

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
 Data File : BP020717.D
 Acq On : 01 Jul 2024 14:03
 Operator : MA/JU
 Sample : P2871-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4C53

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: BP020717.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.016	3	5	20	rVB	1878691	2143045	64.67%	3.930%
2	3.275	46	49	57	rVB	38944	49707	1.50%	0.091%
3	3.546	91	95	107	rVB	115890	164013	4.95%	0.301%
4	3.693	116	120	133	rBV	304627	469539	14.17%	0.861%
5	4.005	169	173	177	rVB	15362	21411	0.65%	0.039%
6	4.363	231	234	238	rVB5	5649	7467	0.23%	0.014%
7	4.505	254	258	261	rBV4	4454	7271	0.22%	0.013%
8	4.552	263	266	271	rVB4	8573	11503	0.35%	0.021%
9	4.781	304	305	310	rBV5	3040	4956	0.15%	0.009%
10	4.905	320	326	332	rBV	28622	45920	1.39%	0.084%
11	5.146	363	367	373	rVB	4900	9035	0.27%	0.017%
12	5.863	483	489	503	rVB	22469	39021	1.18%	0.072%
13	6.369	569	575	580	rBV8	4705	10280	0.31%	0.019%
14	6.775	639	644	651	rVB5	8674	18019	0.54%	0.033%
15	6.928	662	670	681	rBV	754346	1155339	34.86%	2.118%
16	7.087	690	697	711	rVB	525581	854571	25.79%	1.567%
17	7.287	724	731	739	rBV	779475	1154934	34.85%	2.118%
18	7.357	739	743	753	rVB	38974	74553	2.25%	0.137%
19	7.740	801	808	816	rBV	798421	1268925	38.29%	2.327%
20	7.810	816	820	829	rVB2	13904	25524	0.77%	0.047%
21	8.328	904	908	914	rBV9	3344	5589	0.17%	0.010%
22	8.440	920	927	938	rBV	877003	1468664	44.32%	2.693%
23	8.893	996	1004	1013	rBV	672141	1124560	33.93%	2.062%
24	9.010	1021	1024	1027	rBV5	4632	6541	0.20%	0.012%
25	9.051	1027	1031	1037	rVB8	9510	15809	0.48%	0.029%
26	9.451	1091	1099	1108	rBV	87968	135880	4.10%	0.249%
27	9.604	1116	1125	1137	rBV	565848	965258	29.13%	1.770%
28	9.710	1140	1143	1147	rVV6	3685	5564	0.17%	0.010%
29	9.781	1151	1155	1160	rVV7	6954	15936	0.48%	0.029%
30	9.840	1160	1165	1174	rVV7	10853	24434	0.74%	0.045%
31	10.140	1207	1216	1227	rBV	892534	1547509	46.70%	2.838%
32	10.292	1239	1242	1249	rVB8	4420	8803	0.27%	0.016%
33	10.516	1270	1280	1287	rBV	1156343	1800199	54.32%	3.301%
34	10.651	1296	1303	1318	rVV	515996	944256	28.49%	1.731%
35	11.045	1360	1370	1378	rBV2	34552	67346	2.03%	0.123%
36	11.257	1398	1406	1413	rBV	175069	324332	9.79%	0.595%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
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37	12.004	1529	1533	1538	rBV4	6394	10597	0.32%	0.019%
38	12.092	1543	1548	1554	rVB2	16644	25758	0.78%	0.047%
39	12.169	1556	1561	1564	rBV7	2948	5316	0.16%	0.010%
40	12.234	1566	1572	1580	rBV2	23880	51991	1.57%	0.095%
41	12.639	1634	1641	1647	rBV8	8435	17680	0.53%	0.032%
42	12.722	1650	1655	1660	rBV8	4558	9414	0.28%	0.017%
43	12.969	1692	1697	1705	rBV9	4800	11505	0.35%	0.021%
44	13.145	1722	1727	1728	rBV5	3437	5499	0.17%	0.010%
45	13.322	1755	1757	1763	rVB7	6157	8558	0.26%	0.016%
46	13.545	1786	1795	1798	rBV9	7752	15850	0.48%	0.029%
47	13.675	1813	1817	1820	rBV6	3115	4874	0.15%	0.009%
48	13.775	1826	1834	1844	rBV	1701534	2320647	70.03%	4.255%
49	13.839	1844	1845	1849	rVB3	6434	7269	0.22%	0.013%
50	14.022	1870	1876	1877	rBV2	21880	35676	1.08%	0.065%
51	14.063	1877	1883	1897	rVB	1899546	2760197	83.29%	5.061%
52	14.245	1908	1914	1915	rBV5	6531	10684	0.32%	0.020%
53	14.375	1927	1936	1942	rBV	1561220	2291628	69.15%	4.202%
54	14.433	1942	1946	1951	rVB7	4539	9893	0.30%	0.018%
55	14.592	1965	1973	1988	rVV	693360	1360920	41.07%	2.495%
56	14.739	1994	1998	2003	rBV3	13941	22025	0.66%	0.040%
57	14.798	2004	2008	2013	rVB5	20785	28688	0.87%	0.053%
58	14.963	2033	2036	2041	rVB6	4686	7463	0.23%	0.014%
59	15.186	2068	2074	2077	rBV6	9592	20129	0.61%	0.037%
60	15.222	2077	2080	2081	rBV3	4639	4795	0.14%	0.009%
61	15.375	2100	2106	2121	rVB	2233451	3250890	98.10%	5.961%
62	15.504	2122	2128	2137	rBV	549257	779484	23.52%	1.429%
63	15.628	2146	2149	2159	rVB7	15017	35449	1.07%	0.065%
64	15.980	2202	2209	2211	rBV7	12940	29662	0.90%	0.054%
65	16.051	2213	2221	2228	rVV10	29972	87080	2.63%	0.160%
66	16.122	2228	2233	2236	rVB5	13151	22770	0.69%	0.042%
67	16.298	2257	2263	2268	rBV9	14711	27741	0.84%	0.051%
68	16.357	2269	2273	2275	rBV4	15885	17193	0.52%	0.032%
69	16.569	2303	2309	2316	rBV5	56301	111431	3.36%	0.204%
70	16.951	2367	2374	2388	rVB9	50982	189217	5.71%	0.347%
71	17.086	2393	2397	2399	rBV3	31936	35658	1.08%	0.065%
72	17.180	2405	2413	2418	rBV	1576970	2269074	68.47%	4.161%
73	17.222	2418	2420	2424	rVV	91715	120268	3.63%	0.221%
74	17.280	2424	2430	2442	rVB	1968879	3025037	91.28%	5.547%
75	17.374	2442	2446	2451	rBV6	23118	47880	1.44%	0.088%
76	17.586	2477	2482	2490	rBV2	113198	184353	5.56%	0.338%

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77	17.710	2500	2503	2507	rVB4	19654	24626	0.74%	0.045%
78	17.851	2525	2527	2529	rBV3	18873	20048	0.60%	0.037%
79	18.127	2569	2574	2583	rBV	570935	874106	26.38%	1.603%
80	18.292	2598	2602	2604	rBV5	24235	39175	1.18%	0.072%
81	18.333	2605	2609	2620	rVB2	83647	196202	5.92%	0.360%
82	18.610	2653	2656	2662	rBV6	39308	72977	2.20%	0.134%
83	19.027	2723	2727	2730	rBV4	40795	63642	1.92%	0.117%
84	19.086	2734	2737	2741	rVV3	31209	43781	1.32%	0.080%
85	19.321	2772	2777	2781	rBV	162259	286353	8.64%	0.525%
86	19.398	2785	2790	2792	rVV3	152707	221329	6.68%	0.406%
87	19.416	2792	2793	2795	rVV	166434	125745	3.79%	0.231%
88	19.433	2795	2796	2797	rVV	154367	100106	3.02%	0.184%
89	19.457	2797	2800	2809	rVB2	284144	462462	13.96%	0.848%
90	19.668	2830	2836	2845	rVB	2163797	3313877	100.00%	6.077%
91	20.139	2912	2916	2917	rBV3	71409	90163	2.72%	0.165%
92	20.921	3044	3049	3054	rBV	312522	481320	14.52%	0.883%
93	21.315	3112	3116	3122	rBV	1077879	1676371	50.59%	3.074%
94	21.645	3166	3172	3178	rBV2	1166031	2209428	66.67%	4.051%
95	21.733	3180	3187	3188	rBV2	678641	940262	28.37%	1.724%
96	22.568	3321	3329	3338	rVB3	563997	1454386	43.89%	2.667%
97	24.156	3593	3599	3604	rBV	358594	730488	22.04%	1.339%
98	24.221	3605	3610	3618	rVB4	226492	542583	16.37%	0.995%
99	24.780	3698	3705	3715	rVB2	1160670	2966758	89.53%	5.440%
100	25.009	3737	3744	3757	rVB	853390	2323626	70.12%	4.261%

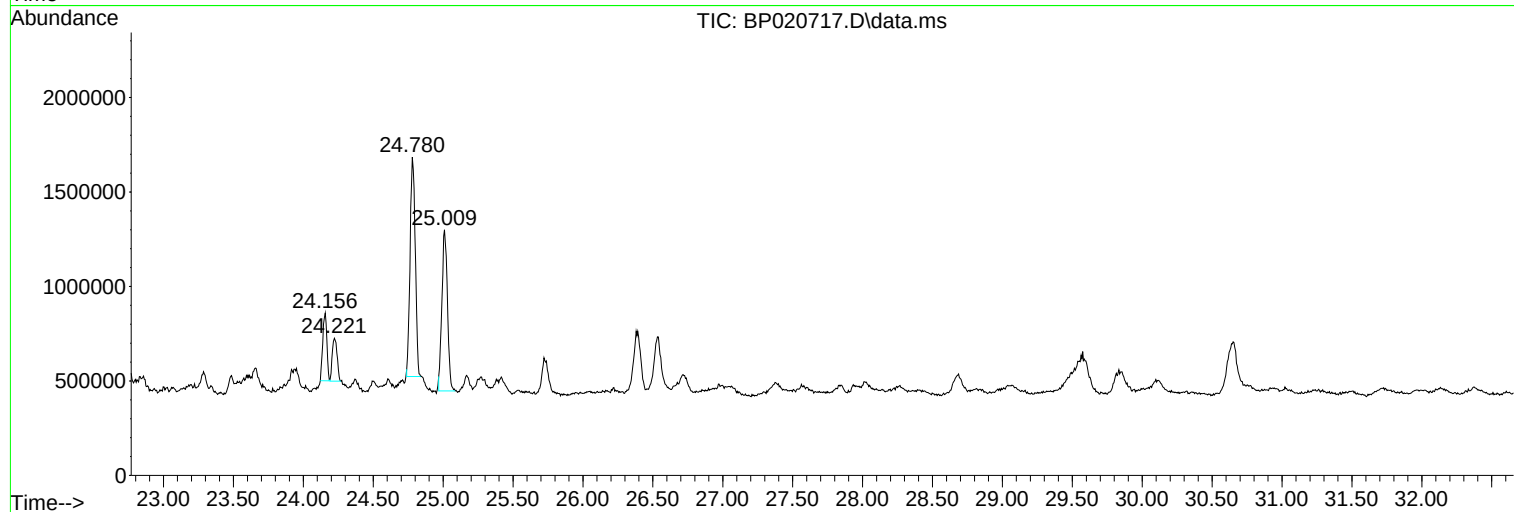
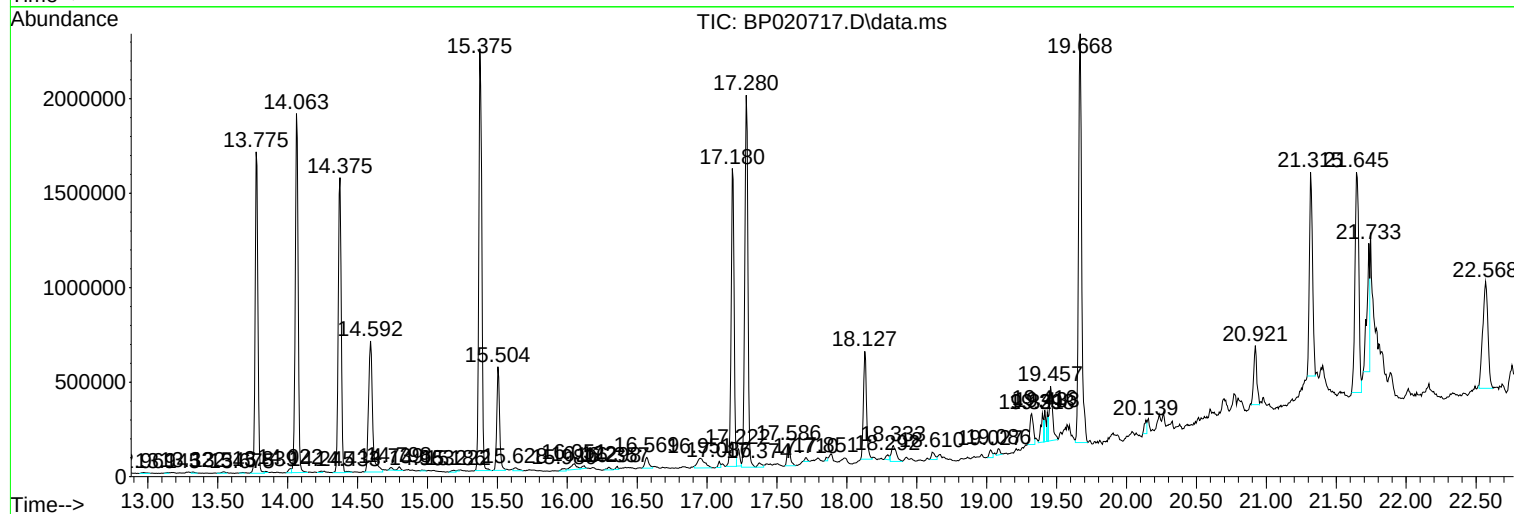
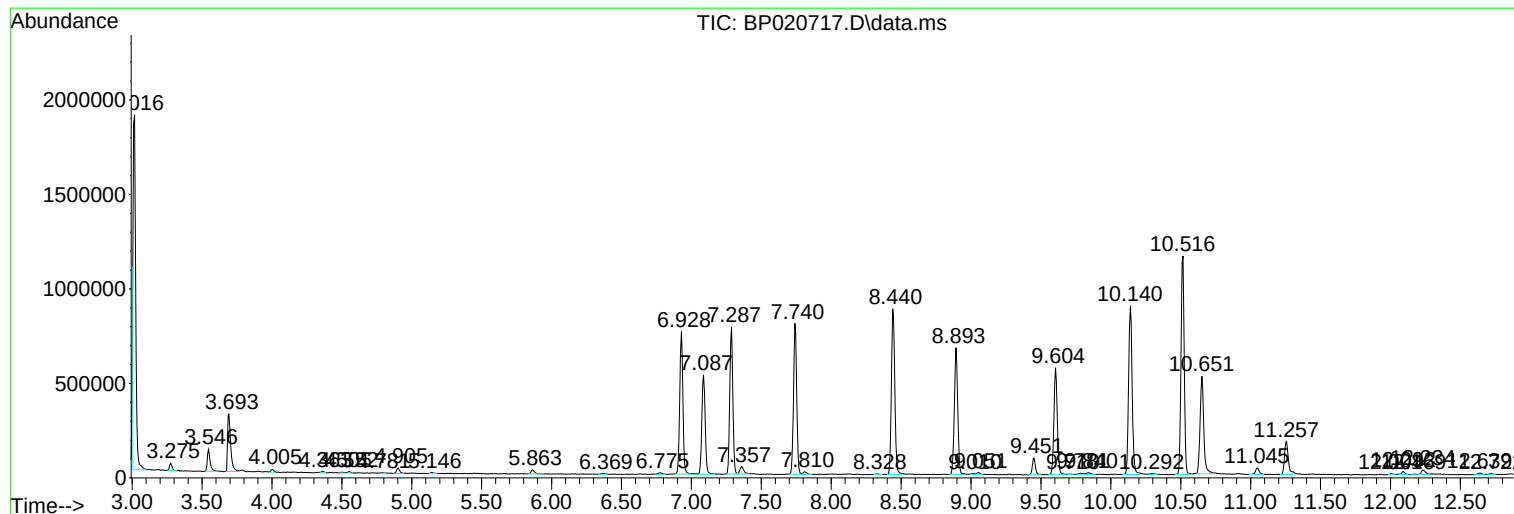
Sum of corrected areas: 54535770

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ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4C53

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



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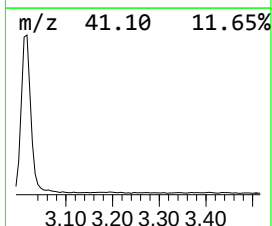
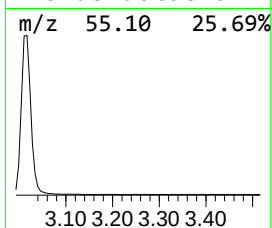
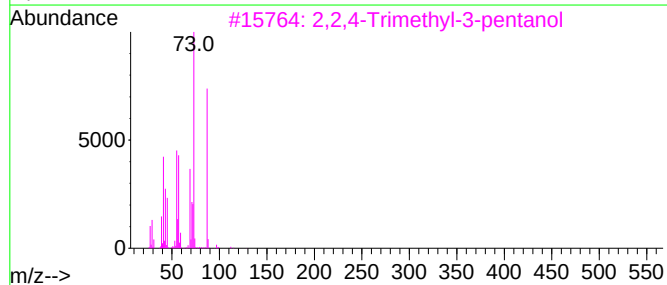
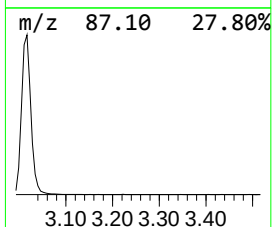
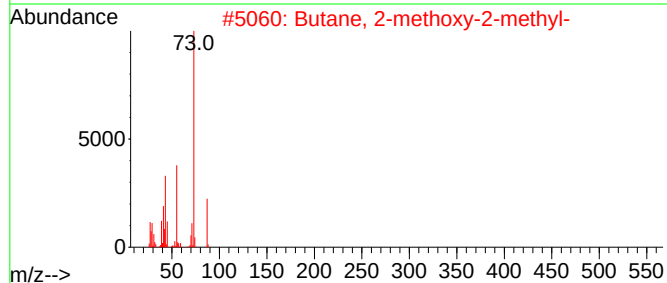
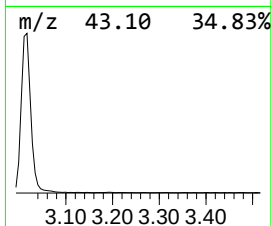
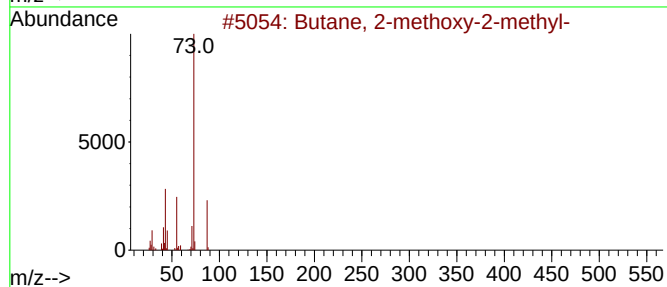
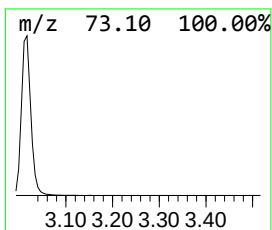
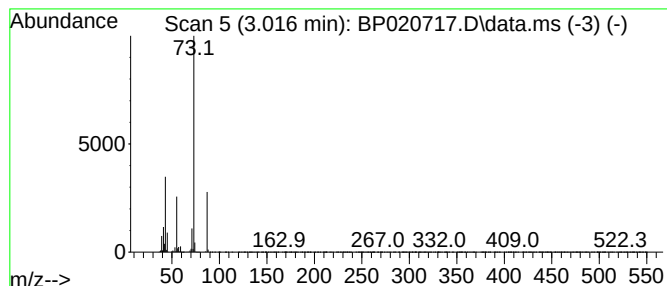
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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.
3.016	33.78 ng/ul	2143050	1,4-Dichlorobenzene-d4	7.740

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	25
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23



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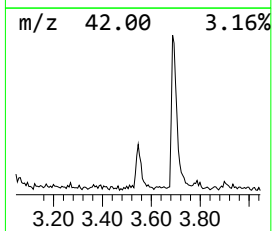
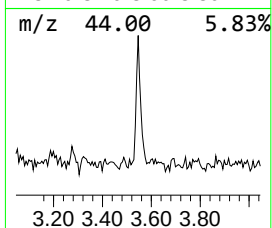
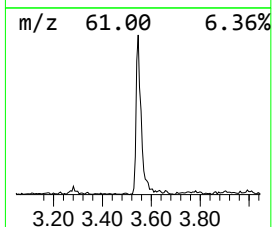
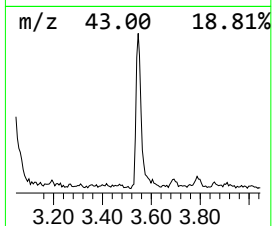
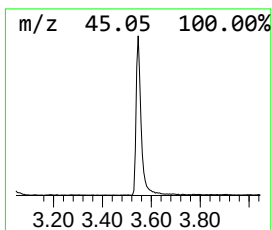
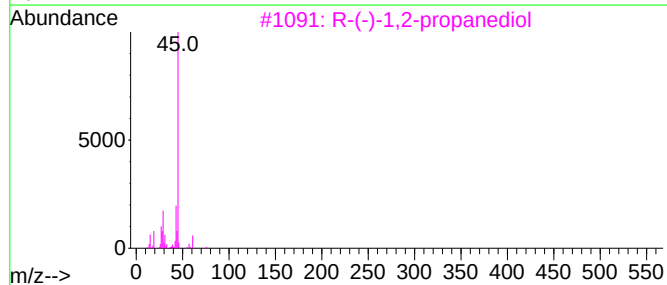
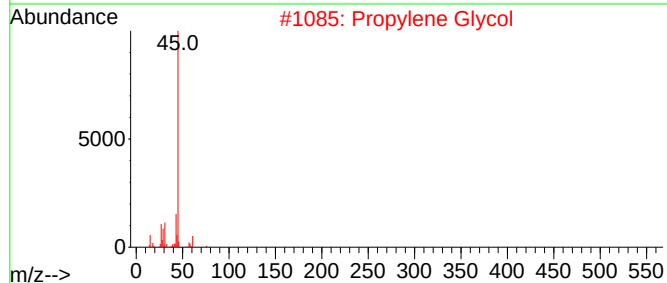
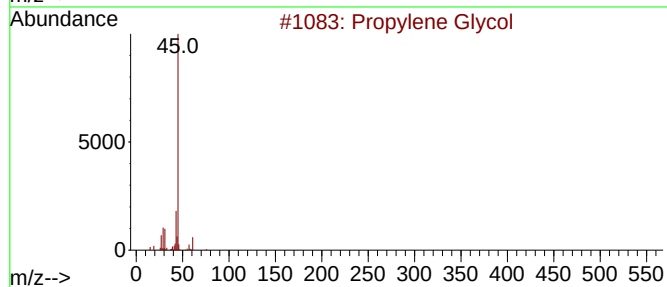
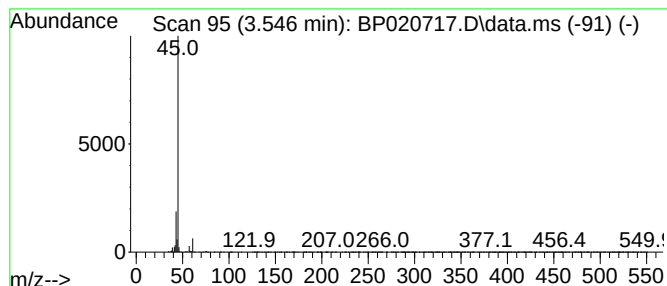
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Propylene Glycol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.546	2.59 ng/ul	164013	1,4-Dichlorobenzene-d4	7.740

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	86
2			Propylene Glycol	76	C3H8O2	000057-55-6	78
3			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	72
4			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	56
5			Isopropyl Alcohol	60	C3H8O	000067-63-0	50



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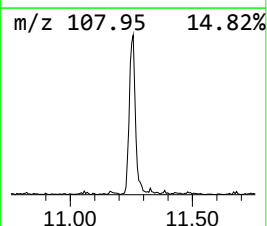
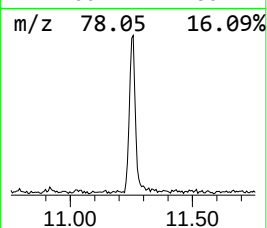
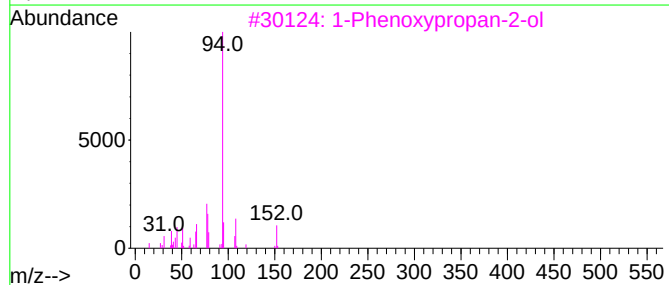
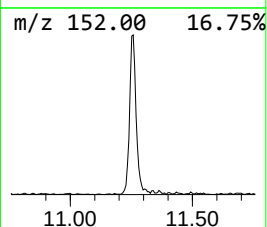
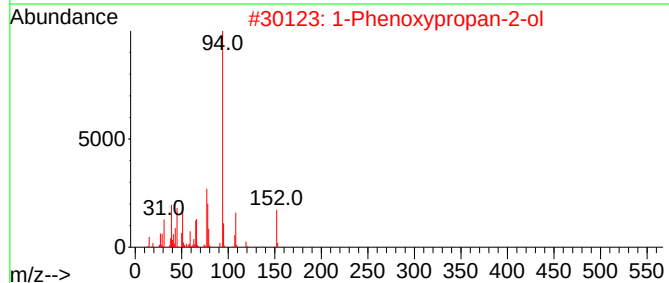
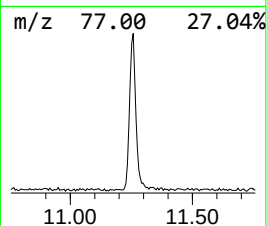
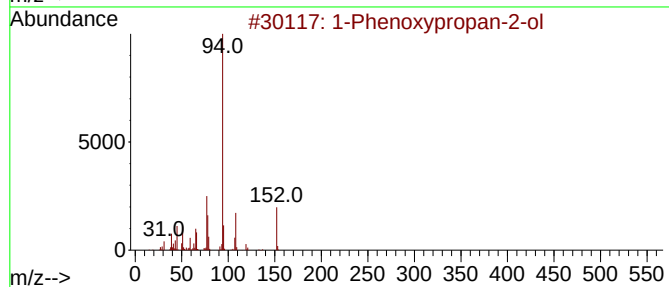
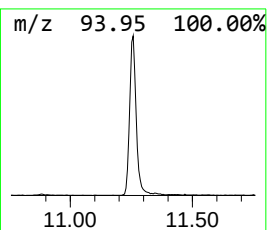
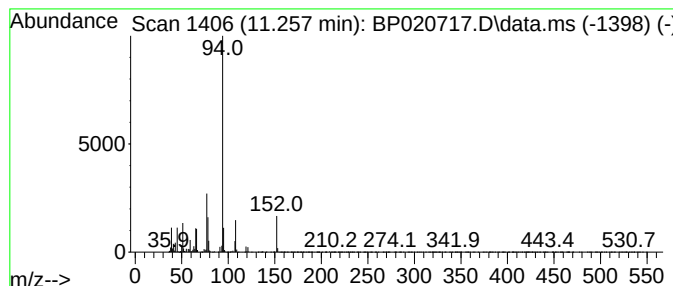
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.257	3.60 ng/ul	324332	Naphthalene-d8	10.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020717.D
Acq On : 01 Jul 2024 14:03
Operator : MA/JU
Sample : P2871-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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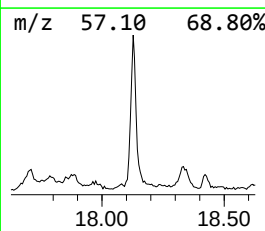
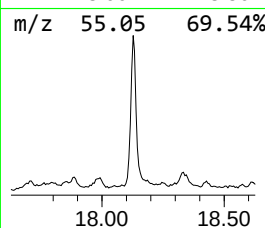
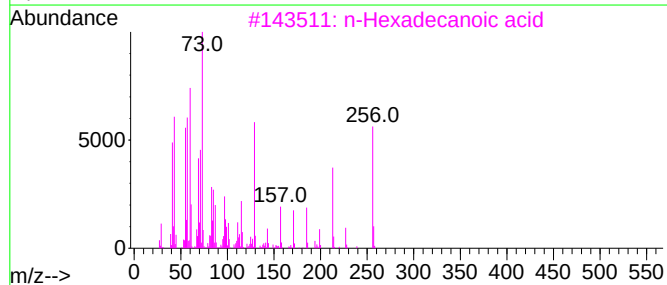
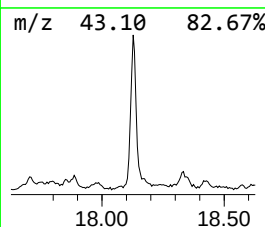
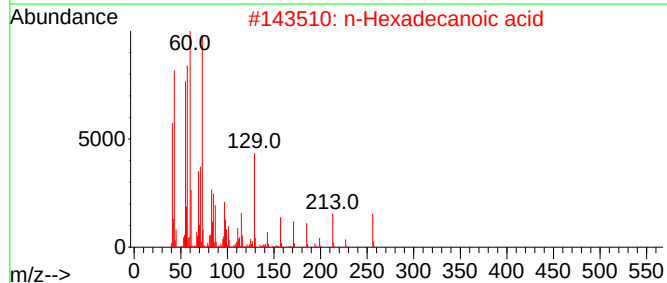
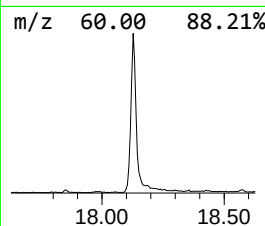
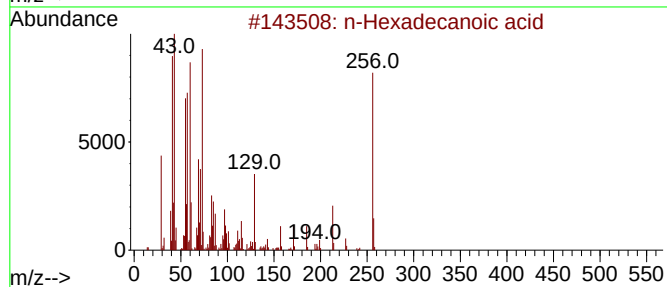
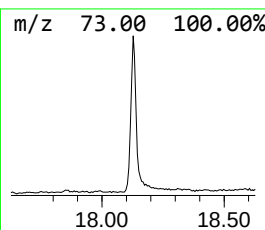
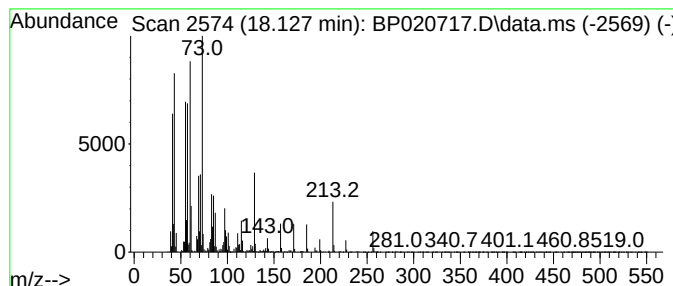
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.127	7.70 ng/ul	874106	Phenanthrene-d10	17.180

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			Tridecanoic acid	214	C13H26O2	000638-53-9	95
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020717.D
Acq On : 01 Jul 2024 14:03
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Sample : P2871-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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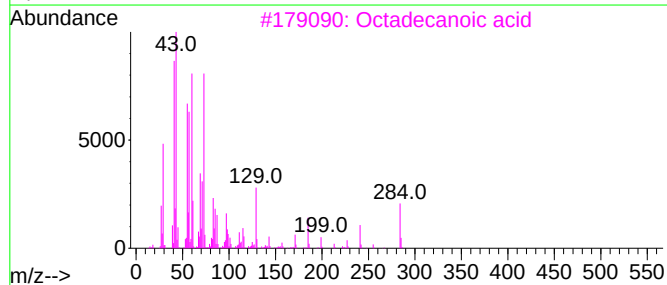
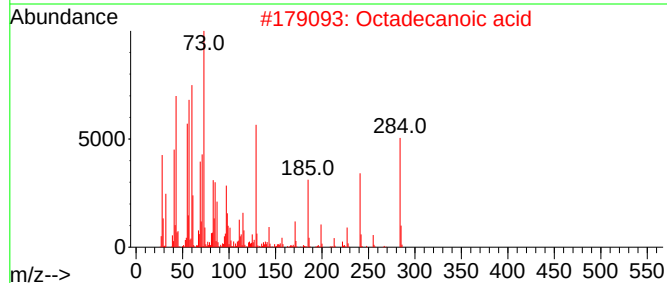
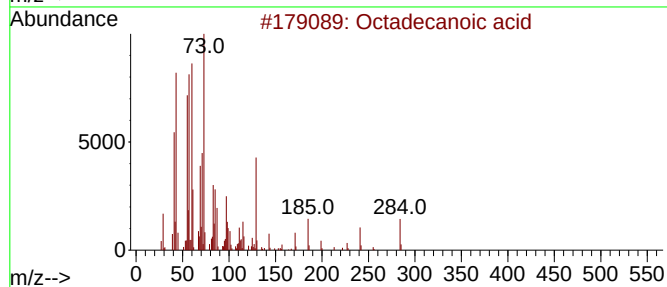
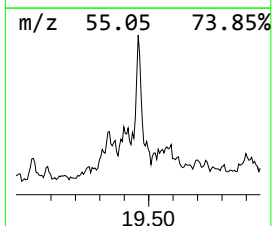
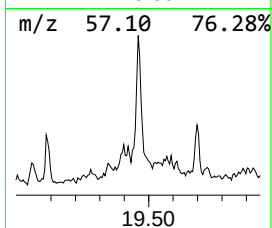
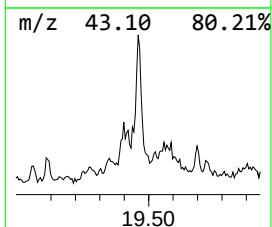
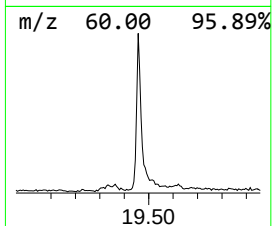
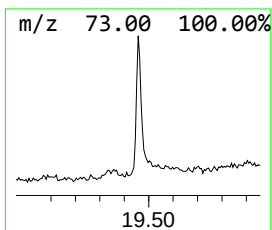
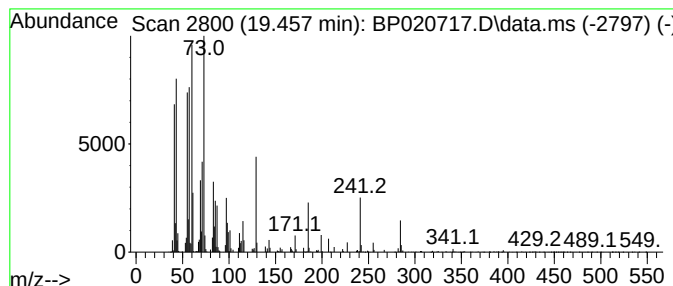
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.457	4.19 ng/ul	462462	Chrysene-d12	21.645

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	98
3			Octadecanoic acid	284	C18H36O2	000057-11-4	95
4			Octadecanoic acid	284	C18H36O2	000057-11-4	93
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020717.D
Acq On : 01 Jul 2024 14:03
Operator : MA/JU
Sample : P2871-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

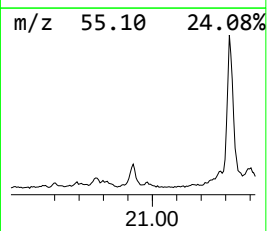
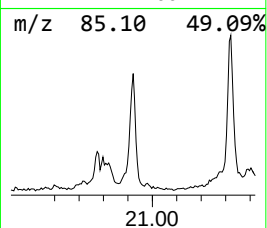
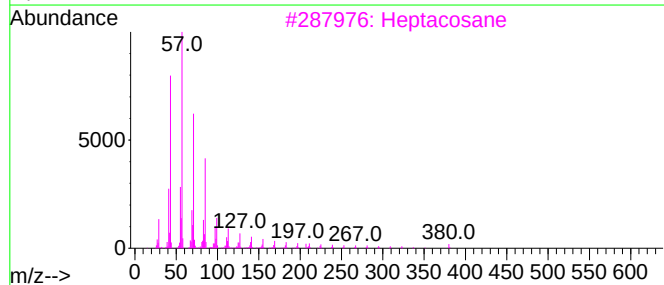
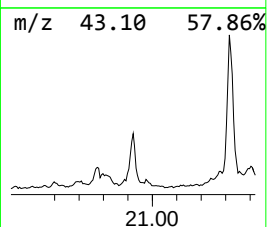
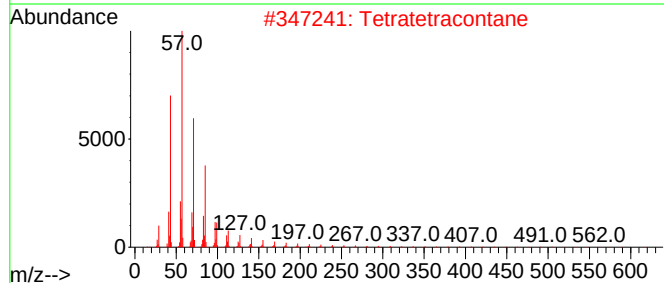
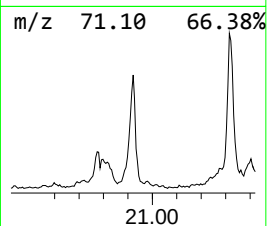
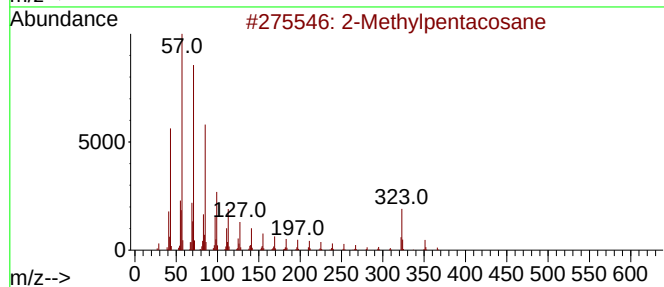
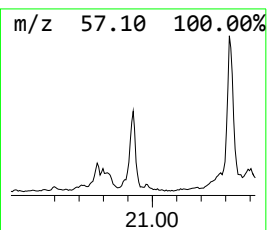
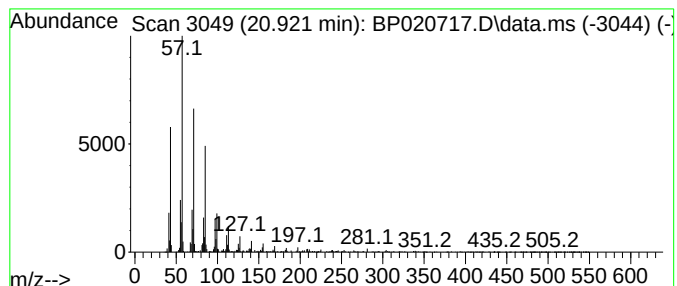
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.921	4.36 ng/ul	481320	Chrysene-d12	21.645	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylpentacosane	366	C26H54	000629-87-8	93
2	Tetratetracontane	619	C44H90	007098-22-8	91
3	Heptacosane	380	C27H56	000593-49-7	91
4	Hexacosane, 9-octyl-	479	C34H70	055429-83-9	91
5	Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020717.D
Acq On : 01 Jul 2024 14:03
Operator : MA/JU
Sample : P2871-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

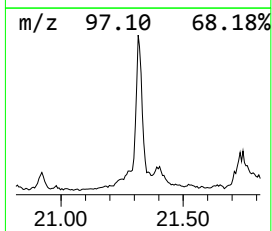
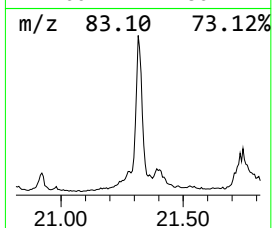
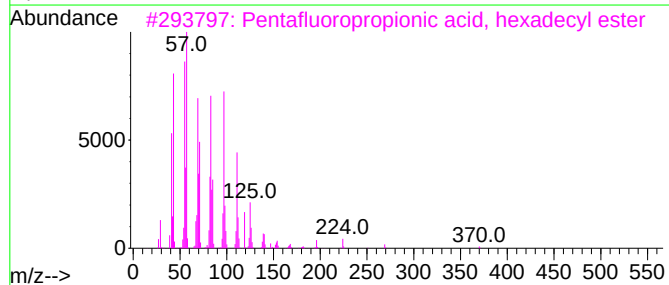
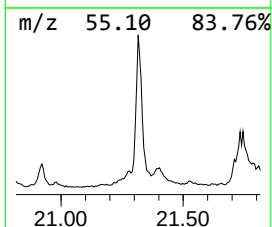
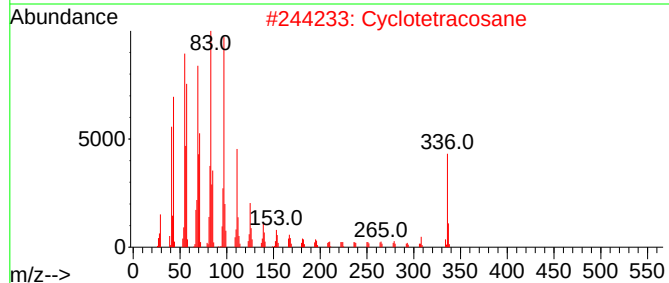
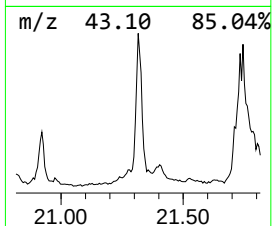
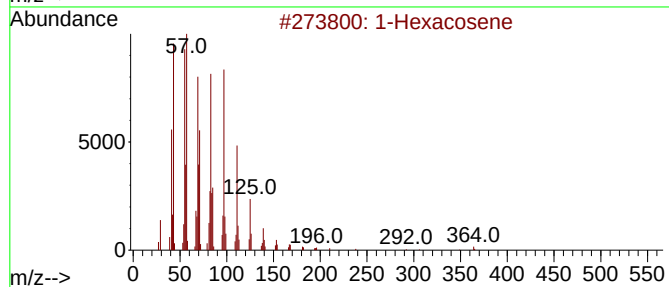
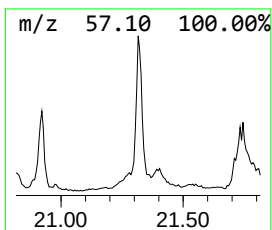
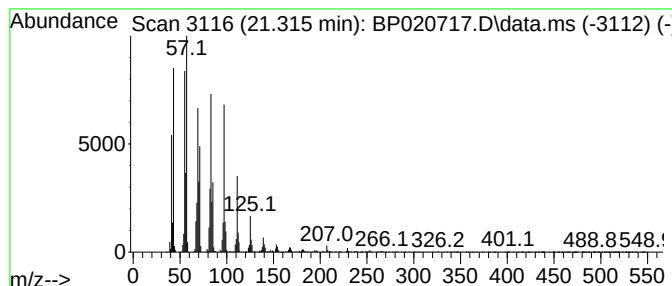
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.315	15.17 ng/ul	1676370	Chrysene-d12		21.645	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Hexacosene		364	C26H52	018835-33-1	95
2	Cyclotetracosane		336	C24H48	000297-03-0	95
3	Pentafluoropropionic acid, hexadecyl ester		388	C19H33F5O2	006222-07-7	94
4	Heptadecyl heptafluorobutyrate		452	C21H35F7O2	959085-66-6	94
5	Eicosyl pentafluoropropionate		444	C23H41F5O2	1000351-80-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020717.D
Acq On : 01 Jul 2024 14:03
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Sample : P2871-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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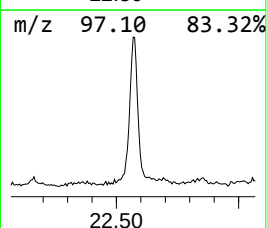
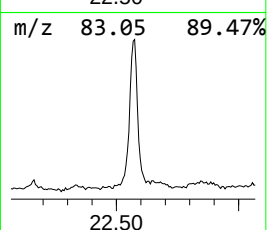
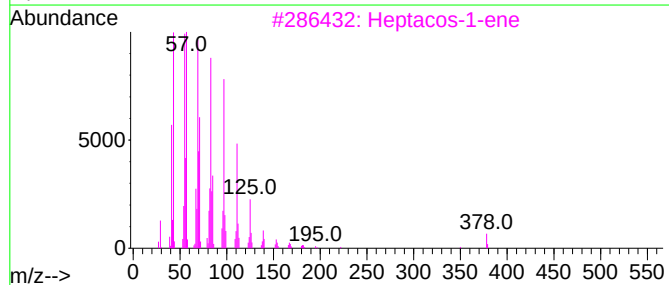
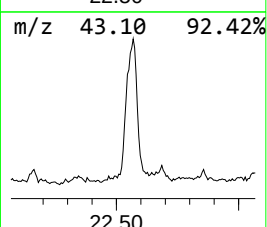
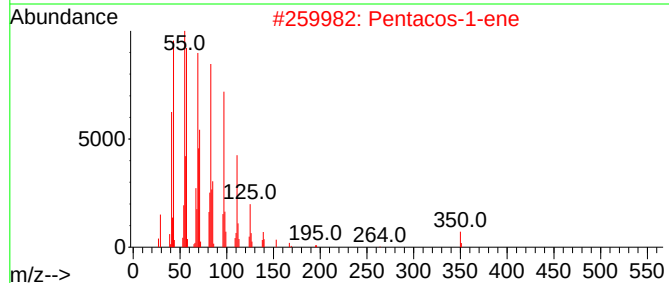
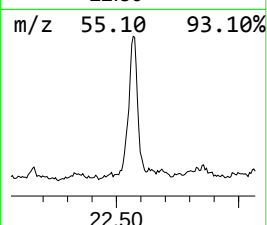
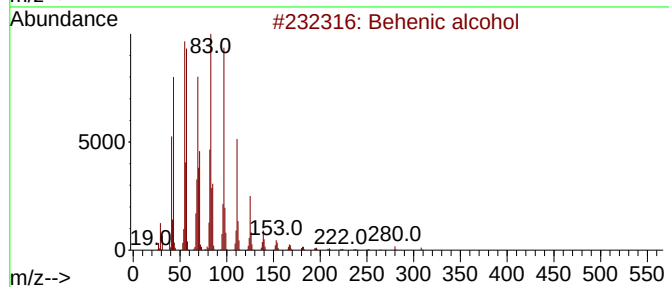
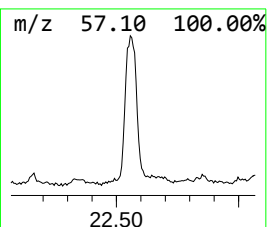
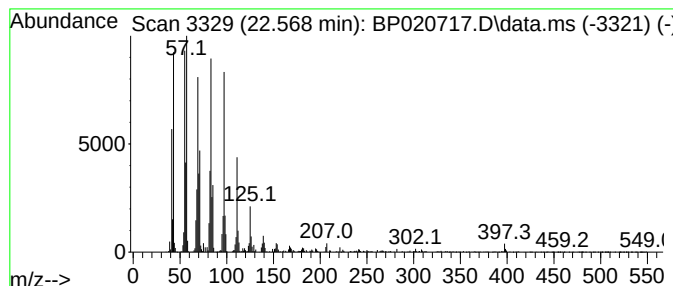
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Behenic alcohol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.568	13.17 ng/ul	1454390	Chrysene-d12	21.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Behenic alcohol	326	C22H46O	000661-19-8	95
2			Pentacos-1-ene	350	C25H50	016980-85-1	94
3			Heptacos-1-ene	378	C27H54	015306-27-1	94
4			1-Tetracosene	336	C24H48	010192-32-2	94
5			1-Hexacosene	364	C26H52	018835-33-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020717.D
Acq On : 01 Jul 2024 14:03
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Sample : P2871-05
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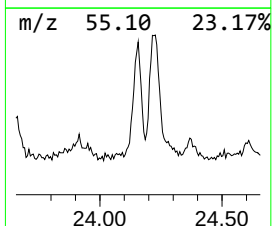
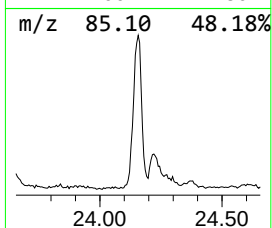
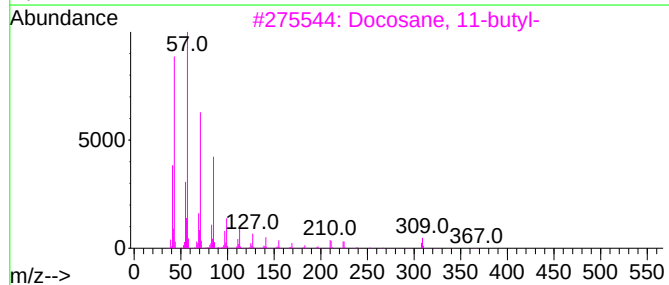
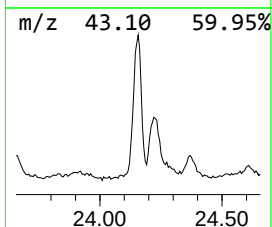
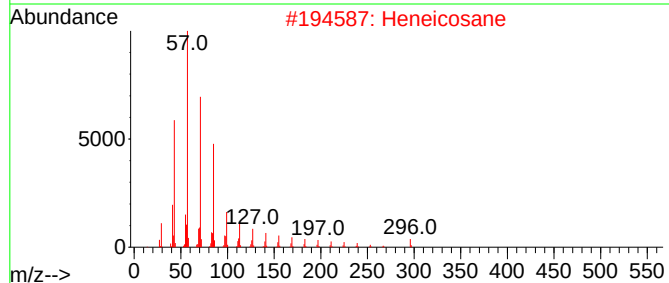
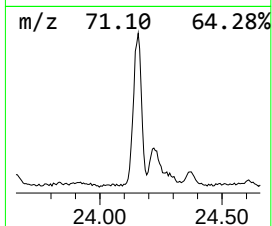
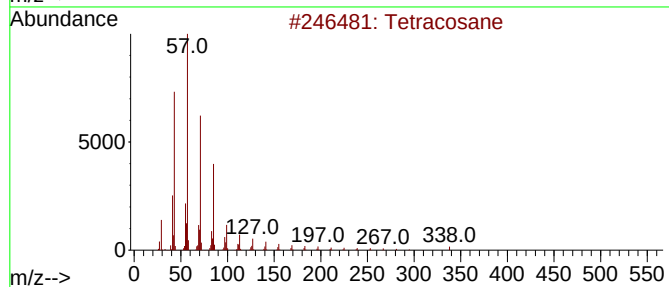
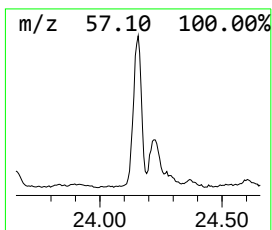
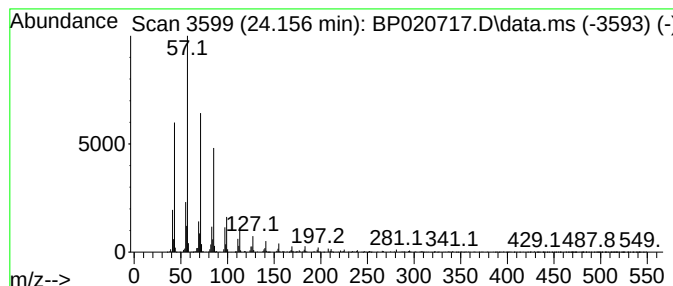
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.157	6.29 ng/ul	730488	Perylene-d12	25.009

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	97
2		Heneicosane	296	C21H44	000629-94-7	96
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	93
4		Hexacosane	366	C26H54	000630-01-3	93
5		Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020717.D
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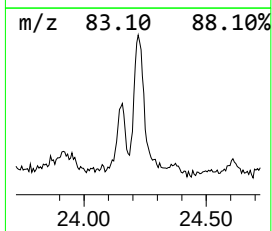
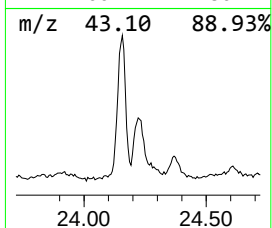
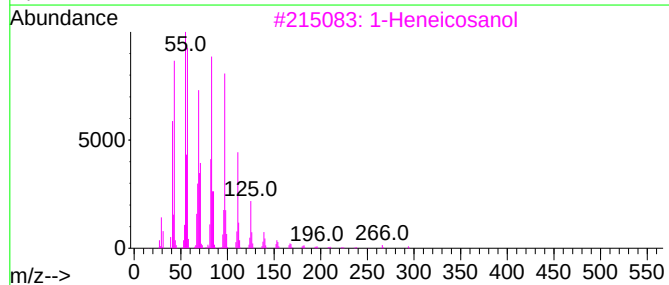
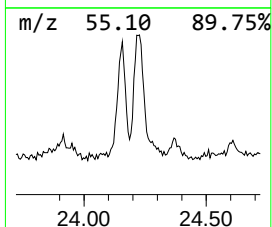
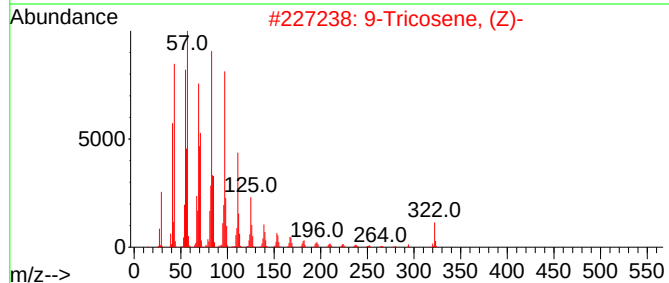
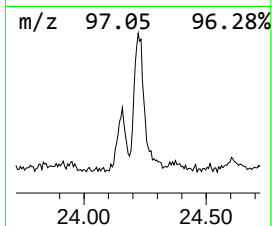
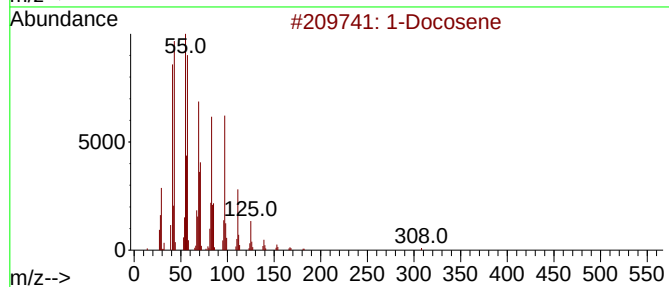
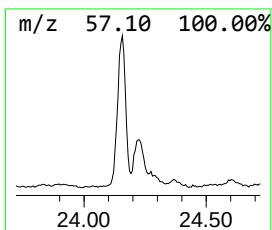
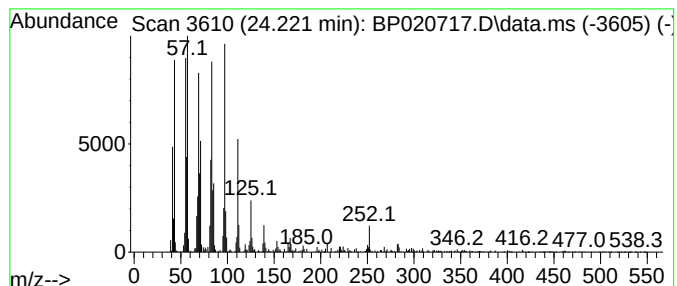
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.221	4.67 ng/ul	542583	Perylene-d12	25.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene			308	C22H44	001599-67-3	99
2	9-Tricosene, (Z)-			322	C23H46	027519-02-4	96
3	1-Heneicosanol			312	C21H44O	015594-90-8	95
4	Trifluoroacetoxy hexadecane			338	C18H33F3O2	006222-03-3	94
5	n-Tetracosanol-1			354	C24H50O	000506-51-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020717.D
Acq On : 01 Jul 2024 14:03
Operator : MA/JU
Sample : P2871-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4C53

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

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.016	33.8	ng/ul	2143050	1	7.740	1268930	20.0
Propylene Glycol	3.546	2.6	ng/ul	164013	1	7.740	1268930	20.0
1-Phenoxypropan...	11.257	3.6	ng/ul	324332	2	10.516	1800200	20.0
n-Hexadecanoic ...	18.127	7.7	ng/ul	874106	4	17.180	2269070	20.0
Octadecanoic acid	19.457	4.2	ng/ul	462462	5	21.645	2209430	20.0
(DEL) Alkane: S...	20.921	4.4	ng/ul	481320	5	21.645	2209430	20.0
(DEL) Alkane: S...	21.315	15.2	ng/ul	1676370	5	21.645	2209430	20.0
Behenic alcohol	22.568	13.2	ng/ul	1454390	5	21.645	2209430	20.0
(DEL) Alkane: S...	24.157	6.3	ng/ul	730488	6	25.009	2323630	20.0
(DEL) Alkane: S...	24.221	4.7	ng/ul	542583	6	25.009	2323630	20.0