

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
 Data File : BP006346.D
 Acq On : 26 Jul 2021 19:42
 Operator : CG/JU
 Sample : M3063-02DL 20X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AT9DL

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-BP072421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP006346.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.711	103	106	117	rBV	58188	102977	0.63%	0.203%
2	4.022	154	159	168	rVB	55515	80234	0.49%	0.158%
3	5.810	458	463	470	rBV	32733	52886	0.32%	0.104%
4	6.287	536	544	550	rBV2	18333	32862	0.20%	0.065%
5	6.775	622	627	632	rBV2	12210	23972	0.15%	0.047%
6	6.957	650	658	663	rBV2	63340	109859	0.67%	0.216%
7	7.128	681	687	692	rBV	220945	347530	2.13%	0.684%
8	7.181	692	696	703	rVB	36840	62196	0.38%	0.122%
9	7.316	714	719	725	rVB	200730	317657	1.95%	0.625%
10	7.440	734	740	746	rVB	17718	29511	0.18%	0.058%
11	7.505	746	751	760	rBV6	11859	28175	0.17%	0.055%
12	7.687	779	782	789	rVB	14468	22407	0.14%	0.044%
13	7.787	789	799	806	rBV	835962	1325609	8.12%	2.609%
14	7.899	814	818	826	rVB	73445	115916	0.71%	0.228%
15	8.099	844	852	856	rBV5	12013	23345	0.14%	0.046%
16	8.157	856	862	869	rBV	85047	144545	0.89%	0.285%
17	8.275	877	882	886	rBV2	26153	40129	0.25%	0.079%
18	8.322	886	890	897	rVB3	13597	24678	0.15%	0.049%
19	8.428	903	908	912	rBV6	23229	38586	0.24%	0.076%
20	8.493	912	919	927	rVB	176499	293315	1.80%	0.577%
21	8.587	927	935	937	rBV3	17525	28337	0.17%	0.056%
22	8.628	937	942	948	rVB2	39413	70585	0.43%	0.139%
23	8.893	980	987	990	rBV	35585	57193	0.35%	0.113%
24	8.940	990	995	1009	rVV2	248965	508764	3.12%	1.001%
25	9.134	1021	1028	1035	rVB5	17447	37136	0.23%	0.073%
26	9.228	1035	1044	1058	rBV6	23283	65994	0.40%	0.130%
27	9.410	1070	1075	1080	rVB	16211	25463	0.16%	0.050%
28	9.469	1080	1085	1090	rBV2	13835	25229	0.15%	0.050%
29	9.657	1111	1117	1125	rBV	174825	302879	1.85%	0.596%
30	9.769	1130	1136	1141	rVV3	17401	31972	0.20%	0.063%
31	9.822	1141	1145	1150	rVV	41173	69153	0.42%	0.136%
32	9.975	1164	1171	1180	rBV3	55201	106679	0.65%	0.210%
33	10.193	1202	1208	1217	rVV	288622	475232	2.91%	0.935%
34	10.275	1217	1222	1229	rVB7	15957	29472	0.18%	0.058%
35	10.363	1230	1237	1244	rBV	149385	237339	1.45%	0.467%
36	10.569	1263	1272	1279	rVV	1172733	1957998	11.99%	3.854%

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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-BP072421.M
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37	10.622	1279	1281	1287	rVV3	20980	37157	0.23%	0.073%
38	10.710	1287	1296	1307	rVB2	292052	644227	3.95%	1.268%
39	10.951	1330	1337	1346	rBV2	6417	21803	0.13%	0.043%
40	11.040	1346	1352	1360	rVB4	20633	40052	0.25%	0.079%
41	11.151	1360	1371	1374	rBV6	18148	37283	0.23%	0.073%
42	11.487	1423	1428	1434	rVB4	18739	32404	0.20%	0.064%
43	11.681	1455	1461	1468	rBV3	46720	83859	0.51%	0.165%
44	11.969	1502	1510	1517	rVB7	31511	72573	0.44%	0.143%
45	12.128	1532	1537	1542	rBV	62722	100600	0.62%	0.198%
46	12.451	1585	1592	1597	rVV3	80723	154441	0.95%	0.304%
47	12.540	1601	1607	1613	rVV7	18991	49226	0.30%	0.097%
48	12.598	1613	1617	1624	rVB10	12036	23724	0.15%	0.047%
49	12.840	1649	1658	1663	rBV3	92373	185734	1.14%	0.366%
50	12.904	1665	1669	1675	rVB	84022	126808	0.78%	0.250%
51	13.116	1695	1705	1718	rBV2	229753	465767	2.85%	0.917%
52	13.251	1720	1728	1733	rVB	102542	165406	1.01%	0.326%
53	13.387	1745	1751	1753	rBV2	37369	59730	0.37%	0.118%
54	13.428	1753	1758	1762	rVV	184862	285486	1.75%	0.562%
55	13.469	1762	1765	1772	rVB2	60603	92233	0.56%	0.182%
56	13.545	1773	1778	1782	rBV	57351	93281	0.57%	0.184%
57	13.587	1782	1785	1789	rVB4	32415	46945	0.29%	0.092%
58	13.834	1821	1827	1833	rVV	671412	905562	5.55%	1.783%
59	13.969	1846	1850	1853	rBV	25715	29596	0.18%	0.058%
60	14.016	1853	1858	1865	rVB6	21583	44929	0.28%	0.088%
61	14.104	1866	1873	1884	rBV	671320	1036410	6.35%	2.040%
62	14.187	1884	1887	1892	rBV2	15479	25314	0.16%	0.050%
63	14.410	1915	1925	1934	rBV	1694025	2572897	15.76%	5.065%
64	14.504	1938	1941	1945	rVV2	30397	45171	0.28%	0.089%
65	14.551	1945	1949	1952	rVV6	21589	41205	0.25%	0.081%
66	14.604	1952	1958	1968	rVB4	61376	162562	1.00%	0.320%
67	14.810	1989	1993	2000	rVB	45980	75740	0.46%	0.149%
68	14.886	2001	2006	2010	rVB4	32630	49575	0.30%	0.098%
69	14.951	2010	2017	2021	rBV2	107180	187516	1.15%	0.369%
70	14.992	2021	2024	2027	rVV2	54023	74868	0.46%	0.147%
71	15.034	2027	2031	2038	rVB	339433	517071	3.17%	1.018%
72	15.151	2048	2051	2055	rBV3	18273	22894	0.14%	0.045%
73	15.245	2061	2067	2072	rVB9	26841	57298	0.35%	0.113%
74	15.345	2076	2084	2088	rBV7	25095	57312	0.35%	0.113%
75	15.410	2089	2095	2100	rVV	903238	1294846	7.93%	2.549%
76	15.463	2100	2104	2108	rVB	139068	190840	1.17%	0.376%

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 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

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77	15.539	2108	2117	2124	rBV	1528481	2375138	14.55%	4.675%
78	15.622	2128	2131	2135	rVB3	36008	43813	0.27%	0.086%
79	15.751	2149	2153	2161	rVB4	37704	69214	0.42%	0.136%
80	15.833	2162	2167	2171	rBV4	25120	45396	0.28%	0.089%
81	16.151	2217	2221	2229	rVB9	16849	35628	0.22%	0.070%
82	16.228	2229	2234	2238	rBV8	19970	35145	0.22%	0.069%
83	16.451	2267	2272	2278	rBV3	41236	85586	0.52%	0.168%
84	16.516	2279	2283	2287	rVB	52988	68810	0.42%	0.135%
85	16.675	2306	2310	2314	rBV6	19696	33639	0.21%	0.066%
86	16.816	2327	2334	2349	rBV	11360639	16329080	100.00%	32.144%
87	16.933	2350	2354	2360	rVV6	71867	151438	0.93%	0.298%
88	16.980	2360	2362	2368	rVV	36825	64430	0.39%	0.127%
89	17.039	2368	2372	2380	rVB	78956	127240	0.78%	0.250%
90	17.163	2387	2393	2399	rBV	1955488	2778332	17.01%	5.469%
91	17.263	2404	2410	2422	rBV	997070	1450308	8.88%	2.855%
92	17.369	2423	2428	2437	rVB	75760	122674	0.75%	0.241%
93	17.539	2454	2457	2461	rBV3	23357	25751	0.16%	0.051%
94	18.069	2543	2547	2553	rBV2	97447	133778	0.82%	0.263%
95	19.251	2744	2748	2754	rBV	378440	417811	2.56%	0.822%
96	19.304	2754	2757	2766	rVB2	71381	97045	0.59%	0.191%
97	19.504	2786	2791	2797	rBV	1178110	1553455	9.51%	3.058%
98	21.257	3085	3089	3097	rBV	2278562	2910088	17.82%	5.728%
99	23.451	3457	3462	3470	rVB2	768548	1416983	8.68%	2.789%
100	23.604	3482	3488	3498	rVB2	1599324	2973445	18.21%	5.853%

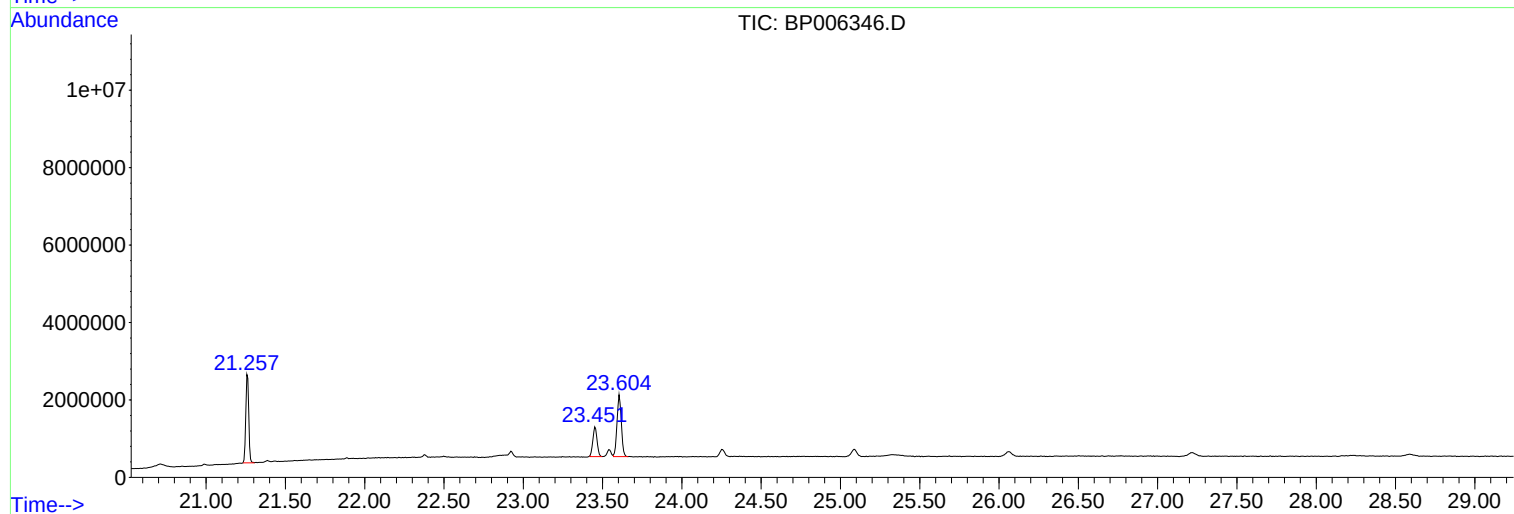
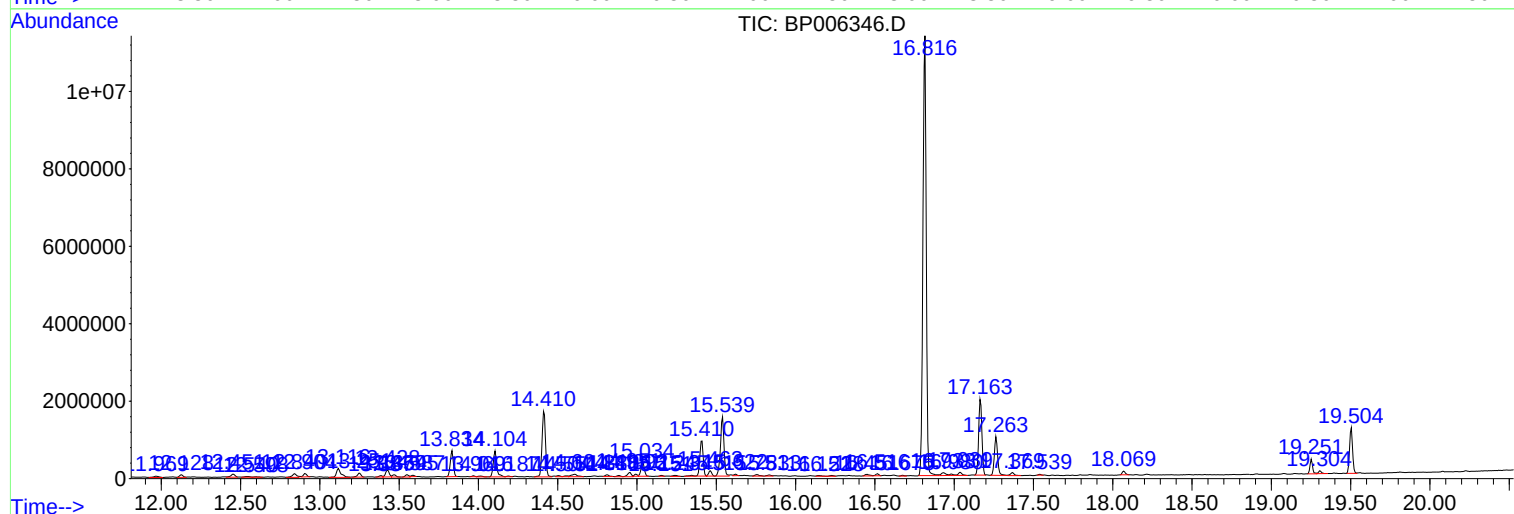
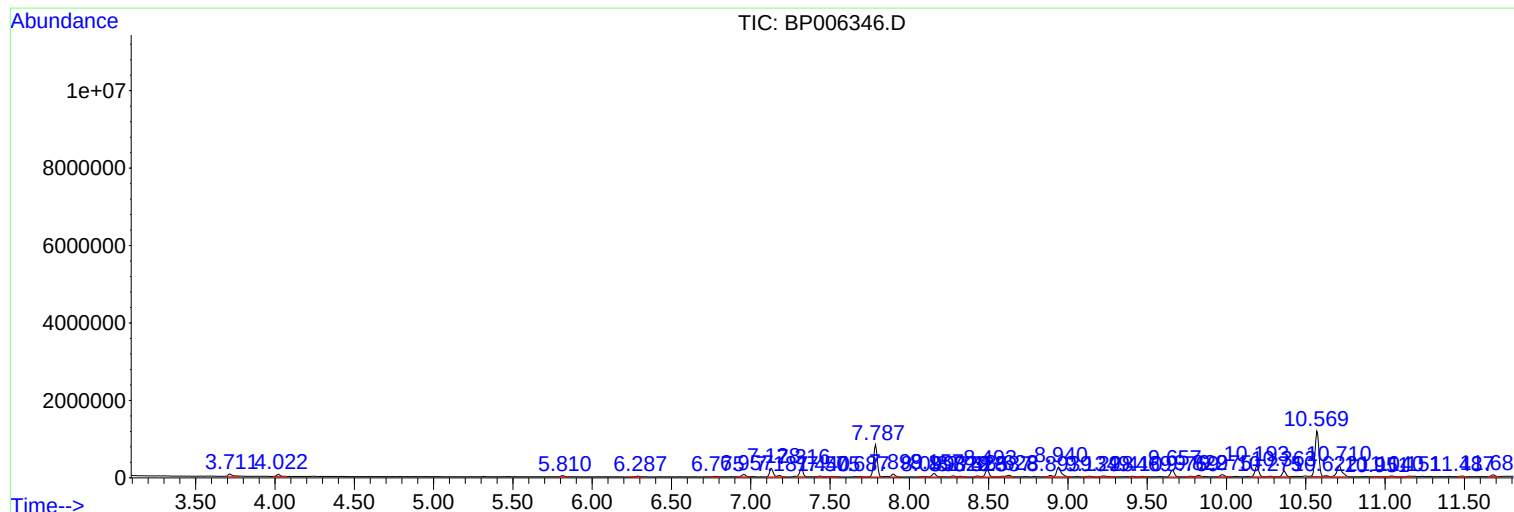
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ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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



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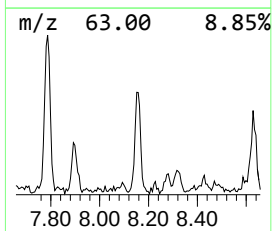
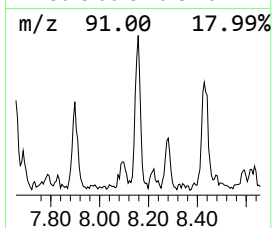
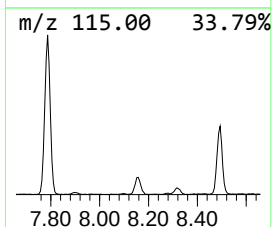
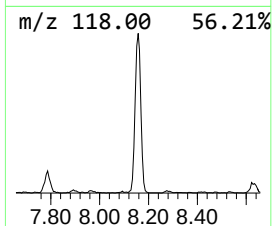
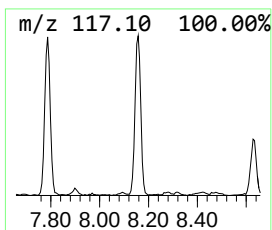
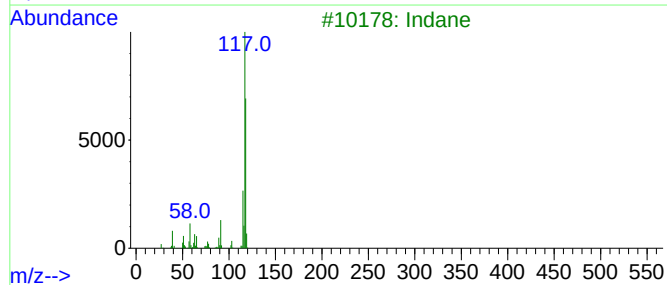
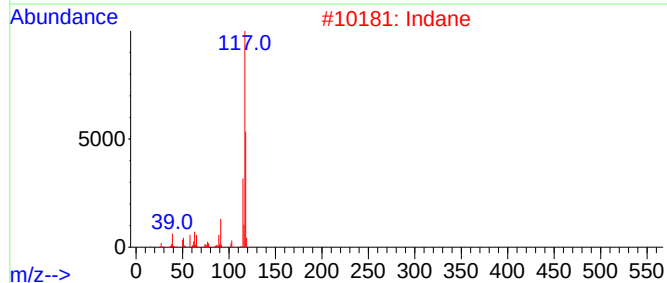
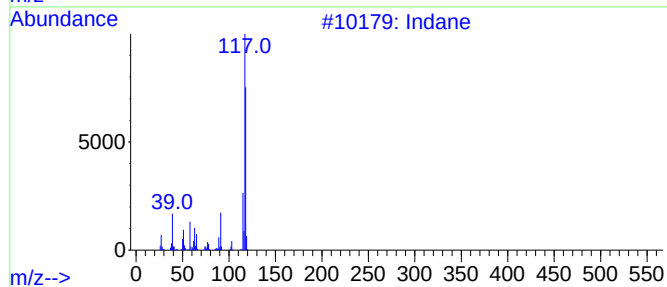
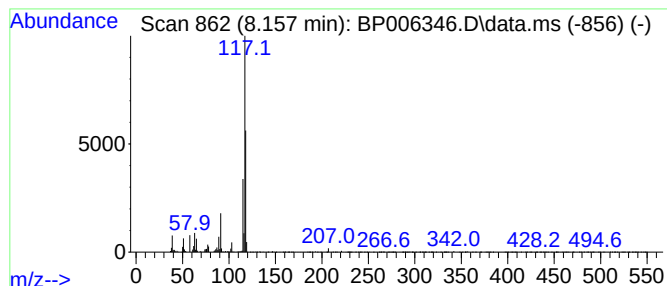
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.160	2.18 ng/ul	144545	1,4-Dichlorobenzene-d4	7.787

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	93
2	Indane			118	C9H10	000496-11-7	93
3	Indane			118	C9H10	000496-11-7	91
4	Benzene, cyclopropyl-			118	C9H10	000873-49-4	83
5	Benzene, 1-propenyl-			118	C9H10	000637-50-3	83



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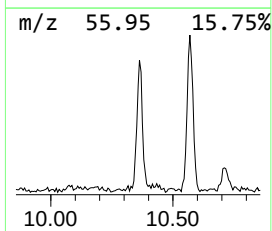
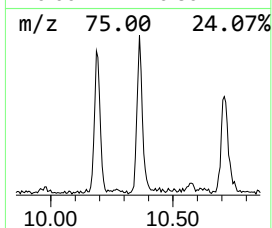
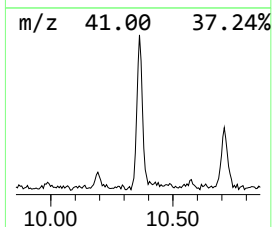
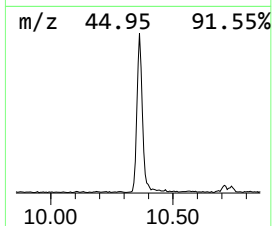
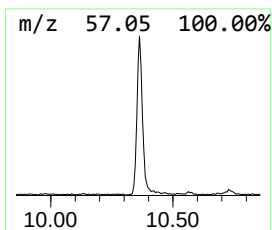
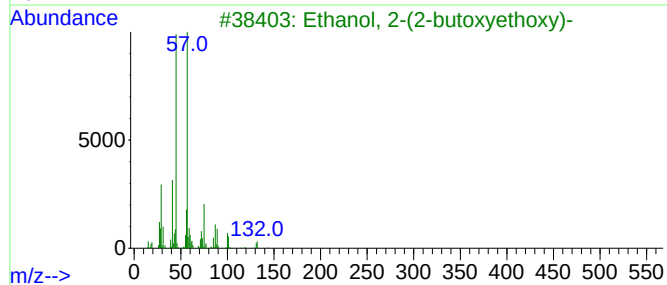
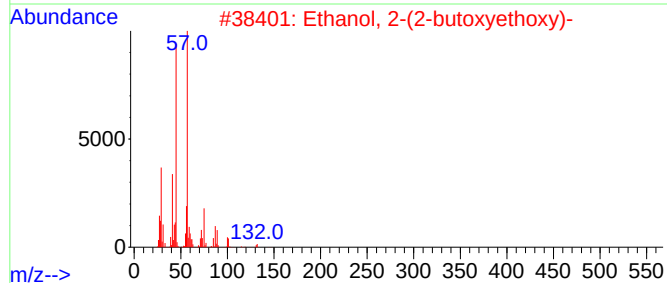
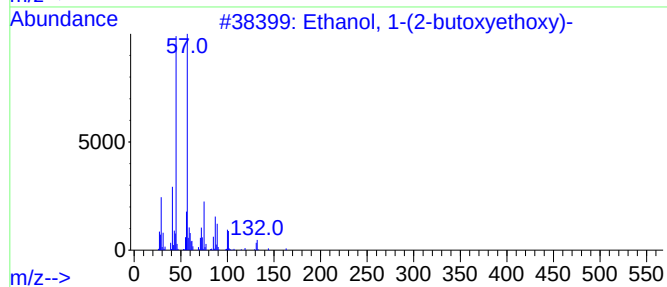
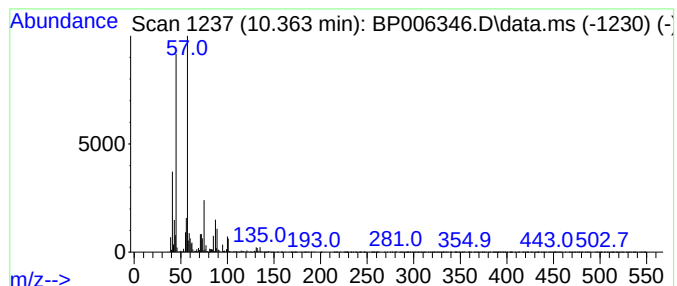
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethanol, 1-(2-butoxyethoxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.360	2.42 ng/ul	237339	Naphthalene-d8	10.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	78



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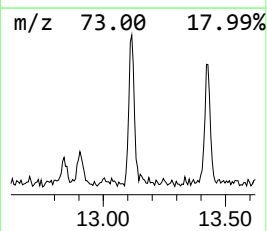
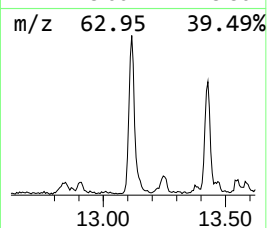
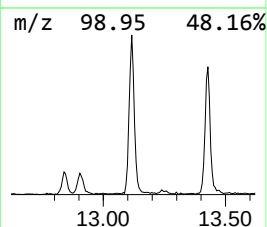
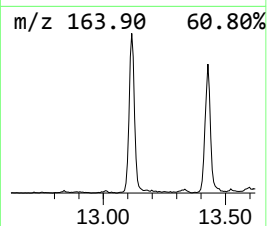
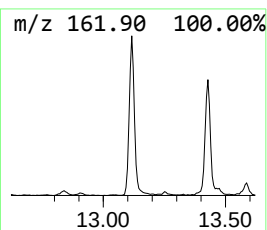
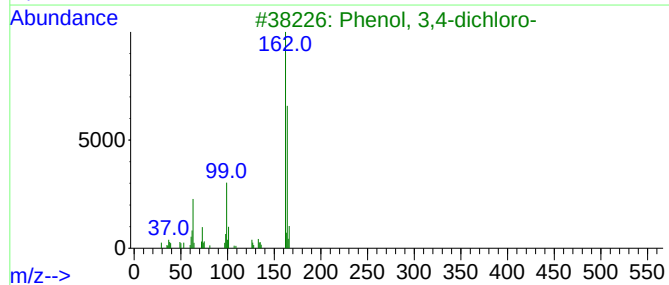
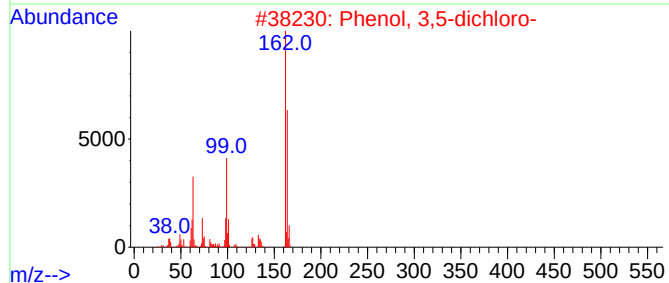
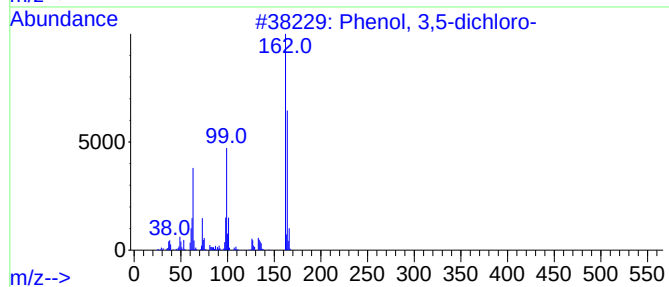
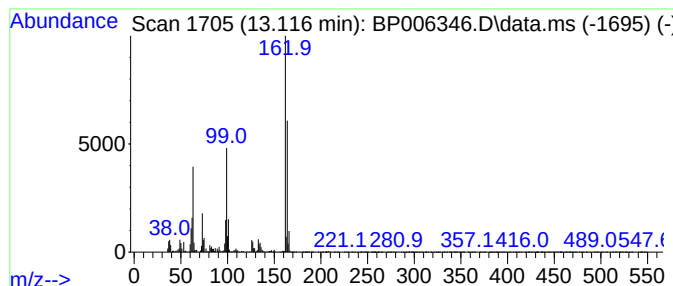
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Phenol, 3,5-dichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.120	3.62 ng/ul	465767	Acenaphthene-d10	14.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	98
2			Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	98
3			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97
4			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	96
5			Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
Data File : BP006346.D
Acq On : 26 Jul 2021 19:42
Operator : CG/JU
Sample : M3063-02DL 20X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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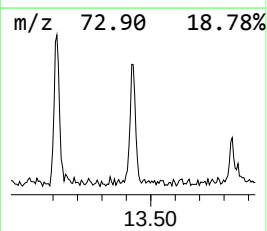
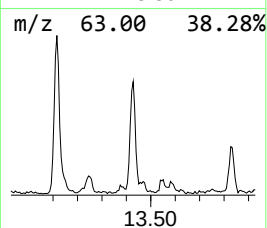
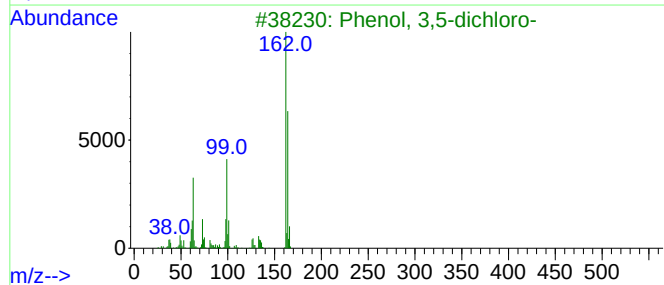
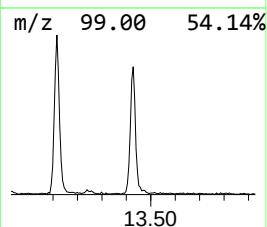
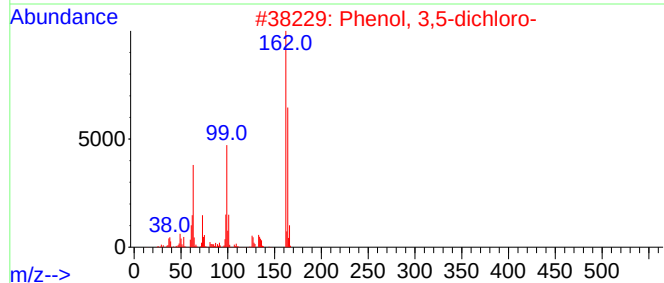
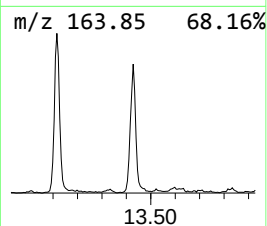
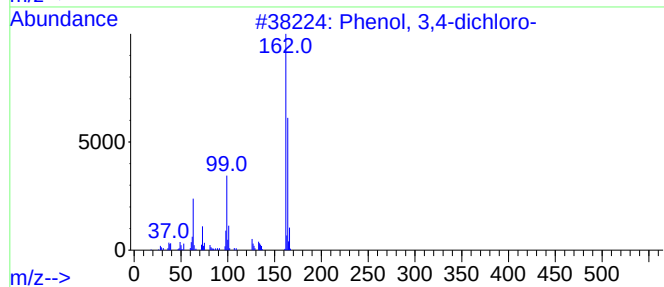
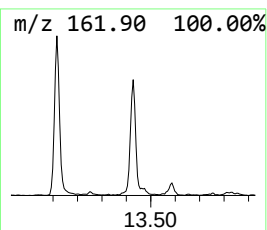
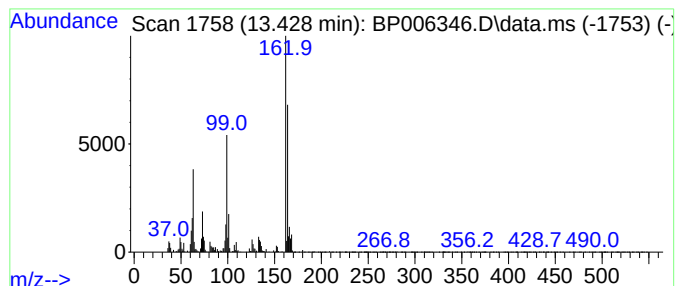
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenol, 3,4-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.430	2.22 ng/ul	285486	Acenaphthene-d10	14.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97
2			Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	95
3			Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	95
4			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	95
5			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
Data File : BP006346.D
Acq On : 26 Jul 2021 19:42
Operator : CG/JU
Sample : M3063-02DL 20X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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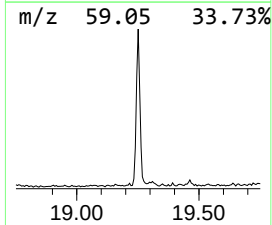
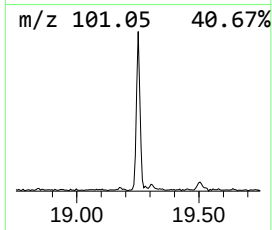
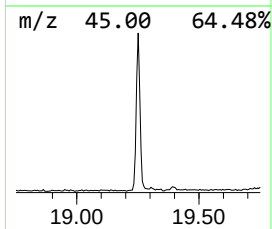
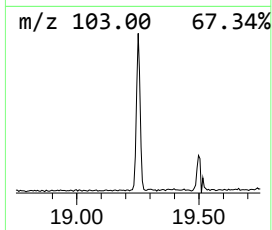
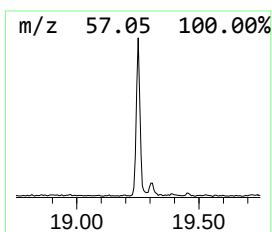
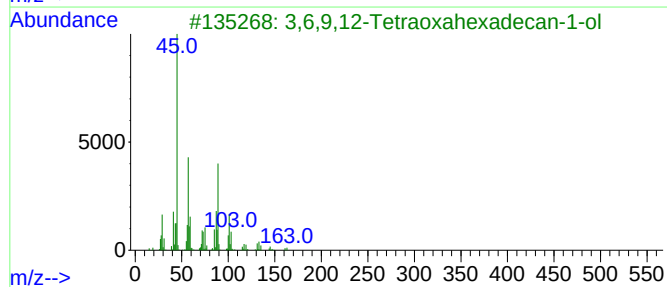
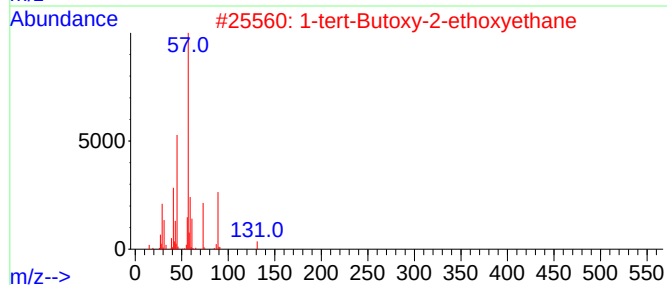
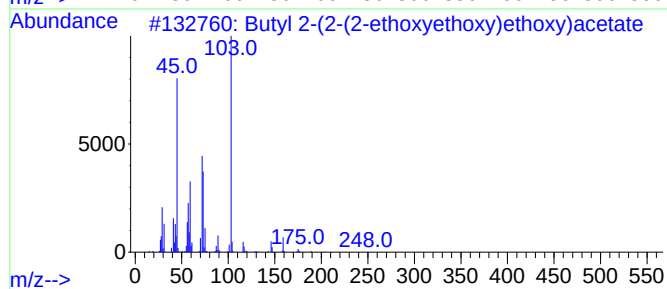
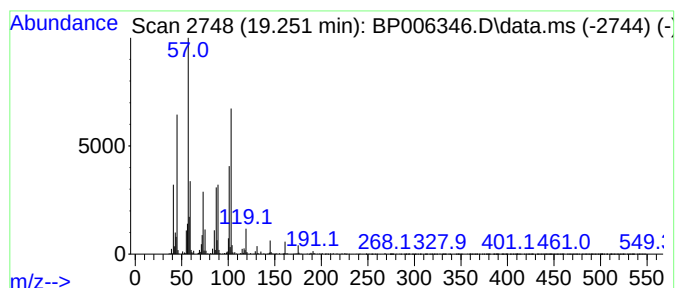
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.250	2.87 ng/ul	417811	Chrysene-d12	21.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl 2-(2-(2-ethoxyethoxy)ethox...	248	C12H24O5	1000366-77-5	38		
2	1-tert-Butoxy-2-ethoxyethane	146	C8H18O2	051422-54-9	38		
3	3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	35		
4	3,6,9,12,15-Pentaoxanonadecan-1-ol	294	C14H30O6	001786-94-3	35		
5	Ethanol, 2-[2-(2-methoxyethoxy)e...	164	C7H16O4	000112-35-6	27		



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Operator : CG/JU
Sample : M3063-02DL 20X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AT9DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Indane	8.160	2.2	ng/ul	144545	1	7.787	1325610	20.0
Ethanol, 1-(2-b...	10.360	2.4	ng/ul	237339	2	10.569	1958000	20.0
Phenol, 3,5-dic...	13.120	3.6	ng/ul	465767	3	14.410	2572900	20.0
Phenol, 3,4-dic...	13.430	2.2	ng/ul	285486	3	14.410	2572900	20.0
unknown-01	19.250	2.9	ng/ul	417811	5	21.257	2910090	20.0