%
20	11.557	1417	1423	1433	rBV	13459	34810	0.10%	0.018%
21	12.128	1514	1520	1524	rBV9	22514	38331	0.11%	0.020%
22	12.351	1553	1558	1564	rVB	62306	95895	0.28%	0.049%
23	13.410	1734	1738	1744	rVB	40315	61789	0.18%	0.032%
24	13.587	1760	1768	1772	rVB5	20228	45871	0.13%	0.023%
25	14.004	1831	1839	1848	rBV	4973163	6995102	20.40%	3.575%
26	14.293	1880	1888	1897	rVV	6158492	9019462	26.31%	4.610%
27	14.598	1932	1940	1947	rBV	2904039	4399001	12.83%	2.248%
28	14.792	1966	1973	1988	rBV	1895123	3029969	8.84%	1.549%
29	14.945	1994	1999	2004	rBV2	72816	107035	0.31%	0.055%
30	15.010	2004	2010	2016	rVB5	69186	123241	0.36%	0.063%
31	15.592	2101	2109	2115	rBV	7226593	10581947	30.87%	5.408%
32	15.640	2115	2117	2122	rVV3	40719	62624	0.18%	0.032%
33	15.704	2122	2128	2145	rVV	2164513	2970331	8.66%	1.518%
34	15.828	2145	2149	2156	rVB2	37593	69554	0.20%	0.036%
35	15.957	2166	2171	2180	rVB	77730	132196	0.39%	0.068%
36	16.734	2300	2303	2307	rVB5	32787	38655	0.11%	0.020%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013160.D
Acq On : 20 Dec 2022 11:22
Operator : CG/JU
Sample : N6003-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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\SFAM-EPA-BP120722.M
Title : SVOA CALIBRATION

37	16.886	2324	2329	2334	rVB2	37165	52143	0.15%	0.027%
38	17.234	2384	2388	2390	rBV3	53401	69773	0.20%	0.036%
39	17.257	2390	2392	2396	rVB3	43154	44153	0.13%	0.023%
40	17.351	2402	2408	2413	rBV	3300746	4605442	13.43%	2.354%
41	17.451	2419	2425	2433	rBV	6870303	9623457	28.07%	4.918%
42	17.534	2436	2439	2444	rVB	97836	132386	0.39%	0.068%
43	17.728	2468	2472	2485	rVB3	137462	259410	0.76%	0.133%
44	18.010	2517	2520	2524	rBV4	34309	43091	0.13%	0.022%
45	18.110	2534	2537	2542	rVB3	52326	68833	0.20%	0.035%
46	18.163	2543	2546	2551	rBV2	102077	128157	0.37%	0.065%
47	18.222	2551	2556	2566	rBV	1258894	1734681	5.06%	0.887%
48	18.557	2609	2613	2617	rBV2	54808	92557	0.27%	0.047%
49	18.633	2622	2626	2630	rBV	734309	806040	2.35%	0.412%
50	18.704	2634	2638	2643	rVV4	60894	101513	0.30%	0.052%
51	19.051	2693	2697	2702	rBV	203514	283751	0.83%	0.145%
52	19.104	2702	2706	2715	rVB5	98833	192559	0.56%	0.098%
53	19.204	2721	2723	2727	rVB2	36411	35286	0.10%	0.018%
54	19.257	2729	2732	2737	rVB	125301	168529	0.49%	0.086%
55	19.357	2738	2749	2754	rBV2	372358	1035733	3.02%	0.529%
56	19.451	2761	2765	2777	rBV	482361	654254	1.91%	0.334%
57	19.680	2799	2804	2816	rBV	9041369	12874245	37.55%	6.580%
58	20.192	2883	2891	2893	rBV3	398412	803774	2.34%	0.411%
59	21.127	3046	3050	3061	rBV	1644268	2313242	6.75%	1.182%
60	21.439	3098	3103	3108	rBV	4246220	5563011	16.23%	2.843%
61	22.045	3204	3206	3219	rVB3	334873	759501	2.22%	0.388%
62	23.133	3387	3391	3397	rBV2	820197	1388803	4.05%	0.710%
63	23.763	3491	3498	3512	rVB2	7440351	15920148	46.44%	8.136%
64	23.921	3518	3525	3533	rVB	3004355	6103436	17.80%	3.119%
65	24.539	3625	3630	3640	rVB	1012844	2134756	6.23%	1.091%
66	26.457	3951	3956	3977	rVB3	1071826	3281715	9.57%	1.677%
67	27.468	4120	4128	4149	rVB3	1882545	7010502	20.45%	3.583%
68	28.515	4294	4306	4327	rVB3	8658043	34282224	100.00%	17.520%

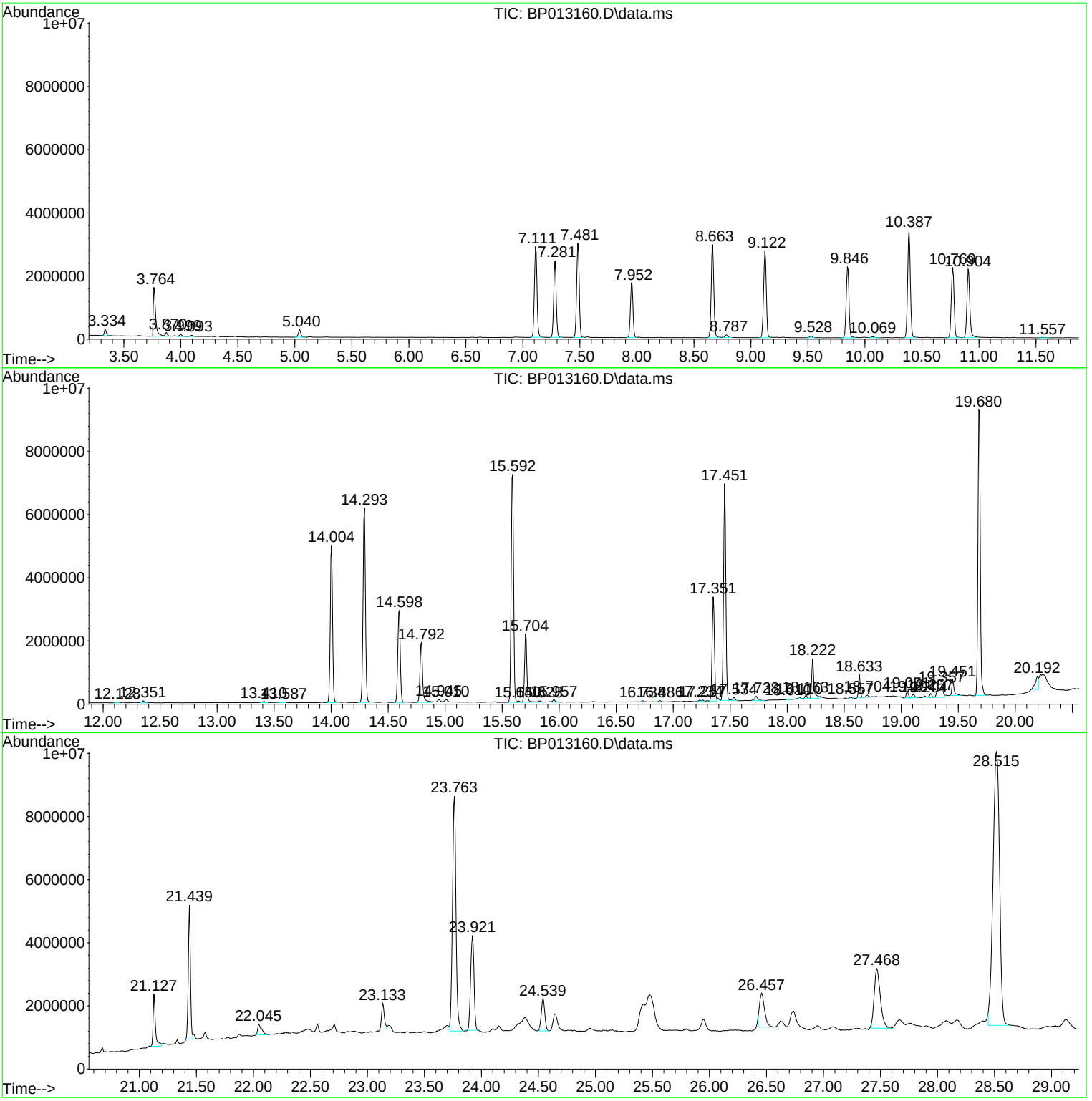
Sum of corrected areas: 195671085

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013160.D
Acq On : 20 Dec 2022 11:22
Operator : CG/JU
Sample : N6003-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013160.D
Acq On : 20 Dec 2022 11:22
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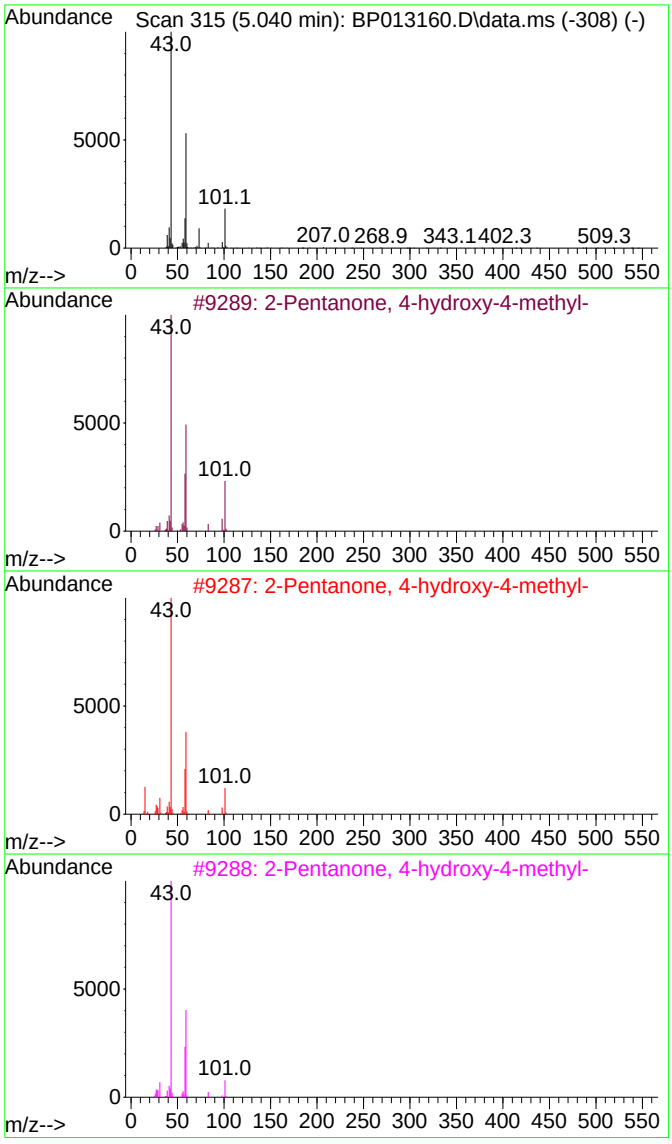
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

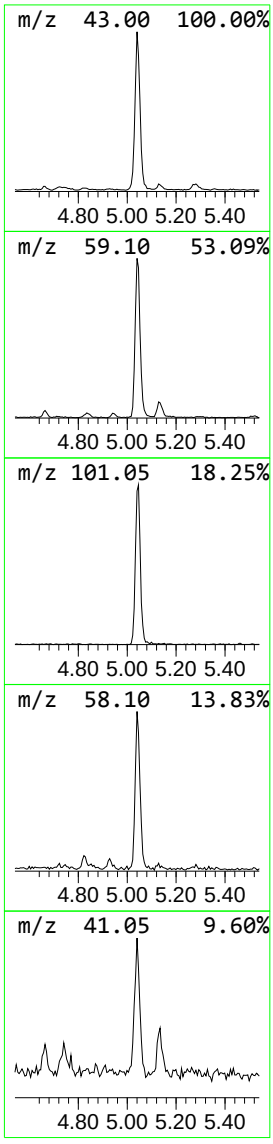
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.040	2.68 ng/ul	364810	1,4-Dichlorobenzene-d4	7.958

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
5	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013160.D
Acq On : 20 Dec 2022 11:22
Operator : CG/JU
Sample : N6003-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

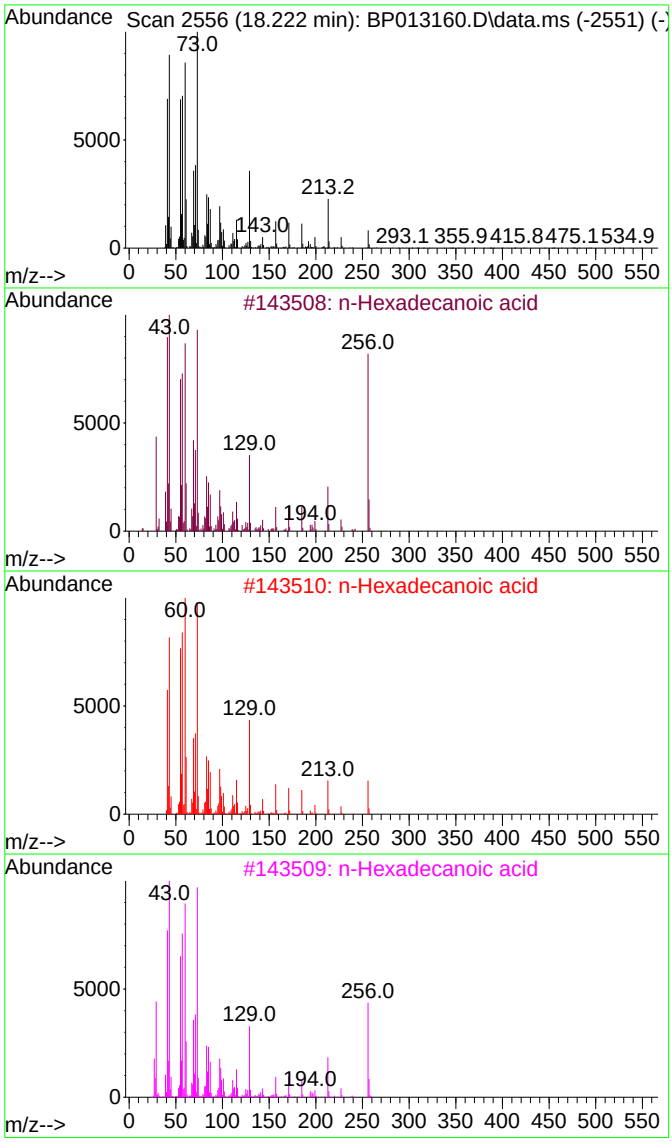
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

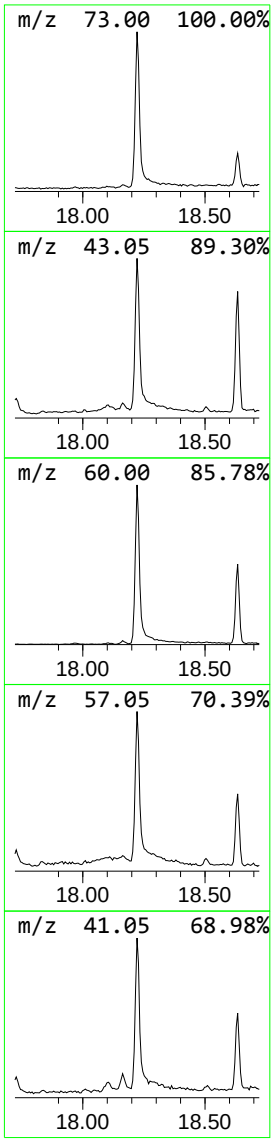
****

Peak Number 2 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.222	7.53 ng/ul	1734680	Phenanthrene-d10	17.351

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013160.D
Acq On : 20 Dec 2022 11:22
Operator : CG/JU
Sample : N6003-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

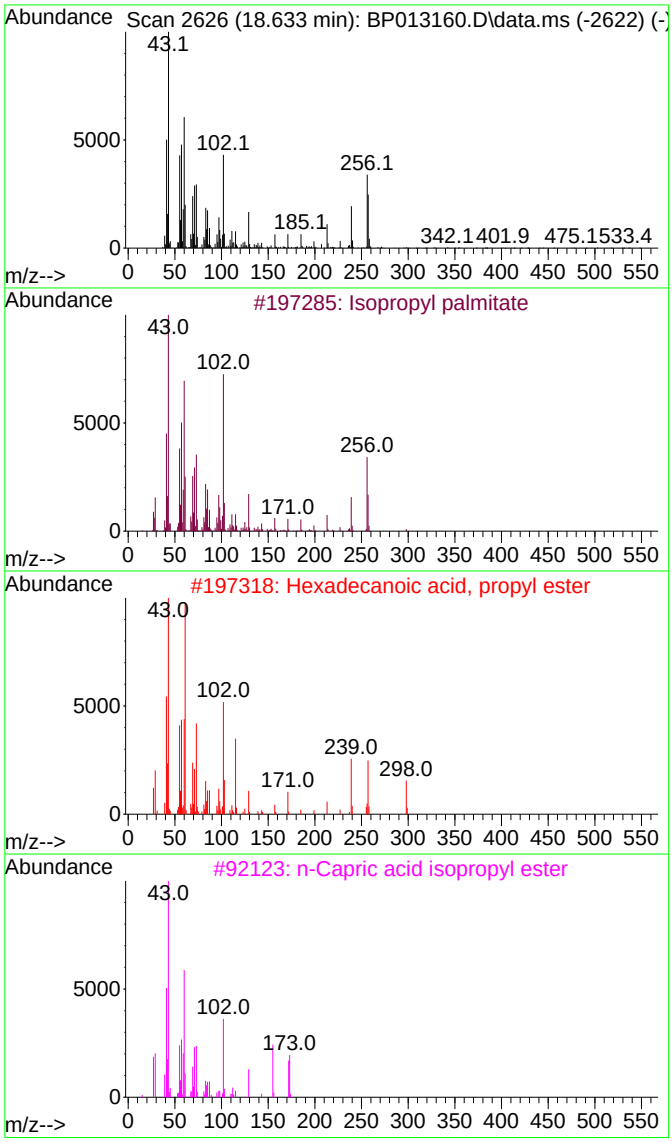
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

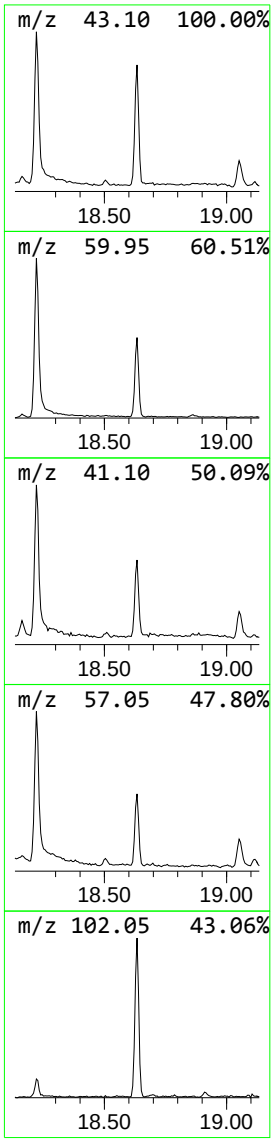
****

Peak Number 3 Isopropyl palmitate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.634	3.50 ng/ul	806040	Phenanthrene-d10	17.351

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isopropyl palmitate	298	C19H38O2	000142-91-6	72
2	Hexadecanoic acid, propyl ester	298	C19H38O2	002239-78-3	50
3	n-Capric acid isopropyl ester	214	C13H26O2	002311-59-3	43
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	42
5	Dodecanoic acid, 1-methylethyl e...	242	C15H30O2	010233-13-3	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013160.D
Acq On : 20 Dec 2022 11:22
Operator : CG/JU
Sample : N6003-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

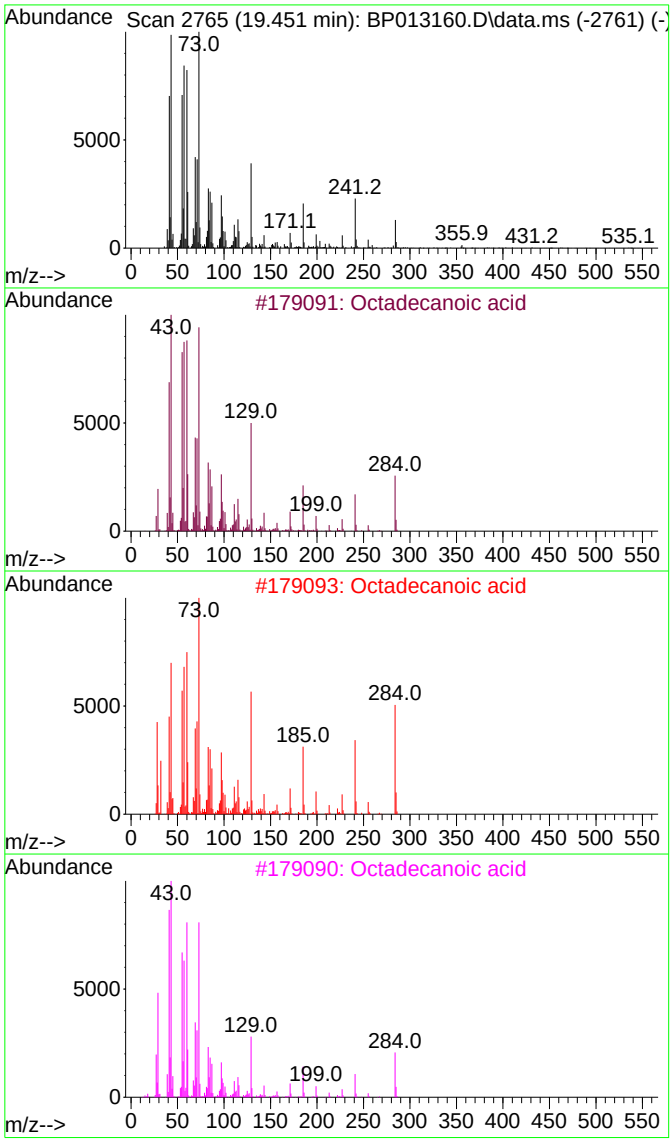
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

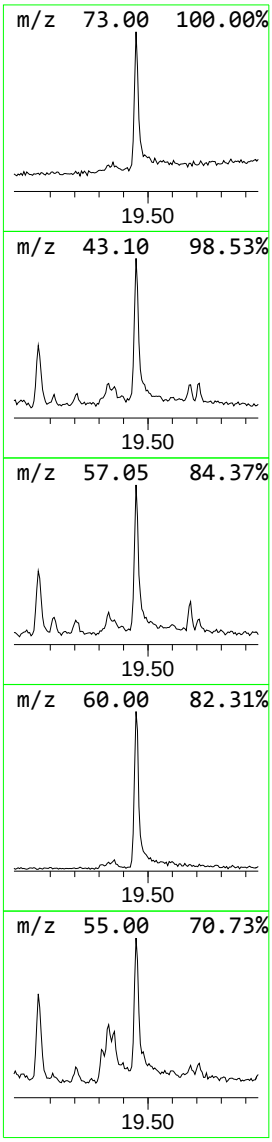
****

Peak Number 4 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.451	2.35 ng/ul	654254	Chrysene-d12	21.439

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	98
4	Octadecanoic acid	284	C18H36O2	000057-11-4	95
5	Tridecanoic acid	214	C13H26O2	000638-53-9	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013160.D
Acq On : 20 Dec 2022 11:22
Operator : CG/JU
Sample : N6003-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ1

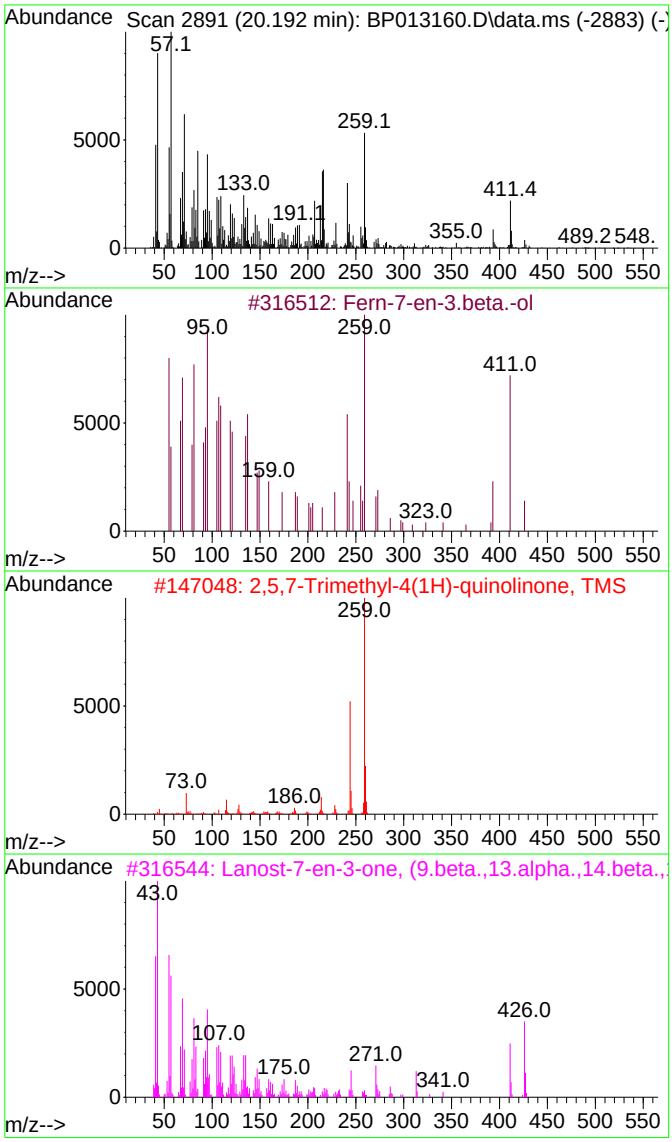
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

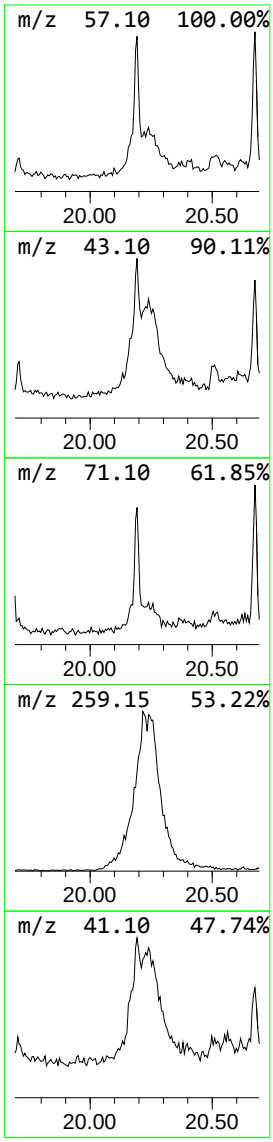
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Fern-7-en-3.beta.-ol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.192	2.89 ng/ul	803774	Chrysene-d12	21.439

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fern-7-en-3.beta.-ol	426	C30H50O	004966-00-1	55
2	2,5,7-Trimethyl-4(1H)-quinolinon...	259	C15H21NOSi	1000497-22-6	25
3	Lanost-7-en-3-one, (9.beta.,13.a...	426	C30H50O	055515-18-9	18
4	6-Fluoro-7-(4-methylpiperazin-1-...	259	C15H18FN3	1000409-37-0	15
5	4-(4-Fluorophenoxy)benzylamine,4...	259	C15H14FNO2	949201-55-2	14





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Operator : CG/JU
Sample : N6003-10
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ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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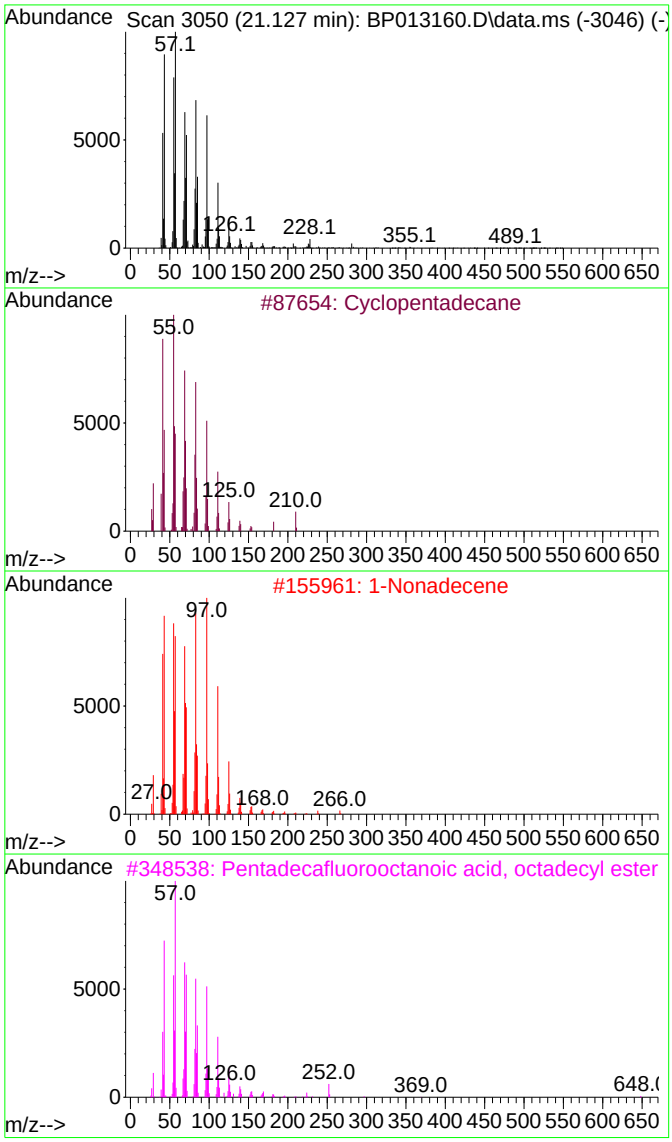
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

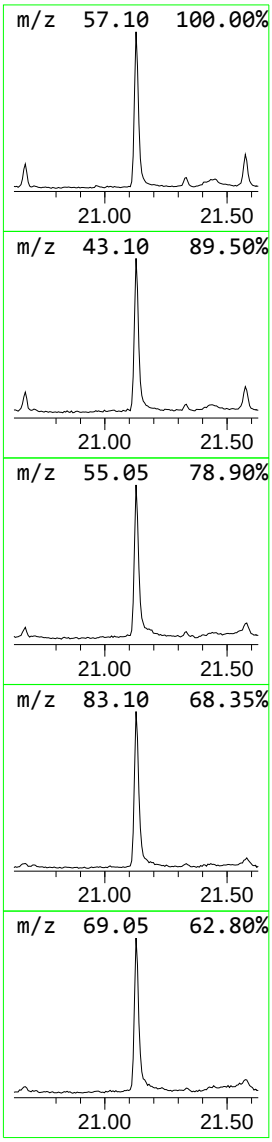
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Cyclic21.127 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.127	8.32 ng/ul	2313240	Chrysene-d12	21.439

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentadecane	210	C15H30	000295-48-7	97
2	1-Nonadecene	266	C19H38	018435-45-5	97
3	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
4	1-Hexadecanol, 2-methyl-	256	C17H36O	002490-48-4	93
5	Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013160.D
Acq On : 20 Dec 2022 11:22
Operator : CG/JU
Sample : N6003-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ1

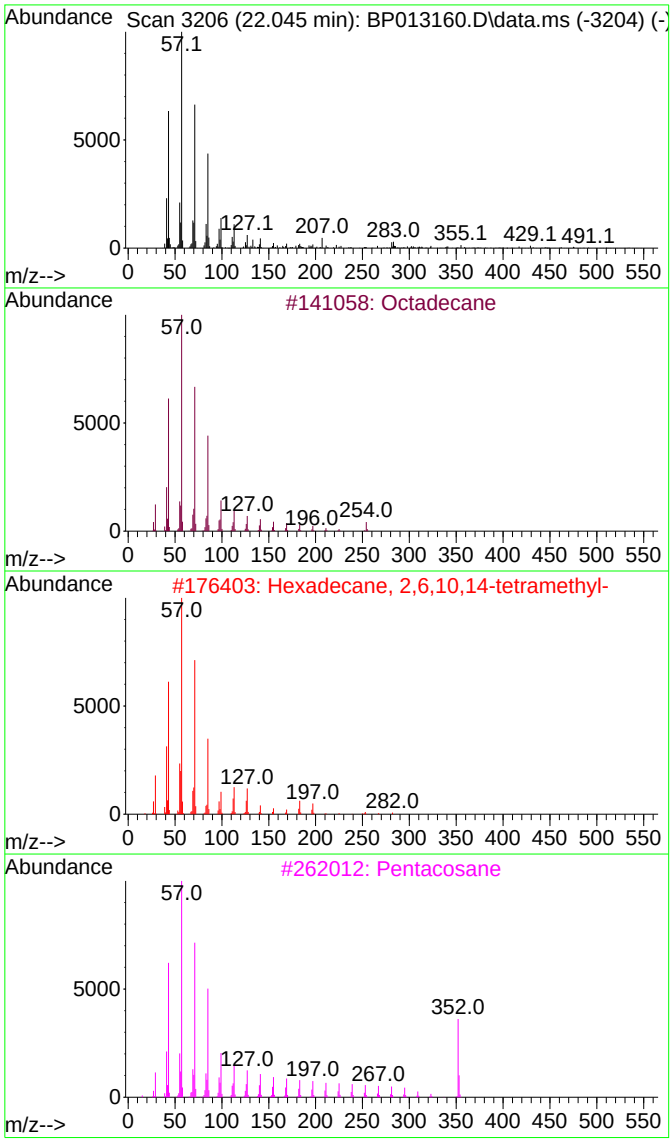
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

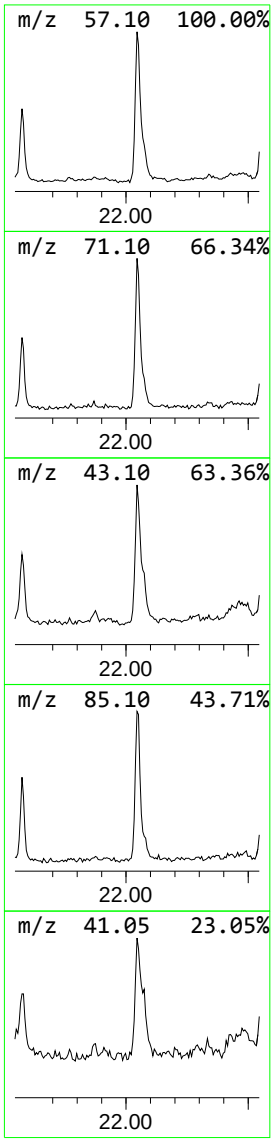
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.045	2.73 ng/ul	759501	Chrysene-d12	21.439

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	95
2	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	93
3	Pentacosane	352	C25H52	000629-99-2	90
4	Hexacosane	366	C26H54	000630-01-3	90
5	Eicosane	282	C20H42	000112-95-8	90





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Data File : BP013160.D
Acq On : 20 Dec 2022 11:22
Operator : CG/JU
Sample : N6003-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ1

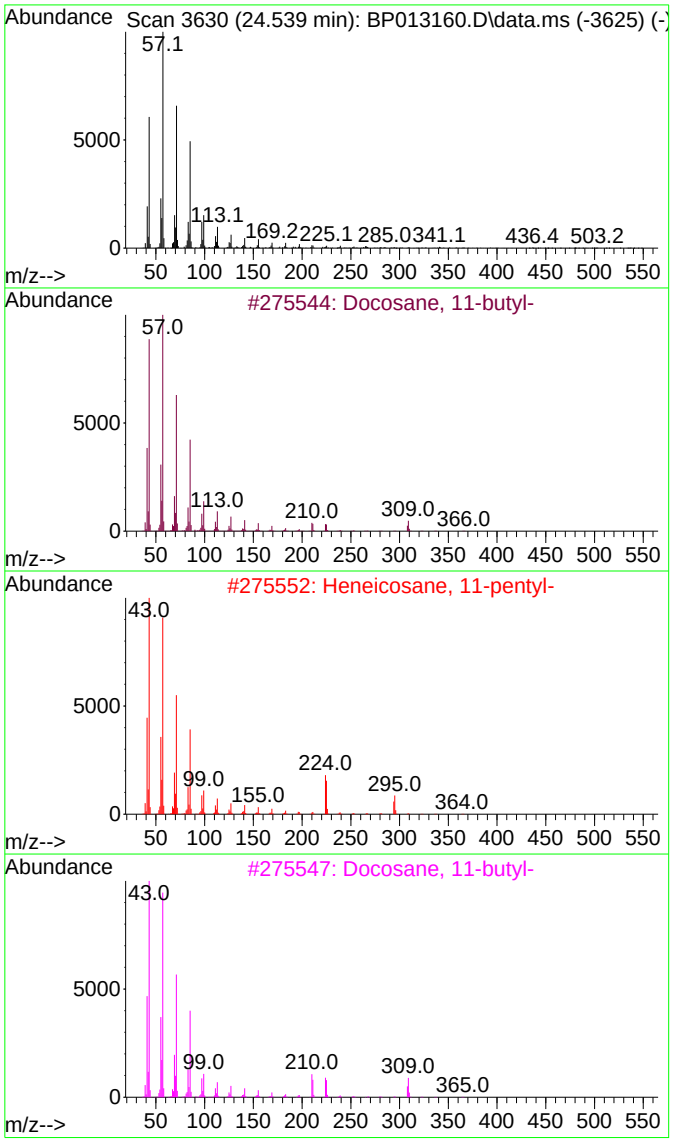
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

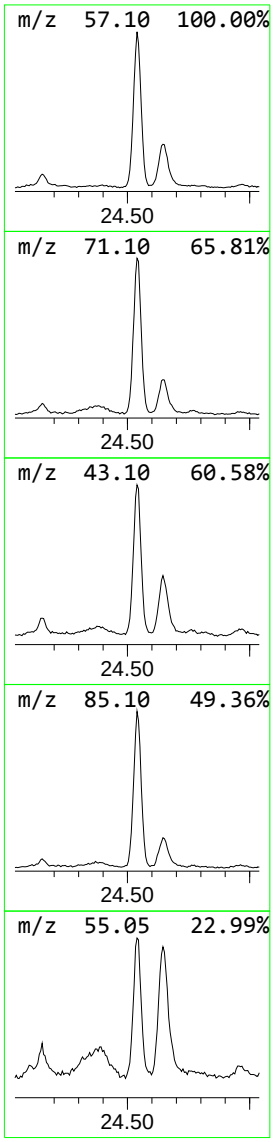
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.539	7.00 ng/ul	2134760	Perylene-d12	23.921

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-	366	C26H54	013475-76-8	95
2	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	94
3	Docosane, 11-butyl-	366	C26H54	013475-76-8	94
4	Octadecane, 2-methyl-	268	C19H40	001560-88-9	94
5	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94





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Acq On : 20 Dec 2022 11:22
Operator : CG/JU
Sample : N6003-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

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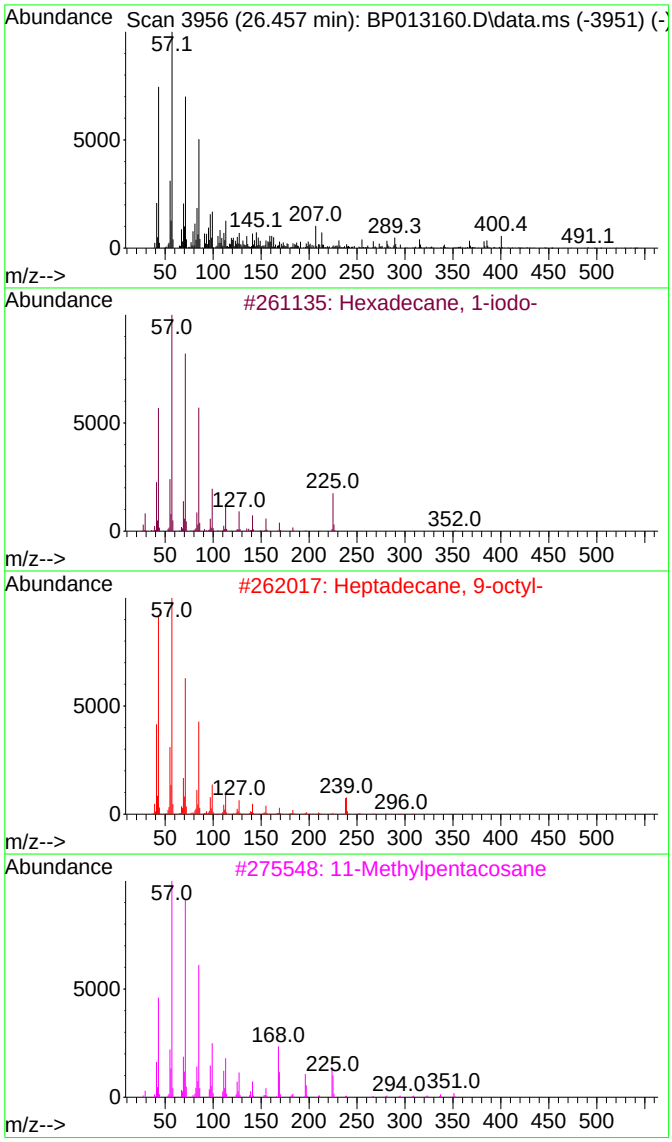
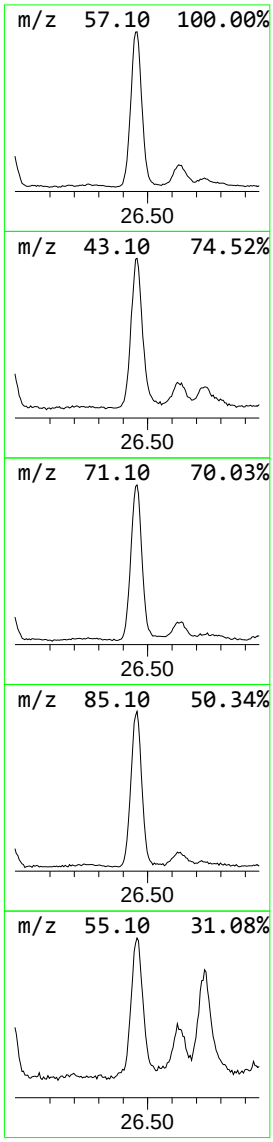
Peak Number 9 Hexadecane, 1-iodo- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.457	10.75 ng/ul	3281720	Perylene-d12	23.921

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	94
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3	11-Methylpentacosane	366	C26H54	015689-71-1	89
4	Eicosane, 10-methyl-	296	C21H44	054833-23-7	89
5	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	89





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Data File : BP013160.D
Acq On : 20 Dec 2022 11:22
Operator : CG/JU
Sample : N6003-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

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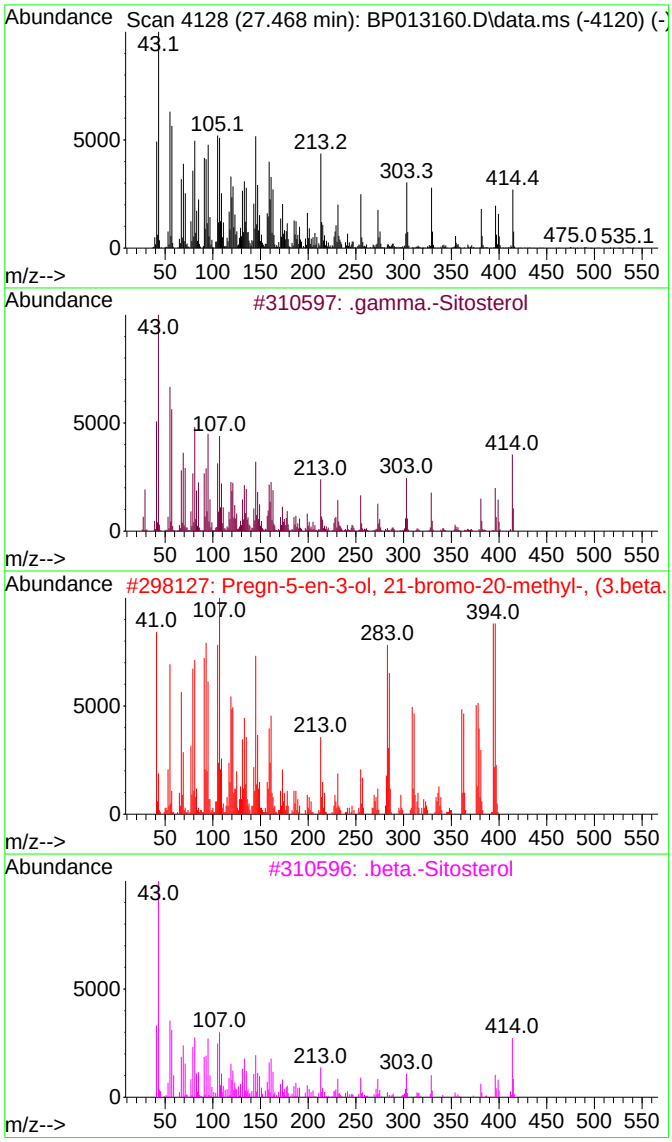
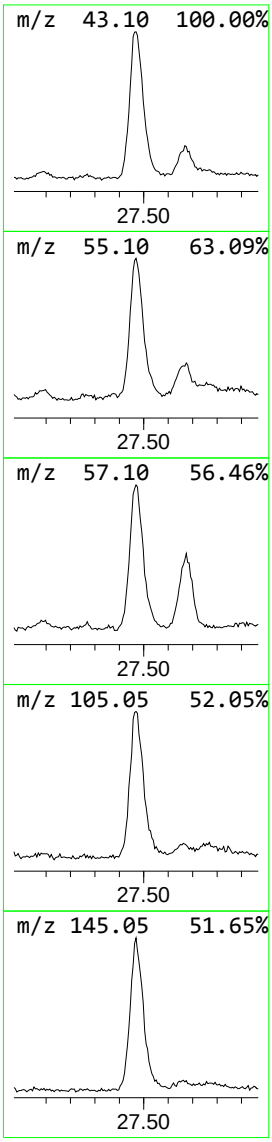
Peak Number 10 .gamma.-Sitosterol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.468	22.97 ng/ul	7010500	Perylene-d12	23.921

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.gamma.-Sitosterol	414	C29H50O	000083-47-6	99
2	Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	95
3	.beta.-Sitosterol	414	C29H50O	000083-46-5	91
4	.beta.-Sitosterol	414	C29H50O	000083-46-5	78
5	Campesterol	400	C28H48O	000474-62-4	53





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Data File : BP013160.D
Acq On : 20 Dec 2022 11:22
Operator : CG/JU
Sample : N6003-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ1

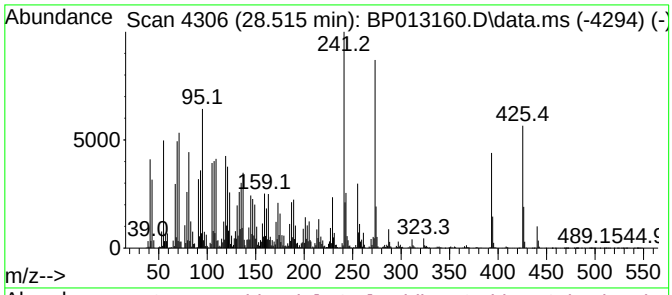
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

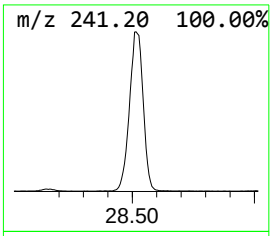
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

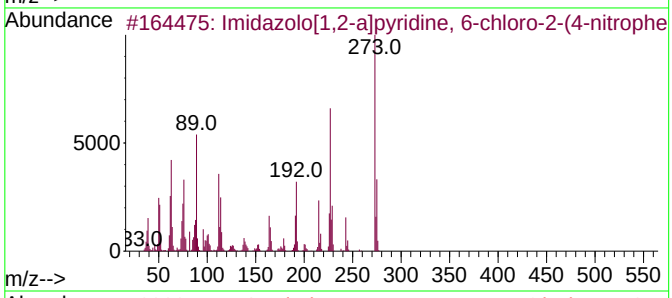
****
Peak Number 11 Imidazolo[1,2-a]pyridine, 6... Concentration Rank 1

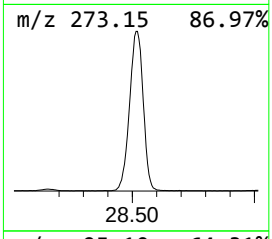
R.T.	EstConc	Area	Relative to ISTD	R.T.
28.515	112.34 ng/ul	34282200	Perylene-d12	23.921

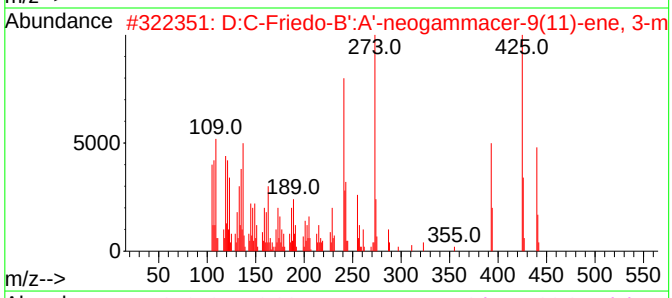
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Imidazolo[1,2-a]pyridine, 6-chlo...	273	C13H8ClN3O2	118000-62-7	56
2	D:C-Friedo-B':A'-neogammacer-9(1...	440	C31H52O	004555-56-0	53
3	16,22-Epoxycoprostan-3.beta.,26,...	434	C27H46O4	122337-91-1	30
4	2-(1H-Indol-3-ylmethylene)-1H-in...	273	C18H11NO2	031060-64-7	25
5	Lupeol	426	C30H50O	000545-47-1	15

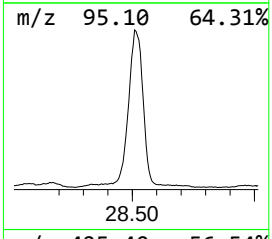


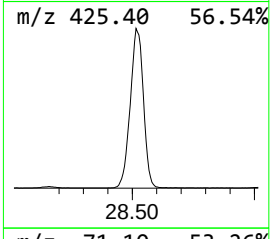


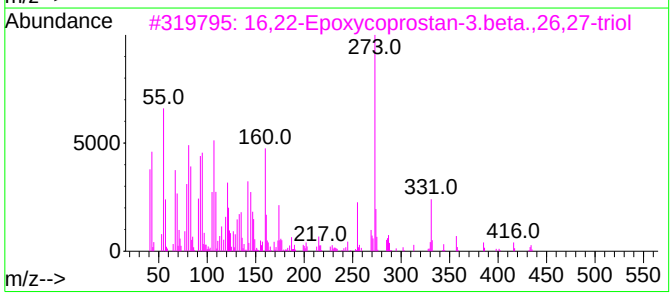


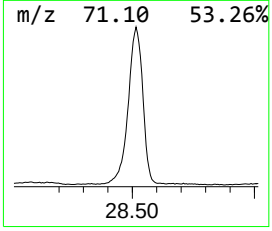












Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013160.D
Acq On : 20 Dec 2022 11:22
Operator : CG/JU
Sample : N6003-10
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA 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
2-Pentanone, 4-...	5.040	2.7	ng/ul	364810	1	7.958	2718280	20.0
n-Hexadecanoic ...	18.222	7.5	ng/ul	1734680	4	17.351	4605440	20.0
Isopropyl palmi...	18.634	3.5	ng/ul	806040	4	17.351	4605440	20.0
Octadecanoic acid	19.451	2.4	ng/ul	654254	5	21.439	5563010	20.0
Fern-7-en-3.bet...	20.192	2.9	ng/ul	803774	5	21.439	5563010	20.0
(DEL) Alkane: C...	21.127	8.3	ng/ul	2313240	5	21.439	5563010	20.0
(DEL) Alkane: S...	22.045	2.7	ng/ul	759501	5	21.439	5563010	20.0
(DEL) Alkane: S...	24.539	7.0	ng/ul	2134760	6	23.921	6103440	20.0
Hexadecane, 1-i...	26.457	10.8	ng/ul	3281720	6	23.921	6103440	20.0
.gamma.-Sitosterol	27.468	23.0	ng/ul	7010500	6	23.921	6103440	20.0
Imidazolo[1,2-a...	28.515	112.3	ng/ul	34282200	6	23.921	6103440	20.0