

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
 Data File : BP019400.D
 Acq On : 19 Jan 2024 12:43
 Operator : MA/JU
 Sample : P1160-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 XA313

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019400.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.210	18	21	27	rVB	45541	54097	0.33%	0.028%
2	3.569	76	82	88	rBV	1402107	1860315	11.32%	0.950%
3	3.628	88	92	110	rVV	579776	893099	5.43%	0.456%
4	3.763	113	115	123	rVB3	16073	23667	0.14%	0.012%
5	3.946	141	146	157	rVB	210349	275909	1.68%	0.141%
6	4.069	164	167	182	rVB	28942	61171	0.37%	0.031%
7	4.787	285	289	295	rVB	60132	77249	0.47%	0.039%
8	4.993	318	324	327	rBV5	9252	21773	0.13%	0.011%
9	6.957	652	658	672	rVB	737754	1220010	7.42%	0.623%
10	7.110	677	684	692	rBV	555622	873930	5.32%	0.446%
11	7.204	693	700	708	rBV	1390877	2050906	12.48%	1.047%
12	7.304	710	717	726	rVB	702122	1113498	6.77%	0.569%
13	7.651	769	776	784	rBV	960186	1470083	8.94%	0.751%
14	7.763	788	795	810	rVB	526937	844914	5.14%	0.431%
15	8.116	846	855	866	rBV	1191307	1897216	11.54%	0.969%
16	8.504	914	921	928	rBV	825585	1380650	8.40%	0.705%
17	8.581	928	934	943	rVB	622600	930711	5.66%	0.475%
18	8.834	969	977	987	rVB	1409662	2240293	13.63%	1.144%
19	8.934	987	994	998	rBV	675987	1188102	7.23%	0.607%
20	8.981	998	1002	1013	rVB	1459991	2229842	13.56%	1.138%
21	9.510	1080	1092	1101	rBV	4197454	6872057	41.80%	3.509%
22	9.657	1109	1117	1127	rBV	602337	979702	5.96%	0.500%
23	9.816	1138	1144	1152	rVB	57143	92510	0.56%	0.047%
24	9.998	1166	1175	1184	rBV	3052462	4978211	30.28%	2.542%
25	10.198	1201	1209	1223	rBV	909345	1555972	9.47%	0.794%
26	10.322	1225	1230	1235	rVB3	20168	32201	0.20%	0.016%
27	10.434	1240	1249	1256	rBV	3522118	5778165	35.15%	2.950%
28	10.563	1259	1271	1275	rBV	629827	1197939	7.29%	0.612%
29	10.616	1275	1280	1289	rVB	4132507	6751082	41.07%	3.447%
30	10.716	1289	1297	1313	rBV	720300	1323018	8.05%	0.675%
31	10.892	1319	1327	1334	rBV	5295660	8153309	49.60%	4.163%
32	10.951	1334	1337	1341	rVB3	12210	17110	0.10%	0.009%
33	11.263	1382	1390	1394	rBV	29210	52054	0.32%	0.027%
34	11.316	1394	1399	1404	rVB2	26341	41872	0.25%	0.021%
35	11.798	1476	1481	1489	rVB	33889	58603	0.36%	0.030%
36	12.057	1519	1525	1528	rBV5	19727	38205	0.23%	0.020%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoother: OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
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37	12.104	1528	1533	1539	rVV2	82712	143529	0.87%	0.073%
38	12.157	1539	1542	1547	rVB	16223	23468	0.14%	0.012%
39	12.234	1547	1555	1562	rBV	2783713	4352510	26.48%	2.222%
40	12.363	1570	1577	1587	rBV2	88449	187091	1.14%	0.096%
41	12.451	1587	1592	1598	rVB	40505	70736	0.43%	0.036%
42	12.569	1605	1612	1621	rVB	2625213	3989392	24.27%	2.037%
43	12.663	1624	1628	1639	rVB	20614	31543	0.19%	0.016%
44	13.292	1728	1735	1747	rVB	4365480	6498710	39.53%	3.318%
45	13.475	1761	1766	1771	rBV	69958	98687	0.60%	0.050%
46	13.528	1771	1775	1781	rVB2	35752	54040	0.33%	0.028%
47	13.592	1781	1786	1793	rVB	27726	41692	0.25%	0.021%
48	13.839	1821	1828	1832	rBV	1708766	2367570	14.40%	1.209%
49	13.886	1832	1836	1851	rVB	2422490	3217334	19.57%	1.643%
50	14.016	1851	1858	1865	rBV	668606	910427	5.54%	0.465%
51	14.110	1867	1874	1877	rVV2	1803664	2941451	17.89%	1.502%
52	14.139	1877	1879	1890	rVB	2090960	2607706	15.86%	1.331%
53	14.416	1917	1926	1934	rBV	1074932	1632788	9.93%	0.834%
54	14.669	1963	1969	1977	rBV	564290	990605	6.03%	0.506%
55	14.739	1977	1981	1991	rVV	54279	106010	0.64%	0.054%
56	14.822	1991	1995	1999	rVV	23823	40269	0.24%	0.021%
57	14.863	1999	2002	2009	rVB2	26516	39419	0.24%	0.020%
58	15.257	2062	2069	2076	rBV	3076634	3945783	24.00%	2.015%
59	15.416	2086	2096	2100	rBV	2431053	3467605	21.09%	1.770%
60	15.475	2100	2106	2114	rVV	12518658	16439175	100.00%	8.393%
61	15.557	2114	2120	2131	rVV	884164	1181716	7.19%	0.603%
62	15.657	2131	2137	2143	rVV4	12950	27899	0.17%	0.014%
63	15.722	2143	2148	2154	rVB2	38253	57484	0.35%	0.029%
64	15.933	2179	2184	2190	rBV2	27407	38666	0.24%	0.020%
65	16.010	2190	2197	2202	rVV7	11108	18080	0.11%	0.009%
66	16.069	2202	2207	2213	rVV	52583	72811	0.44%	0.037%
67	16.369	2248	2258	2264	rBV	4883361	6418194	39.04%	3.277%
68	16.480	2269	2277	2284	rVV	4980064	6956382	42.32%	3.552%
69	16.845	2330	2339	2347	rBV4	11636	27476	0.17%	0.014%
70	16.974	2356	2361	2363	rBV2	20909	31047	0.19%	0.016%
71	17.180	2388	2396	2399	rBV	1504005	2127753	12.94%	1.086%
72	17.222	2399	2403	2408	rVV	3070263	4268507	25.97%	2.179%
73	17.280	2408	2413	2416	rVV	2675010	3854132	23.44%	1.968%
74	17.316	2416	2419	2430	rVB	3584897	4663629	28.37%	2.381%
75	17.463	2438	2444	2449	rVB	10435	21959	0.13%	0.011%
76	17.610	2465	2469	2477	rVB	17422	25656	0.16%	0.013%

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77	17.792	2496	2500	2503	rBV2	20581	28649	0.17%	0.015%
78	18.151	2555	2561	2566	rBV	4239295	5275924	32.09%	2.694%
79	18.551	2626	2629	2631	rBV3	16898	18020	0.11%	0.009%
80	18.586	2631	2635	2652	rVB	238850	363378	2.21%	0.186%
81	18.810	2670	2673	2676	rBV	17351	21583	0.13%	0.011%
82	19.104	2718	2723	2729	rBV	40573	60260	0.37%	0.031%
83	19.174	2731	2735	2738	rBV	32355	45064	0.27%	0.023%
84	19.527	2789	2795	2803	rBV	3798384	4742843	28.85%	2.422%
85	21.292	3089	3095	3097	rBV	1885179	2357387	14.34%	1.204%
86	21.327	3097	3101	3108	rVB	4637461	5452499	33.17%	2.784%
87	22.068	3224	3227	3229	rBV	172665	222042	1.35%	0.113%
88	22.109	3229	3234	3240	rVB	2281788	2971644	18.08%	1.517%
89	22.927	3366	3373	3377	rBV	2330167	4234863	25.76%	2.162%
90	22.980	3377	3382	3394	rVB	4654078	7752175	47.16%	3.958%
91	23.498	3463	3470	3474	rBV	2121948	4216517	25.65%	2.153%
92	23.545	3474	3478	3488	rVB	2423507	4287385	26.08%	2.189%
93	23.645	3489	3495	3503	rVB	1168683	2170057	13.20%	1.108%
94	26.109	3902	3914	3927	rVB2	2981129	9019303	54.86%	4.605%

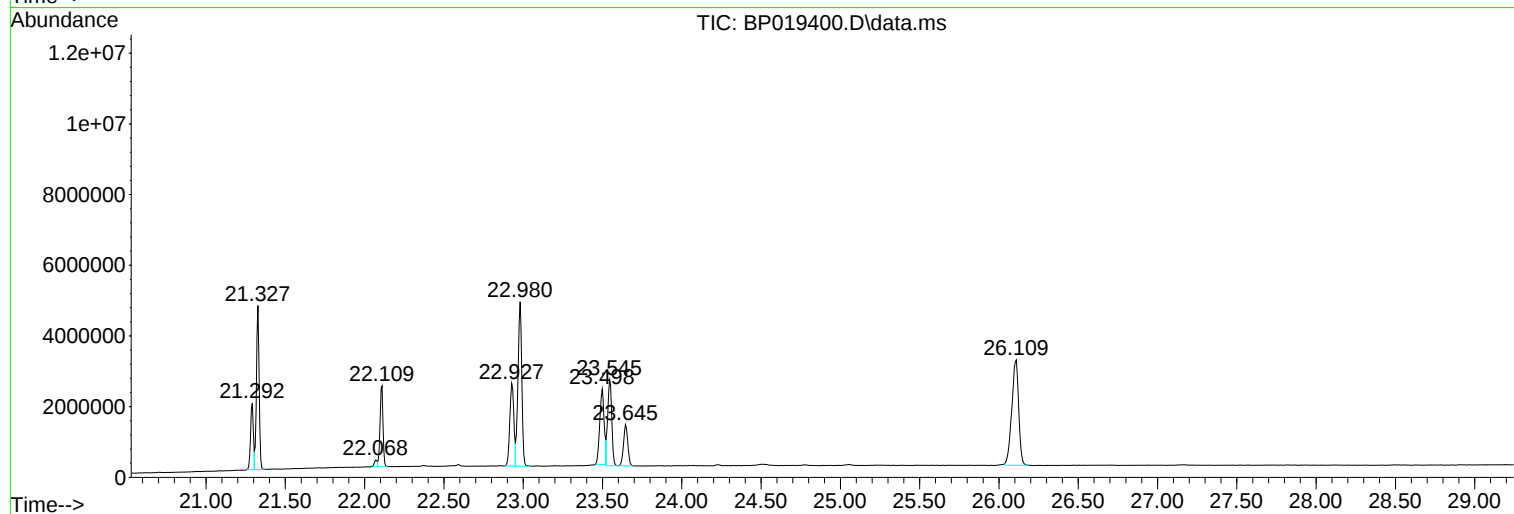
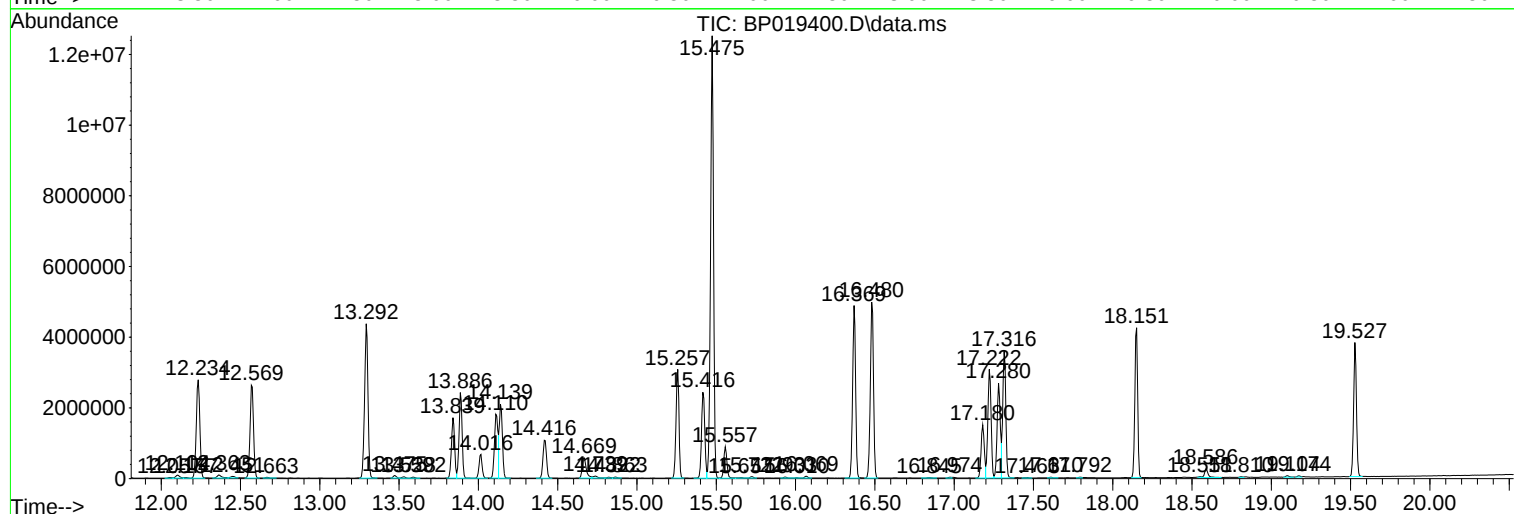
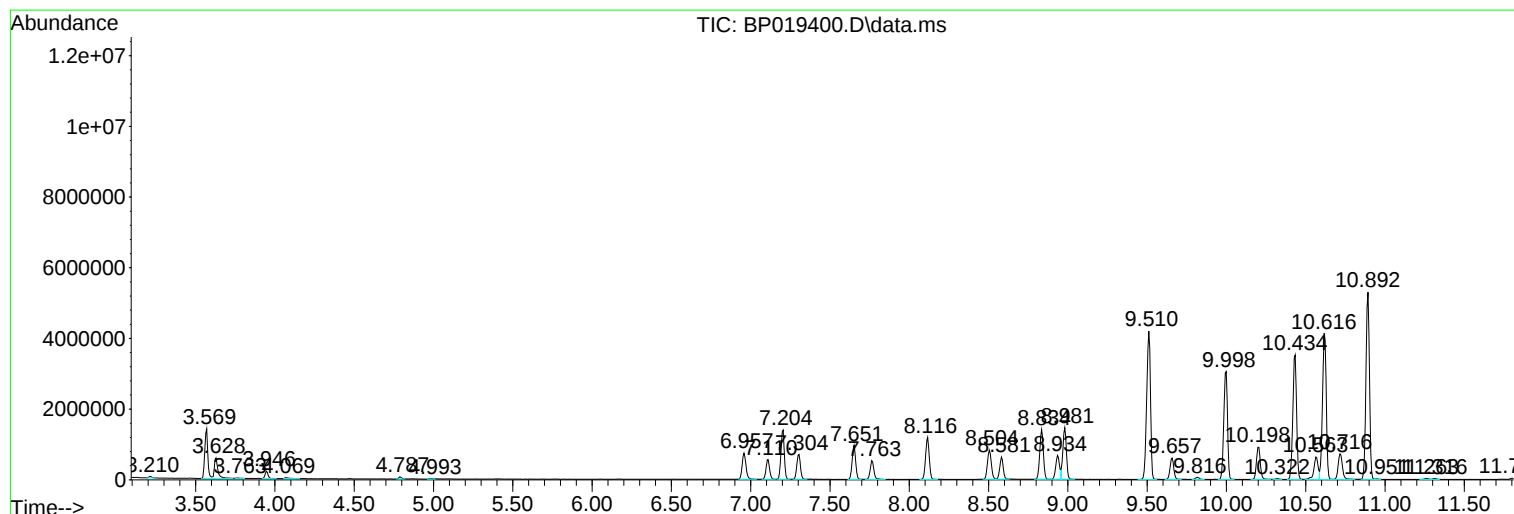
Sum of corrected areas: 195857969

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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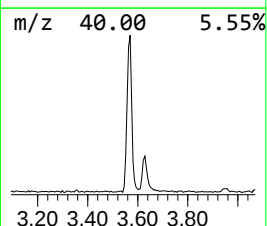
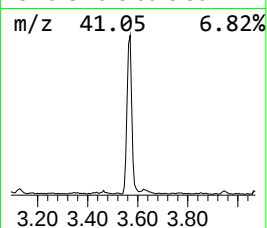
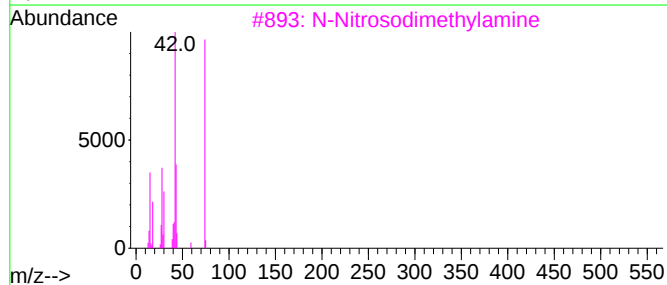
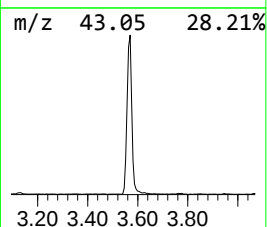
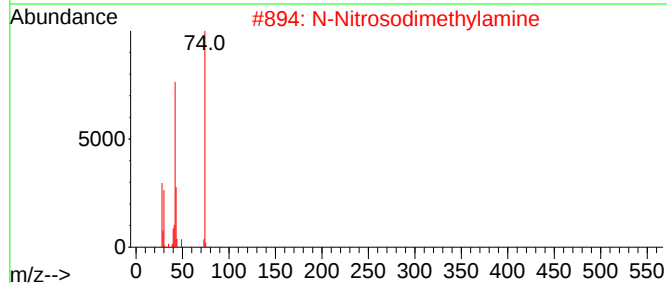
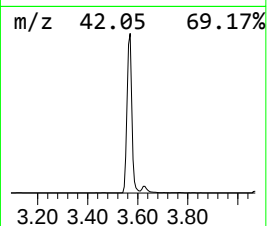
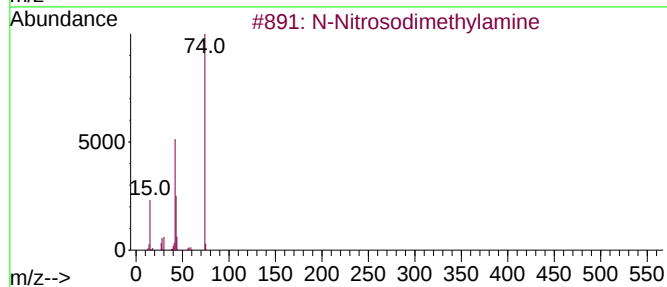
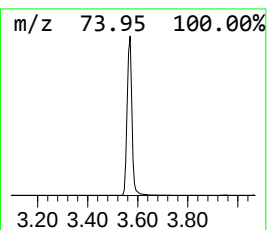
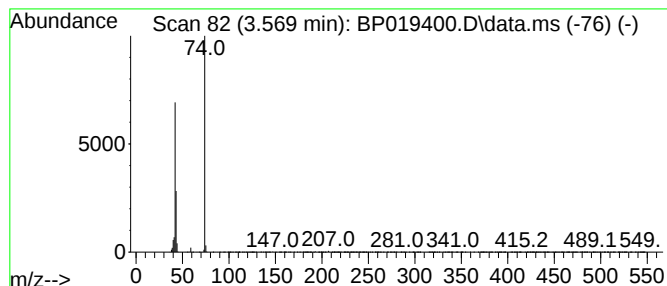
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 N-Nitrosodimethylamine Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.569	44.04 ng/ul	1860320	1,4-Dichlorobenzene-d4	7.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-Nitrosodimethylamine	74	C2H6N2O	000062-75-9	78
2			N-Nitrosodimethylamine	74	C2H6N2O	000062-75-9	9
3			N-Nitrosodimethylamine	74	C2H6N2O	000062-75-9	5
4			N-Nitrosodimethylamine	74	C2H6N2O	000062-75-9	4
5			Aminoguanidine	74	CH6N4	000079-17-4	4



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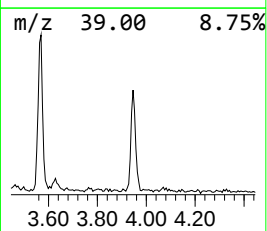
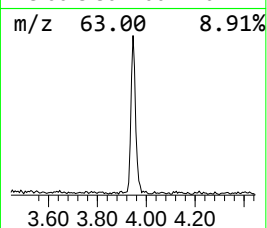
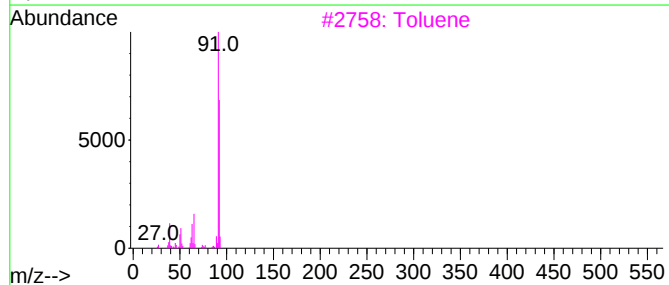
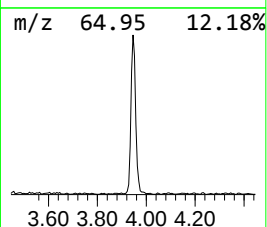
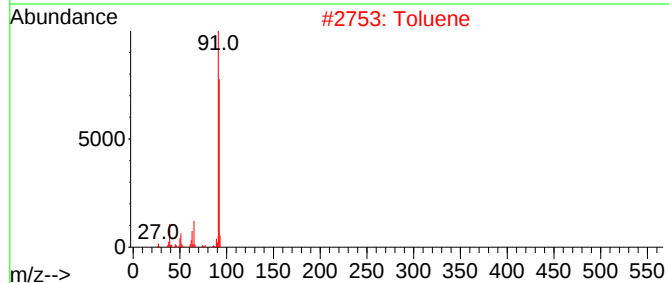
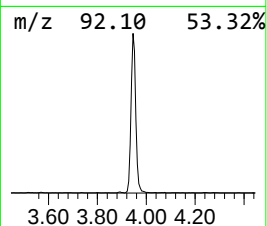
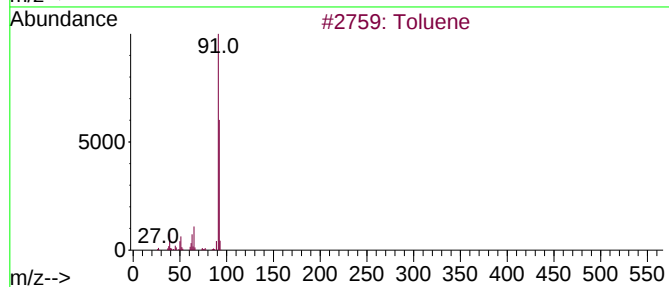
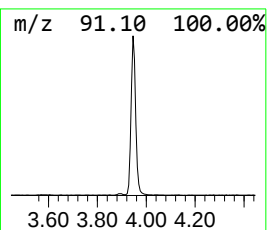
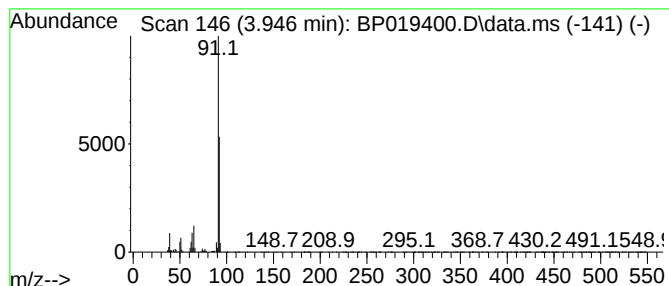
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Toluene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.946	6.53 ng/ul	275909	1,4-Dichlorobenzene-d4	7.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene			92	C7H8	000108-88-3	95
2	Toluene			92	C7H8	000108-88-3	95
3	Toluene			92	C7H8	000108-88-3	93
4	Toluene			92	C7H8	000108-88-3	91
5	1,3,5-Cycloheptatriene			92	C7H8	000544-25-2	90



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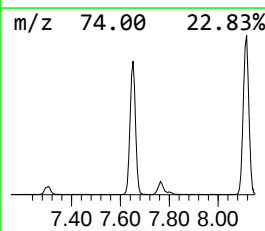
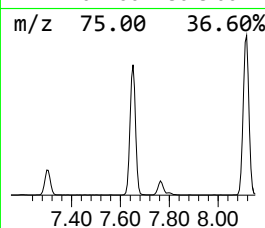
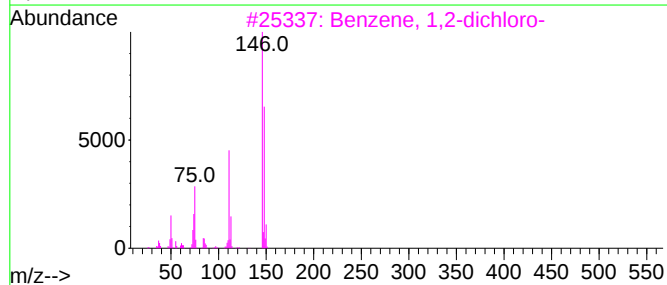
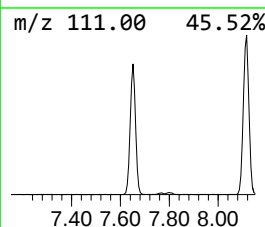
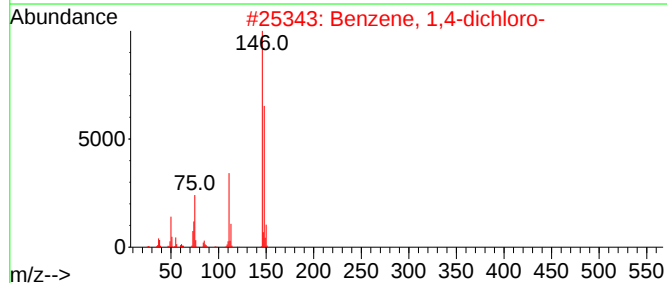
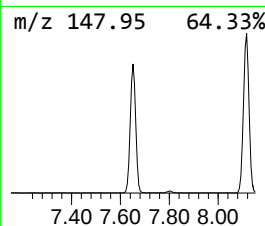
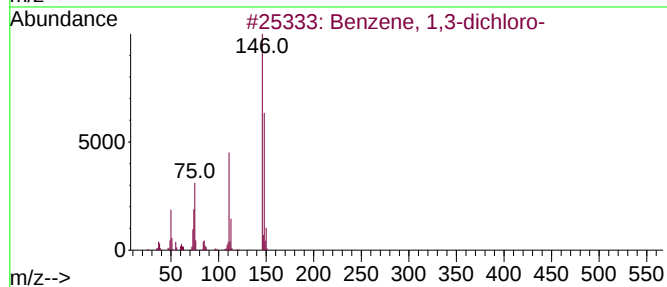
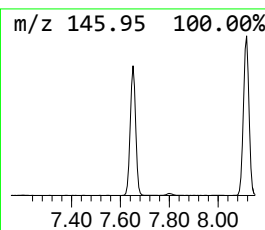
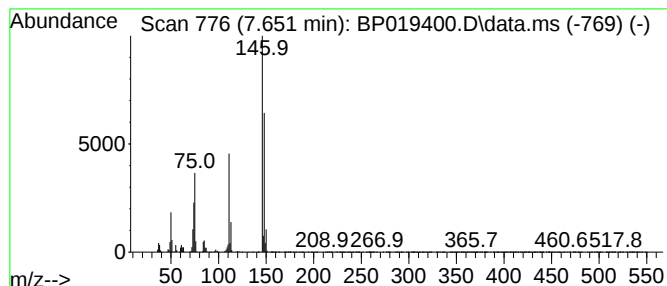
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,3-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.651	34.80 ng/ul	1470080	1,4-Dichlorobenzene-d4	7.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	96
4			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
Data File : BP019400.D
Acq On : 19 Jan 2024 12:43
Operator : MA/JU
Sample : P1160-04
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
XA313

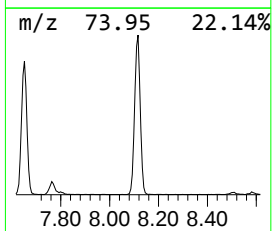
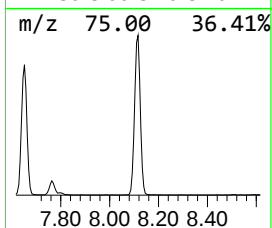
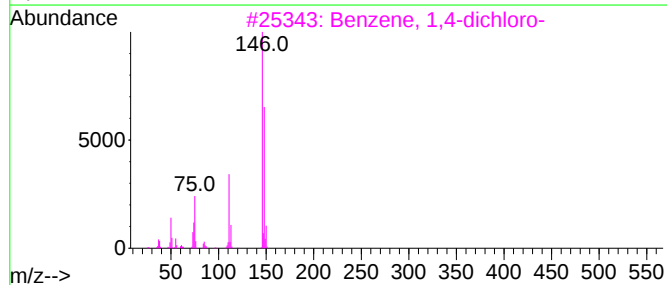
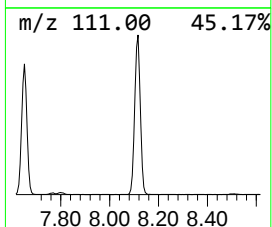
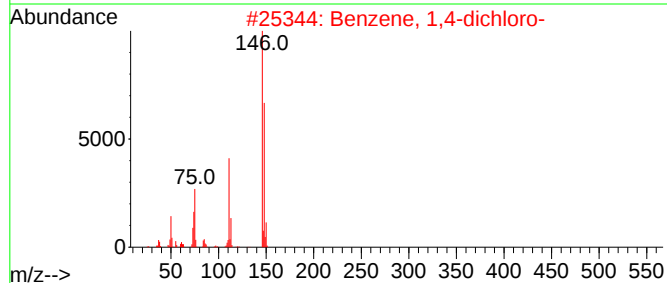
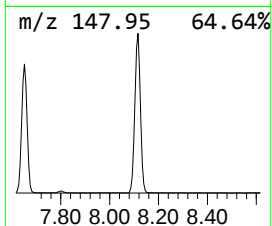
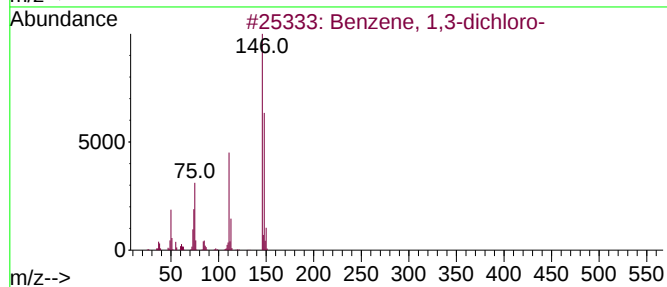
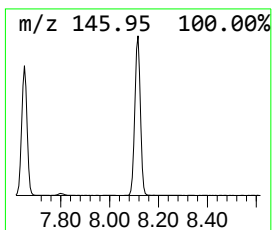
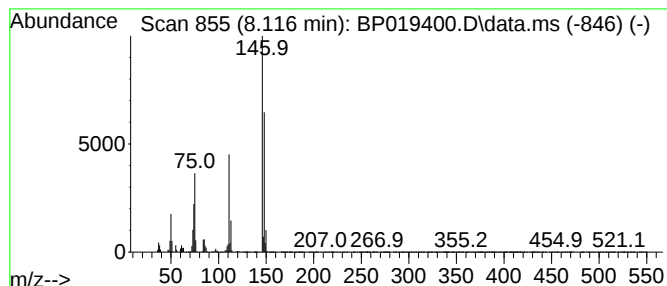
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1,4-dichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.116	44.91 ng/ul	1897220	1,4-Dichlorobenzene-d4	7.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
3			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	96
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
Data File : BP019400.D
Acq On : 19 Jan 2024 12:43
Operator : MA/JU
Sample : P1160-04
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
XA313

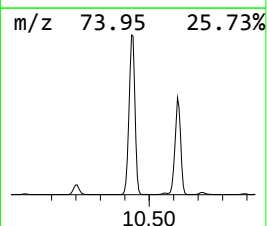
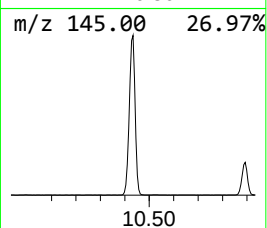
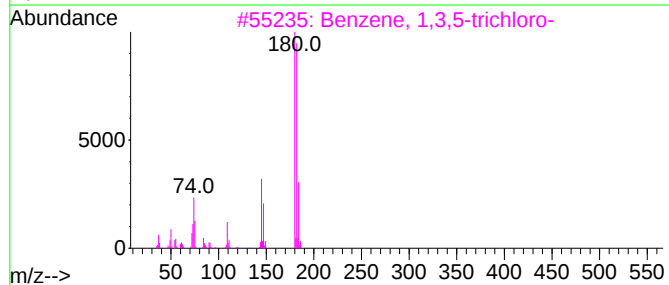
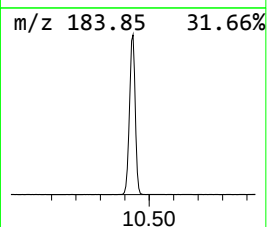
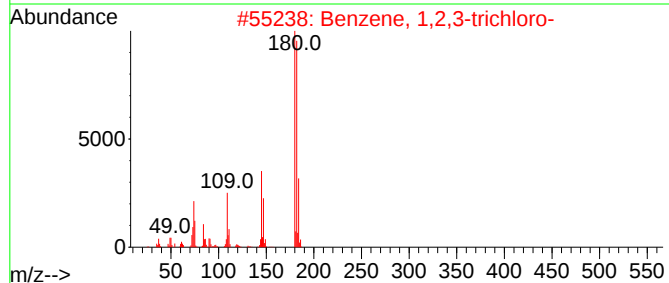
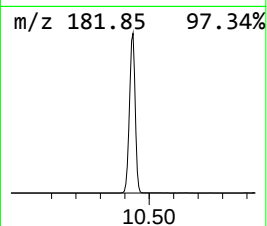
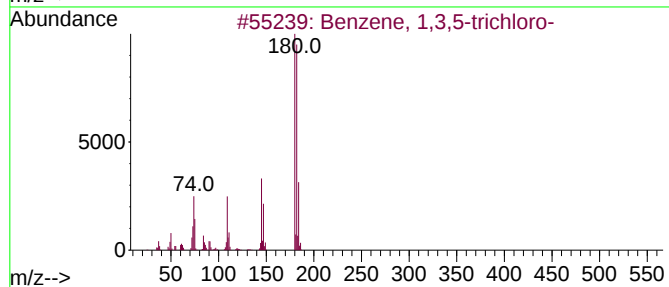
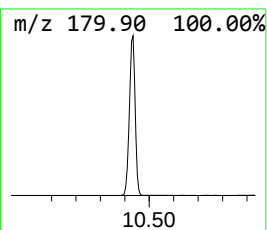
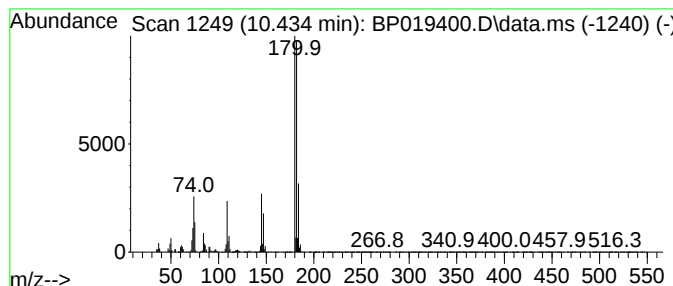
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.434	96.47 ng/ul	5778170	Naphthalene-d8	10.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
2			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
3			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	96
4			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	95
5			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
Data File : BP019400.D
Acq On : 19 Jan 2024 12:43
Operator : MA/JU
Sample : P1160-04
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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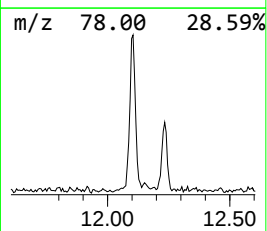
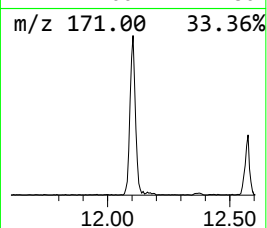
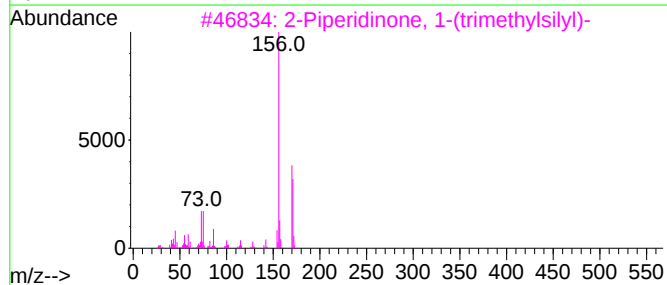
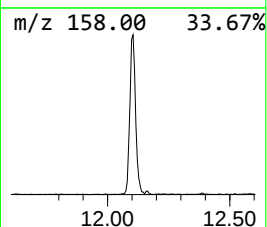
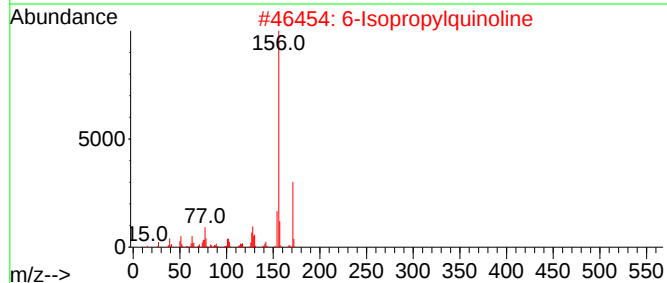
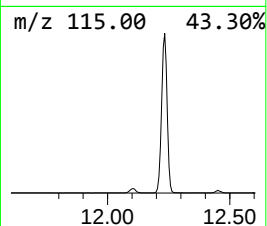
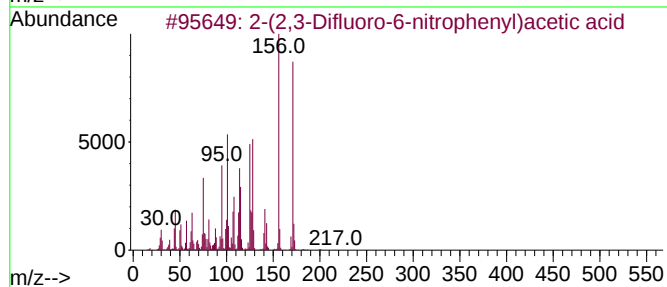
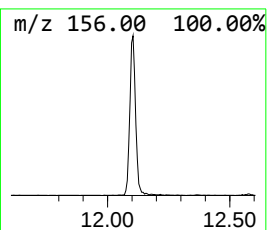
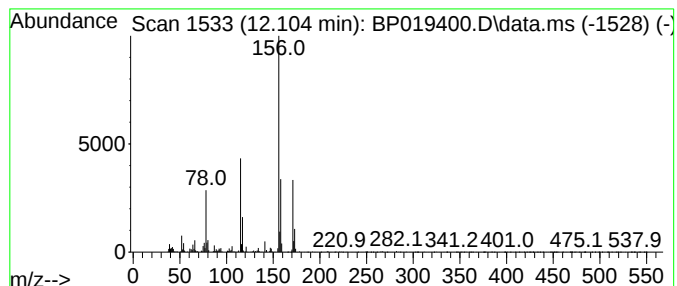
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.104	2.40 ng/ul	143529	Naphthalene-d8	10.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(2,3-Difluoro-6-nitrophenyl)ac...	217	C8H5F2NO4	141428-47-9	49		
2	6-Isopropylquinoline	171	C12H13N	000135-79-5	43		
3	2-Piperidinone, 1-(trimethylsilyl)-	171	C8H17NOSi	003553-93-3	37		
4	1,2,3,4-Tetrahydropentalene, 1,1...	171	C12H13N	1000217-18-6	37		
5	4-Chloro-3,5-dimethylphenol, 0-i...	242	C12H15ClO3	1000487-39-3	32		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
Data File : BP019400.D
Acq On : 19 Jan 2024 12:43
Operator : MA/JU
Sample : P1160-04
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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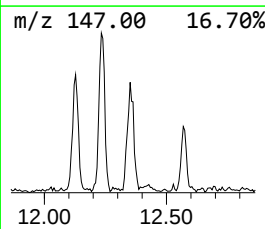
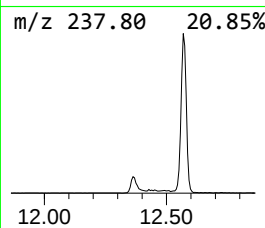
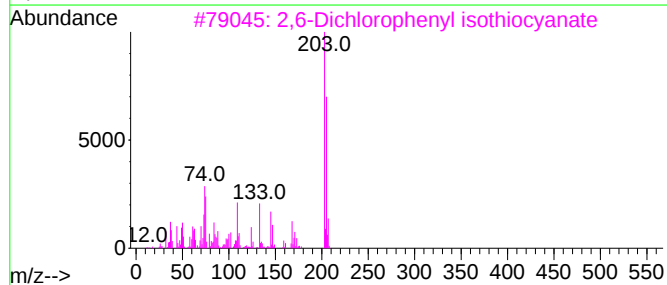
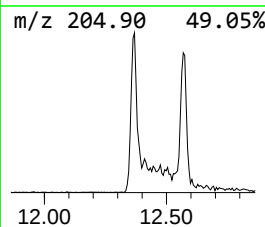
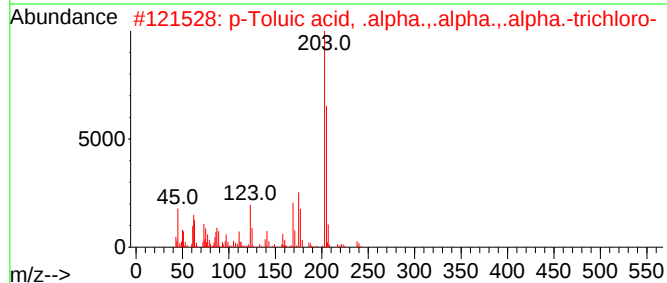
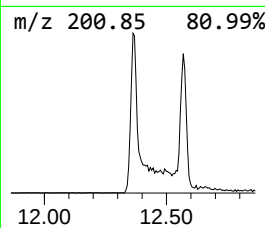
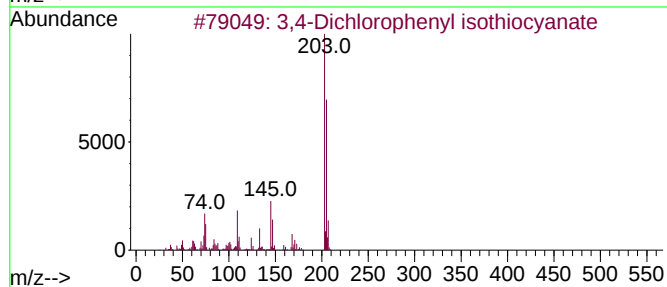
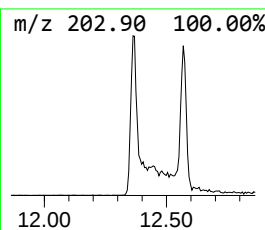
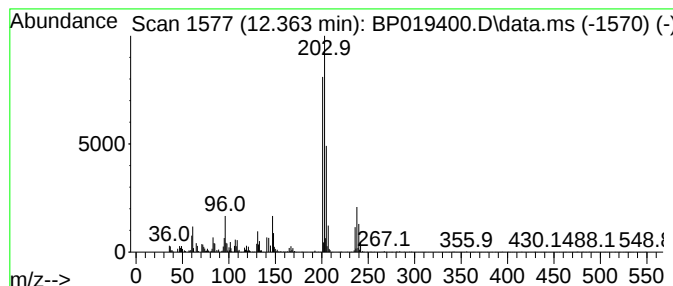
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.363	3.12 ng/ul	187091	Naphthalene-d8	10.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,4-Dichlorophenyl isothiocyanate	203	C7H3Cl2NS	006590-94-9	43
2			p-Toluic acid, .alpha.,.alpha.,....	238	C8H5Cl3O2	005264-40-4	38
3			2,6-Dichlorophenyl isothiocyanate	203	C7H3Cl2NS	006590-95-0	35
4			1-Benzenecarbonyl chloride, 2,4-...	238	C8H5Cl3O2	1000196-85-1	30
5			5-Chlorosulfonyl-2-fluoro-benzoi...	238	C7H4ClFO4S	1000297-26-8	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
Data File : BP019400.D
Acq On : 19 Jan 2024 12:43
Operator : MA/JU
Sample : P1160-04
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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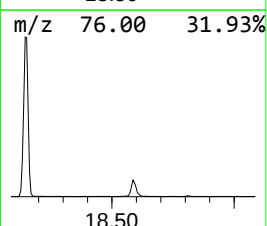
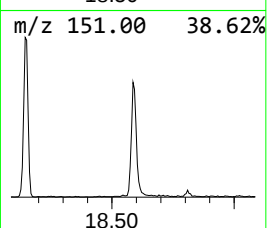
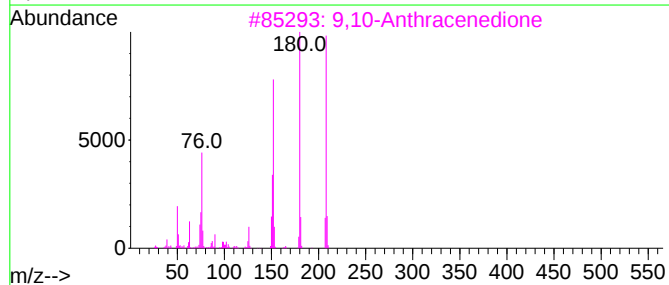
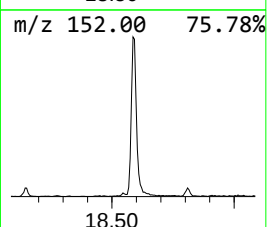
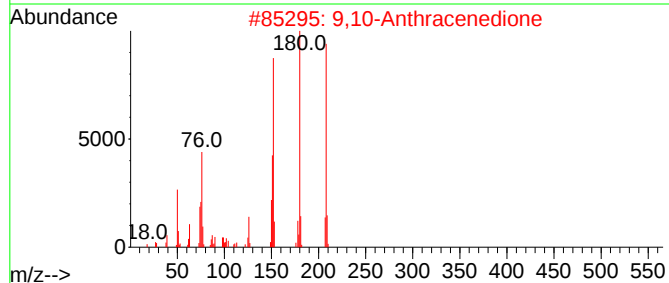
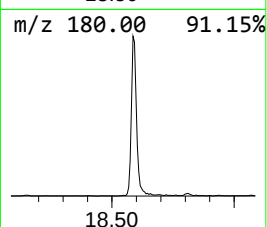
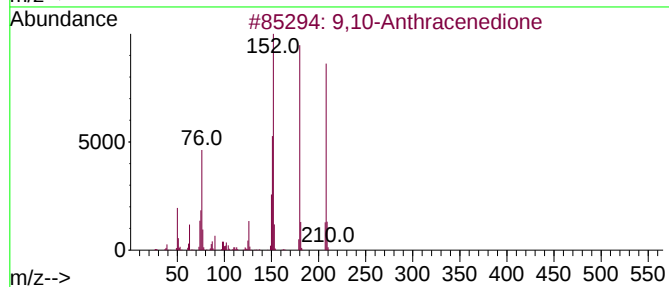
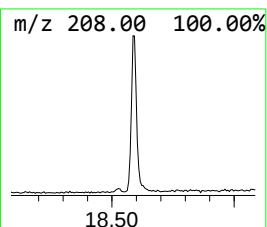
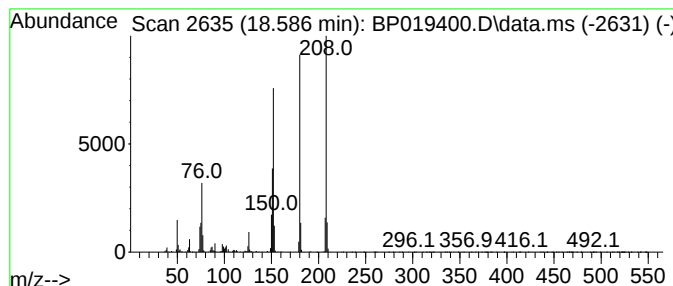
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 9,10-Anthracenedione Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.586	3.42 ng/ul	363378	Phenanthrene-d10	17.180

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
2			9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
3			9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
4			9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
5			Benzo[c]cinnoline	180	C12H8N2	000230-17-1	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
Data File : BP019400.D
Acq On : 19 Jan 2024 12:43
Operator : MA/JU
Sample : P1160-04
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
N-Nitrosodimeth...	3.569	44.0	ng/u1	1860320	1	7.763	844914	20.0
Toluene	3.946	6.5	ng/u1	275909	1	7.763	844914	20.0
Benzene, 1,3-di...	7.651	34.8	ng/u1	1470080	1	7.763	844914	20.0
Benzene, 1,4-di...	8.116	44.9	ng/u1	1897220	1	7.763	844914	20.0
Benzene, 1,3,5-...	10.434	96.5	ng/u1	5778170	2	10.563	1197940	20.0
unknown-01	12.104	2.4	ng/u1	143529	2	10.563	1197940	20.0
unknown-02	12.363	3.1	ng/u1	187091	2	10.563	1197940	20.0
9,10-Anthracene...	18.586	3.4	ng/u1	363378	4	17.180	2127750	20.0