

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
 Data File : BP020994.D
 Acq On : 16 Jul 2024 17:16
 Operator : MA/JU
 Sample : P3178-13
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4BT8

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-BP071024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020994.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.246	40	44	51	rVB	48508	63242	1.92%	0.117%
2	3.522	87	91	102	rBV	100368	154518	4.69%	0.286%
3	3.669	114	116	126	rBV	60565	98404	2.99%	0.182%
4	3.752	128	130	136	rVB	16857	20390	0.62%	0.038%
5	3.964	162	166	172	rVB	21229	28953	0.88%	0.054%
6	4.328	225	228	231	rVB5	7460	8701	0.26%	0.016%
7	4.369	231	235	242	rVV	13311	19553	0.59%	0.036%
8	4.458	246	250	254	rVV2	8154	11527	0.35%	0.021%
9	4.505	254	258	263	rVB	14683	21460	0.65%	0.040%
10	4.740	293	298	306	rBV9	7836	15528	0.47%	0.029%
11	4.852	312	317	327	rVB	65918	106923	3.24%	0.198%
12	5.822	474	482	492	rVB2	18109	43163	1.31%	0.080%
13	6.193	540	545	555	rBV7	7122	16597	0.50%	0.031%
14	6.587	605	612	615	rBV9	3128	7397	0.22%	0.014%
15	6.763	634	642	646	rBV2	12122	28679	0.87%	0.053%
16	6.805	646	649	655	rVB4	9674	15633	0.47%	0.029%
17	6.893	655	664	682	rBV	802433	1330393	40.36%	2.465%
18	7.040	682	689	701	rVB	641142	1036591	31.45%	1.920%
19	7.240	715	723	731	rBV	901906	1391630	42.22%	2.578%
20	7.316	731	736	749	rVB2	22912	63972	1.94%	0.119%
21	7.428	749	755	758	rBV	51911	82186	2.49%	0.152%
22	7.687	791	799	807	rBV	951055	1384349	42.00%	2.565%
23	7.775	807	814	827	rVB	78126	162902	4.94%	0.302%
24	7.957	839	845	853	rBV4	13137	31834	0.97%	0.059%
25	8.152	871	878	882	rBV9	4636	9384	0.28%	0.017%
26	8.204	882	887	898	rVB2	29136	55080	1.67%	0.102%
27	8.304	898	904	912	rVB2	4778	12742	0.39%	0.024%
28	8.404	912	921	929	rBV	894061	1558698	47.29%	2.888%
29	8.463	929	931	936	rVV5	27185	51273	1.56%	0.095%
30	8.510	936	939	946	rVB	17225	28503	0.86%	0.053%
31	8.840	986	995	1012	rBV	673116	1285422	39.00%	2.382%
32	8.981	1015	1019	1024	rVV8	6237	9755	0.30%	0.018%
33	9.193	1048	1055	1060	rBV3	19355	30337	0.92%	0.056%
34	9.393	1081	1089	1098	rBV	89308	145589	4.42%	0.270%
35	9.551	1106	1116	1127	rBV	651672	1083688	32.88%	2.008%
36	9.646	1129	1132	1138	rVV3	10490	16940	0.51%	0.031%

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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-BP071024.MA.M
 Title : SVOA CALIBRATION

37	9.716	1140	1144	1149	rBV7	4836	8264	0.25%	0.015%
38	9.775	1149	1154	1156	rBV4	6540	11071	0.34%	0.021%
39	9.804	1156	1159	1165	rVB7	11604	20105	0.61%	0.037%
40	10.093	1200	1208	1227	rBV	823395	1622753	49.23%	3.006%
41	10.298	1237	1243	1256	rBV	65461	174121	5.28%	0.323%
42	10.451	1262	1269	1282	rVB	1023361	1800558	54.63%	3.336%
43	10.604	1287	1295	1312	rBV	335162	789553	23.95%	1.463%
44	10.875	1336	1341	1349	rVB	29026	48352	1.47%	0.090%
45	10.987	1354	1360	1366	rBV2	44878	71745	2.18%	0.133%
46	11.198	1389	1396	1404	rBV	166332	324399	9.84%	0.601%
47	11.434	1432	1436	1440	rBV3	8915	17004	0.52%	0.032%
48	11.493	1441	1446	1451	rVV2	25408	56645	1.72%	0.105%
49	11.545	1451	1455	1463	rVB2	39576	72619	2.20%	0.135%
50	11.669	1473	1476	1483	rVB4	6025	10175	0.31%	0.019%
51	11.845	1501	1506	1512	rVB8	5109	9728	0.30%	0.018%
52	12.040	1534	1539	1545	rVV2	13721	24095	0.73%	0.045%
53	12.369	1591	1595	1604	rVV4	6991	15623	0.47%	0.029%
54	12.892	1681	1684	1692	rVB6	5904	12512	0.38%	0.023%
55	13.110	1715	1721	1725	rBV5	10697	20339	0.62%	0.038%
56	13.257	1740	1746	1753	rVB10	8456	18358	0.56%	0.034%
57	13.487	1780	1785	1790	rBV9	5551	12638	0.38%	0.023%
58	13.616	1802	1807	1814	rVB9	7649	14093	0.43%	0.026%
59	13.716	1814	1824	1834	rBV	1392951	2005958	60.86%	3.716%
60	13.822	1840	1842	1849	rVB8	8569	11837	0.36%	0.022%
61	13.963	1858	1866	1867	rBV6	15075	23737	0.72%	0.044%
62	14.004	1867	1873	1884	rVB	1829030	2430889	73.75%	4.504%
63	14.198	1902	1906	1910	rVB7	8927	11710	0.36%	0.022%
64	14.316	1919	1926	1932	rBV	1435877	2050181	62.20%	3.798%
65	14.375	1932	1936	1944	rVB3	18691	38549	1.17%	0.071%
66	14.539	1961	1964	1967	rVV4	24109	31257	0.95%	0.058%
67	14.581	1967	1971	1985	rVB	258932	601355	18.24%	1.114%
68	14.722	1991	1995	2001	rBV6	20091	40689	1.23%	0.075%
69	15.110	2058	2061	2065	rVB4	10409	13917	0.42%	0.026%
70	15.163	2065	2070	2073	rVB5	12466	18403	0.56%	0.034%
71	15.328	2092	2098	2104	rBV	2089110	2921672	88.64%	5.413%
72	15.469	2116	2122	2133	rBV	517217	797182	24.19%	1.477%
73	15.881	2188	2192	2204	rVB9	20544	40980	1.24%	0.076%
74	16.063	2217	2223	2231	rBV	72507	157735	4.79%	0.292%
75	16.516	2296	2300	2309	rBV6	43507	92030	2.79%	0.171%
76	16.875	2357	2361	2364	rBV5	22037	34020	1.03%	0.063%

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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-BP071024.MA.M
 Title : SVOA CALIBRATION

77	17.022	2383	2386	2392	rVB7	19549	34967	1.06%	0.065%
78	17.128	2397	2404	2409	rVV	1399273	2304453	69.91%	4.269%
79	17.175	2409	2412	2416	rVV	164530	230037	6.98%	0.426%
80	17.228	2416	2421	2431	rVV	2133878	3020426	91.64%	5.596%
81	17.304	2431	2434	2445	rVB	24905	61745	1.87%	0.114%
82	17.533	2469	2473	2476	rBV5	13672	25128	0.76%	0.047%
83	17.810	2516	2520	2524	rVB2	21214	27501	0.83%	0.051%
84	17.933	2538	2541	2547	rBV7	18166	33376	1.01%	0.062%
85	18.051	2555	2561	2571	rBV	995174	1444701	43.83%	2.677%
86	18.222	2587	2590	2595	rBV3	31835	49461	1.50%	0.092%
87	18.357	2610	2613	2617	rBV3	29468	37998	1.15%	0.070%
88	19.257	2758	2766	2773	rBV	490395	791358	24.01%	1.466%
89	19.386	2784	2788	2798	rBV	315356	425205	12.90%	0.788%
90	19.604	2819	2825	2840	rBV3	2003900	3296081	100.00%	6.107%
91	20.151	2915	2918	2919	rBV2	114348	120355	3.65%	0.223%
92	20.669	3003	3006	3008	rBV3	151660	191615	5.81%	0.355%
93	21.063	3070	3073	3079	rVB	321126	396421	12.03%	0.734%
94	21.216	3094	3099	3108	rVB	999399	1560000	47.33%	2.890%
95	21.557	3151	3157	3163	rBV2	1618973	2994867	90.86%	5.549%
96	22.404	3298	3301	3305	rBV	271858	353969	10.74%	0.656%
97	23.957	3558	3565	3570	rVB	1016273	2086047	63.29%	3.865%
98	24.627	3671	3679	3689	rBV2	842252	2420369	73.43%	4.484%
99	24.868	3712	3720	3730	rBV	926468	2775498	84.21%	5.142%
100	26.092	3921	3928	3944	rVB	710977	2384772	72.35%	4.418%

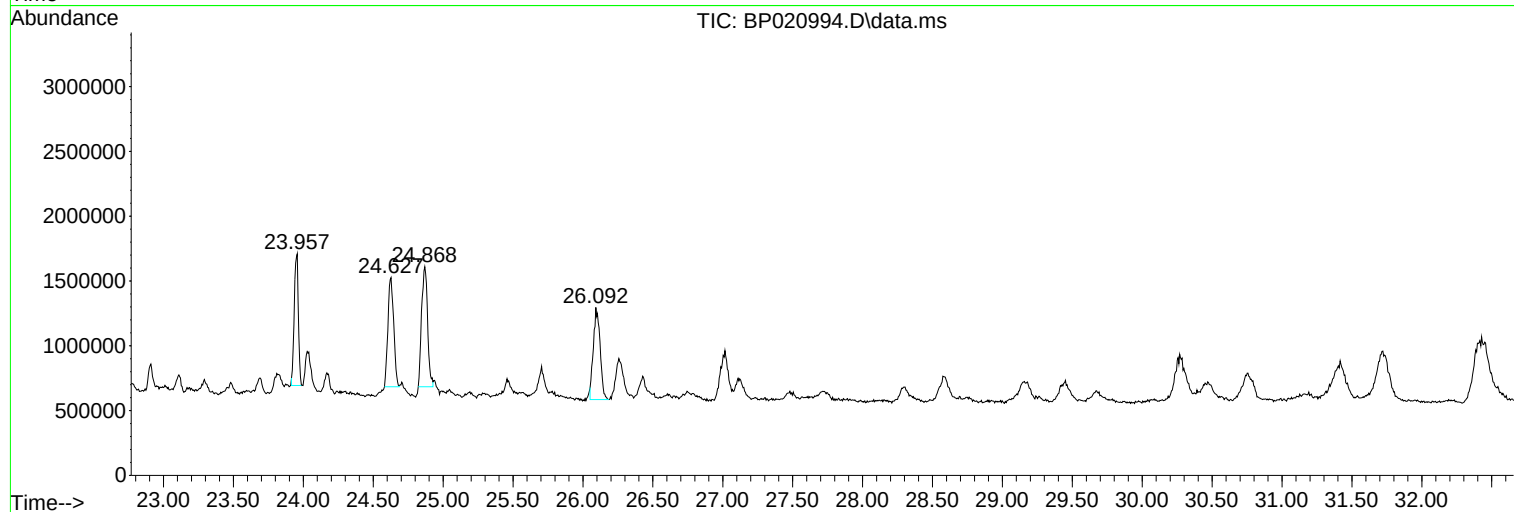
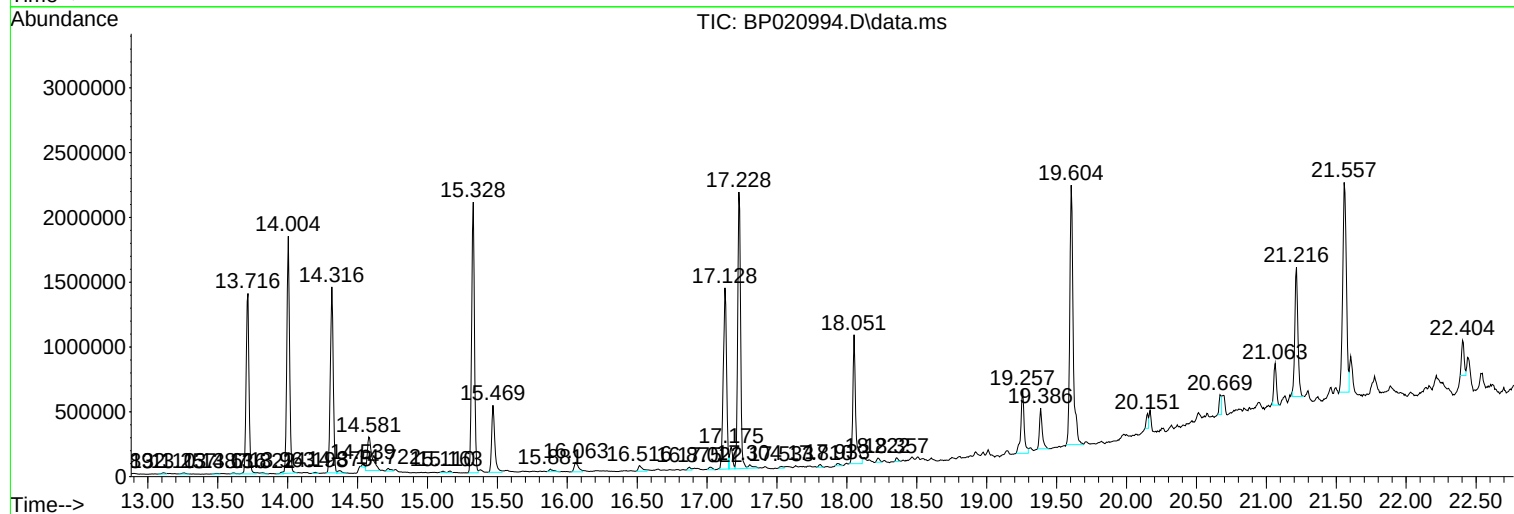
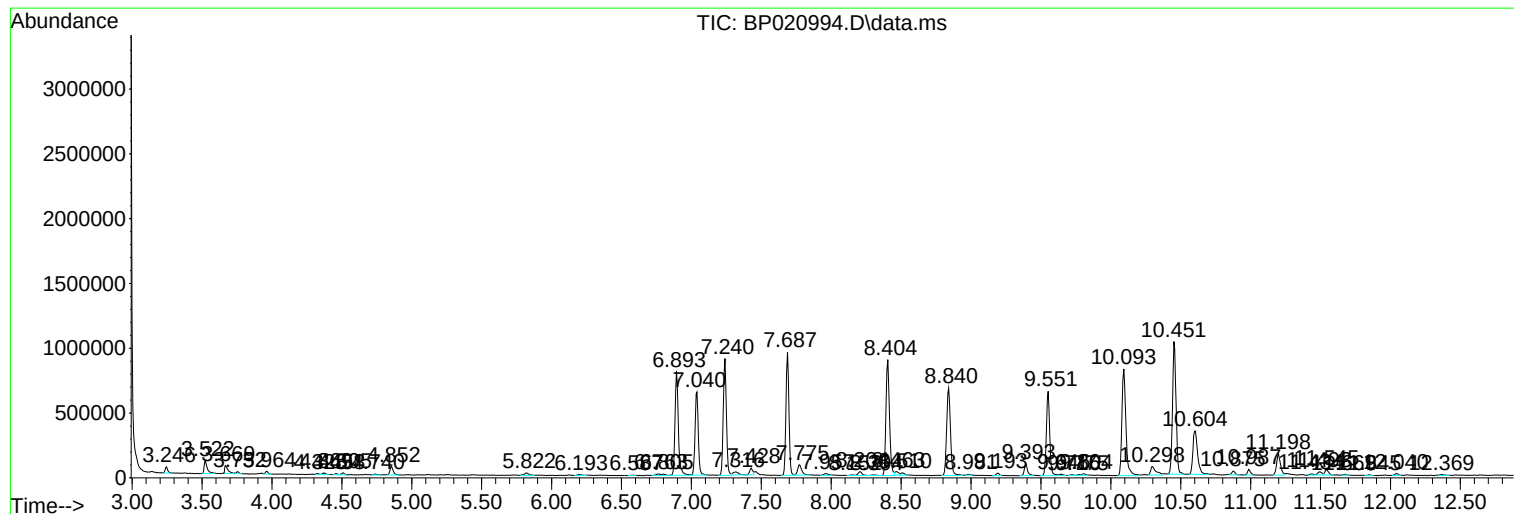
Sum of corrected areas: 53975137

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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



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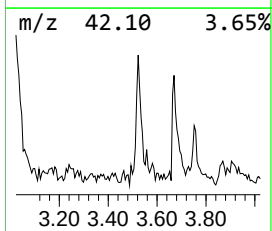
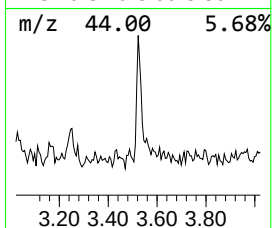
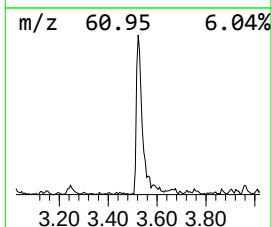
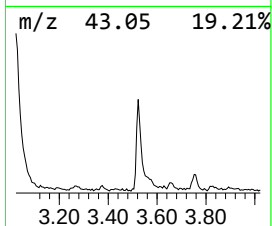
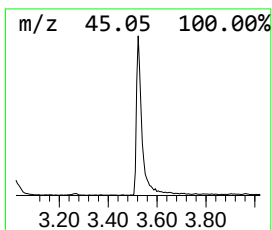
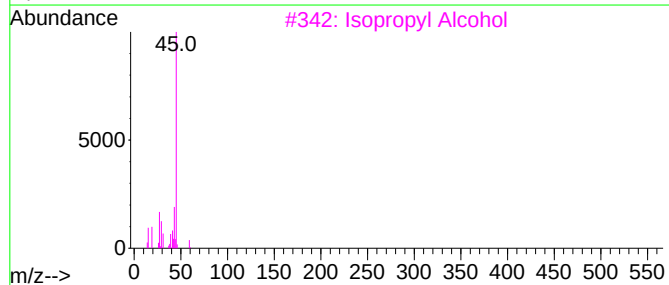
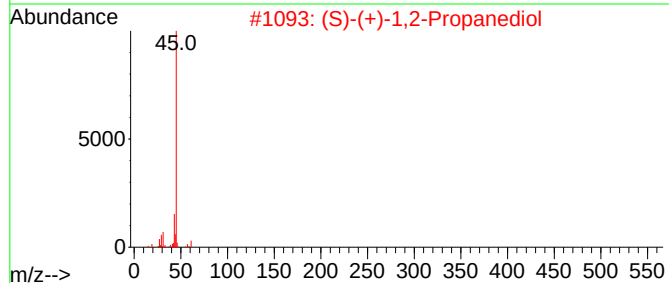
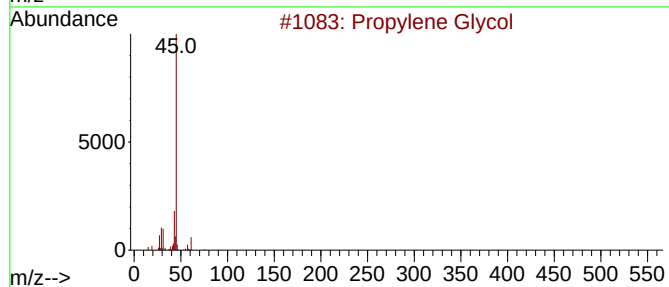
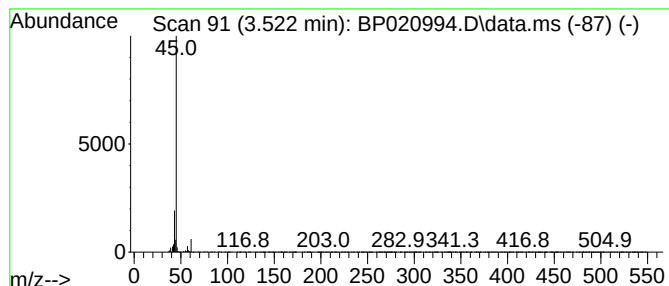
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.522	2.23 ng/ul	154518	1,4-Dichlorobenzene-d4	7.687

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	83
2			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	64
3			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
4			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	43
5			Isopropyl Alcohol	60	C3H8O	000067-63-0	9



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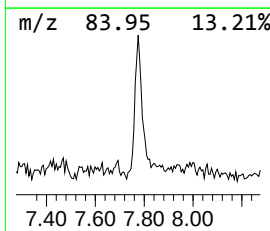
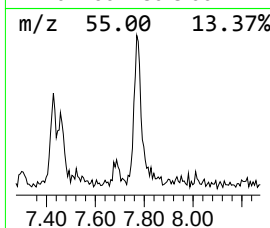
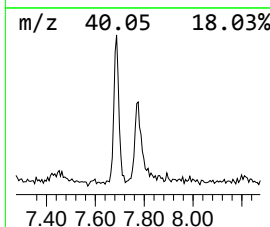
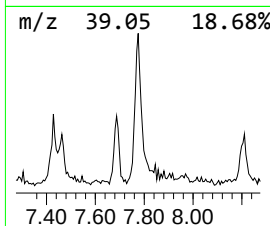
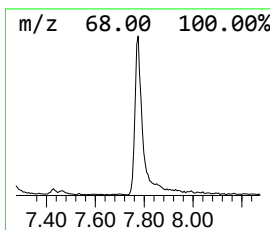
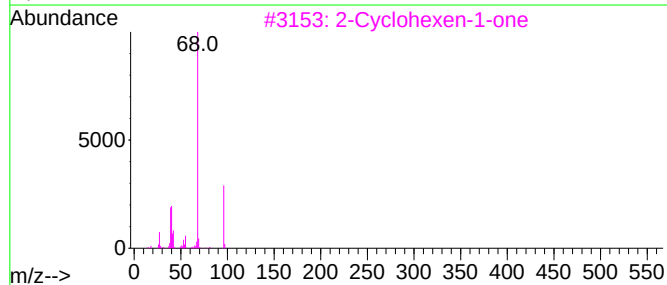
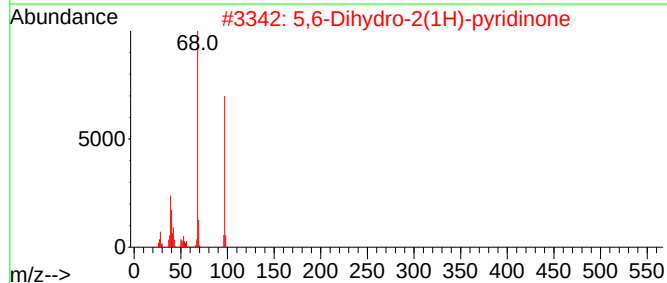
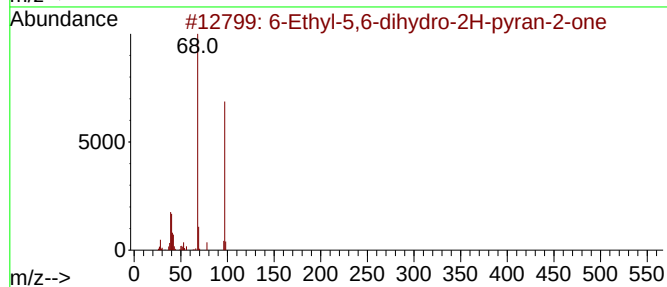
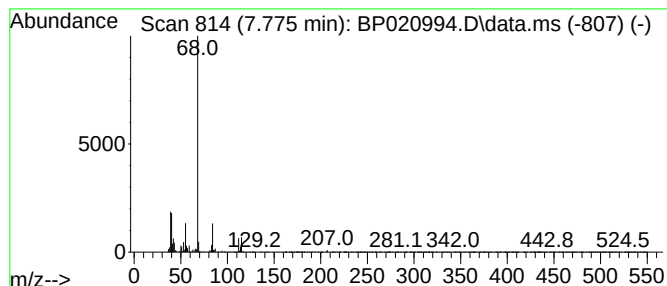
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 6-Ethyl-5,6-dihydro-2H-pyra... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.775	2.35 ng/ul	162902	1,4-Dichlorobenzene-d4	7.687

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Ethyl-5,6-dihydro-2H-pyran-2-one	126	C7H10O2	019895-35-3	50
2			5,6-Dihydro-2(1H)-pyridinone	97	C5H7NO	006052-73-9	40
3			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	36
4			Dihydro-3-methylene-5-methyl-2-f...	112	C6H8O2	062873-16-9	36
5			1H-Pyrazole	68	C3H4N2	000288-13-1	9



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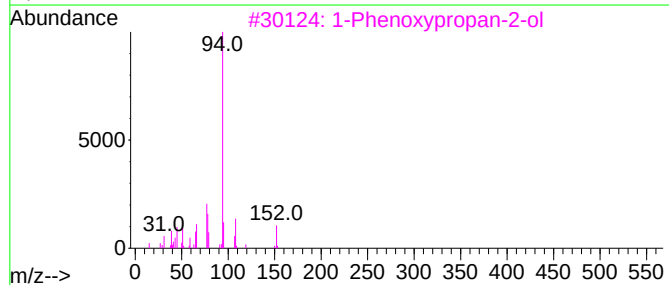
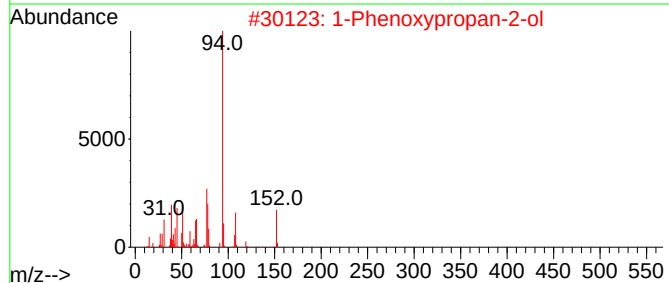
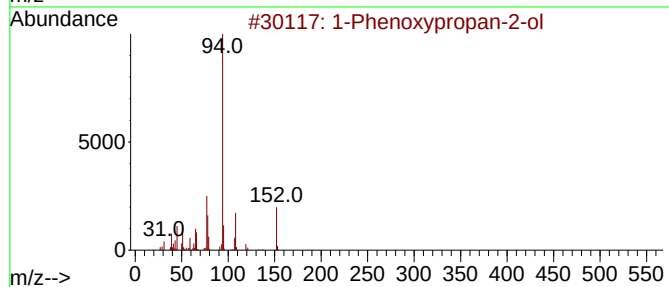
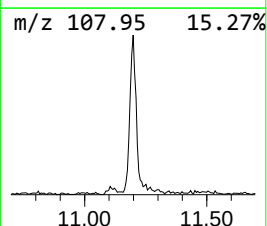
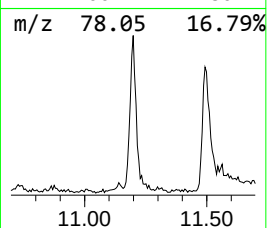
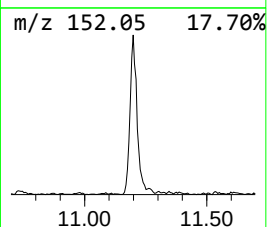
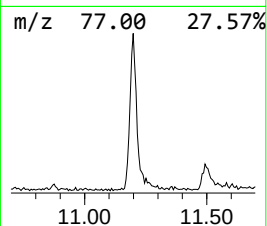
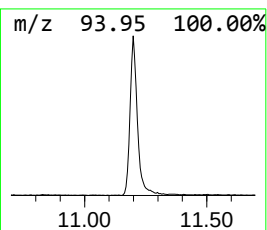
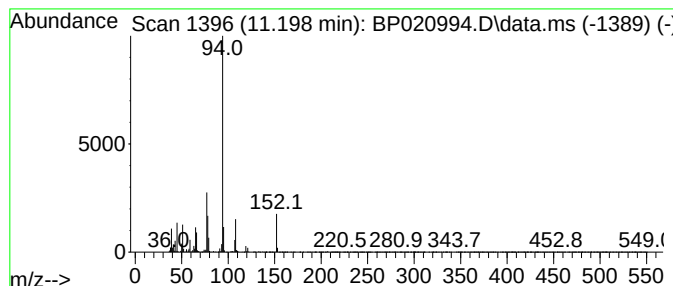
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.198	3.60 ng/ul	324399	Naphthalene-d8	10.451

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	96
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020994.D
Acq On : 16 Jul 2024 17:16
Operator : MA/JU
Sample : P3178-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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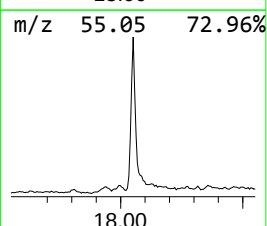
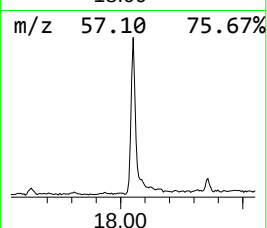
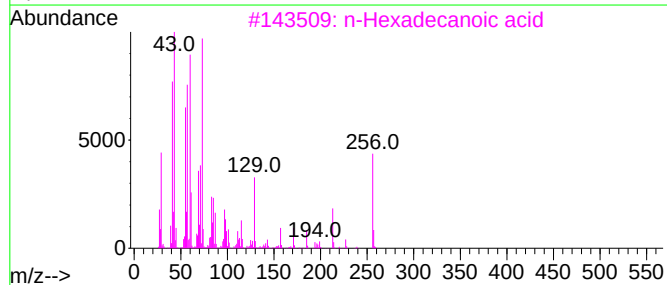
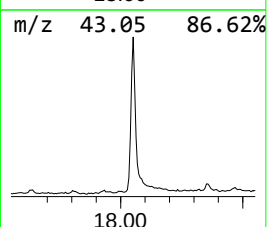
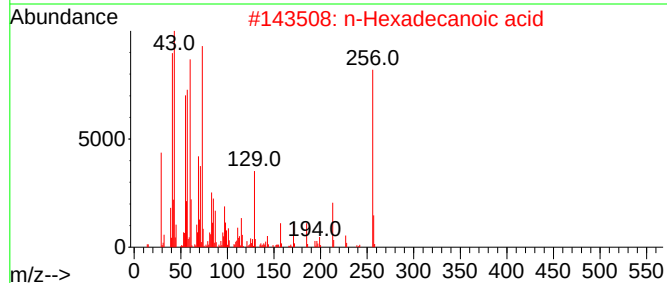
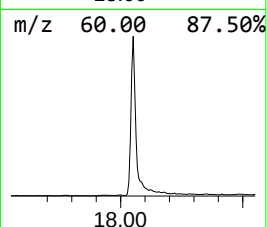
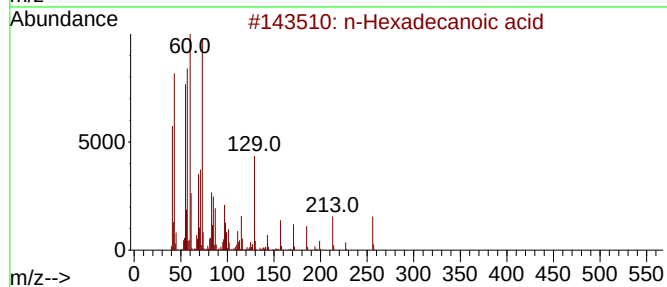
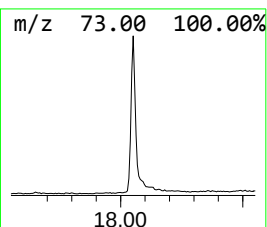
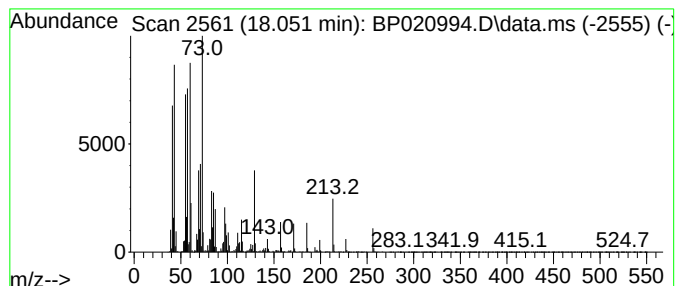
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.051	12.54 ng/ul	1444700	Phenanthrene-d10	17.128

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020994.D
Acq On : 16 Jul 2024 17:16
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Sample : P3178-13
Misc :
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Instrument :
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ClientSampleId :
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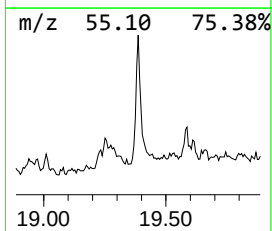
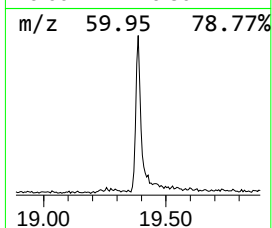
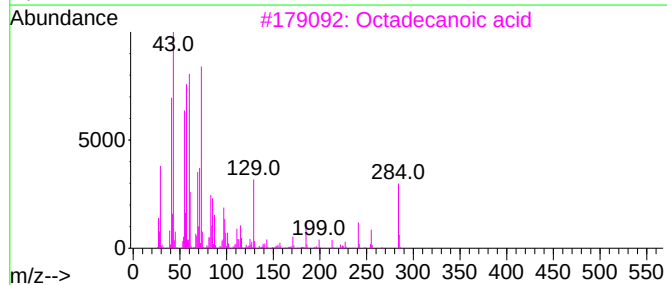
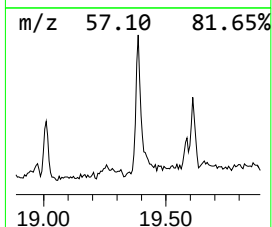
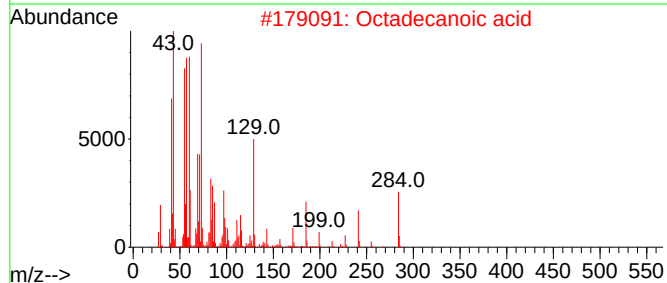
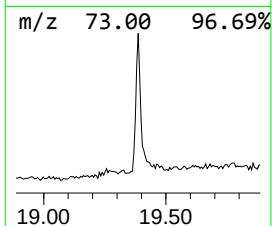
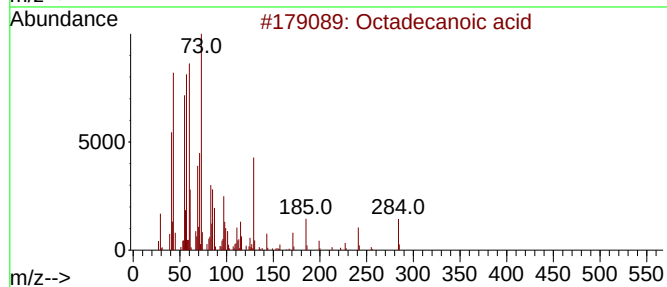
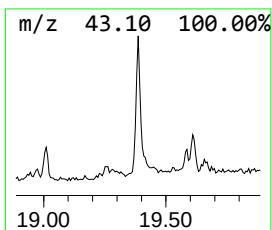
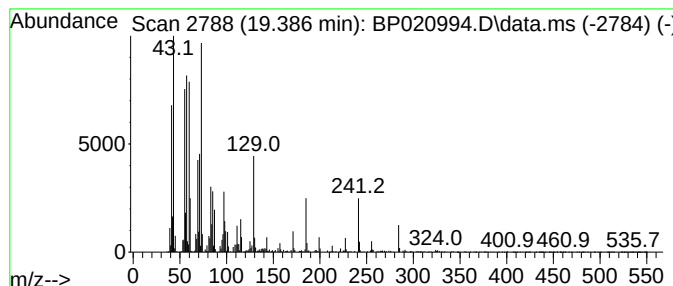
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.386	2.84 ng/ul	425205	Chrysene-d12	21.557

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	95
3			Octadecanoic acid	284	C18H36O2	000057-11-4	94
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020994.D
Acq On : 16 Jul 2024 17:16
Operator : MA/JU
Sample : P3178-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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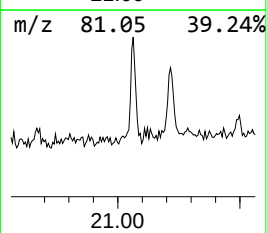
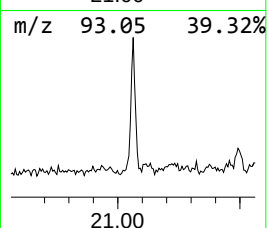
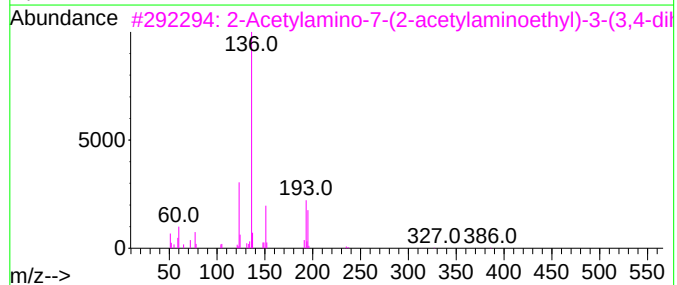
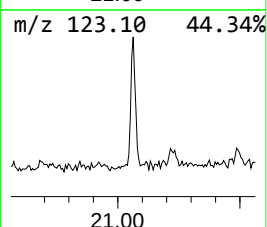
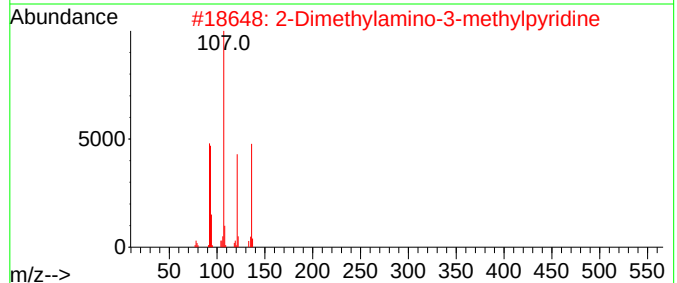
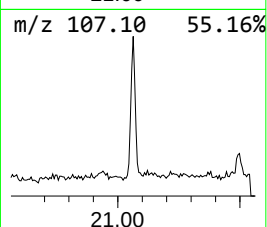
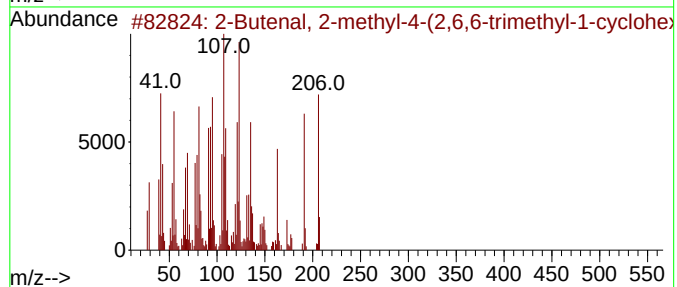
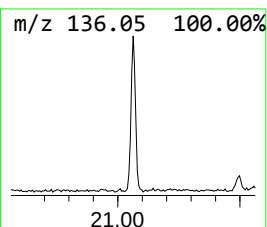
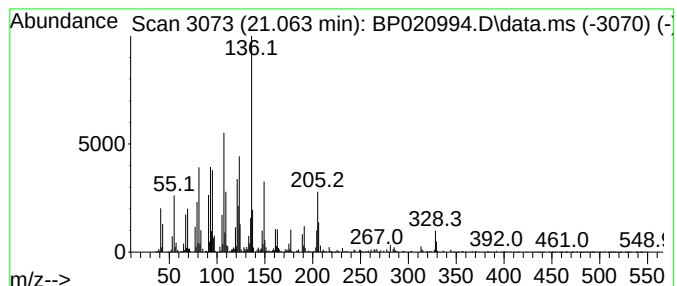
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2-Butenal, 2-methyl-4-(2,6,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.063	2.65 ng/ul	396421	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butenal, 2-methyl-4-(2,6,6-tri...	206	C14H22O	003155-71-3	55		
2	2-Dimethylamino-3-methylpyridine	136	C8H12N2	061713-46-0	38		
3	2-Acetylamino-7-(2-acetylaminoet...	386	C20H22N2O6	076734-99-1	38		
4	(3aS,8aS)-6,8a-Dimethyl-3-(propa...	204	C15H24	395070-76-5	25		
5	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	22		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020994.D
Acq On : 16 Jul 2024 17:16
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Sample : P3178-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

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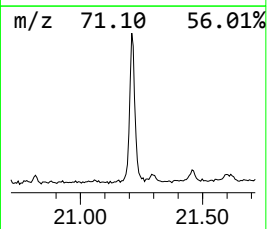
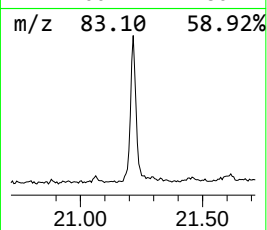
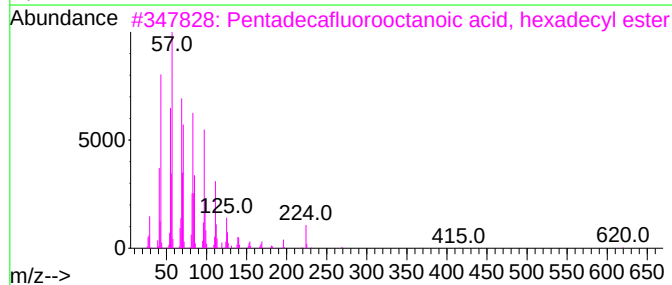
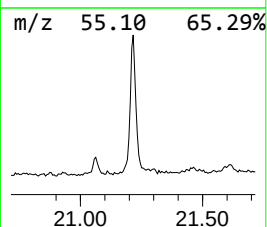
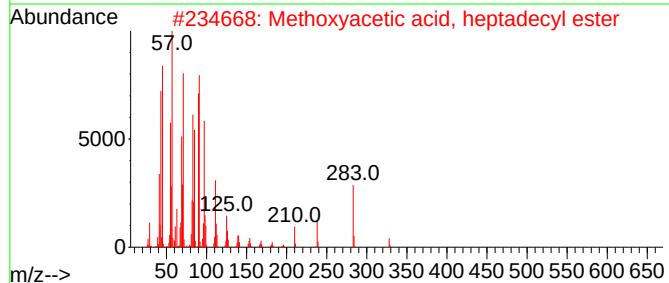
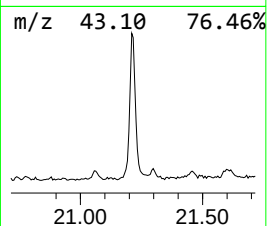
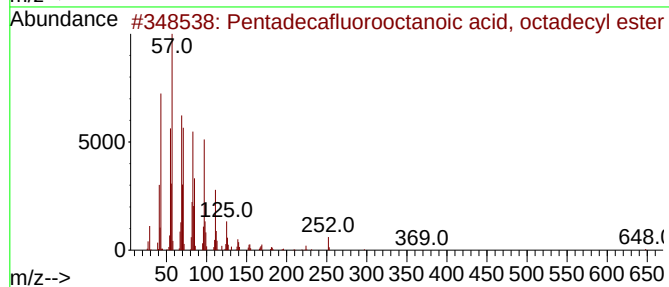
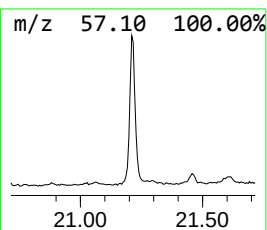
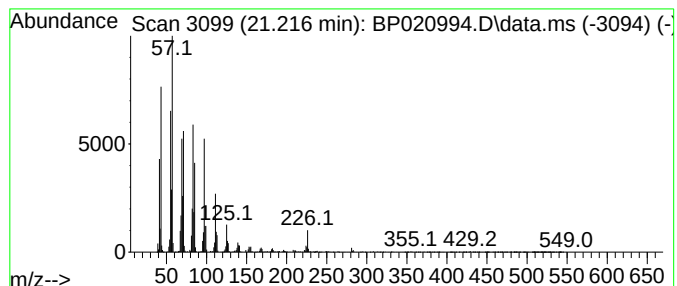
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Pentadecafluorooctanoic aci... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.216	10.42 ng/ul	1560000	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	95
2			Methoxyacetic acid, heptadecyl e...	328	C20H40O3	1000282-99-1	91
3			Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	87
4			Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	87
5			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020994.D
Acq On : 16 Jul 2024 17:16
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Sample : P3178-13
Misc :
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Instrument :
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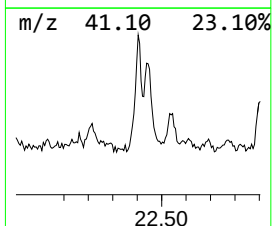
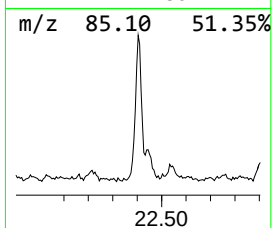
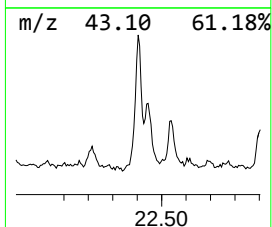
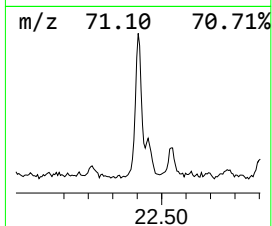
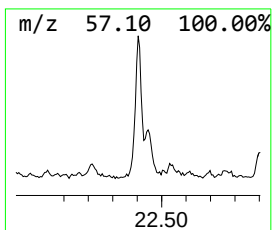
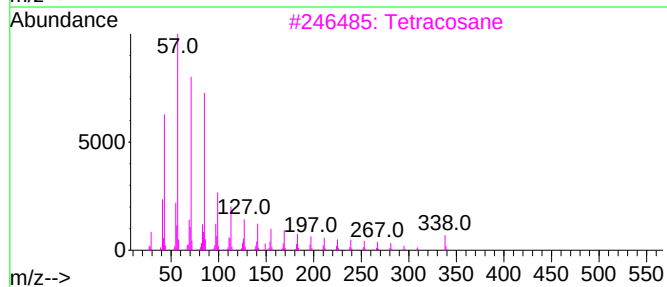
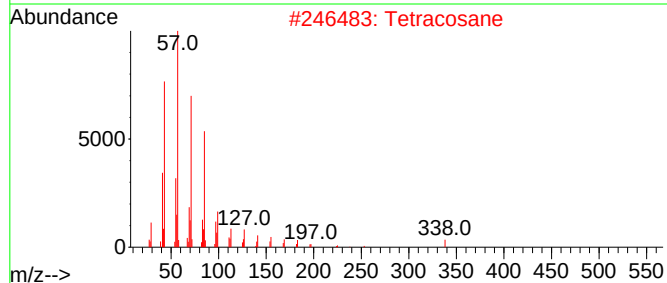
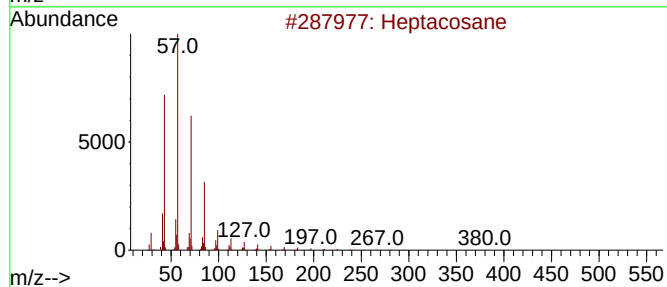
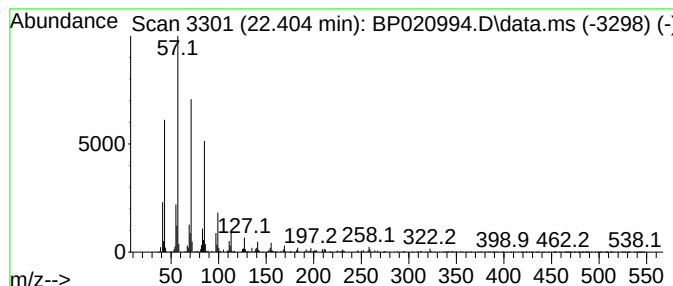
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.404	2.36 ng/ul	353969	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	95
2			Tetracosane	338	C24H50	000646-31-1	93
3			Tetracosane	338	C24H50	000646-31-1	90
4			2-Methylheptacosane	394	C28H58	001561-00-8	90
5			Tetracosane, 1-iodo-	464	C24H49I	1000406-32-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
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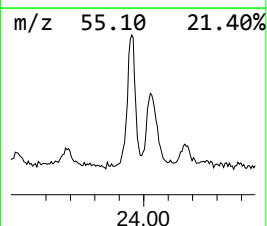
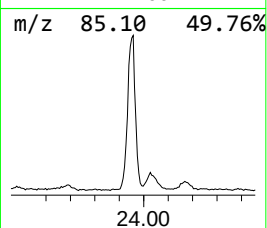
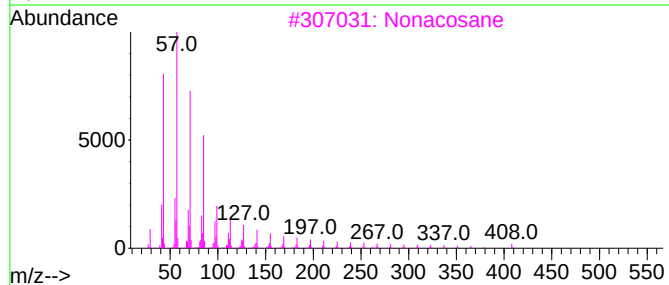
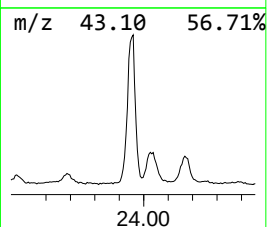
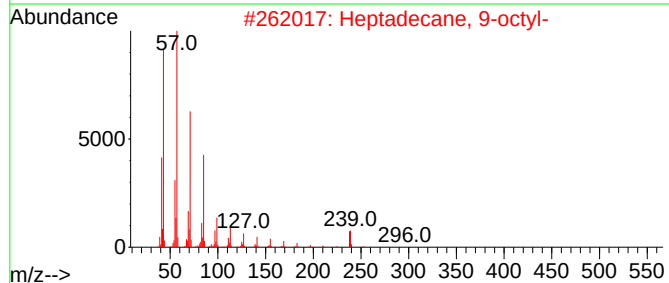
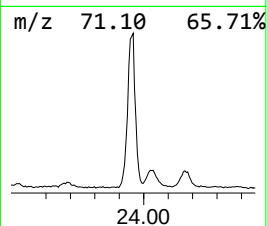
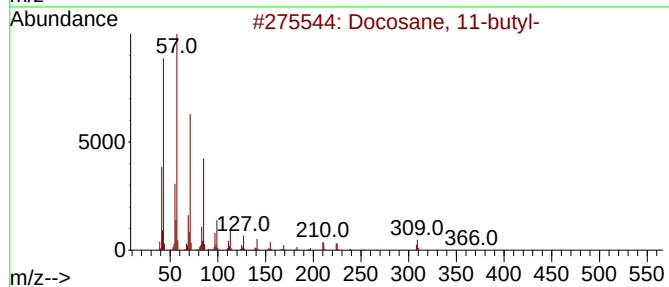
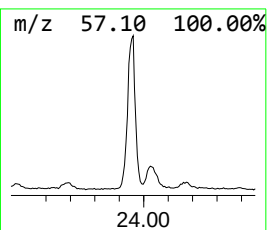
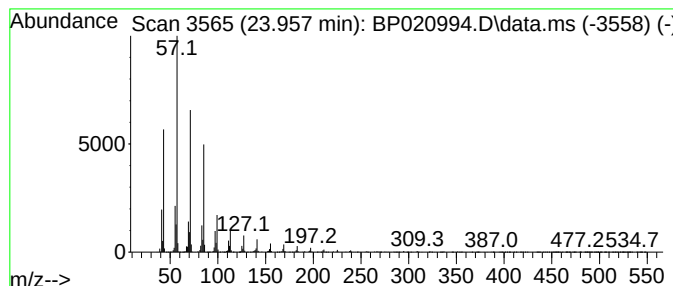
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.957	15.03 ng/ul	2086050	Perylene-d12	24.868	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-	366	C26H54	013475-76-8	95
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
3	Nonacosane	408	C29H60	000630-03-5	93
4	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93
5	2-Methylpentacosane	366	C26H54	000629-87-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020994.D
Acq On : 16 Jul 2024 17:16
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Misc :
ALS Vial : 5 Sample Multiplier: 1

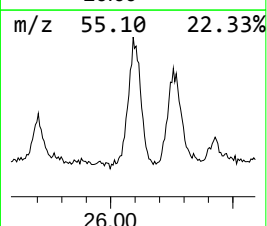
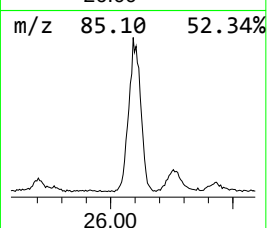
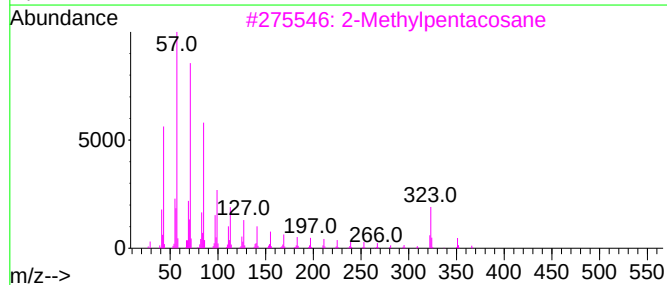
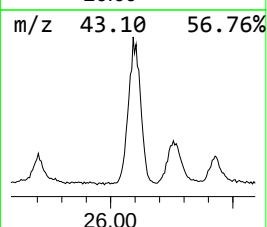
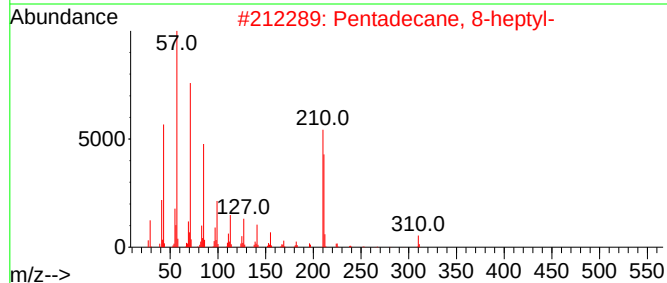
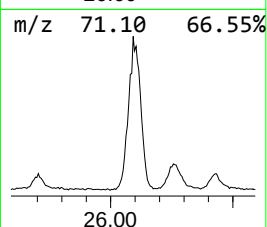
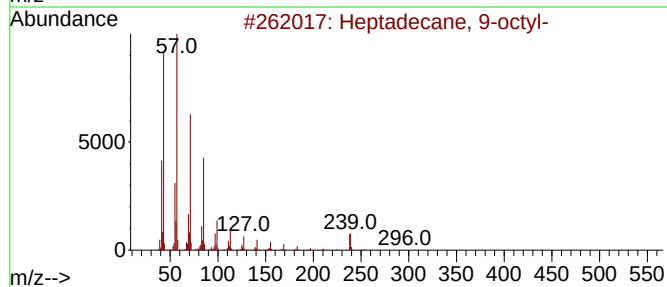
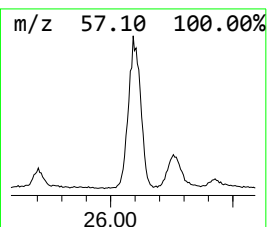
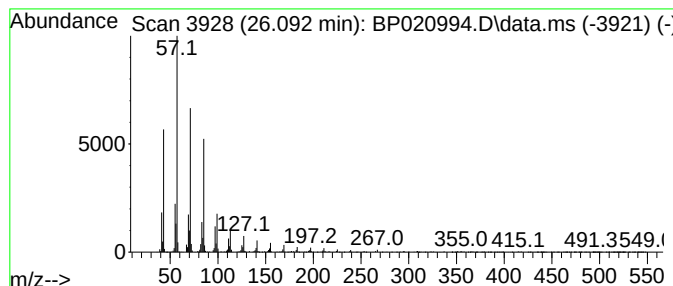
Instrument :
BNA_P
ClientSampleId :
A4BT8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.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 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
26.092	17.18 ng/ul	2384770	Perylene-d12	24.868	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	94
3	2-Methylpentacosane	366	C26H54	000629-87-8	94
4	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
5	Tricosane	324	C23H48	000638-67-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020994.D
Acq On : 16 Jul 2024 17:16
Operator : MA/JU
Sample : P3178-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BT8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.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
Propylene Glycol	3.522	2.2	ng/ul	154518	1	7.687	1384350	20.0
6-Ethyl-5,6-dih...	7.775	2.4	ng/ul	162902	1	7.687	1384350	20.0
1-Phenoxypropan...	11.198	3.6	ng/ul	324399	2	10.451	1800560	20.0
n-Hexadecanoic ...	18.051	12.5	ng/ul	1444700	4	17.128	2304450	20.0
Octadecanoic acid	19.386	2.8	ng/ul	425205	5	21.557	2994870	20.0
2-Butenal, 2-me...	21.063	2.6	ng/ul	396421	5	21.557	2994870	20.0
Pentadecafluoro...	21.216	10.4	ng/ul	1560000	5	21.557	2994870	20.0
(DEL) Alkane: S...	22.404	2.4	ng/ul	353969	5	21.557	2994870	20.0
(DEL) Alkane: S...	23.957	15.0	ng/ul	2086050	6	24.868	2775500	20.0
(DEL) Alkane: S...	26.092	17.2	ng/ul	2384770	6	24.868	2775500	20.0