LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : EY0C3

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-BP062322.M
Title : SVOA CALIBRATION

Signal : TIC: BP010831.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.617	87	90	99	rBV	577296	780974	5.72%	0.504%
2	3.693	99	103	114	rVB2	94014	181679	1.33%	0.117%
3	3.787	114	119	124	rBV	89522	126360	0.93%	0.082%
4	3.840	124	128	134	rVB2	56033	88679	0.65%	0.057%
5	3.934	140	144	148	rBV	158157	200690	1.47%	0.129%
6	4.858	294	301	305	rBV2	162065	260705	1.91%	0.168%
7	5.016	322	328	335	rVB2	34473	70667	0.52%	0.046%
8	5.240	358	366	373	rVB	334157	501788	3.68%	0.324%
9	5.393	383	392	397	rBV	138742	248488	1.82%	0.160%
10	5.734	443	450	455	rVV	102222	171192	1.25%	0.110%
11	6.211	524	531	540	rBV	444109	671570	4.92%	0.433%
12	6.699	608	614	624	rVB	1301474	1920232	14.07%	1.239%
13	6.810	624	633	638	rBV	553410	808695	5.93%	0.522%
14	6.869	638	643	646	rVV	941874	1437725	10.54%	0.928%
15	6.899	646	648	652	rVV	441256	557063	4.08%	0.359%
16	6.946	652	656	665	rVB	418539	619979	4.54%	0.400%
17	7.063	668	676	679	rBV	1548312	2397625	17.57%	1.547%
18	7.110	679	684	698	rVB	2718834	4116056	30.16%	2.655%
19	7.252	700	708	716	rBV	1523365	2308527	16.92%	1.489%
20	7.369	720	728	736	rVB	9257719	13645808	100.00%	8.803%
21	7.587	760	765	768	rVV	111536	176447	1.29%	0.114%
22	7.622	768	771	780	rVV	137602	213899	1.57%	0.138%
23	7.716	780	787	792	rVV	1497339	2331763	17.09%	1.504%
24	7.757	792	794	799	rVV	141294	193224	1.42%	0.125%
25	7.828	800	806	815	rVV	1896406	3078128	22.56%	1.986%
26	8.022	824	839	843	rBV5	30995	95278	0.70%	0.061%
27	8.087	843	850	860	rVB	1030388	1577321	11.56%	1.018%
28	8.210	865	871	876	rVV	722797	1050984	7.70%	0.678%
29	8.269	876	881	886	rVV	323253	499664	3.66%	0.322%
30	8.357	889	896	902	rVV2	1150090	1843090	13.51%	1.189%
31	8.434	902	909	918	rVV	1248491	2145563	15.72%	1.384%
32	8.522	918	924	933	rVB	504628	760651	5.57%	0.491%
33	8.669	943	949	954	rBV	1088706	1696853	12.43%	1.095%
34	8.722	954	958	965	rVB	1058258	1565912	11.48%	1.010%
35	8.822	965	975	980	rBV	2178441	3316015	24.30%	2.139%
36	8.881	980	985	999	rVV2	1849325	4210923	30.86%	2.717%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

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-BP062322.M
Title : SVOA CALIBRATION

37	9.069	1010	1017	1025	rVB4	97726	233624	1.71%	0.151%
38	9.152	1025	1031	1038	rBV	394082	650435	4.77%	0.420%
39	9.216	1038	1042	1050	rVB3	42335	71936	0.53%	0.046%
40	9.346	1050	1064	1069	rBV	1203246	1888484	13.84%	1.218%
41	9.404	1069	1074	1081	rVB	1582516	2352453	17.24%	1.518%
42	9.599	1098	1107	1113	rBV2	1309219	2143559	15.71%	1.383%
43	9.657	1113	1117	1121	rVV	129577	196496	1.44%	0.127%
44	9.710	1121	1126	1130	rVV3	103373	194092	1.42%	0.125%
45	9.763	1130	1135	1147	rVV	900090	1534397	11.24%	0.990%
46	9.910	1147	1160	1169	rVV2	1818412	3716269	27.23%	2.398%
47	9.987	1169	1173	1182	rVV3	258072	604059	4.43%	0.390%
48	10.075	1182	1188	1190	rVV2	114033	234116	1.72%	0.151%
49	10.140	1190	1199	1207	rVV2	2002185	3520564	25.80%	2.271%
50	10.222	1207	1213	1220	rVB	189341	306307	2.24%	0.198%
51	10.404	1231	1244	1246	rBV3	86041	186195	1.36%	0.120%
52	10.434	1246	1249	1252	rVV	141186	227875	1.67%	0.147%
53	10.516	1252	1263	1267	rVV2	2030971	4285800	31.41%	2.765%
54	10.563	1267	1271	1278	rVV	1392837	2503142	18.34%	1.615%
55	10.657	1278	1287	1295	rVV	1809670	3526590	25.84%	2.275%
56	10.734	1295	1300	1310	rVB	180716	285058	2.09%	0.184%
57	10.887	1320	1326	1330	rBV2	51411	84175	0.62%	0.054%
58	10.981	1337	1342	1350	rVB2	84003	139910	1.03%	0.090%
59	11.146	1365	1370	1371	rBV2	75368	105095	0.77%	0.068%
60	11.181	1371	1376	1381	rVB	284718	449623	3.29%	0.290%
61	11.434	1413	1419	1425	rBV	372513	574669	4.21%	0.371%
62	11.628	1445	1452	1459	rBV	267021	451347	3.31%	0.291%
63	11.710	1461	1466	1475	rVB	187217	337649	2.47%	0.218%
64	11.846	1482	1489	1494	rBV	108584	176445	1.29%	0.114%
65	11.904	1494	1499	1510	rVB	198042	350792	2.57%	0.226%
66	12.028	1510	1520	1522	rBV4	47420	117186	0.86%	0.076%
67	12.075	1522	1528	1533	rVB4	109770	185488	1.36%	0.120%
68	12.251	1552	1558	1562	rBV	56636	91145	0.67%	0.059%
69	12.404	1572	1584	1590	rBV	1219081	1986089	14.55%	1.281%
70	12.498	1595	1600	1607	rVB3	72839	134879	0.99%	0.087%
71	12.716	1630	1637	1643	rBV	45849	83995	0.62%	0.054%
72	12.828	1651	1656	1662	rVB	105014	163484	1.20%	0.105%
73	13.245	1723	1727	1732	rVV	72163	125042	0.92%	0.081%
74	13.340	1738	1743	1750	rVV	228083	369822	2.71%	0.239%
75	13.428	1750	1758	1766	rVB	54203	122357	0.90%	0.079%
76	13.540	1766	1777	1784	rBV4	213101	641496	4.70%	0.414%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

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-BP062322.M
Title : SVOA CALIBRATION

77	13.616	1786	1790	1793	rVV	48855	74232	0.54%	0.048%
78	13.692	1797	1803	1808	rVV	319619	589561	4.32%	0.380%
79	13.745	1808	1812	1815	rVV2	213651	327835	2.40%	0.211%
80	13.792	1815	1820	1827	rVV	3409738	4849715	35.54%	3.129%
81	13.857	1827	1831	1839	rVV2	41306	94084	0.69%	0.061%
82	13.928	1839	1843	1847	rVV	100760	148813	1.09%	0.096%
83	14.063	1857	1866	1882	rBV	3868959	5923474	43.41%	3.821%
84	14.198	1884	1889	1896	rVB	57623	89050	0.65%	0.057%
85	14.375	1910	1919	1929	rBV	2621089	3899542	28.58%	2.516%
86	14.592	1951	1956	1965	rVB	296652	442407	3.24%	0.285%
87	14.698	1969	1974	1979	rBV	97719	132695	0.97%	0.086%
88	15.004	2019	2026	2031	rBV	55465	85590	0.63%	0.055%
89	15.375	2080	2089	2095	rBV	5241253	7264393	53.24%	4.687%
90	15.492	2102	2109	2117	rBV	943862	1314993	9.64%	0.848%
91	15.616	2126	2130	2134	rVB	56627	82257	0.60%	0.053%
92	17.139	2382	2389	2395	rBV	2724940	3878884	28.43%	2.502%
93	17.239	2399	2406	2419	rBV	5309982	7517322	55.09%	4.850%
94	19.063	2712	2716	2721	rBV2	67143	88155	0.65%	0.057%
95	19.127	2724	2727	2733	rVV	62147	81273	0.60%	0.052%
96	19.486	2782	2788	2796	rBV	6755796	8525176	62.47%	5.500%
97	20.980	3039	3042	3046	rBV	178127	194287	1.42%	0.125%
98	21.251	3083	3088	3093	rBV	3444549	4252708	31.16%	2.744%
99	23.468	3458	3465	3473	rVB2	5044197	9420904	69.04%	6.078%
100	23.621	3485	3491	3502	rVB	2257391	4505854	33.02%	2.907%

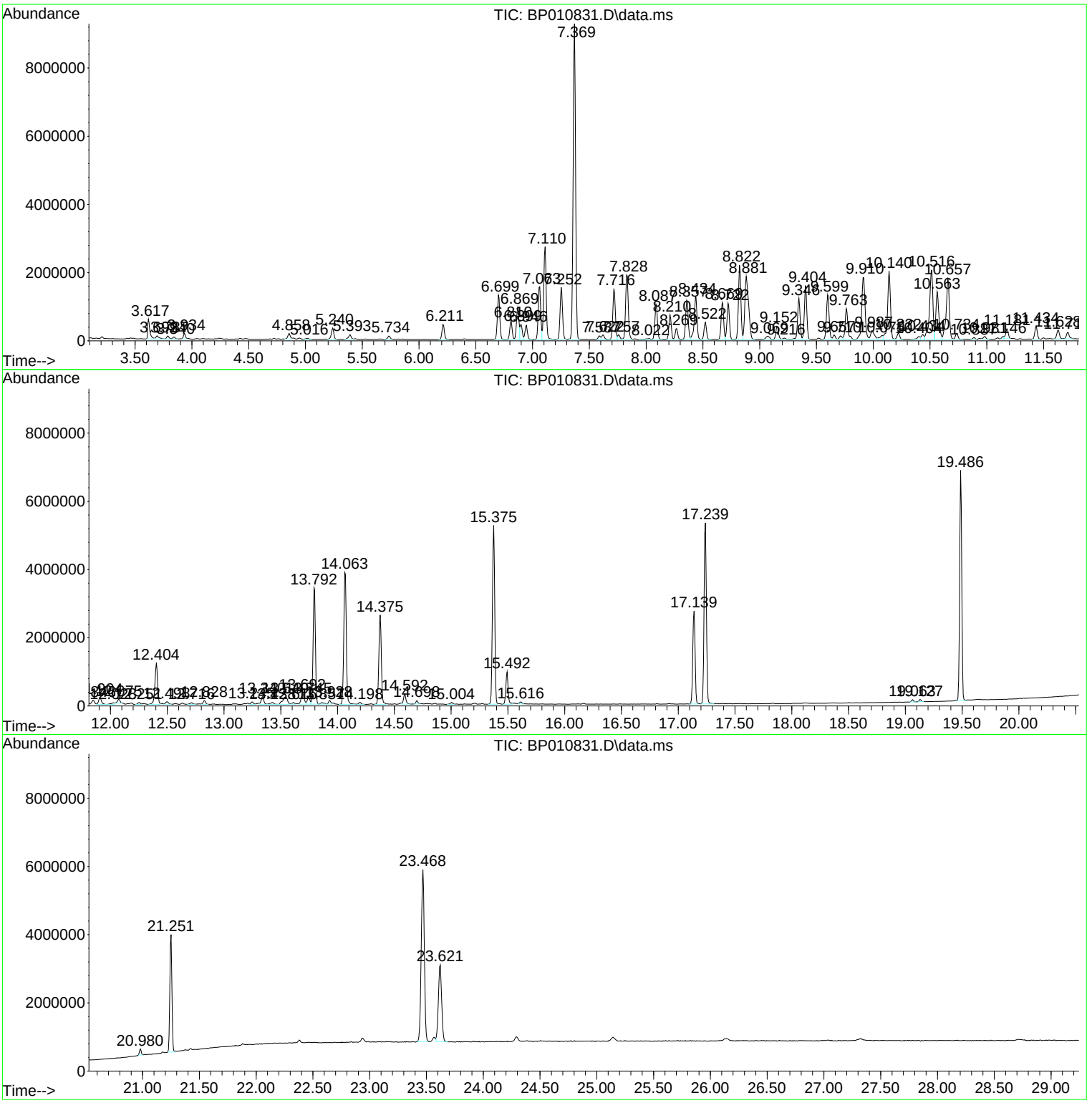
Sum of corrected areas: 155005560

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

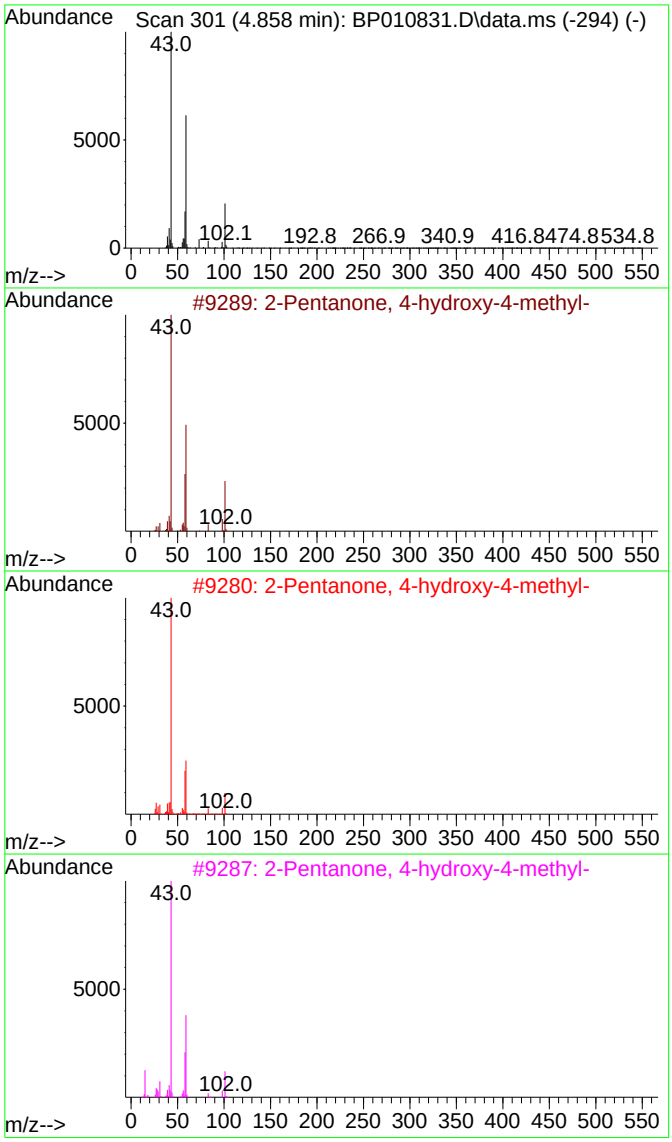
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

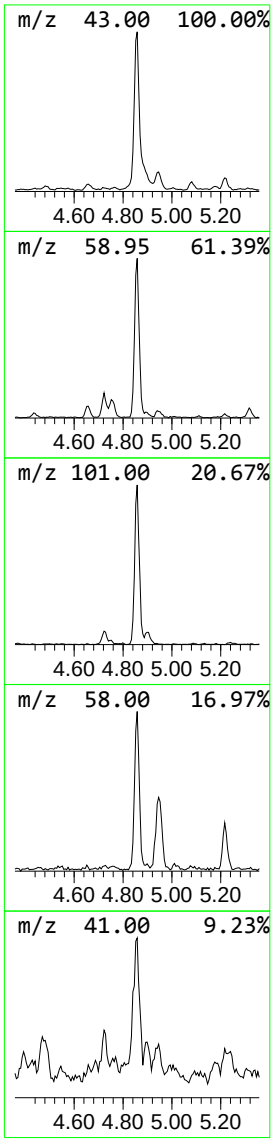
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.858	2.24 ng/ul	260705	1,4-Dichlorobenzene-d4	7.716

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
5	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

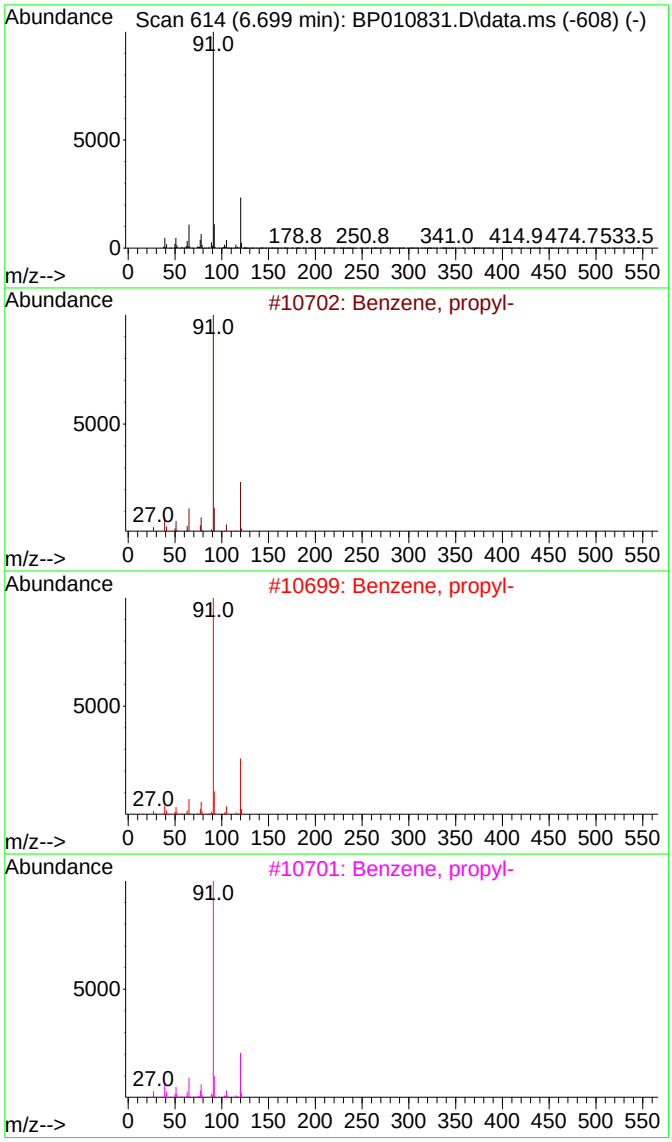
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

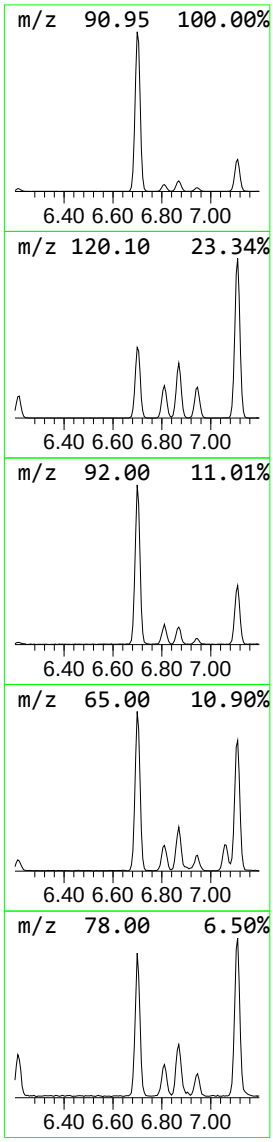
****

Peak Number 5 Benzene, propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.699	16.47 ng/ul	1920230	1,4-Dichlorobenzene-d4	7.716

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	94
2	Benzene, propyl-	120	C9H12	000103-65-1	90
3	Benzene, propyl-	120	C9H12	000103-65-1	90
4	Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	72
5	Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

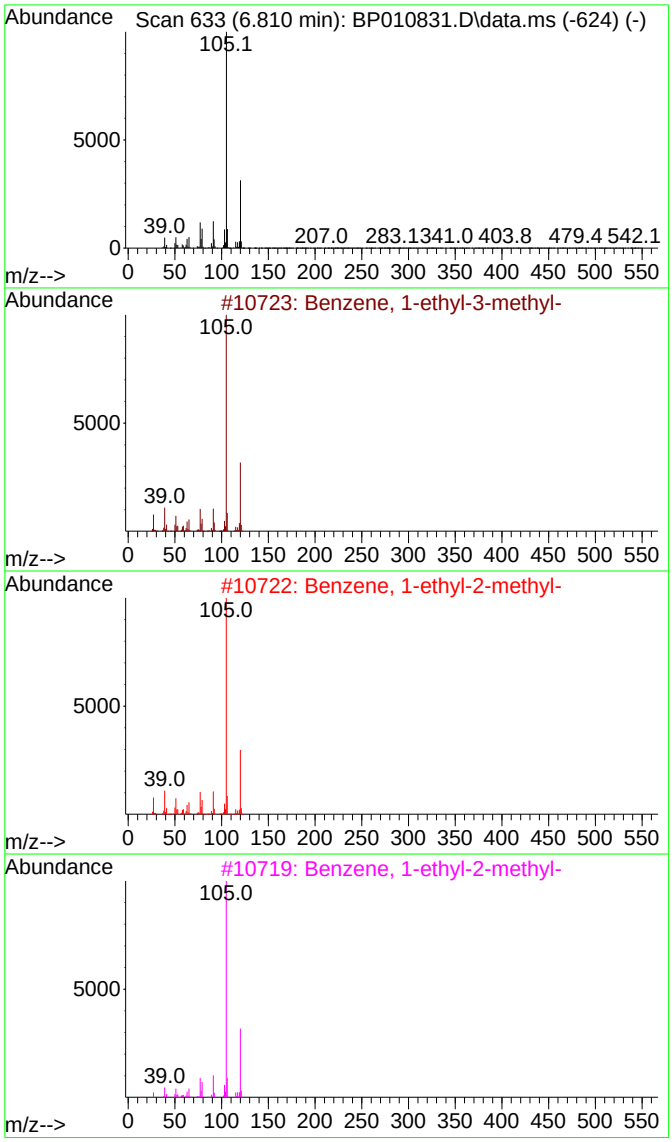
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

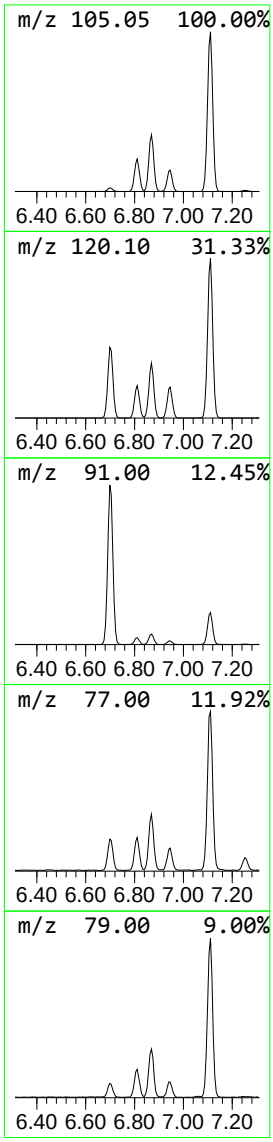
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1-ethyl-3-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.810	6.94 ng/ul	808695	1,4-Dichlorobenzene-d4	7.716

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

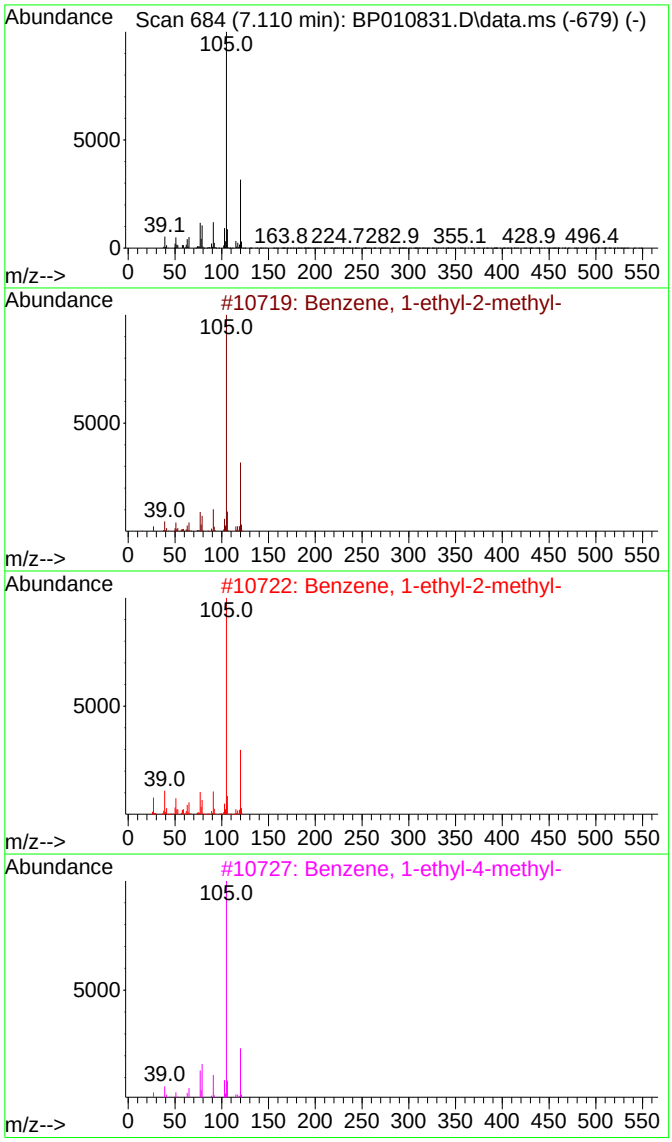
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

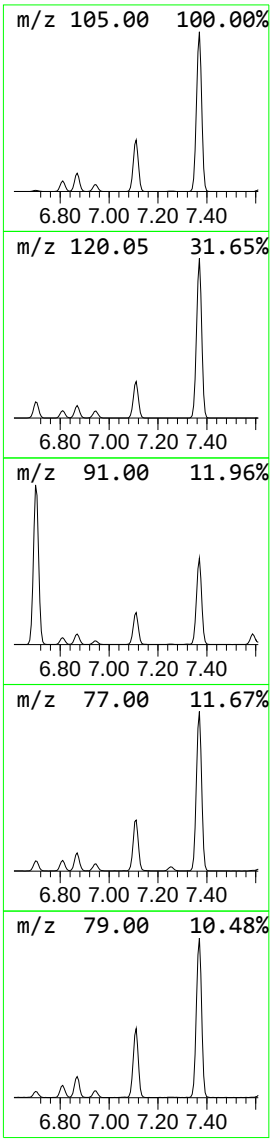
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.110	35.30 ng/ul	4116060	1,4-Dichlorobenzene-d4	7.716

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

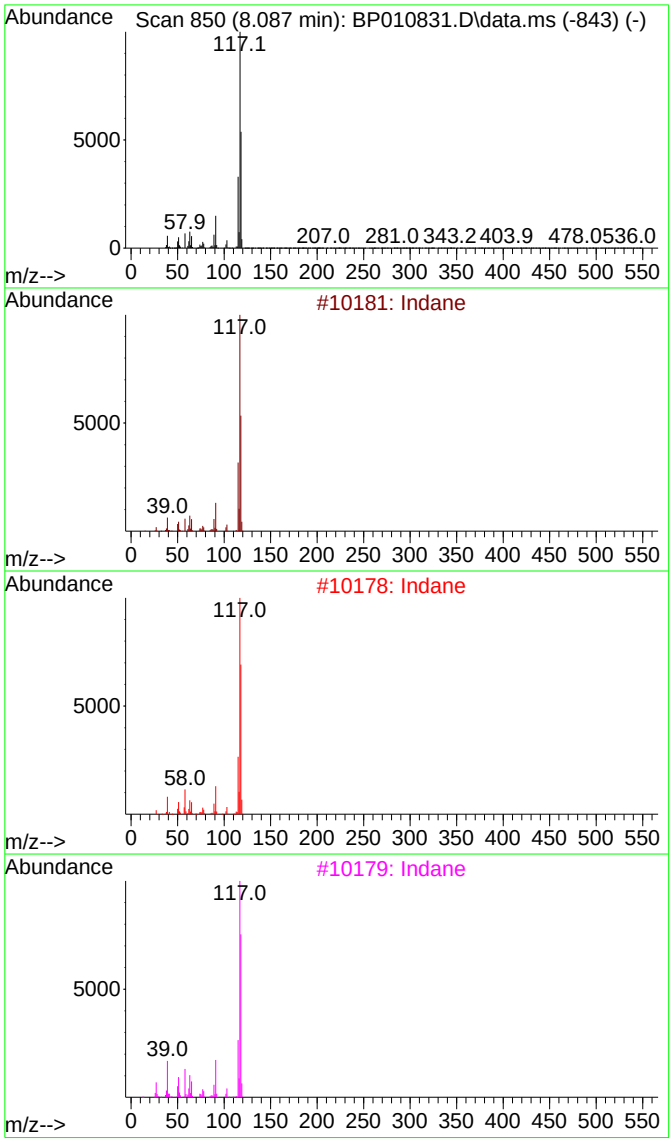
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

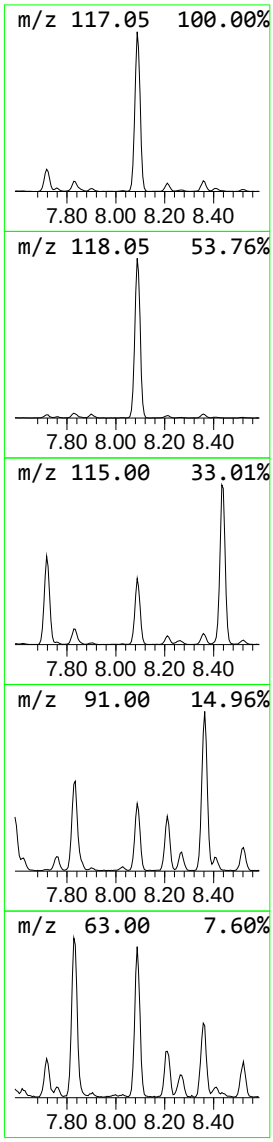
****

Peak Number 10 Indane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.087	13.53 ng/ul	1577320	1,4-Dichlorobenzene-d4	7.716

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	91
3	Indane	118	C9H10	000496-11-7	91
4	Deltacyclene	118	C9H10	007785-10-6	72
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64





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Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

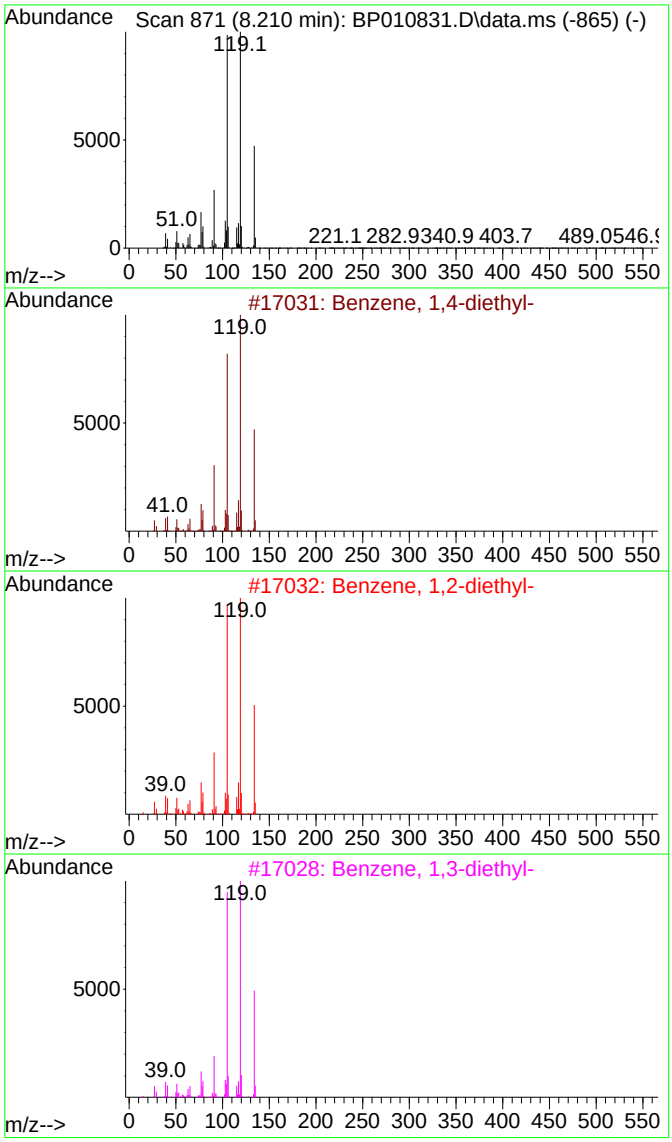
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

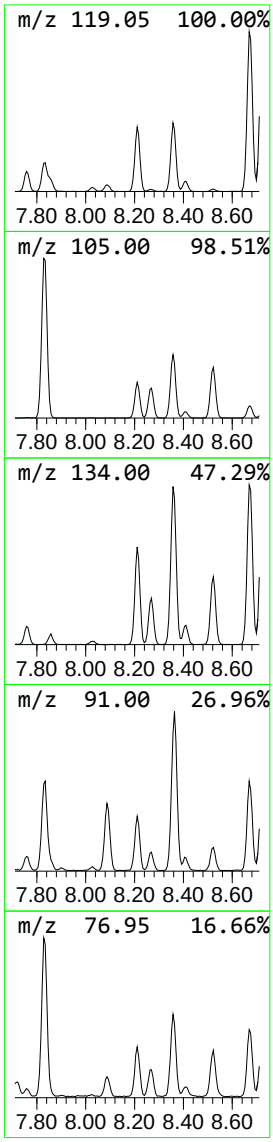
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 1,4-diethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.210	9.01 ng/ul	1050980	1,4-Dichlorobenzene-d4	7.716

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
4	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94





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Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

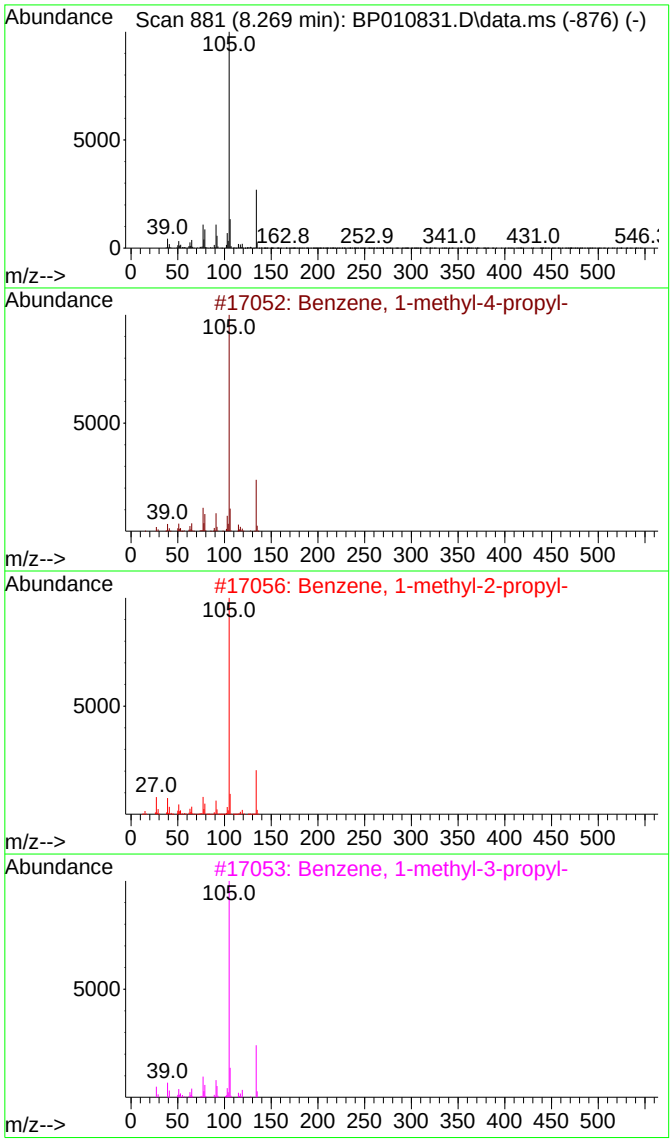
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

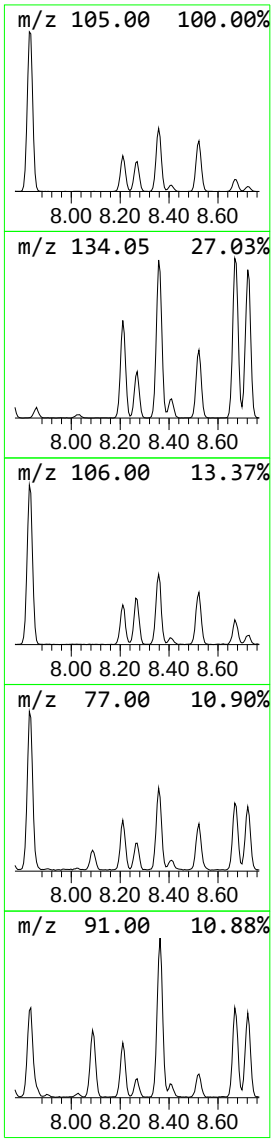
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzene, 1-methyl-4-propyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.269	4.29 ng/ul	499664	1,4-Dichlorobenzene-d4	7.716

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	94
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
3	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	91
4	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91





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Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

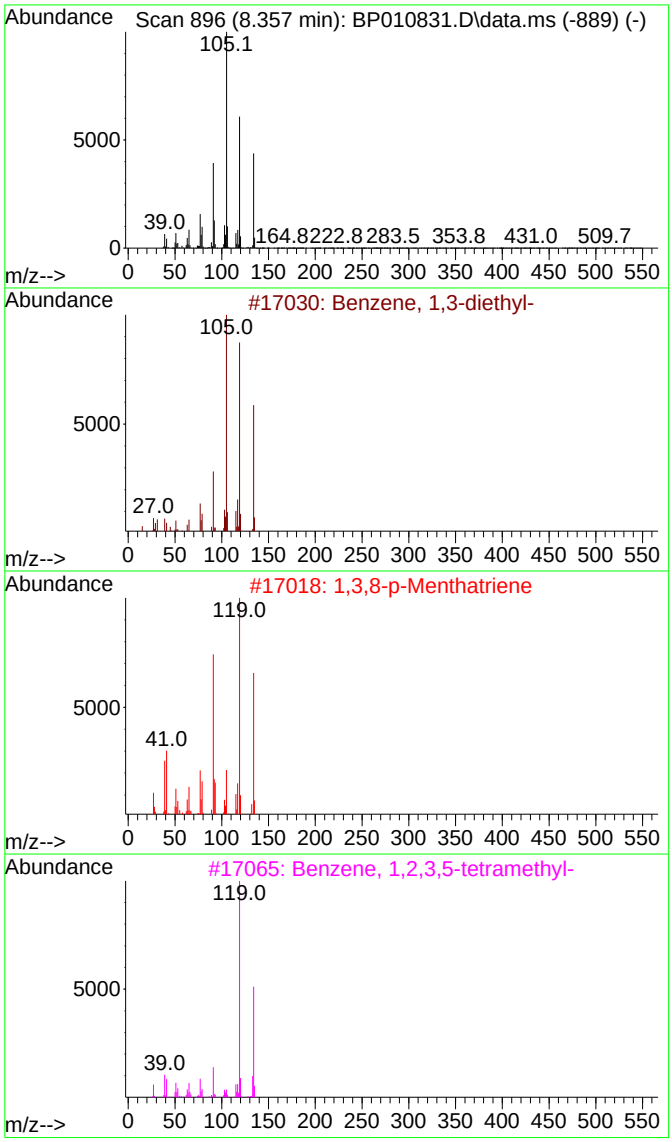
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

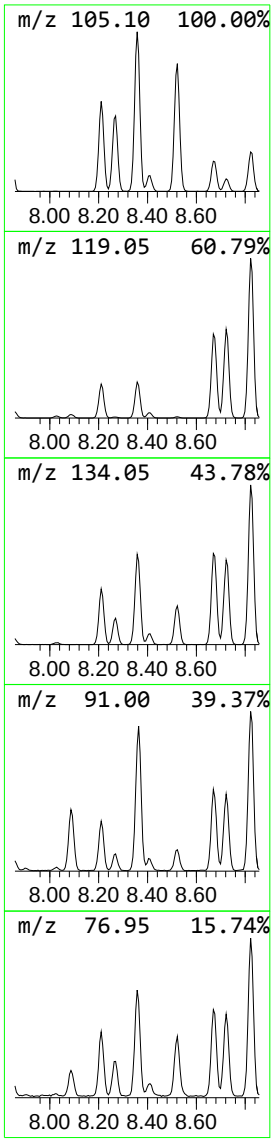
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 1,3-diethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.357	15.81 ng/ul	1843090	1,4-Dichlorobenzene-d4	7.716

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	93
2	1,3,8-p-Menthatriene	134	C10H14	018368-95-1	81
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	81
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	81
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	81





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Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

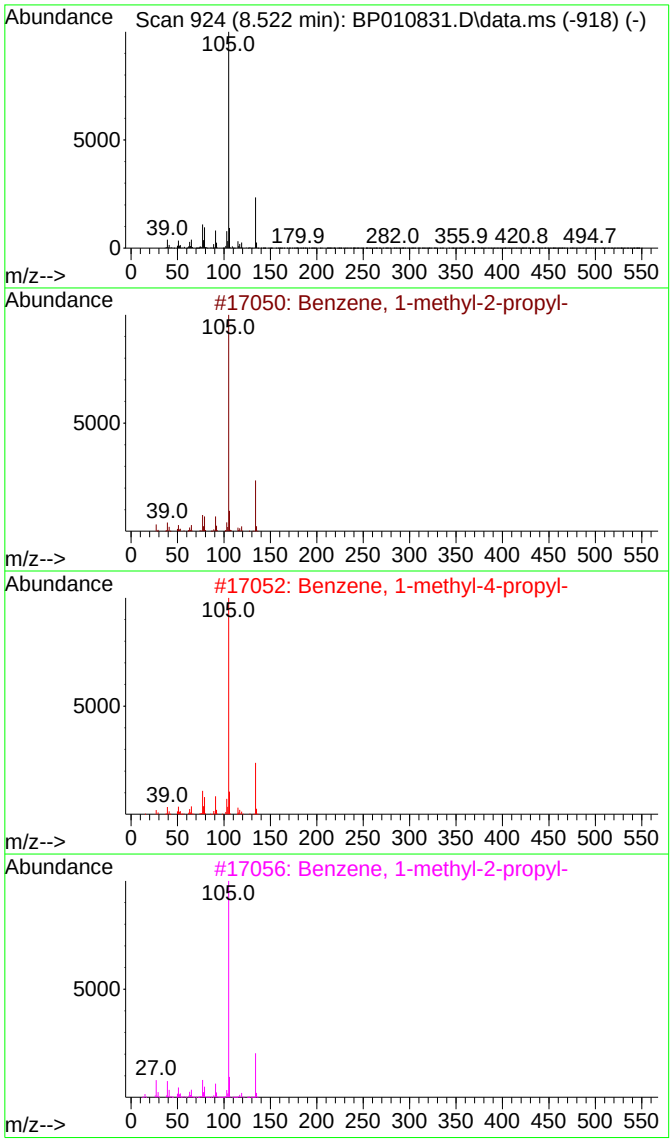
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

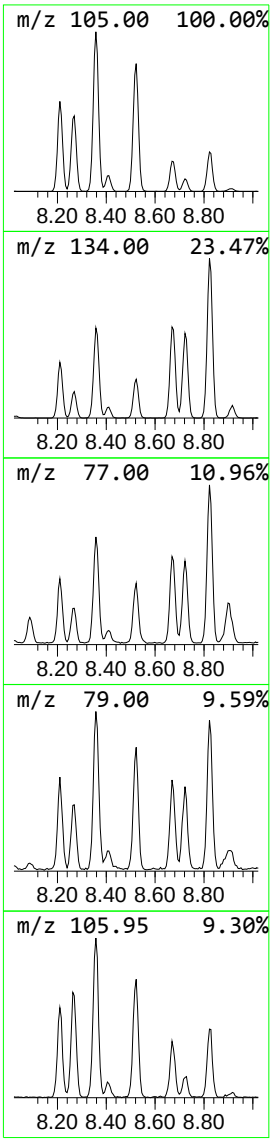
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Benzene, 1-methyl-2-propyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.522	6.52 ng/ul	760651	1,4-Dichlorobenzene-d4	7.716

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	94
2	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	94
3	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
4	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91





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Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

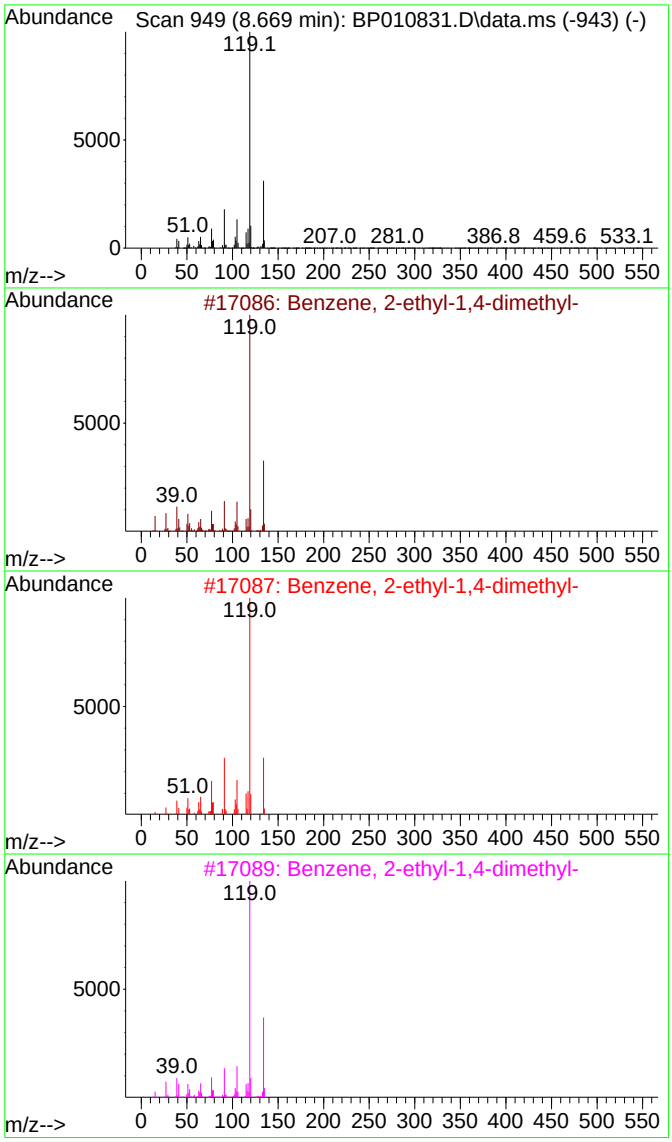
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

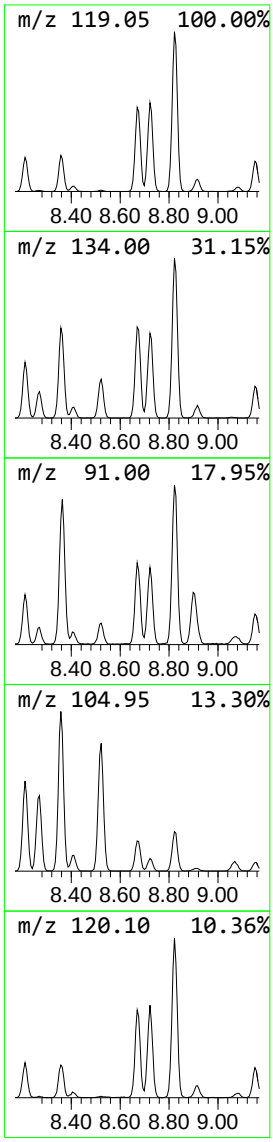
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.669	14.55 ng/ul	1696850	1,4-Dichlorobenzene-d4	7.716

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95





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Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

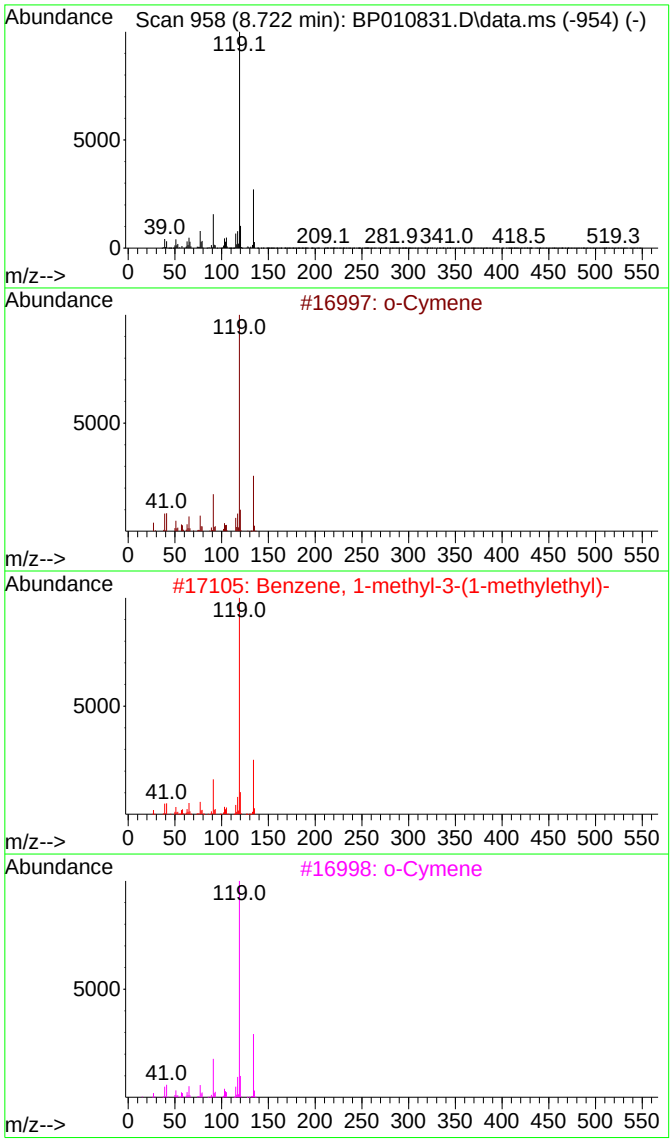
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

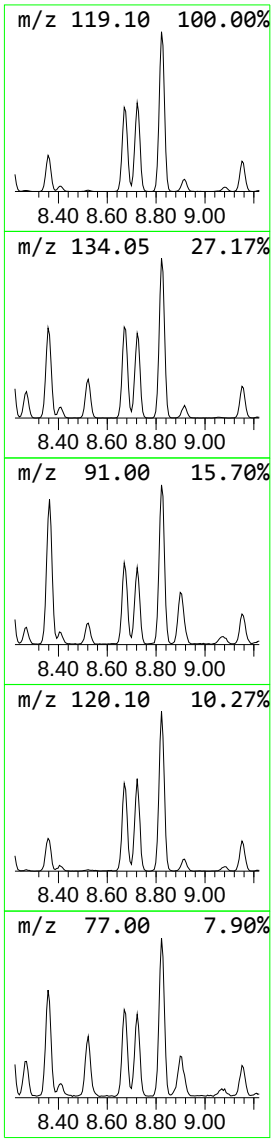
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 o-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.722	13.43 ng/ul	1565910	1,4-Dichlorobenzene-d4	7.716

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	97
3	o-Cymene	134	C10H14	000527-84-4	97
4	o-Cymene	134	C10H14	000527-84-4	97
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95





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Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

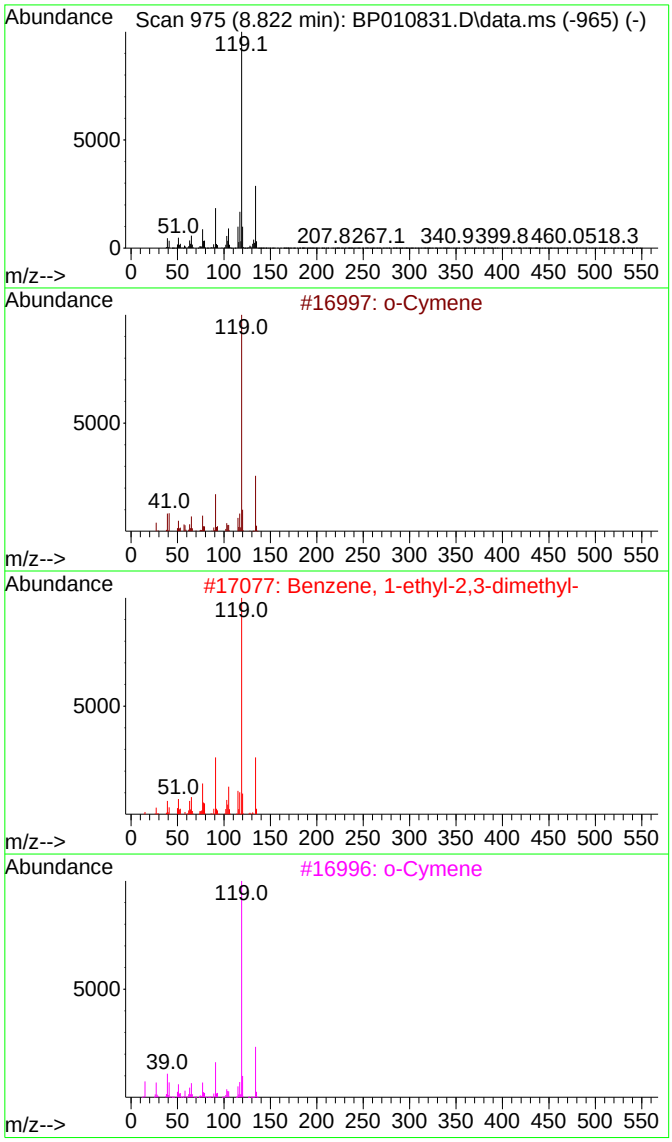
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

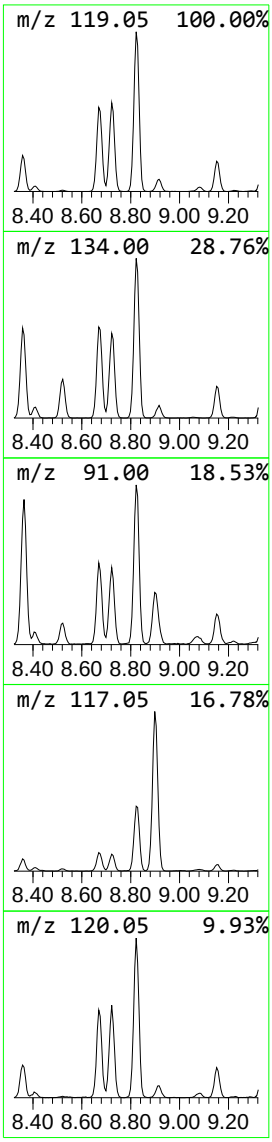
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.822	28.44 ng/ul	3316020	1,4-Dichlorobenzene-d4	7.716

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	95
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
3	o-Cymene	134	C10H14	000527-84-4	94
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
5	p-Cymene	134	C10H14	000099-87-6	93





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Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

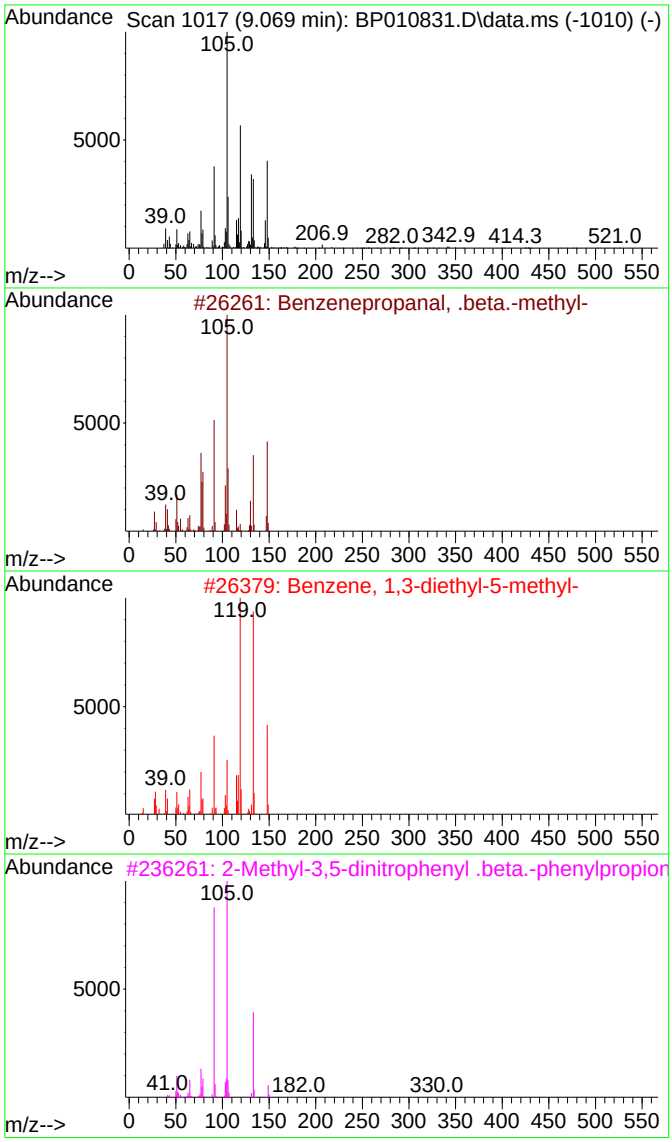
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

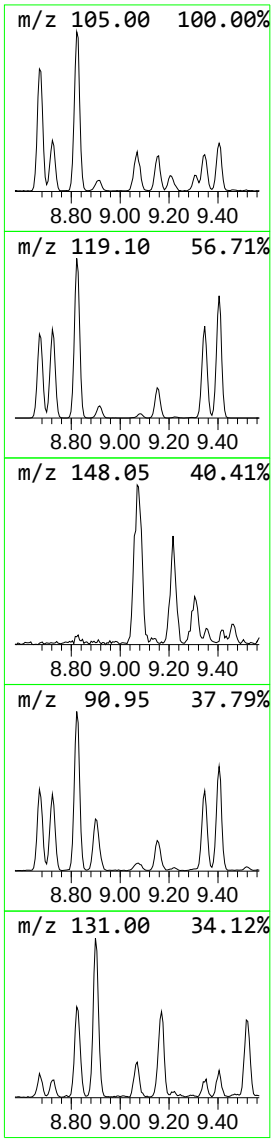
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzenepropanal, .beta.-met... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.069	2.00 ng/ul	233624	1,4-Dichlorobenzene-d4	7.716

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenepropanal, .beta.-methyl-	148	C10H12O	016251-77-7	50
2	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	49
3	2-Methyl-3,5-dinitrophenyl .beta...	330	C16H14N2O6	1000129-40-5	48
4	Benzenepropanal, .beta.-methyl-	148	C10H12O	016251-77-7	46
5	2-Methylindan-2-ol	148	C10H12O	033223-84-6	43





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Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

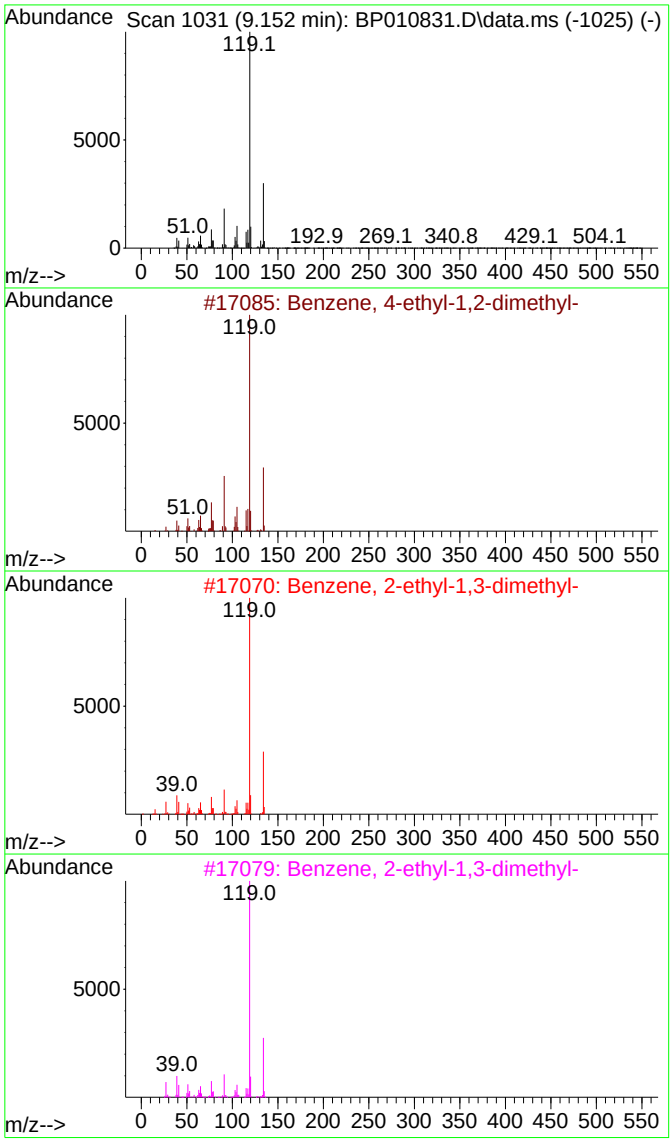
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

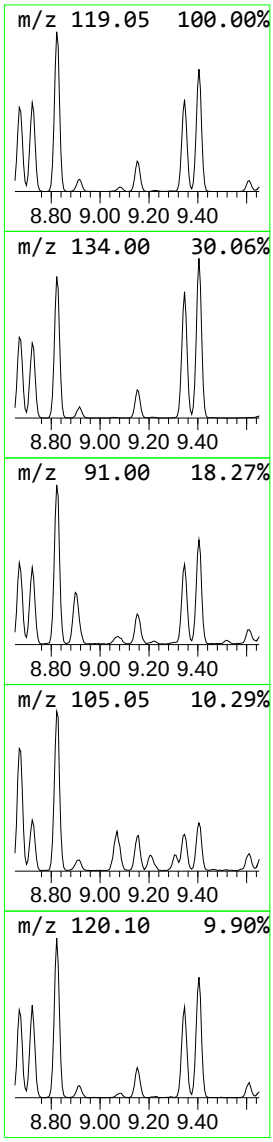
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.152	3.04 ng/ul	650435	Naphthalene-d8	10.516

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	o-Cymene	134	C10H14	000527-84-4	95
5	o-Cymene	134	C10H14	000527-84-4	95





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Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

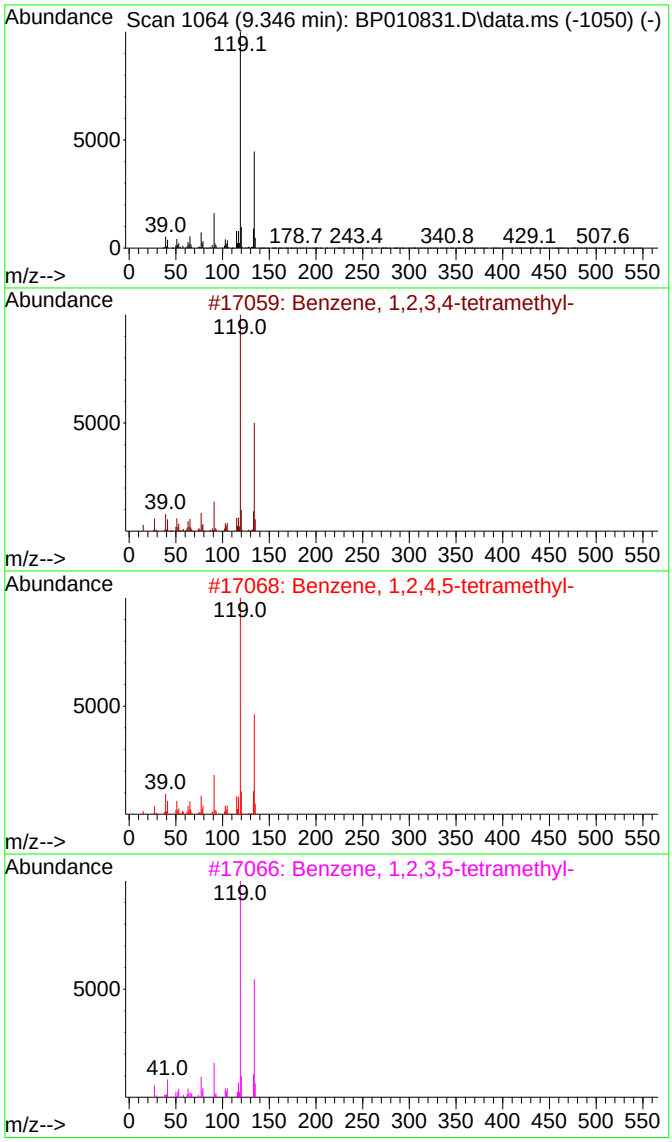
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

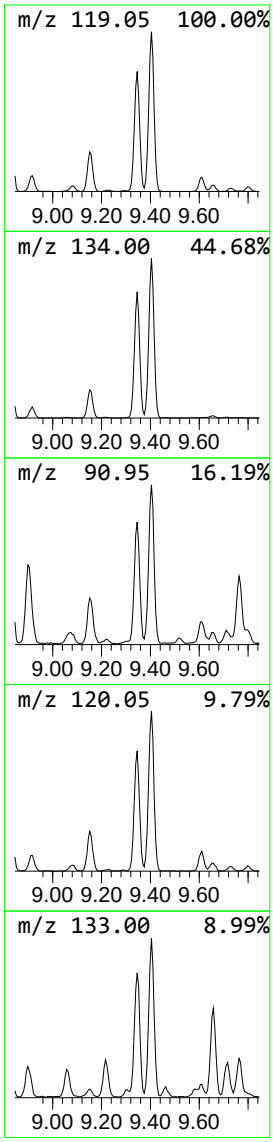
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.346	8.81 ng/ul	1888480	Naphthalene-d8	10.516

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95





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Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

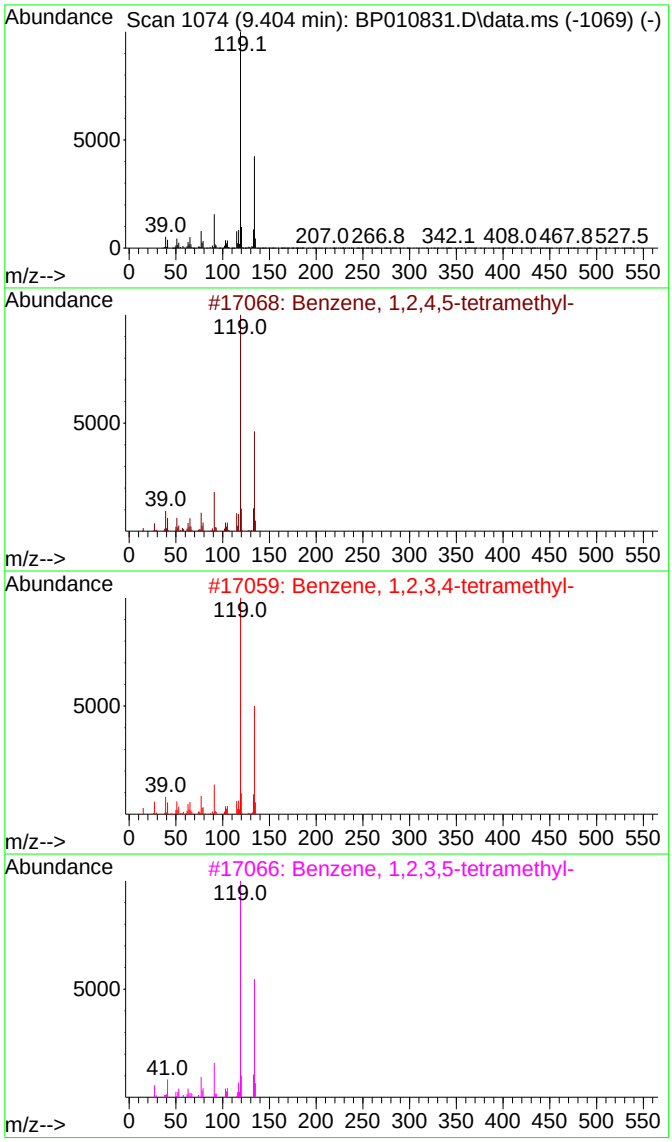
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

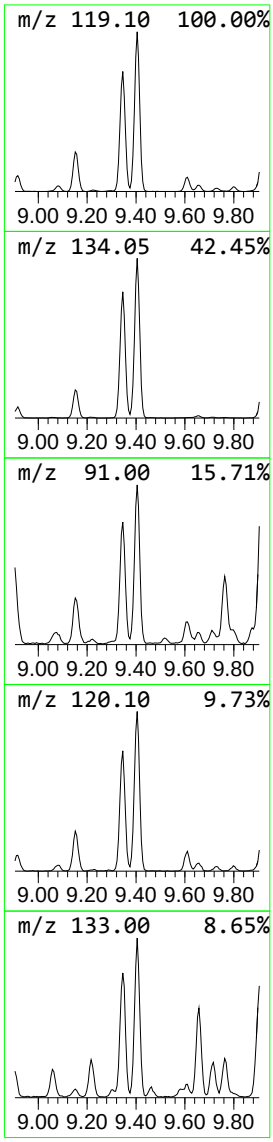
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.404	10.98 ng/ul	2352450	Naphthalene-d8	10.516

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95





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Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

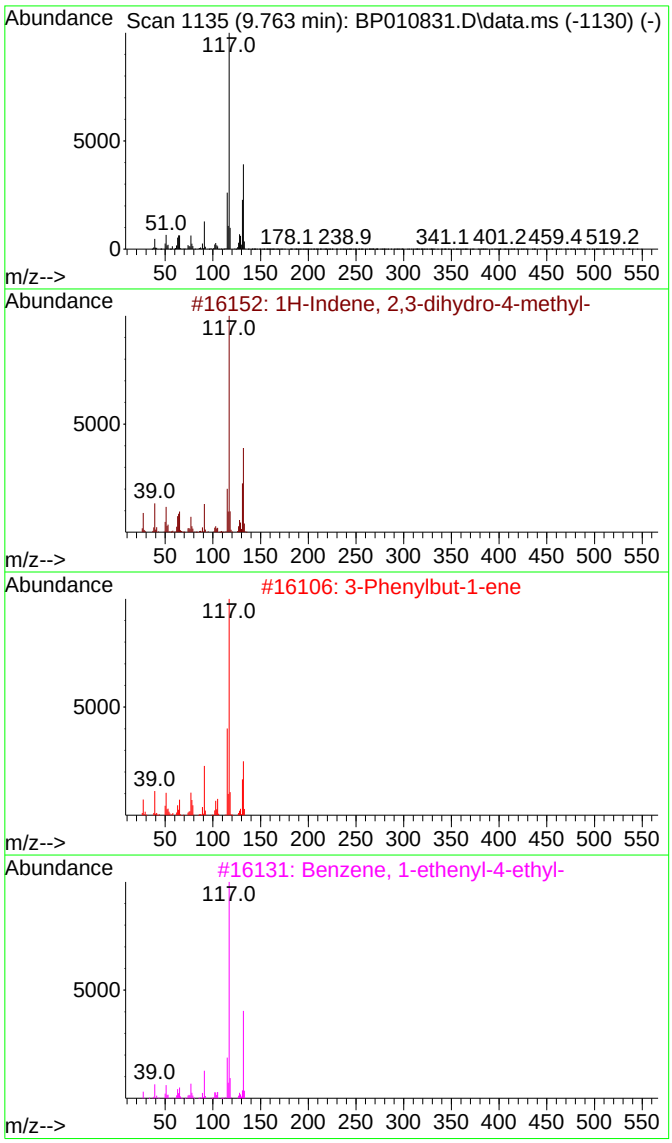
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

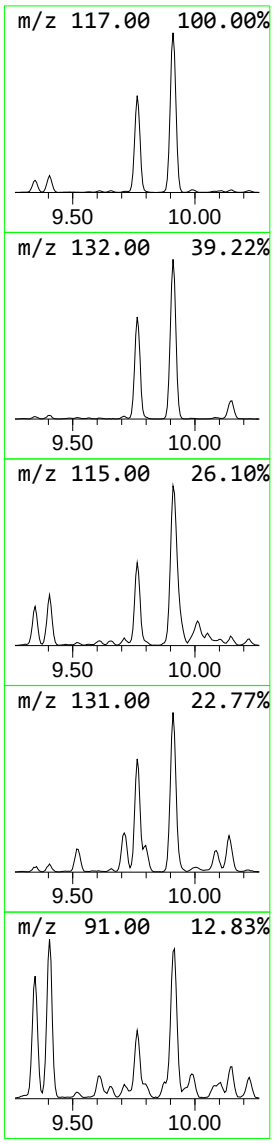
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.763	7.16 ng/ul	1534400	Naphthalene-d8	10.516

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
5	Indan, 1-methyl-	132	C10H12	000767-58-8	83





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Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

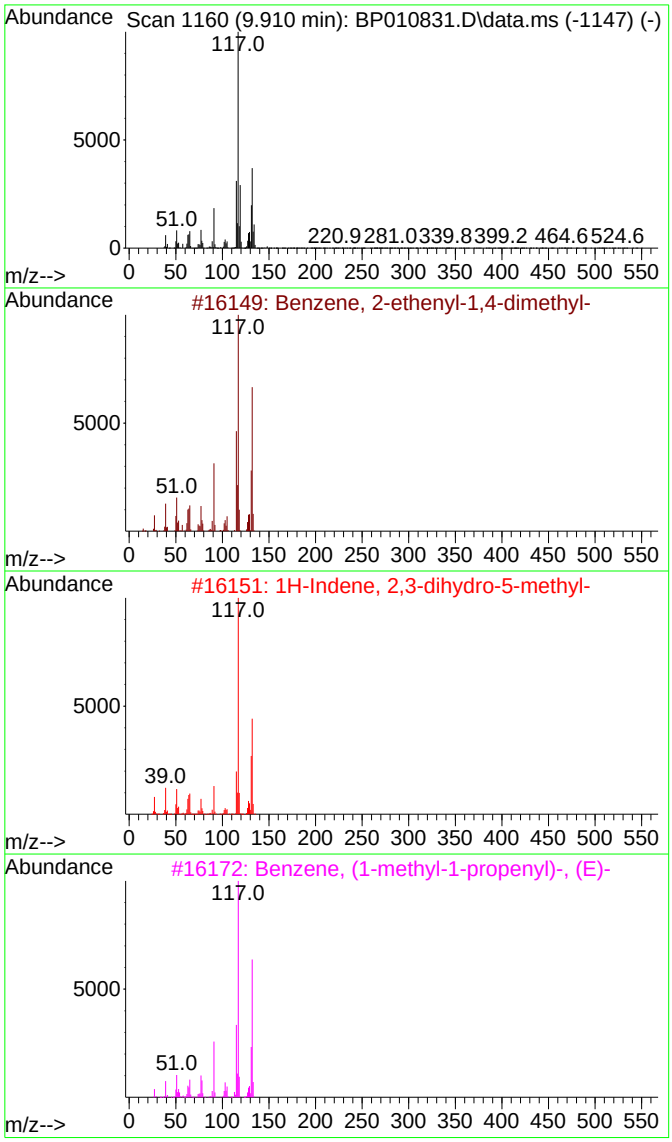
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

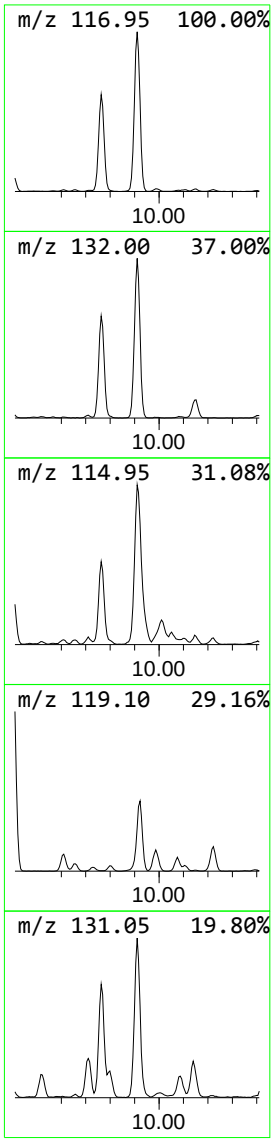
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.910	17.34 ng/ul	3716270	Naphthalene-d8	10.516

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	81
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	76
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76





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Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

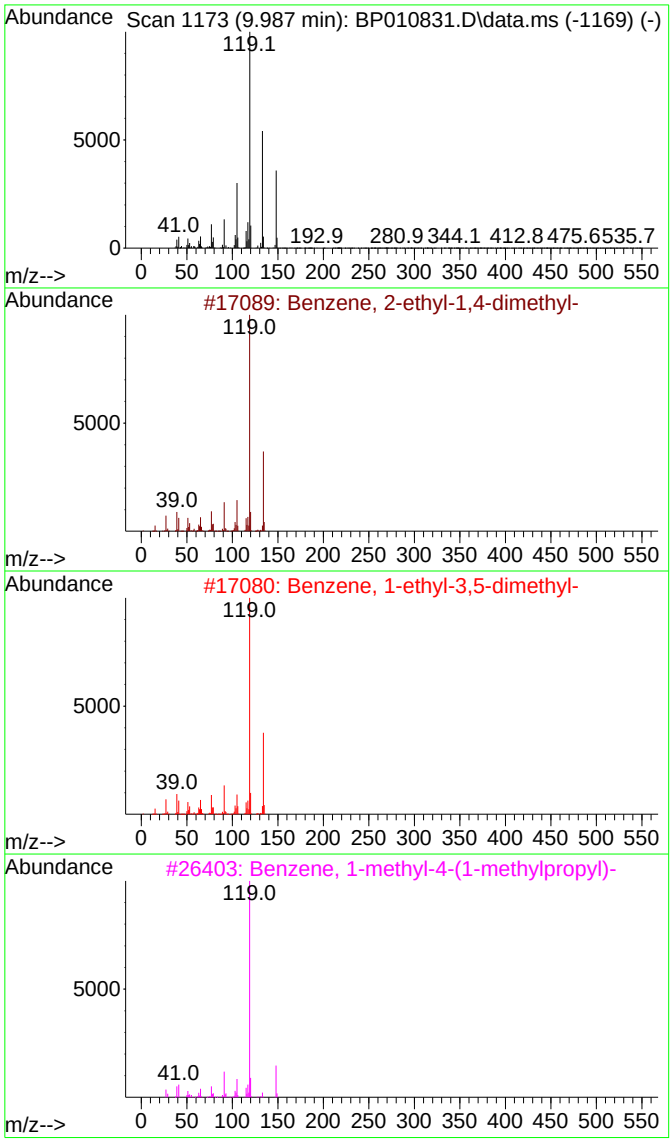
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

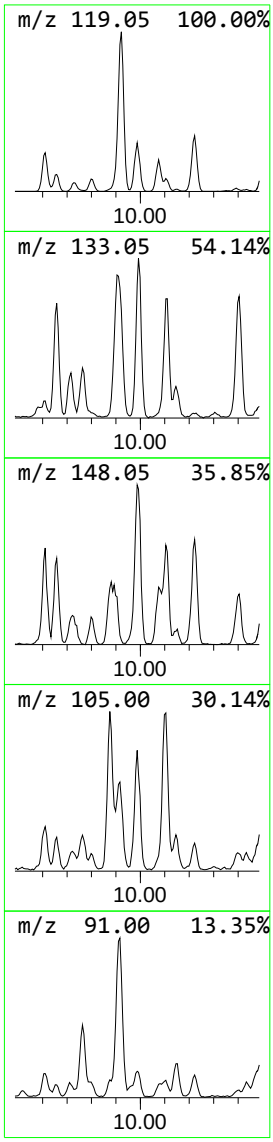
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.987	2.82 ng/ul	604059	Naphthalene-d8	10.516

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	60
2	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	60
3	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	58
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	53
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

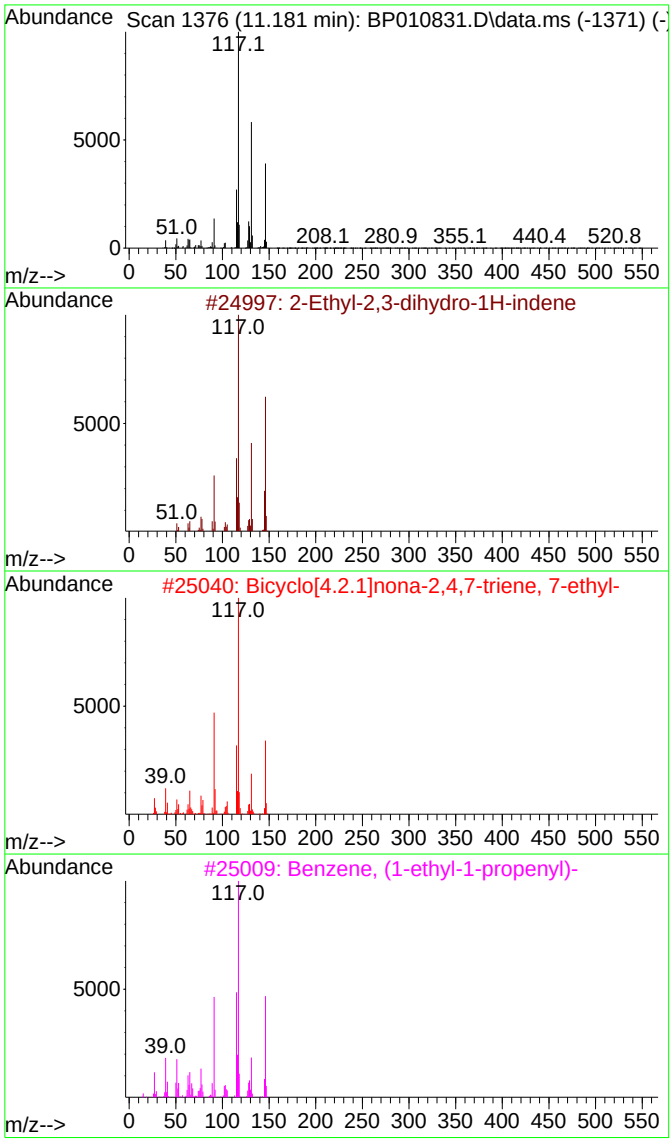
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

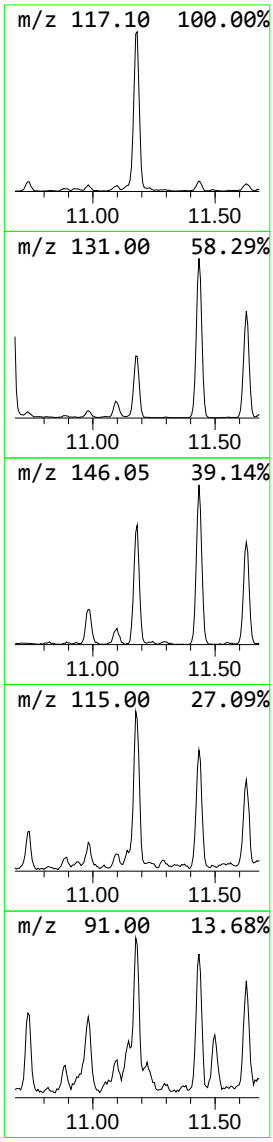
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 2-Ethyl-2,3-dihydro-1H-indene Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.181	2.10 ng/ul	449623	Naphthalene-d8	10.516

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Ethyl-2,3-dihydro-1H-indene	146	C11H14	056147-63-8	83
2	Bicyclo[4.2.1]nona-2,4,7-triene,...	146	C11H14	1000164-42-6	58
3	Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	53
4	Benzene, 1-pentenyl-	146	C11H14	000826-18-6	53
5	1H-Indene, 1-ethyl-2,3-dihydro-	146	C11H14	004830-99-3	52





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Data File : BP010831.D
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Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

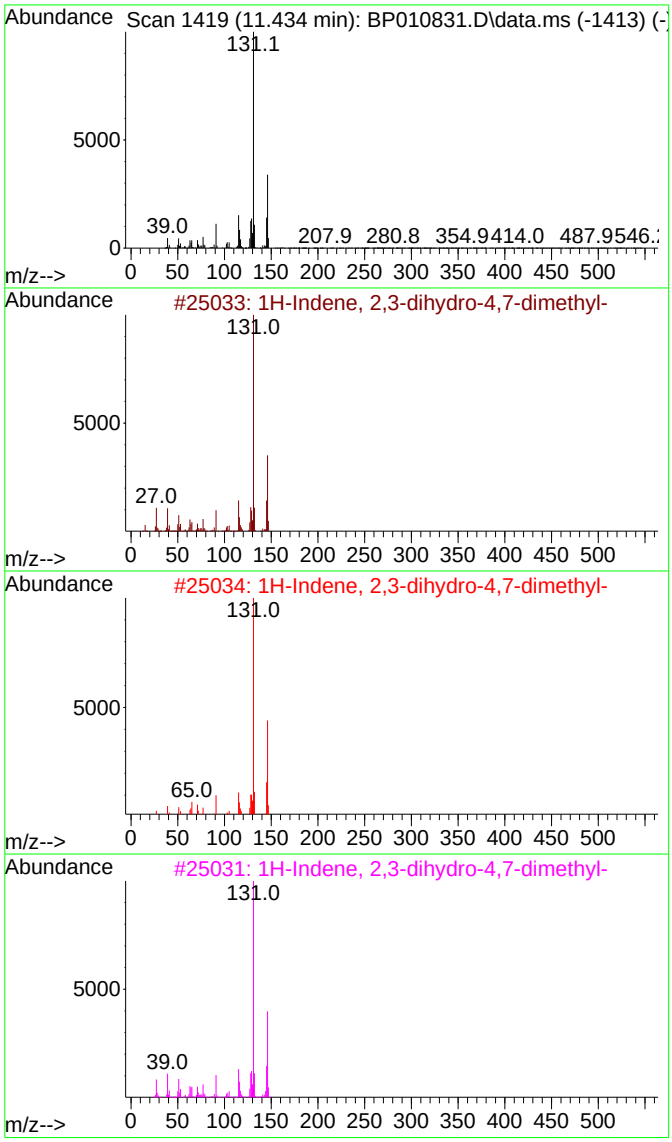
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

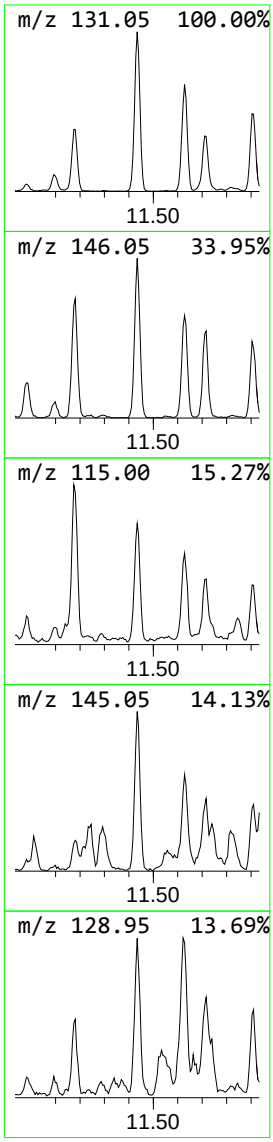
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.434	2.68 ng/ul	574669	Naphthalene-d8	10.516

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	95
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
4	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	94
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	94





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Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

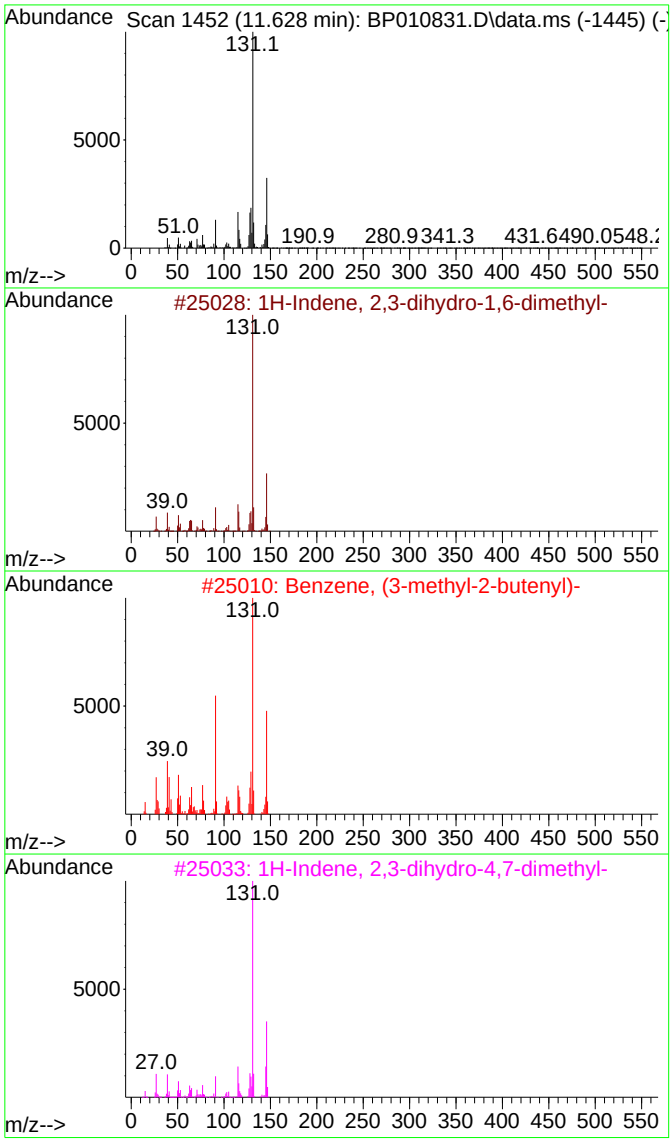
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

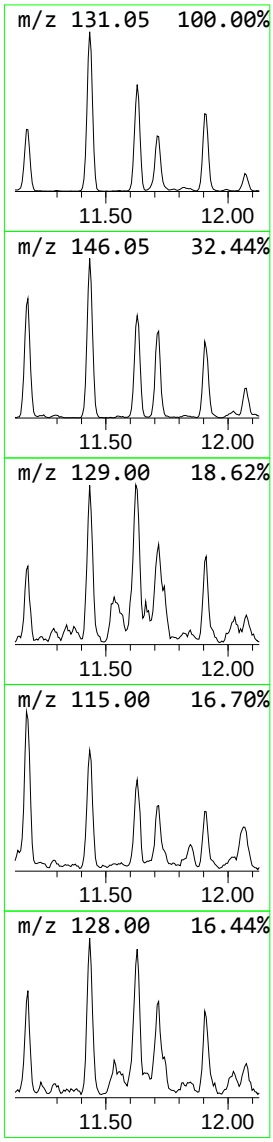
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.628	2.11 ng/ul	451347	Naphthalene-d8	10.516

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94
2	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	94
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
4	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
5	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	91





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Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

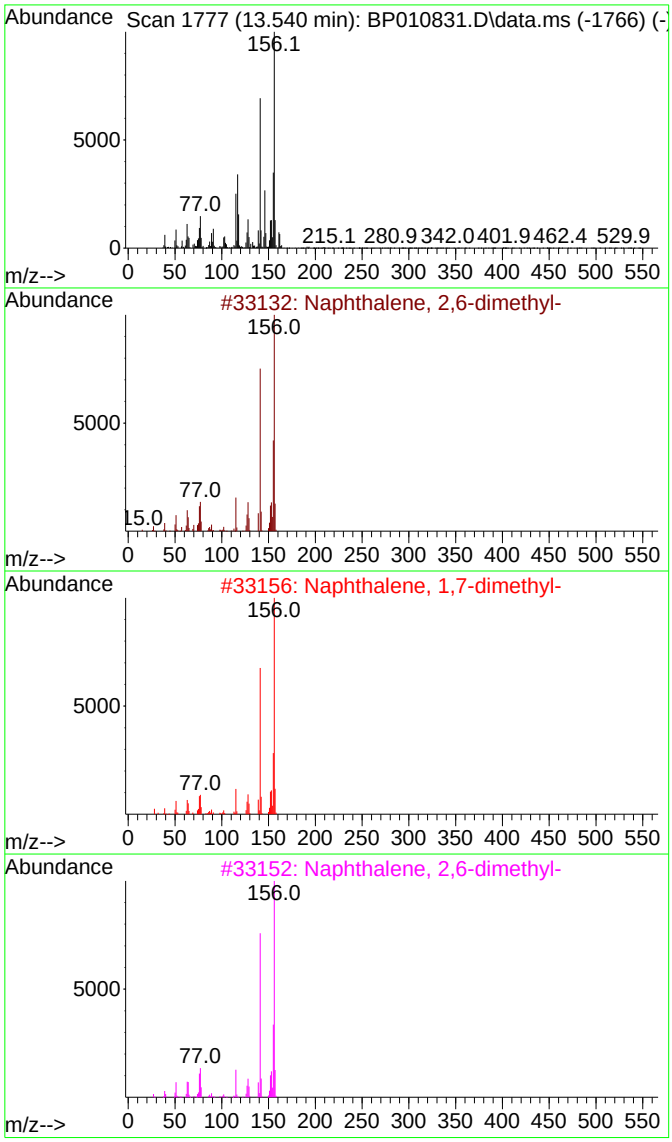
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

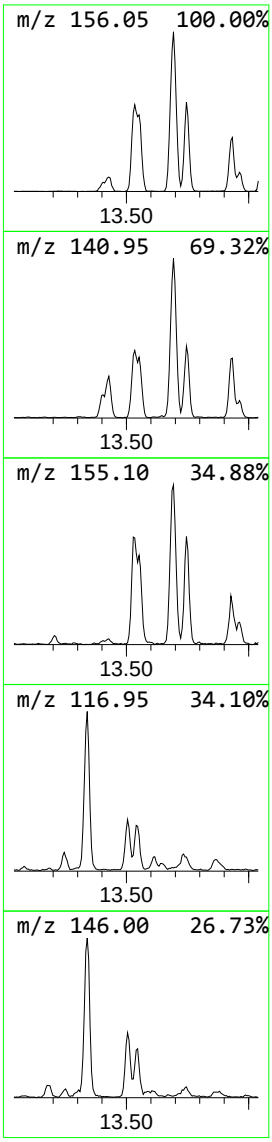
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 Naphthalene, 2,6-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.540	3.29 ng/ul	641496	Acenaphthene-d10	14.375

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
2	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
3	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96





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Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

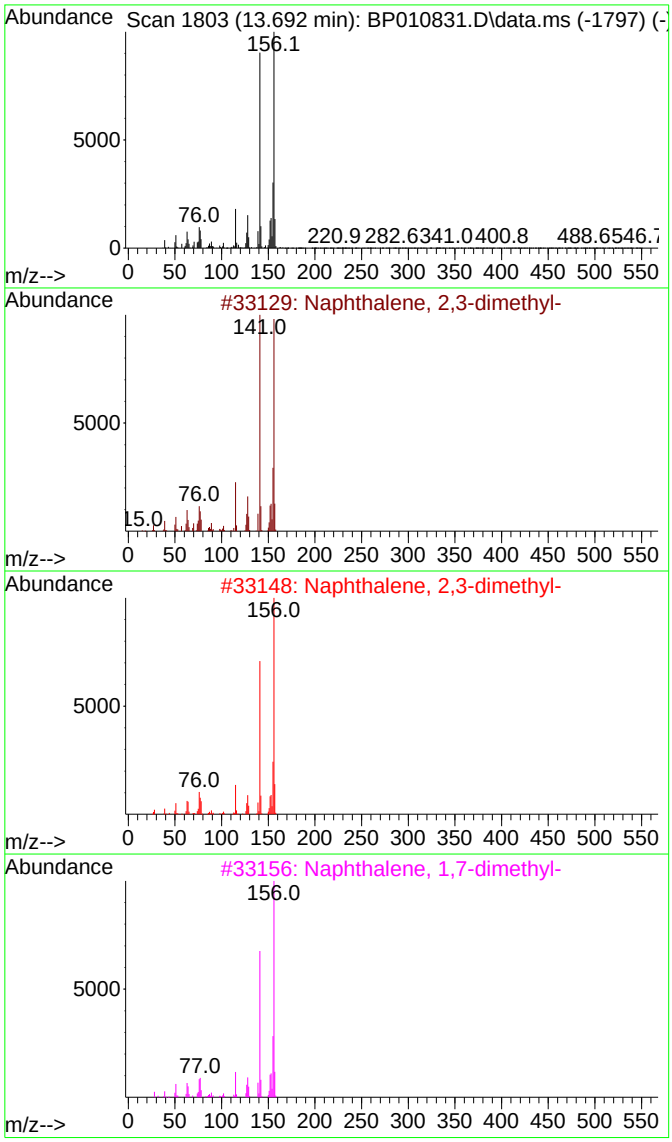
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

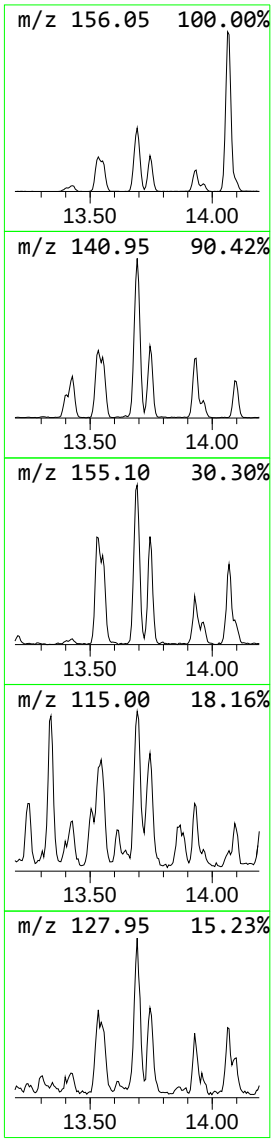
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 Naphthalene, 2,3-dimethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.693	3.02 ng/ul	589561	Acenaphthene-d10	14.375

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
4	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
Data File : BP010831.D
Acq On : 25 Jun 2022 16:44
Operator : CG/JU
Sample : N3355-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0C3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.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
2-Pentanone, 4-...	4.858	2.2	ng/ul	260705	1	7.716	2331760	20.0
Benzene, propyl-	6.699	16.5	ng/ul	1920230	1	7.716	2331760	20.0
Benzene, 1-ethy...	6.810	6.9	ng/ul	808695	1	7.716	2331760	20.0
Benzene, 1-ethy...	7.110	35.3	ng/ul	4116060	1	7.716	2331760	20.0
Indane	8.087	13.5	ng/ul	1577320	1	7.716	2331760	20.0
Benzene, 1,4-di...	8.210	9.0	ng/ul	1050980	1	7.716	2331760	20.0
Benzene, 1-meth...	8.269	4.3	ng/ul	499664	1	7.716	2331760	20.0
Benzene, 1,3-di...	8.357	15.8	ng/ul	1843090	1	7.716	2331760	20.0
Benzene, 1-meth...	8.522	6.5	ng/ul	760651	1	7.716	2331760	20.0
Benzene, 2-ethy...	8.669	14.6	ng/ul	1696850	1	7.716	2331760	20.0
o-Cymene	8.722	13.4	ng/ul	1565910	1	7.716	2331760	20.0
Benzene, 1-ethy...	8.822	28.4	ng/ul	3316020	1	7.716	2331760	20.0
Benzenepropanal...	9.069	2.0	ng/ul	233624	1	7.716	2331760	20.0
Benzene, 4-ethy...	9.152	3.0	ng/ul	650435	2	10.516	4285800	20.0
Benzene, 1,2,3,...	9.346	8.8	ng/ul	1888480	2	10.516	4285800	20.0
Benzene, 1,2,4,...	9.404	11.0	ng/ul	2352450	2	10.516	4285800	20.0
1H-Indene, 2,3-...	9.763	7.2	ng/ul	1534400	2	10.516	4285800	20.0
Benzene, 2-ethe...	9.910	17.3	ng/ul	3716270	2	10.516	4285800	20.0
Benzene, 1-ethy...	9.987	2.8	ng/ul	604059	2	10.516	4285800	20.0
2-Ethyl-2,3-dih...	11.181	2.1	ng/ul	449623	2	10.516	4285800	20.0
1H-Indene, 2,3-...	11.434	2.7	ng/ul	574669	2	10.516	4285800	20.0
1H-Indene, 2,3-...	11.628	2.1	ng/ul	451347	2	10.516	4285800	20.0
Naphthalene, 2,...	13.540	3.3	ng/ul	641496	3	14.375	3899540	20.0
Naphthalene, 2,...	13.693	3.0	ng/ul	589561	3	14.375	3899540	20.0