

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
 Data File : BN029730.D
 Acq On : 13 Feb 2024 12:27
 Operator : MA/JU
 Sample : P1305-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0FP3

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_N\Methods\SFAM-EPA-BN020724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN029730.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.311	34	38	46	rVB	85723	116025	1.12%	0.097%
2	3.599	84	87	93	rBV	16514	24956	0.24%	0.021%
3	3.722	105	108	124	rBV	503518	840841	8.11%	0.701%
4	4.046	158	163	168	rBV	26616	39108	0.38%	0.033%
5	4.905	274	309	311	rBV3	235545	1459428	14.08%	1.216%
6	6.846	636	639	645	rVB6	6642	11622	0.11%	0.010%
7	7.028	662	670	683	rBV	1310351	2130674	20.56%	1.775%
8	7.199	693	699	709	rBV	1051096	1698162	16.39%	1.415%
9	7.387	724	731	742	rBV	1291645	1996803	19.27%	1.664%
10	7.475	742	746	756	rVB	30625	62247	0.60%	0.052%
11	7.857	803	811	818	rBV	1047828	1632425	15.75%	1.360%
12	7.922	818	822	828	rVB	18464	27954	0.27%	0.023%
13	8.569	925	932	941	rBV	1531118	2442438	23.57%	2.035%
14	9.028	1002	1010	1017	rBV	1221118	1987956	19.18%	1.656%
15	9.422	1070	1077	1082	rVB3	11004	19661	0.19%	0.016%
16	9.745	1123	1132	1142	rBV	935470	1532061	14.78%	1.276%
17	9.928	1156	1163	1164	rBV6	15266	28087	0.27%	0.023%
18	10.281	1212	1223	1234	rBV	1424111	2323844	22.42%	1.936%
19	10.657	1279	1287	1296	rBV	1358579	2293649	22.13%	1.911%
20	10.804	1303	1312	1330	rBV	1216346	2111538	20.38%	1.759%
21	12.087	1526	1530	1535	rVB4	7345	13287	0.13%	0.011%
22	12.251	1552	1558	1565	rVB	28324	50220	0.48%	0.042%
23	12.569	1605	1612	1618	rVB6	12441	26584	0.26%	0.022%
24	12.692	1624	1633	1641	rBV	85625	202678	1.96%	0.169%
25	13.181	1712	1716	1721	rVB6	8080	12144	0.12%	0.010%
26	13.328	1730	1741	1750	rBV	2400107	4248769	41.00%	3.540%
27	13.410	1750	1755	1761	rVV	89493	148672	1.43%	0.124%
28	13.492	1761	1769	1782	rVV2	2807302	5167411	49.86%	4.305%
29	13.675	1792	1800	1801	rBV7	6249	13977	0.13%	0.012%
30	13.733	1806	1810	1815	rVV6	12535	24264	0.23%	0.020%
31	13.933	1836	1844	1853	rBV	4346609	6677073	64.43%	5.563%
32	14.063	1863	1866	1872	rVB7	8862	13464	0.13%	0.011%
33	14.204	1874	1890	1903	rBV	2686266	3956296	38.18%	3.296%
34	14.398	1916	1923	1935	rBV	965886	1665422	16.07%	1.388%
35	14.510	1935	1942	1953	rVB	1824351	2697078	26.03%	2.247%
36	14.598	1953	1957	1967	rVB10	18646	47162	0.46%	0.039%

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.MA.M
 Title : SVOA CALIBRATION

37	14.716	1970	1977	1992	rBV	899630	1472091	14.21%	1.227%
38	14.869	1999	2003	2010	rVV9	7992	15136	0.15%	0.013%
39	14.933	2010	2014	2021	rVB9	15348	32015	0.31%	0.027%
40	15.116	2040	2045	2050	rBV5	7454	15557	0.15%	0.013%
41	15.210	2053	2061	2075	rBV	3755611	6006333	57.96%	5.004%
42	15.328	2075	2081	2082	rBV6	7727	14871	0.14%	0.012%
43	15.404	2086	2094	2105	rVV	1054017	1605547	15.49%	1.338%
44	15.510	2105	2112	2119	rVB	3083314	4613096	44.52%	3.843%
45	15.627	2125	2132	2148	rBV	933913	1374503	13.26%	1.145%
46	15.751	2148	2153	2156	rVV3	16175	22516	0.22%	0.019%
47	16.127	2213	2217	2219	rBV4	13002	18762	0.18%	0.016%
48	16.298	2242	2246	2253	rBV8	12833	25819	0.25%	0.022%
49	16.616	2297	2300	2306	rBV5	21753	32391	0.31%	0.027%
50	16.675	2306	2310	2317	rVB4	23360	33222	0.32%	0.028%
51	16.827	2332	2336	2340	rVB7	7844	11065	0.11%	0.009%
52	16.945	2349	2356	2360	rBV8	8685	17633	0.17%	0.015%
53	16.992	2361	2364	2368	rVB6	11750	13925	0.13%	0.012%
54	17.045	2368	2373	2376	rBV6	10487	19273	0.19%	0.016%
55	17.169	2386	2394	2404	rBV	2360405	3543211	34.19%	2.952%
56	17.274	2404	2412	2421	rVV2	5035924	8342484	80.50%	6.951%
57	17.363	2421	2427	2434	rVV	3046246	4391170	42.37%	3.659%
58	17.422	2434	2437	2440	rVB4	16121	19494	0.19%	0.016%
59	17.945	2518	2526	2532	rVB5	28866	72052	0.70%	0.060%
60	18.033	2537	2541	2542	rBV3	15289	18432	0.18%	0.015%
61	18.180	2557	2566	2583	rBV	1314636	2760393	26.64%	2.300%
62	18.469	2611	2615	2624	rVB10	22822	44515	0.43%	0.037%
63	18.604	2634	2638	2644	rBV4	26675	47676	0.46%	0.040%
64	18.845	2675	2679	2685	rBV6	29837	42027	0.41%	0.035%
65	19.045	2709	2713	2715	rBV5	13361	21617	0.21%	0.018%
66	19.068	2715	2717	2722	rVV6	16298	23897	0.23%	0.020%
67	19.110	2722	2724	2727	rVV4	8779	10365	0.10%	0.009%
68	19.204	2735	2740	2743	rBV4	45143	76569	0.74%	0.064%
69	19.280	2748	2753	2755	rBV3	42496	68516	0.66%	0.057%
70	19.339	2761	2763	2778	rVB2	67871	135761	1.31%	0.113%
71	19.463	2779	2784	2798	rBV	548530	882728	8.52%	0.735%
72	19.651	2806	2816	2828	rBV2	4654340	10362962	100.00%	8.634%
73	19.933	2860	2864	2867	rBV4	16770	20900	0.20%	0.017%
74	20.204	2906	2910	2913	rBV5	43994	65551	0.63%	0.055%
75	20.239	2913	2916	2921	rVB	64176	75911	0.73%	0.063%
76	20.563	2967	2971	2974	rBV5	26143	39534	0.38%	0.033%

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 Title : SVOA CALIBRATION

77	20.615	2977	2980	2983	rVB3	41898	46365	0.45%	0.039%
78	20.739	2996	3001	3005	rVB2	82289	105182	1.01%	0.088%
79	21.204	3075	3080	3097	rBV	1469817	2084634	20.12%	1.737%
80	21.468	3120	3125	3143	rBV2	3929563	5467738	52.76%	4.556%
81	21.598	3144	3147	3153	rVV2	132617	169050	1.63%	0.141%
82	21.657	3153	3157	3167	rVV	145705	205842	1.99%	0.172%
83	22.127	3232	3237	3244	rBV2	158729	327459	3.16%	0.273%
84	22.615	3317	3320	3325	rVB2	100152	140137	1.35%	0.117%
85	22.751	3339	3343	3349	rVB	168626	219983	2.12%	0.183%
86	23.174	3411	3415	3419	rBV	147330	213587	2.06%	0.178%
87	23.698	3495	3504	3516	rBV	4532229	9260199	89.36%	7.715%
88	23.851	3518	3530	3540	rVB	2838255	5880164	56.74%	4.899%
89	24.527	3641	3645	3651	rVB2	128794	256266	2.47%	0.214%
90	24.609	3653	3659	3672	rVB2	306151	751061	7.25%	0.626%
91	27.039	4064	4072	4084	rVB3	238804	742159	7.16%	0.618%

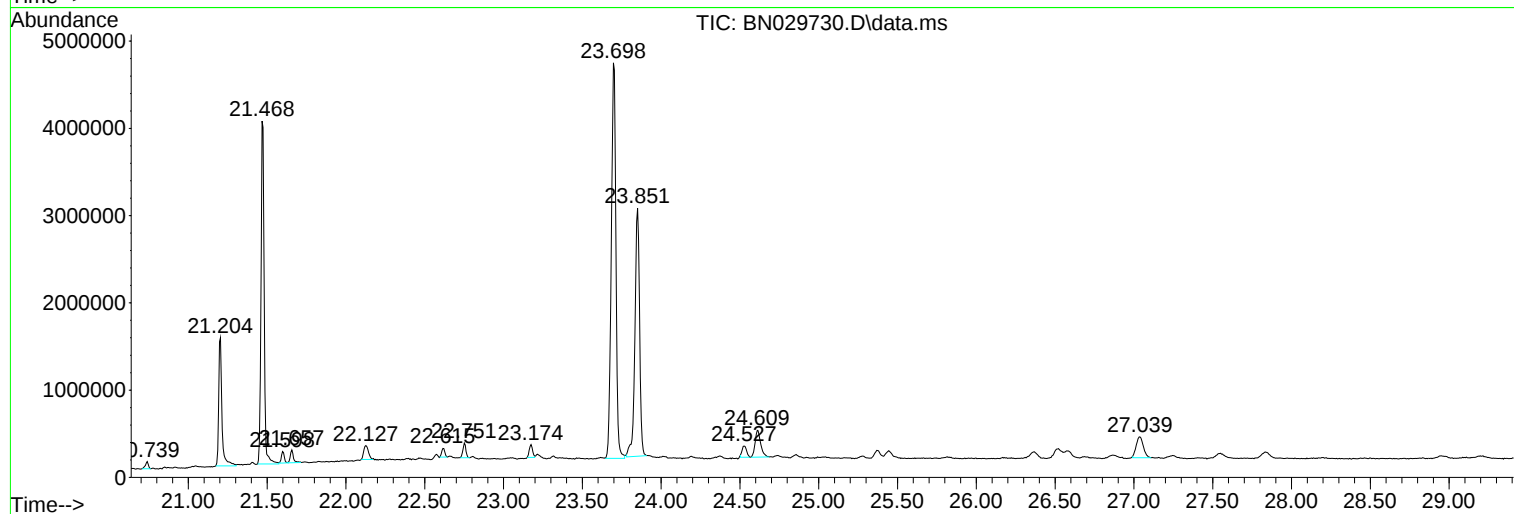
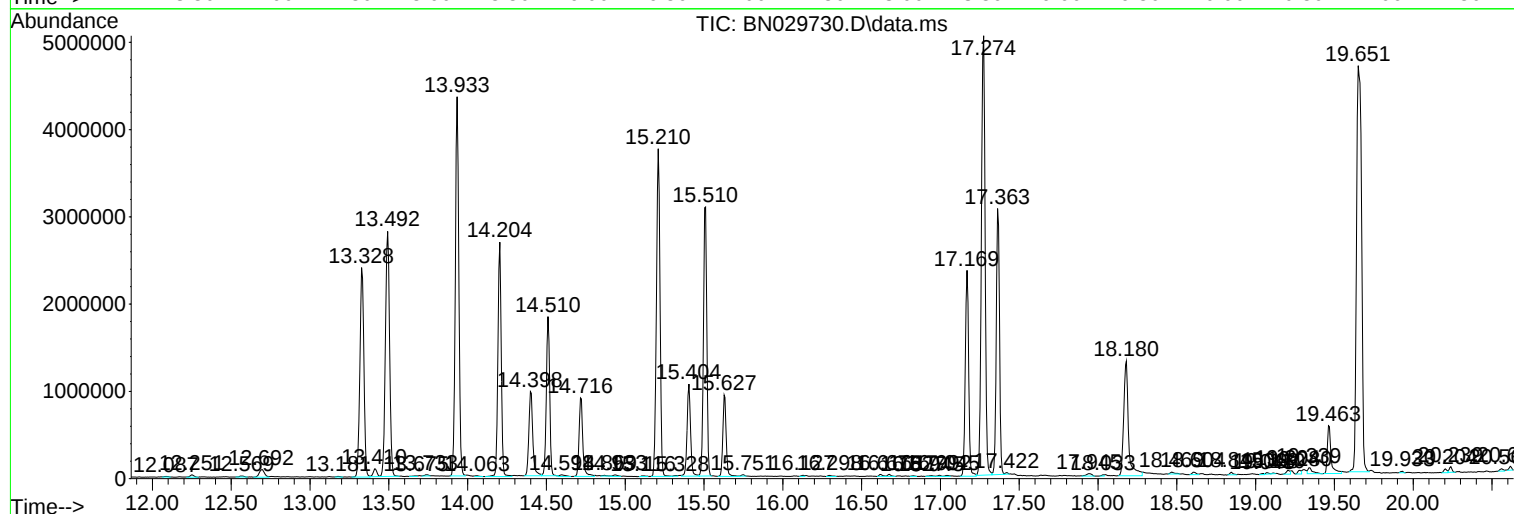
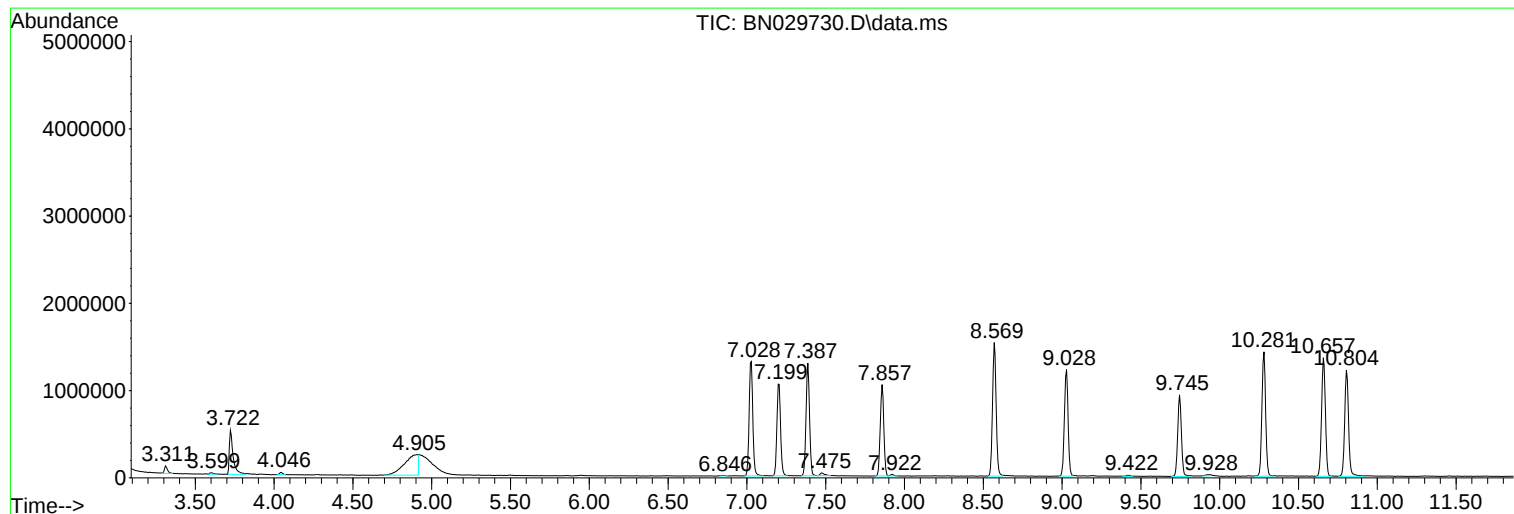
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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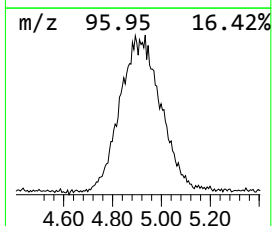
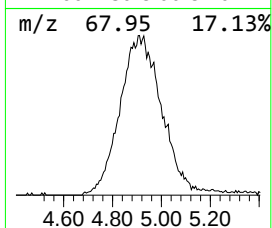
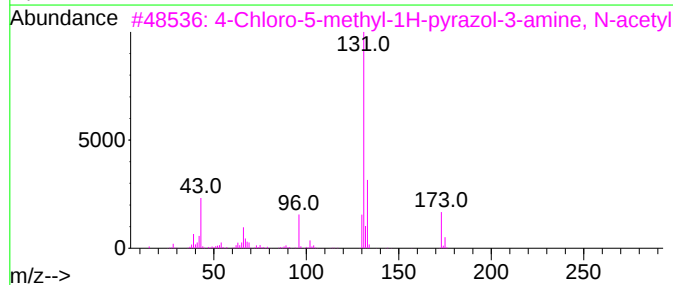
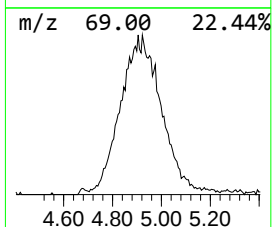
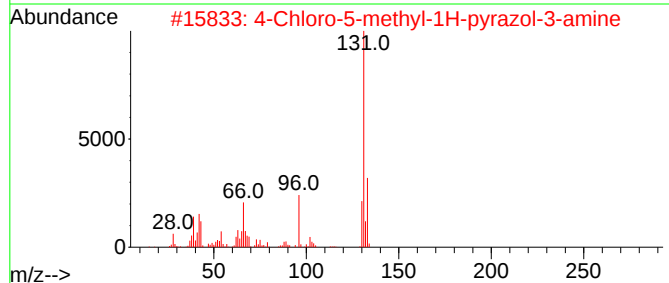
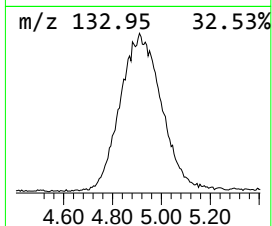
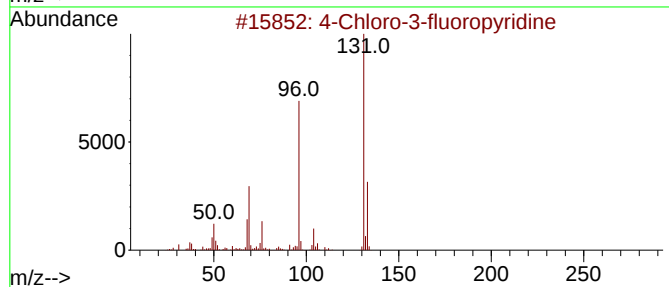
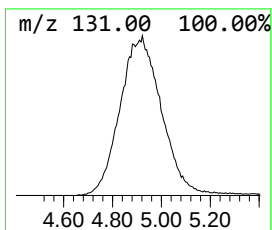
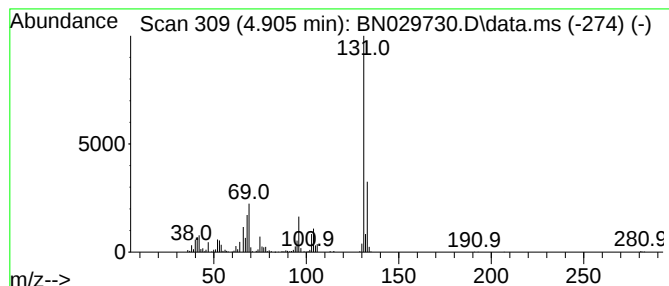
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 4-Chloro-3-fluoropyridine Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.905	17.88 ng/ul	1459430	1,4-Dichlorobenzene-d4	7.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Chloro-3-fluoropyridine	131	C5H3ClFN	002546-56-7	59
2			4-Chloro-5-methyl-1H-pyrazol-3-a...	131	C4H6ClN3	110580-44-4	50
3			4-Chloro-5-methyl-1H-pyrazol-3-a...	173	C6H8ClN3O	1000511-56-8	50
4			4-Chloro-5-methyl-1H-pyrazol-3-a...	215	C8H10ClN3O2	1010511-78-3	42
5			1H-Benzimidazole, 2-(4-methoxyp...	254	C15H14N2O2	1000303-90-9	38



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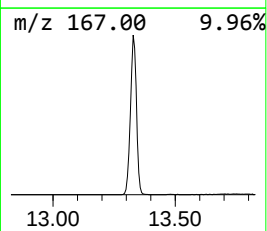
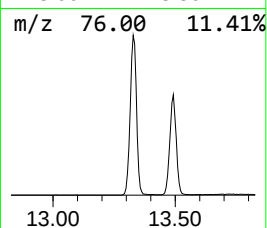
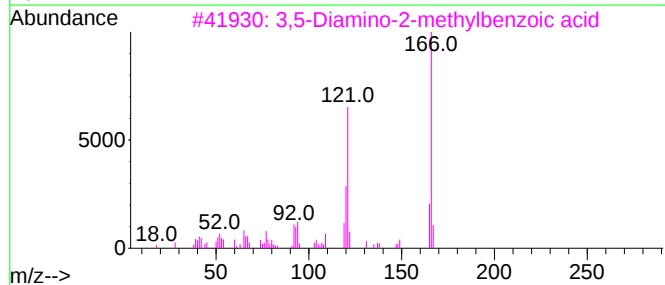
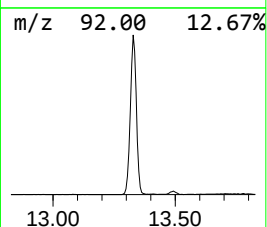
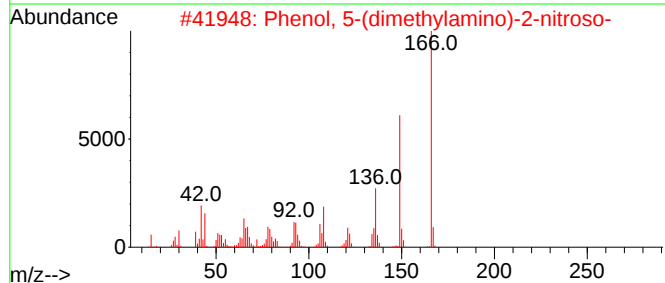
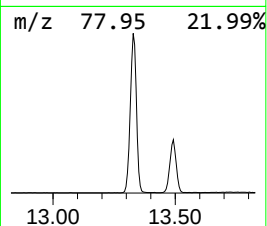
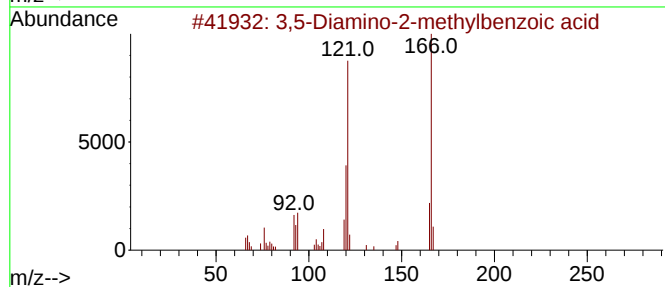
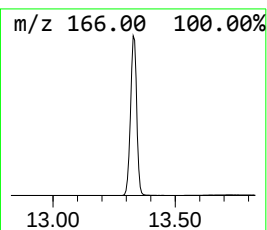
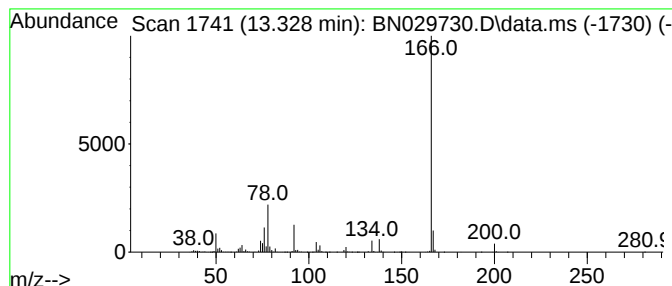
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 3,5-Diamino-2-methylbenzoic... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.328	31.51 ng/ul	4248770	Acenaphthene-d10	14.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-Diamino-2-methylbenzoic acid	166	C8H10N2O2	090007-30-0	58
2			Phenol, 5-(dimethylamino)-2-nitr...	166	C8H10N2O2	016761-04-9	33
3			3,5-Diamino-2-methylbenzoic acid	166	C8H10N2O2	090007-30-0	33
4			Benzeneacetic acid, 3-methoxy-	166	C9H10O3	001798-09-0	9
5			Benzo[b]tetrahydrofuran-3-one, 5...	166	C8H6O4	014771-00-7	9



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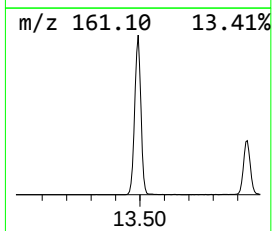
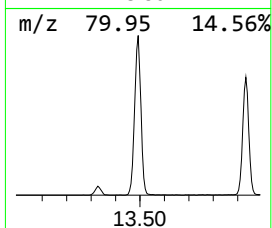
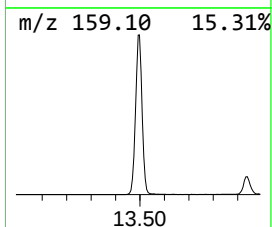
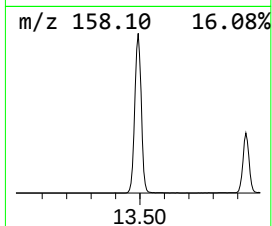
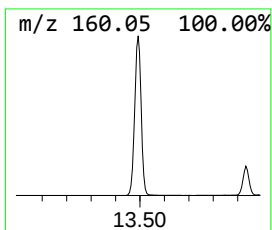
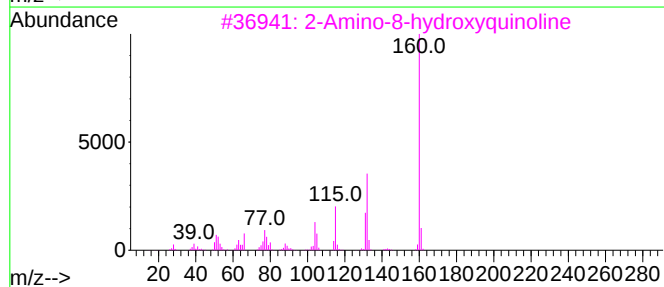
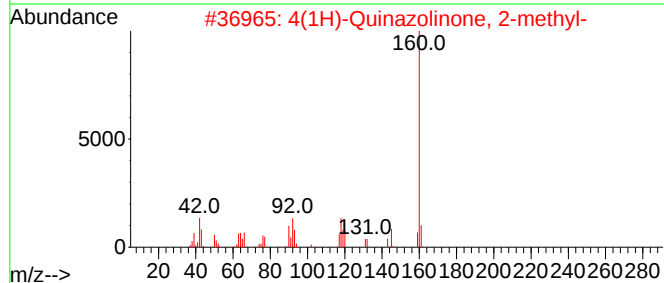
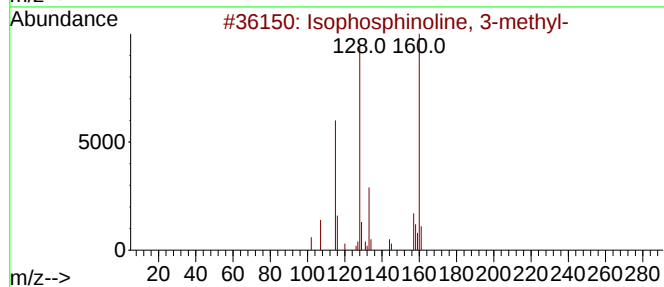
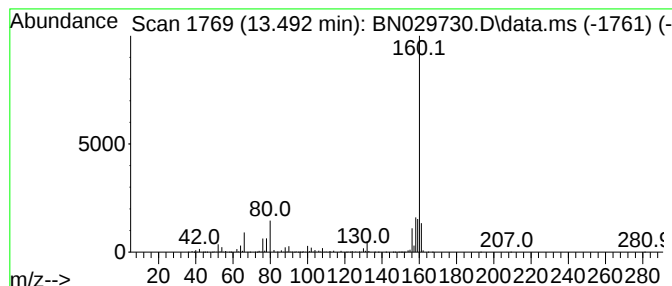
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Isophosphinoline, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.492	38.32 ng/ul	5167410	Acenaphthene-d10	14.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isophosphinoline, 3-methyl-	160	C10H9P	049622-63-1	80
2			4(1H)-Quinazolinone, 2-methyl-	160	C9H8N2O	001769-24-0	59
3			2-Amino-8-hydroxyquinoline	160	C9H8N2O	070125-16-5	58
4			1-Benzosuberone	160	C11H12O	000826-73-3	53
5			Benzonitrile, 4-isothiocyanato-	160	C8H4N2S	002719-32-6	50



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Data File : BN029730.D
Acq On : 13 Feb 2024 12:27
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Sample : P1305-02
Misc :
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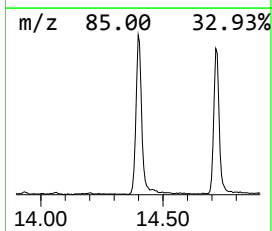
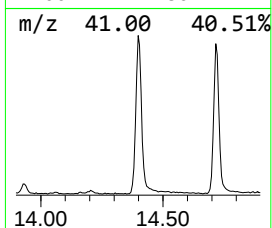
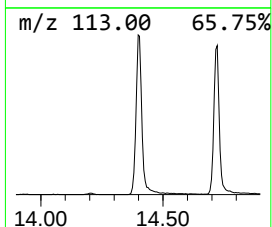
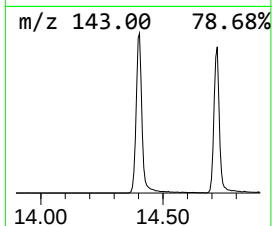
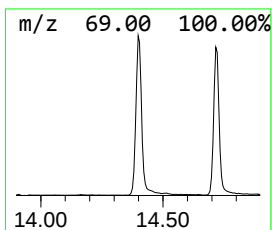
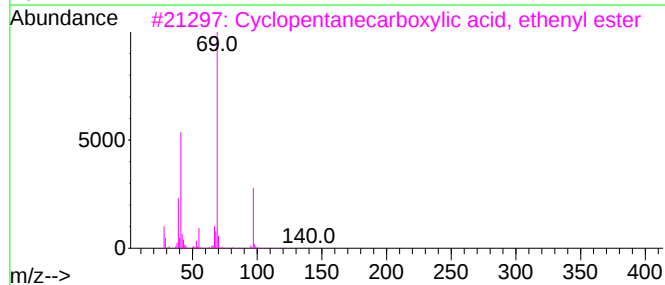
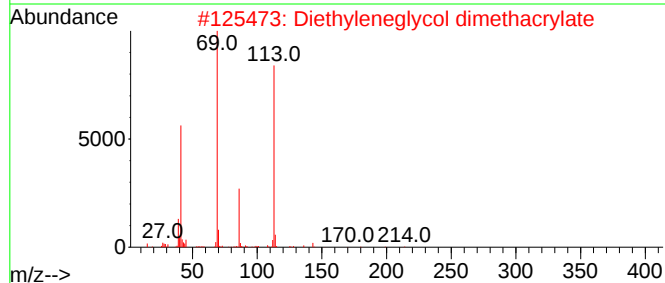
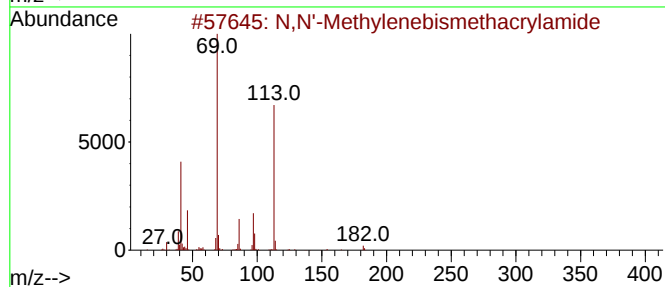
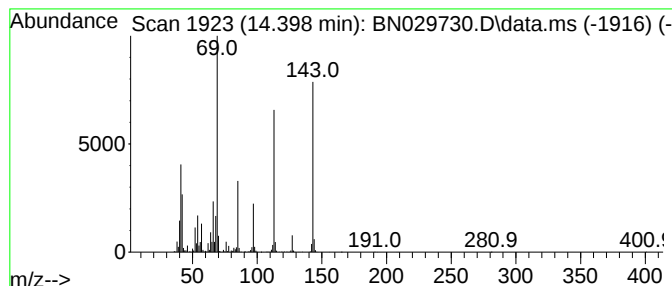
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.398	12.35 ng/ul	1665420	Acenaphthene-d10	14.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N,N'-Methylenebismethacrylamide	182	C9H14N2O2	002359-15-1	22
2			Diethyleneglycol dimethacrylate	242	C12H18O5	002358-84-1	22
3			Cyclopentanecarboxylic acid, eth...	140	C8H12O2	016523-06-1	14
4			Glutaric acid, ethyl 4-heptyl ester	258	C14H26O4	1000359-45-2	10
5			2-(Hydroxymethyl)-3-methoxy-2H-f...	144	C6H8O4	1332747-95-1	10



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Data File : BN029730.D
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Instrument :
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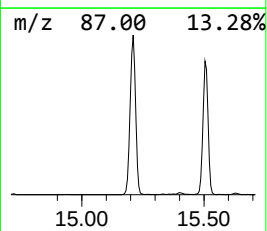
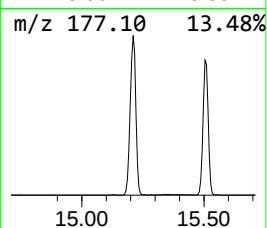
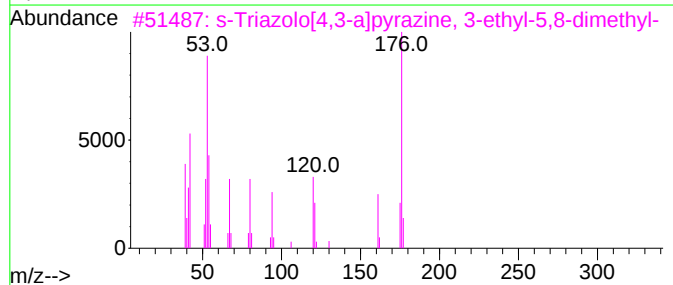
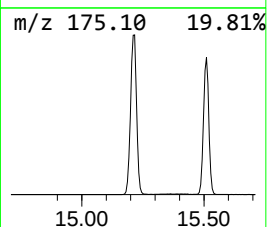
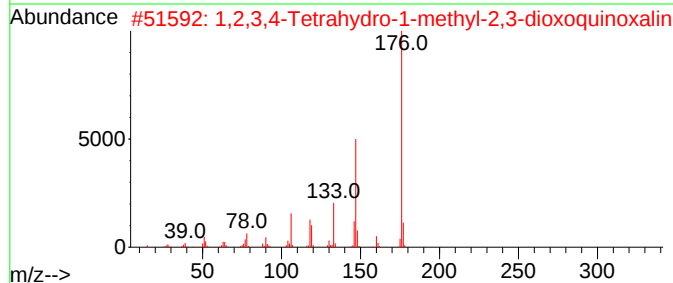
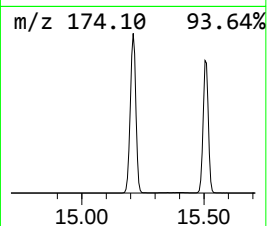
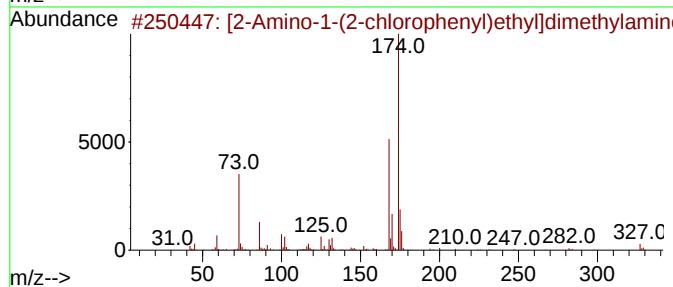
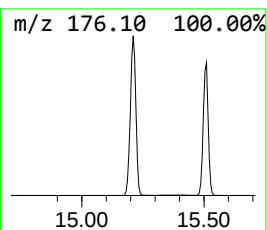
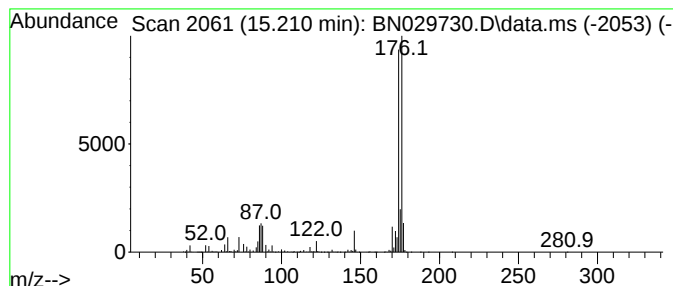
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.210	44.54 ng/ul	6006330	Acenaphthene-d10	14.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			[2-Amino-1-(2-chlorophenyl)ethyl]...	342	C16H31ClN2Si2	1000479-52-3	32
2			1,2,3,4-Tetrahydro-1-methyl-2,3-...	176	C9H8N2O2	020934-51-4	27
3			s-Triazolo[4,3-a]pyrazine, 3-eth...	176	C9H12N4	019848-81-8	25
4			Trifluoromethylbenzene, 3,4-diam...	176	C7H7F3N2	000368-71-8	17
5			N-(trimethylsilyl)-4,5-dihydroth...	174	C6H14N2SSi	1000474-12-9	17



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
Data File : BN029730.D
Acq On : 13 Feb 2024 12:27
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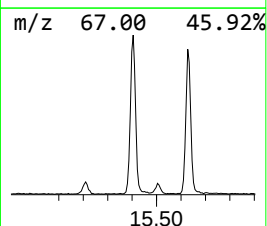
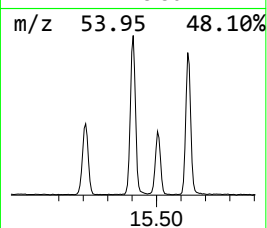
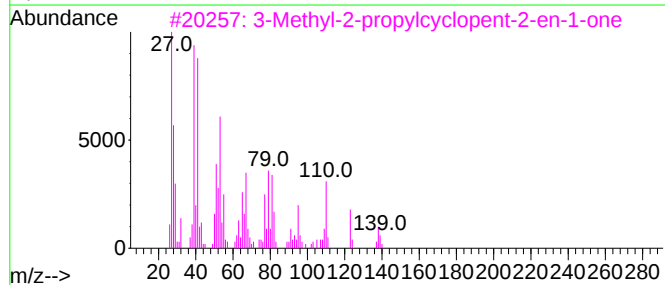
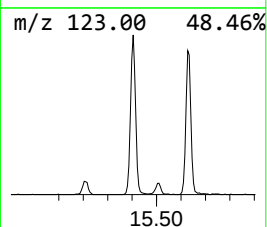
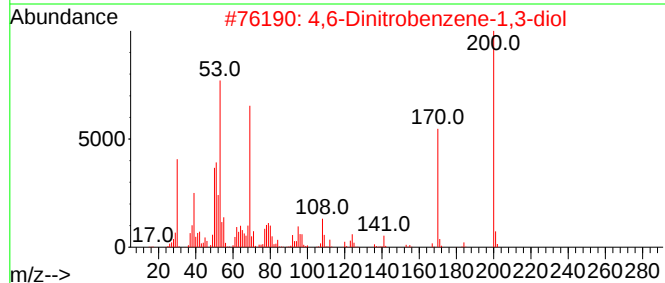
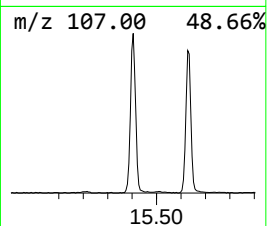
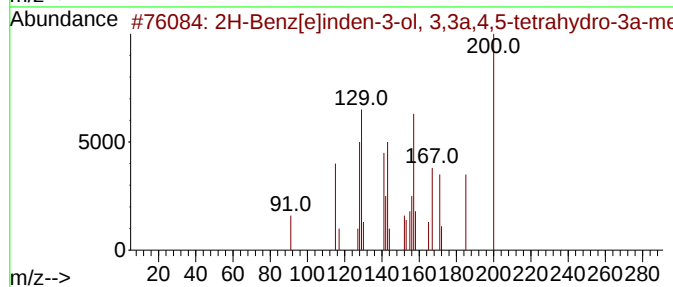
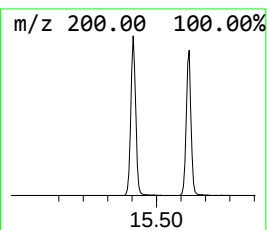
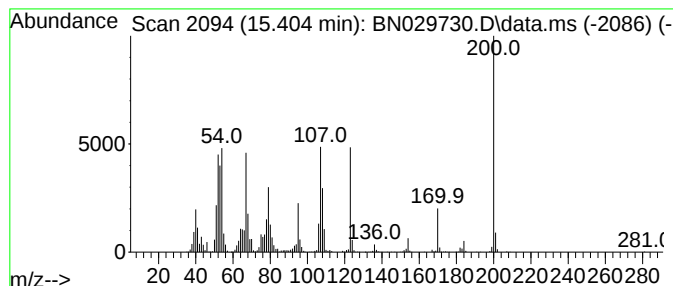
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-03 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.404	11.91 ng/ul	1605550	Acenaphthene-d10	14.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Benz[e]inden-3-ol, 3,3a,4,5-t...	200	C14H16O	071805-91-9	22		
2	4,6-Dinitrobenzene-1,3-diol	200	C6H4N2O6	000616-74-0	16		
3	3-Methyl-2-propylcyclopent-2-en-...	138	C9H14O	1000423-46-9	9		
4	5-Methyl-5-formylamino-6-imino-h...	200	C6H8N4O2S	071500-77-1	4		
5	1,4-benzenediol, 2,5-dinitro-	200	C6H4N2O6	1000400-38-8	4		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
Data File : BN029730.D
Acq On : 13 Feb 2024 12:27
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Sample : P1305-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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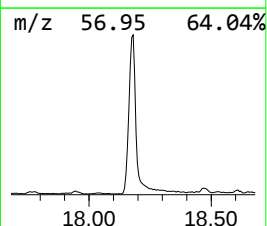
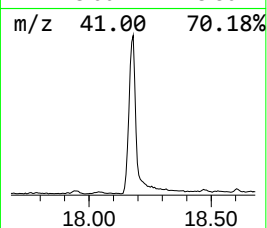
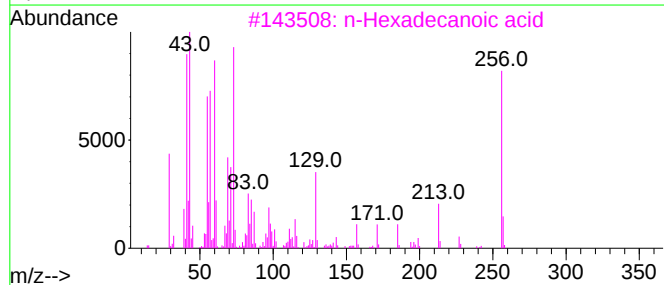
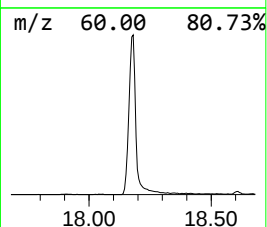
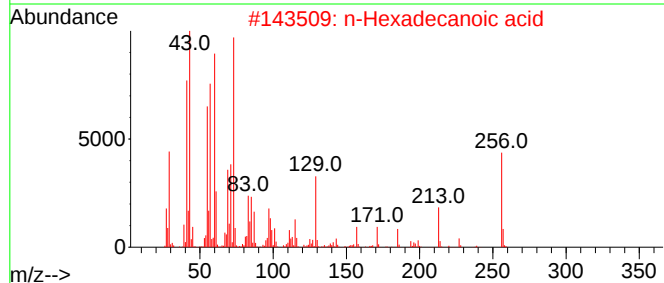
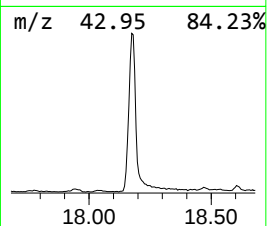
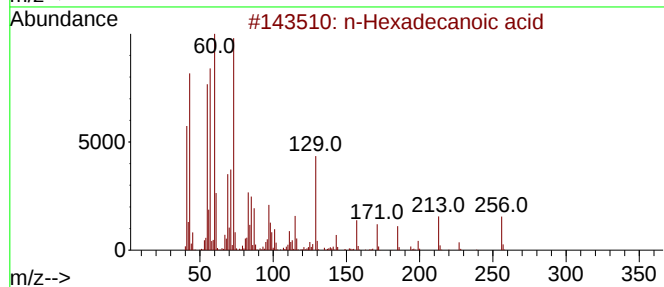
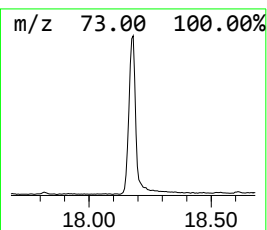
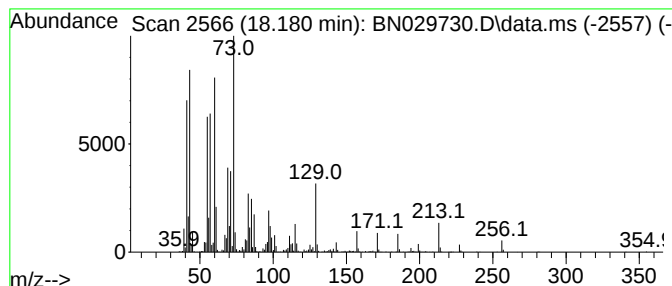
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.180	6.62 ng/ul	2760390	Phenanthrene-d10	17.275

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	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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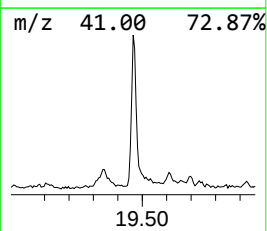
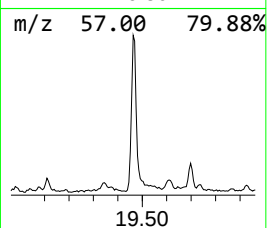
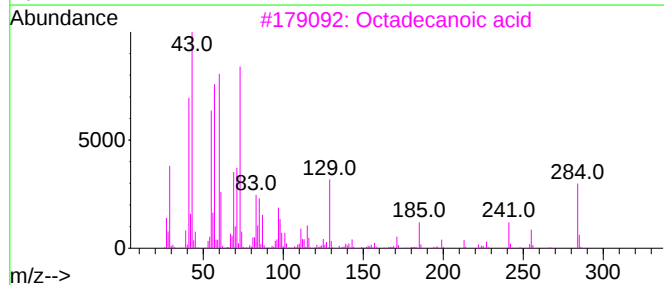
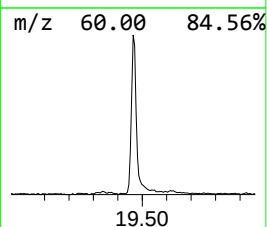
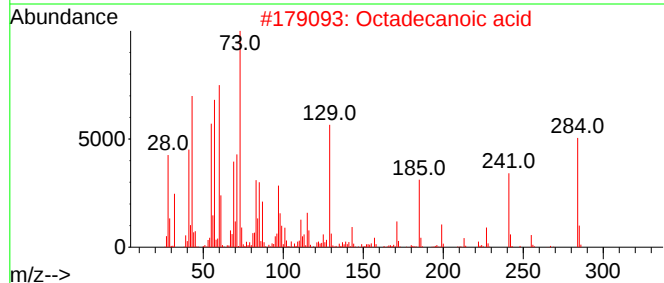
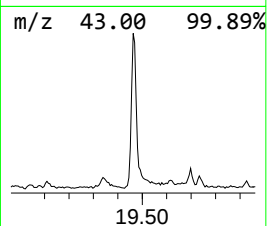
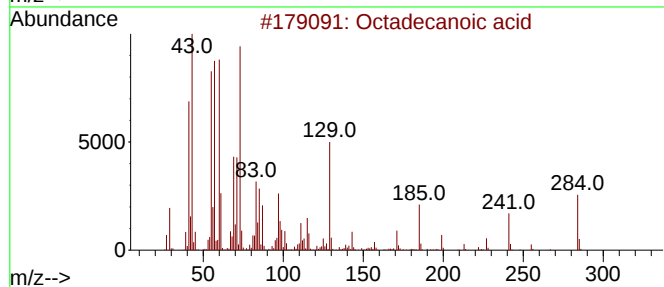
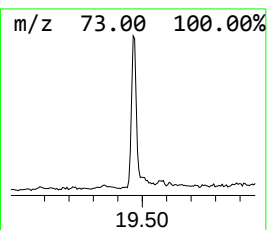
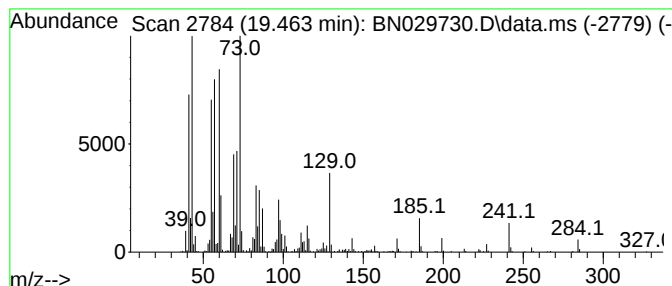
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.463	3.23 ng/ul	882728	Chrysene-d12	21.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	96
4			Octadecanoic acid	284	C18H36O2	000057-11-4	93
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
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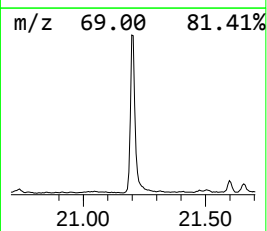
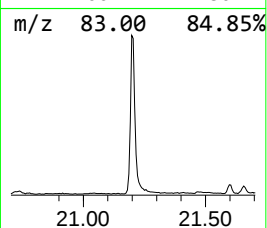
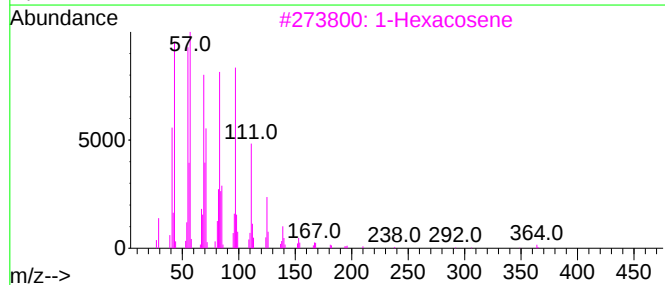
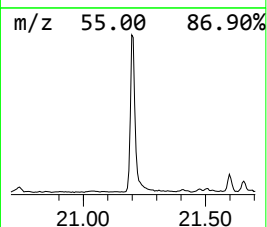
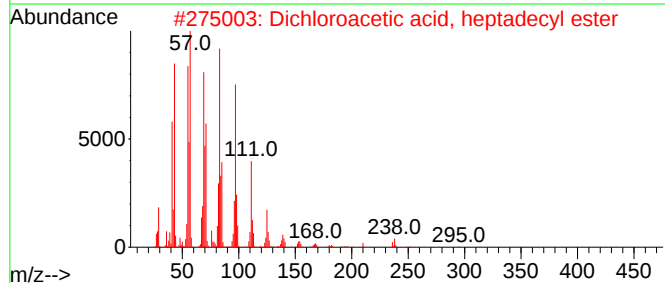
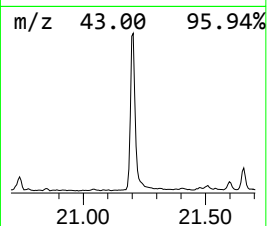
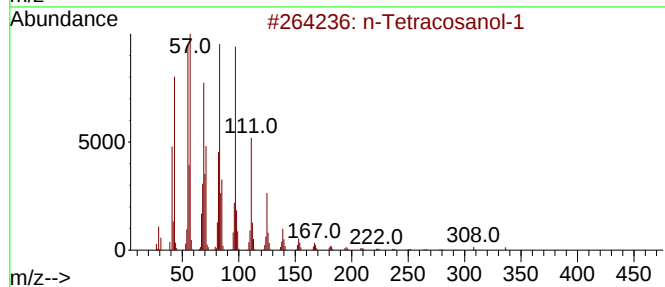
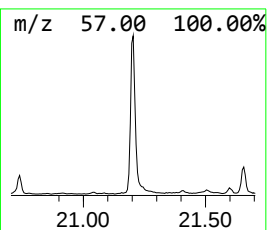
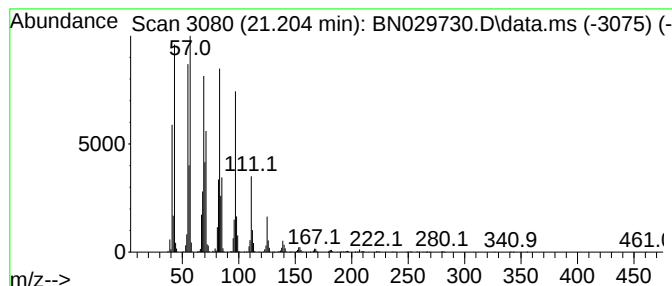
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 n-Tetracosanol-1 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.204	7.63 ng/ul	2084630	Chrysene-d12	21.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
2			Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
3			1-Hexacosene	364	C26H52	018835-33-1	91
4			2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	91
5			1-Heptacosanol	396	C27H56O	002004-39-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
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Sample : P1305-02
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ALS Vial : 5 Sample Multiplier: 1

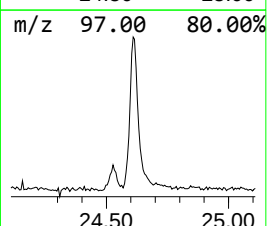
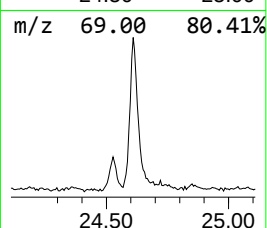
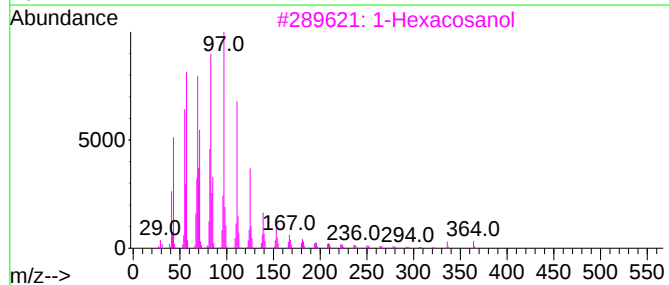
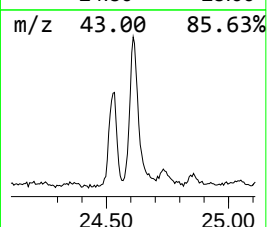
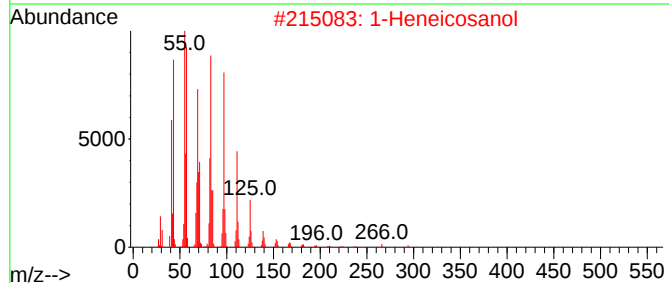
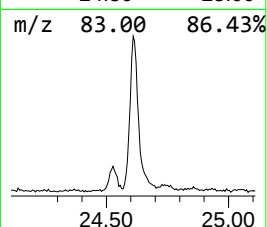
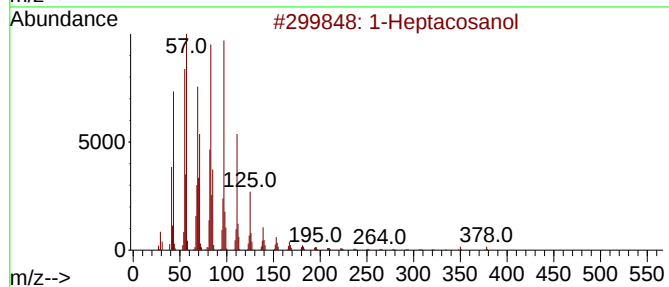
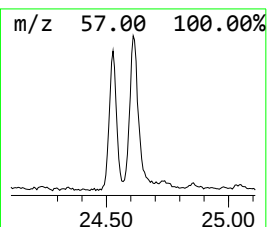
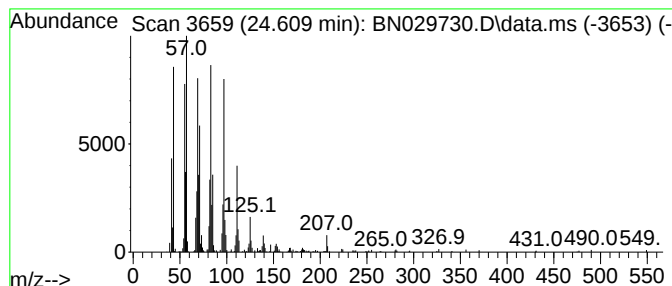
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BNA_N
ClientSampleId :
C0FP3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 1-Heptacosanol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.609	2.55 ng/ul	751061	Perylene-d12	23.851	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptacosanol	396	C27H56O	002004-39-9	94
2	1-Heneicosanol	312	C21H44O	015594-90-8	93
3	1-Hexacosanol	382	C26H54O	000506-52-5	91
4	Triacontan-15-ol	438	C30H62O	031849-26-0	91
5	Octacosanol	410	C28H58O	000557-61-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
Data File : BN029730.D
Acq On : 13 Feb 2024 12:27
Operator : MA/JU
Sample : P1305-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0FP3

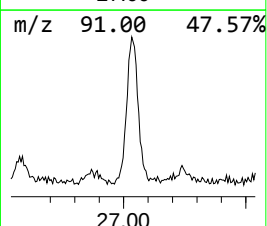
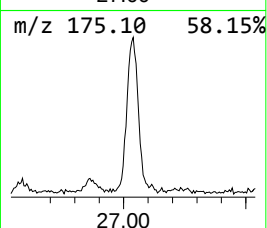
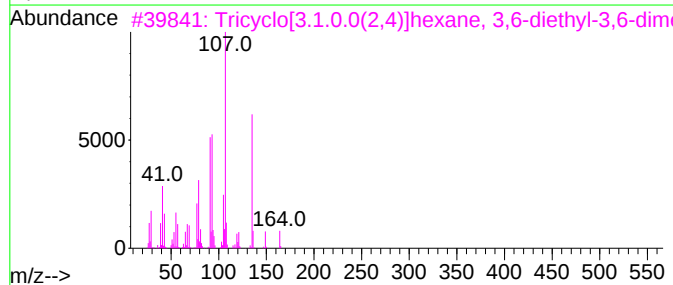
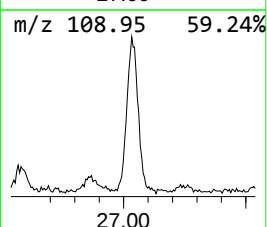
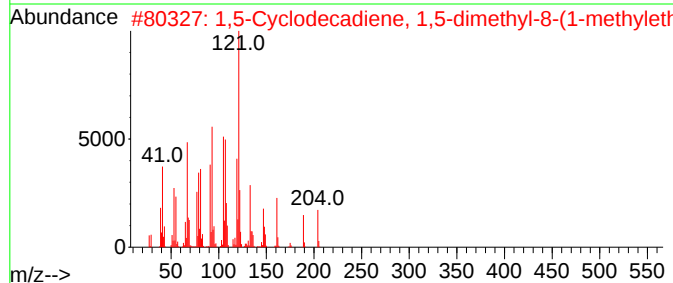
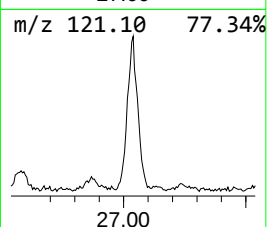
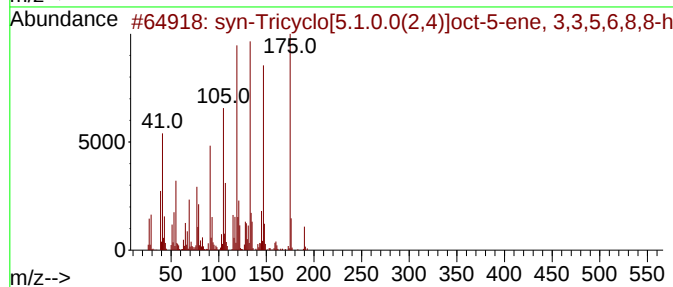
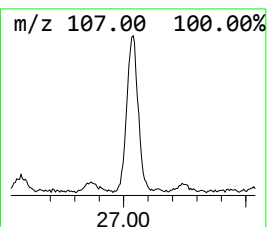
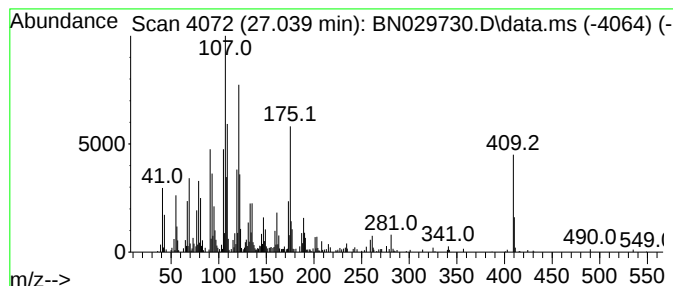
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-04 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.039	2.52 ng/ul	742159	Perylene-d12	23.851

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			syn-Tricyclo[5.1.0.0(2,4)]oct-5-ene, 3,3,5,6,8,8-h	190	C14H22	1000161-99-5	15
2			1,5-Cyclodecadiene, 1,5-dimethyl-8-(1-methylethyl)-3,6-dimethyl-	204	C15H24	015423-57-1	14
3			Tricyclo[3.1.0.0(2,4)]hexane, 3,6-diethyl-3,6-dimethyl-	164	C12H20	1000150-21-2	14
4			Phenol, 2-ethyl-	122	C8H10O	000090-00-6	11
5			Resedine	163	C9H9NO2	060426-44-0	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
Data File : BN029730.D
Acq On : 13 Feb 2024 12:27
Operator : MA/JU
Sample : P1305-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0FP3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.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
4-Chloro-3-fluo...	4.905	17.9	ng/u1	1459430	1	7.857	1632430	20.0
3,5-Diamino-2-m...	13.328	31.5	ng/u1	4248770	3	14.510	2697080	20.0
Isophosphinolin...	13.492	38.3	ng/u1	5167410	3	14.510	2697080	20.0
unknown-01	14.398	12.4	ng/u1	1665420	3	14.510	2697080	20.0
unknown-02	15.210	44.5	ng/u1	6006330	3	14.510	2697080	20.0
unknown-03	15.404	11.9	ng/u1	1605550	3	14.510	2697080	20.0
n-Hexadecanoic ...	18.180	6.6	ng/u1	2760390	4	17.275	8342480	20.0
Octadecanoic acid	19.463	3.2	ng/u1	882728	5	21.468	5467740	20.0
n-Tetracosanol-1	21.204	7.6	ng/u1	2084630	5	21.468	5467740	20.0
1-Heptacosanol	24.609	2.5	ng/u1	751061	6	23.851	5880160	20.0
unknown-04	27.039	2.5	ng/u1	742159	6	23.851	5880160	20.0