LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023444.D
Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

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-BP112624.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP023444.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.016	3	5	10	rVB	2977089	2802456	100.00%	8.451%
2	3.346	58	61	66	rVB	33632	35973	1.28%	0.108%
3	3.816	137	141	146	rVB2	18052	25528	0.91%	0.077%
4	4.234	207	212	219	rVB	13578	22273	0.79%	0.067%
5	4.628	275	279	287	rVB3	12844	18982	0.68%	0.057%
6	5.099	353	359	366	rBV	98597	136519	4.87%	0.412%
7	5.228	374	381	382	rBV	22501	31555	1.13%	0.095%
8	5.587	437	442	450	rVB4	17324	31133	1.11%	0.094%
9	6.046	514	520	527	rBV	163591	262885	9.38%	0.793%
10	6.110	527	531	534	rVV2	13790	21890	0.78%	0.066%
11	6.187	540	544	546	rVV3	10431	16962	0.61%	0.051%
12	6.216	546	549	557	rVV2	11588	18162	0.65%	0.055%
13	6.534	599	603	608	rVV	25834	36967	1.32%	0.111%
14	6.587	608	612	616	rVB2	12839	17818	0.64%	0.054%
15	6.634	616	620	624	rBV3	13853	23181	0.83%	0.070%
16	6.693	624	630	635	rVV4	25651	54442	1.94%	0.164%
17	7.151	702	708	711	rBV	44555	71118	2.54%	0.214%
18	7.181	711	713	720	rVV2	16972	26359	0.94%	0.079%
19	7.340	734	740	746	rVB7	7685	18761	0.67%	0.057%
20	7.434	746	756	761	rBV8	13381	38205	1.36%	0.115%
21	7.528	767	772	787	rVB	619135	1103272	39.37%	3.327%
22	7.640	787	791	793	rVV4	11543	17961	0.64%	0.054%
23	7.698	797	801	805	rVV5	15910	25600	0.91%	0.077%
24	7.769	810	813	815	rVV4	10867	16908	0.60%	0.051%
25	7.793	815	817	824	rVB6	11672	18559	0.66%	0.056%
26	8.004	848	853	856	rBV2	30355	50067	1.79%	0.151%
27	8.057	859	862	866	rVB	26787	36585	1.31%	0.110%
28	8.151	872	878	884	rBV	183413	321109	11.46%	0.968%
29	8.198	884	886	893	rVB3	16188	19040	0.68%	0.057%
30	8.263	893	897	900	rBV	27401	36389	1.30%	0.110%
31	8.310	900	905	911	rVB	53825	84925	3.03%	0.256%
32	8.457	924	930	934	rBV	172982	273781	9.77%	0.826%
33	8.510	934	939	945	rVB	232322	357109	12.74%	1.077%
34	8.604	945	955	963	rBV	762420	1165629	41.59%	3.515%
35	8.687	963	969	972	rBV3	80958	168050	6.00%	0.507%
36	8.722	972	975	985	rVB	96092	164320	5.86%	0.495%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023444.D
Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

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-BP112624.MA.M
Title : SVOA CALIBRATION

37	8.845	985	996	1004	rBV3	164661	374808	13.37%	1.130%
38	8.928	1004	1010	1015	rBV	109018	202191	7.21%	0.610%
39	8.975	1015	1018	1026	rVB5	82439	182218	6.50%	0.549%
40	9.075	1026	1035	1038	rBV5	61014	130058	4.64%	0.392%
41	9.122	1038	1043	1048	rBV	533876	799111	28.51%	2.410%
42	9.181	1048	1053	1060	rVB	811156	1207491	43.09%	3.641%
43	9.245	1060	1064	1068	rVB3	13546	18581	0.66%	0.056%
44	9.369	1075	1085	1089	rBV	407890	731547	26.10%	2.206%
45	9.416	1089	1093	1098	rVB	236263	367291	13.11%	1.108%
46	9.492	1098	1106	1109	rBV	326617	582670	20.79%	1.757%
47	9.534	1109	1113	1115	rBV	200455	264133	9.43%	0.796%
48	9.634	1123	1130	1131	rBV3	176370	262134	9.35%	0.790%
49	9.675	1131	1137	1143	rVV2	708197	1407440	50.22%	4.244%
50	9.745	1143	1149	1155	rVV2	488811	885021	31.58%	2.669%
51	9.840	1155	1165	1166	rVV	214988	376533	13.44%	1.135%
52	9.863	1166	1169	1174	rVV	278394	450930	16.09%	1.360%
53	9.904	1174	1176	1181	rVV2	72006	106293	3.79%	0.321%
54	9.975	1182	1188	1195	rVB	443464	681216	24.31%	2.054%
55	10.045	1195	1200	1205	rBV	27490	42313	1.51%	0.128%
56	10.098	1205	1209	1212	rVV2	22852	35507	1.27%	0.107%
57	10.163	1212	1220	1228	rVV6	70270	237404	8.47%	0.716%
58	10.275	1228	1239	1243	rVV	845749	1677810	59.87%	5.059%
59	10.322	1243	1247	1254	rVV2	365754	810881	28.93%	2.445%
60	10.381	1254	1257	1261	rVV3	40496	67448	2.41%	0.203%
61	10.422	1261	1264	1269	rVV3	21986	35365	1.26%	0.107%
62	10.487	1269	1275	1282	rVV	66503	127840	4.56%	0.385%
63	10.563	1282	1288	1292	rVV8	10367	18877	0.67%	0.057%
64	10.634	1296	1300	1303	rVV3	13306	19838	0.71%	0.060%
65	10.681	1303	1308	1314	rVV	29828	66343	2.37%	0.200%
66	10.928	1346	1350	1353	rVV5	9464	19600	0.70%	0.059%
67	11.169	1386	1391	1398	rVB6	11884	23474	0.84%	0.071%
68	11.281	1405	1410	1414	rBV3	23112	42479	1.52%	0.128%
69	11.686	1474	1479	1485	rBV4	13055	26284	0.94%	0.079%
70	12.328	1582	1588	1592	rBV8	10869	25425	0.91%	0.077%
71	12.539	1621	1624	1630	rVB6	11513	18978	0.68%	0.057%
72	12.651	1639	1643	1649	rVB2	35070	53898	1.92%	0.163%
73	12.892	1679	1684	1687	rBV3	28052	42961	1.53%	0.130%
74	13.539	1790	1794	1799	rVB	72027	98458	3.51%	0.297%
75	13.616	1802	1807	1811	rVB	51497	69540	2.48%	0.210%
76	13.869	1846	1850	1855	rBV7	34247	66085	2.36%	0.199%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023444.D
Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

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-BP112624.MA.M
Title : SVOA CALIBRATION

77	13.951	1861	1864	1872	rVB2	75802	121946	4.35%	0.368%
78	14.033	1874	1878	1884	rBV6	22830	46474	1.66%	0.140%
79	14.104	1885	1890	1891	rBV4	33004	50144	1.79%	0.151%
80	14.151	1892	1898	1905	rVB	1248832	1953940	69.72%	5.892%
81	14.669	1983	1986	1994	rVB6	32367	54376	1.94%	0.164%
82	14.869	2015	2020	2026	rBV9	34723	91931	3.28%	0.277%
83	14.939	2028	2032	2037	rVB2	74032	108740	3.88%	0.328%
84	15.427	2111	2115	2120	rVB	104001	167008	5.96%	0.504%
85	15.922	2190	2199	2205	rBV2	230948	533163	19.02%	1.608%
86	16.751	2336	2340	2345	rVB	170110	244177	8.71%	0.736%
87	16.816	2348	2351	2355	rVB	196486	258139	9.21%	0.778%
88	16.957	2369	2375	2380	rBV2	1451574	2263061	80.75%	6.824%
89	16.998	2380	2382	2387	rVB	186165	267278	9.54%	0.806%
90	17.121	2399	2403	2412	rVB	213172	364763	13.02%	1.100%
91	17.563	2473	2478	2487	rVB2	280175	629824	22.47%	1.899%
92	18.051	2558	2561	2566	rVB2	338710	462256	16.49%	1.394%
93	18.116	2568	2572	2576	rVV4	226271	360500	12.86%	1.087%
94	18.151	2576	2578	2584	rVB2	150534	143148	5.11%	0.432%
95	18.351	2609	2612	2616	rBV	259021	377088	13.46%	1.137%
96	19.092	2735	2738	2742	rBV	300921	406687	14.51%	1.226%
97	19.145	2743	2747	2752	rVB	599339	866235	30.91%	2.612%
98	21.386	3123	3128	3134	rBV	987319	1559476	55.65%	4.703%
99	24.556	3662	3667	3673	rVB	633855	1522824	54.34%	4.592%
100	27.256	4124	4126	4127	rBV	122219	82556	2.95%	0.249%

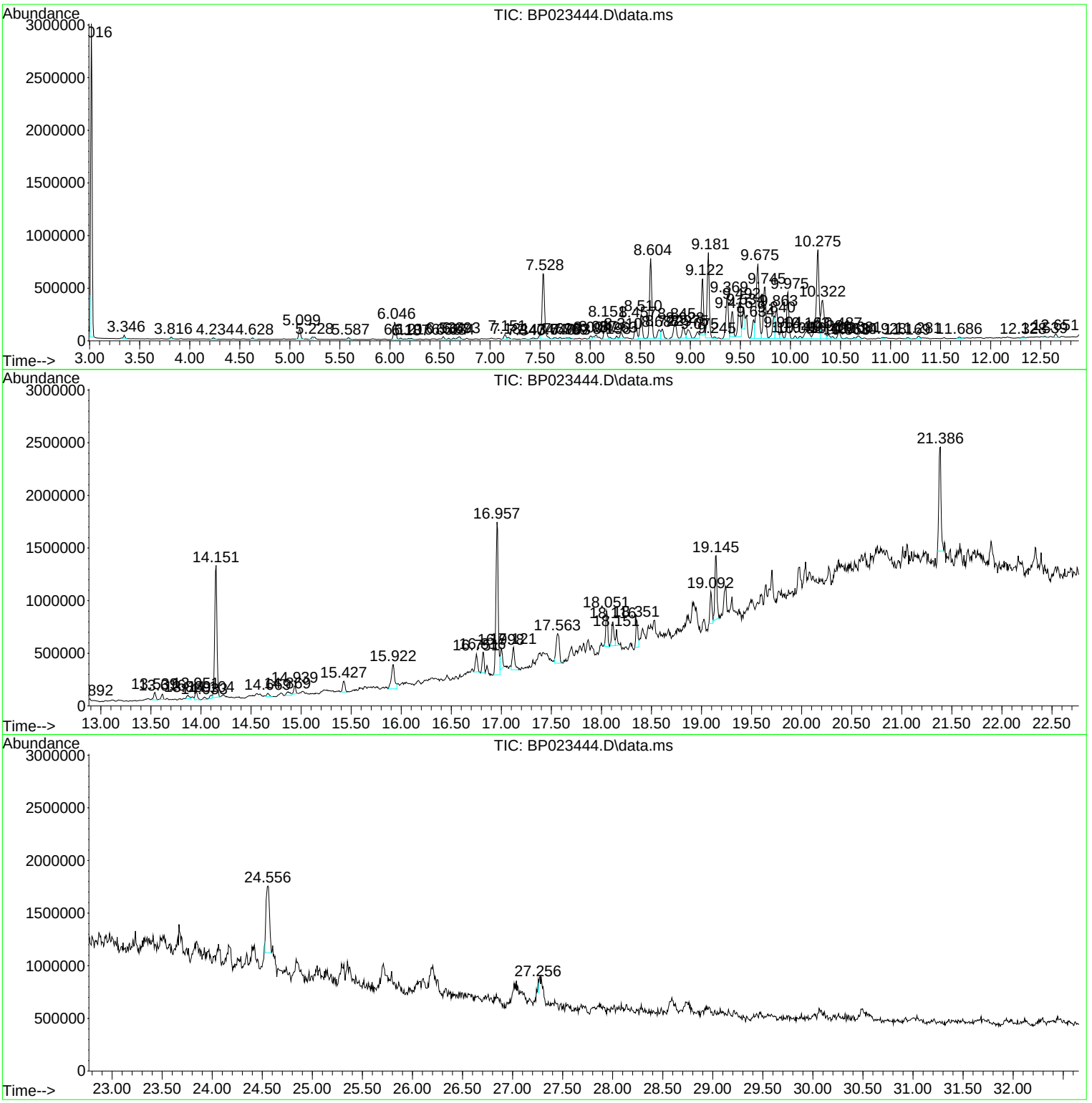
Sum of corrected areas: 33162681

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023444.D
Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023444.D
Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

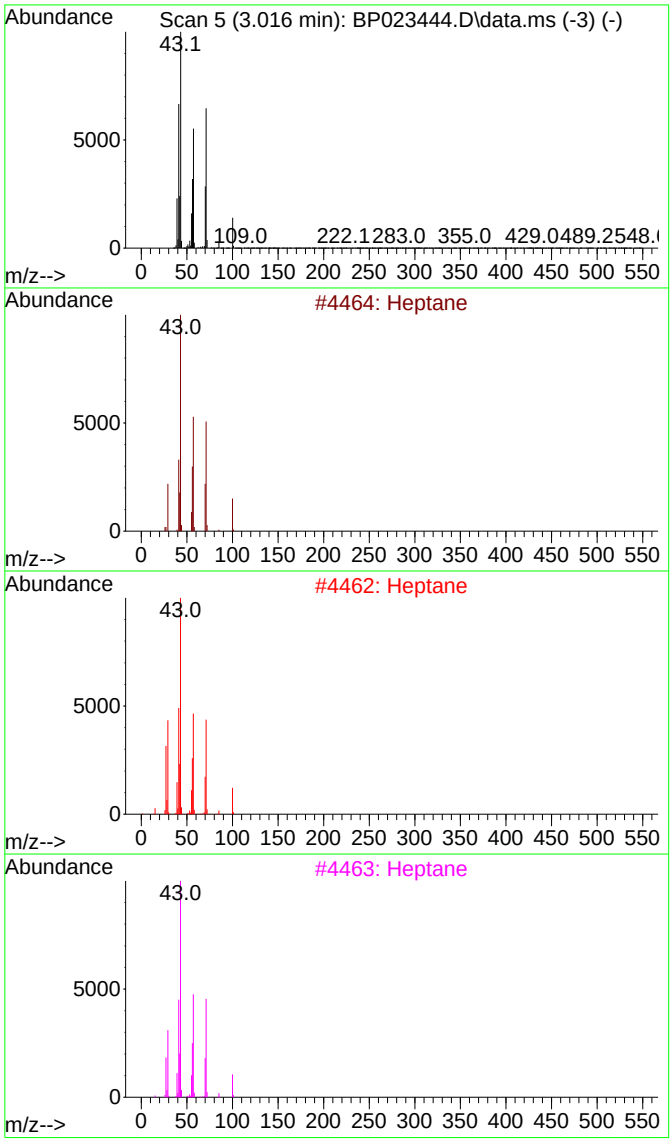
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

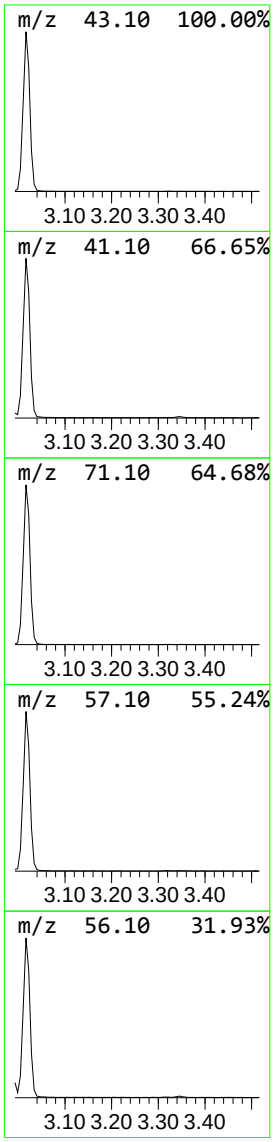
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.016	50.80 ng/ul	2802460	1,4-Dichlorobenzene-d4	7.534

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane	100	C7H16	000142-82-5	91
2	Heptane	100	C7H16	000142-82-5	90
3	Heptane	100	C7H16	000142-82-5	87
4	Heptane	100	C7H16	000142-82-5	86
5	Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	42





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023444.D
Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

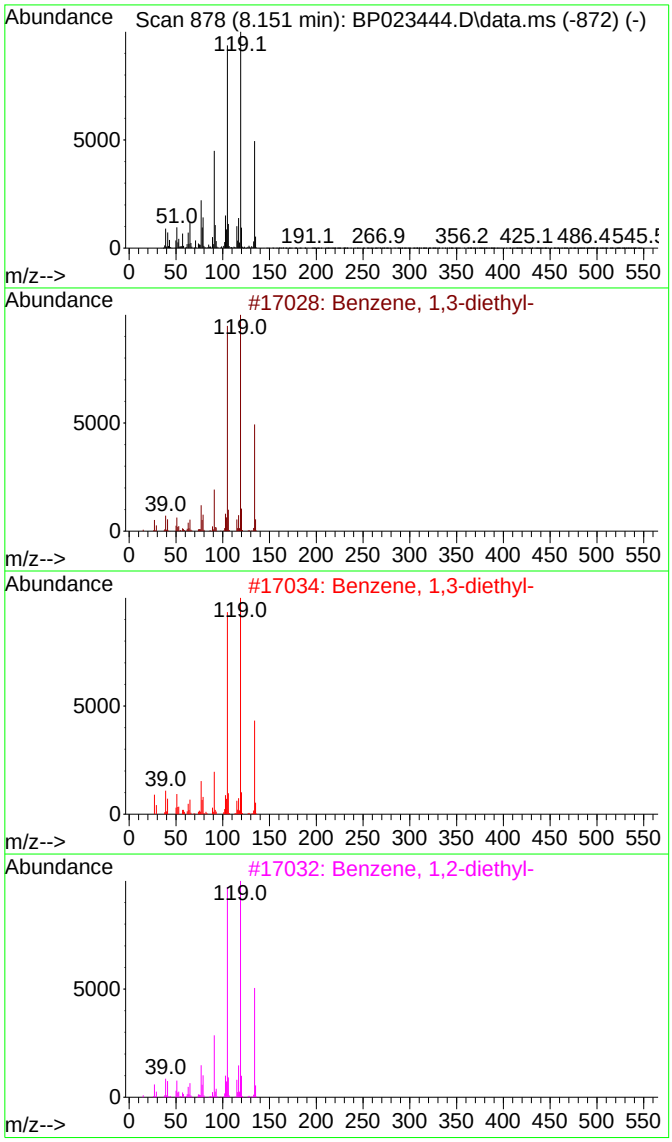
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

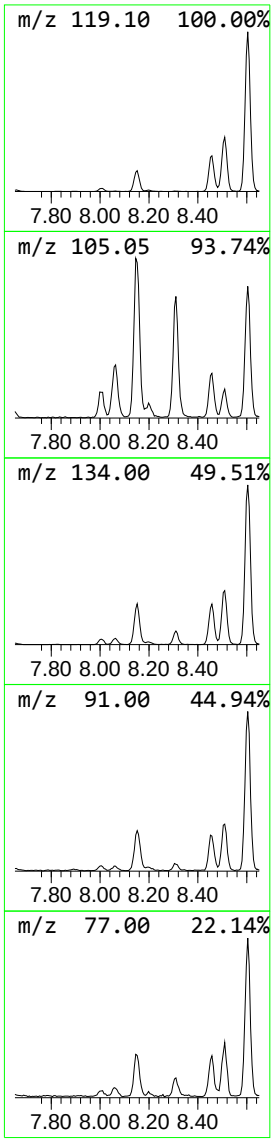
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1,3-diethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.151	5.82 ng/ul	321109	1,4-Dichlorobenzene-d4	7.534

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
2	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
3	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	94
4	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023444.D
Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

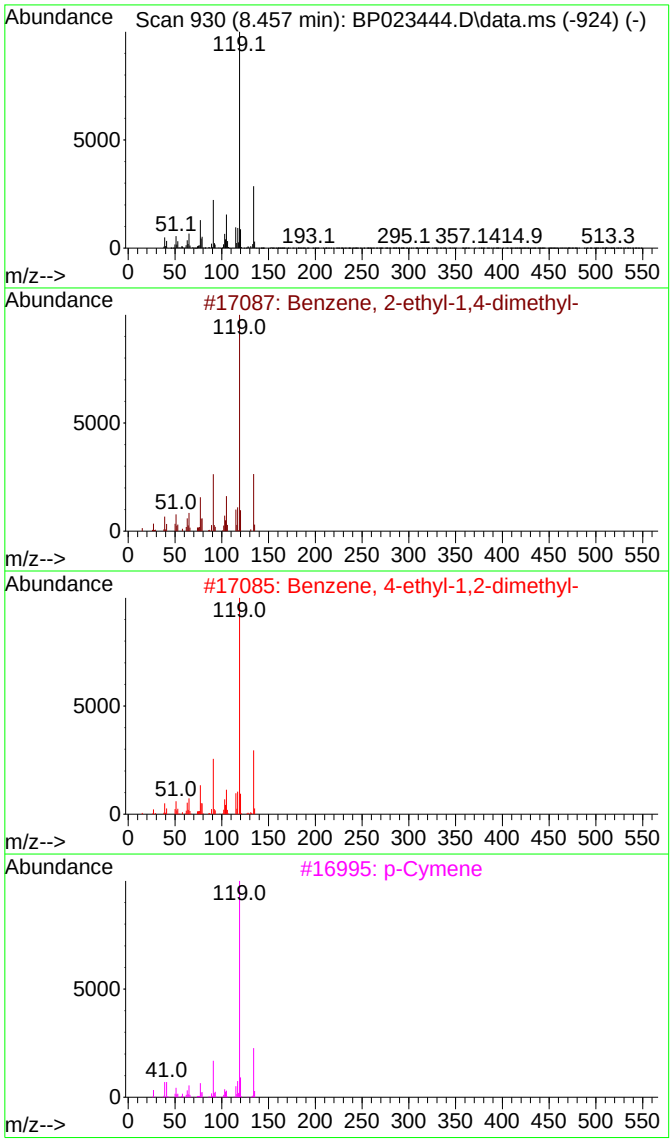
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

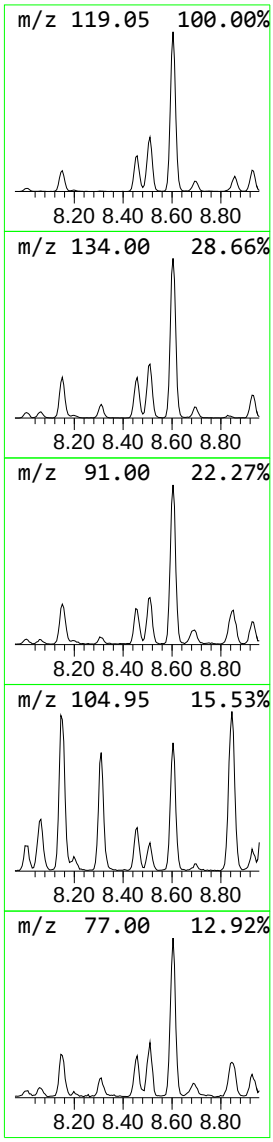
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.457	4.96 ng/ul	273781	1,4-Dichlorobenzene-d4	7.534

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3	p-Cymene	134	C10H14	000099-87-6	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023444.D
Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

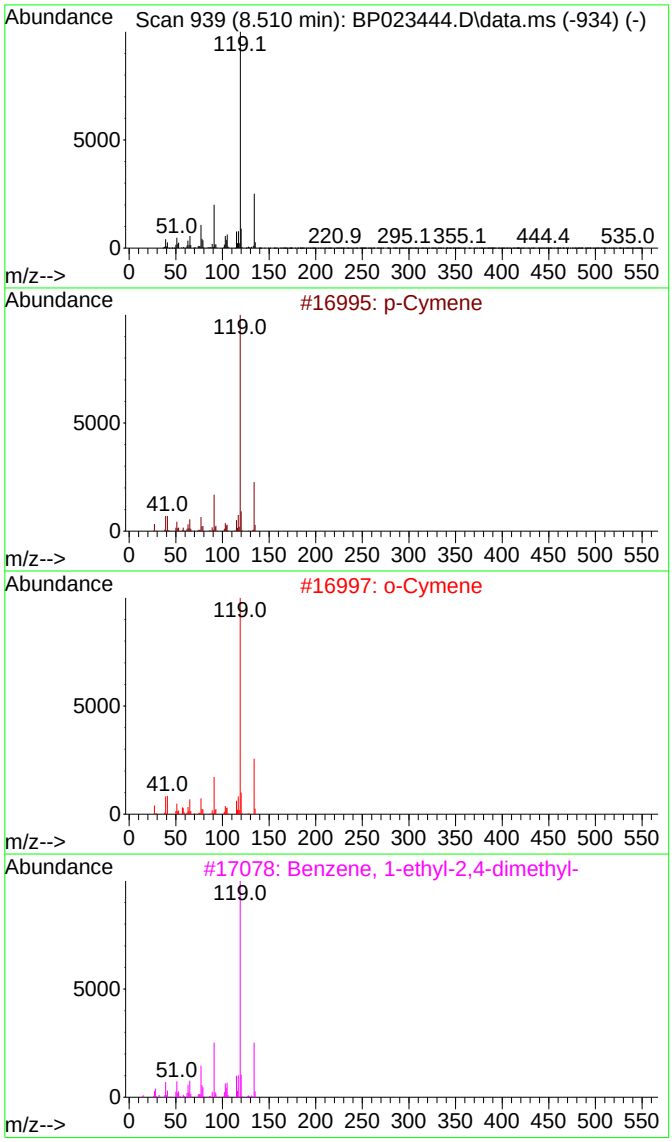
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

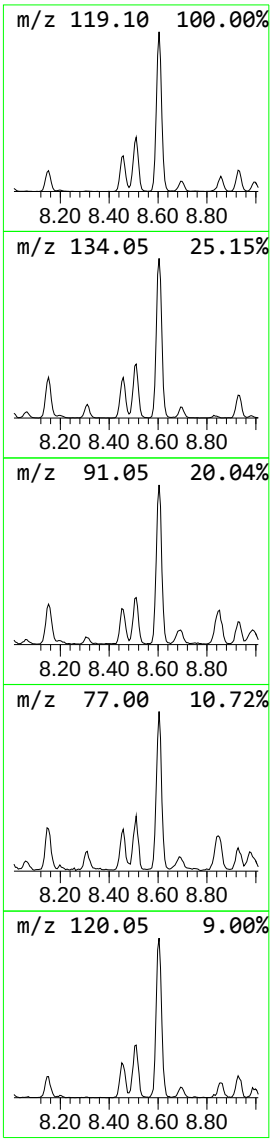
****

Peak Number 6 p-Cymene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.510	6.47 ng/ul	357109	1,4-Dichlorobenzene-d4	7.534

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
4	o-Cymene	134	C10H14	000527-84-4	97
5	o-Cymene	134	C10H14	000527-84-4	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023444.D
Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

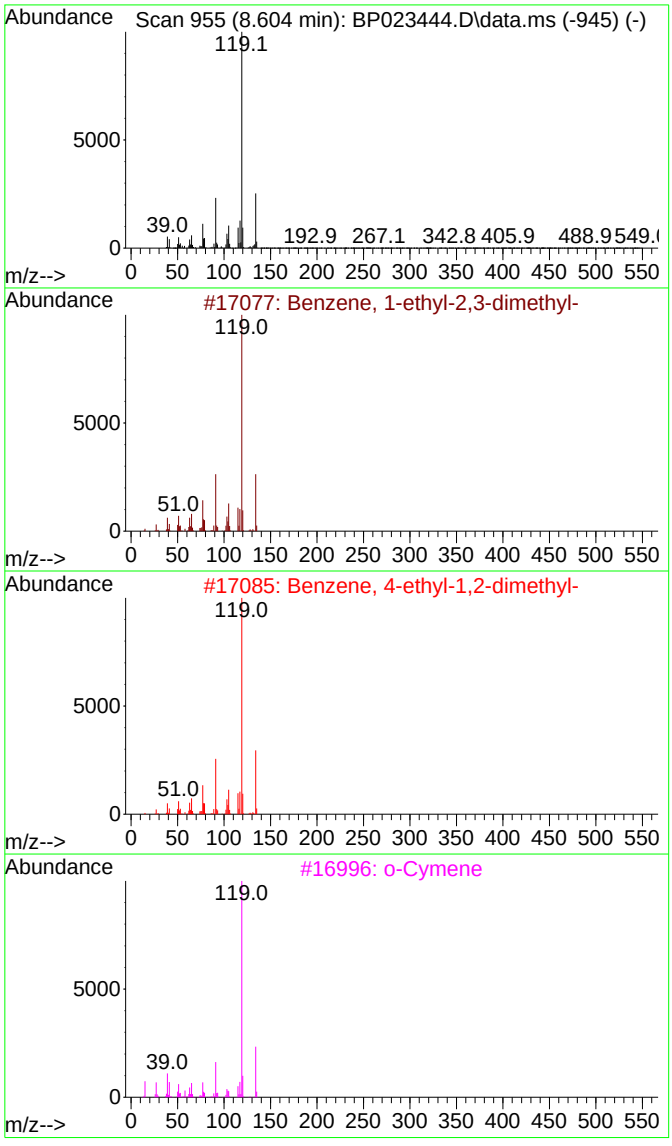
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

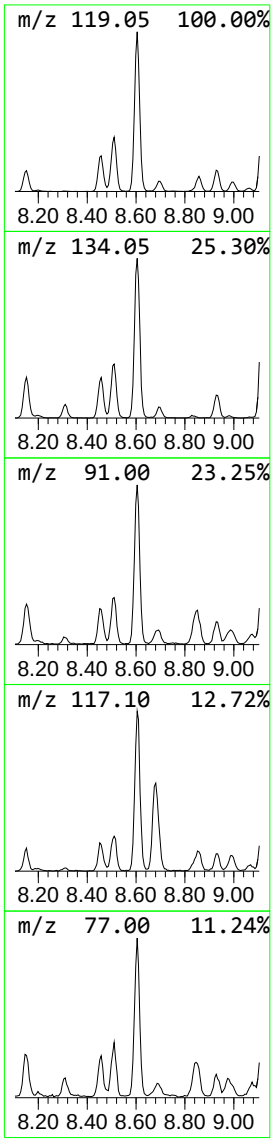
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.604	21.13 ng/ul	1165630	1,4-Dichlorobenzene-d4	7.534

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3	o-Cymene	134	C10H14	000527-84-4	95
4	p-Cymene	134	C10H14	000099-87-6	95
5	o-Cymene	134	C10H14	000527-84-4	95





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Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

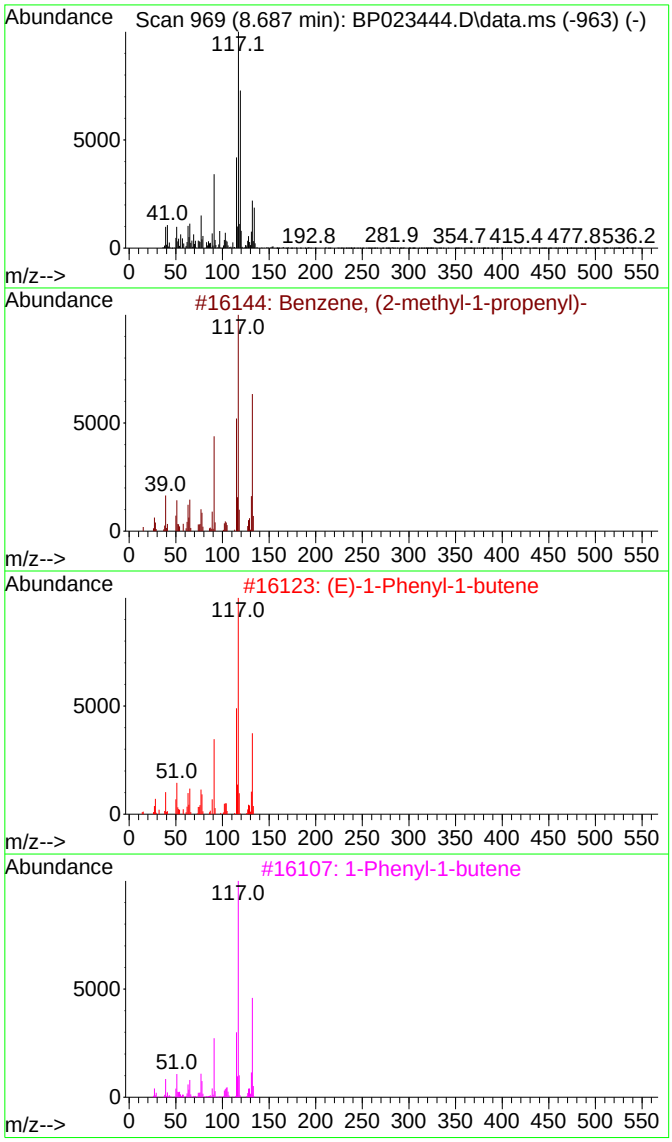
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

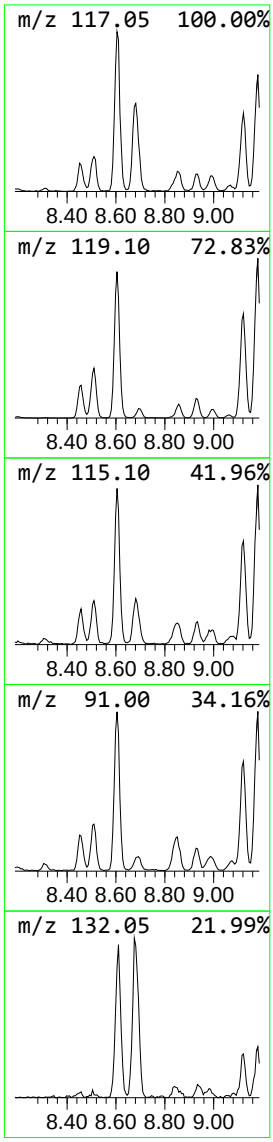
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, (2-methyl-1-propen... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.687	3.05 ng/ul	168050	1,4-Dichlorobenzene-d4	7.534

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
2	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	86
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	86
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	70





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Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

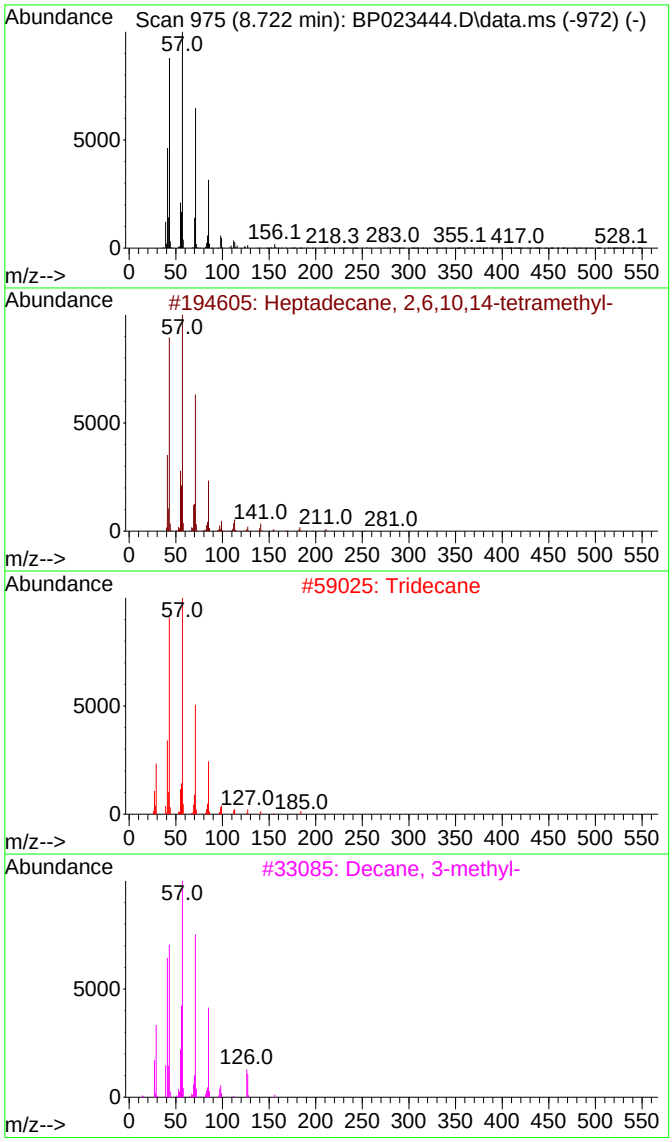
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

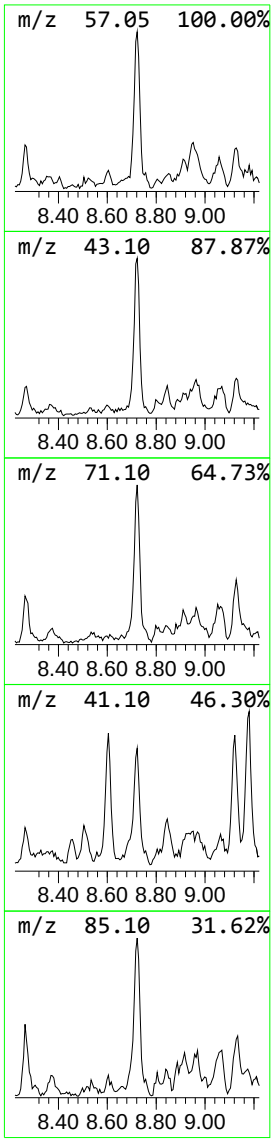
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.722	2.98 ng/ul	164320	1,4-Dichlorobenzene-d4	7.534

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	87
2	Tridecane	184	C13H28	000629-50-5	86
3	Decane, 3-methyl-	156	C11H24	013151-34-3	80
4	Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	78
5	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	72





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Acq On : 16 Dec 2024 21:51
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Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

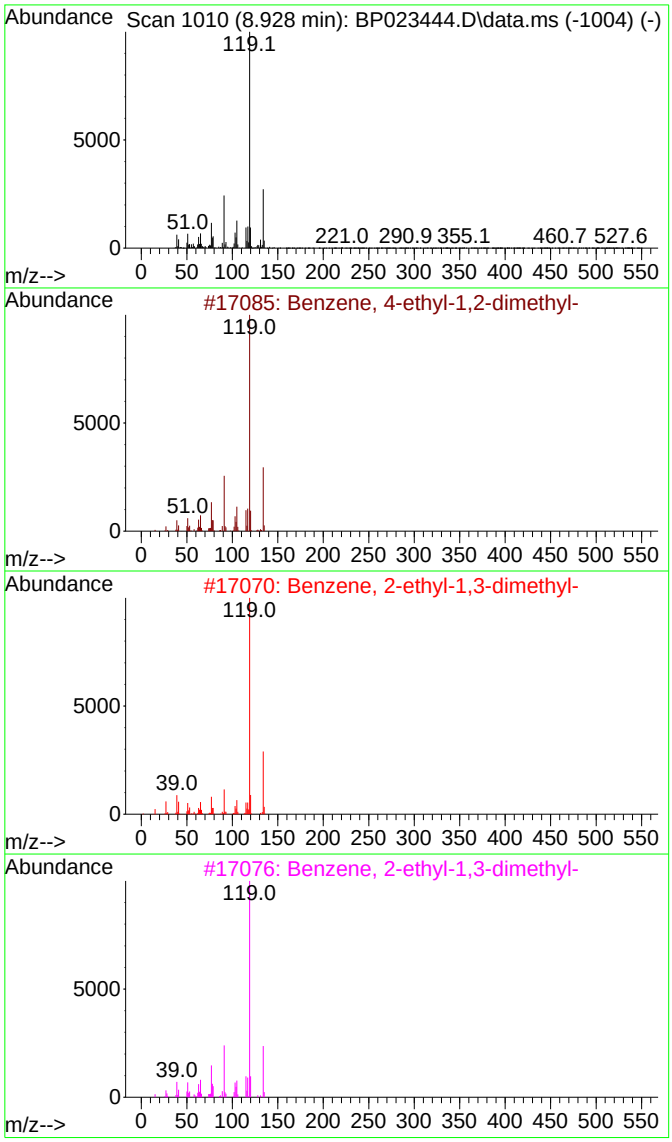
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

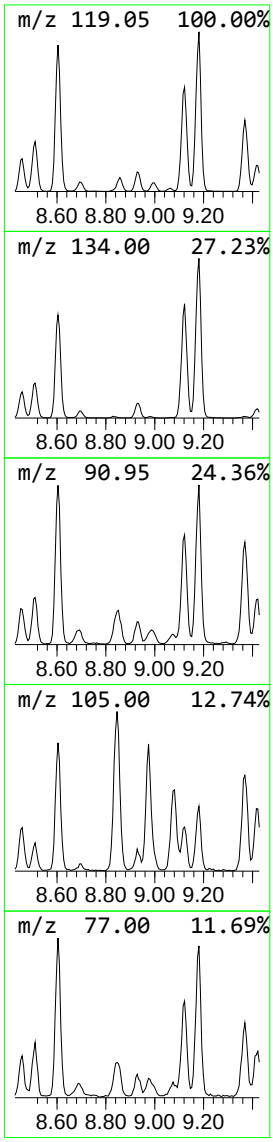
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.928	2.41 ng/ul	202191	Naphthalene-d8	10.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94





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Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

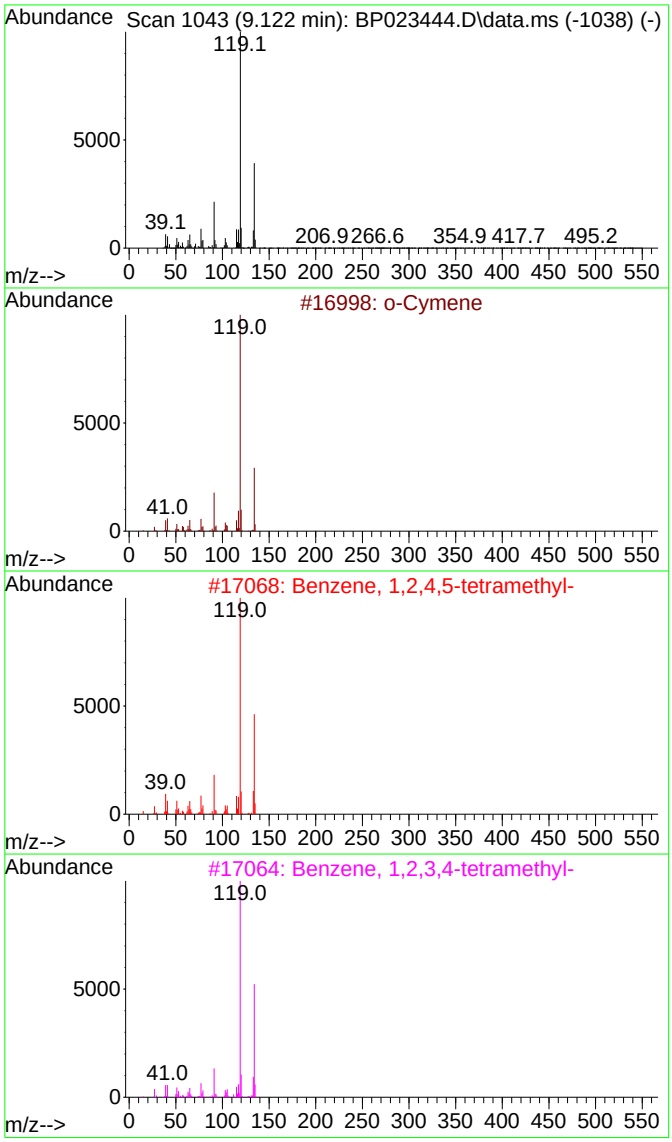
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

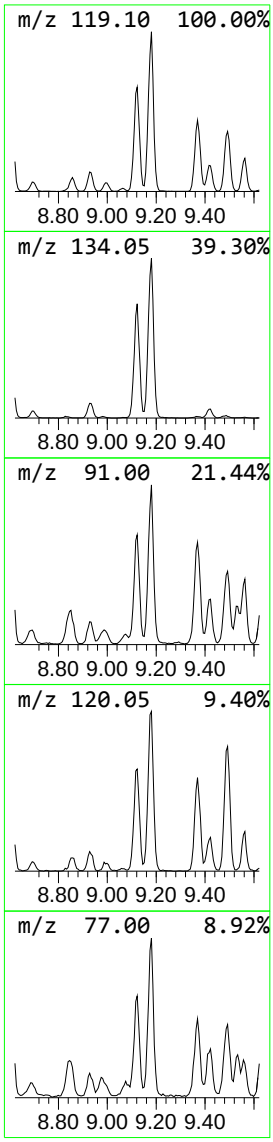
****

Peak Number 13 o-Cymene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.122	9.53 ng/ul	799111	Naphthalene-d8	10.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	95
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91





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Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

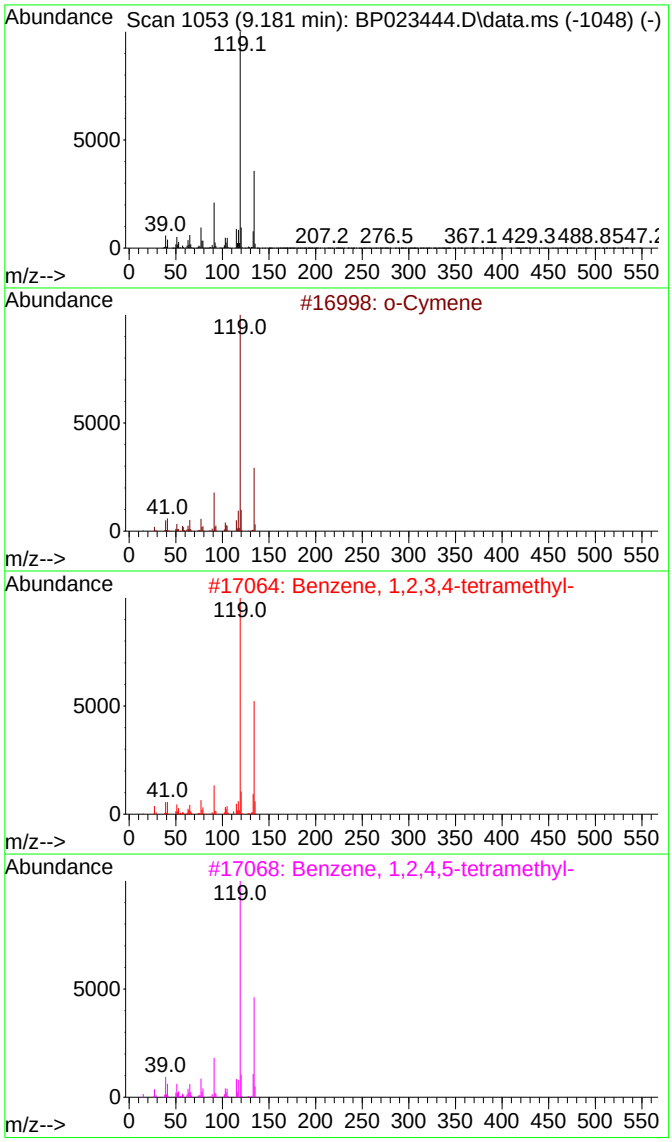
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

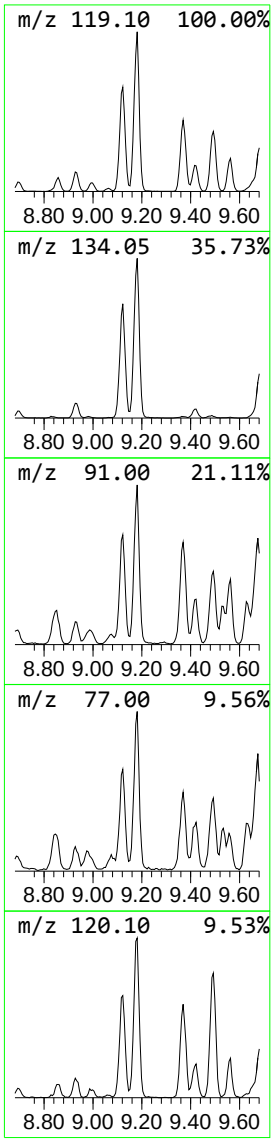
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.181	14.39 ng/ul	1207490	Naphthalene-d8	10.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	95
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
4	p-Cymene	134	C10H14	000099-87-6	91
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91





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Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

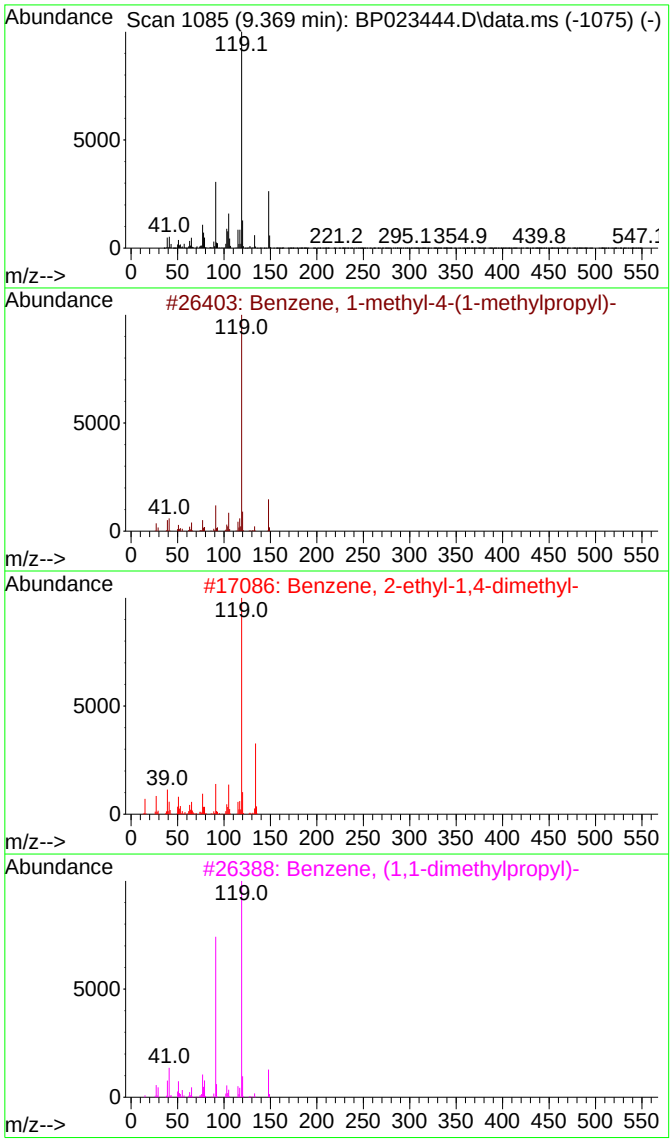
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

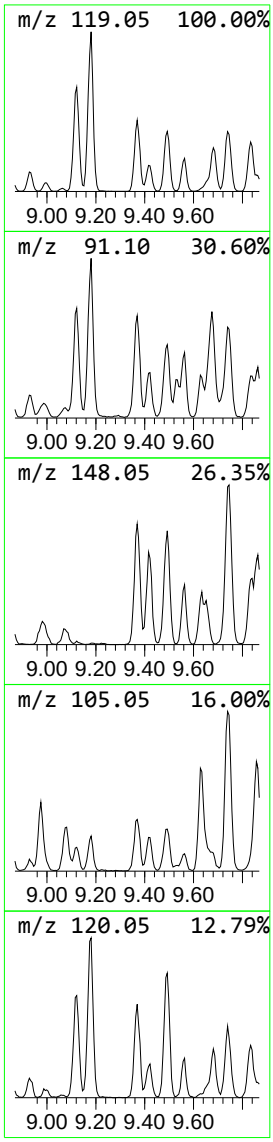
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Benzene, 1-methyl-4-(1-meth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.369	8.72 ng/ul	731547	Naphthalene-d8	10.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	90
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	81
3	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	80
4	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	80
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	74





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Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
C0AD3

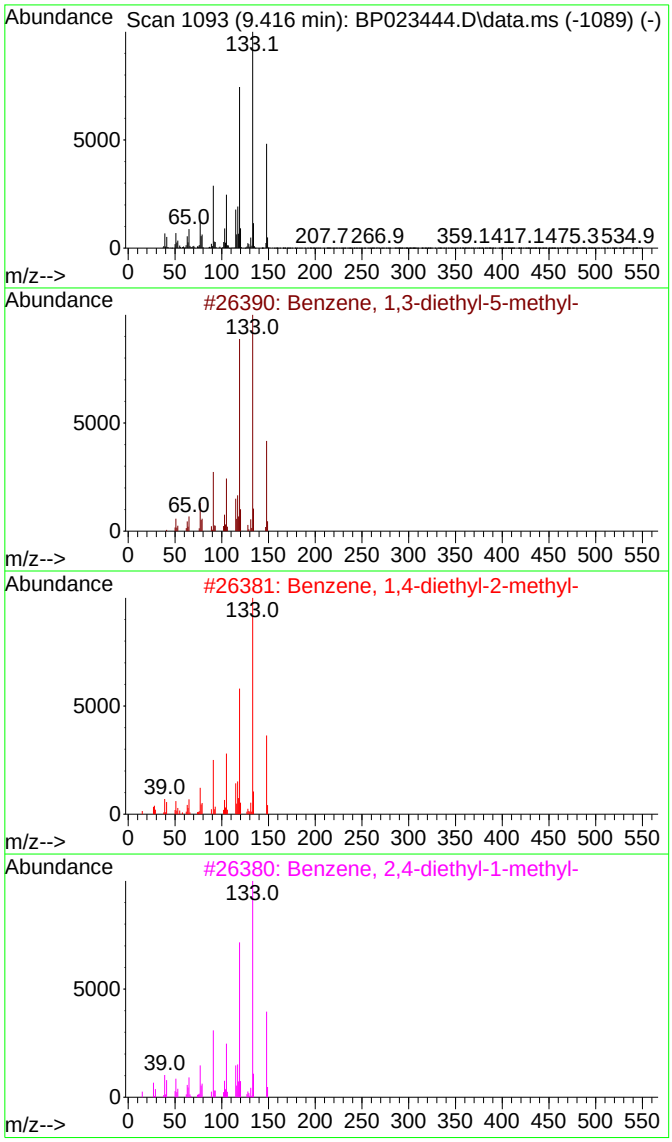
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

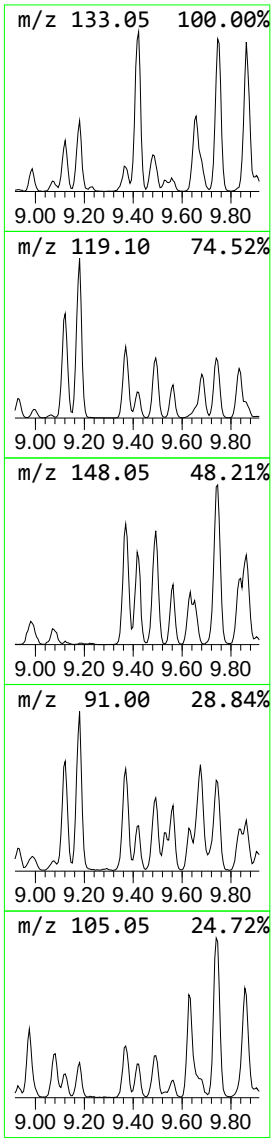
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.416	4.38 ng/ul	367291	Naphthalene-d8	10.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	98
2	Benzene, 1,4-diethyl-2-methyl-	148	C11H16	013632-94-5	97
3	Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	95
4	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	90
5	Benzene, pentamethyl-	148	C11H16	000700-12-9	64





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Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

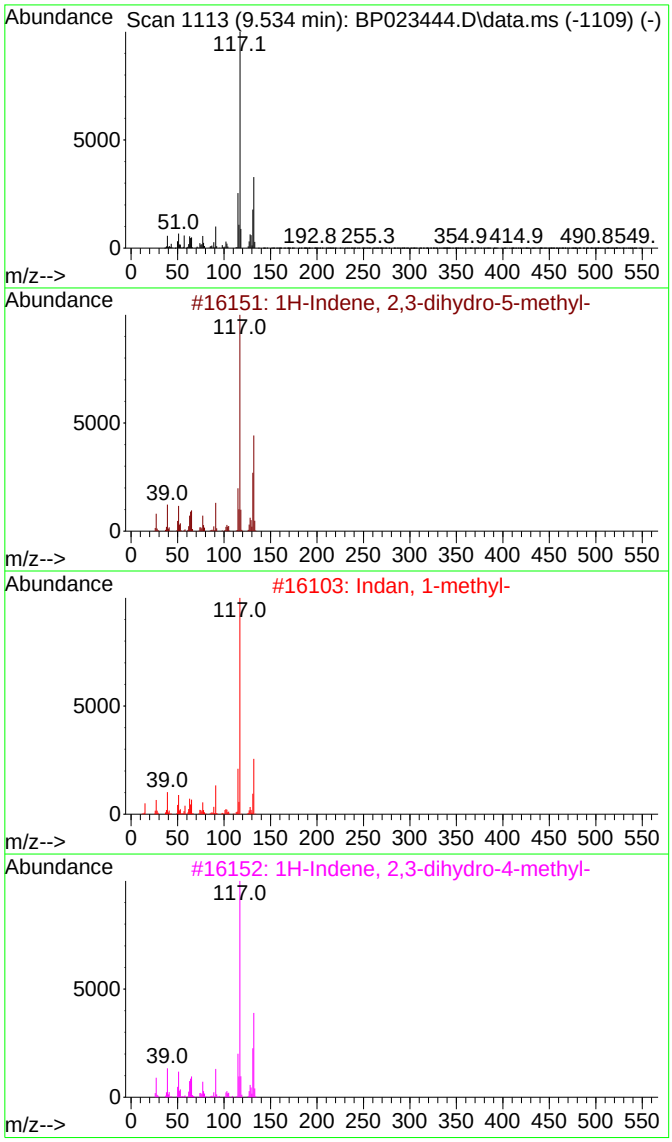
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

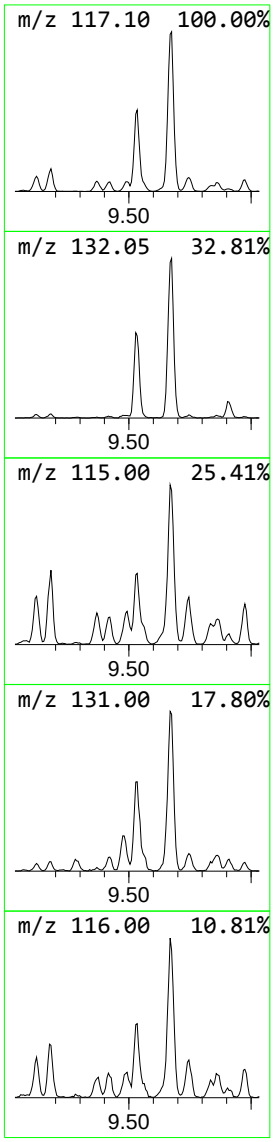
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.534	3.15 ng/ul	264133	Naphthalene-d8	10.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
2	Indan, 1-methyl-	132	C10H12	000767-58-8	90
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	83





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Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

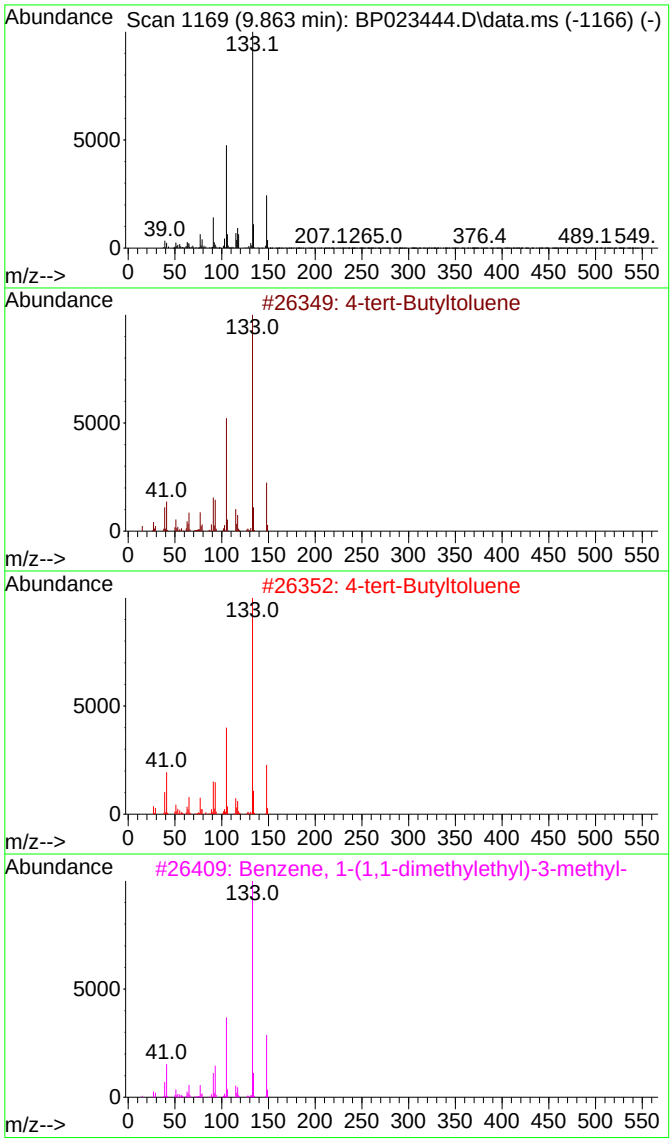
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

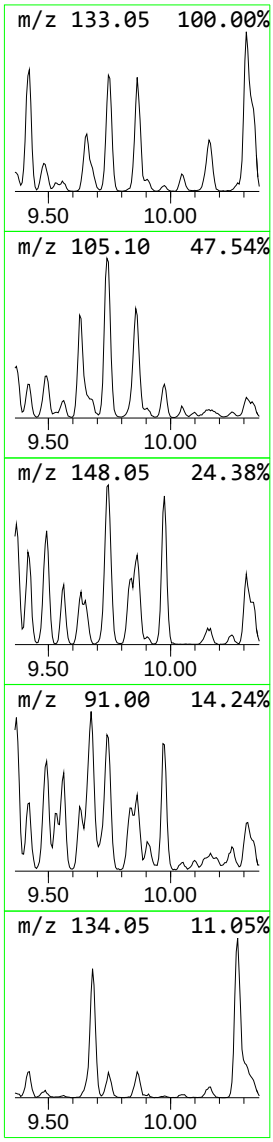
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 4-tert-Butyltoluene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.863	5.38 ng/ul	450930	Naphthalene-d8	10.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-tert-Butyltoluene	148	C11H16	000098-51-1	91
2	4-tert-Butyltoluene	148	C11H16	000098-51-1	91
3	Benzene, 1-(1,1-dimethylethyl)-3...	148	C11H16	001075-38-3	91
4	4-tert-Butyltoluene	148	C11H16	000098-51-1	91
5	2-tert-Butyltoluene	148	C11H16	001074-92-6	91





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Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

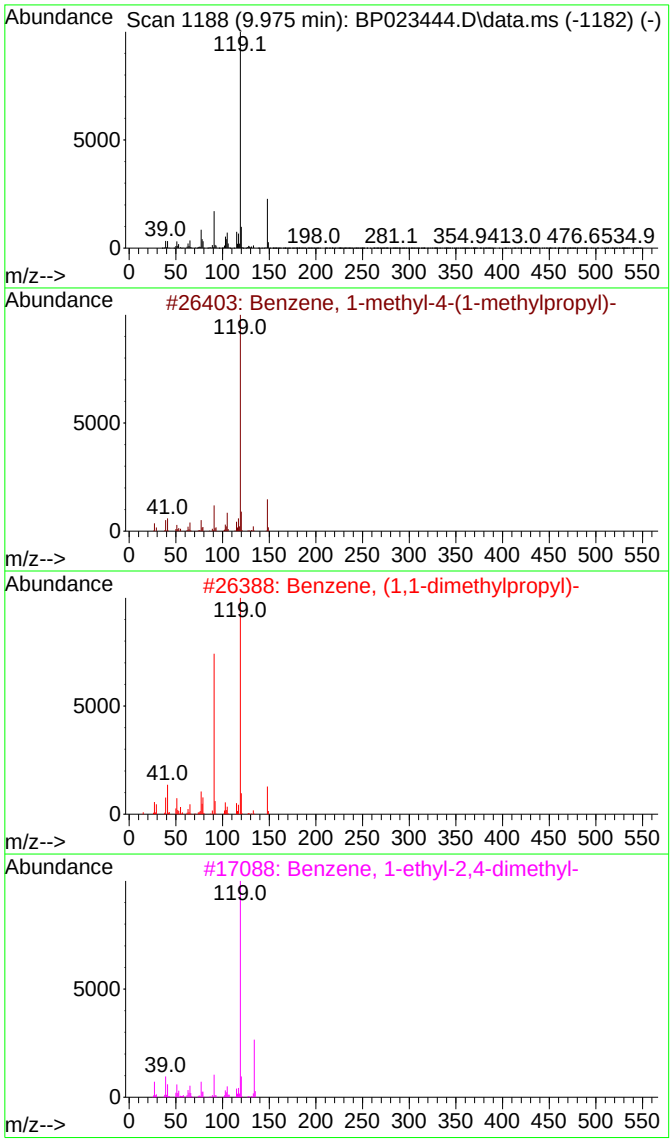
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

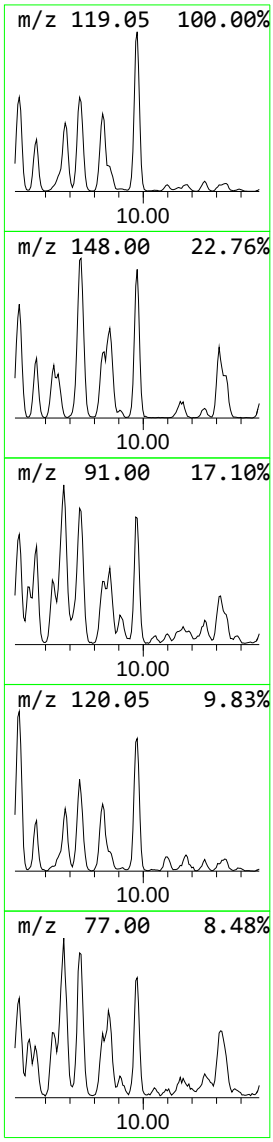
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Benzene, (1,1-dimethylpropyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.975	8.12 ng/ul	681216	Naphthalene-d8	10.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	91
2	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	86
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	81
4	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	78
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	72





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Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

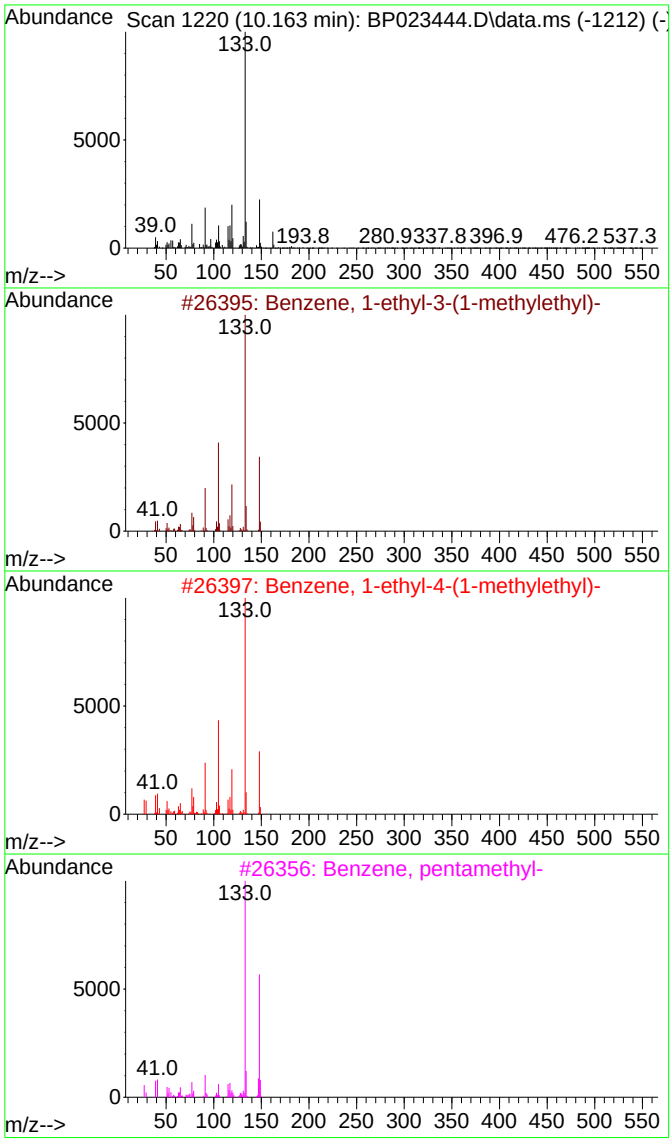
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

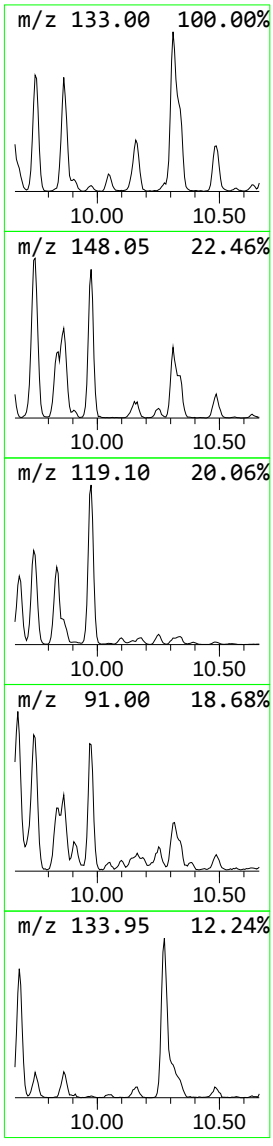
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Benzene, 1-ethyl-3-(1-methy... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.163	2.83 ng/ul	237404	Naphthalene-d8	10.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	91
2	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	90
3	Benzene, pentamethyl-	148	C11H16	000700-12-9	87
4	Benzene, pentamethyl-	148	C11H16	000700-12-9	87
5	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	87





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Data File : BP023444.D
Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

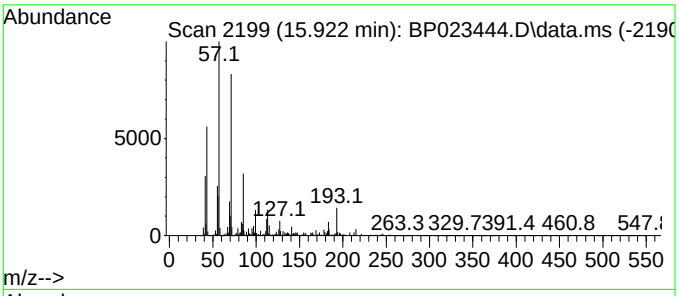
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

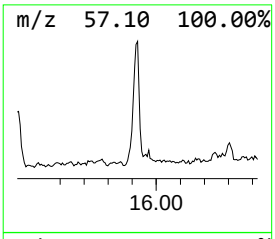
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

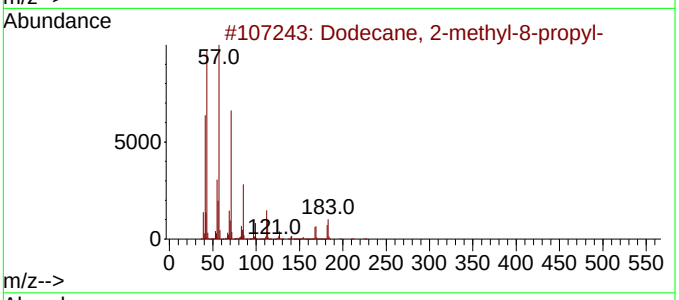
****
Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 18

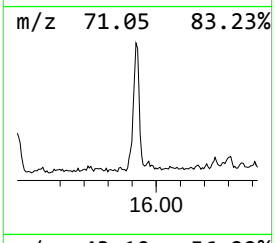
R.T.	EstConc	Area	Relative to ISTD	R.T.
15.922	4.71 ng/ul	533163	Phenanthrene-d10	16.957

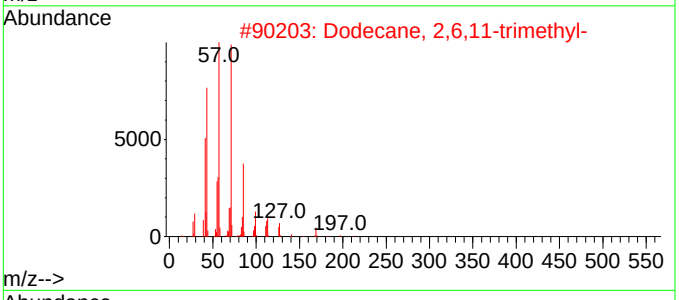
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2-methyl-8-propyl-	226	C16H34	055045-07-3	93
2	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	91
3	Heptadecane, 7-methyl-	254	C18H38	020959-33-5	87
4	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	83
5	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	80

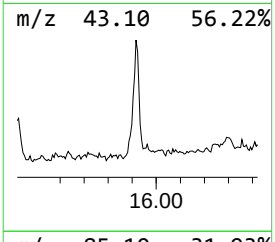


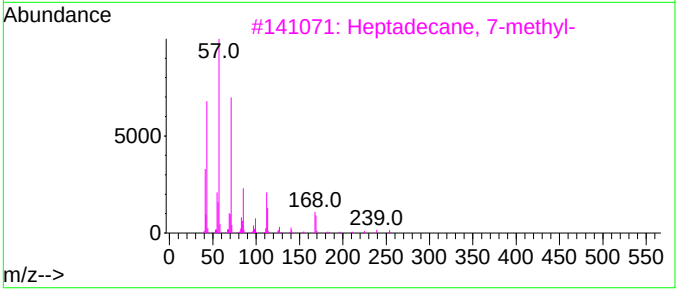


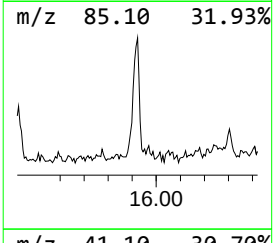


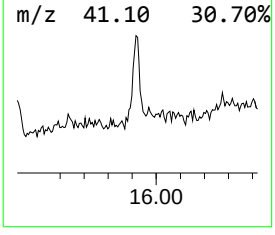












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Data File : BP023444.D
Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

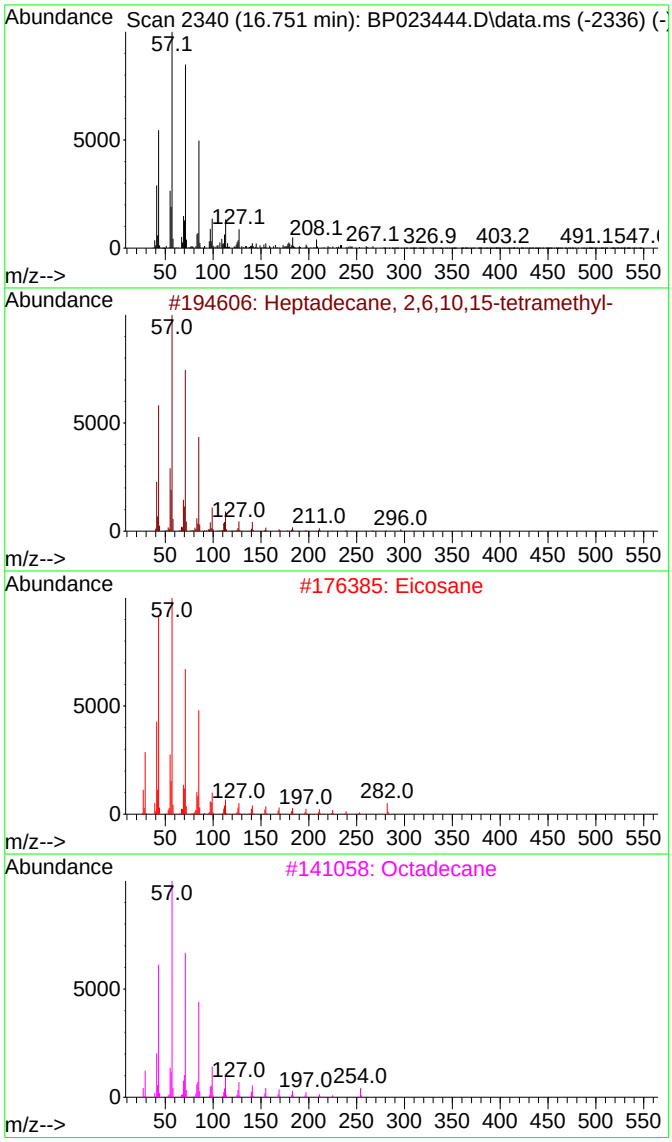
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

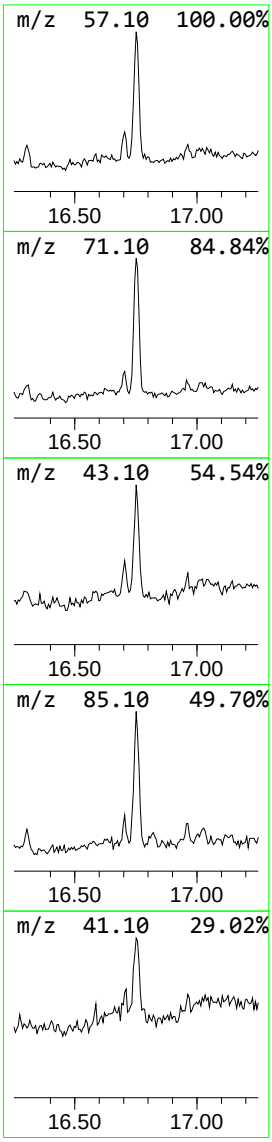
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.751	2.16 ng/ul	244177	Phenanthrene-d10	16.957

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
2	Eicosane	282	C20H42	000112-95-8	86
3	Octadecane	254	C18H38	000593-45-3	80
4	Hexadecane	226	C16H34	000544-76-3	80
5	Heneicosane	296	C21H44	000629-94-7	80





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Data File : BP023444.D
Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

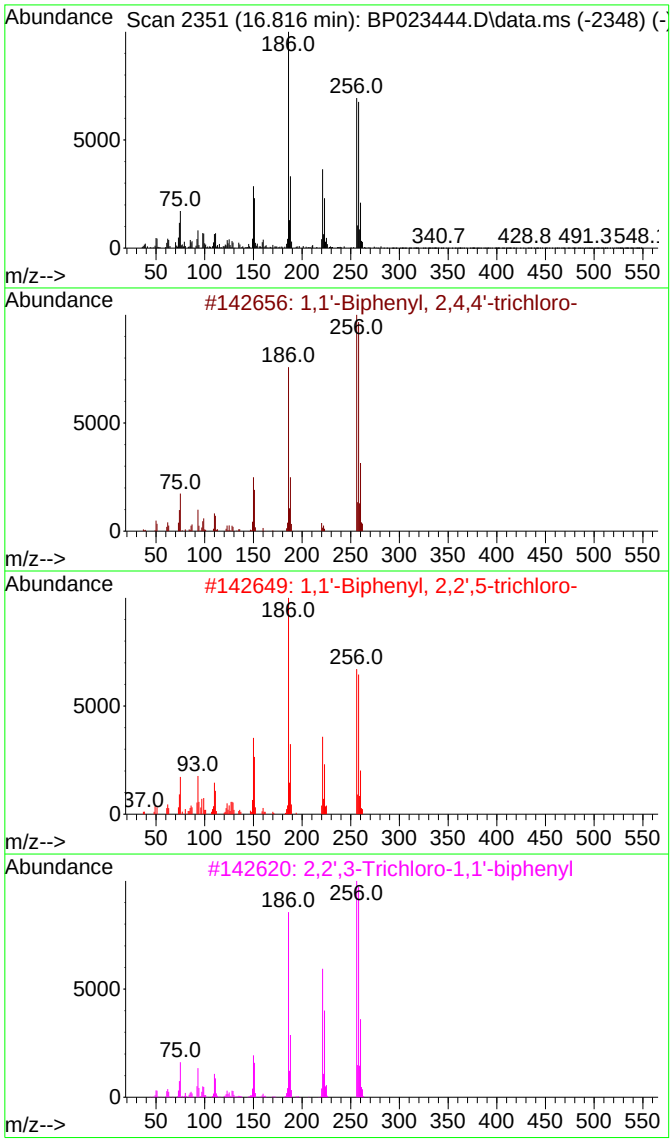
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

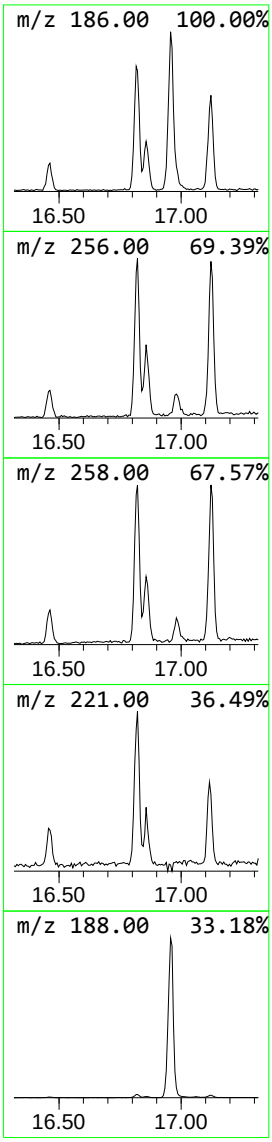
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 1,1'-Biphenyl, 2,4,4'-trich... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.816	2.28 ng/ul	258139	Phenanthrene-d10	16.957

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99
2	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99
3	2,2',3-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-78-9	99
4	2,3',6-Trichlorobiphenyl	256	C12H7Cl3	038444-76-7	97
5	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	96





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Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

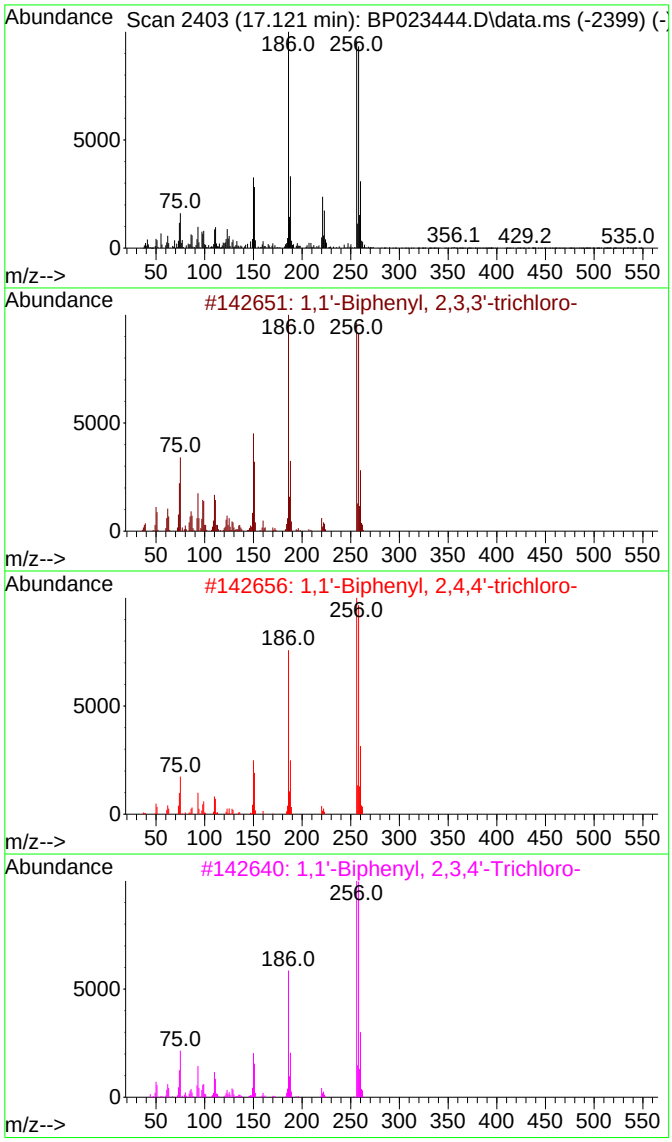
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

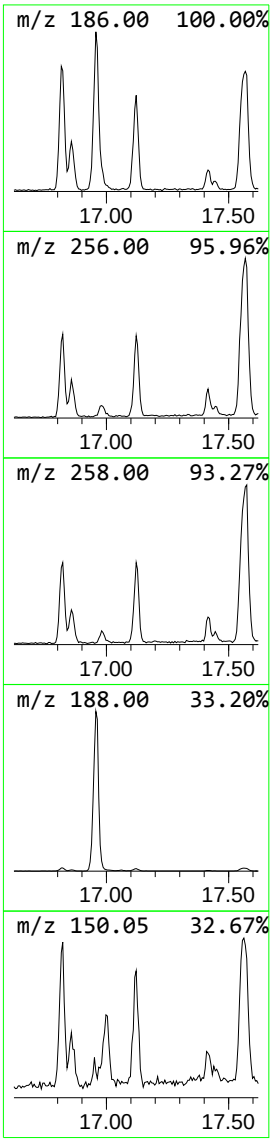
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 1,1'-Biphenyl, 2,3,3'-trich... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.122	3.22 ng/ul	364763	Phenanthrene-d10	16.957

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99
2	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99
3	1,1'-Biphenyl, 2,3,4'-Trichloro-	256	C12H7Cl3	038444-85-8	97
4	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	97
5	1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	97





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Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

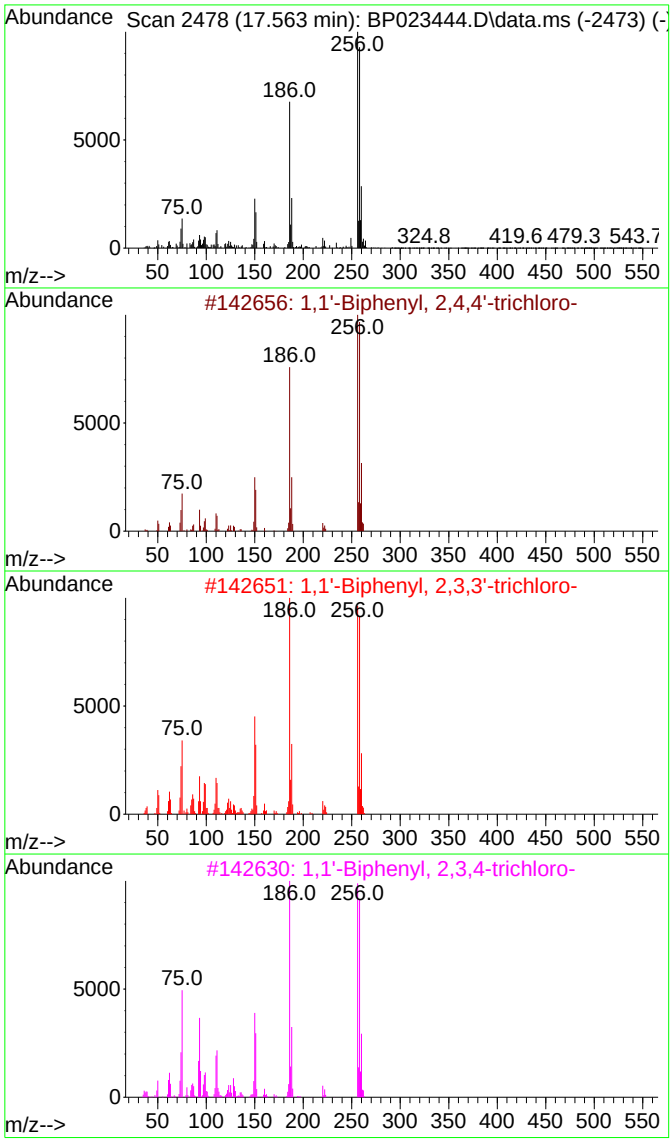
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

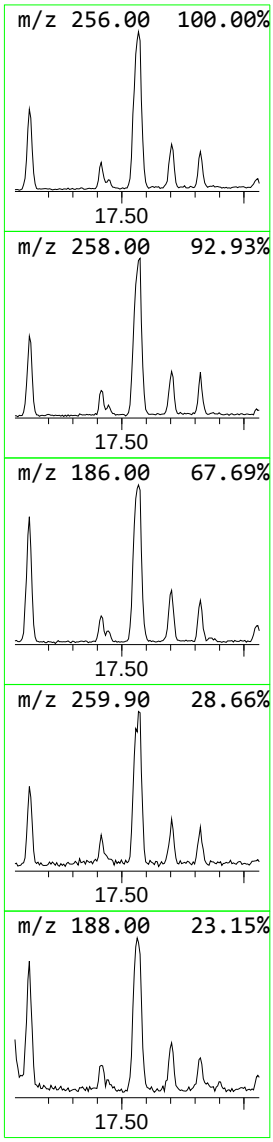
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 30 1,1'-Biphenyl, 2,3,4-trichl... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.563	5.57 ng/ul	629824	Phenanthrene-d10	16.957

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99
2	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99
3	1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	99
4	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	98
5	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	98





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Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

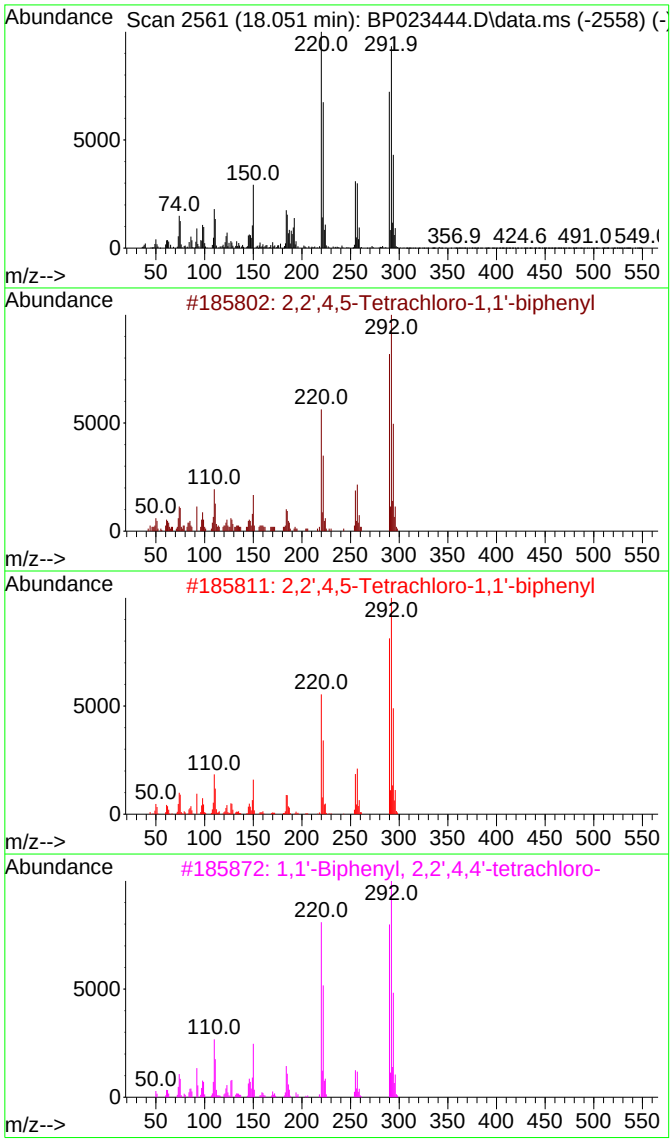
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

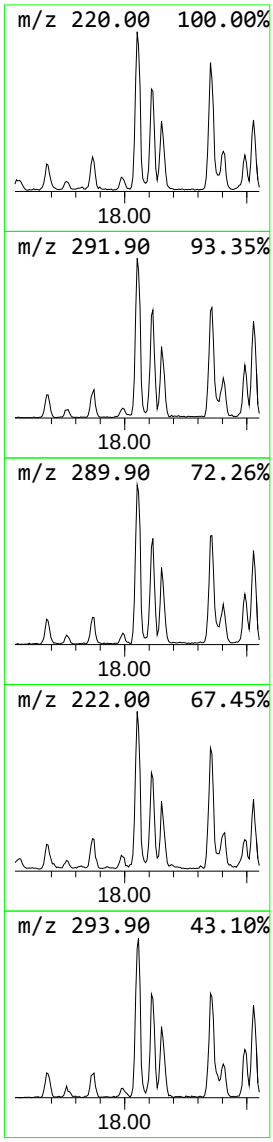
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 31 2,2',4,5-Tetrachloro-1,1'-b... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.051	4.09 ng/ul	462256	Phenanthrene-d10	16.957

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
2	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
3	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
4	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
5	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99





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Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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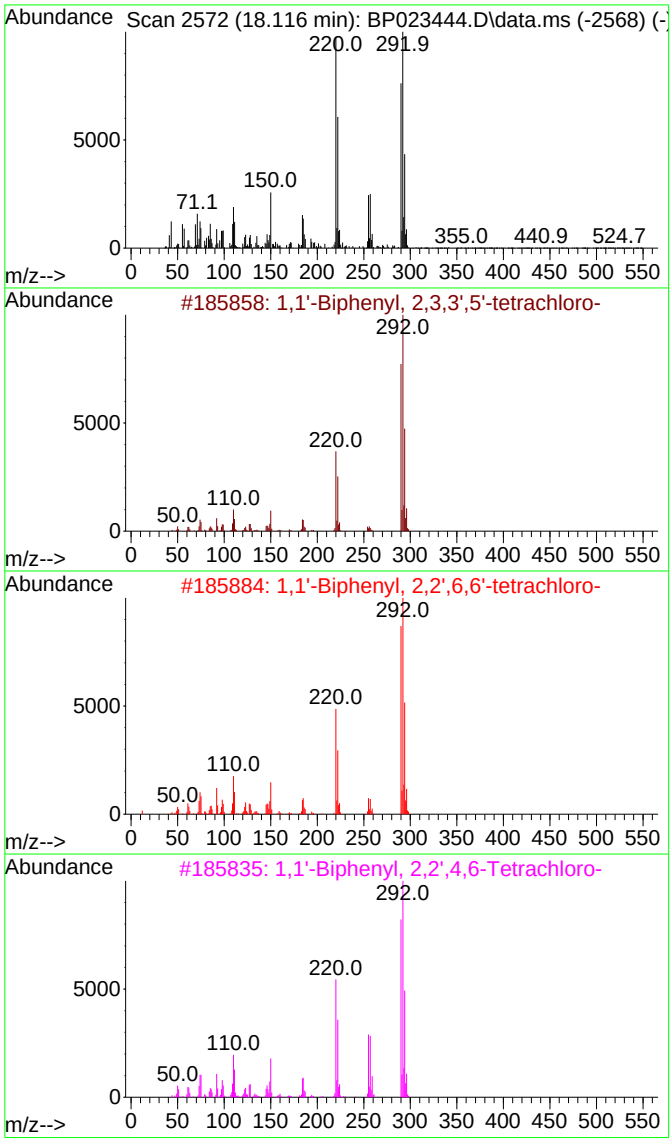
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

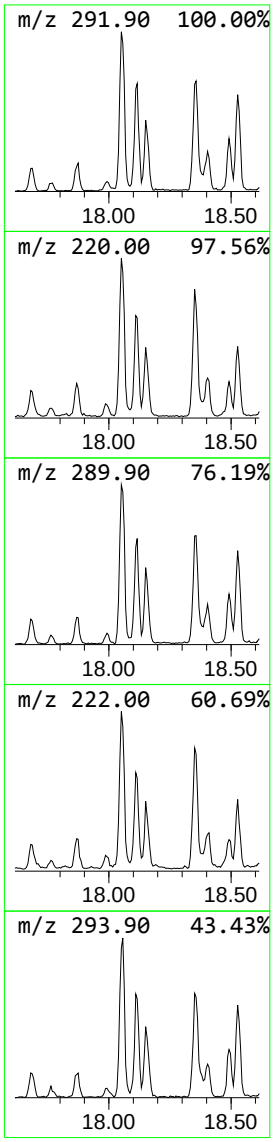
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 32 1,1'-Biphenyl, 2,3,3',5'-te... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.116	3.19 ng/ul	360500	Phenanthrene-d10	16.957

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99
2	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
3	1,1'-Biphenyl, 2,2',4,6-Tetrachl...	290	C12H6Cl4	062796-65-0	99
4	1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	99
5	1,1'-Biphenyl, 3,3',4,4'-tetrach...	290	C12H6Cl4	032598-13-3	98





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Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

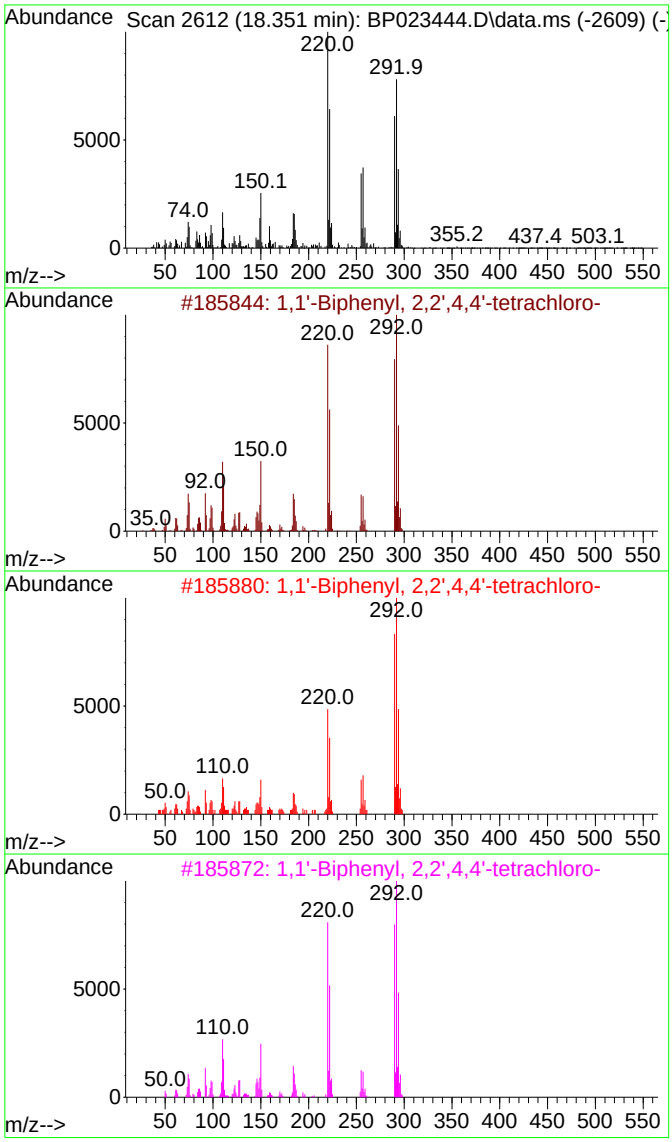
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

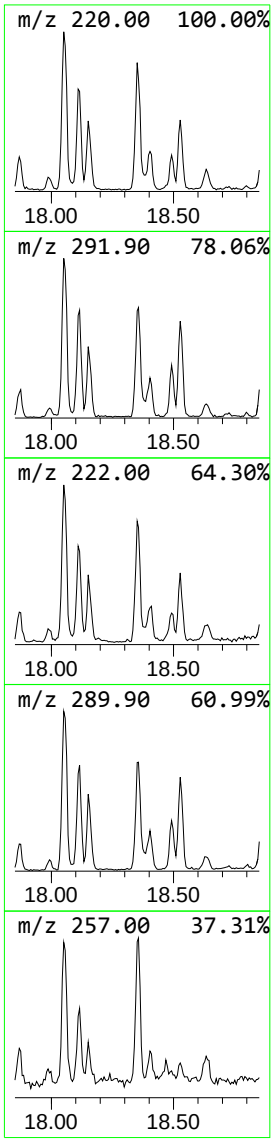
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 33 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.351	3.33 ng/ul	377088	Phenanthrene-d10	16.957

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
2	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
3	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
4	1,1'-Biphenyl, 2,2',5,6'-Tetrach...	290	C12H6Cl4	041464-41-9	99
5	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99





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Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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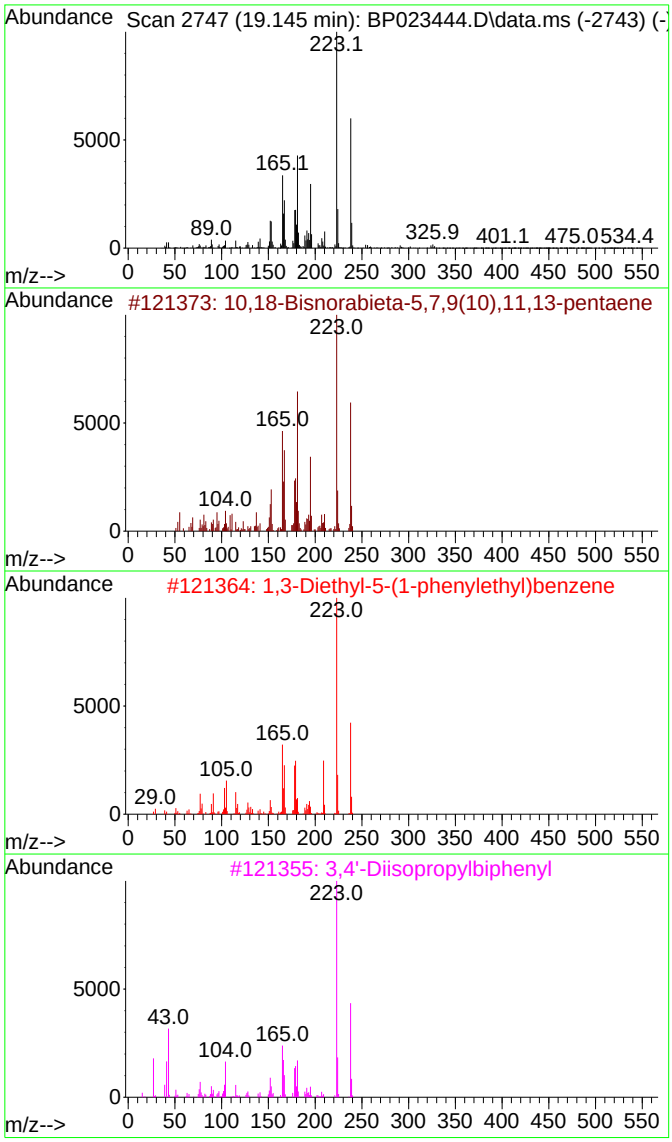
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

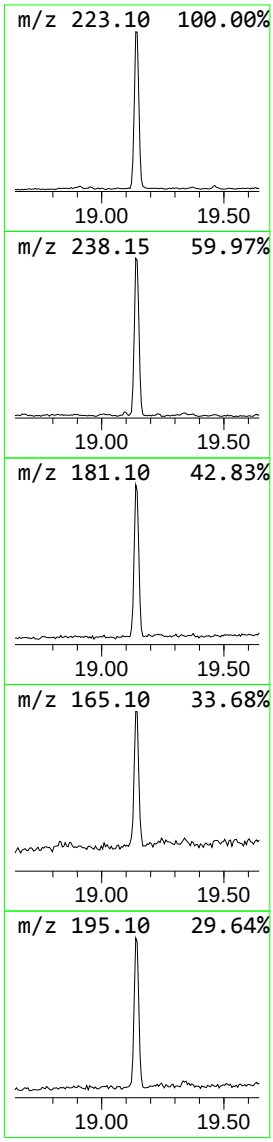
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 34 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.145	7.66 ng/ul	866235	Phenanthrene-d10	16.957

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2	1,3-Diethyl-5-(1-phenylethyl)ben...	238	C18H22	1000482-32-6	76
3	3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	70
4	Anthracene, 1,4-dimethoxy-	238	C16H14O2	013076-29-4	58
5	4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	53





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023444.D
Acq On : 16 Dec 2024 21:51
Operator : RC/JU
Sample : P5234-01
Misc : GCMS CONFIRMATION
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: S...	3.016	50.8	ng/ul	2802460	1	7.534	1103270	20.0
Benzene, 1,3-di...	8.151	5.8	ng/ul	321109	1	7.534	1103270	20.0
Benzene, 2-ethy...	8.457	5.0	ng/ul	273781	1	7.534	1103270	20.0
p-Cymene	8.510	6.5	ng/ul	357109	1	7.534	1103270	20.0
Benzene, 1-ethy...	8.604	21.1	ng/ul	1165630	1	7.534	1103270	20.0
Benzene, (2-met...	8.687	3.0	ng/ul	168050	1	7.534	1103270	20.0
(DEL) Alkane: S...	8.722	3.0	ng/ul	164320	1	7.534	1103270	20.0
Benzene, 4-ethy...	8.928	2.4	ng/ul	202191	2	10.275	1677810	20.0
o-Cymene	9.122	9.5	ng/ul	799111	2	10.275	1677810	20.0
Benzene, 1,2,3,...	9.181	14.4	ng/ul	1207490	2	10.275	1677810	20.0
Benzene, 1-meth...	9.369	8.7	ng/ul	731547	2	10.275	1677810	20.0
Benzene, 1,3-di...	9.416	4.4	ng/ul	367291	2	10.275	1677810	20.0
1H-Indene, 2,3-...	9.534	3.1	ng/ul	264133	2	10.275	1677810	20.0
4-tert-Butyltol...	9.863	5.4	ng/ul	450930	2	10.275	1677810	20.0
Benzene, (1,1-d...	9.975	8.1	ng/ul	681216	2	10.275	1677810	20.0
Benzene, 1-ethy...	10.163	2.8	ng/ul	237404	2	10.275	1677810	20.0
(DEL) Alkane: S...	15.922	4.7	ng/ul	533163	4	16.957	2263060	20.0
(DEL) Alkane: S...	16.751	2.2	ng/ul	244177	4	16.957	2263060	20.0
1,1'-Biphenyl, ...	16.816	2.3	ng/ul	258139	4	16.957	2263060	20.0
1,1'-Biphenyl, ...	17.122	3.2	ng/ul	364763	4	16.957	2263060	20.0
1,1'-Biphenyl, ...	17.563	5.6	ng/ul	629824	4	16.957	2263060	20.0
2,2',4,5-Tetrac...	18.051	4.1	ng/ul	462256	4	16.957	2263060	20.0
1,1'-Biphenyl, ...	18.116	3.2	ng/ul	360500	4	16.957	2263060	20.0
1,1'-Biphenyl, ...	18.351	3.3	ng/ul	377088	4	16.957	2263060	20.0
10,18-Bisnorabi...	19.145	7.7	ng/ul	866235	4	16.957	2263060	20.0