LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020683.D
Acq On : 28 Jun 2024 06:58
Operator : MA/JU
Sample : P2898-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G84

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-BP062524.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP020683.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.011	3	4	23	rVB	5763301	6305437	100.00%	9.306%
2	3.240	40	43	46	rBV	17123	22004	0.35%	0.032%
3	3.275	46	49	58	rVB	47718	65760	1.04%	0.097%
4	3.540	90	94	103	rBV	104176	160241	2.54%	0.236%
5	3.693	116	120	132	rBV	137561	219407	3.48%	0.324%
6	3.999	168	172	177	rVB	29227	37497	0.59%	0.055%
7	4.358	229	233	241	rVB9	10381	19873	0.32%	0.029%
8	4.422	241	244	252	rVB3	5233	10615	0.17%	0.016%
9	4.499	252	257	261	rBV6	5476	7990	0.13%	0.012%
10	4.552	261	266	272	rBV4	8085	12572	0.20%	0.019%
11	4.899	320	325	334	rVB	28432	45434	0.72%	0.067%
12	5.528	428	432	437	rBV	9571	17147	0.27%	0.025%
13	5.569	437	439	445	rVB5	5224	7379	0.12%	0.011%
14	5.834	480	484	487	rBV4	8514	13507	0.21%	0.020%
15	6.664	620	625	632	rVB4	7319	13292	0.21%	0.020%
16	6.775	638	644	654	rVB	23248	39967	0.63%	0.059%
17	6.922	663	669	683	rVB	247540	420060	6.66%	0.620%
18	7.087	690	697	710	rBV	908997	1411736	22.39%	2.084%
19	7.287	723	731	739	rBV	887920	1488213	23.60%	2.196%
20	7.358	739	743	752	rVV	29541	60440	0.96%	0.089%
21	7.440	755	757	763	rVB7	5831	7844	0.12%	0.012%
22	7.752	802	810	815	rBV	795853	1246723	19.77%	1.840%
23	7.805	815	819	834	rVB	389934	623564	9.89%	0.920%
24	8.146	871	877	882	rVB6	3455	7636	0.12%	0.011%
25	8.322	899	907	921	rVB2	15547	36201	0.57%	0.053%
26	8.452	921	929	941	rBV	754049	1220814	19.36%	1.802%
27	8.563	943	948	952	rVB3	6908	12105	0.19%	0.018%
28	8.840	989	995	998	rBV	19568	29396	0.47%	0.043%
29	8.899	998	1005	1016	rVV	1167949	1900569	30.14%	2.805%
30	9.016	1020	1025	1029	rVV	20343	34736	0.55%	0.051%
31	9.052	1029	1031	1038	rVB3	13593	19360	0.31%	0.029%
32	9.287	1062	1071	1073	rBV8	4770	10226	0.16%	0.015%
33	9.452	1093	1099	1105	rBV	37311	62281	0.99%	0.092%
34	9.605	1118	1125	1142	rBV	973784	1585968	25.15%	2.341%
35	9.799	1146	1158	1163	rVV	87659	186512	2.96%	0.275%
36	9.857	1163	1168	1182	rVB	86045	165593	2.63%	0.244%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020683.D
Acq On : 28 Jun 2024 06:58
Operator : MA/JU
Sample : P2898-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G84

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-BP062524.MA.M
Title : SVOA CALIBRATION

37	10.146	1208	1217	1235	rBV	1262411	2308924	36.62%	3.408%
38	10.299	1237	1243	1253	rBV	143553	229380	3.64%	0.339%
39	10.522	1272	1281	1289	rBV	1043976	1825569	28.95%	2.694%
40	10.657	1296	1304	1320	rBV	846631	1365770	21.66%	2.016%
41	10.887	1337	1343	1346	rBV5	5654	9818	0.16%	0.014%
42	10.934	1348	1351	1356	rVB7	4008	7014	0.11%	0.010%
43	11.057	1367	1372	1380	rVB2	24394	42264	0.67%	0.062%
44	11.169	1382	1391	1400	rBV	80922	155016	2.46%	0.229%
45	11.263	1400	1407	1412	rBV2	93263	169618	2.69%	0.250%
46	11.328	1412	1418	1430	rVV	189250	375163	5.95%	0.554%
47	11.416	1430	1433	1438	rVB7	4350	8432	0.13%	0.012%
48	11.504	1446	1448	1454	rVB7	4966	6856	0.11%	0.010%
49	11.775	1489	1494	1498	rVB6	4220	7487	0.12%	0.011%
50	12.104	1545	1550	1557	rVB	30572	48127	0.76%	0.071%
51	12.322	1577	1587	1591	rBV	4924	13123	0.21%	0.019%
52	12.640	1636	1641	1648	rBV5	16012	30364	0.48%	0.045%
53	12.969	1693	1697	1703	rBV9	5035	10099	0.16%	0.015%
54	13.116	1718	1722	1724	rBV5	4816	6500	0.10%	0.010%
55	13.181	1731	1733	1741	rVB8	4633	8346	0.13%	0.012%
56	13.357	1760	1763	1774	rVB8	6206	14423	0.23%	0.021%
57	13.793	1828	1837	1846	rBV	2326320	3748298	59.45%	5.532%
58	13.904	1851	1856	1864	rVB	6931	17409	0.28%	0.026%
59	14.069	1877	1884	1894	rVB2	2664170	4328855	68.65%	6.389%
60	14.287	1916	1921	1926	rVB4	14872	23246	0.37%	0.034%
61	14.375	1928	1936	1946	rBV	1589451	2505763	39.74%	3.698%
62	14.604	1967	1975	1985	rBV	185623	356802	5.66%	0.527%
63	14.745	1995	1999	2004	rBV4	17254	25380	0.40%	0.037%
64	14.804	2004	2009	2019	rVB3	38169	65557	1.04%	0.097%
65	14.981	2035	2039	2043	rBV4	5201	9462	0.15%	0.014%
66	15.134	2059	2065	2070	rBV	22438	32036	0.51%	0.047%
67	15.187	2070	2074	2080	rVV	19099	36181	0.57%	0.053%
68	15.245	2082	2084	2090	rVB3	14303	19515	0.31%	0.029%
69	15.392	2096	2109	2121	rBV	3721386	5479854	86.91%	8.088%
70	15.516	2123	2130	2142	rVV	1099634	1531626	24.29%	2.261%
71	15.628	2146	2149	2162	rVB3	19437	42518	0.67%	0.063%
72	15.734	2163	2167	2171	rBV7	6731	10834	0.17%	0.016%
73	15.957	2202	2205	2209	rBV6	4755	7668	0.12%	0.011%
74	16.098	2224	2229	2236	rBV	36606	72546	1.15%	0.107%
75	16.187	2241	2244	2249	rVB5	7459	11724	0.19%	0.017%
76	16.575	2306	2310	2314	rBV5	23915	31868	0.51%	0.047%

LSC Area Percent Report

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Acq On : 28 Jun 2024 06:58
Operator : MA/JU
Sample : P2898-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G84

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-BP062524.MA.M
Title : SVOA CALIBRATION

77	16.857	2356	2358	2362	rVB5	6237	7058	0.11%	0.010%
78	17.192	2407	2415	2423	rBV	1856681	2859466	45.35%	4.220%
79	17.292	2425	2432	2443	rVV	3907140	5666132	89.86%	8.363%
80	17.386	2443	2448	2453	rVB7	14218	31093	0.49%	0.046%
81	17.498	2462	2467	2476	rBV3	47909	80888	1.28%	0.119%
82	18.122	2568	2573	2583	rBV	222017	408056	6.47%	0.602%
83	18.628	2656	2659	2664	rVB6	14892	20831	0.33%	0.031%
84	19.216	2756	2759	2764	rVB3	55457	74948	1.19%	0.111%
85	19.292	2769	2772	2777	rVV	46940	65267	1.04%	0.096%
86	19.357	2778	2783	2792	rVB2	96815	162293	2.57%	0.240%
87	19.463	2797	2801	2809	rBV	156907	218165	3.46%	0.322%
88	19.610	2822	2826	2831	rBV	168495	250881	3.98%	0.370%
89	19.675	2831	2837	2846	rVB	4003130	5639691	89.44%	8.324%
90	20.592	2990	2993	3002	rBV2	207788	310898	4.93%	0.459%
91	21.645	3166	3172	3179	rBV2	1283127	2227742	35.33%	3.288%
92	24.798	3697	3708	3724	rBV	1699690	4890404	77.56%	7.218%
93	25.016	3737	3745	3761	rVB	853694	2314059	36.70%	3.415%

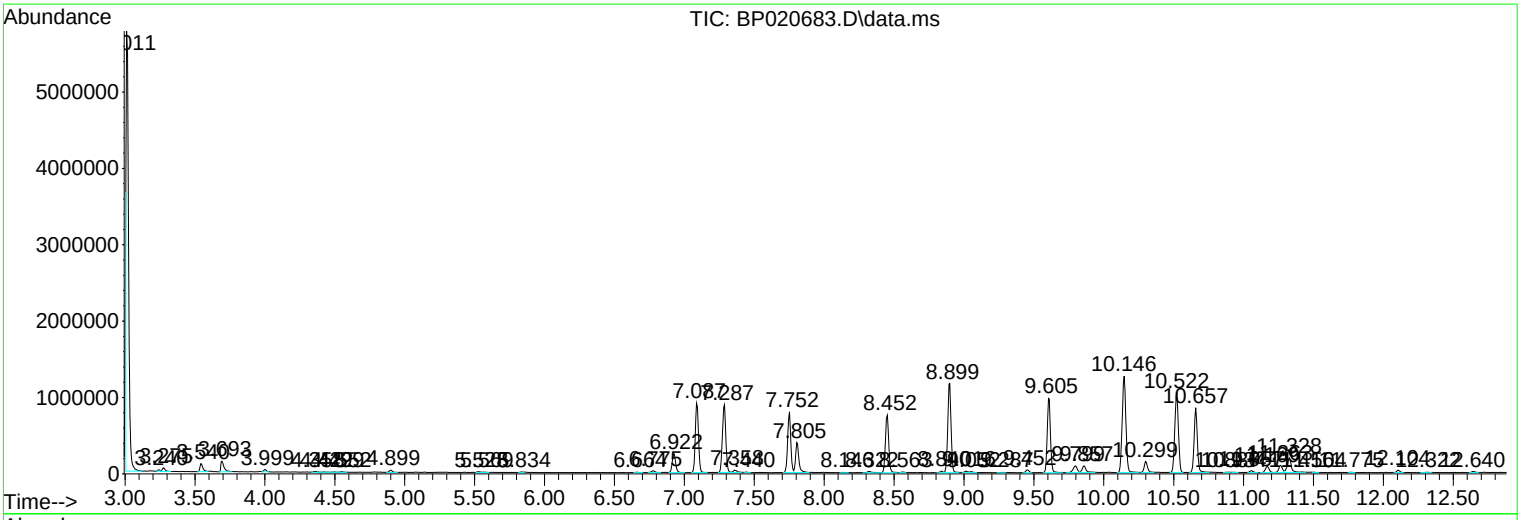
Sum of corrected areas: 67755406

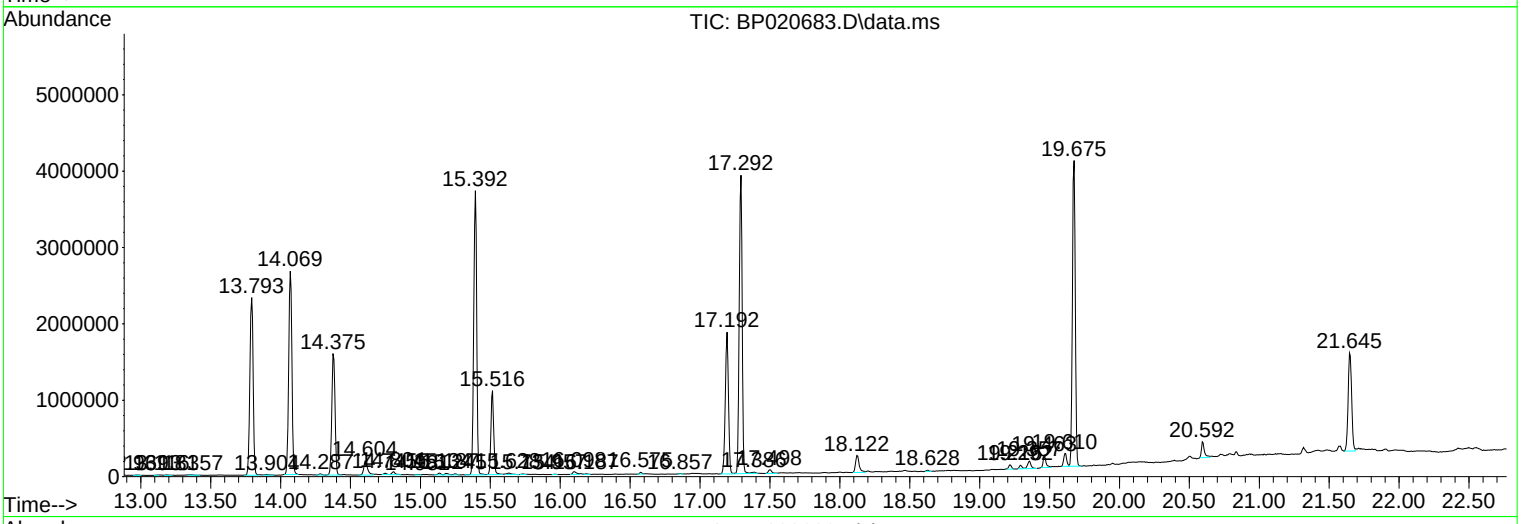
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020683.D
Acq On : 28 Jun 2024 06:58
Operator : MA/JU
Sample : P2898-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

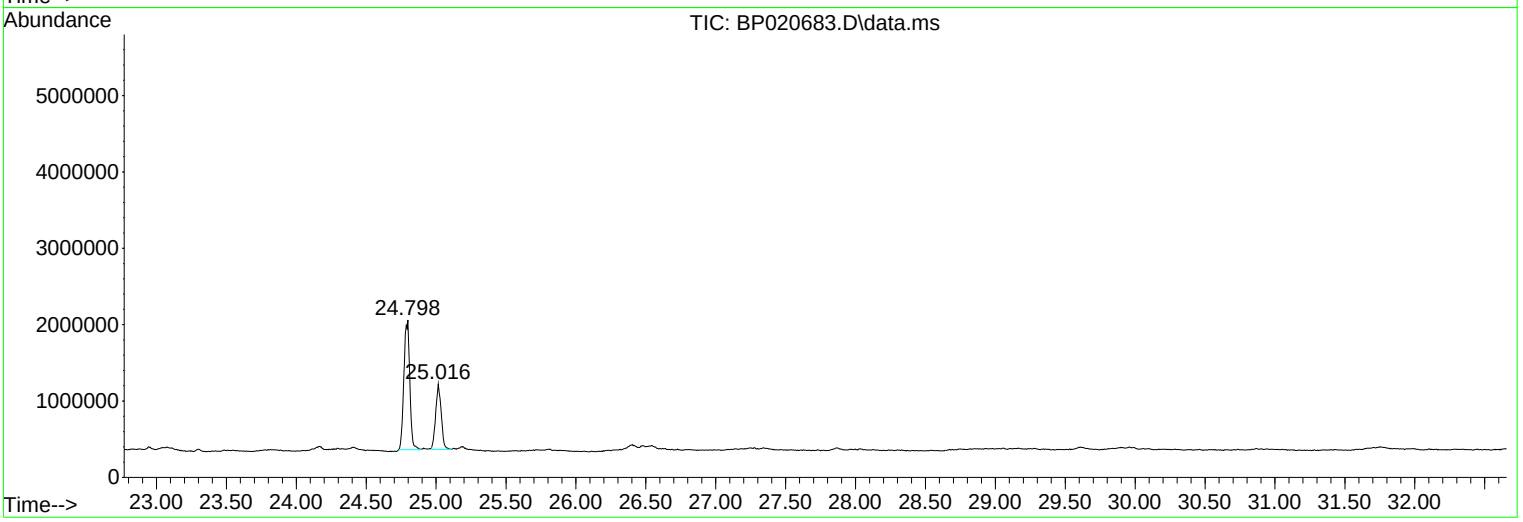
Instrument :
BNA_P
ClientSampleId :
E1G84

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.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\BP062724\
Data File : BP020683.D
Acq On : 28 Jun 2024 06:58
Operator : MA/JU
Sample : P2898-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G84

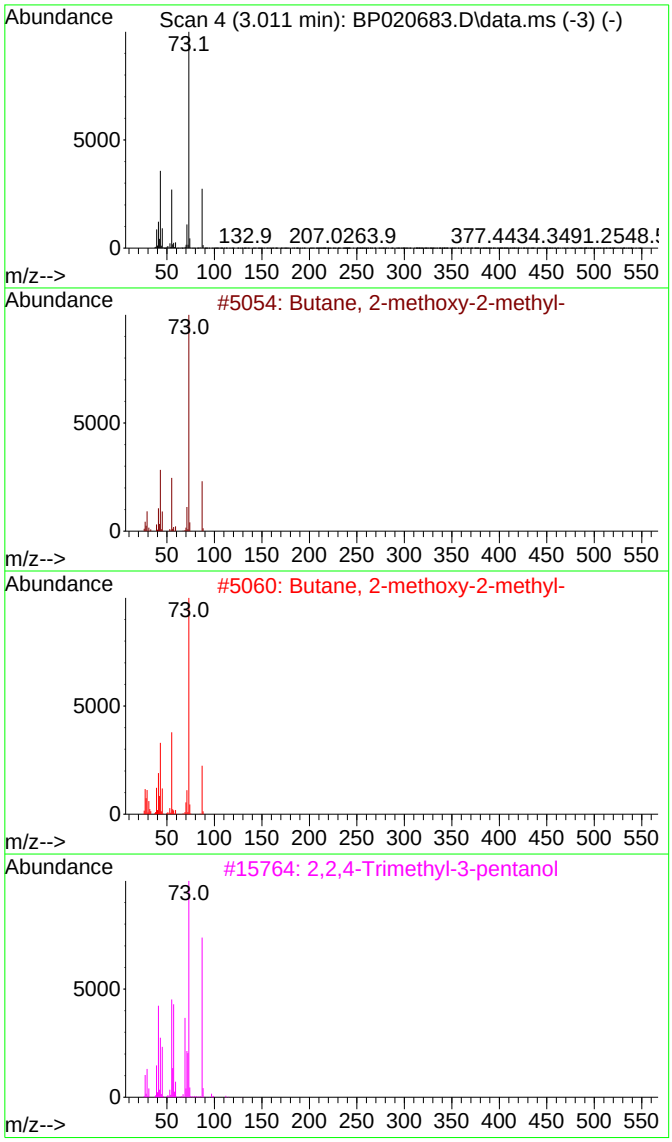
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

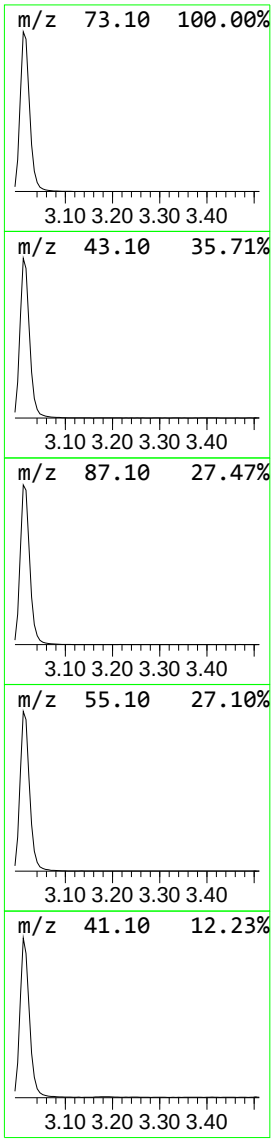
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.011	101.15 ng/ul	6305440	1,4-Dichlorobenzene-d4	7.752

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
5	1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	17





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020683.D
Acq On : 28 Jun 2024 06:58
Operator : MA/JU
Sample : P2898-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G84

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

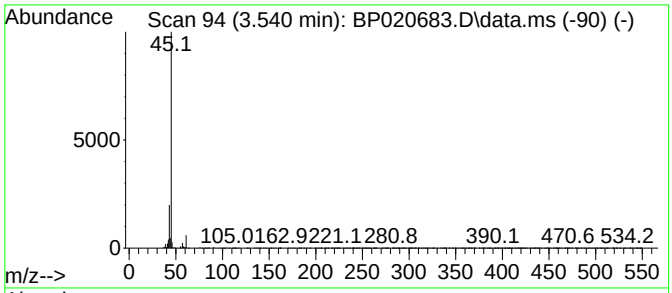
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

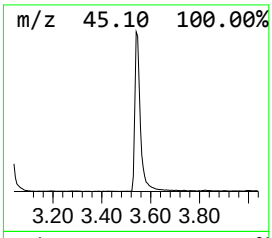
****

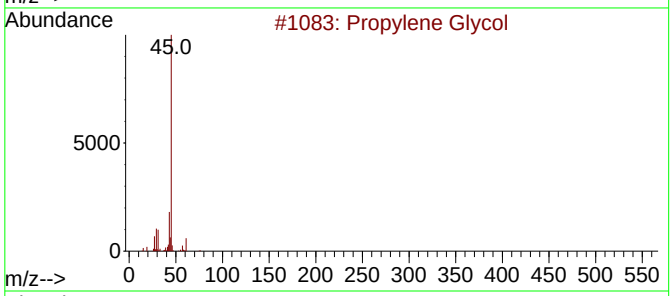
Peak Number 2 Propylene Glycol Concentration Rank 6

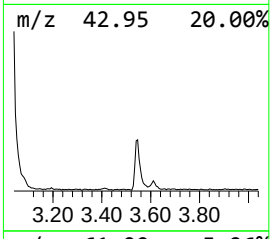
R.T.	EstConc	Area	Relative to ISTD	R.T.
3.540	2.57 ng/ul	160241	1,4-Dichlorobenzene-d4	7.752

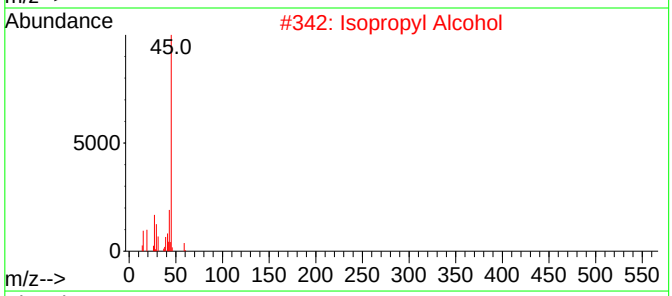
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propylene Glycol	76	C3H8O2	000057-55-6	90
2	Isopropyl Alcohol	60	C3H8O	000067-63-0	56
3	(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	40
4	Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
5	Isopropyl Alcohol	60	C3H8O	000067-63-0	9

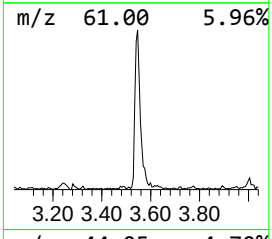


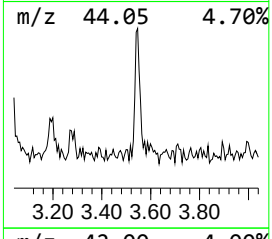


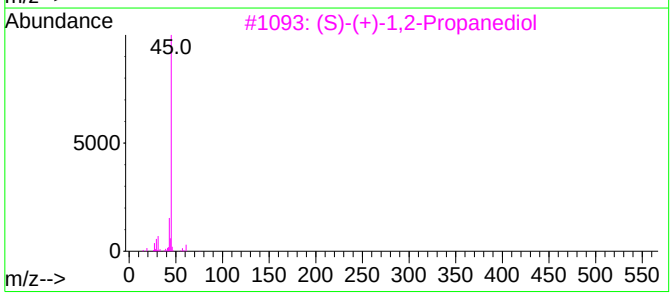


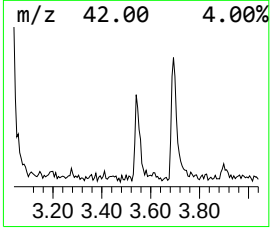












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020683.D
Acq On : 28 Jun 2024 06:58
Operator : MA/JU
Sample : P2898-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G84

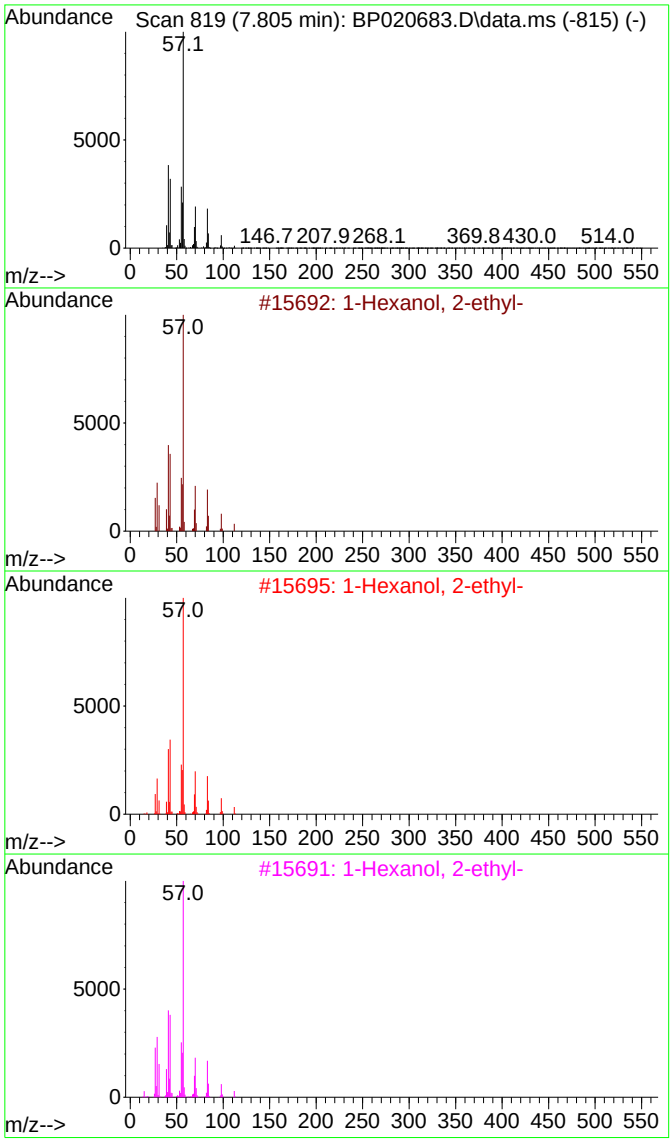
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

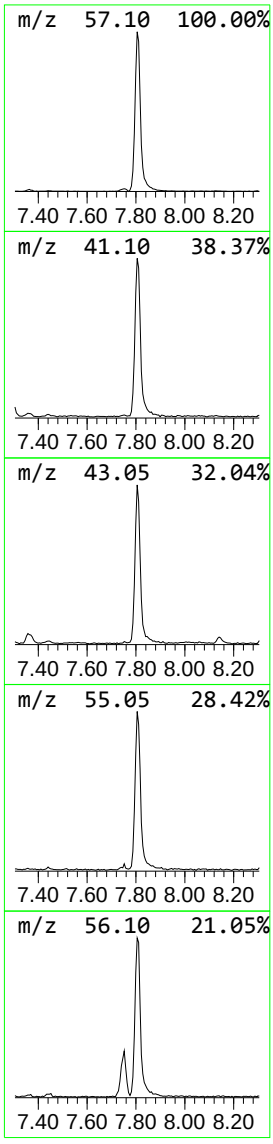
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1-Hexanol, 2-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.805	10.00 ng/ul	623564	1,4-Dichlorobenzene-d4	7.752

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
2	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
3	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72
4	1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	64
5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020683.D
Acq On : 28 Jun 2024 06:58
Operator : MA/JU
Sample : P2898-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G84

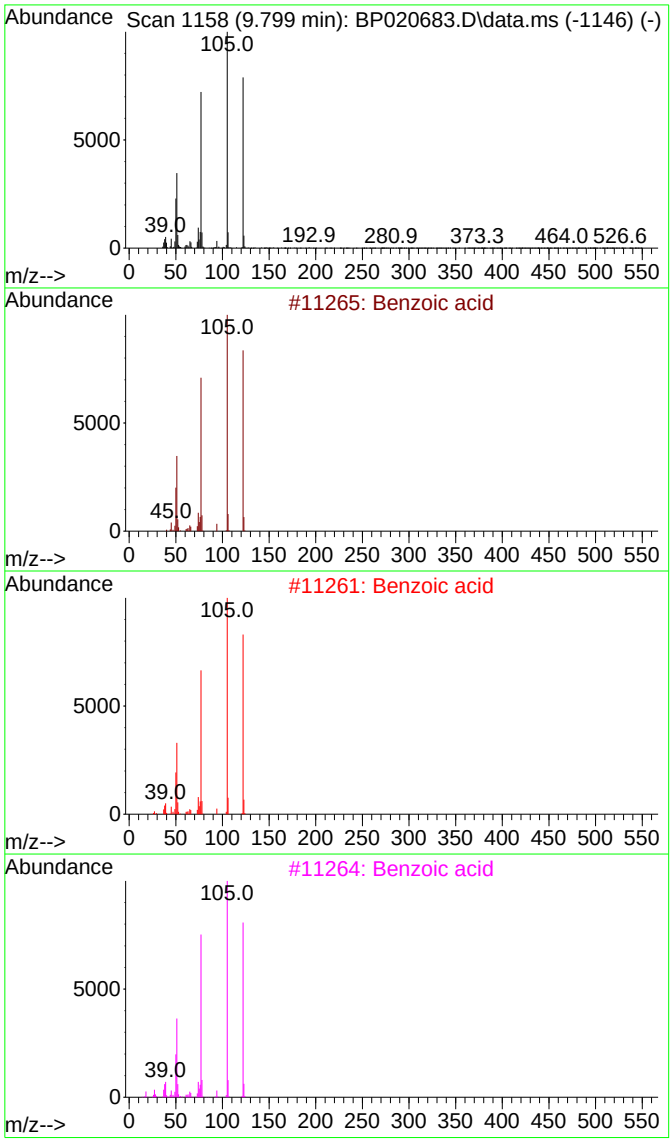
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

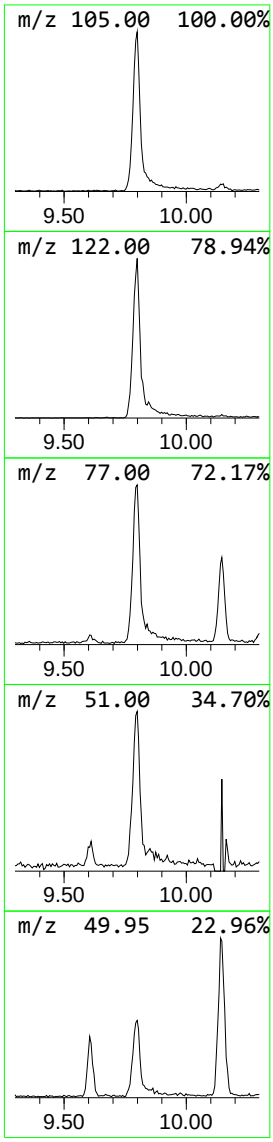
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.799	2.04 ng/ul	186512	Naphthalene-d8	10.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid	122	C7H6O2	000065-85-0	96
2	Benzoic acid	122	C7H6O2	000065-85-0	95
3	Benzoic acid	122	C7H6O2	000065-85-0	94
4	Heptanediamide, N,N'-di-benzoyloxy-	398	C21H22N2O6	1000253-26-4	91
5	Benzoic acid	122	C7H6O2	000065-85-0	90





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Data File : BP020683.D
Acq On : 28 Jun 2024 06:58
Operator : MA/JU
Sample : P2898-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G84

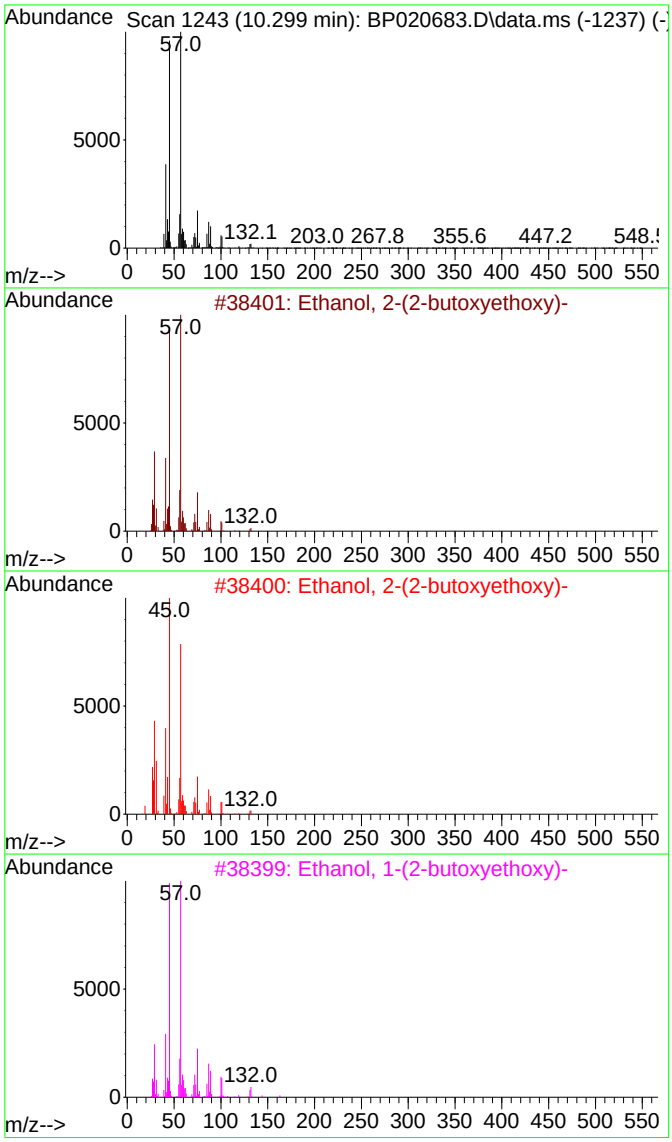
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

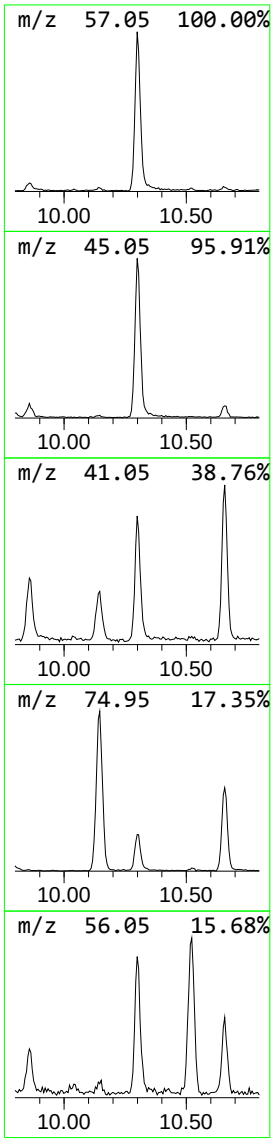
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.299	2.51 ng/ul	229380	Naphthalene-d8	10.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86
3	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
4	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	78
5	Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	78





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Data File : BP020683.D
Acq On : 28 Jun 2024 06:58
Operator : MA/JU
Sample : P2898-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G84

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

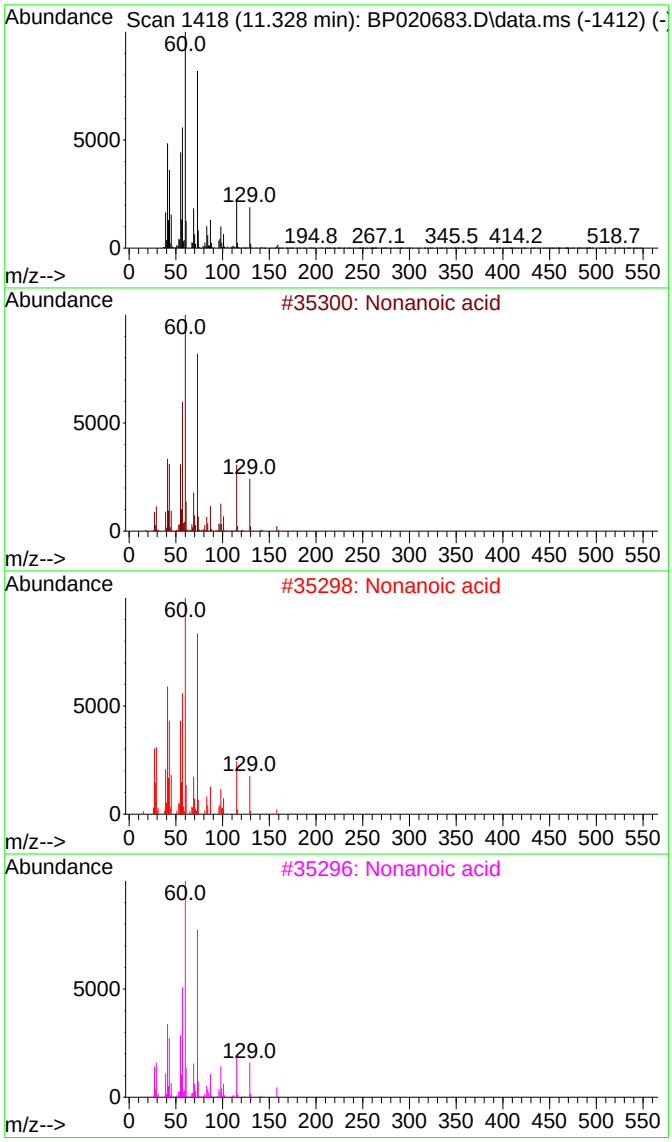
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

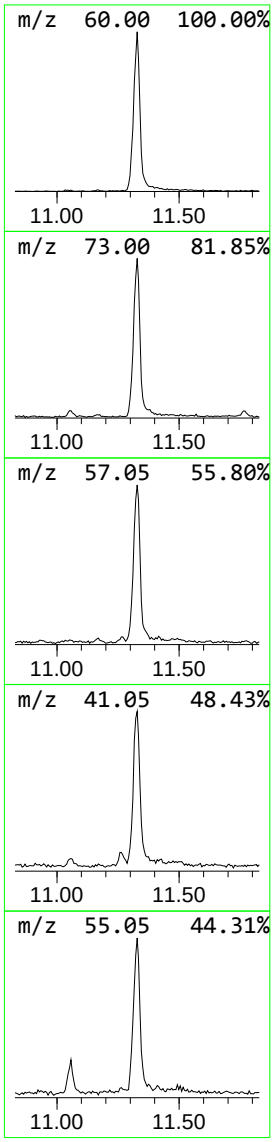
****

Peak Number 6 Nonanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.328	4.11 ng/ul	375163	Naphthalene-d8	10.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanoic acid	158	C9H18O2	000112-05-0	96
2	Nonanoic acid	158	C9H18O2	000112-05-0	91
3	Nonanoic acid	158	C9H18O2	000112-05-0	86
4	Nonanoic acid	158	C9H18O2	000112-05-0	72
5	n-Decanoic acid	172	C10H20O2	000334-48-5	72





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Data File : BP020683.D
Acq On : 28 Jun 2024 06:58
Operator : MA/JU
Sample : P2898-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G84

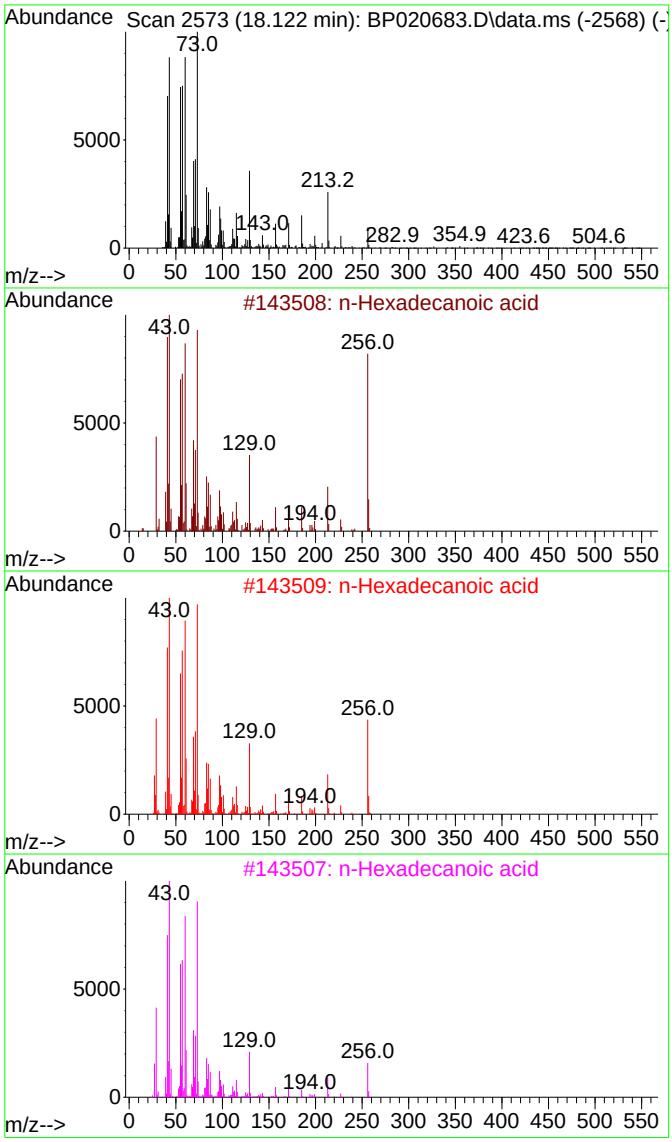
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

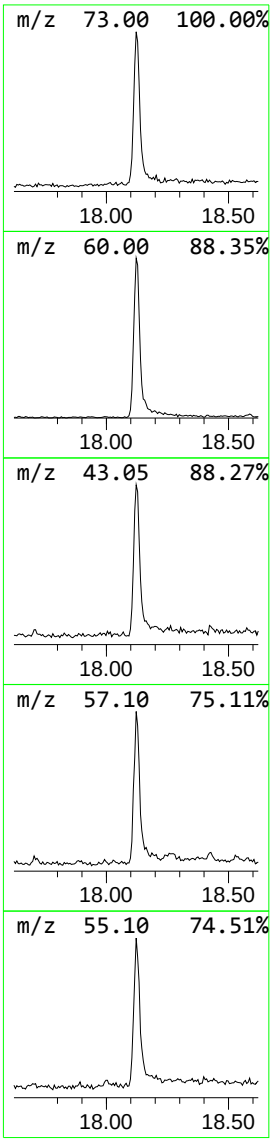
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.122	2.85 ng/ul	408056	Phenanthrene-d10	17.192

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	97
4	Tridecanoic acid	214	C13H26O2	000638-53-9	95
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95





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Data File : BP020683.D
Acq On : 28 Jun 2024 06:58
Operator : MA/JU
Sample : P2898-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G84

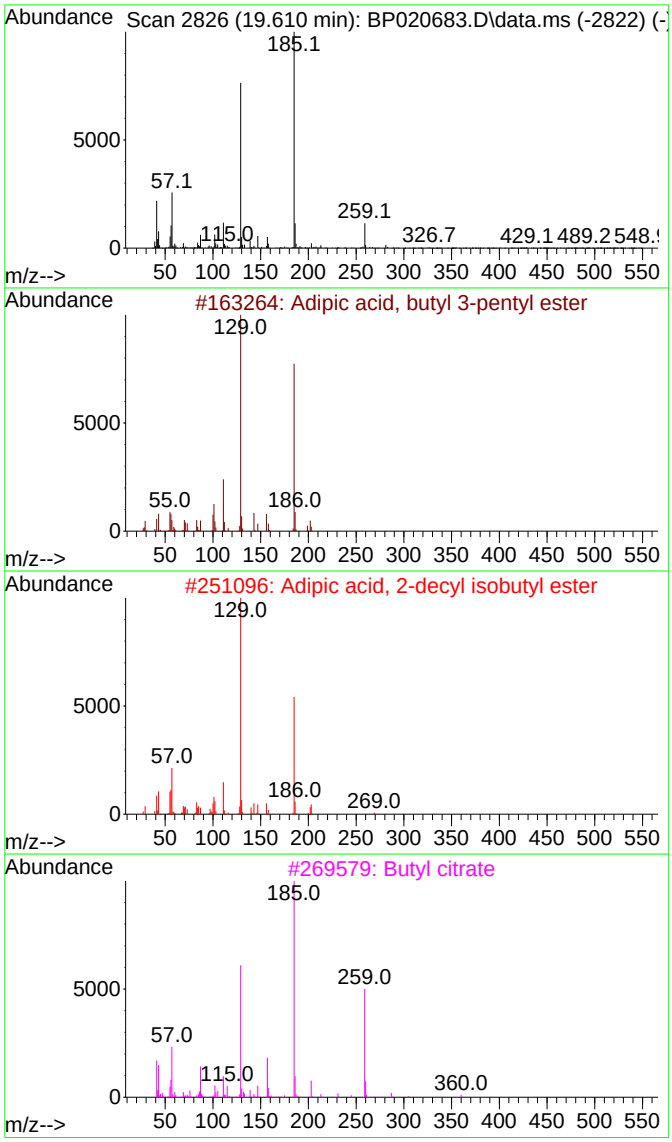
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

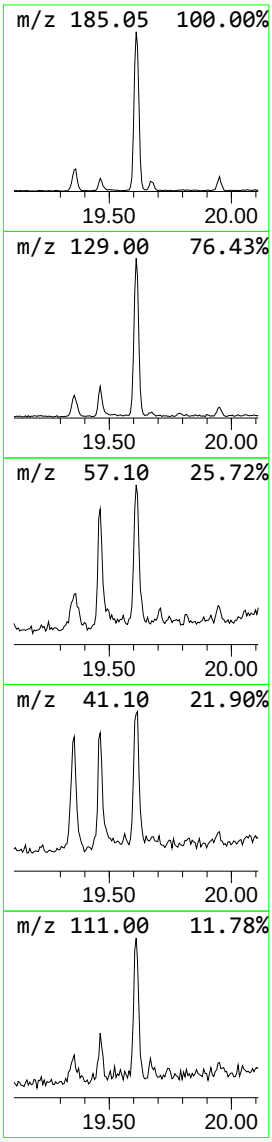
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Adipic acid, butyl 3-pentyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.610	2.25 ng/ul	250881	Chrysene-d12	21.645

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Adipic acid, butyl 3-pentyl ester	272	C15H28O4	1000324-60-5	78
2	Adipic acid, 2-decyl isobutyl ester	342	C20H38O4	1000353-59-6	78
3	Butyl citrate	360	C18H32O7	000077-94-1	74
4	Hexanedioic acid, bis(2-methylpr...	258	C14H26O4	000141-04-8	72
5	Adipic acid, 2-heptyl isobutyl e...	300	C17H32O4	1000353-64-3	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020683.D
Acq On : 28 Jun 2024 06:58
Operator : MA/JU
Sample : P2898-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G84

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

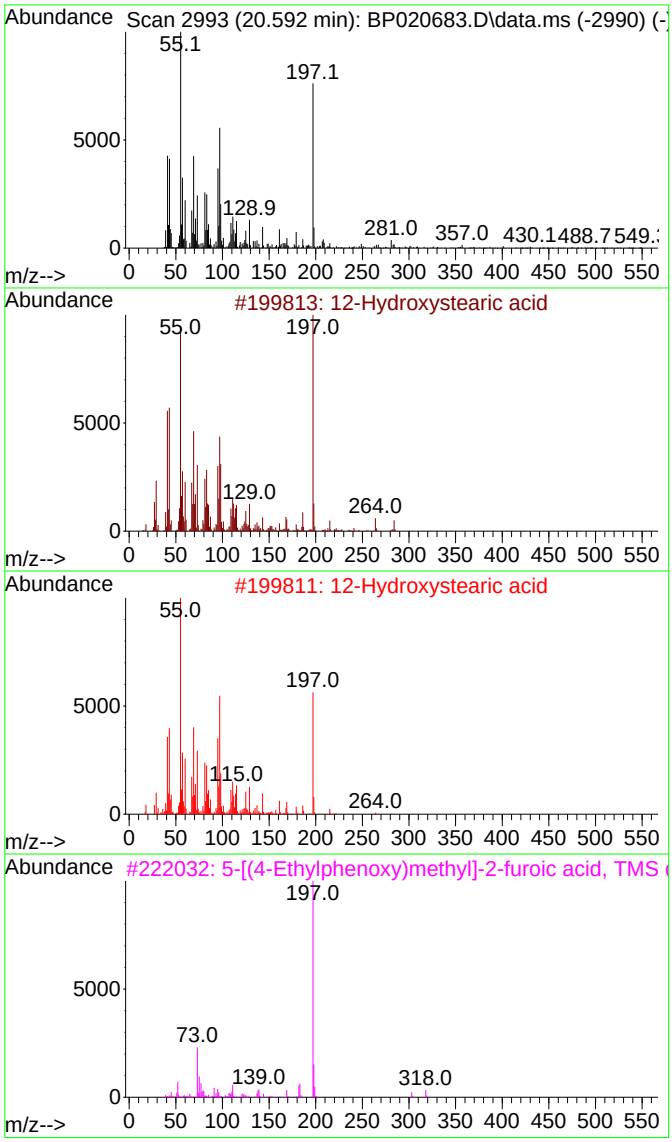
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

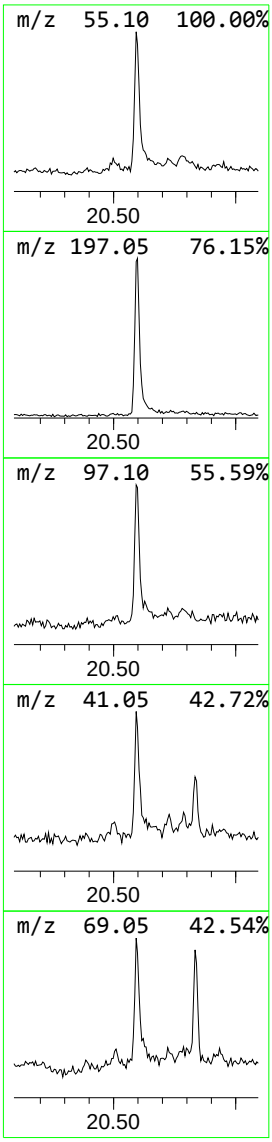
****

Peak Number 9 12-Hydroxystearic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.592	2.79 ng/ul	310898	Chrysene-d12	21.645

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	12-Hydroxystearic acid	300	C18H36O3	000106-14-9	83
2	12-Hydroxystearic acid	300	C18H36O3	000106-14-9	70
3	5-[(4-Ethylphenoxy)methyl]-2-fur...	318	C17H22O4Si	1000474-93-6	38
4	12-Hydroxystearic acid	300	C18H36O3	000106-14-9	32
5	Fumarylacetone diethoxime mono-TMS	314	C14H26N2O4Si	1000101-77-3	32





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020683.D
Acq On : 28 Jun 2024 06:58
Operator : MA/JU
Sample : P2898-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G84

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.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
Butane, 2-metho...	3.011	101.2	ng/ul	6305440	1	7.752	1246720	20.0
Propylene Glycol	3.540	2.6	ng/ul	160241	1	7.752	1246720	20.0
1-Hexanol, 2-et...	7.805	10.0	ng/ul	623564	1	7.752	1246720	20.0
Benzoic acid	9.799	2.0	ng/ul	186512	2	10.522	1825570	20.0
Ethanol, 2-(2-b...	10.299	2.5	ng/ul	229380	2	10.522	1825570	20.0
Nonanoic acid	11.328	4.1	ng/ul	375163	2	10.522	1825570	20.0
n-Hexadecanoic ...	18.122	2.9	ng/ul	408056	4	17.192	2859470	20.0
Adipic acid, bu...	19.610	2.3	ng/ul	250881	5	21.645	2227740	20.0
12-Hydroxystear...	20.592	2.8	ng/ul	310898	5	21.645	2227740	20.0