LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012357.D
Acq On : 27 Oct 2022 23:31
Operator : CG/JU
Sample : N5144-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0BK8

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-BP102522.M
Title : SVOA CALIBRATION

Signal : TIC: BP012357.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.246	41	44	57	rVB	63314	100414	1.08%	0.088%
2	3.676	113	117	119	rBV	20692	27929	0.30%	0.025%
3	3.699	119	121	129	rVV	22258	46276	0.50%	0.041%
4	3.770	129	133	141	rVV	103794	155966	1.67%	0.137%
5	3.846	143	146	148	rVV2	14921	20217	0.22%	0.018%
6	3.887	148	153	161	rVV	76217	128871	1.38%	0.113%
7	3.981	165	169	180	rVB	31595	53510	0.57%	0.047%
8	4.364	229	234	240	rBV3	19341	29908	0.32%	0.026%
9	4.546	260	265	271	rVB2	25378	39147	0.42%	0.034%
10	4.923	321	329	337	rBV2	187795	339799	3.64%	0.299%
11	5.440	411	417	423	rBV	16296	29225	0.31%	0.026%
12	5.811	473	480	486	rBV3	28794	49585	0.53%	0.044%
13	6.975	671	678	692	rVV	1131125	1963327	21.05%	1.727%
14	7.140	700	706	723	rBV	953268	1585754	17.00%	1.395%
15	7.334	731	739	749	rBV	1141757	1892664	20.29%	1.664%
16	7.417	749	753	755	rVV2	18824	28968	0.31%	0.025%
17	7.446	755	758	764	rVB	17378	25817	0.28%	0.023%
18	7.805	808	819	827	rBV	1123162	1816482	19.47%	1.597%
19	8.522	934	941	957	rBV	1118514	1916658	20.55%	1.686%
20	8.640	957	961	967	rVB	38059	63338	0.68%	0.056%
21	8.975	1010	1018	1031	rBV	1132012	1937622	20.77%	1.704%
22	9.358	1078	1083	1091	rVV2	35794	55619	0.60%	0.049%
23	9.699	1132	1141	1153	rBV	925746	1627277	17.44%	1.431%
24	10.234	1224	1232	1246	rVV	1472934	2602251	27.89%	2.288%
25	10.399	1253	1260	1273	rBV2	72697	185472	1.99%	0.163%
26	10.610	1285	1296	1302	rBV	1703957	2923940	31.34%	2.571%
27	10.658	1302	1304	1312	rVB	59323	80300	0.86%	0.071%
28	10.758	1314	1321	1341	rBV	863383	1640449	17.58%	1.443%
29	11.040	1362	1369	1375	rBV	19493	33563	0.36%	0.030%
30	11.628	1464	1469	1480	rBV6	12672	28787	0.31%	0.025%
31	12.022	1532	1536	1543	rVB	18221	27484	0.29%	0.024%
32	12.152	1549	1558	1561	rBV3	11969	28563	0.31%	0.025%
33	12.199	1561	1566	1572	rVV2	28183	49378	0.53%	0.043%
34	12.275	1572	1579	1587	rVB	72256	116091	1.24%	0.102%
35	12.493	1609	1616	1623	rVV	66020	107754	1.16%	0.095%
36	12.981	1693	1699	1706	rVB5	13312	23591	0.25%	0.021%

LSC Area Percent Report

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Operator : CG/JU
Sample : N5144-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BK8

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-BP102522.M
Title : SVOA CALIBRATION

37	13.269	1742	1748	1756	rBV3	43014	83711	0.90%	0.074%
38	13.351	1756	1762	1766	rVV2	14414	23869	0.26%	0.021%
39	13.422	1770	1774	1781	rVB8	10240	20353	0.22%	0.018%
40	13.616	1801	1807	1808	rBV3	16754	24322	0.26%	0.021%
41	13.775	1828	1834	1839	rBV2	46052	85653	0.92%	0.075%
42	13.834	1839	1844	1845	rBV	32642	45148	0.48%	0.040%
43	13.869	1845	1850	1857	rVV	2785788	3923736	42.06%	3.451%
44	14.010	1870	1874	1878	rBV	23634	26721	0.29%	0.023%
45	14.151	1887	1898	1916	rBV	3403557	5124009	54.93%	4.506%
46	14.346	1925	1931	1935	rBV	28400	42565	0.46%	0.037%
47	14.457	1940	1950	1969	rBV	2853250	4298342	46.07%	3.780%
48	14.675	1981	1987	2008	rBV	1096778	1999905	21.44%	1.759%
49	14.846	2010	2016	2017	rVV2	43974	81869	0.88%	0.072%
50	14.887	2020	2023	2028	rVB3	50981	81399	0.87%	0.072%
51	14.963	2035	2036	2047	rVB3	11176	21868	0.23%	0.019%
52	15.075	2051	2055	2061	rVV4	15056	26900	0.29%	0.024%
53	15.134	2061	2065	2070	rVB4	22934	31423	0.34%	0.028%
54	15.246	2078	2084	2087	rBV4	24222	49145	0.53%	0.043%
55	15.298	2091	2093	2099	rVB	33436	41688	0.45%	0.037%
56	15.457	2113	2120	2126	rBV	4373930	6477619	69.43%	5.697%
57	15.587	2136	2142	2153	rBV	1018765	1419712	15.22%	1.249%
58	15.693	2157	2160	2166	rVB5	21613	43185	0.46%	0.038%
59	15.804	2174	2179	2185	rBV3	26858	48439	0.52%	0.043%
60	15.951	2201	2204	2210	rVB2	25967	46789	0.50%	0.041%
61	16.040	2213	2219	2229	rVB7	15686	40490	0.43%	0.036%
62	16.187	2236	2244	2250	rBV3	72622	178062	1.91%	0.157%
63	16.363	2271	2274	2283	rVB10	10404	21924	0.24%	0.019%
64	16.616	2314	2317	2324	rBV5	21149	26542	0.28%	0.023%
65	16.757	2335	2341	2349	rBV3	78092	146748	1.57%	0.129%
66	16.875	2357	2361	2367	rBV5	18558	32675	0.35%	0.029%
67	16.969	2369	2377	2380	rBV2	39005	74634	0.80%	0.066%
68	17.110	2397	2401	2408	rBV4	56966	95304	1.02%	0.084%
69	17.216	2409	2419	2424	rBV	3683869	5361111	57.47%	4.715%
70	17.257	2424	2426	2430	rVV	182274	237176	2.54%	0.209%
71	17.316	2430	2436	2446	rVB2	4650448	6766668	72.53%	5.951%
72	17.710	2500	2503	2506	rBV2	27836	32332	0.35%	0.028%
73	17.851	2524	2527	2530	rBV4	18608	28085	0.30%	0.025%
74	17.887	2530	2533	2538	rVB	54788	67392	0.72%	0.059%
75	17.992	2546	2551	2557	rBV2	139908	221415	2.37%	0.195%
76	18.051	2557	2561	2564	rBV3	93117	133276	1.43%	0.117%

LSC Area Percent Report

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Operator : CG/JU
Sample : N5144-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BK8

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Title : SVOA CALIBRATION

77	18.104	2564	2570	2580	rBV2	1141784	1830813	19.62%	1.610%
78	18.275	2595	2599	2602	rBV2	46823	72775	0.78%	0.064%
79	18.392	2614	2619	2623	rBV5	68071	134952	1.45%	0.119%
80	18.434	2623	2626	2633	rVB2	38412	65634	0.70%	0.058%
81	18.581	2648	2651	2655	rVB2	46990	64068	0.69%	0.056%
82	19.204	2753	2757	2758	rBV	163563	199399	2.14%	0.175%
83	19.228	2758	2761	2775	rVB	544392	1001122	10.73%	0.880%
84	19.339	2776	2780	2794	rBV2	679846	1116670	11.97%	0.982%
85	19.557	2812	2817	2830	rVB	6532666	9329066	100.00%	8.204%
86	20.069	2902	2904	2910	rVB3	200854	247335	2.65%	0.218%
87	20.557	2984	2987	2991	rVB	194020	205338	2.20%	0.181%
88	21.022	3061	3066	3075	rVB3	1257826	2258363	24.21%	1.986%
89	21.216	3096	3099	3103	rBV	298856	310467	3.33%	0.273%
90	21.310	3109	3115	3120	rBV	5047467	6947981	74.48%	6.110%
91	21.445	3136	3138	3150	rVB	271308	390202	4.18%	0.343%
92	21.657	3171	3174	3183	rVB	682144	988545	10.60%	0.869%
93	21.904	3213	3216	3219	rBV	432996	495846	5.32%	0.436%
94	21.939	3220	3222	3231	rVB2	417313	703627	7.54%	0.619%
95	22.651	3339	3343	3350	rVB	1162651	1694808	18.17%	1.490%
96	22.957	3390	3395	3400	rBV2	812955	1479651	15.86%	1.301%
97	23.022	3400	3406	3425	rVB	2869462	7336297	78.64%	6.452%
98	23.551	3489	3496	3511	rVB2	3920641	8849894	94.86%	7.783%
99	23.710	3516	3523	3541	rVB2	2828062	5809350	62.27%	5.109%
100	24.298	3617	3623	3632	rVB	1370734	2865487	30.72%	2.520%

Sum of corrected areas: 113711890

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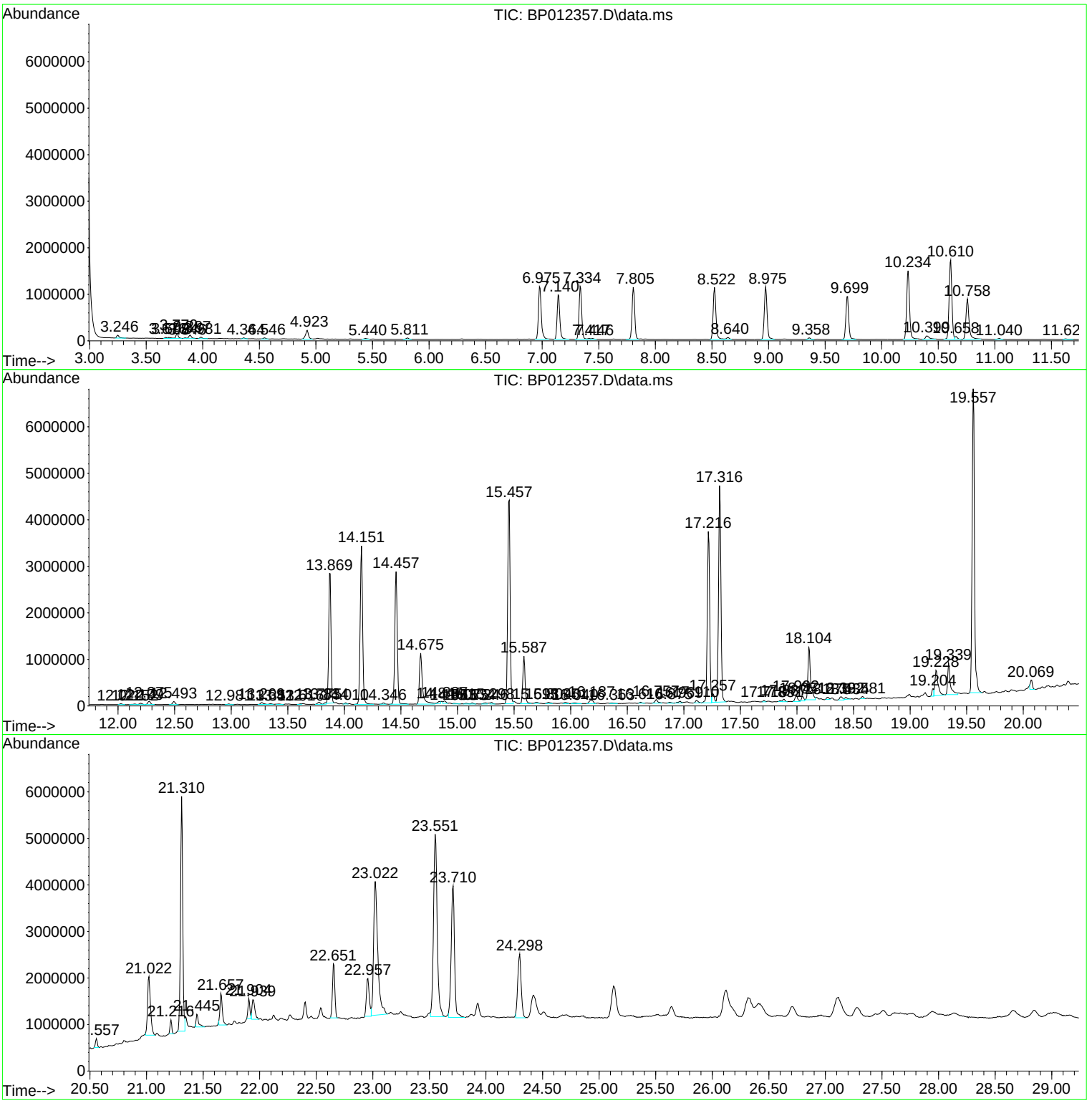
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BK8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012357.D
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Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BK8

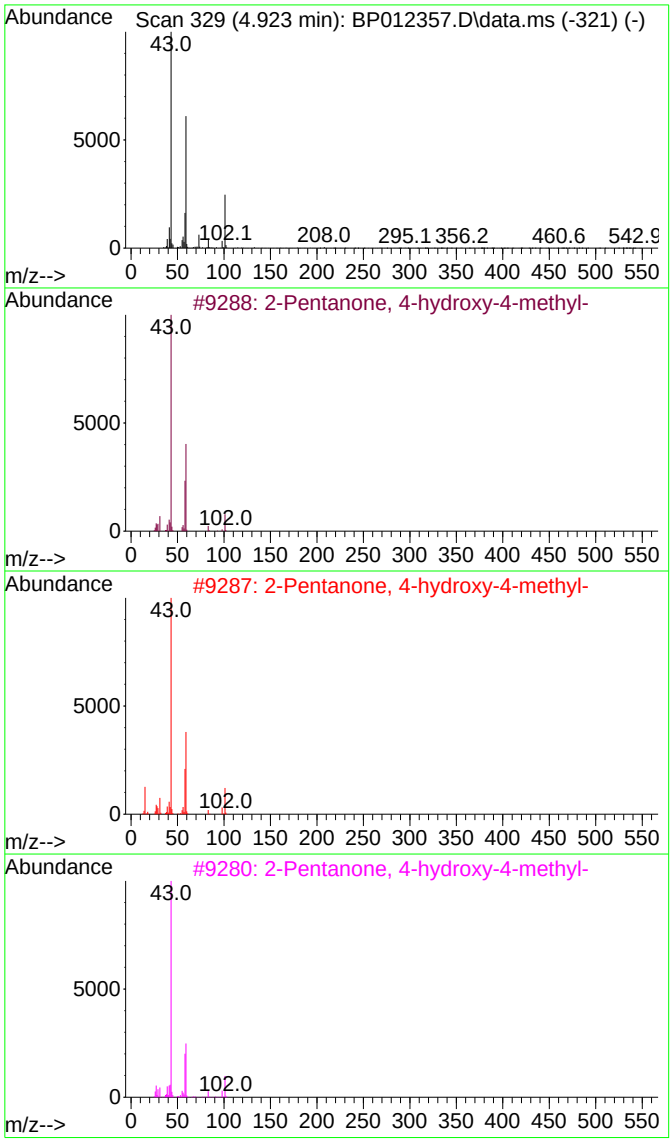
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

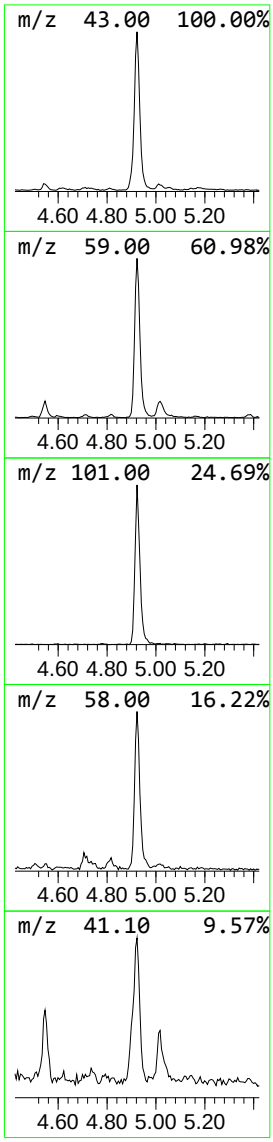
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.923	3.74 ng/ul	339799	1,4-Dichlorobenzene-d4	7.805

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
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	38
5	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	28





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012357.D
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ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

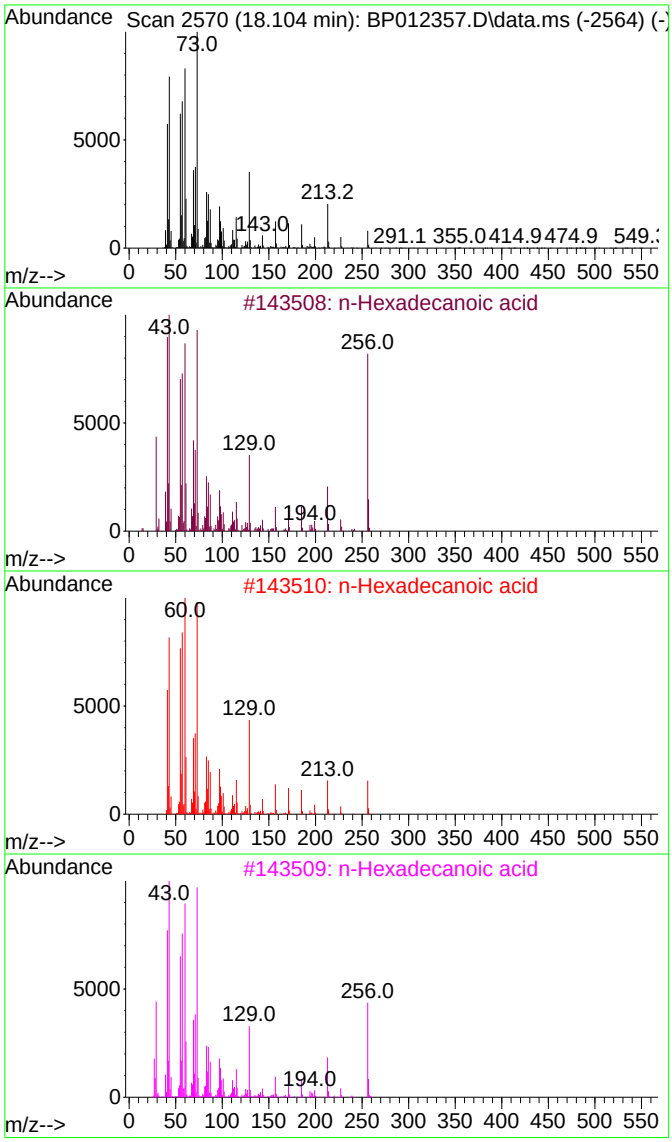
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

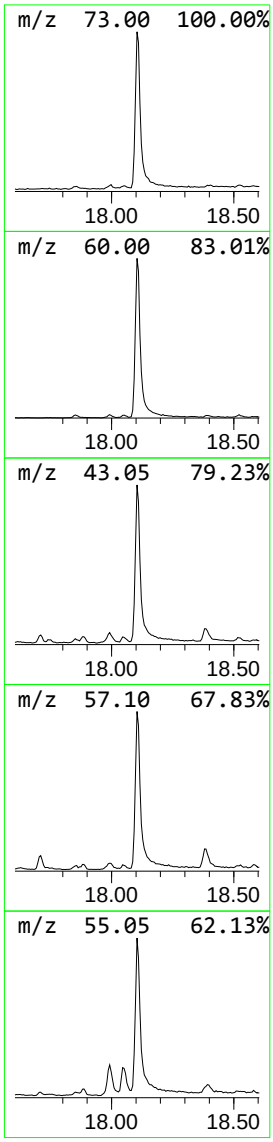
****

Peak Number 2 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.104	6.83 ng/ul	1830810	Phenanthrene-d10	17.216

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	95
5	Tridecanoic acid	214	C13H26O2	000638-53-9	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
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Acq On : 27 Oct 2022 23:31
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

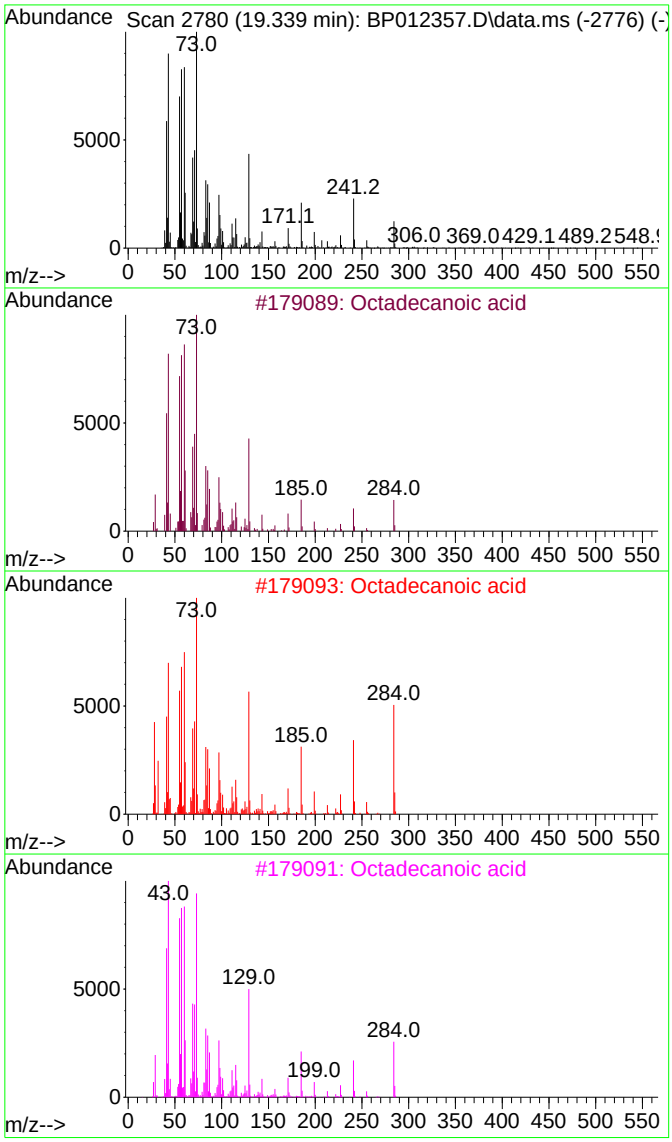
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

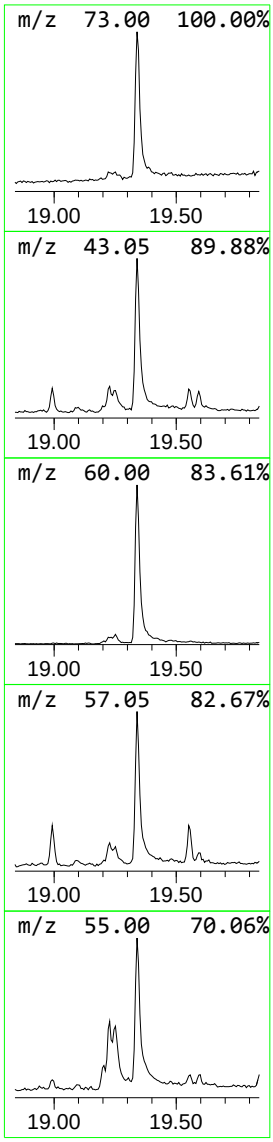
****

Peak Number 3 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.339	3.21 ng/ul	1116670	Chrysene-d12	21.310

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	95
5	Octadecanoic acid	284	C18H36O2	000057-11-4	94





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Sample : N5144-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BK8

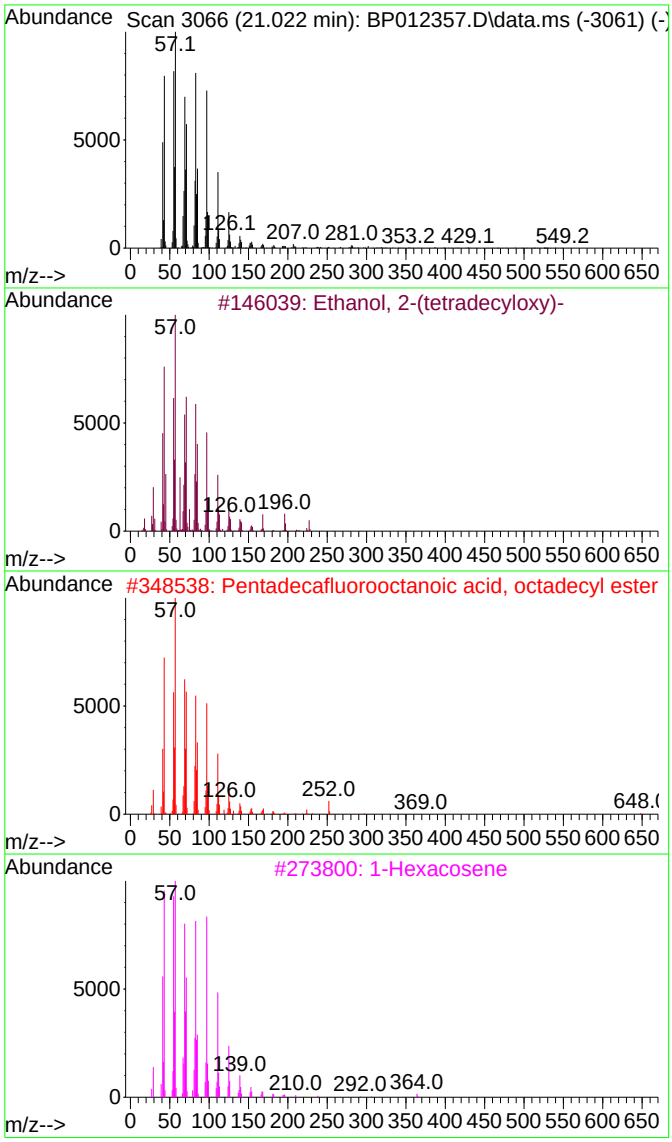
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

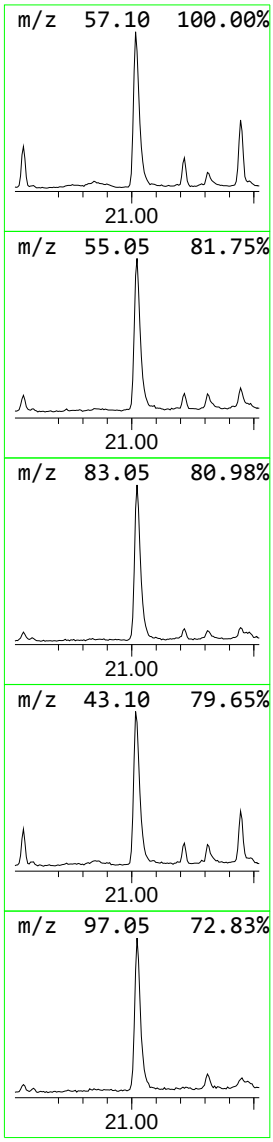
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Ethanol, 2-(tetradecyloxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.022	6.50 ng/ul	2258360	Chrysene-d12	21.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	93
2	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
3	1-Hexacosene	364	C26H52	018835-33-1	91
4	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
5	Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012357.D
Acq On : 27 Oct 2022 23:31
Operator : CG/JU
Sample : N5144-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BK8

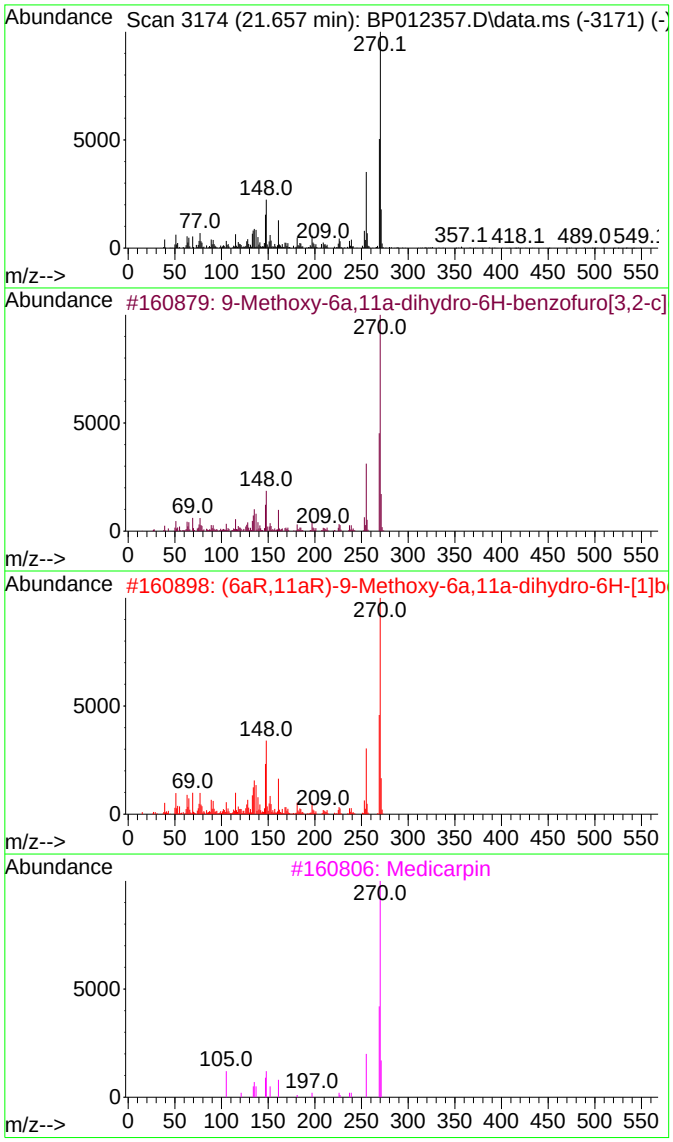
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

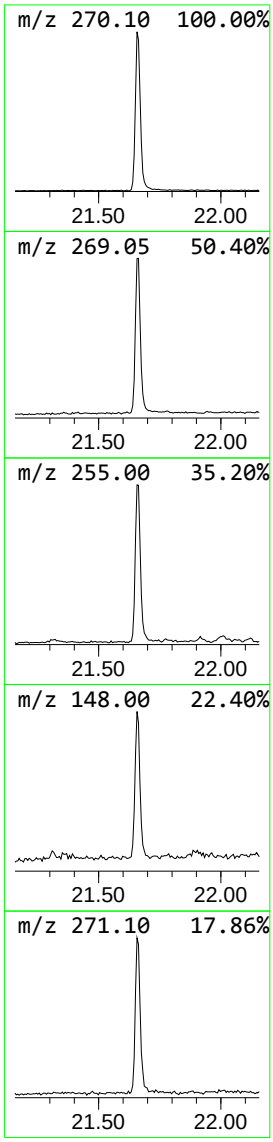
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 9-Methoxy-6a,11a-dihydro-6H... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.657	2.85 ng/ul	988545	Chrysene-d12	21.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Methoxy-6a,11a-dihydro-6H-benz...	270	C16H14O4	607363-34-8	99
2	(6aR,11aR)-9-Methoxy-6a,11a-dihy...	270	C16H14O4	033983-40-3	98
3	Medicarpin	270	C16H14O4	032383-76-9	91
4	3-Acetoxy-9-methoxypterocarpan	312	C18H16O5	001891-12-9	86
5	4-Azafluorenone, 3-methylphenyli...	270	C19H14N2	1000301-74-7	64





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Acq On : 27 Oct 2022 23:31
Operator : CG/JU
Sample : N5144-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BK8

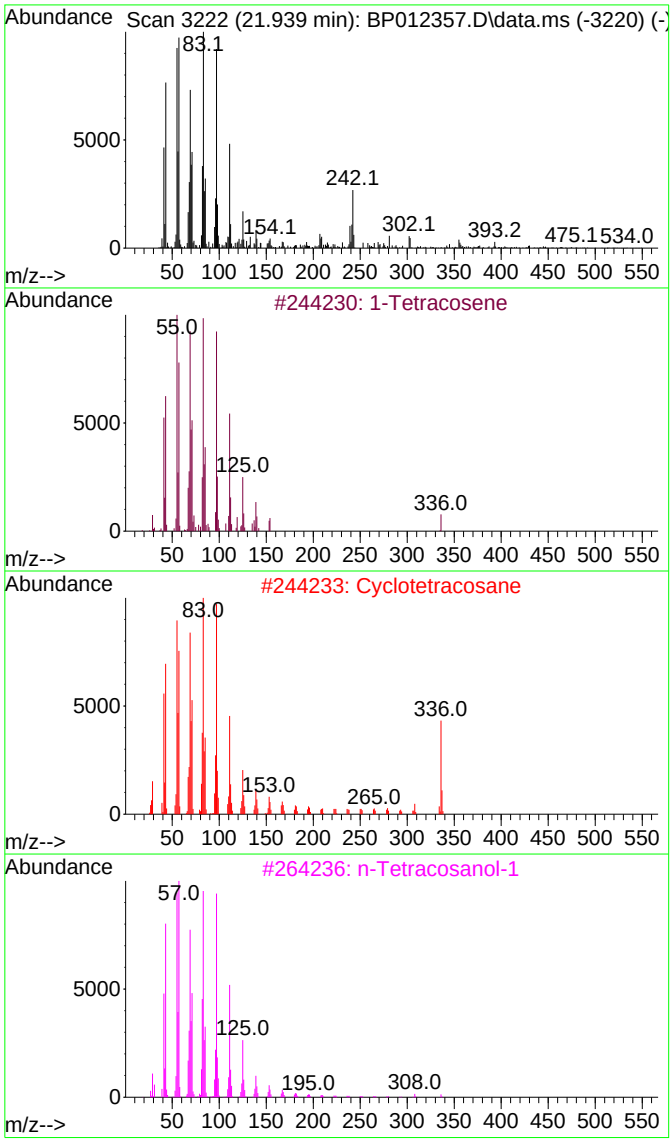
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

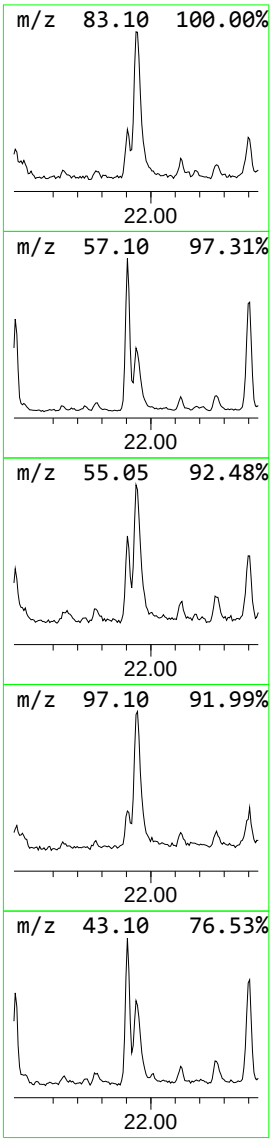
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.939	2.03 ng/ul	703627	Chrysene-d12	21.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene	336	C24H48	010192-32-2	99
2	Cyclotetracosane	336	C24H48	000297-03-0	98
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4	1-Heneicosanol	312	C21H44O	015594-90-8	93
5	1-Heptacosanol	396	C27H56O	002004-39-9	91





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Data File : BP012357.D
Acq On : 27 Oct 2022 23:31
Operator : CG/JU
Sample : N5144-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BK8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

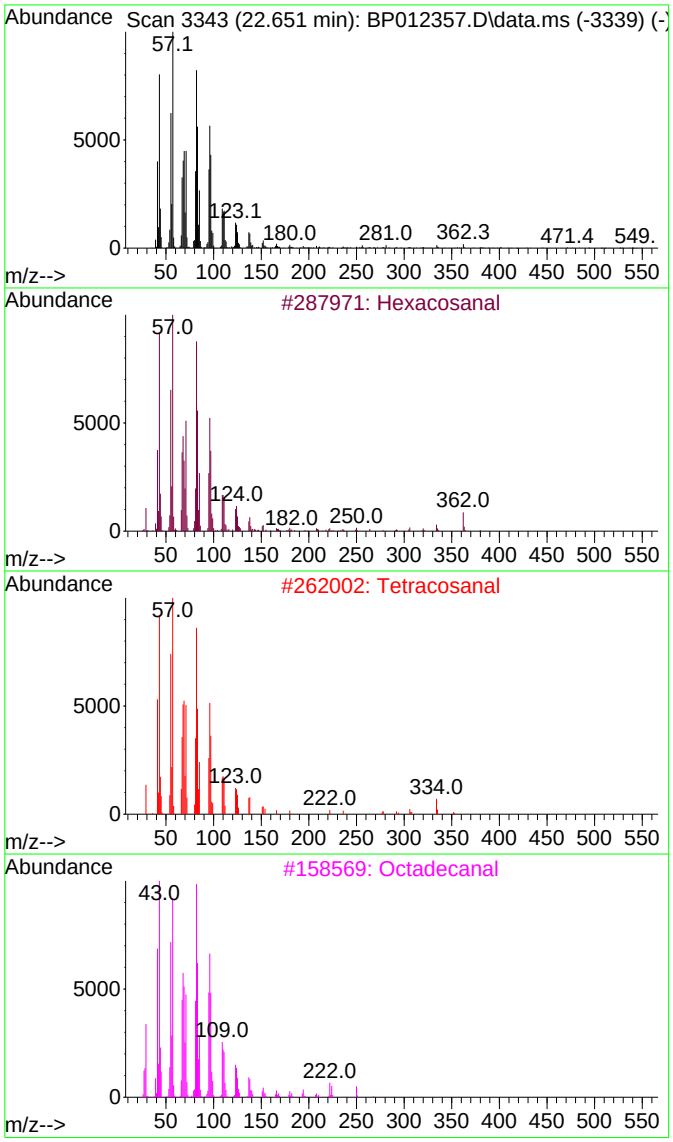
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

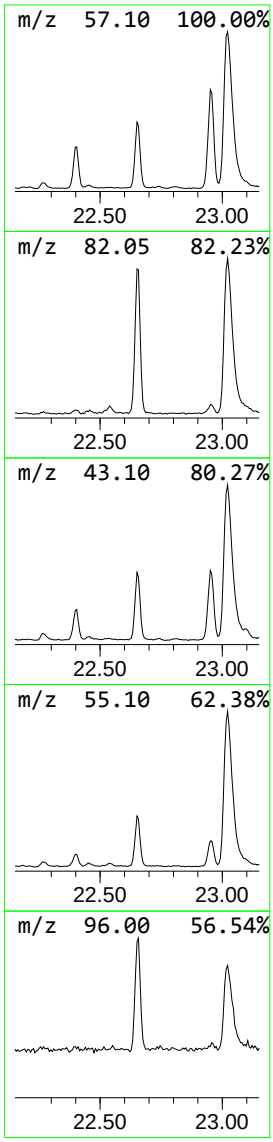
****

Peak Number 7 Hexacosanal Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.651	5.83 ng/ul	1694810	Perylene-d12	23.710

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosanal	380	C26H52O	026627-85-0	97
2	Tetracosanal	352	C24H48O	057866-08-7	94
3	Octadecanal	268	C18H36O	000638-66-4	91
4	Octadecanal	268	C18H36O	000638-66-4	91
5	Octadecanal	268	C18H36O	000638-66-4	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012357.D
Acq On : 27 Oct 2022 23:31
Operator : CG/JU
Sample : N5144-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BK8

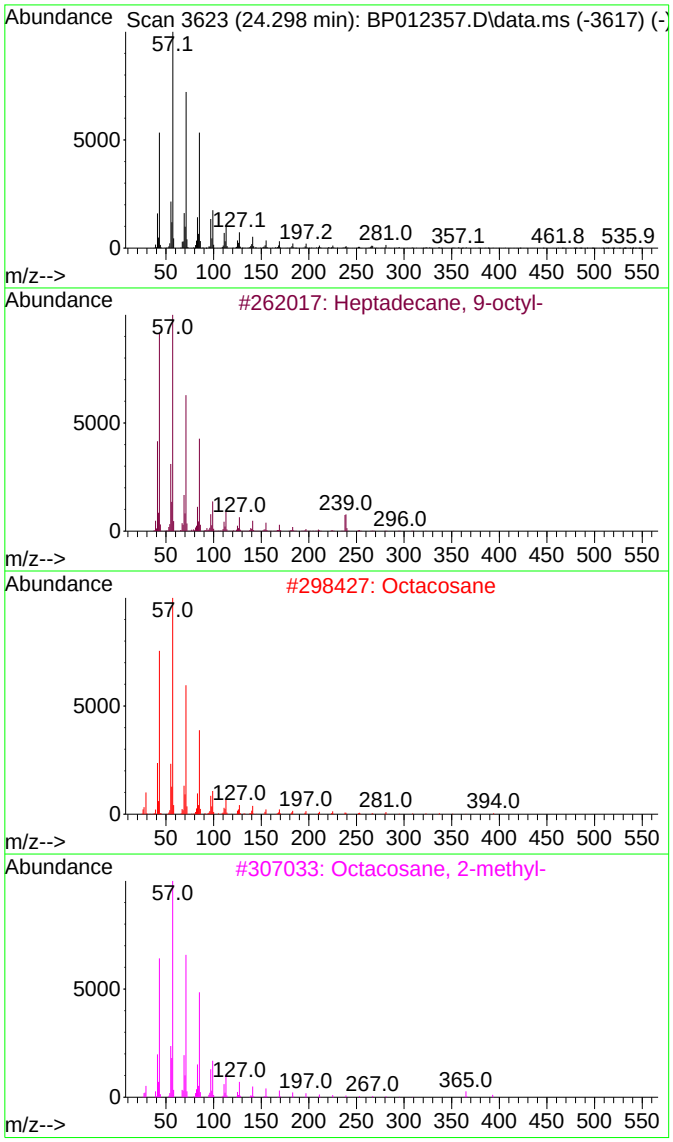
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

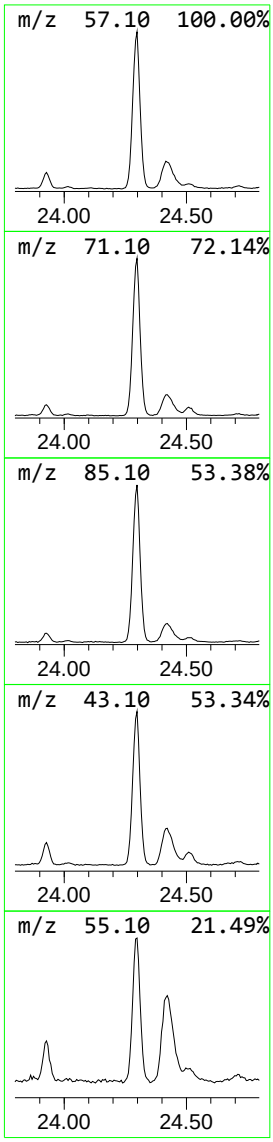
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.298	9.87 ng/ul	2865490	Perylene-d12	23.710

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2	Octacosane	394	C28H58	000630-02-4	91
3	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	4-Methylheneicosane	310	C22H46	025117-29-7	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012357.D
Acq On : 27 Oct 2022 23:31
Operator : CG/JU
Sample : N5144-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BK8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.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-...	4.923	3.7	ng/ul	339799	1	7.805	1816480	20.0
n-Hexadecanoic ...	18.104	6.8	ng/ul	1830810	4	17.216	5361110	20.0
Octadecanoic acid	19.339	3.2	ng/ul	1116670	5	21.310	6947980	20.0
Ethanol, 2-(tet...	21.022	6.5	ng/ul	2258360	5	21.310	6947980	20.0
9-Methoxy-6a,11...	21.657	2.9	ng/ul	988545	5	21.310	6947980	20.0
(DEL) Alkane: S...	21.939	2.0	ng/ul	703627	5	21.310	6947980	20.0
Hexacosanal	22.651	5.8	ng/ul	1694810	6	23.710	5809350	20.0
(DEL) Alkane: S...	24.298	9.9	ng/ul	2865490	6	23.710	5809350	20.0