

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082021\
 Data File : BN015994.D
 Acq On : 20 Aug 2021 20:13
 Operator : CG/JU
 Sample : M3448-14
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBKF3

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-BN081621MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN015994.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.740	108	111	122	rVB2	266638	467811	6.74%	0.595%
2	7.034	665	671	680	rVB	308630	560719	8.08%	0.713%
3	7.205	694	700	708	rBV	824900	1375107	19.82%	1.749%
4	7.405	727	734	742	rBV	958860	1609042	23.19%	2.047%
5	7.469	742	745	751	rVB2	29837	45631	0.66%	0.058%
6	7.875	807	814	821	rVB	809602	1344080	19.37%	1.710%
7	8.252	870	878	887	rBV	1982376	3359297	48.42%	4.273%
8	8.581	925	934	946	rBV	838140	1445435	20.83%	1.839%
9	8.740	957	961	966	rVB5	18499	34209	0.49%	0.044%
10	8.981	997	1002	1004	rVV5	20574	30455	0.44%	0.039%
11	9.034	1004	1011	1021	rVB	1129872	1955331	28.18%	2.487%
12	9.340	1056	1063	1068	rVB9	12237	27285	0.39%	0.035%
13	9.416	1071	1076	1081	rVV5	19457	40291	0.58%	0.051%
14	9.616	1105	1110	1118	rVB4	43478	78013	1.12%	0.099%
15	9.758	1126	1134	1145	rBV	870295	1552566	22.38%	1.975%
16	9.893	1150	1157	1159	rBV7	12607	24877	0.36%	0.032%
17	10.075	1183	1188	1194	rBV3	44496	77239	1.11%	0.098%
18	10.293	1215	1225	1234	rVV	1338464	2332207	33.62%	2.966%
19	10.369	1234	1238	1244	rVV4	36039	63206	0.91%	0.080%
20	10.452	1244	1252	1261	rVV	155611	290711	4.19%	0.370%
21	10.540	1261	1267	1272	rVV8	18066	38613	0.56%	0.049%
22	10.675	1282	1290	1298	rBV	1112107	1967507	28.36%	2.503%
23	10.734	1298	1300	1305	rVV6	25392	39301	0.57%	0.050%
24	10.810	1305	1313	1328	rVV	1078705	1990442	28.69%	2.532%
25	11.128	1359	1367	1380	rBV	132012	260145	3.75%	0.331%
26	11.287	1388	1394	1398	rBV7	23462	45937	0.66%	0.058%
27	11.340	1398	1403	1410	rVB7	20872	43364	0.63%	0.055%
28	11.463	1419	1424	1428	rBV8	12180	24970	0.36%	0.032%
29	11.552	1431	1439	1445	rVV9	28926	85968	1.24%	0.109%
30	11.787	1474	1479	1486	rBV5	41368	86562	1.25%	0.110%
31	11.863	1486	1492	1496	rBV3	77117	133395	1.92%	0.170%
32	11.928	1496	1503	1507	rVV3	114027	287094	4.14%	0.365%
33	11.969	1507	1510	1516	rVB	92012	146648	2.11%	0.187%
34	12.057	1521	1525	1531	rVB2	66821	107342	1.55%	0.137%
35	12.152	1535	1541	1546	rBV5	28278	55044	0.79%	0.070%
36	12.204	1546	1550	1551	rBV2	22233	27067	0.39%	0.034%

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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-BN081621MA.M
 Title : SVOA CALIBRATION

37	12.222	1551	1553	1558	rVB3	27009	41406	0.60%	0.053%
38	12.310	1563	1568	1573	rVB9	18314	40479	0.58%	0.051%
39	12.446	1586	1591	1597	rVB8	17853	37928	0.55%	0.048%
40	12.551	1597	1609	1615	rBV4	139501	260742	3.76%	0.332%
41	12.681	1624	1631	1634	rBV8	40966	90750	1.31%	0.115%
42	12.722	1636	1638	1643	rVB3	27864	33970	0.49%	0.043%
43	12.857	1655	1661	1667	rBV6	57312	117566	1.69%	0.150%
44	12.993	1680	1684	1688	rBV7	28183	43746	0.63%	0.056%
45	13.087	1695	1700	1704	rBV2	46888	81510	1.17%	0.104%
46	13.122	1704	1706	1712	rVB7	21031	34047	0.49%	0.043%
47	13.381	1747	1750	1754	rVV6	19533	30313	0.44%	0.039%
48	13.463	1760	1764	1769	rVV8	20960	45479	0.66%	0.058%
49	13.510	1770	1772	1779	rVV7	20822	44191	0.64%	0.056%
50	13.575	1779	1783	1788	rVB8	29697	55827	0.80%	0.071%
51	13.693	1798	1803	1809	rBV10	32601	56648	0.82%	0.072%
52	13.828	1821	1826	1827	rBV4	19706	30660	0.44%	0.039%
53	13.922	1836	1842	1848	rVB	2543054	3643938	52.52%	4.635%
54	13.993	1849	1854	1863	rVB6	106089	257699	3.71%	0.328%
55	14.110	1869	1874	1879	rVB3	58820	112928	1.63%	0.144%
56	14.204	1879	1890	1902	rBV	2653290	4251032	61.27%	5.407%
57	14.328	1907	1911	1912	rBV4	34357	40013	0.58%	0.051%
58	14.363	1913	1917	1922	rVB	103566	162387	2.34%	0.207%
59	14.422	1922	1927	1931	rBV7	37393	71780	1.03%	0.091%
60	14.510	1934	1942	1948	rBV	1653691	2720174	39.21%	3.460%
61	14.575	1948	1953	1959	rVB	267312	433577	6.25%	0.551%
62	14.704	1970	1975	1982	rBV	311615	467347	6.74%	0.594%
63	14.828	1993	1996	1998	rBV4	25194	31064	0.45%	0.040%
64	14.910	2006	2010	2016	rVB8	39340	75922	1.09%	0.097%
65	14.987	2018	2023	2025	rBV2	76678	101514	1.46%	0.129%
66	15.028	2025	2030	2034	rVV	219695	379957	5.48%	0.483%
67	15.081	2034	2039	2043	rVV2	168519	349414	5.04%	0.444%
68	15.122	2043	2046	2054	rVB3	124199	248345	3.58%	0.316%
69	15.328	2076	2081	2084	rVB5	124960	181165	2.61%	0.230%
70	15.451	2099	2102	2104	rBV4	25311	34068	0.49%	0.043%
71	15.504	2104	2111	2117	rVV	3452702	5185951	74.75%	6.596%
72	15.557	2117	2120	2125	rVB	100362	137181	1.98%	0.174%
73	15.616	2125	2130	2138	rBV	1189522	1752796	25.26%	2.229%
74	15.687	2138	2142	2148	rVB2	172392	250827	3.62%	0.319%
75	15.751	2149	2153	2160	rVB3	93358	178029	2.57%	0.226%
76	15.851	2160	2170	2177	rBV4	100303	303404	4.37%	0.386%

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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-BN081621MA.M
 Title : SVOA CALIBRATION

77	15.992	2190	2194	2198	rVB4	36807	48230	0.70%	0.061%
78	16.234	2231	2235	2242	rVB3	110767	185177	2.67%	0.236%
79	16.557	2286	2290	2292	rBV5	35517	62035	0.89%	0.079%
80	16.587	2293	2295	2300	rVB3	55144	80084	1.15%	0.102%
81	16.787	2324	2329	2333	rBV3	130670	229375	3.31%	0.292%
82	16.898	2344	2348	2353	rVB3	126983	231710	3.34%	0.295%
83	16.963	2356	2359	2367	rVB9	63689	101703	1.47%	0.129%
84	17.069	2374	2377	2386	rVB3	97761	161852	2.33%	0.206%
85	17.257	2403	2409	2415	rBV	1886555	2852229	41.11%	3.628%
86	17.357	2420	2426	2438	rVB	4123300	6101113	87.94%	7.760%
87	18.157	2559	2562	2568	rBV8	44827	76854	1.11%	0.098%
88	18.228	2572	2574	2585	rVV4	64452	145825	2.10%	0.185%
89	18.316	2585	2589	2599	rVB4	95413	213694	3.08%	0.272%
90	18.675	2646	2650	2657	rVB4	88240	152106	2.19%	0.193%
91	19.057	2709	2715	2720	rBV	777276	1117274	16.10%	1.421%
92	19.216	2739	2742	2745	rBV4	58561	81372	1.17%	0.104%
93	19.651	2810	2816	2828	rBV	4885838	6845972	98.68%	8.708%
94	20.651	2983	2986	2992	rVB	624157	744988	10.74%	0.948%
95	21.157	3069	3072	3080	rVB3	194578	252455	3.64%	0.321%
96	21.445	3116	3121	3131	rVB	2446826	3185033	45.91%	4.051%
97	22.116	3231	3235	3242	rVB	513071	743772	10.72%	0.946%
98	22.563	3307	3311	3319	rVB3	212690	353158	5.09%	0.449%
99	23.692	3494	3503	3509	rBV	3292513	6937830	100.00%	8.825%
100	23.845	3521	3529	3538	rVB2	1511835	3246539	46.79%	4.129%

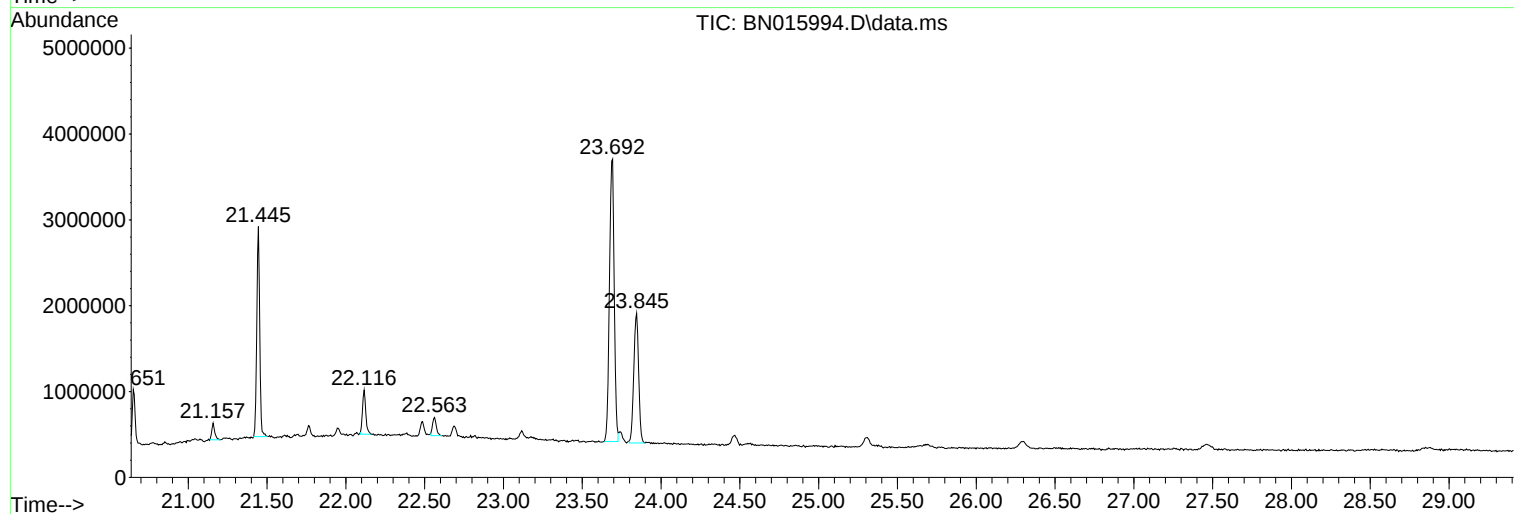
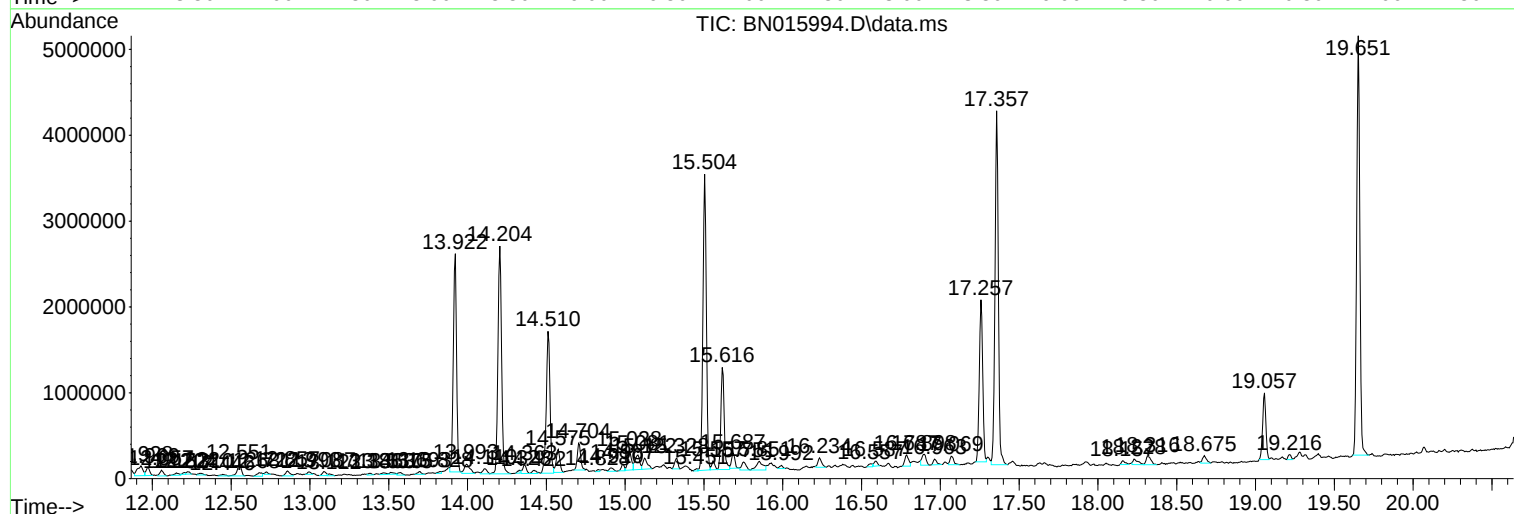
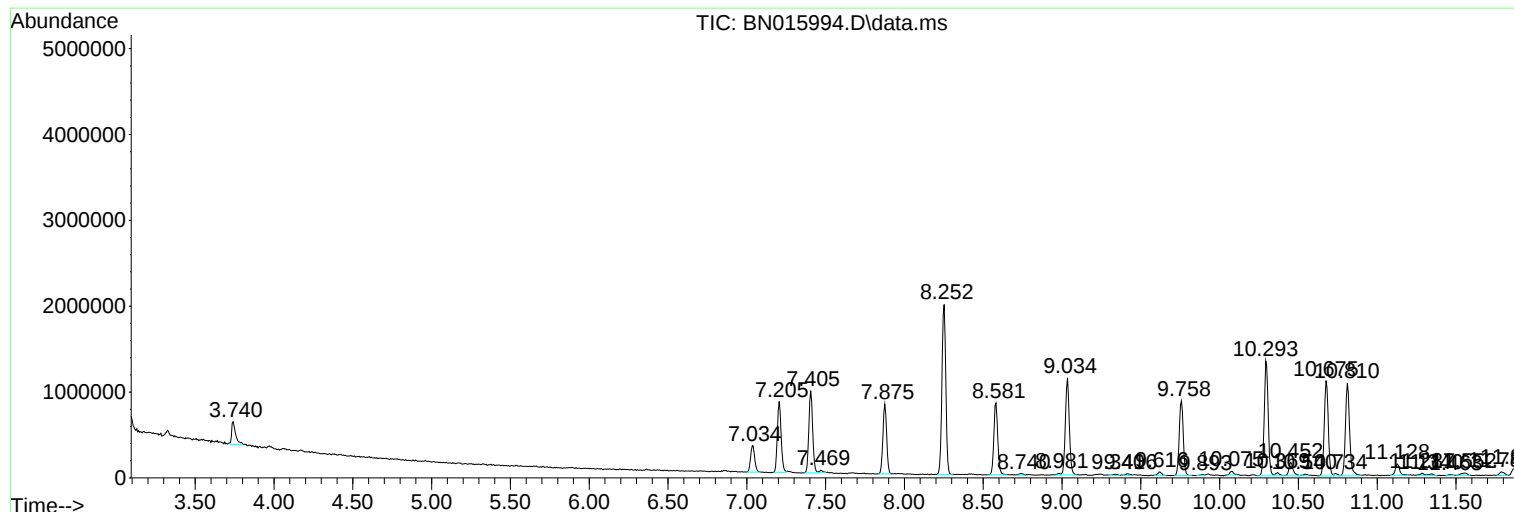
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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



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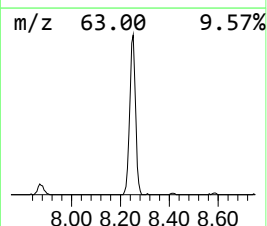
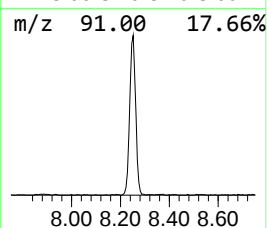
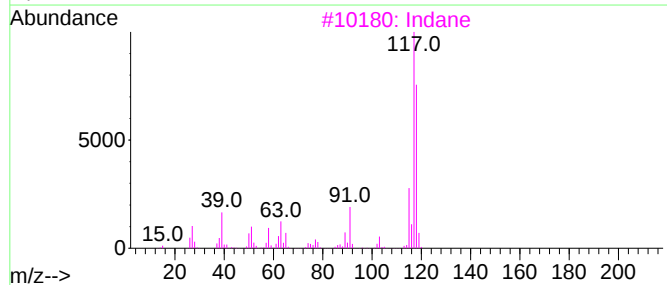
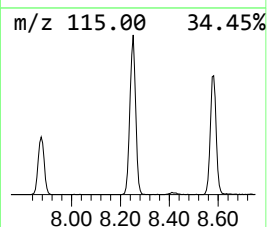
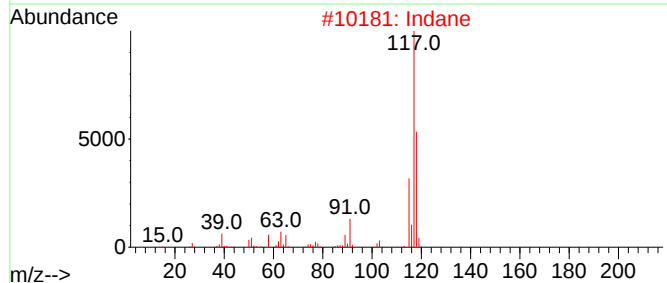
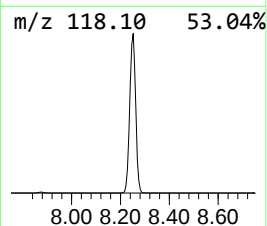
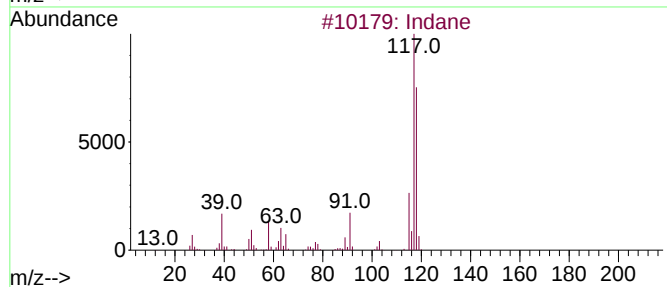
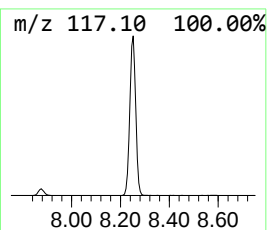
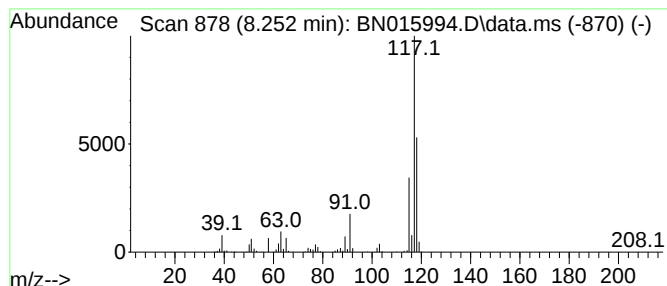
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.252	49.99 ng/ul	3359300	1,4-Dichlorobenzene-d4	7.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	93
2	Indane			118	C9H10	000496-11-7	90
3	Indane			118	C9H10	000496-11-7	90
4	Benzene, 1-ethenyl-2-methyl-			118	C9H10	000611-15-4	60
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...			118	C9H10	1000191-13-7	59



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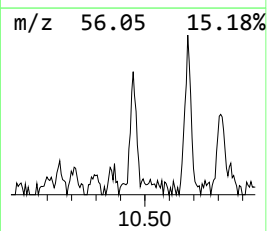
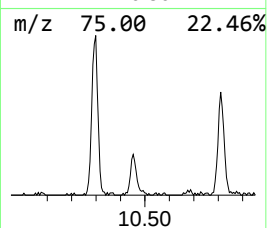
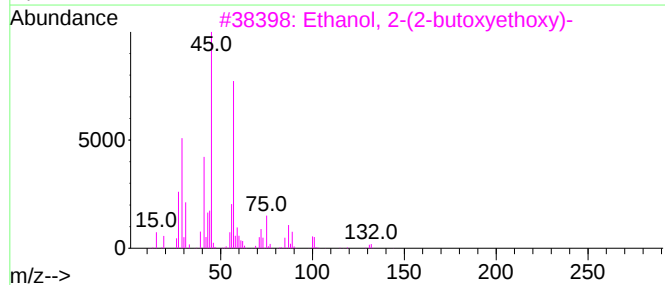
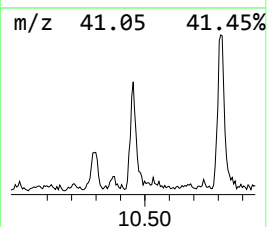
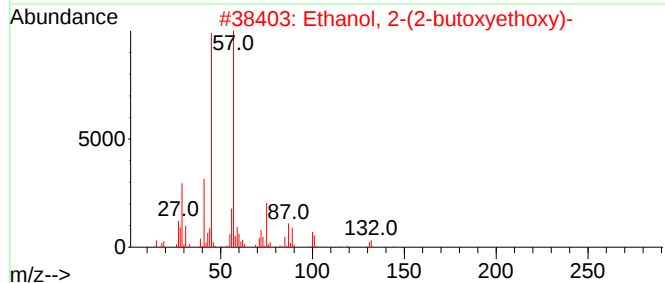
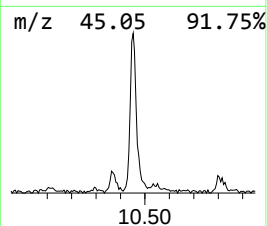
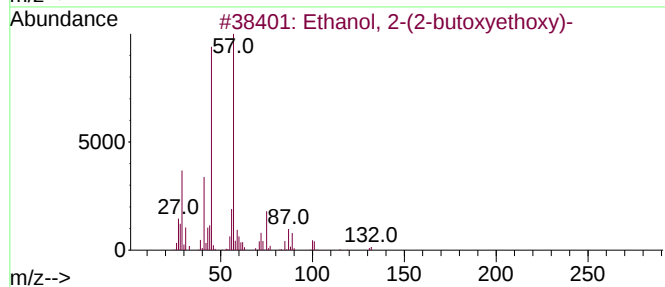
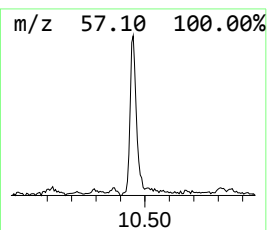
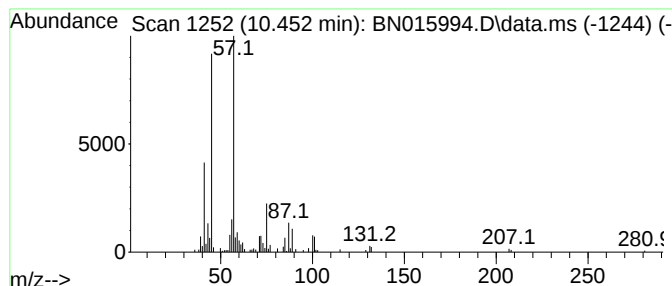
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.452	2.96 ng/ul	290711	Naphthalene-d8	10.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
4			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
5			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64



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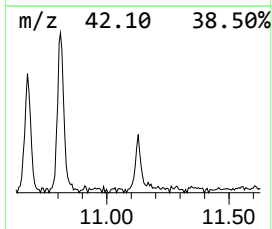
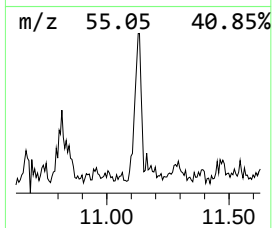
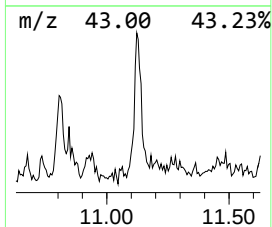
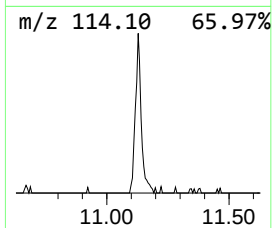
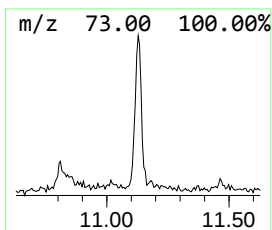
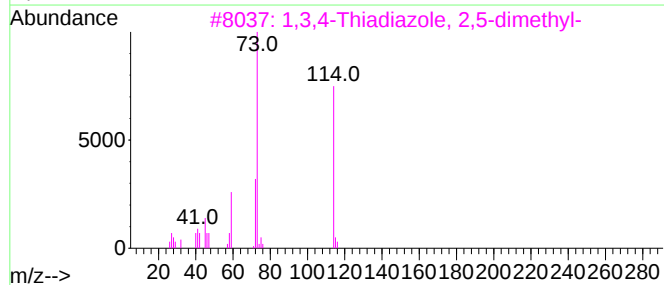
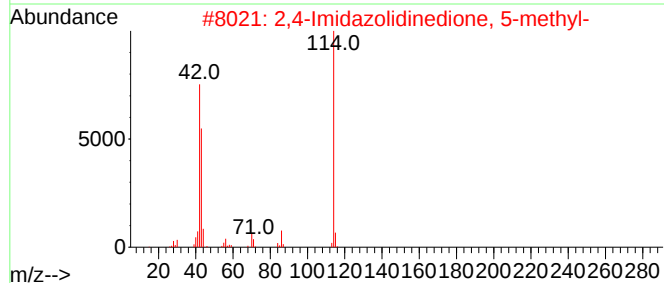
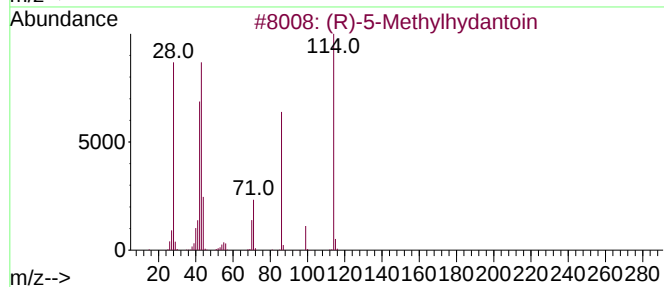
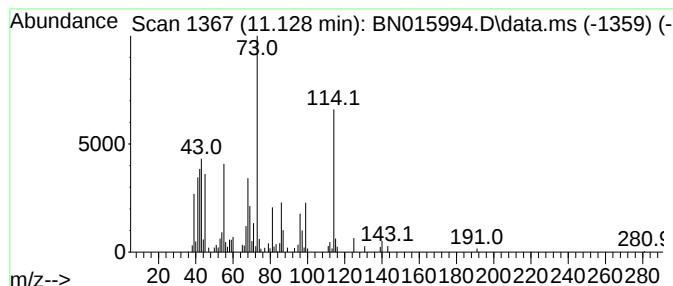
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.128	2.64 ng/ul	260145	Naphthalene-d8	10.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(R)-5-Methylhydantoin			114	C4H6N2O2	055147-68-7	38
2	2,4-Imidazolidinedione, 5-methyl-			114	C4H6N2O2	000616-03-5	27
3	1,3,4-Thiadiazole, 2,5-dimethyl-			114	C4H6N2S	027464-82-0	25
4	2-Methoxythiophene			114	C5H6OS	016839-97-7	22
5	2-Methoxythiophene			114	C5H6OS	016839-97-7	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082021\
Data File : BN015994.D
Acq On : 20 Aug 2021 20:13
Operator : CG/JU
Sample : M3448-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBKF3

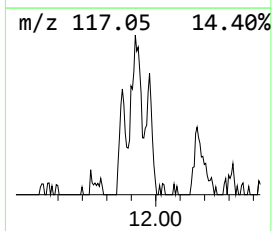
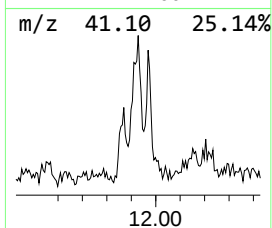
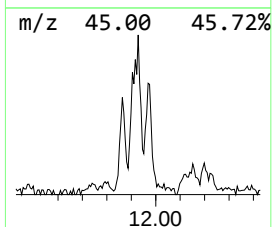
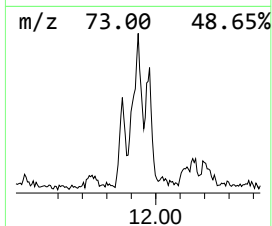
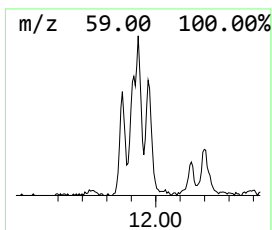
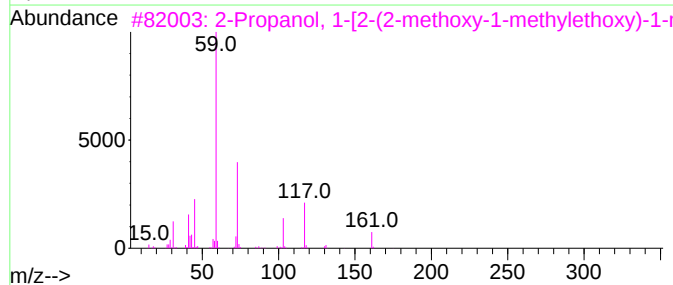
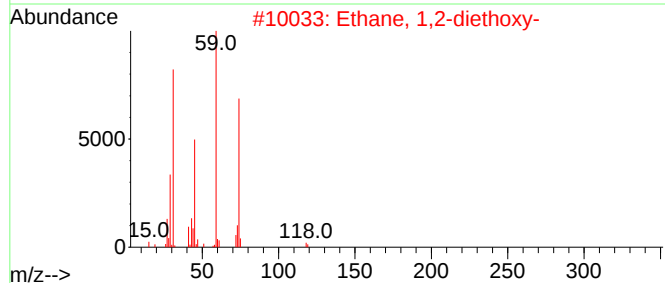
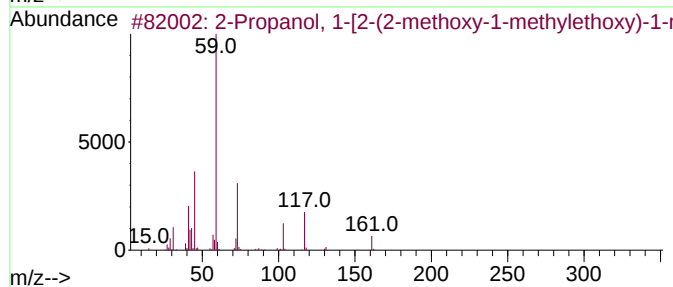
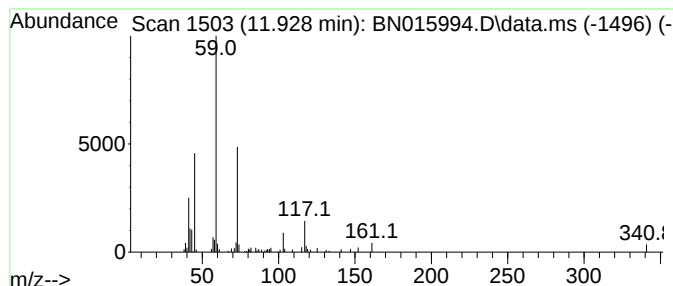
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2-Propanol, 1-[2-(2-methoxy... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.928	2.92 ng/ul	287094	Naphthalene-d8	10.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	78		
2	Ethane, 1,2-diethoxy-	118	C6H14O2	000629-14-1	72		
3	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	64		
4	2-Butanol, 1-methoxy-	104	C5H12O2	053778-73-7	59		
5	Ethanedioic acid, dimethyl ester	118	C4H6O4	000553-90-2	59		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082021\
Data File : BN015994.D
Acq On : 20 Aug 2021 20:13
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Sample : M3448-14
Misc :
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Instrument :
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ClientSampleId :
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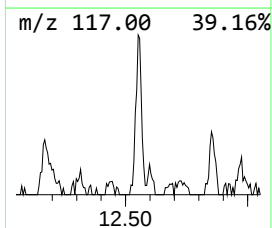
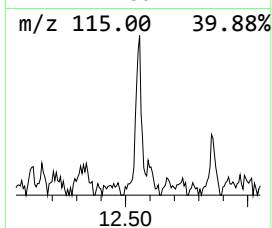
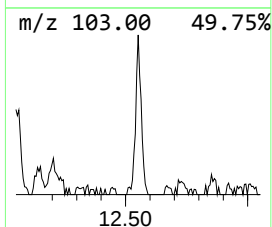
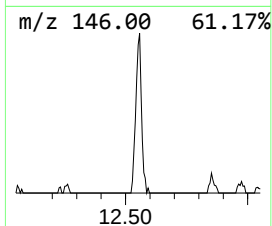
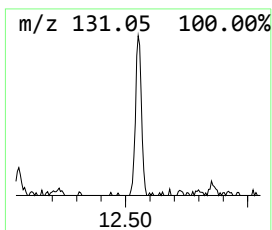
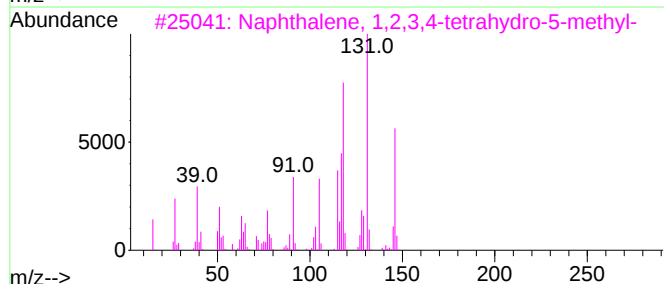
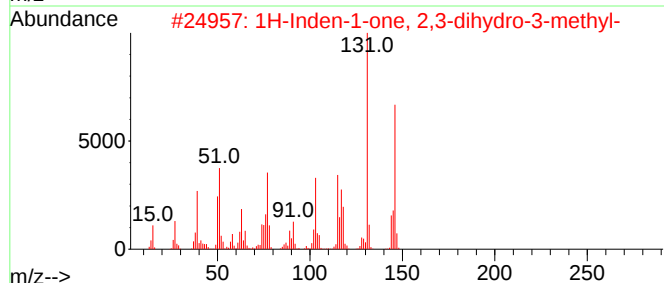
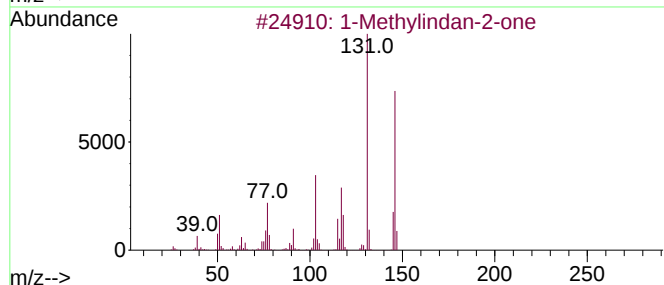
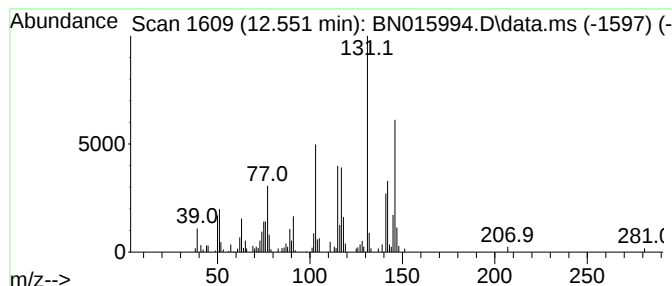
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1-Methylindan-2-one Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.552	2.65 ng/ul	260742	Naphthalene-d8	10.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Methylindan-2-one	146	C10H10O	035587-60-1	95
2			1H-Inden-1-one, 2,3-dihydro-3-me...	146	C10H10O	006072-57-7	86
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	76
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	70
5			1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082021\
Data File : BN015994.D
Acq On : 20 Aug 2021 20:13
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Sample : M3448-14
Misc :
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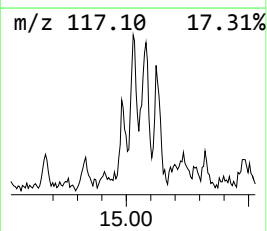
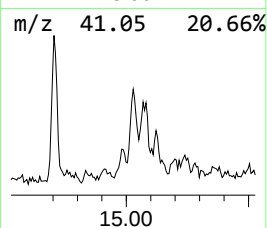
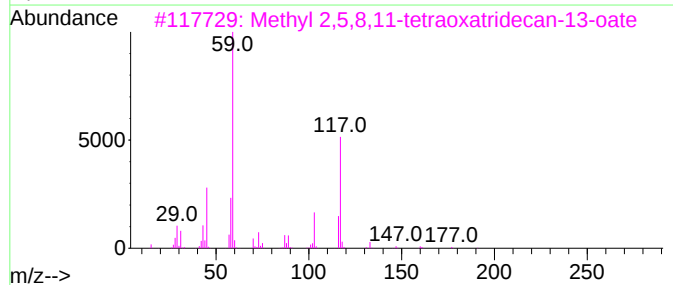
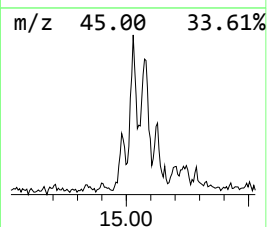
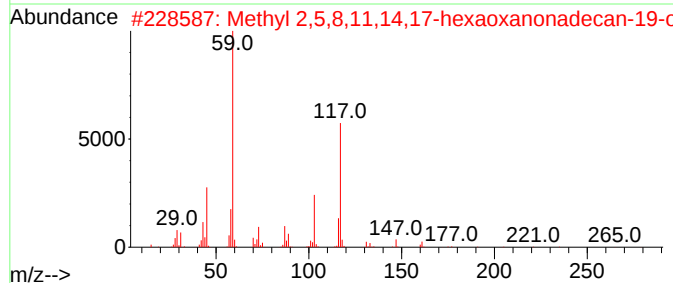
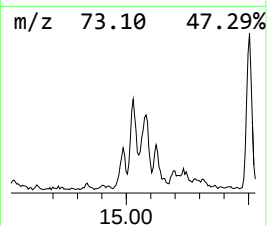
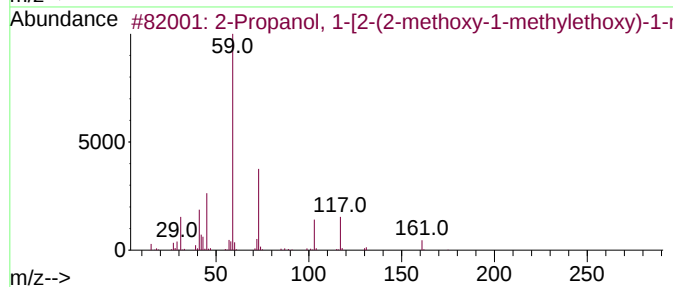
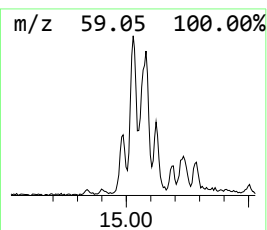
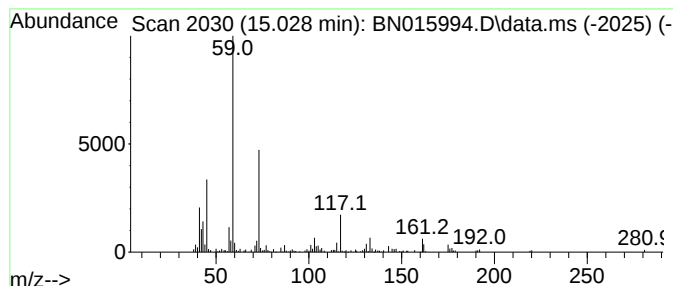
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Methyl 2,5,8,11,14,17-hexao... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.028	2.79 ng/ul	379957	Acenaphthene-d10	14.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	72		
2	Methyl 2,5,8,11,14,17-hexaoxon...	324	C14H28O8	1000366-81-4	64		
3	Methyl 2,5,8,11-tetraoxatridecan...	236	C10H20O6	1000366-78-2	64		
4	2,5,8,11-Tetraoxatetradecan-13-o...	264	C13H28O5	020324-34-9	59		
5	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	59		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082021\
Data File : BN015994.D
Acq On : 20 Aug 2021 20:13
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Sample : M3448-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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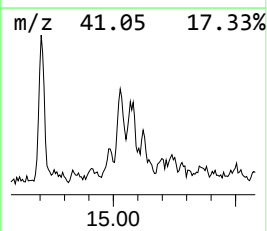
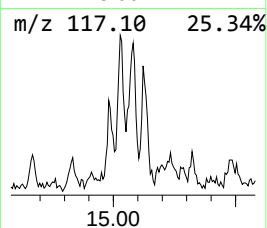
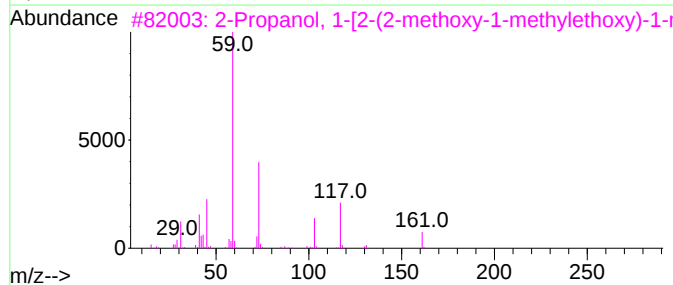
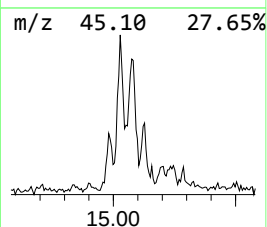
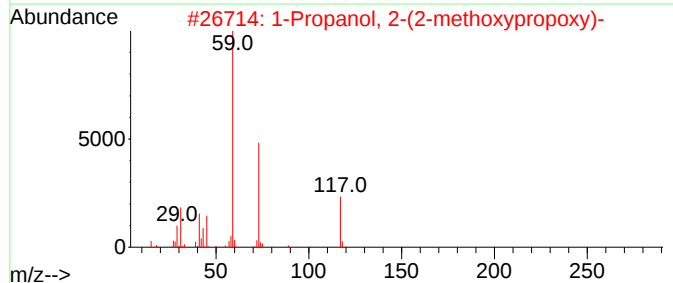
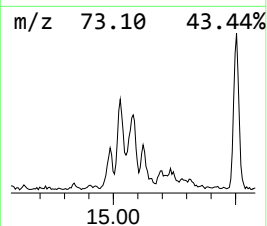
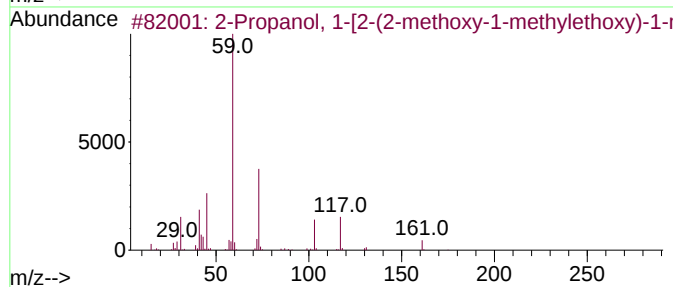
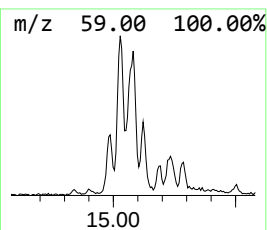
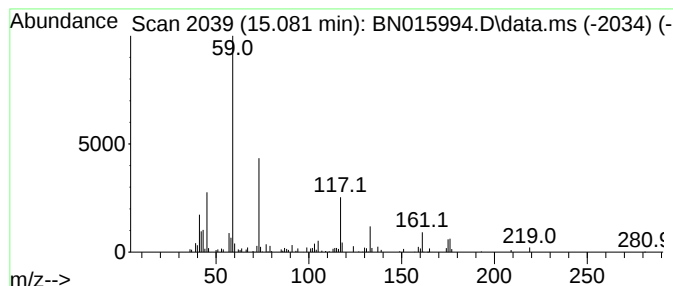
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1-Propanol, 2-(2-methoxypro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.081	2.57 ng/ul	349414	Acenaphthene-d10	14.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	59		
2	1-Propanol, 2-(2-methoxypropoxy)-	148	C7H16O3	013588-28-8	59		
3	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	56		
4	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	50		
5	Butane, 2-methoxy-	88	C5H12O	006795-87-5	47		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082021\
Data File : BN015994.D
Acq On : 20 Aug 2021 20:13
Operator : CG/JU
Sample : M3448-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

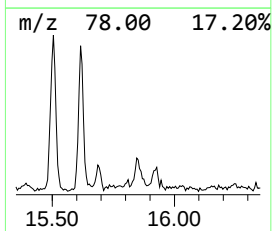
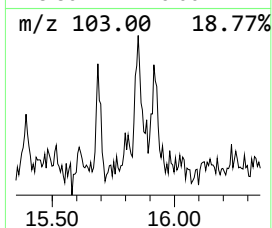
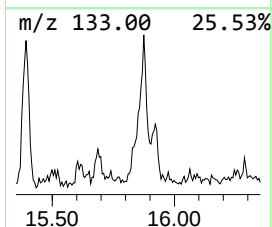
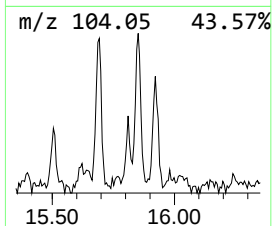
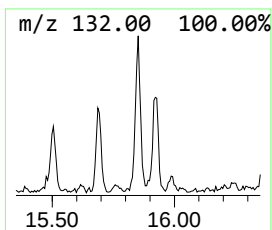
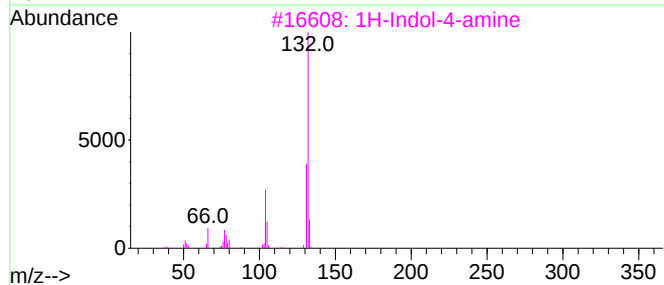
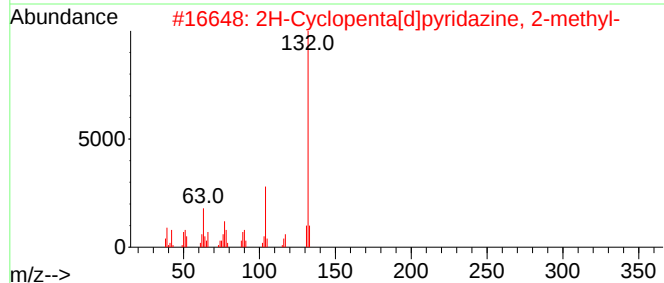
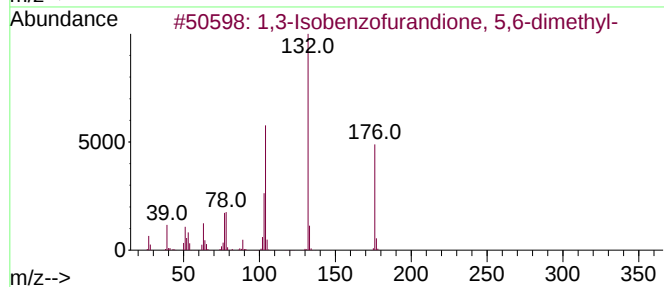
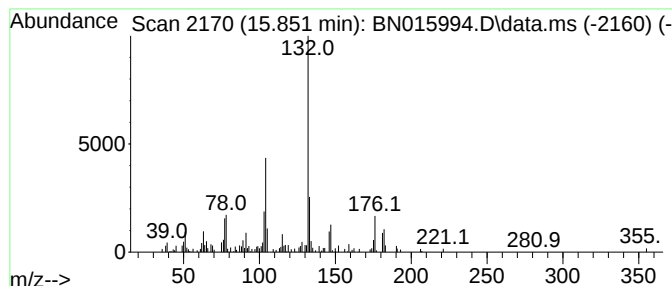
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1,3-Isobenzofurandione, 5,6... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.851	2.23 ng/ul	303404	Acenaphthene-d10	14.510	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Isobenzofurandione, 5,6-dime...	176	C10H8O3	005999-20-2	59
2	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	53
3	1H-Indol-4-amine	132	C8H8N2	005192-23-4	53
4	4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	52
5	(3-Aminophenyl)acetonitrile	132	C8H8N2	004623-24-9	50



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Misc :
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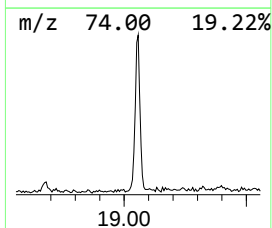
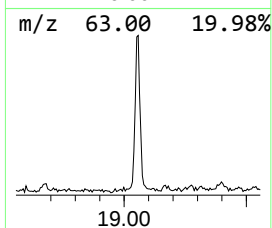
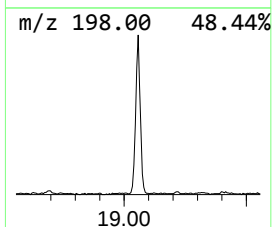
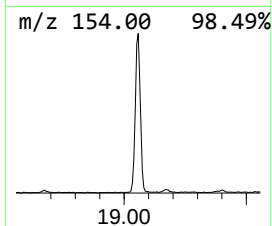
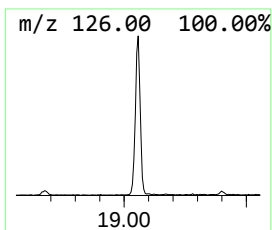
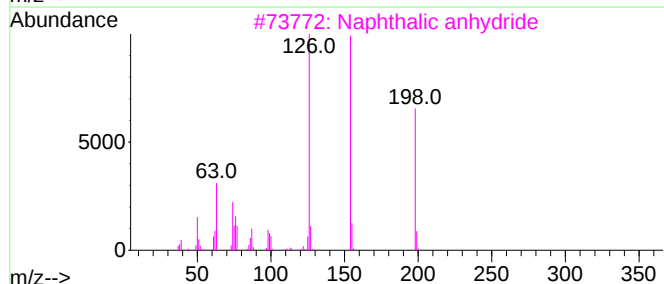
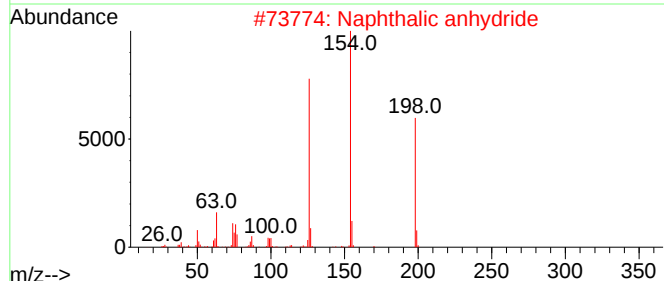
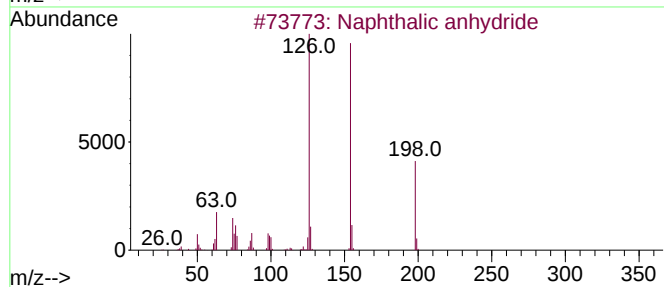
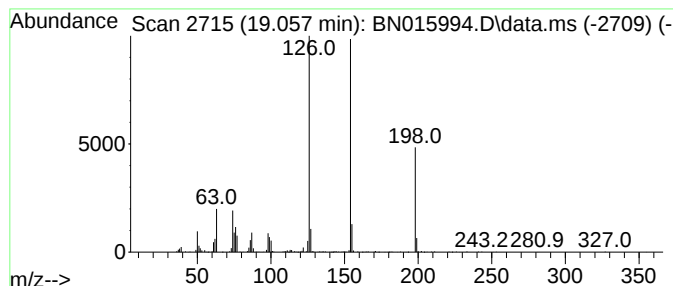
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Naphthalic anhydride Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.057	7.83 ng/ul	1117270	Phenanthrene-d10	17.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalic anhydride	198	C12H6O3	000081-84-5	99
2			Naphthalic anhydride	198	C12H6O3	000081-84-5	98
3			Naphthalic anhydride	198	C12H6O3	000081-84-5	98
4			Naphthalic anhydride	198	C12H6O3	000081-84-5	97
5			Naphthalic anhydride	198	C12H6O3	000081-84-5	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082021\
Data File : BN015994.D
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Misc :
ALS Vial : 8 Sample Multiplier: 1

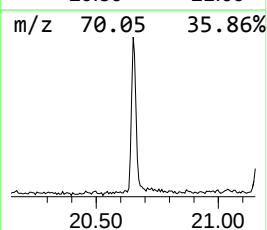
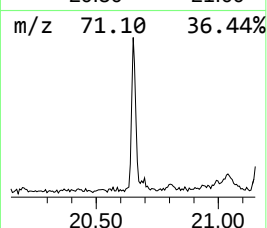
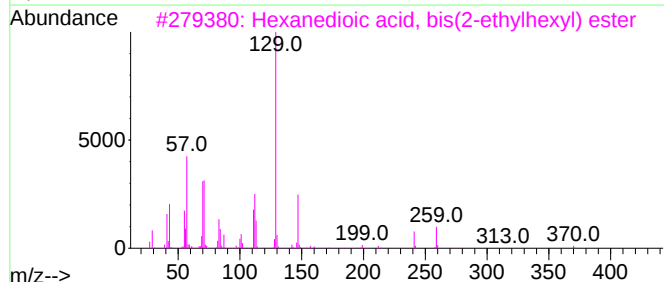
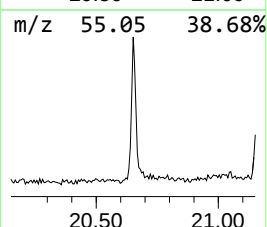
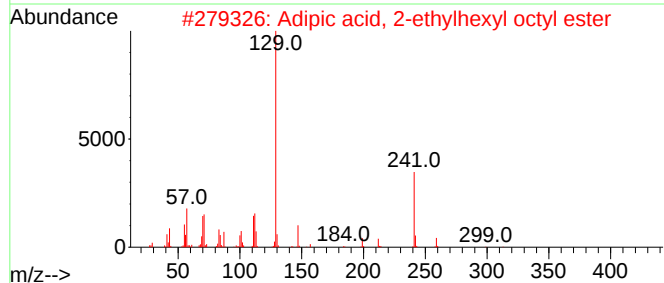
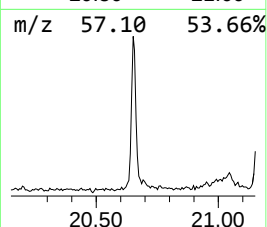
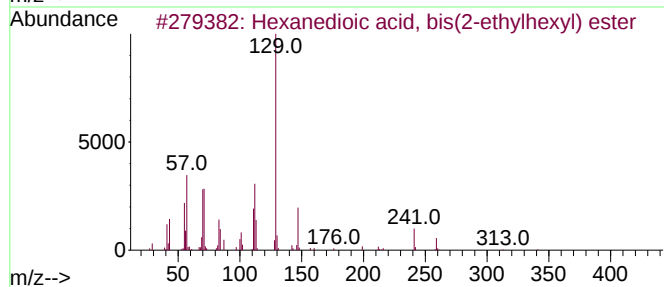
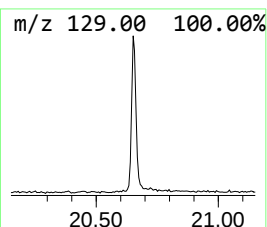
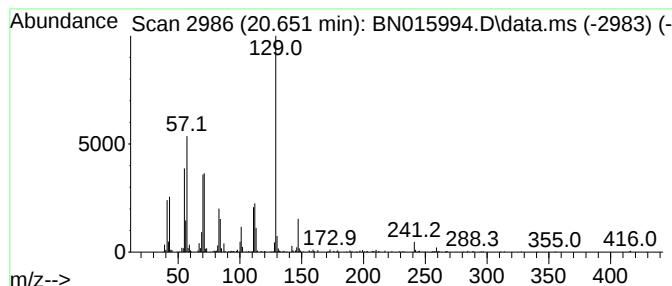
Instrument :
BNA_N
ClientSampleId :
DBKF3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 Hexanedioic acid, bis(2-eth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.651	4.68 ng/ul	744988	Chrysene-d12	21.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	91
2	Adipic acid, 2-ethylhexyl octyl ...	370	C22H42O4	1000324-48-6	90
3	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	78
4	Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	72
5	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082021\
Data File : BN015994.D
Acq On : 20 Aug 2021 20:13
Operator : CG/JU
Sample : M3448-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBKF3

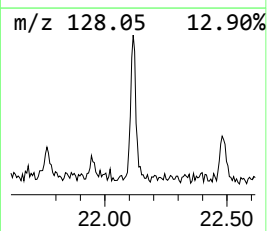
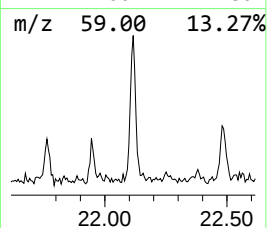
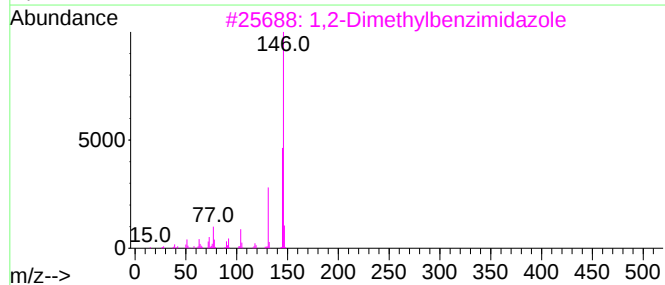
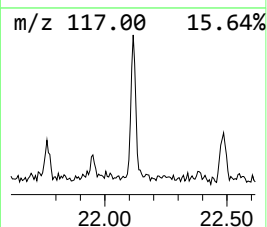
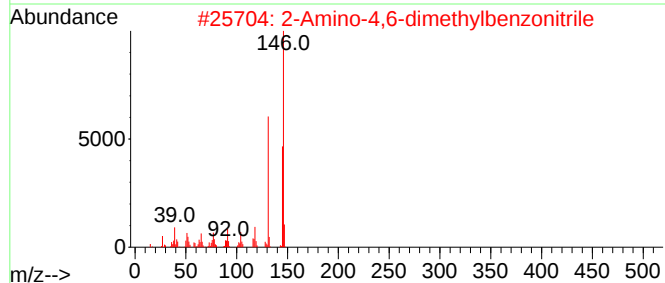
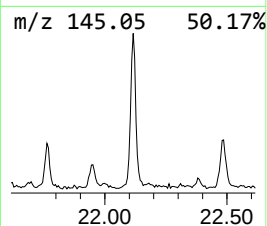
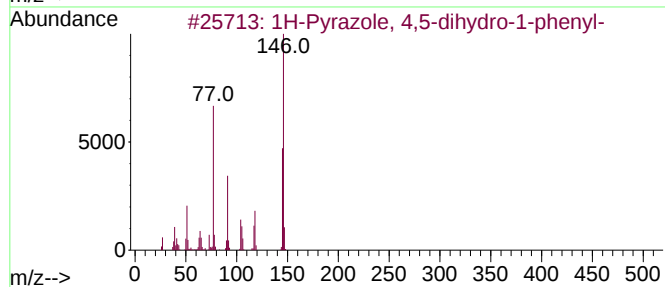
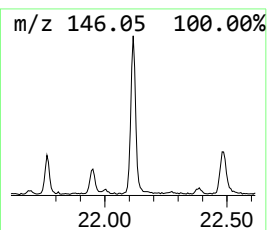
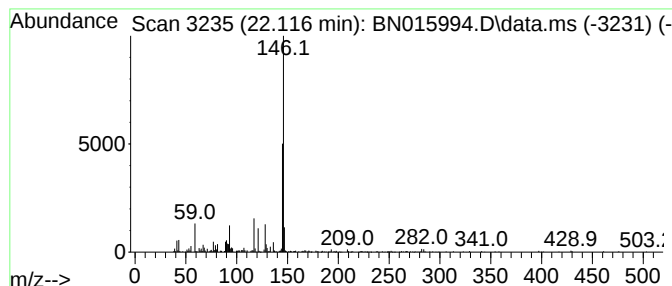
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 1H-Pyrazole, 4,5-dihydro-1-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.116	4.67 ng/ul	743772	Chrysene-d12	21.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Pyrazole, 4,5-dihydro-1-phenyl-	146	C9H10N2	000936-53-8	59
2			2-Amino-4,6-dimethylbenzonitrile	146	C9H10N2	1010461-05-1	59
3			1,2-Dimethylbenzimidazole	146	C9H10N2	002876-08-6	53
4			Acetonitrile, 2-(4-methoxyphenyl...	291	C18H17N3O	1000277-32-9	50
5			1H-Pyrazole, 4,5-dihydro-3-phenyl-	146	C9H10N2	000936-48-1	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082021\
Data File : BN015994.D
Acq On : 20 Aug 2021 20:13
Operator : CG/JU
Sample : M3448-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBKF3

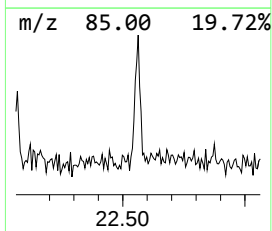
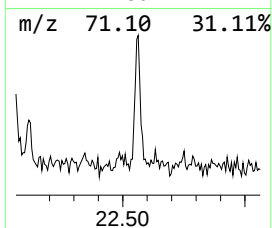
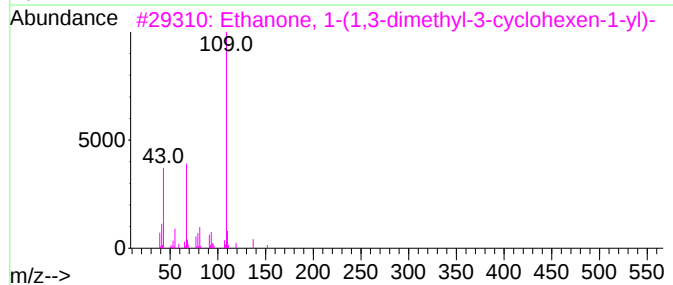
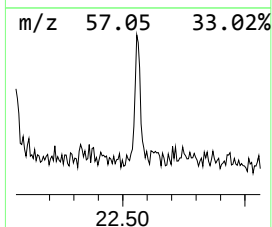
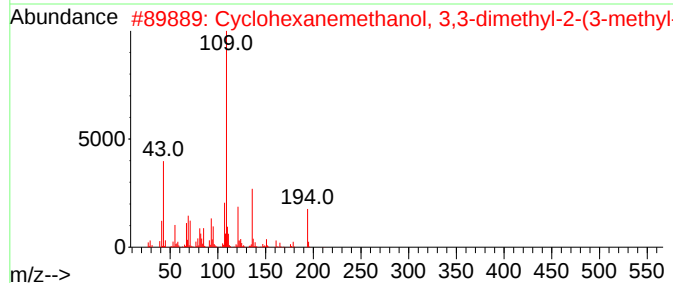
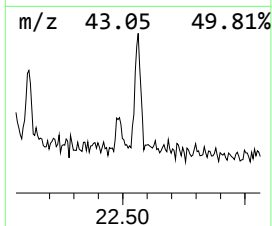
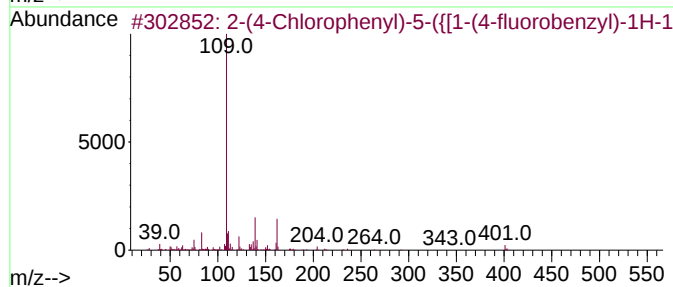
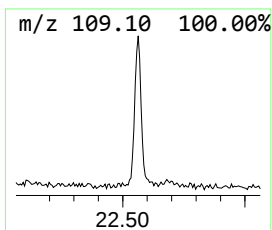
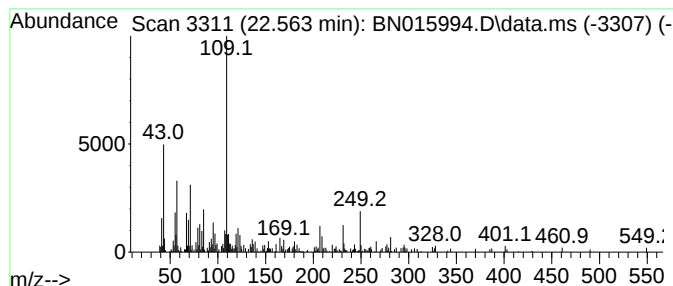
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 2-(4-Chlorophenyl)-5-([1-(... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.563	2.22 ng/ul	353158	Chrysene-d12	21.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(4-Chlorophenyl)-5-([1-(4-flu...	401	C18H13ClFN5OS	1000482-39-7	58		
2	Cyclohexanemethanol, 3,3-dimethy...	212	C13H24O2	072541-28-7	50		
3	Ethanone, 1-(1,3-dimethyl-3-cycl...	152	C10H16O	051733-68-7	49		
4	Disulfide, 3-fluorobenzyl 4-brom...	328	C13H10BrFS2	1000470-64-7	49		
5	6,6-Dimethylhepta-2,4-diene	124	C9H16	1000195-03-3	49		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082021\
Data File : BN015994.D
Acq On : 20 Aug 2021 20:13
Operator : CG/JU
Sample : M3448-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBKF3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.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
Indane	8.252	50.0	ng/ul	3359300	1	7.875	1344080	20.0
Ethanol, 2-(2-b...	10.452	3.0	ng/ul	290711	2	10.675	1967510	20.0
unknown-01	11.128	2.6	ng/ul	260145	2	10.675	1967510	20.0
2-Propanol, 1-[...	11.928	2.9	ng/ul	287094	2	10.675	1967510	20.0
1-Methylindan-2...	12.552	2.6	ng/ul	260742	2	10.675	1967510	20.0
Methyl 2,5,8,11...	15.028	2.8	ng/ul	379957	3	14.510	2720170	20.0
1-Propanol, 2-(...	15.081	2.6	ng/ul	349414	3	14.510	2720170	20.0
1,3-Isobenzofur...	15.851	2.2	ng/ul	303404	3	14.510	2720170	20.0
Naphthalic anhy...	19.057	7.8	ng/ul	1117270	4	17.257	2852230	20.0
Hexanedioic aci...	20.651	4.7	ng/ul	744988	5	21.445	3185030	20.0
1H-Pyrazole, 4,...	22.116	4.7	ng/ul	743772	5	21.445	3185030	20.0
2-(4-Chlorophen...	22.563	2.2	ng/ul	353158	5	21.445	3185030	20.0