

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020725\
 Data File : BP023999.D
 Acq On : 07 Feb 2025 12:28
 Operator : RC/JU
 Sample : Q1229-11DL 20X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E29B5DL

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-BP012925.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023999.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.728	106	109	117	rBV2	16791	31303	0.43%	0.065%
2	4.023	155	159	164	rVB	19902	28258	0.39%	0.058%
3	6.946	653	656	668	rVB4	12348	30960	0.42%	0.064%
4	7.099	677	682	689	rVB	42682	69105	0.95%	0.143%
5	7.305	710	717	725	rBV	37369	67128	0.92%	0.138%
6	7.764	787	795	804	rBV	1303607	1916584	26.26%	3.954%
7	7.834	804	807	815	rVV9	6533	12248	0.17%	0.025%
8	8.469	910	915	924	rBV2	29672	50635	0.69%	0.104%
9	8.911	983	990	999	rBV3	43033	83445	1.14%	0.172%
10	8.999	999	1005	1008	rBV7	7442	12668	0.17%	0.026%
11	9.169	1029	1034	1039	rBV	27665	41934	0.57%	0.087%
12	9.216	1039	1042	1048	rVB4	16631	22991	0.31%	0.047%
13	9.346	1060	1064	1069	rVB8	5319	8663	0.12%	0.018%
14	9.475	1079	1086	1089	rBV6	7535	13791	0.19%	0.028%
15	9.516	1089	1093	1099	rVB	26112	40700	0.56%	0.084%
16	9.622	1106	1111	1118	rBV	25297	42098	0.58%	0.087%
17	9.722	1120	1128	1131	rBV	769693	1331307	18.24%	2.746%
18	9.752	1131	1133	1150	rVB	525076	771064	10.56%	1.591%
19	9.958	1164	1168	1174	rVB7	9997	14940	0.20%	0.031%
20	10.028	1174	1180	1186	rVV	105932	179971	2.47%	0.371%
21	10.087	1188	1190	1195	rVV6	5384	8390	0.11%	0.017%
22	10.169	1197	1204	1213	rVB	49127	87491	1.20%	0.180%
23	10.358	1233	1236	1241	rBV4	6129	9828	0.13%	0.020%
24	10.458	1246	1253	1260	rVV	4124217	6372992	87.31%	13.147%
25	10.534	1260	1266	1273	rVV	1664267	2888127	39.57%	5.958%
26	10.593	1273	1276	1283	rVV6	24463	50254	0.69%	0.104%
27	10.675	1284	1290	1302	rVB2	67332	146325	2.00%	0.302%
28	10.810	1308	1313	1322	rVB	21970	37006	0.51%	0.076%
29	10.905	1322	1329	1338	rVV	84541	152168	2.08%	0.314%
30	10.987	1338	1343	1350	rVV2	39320	74313	1.02%	0.153%
31	11.081	1352	1359	1366	rVV	199479	352493	4.83%	0.727%
32	11.152	1366	1371	1382	rVV	117237	221902	3.04%	0.458%
33	11.234	1382	1385	1390	rVB7	4173	7549	0.10%	0.016%
34	11.369	1402	1408	1415	rVB	34659	61061	0.84%	0.126%
35	11.452	1415	1422	1427	rBV9	4360	9900	0.14%	0.020%
36	11.557	1433	1440	1446	rBV	149867	250690	3.43%	0.517%

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37	11.616	1446	1450	1458	rVV	75907	134944	1.85%	0.278%
38	11.716	1462	1467	1470	rBV2	6990	10848	0.15%	0.022%
39	11.875	1482	1494	1510	rBV	1345704	2344566	32.12%	4.837%
40	11.987	1510	1513	1518	rVB2	12506	19011	0.26%	0.039%
41	12.052	1518	1524	1530	rBV	39275	82986	1.14%	0.171%
42	12.099	1530	1532	1536	rVB	14020	17858	0.24%	0.037%
43	12.175	1543	1545	1552	rVB6	7202	9860	0.14%	0.020%
44	12.234	1552	1555	1558	rBV	20436	28425	0.39%	0.059%
45	12.263	1558	1560	1565	rVV	14584	24166	0.33%	0.050%
46	12.346	1569	1574	1578	rBV	35036	50949	0.70%	0.105%
47	12.416	1583	1586	1592	rVB2	23347	32166	0.44%	0.066%
48	12.469	1592	1595	1600	rVB2	9603	12714	0.17%	0.026%
49	12.528	1600	1605	1609	rBV3	15190	24499	0.34%	0.051%
50	12.575	1609	1613	1618	rBV2	30054	55818	0.76%	0.115%
51	12.622	1618	1621	1623	rBV	9937	11470	0.16%	0.024%
52	12.675	1628	1630	1633	rVV	12739	13215	0.18%	0.027%
53	12.716	1633	1637	1639	rVV2	19342	25609	0.35%	0.053%
54	12.757	1639	1644	1648	rVV3	76876	132808	1.82%	0.274%
55	12.793	1648	1650	1658	rVB2	27811	35187	0.48%	0.073%
56	12.899	1663	1668	1672	rBV7	7372	14708	0.20%	0.030%
57	12.975	1674	1681	1689	rVB4	17329	35440	0.49%	0.073%
58	13.057	1689	1695	1701	rBV5	8392	21341	0.29%	0.044%
59	13.157	1708	1712	1716	rBV6	13144	18739	0.26%	0.039%
60	13.222	1716	1723	1740	rBV	4729960	7299378	100.00%	15.058%
61	13.422	1755	1757	1766	rVB5	11752	23398	0.32%	0.048%
62	13.499	1766	1770	1775	rBV4	10086	16512	0.23%	0.034%
63	13.593	1775	1786	1790	rVV	1691566	2300213	31.51%	4.745%
64	13.634	1790	1793	1801	rVV	510622	804787	11.03%	1.660%
65	13.704	1801	1805	1811	rVV	100214	172143	2.36%	0.355%
66	13.787	1814	1819	1824	rVV	115373	183378	2.51%	0.378%
67	13.840	1826	1828	1838	rVB4	15233	28706	0.39%	0.059%
68	14.075	1855	1868	1876	rBV	118138	207465	2.84%	0.428%
69	14.151	1876	1881	1885	rVB6	8576	11706	0.16%	0.024%
70	14.298	1902	1906	1912	rVB8	9135	13589	0.19%	0.028%
71	14.387	1912	1921	1938	rBV	2393917	3732607	51.14%	7.700%
72	14.628	1957	1962	1975	rVB	12077	42576	0.58%	0.088%
73	15.069	2031	2037	2044	rBV2	27822	41124	0.56%	0.085%
74	15.198	2055	2059	2061	rBV2	6624	9738	0.13%	0.020%
75	15.263	2067	2070	2075	rBV7	7157	11805	0.16%	0.024%
76	15.393	2086	2092	2099	rBV	174294	264491	3.62%	0.546%

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Title : SVOA CALIBRATION

77	15.516	2108	2113	2119	rBV6	18356	33228	0.46%	0.069%
78	15.763	2150	2155	2162	rBV	118188	179151	2.45%	0.370%
79	16.345	2250	2254	2259	rBV3	13898	22279	0.31%	0.046%
80	16.963	2355	2359	2365	rBV3	12137	21737	0.30%	0.045%
81	17.192	2390	2398	2408	rBV2	2680125	4476714	61.33%	9.235%
82	17.292	2409	2415	2428	rVB	205296	323732	4.44%	0.668%
83	18.128	2553	2557	2565	rBV5	22745	39977	0.55%	0.082%
84	19.657	2812	2817	2826	rBV	253176	364988	5.00%	0.753%
85	21.622	3146	3151	3162	rBV2	3134484	4717389	64.63%	9.731%
86	24.986	3715	3723	3737	rVB	1719522	4475133	61.31%	9.232%

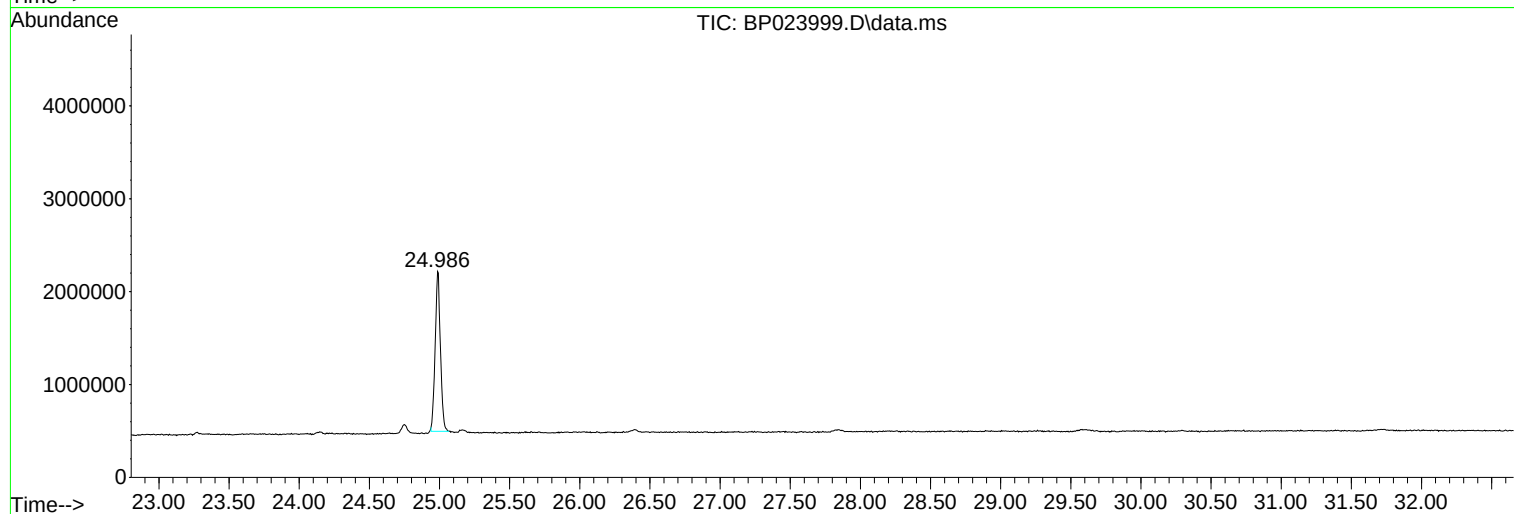
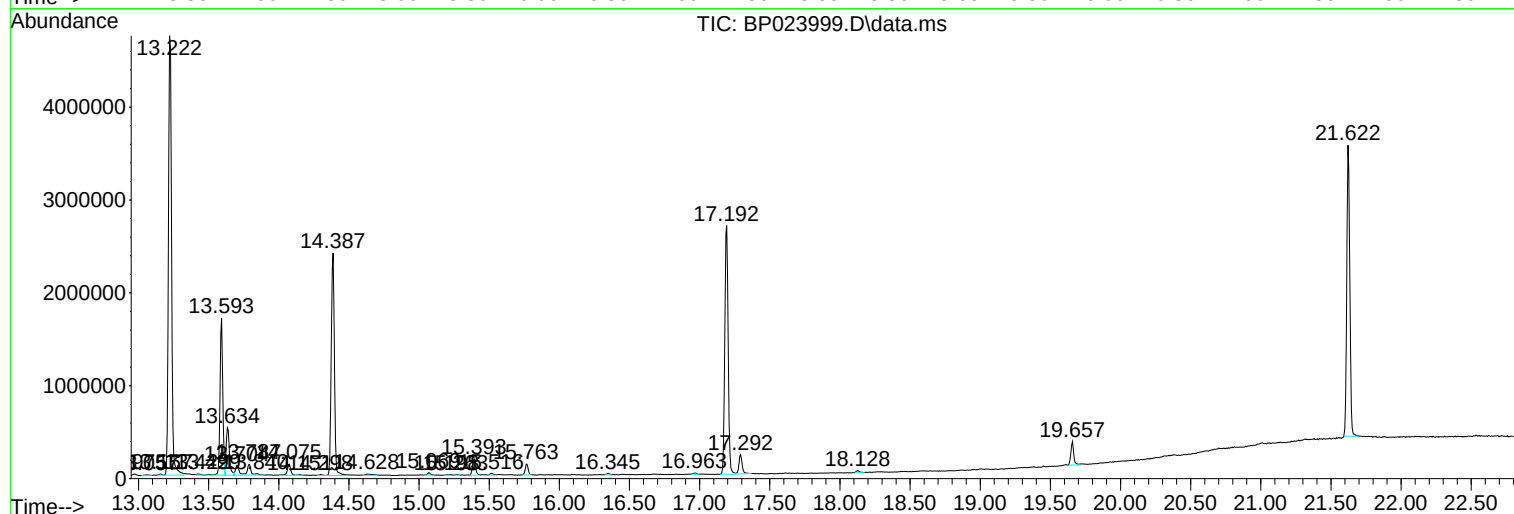
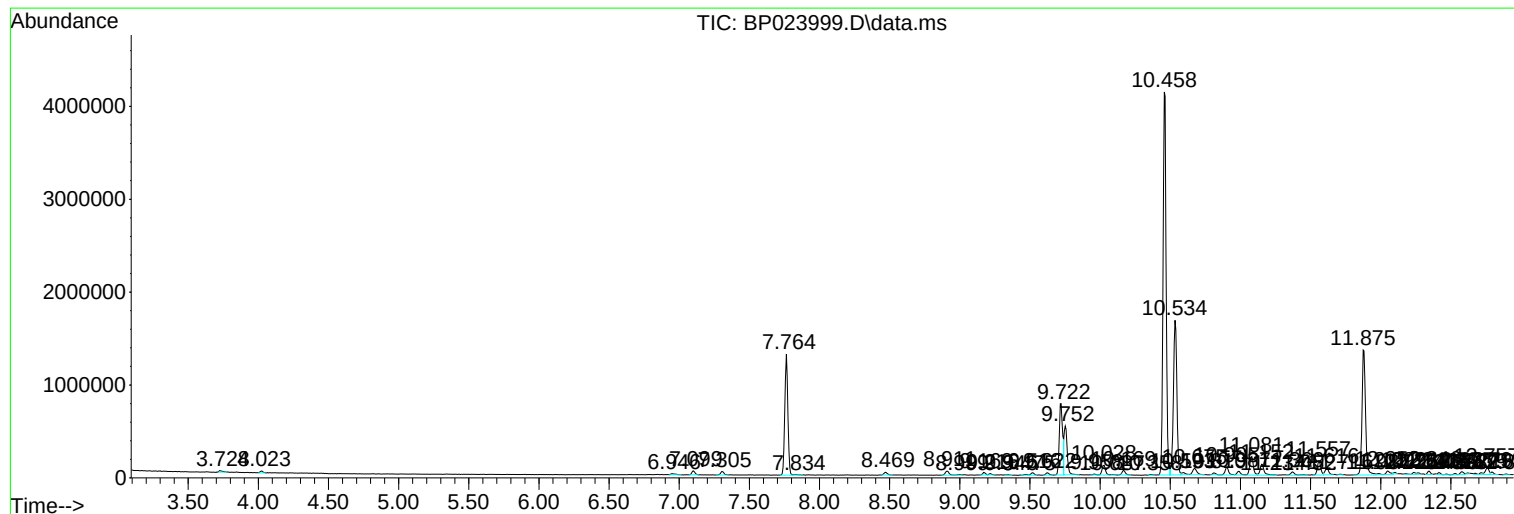
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Instrument :
BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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



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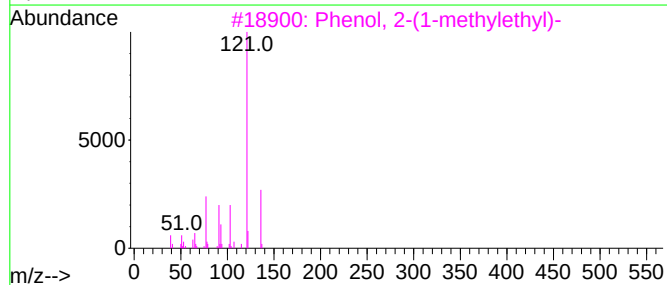
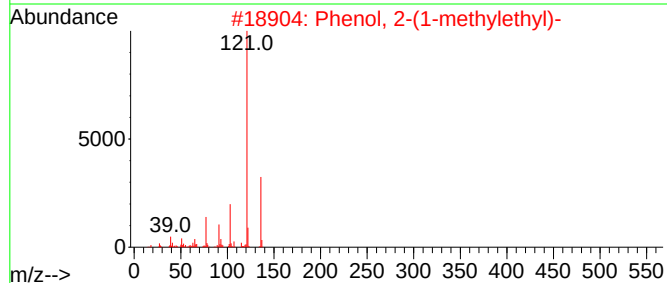
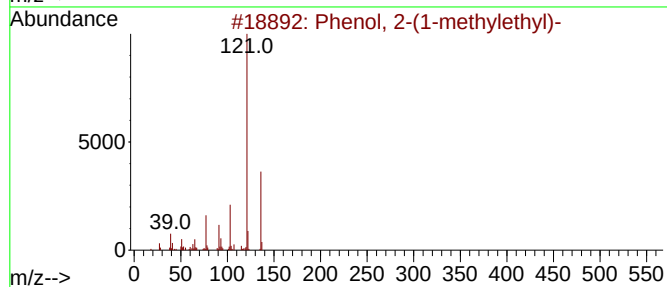
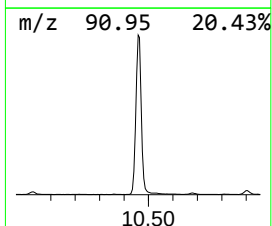
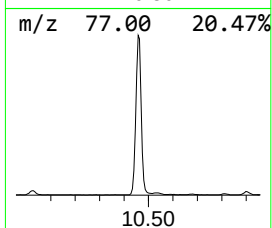
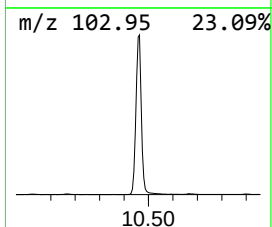
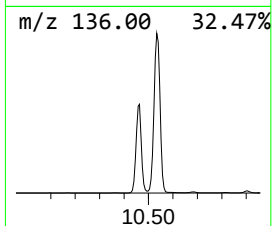
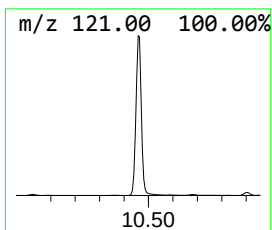
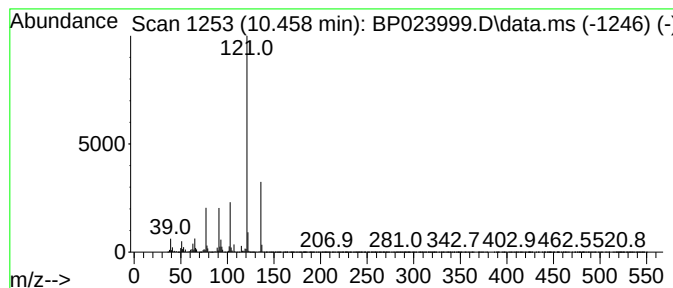
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Phenol, 2-(1-methylethyl)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.458	44.13 ng/ul	6372990	Naphthalene-d8	10.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	95
2			Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	95
3			Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	94
4			Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	90
5			Phenol, 2-ethyl-5-methyl-	136	C9H12O	001687-61-2	87



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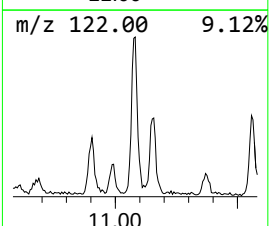
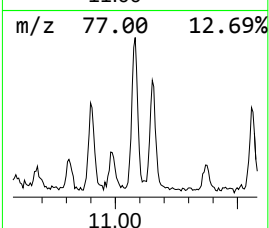
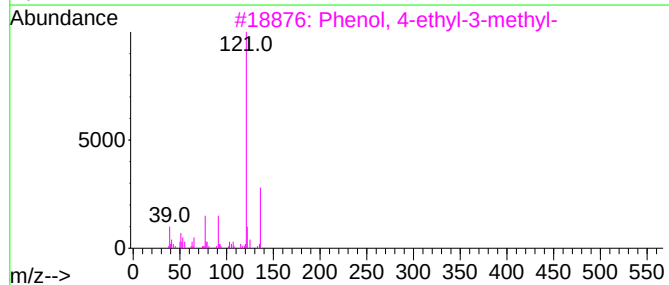
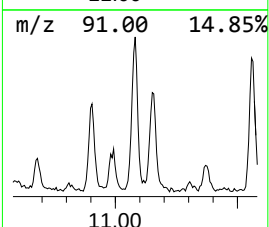
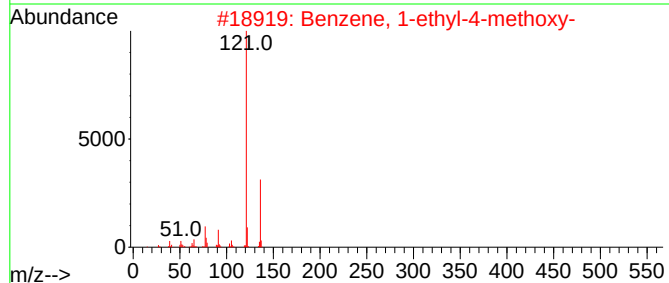
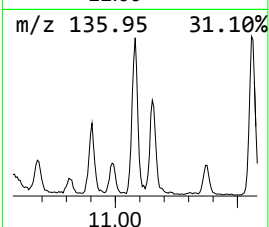
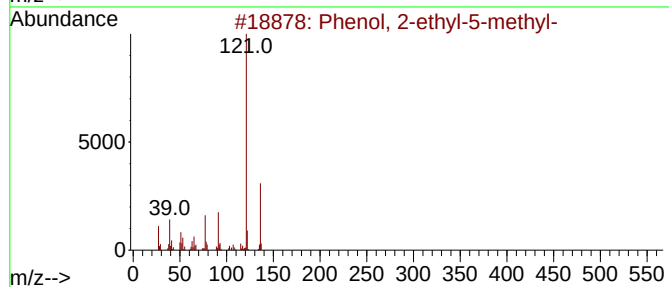
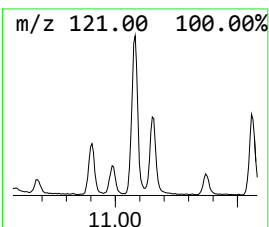
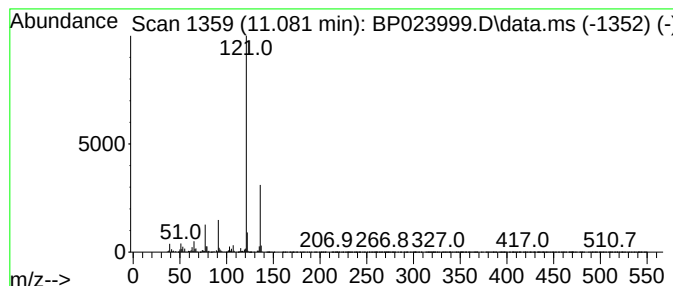
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenol, 2-ethyl-5-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.081	2.44 ng/ul	352493	Naphthalene-d8	10.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2-ethyl-5-methyl-	136	C9H12O	001687-61-2	91
2			Benzene, 1-ethyl-4-methoxy-	136	C9H12O	001515-95-3	91
3			Phenol, 4-ethyl-3-methyl-	136	C9H12O	001123-94-0	91
4			p-Cumenol	136	C9H12O	000099-89-8	91
5			Phenol, 4-ethyl-2-methyl-	136	C9H12O	002219-73-0	91



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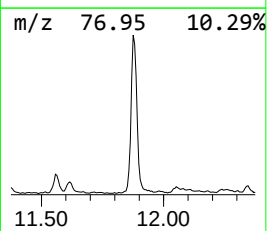
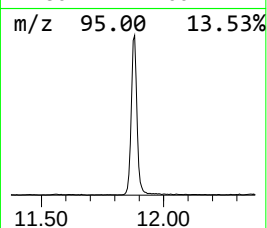
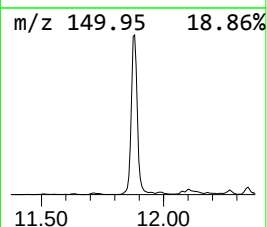
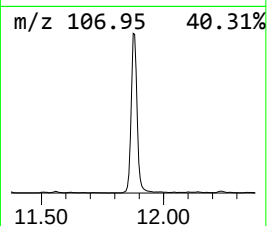
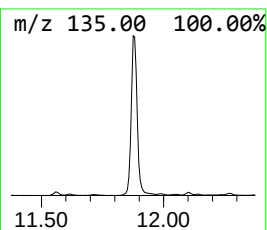
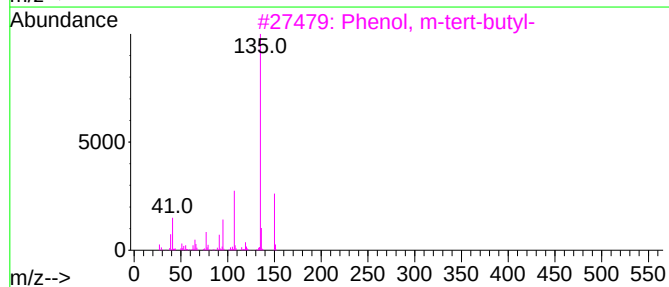
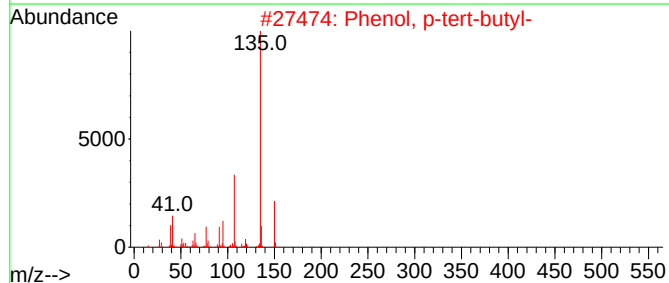
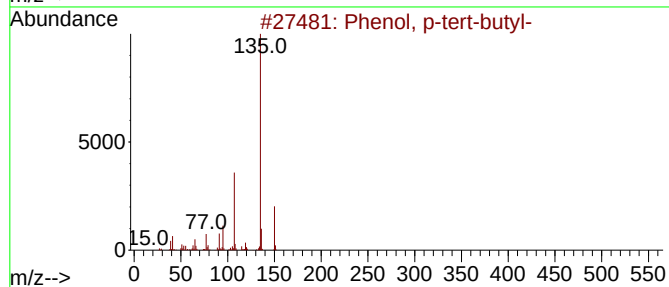
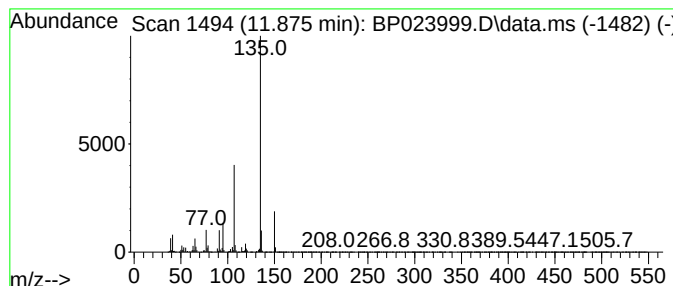
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Phenol, p-tert-butyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.875	16.24 ng/ul	2344570	Naphthalene-d8	10.534

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	96
3			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	94
4			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	94
5			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	87



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E29B5DL

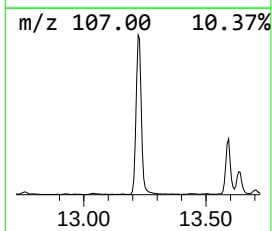
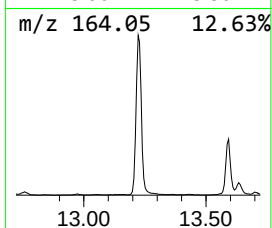
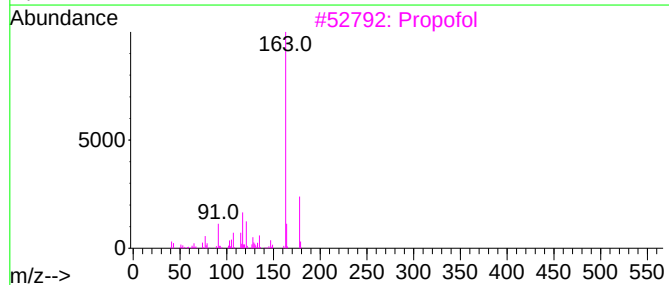
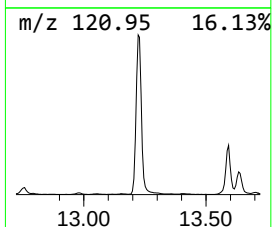
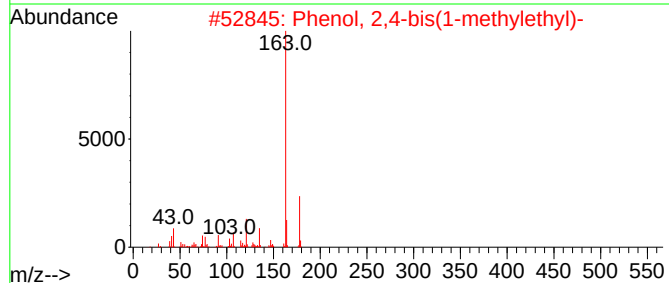
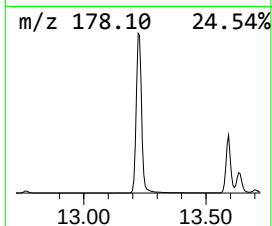
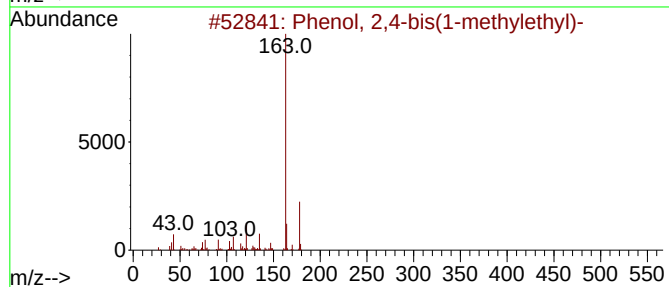
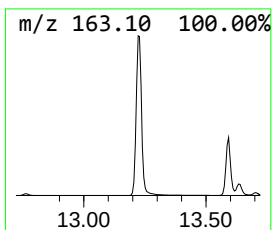
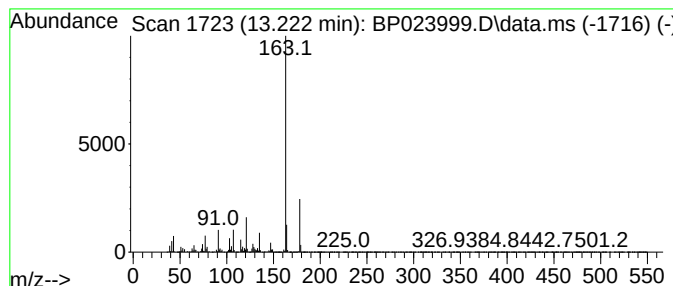
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenol, 2,4-bis(1-methyleth... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.222	39.11 ng/ul	7299380	Acenaphthene-d10	14.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	97
2			Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	94
3			Propofol	178	C12H18O	002078-54-8	94
4			Propofol	178	C12H18O	002078-54-8	93
5			Propofol	178	C12H18O	002078-54-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020725\
Data File : BP023999.D
Acq On : 07 Feb 2025 12:28
Operator : RC/JU
Sample : Q1229-11DL 20X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B5DL

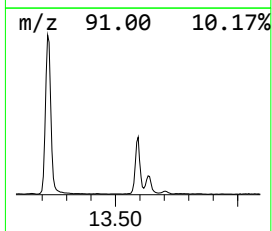
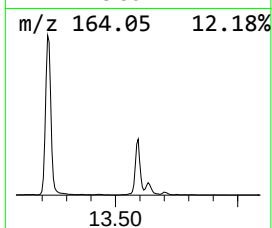
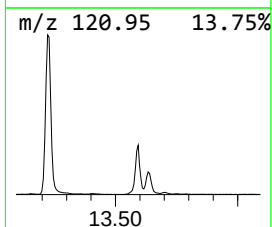
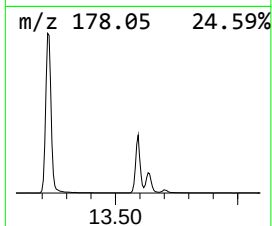
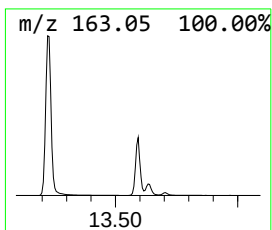
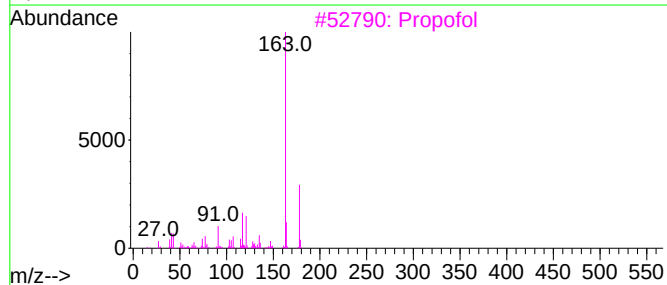
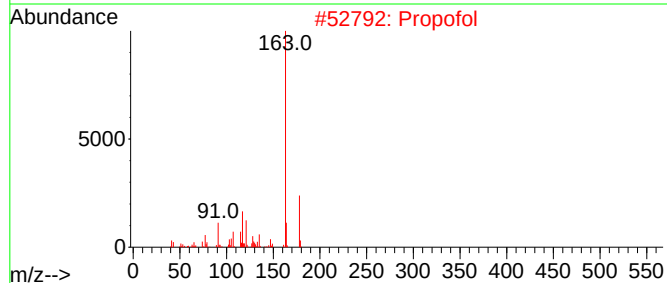
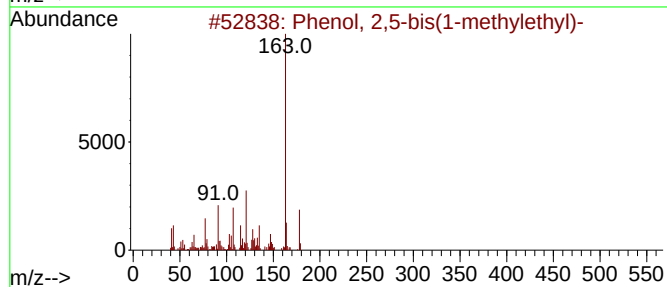
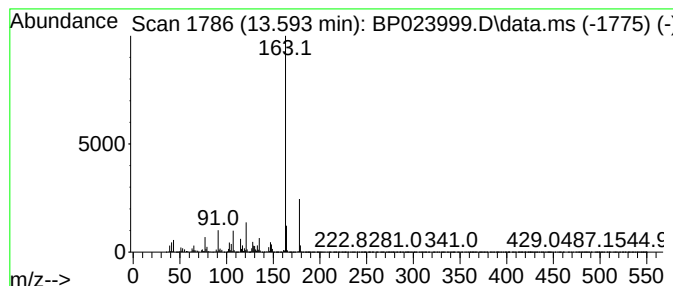
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phenol, 2,5-bis(1-methyleth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.593	12.32 ng/ul	2300210	Acenaphthene-d10	14.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,5-bis(1-methylethyl)-	178	C12H18O	035946-91-9	95
2			Propofol	178	C12H18O	002078-54-8	94
3			Propofol	178	C12H18O	002078-54-8	93
4			Propofol	178	C12H18O	002078-54-8	93
5			Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020725\
Data File : BP023999.D
Acq On : 07 Feb 2025 12:28
Operator : RC/JU
Sample : Q1229-11DL 20X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B5DL

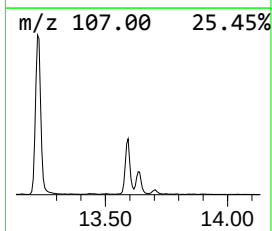
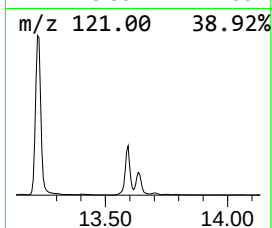
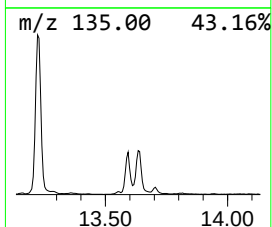
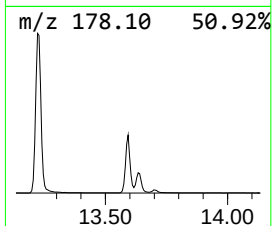
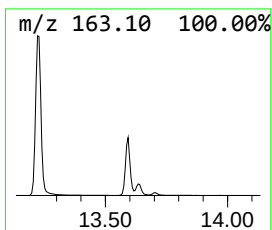
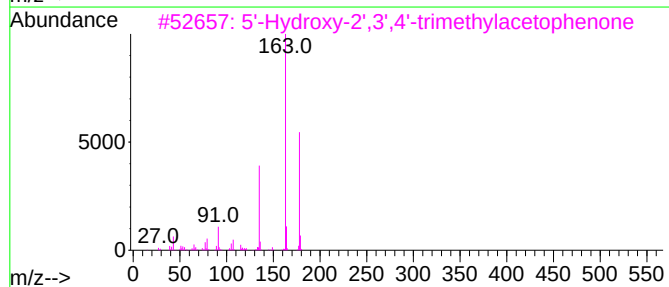
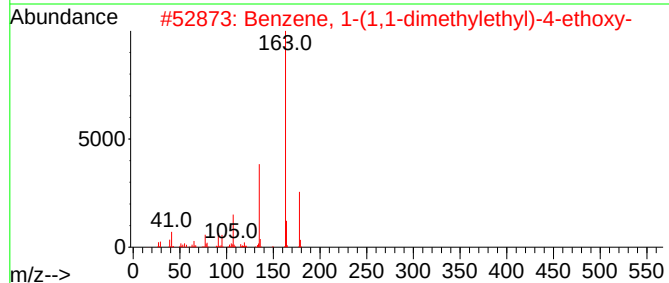
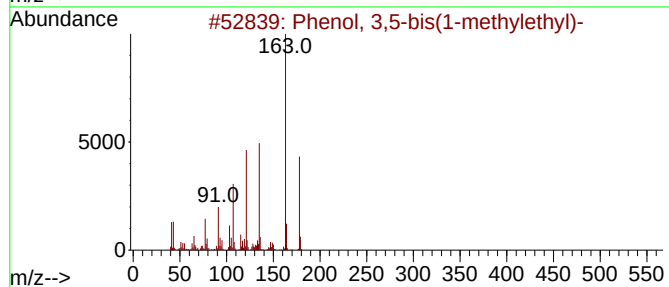
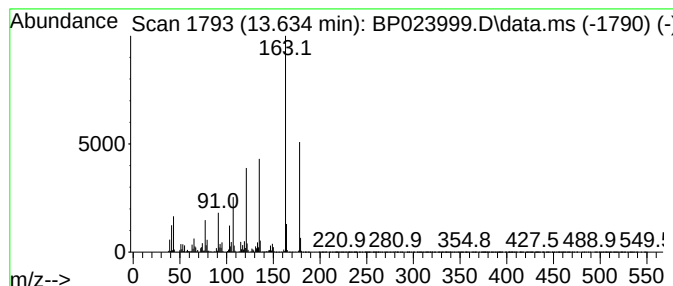
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phenol, 3,5-bis(1-methyleth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.634	4.31 ng/ul	804787	Acenaphthene-d10	14.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 3,5-bis(1-methylethyl)-	178	C12H18O	026886-05-5	96
2			Benzene, 1-(1,1-dimethylethyl)-4...	178	C12H18O	017269-94-2	90
3			5'-Hydroxy-2',3',4'-trimethylace...	178	C11H14O2	013667-28-2	81
4			6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	81
5			Propofol	178	C12H18O	002078-54-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020725\
Data File : BP023999.D
Acq On : 07 Feb 2025 12:28
Operator : RC/JU
Sample : Q1229-11DL 20X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B5DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.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
Phenol, 2-(1-me...	10.458	44.1	ng/u1	6372990	2	10.534	2888130	20.0
Phenol, 2-ethyl...	11.081	2.4	ng/u1	352493	2	10.534	2888130	20.0
Phenol, p-tert-...	11.875	16.2	ng/u1	2344570	2	10.534	2888130	20.0
Phenol, 2,4-bis...	13.222	39.1	ng/u1	7299380	3	14.387	3732610	20.0
Phenol, 2,5-bis...	13.593	12.3	ng/u1	2300210	3	14.387	3732610	20.0
Phenol, 3,5-bis...	13.634	4.3	ng/u1	804787	3	14.387	3732610	20.0