

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
 Data File : BP023160.D
 Acq On : 16 Nov 2024 05:46
 Operator : RC/JU
 Sample : P4803-05
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A0AP8

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-BP110924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023160.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.128	21	24	27	rVV	18090	22455	0.65%	0.070%
2	3.170	27	31	39	rVB	11649	19525	0.56%	0.061%
3	3.552	92	96	117	rBV	98248	193728	5.58%	0.607%
4	4.287	215	221	232	rVB3	19025	34786	1.00%	0.109%
5	4.952	329	334	341	rBV2	16295	26547	0.76%	0.083%
6	6.087	521	527	533	rBV2	7588	14828	0.43%	0.046%
7	6.781	639	645	661	rBV	147647	307204	8.85%	0.962%
8	6.916	662	668	675	rBV	261227	400350	11.53%	1.254%
9	7.122	696	703	718	rVV	322731	574054	16.53%	1.798%
10	7.569	771	779	789	rBV	278895	469303	13.51%	1.470%
11	7.928	834	840	851	rVB	79730	135475	3.90%	0.424%
12	8.052	855	861	865	rBV3	8141	14169	0.41%	0.044%
13	8.211	882	888	890	rBV5	5530	11270	0.32%	0.035%
14	8.281	894	900	921	rVB2	387103	733355	21.12%	2.297%
15	8.487	927	935	943	rVB10	8002	20903	0.60%	0.065%
16	8.593	945	953	957	rBV5	13915	28734	0.83%	0.090%
17	8.646	957	962	967	rVB2	33691	57012	1.64%	0.179%
18	8.716	967	974	984	rBV	342113	623029	17.94%	1.951%
19	8.869	994	1000	1002	rBV4	5897	10903	0.31%	0.034%
20	8.905	1002	1006	1012	rVB9	8906	16717	0.48%	0.052%
21	8.981	1014	1019	1023	rVB7	3961	7984	0.23%	0.025%
22	9.058	1025	1032	1037	rBV5	19104	37390	1.08%	0.117%
23	9.169	1045	1051	1058	rBV3	44361	80781	2.33%	0.253%
24	9.428	1089	1095	1107	rVV	296352	543885	15.66%	1.703%
25	9.522	1107	1111	1117	rVB8	4116	8126	0.23%	0.025%
26	9.575	1117	1120	1123	rBV3	7144	10071	0.29%	0.032%
27	9.728	1135	1146	1157	rBV6	47584	154489	4.45%	0.484%
28	9.969	1180	1187	1202	rBV	456543	945861	27.24%	2.963%
29	10.116	1208	1212	1215	rBV6	4623	8077	0.23%	0.025%
30	10.181	1219	1223	1227	rVB2	13229	19770	0.57%	0.062%
31	10.322	1240	1247	1252	rBV	350494	622449	17.92%	1.950%
32	10.481	1268	1274	1287	rVV	207806	445695	12.83%	1.396%
33	10.599	1289	1294	1303	rVB10	17863	48665	1.40%	0.152%
34	10.746	1315	1319	1323	rVB7	5918	10978	0.32%	0.034%
35	10.899	1340	1345	1348	rBV5	8768	16838	0.48%	0.053%
36	11.052	1366	1371	1374	rBV	26403	38063	1.10%	0.119%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Title : SVOA CALIBRATION

37	11.169	1389	1391	1396	rVB5	7120	9630	0.28%	0.030%
38	11.304	1409	1414	1418	rBV6	9587	18627	0.54%	0.058%
39	11.352	1418	1422	1427	rVV7	15949	30105	0.87%	0.094%
40	11.604	1461	1465	1469	rBV6	5828	10261	0.30%	0.032%
41	11.675	1470	1477	1484	rBV	105811	179399	5.17%	0.562%
42	11.728	1484	1486	1490	rBV5	8246	9682	0.28%	0.030%
43	11.775	1490	1494	1497	rBV6	8641	13234	0.38%	0.041%
44	11.875	1505	1511	1516	rBV9	11981	27234	0.78%	0.085%
45	11.922	1516	1519	1522	rVV5	8447	13071	0.38%	0.041%
46	11.963	1522	1526	1533	rVV6	18761	35341	1.02%	0.111%
47	12.057	1534	1542	1547	rVV6	12191	36769	1.06%	0.115%
48	12.110	1547	1551	1556	rVB6	24210	41394	1.19%	0.130%
49	12.204	1562	1567	1574	rVB2	50147	93208	2.68%	0.292%
50	12.334	1585	1589	1594	rBV5	13100	26016	0.75%	0.081%
51	12.387	1594	1598	1602	rVV6	15416	25650	0.74%	0.080%
52	12.422	1602	1604	1610	rVV7	7707	13778	0.40%	0.043%
53	12.475	1610	1613	1618	rVV7	6696	10787	0.31%	0.034%
54	12.575	1625	1630	1636	rVB2	56151	96191	2.77%	0.301%
55	12.787	1663	1666	1672	rVB7	12689	20280	0.58%	0.064%
56	12.887	1680	1683	1690	rBV9	6862	14296	0.41%	0.045%
57	12.987	1695	1700	1703	rVV7	18562	37112	1.07%	0.116%
58	13.040	1705	1709	1719	rVB3	31737	65766	1.89%	0.206%
59	13.263	1744	1747	1753	rVB5	19746	32462	0.93%	0.102%
60	13.340	1755	1760	1763	rBV6	14142	26594	0.77%	0.083%
61	13.381	1764	1767	1771	rVV6	17204	22872	0.66%	0.072%
62	13.428	1771	1775	1779	rVV4	26609	43402	1.25%	0.136%
63	13.469	1779	1782	1784	rVV	17470	20177	0.58%	0.063%
64	13.499	1785	1787	1792	rVB3	17417	26053	0.75%	0.082%
65	13.593	1797	1803	1813	rVB	908809	1441156	41.50%	4.514%
66	13.710	1814	1823	1827	rBV	17566	58216	1.68%	0.182%
67	13.751	1827	1830	1831	rBV3	7509	8741	0.25%	0.027%
68	13.834	1841	1844	1846	rBV3	15350	17141	0.49%	0.054%
69	13.875	1846	1851	1861	rVB	1085330	1540069	44.35%	4.824%
70	14.016	1872	1875	1884	rVB2	19787	40942	1.18%	0.128%
71	14.122	1891	1893	1898	rVB5	12176	16908	0.49%	0.053%
72	14.193	1898	1905	1912	rBV	635624	960344	27.65%	3.008%
73	14.246	1912	1914	1920	rVB2	25792	38070	1.10%	0.119%
74	14.404	1938	1941	1945	rBV6	8541	12852	0.37%	0.040%
75	14.487	1950	1955	1969	rVV4	87858	265501	7.65%	0.832%
76	14.604	1972	1975	1983	rVB5	37617	69922	2.01%	0.219%

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77	15.122	2060	2063	2068	rBV6	18538	30778	0.89%	0.096%
78	15.210	2073	2078	2083	rVV	1642543	2126773	61.24%	6.661%
79	15.257	2083	2086	2090	rVB2	45484	58177	1.68%	0.182%
80	15.351	2097	2102	2114	rBV	425019	724818	20.87%	2.270%
81	15.451	2115	2119	2122	rVV	62817	87241	2.51%	0.273%
82	15.493	2122	2126	2133	rVV4	35808	70750	2.04%	0.222%
83	15.581	2137	2141	2149	rVB	75154	127929	3.68%	0.401%
84	15.745	2163	2169	2182	rVB3	67366	179159	5.16%	0.561%
85	15.945	2198	2203	2207	rBV	293885	453695	13.06%	1.421%
86	16.275	2254	2259	2266	rBV2	108947	192186	5.53%	0.602%
87	17.004	2377	2383	2388	rBV	862327	1233713	35.53%	3.864%
88	17.092	2393	2398	2412	rVB	1709025	2494376	71.83%	7.813%
89	17.934	2537	2541	2547	rBV4	98515	158554	4.57%	0.497%
90	18.157	2575	2579	2586	rVB	51243	84261	2.43%	0.264%
91	18.345	2608	2611	2614	rBV3	33036	41269	1.19%	0.129%
92	18.734	2673	2677	2681	rBV5	34640	62049	1.79%	0.194%
93	18.916	2704	2708	2713	rVB	68571	107322	3.09%	0.336%
94	19.281	2766	2770	2777	rVB4	60312	87353	2.52%	0.274%
95	19.492	2800	2806	2813	rBV	2195110	3272500	94.24%	10.250%
96	19.563	2813	2818	2828	rVV	759808	1239179	35.68%	3.881%
97	20.586	2989	2992	2995	rVB2	80110	82870	2.39%	0.260%
98	21.439	3132	3137	3147	rVB2	1107903	1645859	47.40%	5.155%
99	24.433	3638	3646	3659	rVB	1406275	3472635	100.00%	10.877%
100	24.657	3676	3684	3703	rVB	552138	1711419	49.28%	5.360%

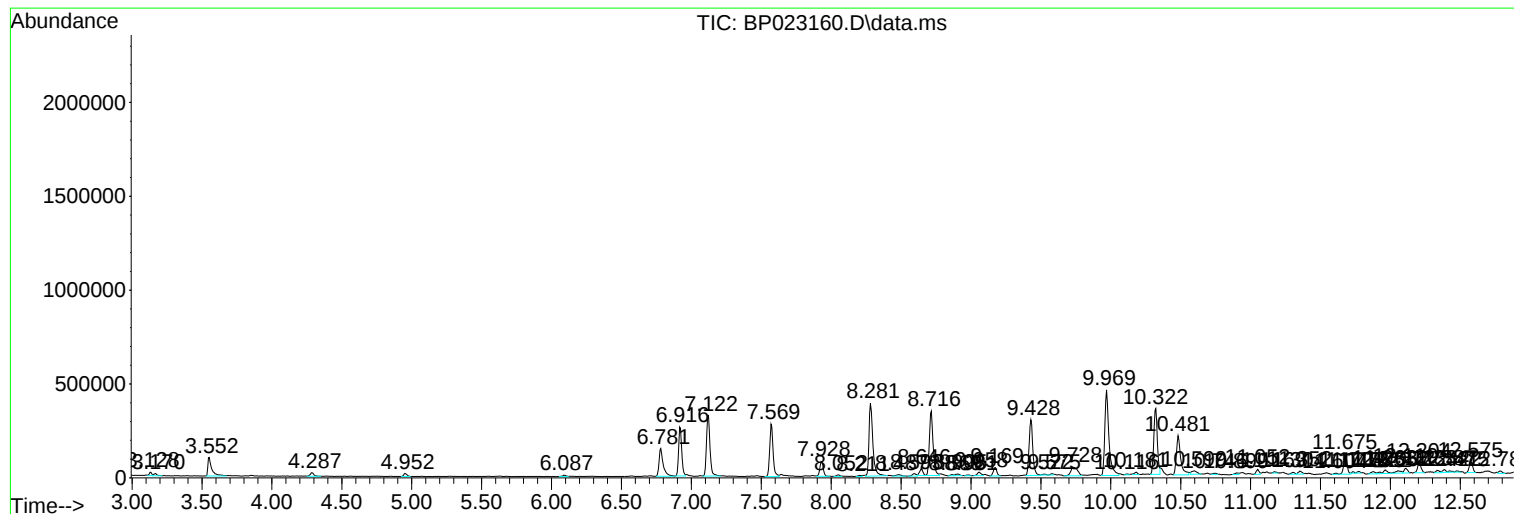
Sum of corrected areas: 31927587

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ALS Vial : 29 Sample Multiplier: 1

Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
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Instrument :
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A0AP8

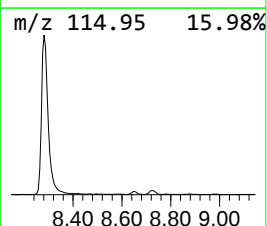
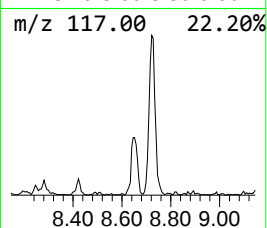
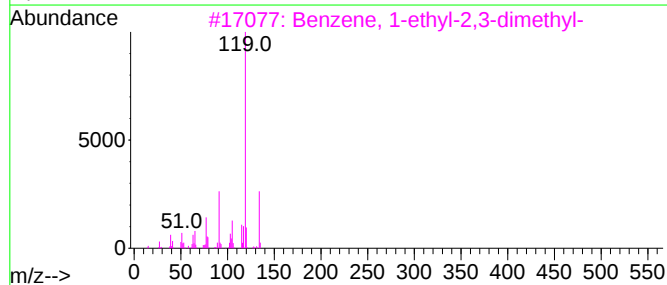
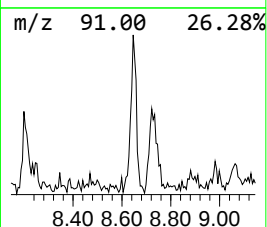
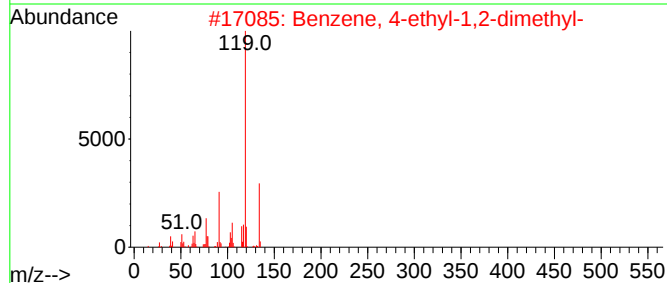
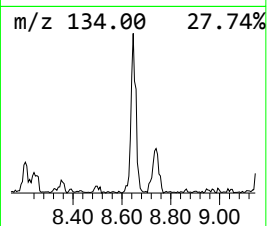
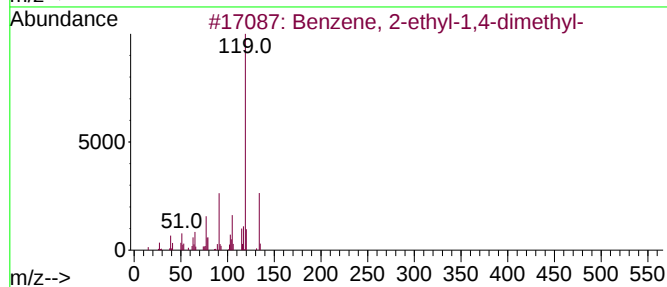
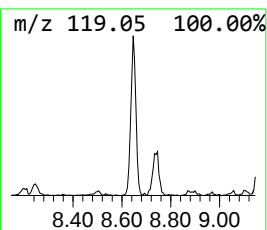
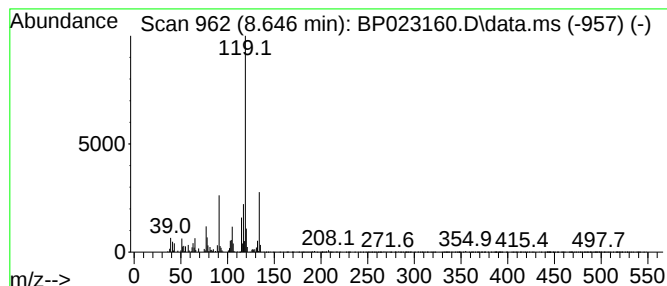
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.646	2.43 ng/ul	57012	1,4-Dichlorobenzene-d4	7.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
4			p-Cymene	134	C10H14	000099-87-6	87
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	87



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ClientSampleId :
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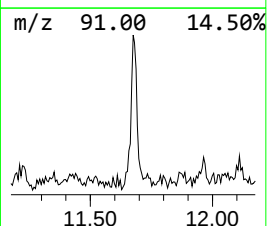
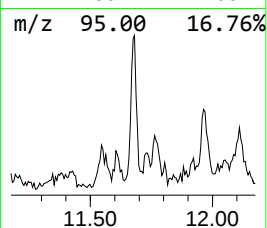
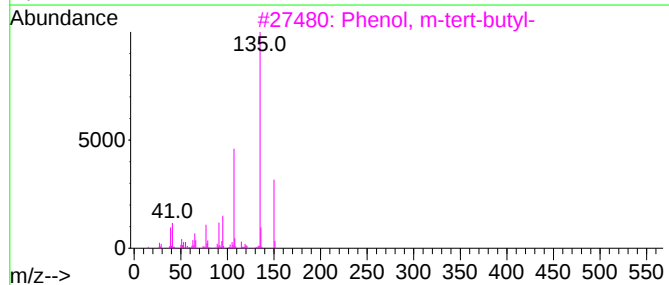
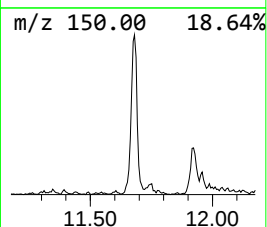
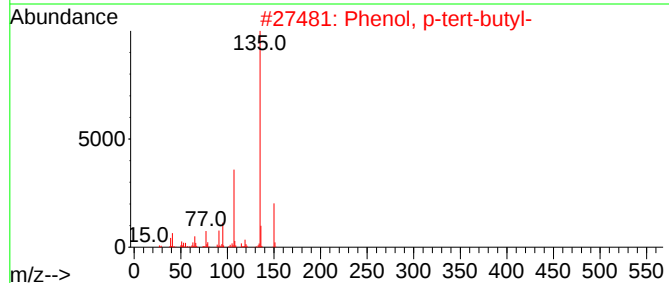
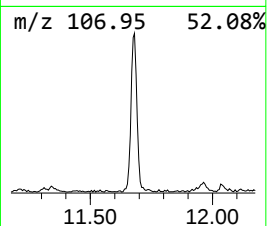
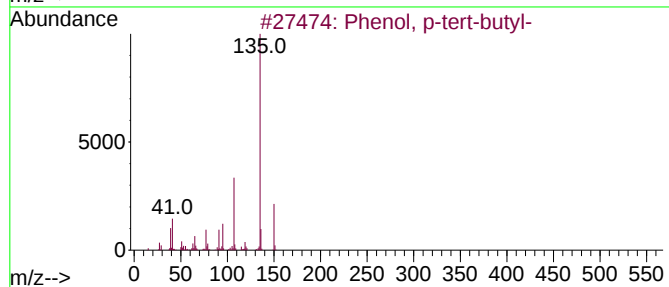
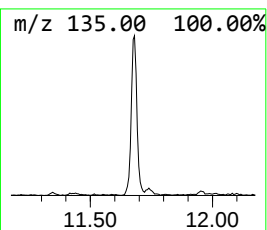
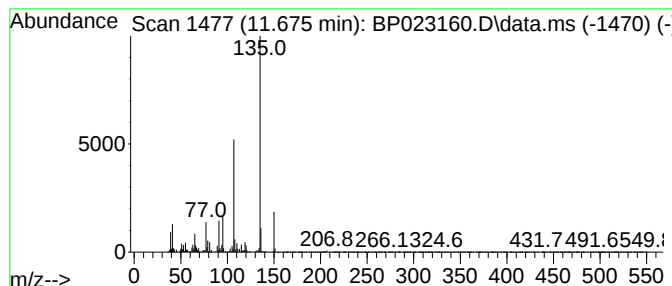
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phenol, p-tert-butyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.675	5.76 ng/ul	179399	Naphthalene-d8	10.322

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	97
2			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	95
3			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
4			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
5			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91



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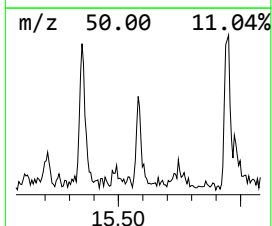
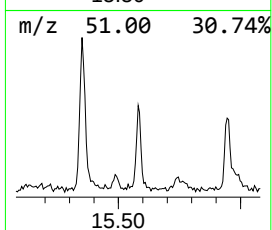
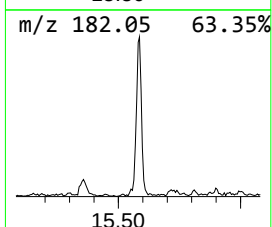
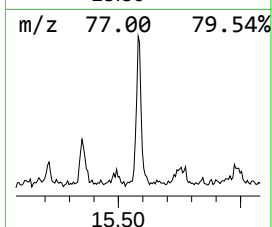
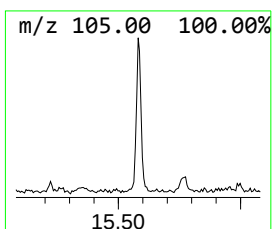
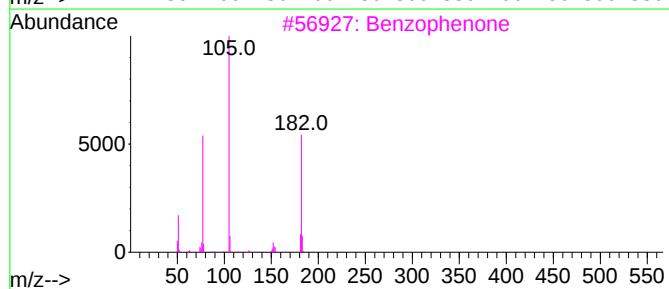
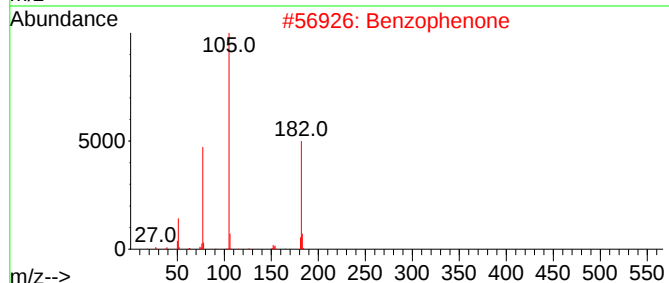
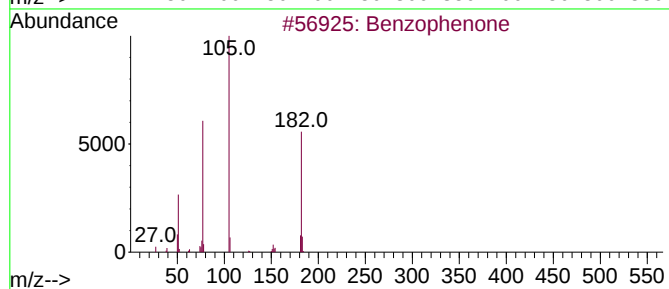
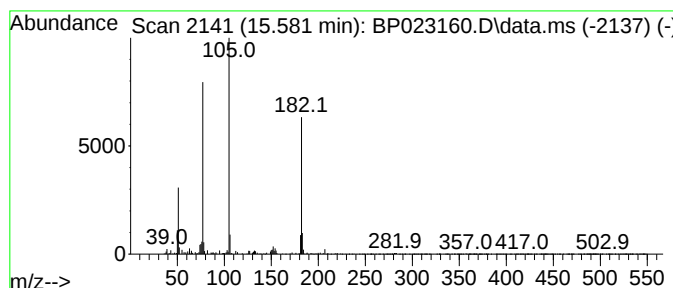
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzophenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.581	2.66 ng/ul	127929	Acenaphthene-d10	14.193

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	91
3			Benzophenone	182	C13H10O	000119-61-9	90
4			Benzophenone	182	C13H10O	000119-61-9	87
5			Benzophenone	182	C13H10O	000119-61-9	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023160.D
Acq On : 16 Nov 2024 05:46
Operator : RC/JU
Sample : P4803-05
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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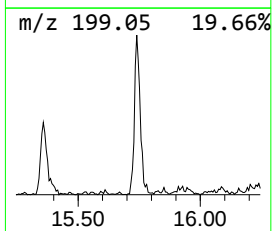
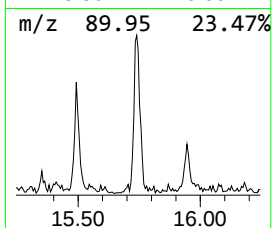
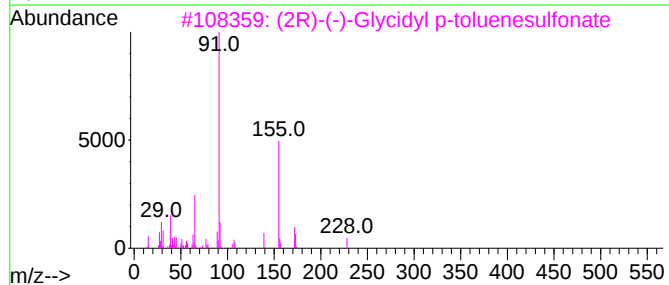
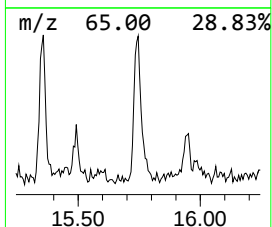
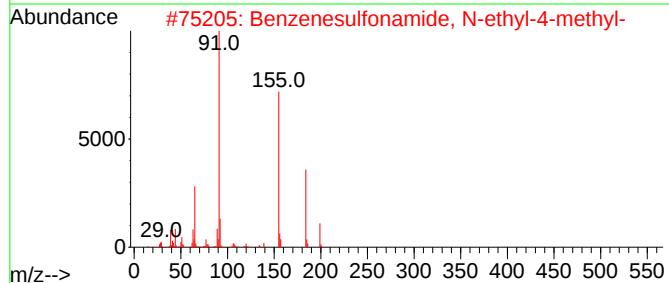
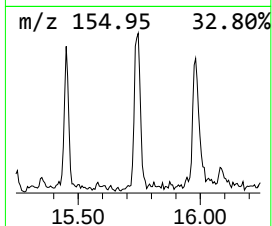
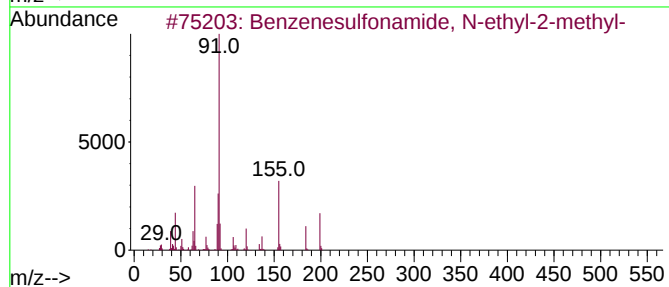
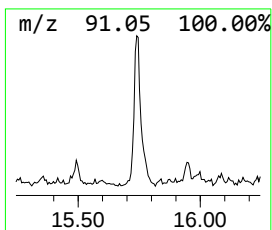
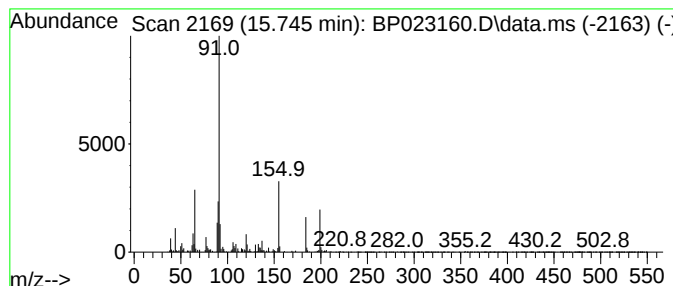
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzenesulfonamide, N-ethyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.745	2.90 ng/ul	179159	Phenanthrene-d10	17.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	96
2			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	53
3			(2R)-(-)-Glycidyl p-toluenesulfo...	228	C10H12O4S	113826-06-5	53
4			Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	52
5			Benzenesulfonamide, N,N'-(dithio...	460	C18H24N2O4S4	023516-74-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023160.D
Acq On : 16 Nov 2024 05:46
Operator : RC/JU
Sample : P4803-05
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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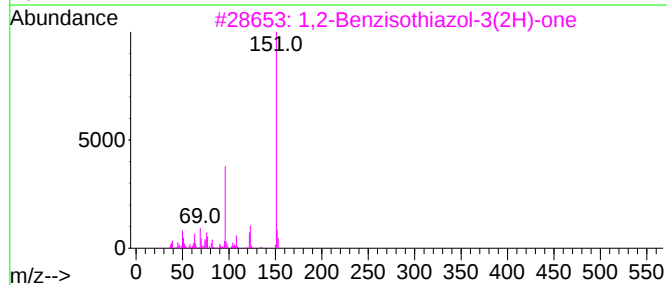
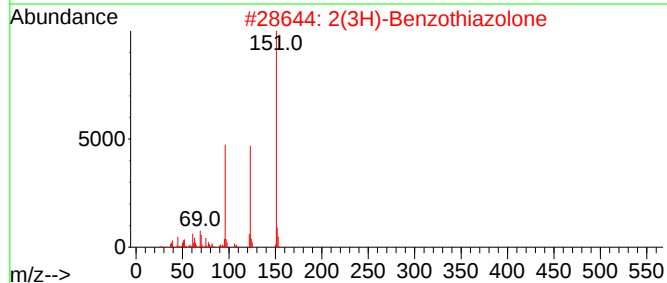
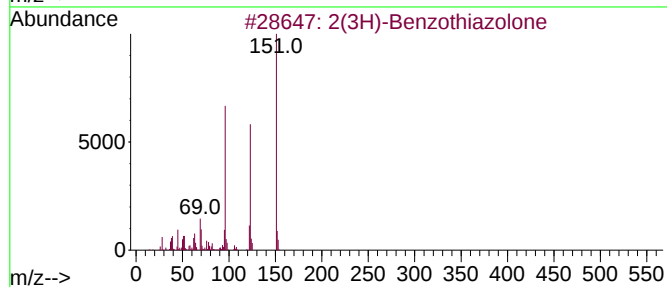
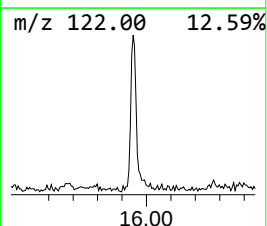
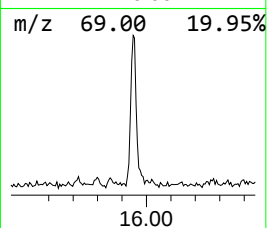
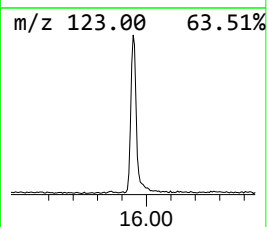
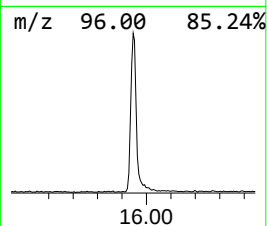
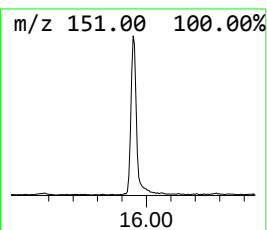
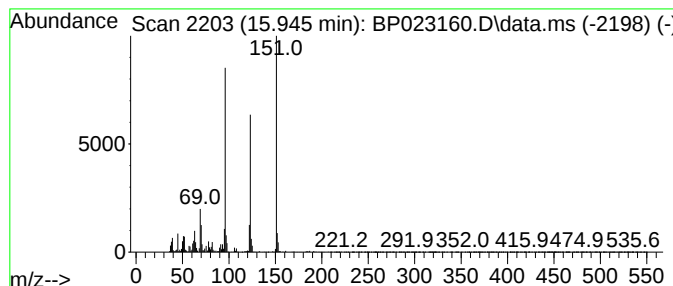
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2(3H)-Benzothiazolone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.945	7.35 ng/ul	453695	Phenanthrene-d10	17.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	96
2			2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	94
3			1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	59
4			Thieno[3,2-b]pyridin-7-ol	151	C7H5NOS	107818-20-2	49
5			Thieno[2,3-b]pyridine-N-oxide	151	C7H5NOS	025557-50-0	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023160.D
Acq On : 16 Nov 2024 05:46
Operator : RC/JU
Sample : P4803-05
Misc :
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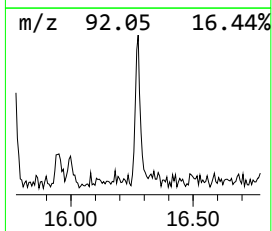
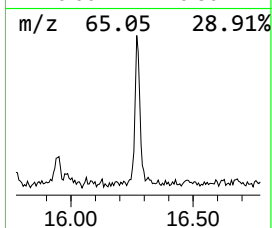
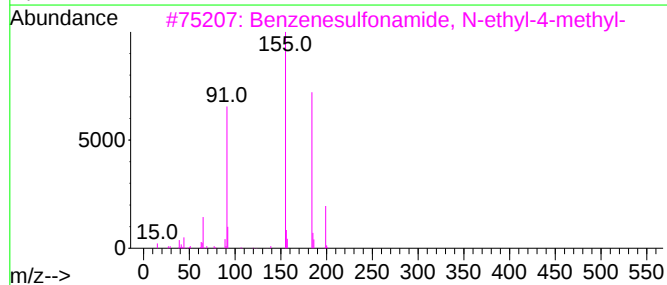
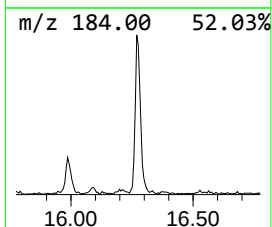
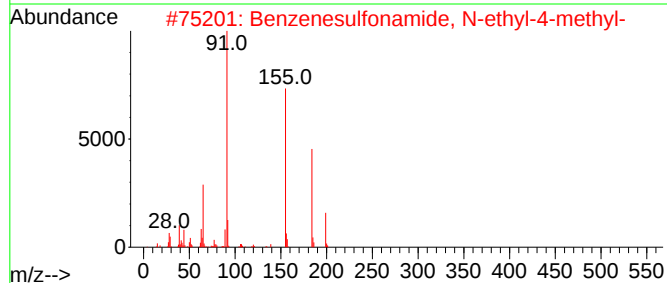
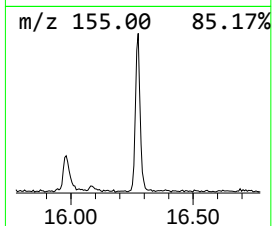
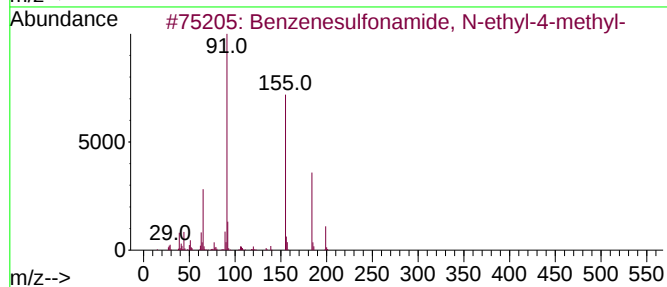
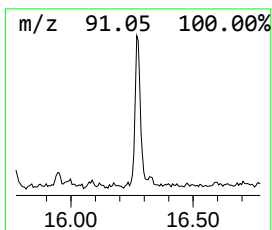
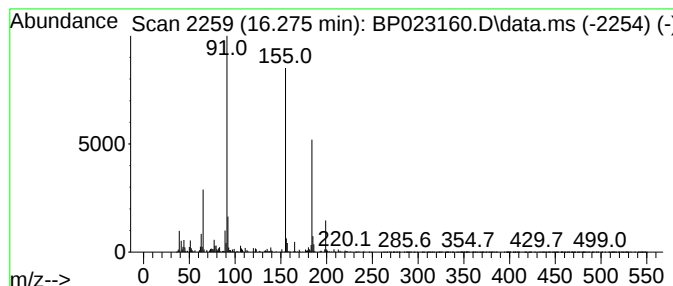
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzenesulfonamide, N-ethyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.275	3.12 ng/ul	192186	Phenanthrene-d10	17.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	95
2			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	95
3			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	91
4			Acetic acid, 2-(4-methylphenylsu...	243	C10H13NO4S	002645-02-5	86
5			N-isobutyl-4-methylbenzenesulfon...	227	C11H17NO2S	023705-38-6	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023160.D
Acq On : 16 Nov 2024 05:46
Operator : RC/JU
Sample : P4803-05
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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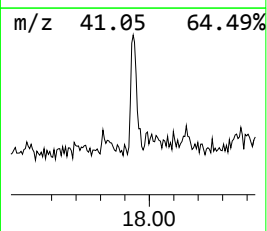
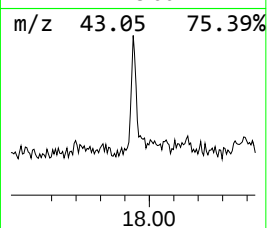
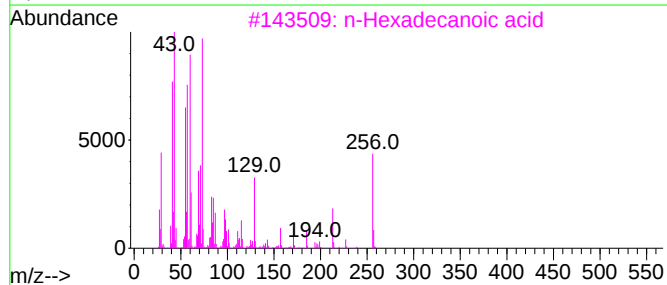
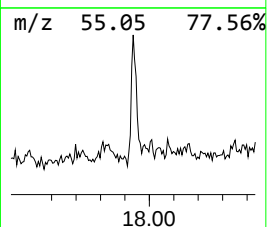
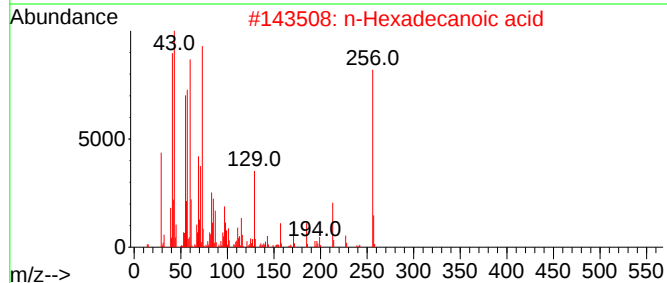
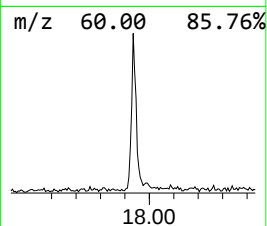
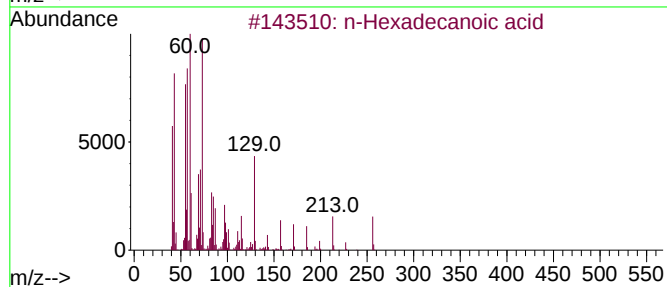
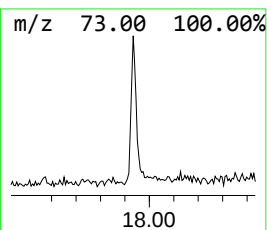
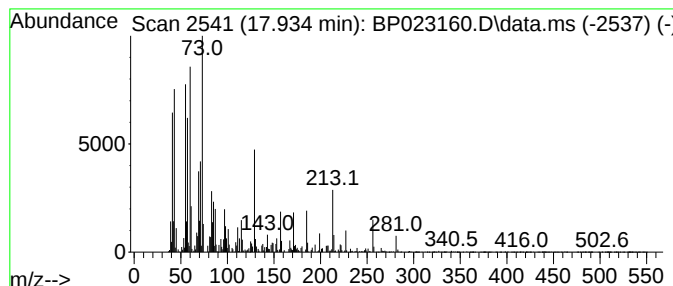
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.934	2.57 ng/ul	158554	Phenanthrene-d10	17.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023160.D
Acq On : 16 Nov 2024 05:46
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Sample : P4803-05
Misc :
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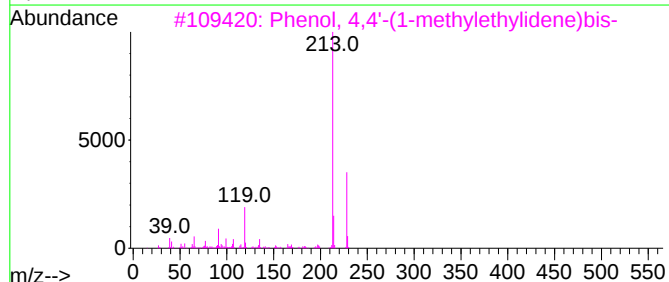
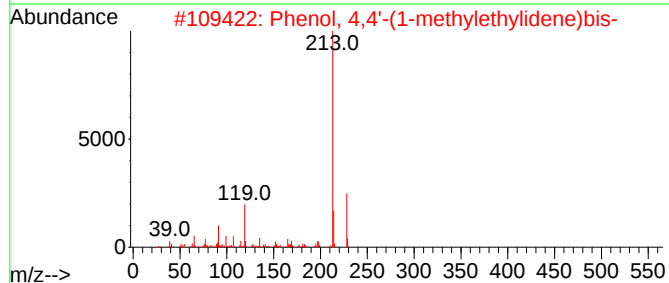
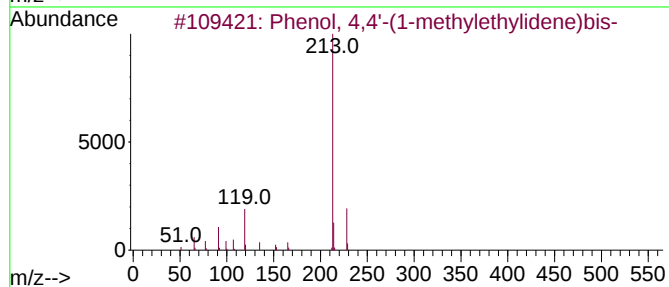
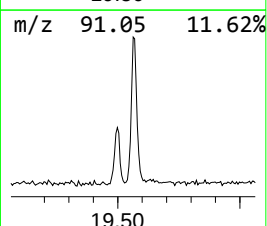
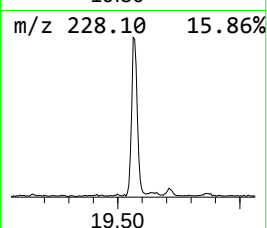
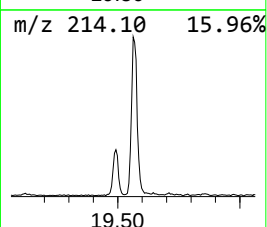
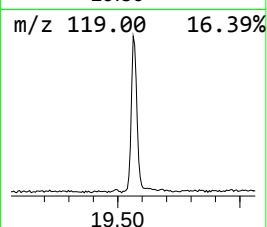
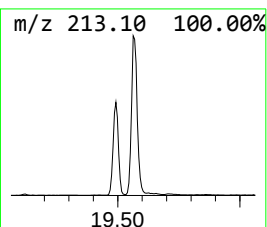
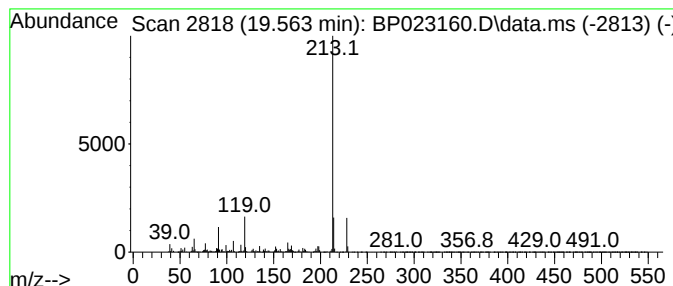
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Phenol, 4,4'-(1-methylethyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.563	15.06 ng/ul	1239180	Chrysene-d12	21.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	97
2			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	93
3			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	90
4			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	87
5			Diphenolic acid	286	C17H18O4	000126-00-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023160.D
Acq On : 16 Nov 2024 05:46
Operator : RC/JU
Sample : P4803-05
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Phenol, p-tert-...	11.675	5.8	ng/ul	179399	2	10.322	622449	20.0
Benzophenone	15.581	2.7	ng/ul	127929	3	14.193	960344	20.0
Benzenesulfonam...	15.745	2.9	ng/ul	179159	4	17.004	1233710	20.0
2(3H)-Benzothia...	15.945	7.3	ng/ul	453695	4	17.004	1233710	20.0
Benzenesulfonam...	16.275	3.1	ng/ul	192186	4	17.004	1233710	20.0
n-Hexadecanoic ...	17.934	2.6	ng/ul	158554	4	17.004	1233710	20.0
Phenol, 4,4'-(1...	19.563	15.1	ng/ul	1239180	5	21.439	1645860	20.0