LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023974.D
Acq On : 06 Feb 2025 17:21
Operator : RC/JU
Sample : Q1202-20DL 30X
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B4DL

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: BP023974.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.981	649	662	668	rBV2	45451	105281	0.16%	0.068%
2	7.105	677	683	691	rVB	74877	120653	0.18%	0.078%
3	7.311	710	718	726	rBV	70577	119975	0.18%	0.077%
4	7.769	788	796	804	rBV	1539221	2401887	3.56%	1.547%
5	8.481	911	917	918	rBV	73629	103528	0.15%	0.067%
6	8.540	922	927	940	rVV	352294	631576	0.94%	0.407%
7	8.911	985	990	999	rBV	78374	126280	0.19%	0.081%
8	9.175	1026	1035	1045	rVB2	99050	207476	0.31%	0.134%
9	9.522	1089	1094	1101	rVB	115919	171427	0.25%	0.110%
10	9.628	1106	1112	1119	rVB2	49905	92263	0.14%	0.059%
11	9.734	1119	1130	1132	rBV	1254914	2071064	3.07%	1.334%
12	9.763	1132	1135	1151	rVB	1824827	2704804	4.01%	1.742%
13	9.993	1164	1174	1178	rBV	930380	2183853	3.23%	1.406%
14	10.034	1178	1181	1192	rVB	910714	1397663	2.07%	0.900%
15	10.187	1199	1207	1218	rVB2	706056	1160448	1.72%	0.747%
16	10.375	1233	1239	1242	rBV2	48374	81804	0.12%	0.053%
17	10.422	1242	1247	1250	rVV	396621	618175	0.92%	0.398%
18	10.475	1250	1256	1262	rVV	11911414	16907352	25.04%	10.886%
19	10.540	1262	1267	1274	rVV	2423724	3807674	5.64%	2.452%
20	10.593	1274	1276	1284	rVB3	47735	94010	0.14%	0.061%
21	10.693	1284	1293	1299	rBV3	77067	165710	0.25%	0.107%
22	10.822	1309	1315	1320	rBV2	83648	150547	0.22%	0.097%
23	10.899	1321	1328	1340	rVV	4673866	7195588	10.66%	4.633%
24	10.999	1340	1345	1349	rVV	81528	125590	0.19%	0.081%
25	11.093	1355	1361	1368	rBV2	190340	347804	0.52%	0.224%
26	11.163	1368	1373	1379	rVB	194132	285182	0.42%	0.184%
27	11.299	1390	1396	1399	rBV3	54143	112322	0.17%	0.072%
28	11.381	1402	1410	1415	rVV	589376	1174513	1.74%	0.756%
29	11.416	1415	1416	1421	rVB2	124499	134612	0.20%	0.087%
30	11.563	1430	1441	1448	rBV2	766657	1562400	2.31%	1.006%
31	11.628	1448	1452	1475	rVB2	284773	905686	1.34%	0.583%
32	11.910	1488	1500	1519	rBV2	31785321	67524626	100.00%	43.478%
33	12.104	1528	1533	1538	rVV6	39887	107782	0.16%	0.069%
34	12.157	1538	1542	1548	rVB3	74633	121514	0.18%	0.078%
35	12.257	1554	1559	1565	rBV3	39217	79096	0.12%	0.051%
36	12.587	1611	1615	1620	rBV3	43343	77857	0.12%	0.050%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023974.D
Acq On : 06 Feb 2025 17:21
Operator : RC/JU
Sample : Q1202-20DL 30X
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B4DL

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	13.240	1718	1726	1741	rBV	5619539	8973079	13.29%	5.778%
38	13.604	1782	1788	1792	rVV	2216114	3119702	4.62%	2.009%
39	13.646	1792	1795	1801	rVV	1290290	1556574	2.31%	1.002%
40	13.710	1801	1806	1813	rVV	267686	412085	0.61%	0.265%
41	13.793	1814	1820	1825	rVB2	211007	324660	0.48%	0.209%
42	14.081	1860	1869	1878	rVB2	212064	389347	0.58%	0.251%
43	14.393	1916	1922	1930	rBV	3380222	5052030	7.48%	3.253%
44	14.646	1958	1965	1970	rBV5	44735	92244	0.14%	0.059%
45	15.040	2026	2032	2036	rBV	175478	243019	0.36%	0.156%
46	15.087	2036	2040	2055	rVB3	43042	104826	0.16%	0.067%
47	15.404	2089	2094	2103	rBV2	321631	536901	0.80%	0.346%
48	17.198	2393	2399	2410	rBV2	4004165	5839734	8.65%	3.760%
49	17.298	2412	2416	2431	rVB	331344	504605	0.75%	0.325%
50	17.592	2461	2466	2477	rBV	245649	394040	0.58%	0.254%
51	19.669	2814	2819	2826	rBV	367802	513540	0.76%	0.331%
52	21.639	3148	3154	3165	rVB2	3607522	6023766	8.92%	3.879%
53	25.016	3720	3728	3745	rVB	2150229	6050853	8.96%	3.896%

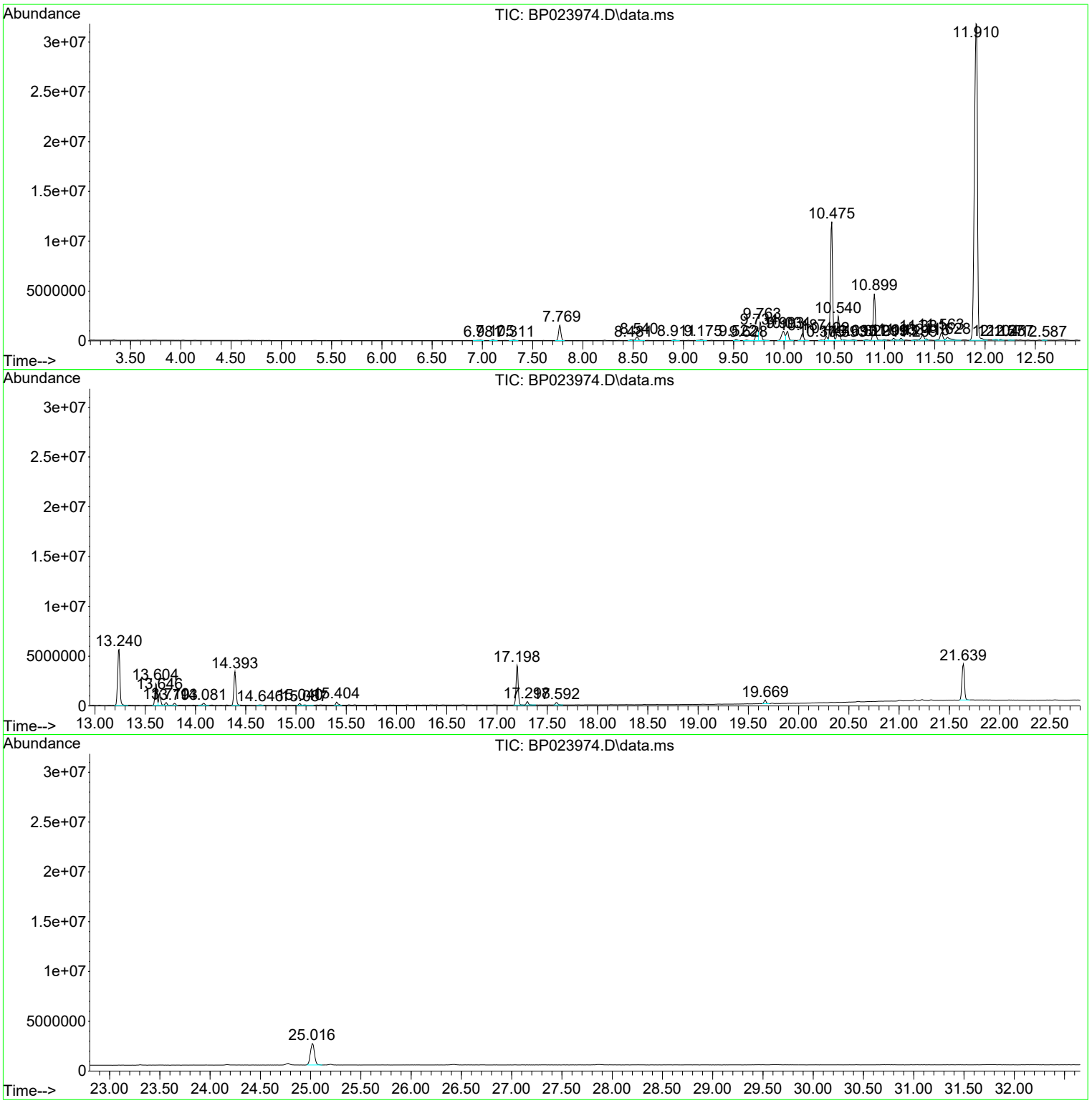
Sum of corrected areas: 155309027

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023974.D
Acq On : 06 Feb 2025 17:21
Operator : RC/JU
Sample : Q1202-20DL 30X
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B4DL

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 : BP023974.D
Acq On : 06 Feb 2025 17:21
Operator : RC/JU
Sample : Q1202-20DL 30X
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B4DL

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 Phenol, 3-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.993	11.47 ng/ul	2183850	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, 3-ethyl-	122	C8H10O	000620-17-7	91
4	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	91
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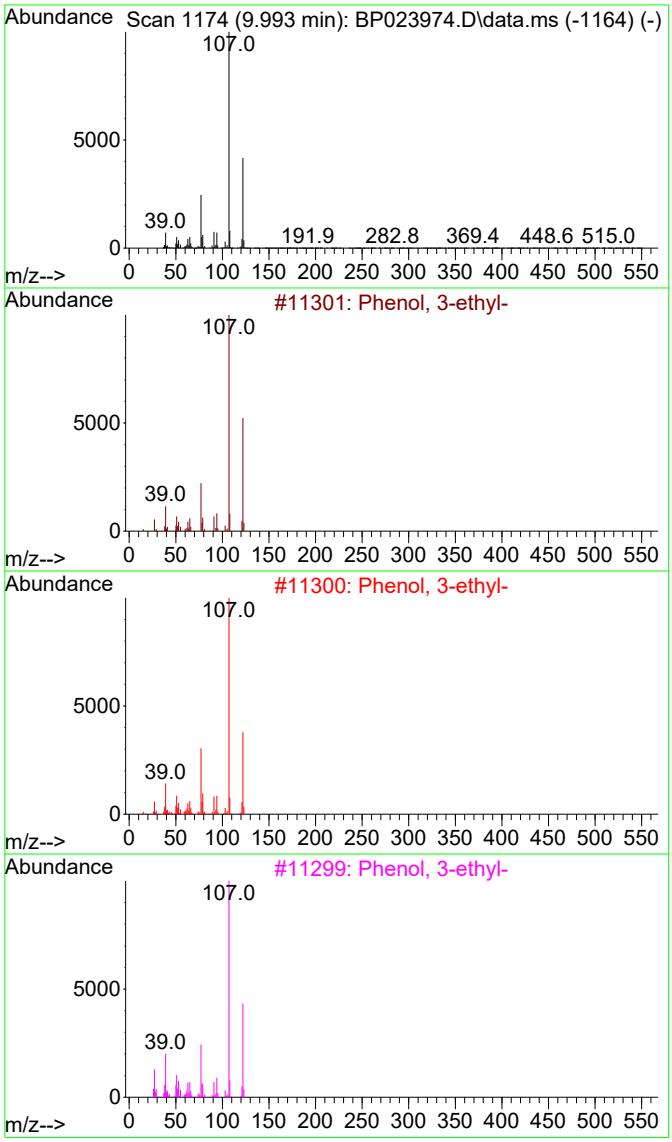
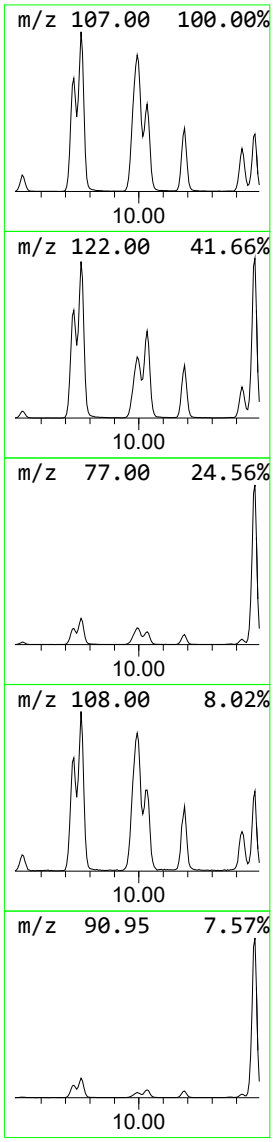




Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023974.D
Acq On : 06 Feb 2025 17:21
Operator : RC/JU
Sample : Q1202-20DL 30X
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B4DL

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

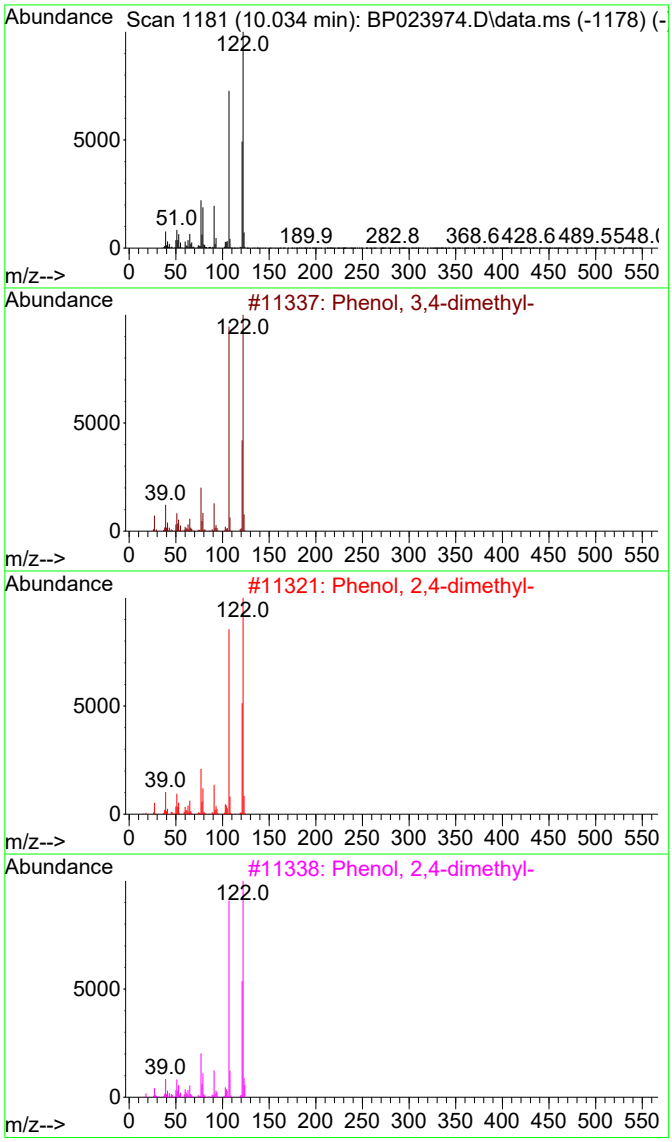
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

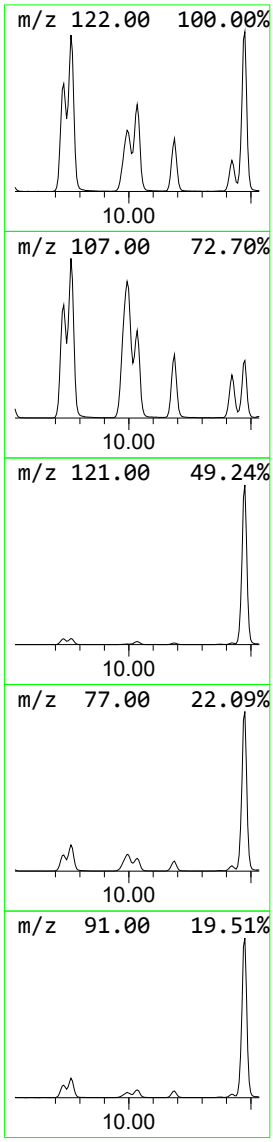
****

Peak Number 2 Phenol, 3,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.034	7.34 ng/ul	1397660	Naphthalene-d8	10.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	91
2	Phenol, 2,4-dimethyl-	122	C8H10O	000105-67-9	90
3	Phenol, 2,4-dimethyl-	122	C8H10O	000105-67-9	90
4	Phenol, 2,5-dimethyl-	122	C8H10O	000095-87-4	90
5	Phenol, 2,4-dimethyl-	122	C8H10O	000105-67-9	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023974.D
Acq On : 06 Feb 2025 17:21
Operator : RC/JU
Sample : Q1202-20DL 30X
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B4DL

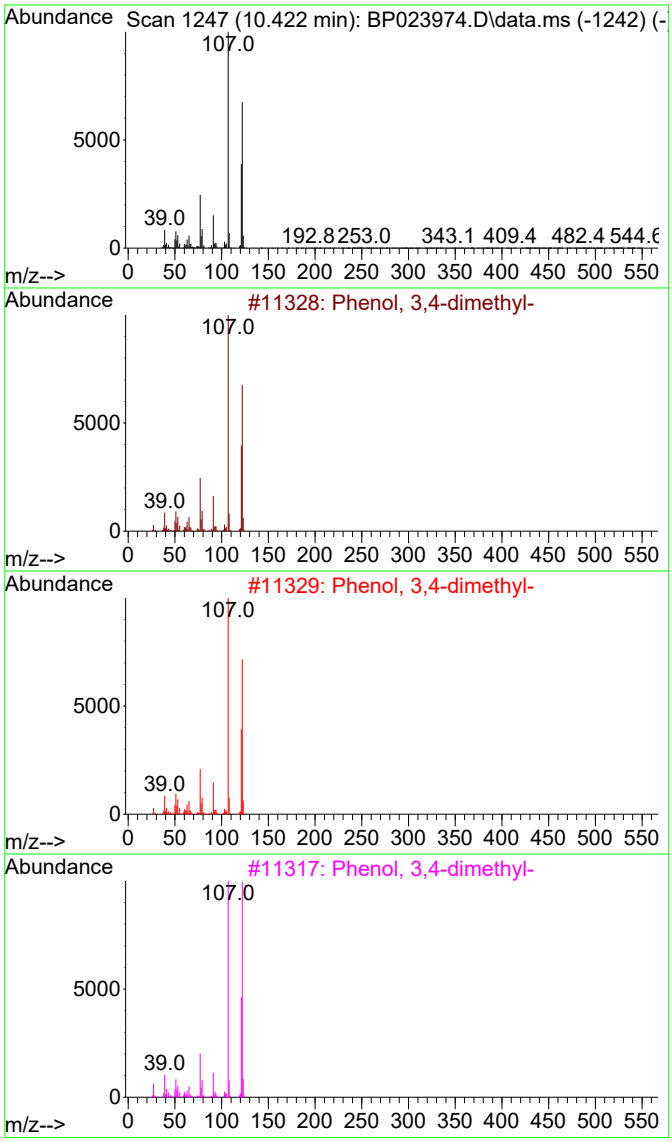
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

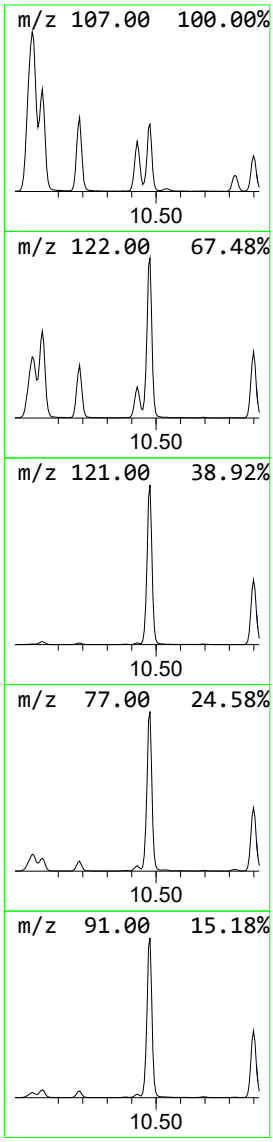
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Phenol, 3,5-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.422	3.25 ng/ul	618175	Naphthalene-d8	10.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	96
2	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	96
3	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	95
4	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	94
5	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023974.D
Acq On : 06 Feb 2025 17:21
Operator : RC/JU
Sample : Q1202-20DL 30X
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B4DL

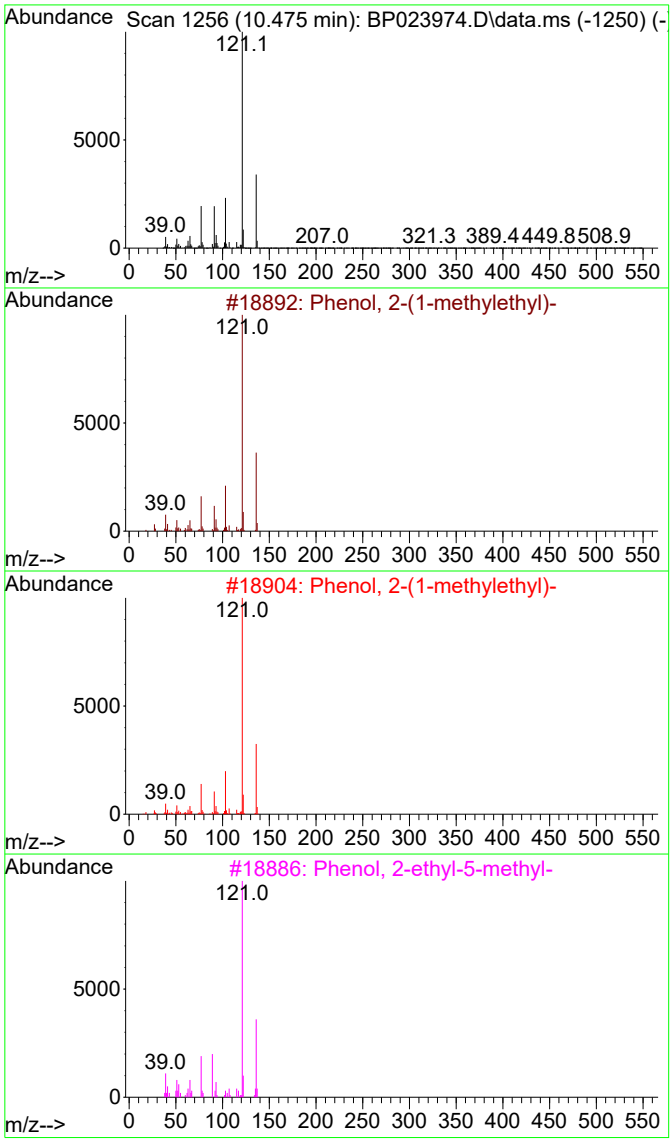
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

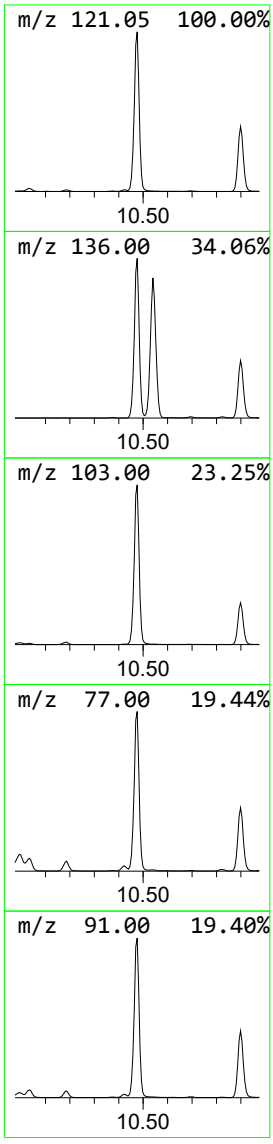
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Phenol, 2-(1-methylethyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.475	88.81 ng/ul	16907400	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	95
3	Phenol, 2-ethyl-5-methyl-	136	C9H12O	001687-61-2	87
4	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	87
5	p-Cumenol	136	C9H12O	000099-89-8	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023974.D
Acq On : 06 Feb 2025 17:21
Operator : RC/JU
Sample : Q1202-20DL 30X
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B4DL

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

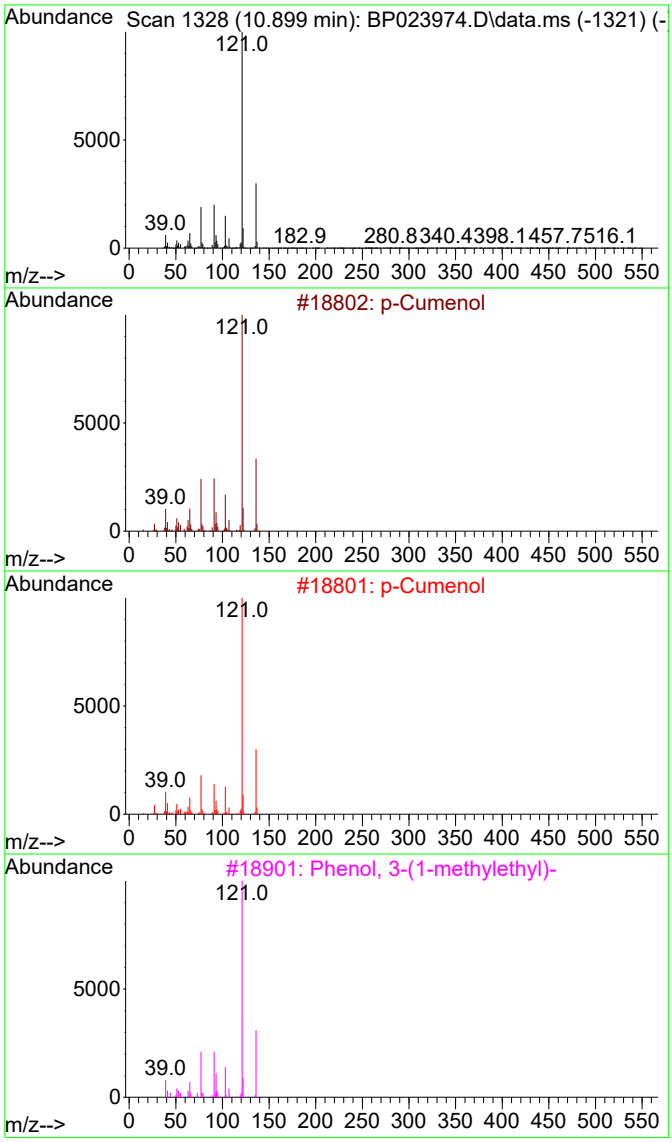
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

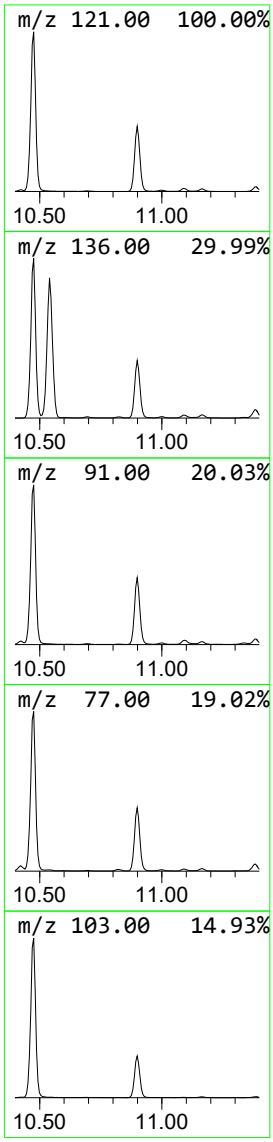
****

Peak Number 5 p-Cumenol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.899	37.80 ng/ul	7195590	Naphthalene-d8	10.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cumenol	136	C9H12O	000099-89-8	96
2	p-Cumenol	136	C9H12O	000099-89-8	95
3	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	94
4	p-Cumenol	136	C9H12O	000099-89-8	91
5	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023974.D
Acq On : 06 Feb 2025 17:21
Operator : RC/JU
Sample : Q1202-20DL 30X
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B4DL

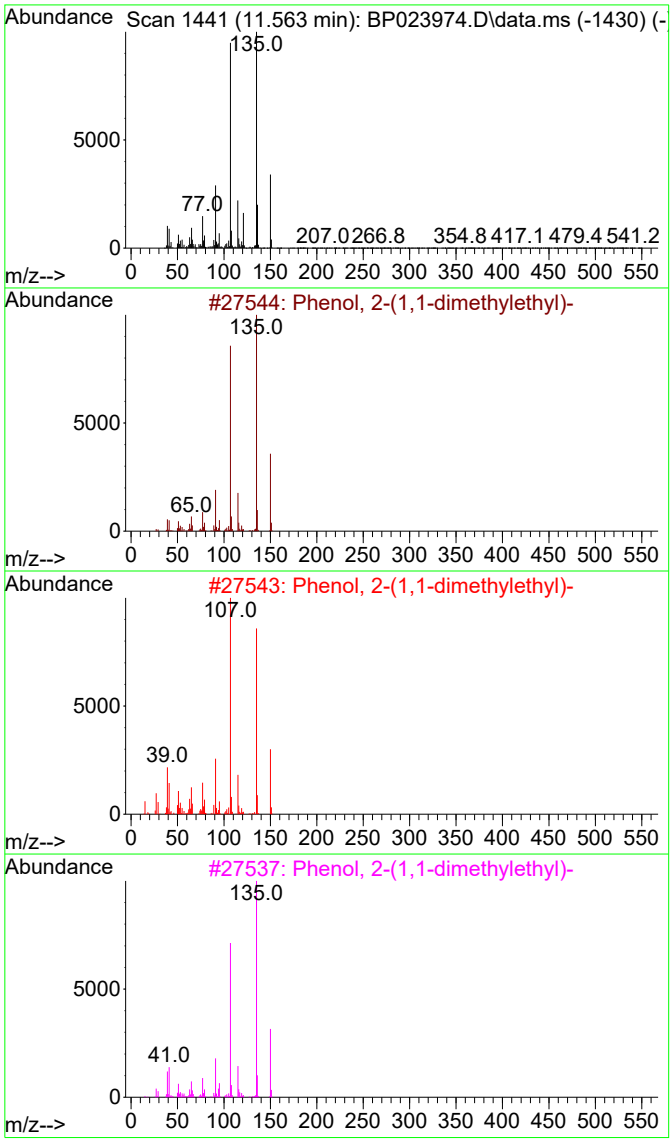
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

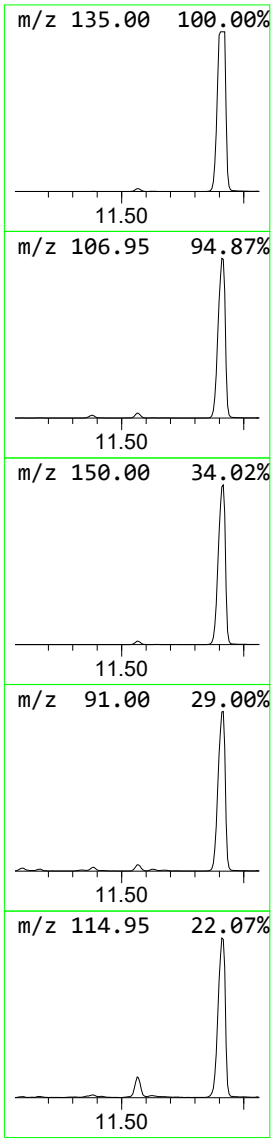
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.563	8.21 ng/ul	1562400	Naphthalene-d8	10.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	94
2	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	83
3	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	81
4	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	58
5	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023974.D
Acq On : 06 Feb 2025 17:21
Operator : RC/JU
Sample : Q1202-20DL 30X
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B4DL

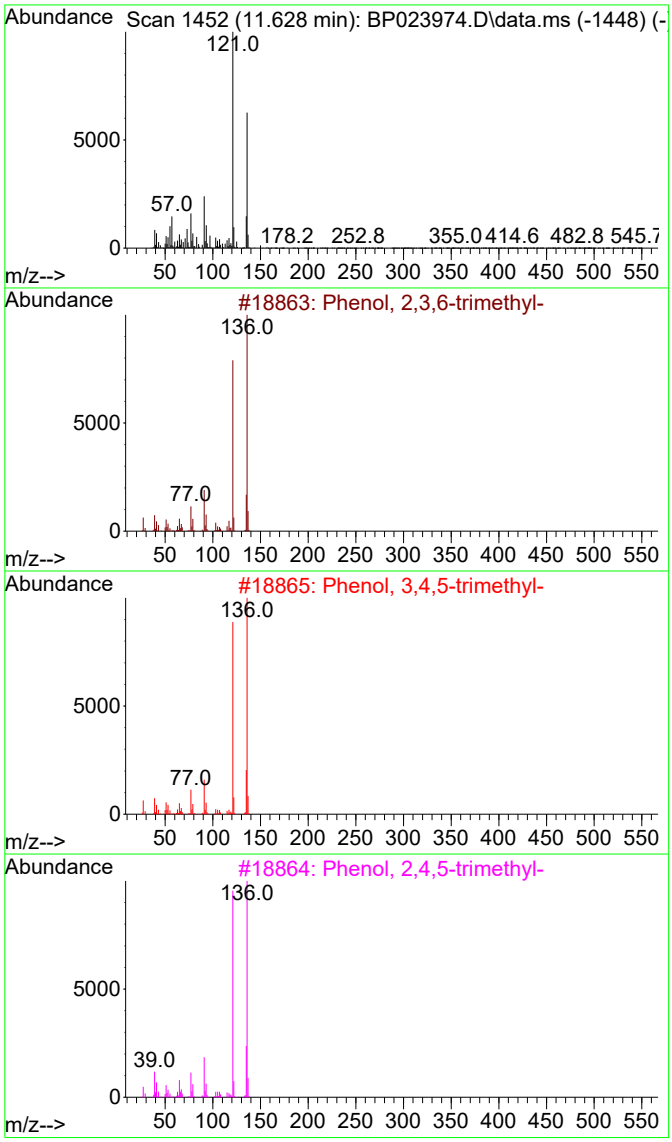
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

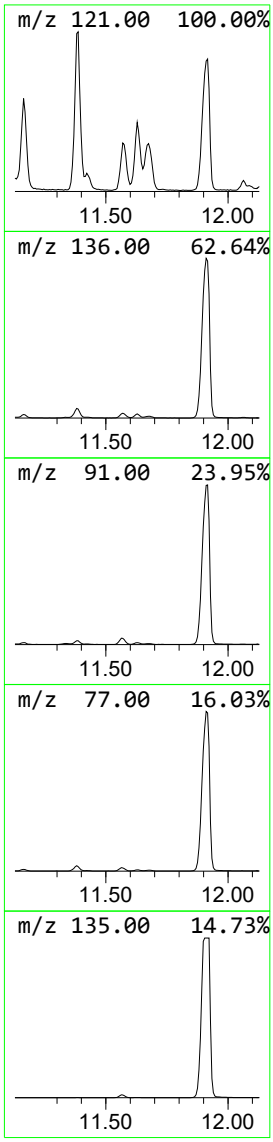
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Phenol, 2,3,6-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.628	4.76 ng/ul	905686	Naphthalene-d8	10.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	96
2	Phenol, 3,4,5-trimethyl-	136	C9H12O	000527-54-8	94
3	Phenol, 2,4,5-trimethyl-	136	C9H12O	000496-78-6	94
4	Phenol, 2,3,5-trimethyl-	136	C9H12O	000697-82-5	91
5	Phenol, 2,3,5-trimethyl-	136	C9H12O	000697-82-5	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023974.D
Acq On : 06 Feb 2025 17:21
Operator : RC/JU
Sample : Q1202-20DL 30X
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B4DL

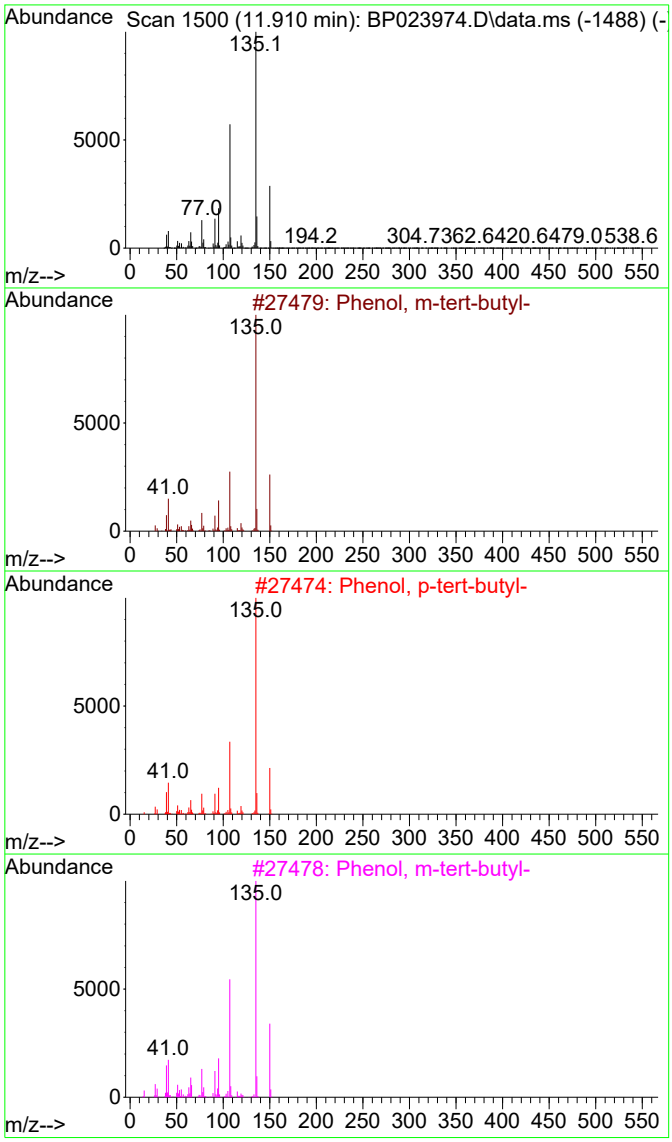
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

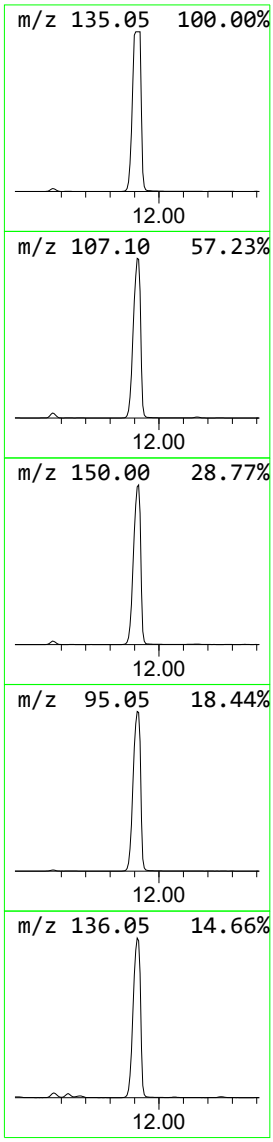
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Phenol, m-tert-butyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.910	354.68 ng/ul	67524600	Naphthalene-d8	10.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	94
2	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	93
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
4	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	90
5	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023974.D
Acq On : 06 Feb 2025 17:21
Operator : RC/JU
Sample : Q1202-20DL 30X
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B4DL

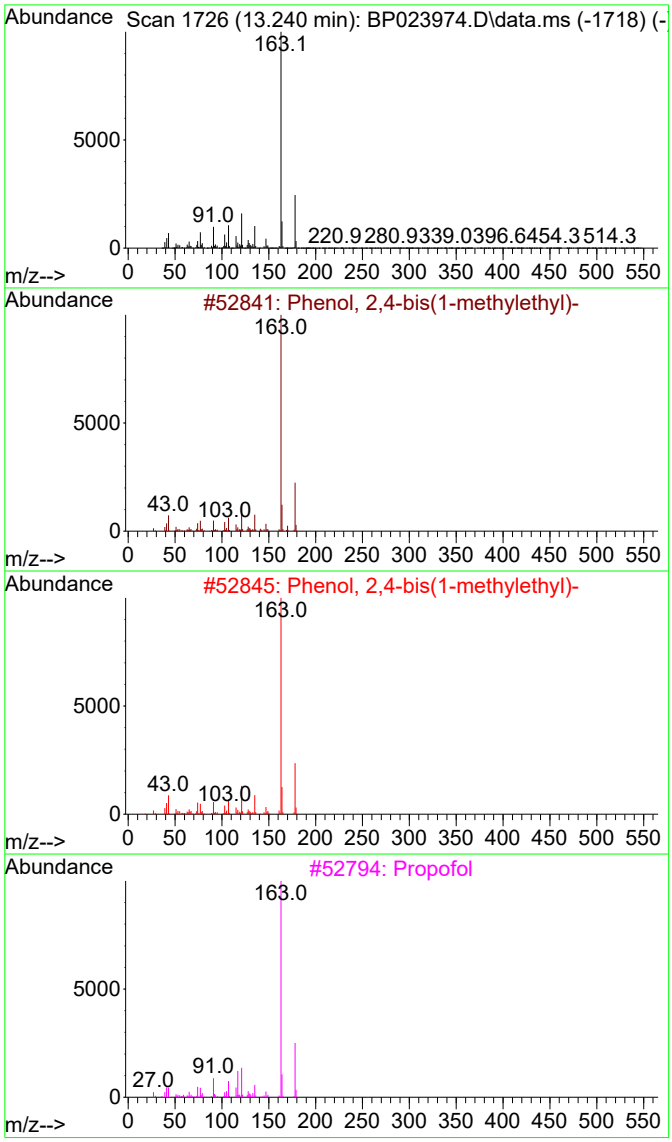
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

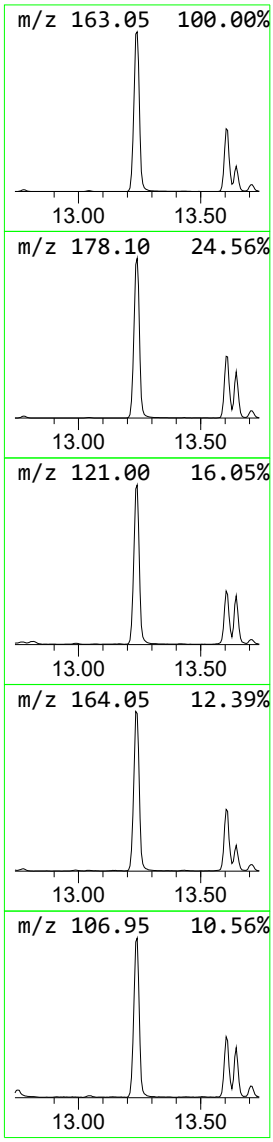
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Phenol, 2,4-bis(1-methyleth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.240	35.52 ng/ul	8973080	Acenaphthene-d10	14.393

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	93
3	Propofol	178	C12H18O	002078-54-8	93
4	Propofol	178	C12H18O	002078-54-8	91
5	Propofol	178	C12H18O	002078-54-8	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023974.D
Acq On : 06 Feb 2025 17:21
Operator : RC/JU
Sample : Q1202-20DL 30X
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B4DL

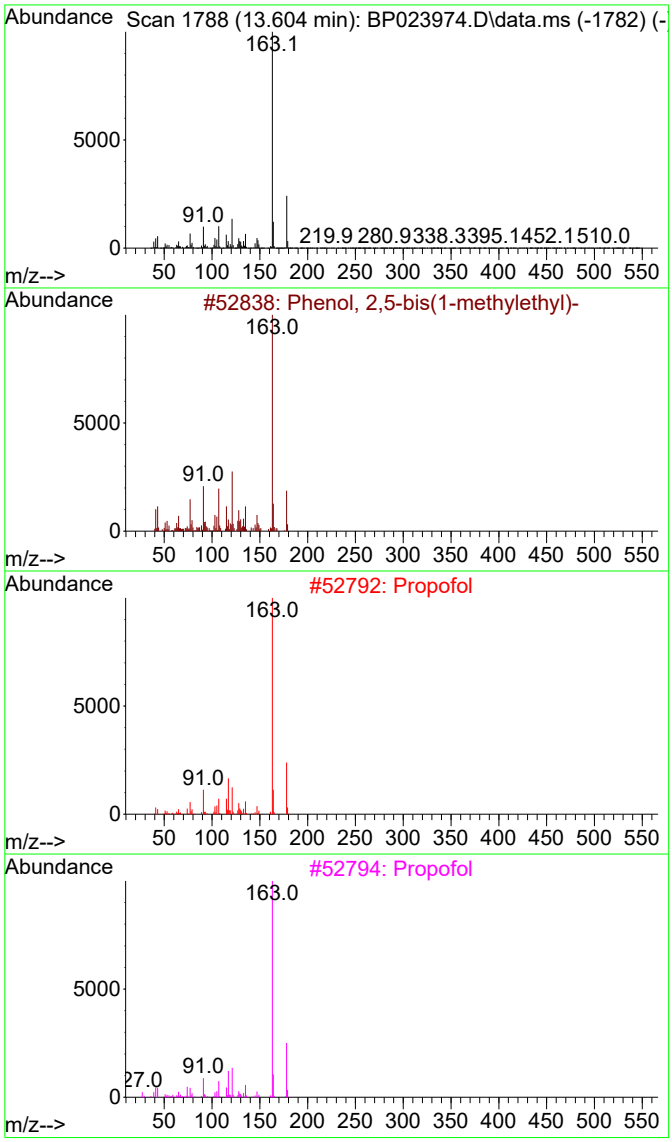
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

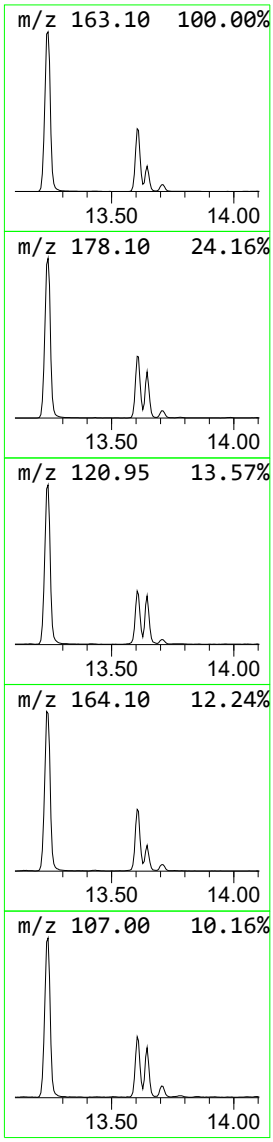
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.604	12.35 ng/ul	3119700	Acenaphthene-d10	14.393

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	93
4	Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	91
5	Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023974.D
Acq On : 06 Feb 2025 17:21
Operator : RC/JU
Sample : Q1202-20DL 30X
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B4DL

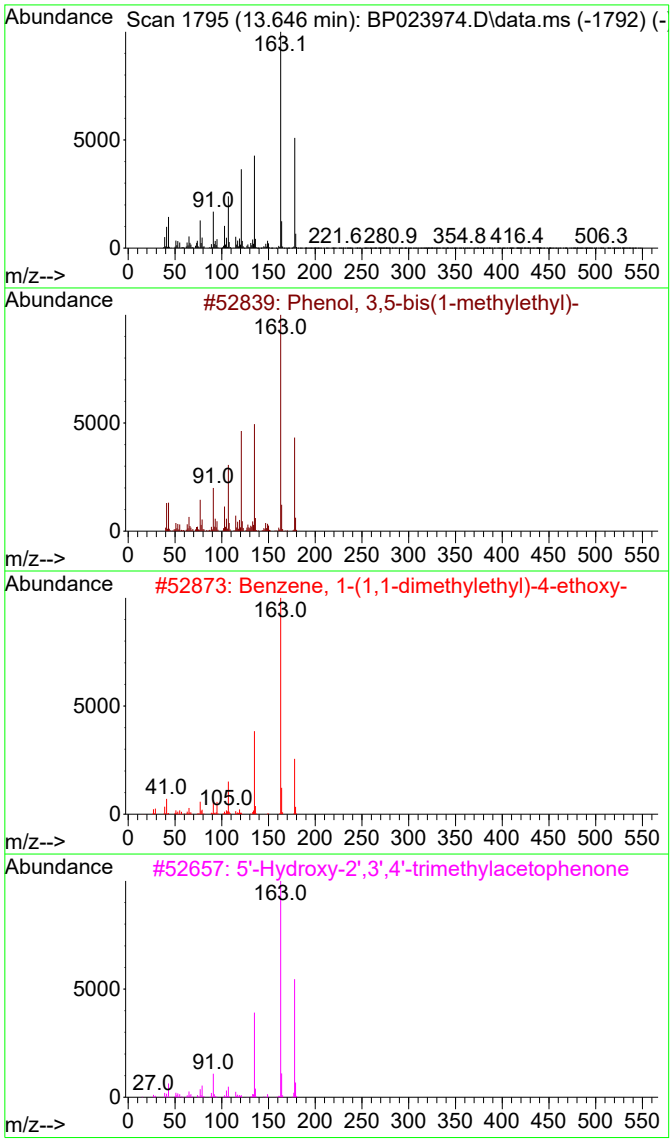
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

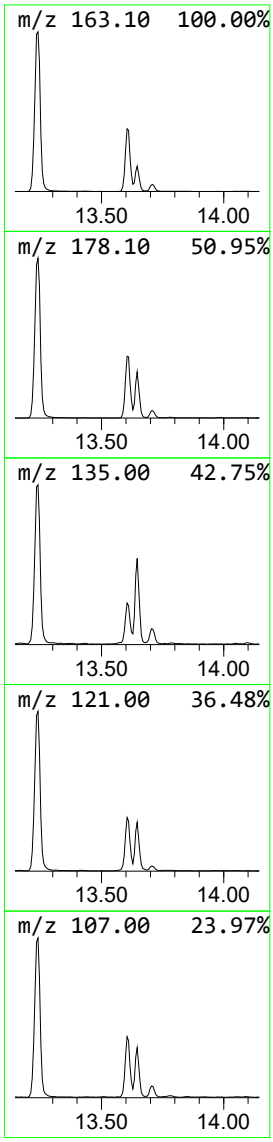
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Phenol, 3,5-bis(1-methyleth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.646	6.16 ng/ul	1556570	Acenaphthene-d10	14.393

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-bis(1-methylethyl)-	178	C12H18O	026886-05-5	94
2	Benzene, 1-(1,1-dimethylethyl)-4...	178	C12H18O	017269-94-2	90
3	5'-Hydroxy-2',3',4'-trimethylace...	178	C11H14O2	013667-28-2	87
4	6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	87
5	6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023974.D
Acq On : 06 Feb 2025 17:21
Operator : RC/JU
Sample : Q1202-20DL 30X
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B4DL

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
Phenol, 3-ethyl-	9.993	11.5	ng/ul	2183850	2	10.540	3807670	20.0
Phenol, 3,4-dim...	10.034	7.3	ng/ul	1397660	2	10.540	3807670	20.0
Phenol, 3,5-dim...	10.422	3.3	ng/ul	618175	2	10.540	3807670	20.0
Phenol, 2-(1-me...	10.475	88.8	ng/ul	16907400	2	10.540	3807670	20.0
p-Cumenol	10.899	37.8	ng/ul	7195590	2	10.540	3807670	20.0
Phenol, 2-(1,1-...	11.563	8.2	ng/ul	1562400	2	10.540	3807670	20.0
Phenol, 2,3,6-t...	11.628	4.8	ng/ul	905686	2	10.540	3807670	20.0
Phenol, m-tert-...	11.910	354.7	ng/ul	67524600	2	10.540	3807670	20.0
Phenol, 2,4-bis...	13.240	35.5	ng/ul	8973080	3	14.393	5052030	20.0
Phenol, 2,5-bis...	13.604	12.4	ng/ul	3119700	3	14.393	5052030	20.0
Phenol, 3,5-bis...	13.646	6.2	ng/ul	1556570	3	14.393	5052030	20.0