LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

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-BP071024.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP021242.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.170	28	31	38	rVB	23551	31227	0.64%	0.045%
2	3.593	100	103	123	rBV	116162	264196	5.42%	0.378%
3	6.134	527	535	542	rBV2	20986	36846	0.76%	0.053%
4	6.299	555	563	570	rBV2	17851	34815	0.71%	0.050%
5	6.852	651	657	672	rBV	182497	411022	8.43%	0.588%
6	6.975	672	678	690	rVB	465420	770571	15.80%	1.102%
7	7.181	706	713	724	rBV	586840	1013622	20.78%	1.450%
8	7.269	724	728	736	rVV3	23246	49045	1.01%	0.070%
9	7.622	782	788	799	rVV	600092	1043693	21.40%	1.493%
10	7.728	799	806	814	rVV	80864	148797	3.05%	0.213%
11	7.981	841	849	864	rBV	2387152	3840916	78.74%	5.495%
12	8.281	895	900	907	rBV	67826	121510	2.49%	0.174%
13	8.358	907	913	928	rVV	491584	1041568	21.35%	1.490%
14	8.481	929	934	937	rVV4	16474	38014	0.78%	0.054%
15	8.528	937	942	950	rVV	64796	116278	2.38%	0.166%
16	8.616	952	957	963	rVV	26629	45666	0.94%	0.065%
17	8.693	964	970	979	rVV3	35303	100161	2.05%	0.143%
18	8.787	979	986	1002	rVB	544306	1057171	21.67%	1.512%
19	8.963	1009	1016	1026	rVV2	92168	189595	3.89%	0.271%
20	9.087	1027	1037	1042	rVV	99996	251224	5.15%	0.359%
21	9.146	1042	1047	1058	rVV2	52786	123998	2.54%	0.177%
22	9.499	1100	1107	1121	rBV	433663	894296	18.33%	1.279%
23	9.640	1123	1131	1138	rVV2	42284	102409	2.10%	0.147%
24	9.710	1138	1143	1149	rVB2	24477	42565	0.87%	0.061%
25	9.781	1149	1155	1168	rBV	87221	185069	3.79%	0.265%
26	10.058	1192	1202	1215	rBV	608990	1429415	29.30%	2.045%
27	10.393	1246	1259	1266	rBV	884914	1519687	31.15%	2.174%
28	10.552	1279	1286	1306	rBV2	651562	1351236	27.70%	1.933%
29	10.987	1353	1360	1372	rVV	897873	1668677	34.21%	2.387%
30	11.081	1373	1376	1382	rVV3	21792	38974	0.80%	0.056%
31	11.163	1384	1390	1397	rVB4	22410	53930	1.11%	0.077%
32	11.293	1409	1412	1423	rVB6	13000	28555	0.59%	0.041%
33	11.434	1425	1436	1438	rBV7	14314	28426	0.58%	0.041%
34	11.469	1438	1442	1446	rVV2	29678	55892	1.15%	0.080%
35	11.510	1446	1449	1456	rVB	42529	75117	1.54%	0.107%
36	11.604	1456	1465	1479	rVB	218588	397955	8.16%	0.569%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

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-BP071024.MA.M
Title : SVOA CALIBRATION

37	11.740	1481	1488	1491	rBV3	23165	43020	0.88%	0.062%
38	11.822	1498	1502	1509	rVB6	36794	71739	1.47%	0.103%
39	11.952	1517	1524	1529	rVV	72328	131057	2.69%	0.188%
40	12.022	1529	1536	1541	rVV	254915	497244	10.19%	0.711%
41	12.069	1541	1544	1550	rVV4	77339	142653	2.92%	0.204%
42	12.187	1550	1564	1574	rVV	2571488	4345262	89.08%	6.217%
43	12.281	1574	1580	1595	rVB7	63127	225383	4.62%	0.322%
44	12.416	1596	1603	1612	rBV	366669	651224	13.35%	0.932%
45	12.528	1619	1622	1627	rVB2	21947	29073	0.60%	0.042%
46	12.599	1629	1634	1639	rBV	387119	519647	10.65%	0.743%
47	12.687	1642	1649	1654	rVV3	352890	697129	14.29%	0.997%
48	12.734	1654	1657	1659	rVV2	64195	96035	1.97%	0.137%
49	12.787	1659	1666	1676	rVB2	380935	784315	16.08%	1.122%
50	12.863	1676	1679	1688	rVB6	13341	29076	0.60%	0.042%
51	12.963	1688	1696	1699	rBV3	23496	51902	1.06%	0.074%
52	13.016	1699	1705	1709	rBV2	230757	383918	7.87%	0.549%
53	13.069	1709	1714	1718	rVV	438350	780765	16.01%	1.117%
54	13.116	1718	1722	1738	rVB2	333757	740775	15.19%	1.060%
55	13.334	1750	1759	1765	rVB2	60832	149179	3.06%	0.213%
56	13.410	1766	1772	1776	rBV	509794	768818	15.76%	1.100%
57	13.463	1776	1781	1792	rVB	279441	550824	11.29%	0.788%
58	13.557	1792	1797	1801	rBV2	70500	124339	2.55%	0.178%
59	13.610	1801	1806	1813	rBV	435556	687007	14.08%	0.983%
60	13.681	1813	1818	1825	rBV	1509554	2080525	42.65%	2.977%
61	13.734	1825	1827	1829	rVV	33308	32220	0.66%	0.046%
62	13.775	1829	1834	1840	rVV	741380	1271328	26.06%	1.819%
63	13.834	1840	1844	1849	rVB4	85553	147147	3.02%	0.211%
64	13.881	1849	1852	1858	rBV	57870	85772	1.76%	0.123%
65	13.969	1858	1867	1874	rBV	1801042	2567569	52.64%	3.673%
66	14.081	1882	1886	1895	rVB2	37785	72187	1.48%	0.103%
67	14.163	1895	1900	1907	rBV2	33877	70746	1.45%	0.101%
68	14.275	1907	1919	1926	rVV	1353148	2292845	47.00%	3.280%
69	14.346	1927	1931	1938	rVB3	70390	133984	2.75%	0.192%
70	14.410	1938	1942	1949	rBV4	62537	136132	2.79%	0.195%
71	14.487	1950	1955	1962	rVB2	60834	116218	2.38%	0.166%
72	14.551	1962	1966	1972	rVB4	23907	37795	0.77%	0.054%
73	14.634	1972	1980	1989	rBV2	108274	298023	6.11%	0.426%
74	14.775	1999	2004	2006	rBV	67495	104720	2.15%	0.150%
75	14.816	2006	2011	2025	rVV	414751	716197	14.68%	1.025%
76	14.928	2026	2030	2036	rVV3	48732	81926	1.68%	0.117%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

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-BP071024.MA.M
Title : SVOA CALIBRATION

77	14.998	2039	2042	2051	rVB5	23152	54241	1.11%	0.078%
78	15.287	2083	2091	2100	rVV	2101149	3219754	66.00%	4.607%
79	15.463	2115	2121	2129	rBV	546902	1005417	20.61%	1.438%
80	15.787	2171	2176	2188	rVB	10606	39183	0.80%	0.056%
81	16.287	2255	2261	2265	rBV3	45218	93116	1.91%	0.133%
82	16.828	2348	2353	2358	rBV	32433	45629	0.94%	0.065%
83	17.010	2380	2384	2386	rBV4	19531	27542	0.56%	0.039%
84	17.092	2392	2398	2405	rBV	1777129	2552829	52.33%	3.652%
85	17.192	2409	2415	2424	rBV2	2152097	3580255	73.40%	5.122%
86	17.592	2480	2483	2490	rVB	36822	50368	1.03%	0.072%
87	17.787	2511	2516	2521	rVB	81501	104627	2.14%	0.150%
88	18.028	2551	2557	2567	rBV2	255941	508637	10.43%	0.728%
89	18.310	2601	2605	2610	rVB2	31274	41799	0.86%	0.060%
90	18.928	2707	2710	2714	rBV5	18229	28831	0.59%	0.041%
91	18.975	2715	2718	2720	rVV2	48248	52995	1.09%	0.076%
92	19.004	2720	2723	2728	rVB	64086	78079	1.60%	0.112%
93	19.116	2739	2742	2745	rBV2	28167	41840	0.86%	0.060%
94	19.145	2745	2747	2751	rVB	40291	46207	0.95%	0.066%
95	19.210	2754	2758	2761	rBV	38045	53852	1.10%	0.077%
96	19.357	2778	2783	2790	rBV2	156502	301809	6.19%	0.432%
97	19.581	2815	2821	2834	rBV	3137009	4797024	98.34%	6.863%
98	21.539	3148	3154	3171	rVB2	1833366	3359263	68.86%	4.806%
99	24.610	3668	3676	3691	rVB	1788971	4878058	100.00%	6.979%
100	24.845	3708	3716	3736	rVB	1232238	3586540	73.52%	5.131%

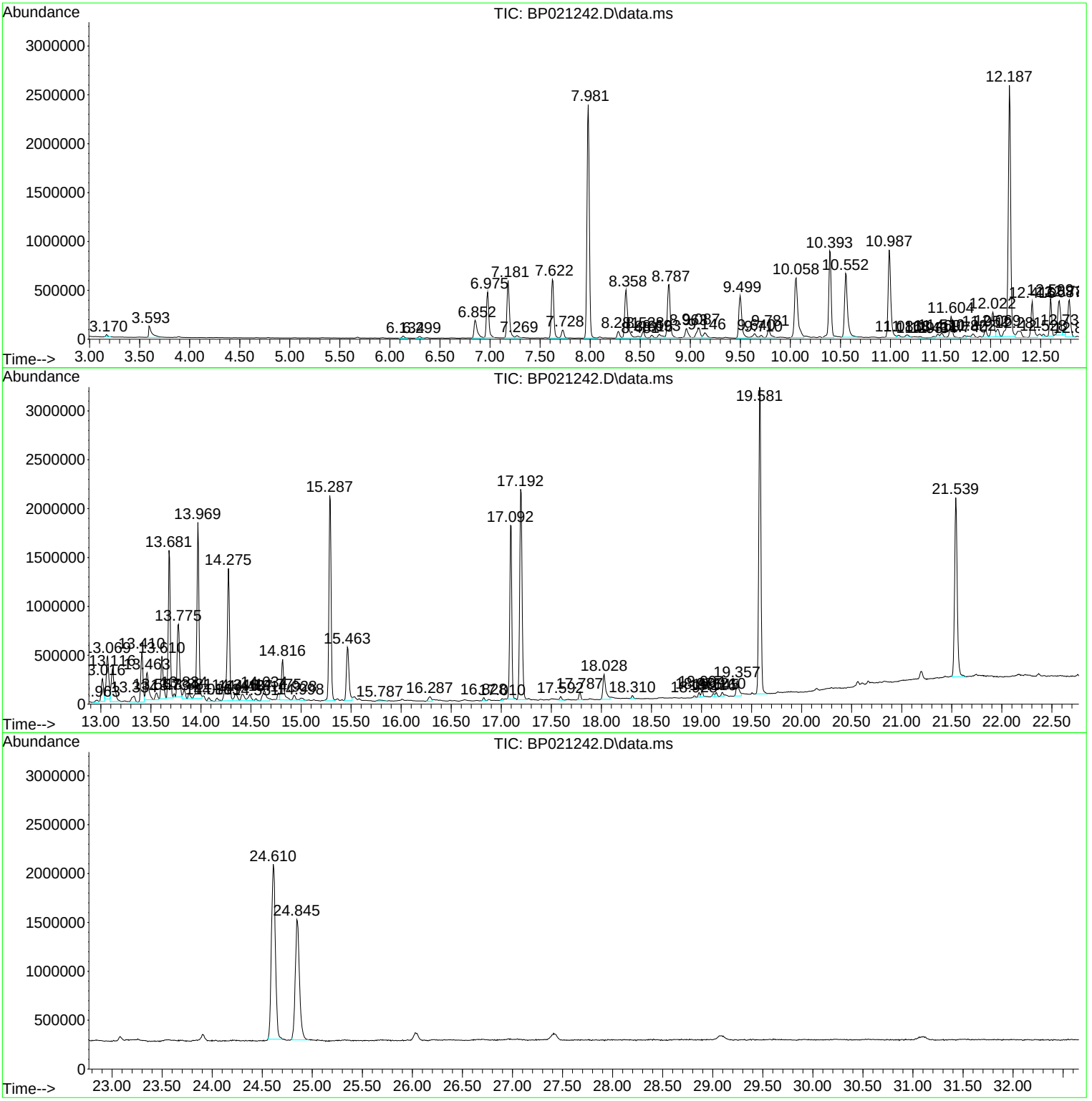
Sum of corrected areas: 69895751

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.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\BP072924\
Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

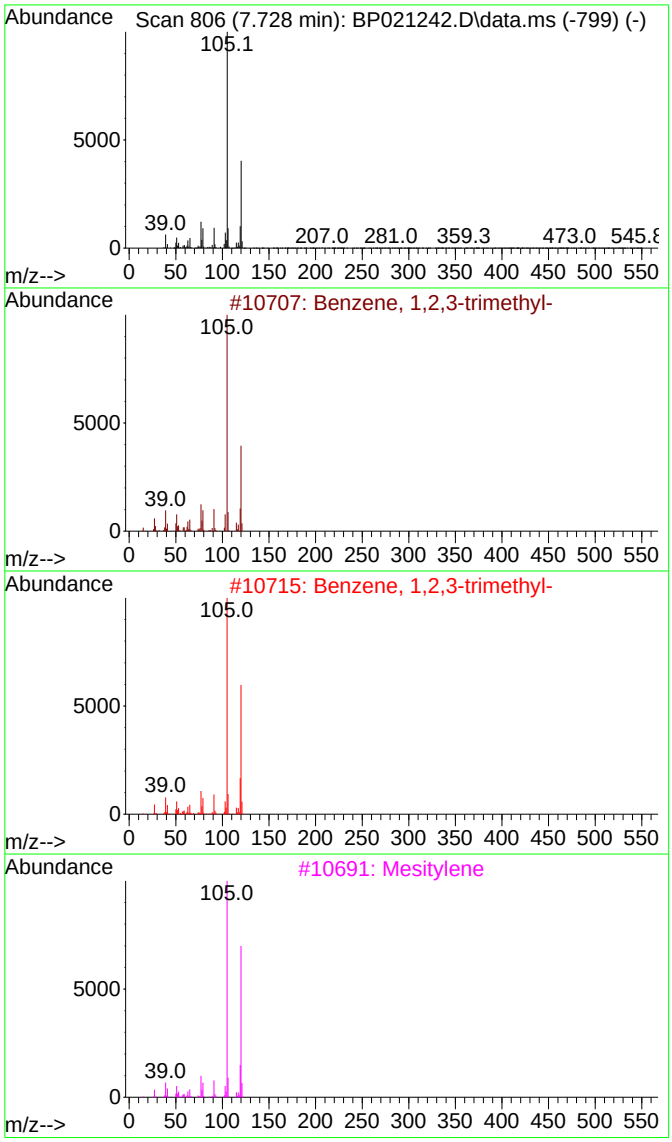
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

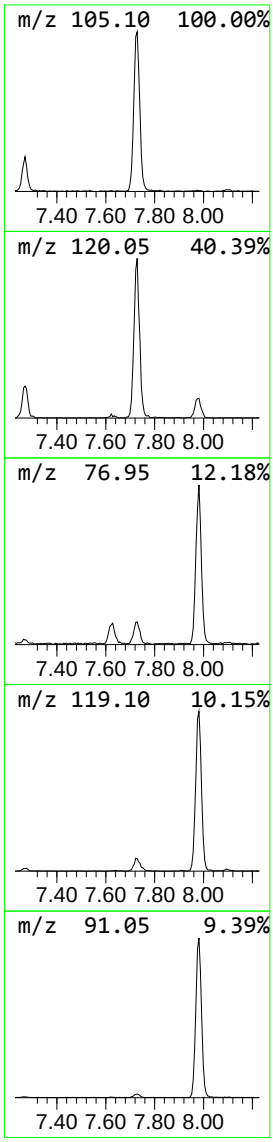
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzene, 1,2,3-trimethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.728	2.85 ng/ul	148797	1,4-Dichlorobenzene-d4	7.622

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Mesitylene	120	C9H12	000108-67-8	94
4	Mesitylene	120	C9H12	000108-67-8	94
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

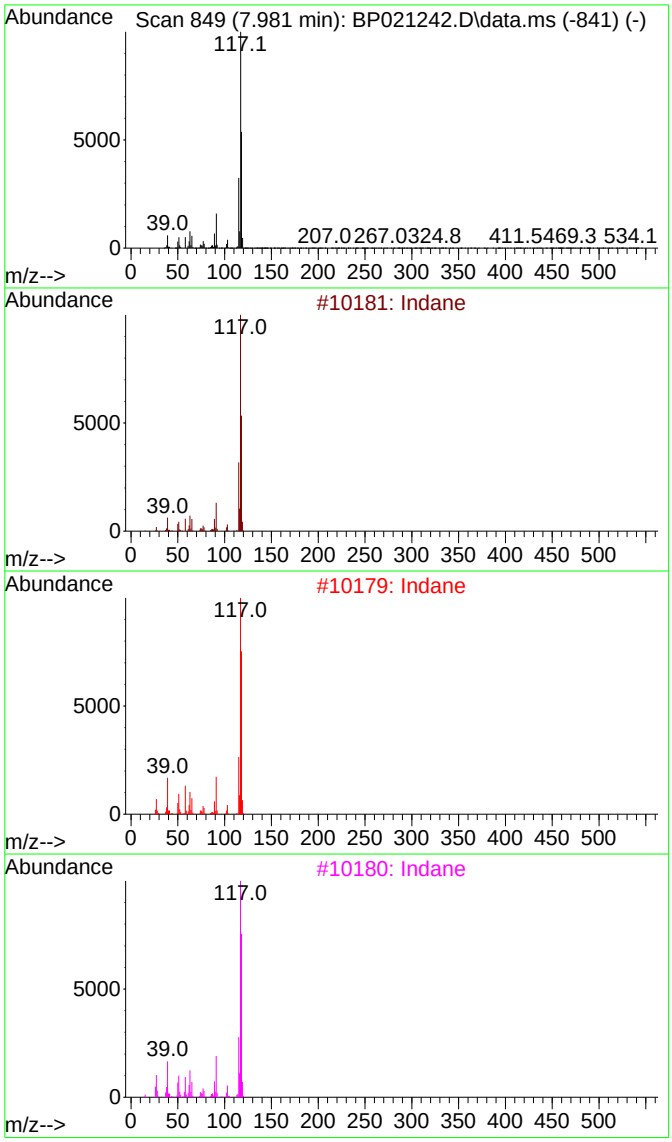
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

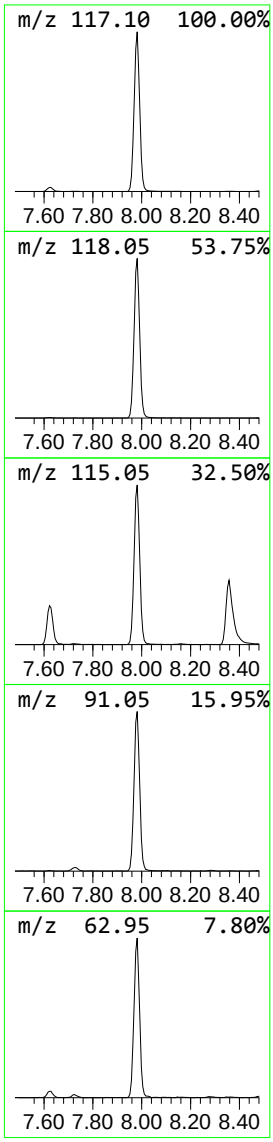
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.981	73.60 ng/ul	3840920	1,4-Dichlorobenzene-d4	7.622

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	90
4	Indane	118	C9H10	000496-11-7	87
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

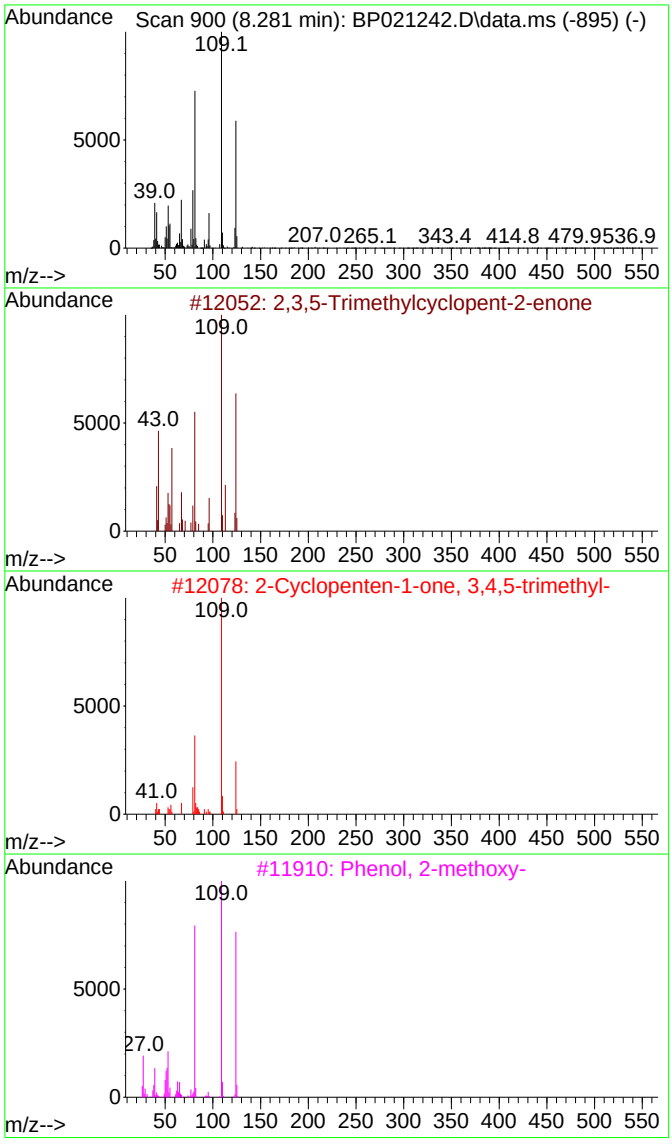
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

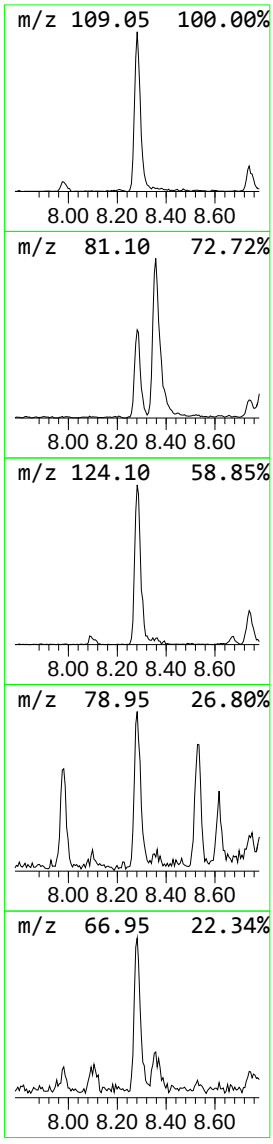
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2,3,5-Trimethylcyclopent-2-... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.281	2.33 ng/ul	121510	1,4-Dichlorobenzene-d4	7.622

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,5-Trimethylcyclopent-2-enone	124	C8H12O	1000427-08-7	91
2	2-Cyclopenten-1-one, 3,4,5-trime...	124	C8H12O	055683-21-1	74
3	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	70
4	2-Cyclopenten-1-one, 3,4,4-trime...	124	C8H12O	030434-65-2	68
5	2-Cyclopenten-1-one, 2,3,4-trime...	124	C8H12O	028790-86-5	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

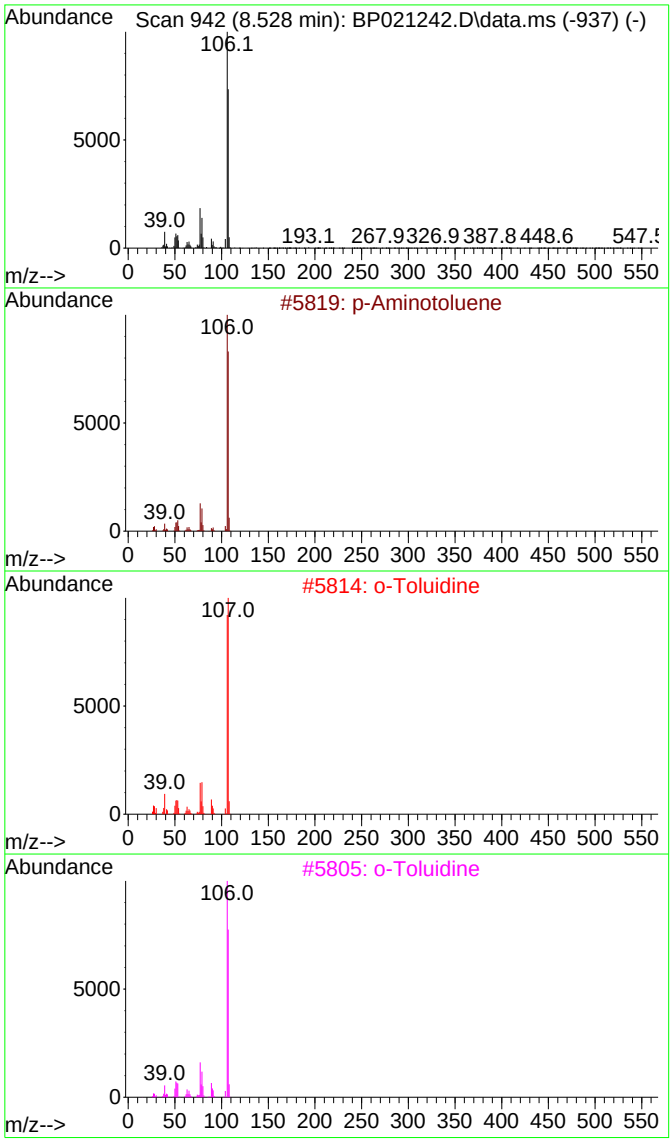
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

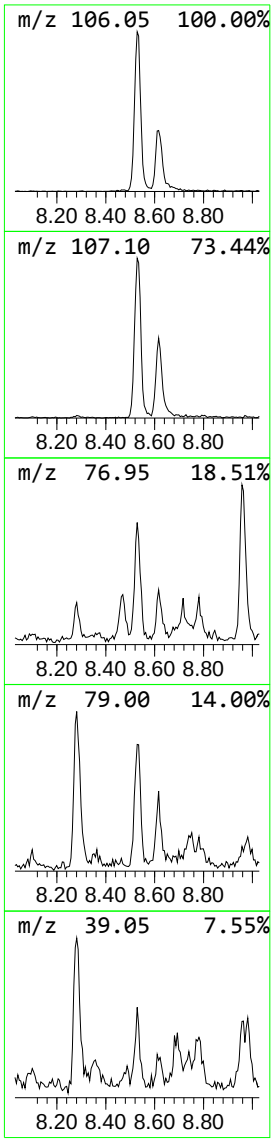
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 p-Aminotoluene Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.528	2.23 ng/ul	116278	1,4-Dichlorobenzene-d4	7.622

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Aminotoluene	107	C7H9N	000106-49-0	94
2	o-Toluidine	107	C7H9N	000095-53-4	94
3	o-Toluidine	107	C7H9N	000095-53-4	94
4	o-Toluidine	107	C7H9N	000095-53-4	91
5	Aniline, N-methyl-	107	C7H9N	000100-61-8	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

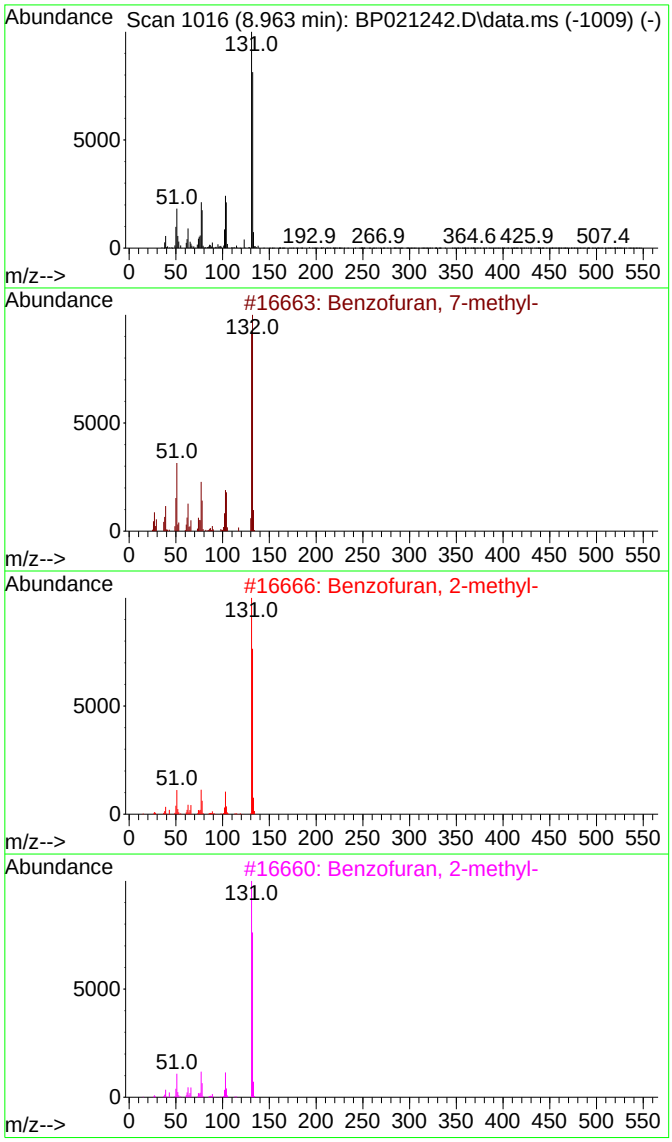
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

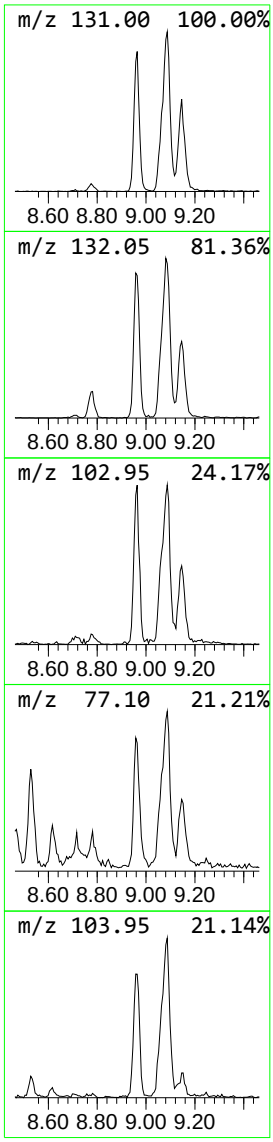
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzofuran, 7-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.963	3.63 ng/ul	189595	1,4-Dichlorobenzene-d4	7.622

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
5	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	90





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Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

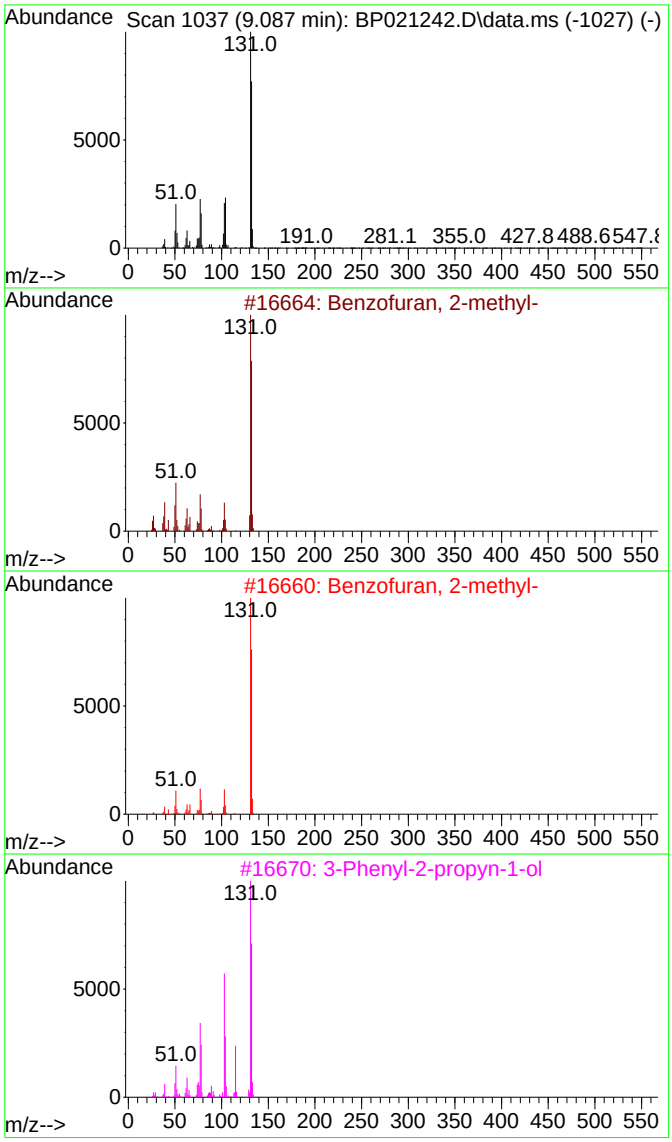
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

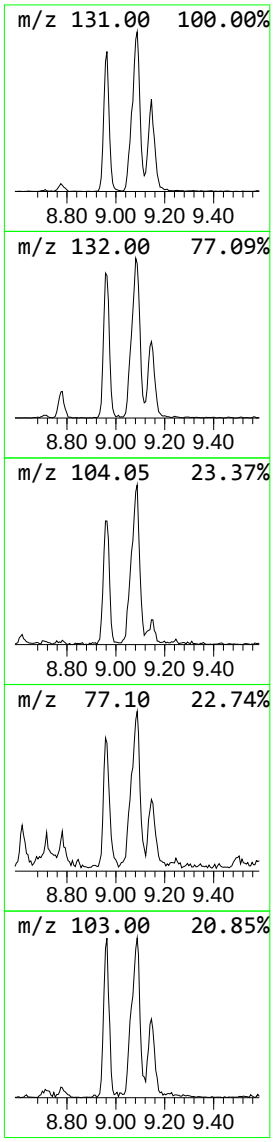
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzofuran, 2-methyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.087	3.31 ng/ul	251224	Naphthalene-d8	10.393

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
3	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	90
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	87
5	2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	87





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Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

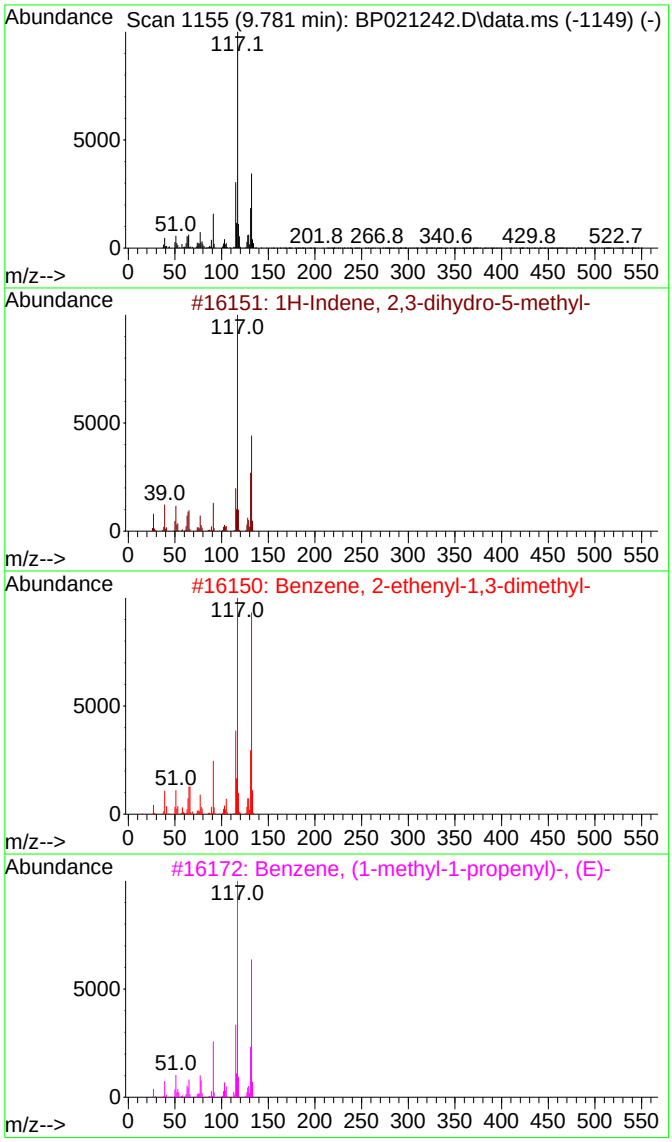
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

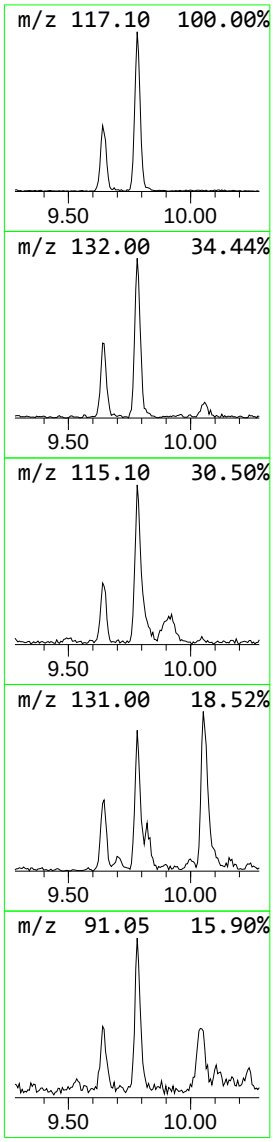
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.781	2.44 ng/ul	185069	Naphthalene-d8	10.393

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
2	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	91
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	90





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Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

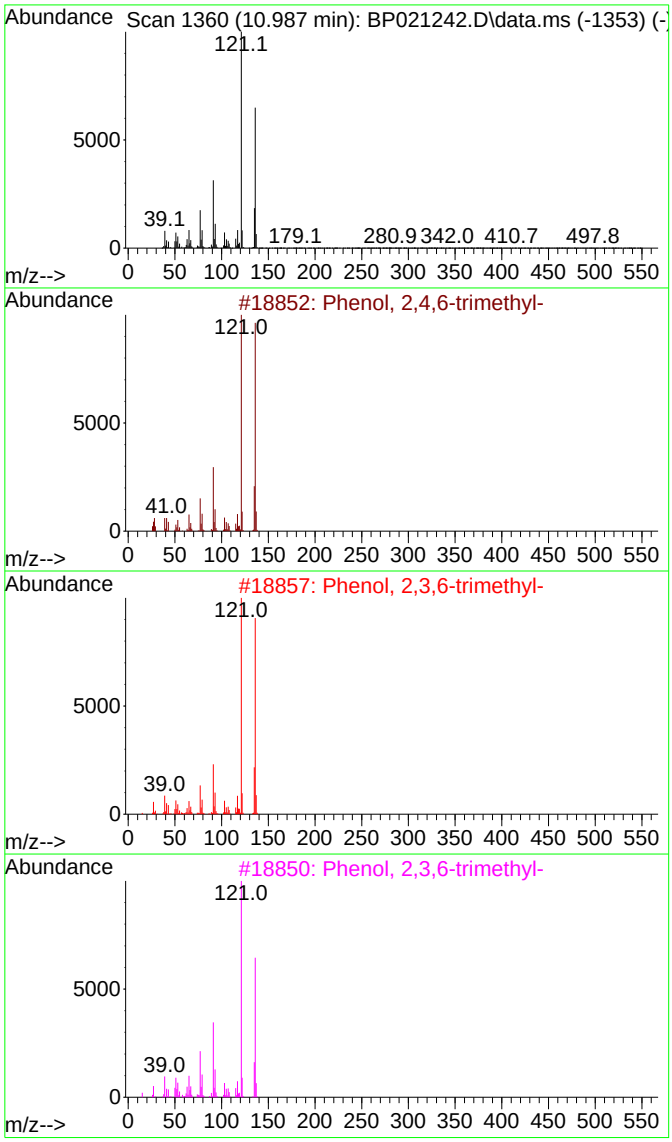
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

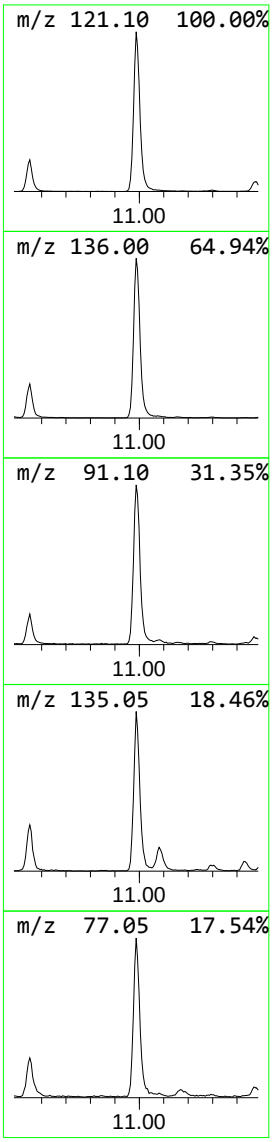
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Phenol, 2,4,6-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.987	21.96 ng/ul	1668680	Naphthalene-d8	10.393

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	97
2	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	97
3	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	96
4	Phenol, 2,3,5-trimethyl-	136	C9H12O	000697-82-5	95
5	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	95





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Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

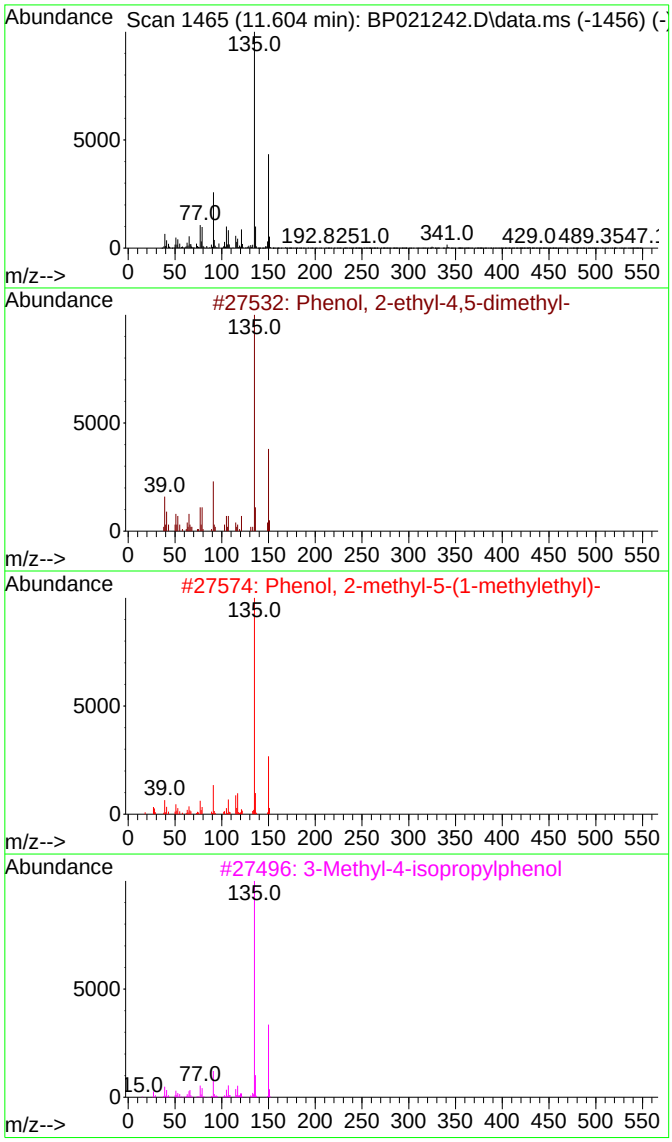
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

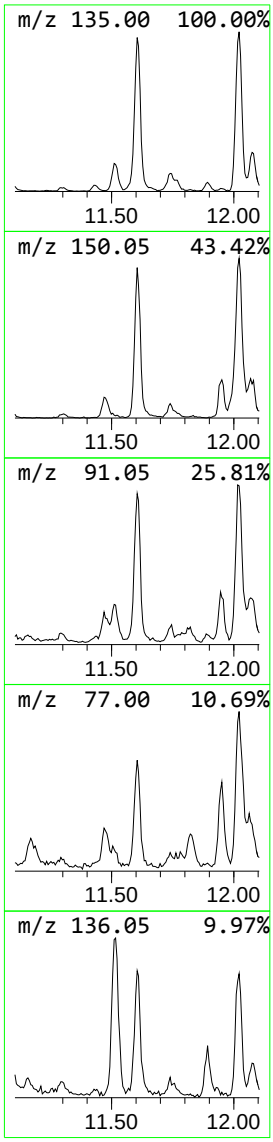
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Phenol, 2-ethyl-4,5-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.605	5.24 ng/ul	397955	Naphthalene-d8	10.393

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-ethyl-4,5-dimethyl-	150	C10H14O	002219-78-5	91
2	Phenol, 2-methyl-5-(1-methylethyl)-	150	C10H14O	000499-75-2	90
3	3-Methyl-4-isopropylphenol	150	C10H14O	003228-02-2	87
4	5-Cyano-1,2,3,4-tetrahydro-4,6-d...	150	C8H10N2O	027036-93-7	86
5	Thymol	150	C10H14O	000089-83-8	86





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Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

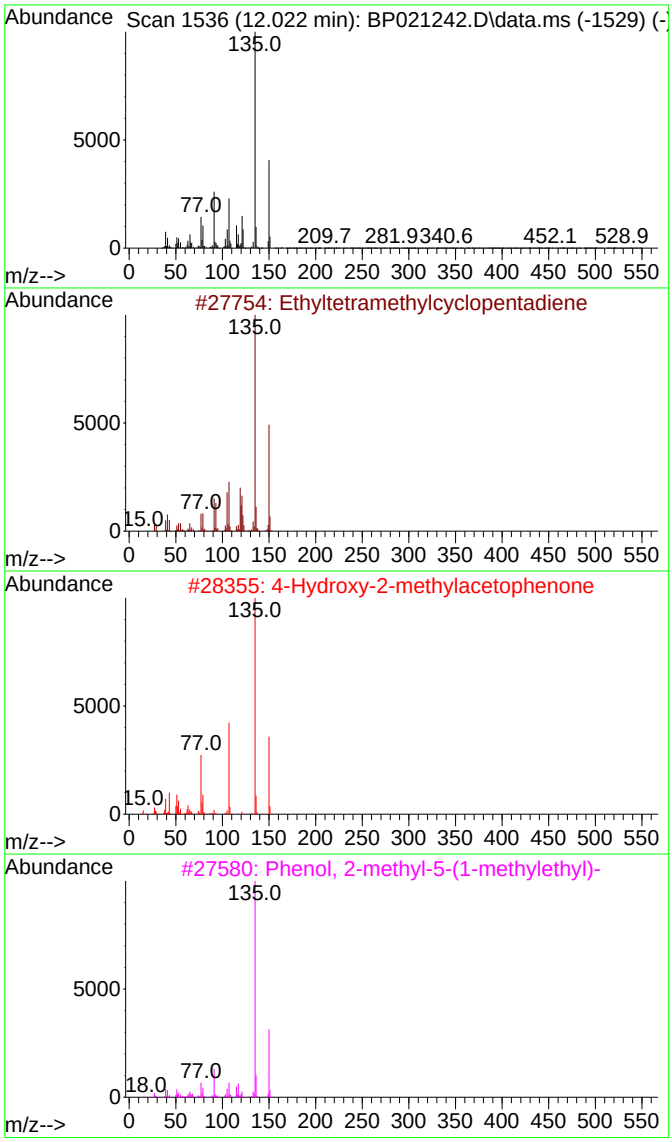
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

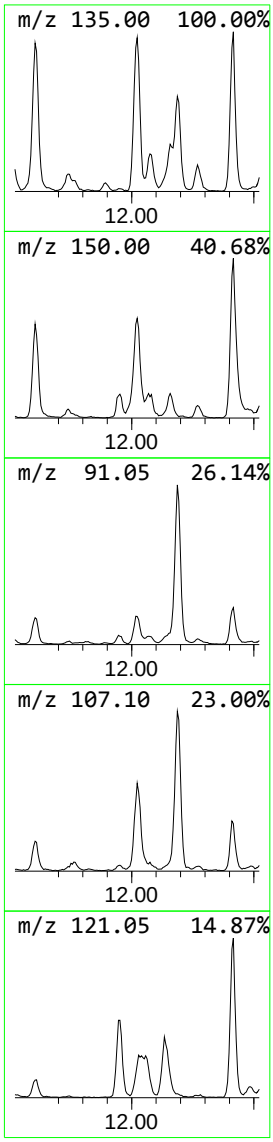
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Ethyltetramethylcyclopentad... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.022	6.54 ng/ul	497244	Naphthalene-d8	10.393

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyltetramethylcyclopentadiene	150	C11H18	057693-77-3	87
2	4-Hydroxy-2-methylacetophenone	150	C9H10O2	000875-59-2	81
3	Phenol, 2-methyl-5-(1-methylethyl)-	150	C10H14O	000499-75-2	76
4	3-Methyl-4-isopropylphenol	150	C10H14O	003228-02-2	76
5	Phenol, 2-methyl-5-(1-methylethyl)-	150	C10H14O	000499-75-2	76





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Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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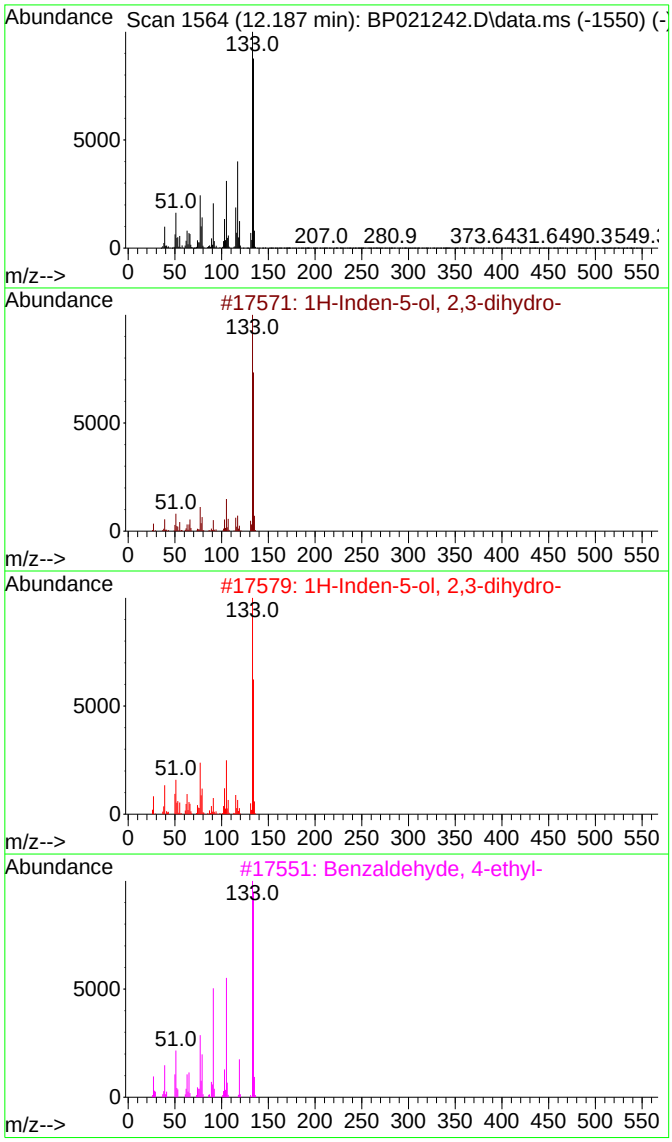
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

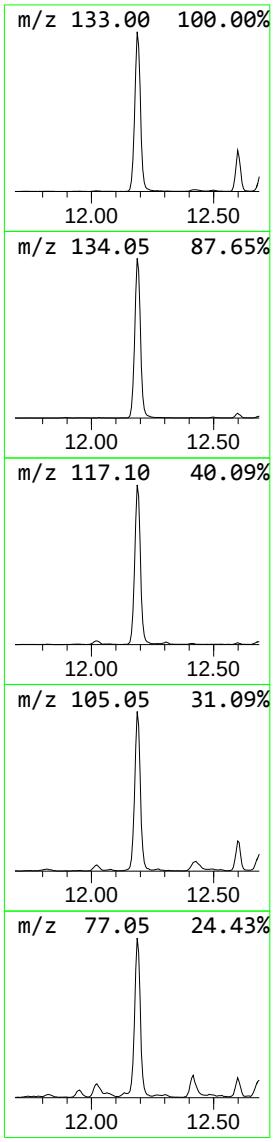
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 1H-Inden-5-ol, 2,3-dihydro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.187	57.19 ng/ul	4345260	Naphthalene-d8	10.393

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Inden-5-ol, 2,3-dihydro-	134	C9H10O	001470-94-6	87
2	1H-Inden-5-ol, 2,3-dihydro-	134	C9H10O	001470-94-6	81
3	Benzaldehyde, 4-ethyl-	134	C9H10O	004748-78-1	78
4	Benzaldehyde, 3,4-dimethyl-	134	C9H10O	005973-71-7	76
5	Benzaldehyde, 3,4-dimethyl-	134	C9H10O	005973-71-7	68





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Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

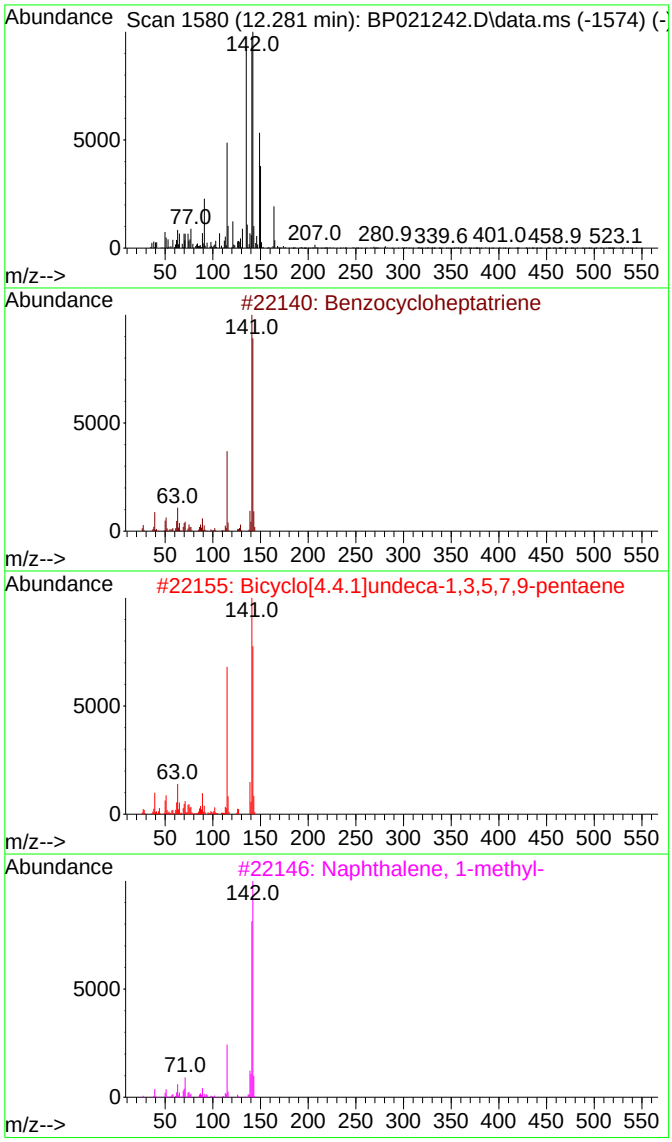
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

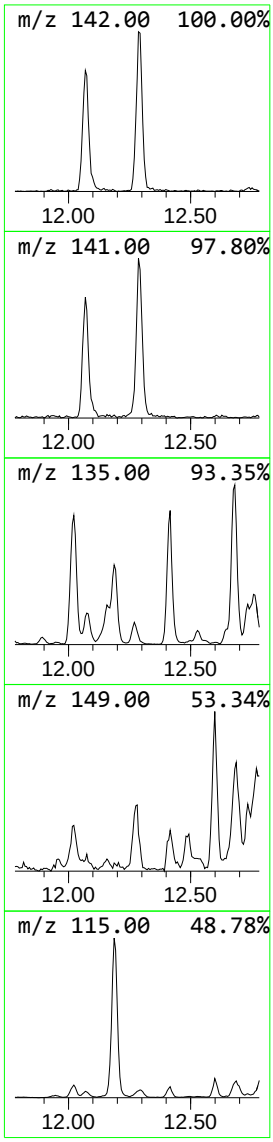
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzocycloheptatriene Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.281	2.97 ng/ul	225383	Naphthalene-d8	10.393

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzocycloheptatriene	142	C11H10	000264-09-5	50
2	Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	50
3	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	45
4	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	45
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	45





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Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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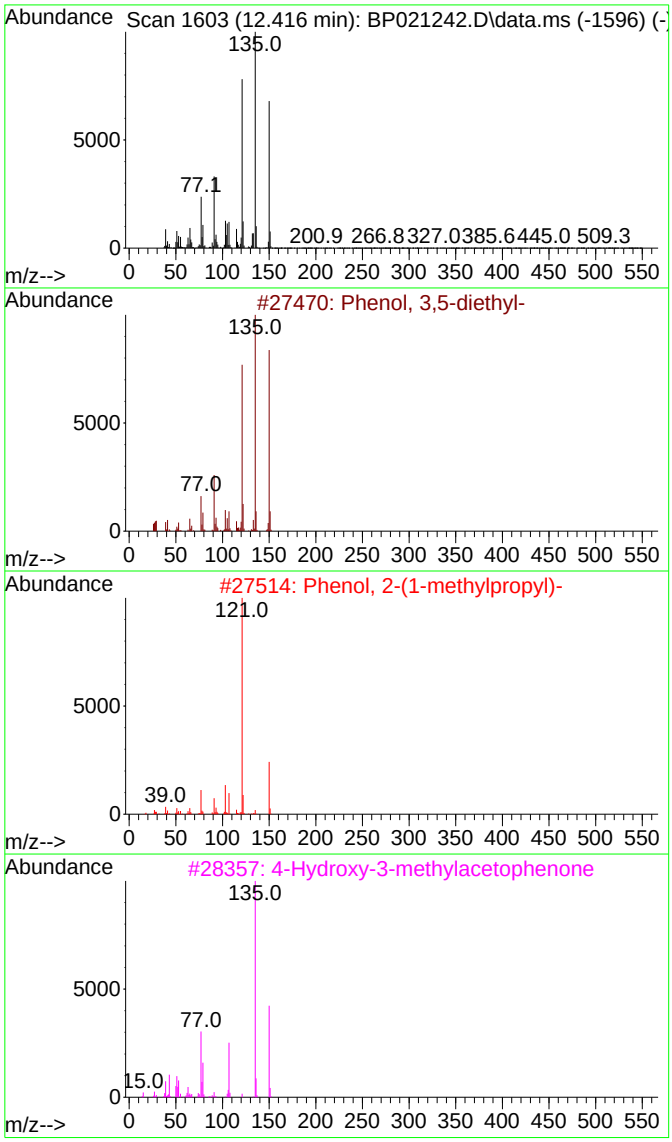
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

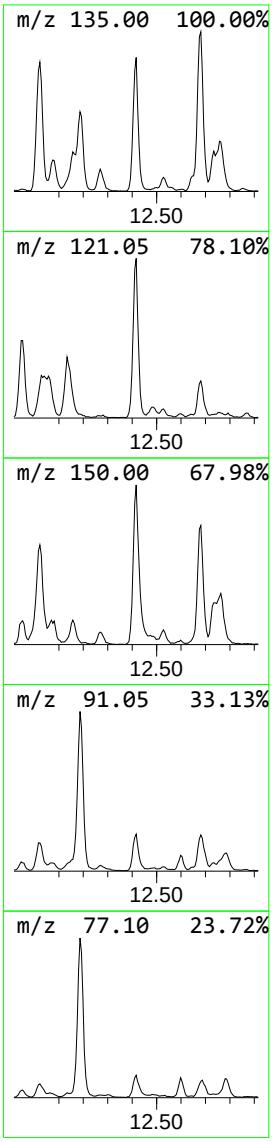
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Phenol, 3,5-diethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.416	5.68 ng/ul	651224	Acenaphthene-d10	14.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-diethyl-	150	C10H14O	001197-34-8	96
2	Phenol, 2-(1-methylpropyl)-	150	C10H14O	000089-72-5	64
3	4-Hydroxy-3-methylacetophenone	150	C9H10O2	000876-02-8	60
4	4-Hydroxy-3-methylacetophenone	150	C9H10O2	000876-02-8	60
5	3-Ethylbenzoic acid	150	C9H10O2	000619-20-5	58





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Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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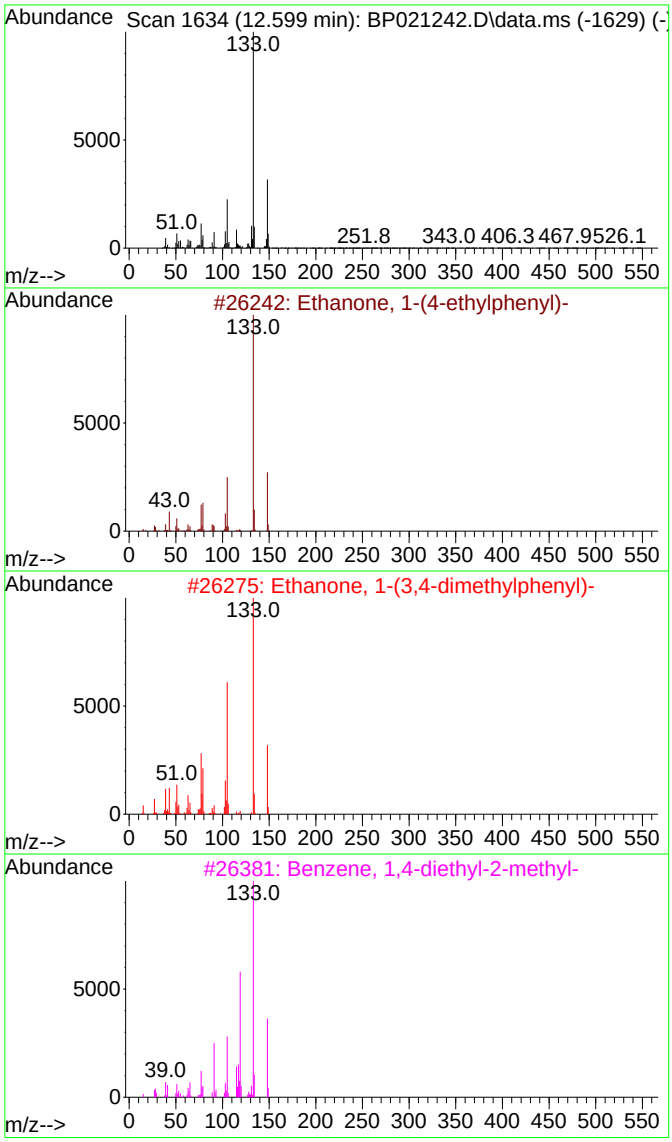
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

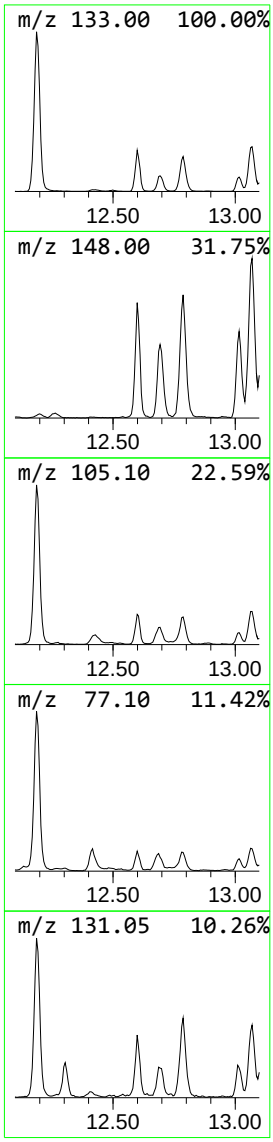
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Ethanone, 1-(4-ethylphenyl)- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.599	4.53 ng/ul	519647	Acenaphthene-d10	14.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 1-(4-ethylphenyl)-	148	C10H12O	000937-30-4	87
2	Ethanone, 1-(3,4-dimethylphenyl)-	148	C10H12O	003637-01-2	87
3	Benzene, 1,4-diethyl-2-methyl-	148	C11H16	013632-94-5	86
4	Ethanone, 1-(4-ethylphenyl)-	148	C10H12O	000937-30-4	83
5	m-Ethylacetophenone	148	C10H12O	022699-70-3	80





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Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

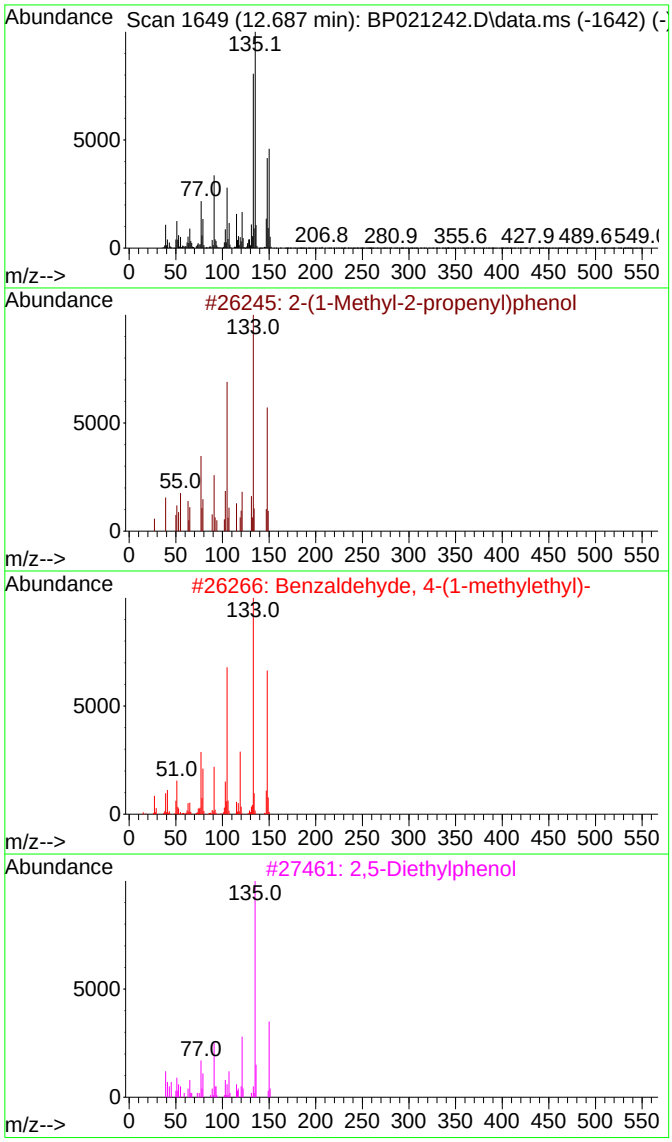
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

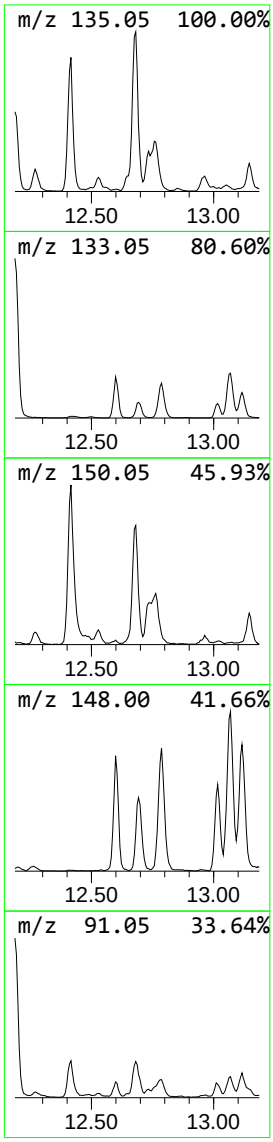
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 2-(1-Methyl-2-propenyl)phenol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.687	6.08 ng/ul	697129	Acenaphthene-d10	14.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(1-Methyl-2-propenyl)phenol	148	C10H12O	1000432-85-2	87
2	Benzaldehyde, 4-(1-methylethyl)-	148	C10H12O	000122-03-2	64
3	2,5-Diethylphenol	150	C10H14O	000876-20-0	52
4	1-Hexene, 2-(O-anisyl)-4-methyl-	204	C14H20O	1000149-68-0	50
5	Ethanone, 1-(2,5-dimethylphenyl)-	148	C10H12O	002142-73-6	46





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ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
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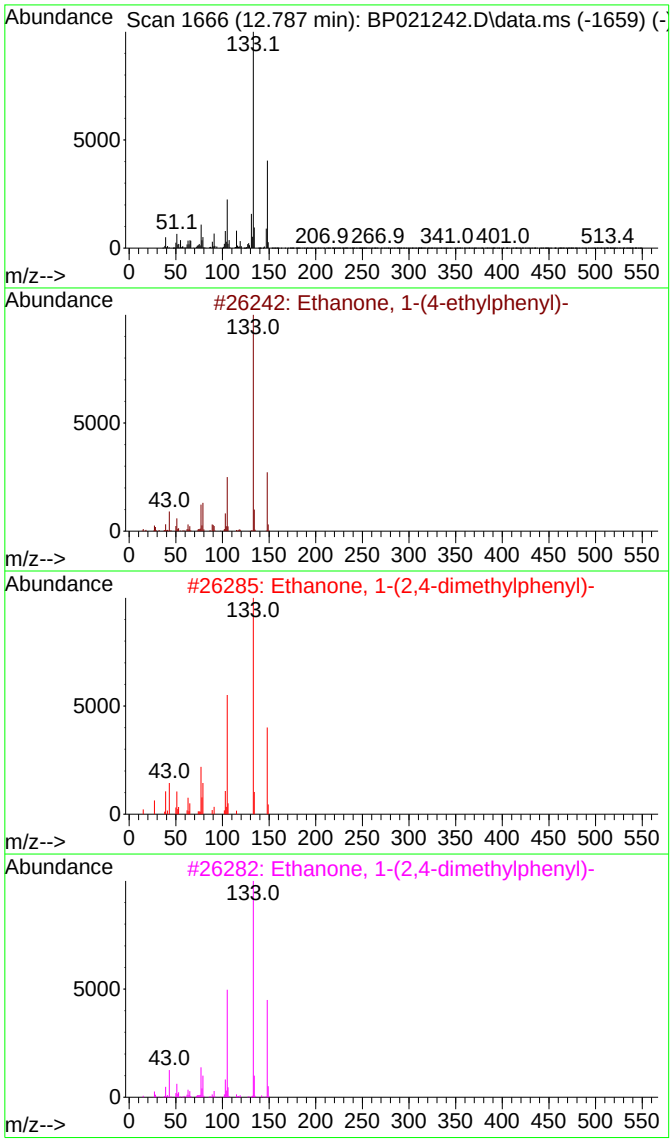
Quant Title : SVOA CALIBRATION

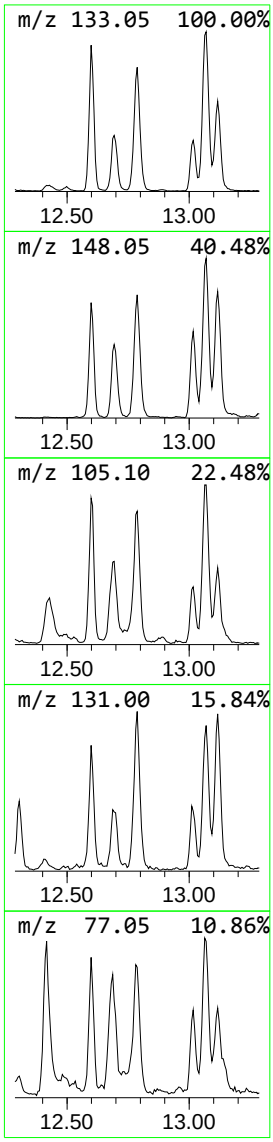
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Ethanone, 1-(2,4-dimethylph... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.787	6.84 ng/ul	784315	Acenaphthene-d10	14.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 1-(4-ethylphenyl)-	148	C10H12O	000937-30-4	87
2	Ethanone, 1-(2,4-dimethylphenyl)-	148	C10H12O	000089-74-7	72
3	Ethanone, 1-(2,4-dimethylphenyl)-	148	C10H12O	000089-74-7	64
4	Benzene, pentamethyl-	148	C11H16	000700-12-9	64
5	Ethanone, 1-(2,5-dimethylphenyl)-	148	C10H12O	002142-73-6	62





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Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

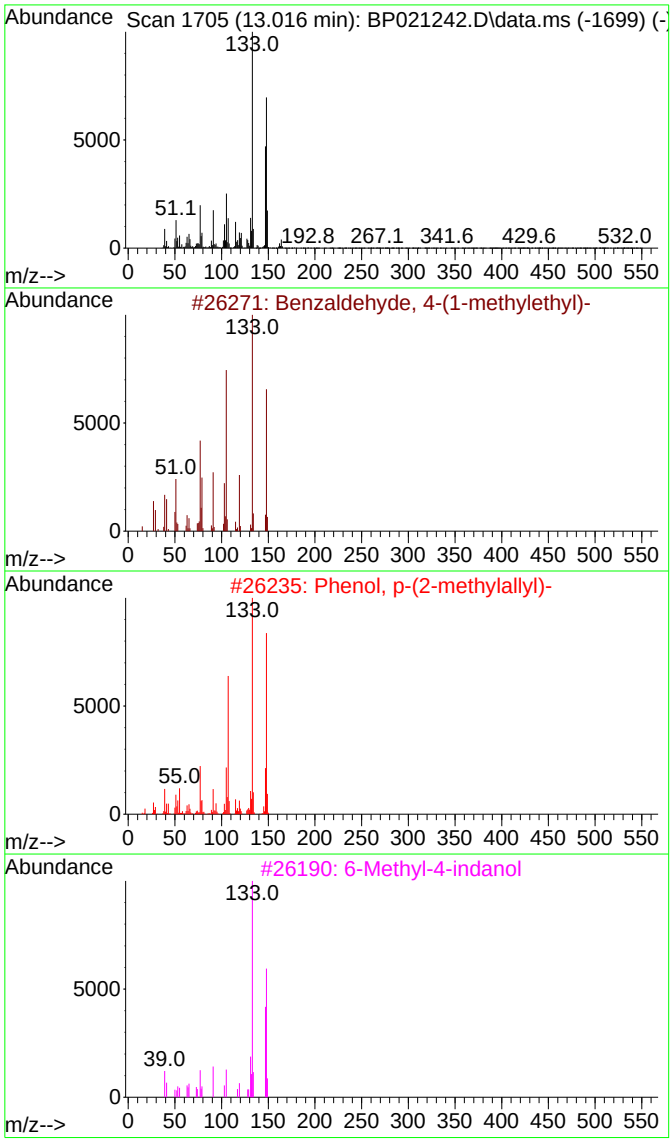
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

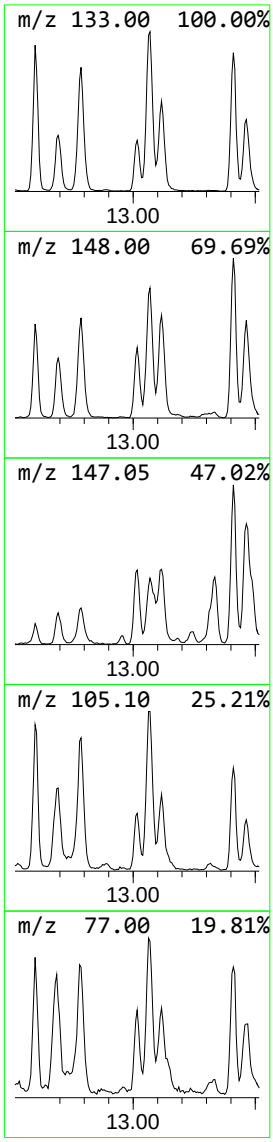
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzaldehyde, 4-(1-methylet... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.016	3.35 ng/ul	383918	Acenaphthene-d10	14.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 4-(1-methylethyl)-	148	C10H12O	000122-03-2	70
2	Phenol, p-(2-methylallyl)-	148	C10H12O	033641-78-0	68
3	6-Methyl-4-indanol	148	C10H12O	020294-32-0	64
4	2-(1-Methyl-2-propenyl)phenol	148	C10H12O	1000432-85-2	59
5	Benzene, 1-(1,1-dimethylethyl)-3...	148	C11H16	001075-38-3	53





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Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

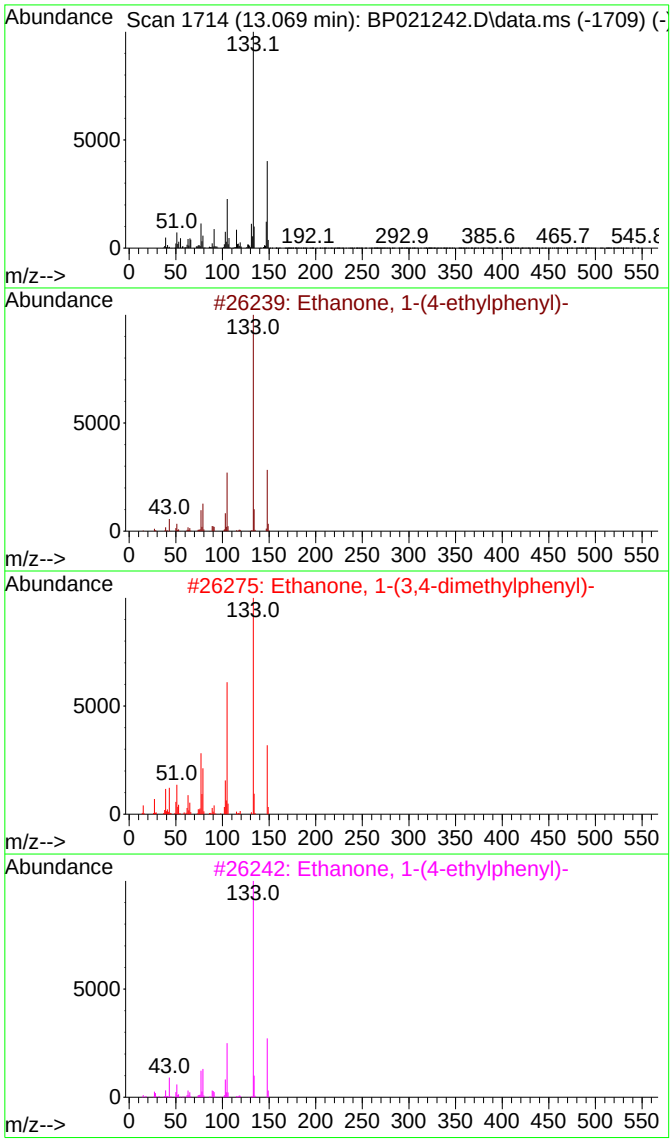
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

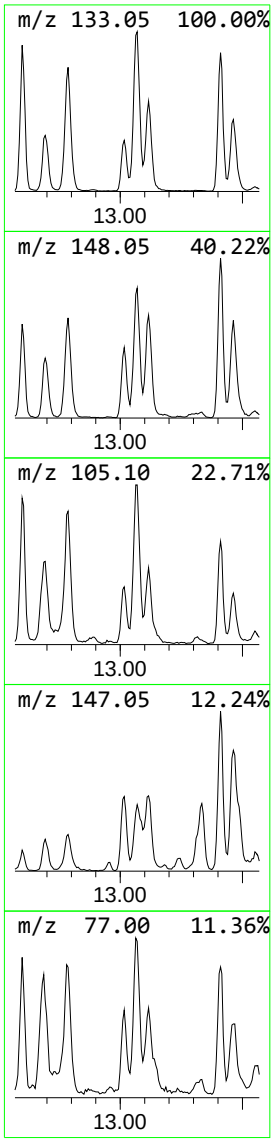
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Ethanone, 1-(3,4-dimethylph... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.069	6.81 ng/ul	780765	Acenaphthene-d10	14.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 1-(4-ethylphenyl)-	148	C10H12O	000937-30-4	87
2	Ethanone, 1-(3,4-dimethylphenyl)-	148	C10H12O	003637-01-2	87
3	Ethanone, 1-(4-ethylphenyl)-	148	C10H12O	000937-30-4	87
4	Ethanone, 1-(4-ethylphenyl)-	148	C10H12O	000937-30-4	83
5	5H-Cyclopentapyrazine, 6,7-dihyd...	148	C9H12N2	038917-61-2	72





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Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

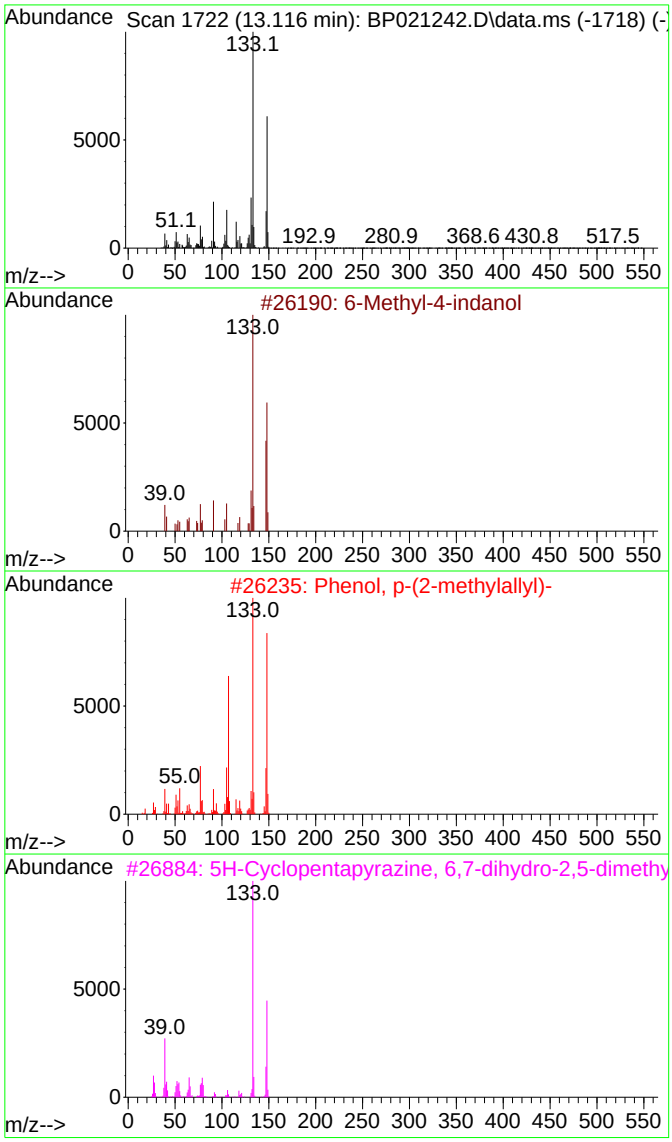
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

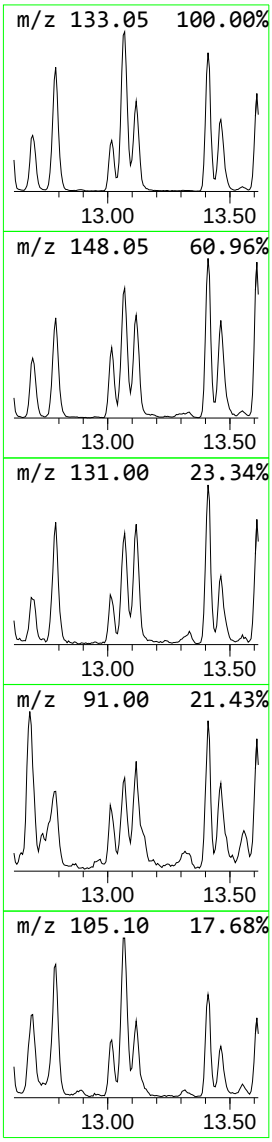
****

Peak Number 19 6-Methyl-4-indanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.116	6.46 ng/ul	740775	Acenaphthene-d10	14.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methyl-4-indanol	148	C10H12O	020294-32-0	93
2	Phenol, p-(2-methylallyl)-	148	C10H12O	033641-78-0	87
3	5H-Cyclopentapyrazine, 6,7-dihyd...	148	C9H12N2	038917-61-2	72
4	Ethanone, 1-(2,4-dimethylphenyl)-	148	C10H12O	000089-74-7	70
5	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	68





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Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

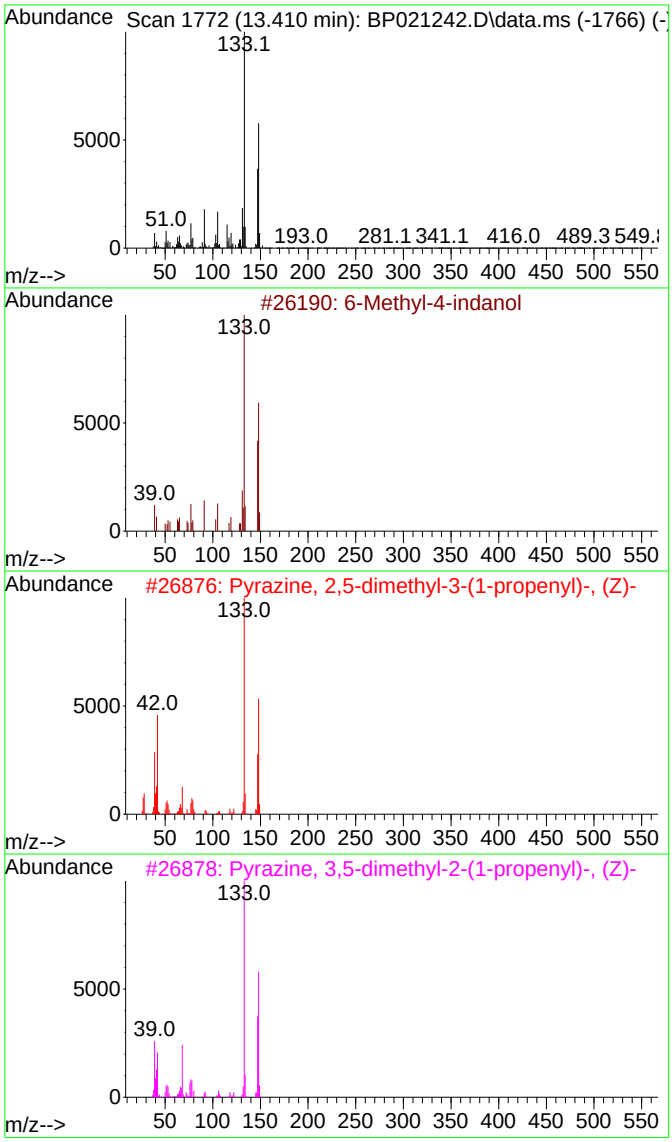
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

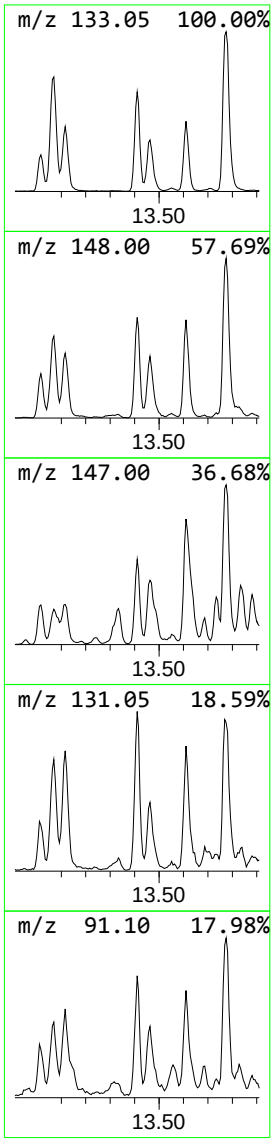
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Pyrazine, 3,5-dimethyl-2-(1... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.410	6.71 ng/ul	768818	Acenaphthene-d10	14.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methyl-4-indanol	148	C10H12O	020294-32-0	94
2	Pyrazine, 2,5-dimethyl-3-(1-prop...	148	C9H12N2	055138-73-3	80
3	Pyrazine, 3,5-dimethyl-2-(1-prop...	148	C9H12N2	055138-74-4	80
4	2-(1-Methyl-2-propenyl)phenol	148	C10H12O	1000432-85-2	80
5	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	70





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Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

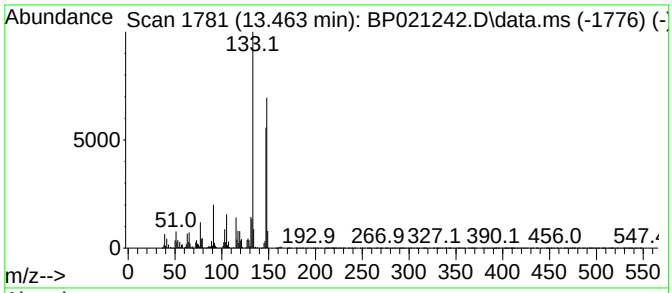
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

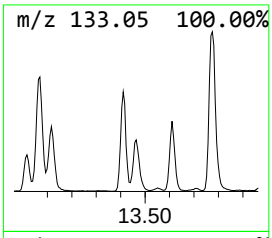
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

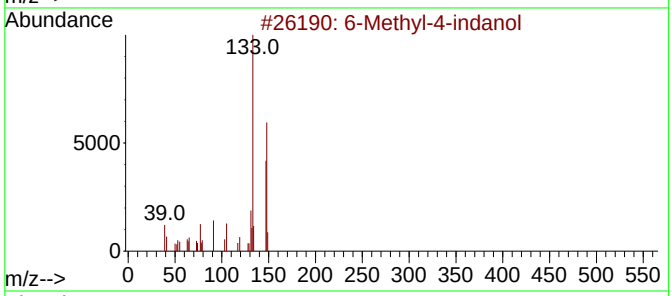
****
Peak Number 21 5H-Cyclopentapyrazine, 6,7-... Concentration Rank 15

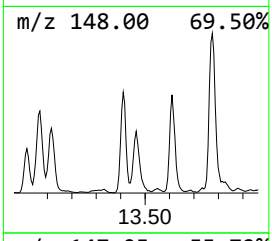
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.463	4.80 ng/ul	550824	Acenaphthene-d10	14.275

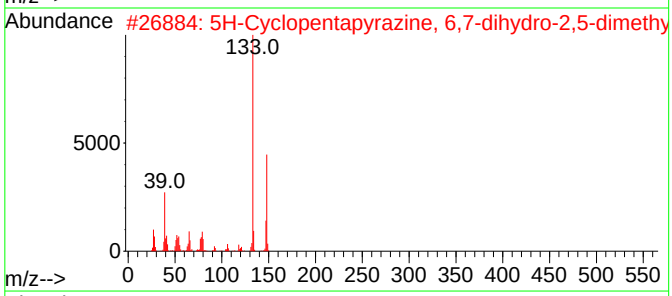
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methyl-4-indanol	148	C10H12O	020294-32-0	80
2	5H-Cyclopentapyrazine, 6,7-dihyd...	148	C9H12N2	038917-61-2	64
3	Benzene, pentamethyl-	148	C11H16	000700-12-9	58
4	Phenol, 2-(2-methyl-2-propenyl)-	148	C10H12O	020944-88-1	58
5	Ethanone, 1-(2,4-dimethylphenyl)-	148	C10H12O	000089-74-7	53

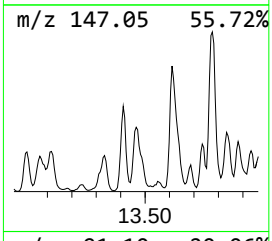


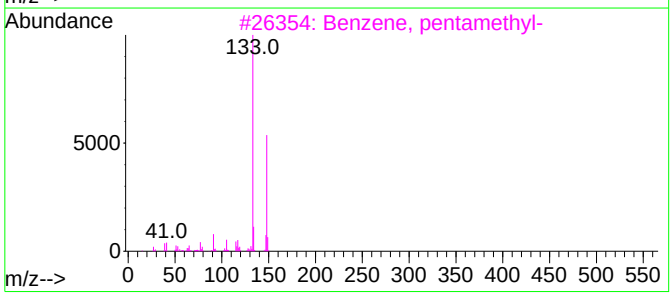


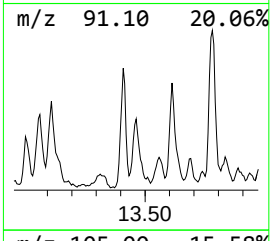


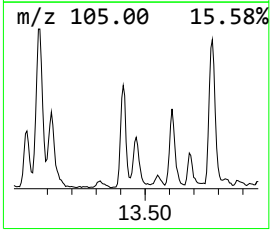












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Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

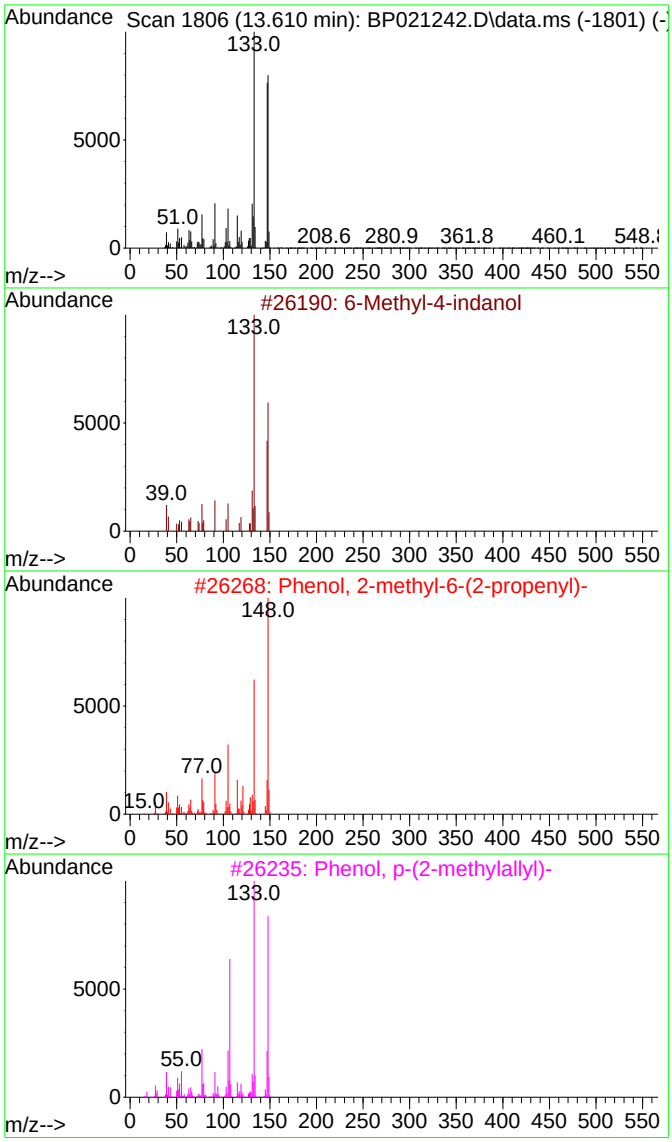
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

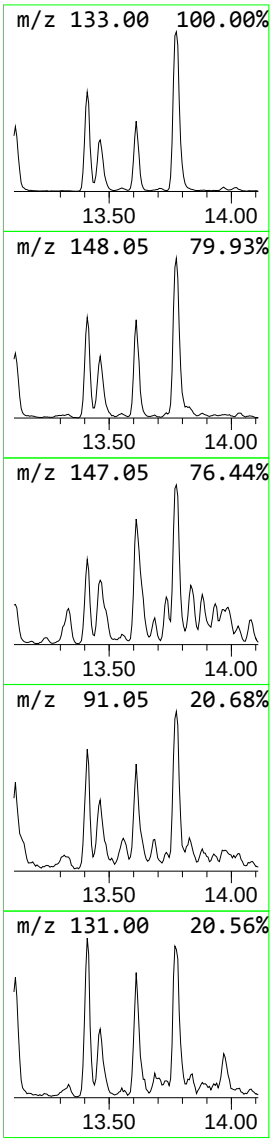
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 Phenol, 2-methyl-6-(2-prope... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.610	5.99 ng/ul	687007	Acenaphthene-d10	14.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methyl-4-indanol	148	C10H12O	020294-32-0	90
2	Phenol, 2-methyl-6-(2-propenyl)-	148	C10H12O	003354-58-3	58
3	Phenol, p-(2-methylallyl)-	148	C10H12O	033641-78-0	58
4	Pyrrole-3-carbonitrile, 5-formyl...	148	C8H8N2O	032487-71-1	55
5	Benzaldehyde, 2,4,5-trimethyl-	148	C10H12O	005779-72-6	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

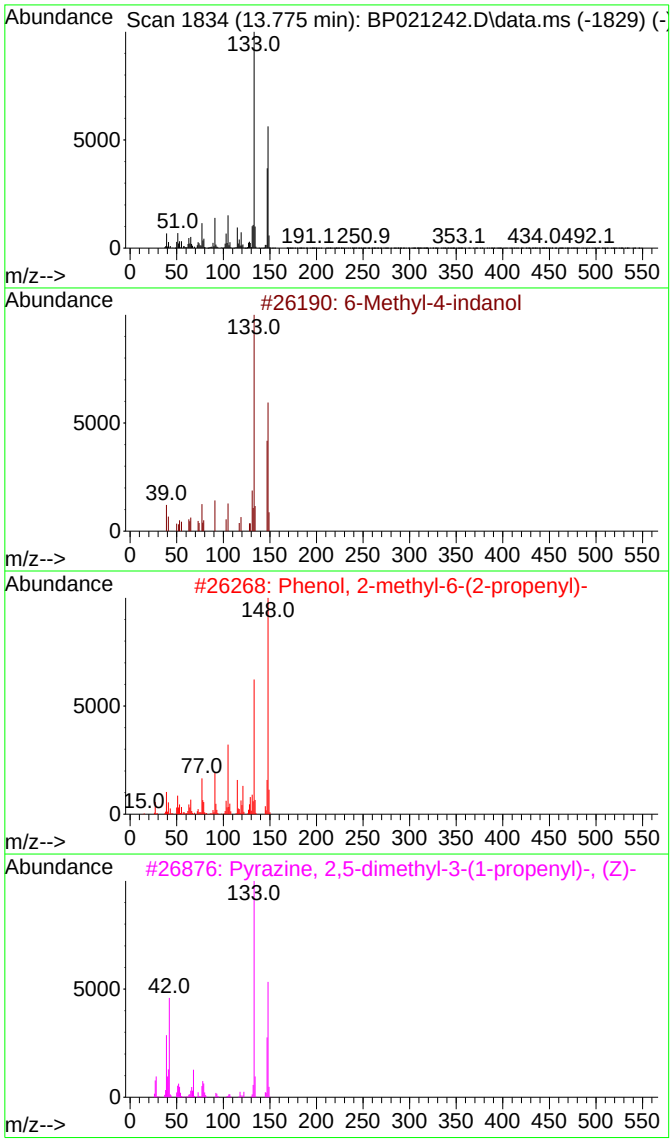
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

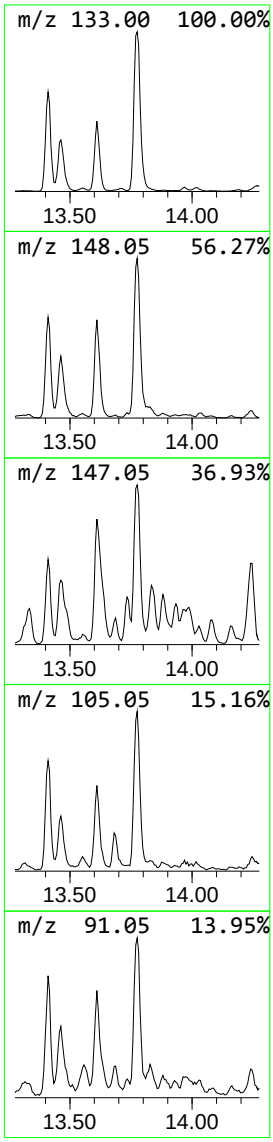
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Pyrazine, 2,5-dimethyl-3-(1... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.775	11.09 ng/ul	1271330	Acenaphthene-d10	14.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methyl-4-indanol	148	C10H12O	020294-32-0	87
2	Phenol, 2-methyl-6-(2-propenyl)-	148	C10H12O	003354-58-3	83
3	Pyrazine, 2,5-dimethyl-3-(1-prop...	148	C9H12N2	055138-73-3	80
4	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	76
5	Benzene, pentamethyl-	148	C11H16	000700-12-9	74





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Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

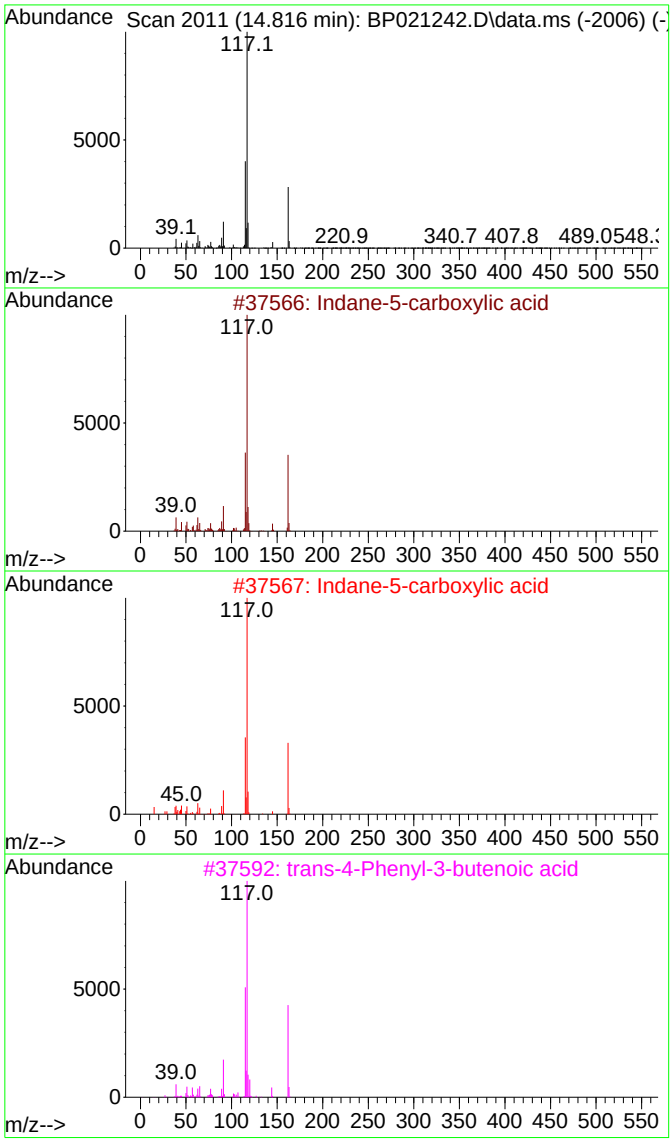
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

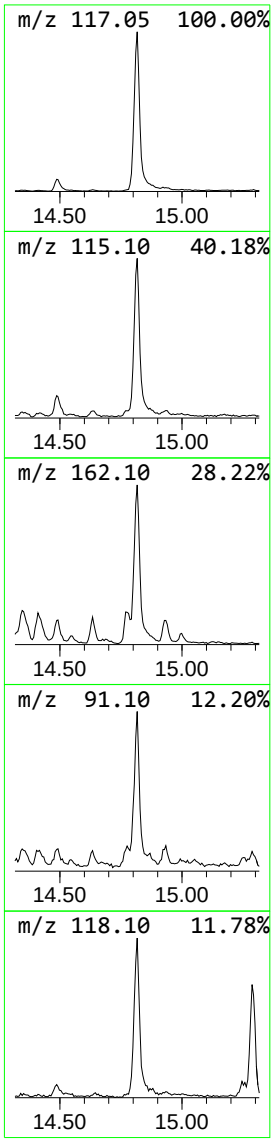
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Indane-5-carboxylic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.816	6.25 ng/ul	716197	Acenaphthene-d10	14.275

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane-5-carboxylic acid	162	C10H10O2	065898-38-6	94
2	Indane-5-carboxylic acid	162	C10H10O2	065898-38-6	91
3	trans-4-Phenyl-3-butenoic acid	162	C10H10O2	001914-58-5	76
4	Benzene, 1-pentenyl-	146	C11H14	000826-18-6	64
5	Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	64





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Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

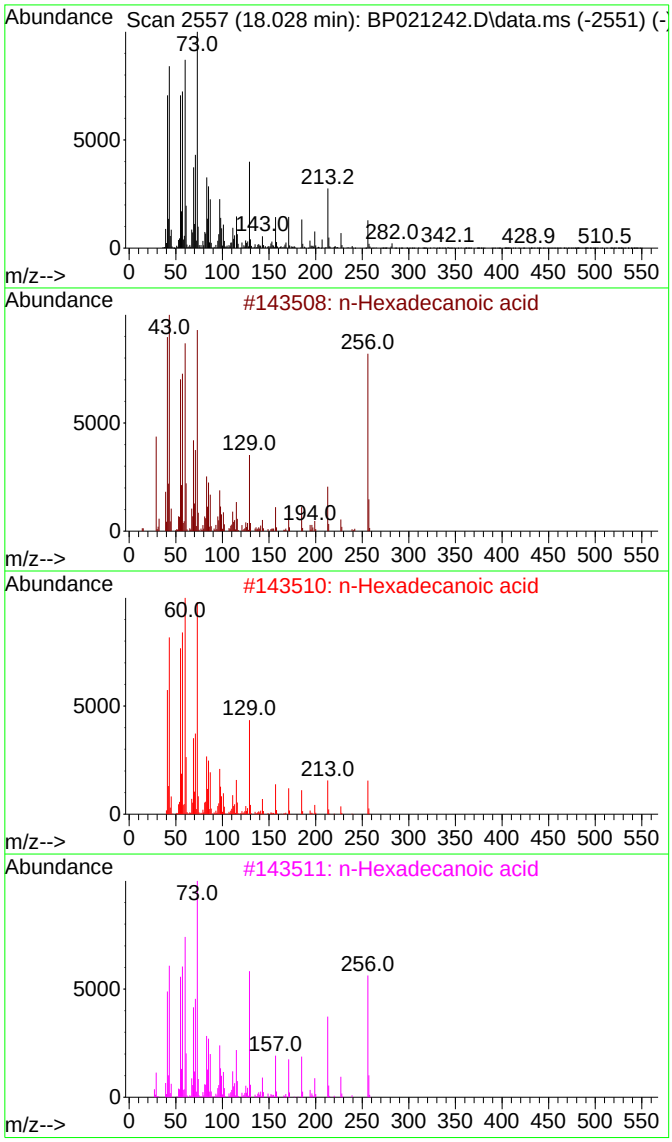
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

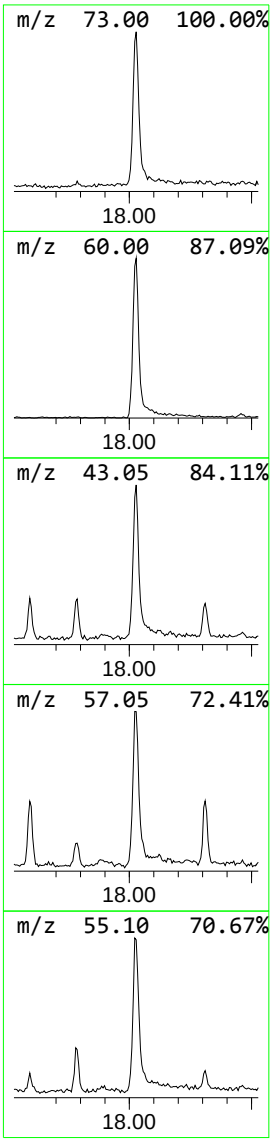
****

Peak Number 25 n-Hexadecanoic acid Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.028	3.98 ng/ul	508637	Phenanthrene-d10	17.092

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	93





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021242.D
Acq On : 30 Jul 2024 19:58
Operator : MA/JU
Sample : P3347-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.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
Benzene, 1,2,3-...	7.728	2.9	ng/ul	148797	1	7.622	1043690	20.0
Indane	7.981	73.6	ng/ul	3840920	1	7.622	1043690	20.0
2,3,5-Trimethyl...	8.281	2.3	ng/ul	121510	1	7.622	1043690	20.0
p-Aminotoluene	8.528	2.2	ng/ul	116278	1	7.622	1043690	20.0
Benzofuran, 7-m...	8.963	3.6	ng/ul	189595	1	7.622	1043690	20.0
Benzofuran, 2-m...	9.087	3.3	ng/ul	251224	2	10.393	1519690	20.0
1H-Indene, 2,3-...	9.781	2.4	ng/ul	185069	2	10.393	1519690	20.0
Phenol, 2,4,6-t...	10.987	22.0	ng/ul	1668680	2	10.393	1519690	20.0
Phenol, 2-ethyl...	11.605	5.2	ng/ul	397955	2	10.393	1519690	20.0
Ethyltetramethy...	12.022	6.5	ng/ul	497244	2	10.393	1519690	20.0
1H-Inden-5-ol, ...	12.187	57.2	ng/ul	4345260	2	10.393	1519690	20.0
Benzocyclohepta...	12.281	3.0	ng/ul	225383	2	10.393	1519690	20.0
Phenol, 3,5-die...	12.416	5.7	ng/ul	651224	3	14.275	2292850	20.0
Ethanone, 1-(4-...	12.599	4.5	ng/ul	519647	3	14.275	2292850	20.0
2-(1-Methyl-2-p...	12.687	6.1	ng/ul	697129	3	14.275	2292850	20.0
Ethanone, 1-(2,...	12.787	6.8	ng/ul	784315	3	14.275	2292850	20.0
Benzaldehyde, 4...	13.016	3.4	ng/ul	383918	3	14.275	2292850	20.0
Ethanone, 1-(3,...	13.069	6.8	ng/ul	780765	3	14.275	2292850	20.0
6-Methyl-4-indanol	13.116	6.5	ng/ul	740775	3	14.275	2292850	20.0
Pyrazine, 3,5-d...	13.410	6.7	ng/ul	768818	3	14.275	2292850	20.0
5H-Cyclopentapy...	13.463	4.8	ng/ul	550824	3	14.275	2292850	20.0
Phenol, 2-methy...	13.610	6.0	ng/ul	687007	3	14.275	2292850	20.0
Pyrazine, 2,5-d...	13.775	11.1	ng/ul	1271330	3	14.275	2292850	20.0
Indane-5-carbox...	14.816	6.3	ng/ul	716197	3	14.275	2292850	20.0
n-Hexadecanoic ...	18.028	4.0	ng/ul	508637	4	17.092	2552830	20.0