

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013125\
 Data File : BP023826.D
 Acq On : 31 Jan 2025 15:20
 Operator : RC/JU
 Sample : Q1197-18
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E2997

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: BP023826.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.299	33	36	50	rVB	137876	224696	1.43%	0.163%
2	3.711	103	106	124	rVB	607332	985099	6.26%	0.715%
3	4.034	158	161	169	rVB3	30459	44245	0.28%	0.032%
4	4.475	233	236	243	rVB3	14953	20701	0.13%	0.015%
5	4.940	312	315	321	rVB4	18052	29558	0.19%	0.021%
6	6.975	650	661	671	rBV	651196	997018	6.33%	0.723%
7	7.116	678	685	704	rVB	1828197	2862736	18.18%	2.076%
8	7.328	714	721	732	rBV	2191858	3231391	20.52%	2.344%
9	7.416	733	736	743	rVB5	10563	17862	0.11%	0.013%
10	7.781	790	798	805	rBV	1500525	2345006	14.89%	1.701%
11	7.852	805	810	819	rVB	98542	166757	1.06%	0.121%
12	8.204	866	870	878	rBV	14741	26233	0.17%	0.019%
13	8.493	909	919	929	rBV	1897018	2992199	19.00%	2.170%
14	8.928	983	993	1005	rBV	2252451	3444920	21.87%	2.499%
15	9.063	1011	1016	1022	rBV	22529	37162	0.24%	0.027%
16	9.369	1062	1068	1072	rVB	27574	43005	0.27%	0.031%
17	9.640	1105	1114	1129	rBV	1798476	2852895	18.12%	2.069%
18	10.187	1200	1207	1222	rBV	2997706	4697491	29.83%	3.407%
19	10.487	1253	1258	1263	rBV2	73551	136449	0.87%	0.099%
20	10.557	1263	1270	1277	rVB	2278926	3558947	22.60%	2.581%
21	10.687	1285	1292	1302	rBV	2519970	4344418	27.59%	3.151%
22	10.916	1326	1331	1337	rVB	12384	20685	0.13%	0.015%
23	11.351	1400	1405	1409	rBV8	7971	16717	0.11%	0.012%
24	11.587	1440	1445	1448	rBV2	13549	23231	0.15%	0.017%
25	11.640	1451	1454	1459	rVB6	16694	25663	0.16%	0.019%
26	11.904	1493	1499	1506	rBV	365764	553727	3.52%	0.402%
27	12.081	1524	1529	1534	rBV3	17707	37406	0.24%	0.027%
28	12.140	1534	1539	1549	rVB2	56154	103668	0.66%	0.075%
29	12.587	1609	1615	1619	rBV8	8188	16060	0.10%	0.012%
30	13.016	1682	1688	1695	rVB6	11507	22979	0.15%	0.017%
31	13.240	1719	1726	1736	rBV	650231	893213	5.67%	0.648%
32	13.363	1743	1747	1752	rVB3	10350	17179	0.11%	0.012%
33	13.610	1777	1789	1793	rBV	177796	264019	1.68%	0.191%
34	13.651	1793	1796	1803	rVB4	17776	23477	0.15%	0.017%
35	13.728	1803	1809	1814	rBV4	17546	29986	0.19%	0.022%
36	13.804	1815	1822	1837	rVV	5989958	7713475	48.98%	5.595%

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37	14.087	1863	1870	1885	rBV	6148395	9291666	59.00%	6.739%
38	14.398	1913	1923	1932	rBV	3198878	5046370	32.04%	3.660%
39	14.475	1932	1936	1943	rVB9	14528	28861	0.18%	0.021%
40	14.628	1954	1962	1974	rBV	568587	965356	6.13%	0.700%
41	14.822	1992	1995	2000	rVB7	14997	22297	0.14%	0.016%
42	15.310	2074	2078	2084	rVB	25570	34770	0.22%	0.025%
43	15.404	2087	2094	2102	rBV	8276839	11965613	75.98%	8.679%
44	15.516	2108	2113	2126	rBV	2494662	3269509	20.76%	2.371%
45	15.645	2130	2135	2144	rVB	38329	74752	0.47%	0.054%
46	15.775	2152	2157	2162	rVB	48980	75076	0.48%	0.054%
47	16.975	2357	2361	2368	rVB	18536	29636	0.19%	0.021%
48	17.198	2392	2399	2408	rBV2	4055856	5981331	37.98%	4.338%
49	17.298	2408	2416	2428	rBV	9082896	13461631	85.48%	9.764%
50	17.580	2461	2464	2476	rVB3	33552	56616	0.36%	0.041%
51	18.133	2554	2558	2564	rBV3	43440	62602	0.40%	0.045%
52	19.210	2738	2741	2747	rVB	108554	147840	0.94%	0.107%
53	19.286	2750	2754	2759	rBV	103360	142433	0.90%	0.103%
54	19.669	2812	2819	2828	rBV	11581919	15748435	100.00%	11.423%
55	21.639	3148	3154	3162	rBV	4439983	6413596	40.73%	4.652%
56	24.792	3680	3690	3701	rVB	5819712	15416058	97.89%	11.182%
57	25.009	3719	3727	3742	rVB2	2500740	6815961	43.28%	4.944%

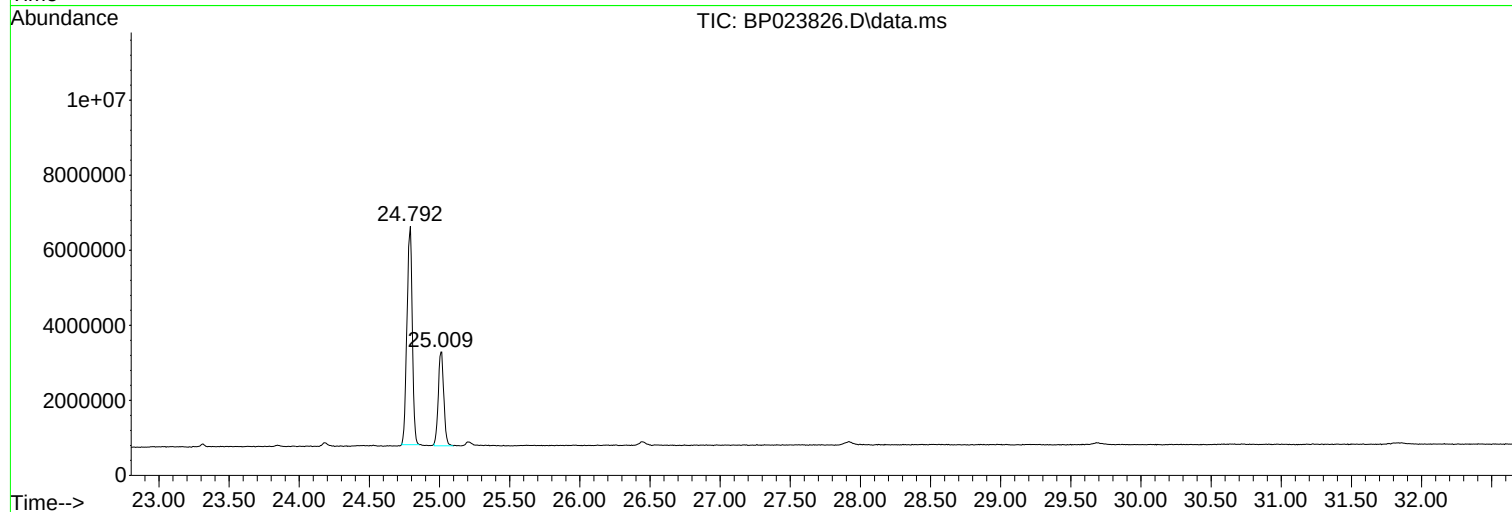
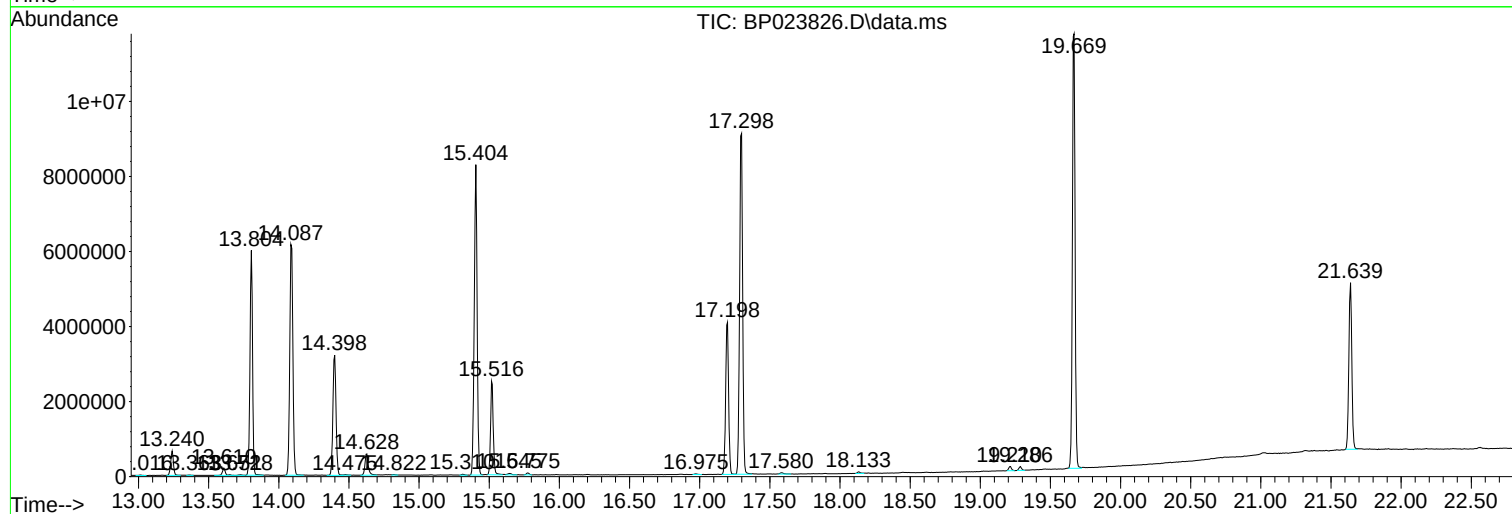
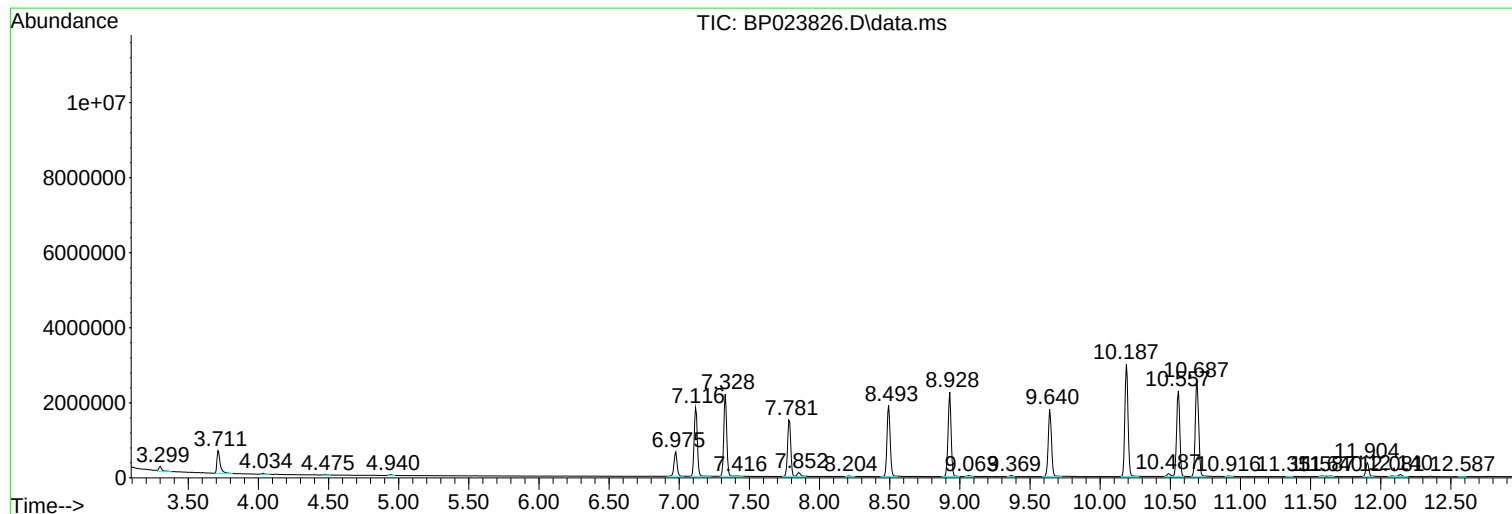
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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



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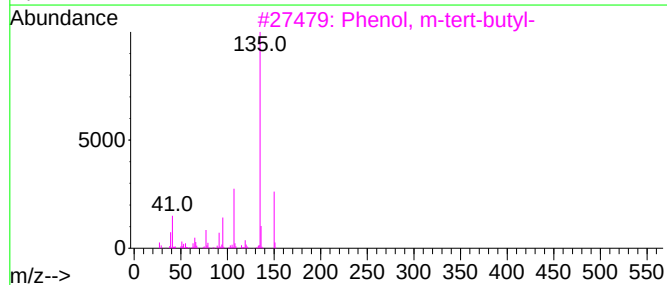
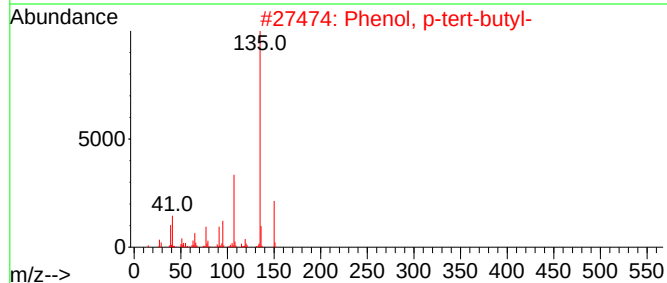
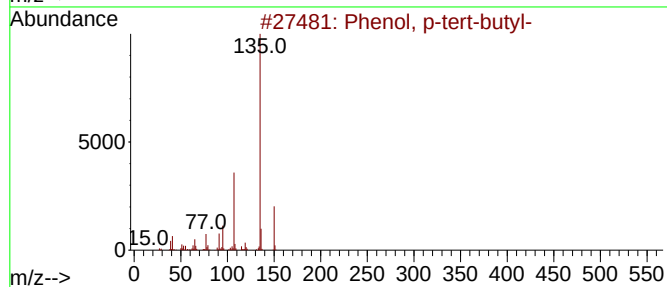
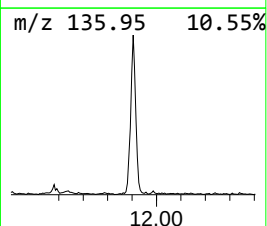
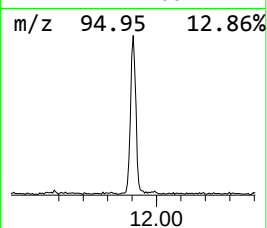
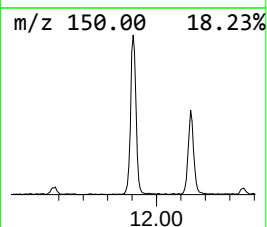
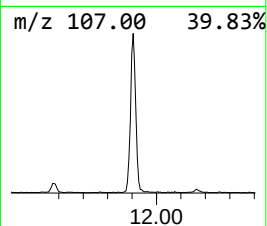
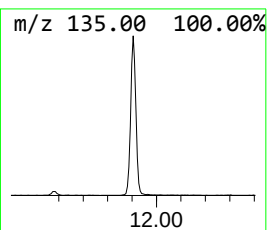
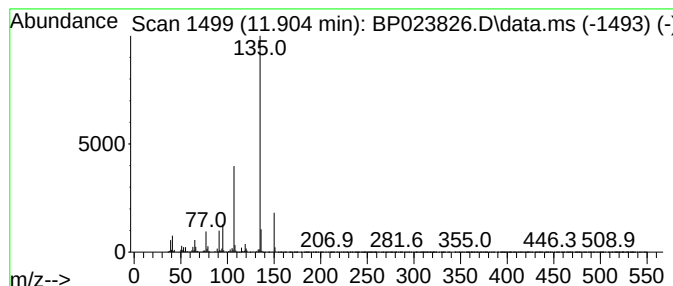
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Phenol, p-tert-butyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.904	3.11 ng/ul	553727	Naphthalene-d8	10.557

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	90



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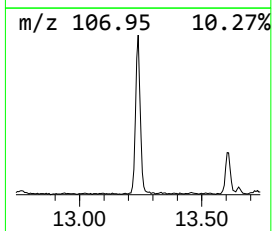
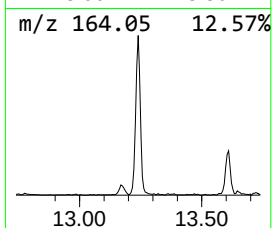
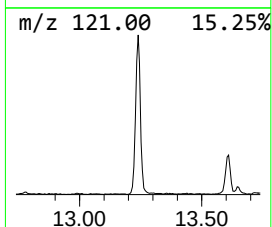
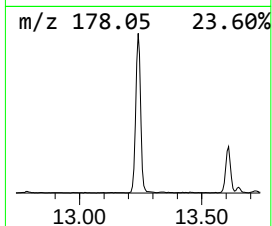
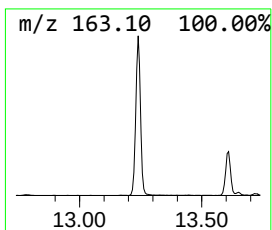
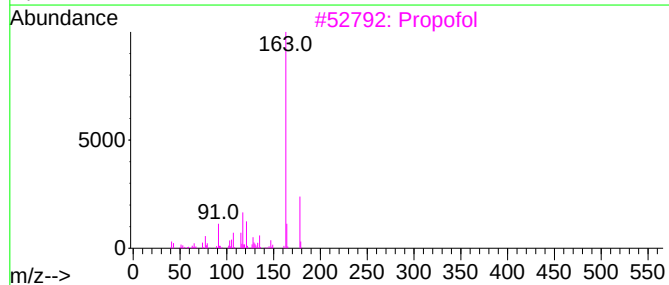
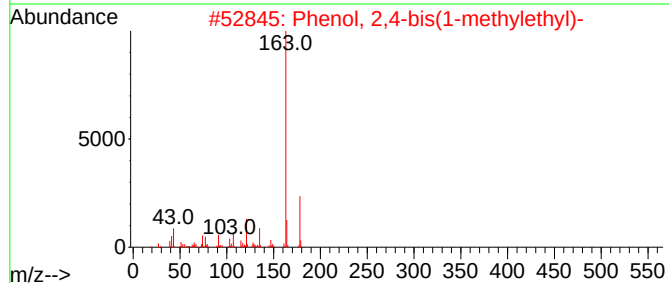
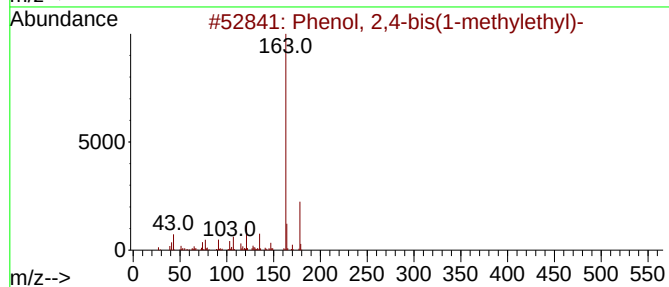
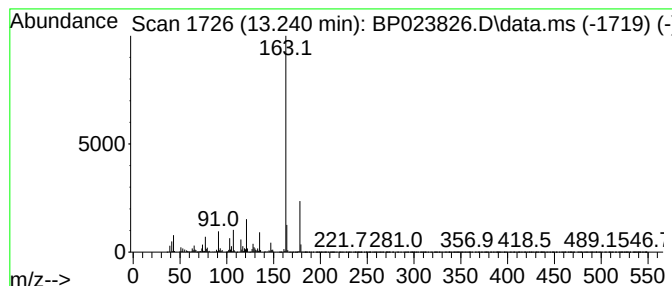
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TIC Library : C:\DATABASE\NIST20.L
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Peak Number 2 Phenol, 2,4-bis(1-methyleth... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.240	3.54 ng/ul	893213	Acenaphthene-d10	14.398

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	95
3			Propofol	178	C12H18O	002078-54-8	94
4			Propofol	178	C12H18O	002078-54-8	94
5			Propofol	178	C12H18O	002078-54-8	93



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, p-tert-...	11.904	3.1	ng/ul	553727	2	10.557	3558950	20.0
Phenol, 2,4-bis...	13.240	3.5	ng/ul	893213	3	14.398	5046370	20.0