LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013159.D
Acq On : 20 Dec 2022 10:48
Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

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: BP013159.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	45	rVB	180686	276534	0.42%	0.117%
2	3.764	95	98	114	rBV	1240158	1833962	2.77%	0.778%
3	3.999	133	138	144	rVB	69702	102836	0.16%	0.044%
4	4.322	189	193	198	rVB	74895	95399	0.14%	0.040%
5	4.740	260	264	275	rVB2	51879	84426	0.13%	0.036%
6	5.040	309	315	325	rVB2	289690	439951	0.66%	0.187%
7	6.628	578	585	593	rVB	1468327	2255671	3.40%	0.957%
8	6.934	628	637	642	rBV7	43792	88731	0.13%	0.038%
9	7.110	659	667	683	rVB	2420288	3814797	5.76%	1.618%
10	7.281	688	696	705	rVV	1972496	3012107	4.55%	1.277%
11	7.387	708	714	721	rVV	346168	533041	0.80%	0.226%
12	7.481	723	730	740	rVV	2511833	3881832	5.86%	1.646%
13	7.952	803	810	818	rBV	1906912	3031298	4.58%	1.285%
14	8.663	923	931	943	rBV	2439248	3918038	5.91%	1.661%
15	8.781	946	951	961	rVB	101686	177148	0.27%	0.075%
16	9.122	1001	1009	1016	rBV	2180651	3542191	5.35%	1.502%
17	9.528	1071	1078	1085	rBV2	59371	113990	0.17%	0.048%
18	9.846	1123	1132	1151	rBV	1817162	2999298	4.53%	1.272%
19	10.075	1165	1171	1182	rVB4	74202	159673	0.24%	0.068%
20	10.387	1216	1224	1241	rBV	2739045	4577232	6.91%	1.941%
21	10.769	1280	1289	1296	rBV	2425025	4025540	6.08%	1.707%
22	10.904	1305	1312	1331	rVB	1605336	2842915	4.29%	1.206%
23	12.351	1552	1558	1565	rVB2	50540	82365	0.12%	0.035%
24	12.869	1639	1646	1652	rBV	172552	266015	0.40%	0.113%
25	13.587	1762	1768	1773	rVB2	39200	66474	0.10%	0.028%
26	13.916	1819	1824	1831	rVV2	199720	295215	0.45%	0.125%
27	14.004	1831	1839	1854	rVV	3911986	5638318	8.51%	2.391%
28	14.292	1881	1888	1903	rVV	4838635	7216320	10.89%	3.060%
29	14.598	1933	1940	1946	rVB	3110614	4633392	6.99%	1.965%
30	14.657	1946	1950	1955	rBV3	67541	103621	0.16%	0.044%
31	14.792	1966	1973	1978	rBV	1406177	2229781	3.37%	0.946%
32	14.839	1978	1981	1990	rVB2	162488	293822	0.44%	0.125%
33	14.945	1995	1999	2004	rVV2	84588	131355	0.20%	0.056%
34	14.998	2004	2008	2014	rVV3	104993	192981	0.29%	0.082%
35	15.504	2086	2094	2099	rBV4	94006	180414	0.27%	0.077%
36	15.592	2101	2109	2122	rBV	5853050	8570665	12.94%	3.634%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013159.D
Acq On : 20 Dec 2022 10:48
Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

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	15.704	2122	2128	2141	rBV	1619975	2213293	3.34%	0.939%
38	15.834	2146	2150	2156	rVB4	43541	74510	0.11%	0.032%
39	15.951	2166	2170	2176	rVB	66772	98941	0.15%	0.042%
40	16.163	2201	2206	2211	rVB2	94374	130979	0.20%	0.056%
41	16.733	2299	2303	2308	rBV	96343	133654	0.20%	0.057%
42	17.351	2402	2408	2413	rVV	3546173	4933282	7.45%	2.092%
43	17.392	2413	2415	2419	rVV	119135	151498	0.23%	0.064%
44	17.451	2419	2425	2433	rVV	5495902	7788581	11.76%	3.303%
45	17.533	2436	2439	2444	rVB2	82796	120202	0.18%	0.051%
46	18.110	2533	2537	2541	rBV4	45080	70867	0.11%	0.030%
47	18.163	2542	2546	2551	rBV	473099	652349	0.98%	0.277%
48	18.227	2551	2557	2573	rVV	2136446	3160181	4.77%	1.340%
49	18.557	2609	2613	2621	rVB4	74508	112880	0.17%	0.048%
50	18.633	2623	2626	2633	rBV3	119501	190938	0.29%	0.081%
51	19.051	2694	2697	2701	rBV4	70794	91363	0.14%	0.039%
52	19.110	2702	2707	2712	rVV7	99219	192233	0.29%	0.082%
53	19.263	2730	2733	2738	rVB3	116829	163262	0.25%	0.069%
54	19.457	2761	2766	2786	rBV	6741992	9948655	15.02%	4.219%
55	19.680	2799	2804	2819	rVB	7534740	10654768	16.08%	4.518%
56	20.110	2874	2877	2881	rBV2	183389	248294	0.37%	0.105%
57	20.627	2961	2965	2968	rBV	308460	346845	0.52%	0.147%
58	21.127	3045	3050	3055	rBV2	1711687	2785402	4.20%	1.181%
59	21.439	3097	3103	3107	rBV	4479174	6299941	9.51%	2.671%
60	22.068	3208	3210	3219	rVB3	552338	853321	1.29%	0.362%
61	23.133	3388	3391	3398	rBV2	381268	779949	1.18%	0.331%
62	23.757	3490	3497	3511	rVB2	5814480	12327076	18.61%	5.227%
63	23.921	3518	3525	3533	rVB2	3059108	6316396	9.53%	2.678%
64	24.539	3625	3630	3638	rVB	967090	1999020	3.02%	0.848%
65	24.639	3640	3647	3665	rVB2	2892802	7468509	11.27%	3.167%
66	26.456	3949	3956	3977	rVB2	1777484	6213941	9.38%	2.635%
67	27.480	4120	4130	4150	rVB2	3087744	11345222	17.13%	4.811%
68	28.533	4294	4309	4327	rVB2	16174982	66247963	100.00%	28.092%

Sum of corrected areas: 235821663

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
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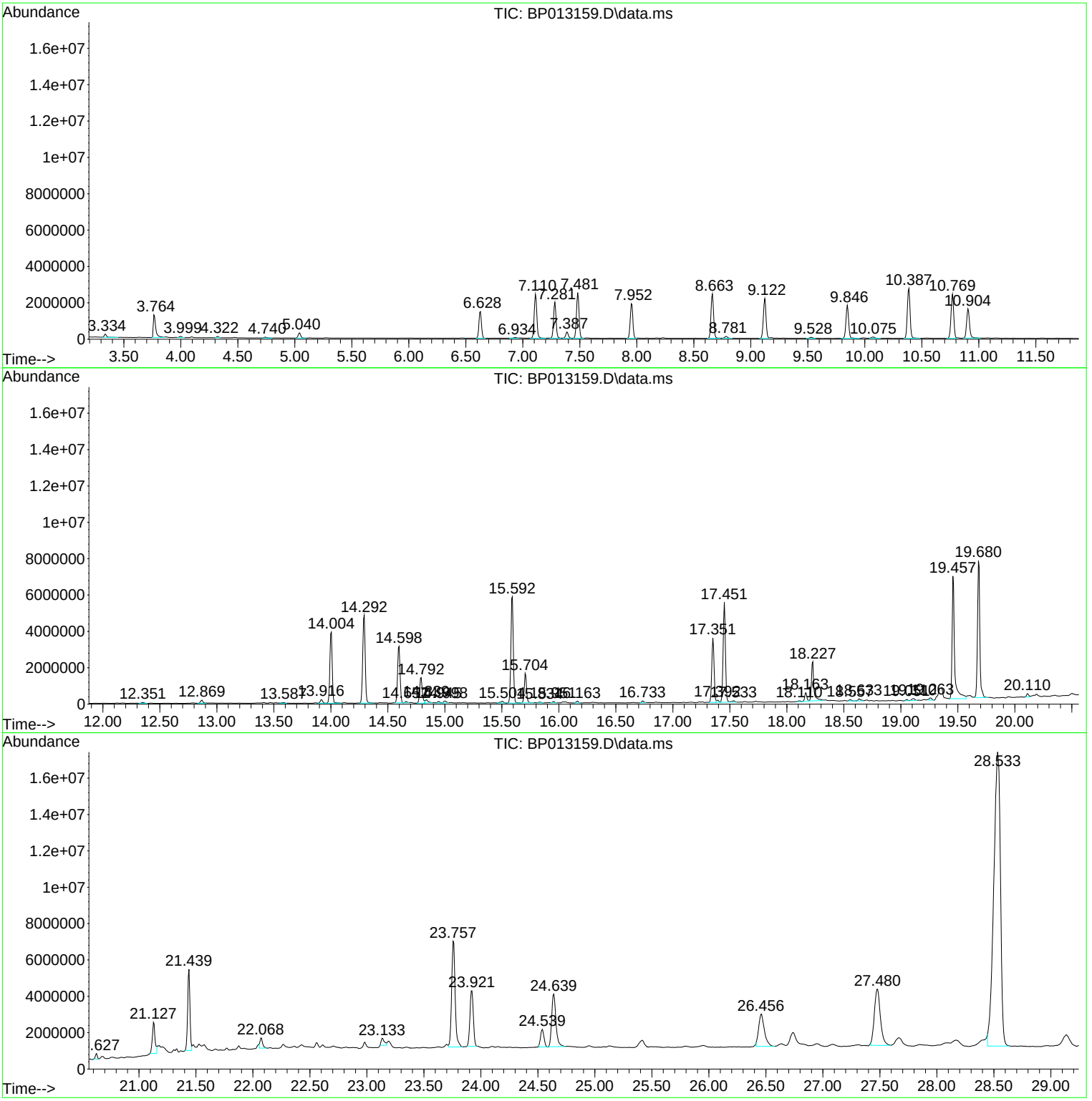
Acq On : 20 Dec 2022 10:48
Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

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 : BP013159.D
Acq On : 20 Dec 2022 10:48
Operator : CG/JU
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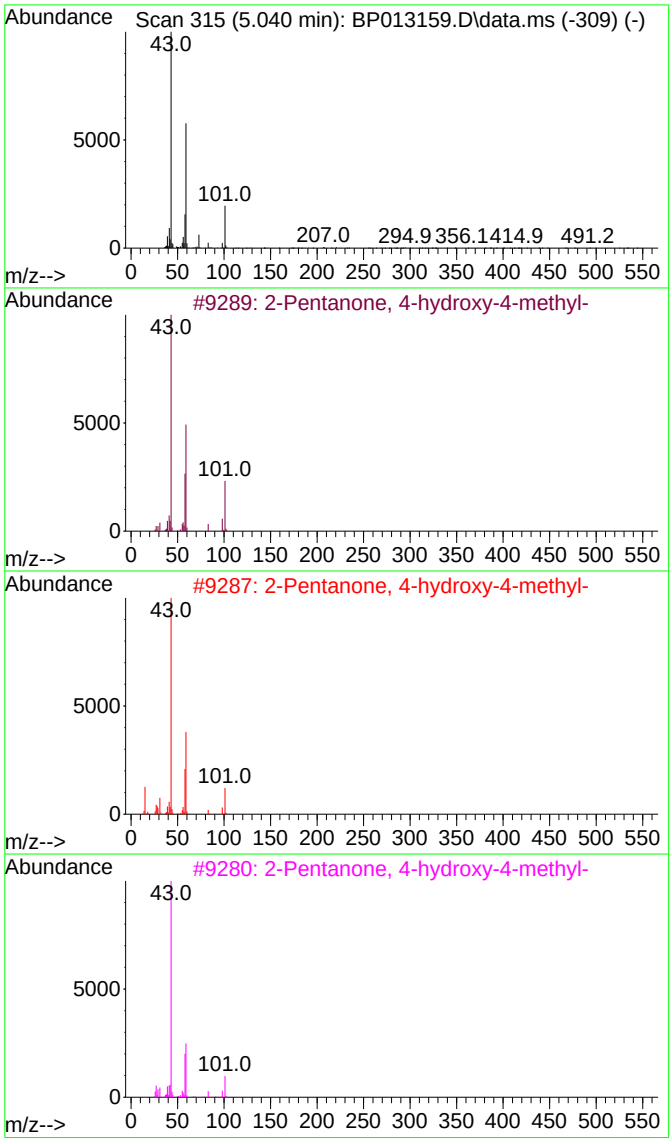
Quant Title : SVOA CALIBRATION

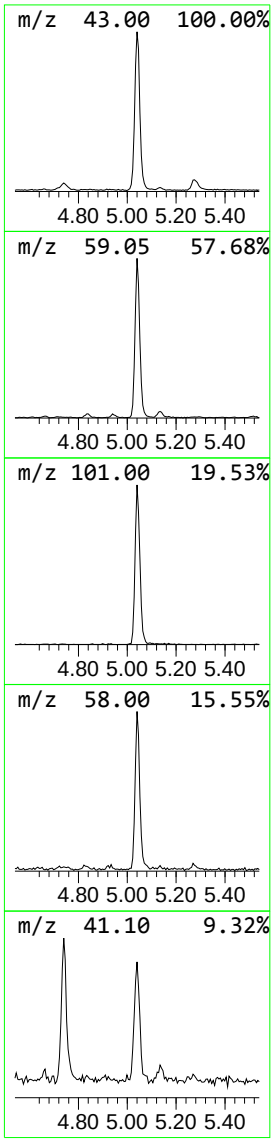
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.040	2.90 ng/ul	439951	1,4-Dichlorobenzene-d4	7.952

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	50
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	45
5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013159.D
Acq On : 20 Dec 2022 10:48
Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

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

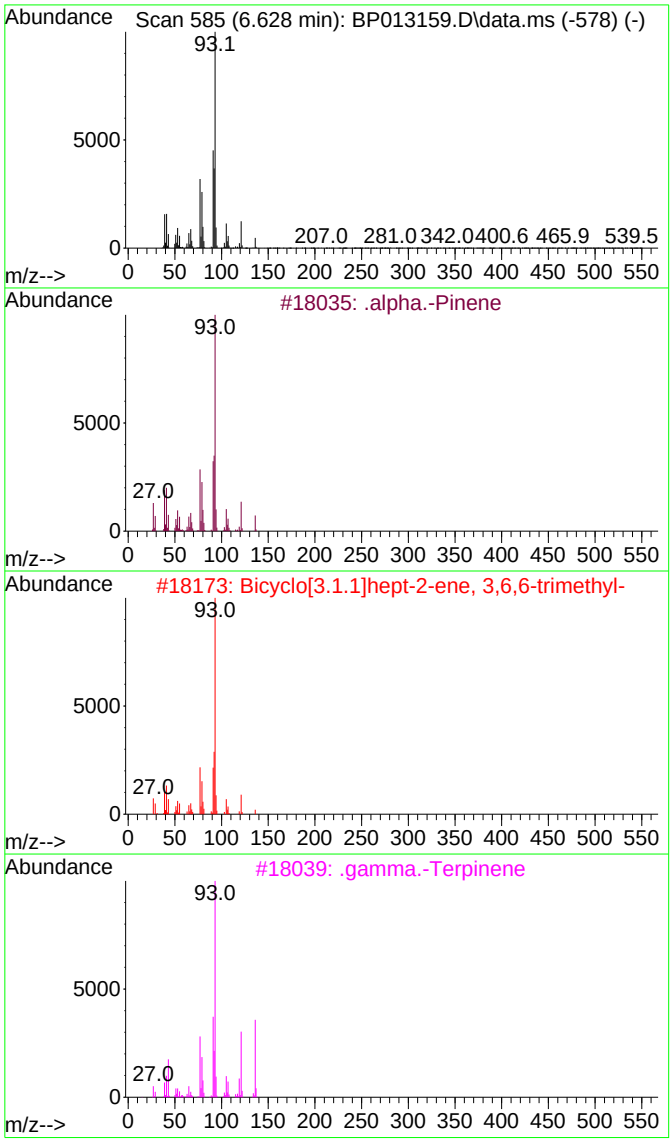
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

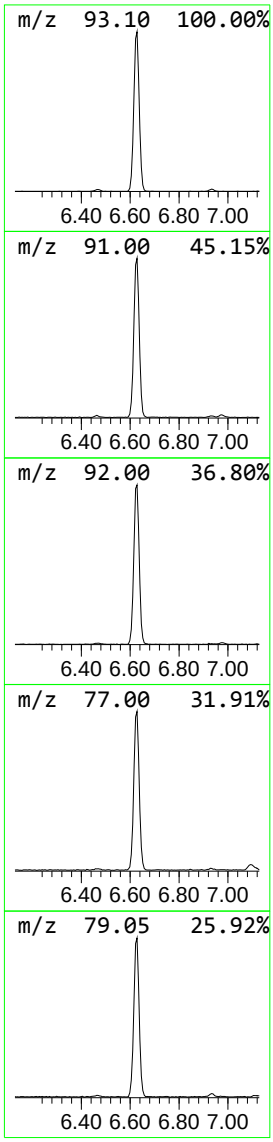
****

Peak Number 2 .alpha.-Pinene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.628	14.88 ng/ul	2255670	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Pinene	136	C10H16	000080-56-8	94
2	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	91
3	.gamma.-Terpinene	136	C10H16	000099-85-4	91
4	3-Carene	136	C10H16	013466-78-9	91
5	(+)-3-Carene	136	C10H16	000498-15-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013159.D
Acq On : 20 Dec 2022 10:48
Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

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

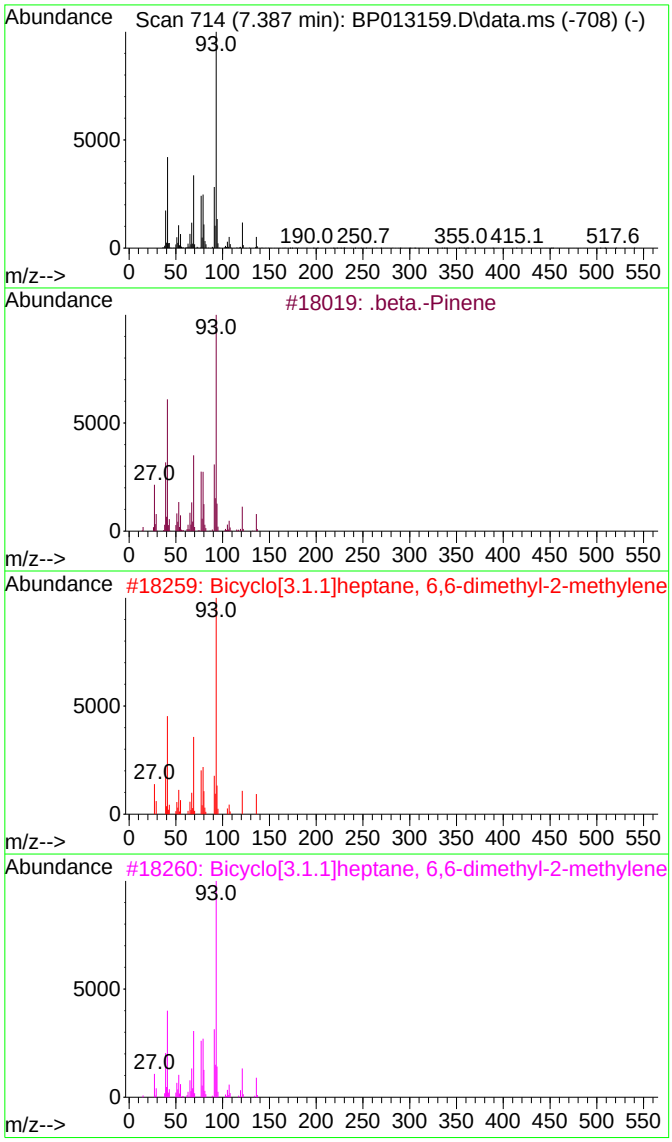
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

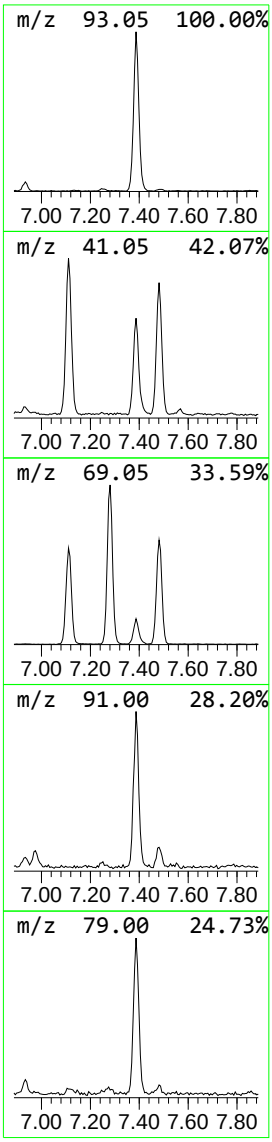
****

Peak Number 3 .beta.-Pinene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.387	3.52 ng/ul	533041	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-Pinene	136	C10H16	000127-91-3	93
2	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	91
3	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	91
4	.beta.-Pinene	136	C10H16	000127-91-3	91
5	.beta.-Pinene	136	C10H16	000127-91-3	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013159.D
Acq On : 20 Dec 2022 10:48
Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

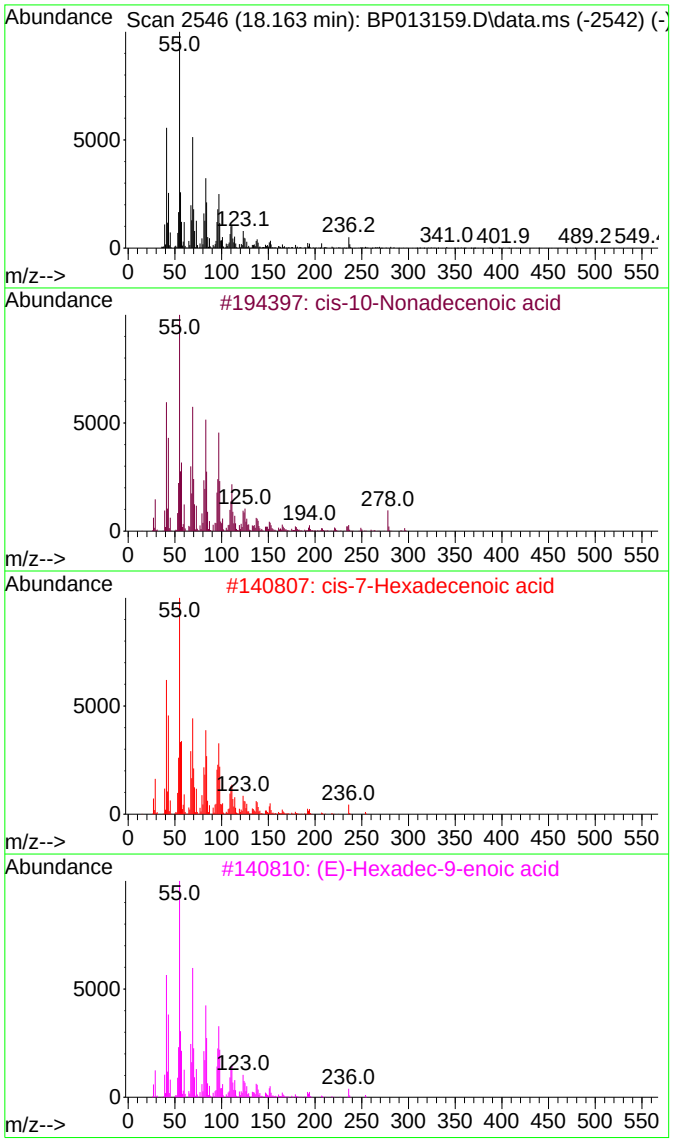
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

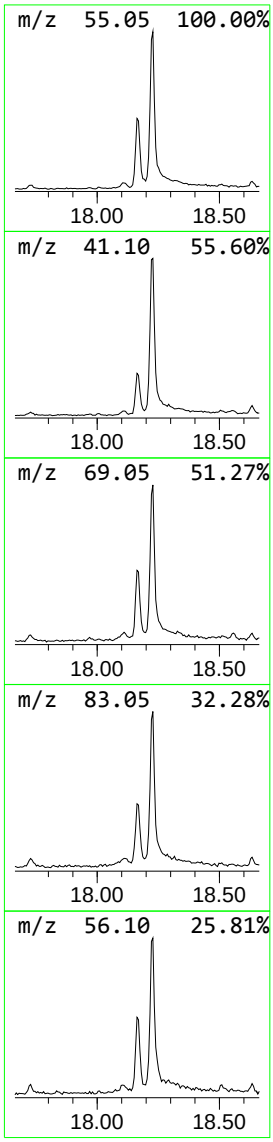
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 cis-10-Nonadecenoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.163	2.64 ng/ul	652349	Phenanthrene-d10	17.351

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	cis-10-Nonadecenoic acid	296	C19H36O2	073033-09-7	99
2	cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	98
3	(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	98
4	cis-11-Eicosenoic acid	310	C20H38O2	005561-99-9	91
5	cis-Vaccenic acid	282	C18H34O2	000506-17-2	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013159.D
Acq On : 20 Dec 2022 10:48
Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

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

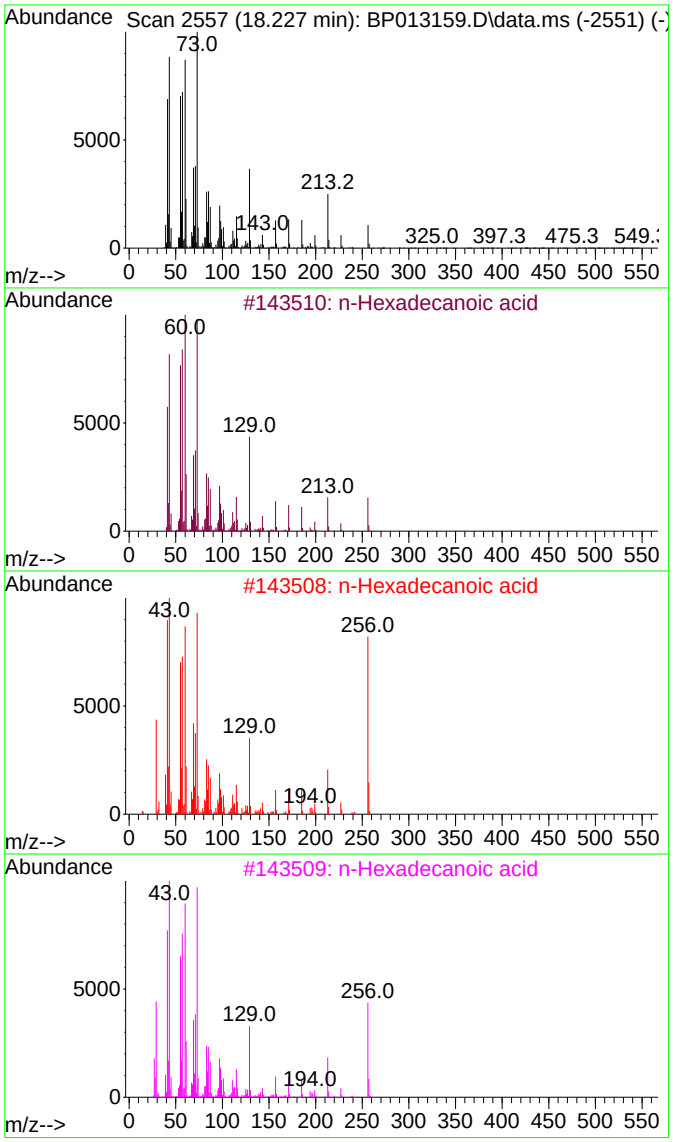
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

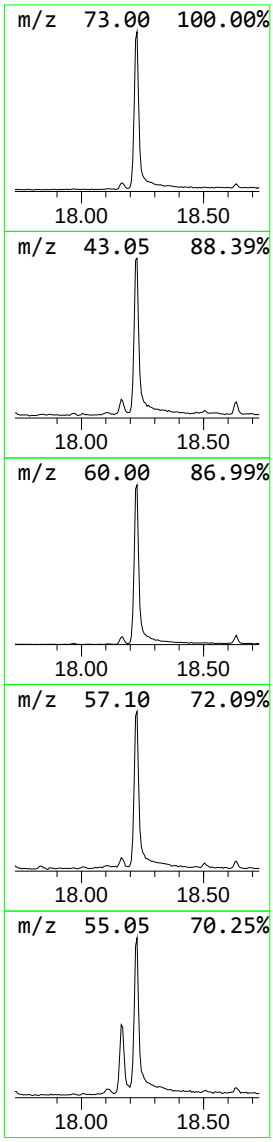
****

Peak Number 5 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.227	12.81 ng/ul	3160180	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	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
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Acq On : 20 Dec 2022 10:48
Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

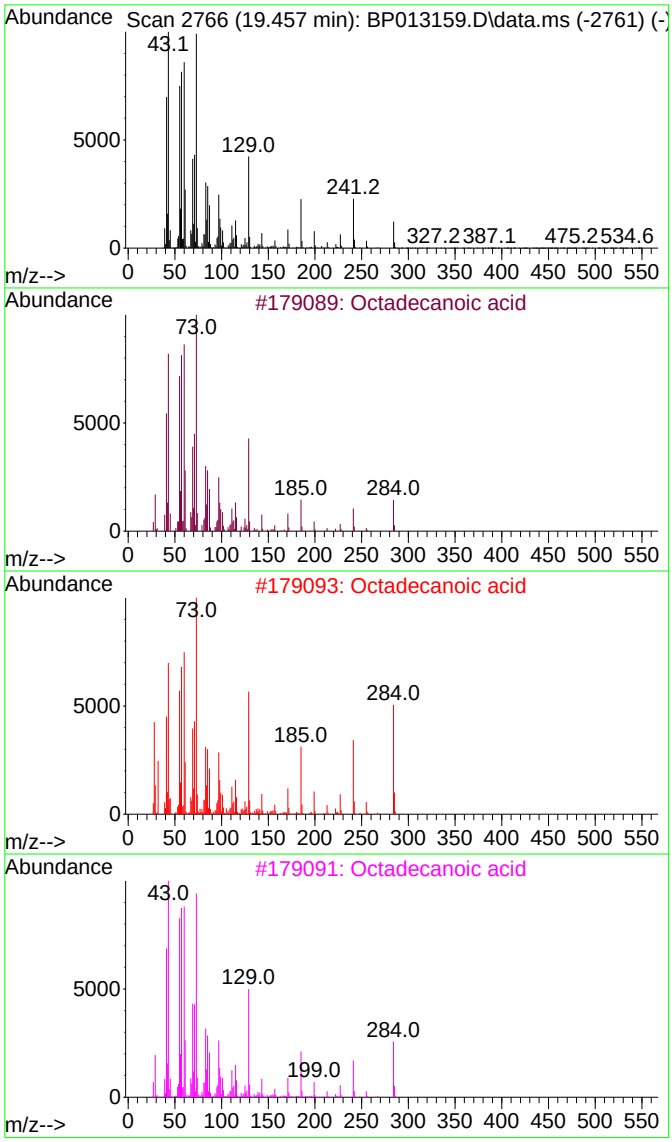
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

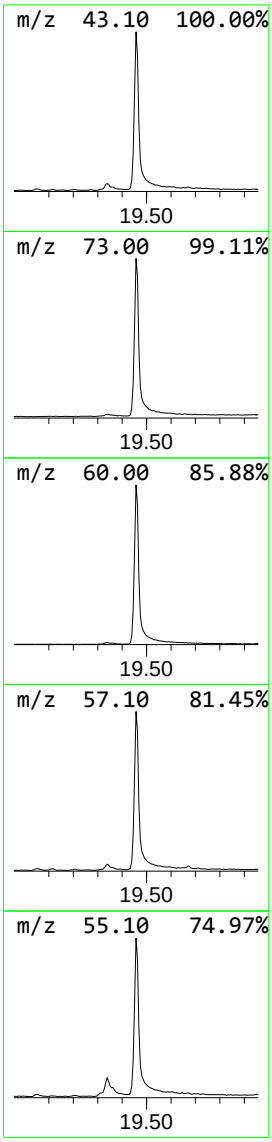
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.457	31.58 ng/ul	9948660	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	99
4	Octadecanoic acid	284	C18H36O2	000057-11-4	98
5	Octadecanoic acid	284	C18H36O2	000057-11-4	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013159.D
Acq On : 20 Dec 2022 10:48
Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

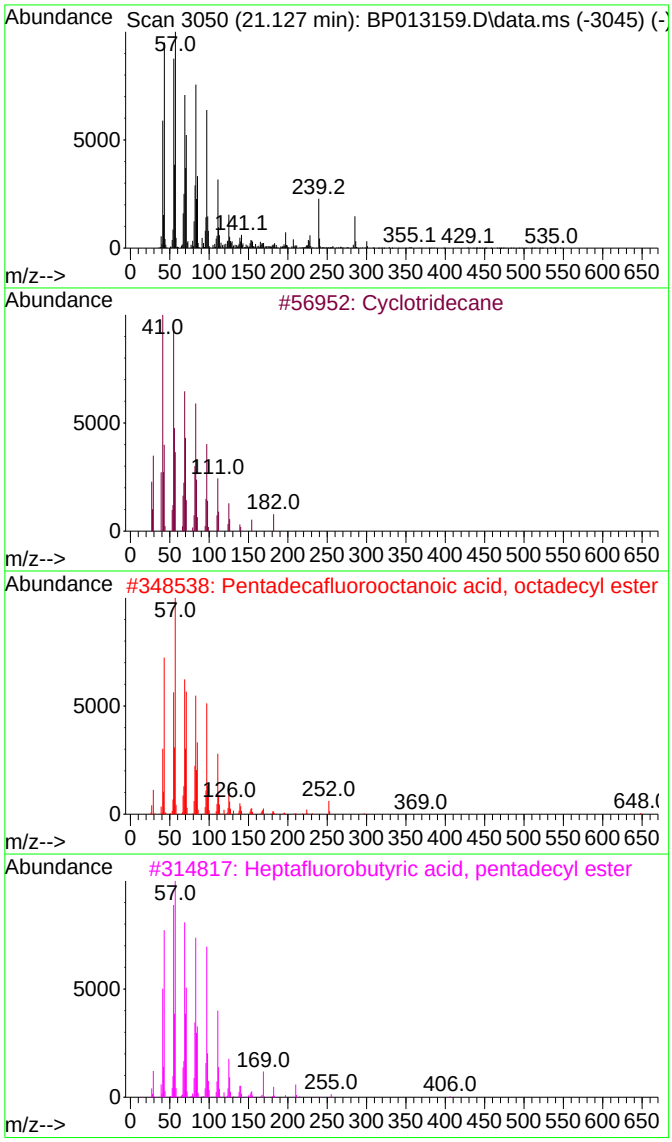
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

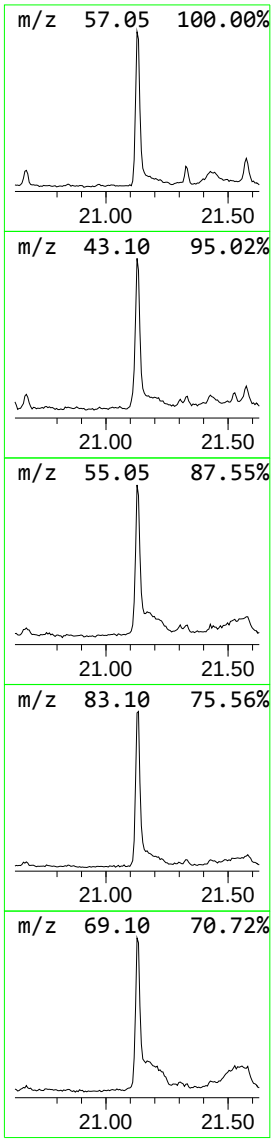
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Cyclic21.127 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.127	8.84 ng/ul	2785400	Chrysene-d12	21.439

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotridecane	182	C13H26	000295-02-3	97
2	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91
3	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	91
4	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
5	1-Hexacosene	364	C26H52	018835-33-1	91





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Data File : BP013159.D
Acq On : 20 Dec 2022 10:48
Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

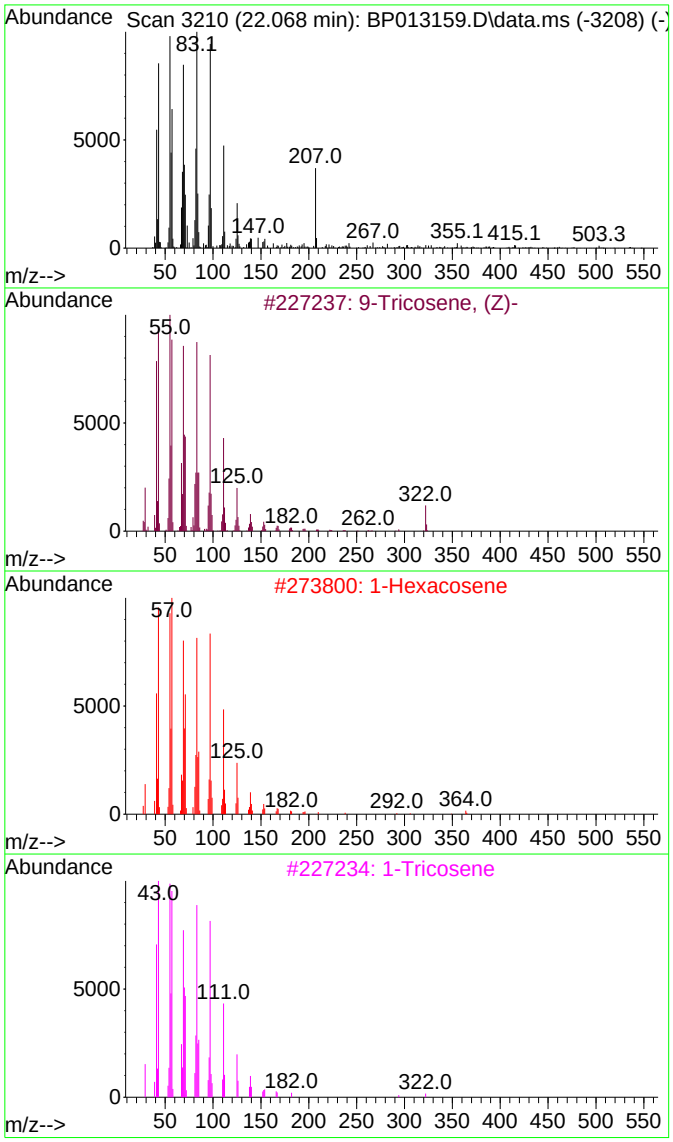
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

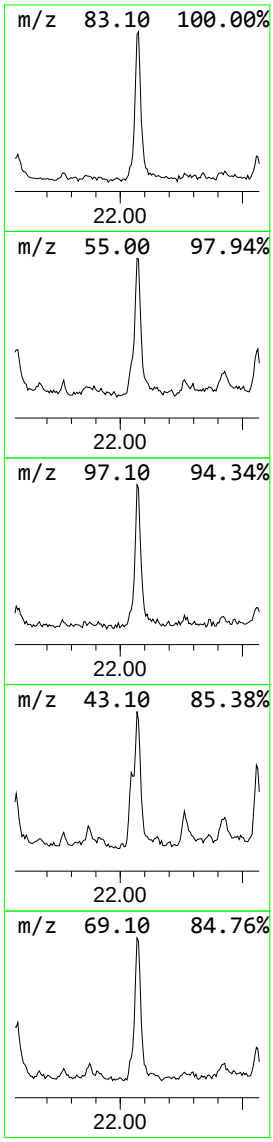
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.069	2.71 ng/ul	853321	Chrysene-d12	21.439

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Tricosene, (Z)-	322	C23H46	027519-02-4	87
2	1-Hexacosene	364	C26H52	018835-33-1	83
3	1-Tricosene	322	C23H46	018835-32-0	83
4	Dodecane, 1-cyclopentyl-4-(3-cyc...	348	C25H48	007225-68-5	72
5	n-Tetracosanol-1	354	C24H50O	000506-51-4	72





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Acq On : 20 Dec 2022 10:48
Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

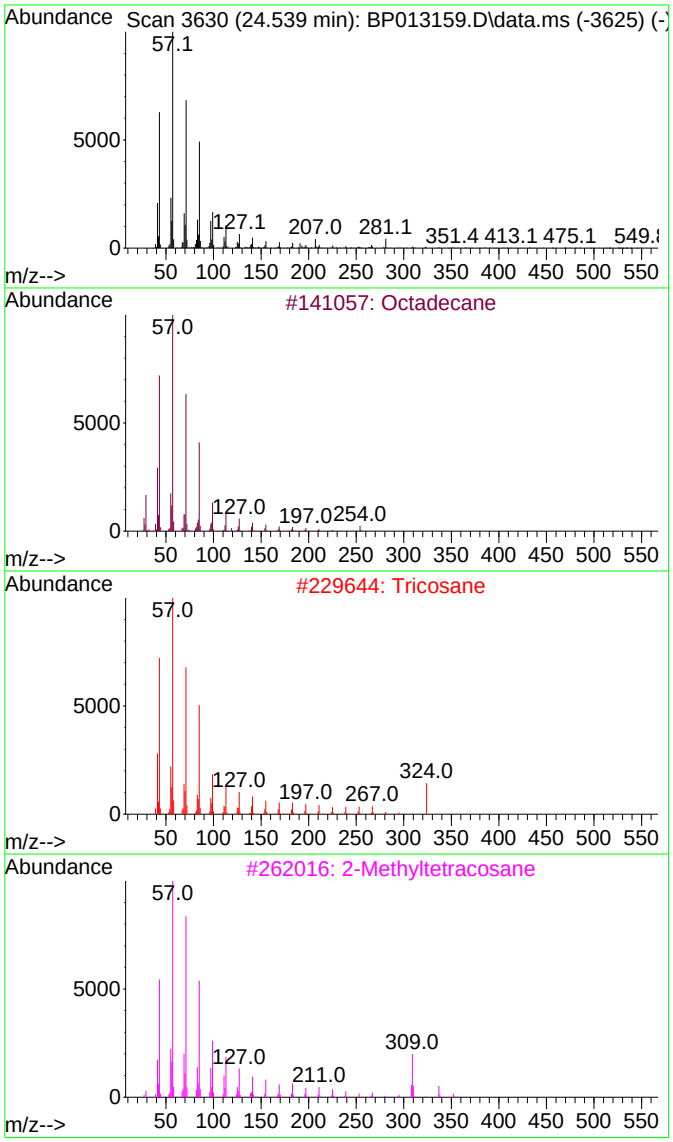
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

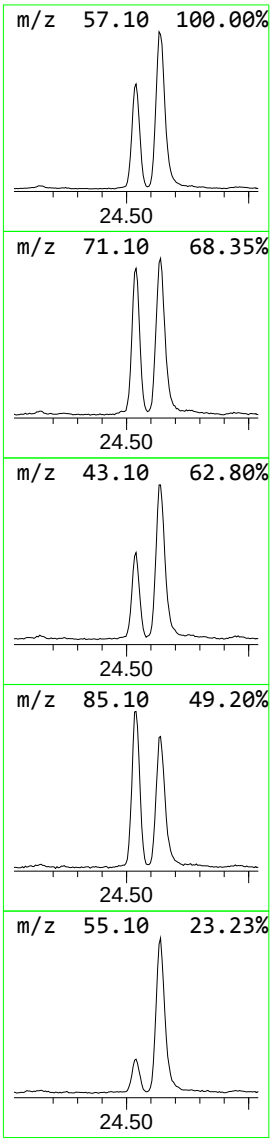
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.539	6.33 ng/ul	1999020	Perylene-d12	23.921

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	95
2	Tricosane	324	C23H48	000638-67-5	94
3	2-Methyltetracosane	352	C25H52	001560-78-7	94
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
5	Tetracosane	338	C24H50	000646-31-1	93





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Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

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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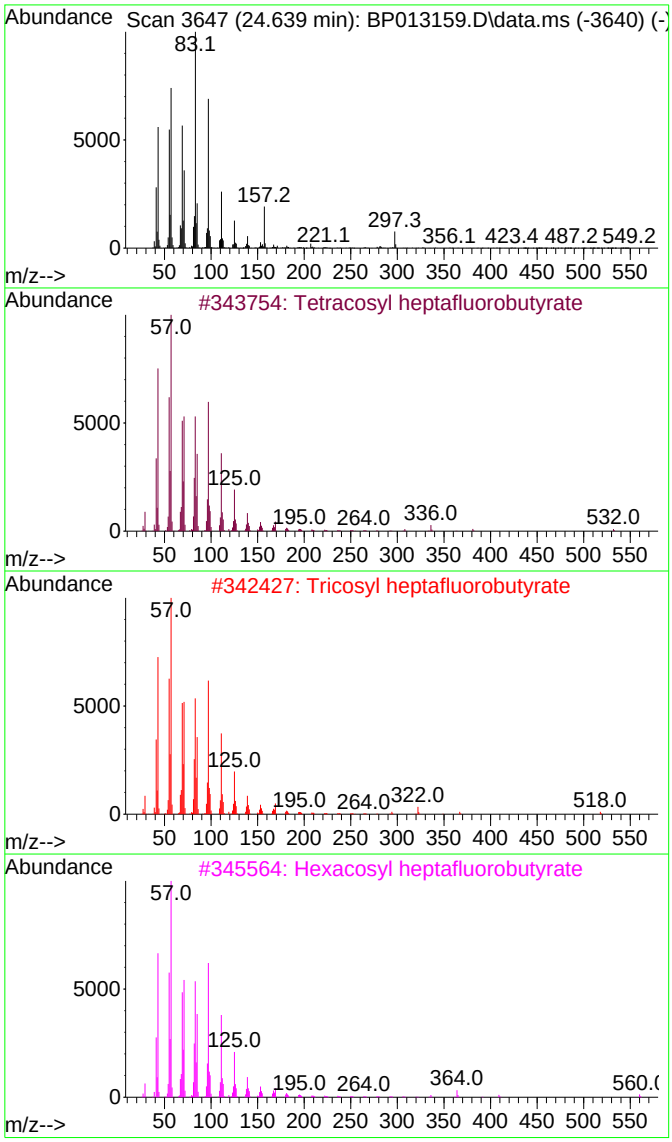
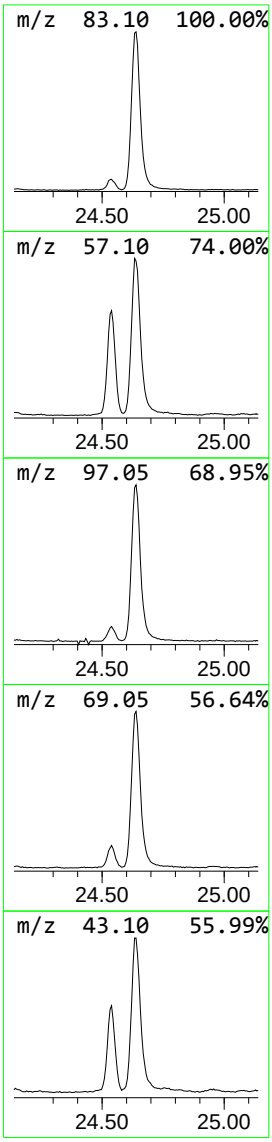
Peak Number 10 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.639	23.65 ng/ul	7468510	Perylene-d12	23.921

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosyl heptafluorobutyrate	550	C28H49F7O2	1000351-83-7	49
2	Tricosyl heptafluorobutyrate	536	C27H47F7O2	1000351-83-4	49
3	Hexacosyl heptafluorobutyrate	578	C30H53F7O2	1000351-83-3	49
4	Hexacosyl pentafluoropropionate	528	C29H53F5O2	1000351-81-2	49
5	Hexacosyl trifluoroacetate	478	C28H53F3O2	1000351-75-0	49





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Data File : BP013159.D
Acq On : 20 Dec 2022 10:48
Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

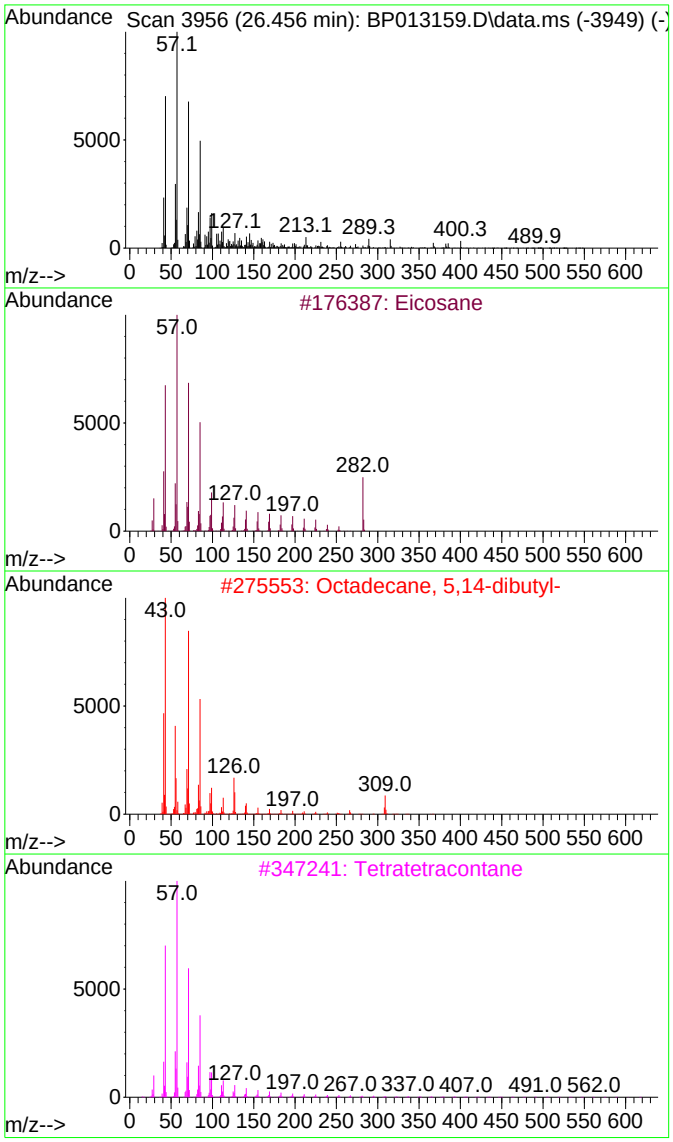
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

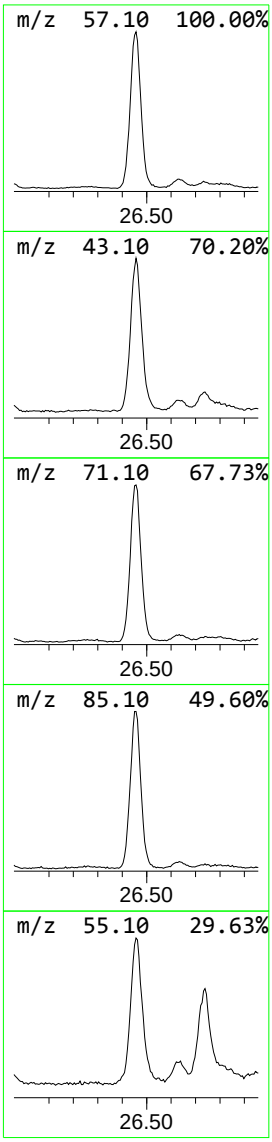
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.456	19.68 ng/ul	6213940	Perylene-d12	23.921

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	94
2	Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	93
3	Tetratetracontane	619	C44H90	007098-22-8	93
4	Docosane, 9-butyl-	366	C26H54	055282-14-9	93
5	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93





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Data File : BP013159.D
Acq On : 20 Dec 2022 10:48
Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

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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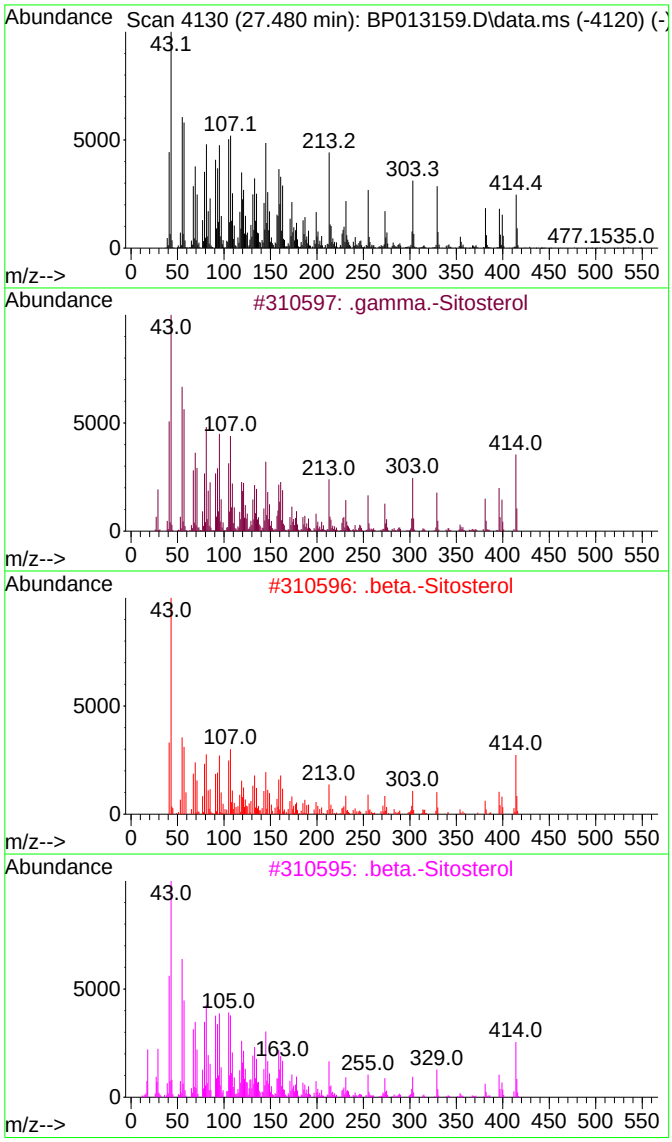
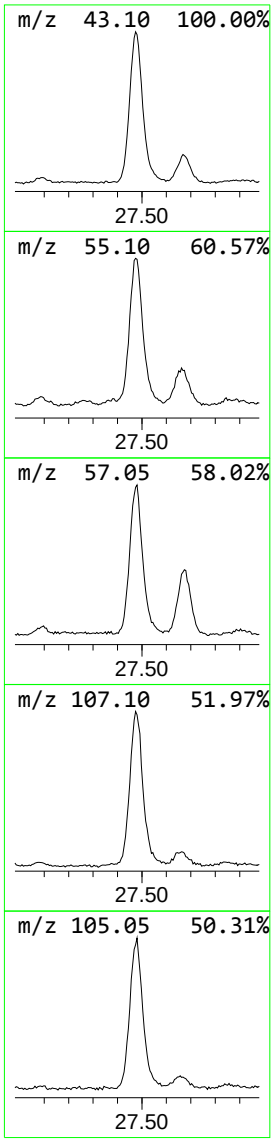
Peak Number 12 .gamma.-Sitosterol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.480	35.92 ng/ul	11345200	Perylene-d12	23.921

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.gamma.-Sitosterol	414	C29H50O	000083-47-6	99
2	.beta.-Sitosterol	414	C29H50O	000083-46-5	95
3	.beta.-Sitosterol	414	C29H50O	000083-46-5	91
4	.gamma.-Sitosterol	414	C29H50O	000083-47-6	83
5	Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	74





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Data File : BP013159.D
Acq On : 20 Dec 2022 10:48
Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

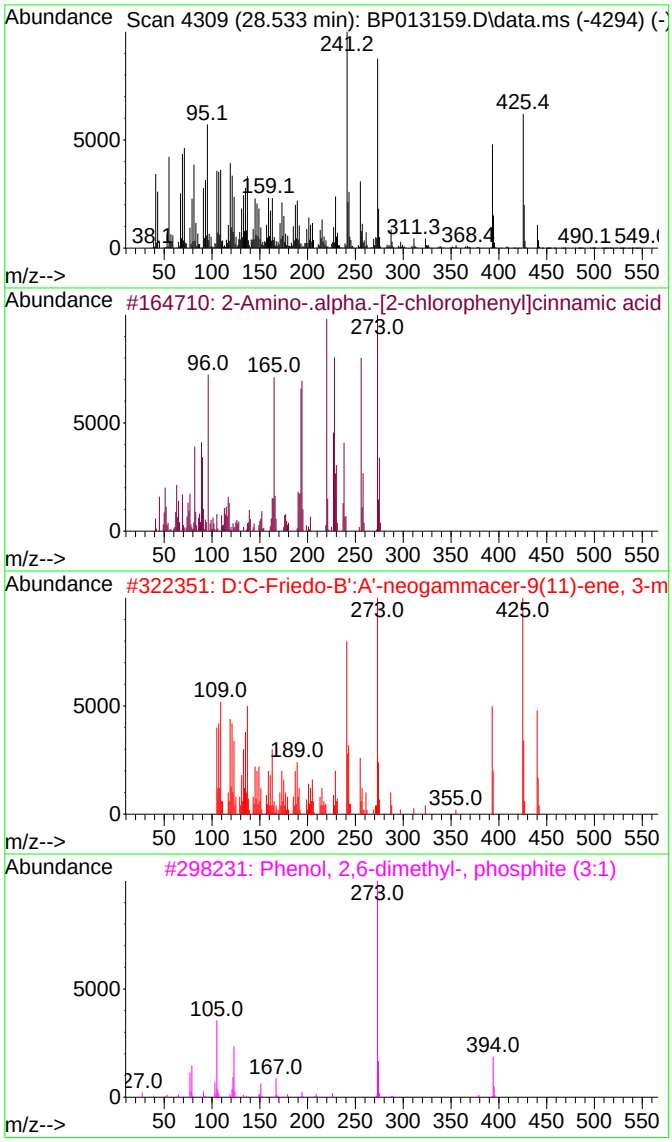
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

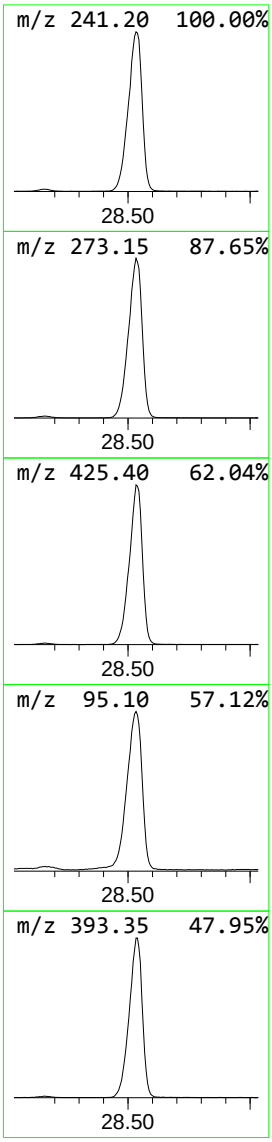
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 2-Amino-.alpha.-[2-chloroph... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.533	209.76 ng/ul	66248000	Perylene-d12	23.921

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Amino-.alpha.-[2-chlorophenyl]...	273	C15H12ClNO2	1000254-08-5	59
2	D:C-Friedo-B':A'-neogammacer-9(1...	440	C31H52O	004555-56-0	58
3	Phenol, 2,6-dimethyl-, phosphite...	394	C24H27O3P	052830-49-6	25
4	Benzamide, 2,4,6-tris(1,1-dimeth...	393	C27H39NO	016420-52-3	25
5	2-(1H-Indol-3-ylmethylene)-1H-in...	273	C18H11NO2	031060-64-7	25





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013159.D
Acq On : 20 Dec 2022 10:48
Operator : CG/JU
Sample : N6003-09
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ0

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	2.9	ng/ul	439951	1	7.952	3031300	20.0
.alpha.-Pinene	6.628	14.9	ng/ul	2255670	1	7.952	3031300	20.0
.beta.-Pinene	7.387	3.5	ng/ul	533041	1	7.952	3031300	20.0
cis-10-Nonadece...	18.163	2.6	ng/ul	652349	4	17.351	4933280	20.0
n-Hexadecanoic ...	18.227	12.8	ng/ul	3160180	4	17.351	4933280	20.0
Octadecanoic acid	19.457	31.6	ng/ul	9948660	5	21.439	6299940	20.0
(DEL) Alkane: C...	21.127	8.8	ng/ul	2785400	5	21.439	6299940	20.0
(DEL) Alkane: S...	22.069	2.7	ng/ul	853321	5	21.439	6299940	20.0
(DEL) Alkane: S...	24.539	6.3	ng/ul	1999020	6	23.921	6316400	20.0
unknown-01	24.639	23.6	ng/ul	7468510	6	23.921	6316400	20.0
(DEL) Alkane: S...	26.456	19.7	ng/ul	6213940	6	23.921	6316400	20.0
.gamma.-Sitosterol	27.480	35.9	ng/ul	11345200	6	23.921	6316400	20.0
2-Amino-.alpha....	28.533	209.8	ng/ul	66248000	6	23.921	6316400	20.0