

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012825\
 Data File : BM049493.D
 Acq On : 29 Jan 2025 07:47
 Operator : RC/JU
 Sample : Q1184-05
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E2957

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-BM012825.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM049493.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.116	19	22	28	rVB	63220	73284	1.21%	0.125%
2	3.510	85	89	107	rVB	299694	480710	7.92%	0.818%
3	3.816	138	141	144	rVB4	14358	17754	0.29%	0.030%
4	3.910	154	157	160	rBV5	16717	22924	0.38%	0.039%
5	4.240	210	213	218	rVB3	8931	12435	0.20%	0.021%
6	4.704	287	292	298	rVB2	38717	59717	0.98%	0.102%
7	5.940	498	502	506	rVB2	15014	21615	0.36%	0.037%
8	6.704	626	632	644	rVB	266036	429103	7.07%	0.730%
9	6.863	652	659	672	rVB	806521	1256945	20.72%	2.139%
10	7.051	685	691	704	rBV	906855	1385050	22.83%	2.357%
11	7.510	762	769	778	rBV	823221	1264535	20.85%	2.152%
12	7.593	779	783	787	rVB3	8858	14519	0.24%	0.025%
13	8.222	884	890	901	rBV	733754	1151992	18.99%	1.961%
14	8.651	956	963	974	rBV	864225	1411158	23.26%	2.402%
15	8.798	984	988	991	rVB4	5364	7158	0.12%	0.012%
16	9.098	1034	1039	1045	rBV3	13927	18838	0.31%	0.032%
17	9.369	1078	1085	1094	rBV	672388	1108728	18.28%	1.887%
18	9.475	1098	1103	1107	rBV4	10641	19565	0.32%	0.033%
19	9.781	1152	1155	1160	rVB5	7653	8911	0.15%	0.015%
20	9.904	1168	1176	1190	rBV	1039295	1780785	29.36%	3.031%
21	10.269	1231	1238	1248	rBV	1009580	1617714	26.67%	2.753%
22	10.416	1256	1263	1278	rBV	779455	1392456	22.95%	2.370%
23	10.657	1297	1304	1307	rBV7	6620	12639	0.21%	0.022%
24	11.633	1459	1470	1491	rBV	3763497	6066055	100.00%	10.325%
25	11.869	1505	1510	1518	rVB2	12215	24813	0.41%	0.042%
26	12.998	1696	1702	1714	rBV	182380	261640	4.31%	0.445%
27	13.180	1727	1733	1736	rVB7	5250	8393	0.14%	0.014%
28	13.374	1761	1766	1769	rBV	46702	62842	1.04%	0.107%
29	13.416	1769	1773	1780	rVB	65649	93198	1.54%	0.159%
30	13.486	1780	1785	1793	rVB3	11461	18000	0.30%	0.031%
31	13.574	1793	1800	1812	rBV	1982408	2688536	44.32%	4.576%
32	13.839	1838	1845	1860	rBV	2259462	3139380	51.75%	5.343%
33	14.151	1891	1898	1908	rBV	1493709	2087760	34.42%	3.554%
34	14.392	1933	1939	1955	rBV2	75714	246116	4.06%	0.419%
35	14.992	2035	2041	2046	rVB6	7047	13599	0.22%	0.023%
36	15.151	2061	2068	2083	rBV	2936447	4116339	67.86%	7.006%

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 Title : SVOA CALIBRATION

37	15.280	2083	2090	2098	rVV	839409	1136489	18.74%	1.934%
38	15.339	2098	2100	2106	rVV6	15259	28819	0.48%	0.049%
39	15.392	2106	2109	2119	rVB5	13730	25550	0.42%	0.043%
40	15.527	2127	2132	2141	rVB	60453	89590	1.48%	0.152%
41	16.086	2222	2227	2235	rVB5	10626	19095	0.31%	0.033%
42	16.692	2326	2330	2333	rBV2	7930	9480	0.16%	0.016%
43	16.904	2360	2366	2376	rBV	1847745	2408495	39.70%	4.099%
44	17.004	2376	2383	2399	rVV	3363073	4637137	76.44%	7.893%
45	17.827	2518	2523	2529	rBV5	28764	48069	0.79%	0.082%
46	17.898	2531	2535	2542	rVB4	6280	13098	0.22%	0.022%
47	18.298	2600	2603	2606	rVV4	6584	6976	0.12%	0.012%
48	18.668	2661	2666	2669	rBV7	6223	8870	0.15%	0.015%
49	18.868	2696	2700	2705	rVB2	40256	55109	0.91%	0.094%
50	18.945	2708	2713	2717	rBV	32550	46156	0.76%	0.079%
51	19.121	2739	2743	2748	rBV8	12539	21619	0.36%	0.037%
52	19.309	2767	2775	2784	rBV	4796532	5872954	96.82%	9.996%
53	21.133	3080	3085	3099	rBV	1966237	2772278	45.70%	4.719%
54	23.727	3517	3526	3539	rBV	2797587	6034598	99.48%	10.271%
55	23.915	3550	3558	3576	rVB	1358989	3122417	51.47%	5.315%

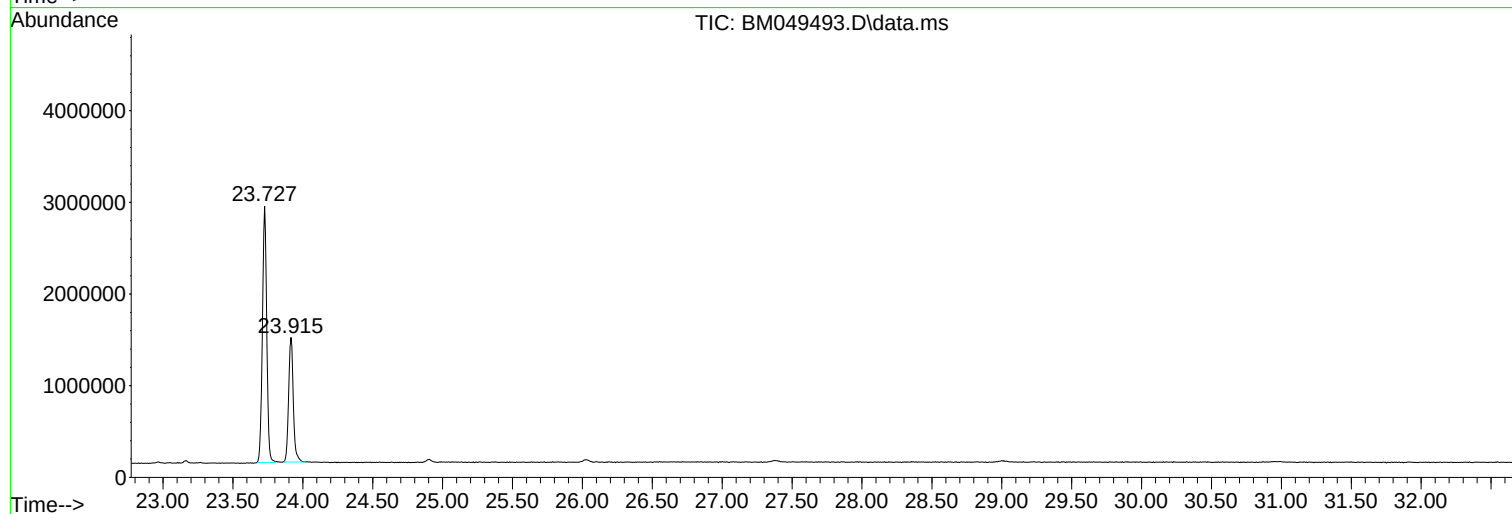
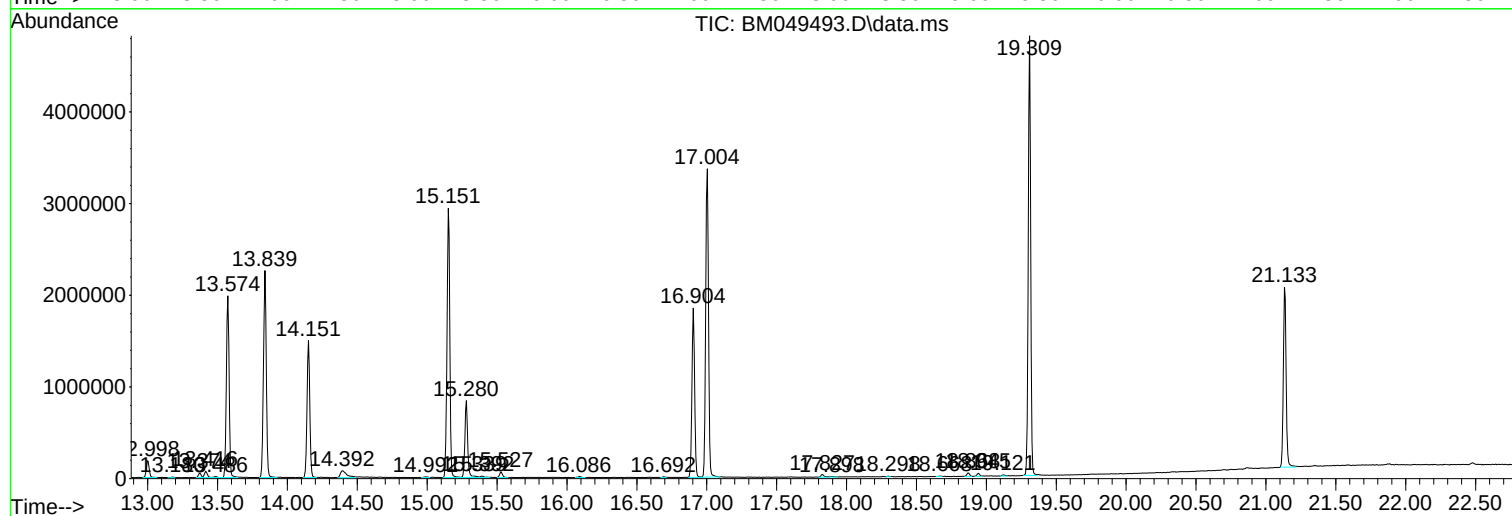
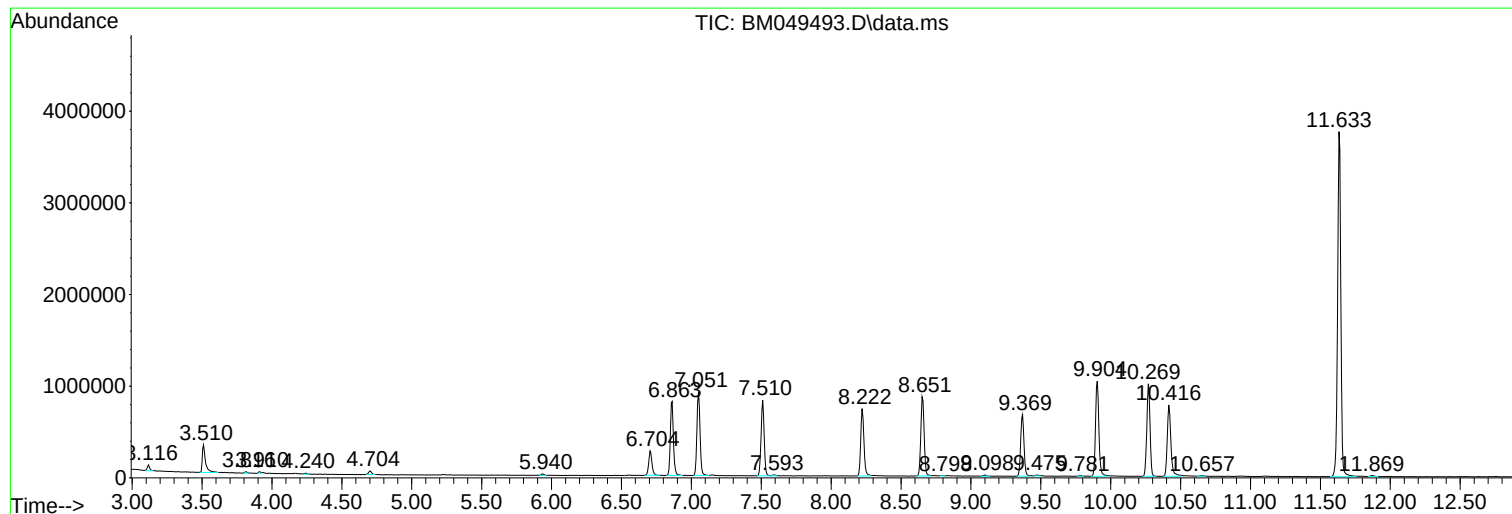
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM012825.MA.M
Quant Title : SVOA CALIBRATION

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



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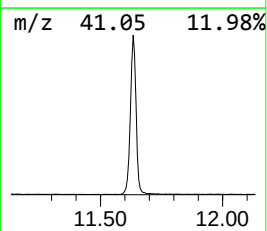
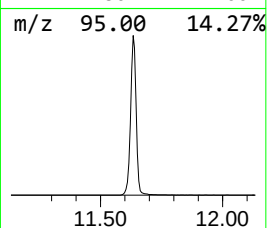
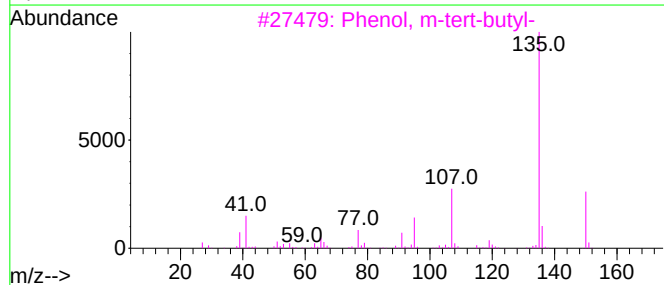
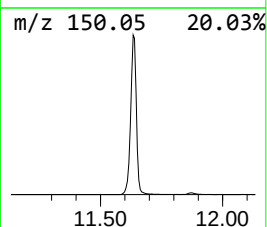
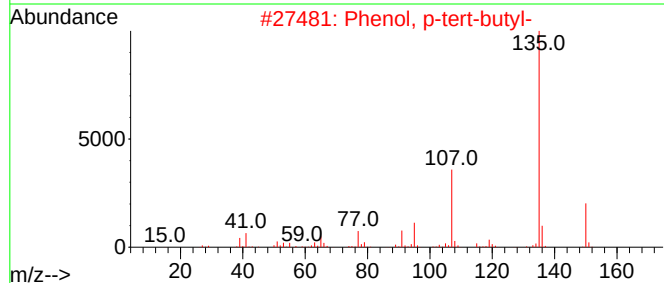
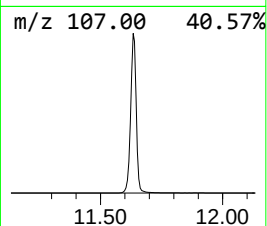
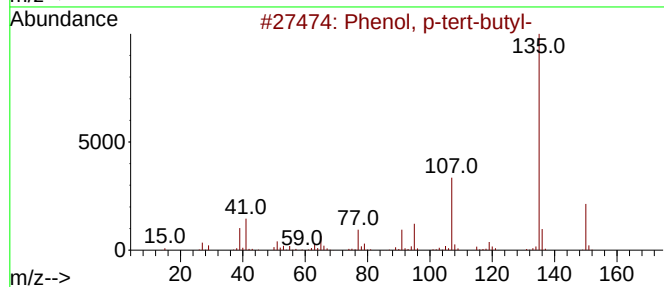
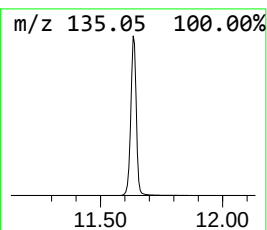
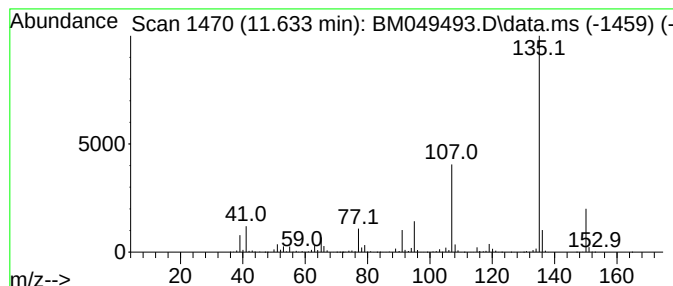
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM012825.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.633	75.00 ng/ul	6066060	Naphthalene-d8	10.269

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	96
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	91



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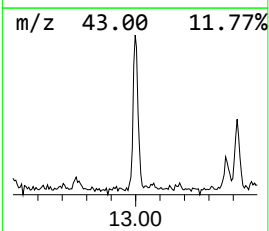
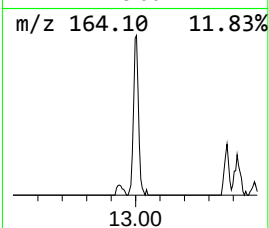
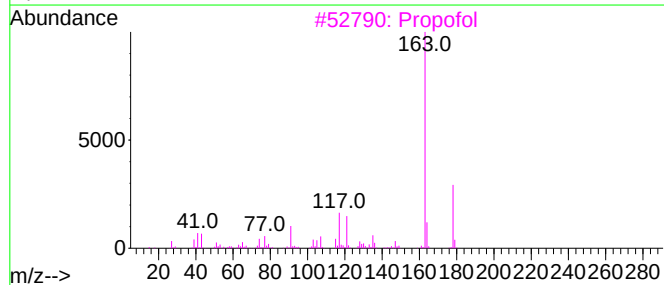
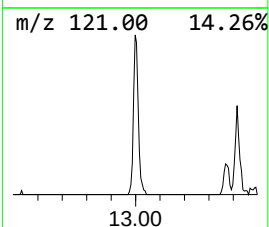
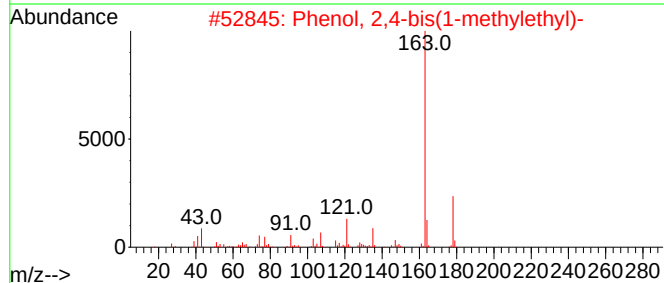
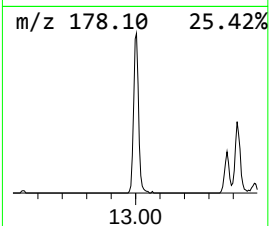
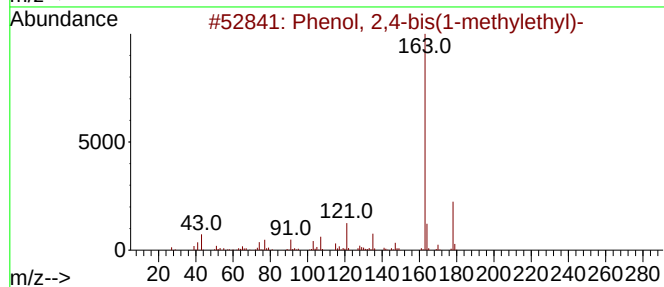
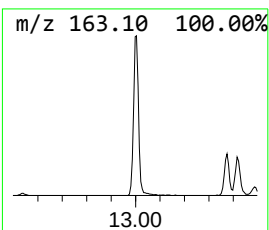
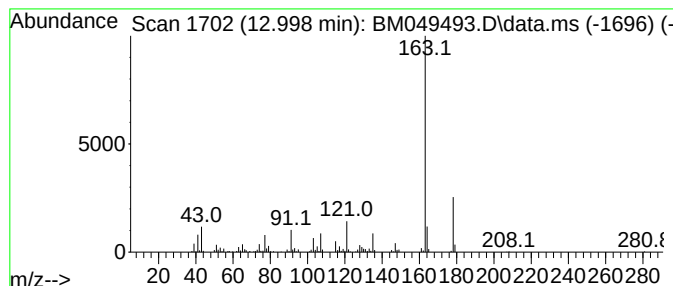
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM012825.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.998	2.51 ng/ul	261640	Acenaphthene-d10	14.151

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	96
3			Propofol	178	C12H18O	002078-54-8	95
4			Propofol	178	C12H18O	002078-54-8	93
5			Propofol	178	C12H18O	002078-54-8	91



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, p-tert-...	11.633	75.0	ng/u1	6066060	2	10.269	1617710	20.0
Phenol, 2,4-bis...	12.998	2.5	ng/u1	261640	3	14.151	2087760	20.0