

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062121\
 Data File : BM030507.D
 Acq On : 21 Jun 2021 16:42
 Operator : CG/JU
 Sample : M2697-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JLX82

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_M\METHODS\SFAM-EPA-BM060621.M
 Title : SVOA CALIBRATION

Signal : TIC: BM030507.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.746	107	112	113	rBV3	11655	18021	0.85%	0.061%
2	3.769	113	116	124	rVV	59485	156216	7.34%	0.530%
3	3.840	124	128	138	rVB	46820	105943	4.98%	0.360%
4	3.963	144	149	155	rVV2	22760	47547	2.23%	0.161%
5	4.051	158	164	176	rVB	61737	151904	7.14%	0.516%
6	4.216	188	192	196	rBV5	7972	13547	0.64%	0.046%
7	4.269	196	201	215	rVB	105215	212129	9.97%	0.720%
8	4.422	221	227	233	rVB2	27858	51175	2.41%	0.174%
9	4.551	244	249	252	rBV	22642	42044	1.98%	0.143%
10	4.598	252	257	264	rVV	94796	179151	8.42%	0.608%
11	4.675	264	270	277	rVB	75561	132500	6.23%	0.450%
12	4.957	312	318	329	rVB	35343	70354	3.31%	0.239%
13	5.057	331	335	343	rBV	13145	28738	1.35%	0.098%
14	5.357	379	386	394	rVB	54299	102509	4.82%	0.348%
15	5.487	401	408	420	rBV	165006	397909	18.70%	1.351%
16	5.851	462	470	480	rVB	83190	155754	7.32%	0.529%
17	6.222	528	533	539	rVB2	51200	84640	3.98%	0.287%
18	6.822	630	635	642	rBV4	8336	16627	0.78%	0.056%
19	6.992	654	664	678	rBV	374846	741133	34.83%	2.516%
20	7.157	686	692	703	rVB	291577	483994	22.75%	1.643%
21	7.251	703	708	716	rVV7	8319	16822	0.79%	0.057%
22	7.357	719	726	736	rVV	386331	661620	31.09%	2.246%
23	7.645	773	775	784	rVB8	4319	9440	0.44%	0.032%
24	7.822	798	805	812	rVV	652604	1087753	51.12%	3.693%
25	7.887	812	816	820	rVV3	6688	11292	0.53%	0.038%
26	7.945	823	826	831	rVB5	5645	8883	0.42%	0.030%
27	8.051	837	844	852	rBV	14511	29975	1.41%	0.102%
28	8.357	891	896	902	rBV2	10637	20282	0.95%	0.069%
29	8.522	914	924	939	rBV	337456	594541	27.94%	2.019%
30	8.645	939	945	951	rVV	106513	180932	8.50%	0.614%
31	8.916	985	991	995	rVV	23293	39197	1.84%	0.133%
32	8.975	995	1001	1011	rVV	341928	580959	27.30%	1.972%
33	9.081	1014	1019	1025	rVV2	10241	20901	0.98%	0.071%
34	9.139	1025	1029	1038	rVB8	5784	12543	0.59%	0.043%
35	9.386	1067	1071	1076	rVB2	10876	18558	0.87%	0.063%
36	9.692	1115	1123	1131	rBV	258342	446574	20.99%	1.516%

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Integrator: RTE

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Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
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37	9.869	1147	1153	1154	rBV5	7667	11147	0.52%	0.038%
38	10.051	1174	1184	1188	rBV3	17007	28201	1.33%	0.096%
39	10.233	1208	1215	1232	rBV	392010	698000	32.80%	2.370%
40	10.404	1237	1244	1259	rBV	31063	88328	4.15%	0.300%
41	10.610	1270	1279	1292	rVB	795723	1390582	65.35%	4.721%
42	10.745	1294	1302	1313	rBV	246517	460451	21.64%	1.563%
43	11.151	1365	1371	1376	rVV6	7032	12182	0.57%	0.041%
44	11.410	1409	1415	1421	rBV2	11111	25959	1.22%	0.088%
45	11.633	1448	1453	1457	rBV6	6789	13929	0.65%	0.047%
46	11.710	1460	1466	1467	rBV4	7156	11670	0.55%	0.040%
47	11.969	1506	1510	1516	rBV3	4527	8894	0.42%	0.030%
48	12.392	1577	1582	1588	rBV9	4008	8512	0.40%	0.029%
49	12.486	1592	1598	1603	rBV5	9014	13754	0.65%	0.047%
50	12.986	1677	1683	1687	rBV2	13236	21005	0.99%	0.071%
51	13.057	1690	1695	1705	rVB7	5876	12940	0.61%	0.044%
52	13.269	1724	1731	1738	rVV2	28131	47044	2.21%	0.160%
53	13.351	1738	1745	1749	rVV10	3262	8669	0.41%	0.029%
54	13.774	1812	1817	1822	rBV5	13851	29325	1.38%	0.100%
55	13.857	1822	1831	1840	rVV	679448	1022042	48.03%	3.470%
56	13.927	1840	1843	1847	rVB	22368	30127	1.42%	0.102%
57	14.139	1868	1879	1887	rBV	758194	1155207	54.29%	3.922%
58	14.345	1907	1914	1919	rBV	34548	48405	2.27%	0.164%
59	14.445	1922	1931	1939	rVV	1159060	1674574	78.70%	5.685%
60	14.510	1939	1942	1946	rVV3	9916	14525	0.68%	0.049%
61	14.563	1947	1951	1956	rVB5	8838	13112	0.62%	0.045%
62	14.651	1960	1966	1986	rVV	121058	281299	13.22%	0.955%
63	14.810	1988	1993	1996	rVV5	13736	25482	1.20%	0.087%
64	14.868	1996	2003	2009	rVV5	55660	103106	4.85%	0.350%
65	14.921	2009	2012	2016	rVB5	6916	10023	0.47%	0.034%
66	14.968	2016	2020	2026	rVB5	7534	12492	0.59%	0.042%
67	15.292	2072	2075	2081	rVB2	20314	27661	1.30%	0.094%
68	15.439	2093	2100	2106	rBV	964967	1430631	67.24%	4.857%
69	15.486	2106	2108	2114	rVB5	10703	15753	0.74%	0.053%
70	15.557	2114	2120	2125	rBV	150589	253410	11.91%	0.860%
71	15.698	2136	2144	2151	rVB3	24831	52388	2.46%	0.178%
72	15.857	2165	2171	2177	rBV	265455	332171	15.61%	1.128%
73	15.910	2177	2180	2185	rVB6	6553	9291	0.44%	0.032%
74	16.180	2218	2226	2230	rVV	38970	75006	3.53%	0.255%
75	16.498	2273	2280	2283	rBV7	10481	21338	1.00%	0.072%
76	16.557	2283	2290	2294	rVV8	6462	18535	0.87%	0.063%

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 Sampling : 1 Min Area: 0.1 % of largest Peak
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77	16.668	2307	2309	2316	rVB8	6453	9461	0.44%	0.032%
78	16.727	2316	2319	2323	rBV2	16740	24438	1.15%	0.083%
79	16.945	2350	2356	2359	rBV	32615	52212	2.45%	0.177%
80	16.998	2361	2365	2369	rVB2	34497	52308	2.46%	0.178%
81	17.186	2390	2397	2402	rBV	1276802	1830335	86.02%	6.214%
82	17.227	2402	2404	2408	rVV	47409	59765	2.81%	0.203%
83	17.280	2408	2413	2423	rVV	984222	1472279	69.19%	4.999%
84	17.374	2426	2429	2433	rVB4	29428	39231	1.84%	0.133%
85	17.539	2454	2457	2461	rBV4	11993	17923	0.84%	0.061%
86	17.592	2464	2466	2473	rVB7	14123	22472	1.06%	0.076%
87	17.680	2477	2481	2485	rBV	49038	60152	2.83%	0.204%
88	17.851	2507	2510	2517	rVB5	32122	52534	2.47%	0.178%
89	18.074	2544	2548	2551	rBV3	55883	78360	3.68%	0.266%
90	18.151	2558	2561	2564	rVB2	25369	31768	1.49%	0.108%
91	18.368	2595	2598	2604	rBV	46512	65384	3.07%	0.222%
92	19.003	2703	2706	2713	rBV	58259	91959	4.32%	0.312%
93	19.574	2798	2803	2816	rVB	1368811	1943354	91.33%	6.598%
94	20.568	2968	2972	2975	rBV	145901	170693	8.02%	0.580%
95	21.068	3054	3057	3060	rBV	112566	132264	6.22%	0.449%
96	21.268	3088	3091	3096	rVB	338858	402438	18.91%	1.366%
97	21.356	3101	3106	3116	rBV	1585452	2127789	100.00%	7.224%
98	22.209	3248	3251	3259	rVB	230400	302249	14.20%	1.026%
99	23.491	3463	3469	3479	rVB	957422	1958338	92.04%	6.649%
100	23.638	3488	3494	3501	rVB	1103733	2007295	94.34%	6.815%

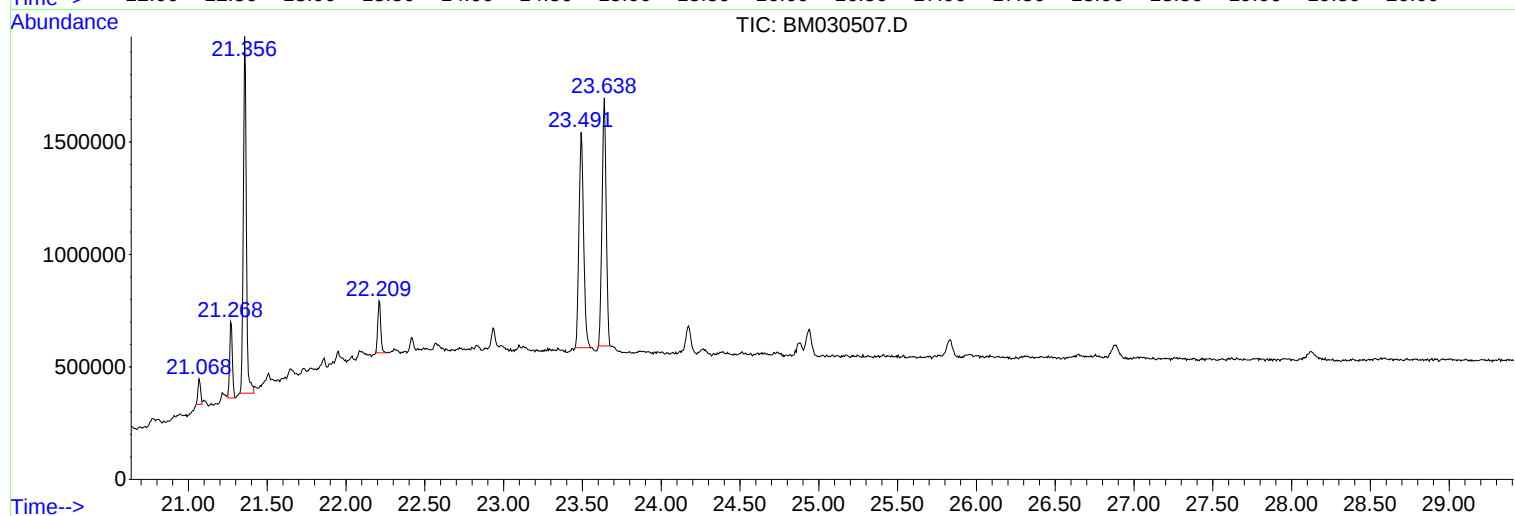
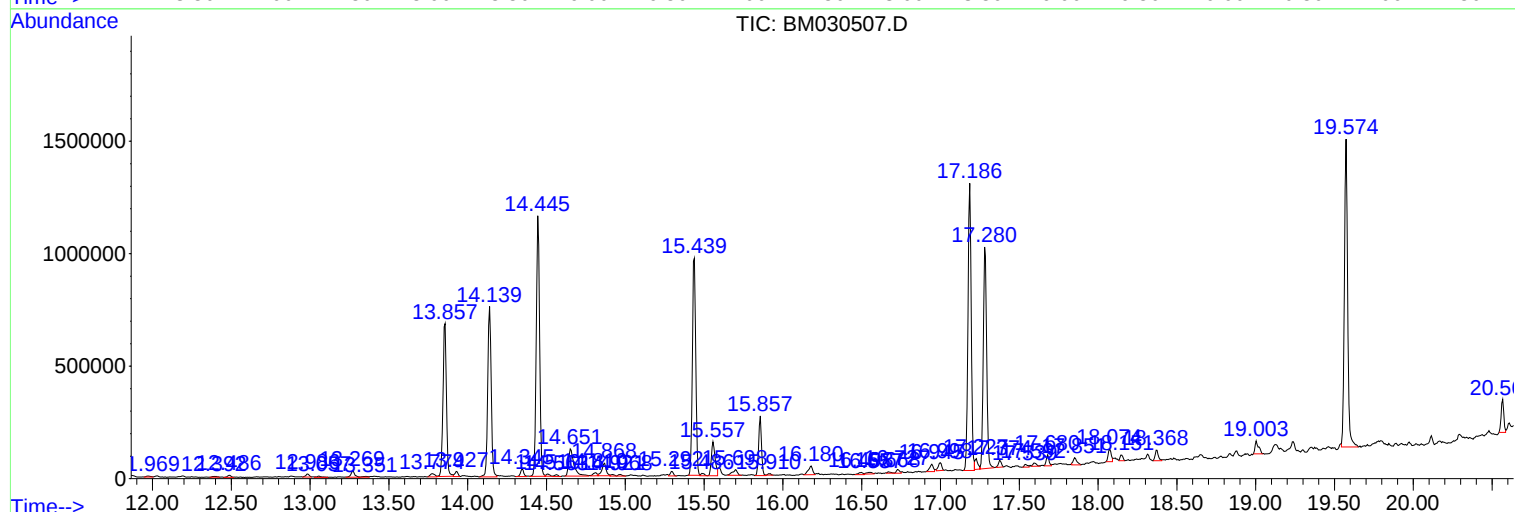
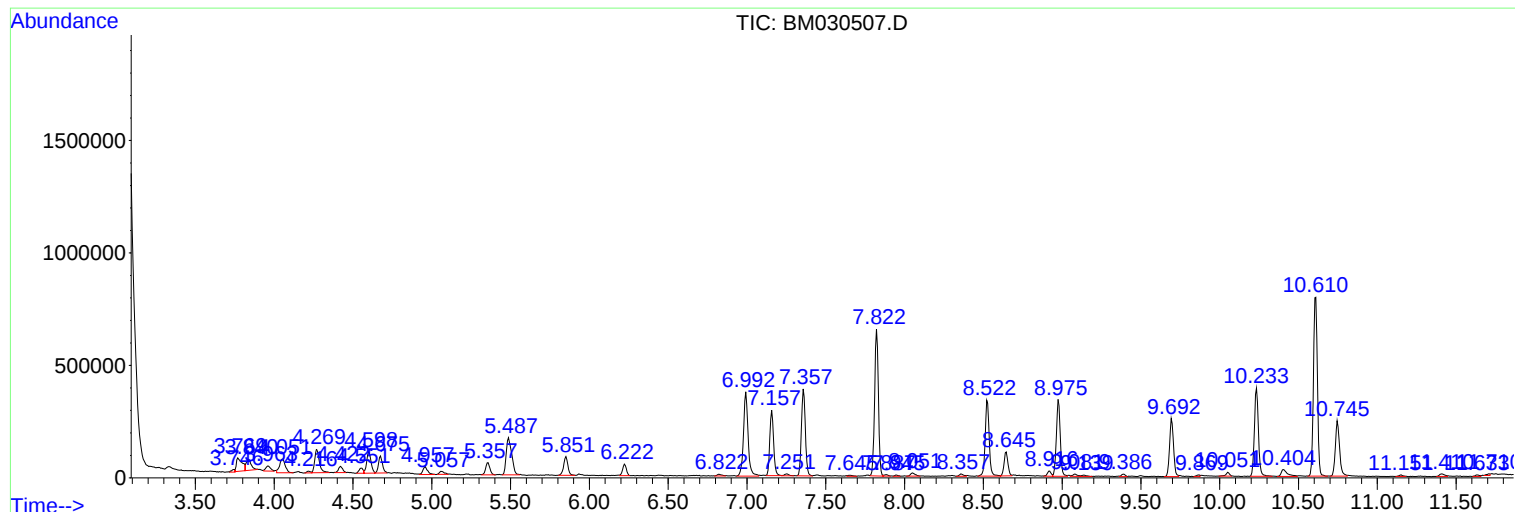
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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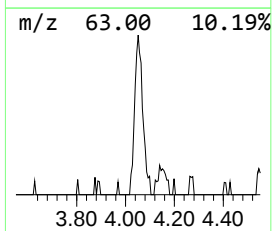
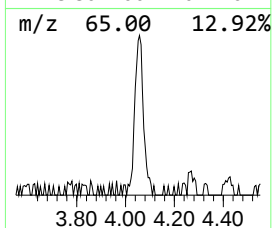
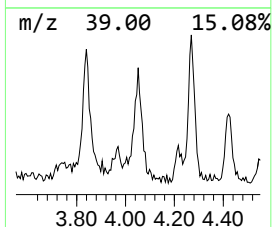
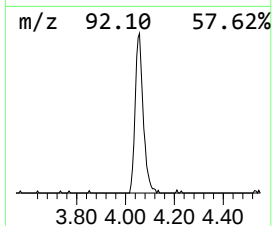
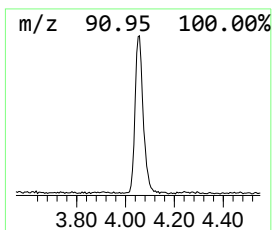
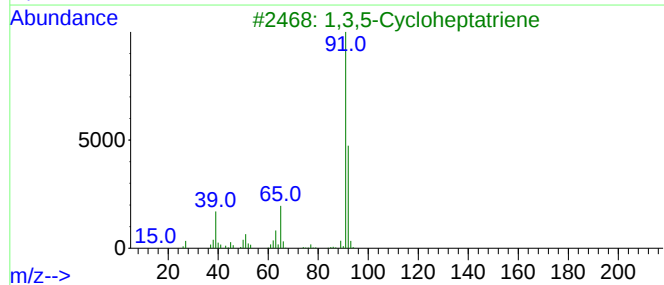
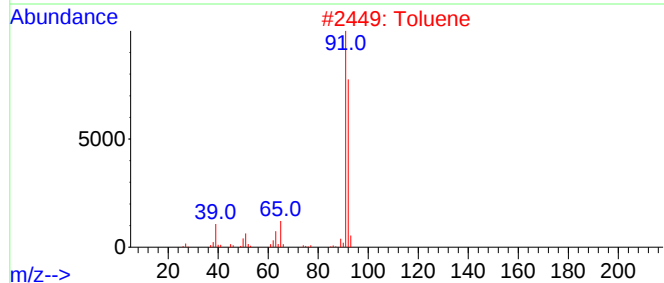
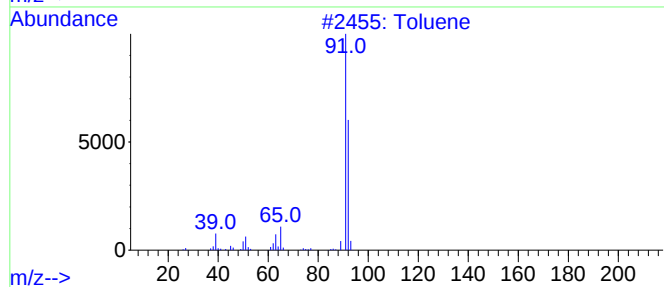
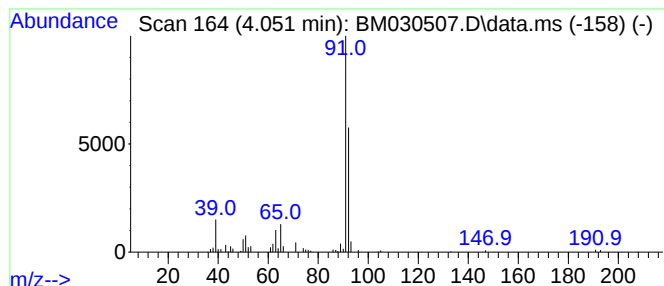
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Toluene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.050	2.79 ng/ul	151904	1,4-Dichlorobenzene-d4	7.822

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Toluene	92	C7H8	000108-88-3	94
2			Toluene	92	C7H8	000108-88-3	91
3			1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	90
4			Toluene	92	C7H8	000108-88-3	90
5			Toluene	92	C7H8	000108-88-3	87



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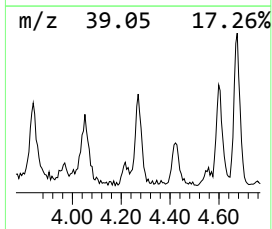
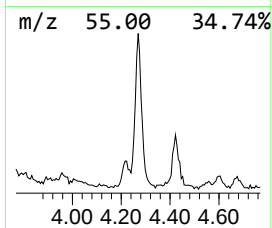
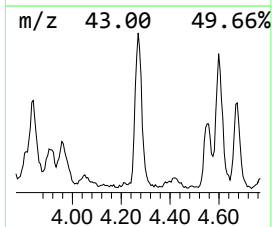
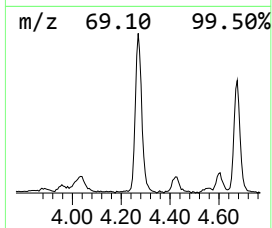
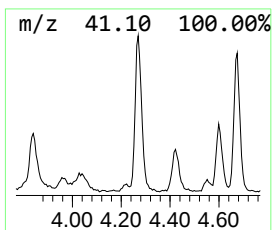
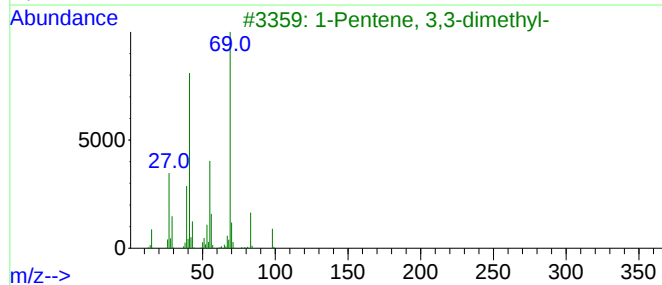
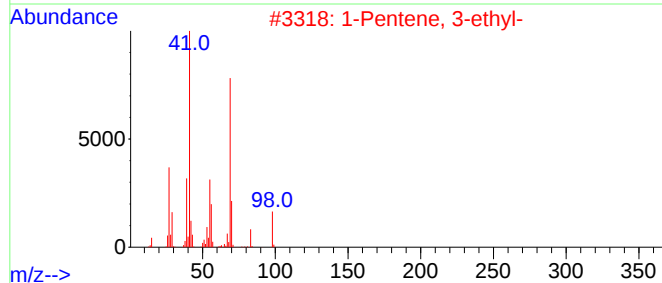
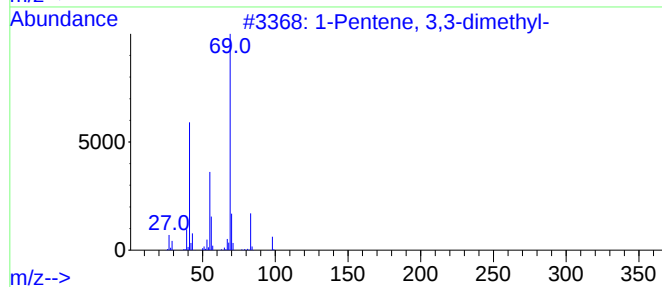
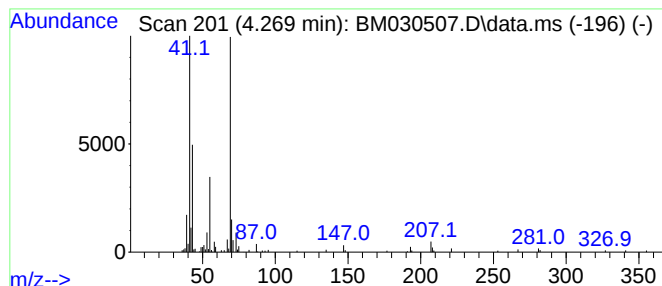
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.270	3.90 ng/ul	212129	1,4-Dichlorobenzene-d4	7.822

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene, 3,3-dimethyl-	98	C7H14	003404-73-7	45	
2	1-Pentene, 3-ethyl-	98	C7H14	004038-04-4	42	
3	1-Pentene, 3,3-dimethyl-	98	C7H14	003404-73-7	38	
4	1-Pentyn-3-ol, 3,4-dimethyl-	112	C7H12O	001482-15-1	38	
5	1-Pentene, 2,3-dimethyl-	98	C7H14	003404-72-6	36	



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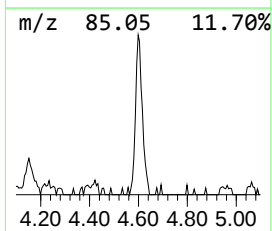
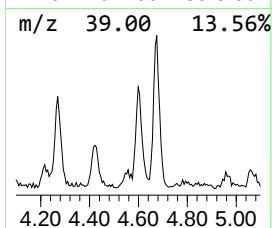
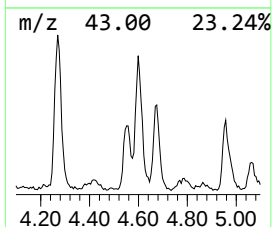
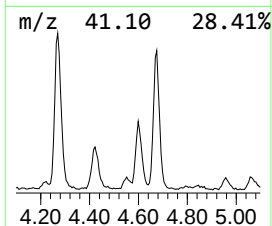
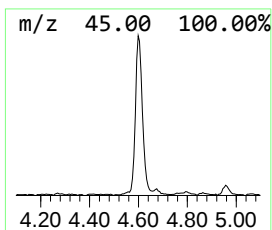
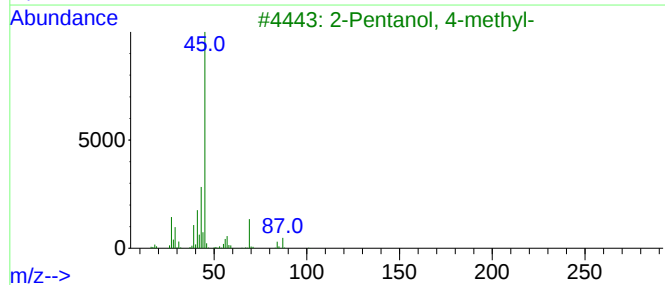
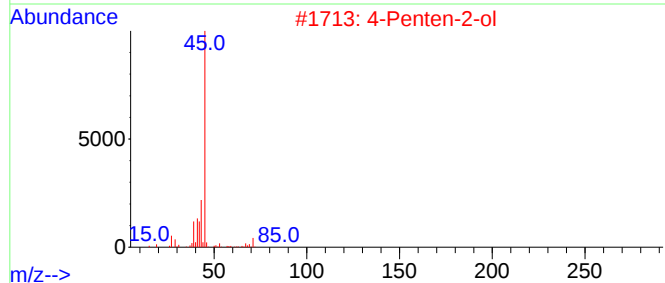
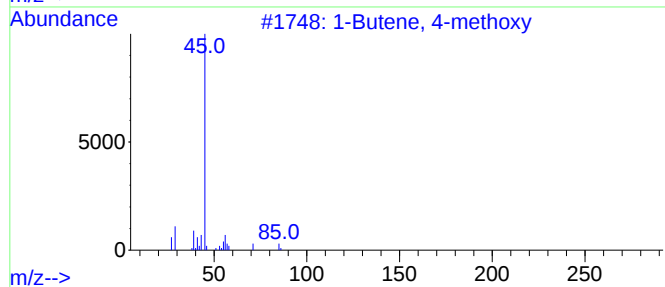
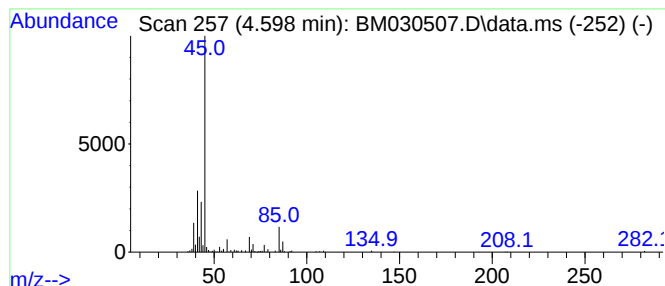
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.600	3.29 ng/ul	179151	1,4-Dichlorobenzene-d4	7.822

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Butene, 4-methoxy	86	C5H10O	004696-30-4	45
2			4-Penten-2-ol	86	C5H10O	000625-31-0	45
3			2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	45
4			2-Hexanol, 3-methyl-	116	C7H16O	002313-65-7	45
5			2-Pentanol, 3-methyl-	102	C6H14O	000565-60-6	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062121\
Data File : BM030507.D
Acq On : 21 Jun 2021 16:42
Operator : CG/JU
Sample : M2697-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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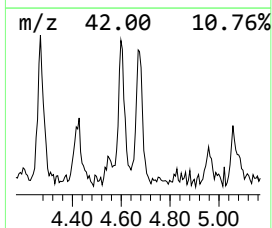
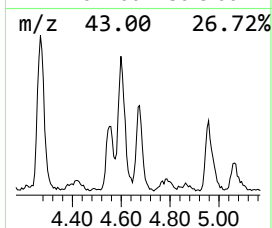
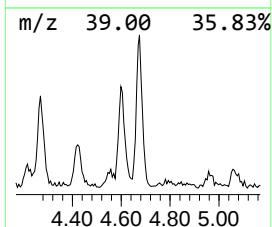
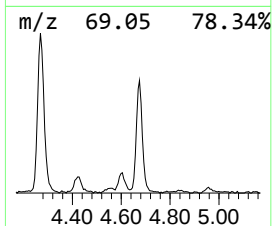
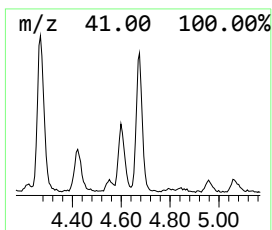
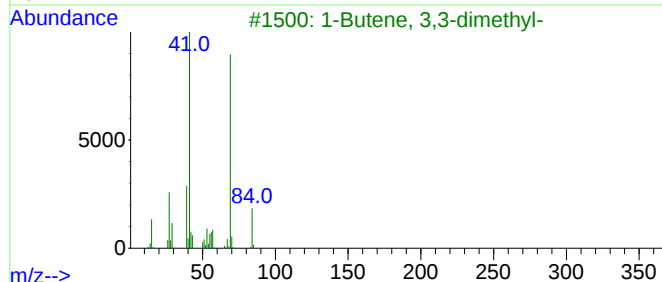
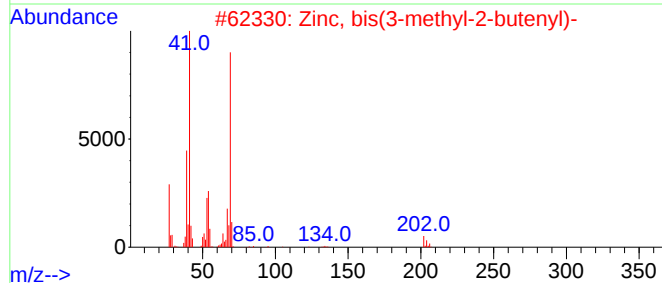
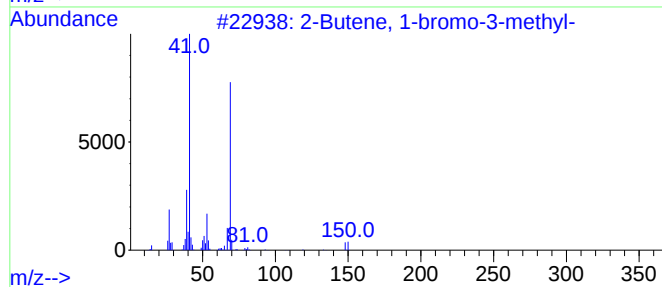
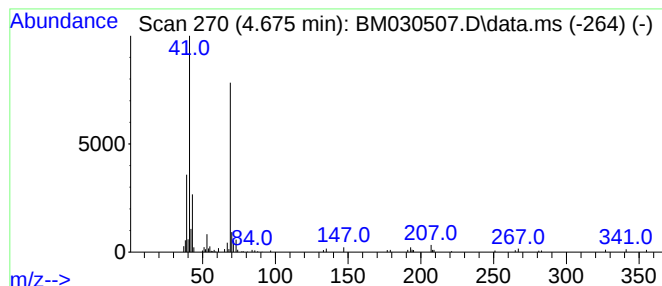
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.670	2.44 ng/ul	132500	1,4-Dichlorobenzene-d4	7.822

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butene, 1-bromo-3-methyl-	148	C5H9Br	000870-63-3	40		
2	Zinc, bis(3-methyl-2-butenyl)-	202	C10H18Zn	066094-28-8	40		
3	1-Butene, 3,3-dimethyl-	84	C6H12	000558-37-2	40		
4	1-Pentene, 2,3-dimethyl-	98	C7H14	003404-72-6	39		
5	Cyclopentane, bromo-	148	C5H9Br	000137-43-9	39		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062121\
Data File : BM030507.D
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Operator : CG/JU
Sample : M2697-01
Misc :
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Instrument :
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ClientSampleId :
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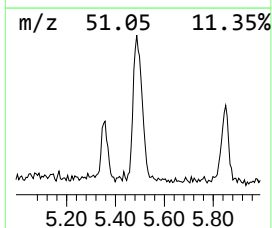
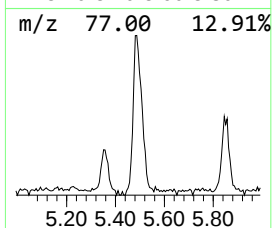
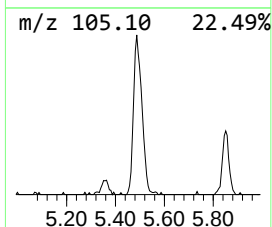
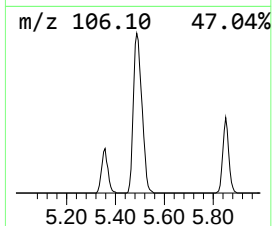
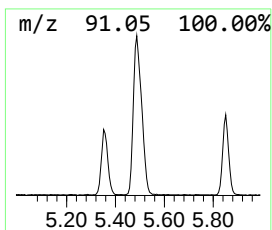
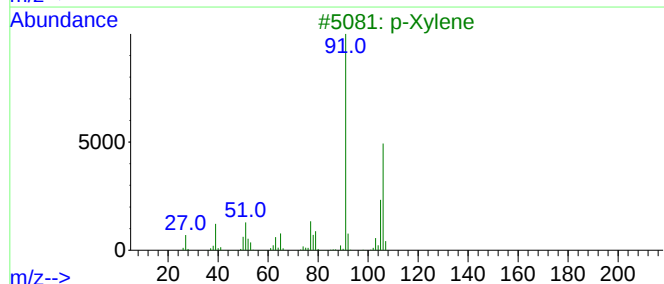
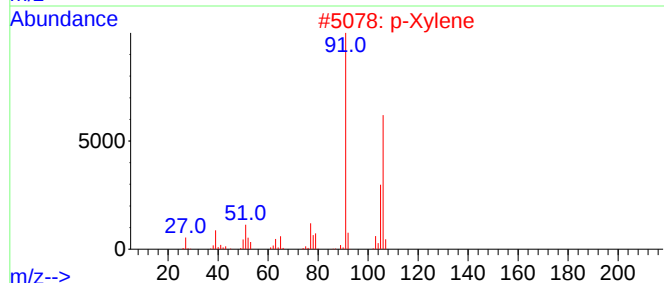
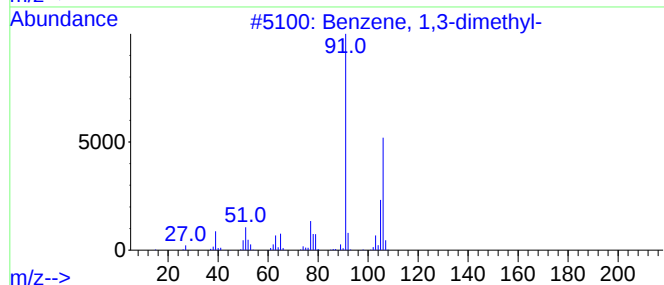
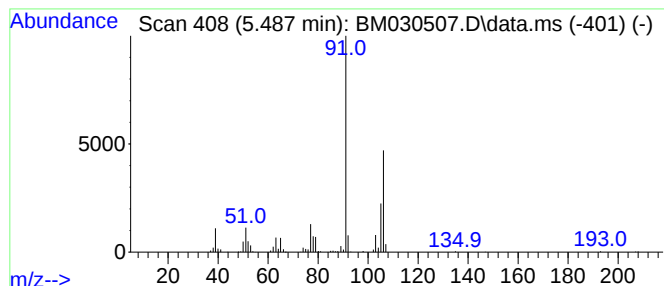
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.490	7.32 ng/ul	397909	1,4-Dichlorobenzene-d4	7.822

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2			p-Xylene	106	C8H10	000106-42-3	97
3			p-Xylene	106	C8H10	000106-42-3	97
4			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
5			o-Xylene	106	C8H10	000095-47-6	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062121\
Data File : BM030507.D
Acq On : 21 Jun 2021 16:42
Operator : CG/JU
Sample : M2697-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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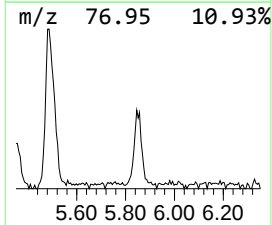
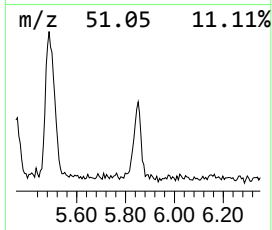
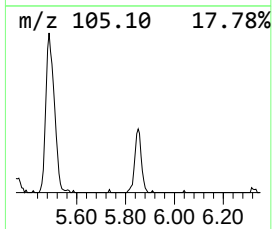
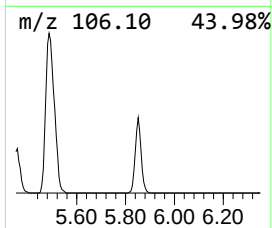
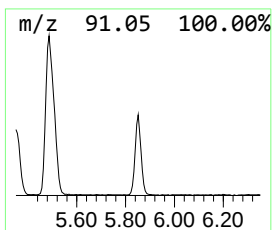
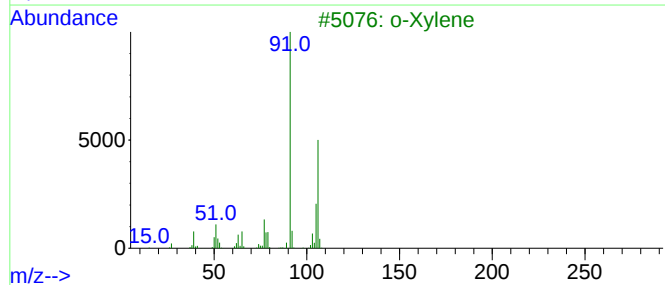
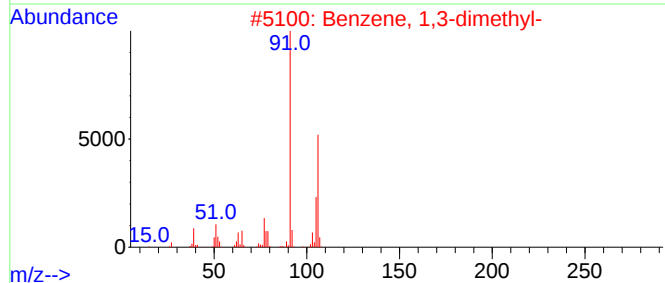
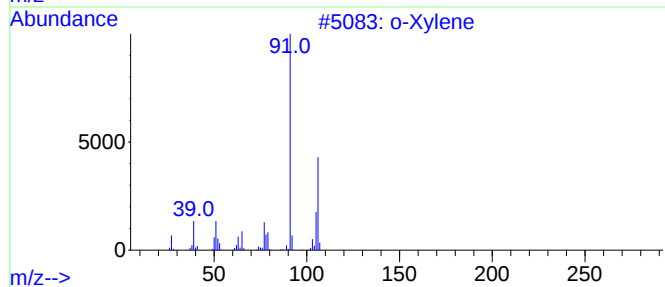
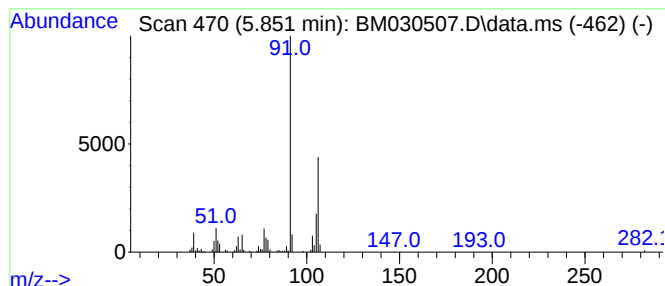
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 o-Xylene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.850	2.86 ng/ul	155754	1,4-Dichlorobenzene-d4	7.822

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Xylene	106	C8H10	000095-47-6	97
2			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
3			o-Xylene	106	C8H10	000095-47-6	97
4			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
5			o-Xylene	106	C8H10	000095-47-6	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062121\
Data File : BM030507.D
Acq On : 21 Jun 2021 16:42
Operator : CG/JU
Sample : M2697-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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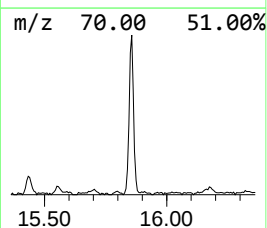
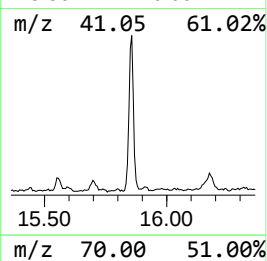
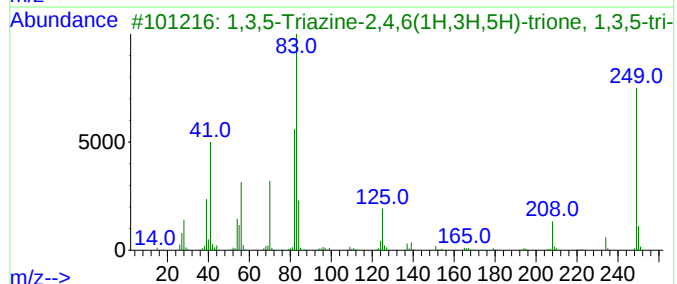
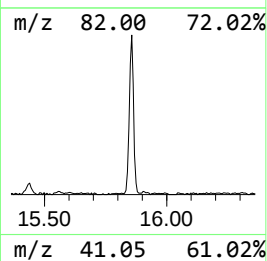
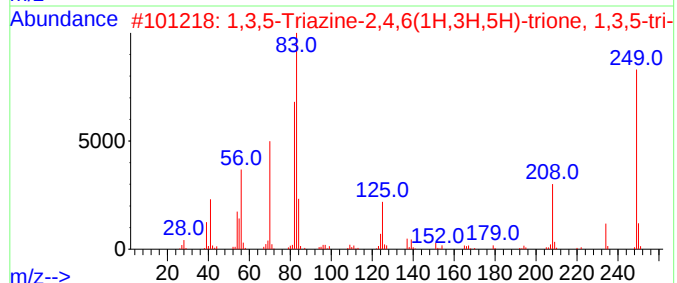
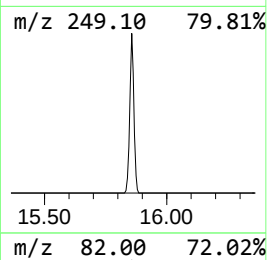
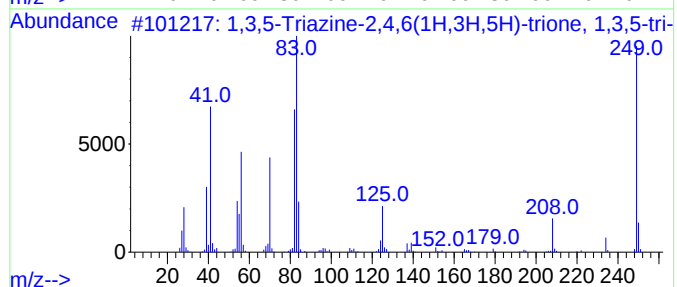
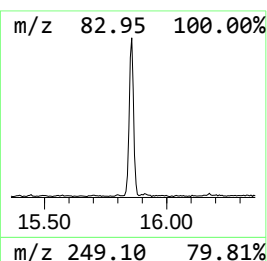
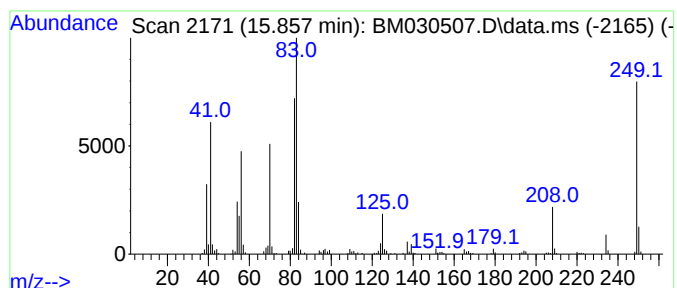
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1,3,5-Triazine-2,4,6(1H,3H,... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.860	3.63 ng/ul	332171	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	99		
2	1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	99		
3	1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	95		
4	4H-1-Benzopyran-2-carboxylic aci...	249	C12H11NO5	032142-43-1	53		
5	2-Methylphenanthro[3,4-d][1,3]ox...	249	C16H11NO2	098033-24-0	45		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062121\
Data File : BM030507.D
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Sample : M2697-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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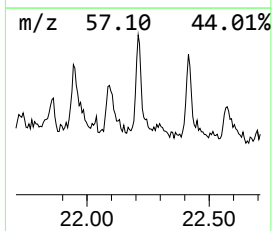
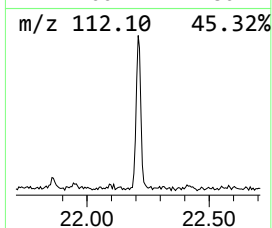
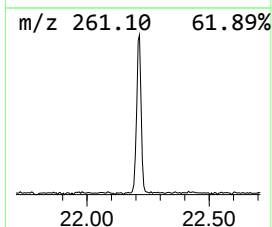
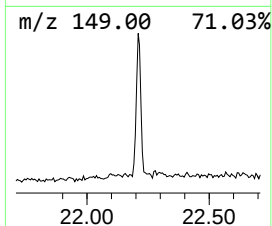
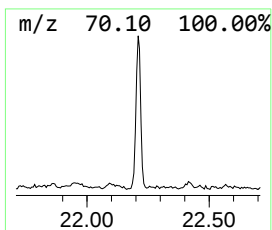
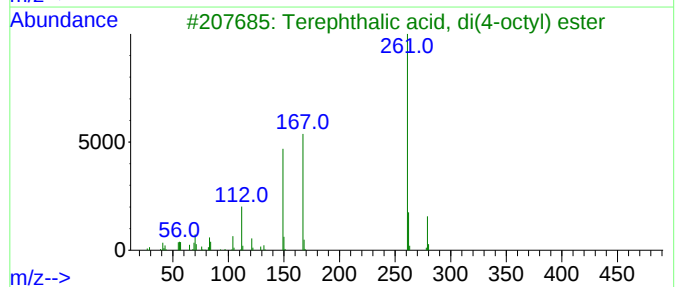
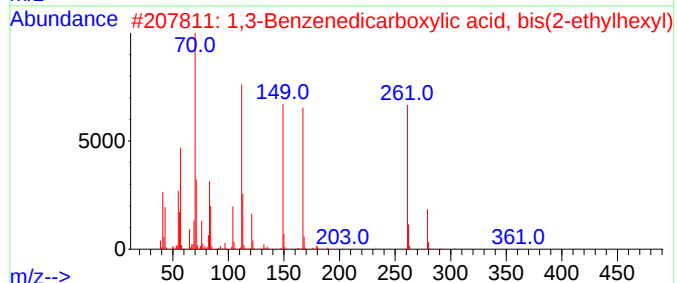
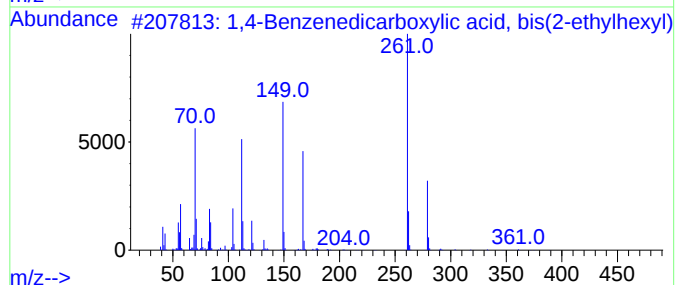
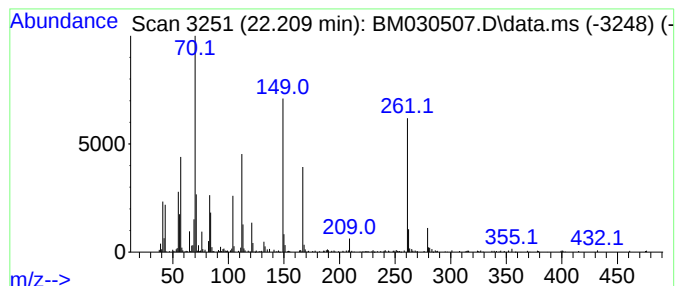
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1,4-Benzenedicarboxylic aci... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.210	2.84 ng/ul	302249	Chrysene-d12	21.356

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	64
2			1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	58
3			Terephthalic acid, di(4-octyl) e...	390	C24H38O4	1000323-74-2	43
4			Terephthalic acid, 3-methylbut-2...	348	C21H32O4	1000356-27-6	43
5			Terephthalic acid, 2-ethylhexyl ...	390	C24H38O4	1000324-00-5	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062121\
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 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Toluene	4.050	2.8	ng/ul	151904	1	7.822	1087750	20.0
(DEL) Alkane: S...	4.270	3.9	ng/ul	212129	1	7.822	1087750	20.0
unknown-01	4.600	3.3	ng/ul	179151	1	7.822	1087750	20.0
unknown-02	4.670	2.4	ng/ul	132500	1	7.822	1087750	20.0
Benzene, 1,3-di...	5.490	7.3	ng/ul	397909	1	7.822	1087750	20.0
o-Xylene	5.850	2.9	ng/ul	155754	1	7.822	1087750	20.0
1,3,5-Triazine-...	15.860	3.6	ng/ul	332171	4	17.186	1830340	20.0
1,4-Benzenedica...	22.210	2.8	ng/ul	302249	5	21.356	2127790	20.0