

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061522\
 Data File : BN020215.D
 Acq On : 16 Jun 2022 21:32
 Operator : CG/JU
 Sample : N3285-16
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EW4Q8

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_N\Methods\SFAM-EPA-BN060622.M
 Title : SVOA CALIBRATION

Signal : TIC: BN020215.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.258	25	29	35	rBV	30364	41444	0.83%	0.074%
2	3.670	96	99	115	rBV	186744	336699	6.77%	0.603%
3	3.905	135	139	146	rVB	15451	25972	0.52%	0.046%
4	3.999	151	155	160	rVB	12623	17480	0.35%	0.031%
5	4.370	215	218	225	rVB8	3638	7664	0.15%	0.014%
6	4.670	266	269	274	rBV4	4894	8442	0.17%	0.015%
7	4.905	306	309	315	rVB2	7368	11159	0.22%	0.020%
8	5.340	377	383	389	rBV4	5141	10156	0.20%	0.018%
9	6.817	630	634	642	rBV6	3163	6652	0.13%	0.012%
10	6.987	655	663	678	rBV2	183845	483935	9.72%	0.866%
11	7.122	680	686	699	rVV	580736	966755	19.43%	1.730%
12	7.334	714	722	733	rVV	625765	1023536	20.57%	1.832%
13	7.793	792	800	808	rVV	629732	1037004	20.84%	1.856%
14	7.869	808	813	821	rVB4	5426	12853	0.26%	0.023%
15	8.516	916	923	943	rBV	562583	964043	19.37%	1.725%
16	8.946	988	996	1010	rBV	765141	1323655	26.60%	2.369%
17	9.087	1016	1020	1026	rBV4	7701	12591	0.25%	0.023%
18	9.193	1033	1038	1046	rVB7	7844	18542	0.37%	0.033%
19	9.322	1057	1060	1067	rBV6	4435	6673	0.13%	0.012%
20	9.663	1110	1118	1131	rBV	626520	1100097	22.11%	1.969%
21	9.775	1131	1137	1140	rVV3	16189	29337	0.59%	0.052%
22	9.810	1140	1143	1148	rVB3	13727	21301	0.43%	0.038%
23	9.928	1159	1163	1171	rVB7	4032	8962	0.18%	0.016%
24	10.052	1176	1184	1186	rVV2	14556	30914	0.62%	0.055%
25	10.081	1186	1189	1194	rVV	20242	36938	0.74%	0.066%
26	10.116	1194	1195	1203	rVB6	7223	9121	0.18%	0.016%
27	10.222	1205	1213	1228	rVV	913897	1602465	32.20%	2.868%
28	10.475	1250	1256	1257	rBV3	8326	13146	0.26%	0.024%
29	10.522	1257	1264	1269	rVV	727365	1201601	24.15%	2.150%
30	10.581	1269	1274	1283	rVB	988826	1709470	34.35%	3.059%
31	10.716	1289	1297	1309	rBV	687176	1232462	24.77%	2.205%
32	10.952	1329	1337	1349	rVV	179325	301783	6.06%	0.540%
33	11.069	1352	1357	1365	rVB6	10572	17853	0.36%	0.032%
34	11.146	1365	1370	1375	rVB7	7015	12202	0.25%	0.022%
35	11.222	1375	1383	1385	rBV7	6064	14134	0.28%	0.025%
36	11.246	1385	1387	1395	rVV7	5779	11169	0.22%	0.020%

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
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37	11.334	1397	1402	1404	rBV6	4970	8041	0.16%	0.014%
38	11.393	1407	1412	1417	rVV6	20131	47745	0.96%	0.085%
39	11.446	1417	1421	1427	rVV5	14161	26488	0.53%	0.047%
40	11.505	1427	1431	1435	rVV7	5409	8776	0.18%	0.016%
41	11.563	1435	1441	1446	rVV6	19178	38558	0.77%	0.069%
42	11.622	1446	1451	1455	rVV4	22978	54328	1.09%	0.097%
43	11.681	1459	1461	1467	rVV3	19412	32646	0.66%	0.058%
44	11.728	1467	1469	1475	rVB6	5808	9296	0.19%	0.017%
45	11.822	1482	1485	1487	rBV3	5928	7077	0.14%	0.013%
46	11.940	1494	1505	1522	rBV	2348263	3901441	78.40%	6.982%
47	12.169	1534	1544	1552	rVB3	21562	44331	0.89%	0.079%
48	13.275	1725	1732	1741	rBV	244027	350831	7.05%	0.628%
49	13.640	1788	1794	1798	rBV	44090	61983	1.25%	0.111%
50	13.687	1798	1802	1808	rVB	32519	48796	0.98%	0.087%
51	13.757	1808	1814	1817	rVB2	9318	13236	0.27%	0.024%
52	13.834	1820	1827	1836	rBV	1986769	2775231	55.77%	4.966%
53	13.969	1845	1850	1856	rVB8	4783	8876	0.18%	0.016%
54	14.116	1865	1875	1883	rBV	2027849	3168609	63.67%	5.670%
55	14.216	1885	1892	1900	rVB10	12339	37355	0.75%	0.067%
56	14.334	1908	1912	1918	rVB3	14135	20301	0.41%	0.036%
57	14.422	1920	1927	1935	rBV	1593476	2420947	48.65%	4.332%
58	14.675	1961	1970	1982	rBV	164013	327244	6.58%	0.586%
59	14.781	1984	1988	1995	rVV3	19666	37051	0.74%	0.066%
60	14.840	1995	1998	2004	rVB6	8139	14200	0.29%	0.025%
61	14.910	2004	2010	2014	rBV2	12107	18262	0.37%	0.033%
62	15.240	2060	2066	2069	rBV7	6837	13744	0.28%	0.025%
63	15.287	2069	2074	2079	rVB	35343	48363	0.97%	0.087%
64	15.416	2089	2096	2108	rBV	2682176	4012390	80.63%	7.180%
65	15.534	2110	2116	2132	rBV	1149752	1570568	31.56%	2.810%
66	15.651	2132	2136	2142	rBV	34537	57951	1.16%	0.104%
67	15.822	2161	2165	2174	rVB5	8810	17744	0.36%	0.032%
68	16.145	2216	2220	2223	rVV	39832	55805	1.12%	0.100%
69	16.198	2227	2229	2234	rVV4	14224	17273	0.35%	0.031%
70	16.487	2274	2278	2280	rBV4	5195	8164	0.16%	0.015%
71	16.657	2304	2307	2308	rBV3	8763	9050	0.18%	0.016%
72	16.845	2334	2339	2343	rBV7	7834	13521	0.27%	0.024%
73	16.934	2350	2354	2359	rBV2	40434	58834	1.18%	0.105%
74	16.987	2359	2363	2368	rVB3	14602	21526	0.43%	0.039%
75	17.092	2379	2381	2386	rVB5	7354	8636	0.17%	0.015%
76	17.163	2386	2393	2402	rBV	2018063	2877456	57.82%	5.149%

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 Title : SVOA CALIBRATION

77	17.263	2403	2410	2424	rVV	2917063	4401609	88.45%	7.877%
78	17.675	2476	2480	2485	rVB	30238	36651	0.74%	0.066%
79	17.787	2496	2499	2503	rVB4	19726	24649	0.50%	0.044%
80	17.839	2506	2508	2512	rVB5	8664	10381	0.21%	0.019%
81	18.063	2542	2546	2553	rBV2	57309	89973	1.81%	0.161%
82	18.369	2594	2598	2606	rBV	29380	48920	0.98%	0.088%
83	18.822	2671	2675	2680	rBV5	21694	38920	0.78%	0.070%
84	19.004	2702	2706	2711	rBV	20943	27936	0.56%	0.050%
85	19.122	2722	2726	2730	rVB	44959	61350	1.23%	0.110%
86	19.192	2734	2738	2743	rBV	43363	61681	1.24%	0.110%
87	19.351	2760	2765	2769	rBV2	52363	85341	1.71%	0.153%
88	19.498	2787	2790	2793	rBV4	24745	29456	0.59%	0.053%
89	19.557	2794	2800	2807	rBV	3494703	4976383	100.00%	8.905%
90	20.116	2893	2895	2899	rVB2	20431	21011	0.42%	0.038%
91	20.545	2966	2968	2969	rBV2	17706	14869	0.30%	0.027%
92	20.569	2970	2972	2977	rVV2	44307	49328	0.99%	0.088%
93	20.610	2977	2979	2983	rVB2	17551	19013	0.38%	0.034%
94	21.351	3099	3105	3112	rBV	2108006	2847202	57.21%	5.095%
95	22.439	3288	3290	3297	rVB	72312	102304	2.06%	0.183%
96	22.975	3377	3381	3387	rVB2	97712	145597	2.93%	0.261%
97	23.516	3466	3473	3480	rBV	1930539	3970056	79.78%	7.104%
98	23.669	3491	3499	3509	rVB2	1157284	2394200	48.11%	4.284%
99	24.257	3595	3599	3607	rVB	121079	232826	4.68%	0.417%
100	25.057	3730	3735	3745	rVB2	107518	251642	5.06%	0.450%

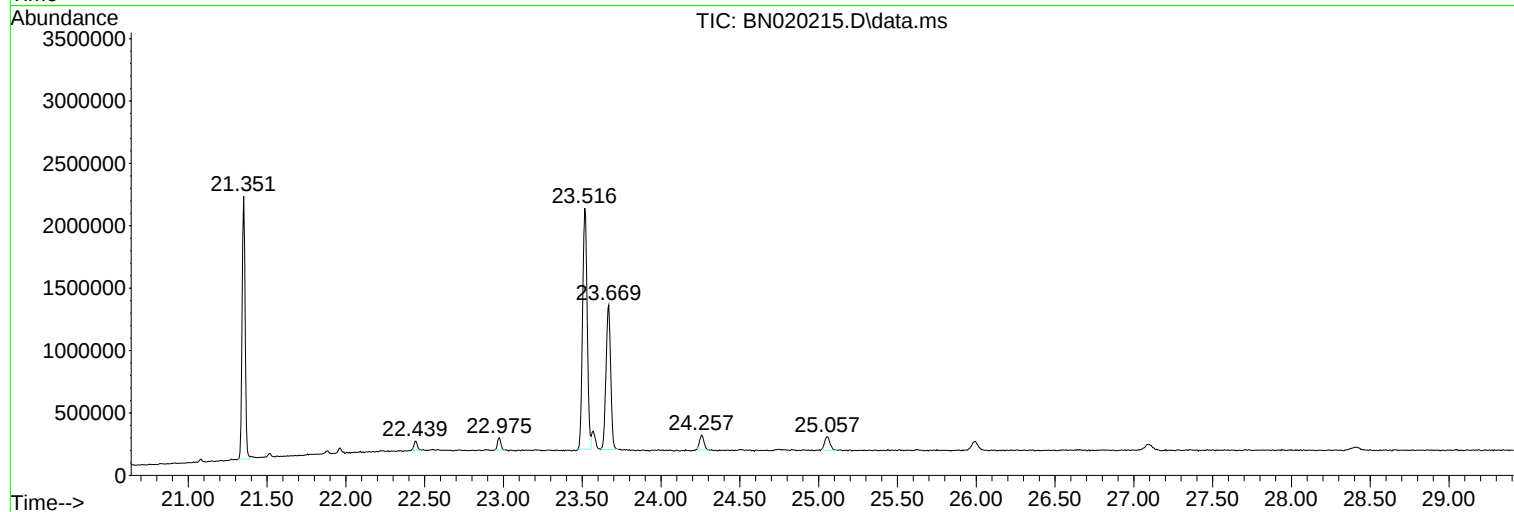
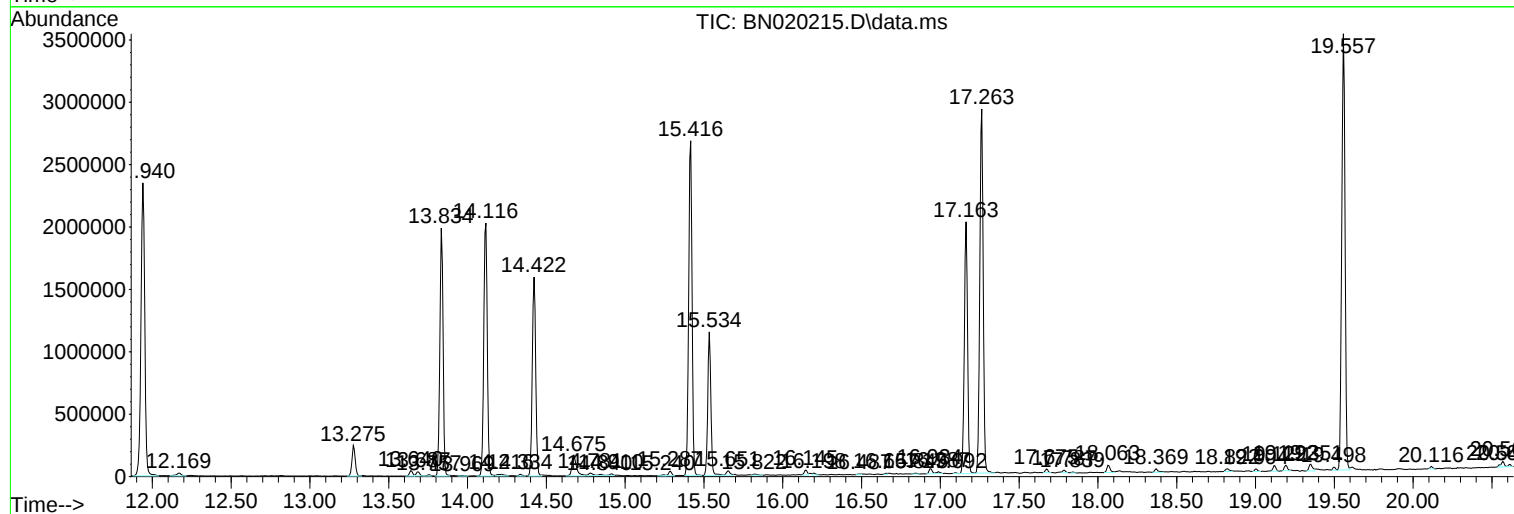
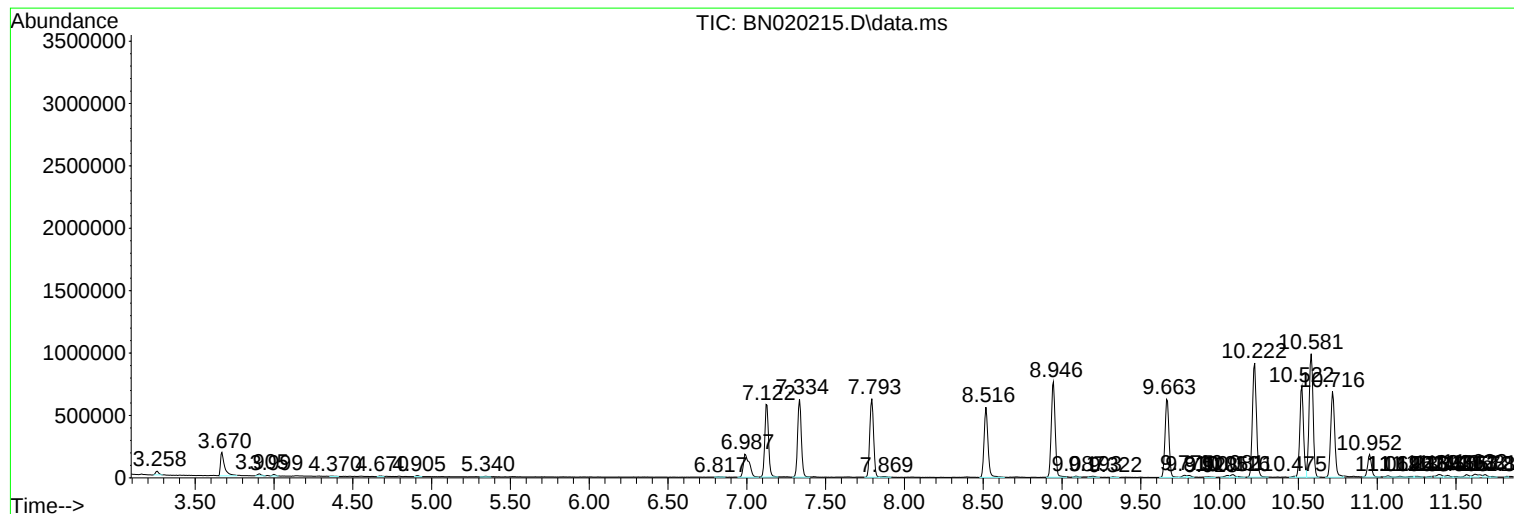
Sum of corrected areas: 55882182

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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



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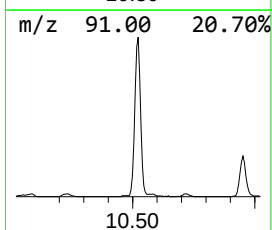
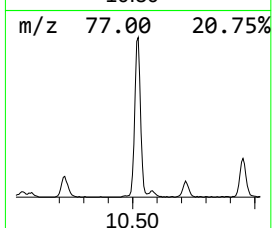
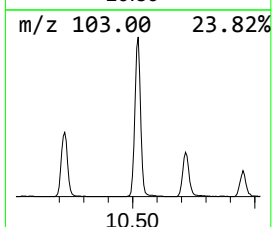
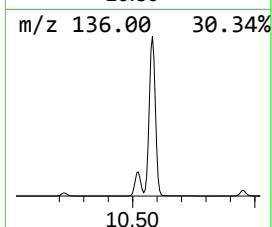
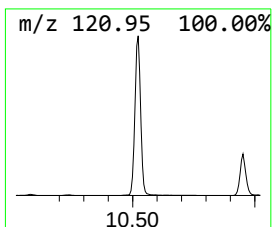
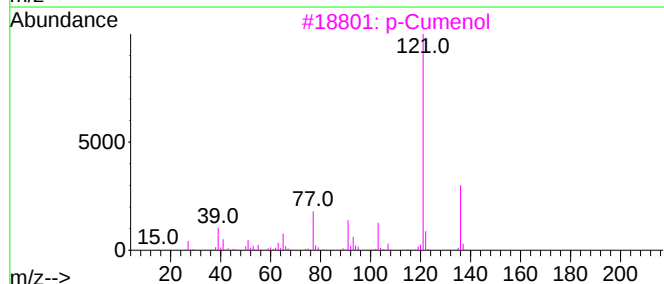
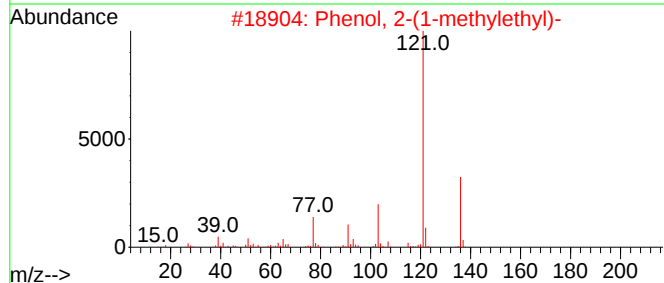
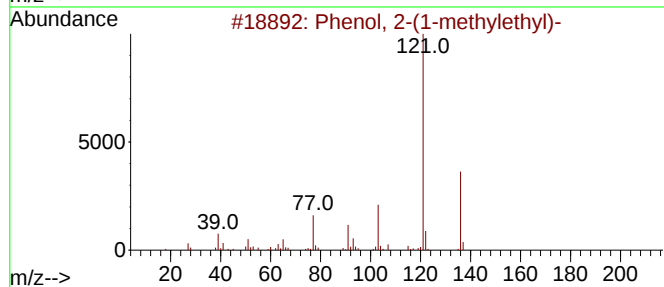
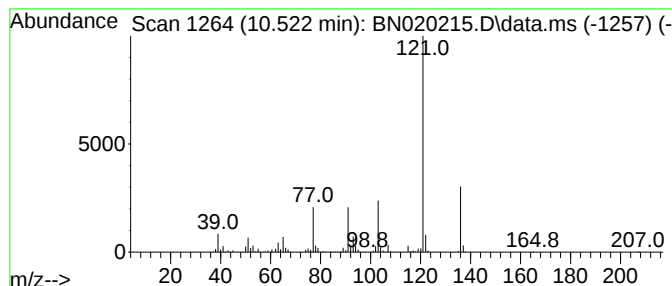
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.522	14.06 ng/ul	1201600	Naphthalene-d8	10.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	93
2			Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	91
3			p-Cumenol	136	C9H12O	000099-89-8	90
4			Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	90
5			Phenol, 2-ethyl-6-methyl-	136	C9H12O	001687-64-5	87



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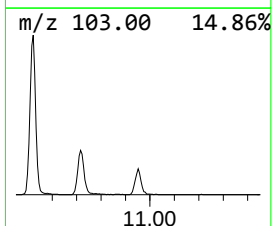
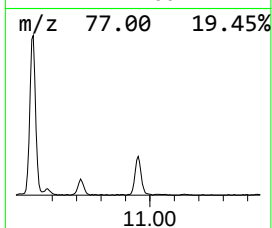
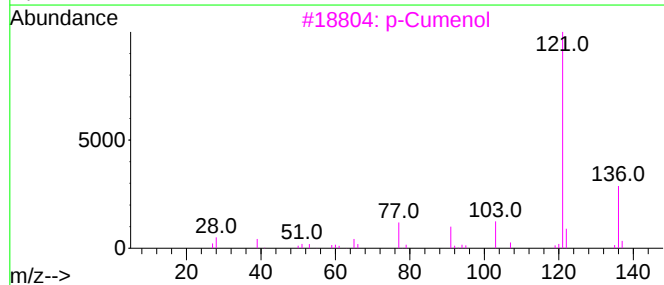
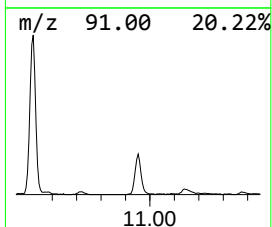
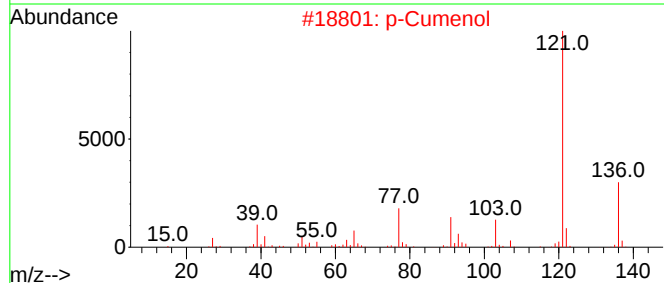
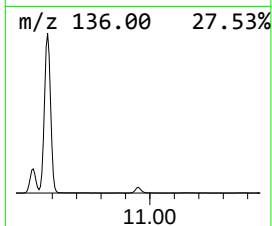
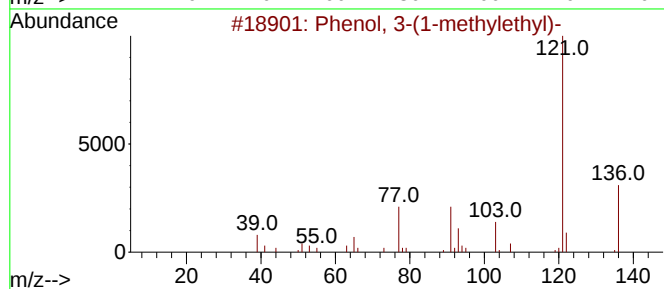
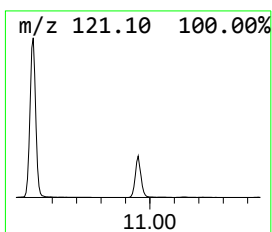
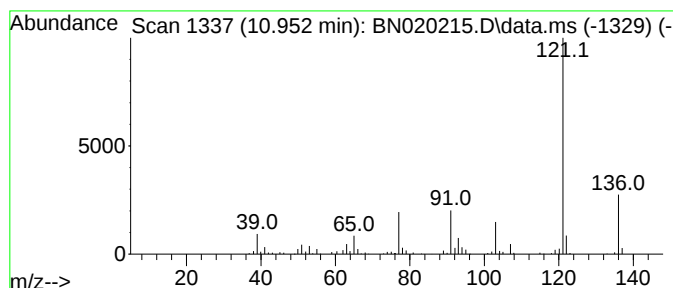
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenol, 3-(1-methylethyl)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.952	3.53 ng/ul	301783	Naphthalene-d8	10.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	94
2			p-Cumenol	136	C9H12O	000099-89-8	94
3			p-Cumenol	136	C9H12O	000099-89-8	91
4			Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	91
5			Phenol, 2-ethyl-6-methyl-	136	C9H12O	001687-64-5	90



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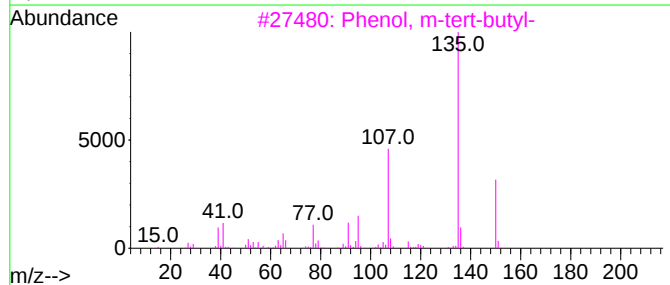
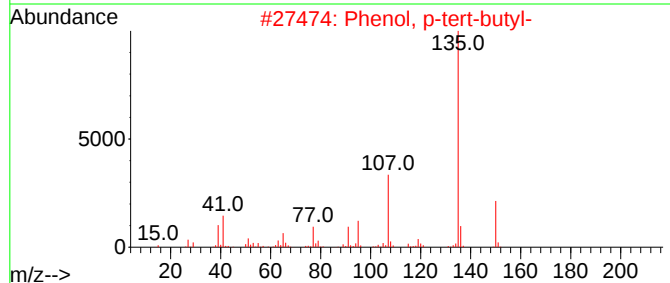
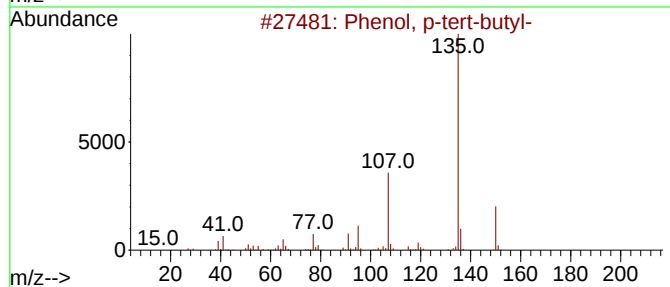
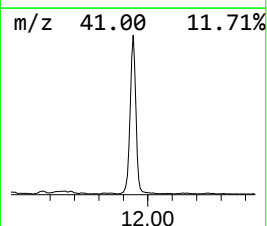
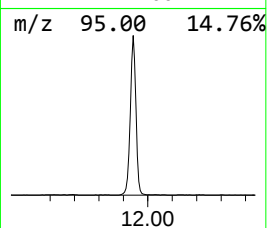
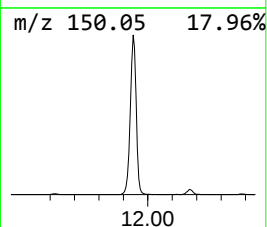
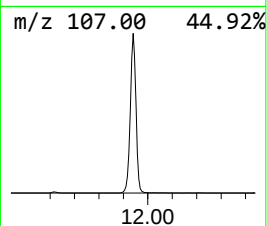
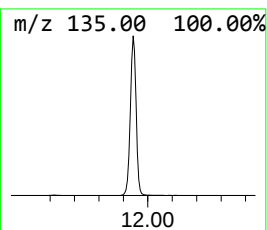
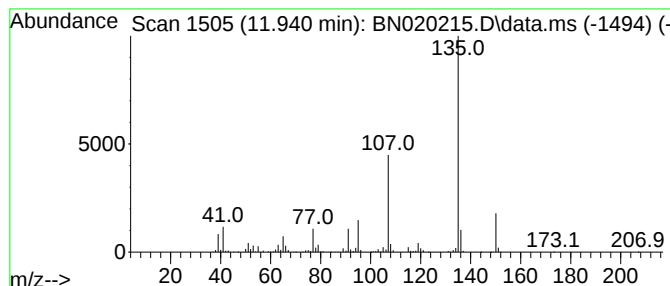
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Phenol, p-tert-butyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.940	45.65 ng/ul	3901440	Naphthalene-d8	10.581

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	94
3			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
4			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
5			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061522\
Data File : BN020215.D
Acq On : 16 Jun 2022 21:32
Operator : CG/JU
Sample : N3285-16
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
EW4Q8

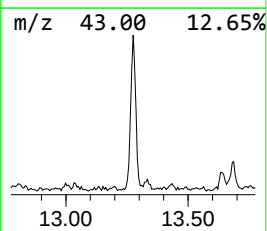
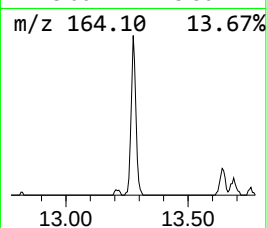
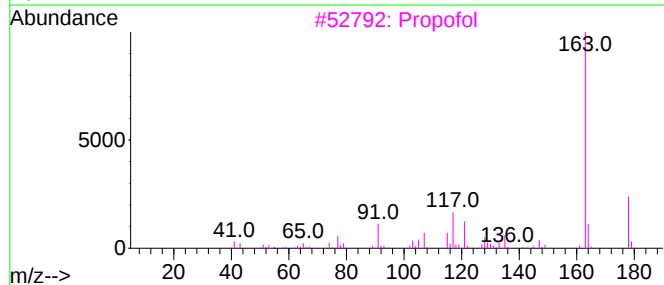
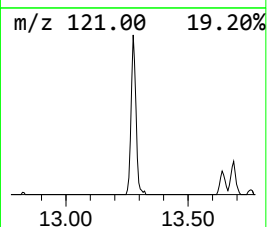
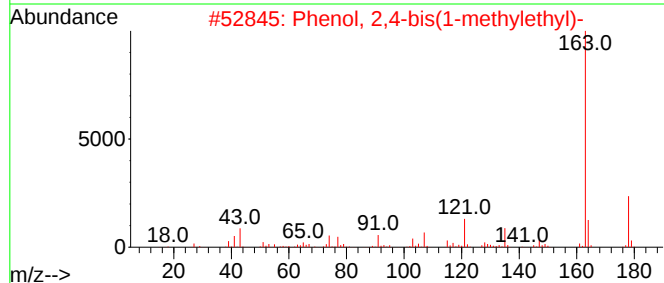
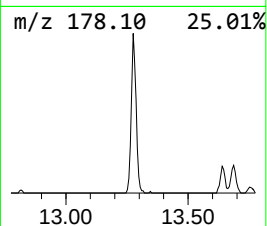
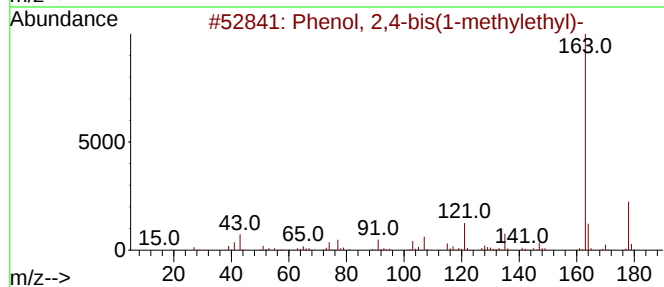
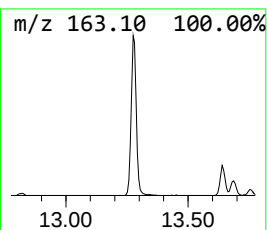
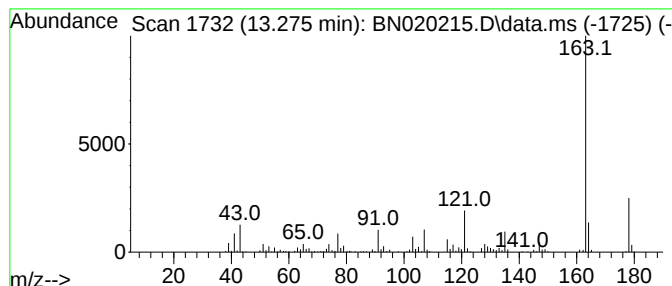
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.275	2.90 ng/ul	350831	Acenaphthene-d10	14.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	95
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_N\Data\BN061522\
Data File : BN020215.D
Acq On : 16 Jun 2022 21:32
Operator : CG/JU
Sample : N3285-16
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
EW4Q8

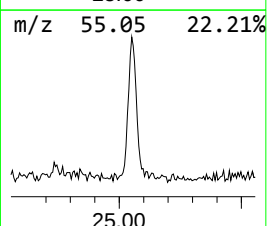
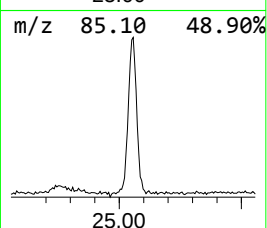
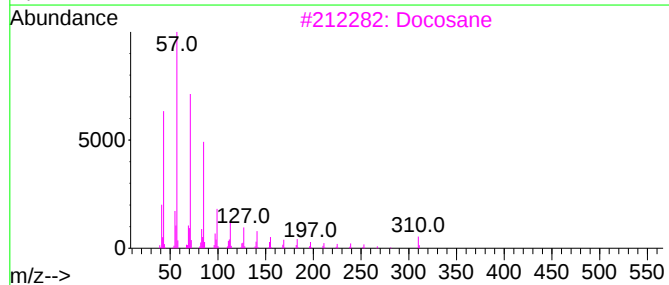
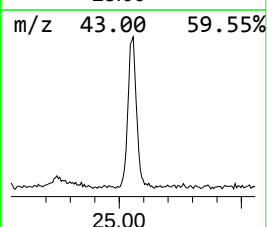
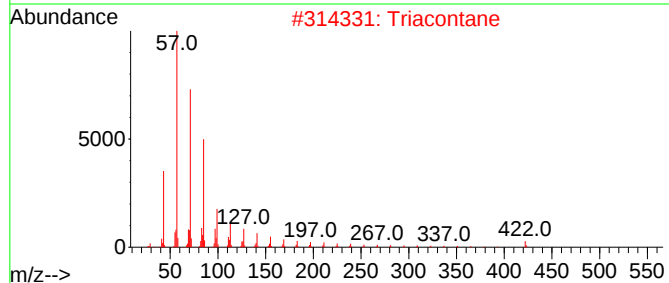
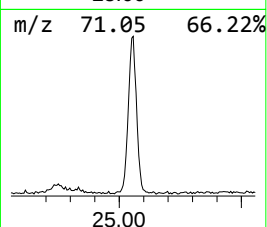
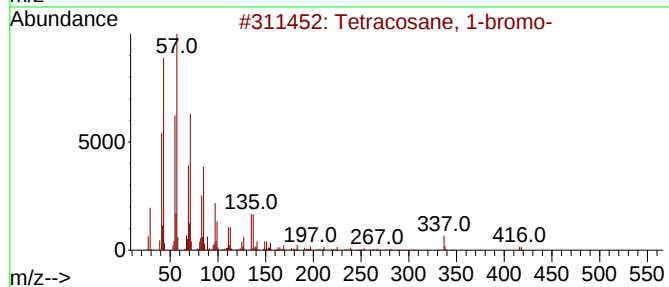
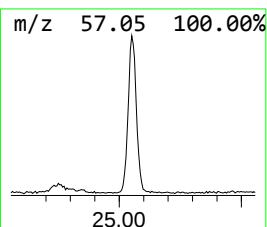
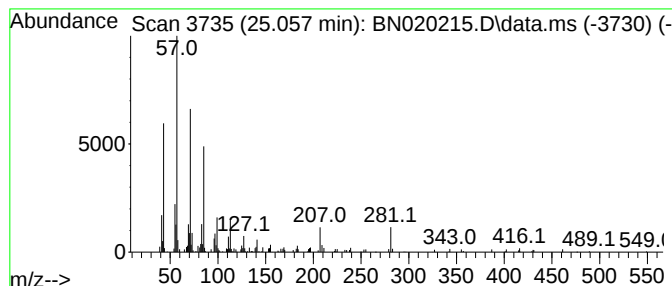
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Tetracosane, 1-bromo- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.057	2.10 ng/ul	251642	Perylene-d12	23.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane, 1-bromo-	416	C24H49Br	006946-24-3	64
2			Triacontane	422	C30H62	000638-68-6	62
3			Docosane	310	C22H46	000629-97-0	62
4			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	62
5			Hexacosane	366	C26H54	000630-01-3	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061522\
Data File : BN020215.D
Acq On : 16 Jun 2022 21:32
Operator : CG/JU
Sample : N3285-16
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
EW4Q8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.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...	10.522	14.1	ng/ul	1201600	2	10.581	1709470	20.0
Phenol, 3-(1-me...	10.952	3.5	ng/ul	301783	2	10.581	1709470	20.0
Phenol, p-tert-...	11.940	45.6	ng/ul	3901440	2	10.581	1709470	20.0
Phenol, 2,4-bis...	13.275	2.9	ng/ul	350831	3	14.422	2420950	20.0
Tetracosane, 1-...	25.057	2.1	ng/ul	251642	6	23.669	2394200	20.0