LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

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-BP092424.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP022153.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.381	400	407	417	rBV	52488	109931	1.32%	0.340%
2	5.746	463	469	476	rBV	55276	89569	1.08%	0.277%
3	6.222	540	550	558	rVB4	21415	47650	0.57%	0.148%
4	6.699	624	631	637	rBV	36047	56145	0.67%	0.174%
5	6.805	642	649	654	rVV	162149	252714	3.04%	0.783%
6	6.863	654	659	666	rVV	131907	208654	2.51%	0.646%
7	6.934	666	671	676	rVV	123892	189070	2.27%	0.586%
8	6.981	676	679	684	rVV2	22747	32943	0.40%	0.102%
9	7.099	692	699	701	rVV	109652	180341	2.17%	0.559%
10	7.134	701	705	715	rVV	253138	386823	4.65%	1.198%
11	7.228	715	721	731	rVV2	49158	96137	1.15%	0.298%
12	7.352	734	742	748	rVV	287180	478622	5.75%	1.482%
13	7.699	791	801	806	rBV	439023	714989	8.59%	2.215%
14	7.763	806	812	817	rVV	379550	598235	7.18%	1.853%
15	7.816	817	821	832	rVV	142217	241338	2.90%	0.747%
16	7.916	832	838	843	rVV	84115	127267	1.53%	0.394%
17	8.069	857	864	868	rBV	46842	78375	0.94%	0.243%
18	8.122	868	873	877	rVV	86141	147515	1.77%	0.457%
19	8.169	877	881	888	rVB2	155838	236317	2.84%	0.732%
20	8.240	888	893	897	rBV	27088	38558	0.46%	0.119%
21	8.328	902	908	913	rBV	58504	94944	1.14%	0.294%
22	8.487	928	935	945	rVB3	46564	83776	1.01%	0.259%
23	8.634	954	960	964	rBV	44250	72491	0.87%	0.225%
24	8.681	964	968	975	rVB2	65057	117326	1.41%	0.363%
25	8.787	980	986	992	rBV3	56483	104484	1.25%	0.324%
26	8.969	1012	1017	1021	rBV2	23662	38797	0.47%	0.120%
27	9.110	1035	1041	1048	rVB2	55782	99471	1.19%	0.308%
28	9.263	1062	1067	1072	rBV2	104546	164843	1.98%	0.511%
29	9.363	1078	1084	1095	rVB2	37297	85825	1.03%	0.266%
30	9.469	1095	1102	1108	rVB	81432	124466	1.49%	0.386%
31	9.616	1122	1127	1132	rVB	63407	95265	1.14%	0.295%
32	9.716	1139	1144	1152	rVB2	34301	57315	0.69%	0.178%
33	9.869	1163	1170	1174	rBV2	35053	70946	0.85%	0.220%
34	9.969	1182	1187	1194	rVB2	24954	41413	0.50%	0.128%
35	10.093	1203	1208	1214	rVB3	38328	62319	0.75%	0.193%
36	10.151	1214	1218	1225	rVB2	40464	61831	0.74%	0.192%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

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-BP092424.MA.M
Title : SVOA CALIBRATION

37	10.222	1225	1230	1238	rVB3	18796	34111	0.41%	0.106%
38	10.451	1262	1269	1275	rVV2	816758	1433149	17.21%	4.439%
39	10.504	1275	1278	1285	rVB	90389	149355	1.79%	0.463%
40	10.587	1285	1292	1298	rBV4	123724	284856	3.42%	0.882%
41	10.640	1298	1301	1311	rVB4	34338	78012	0.94%	0.242%
42	10.728	1311	1316	1320	rBV3	41630	79011	0.95%	0.245%
43	10.828	1327	1333	1341	rVB	262393	392358	4.71%	1.215%
44	10.957	1347	1355	1359	rBV5	31118	71329	0.86%	0.221%
45	11.034	1364	1368	1378	rVB7	21053	60007	0.72%	0.186%
46	11.245	1400	1404	1409	rVV4	24395	48115	0.58%	0.149%
47	11.398	1425	1430	1439	rVB3	65657	120368	1.45%	0.373%
48	11.551	1451	1456	1461	rVB2	24684	36870	0.44%	0.114%
49	11.716	1475	1484	1495	rVB4	100328	214576	2.58%	0.665%
50	11.869	1505	1510	1513	rBV2	39663	55562	0.67%	0.172%
51	11.910	1513	1517	1523	rVB	107647	151857	1.82%	0.470%
52	11.998	1525	1532	1536	rBV5	52480	98658	1.18%	0.306%
53	12.051	1536	1541	1546	rVV3	39360	67353	0.81%	0.209%
54	12.110	1546	1551	1557	rVB3	78191	135877	1.63%	0.421%
55	12.216	1563	1569	1574	rBV	43724	77638	0.93%	0.240%
56	12.281	1574	1580	1585	rVV4	68891	131755	1.58%	0.408%
57	12.334	1585	1589	1594	rVB	48814	75443	0.91%	0.234%
58	12.469	1608	1612	1615	rBV3	26821	42283	0.51%	0.131%
59	12.716	1645	1654	1659	rBV4	52519	111186	1.34%	0.344%
60	12.816	1667	1671	1676	rVB5	21582	34938	0.42%	0.108%
61	12.875	1676	1681	1685	rVB	63209	88134	1.06%	0.273%
62	13.051	1703	1711	1718	rBV	78146	128504	1.54%	0.398%
63	13.128	1720	1724	1730	rVB5	19221	36680	0.44%	0.114%
64	13.257	1741	1746	1751	rBV4	29183	49942	0.60%	0.155%
65	13.351	1756	1762	1767	rVB2	77741	129329	1.55%	0.401%
66	13.481	1773	1784	1791	rBV4	41606	112002	1.35%	0.347%
67	13.563	1794	1798	1803	rVV	57665	98273	1.18%	0.304%
68	13.616	1803	1807	1810	rVV2	28472	52738	0.63%	0.163%
69	13.657	1810	1814	1819	rVV	138760	226576	2.72%	0.702%
70	13.710	1819	1823	1828	rVB	53113	79023	0.95%	0.245%
71	13.992	1867	1871	1874	rVV	53854	89152	1.07%	0.276%
72	14.028	1874	1877	1883	rVV	82768	134416	1.61%	0.416%
73	14.304	1918	1924	1931	rVV	946880	1397904	16.79%	4.330%
74	14.369	1931	1935	1940	rVV2	33373	62712	0.75%	0.194%
75	14.416	1940	1943	1951	rVV5	31813	57723	0.69%	0.179%
76	14.522	1957	1961	1970	rVV9	28886	68972	0.83%	0.214%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DDC95DL

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-BP092424.MA.M
Title : SVOA CALIBRATION

77	14.622	1972	1978	1982	rVV	58526	113118	1.36%	0.350%
78	14.698	1982	1991	1996	rVV8	49453	139788	1.68%	0.433%
79	14.751	1996	2000	2007	rVV5	32234	52421	0.63%	0.162%
80	14.816	2007	2011	2014	rVV4	31067	47282	0.57%	0.146%
81	14.857	2014	2018	2022	rVV2	51285	86323	1.04%	0.267%
82	14.945	2022	2033	2049	rVV	1997862	3143481	37.75%	9.736%
83	15.069	2049	2054	2057	rVV5	20168	36199	0.43%	0.112%
84	15.110	2057	2061	2067	rVV3	31980	59749	0.72%	0.185%
85	15.181	2069	2073	2079	rVV6	21998	39159	0.47%	0.121%
86	15.310	2090	2095	2101	rVB	83094	126602	1.52%	0.392%
87	15.422	2107	2114	2118	rBV	74078	116864	1.40%	0.362%
88	15.463	2118	2121	2132	rVB3	46751	73667	0.88%	0.228%
89	16.151	2233	2238	2246	rBV5	25063	48190	0.58%	0.149%
90	16.592	2307	2313	2319	rVV3	21541	48299	0.58%	0.150%
91	16.745	2332	2339	2349	rBV	5880491	8326601	100.00%	25.790%
92	17.098	2393	2399	2406	rBV	1218164	1857105	22.30%	5.752%
93	17.204	2412	2417	2424	rBV2	84850	144479	1.74%	0.447%
94	19.516	2807	2810	2815	rVB3	26477	34751	0.42%	0.108%
95	19.574	2816	2820	2828	rVB	118269	175734	2.11%	0.544%
96	20.245	2931	2934	2939	rBV3	53736	58621	0.70%	0.182%
97	20.916	3044	3048	3053	rBV2	77015	106616	1.28%	0.330%
98	21.533	3148	3153	3161	rBV2	1599209	2379942	28.58%	7.371%
99	21.616	3163	3167	3173	rVB2	103168	160850	1.93%	0.498%
100	24.815	3703	3711	3725	rVB	920976	2426869	29.15%	7.517%

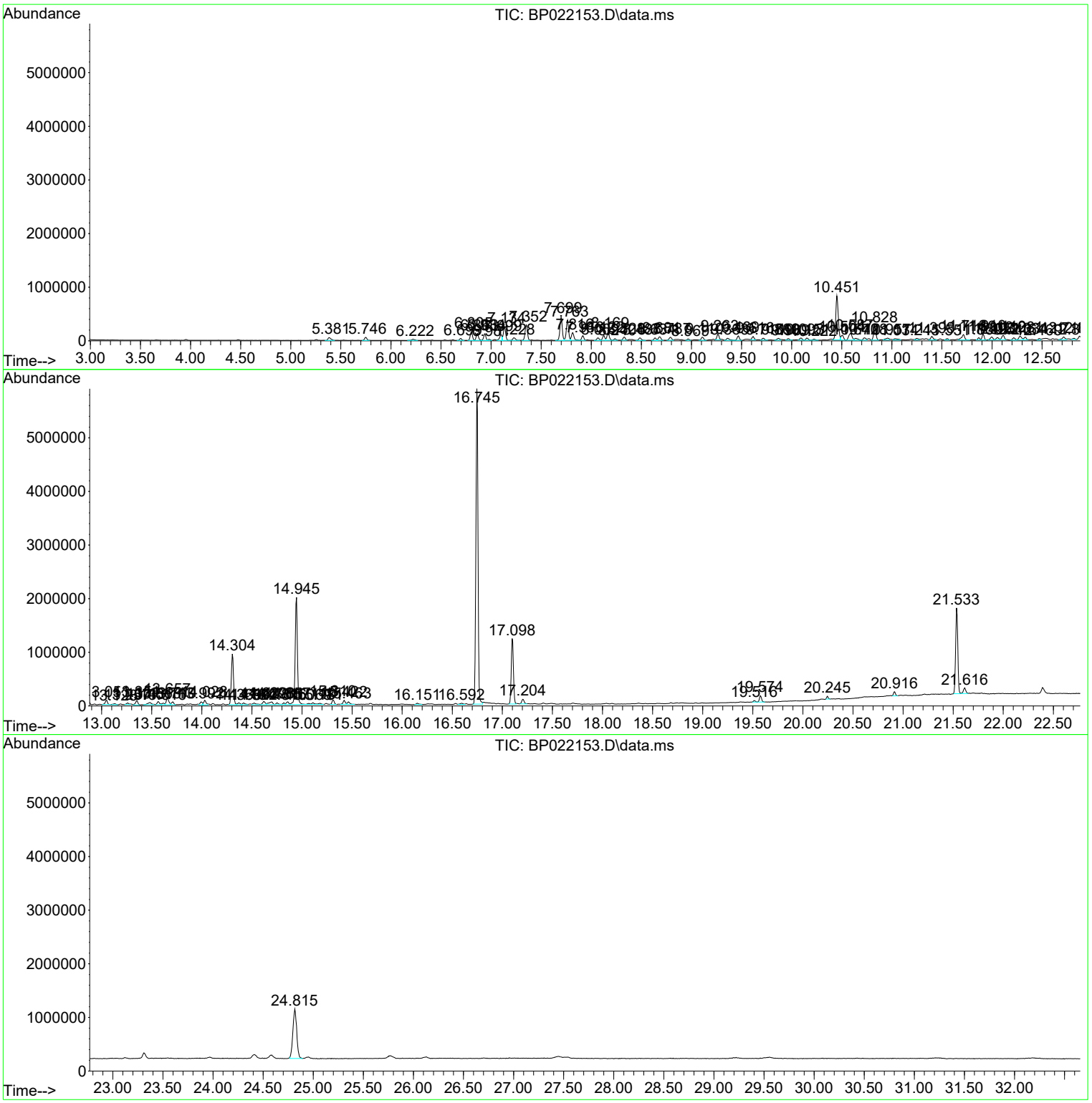
Sum of corrected areas: 32286162

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.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\BP100124\
Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

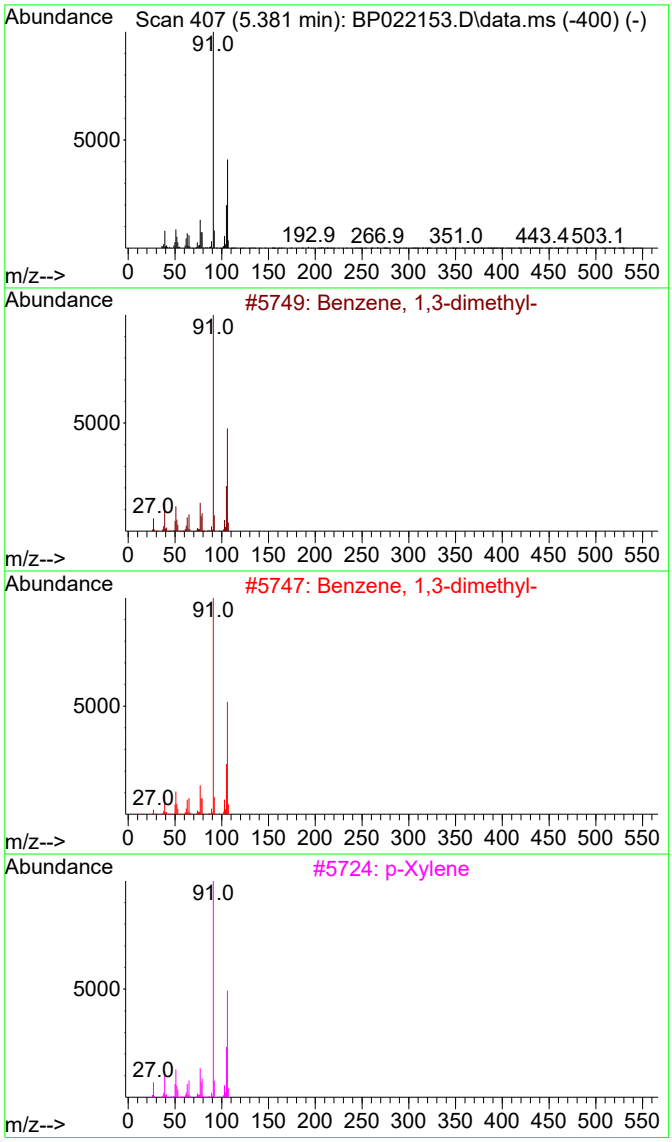
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

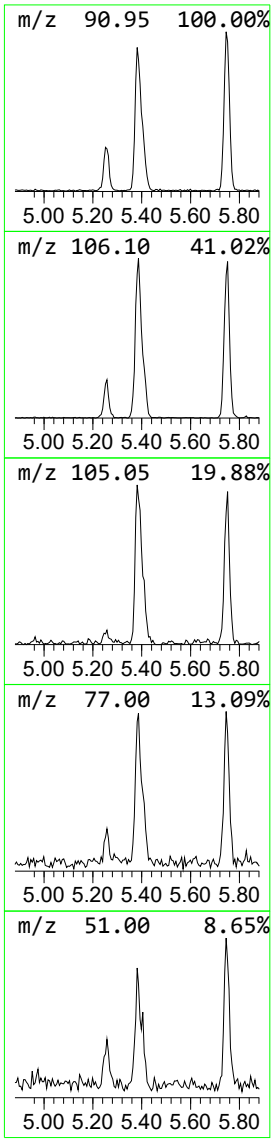
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzene, 1,3-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.381	3.08 ng/ul	109931	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
2	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
3	p-Xylene	106	C8H10	000106-42-3	95
4	p-Xylene	106	C8H10	000106-42-3	94
5	p-Xylene	106	C8H10	000106-42-3	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

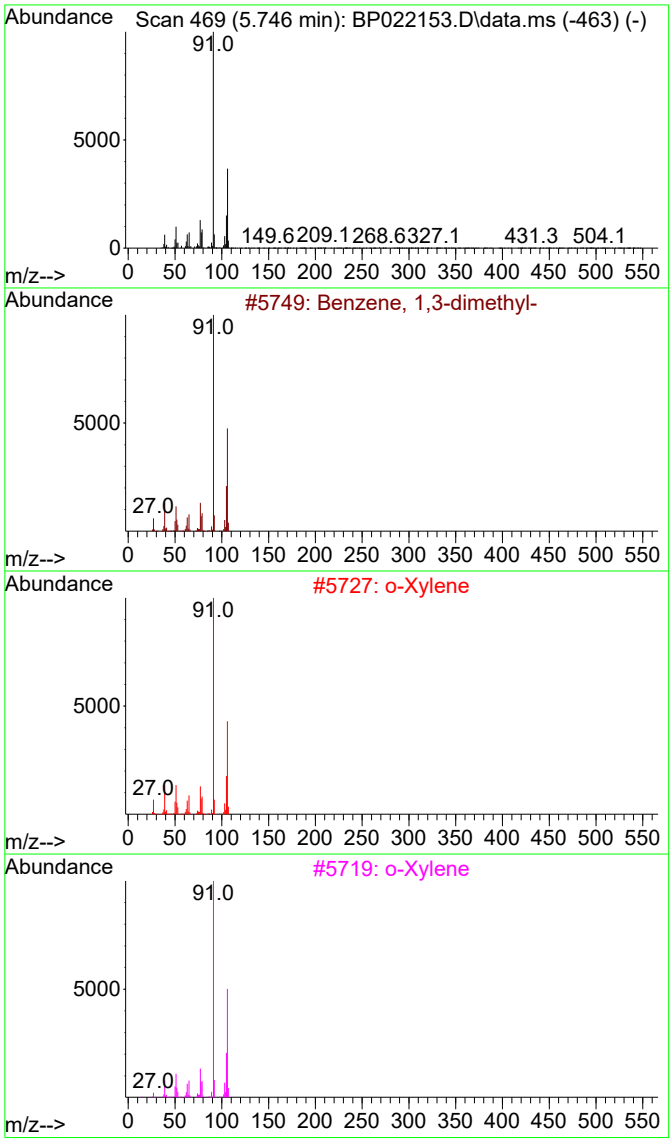
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

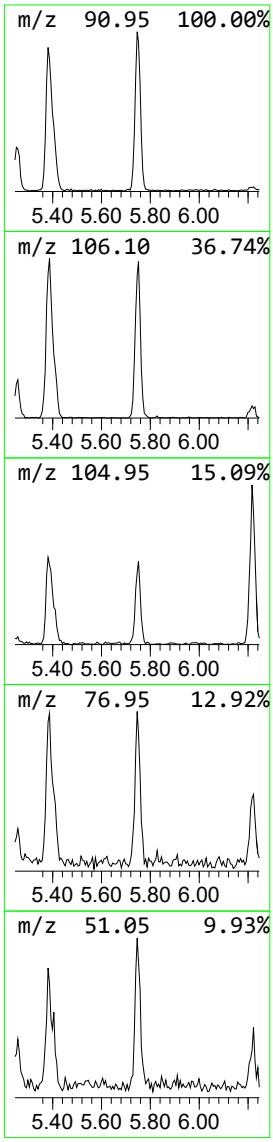
****

Peak Number 2 o-Xylene Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.746	2.51 ng/ul	89569	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2	o-Xylene	106	C8H10	000095-47-6	97
3	o-Xylene	106	C8H10	000095-47-6	97
4	o-Xylene	106	C8H10	000095-47-6	95
5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

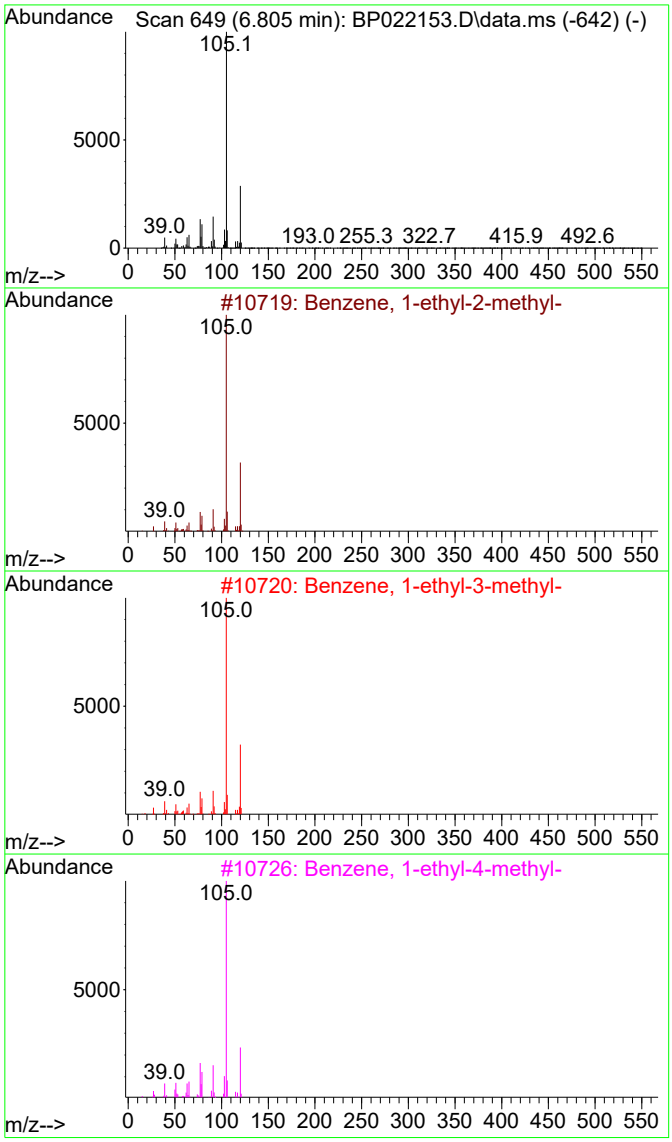
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

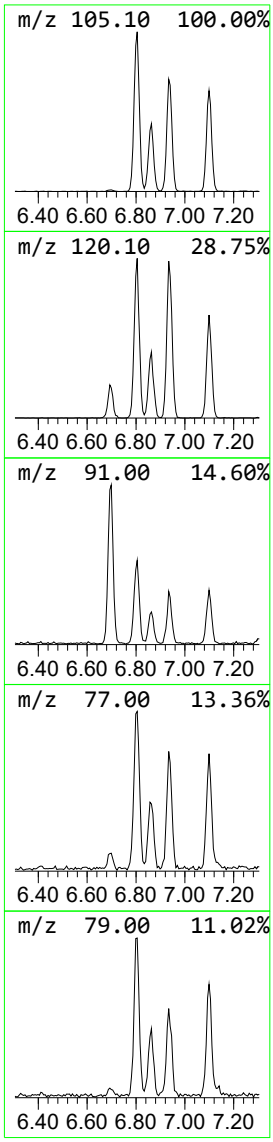
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.805	7.07 ng/ul	252714	1,4-Dichlorobenzene-d4	7.699

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-3-methyl-	120	C9H12	000620-14-4	91
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

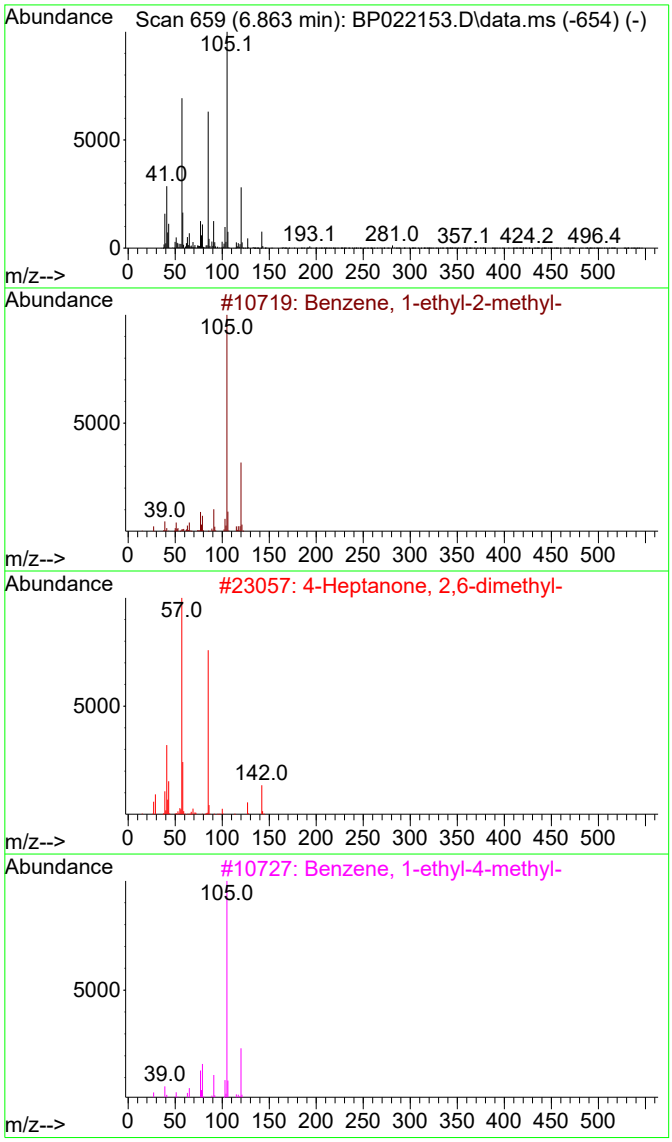
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

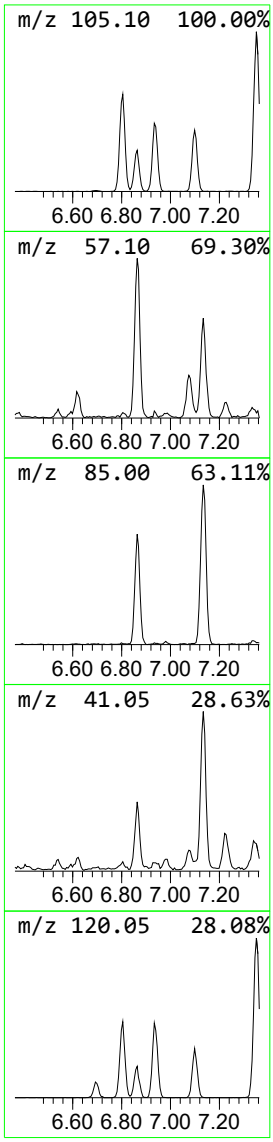
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 4-Heptanone, 2,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.863	5.84 ng/ul	208654	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	86
2	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	83
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	64
4	Mesitylene	120	C9H12	000108-67-8	50
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

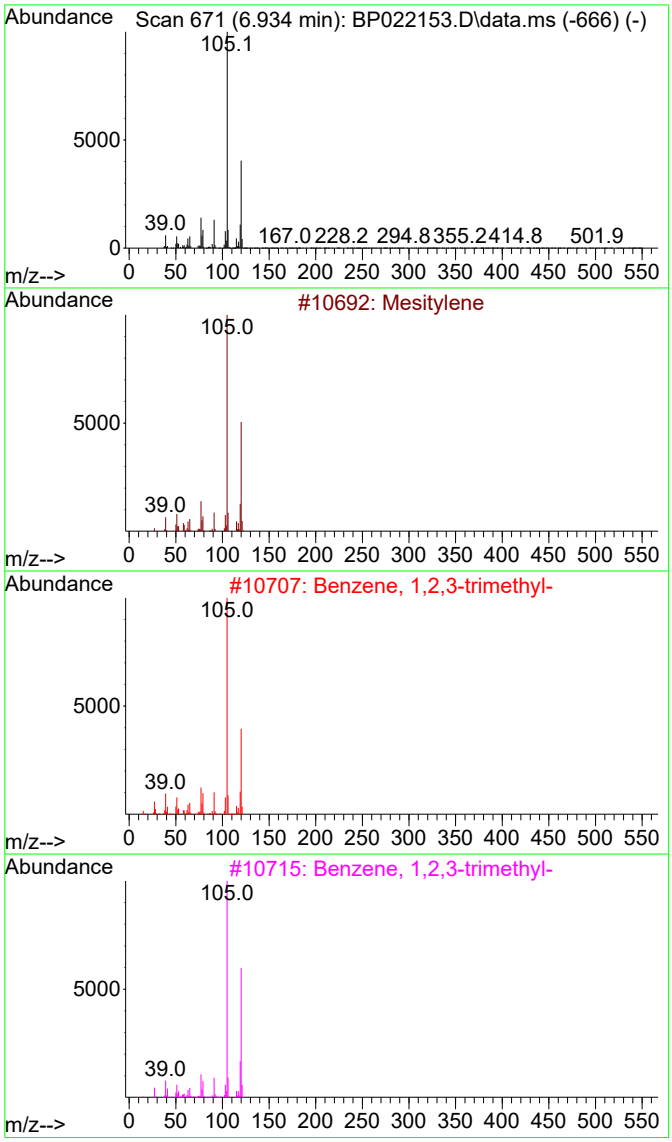
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

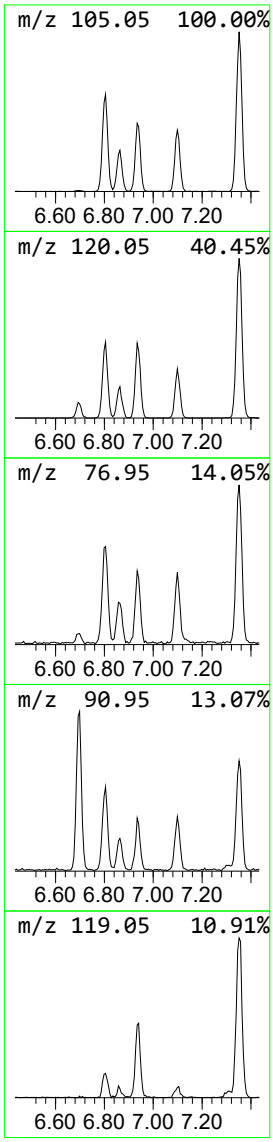
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Mesitylene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.934	5.29 ng/ul	189070	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

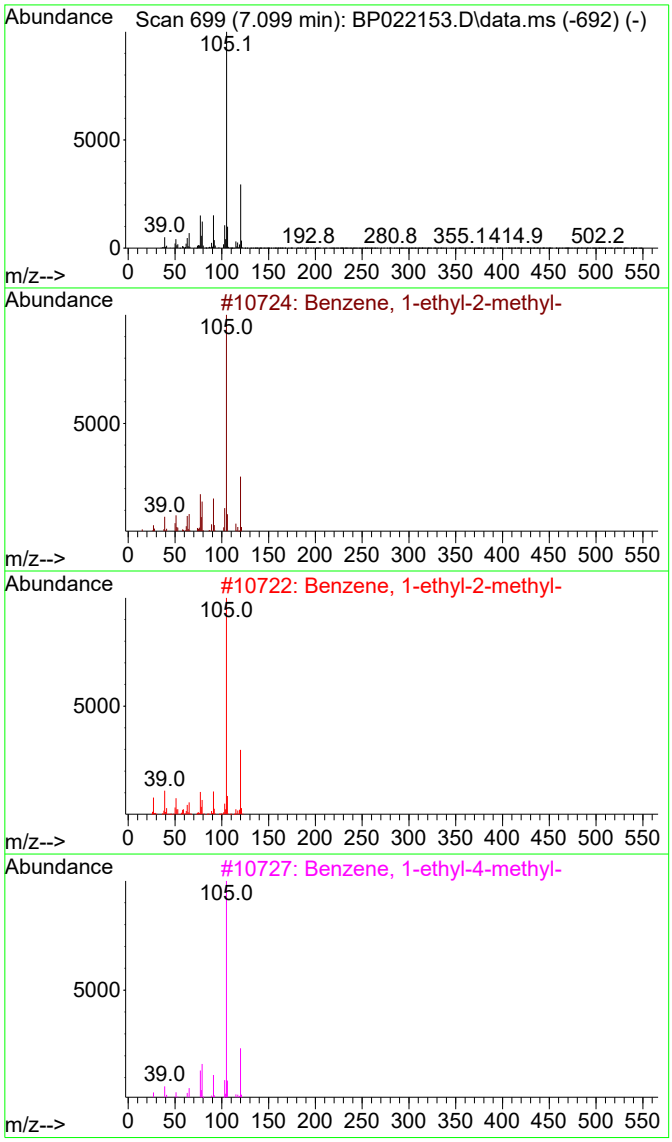
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

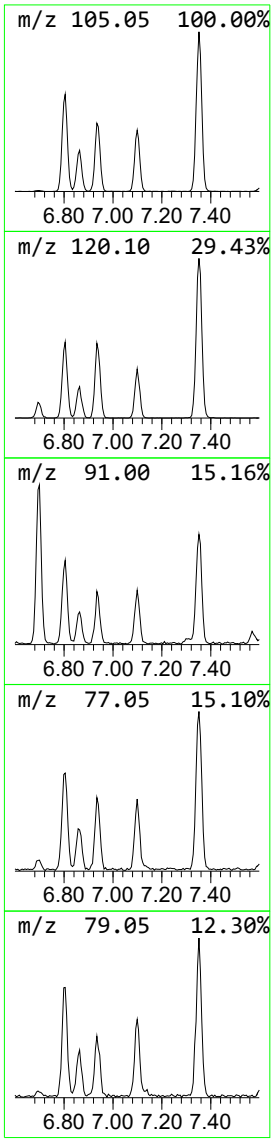
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1-ethyl-4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.099	5.04 ng/ul	180341	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





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Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

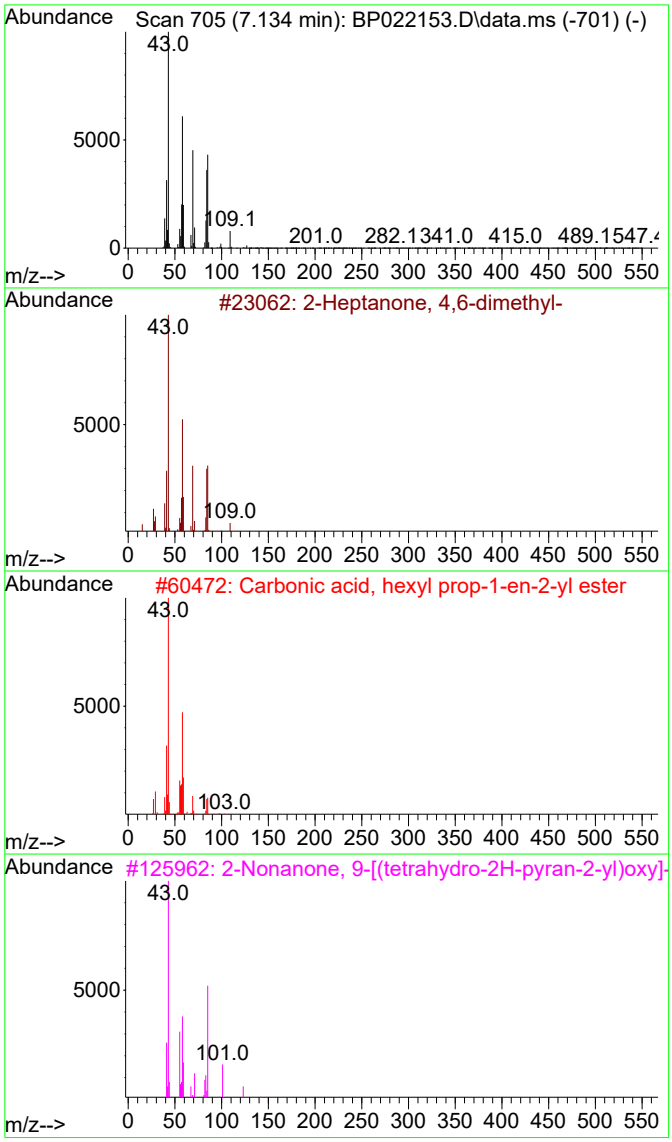
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

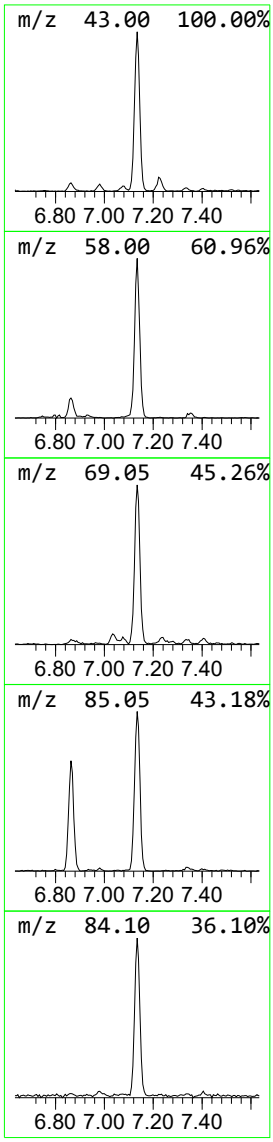
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 2-Heptanone, 4,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.134	10.82 ng/ul	386823	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	90
2	Carbonic acid, hexyl prop-1-en-2...	186	C10H18O3	1000382-54-2	78
3	2-Nonanone, 9-[(tetrahydro-2H-py...	242	C14H26O3	054699-41-1	59
4	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	50
5	Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	43





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Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

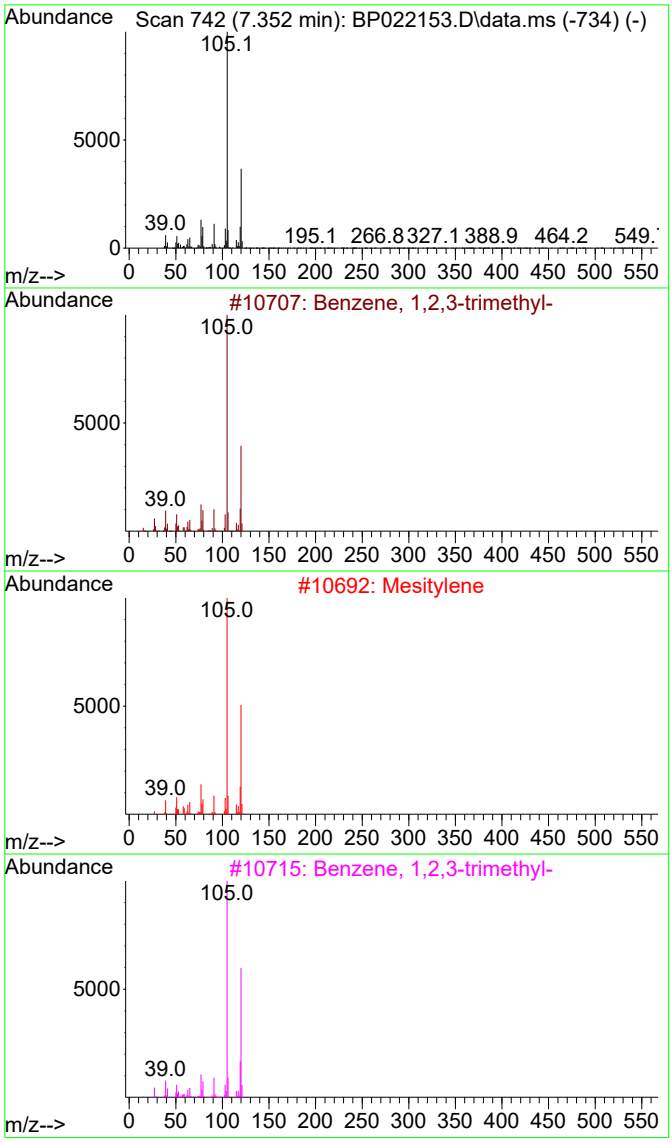
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

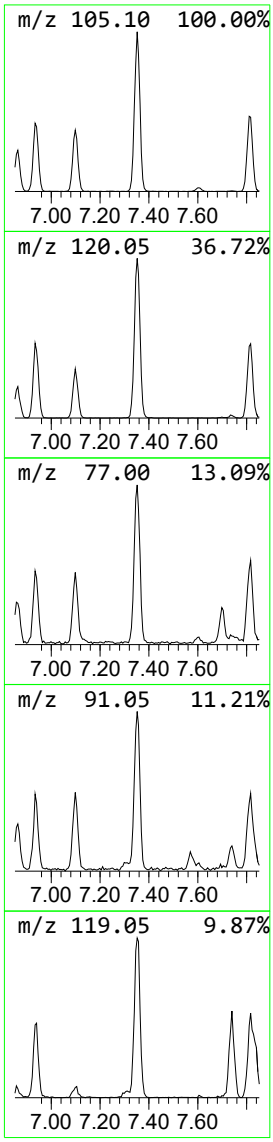
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.352	13.39 ng/ul	478622	1,4-Dichlorobenzene-d4	7.699

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





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Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

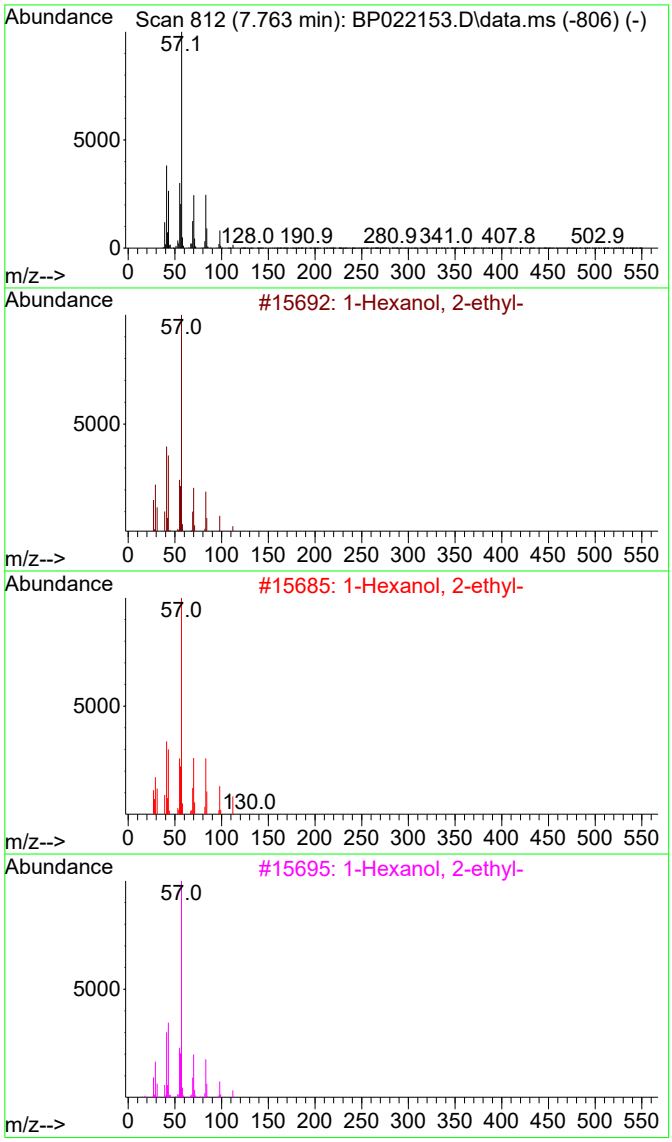
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

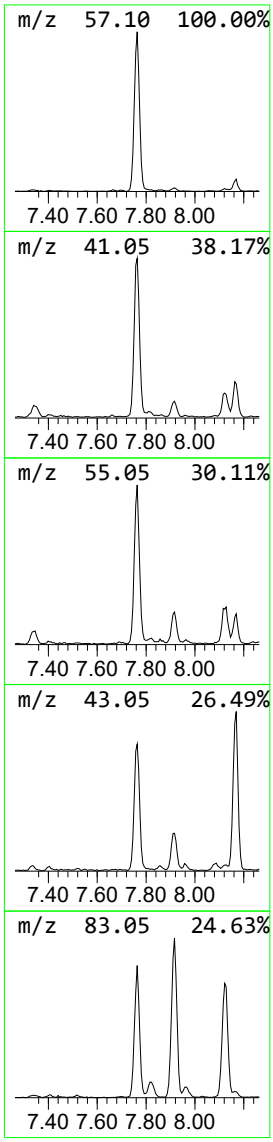
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1-Hexanol, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.763	16.73 ng/ul	598235	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	90
2	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
3	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
4	2-Propyl-1-pentanol	130	C8H18O	058175-57-8	64
5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	59





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Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

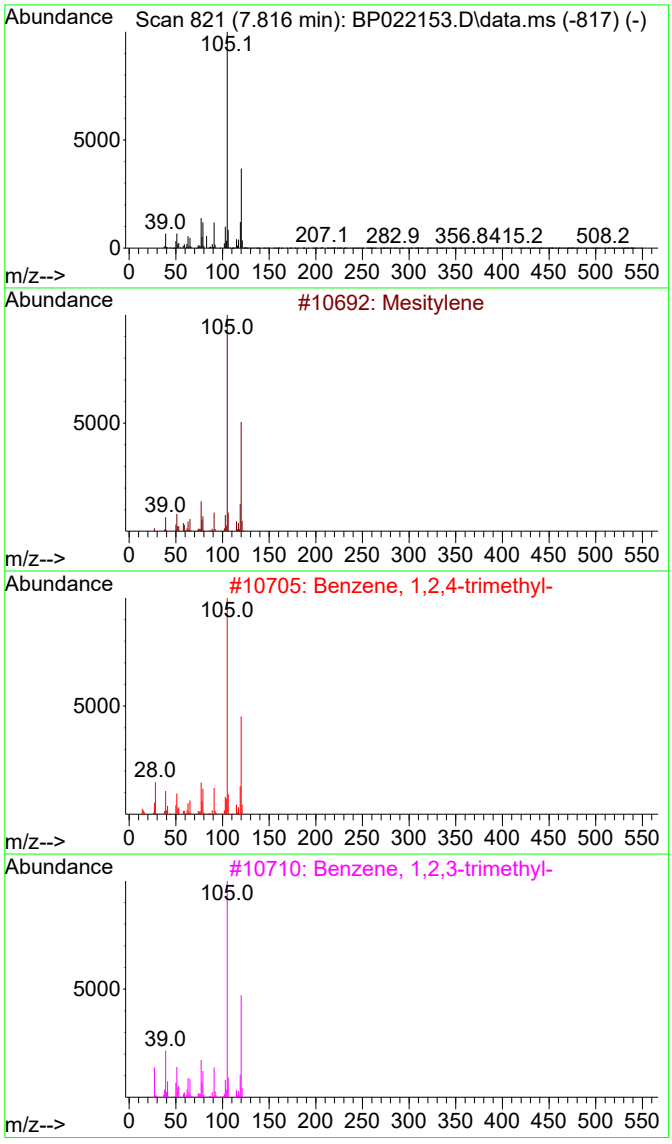
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

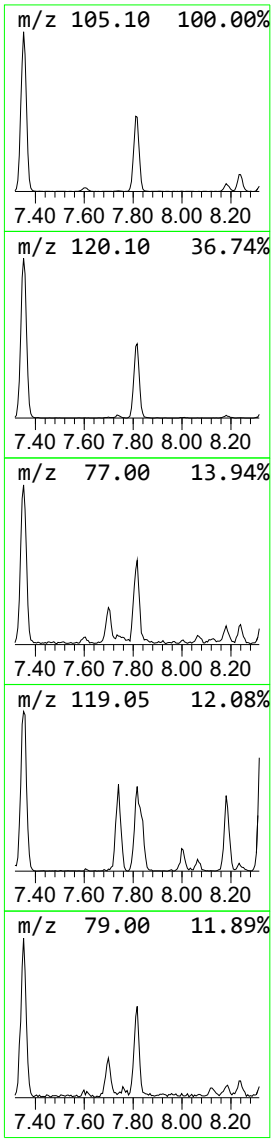
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1,2,4-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.816	6.75 ng/ul	241338	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	94
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





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Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

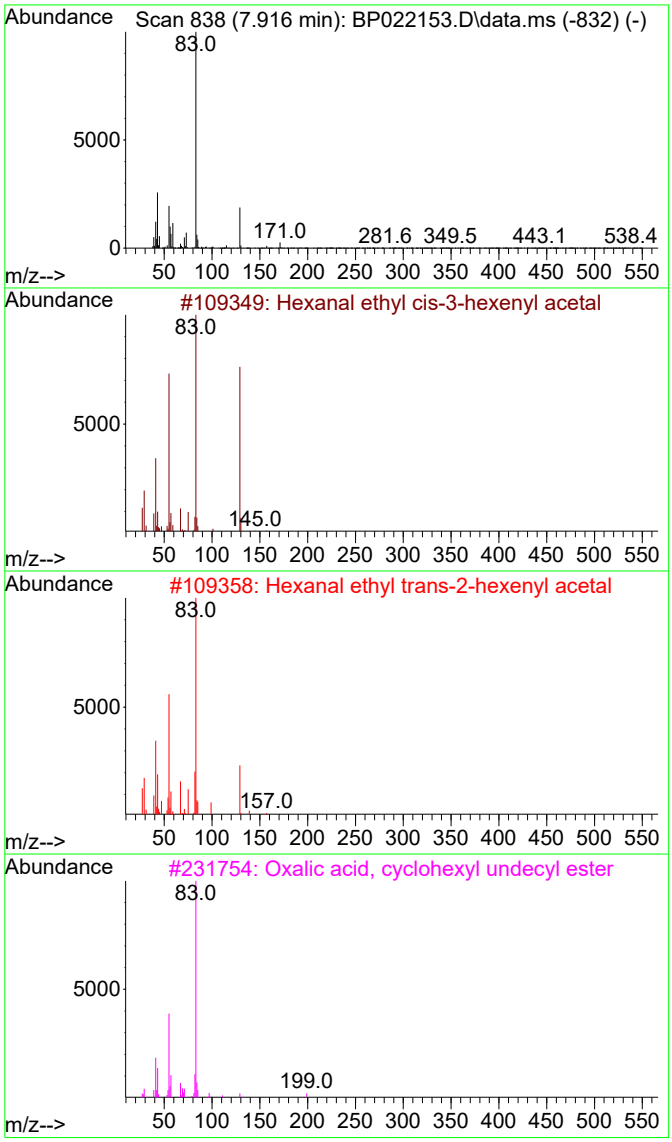
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

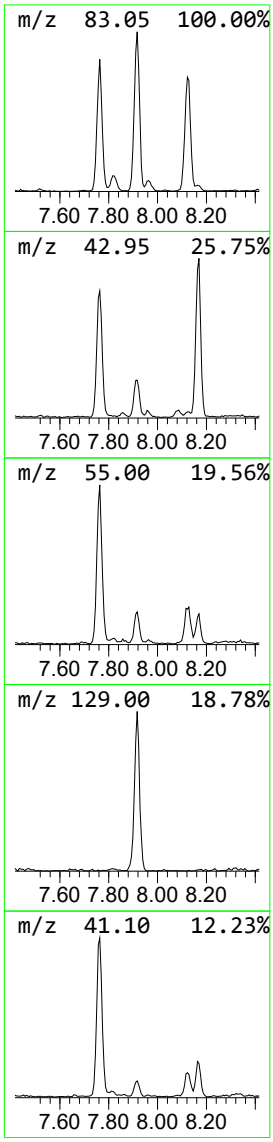
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Hexanal ethyl cis-3-hexenyl... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.916	3.56 ng/ul	127267	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanal ethyl cis-3-hexenyl acetal	228	C14H28O2	1000431-42-9	64
2	Hexanal ethyl trans-2-hexenyl ac...	228	C14H28O2	1000431-42-6	53
3	Oxalic acid, cyclohexyl undecyl ...	326	C19H34O4	1000309-31-3	45
4	Pentanoic acid, 2,4-dimethyl-4-n...	189	C8H15NO4	005762-40-3	39
5	Oxalic acid, cyclohexyl dodecyl ...	340	C20H36O4	1000309-31-4	38





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Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

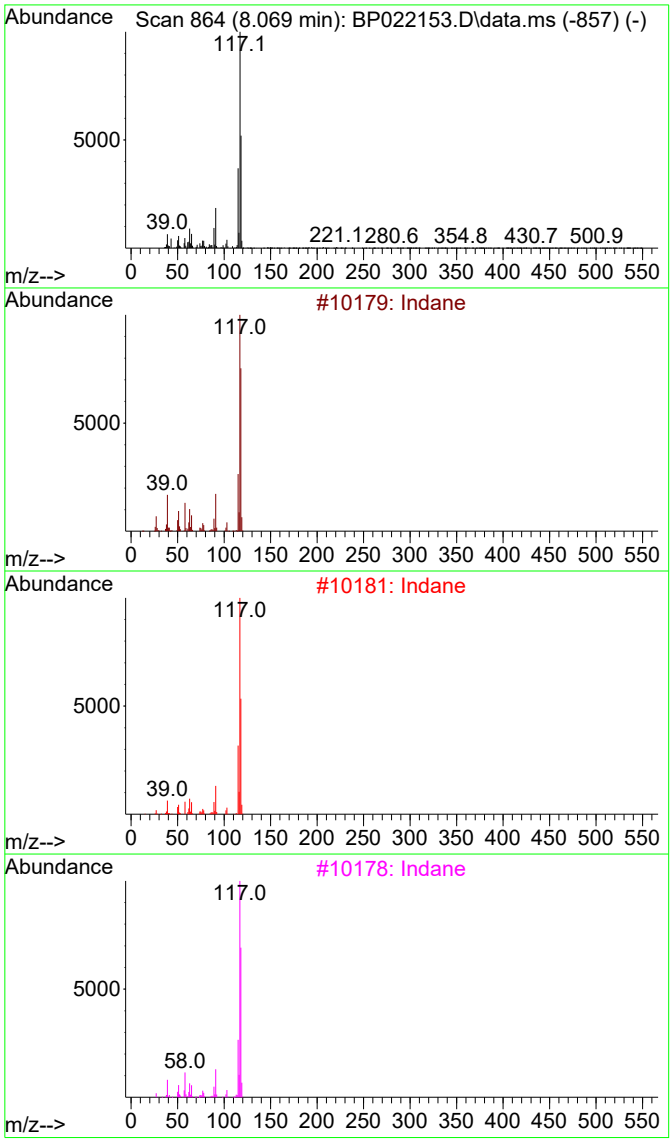
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

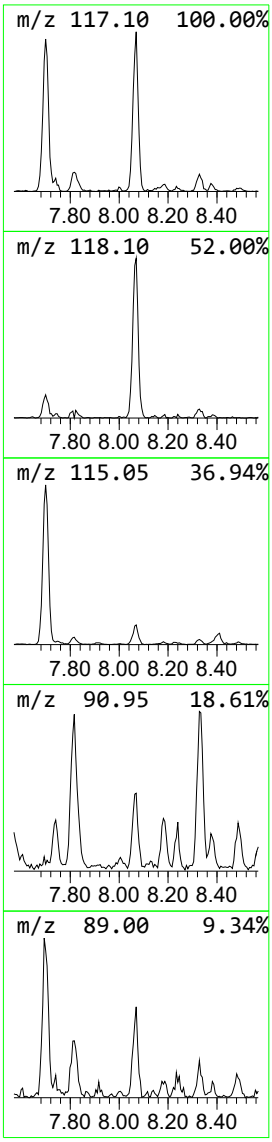
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Indane Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.069	2.19 ng/ul	78375	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	83
2	Indane	118	C9H10	000496-11-7	81
3	Indane	118	C9H10	000496-11-7	68
4	Deltacyclene	118	C9H10	007785-10-6	68
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64





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Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

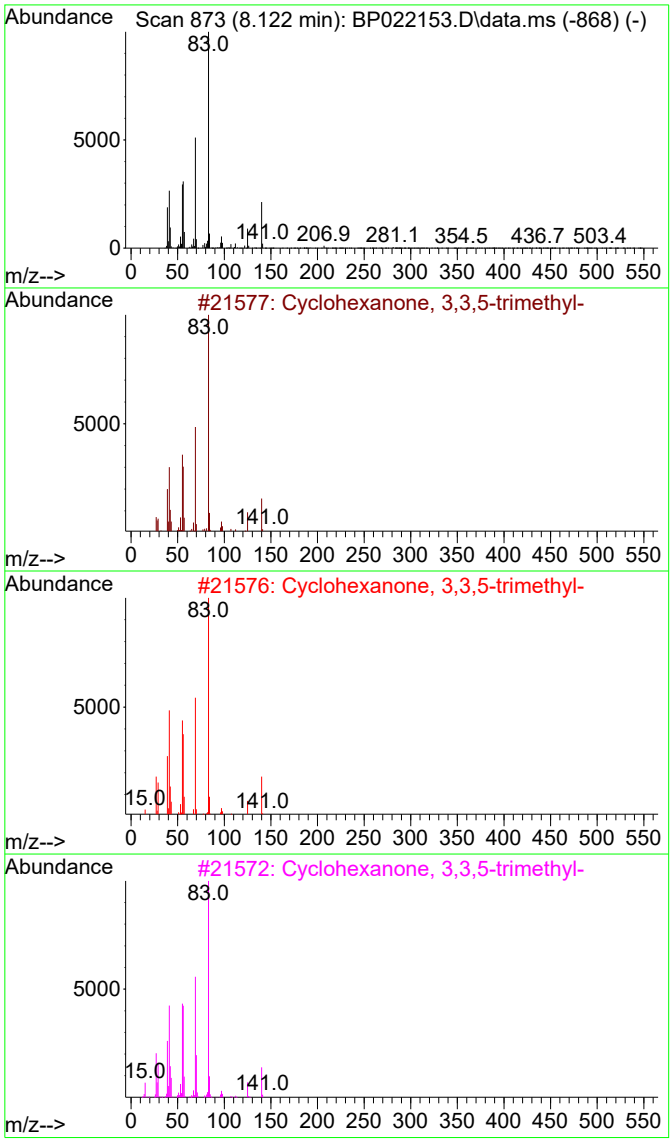
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

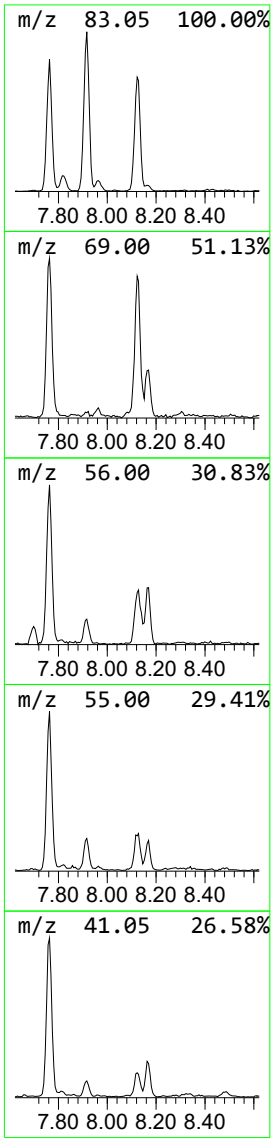
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Cyclohexanone, 3,3,5-trimet... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.122	4.13 ng/ul	147515	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	94
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	90
3	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
4	Cyclopentanone, 3-butyl-	140	C9H16O	057283-81-5	47
5	Methanone, dicyclohexyl-	194	C13H22O	000119-60-8	43





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Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

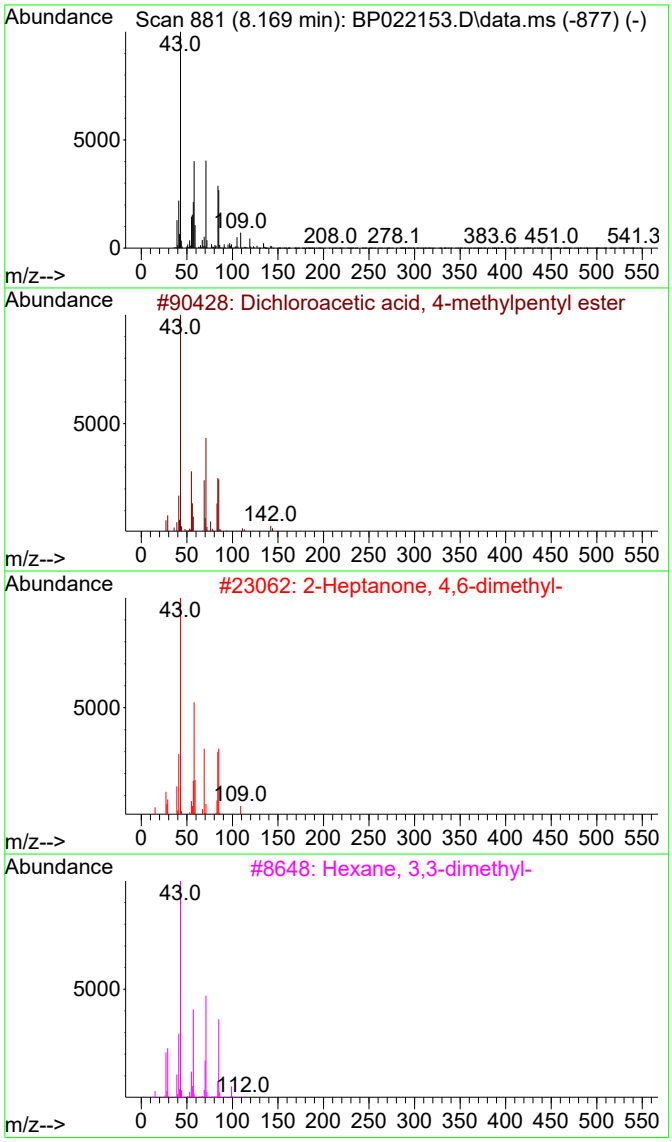
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

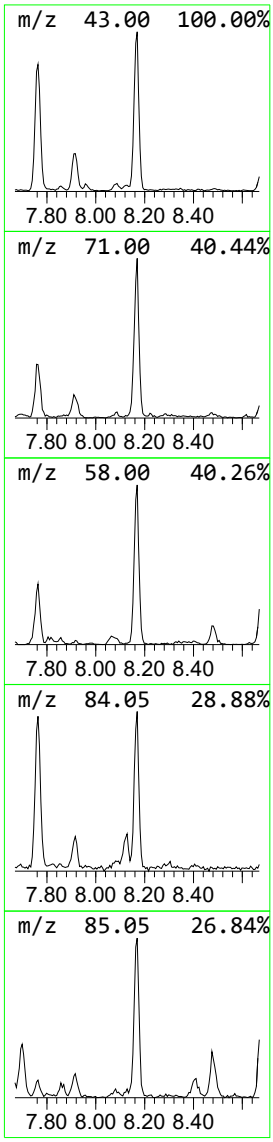
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Dichloroacetic acid, 4-meth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.169	6.61 ng/ul	236317	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dichloroacetic acid, 4-methylpen...	212	C8H14Cl2O2	1000282-43-7	50
2	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	47
3	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	43
4	2-Pentanone, 5,5'-oxybis-	186	C10H18O3	093677-66-8	38
5	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	38





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Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

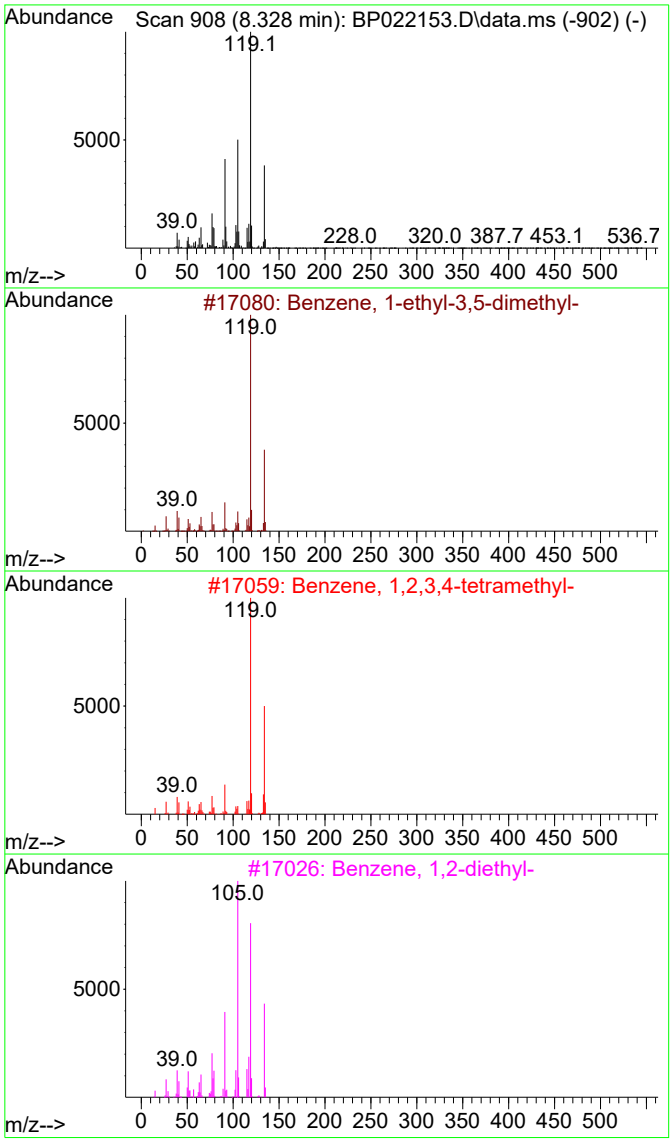
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

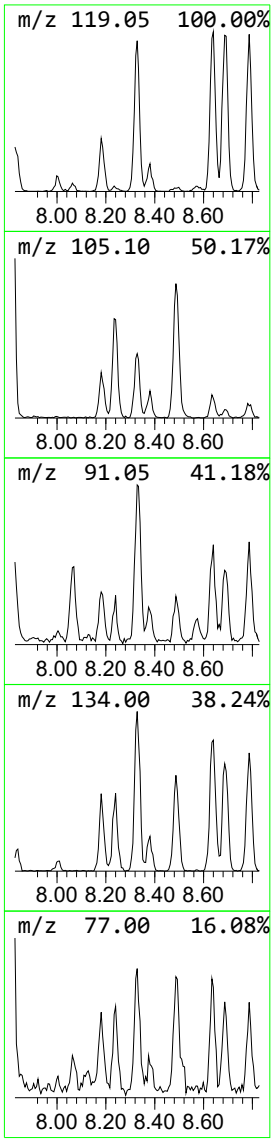
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.328	2.66 ng/ul	94944	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	95
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
3	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91





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Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

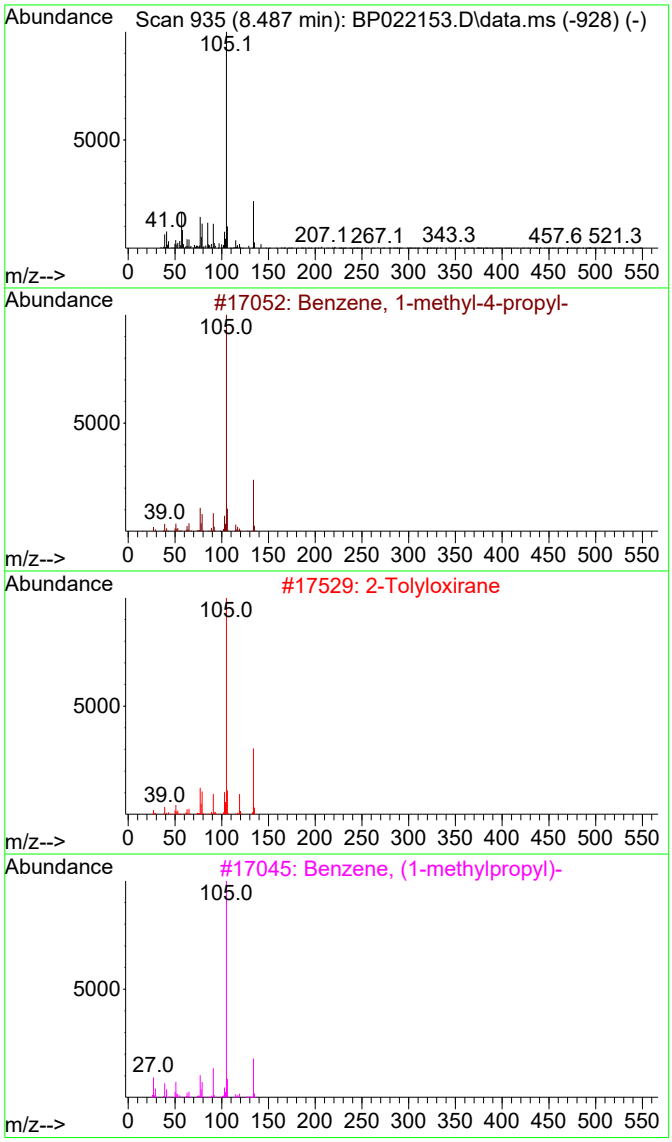
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

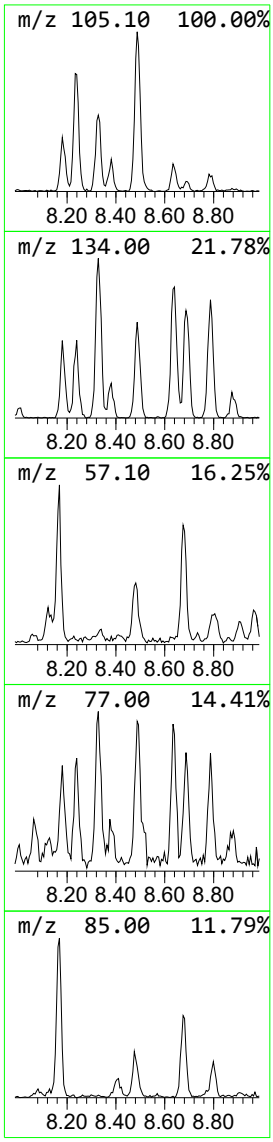
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Benzene, 1-methyl-4-propyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.487	2.34 ng/ul	83776	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	81
2	2-Tolyloxirane	134	C9H10O	002783-26-8	81
3	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	81
4	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	81
5	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	81





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Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

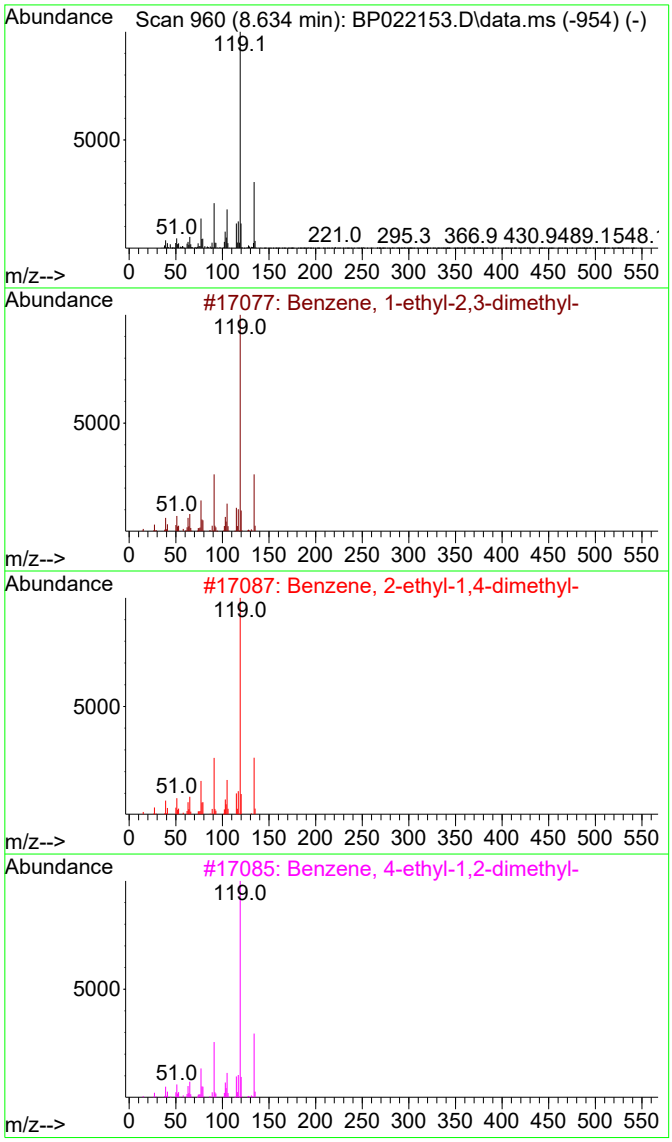
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

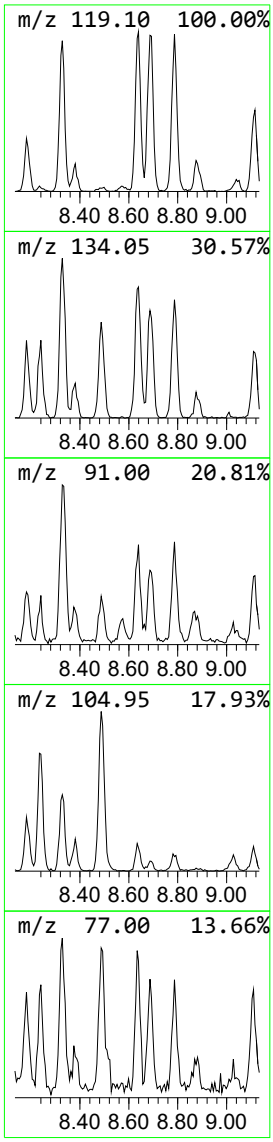
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.634	2.03 ng/ul	72491	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	97
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5	o-Cymene	134	C10H14	000527-84-4	94





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Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

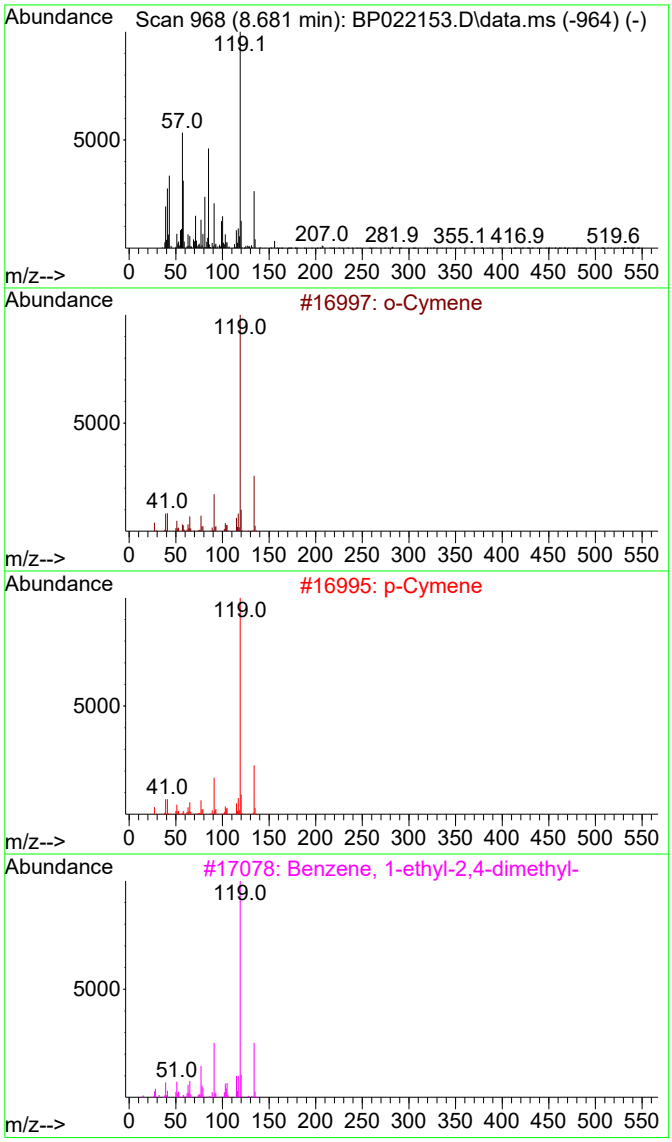
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

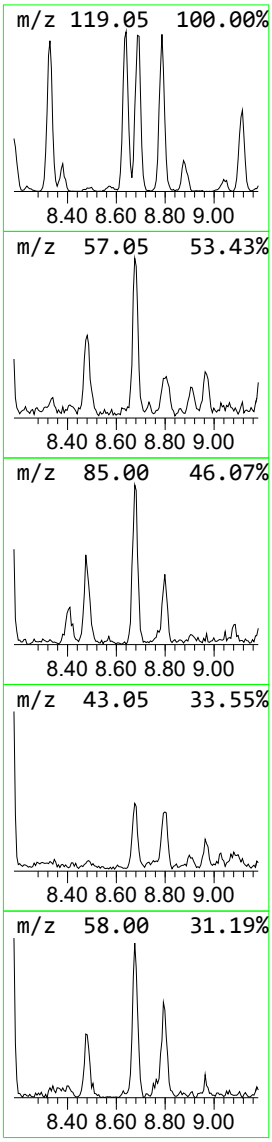
****

Peak Number 18 o-Cymene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.681	3.28 ng/ul	117326	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	87
2	p-Cymene	134	C10H14	000099-87-6	87
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	70
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	55
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	55





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Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

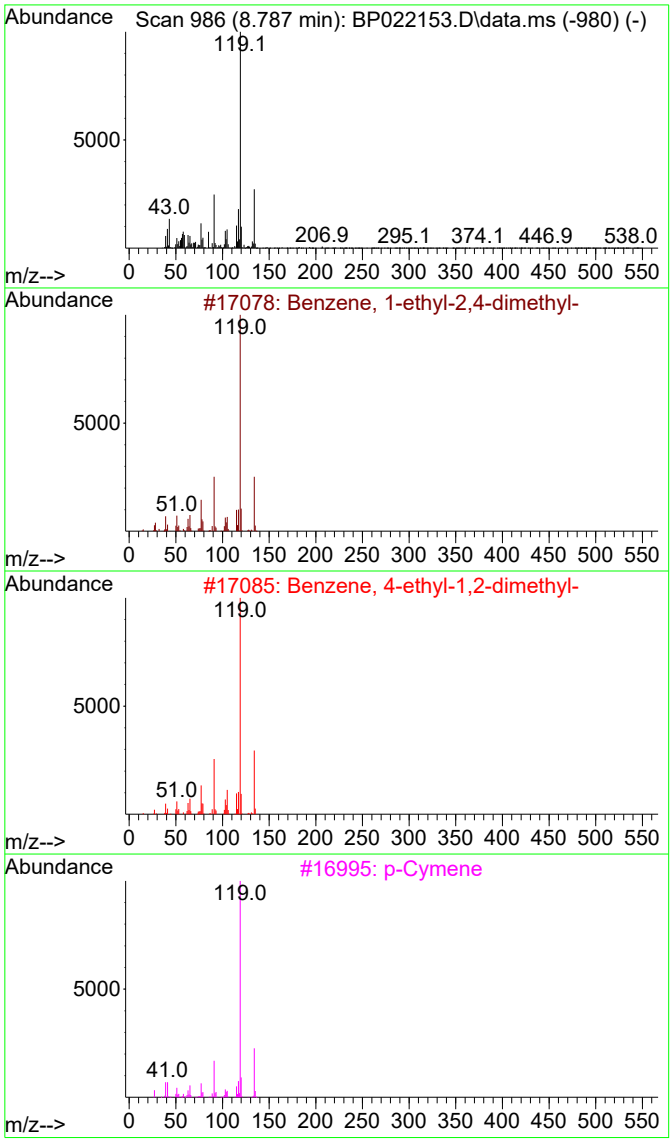
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

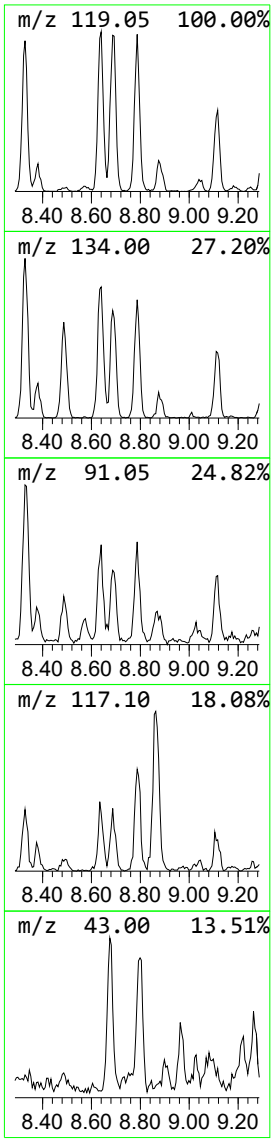
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.787	2.92 ng/ul	104484	1,4-Dichlorobenzene-d4	7.699

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
3	p-Cymene	134	C10H14	000099-87-6	93
4	o-Cymene	134	C10H14	000527-84-4	93
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93





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Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

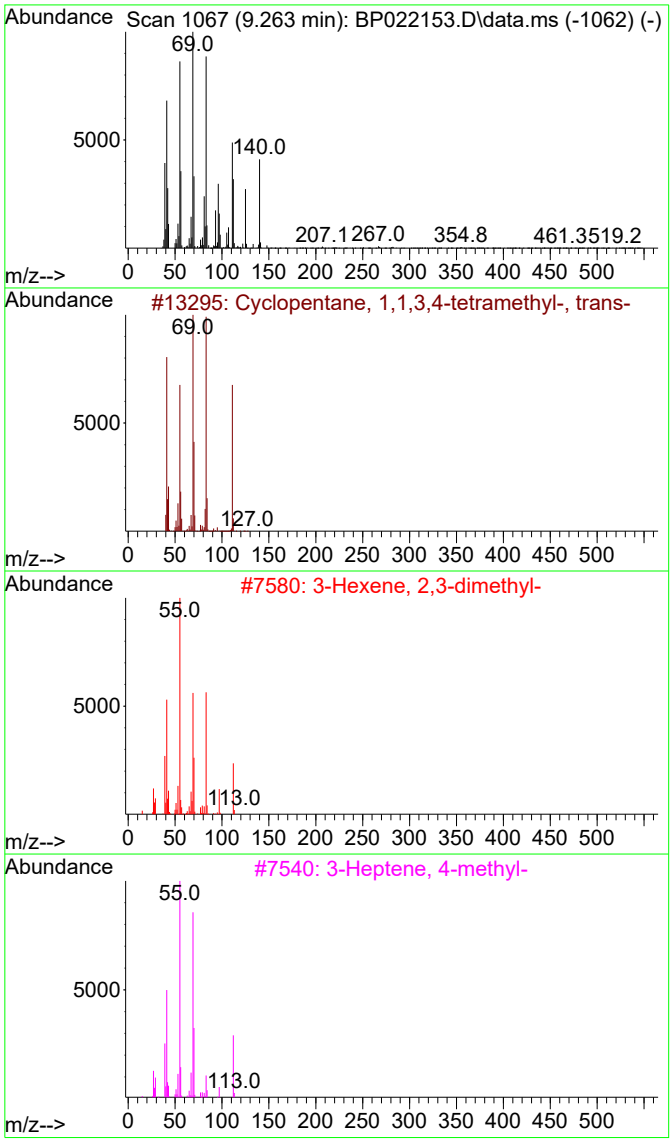
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

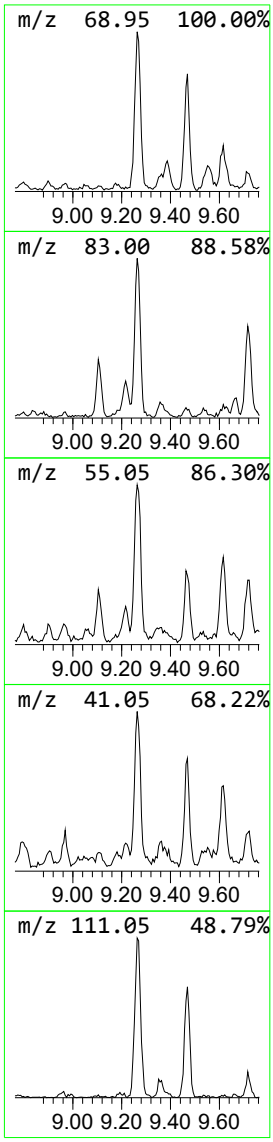
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 (DEL) Alkane: Cyclic9.263 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.263	2.30 ng/ul	164843	Naphthalene-d8	10.457

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,1,3,4-tetramethy...	126	C9H18	020309-77-7	50
2	3-Hexene, 2,3-dimethyl-	112	C8H16	007145-23-5	49
3	3-Heptene, 4-methyl-	112	C8H16	004485-16-9	38
4	3-Ethyl-4-methyl-2-pentene	112	C8H16	019780-68-8	38
5	3-Hexene, 3,4-dimethyl-	112	C8H16	030951-95-2	30





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Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

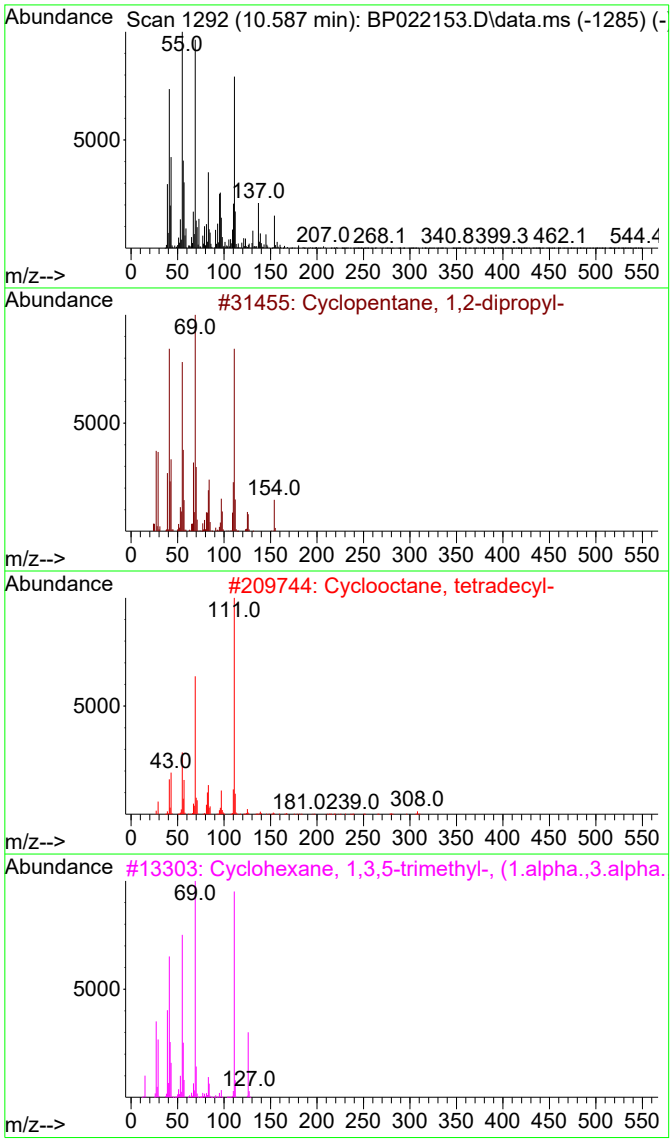
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

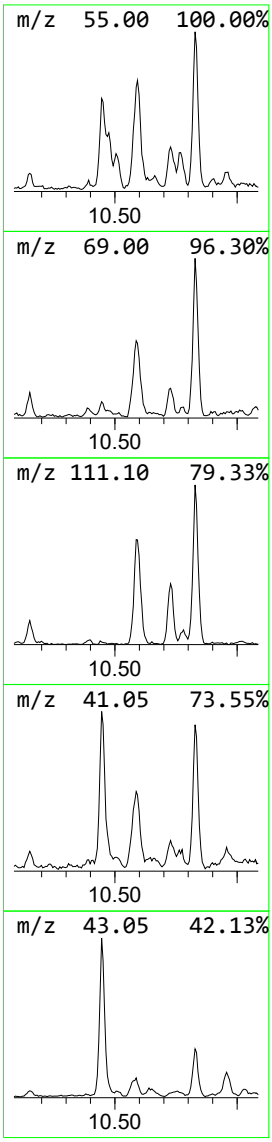
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 (DEL) Alkane: Cyclic10.587 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.587	3.98 ng/ul	284856	Naphthalene-d8	10.457

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,2-dipropyl-	154	C11H22	091242-57-8	81
2	Cyclooctane, tetradecyl-	308	C22H44	149003-36-1	49
3	Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	43
4	3-Heptene, 2-methyl-, (E)-	112	C8H16	000692-96-6	41
5	Cyclohexanone, 2-methyl-5-(1-met...	154	C10H18O	059471-80-6	38





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Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

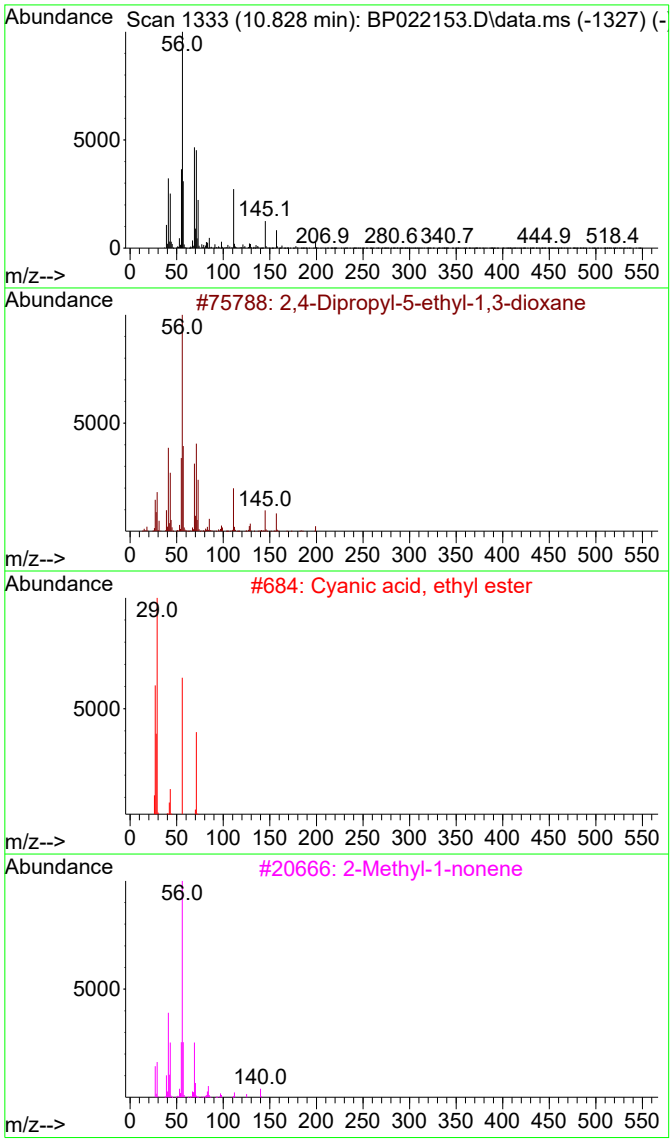
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

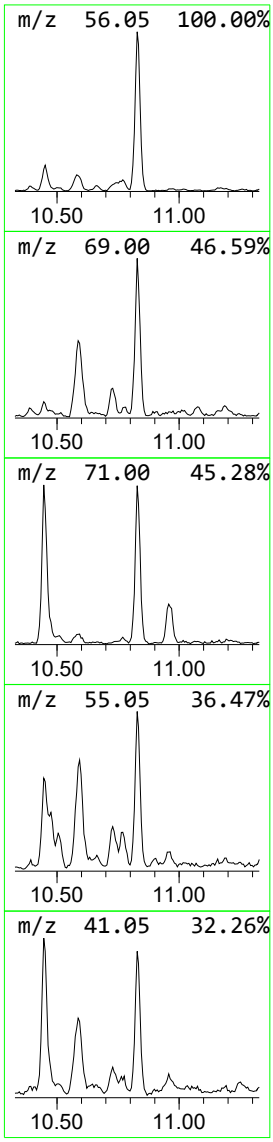
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 2,4-Dipropyl-5-ethyl-1,3-di... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.828	5.48 ng/ul	392358	Naphthalene-d8	10.457

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	72
2	Cyanic acid, ethyl ester	71	C3H5NO	000627-48-5	40
3	2-Methyl-1-nonene	140	C10H20	002980-71-4	35
4	2-Chloro-2-methylnonane	176	C10H21Cl	004325-50-2	32
5	3,4,5-Trimethyldihydrofuran-2-one	128	C7H12O2	1000194-05-2	25





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Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

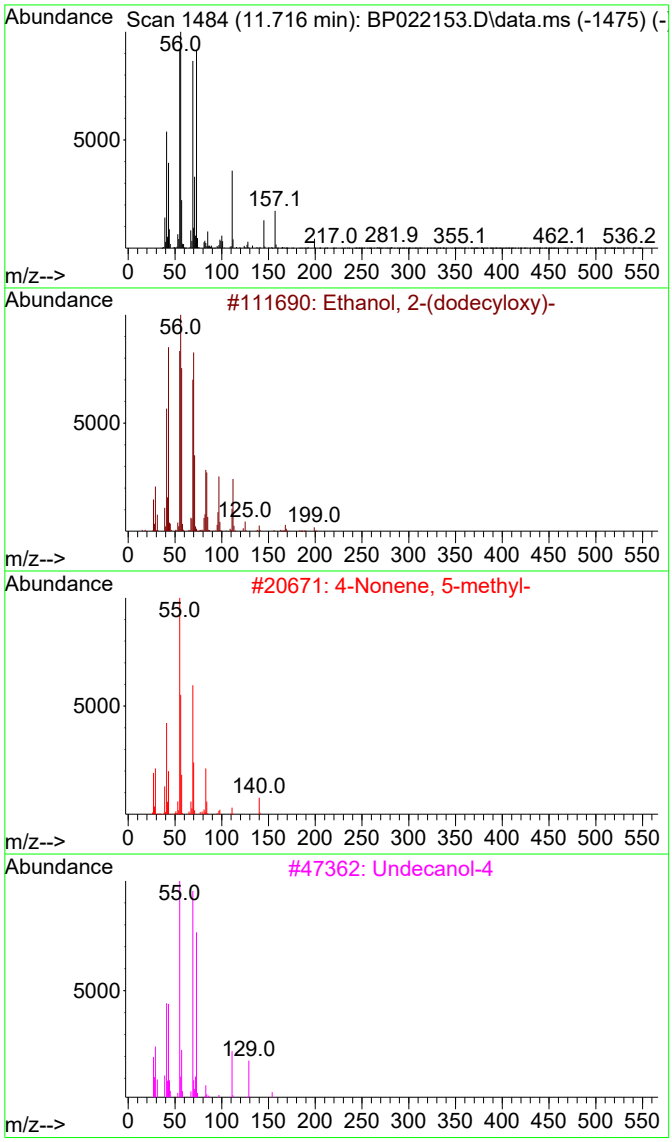
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

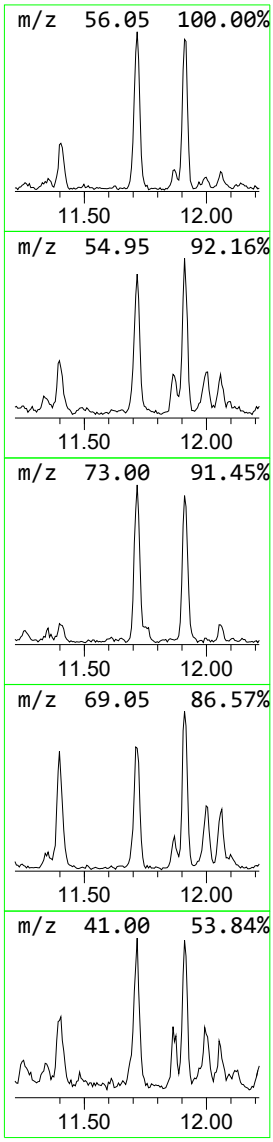
****

Peak Number 23 unknown-01 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.716	2.99 ng/ul	214576	Naphthalene-d8	10.457

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(dodecyloxy)-	230	C14H30O2	004536-30-5	45
2	4-Nonene, 5-methyl-	140	C10H20	015918-07-7	38
3	Undecanol-4	172	C11H24O	004272-06-4	37
4	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	37
5	1H-1,2,4-Triazole-3-carboxaldehy...	111	C4H5N3O	056804-98-9	35





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Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

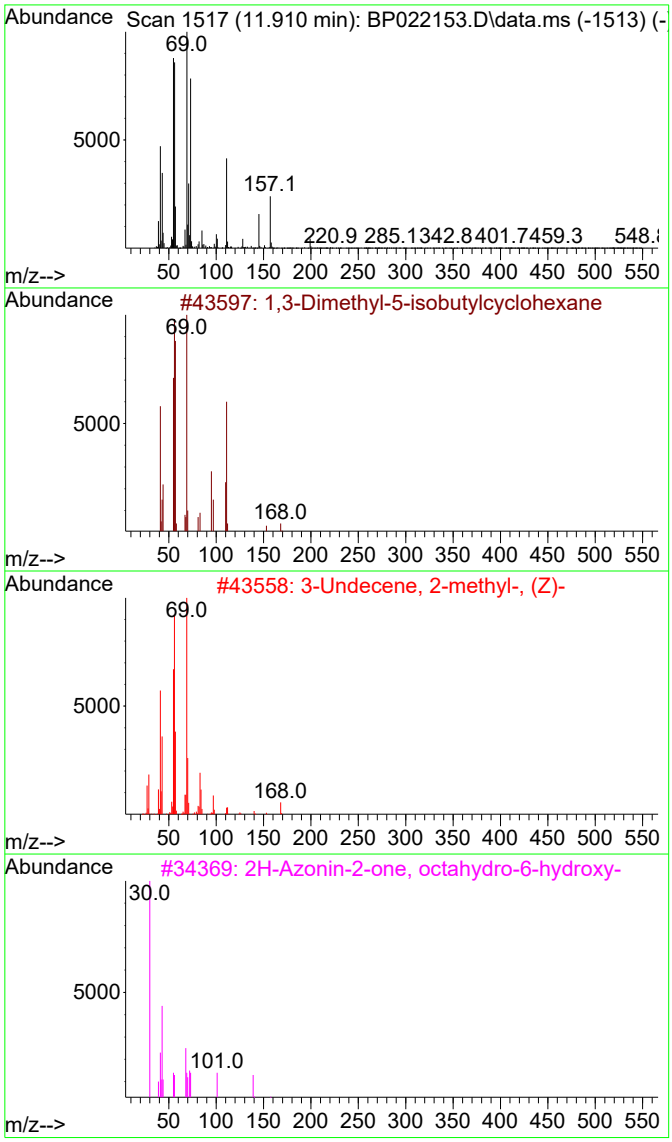
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

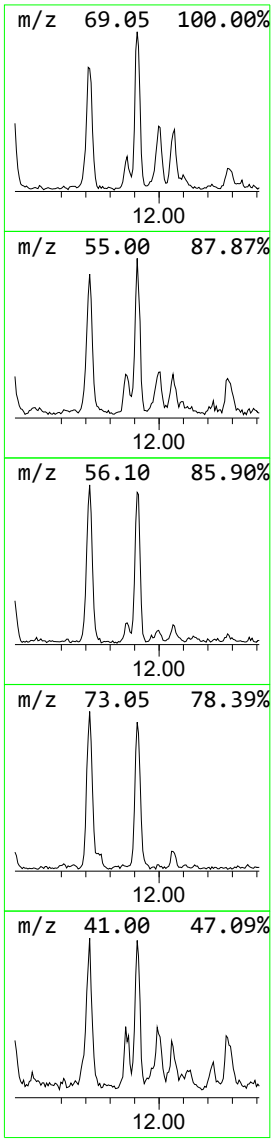
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 (DEL) Alkane: Cyclic11.910 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.910	2.12 ng/ul	151857	Naphthalene-d8	10.457

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dimethyl-5-isobutylcyclohexane	168	C12H24	013131-76-5	53
2	3-Undecene, 2-methyl-, (Z)-	168	C12H24	074630-48-1	38
3	2H-Azonin-2-one, octahydro-6-hyd...	157	C8H15NO2	023435-07-6	36
4	2-Pyrrolidinone, 1-ethenyl-	111	C6H9NO	000088-12-0	27
5	2,5-Dimethyl-5-hexen-3-ol	128	C8H16O	067760-91-2	25





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Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

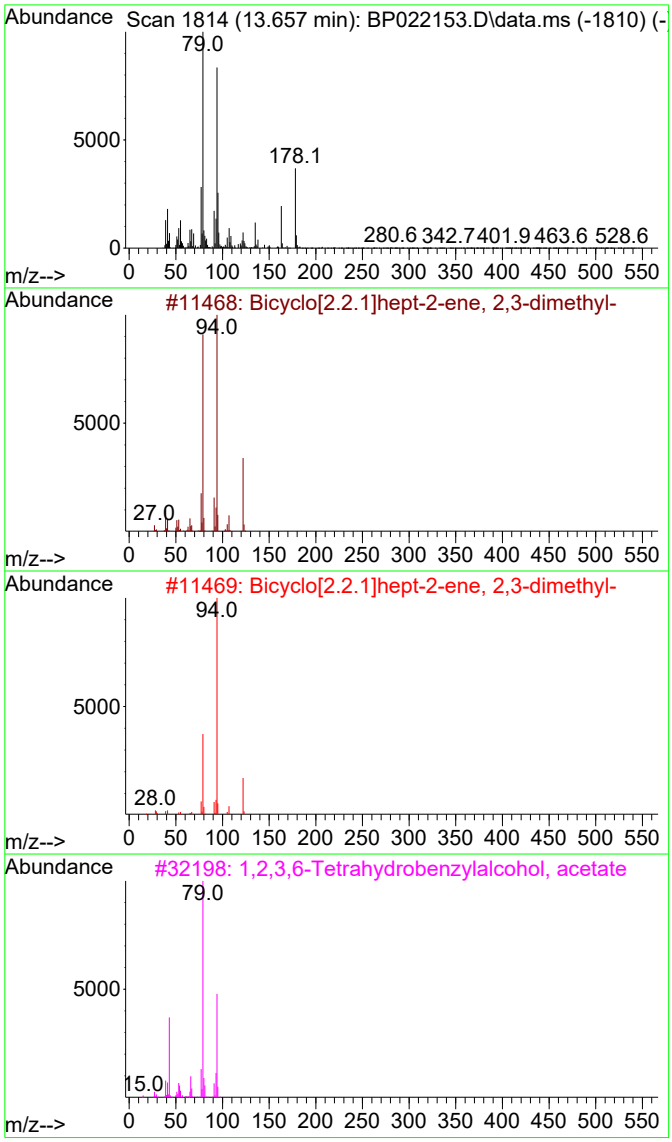
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

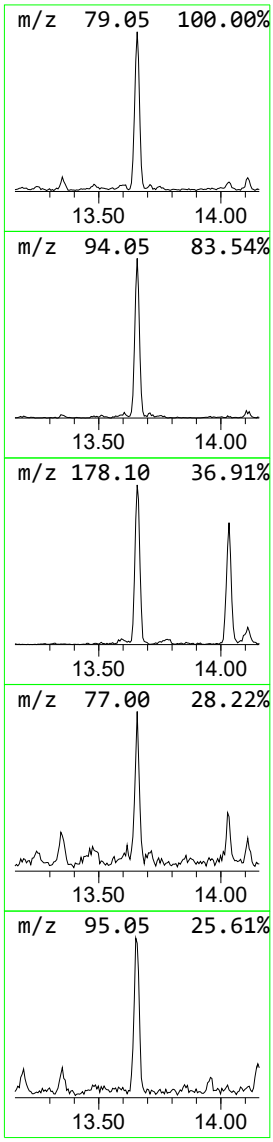
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Bicyclo[2.2.1]hept-2-ene, 2... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.657	3.24 ng/ul	226576	Acenaphthene-d10	14.304

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]hept-2-ene, 2,3-di...	122	C9H14	000529-16-8	58
2	Bicyclo[2.2.1]hept-2-ene, 2,3-di...	122	C9H14	000529-16-8	50
3	1,2,3,6-Tetrahydrobenzylalcohol,...	154	C9H14O2	1000365-58-6	50
4	1,3-Cyclopentadiene, 1,2-dimethyl-	94	C7H10	004784-86-5	49
5	1,3-Cyclopentadiene, 5,5-dimethyl-	94	C7H10	004125-18-2	47





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
Data File : BP022153.D
Acq On : 02 Oct 2024 03:22
Operator : RC/JU
Sample : P4022-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC95DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.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, 1,3-di...	5.381	3.1	ng/ul	109931	1	7.699	714989	20.0
o-Xylene	5.746	2.5	ng/ul	89569	1	7.699	714989	20.0
Benzene, 1-ethy...	6.805	7.1	ng/ul	252714	1	7.699	714989	20.0
4-Heptanone, 2,...	6.863	5.8	ng/ul	208654	1	7.699	714989	20.0
Mesitylene	6.934	5.3	ng/ul	189070	1	7.699	714989	20.0
Benzene, 1-ethy...	7.099	5.0	ng/ul	180341	1	7.699	714989	20.0
2-Heptanone, 4,...	7.134	10.8	ng/ul	386823	1	7.699	714989	20.0
Benzene, 1,2,3-...	7.352	13.4	ng/ul	478622	1	7.699	714989	20.0
1-Hexanol, 2-et...	7.763	16.7	ng/ul	598235	1	7.699	714989	20.0
Benzene, 1,2,4-...	7.816	6.8	ng/ul	241338	1	7.699	714989	20.0
Hexanal ethyl c...	7.916	3.6	ng/ul	127267	1	7.699	714989	20.0
Indane	8.069	2.2	ng/ul	78375	1	7.699	714989	20.0
Cyclohexanone, ...	8.122	4.1	ng/ul	147515	1	7.699	714989	20.0
Dichloroacetic ...	8.169	6.6	ng/ul	236317	1	7.699	714989	20.0
Benzene, 1-ethy...	8.328	2.7	ng/ul	94944	1	7.699	714989	20.0
Benzene, 1-meth...	8.487	2.3	ng/ul	83776	1	7.699	714989	20.0
Benzene, 1-ethy...	8.634	2.0	ng/ul	72491	1	7.699	714989	20.0
o-Cymene	8.681	3.3	ng/ul	117326	1	7.699	714989	20.0
Benzene, 1-ethy...	8.787	2.9	ng/ul	104484	1	7.699	714989	20.0
(DEL) Alkane: C...	9.263	2.3	ng/ul	164843	2	10.457	1433150	20.0
(DEL) Alkane: C...	10.587	4.0	ng/ul	284856	2	10.457	1433150	20.0
2,4-Dipropyl-5-...	10.828	5.5	ng/ul	392358	2	10.457	1433150	20.0
unknown-01	11.716	3.0	ng/ul	214576	2	10.457	1433150	20.0
(DEL) Alkane: C...	11.910	2.1	ng/ul	151857	2	10.457	1433150	20.0
Bicyclo[2.2.1]h...	13.657	3.2	ng/ul	226576	3	14.304	1397900	20.0