LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023977.D
Acq On : 06 Feb 2025 19:22
Operator : RC/JU
Sample : Q1202-19DL2 150X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B3DL2

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-BP012925.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP023977.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.681	266	271	280	rBV	40510	67097	0.23%	0.081%
2	4.816	290	294	302	rVB2	23268	36849	0.13%	0.044%
3	6.981	654	662	672	rBV	424605	680003	2.32%	0.820%
4	7.769	787	796	802	rBV	1598438	2366989	8.07%	2.855%
5	7.834	802	807	833	rVB	402309	858340	2.93%	1.035%
6	8.357	889	896	911	rBV	163184	309888	1.06%	0.374%
7	8.540	921	927	936	rVB	333129	590394	2.01%	0.712%
8	9.063	1008	1016	1027	rBV	333726	622145	2.12%	0.750%
9	9.175	1029	1035	1044	rVB2	36040	80663	0.28%	0.097%
10	9.522	1089	1094	1107	rVB	85638	146405	0.50%	0.177%
11	9.728	1121	1129	1131	rBV	954759	1518907	5.18%	1.832%
12	9.757	1131	1134	1148	rVV	811241	1280424	4.37%	1.544%
13	9.869	1148	1153	1163	rVB3	31088	53855	0.18%	0.065%
14	9.993	1165	1174	1177	rBV	373256	795887	2.71%	0.960%
15	10.028	1177	1180	1188	rVB	512212	764281	2.61%	0.922%
16	10.181	1198	1206	1221	rVB	395260	649094	2.21%	0.783%
17	10.369	1231	1238	1241	rBV3	20808	34476	0.12%	0.042%
18	10.416	1241	1246	1249	rVV	203415	335050	1.14%	0.404%
19	10.463	1249	1254	1261	rVV	4360544	6818082	23.25%	8.223%
20	10.540	1261	1267	1274	rVV	2158021	3663482	12.49%	4.419%
21	10.593	1274	1276	1284	rVV	43884	70380	0.24%	0.085%
22	10.681	1285	1291	1300	rVB5	27784	58788	0.20%	0.071%
23	10.816	1309	1314	1320	rBV2	34806	53485	0.18%	0.065%
24	10.898	1320	1328	1340	rVV	1366113	2309192	7.88%	2.785%
25	10.993	1340	1344	1351	rVV	44479	72794	0.25%	0.088%
26	11.075	1351	1358	1367	rVV	140259	272480	0.93%	0.329%
27	11.157	1367	1372	1384	rVB	67219	127367	0.43%	0.154%
28	11.369	1394	1408	1426	rBV3	209870	496396	1.69%	0.599%
29	11.563	1436	1441	1446	rBV2	106461	169615	0.58%	0.205%
30	11.622	1446	1451	1455	rVV	89573	156456	0.53%	0.189%
31	11.663	1455	1458	1464	rVB2	39498	65821	0.22%	0.079%
32	11.892	1483	1497	1517	rBV	17620473	29322589	100.00%	35.366%
33	12.034	1517	1521	1530	rVB3	32413	84877	0.29%	0.102%
34	12.151	1537	1541	1547	rVB2	32467	50419	0.17%	0.061%
35	12.340	1569	1573	1578	rVB2	21346	32543	0.11%	0.039%
36	12.587	1610	1615	1619	rBV6	18564	37908	0.13%	0.046%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023977.D
Acq On : 06 Feb 2025 19:22
Operator : RC/JU
Sample : Q1202-19DL2 150X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B3DL2

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-BP012925.MA.M
Title : SVOA CALIBRATION

37	12.663	1624	1628	1633	rVV4	17650	33947	0.12%	0.041%
38	12.763	1633	1645	1649	rVV10	20417	78094	0.27%	0.094%
39	13.228	1716	1724	1732	rBV	1885249	2689930	9.17%	3.244%
40	13.298	1732	1736	1743	rVB	81770	142019	0.48%	0.171%
41	13.592	1780	1786	1790	rBV	487061	662761	2.26%	0.799%
42	13.639	1790	1794	1801	rVV	197271	301138	1.03%	0.363%
43	13.710	1801	1806	1812	rVB	43544	74794	0.26%	0.090%
44	13.792	1814	1820	1825	rVB2	40938	61706	0.21%	0.074%
45	14.081	1862	1869	1877	rVB2	40364	67448	0.23%	0.081%
46	14.245	1893	1897	1903	rBV	42864	62774	0.21%	0.076%
47	14.392	1913	1922	1934	rBV	3587930	5026641	17.14%	6.063%
48	15.075	2033	2038	2043	rBV	26981	45218	0.15%	0.055%
49	15.398	2087	2093	2099	rBV2	52730	93379	0.32%	0.113%
50	15.851	2166	2170	2176	rVB	44781	61363	0.21%	0.074%
51	16.463	2265	2274	2282	rBV3	30816	69982	0.24%	0.084%
52	17.180	2389	2396	2408	rBV2	3715166	5869495	20.02%	7.079%
53	17.280	2408	2413	2418	rVB2	62999	95276	0.32%	0.115%
54	19.651	2812	2816	2821	rBV	65721	90509	0.31%	0.109%
55	21.627	3146	3152	3163	rVB2	3775596	6152082	20.98%	7.420%
56	24.998	3717	3725	3740	rVB2	2269150	6178666	21.07%	7.452%

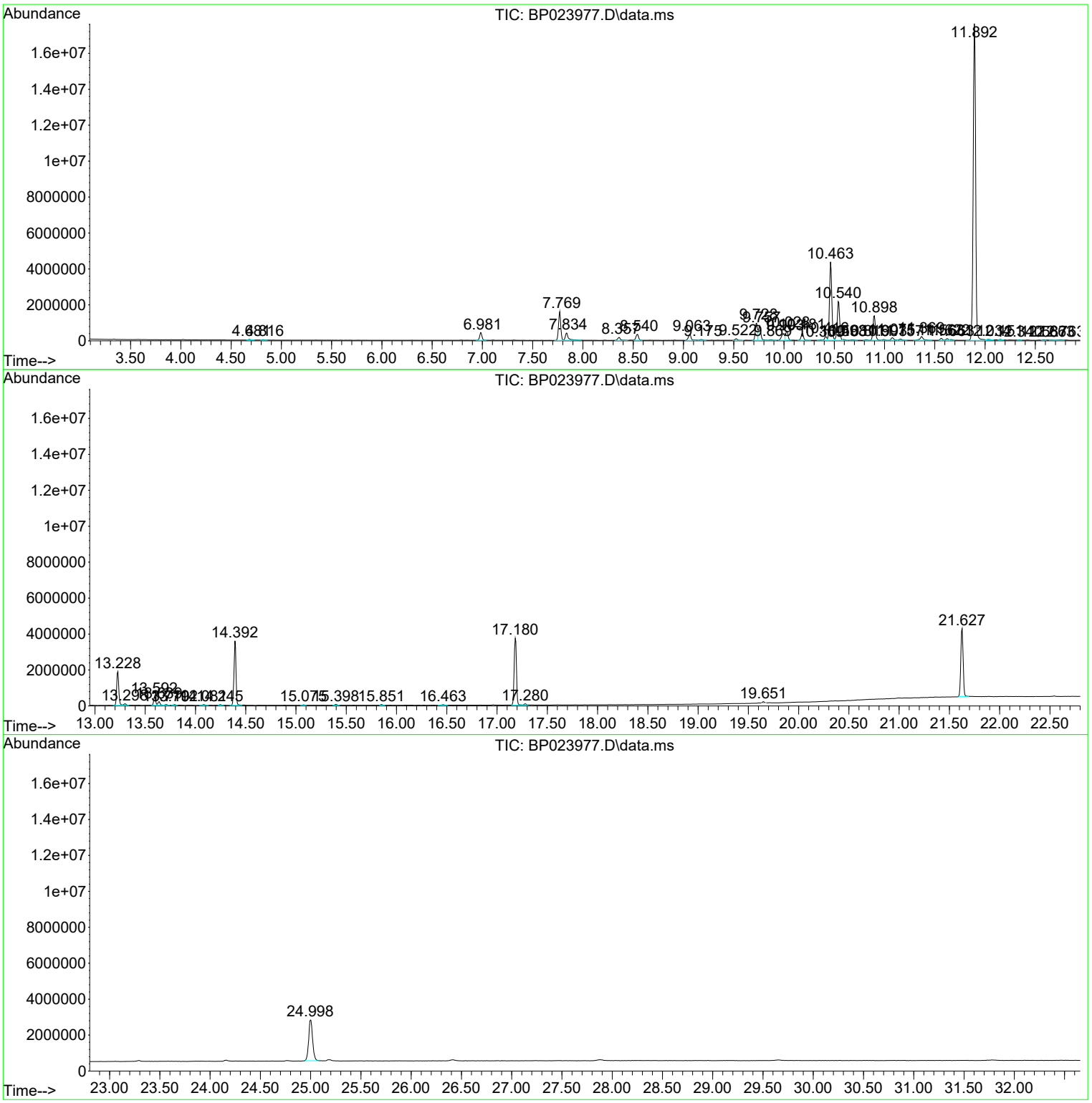
Sum of corrected areas: 82910643

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023977.D
Acq On : 06 Feb 2025 19:22
Operator : RC/JU
Sample : Q1202-19DL2 150X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B3DL2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.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\BP020525\
Data File : BP023977.D
Acq On : 06 Feb 2025 19:22
Operator : RC/JU
Sample : Q1202-19DL2 150X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B3DL2

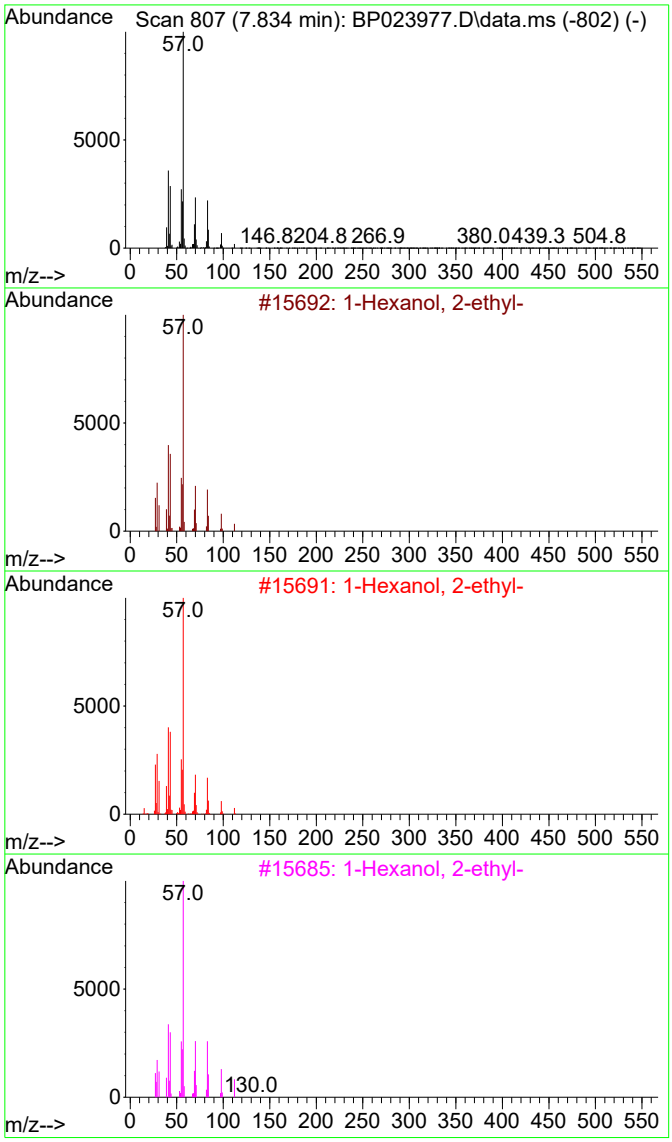
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

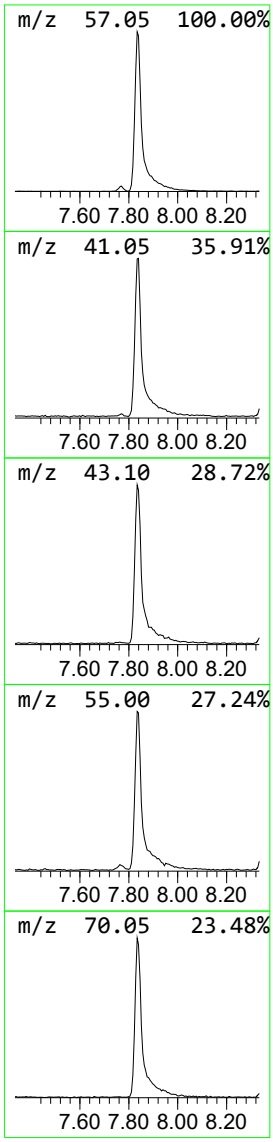
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.834	7.25 ng/ul	858340	1,4-Dichlorobenzene-d4	7.769

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	83
4	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	74
5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023977.D
Acq On : 06 Feb 2025 19:22
Operator : RC/JU
Sample : Q1202-19DL2 150X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B3DL2

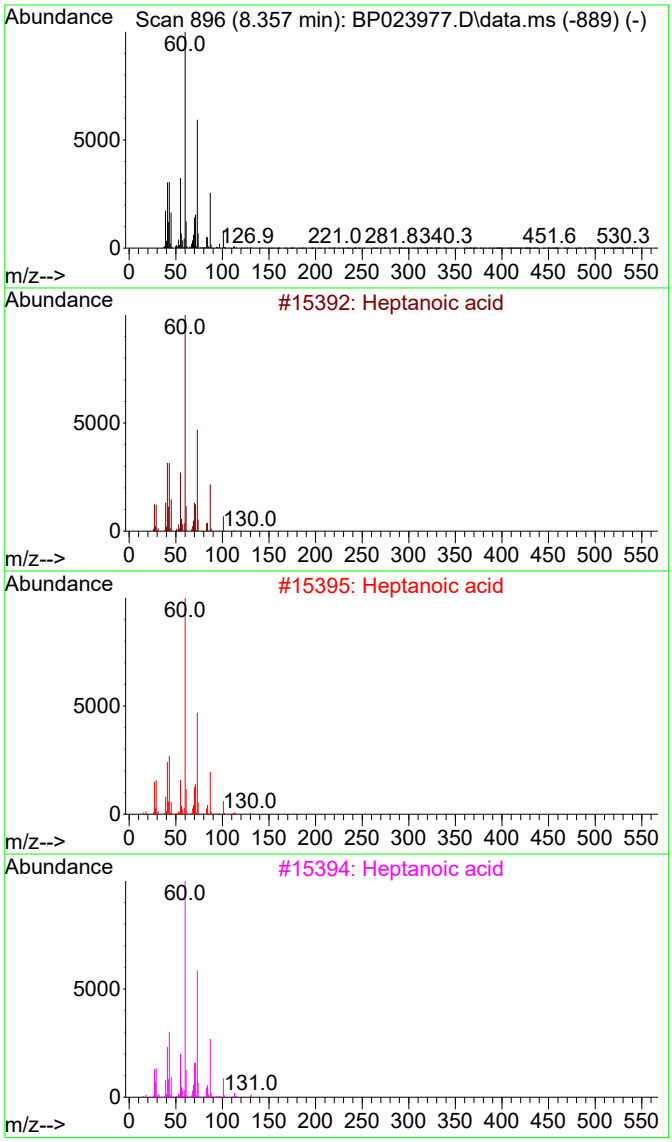
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

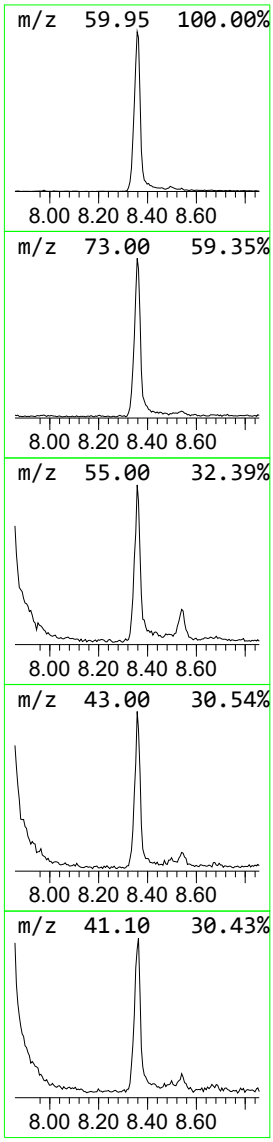
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Heptanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.357	2.62 ng/ul	309888	1,4-Dichlorobenzene-d4	7.769

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptanoic acid	130	C7H14O2	000111-14-8	91
2	Heptanoic acid	130	C7H14O2	000111-14-8	86
3	Heptanoic acid	130	C7H14O2	000111-14-8	78
4	Heptanoic acid	130	C7H14O2	000111-14-8	78
5	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023977.D
Acq On : 06 Feb 2025 19:22
Operator : RC/JU
Sample : Q1202-19DL2 150X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B3DL2

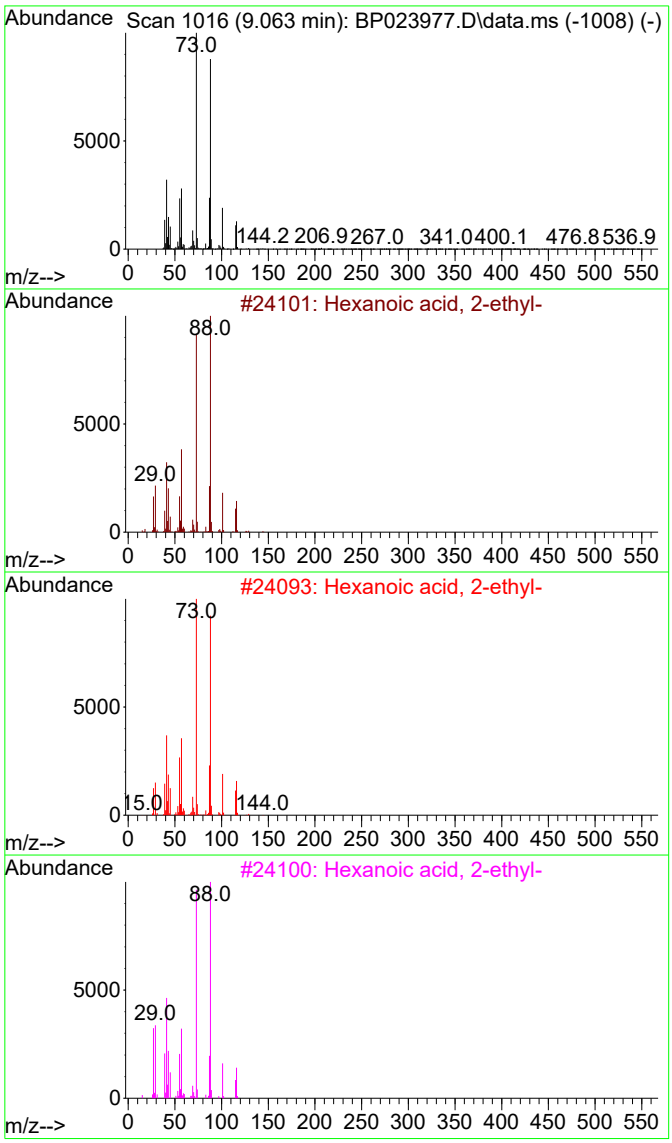
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

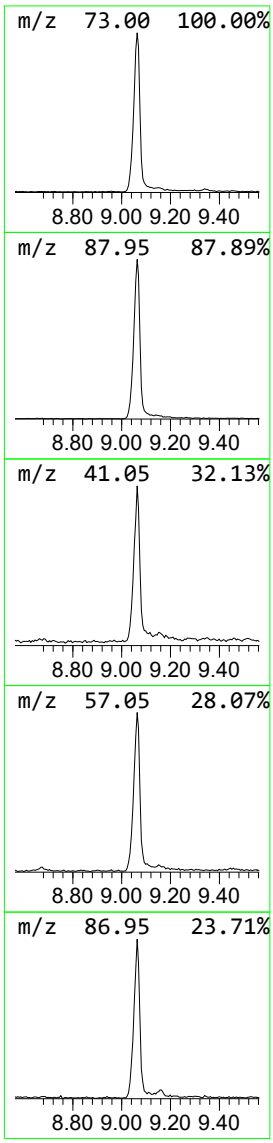
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Hexanoic acid, 2-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.063	5.26 ng/ul	622145	1,4-Dichlorobenzene-d4	7.769

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	93
2	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	91
3	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
4	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	83
5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023977.D
Acq On : 06 Feb 2025 19:22
Operator : RC/JU
Sample : Q1202-19DL2 150X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B3DL2

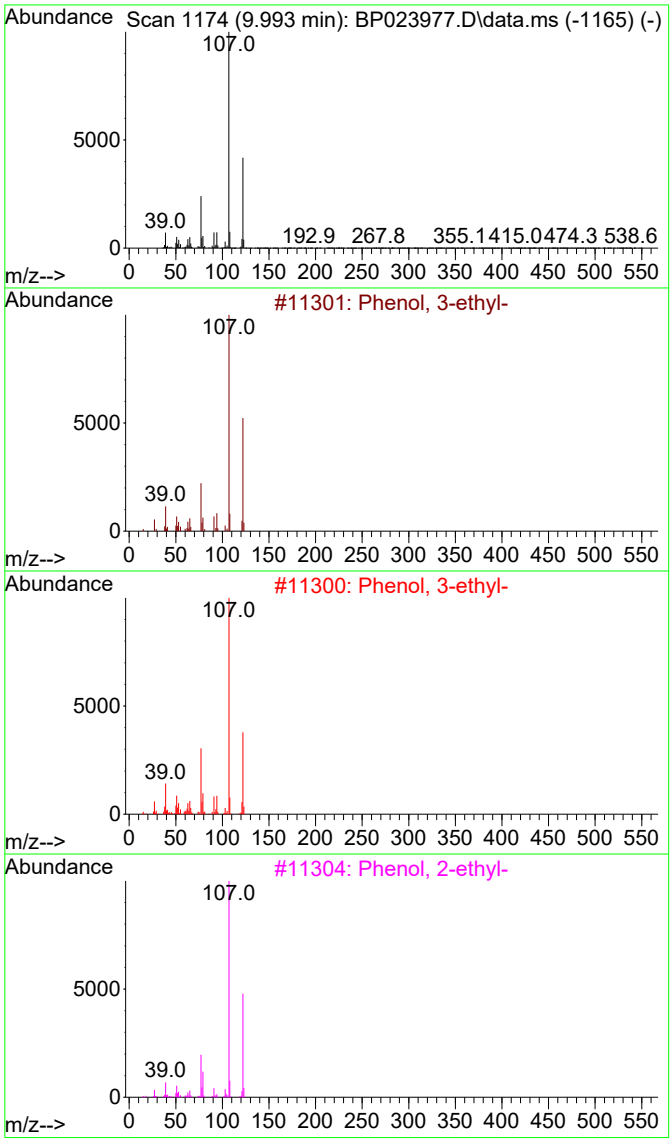
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

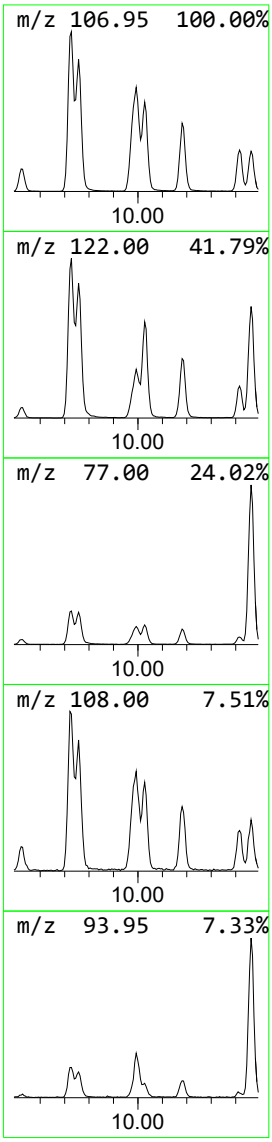
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Phenol, 3-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.993	4.34 ng/ul	795887	Naphthalene-d8	10.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	97
2	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	94
3	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	91
4	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	91
5	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023977.D
Acq On : 06 Feb 2025 19:22
Operator : RC/JU
Sample : Q1202-19DL2 150X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B3DL2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

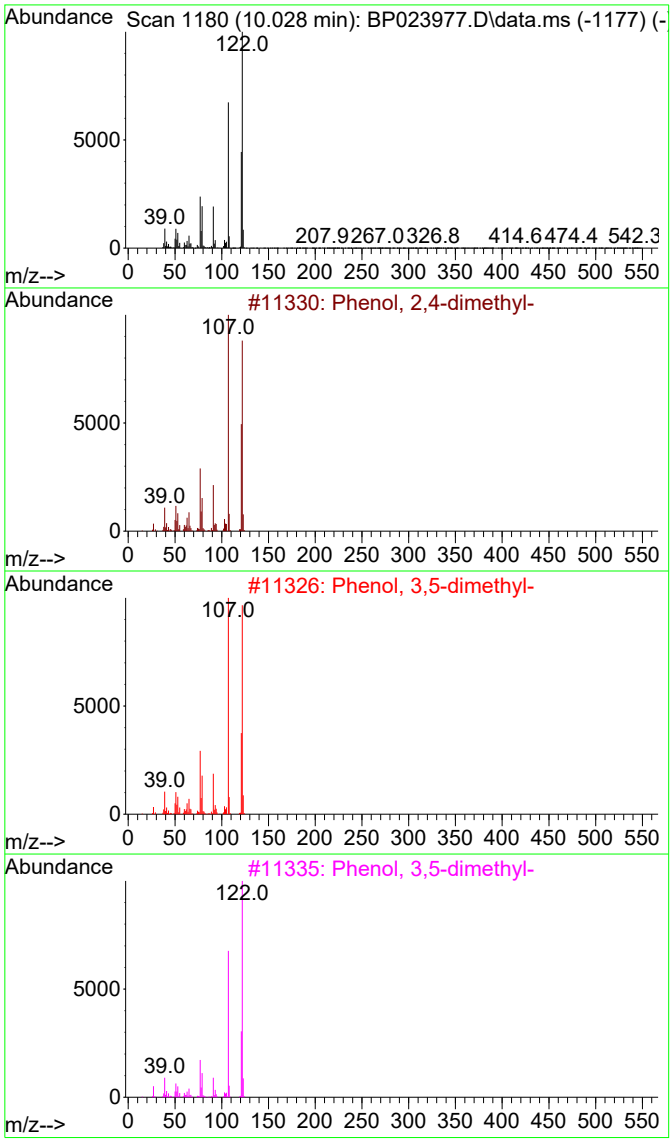
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

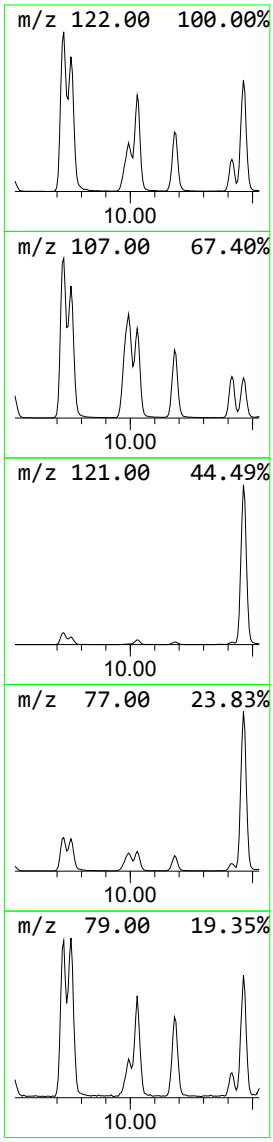
****

Peak Number 5 Phenol, 3,5-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.028	4.17 ng/ul	764281	Naphthalene-d8	10.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,4-dimethyl-	122	C8H10O	000105-67-9	91
2	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	91
3	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	90
4	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	87
5	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023977.D
Acq On : 06 Feb 2025 19:22
Operator : RC/JU
Sample : Q1202-19DL2 150X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B3DL2

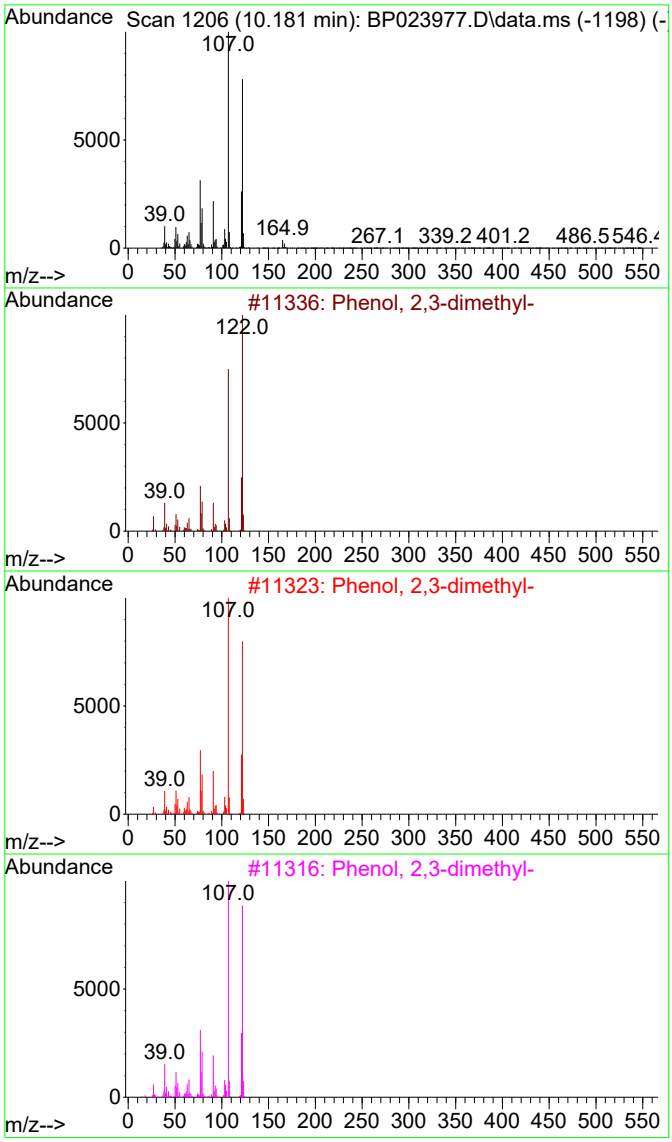
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

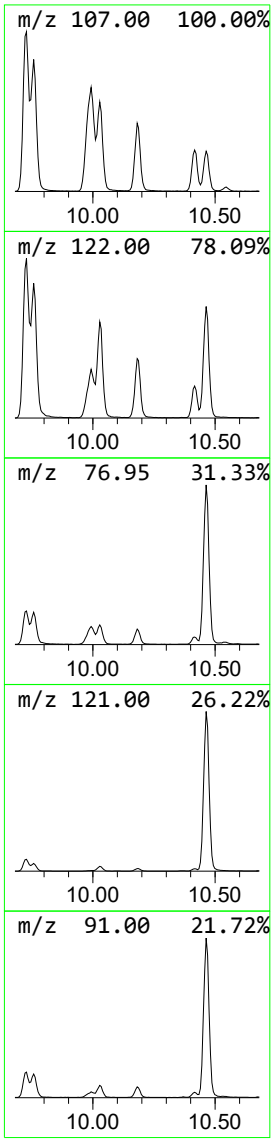
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Phenol, 2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.181	3.54 ng/ul	649094	Naphthalene-d8	10.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	97
2	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	96
3	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	96
4	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	94
5	Phenol, 2,4-dimethyl-	122	C8H10O	000105-67-9	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023977.D
Acq On : 06 Feb 2025 19:22
Operator : RC/JU
Sample : Q1202-19DL2 150X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B3DL2

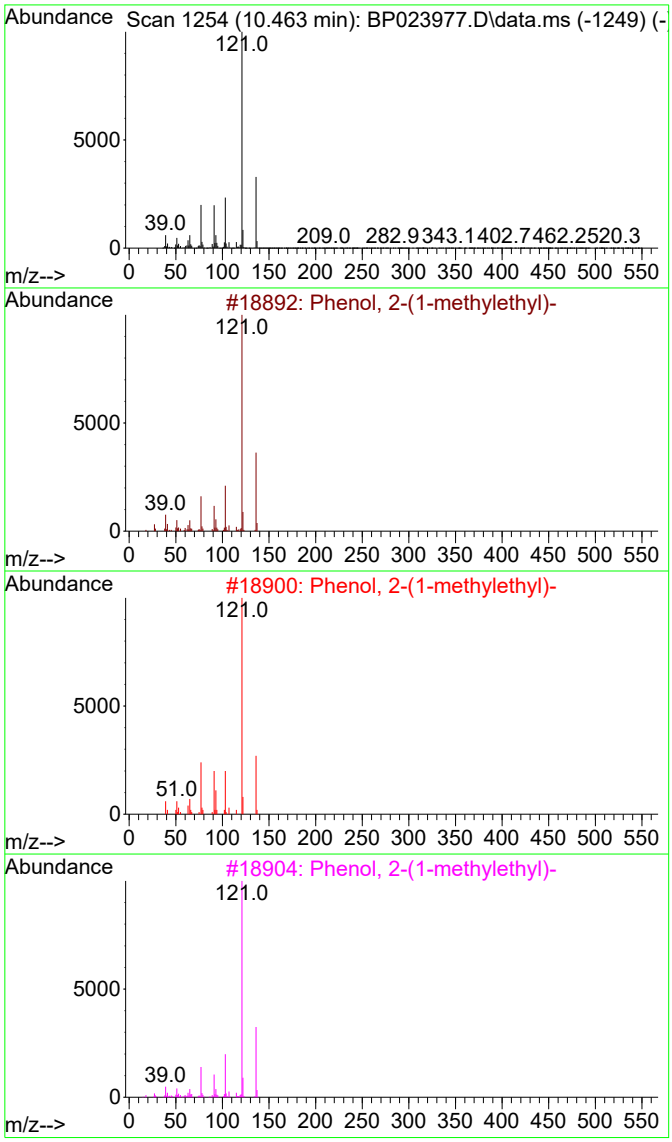
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

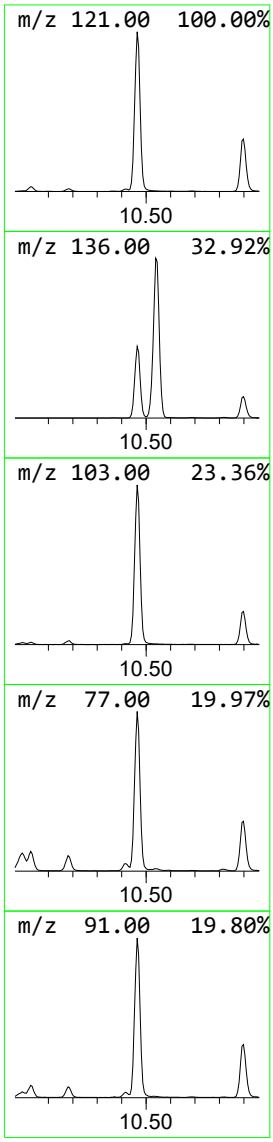
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Phenol, 2-(1-methylethyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.463	37.22 ng/ul	6818080	Naphthalene-d8	10.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	95
2	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	94
3	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	91
4	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	86
5	p-Cumenol	136	C9H12O	000099-89-8	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023977.D
Acq On : 06 Feb 2025 19:22
Operator : RC/JU
Sample : Q1202-19DL2 150X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B3DL2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

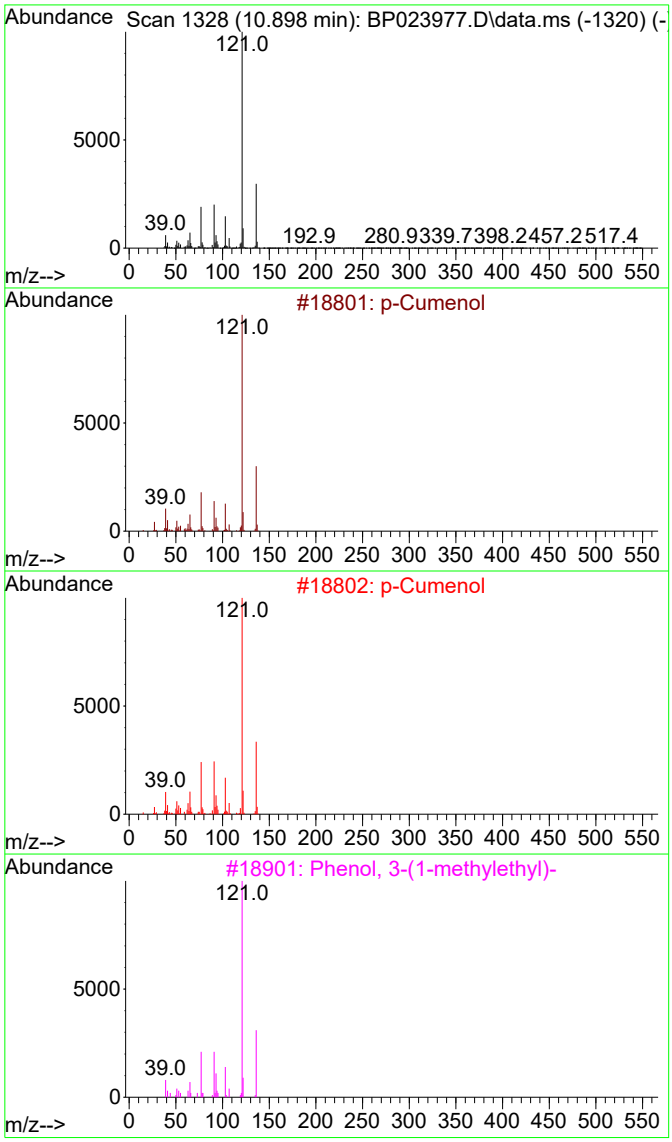
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

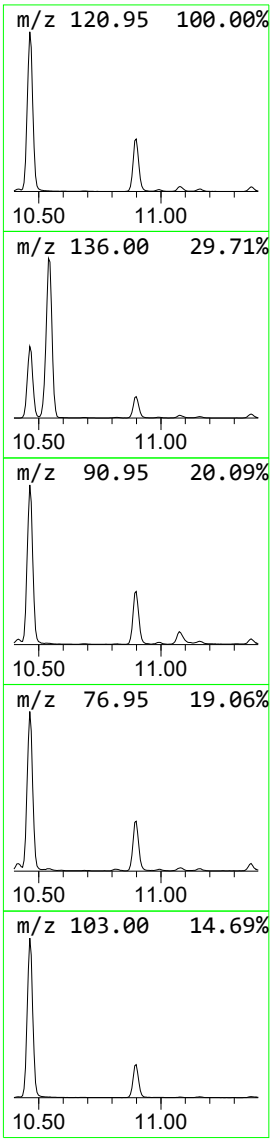
****

Peak Number 8 p-Cumenol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.898	12.61 ng/ul	2309190	Naphthalene-d8	10.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cumenol	136	C9H12O	000099-89-8	97
2	p-Cumenol	136	C9H12O	000099-89-8	96
3	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	94
4	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	91
5	p-Cumenol	136	C9H12O	000099-89-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023977.D
Acq On : 06 Feb 2025 19:22
Operator : RC/JU
Sample : Q1202-19DL2 150X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B3DL2

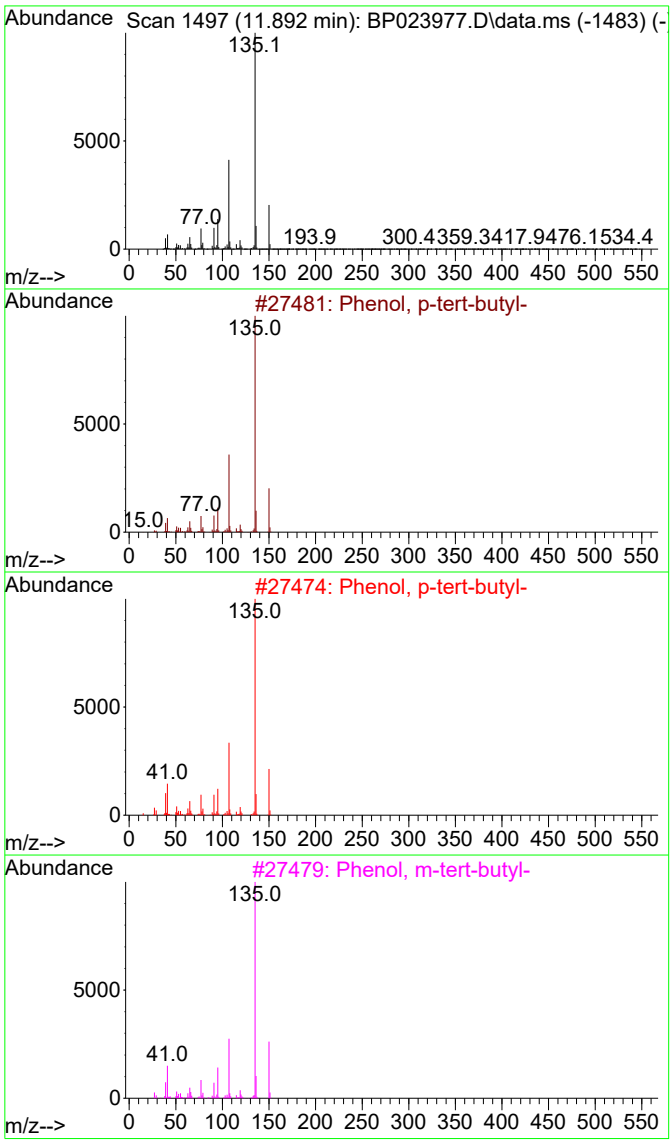
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

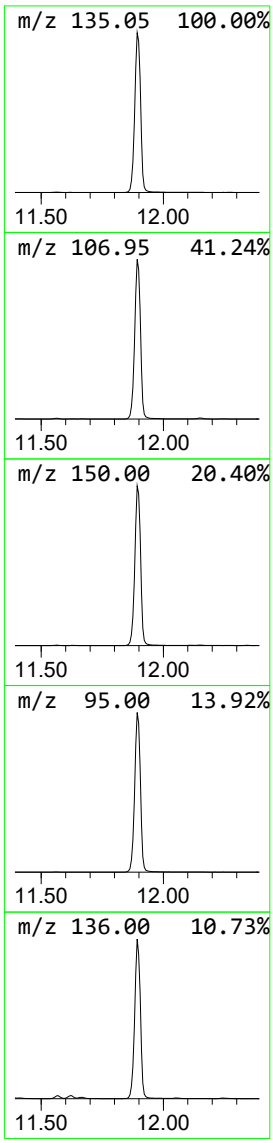
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Phenol, p-tert-butyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.893	160.08 ng/ul	29322600	Naphthalene-d8	10.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	96
2	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	96
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	94
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	94
5	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023977.D
Acq On : 06 Feb 2025 19:22
Operator : RC/JU
Sample : Q1202-19DL2 150X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B3DL2

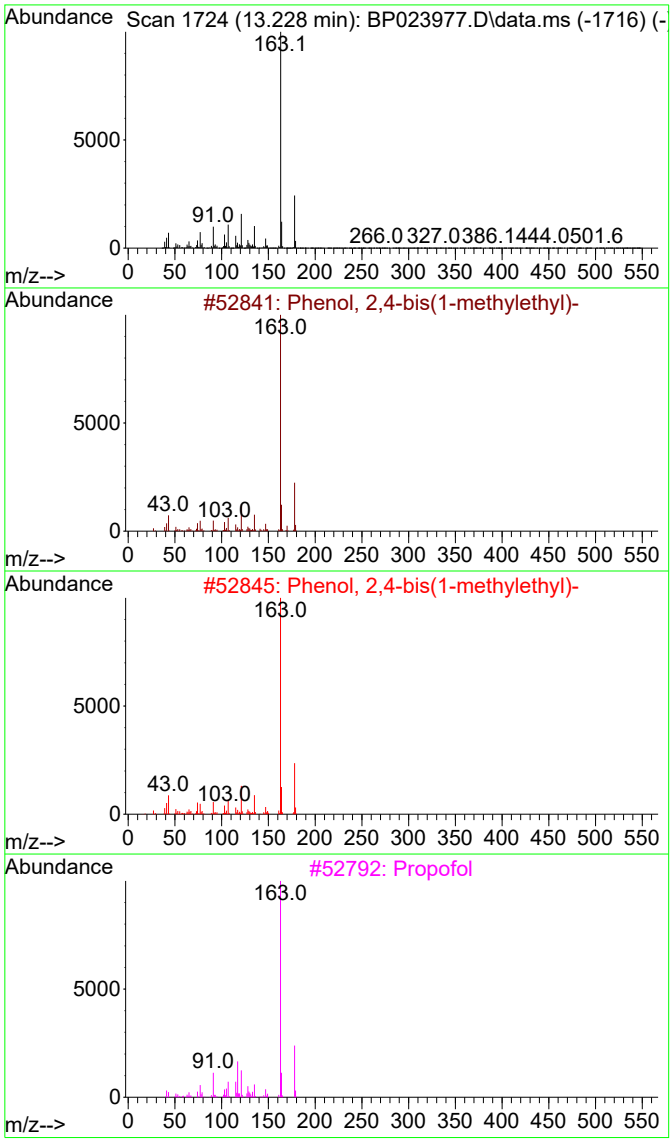
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

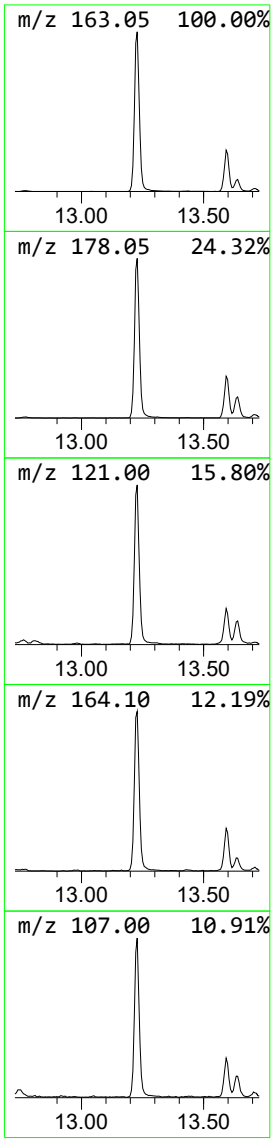
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Phenol, 2,4-bis(1-methyleth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.228	10.70 ng/ul	2689930	Acenaphthene-d10	14.392

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	97
2	Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	95
3	Propofol	178	C12H18O	002078-54-8	94
4	Propofol	178	C12H18O	002078-54-8	93
5	Propofol	178	C12H18O	002078-54-8	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023977.D
Acq On : 06 Feb 2025 19:22
Operator : RC/JU
Sample : Q1202-19DL2 150X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B3DL2

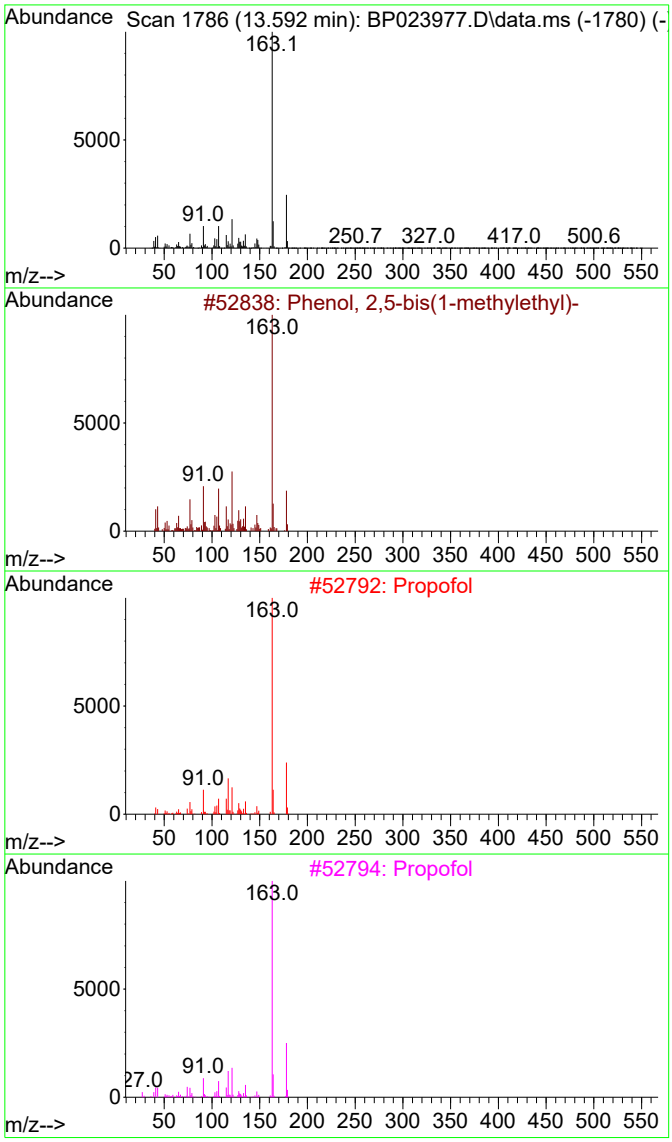
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

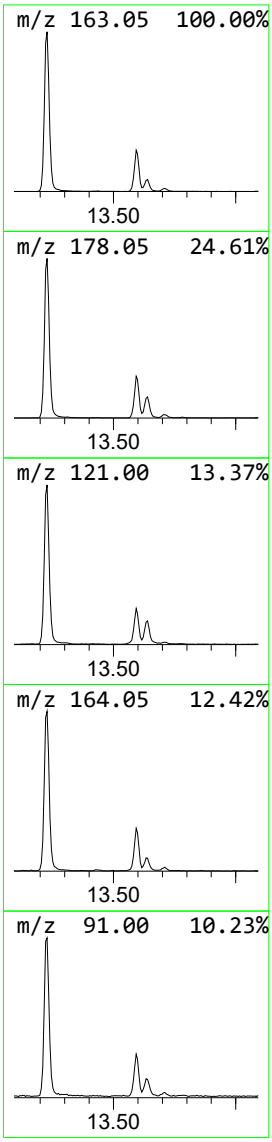
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Phenol, 2,5-bis(1-methyleth... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.592	2.64 ng/ul	662761	Acenaphthene-d10	14.392

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,5-bis(1-methylethyl)-	178	C12H18O	035946-91-9	95
2	Propofol	178	C12H18O	002078-54-8	94
3	Propofol	178	C12H18O	002078-54-8	94
4	Propofol	178	C12H18O	002078-54-8	90
5	Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023977.D
Acq On : 06 Feb 2025 19:22
Operator : RC/JU
Sample : Q1202-19DL2 150X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B3DL2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.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
1-Hexanol, 2-et...	7.834	7.3	ng/ul	858340	1	7.769	2366990	20.0
Heptanoic acid	8.357	2.6	ng/ul	309888	1	7.769	2366990	20.0
Hexanoic acid, ...	9.063	5.3	ng/ul	622145	1	7.769	2366990	20.0
Phenol, 3-ethyl-	9.993	4.3	ng/ul	795887	2	10.540	3663480	20.0
Phenol, 3,5-dim...	10.028	4.2	ng/ul	764281	2	10.540	3663480	20.0
Phenol, 2,3-dim...	10.181	3.5	ng/ul	649094	2	10.540	3663480	20.0
Phenol, 2-(1-me...	10.463	37.2	ng/ul	6818080	2	10.540	3663480	20.0
p-Cumenol	10.898	12.6	ng/ul	2309190	2	10.540	3663480	20.0
Phenol, p-tert-...	11.893	160.1	ng/ul	29322600	2	10.540	3663480	20.0
Phenol, 2,4-bis...	13.228	10.7	ng/ul	2689930	3	14.392	5026640	20.0
Phenol, 2,5-bis...	13.592	2.6	ng/ul	662761	3	14.392	5026640	20.0