LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022138.D
Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DDCG1

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-BP092424.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP022138.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.222	36	40	47	rVB	22918	34765	0.82%	0.054%
2	3.546	85	95	104	rVB	51000	93933	2.22%	0.145%
3	3.640	106	111	122	rBV	223217	365063	8.63%	0.563%
4	3.928	143	160	163	rBV	158396	488069	11.54%	0.753%
5	4.869	314	320	327	rVV	423210	581776	13.76%	0.897%
6	5.634	443	450	455	rVB4	16734	31079	0.73%	0.048%
7	5.787	470	476	481	rBV2	17460	26262	0.62%	0.040%
8	6.075	515	525	530	rBV	1205761	2051522	48.52%	3.163%
9	6.116	530	532	535	rVV	62580	77560	1.83%	0.120%
10	6.145	535	537	546	rVV	29639	39169	0.93%	0.060%
11	6.234	546	552	560	rVB3	14209	26370	0.62%	0.041%
12	6.745	633	639	644	rVB	20321	35873	0.85%	0.055%
13	6.892	655	664	674	rBV	471841	823015	19.46%	1.269%
14	7.040	677	689	695	rVV	327107	539889	12.77%	0.832%
15	7.245	717	724	733	rVB	478451	744647	17.61%	1.148%
16	7.440	747	757	761	rBV5	11870	31102	0.74%	0.048%
17	7.504	761	768	774	rVV	54734	103294	2.44%	0.159%
18	7.698	785	801	808	rBV2	512328	1082061	25.59%	1.668%
19	7.769	808	813	819	rVB2	45334	74937	1.77%	0.116%
20	7.934	828	841	853	rVB	376400	761647	18.01%	1.174%
21	8.039	853	859	868	rBV3	14523	30996	0.73%	0.048%
22	8.128	868	874	878	rVV2	59308	101534	2.40%	0.157%
23	8.163	878	880	886	rVB2	16775	25441	0.60%	0.039%
24	8.328	904	908	914	rVB	46739	75961	1.80%	0.117%
25	8.404	914	921	932	rBV	561343	984959	23.29%	1.519%
26	8.504	932	938	944	rVB	36893	62529	1.48%	0.096%
27	8.581	944	951	960	rBV	49304	85810	2.03%	0.132%
28	8.839	985	995	1007	rBV	468943	843174	19.94%	1.300%
29	9.075	1015	1035	1039	rBV	694405	2438330	57.66%	3.760%
30	9.116	1039	1042	1050	rVV3	27568	57767	1.37%	0.089%
31	9.357	1079	1083	1086	rVV2	24694	38895	0.92%	0.060%
32	9.404	1086	1091	1101	rVB3	81844	160706	3.80%	0.248%
33	9.551	1108	1116	1123	rBV3	428230	835750	19.76%	1.289%
34	9.628	1123	1129	1134	rBV	139764	220854	5.22%	0.341%
35	9.722	1140	1145	1150	rBV	35175	54767	1.30%	0.084%
36	9.975	1182	1188	1199	rVB	298330	514026	12.16%	0.793%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022138.D
Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

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-BP092424.MA.M
Title : SVOA CALIBRATION

37	10.098	1202	1209	1215	rBV	696666	1088278	25.74%	1.678%
38	10.233	1223	1232	1239	rVB4	35822	97218	2.30%	0.150%
39	10.351	1244	1252	1259	rBV	350955	568584	13.45%	0.877%
40	10.469	1265	1272	1278	rVV	702855	1139090	26.94%	1.756%
41	10.598	1287	1294	1301	rBV	410324	790452	18.69%	1.219%
42	10.669	1301	1306	1313	rVV	96405	181103	4.28%	0.279%
43	10.769	1313	1323	1333	rVV	138158	276888	6.55%	0.427%
44	10.998	1357	1362	1367	rVB4	21976	36875	0.87%	0.057%
45	11.369	1418	1425	1431	rBV2	57879	97349	2.30%	0.150%
46	11.586	1452	1462	1465	rBV4	24513	55751	1.32%	0.086%
47	11.622	1465	1468	1473	rVB3	32213	48271	1.14%	0.074%
48	12.004	1523	1533	1538	rBV5	44950	122895	2.91%	0.189%
49	12.222	1565	1570	1578	rVB4	45531	74580	1.76%	0.115%
50	12.392	1588	1599	1604	rBV5	25025	57374	1.36%	0.088%
51	12.563	1620	1628	1629	rBV3	18724	31810	0.75%	0.049%
52	12.592	1629	1633	1643	rVV3	43781	88135	2.08%	0.136%
53	12.722	1650	1655	1664	rVB	59274	94850	2.24%	0.146%
54	12.810	1664	1670	1677	rBV	11430	28628	0.68%	0.044%
55	13.057	1704	1712	1720	rVB2	91521	156477	3.70%	0.241%
56	13.163	1722	1730	1735	rVB4	25462	53322	1.26%	0.082%
57	13.263	1735	1747	1751	rBV	85551	154509	3.65%	0.238%
58	13.345	1757	1761	1770	rVB5	23355	48966	1.16%	0.075%
59	13.569	1794	1799	1809	rVB	157527	252190	5.96%	0.389%
60	13.727	1819	1826	1834	rVB	992583	1545031	36.54%	2.382%
61	13.822	1838	1842	1847	rBV	28526	41739	0.99%	0.064%
62	14.004	1866	1873	1880	rBV	1263836	1830027	43.28%	2.822%
63	14.074	1880	1885	1893	rVB2	38097	83689	1.98%	0.129%
64	14.310	1918	1925	1932	rBV	1151103	1662918	39.33%	2.564%
65	14.533	1955	1963	1971	rBV	560562	802575	18.98%	1.237%
66	14.627	1971	1979	1984	rVV3	71478	170435	4.03%	0.263%
67	14.869	2015	2020	2025	rBV4	32449	57554	1.36%	0.089%
68	14.945	2025	2033	2044	rVB	227820	384556	9.09%	0.593%
69	15.316	2090	2096	2106	rVB	1515350	2350735	55.59%	3.624%
70	15.445	2111	2118	2129	rBV2	341693	780290	18.45%	1.203%
71	15.680	2152	2158	2165	rVB	725268	1001009	23.67%	1.543%
72	16.163	2236	2240	2247	rVB	98574	137221	3.25%	0.212%
73	16.274	2254	2259	2262	rBV	76845	126181	2.98%	0.195%
74	16.516	2295	2300	2303	rVV5	18687	29494	0.70%	0.045%
75	16.751	2334	2340	2351	rBV	2999666	4228508	100.00%	6.520%
76	17.110	2395	2401	2411	rVB	1451469	2160178	51.09%	3.331%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022138.D
Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

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-BP092424.MA.M
Title : SVOA CALIBRATION

77	17.210	2411	2418	2426	rVB	1808688	2604210	61.59%	4.015%
78	17.745	2505	2509	2518	rVB	135792	199361	4.71%	0.307%
79	18.039	2554	2559	2568	rBV	608974	990497	23.42%	1.527%
80	18.139	2571	2576	2589	rVB2	230963	440429	10.42%	0.679%
81	18.333	2599	2609	2612	rBV2	680544	1351599	31.96%	2.084%
82	18.362	2612	2614	2616	rVV2	242245	299742	7.09%	0.462%
83	18.562	2646	2648	2654	rBV5	51051	87506	2.07%	0.135%
84	19.139	2743	2746	2749	rBV3	81876	98911	2.34%	0.153%
85	19.204	2752	2757	2759	rBV3	88120	139703	3.30%	0.215%
86	19.233	2760	2762	2765	rBV3	56031	59277	1.40%	0.091%
87	19.374	2782	2786	2794	rVB	193464	273702	6.47%	0.422%
88	19.586	2817	2822	2832	rBV	2180633	3124927	73.90%	4.818%
89	20.133	2913	2915	2918	rVB2	81665	65586	1.55%	0.101%
90	20.574	2987	2990	2992	rBV	81920	99649	2.36%	0.154%
91	20.674	3002	3007	3014	rBV	1167178	2674552	63.25%	4.124%
92	21.104	3076	3080	3082	rBV3	129472	185463	4.39%	0.286%
93	21.156	3084	3089	3095	rVV	419403	910210	21.53%	1.403%
94	21.545	3149	3155	3161	rBV	1573936	2451424	57.97%	3.780%
95	21.827	3201	3203	3212	rVB2	435648	668228	15.80%	1.030%
96	22.203	3260	3267	3271	rBV6	423209	1071585	25.34%	1.652%
97	22.356	3285	3293	3299	rBV	1166177	3065188	72.49%	4.726%
98	23.380	3463	3467	3471	rBV	258898	348375	8.24%	0.537%
99	24.609	3667	3676	3687	rVB	1209057	3467098	81.99%	5.346%
100	24.833	3707	3714	3725	rVB	983451	2700570	63.87%	4.164%

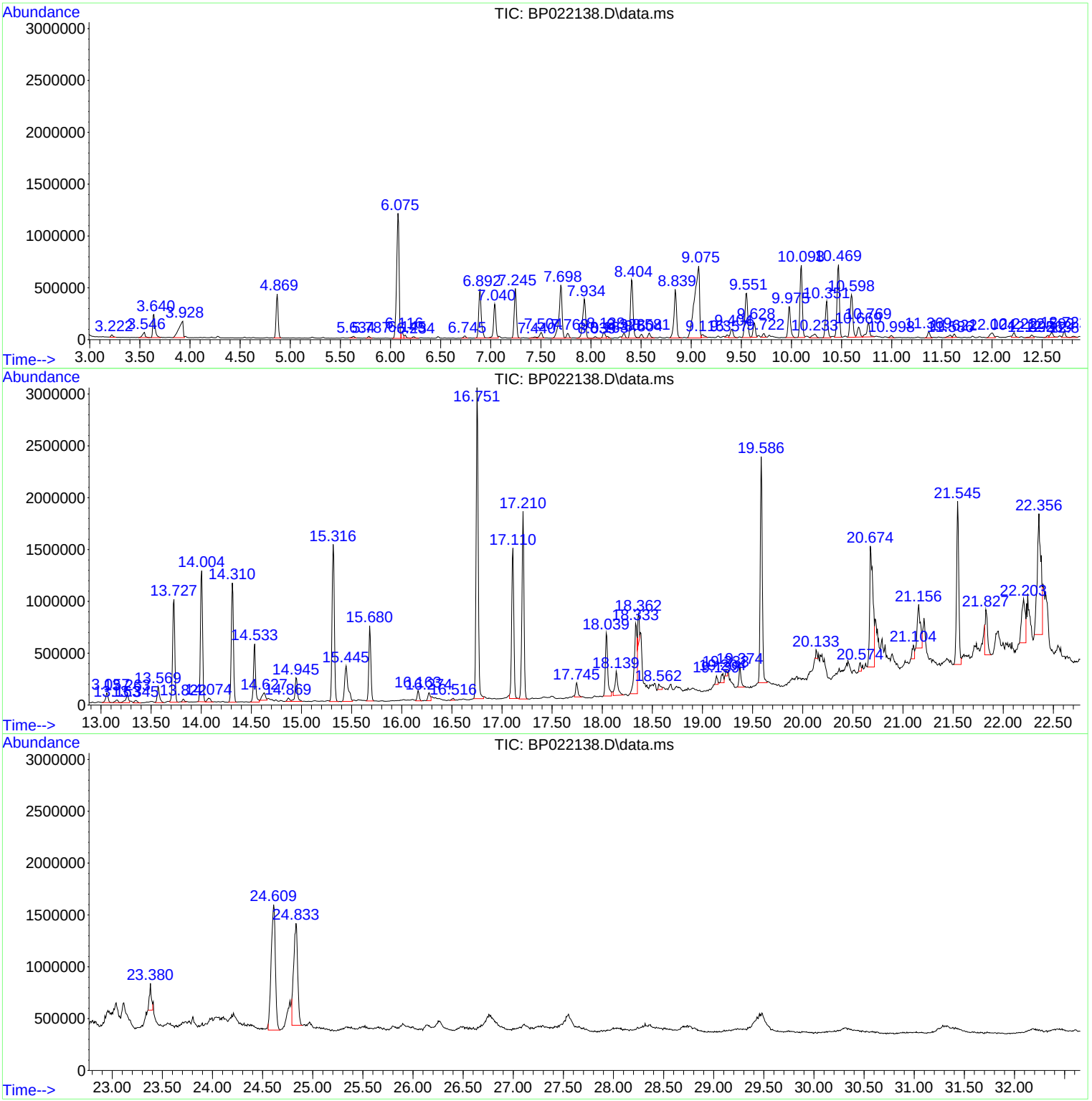
Sum of corrected areas: 64857059

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022138.D
Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022138.D
Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

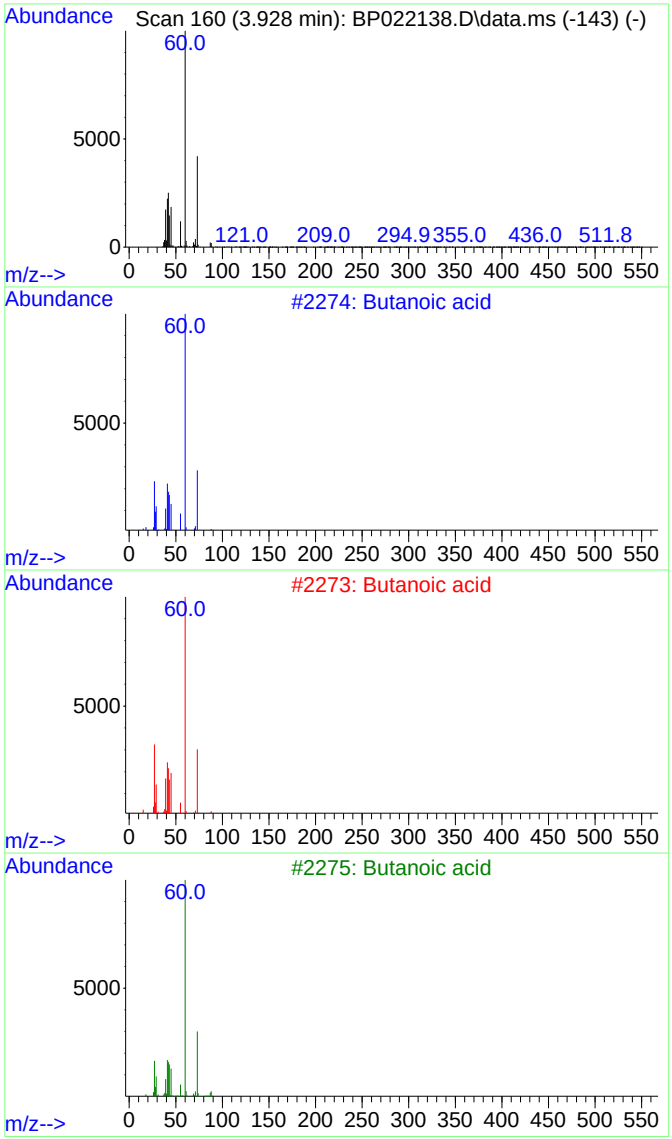
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

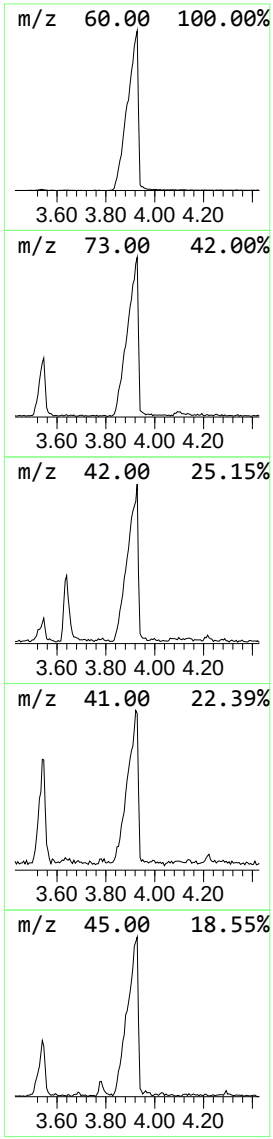
****

Peak Number 1 Butanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.928	9.02 ng/ul	488069	1,4-Dichlorobenzene-d4	7.698

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid	88	C4H8O2	000107-92-6	91
2	Butanoic acid	88	C4H8O2	000107-92-6	91
3	Butanoic acid	88	C4H8O2	000107-92-6	83
4	Butanoic acid	88	C4H8O2	000107-92-6	78
5	Pentanoic acid	102	C5H10O2	000109-52-4	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022138.D
Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

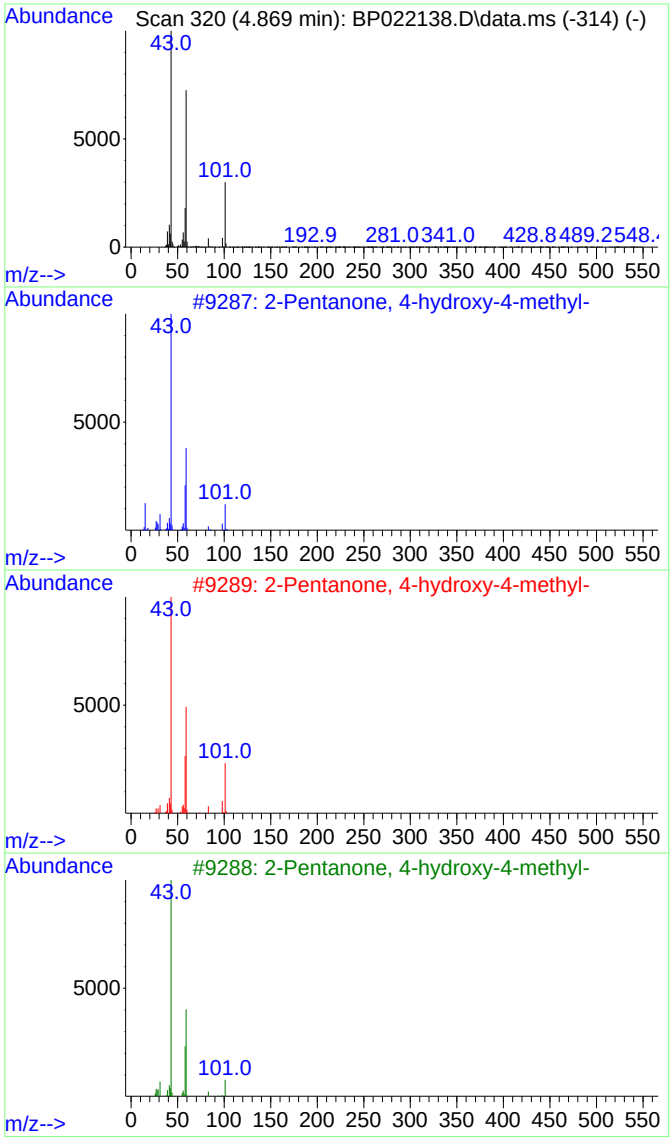
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

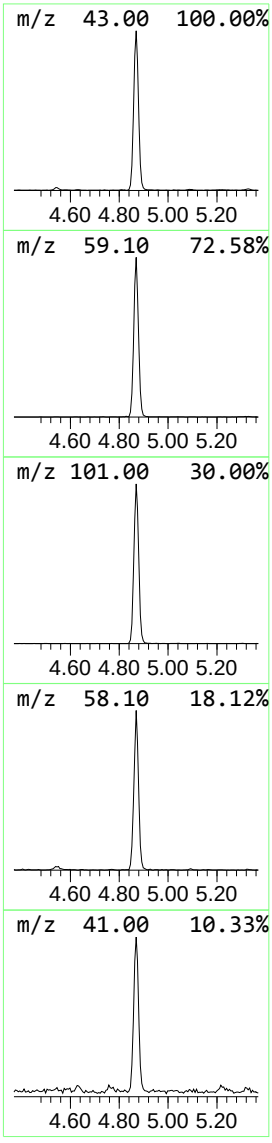
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.869	10.75 ng/ul	581776	1,4-Dichlorobenzene-d4	7.698

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
4	(+-)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	47
5	Guanidine	59	CH5N3	000113-00-8	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022138.D
Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

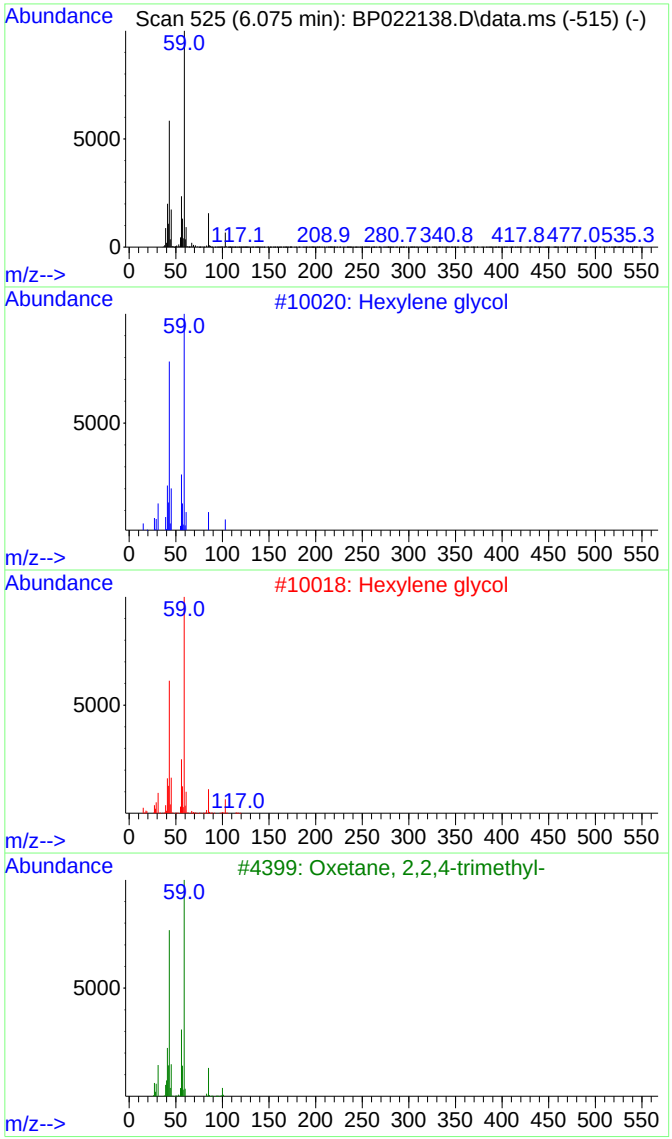
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

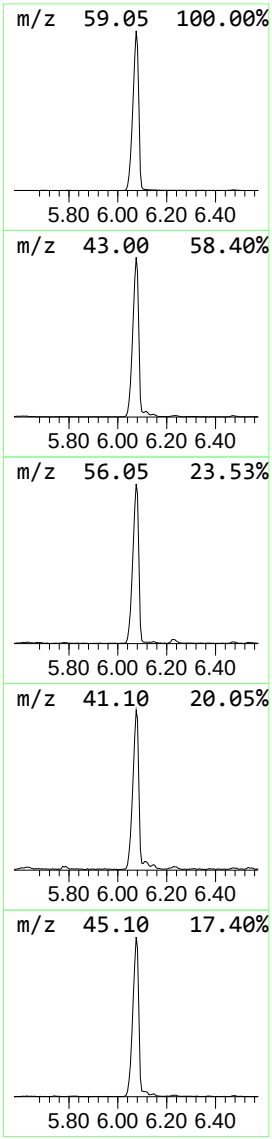
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Hexylene glycol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.075	37.92 ng/ul	2051520	1,4-Dichlorobenzene-d4	7.698

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexylene glycol	118	C6H14O2	000107-41-5	90
2	Hexylene glycol	118	C6H14O2	000107-41-5	83
3	Oxetane, 2,2,4-trimethyl-	100	C6H12O	023120-44-7	59
4	Hexylene glycol	118	C6H14O2	000107-41-5	53
5	(R)-(-)-2-Methyl-2,4-pentanediol	118	C6H14O2	099210-90-9	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022138.D
Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

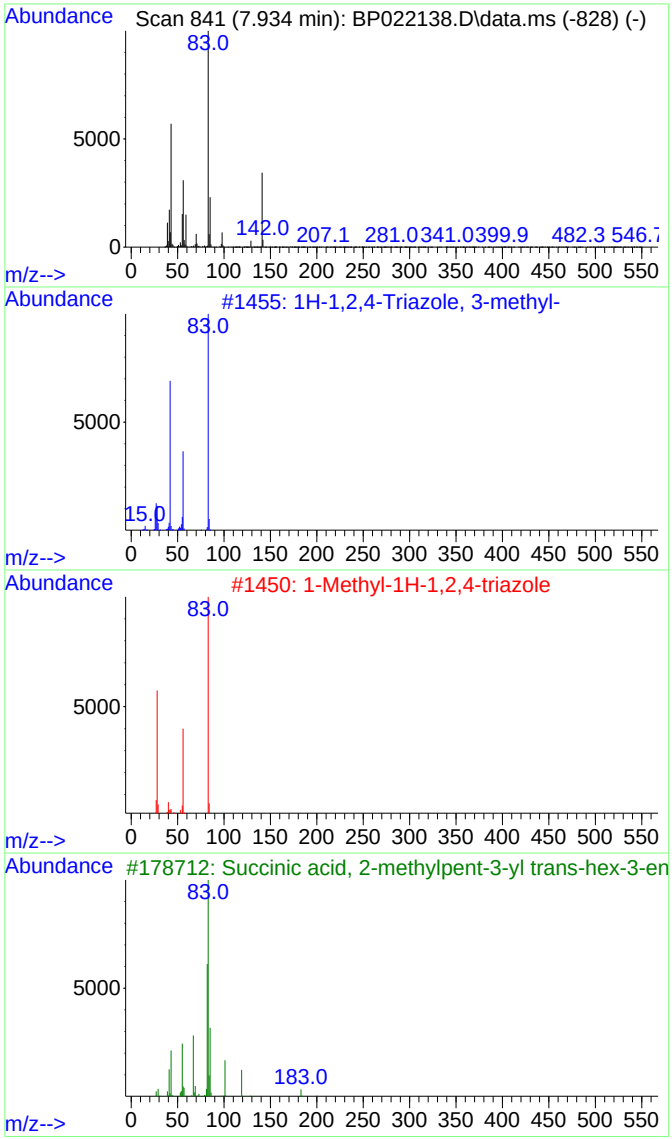
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

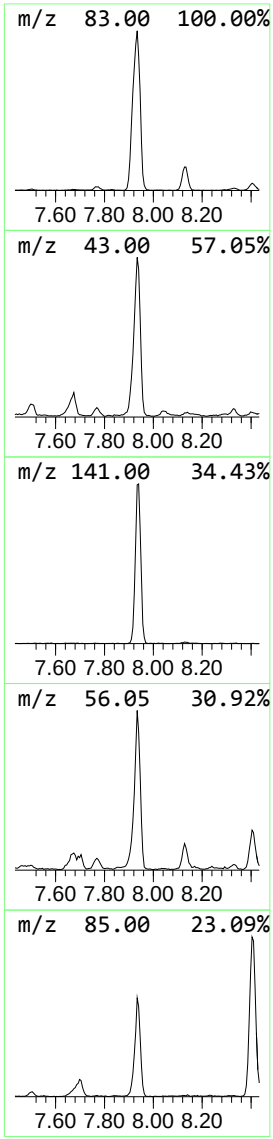
****

Peak Number 4 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.934	14.08 ng/ul	761647	1,4-Dichlorobenzene-d4	7.698

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-1,2,4-Triazole, 3-methyl-	83	C3H5N3	007170-01-6	38
2	1-Methyl-1H-1,2,4-triazole	83	C3H5N3	006086-21-1	38
3	Succinic acid, 2-methylpent-3-yl...	284	C16H28O4	1000391-11-2	12
4	1-Pentanol, 2,3-dimethyl-	116	C7H16O	010143-23-4	10
5	(Trimethylsilyl)acetylene	98	C5H10Si	001066-54-2	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022138.D
Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

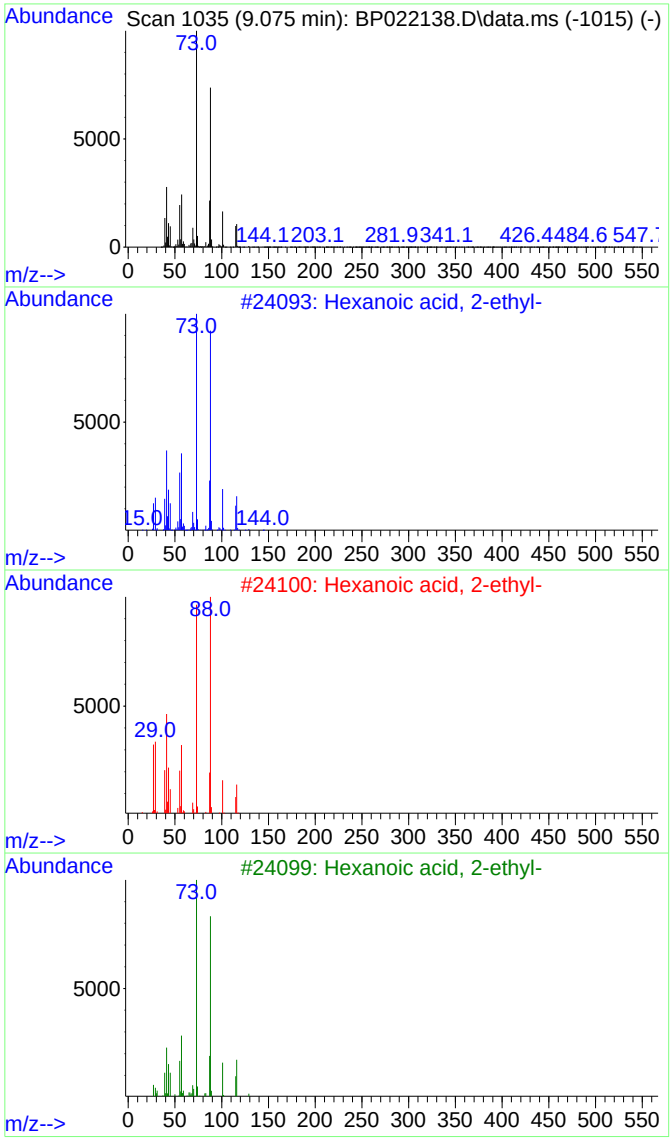
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

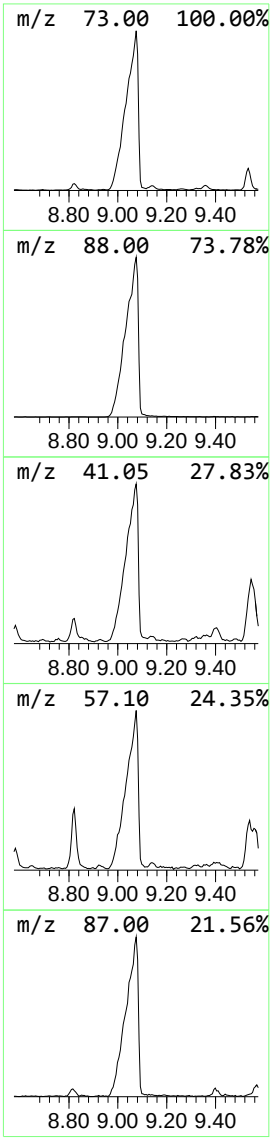
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Hexanoic acid, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.075	45.07 ng/ul	2438330	1,4-Dichlorobenzene-d4	7.698

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
3	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
4	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	72
5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	53





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Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

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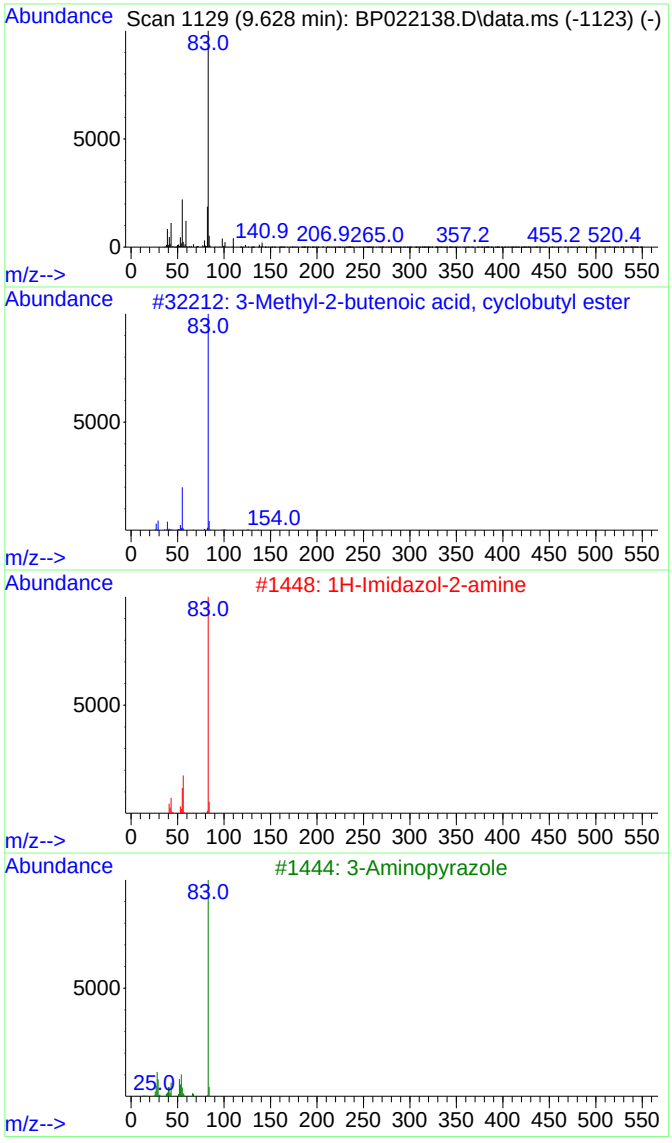
Quant Title : SVOA CALIBRATION

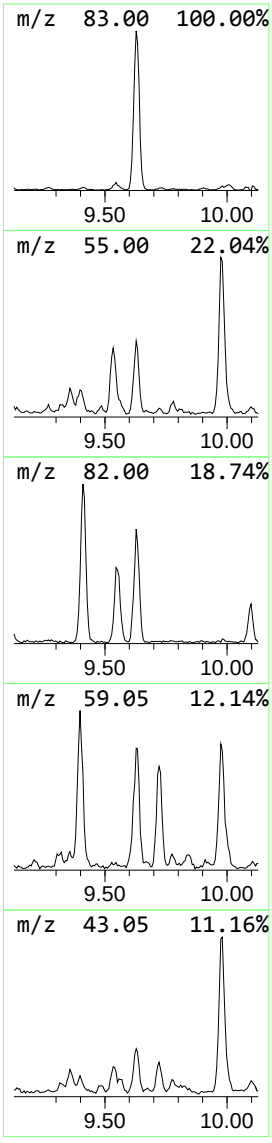
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 3-Methyl-2-butenoic acid, c... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.628	3.88 ng/ul	220854	Naphthalene-d8	10.469

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methyl-2-butenoic acid, cyclob...	154	C9H14O2	1000282-89-1	58
2	1H-Imidazol-2-amine	83	C3H5N3	007720-39-0	53
3	3-Aminopyrazole	83	C3H5N3	001820-80-0	50
4	2,6-Dimethyl-6-nitro-2-hepten-4-one	185	C9H15NO3	073583-56-9	45
5	N,N'-Bis(2,6-dimethyl-6-nitrosoh...	338	C18H30N2O4	066737-12-0	42





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Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

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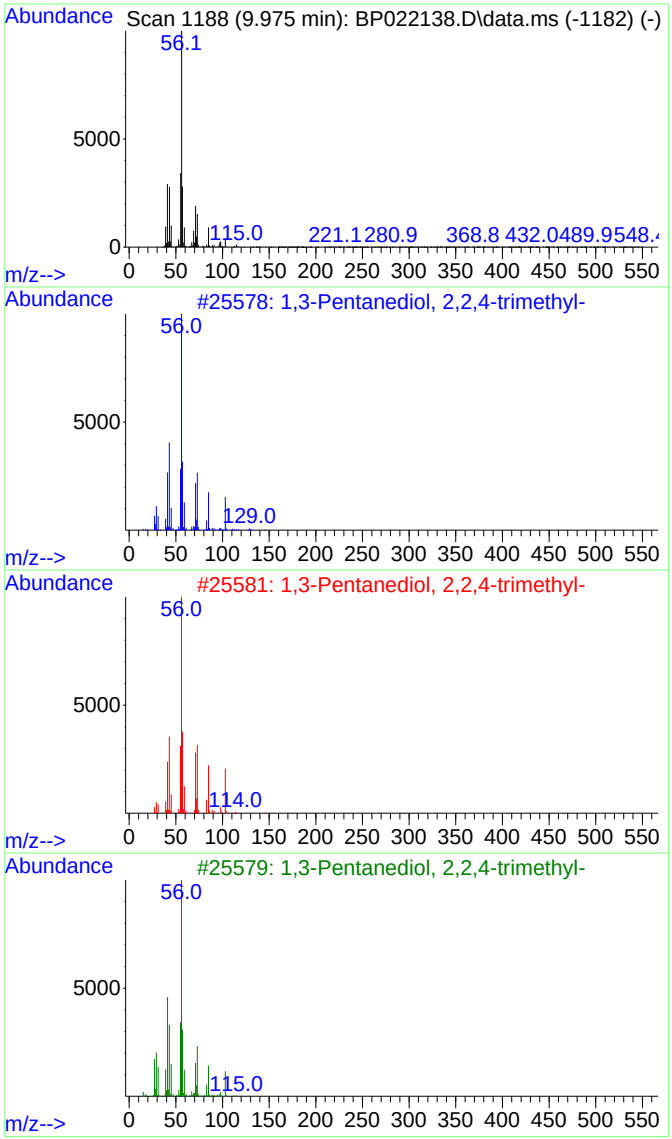
Quant Title : SVOA CALIBRATION

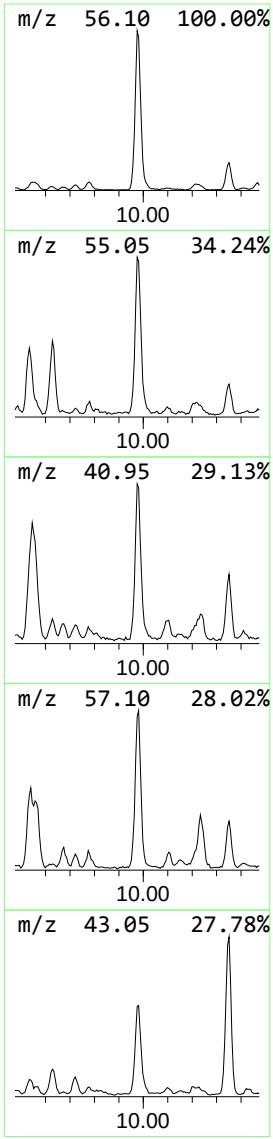
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1,3-Pentanediol, 2,2,4-trim... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.975	9.03 ng/ul	514026	Naphthalene-d8	10.469

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	74
2	1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	64
3	1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	50
4	Propane, 1-isocyanato-	85	C4H7NO	000110-78-1	47
5	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	43





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Misc :
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BNA_P
ClientSampleId :
DDCG1

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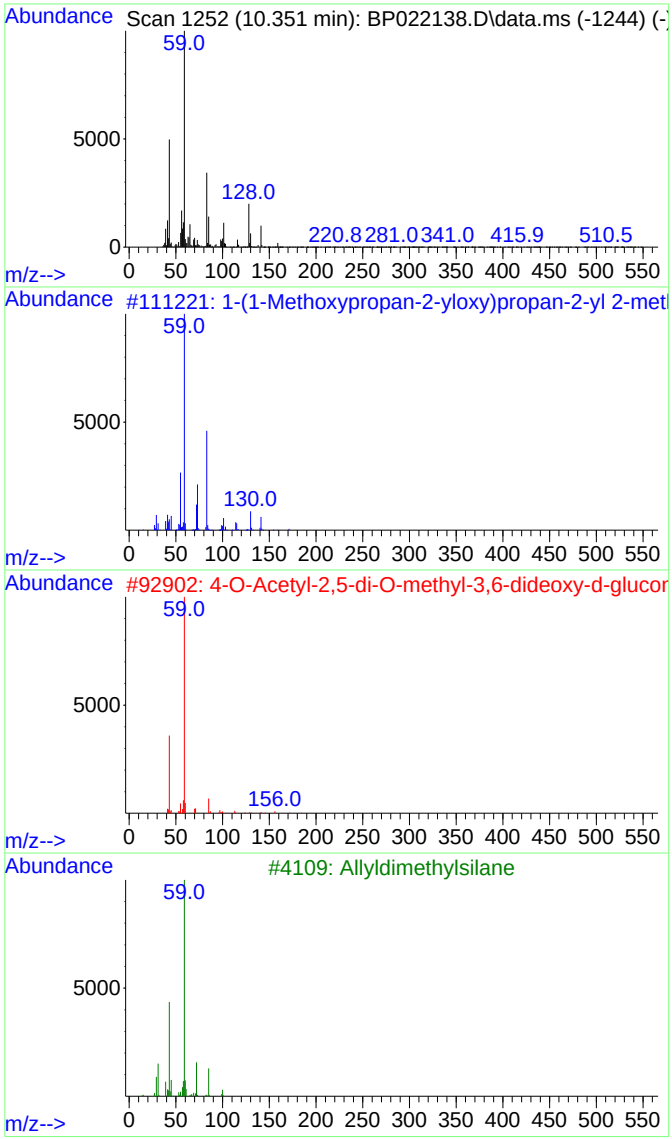
Quant Title : SVOA CALIBRATION

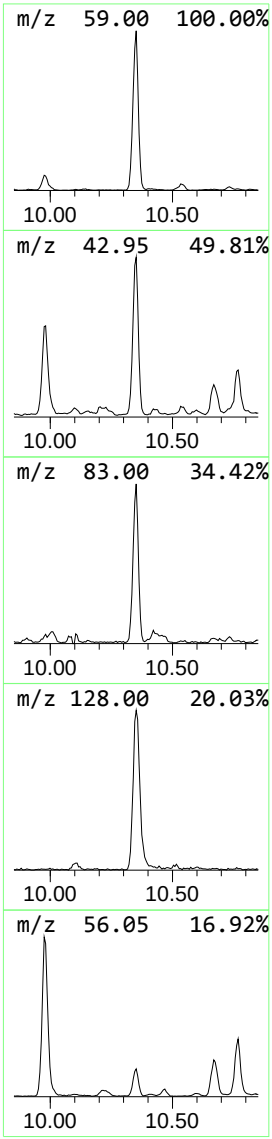
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1-(1-Methoxypropan-2-yloxy)... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.351	9.98 ng/ul	568584	Naphthalene-d8	10.469

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(1-Methoxypropan-2-yloxy)propa...	230	C12H22O4	1000367-11-3	53
2	4-O-Acetyl-2,5-di-O-methyl-3,6-d...	215	C10H17NO4	1010101-80-3	43
3	Allyldimethylsilane	100	C5H12Si	003937-30-2	43
4	4-Pentene-2-ol, 2-methyl	100	C6H12O	000624-97-5	43
5	1-(1-Methoxypropan-2-yloxy)propa...	232	C12H24O4	1000367-11-4	37





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Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

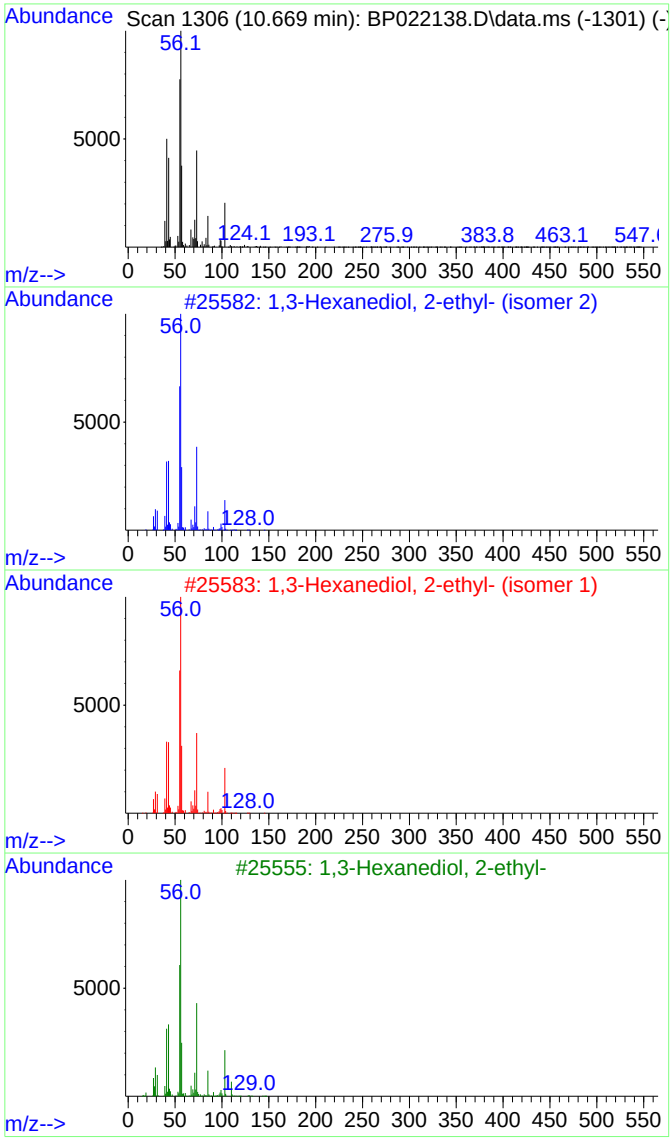
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

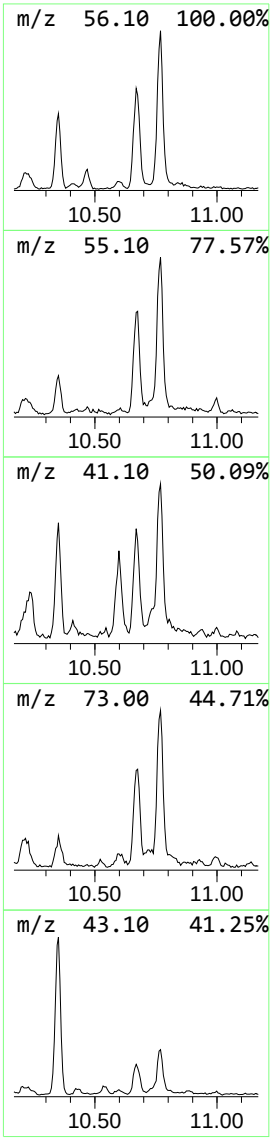
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1,3-Hexanediol, 2-ethyl- (i... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.669	3.18 ng/ul	181103	Naphthalene-d8	10.469

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Hexanediol, 2-ethyl- (isomer 2)	146	C8H18O2	1000484-18-1	90
2	1,3-Hexanediol, 2-ethyl- (isomer 1)	146	C8H18O2	1000484-18-2	90
3	1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	83
4	1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	74
5	1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	72





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Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

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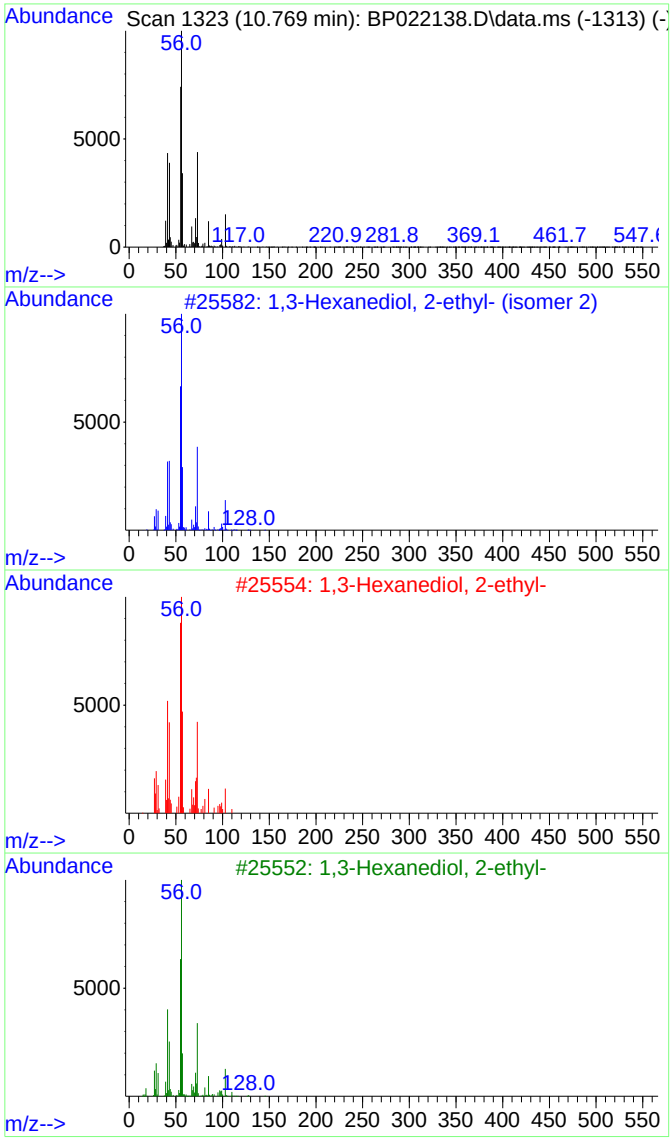
Quant Title : SVOA CALIBRATION

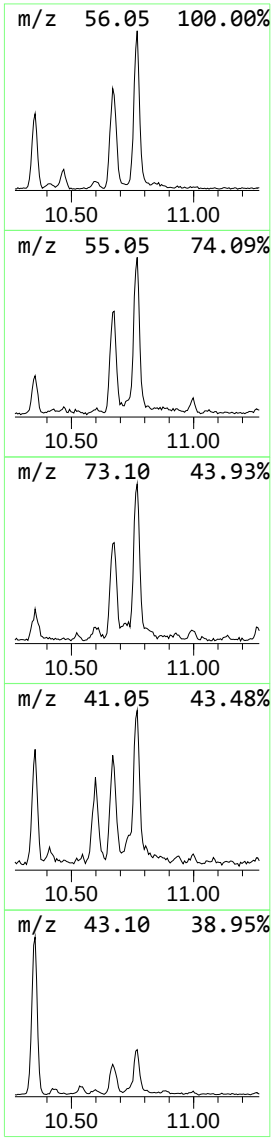
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1,3-Hexanediol, 2-ethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.769	4.86 ng/ul	276888	Naphthalene-d8	10.469

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Hexanediol, 2-ethyl- (isomer 2)	146	C8H18O2	1000484-18-1	90
2	1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	83
3	1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	64
4	1,3-Hexanediol, 2-ethyl- (isomer 1)	146	C8H18O2	1000484-18-2	64
5	Neopentyl glycol	104	C5H12O2	000126-30-7	59





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Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

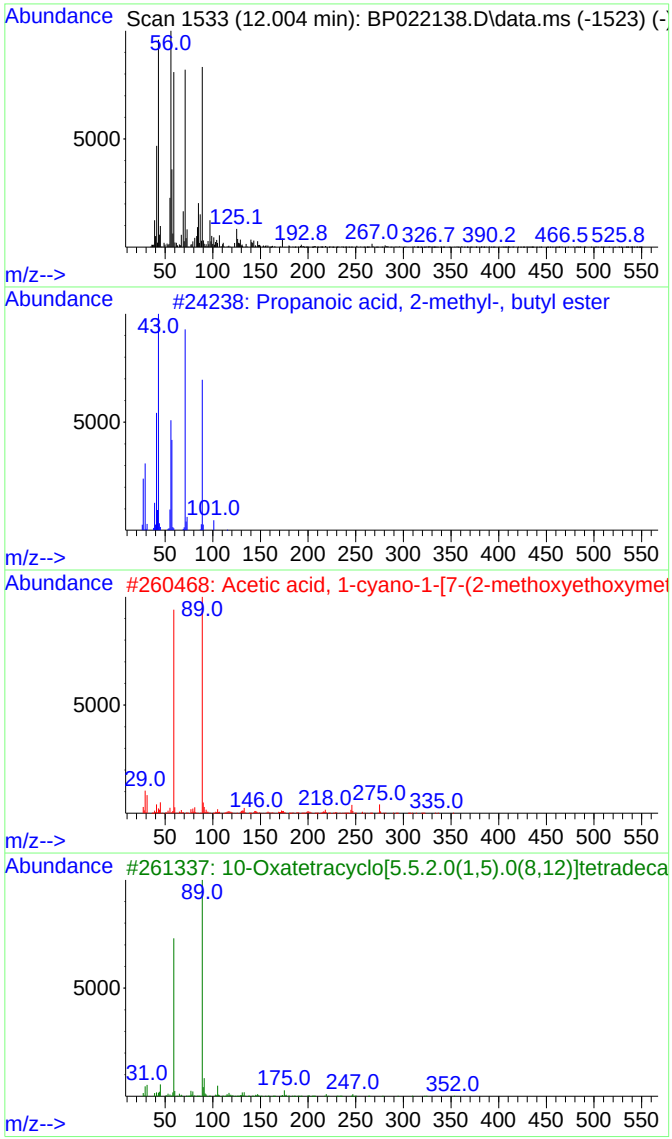
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

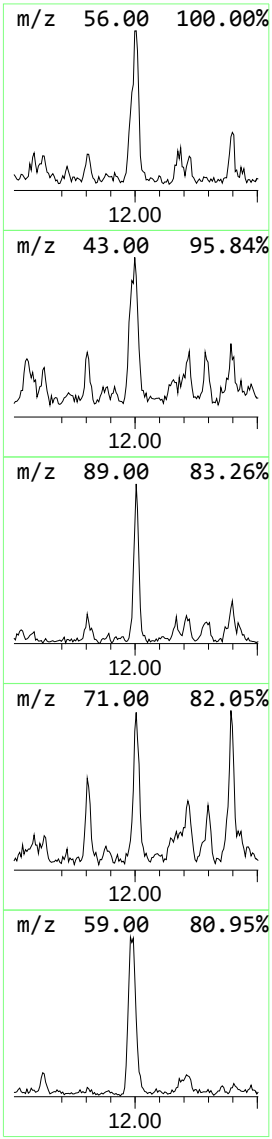
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Propanoic acid, 2-methyl-, ... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.004	2.16 ng/ul	122895	Naphthalene-d8	10.469

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	53
2	Acetic acid, 1-cyano-1-[7-(2-met...	351	C19H29NO5	1000155-83-6	52
3	10-Oxatetracyclo[5.5.2.0(1,5).0(...	352	C18H24O7	1000154-01-5	50
4	Spiro[3,5-dioxatricyclo[6.3.0.0(...	410	C23H38O6	1000157-52-8	50
5	Isobutyl 2,5,8,11-tetraoxatridec...	308	C14H28O7	1000378-27-9	47





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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

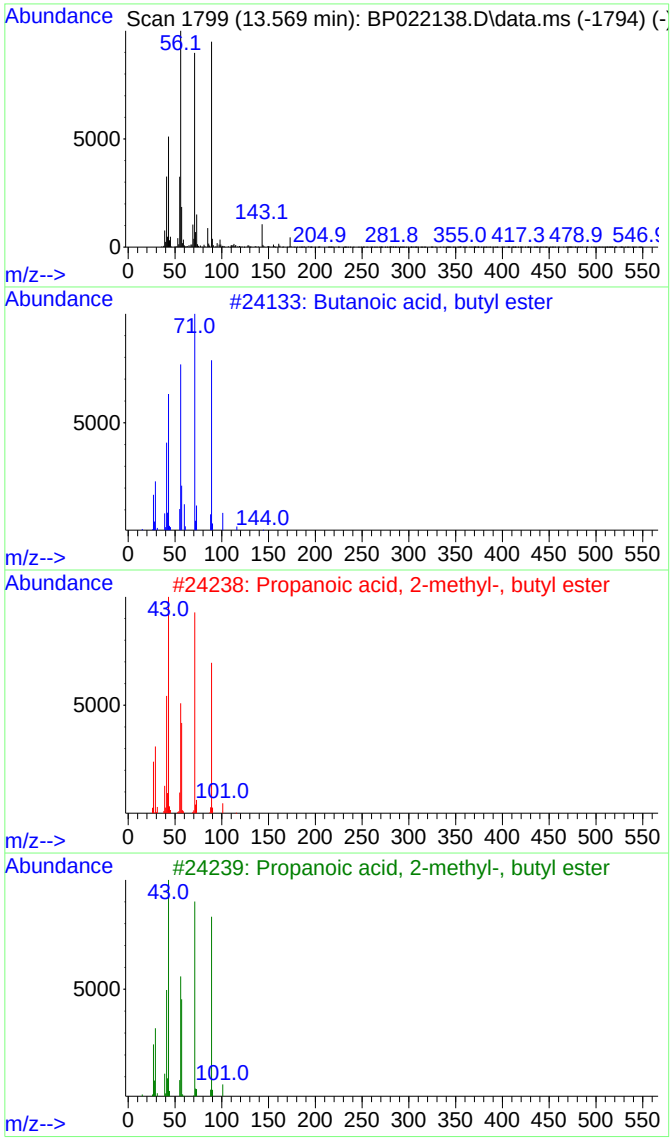
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

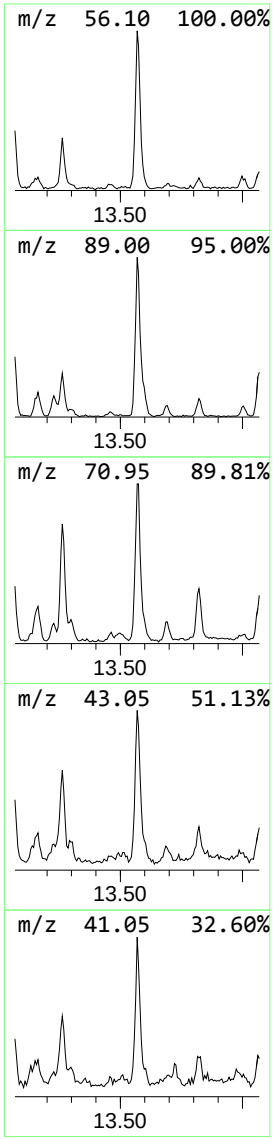
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Butanoic acid, butyl ester Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.569	3.03 ng/ul	252190	Acenaphthene-d10	14.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	80
2	Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	72
3	Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	72
4	Propanoic acid, 2-methyl-, hexyl...	172	C10H20O2	002349-07-7	56
5	Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	50





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Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

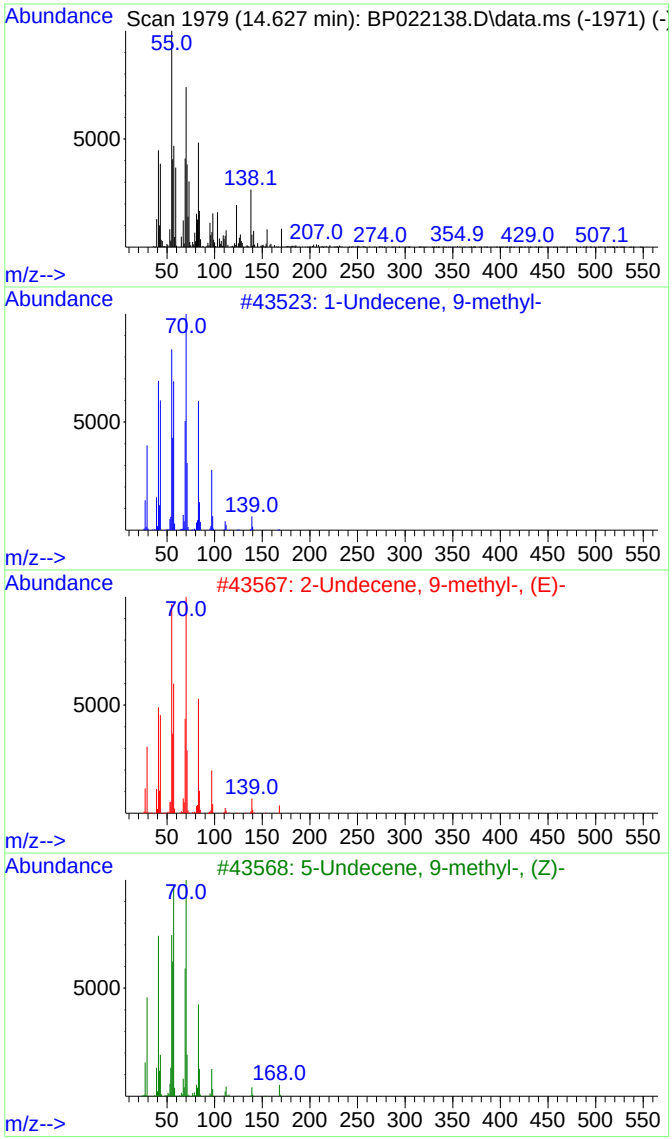
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

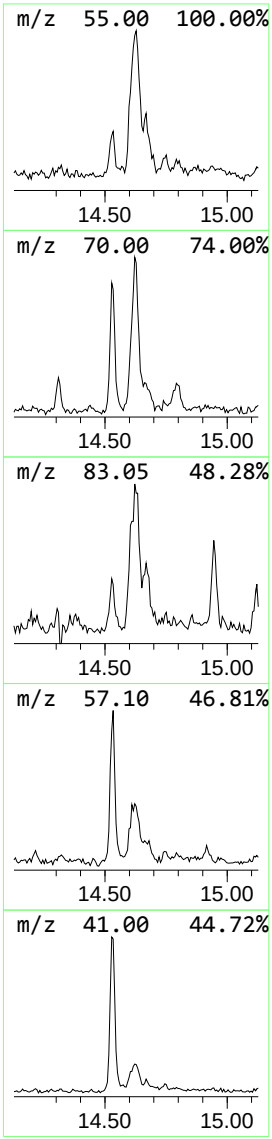
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.627	2.05 ng/ul	170435	Acenaphthene-d10	14.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Undecene, 9-methyl-	168	C12H24	074630-41-4	47
2	2-Undecene, 9-methyl-, (E)-	168	C12H24	074630-46-9	47
3	5-Undecene, 9-methyl-, (Z)-	168	C12H24	074630-65-2	38
4	Cyclopentane, 1,2,3-trimethyl-, ...	112	C8H16	015890-40-1	35
5	1-Decene, 3,3,4-trimethyl-	182	C13H26	049622-17-5	30





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ALS Vial : 11 Sample Multiplier: 1

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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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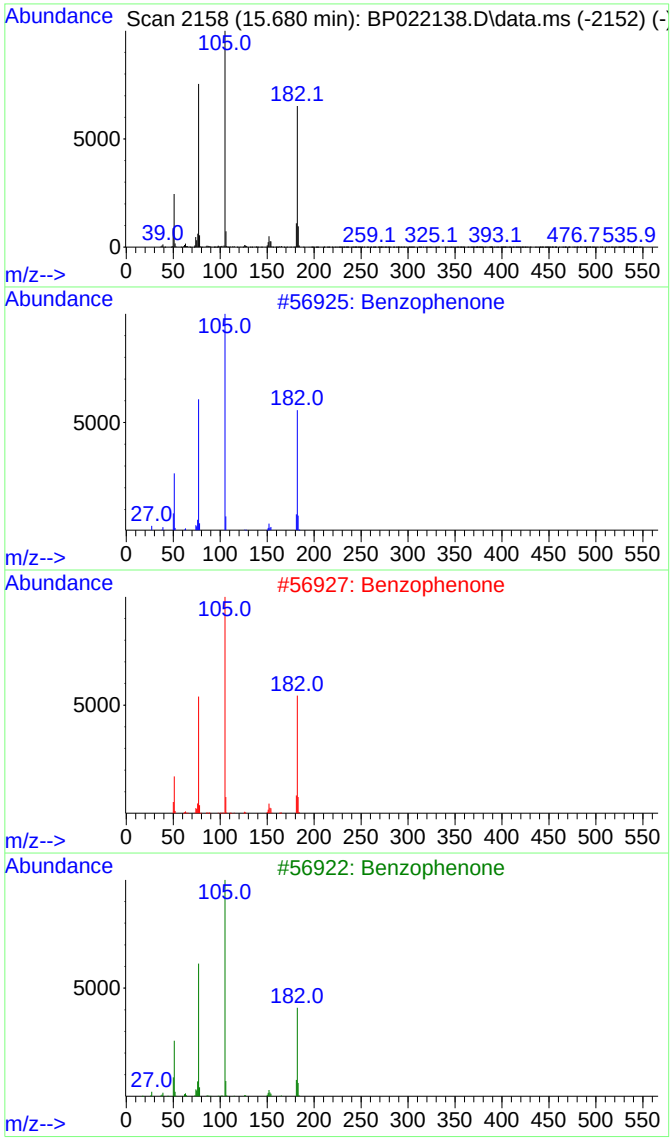
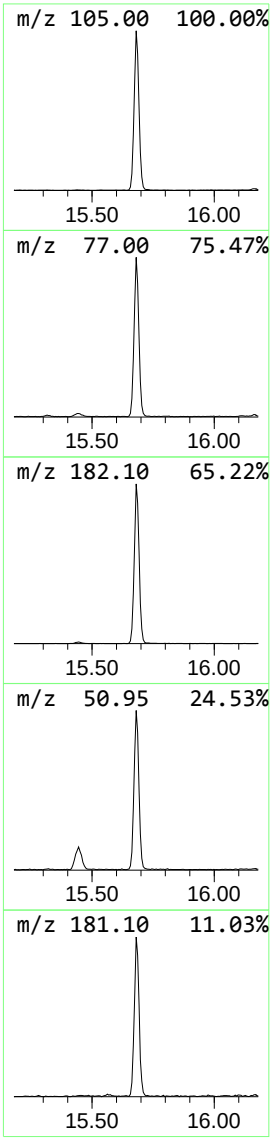
Peak Number 14 Benzophenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.680	12.04 ng/ul	1001010	Acenaphthene-d10	14.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	Benzophenone	182	C13H10O	000119-61-9	96
3	Benzophenone	182	C13H10O	000119-61-9	94
4	Benzophenone	182	C13H10O	000119-61-9	89
5	Benzophenone	182	C13H10O	000119-61-9	87





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ClientSampleId :
DDCG1

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Quant Title : SVOA CALIBRATION

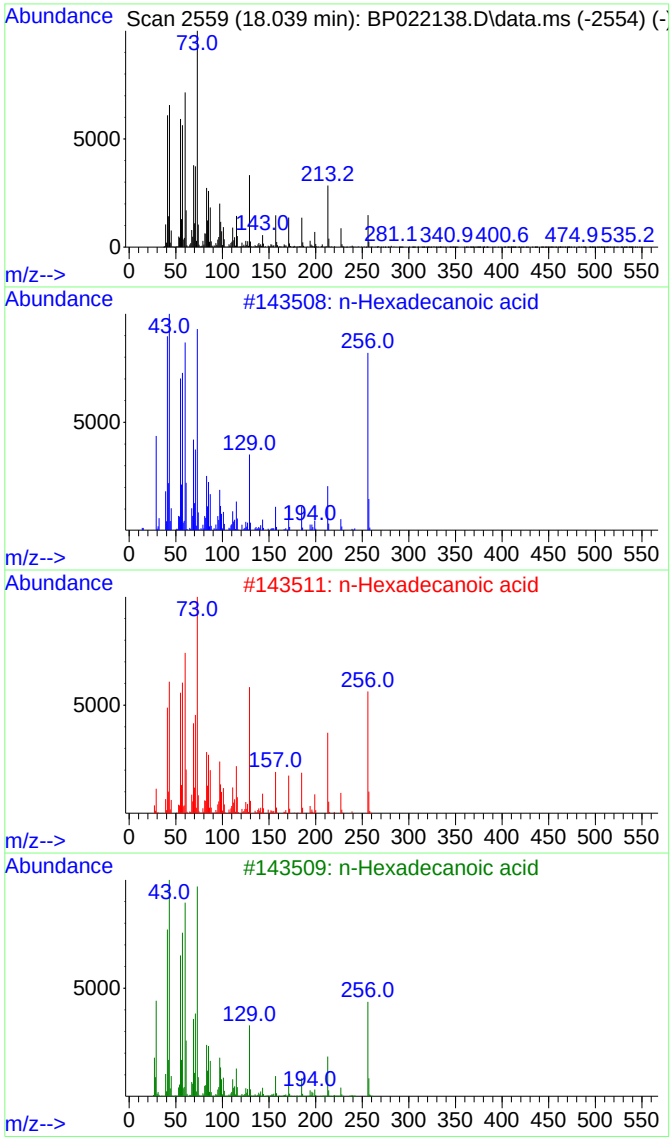
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

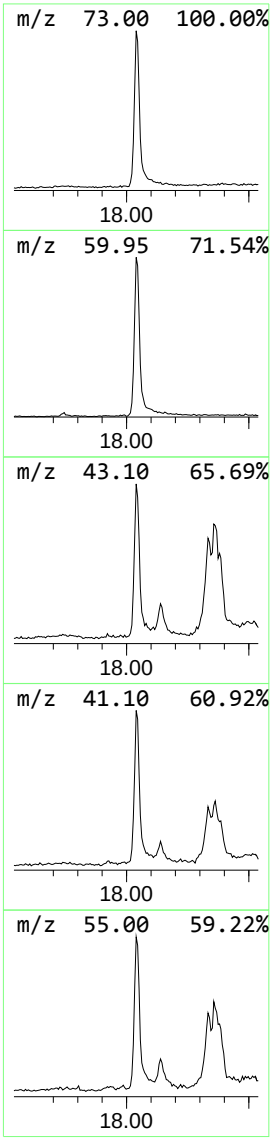
****

Peak Number 15 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.039	9.17 ng/ul	990497	Phenanthrene-d10	17.110

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	98
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96





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ALS Vial : 11 Sample Multiplier: 1

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Quant Title : SVOA CALIBRATION

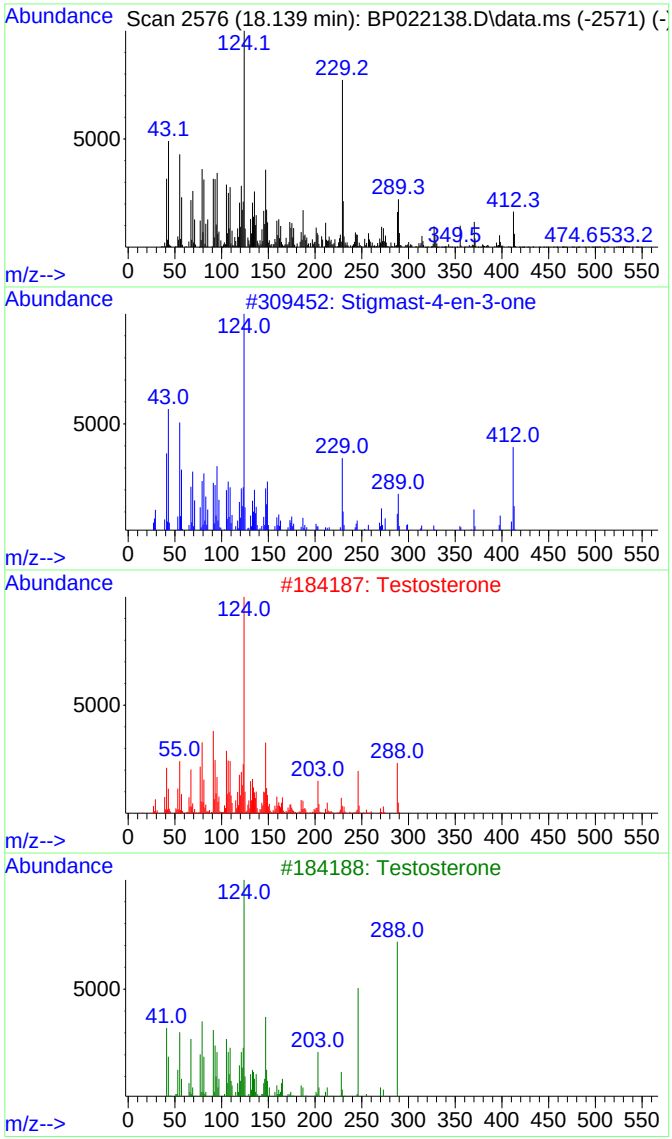
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

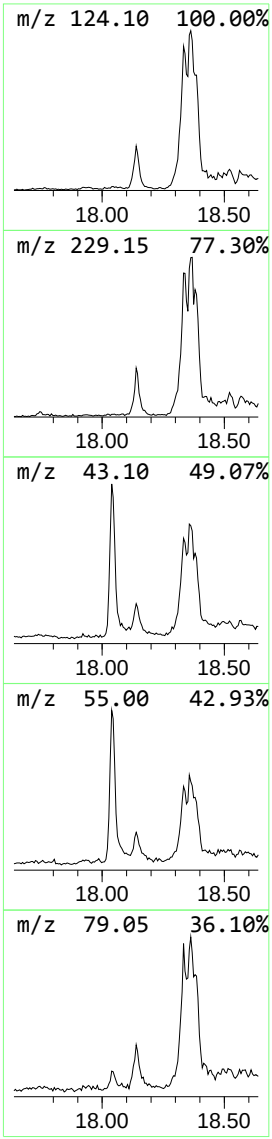
****

Peak Number 16 Stigmast-4-en-3-one Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.139	4.08 ng/ul	440429	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Stigmast-4-en-3-one	412	C29H48O	001058-61-3	83
2	Testosterone	288	C19H28O2	000058-22-0	70
3	Testosterone	288	C19H28O2	000058-22-0	70
4	.gamma.-Sitostenone	412	C29H48O	084924-96-9	60
5	Testosterone	288	C19H28O2	000058-22-0	47





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Data File : BP022138.D
Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

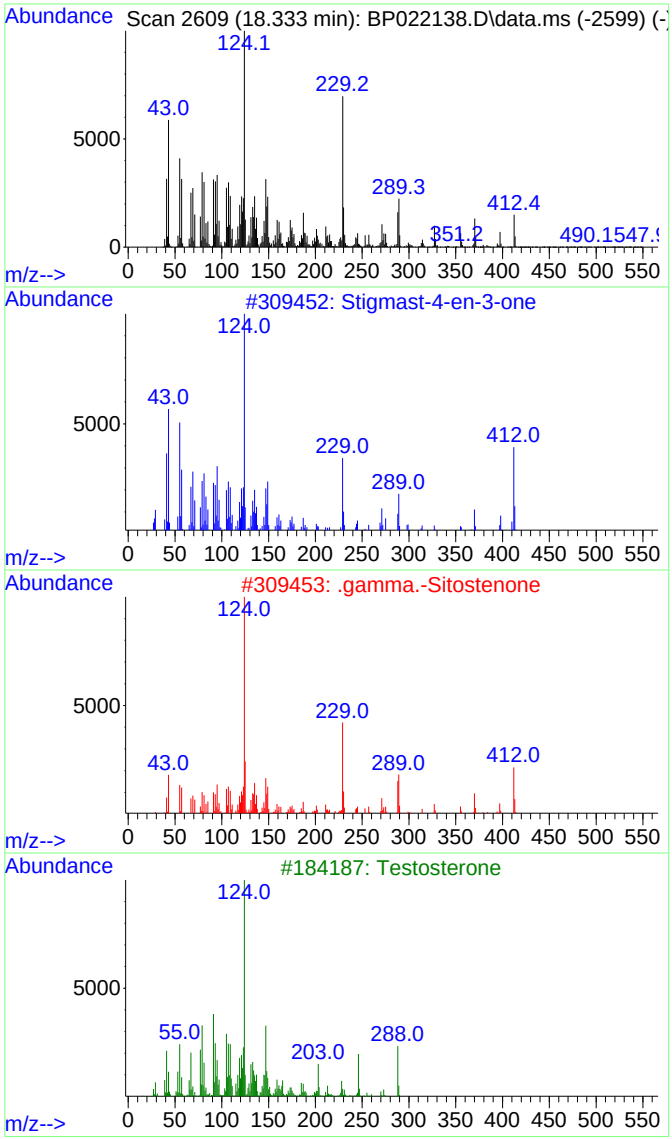
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

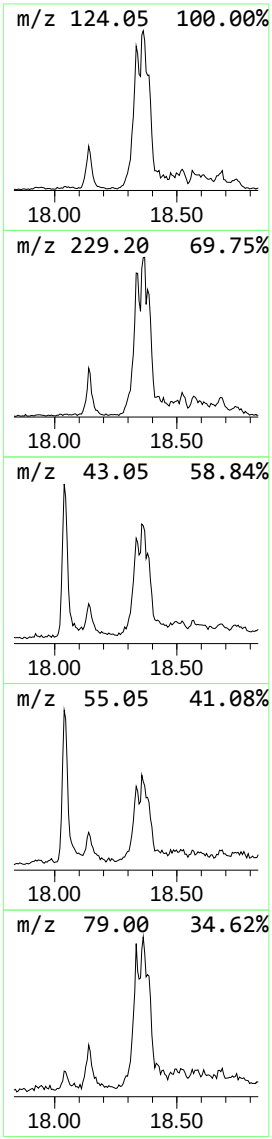
****

Peak Number 17 .gamma.-Sitostenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.333	12.51 ng/ul	1351600	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Stigmast-4-en-3-one	412	C29H48O	001058-61-3	93
2	.gamma.-Sitostenone	412	C29H48O	084924-96-9	93
3	Testosterone	288	C19H28O2	000058-22-0	80
4	Testosterone cypionate	412	C27H40O3	000058-20-8	53
5	Androst-4-en-3-one, 17-hydroxy-,...	288	C19H28O2	000481-30-1	45





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Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

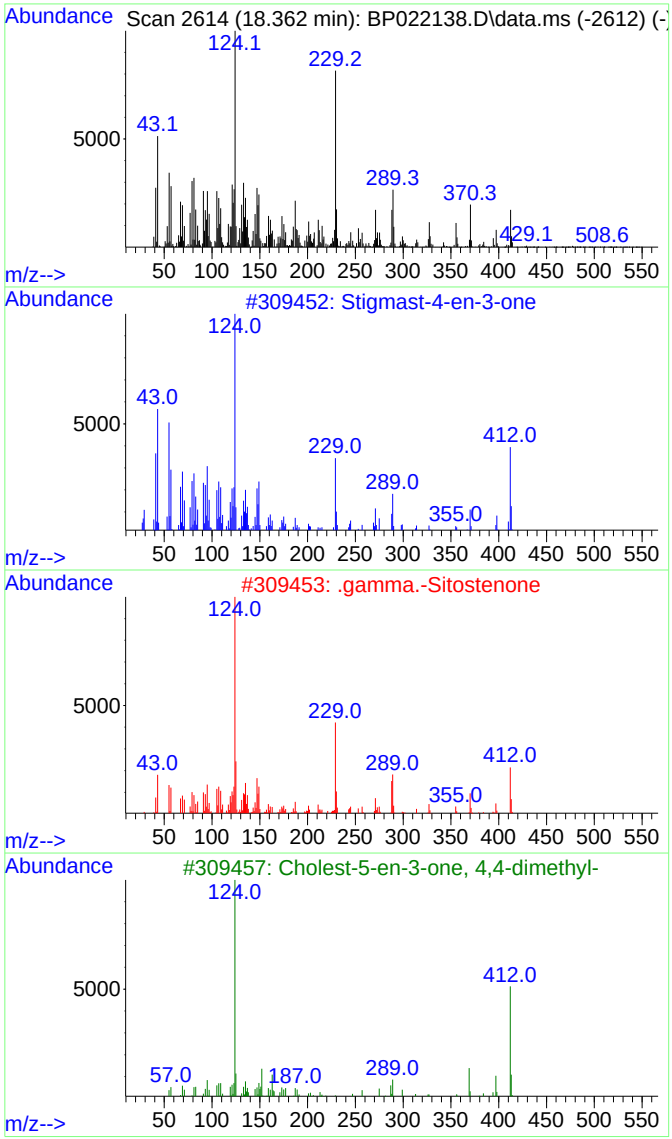
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

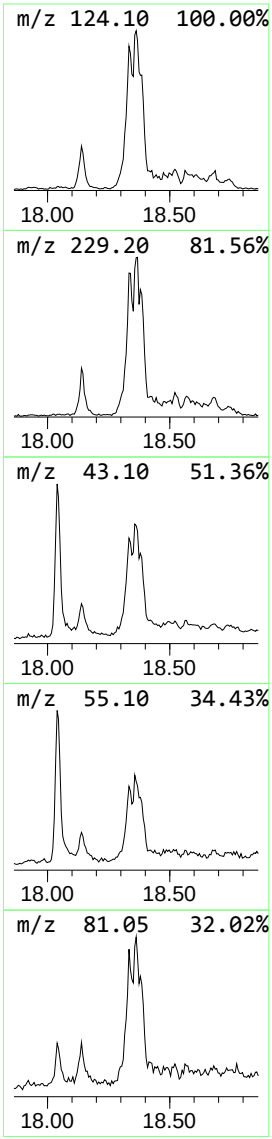
****

Peak Number 18 unknown-02 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.363	2.78 ng/ul	299742	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Stigmast-4-en-3-one	412	C29H48O	001058-61-3	92
2	.gamma.-Sitostenone	412	C29H48O	084924-96-9	90
3	Cholest-5-en-3-one, 4,4-dimethyl-	412	C29H48O	002220-42-0	46
4	Testosterone cypionate	412	C27H40O3	000058-20-8	41
5	Littorine	289	C17H23NO3	021956-47-8	30





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Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

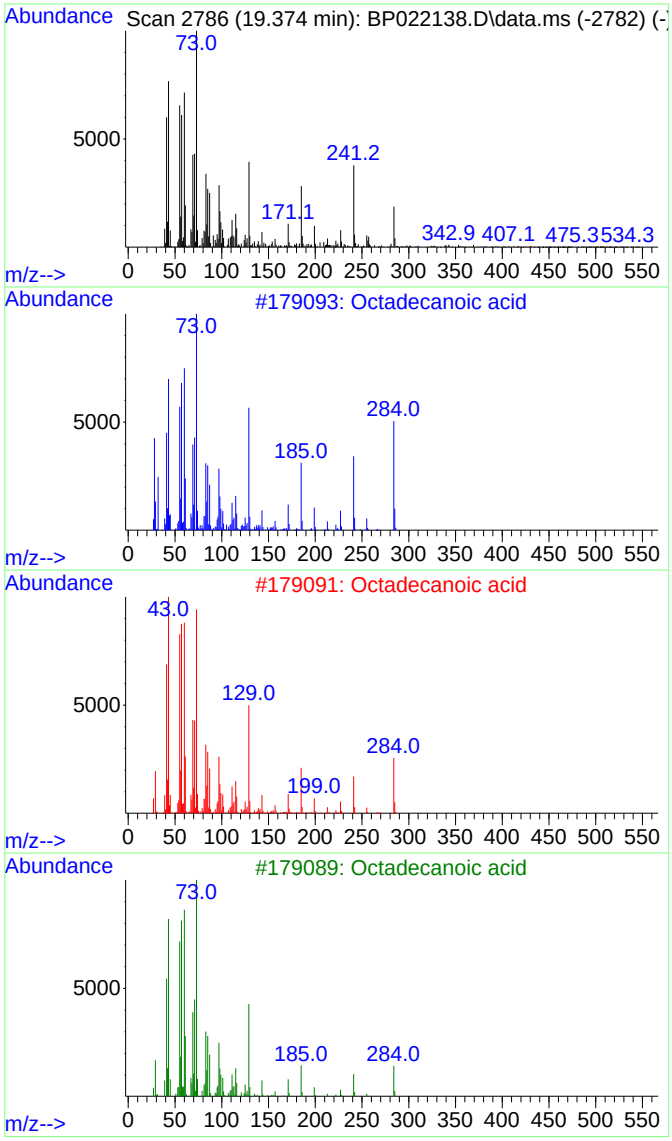
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

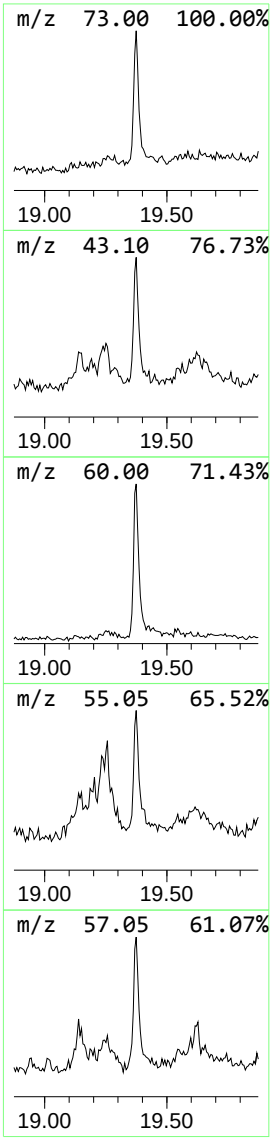
****

Peak Number 19 Octadecanoic acid Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.374	2.23 ng/ul	273702	Chrysene-d12	21.545

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	97
3	Octadecanoic acid	284	C18H36O2	000057-11-4	96
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	86
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86





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Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

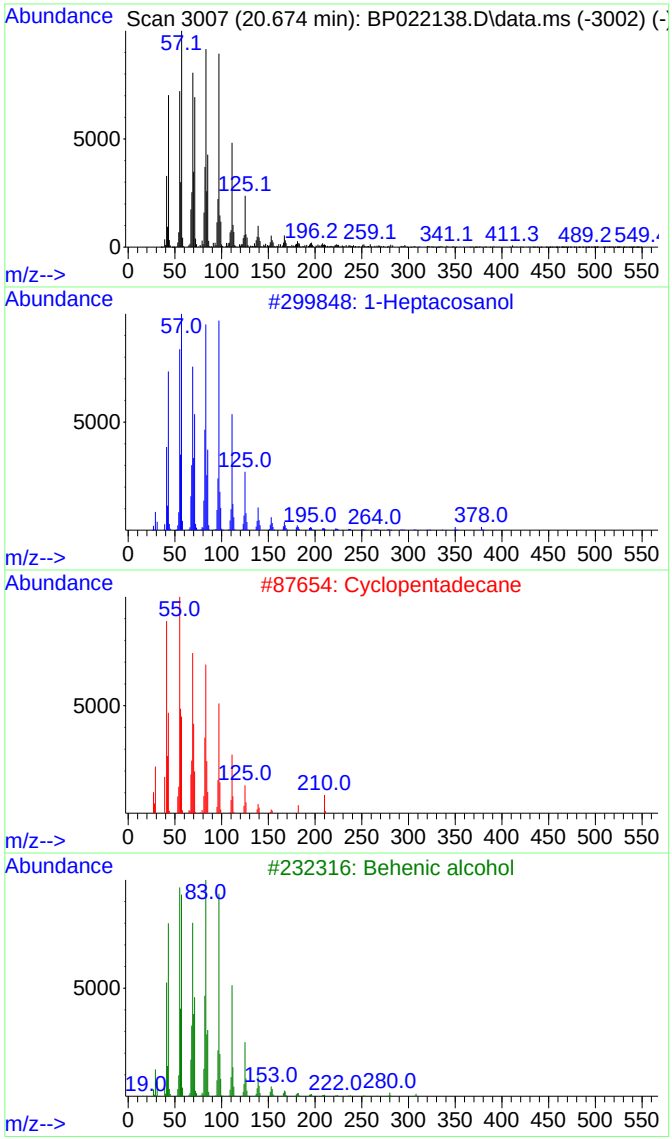
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

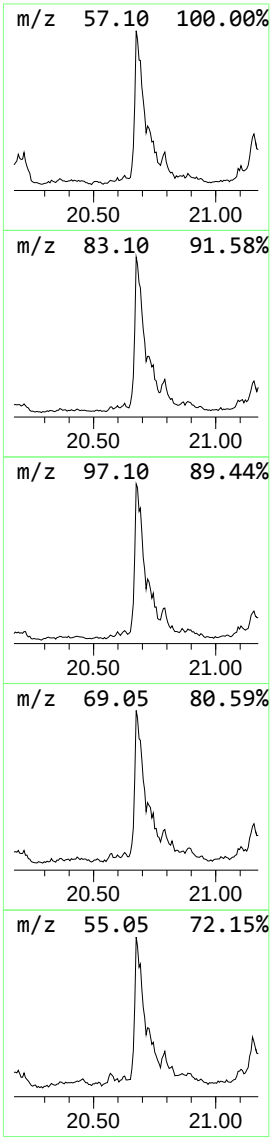
****

Peak Number 20 1-Heptacosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.674	21.82 ng/ul	2674550	Chrysene-d12	21.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptacosanol	396	C27H56O	002004-39-9	95
2	Cyclopentadecane	210	C15H30	000295-48-7	93
3	Behenic alcohol	326	C22H46O	000661-19-8	93
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	93
5	Triacontan-15-ol	438	C30H62O	031849-26-0	93





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Data File : BP022138.D
Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

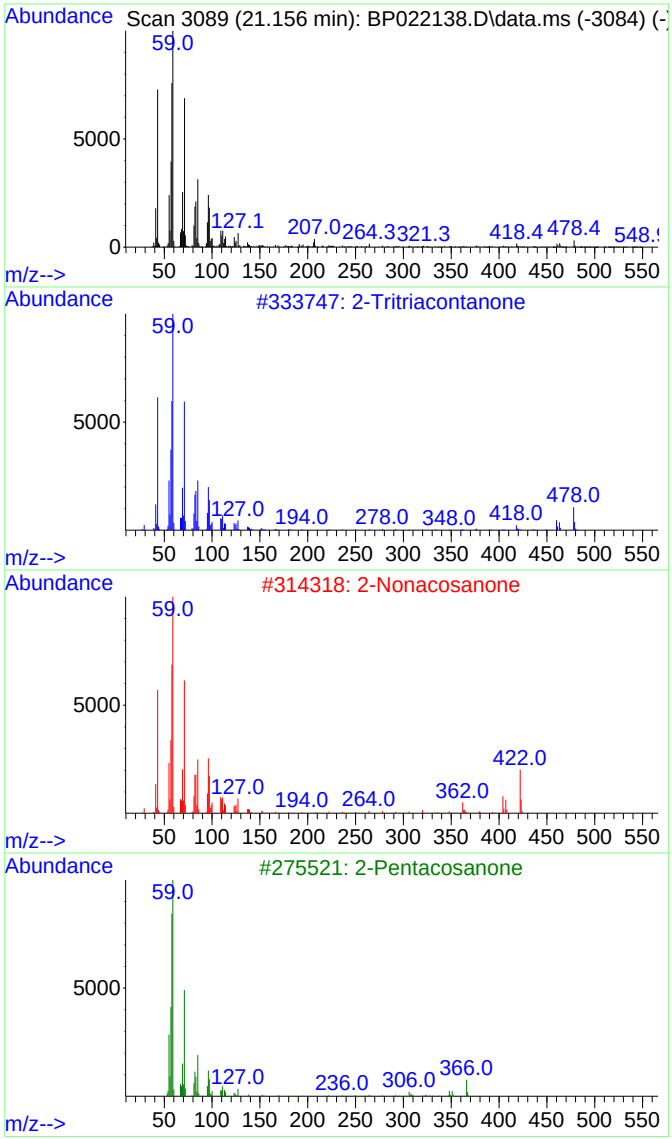
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

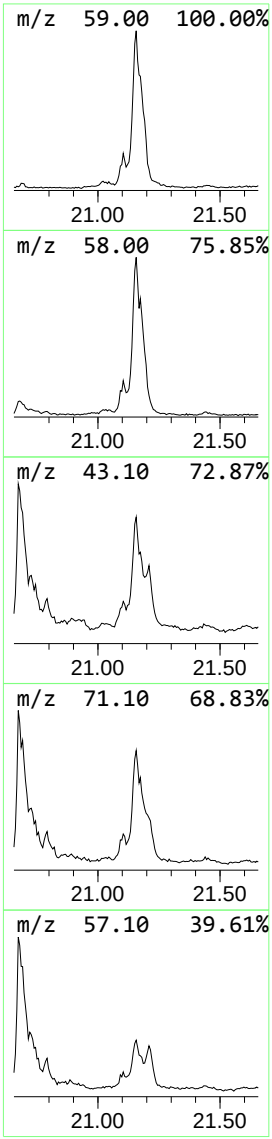
****

Peak Number 21 2-Tritriacontanone Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.157	7.43 ng/ul	910210	Chrysene-d12	21.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Tritriacontanone	479	C33H66O	075207-55-5	98
2	2-Nonacosanone	422	C29H58O	017600-99-6	90
3	2-Pentacosanone	366	C25H50O	075207-54-4	83
4	2-Pentacosanone	366	C25H50O	075207-54-4	83
5	Oxirane, tetramethyl-	100	C6H12O	005076-20-0	47





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Data File : BP022138.D
Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

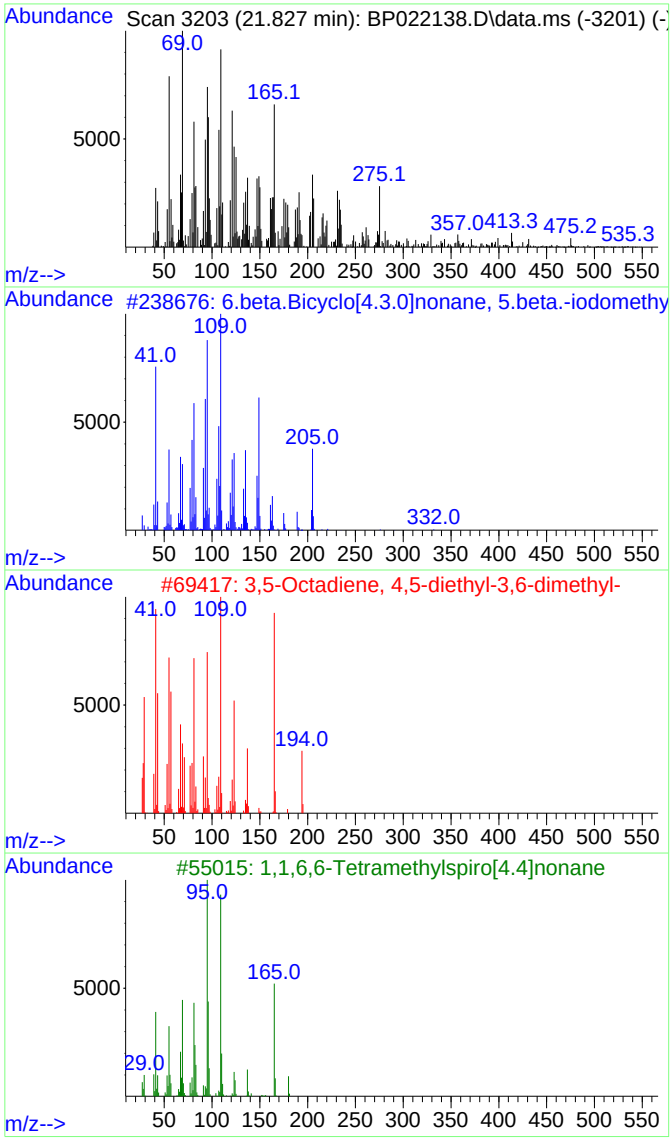
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

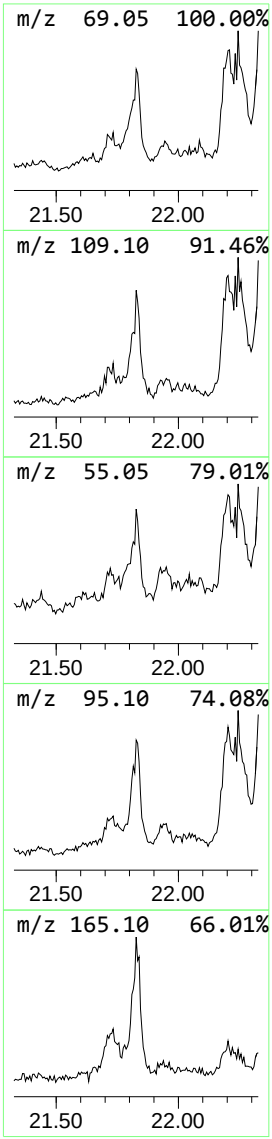
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 6.beta.Bicyclo[4.3.0]nonane... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.827	5.45 ng/ul	668228	Chrysene-d12	21.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6.beta.Bicyclo[4.3.0]nonane, 5.b...	332	C15H25I	1000195-85-9	64
2	3,5-Octadiene, 4,5-diethyl-3,6-d...	194	C14H26	061233-79-2	49
3	1,1,6,6-Tetramethylspiro[4.4]nonane	180	C13H24	074054-92-5	43
4	(1R,2R,4aS,8aS)-1-(2-(Furan-3-yl...	304	C20H32O2	134551-51-2	35
5	2,5,9-Trimethylcycloundeca-4,8-d...	206	C14H22O	1000195-33-9	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
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Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

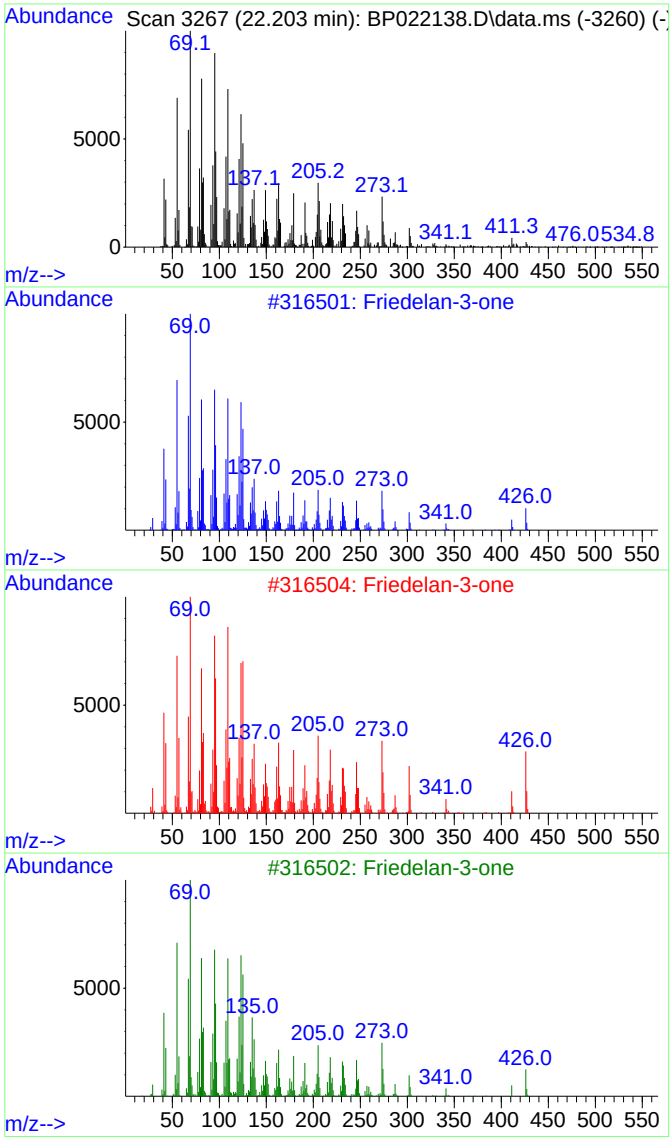
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

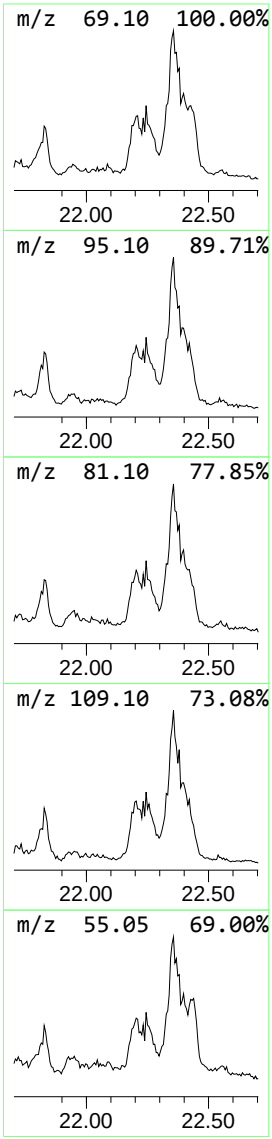
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 2,3,3-Trimethyl-2-(3-methyl... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.203	8.74 ng/ul	1071590	Chrysene-d12	21.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Friedelan-3-one	426	C30H50O	000559-74-0	81
2	Friedelan-3-one	426	C30H50O	000559-74-0	64
3	Friedelan-3-one	426	C30H50O	000559-74-0	62
4	2,3,3-Trimethyl-2-(3-methyl-buta...	206	C14H22O	1000193-60-6	52
5	(E)-3-Methyl-5-((1R,4aR,8aR)-5,5...	290	C20H34O	021738-29-4	46





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Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

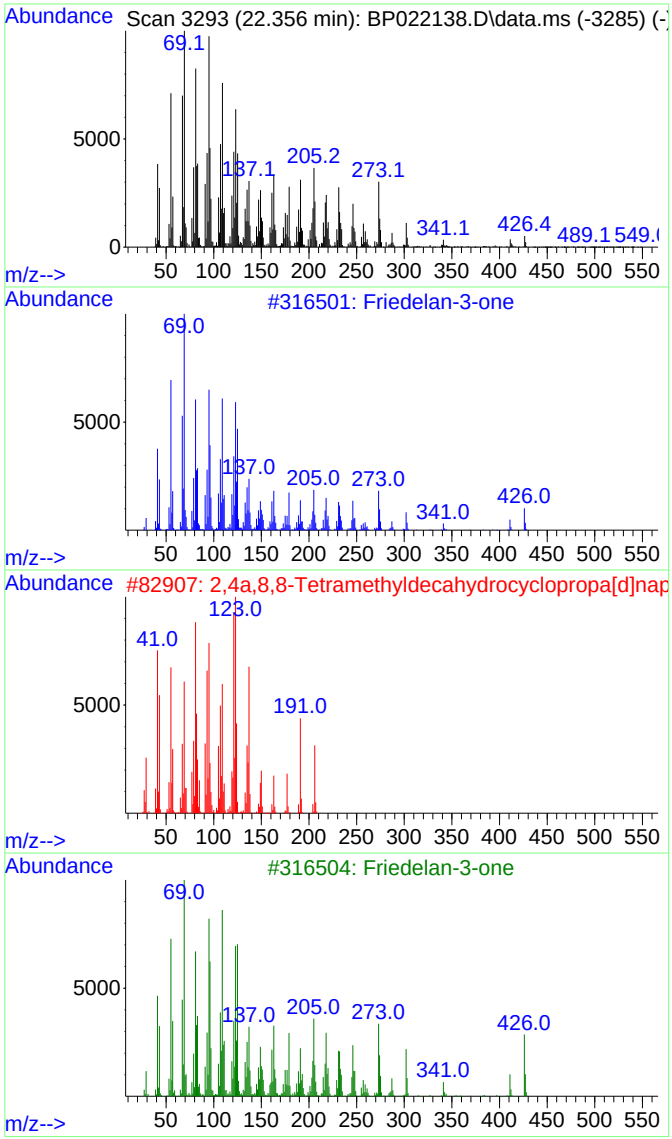
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

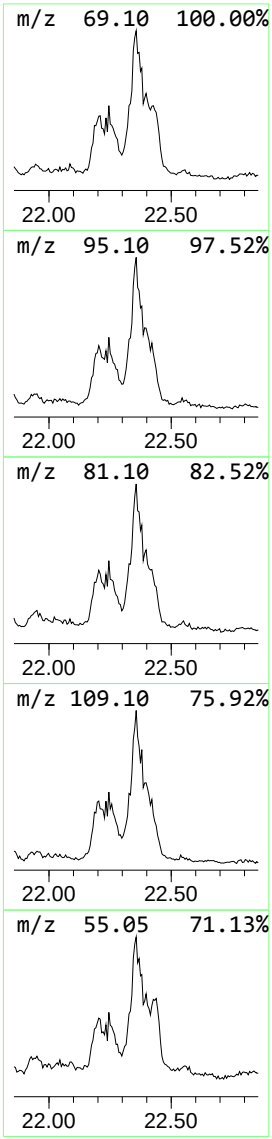
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 2,4a,8,8-Tetramethyldecahyd... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.356	25.01 ng/ul	3065190	Chrysene-d12	21.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Friedelan-3-one	426	C30H50O	000559-74-0	97
2	2,4a,8,8-Tetramethyldecahydrocyc...	206	C15H26	074022-04-1	90
3	Friedelan-3-one	426	C30H50O	000559-74-0	89
4	5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	56
5	2,2,6-Trimethyl-1-(2-methyl-cycl...	218	C15H22O	1000188-72-8	56





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Data File : BP022138.D
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Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

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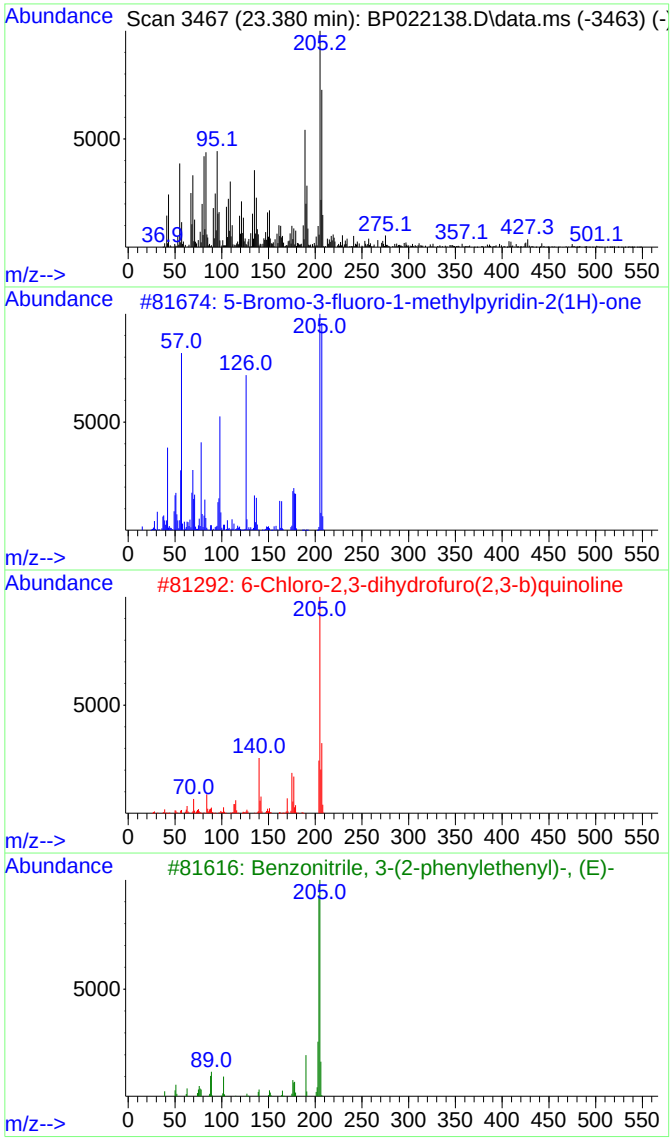
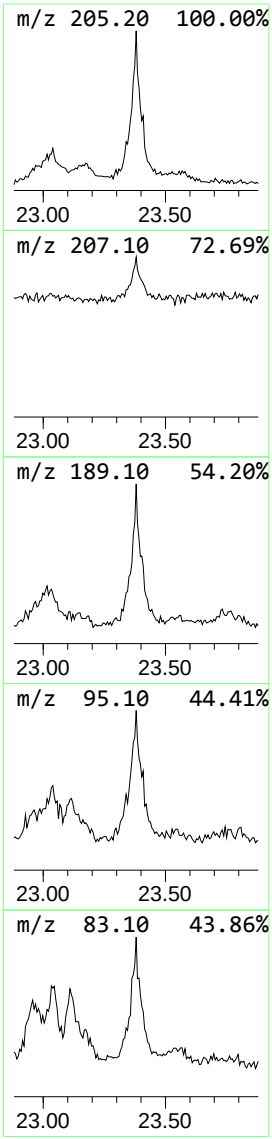
Peak Number 25 unknown-03 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.380	2.58 ng/ul	348375	Perylene-d12	24.833

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Bromo-3-fluoro-1-methylpyridin...	205	C6H5BrFNO	1352152-46-5	35
2	6-Chloro-2,3-dihydrofuro(2,3-b)q...	205	C11H8ClNO	064124-84-1	27
3	Benzonitrile, 3-(2-phenylethenyl...	205	C15H11N	014064-35-8	27
4	1-(3-Chlorophenyl)-3-methyl-1H-p...	207	C10H10ClN3	040401-41-0	25
5	N-(1-Methyl-3-oxobutylidene)-4-m...	205	C12H15NO2	020010-38-2	25





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022138.D
Acq On : 01 Oct 2024 16:31
Operator : RC/JU
Sample : P4021-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCG1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.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
Butanoic acid	3.928	9.0	ng/ul	488069	1	7.698	1082060	20.0
2-Pentanone, 4-...	4.869	10.8	ng/ul	581776	1	7.698	1082060	20.0
Hexylene glycol	6.075	37.9	ng/ul	2051520	1	7.698	1082060	20.0
unknown-01	7.934	14.1	ng/ul	761647	1	7.698	1082060	20.0
Hexanoic acid, ...	9.075	45.1	ng/ul	2438330	1	7.698	1082060	20.0
3-Methyl-2-bute...	9.628	3.9	ng/ul	220854	2	10.469	1139090	20.0
1,3-Pentanediol...	9.975	9.0	ng/ul	514026	2	10.469	1139090	20.0
1-(1-Methoxypro...	10.351	10.0	ng/ul	568584	2	10.469	1139090	20.0
1,3-Hexanediol,...	10.669	3.2	ng/ul	181103	2	10.469	1139090	20.0
1,3-Hexanediol,...	10.769	4.9	ng/ul	276888	2	10.469	1139090	20.0
Propanoic acid,...	12.004	2.2	ng/ul	122895	2	10.469	1139090	20.0
Butanoic acid, ...	13.569	3.0	ng/ul	252190	3	14.310	1662920	20.0
(DEL) Alkane: S...	14.627	2.0	ng/ul	170435	3	14.310	1662920	20.0
Benzophenone	15.680	12.0	ng/ul	1001010	3	14.310	1662920	20.0
n-Hexadecanoic ...	18.039	9.2	ng/ul	990497	4	17.110	2160180	20.0
Stigmast-4-en-3...	18.139	4.1	ng/ul	440429	4	17.110	2160180	20.0
.gamma.-Sitoste...	18.333	12.5	ng/ul	1351600	4	17.110	2160180	20.0
unknown-02	18.363	2.8	ng/ul	299742	4	17.110	2160180	20.0
Octadecanoic acid	19.374	2.2	ng/ul	273702	5	21.545	2451420	20.0
1-Heptacosanol	20.674	21.8	ng/ul	2674550	5	21.545	2451420	20.0
2-Tritriacontanone	21.157	7.4	ng/ul	910210	5	21.545	2451420	20.0
6.beta.Bicyclo[...	21.827	5.5	ng/ul	668228	5	21.545	2451420	20.0
2,3,3-Trimethyl...	22.203	8.7	ng/ul	1071590	5	21.545	2451420	20.0
2,4a,8,8-Tetram...	22.356	25.0	ng/ul	3065190	5	21.545	2451420	20.0
unknown-03	23.380	2.6	ng/ul	348375	6	24.833	2700570	20.0