

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
 Data File : BN032170.D
 Acq On : 22 Jun 2024 17:14
 Operator : MA/JU
 Sample : P2899-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AA7

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-BN052924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN032170.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.163	26	30	39	rVB	47891	67996	1.66%	0.096%
2	3.434	73	76	85	rVB	35746	53744	1.31%	0.076%
3	3.581	100	101	111	rVV	48030	67950	1.66%	0.096%
4	3.669	111	116	122	rVB	47221	63998	1.56%	0.091%
5	3.875	147	151	157	rVB	22011	32169	0.79%	0.046%
6	4.422	239	244	248	rVV	48784	73178	1.79%	0.104%
7	4.769	297	303	312	rBV	282929	431689	10.55%	0.612%
8	4.857	312	318	331	rVB2	21471	56867	1.39%	0.081%
9	5.010	336	344	353	rVB2	14727	30078	0.74%	0.043%
10	5.293	387	392	400	rBV	16649	32236	0.79%	0.046%
11	5.657	448	454	460	rVV4	27064	47577	1.16%	0.067%
12	5.734	460	467	477	rVB2	33128	68914	1.68%	0.098%
13	6.151	525	538	542	rBV4	17313	44344	1.08%	0.063%
14	6.310	559	565	573	rVB	174907	295996	7.23%	0.420%
15	6.657	619	624	631	rVB5	23839	54014	1.32%	0.077%
16	6.810	638	650	664	rBV	830912	1438793	35.16%	2.040%
17	6.975	672	678	688	rVB	703937	1107842	27.07%	1.571%
18	7.063	688	693	703	rVB	120398	200734	4.91%	0.285%
19	7.163	703	710	720	rBV	908617	1471276	35.95%	2.086%
20	7.246	720	724	727	rBV2	22738	32743	0.80%	0.046%
21	7.628	776	789	798	rBV	1291064	2103713	51.41%	2.982%
22	7.898	830	835	841	rVB	56635	90421	2.21%	0.128%
23	8.169	871	881	884	rBV2	18364	48346	1.18%	0.069%
24	8.340	903	910	924	rVV	985717	1685424	41.19%	2.389%
25	8.451	924	929	936	rVV	102105	184840	4.52%	0.262%
26	8.787	979	986	993	rVV	823872	1386035	33.87%	1.965%
27	8.834	993	994	1002	rVV	32648	38810	0.95%	0.055%
28	9.351	1068	1082	1090	rBV	88361	163655	4.00%	0.232%
29	9.504	1100	1108	1115	rBV	652234	1134203	27.72%	1.608%
30	9.681	1131	1138	1140	rBV3	31938	72485	1.77%	0.103%
31	9.722	1141	1145	1156	rVV8	43063	122667	3.00%	0.174%
32	10.040	1191	1199	1210	rBV	1027433	1900101	46.43%	2.694%
33	10.410	1251	1262	1268	rVV	1764475	3091414	75.54%	4.383%
34	10.457	1268	1270	1275	rVV	53054	75671	1.85%	0.107%
35	10.516	1275	1280	1282	rVV4	22368	42397	1.04%	0.060%
36	10.563	1282	1288	1304	rVB2	72283	190245	4.65%	0.270%

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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_N\Methods\SFAM-EPA-BN052924.MA.M
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37	10.975	1349	1358	1366	rBV2	85573	318535	7.78%	0.452%
38	11.157	1384	1389	1395	rBV	94149	182626	4.46%	0.259%
39	11.422	1429	1434	1444	rVB2	32507	58095	1.42%	0.082%
40	11.822	1495	1502	1508	rVB5	51711	115051	2.81%	0.163%
41	11.998	1526	1532	1540	rVV3	25219	51169	1.25%	0.073%
42	12.075	1540	1545	1555	rVB	79879	134227	3.28%	0.190%
43	12.298	1578	1583	1592	rVB	66365	109331	2.67%	0.155%
44	12.392	1593	1599	1605	rBV2	20513	43591	1.07%	0.062%
45	12.798	1664	1668	1676	rVV3	21831	43868	1.07%	0.062%
46	13.092	1713	1718	1723	rBV2	56357	95678	2.34%	0.136%
47	13.139	1723	1726	1733	rVV9	17385	31704	0.77%	0.045%
48	13.269	1744	1748	1754	rVB4	26385	36810	0.90%	0.052%
49	13.433	1766	1776	1777	rBV4	25343	52377	1.28%	0.074%
50	13.451	1777	1779	1785	rVV3	31086	46231	1.13%	0.066%
51	13.533	1787	1793	1798	rVV	205003	327066	7.99%	0.464%
52	13.586	1798	1802	1808	rVV2	67410	120729	2.95%	0.171%
53	13.645	1808	1812	1814	rVV	47353	64508	1.58%	0.091%
54	13.692	1814	1820	1826	rVV	1958711	2752719	67.27%	3.902%
55	13.751	1826	1830	1834	rVV2	42939	62202	1.52%	0.088%
56	13.828	1840	1843	1847	rVV	28809	39895	0.97%	0.057%
57	13.969	1856	1867	1882	rBV	2221033	3467800	84.74%	4.916%
58	14.110	1884	1891	1895	rVB10	17853	38278	0.94%	0.054%
59	14.169	1895	1901	1906	rVB	63298	98841	2.42%	0.140%
60	14.275	1909	1919	1926	rBV	2608376	3989912	97.50%	5.656%
61	14.504	1949	1958	1970	rBV	606375	1267492	30.97%	1.797%
62	14.657	1979	1984	1985	rBV	53113	75062	1.83%	0.106%
63	14.798	2004	2008	2017	rVB10	18024	37182	0.91%	0.053%
64	14.892	2017	2024	2030	rBV5	20522	39283	0.96%	0.056%
65	15.069	2048	2054	2057	rBV4	39569	72493	1.77%	0.103%
66	15.127	2061	2064	2069	rVB	61463	71277	1.74%	0.101%
67	15.269	2081	2088	2094	rBV	2714411	4092166	100.00%	5.801%
68	15.404	2106	2111	2117	rBV	625436	933158	22.80%	1.323%
69	15.451	2117	2119	2125	rVB3	28509	40133	0.98%	0.057%
70	15.622	2143	2148	2155	rBV	39258	72912	1.78%	0.103%
71	15.769	2169	2173	2182	rVB	38094	77525	1.89%	0.110%
72	15.851	2182	2187	2192	rVB5	24984	44765	1.09%	0.063%
73	16.010	2206	2214	2220	rBV2	110410	252754	6.18%	0.358%
74	16.157	2232	2239	2240	rBV6	20110	31517	0.77%	0.045%
75	16.439	2282	2287	2294	rBV6	25887	58726	1.44%	0.083%
76	16.569	2301	2309	2313	rBV5	32801	84190	2.06%	0.119%

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77	16.686	2325	2329	2336	rBV5	24900	44988	1.10%	0.064%
78	16.786	2342	2346	2349	rVV	63197	76683	1.87%	0.109%
79	16.927	2367	2370	2376	rBV4	32101	56761	1.39%	0.080%
80	17.027	2380	2387	2392	rVV	2728766	4077225	99.63%	5.780%
81	17.069	2392	2394	2397	rVV	84761	98596	2.41%	0.140%
82	17.121	2397	2403	2412	rVV2	2602942	3871475	94.61%	5.488%
83	17.204	2415	2417	2423	rVB6	24905	39686	0.97%	0.056%
84	17.521	2468	2471	2481	rVB	66208	89625	2.19%	0.127%
85	17.804	2515	2519	2525	rBV6	44177	80607	1.97%	0.114%
86	17.927	2535	2540	2550	rBV	657100	1050235	25.66%	1.489%
87	18.192	2581	2585	2588	rBV	152064	222163	5.43%	0.315%
88	18.263	2594	2597	2601	rVB4	85243	111247	2.72%	0.158%
89	18.545	2640	2645	2650	rVB	752695	964176	23.56%	1.367%
90	18.604	2652	2655	2657	rBV3	47431	60286	1.47%	0.085%
91	19.216	2755	2759	2770	rBV2	576455	1041947	25.46%	1.477%
92	19.427	2790	2795	2801	rBV	2581373	3565382	87.13%	5.054%
93	20.310	2941	2945	2949	rBV	1664362	2067092	50.51%	2.930%
94	20.351	2949	2952	2958	rVB	990571	1324098	32.36%	1.877%
95	20.921	3045	3049	3051	rBV	1011241	1319841	32.25%	1.871%
96	20.951	3051	3054	3065	rVB	892276	1368028	33.43%	1.939%
97	21.262	3102	3107	3117	rVB2	2301169	3683107	90.00%	5.221%
98	22.033	3234	3238	3248	rVB	808453	1484835	36.28%	2.105%
99	23.968	3561	3567	3576	rVB3	979844	2355887	57.57%	3.340%
100	24.174	3594	3602	3615	rVB	1570428	3951257	96.56%	5.601%

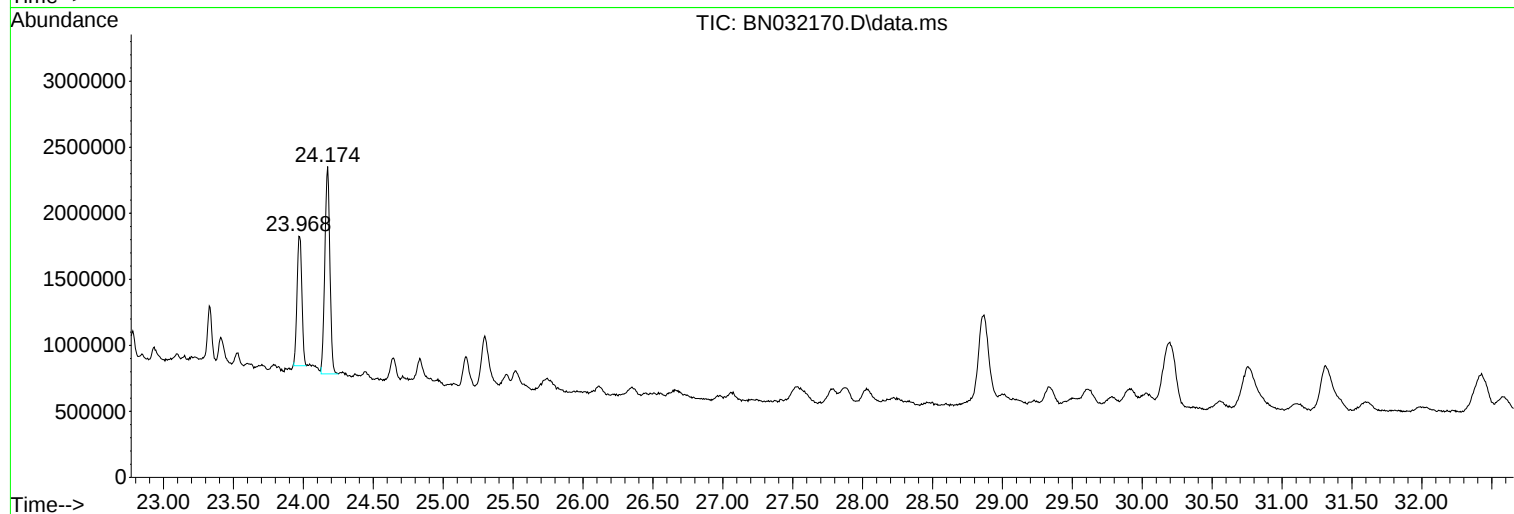
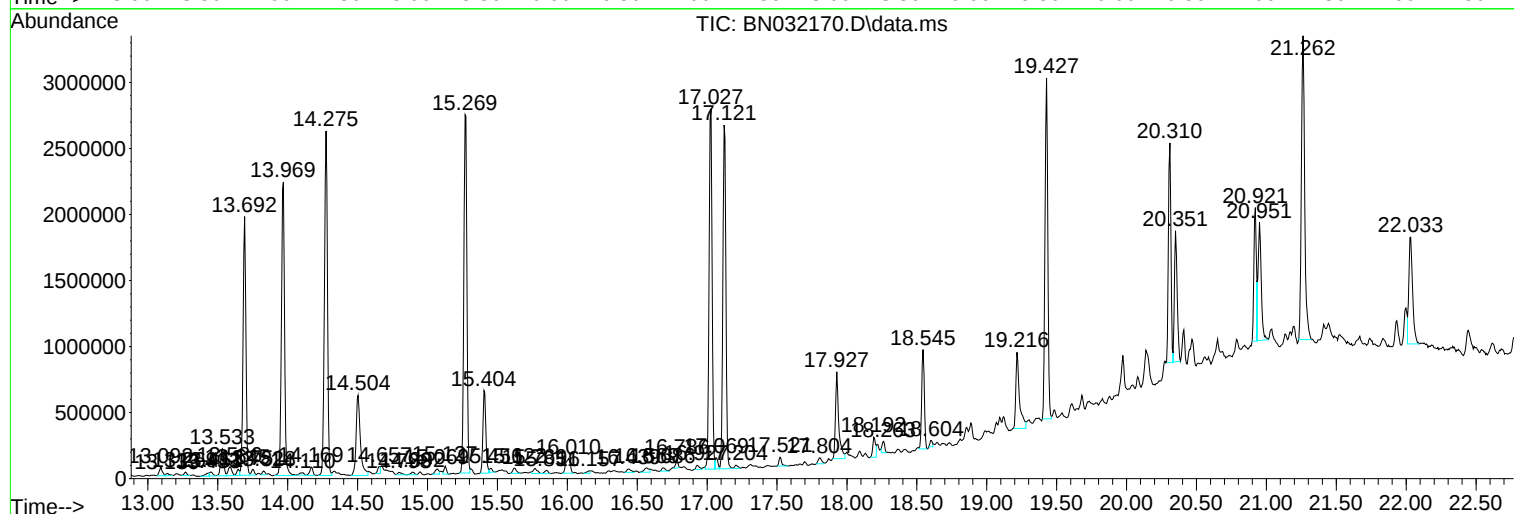
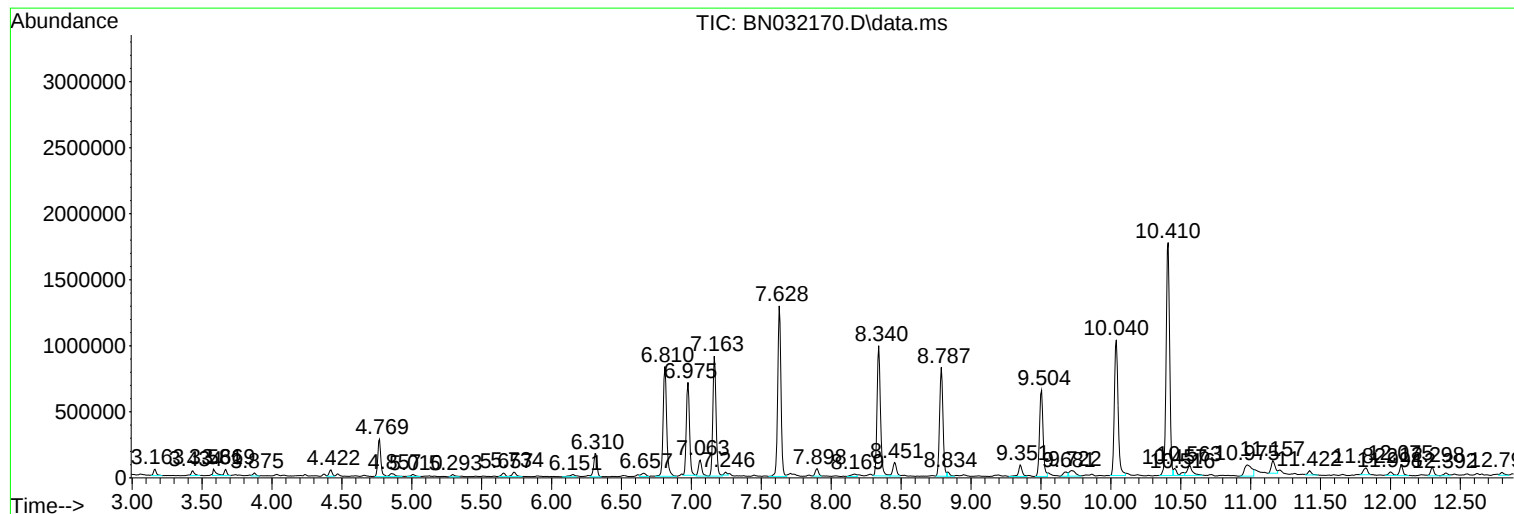
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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



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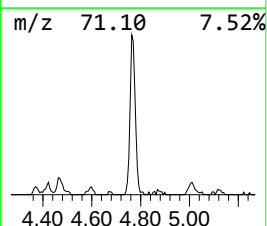
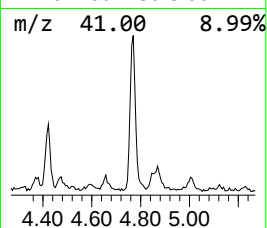
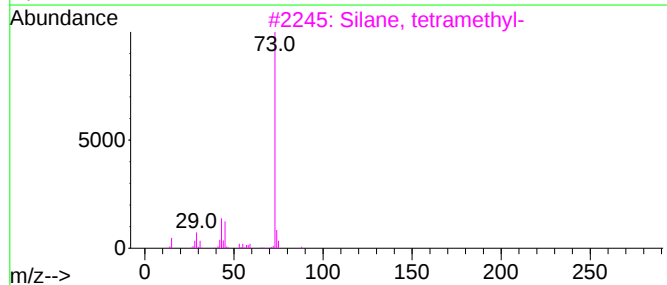
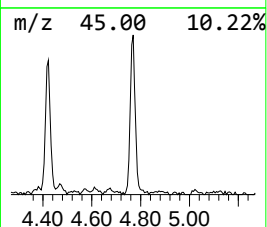
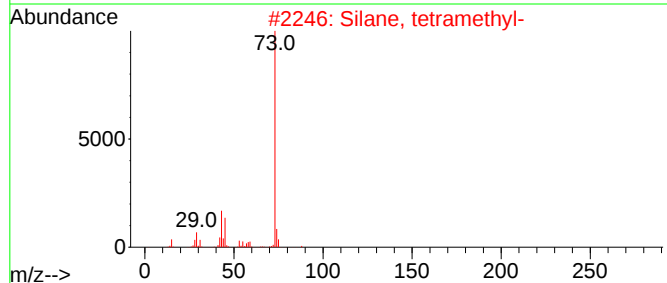
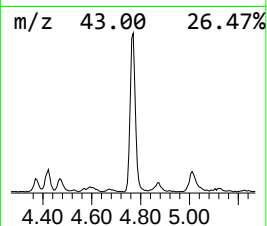
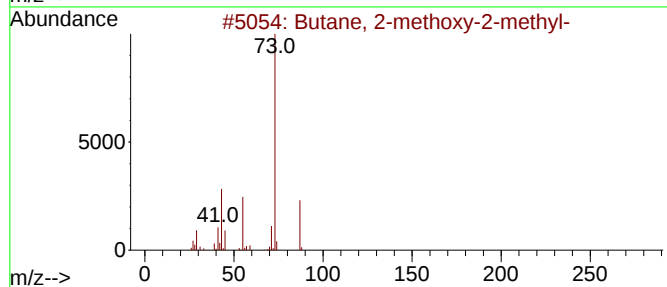
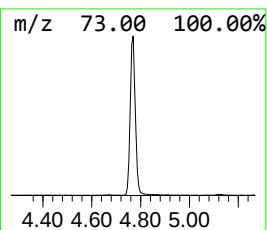
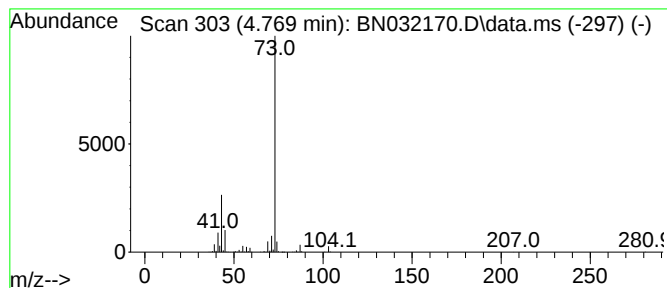
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.769	4.10 ng/ul	431689	1,4-Dichlorobenzene-d4	7.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	45
2			Silane, tetramethyl-	88	C4H12Si	000075-76-3	43
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	37
4			Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	1000186-02-3	33
5			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	28



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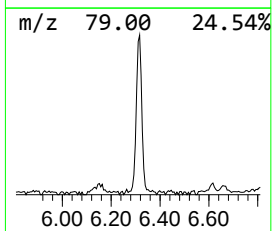
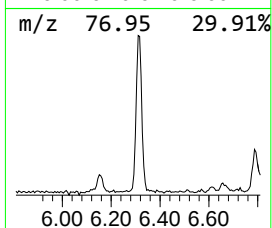
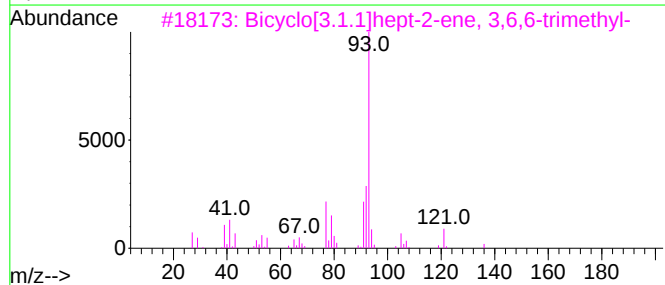
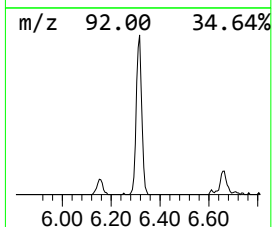
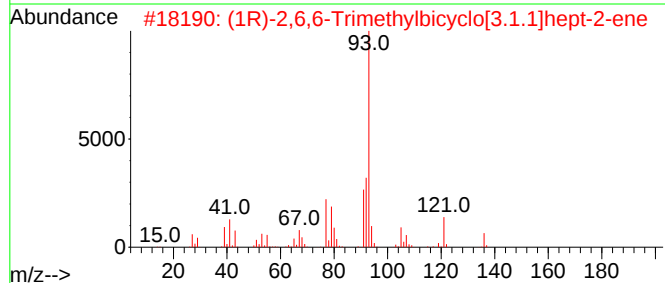
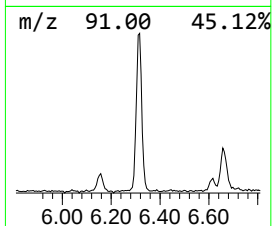
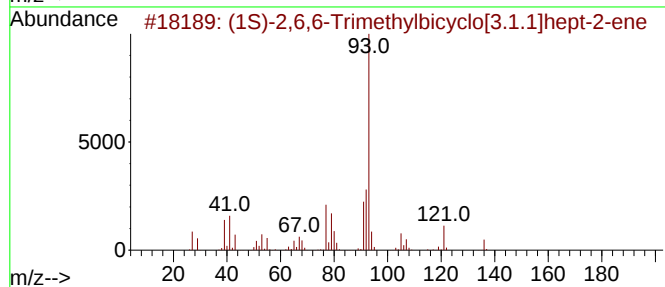
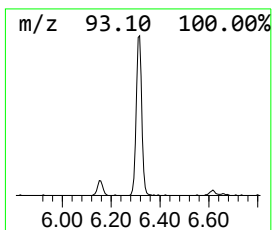
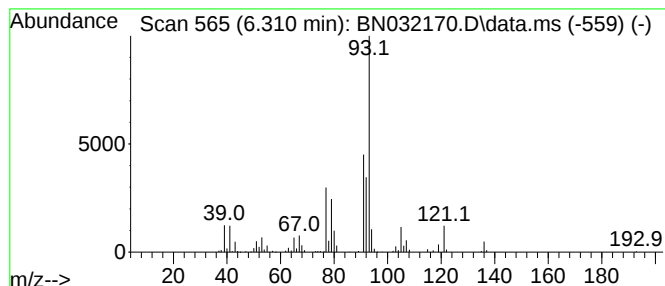
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (1S)-2,6,6-Trimethylbicyclo... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.310	2.81 ng/ul	295996	1,4-Dichlorobenzene-d4	7.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1S)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-26-4	95		
2	(1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-70-8	94		
3	Bicyclo[3.1.1]hept-2-ene, 3,6,6-trimethyl-	136	C10H16	004889-83-2	94		
4	.alpha.-Pinene	136	C10H16	000080-56-8	94		
5	(+)-3-Carene	136	C10H16	000498-15-7	91		



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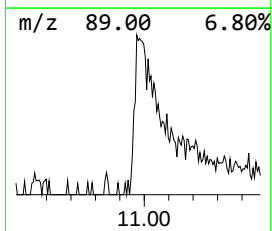
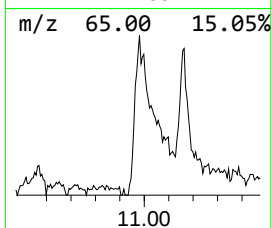
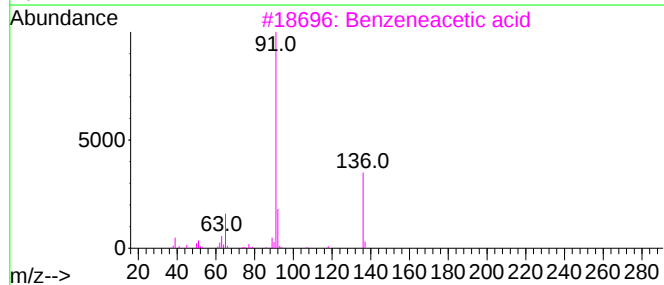
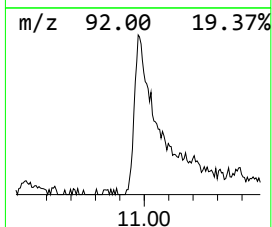
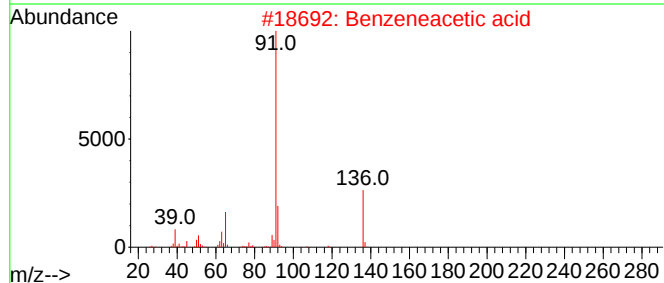
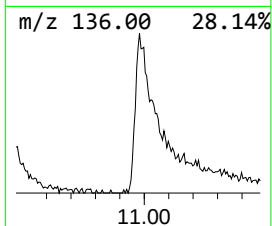
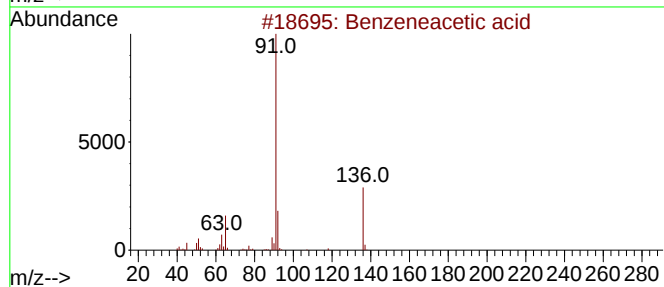
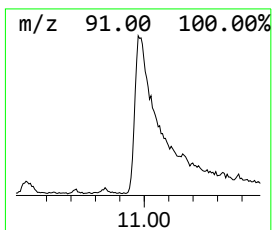
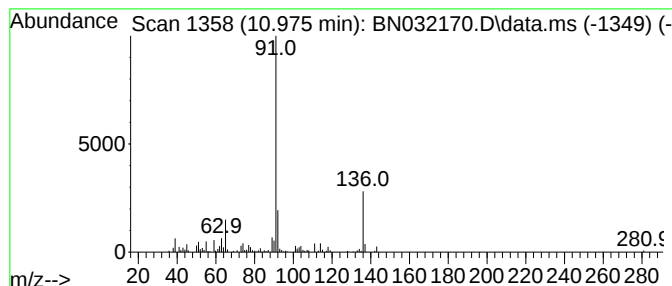
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzeneacetic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.975	2.06 ng/ul	318535	Naphthalene-d8	10.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneacetic acid	136	C8H8O2	000103-82-2	93
2			Benzeneacetic acid	136	C8H8O2	000103-82-2	93
3			Benzeneacetic acid	136	C8H8O2	000103-82-2	81
4			Propanedioic acid, phenyl-	180	C9H8O4	002613-89-0	80
5			Benzeneacetic acid	136	C8H8O2	000103-82-2	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
Data File : BN032170.D
Acq On : 22 Jun 2024 17:14
Operator : MA/JU
Sample : P2899-05
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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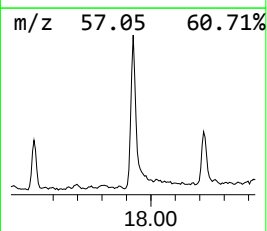
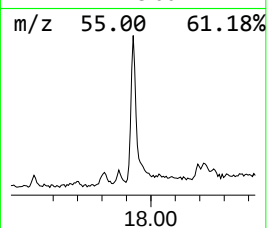
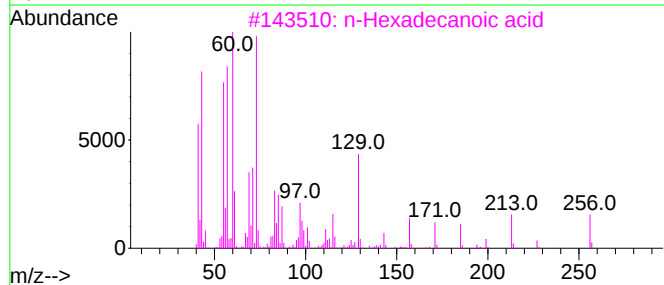
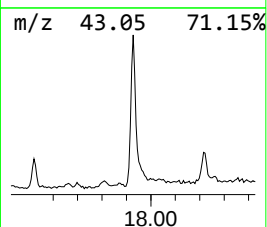
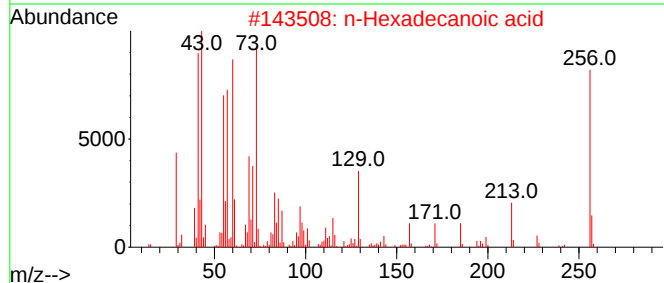
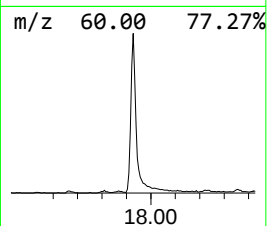
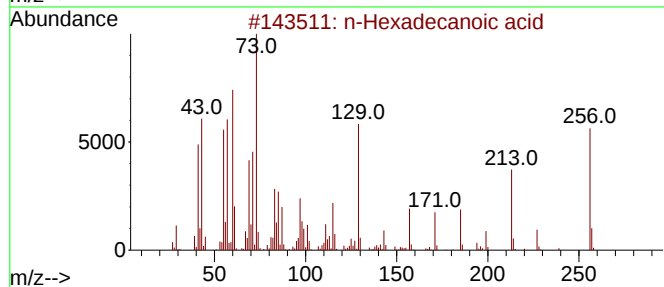
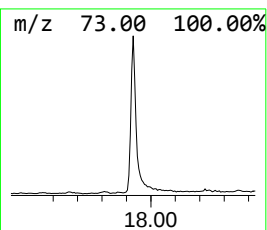
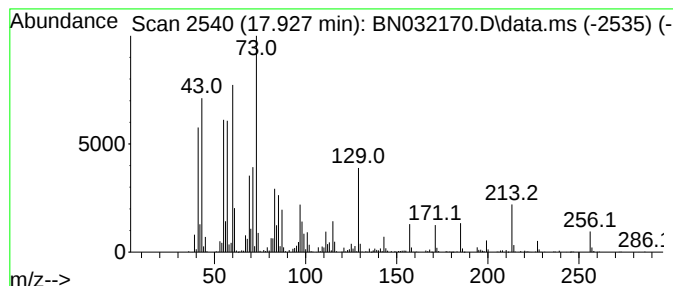
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.927	5.15 ng/ul	1050240	Phenanthrene-d10	17.027

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	99
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
Data File : BN032170.D
Acq On : 22 Jun 2024 17:14
Operator : MA/JU
Sample : P2899-05
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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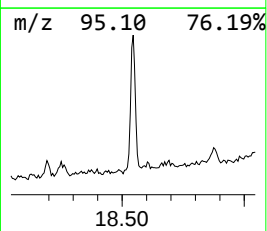
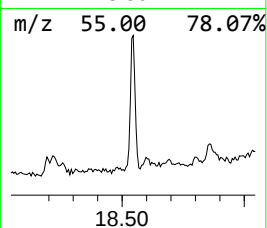
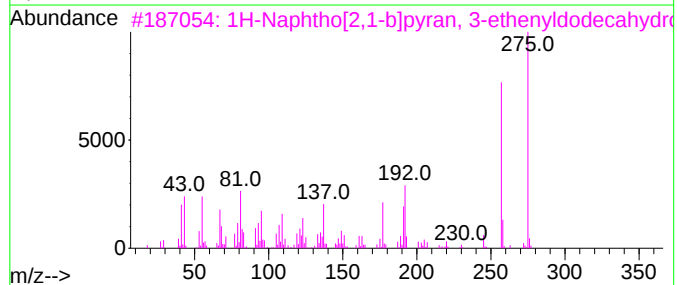
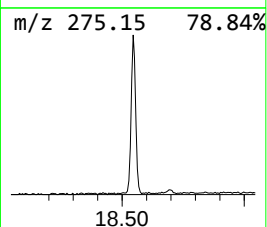
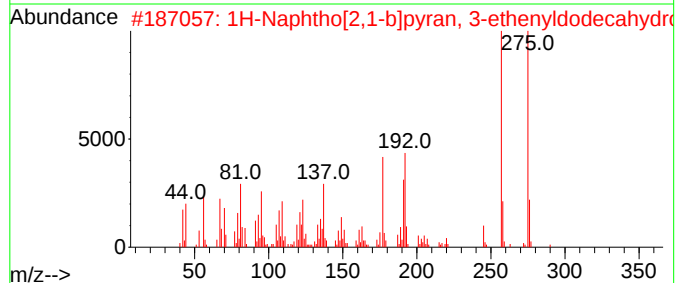
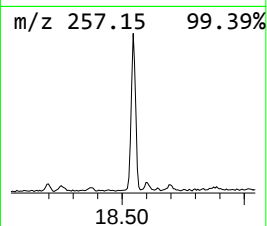
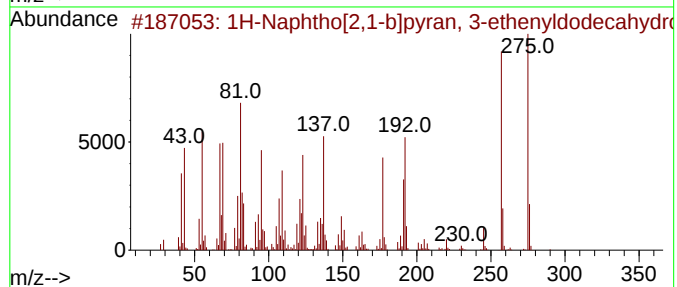
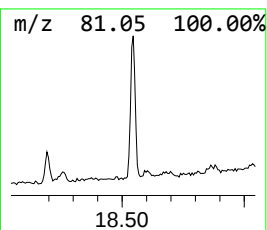
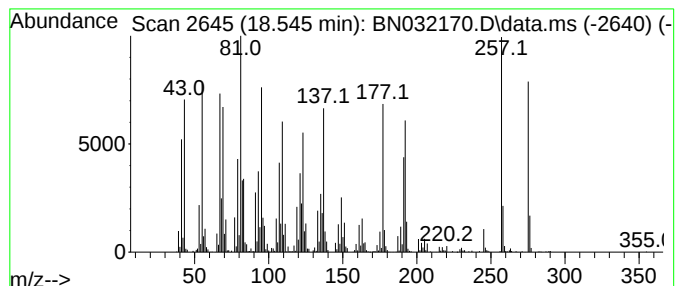
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1H-Naphtho[2,1-b]pyran, 3-e... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.545	4.73 ng/ul	964176	Phenanthrene-d10	17.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Naphtho[2,1-b]pyran, 3-etheny...	290	C20H34O	000596-84-9	96
2			1H-Naphtho[2,1-b]pyran, 3-etheny...	290	C20H34O	000596-84-9	91
3			1H-Naphtho[2,1-b]pyran, 3-etheny...	290	C20H34O	001227-93-6	78
4			4a,7,7,10a-Tetramethyl-dodecahyd...	264	C17H28O2	938512-84-6	60
5			1H-Naphtho[2,1-b]pyran, 3-etheny...	290	C20H34O	000596-84-9	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
Data File : BN032170.D
Acq On : 22 Jun 2024 17:14
Operator : MA/JU
Sample : P2899-05
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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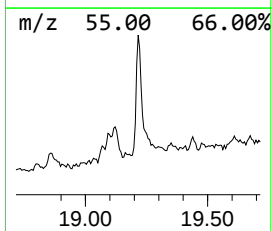
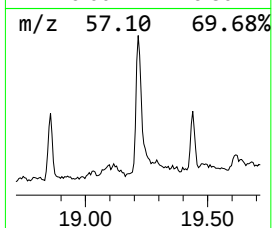
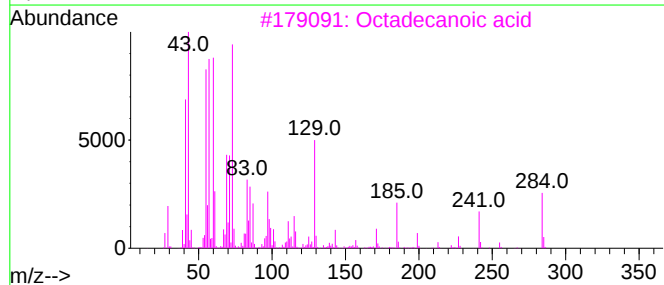
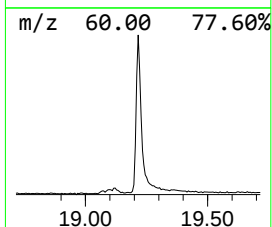
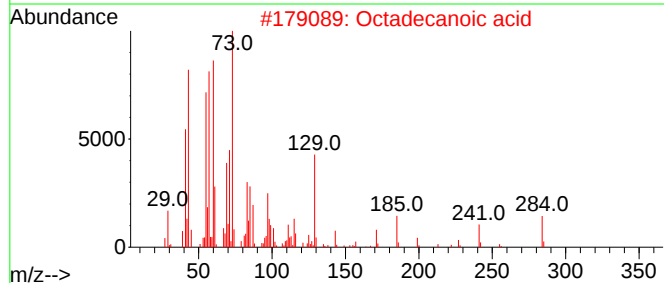
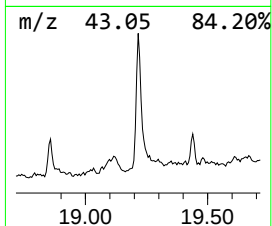
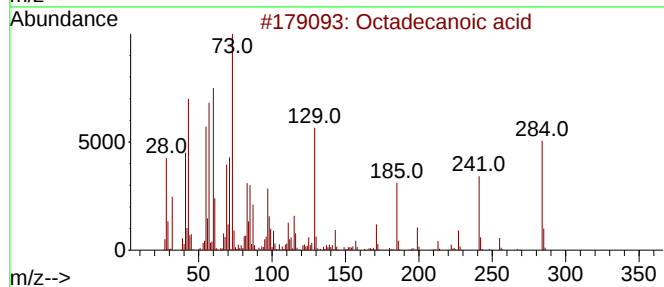
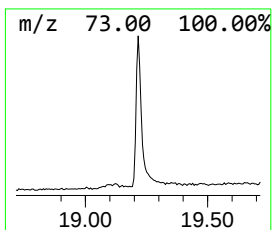
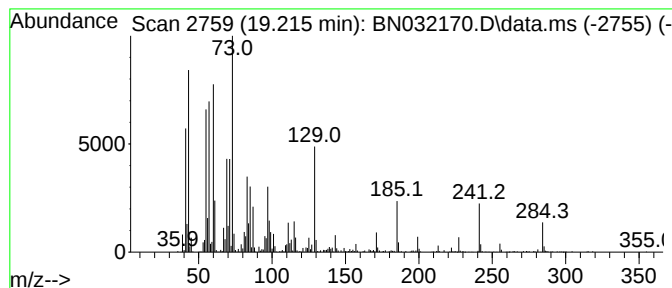
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.215	5.66 ng/ul	1041950	Chrysene-d12	21.262

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	99
4	Octadecanoic acid			284	C18H36O2	000057-11-4	95
5	Octadecanoic acid			284	C18H36O2	000057-11-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
Data File : BN032170.D
Acq On : 22 Jun 2024 17:14
Operator : MA/JU
Sample : P2899-05
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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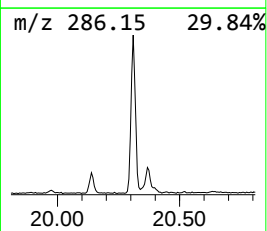
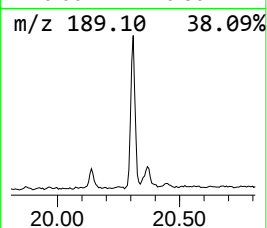
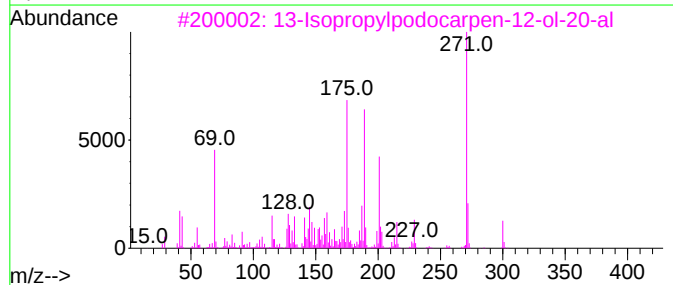
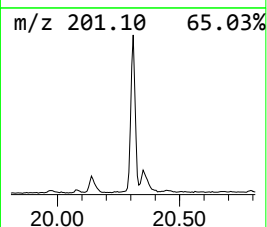
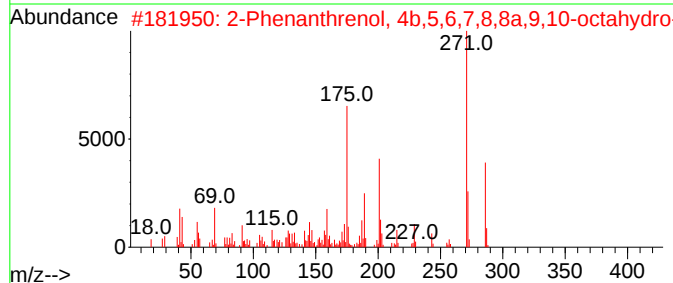
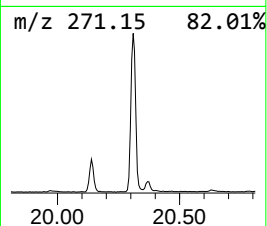
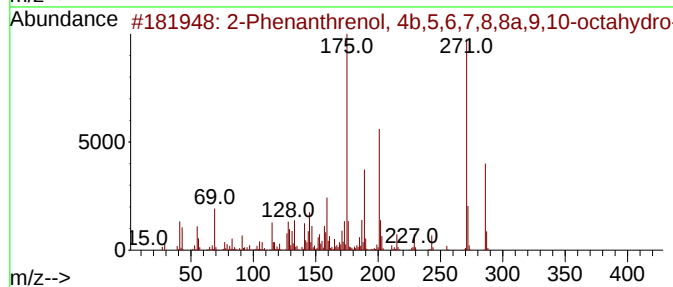
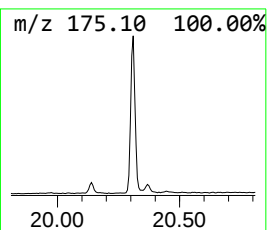
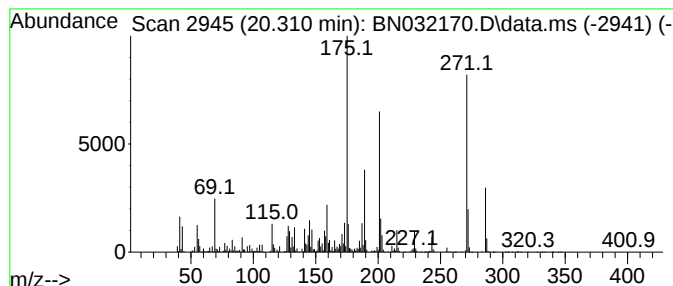
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.310	11.22 ng/ul	2067090	Chrysene-d12	21.262

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	99		
2	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	92		
3	13-Isopropylpodocarpin-12-ol-20-al	300	C20H28O2	024035-37-8	58		
4	13-Isopropylpodocarpin-12-ol-20-al	300	C20H28O2	024035-37-8	47		
5	Pisiferol	302	C20H30O2	024035-36-7	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
Data File : BN032170.D
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Misc :
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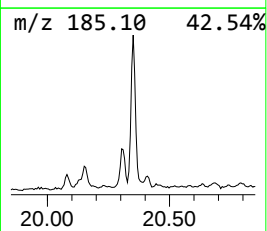
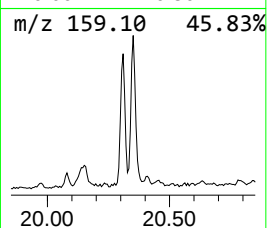
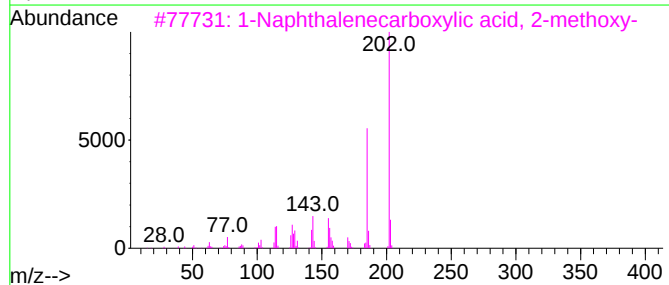
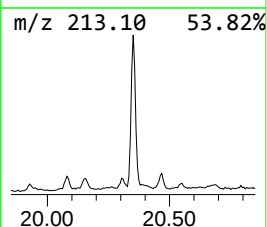
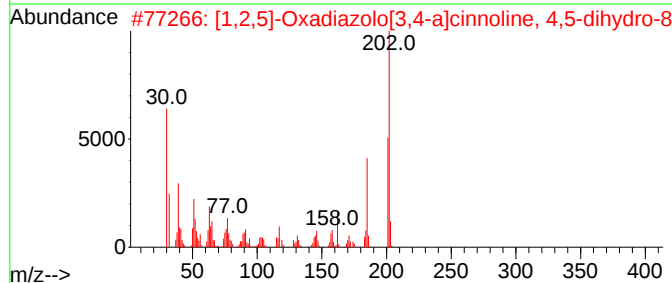
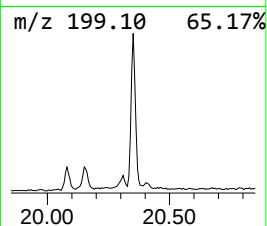
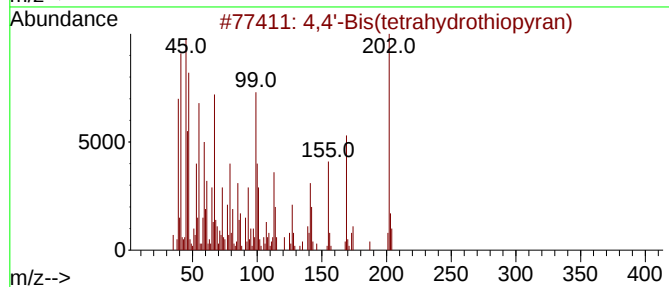
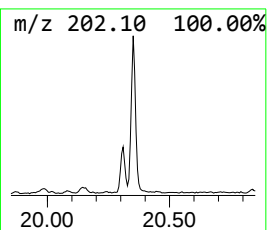
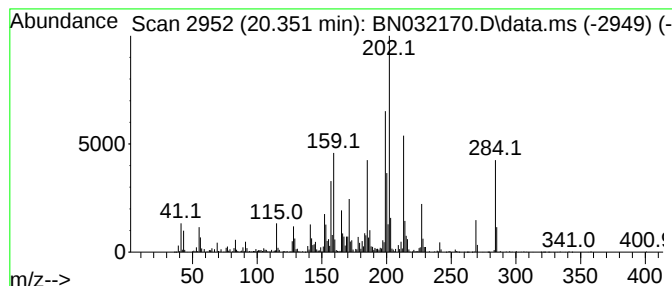
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 4,4'-Bis(tetrahydrothiopyran) Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.351	7.19 ng/ul	1324100	Chrysene-d12	21.262

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4,4'-Bis(tetrahydrothiopyran)	202	C10H18S2	116196-83-9	91
2			[1,2,5]-Oxadiazolo[3,4-a]cinnoli...	202	C10H10N4O	1000260-59-4	38
3			1-Naphthalenecarboxylic acid, 2-...	202	C12H10O3	000947-62-6	22
4			4-Hydroxyamino-6-phenylpyrimidin...	202	C10H10N4O	1000111-25-2	14
5			Pyrene	202	C16H10	000129-00-0	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
Data File : BN032170.D
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Sample : P2899-05
Misc :
ALS Vial : 12 Sample Multiplier: 1

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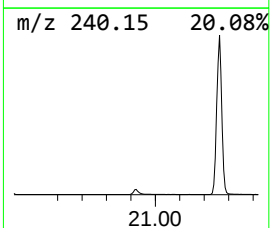
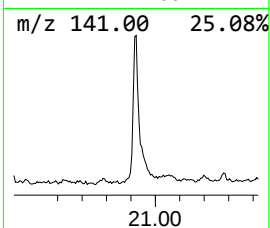
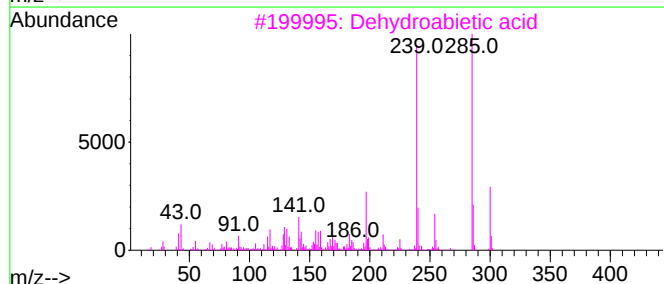
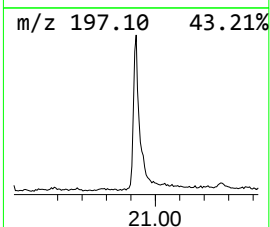
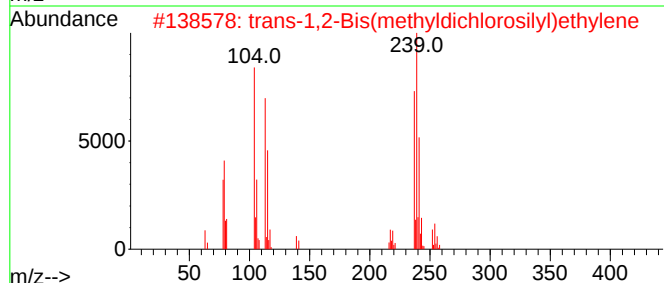
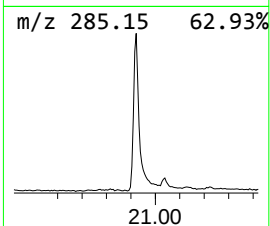
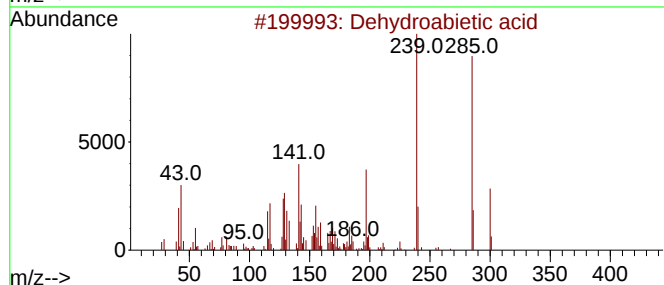
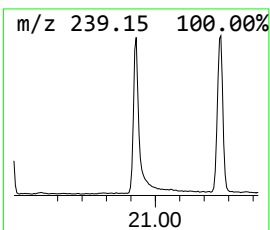
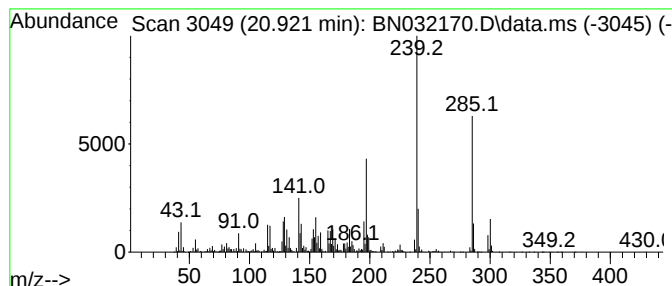
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Dehydroabietic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.921	7.17 ng/ul	1319840	Chrysene-d12	21.262

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dehydroabietic acid	300	C20H28O2	001740-19-8	99
2			trans-1,2-Bis(methyldichlorosilyl)ethylen...	252	C4H8Cl4Si2	065899-10-7	94
3			Dehydroabietic acid	300	C20H28O2	001740-19-8	94
4			Dehydroabietic acid	300	C20H28O2	001740-19-8	91
5			1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
Data File : BN032170.D
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Sample : P2899-05
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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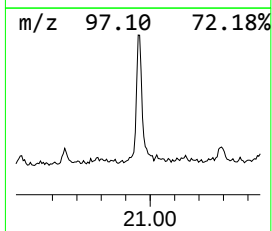
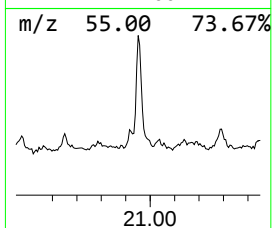
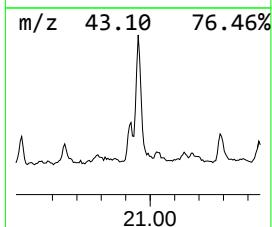
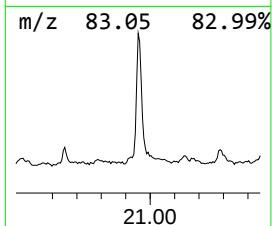
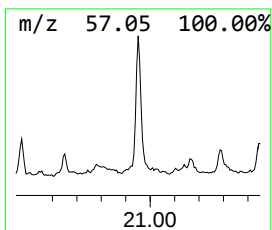
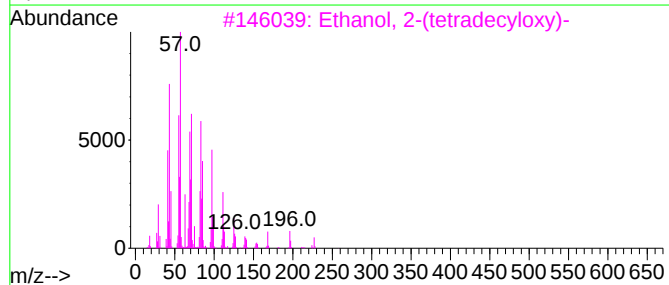
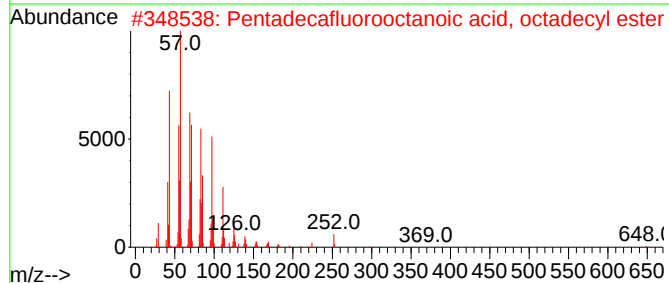
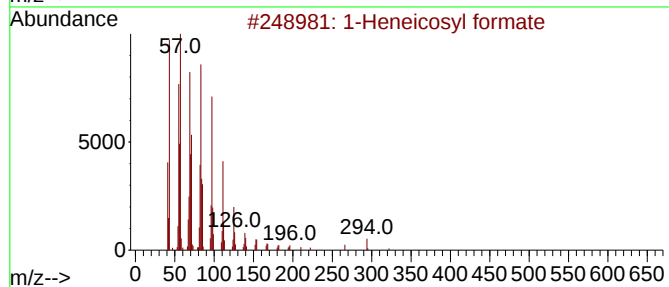
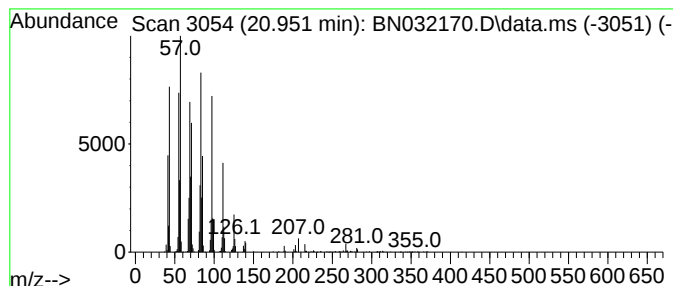
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 1-Heneicosyl formate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.951	7.43 ng/ul	1368030	Chrysene-d12	21.262

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosyl formate	340	C22H44O2	077899-03-7	95
2			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91
3			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	90
4			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	90
5			1-Docosene	308	C22H44	001599-67-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
Data File : BN032170.D
Acq On : 22 Jun 2024 17:14
Operator : MA/JU
Sample : P2899-05
Misc :
ALS Vial : 12 Sample Multiplier: 1

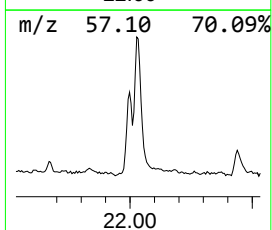
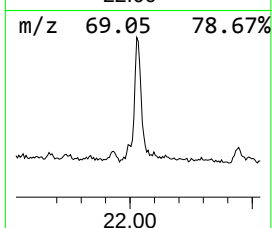
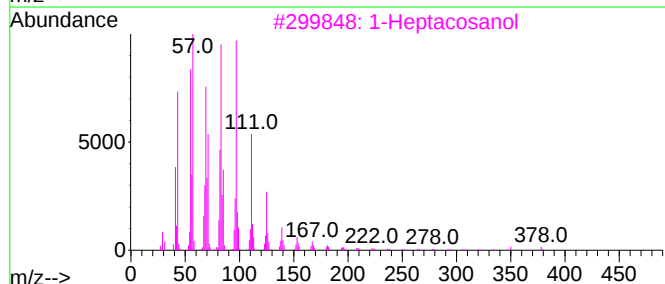
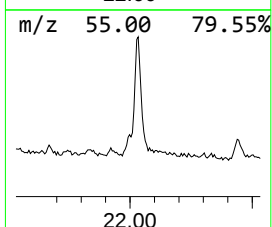
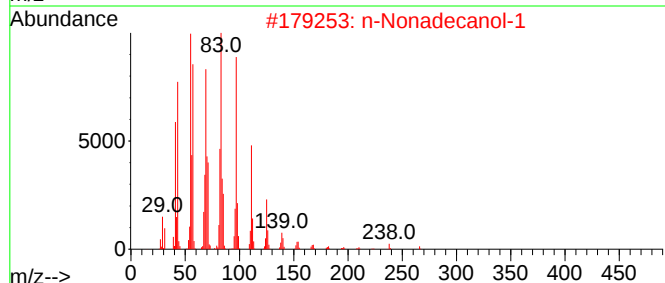
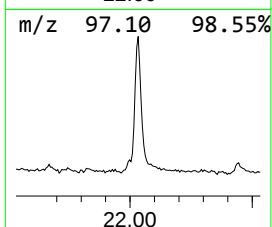
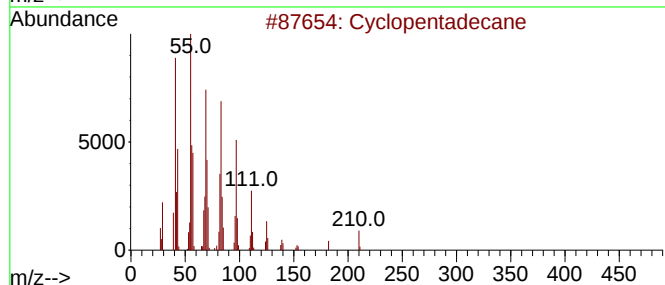
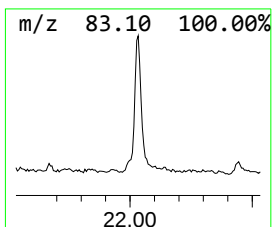
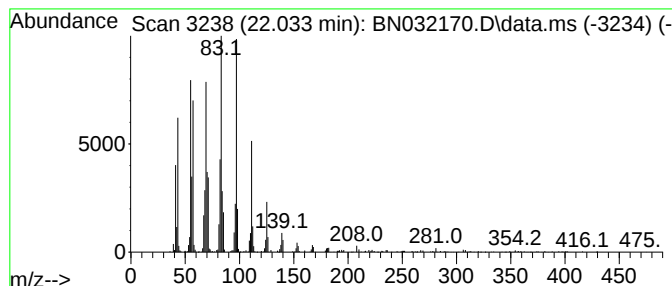
Instrument :
BNA_N
ClientSampleId :
C0AA7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Cyclic22.033 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.033	8.06 ng/ul	1484840	Chrysene-d12	21.262	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentadecane	210	C15H30	000295-48-7	93
2	n-Nonadecanol-1	284	C19H40O	001454-84-8	93
3	1-Heptacosanol	396	C27H56O	002004-39-9	91
4	1-Octadecanol	270	C18H38O	000112-92-5	91
5	1-Nonadecene	266	C19H38	018435-45-5	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
Data File : BN032170.D
Acq On : 22 Jun 2024 17:14
Operator : MA/JU
Sample : P2899-05
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0AA7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.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
unknown-01	4.769	4.1	ng/u1	431689	1	7.628	2103710	20.0
(1S)-2,6,6-Trim...	6.310	2.8	ng/u1	295996	1	7.628	2103710	20.0
Benzeneacetic acid	10.975	2.1	ng/u1	318535	2	10.410	3091410	20.0
n-Hexadecanoic ...	17.927	5.2	ng/u1	1050240	4	17.027	4077230	20.0
1H-Naphtho[2,1-...	18.545	4.7	ng/u1	964176	4	17.027	4077230	20.0
Octadecanoic acid	19.215	5.7	ng/u1	1041950	5	21.262	3683110	20.0
2-Phenanthrenol...	20.310	11.2	ng/u1	2067090	5	21.262	3683110	20.0
4,4'-Bis(tetra...	20.351	7.2	ng/u1	1324100	5	21.262	3683110	20.0
Dehydroabietic ...	20.921	7.2	ng/u1	1319840	5	21.262	3683110	20.0
1-Heneicosyl fo...	20.951	7.4	ng/u1	1368030	5	21.262	3683110	20.0
(DEL) Alkane: C...	22.033	8.1	ng/u1	1484840	5	21.262	3683110	20.0