LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023975.D
Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : E29B0DL

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: BP023975.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.728	105	109	119	rBV	24701	49565	0.17%	0.047%
2	6.958	646	658	671	rVB2	30019	88047	0.30%	0.083%
3	7.111	677	684	691	rBV	88012	141361	0.48%	0.133%
4	7.311	713	718	725	rVB	83698	136059	0.46%	0.128%
5	7.769	789	796	803	rBV	1456674	2261537	7.67%	2.133%
6	8.475	909	916	921	rBV	82606	149435	0.51%	0.141%
7	8.540	921	927	935	rVB	301555	532655	1.81%	0.502%
8	8.910	983	990	996	rBV2	101203	179134	0.61%	0.169%
9	9.046	1009	1013	1020	rVB	34712	47302	0.16%	0.045%
10	9.175	1025	1035	1047	rVB	129460	253317	0.86%	0.239%
11	9.287	1047	1054	1060	rBV	11834	32242	0.11%	0.030%
12	9.528	1089	1095	1106	rBV	85072	144149	0.49%	0.136%
13	9.628	1106	1112	1119	rBV	64409	105267	0.36%	0.099%
14	9.728	1123	1129	1132	rVV	903988	1489128	5.05%	1.405%
15	9.757	1132	1134	1151	rVB	938142	1343419	4.55%	1.267%
16	9.993	1164	1174	1177	rBV	516417	1132131	3.84%	1.068%
17	10.034	1177	1181	1190	rVB	659391	1031857	3.50%	0.973%
18	10.181	1197	1206	1217	rBV2	681497	1073396	3.64%	1.012%
19	10.369	1233	1238	1241	rBV	25811	39905	0.14%	0.038%
20	10.416	1241	1246	1249	rVV	260188	402421	1.36%	0.380%
21	10.469	1249	1255	1261	rVV	7811004	11981588	40.62%	11.302%
22	10.546	1261	1268	1274	rVV	2110734	3590268	12.17%	3.387%
23	10.599	1274	1277	1284	rVV4	37974	71457	0.24%	0.067%
24	10.687	1284	1292	1301	rVB2	88502	192893	0.65%	0.182%
25	10.816	1307	1314	1320	rBV	69559	106969	0.36%	0.101%
26	10.899	1320	1328	1339	rVV	2031797	3469264	11.76%	3.272%
27	10.993	1339	1344	1354	rVV	97603	172504	0.58%	0.163%
28	11.087	1354	1360	1365	rVV	257548	443060	1.50%	0.418%
29	11.157	1365	1372	1386	rVB	165739	336924	1.14%	0.318%
30	11.310	1392	1398	1403	rBV2	23756	50157	0.17%	0.047%
31	11.375	1403	1409	1414	rVV	285939	491956	1.67%	0.464%
32	11.422	1414	1417	1425	rVB	46247	79134	0.27%	0.075%
33	11.563	1435	1441	1446	rBV	177096	305897	1.04%	0.289%
34	11.622	1446	1451	1456	rVV	204442	366706	1.24%	0.346%
35	11.669	1456	1459	1464	rVV2	59532	104432	0.35%	0.099%
36	11.887	1486	1496	1510	rBV	2559725	4061672	13.77%	3.831%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023975.D
Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

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	12.057	1520	1525	1528	rBV	37609	59090	0.20%	0.056%
38	12.110	1531	1534	1537	rVV	39022	55361	0.19%	0.052%
39	12.151	1537	1541	1545	rVV2	37090	56390	0.19%	0.053%
40	12.246	1553	1557	1560	rBV	24989	39532	0.13%	0.037%
41	12.346	1570	1574	1580	rBV	86735	124350	0.42%	0.117%
42	12.422	1584	1587	1592	rVB	48442	65712	0.22%	0.062%
43	12.534	1601	1606	1610	rBV4	24518	39527	0.13%	0.037%
44	12.593	1610	1616	1619	rBV2	59760	111346	0.38%	0.105%
45	12.681	1628	1631	1634	rVB2	36557	45446	0.15%	0.043%
46	12.769	1634	1646	1650	rVV6	179870	441451	1.50%	0.416%
47	12.798	1650	1651	1657	rVB2	45870	57987	0.20%	0.055%
48	12.975	1676	1681	1688	rVB3	54510	112142	0.38%	0.106%
49	13.057	1690	1695	1701	rBV5	13332	36104	0.12%	0.034%
50	13.234	1717	1725	1745	rBV	19205883	29499352	100.00%	27.825%
51	13.504	1766	1771	1775	rBV	43671	73855	0.25%	0.070%
52	13.598	1776	1787	1791	rVV	7018506	9641870	32.69%	9.095%
53	13.640	1791	1794	1800	rVV	1583070	1976538	6.70%	1.864%
54	13.704	1801	1805	1811	rVV	377132	560339	1.90%	0.529%
55	13.793	1815	1820	1826	rVV	250901	363215	1.23%	0.343%
56	13.846	1826	1829	1839	rVB	57596	97421	0.33%	0.092%
57	14.075	1862	1868	1877	rVB2	265204	432410	1.47%	0.408%
58	14.151	1877	1881	1885	rBV	91399	133168	0.45%	0.126%
59	14.387	1914	1921	1927	rBV	3068892	4724481	16.02%	4.456%
60	14.434	1927	1929	1938	rVB3	61547	94927	0.32%	0.090%
61	14.640	1956	1964	1977	rVB6	50169	151979	0.52%	0.143%
62	15.075	2033	2038	2046	rBV2	146282	232750	0.79%	0.220%
63	15.198	2055	2059	2062	rBV	39569	54484	0.18%	0.051%
64	15.263	2067	2070	2078	rVB5	34993	62919	0.21%	0.059%
65	15.392	2086	2092	2099	rBV	386159	597983	2.03%	0.564%
66	15.475	2102	2106	2108	rVV2	31817	45868	0.16%	0.043%
67	15.510	2108	2112	2118	rVV2	59544	106105	0.36%	0.100%
68	15.569	2118	2122	2127	rVB	92880	128236	0.43%	0.121%
69	16.087	2206	2210	2216	rBV3	23597	44619	0.15%	0.042%
70	16.445	2267	2271	2275	rBV2	24809	37671	0.13%	0.036%
71	17.181	2390	2396	2407	rBV2	3812354	5643335	19.13%	5.323%
72	17.281	2408	2413	2422	rVB2	384269	610738	2.07%	0.576%
73	19.222	2739	2743	2748	rVB	147486	209792	0.71%	0.198%
74	19.657	2812	2817	2823	rBV	528594	726253	2.46%	0.685%
75	21.627	3147	3152	3160	rVB2	3477649	5940703	20.14%	5.604%
76	25.010	3719	3727	3742	rVB	2602439	6123193	20.76%	5.776%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023975.D
Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

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

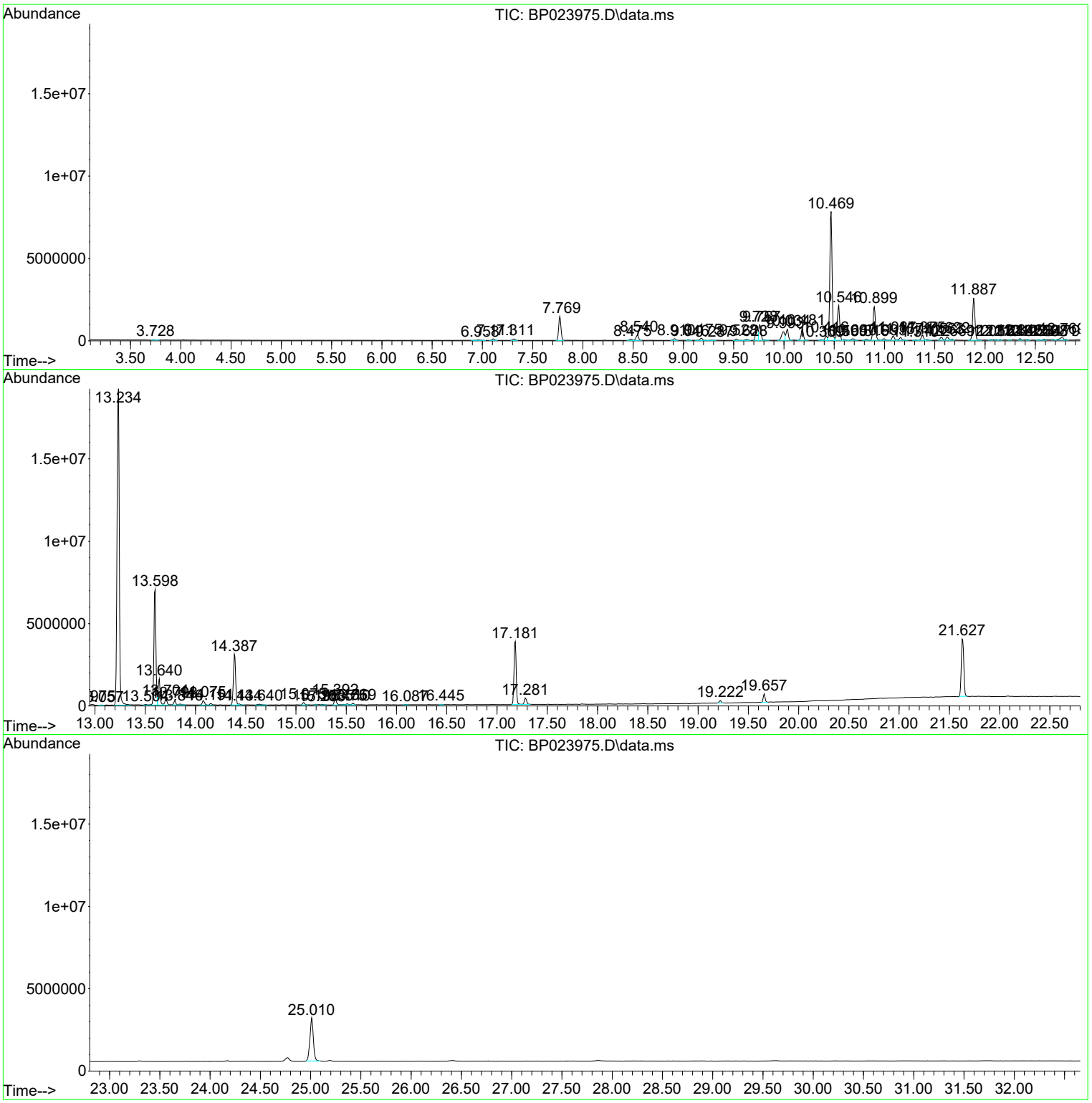
Sum of corrected areas: 106015850

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023975.D
Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

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 : BP023975.D
Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

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

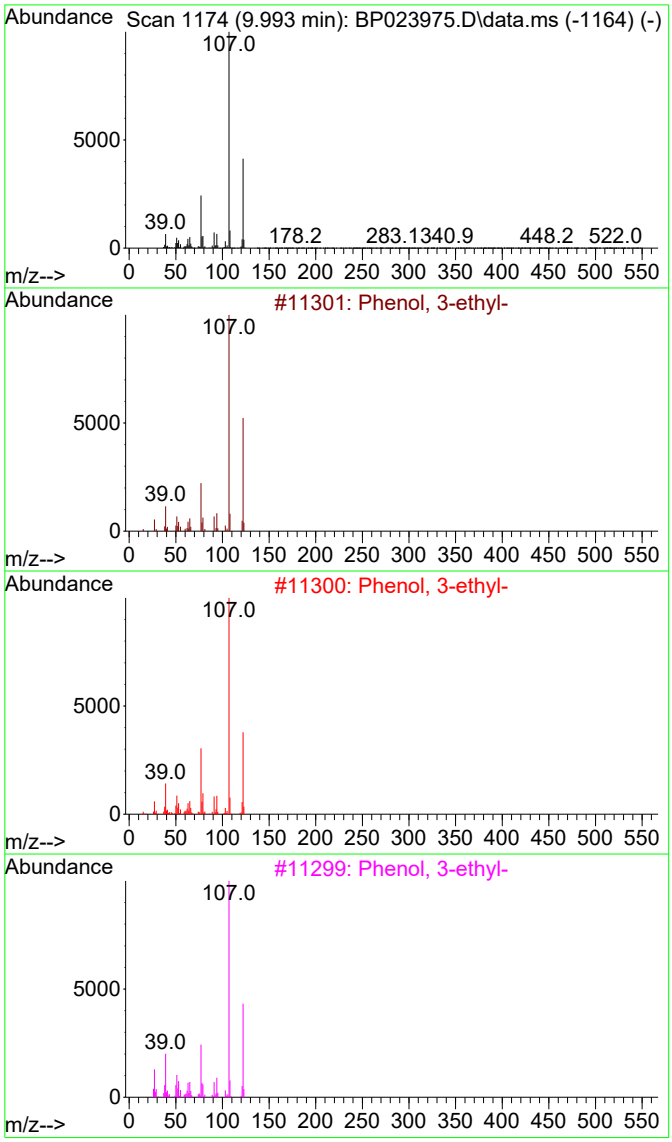
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

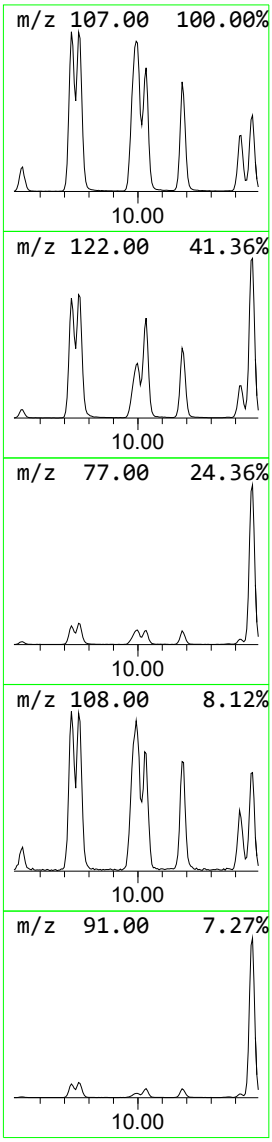
****

Peak Number 1 Phenol, 3-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.993	6.31 ng/ul	1132130	Naphthalene-d8	10.546

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
5	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023975.D
Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

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

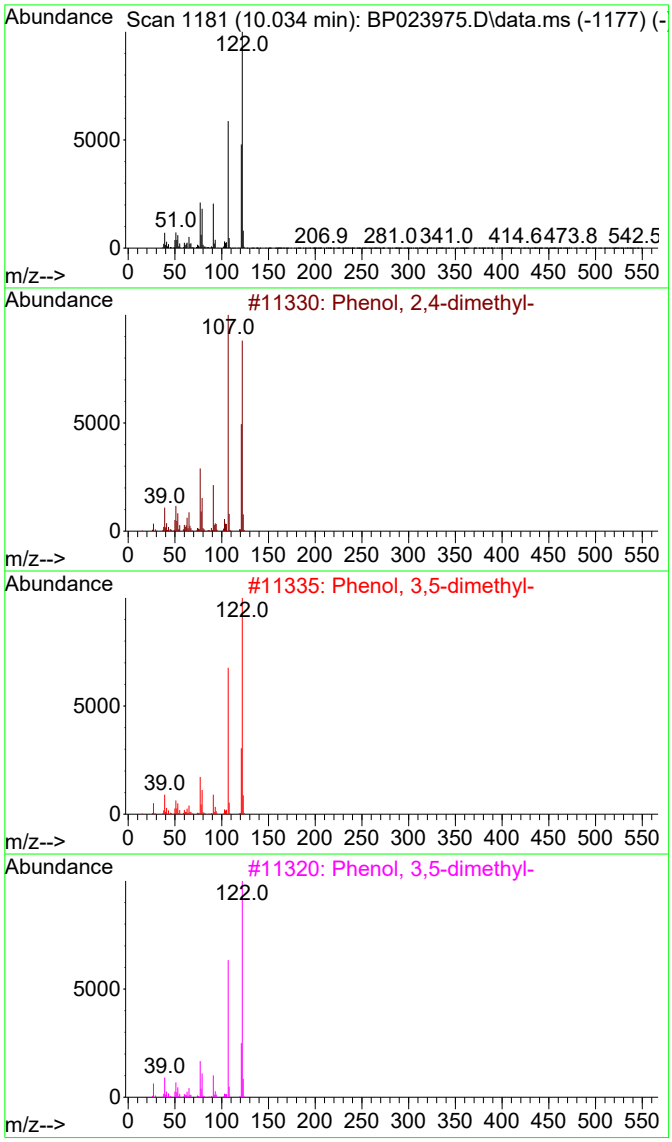
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

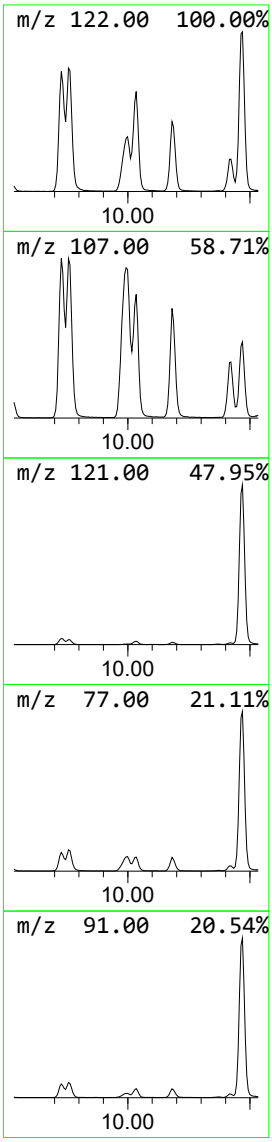
****

Peak Number 2 Phenol, 3,5-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.034	5.75 ng/ul	1031860	Naphthalene-d8	10.546

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	91
4	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	91
5	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023975.D
Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

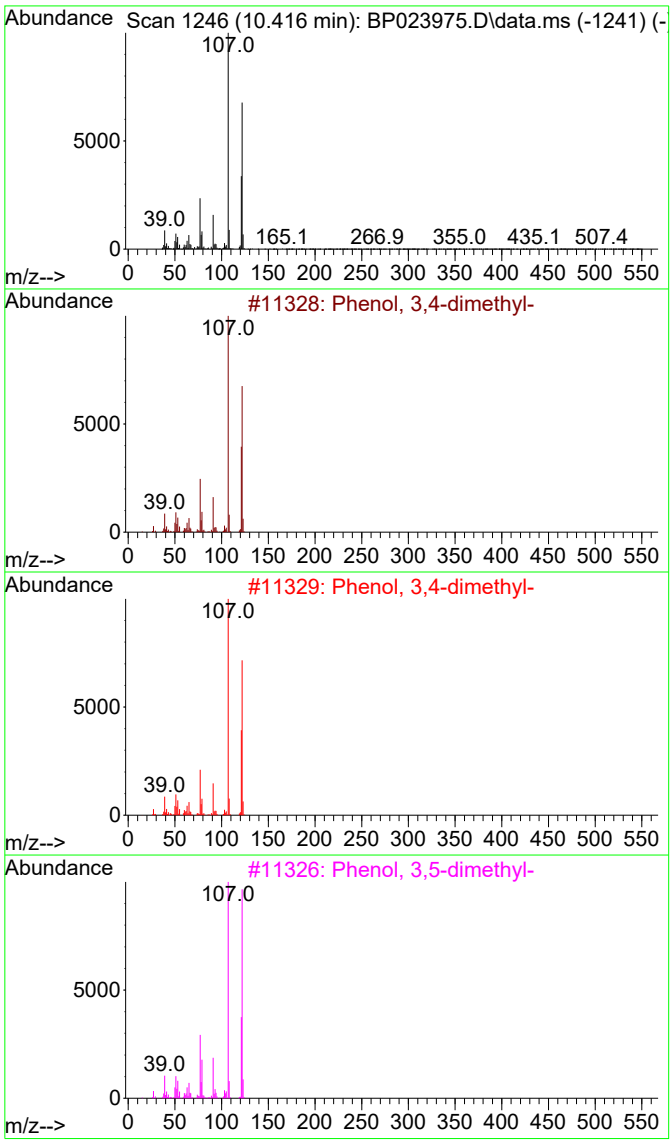
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

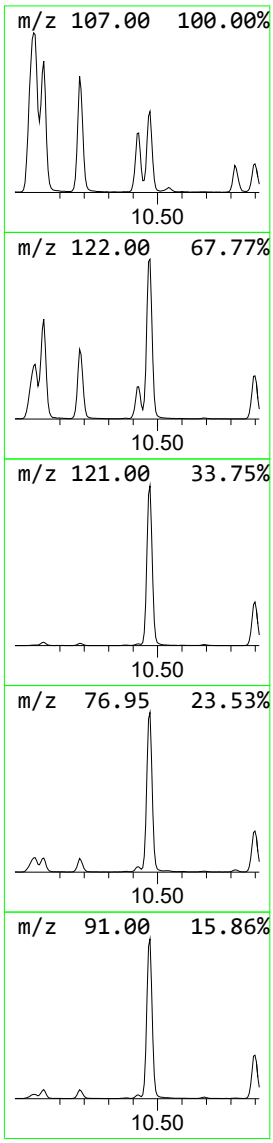
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Phenol, 3,4-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.416	2.24 ng/ul	402421	Naphthalene-d8	10.546

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, 3,5-dimethyl-	122	C8H10O	000108-68-9	94
4	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	94
5	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023975.D
Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

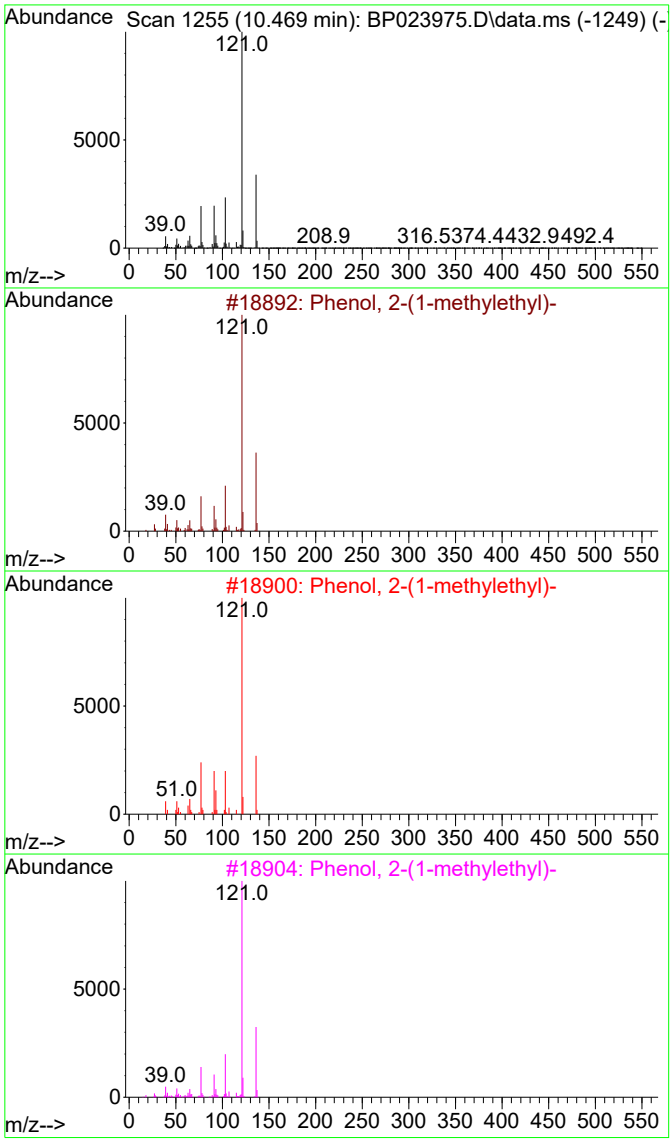
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

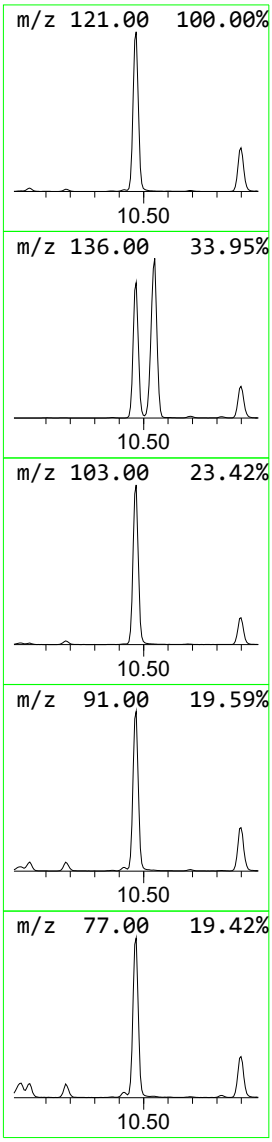
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.469	66.74 ng/ul	11981600	Naphthalene-d8	10.546

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	91
3	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	90
4	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	87
5	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023975.D
Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

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

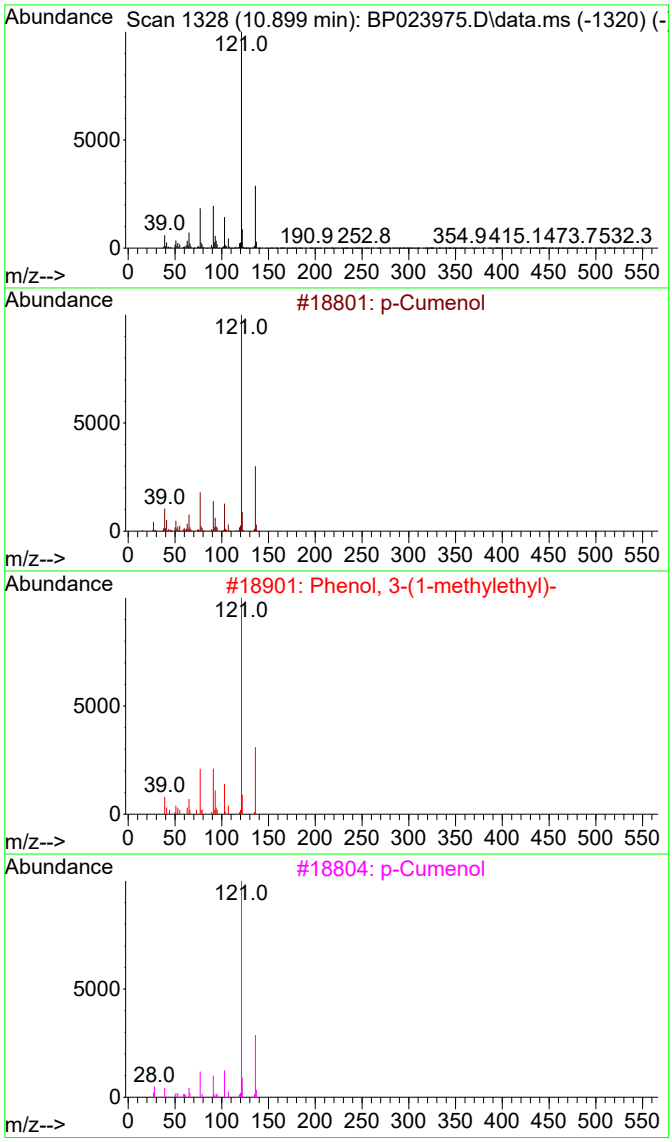
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

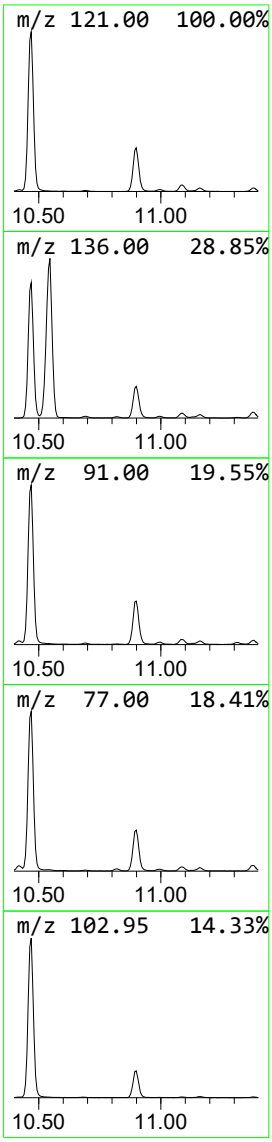
****

Peak Number 5 p-Cumenol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.899	19.33 ng/ul	3469260	Naphthalene-d8	10.546

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





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023975.D
Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

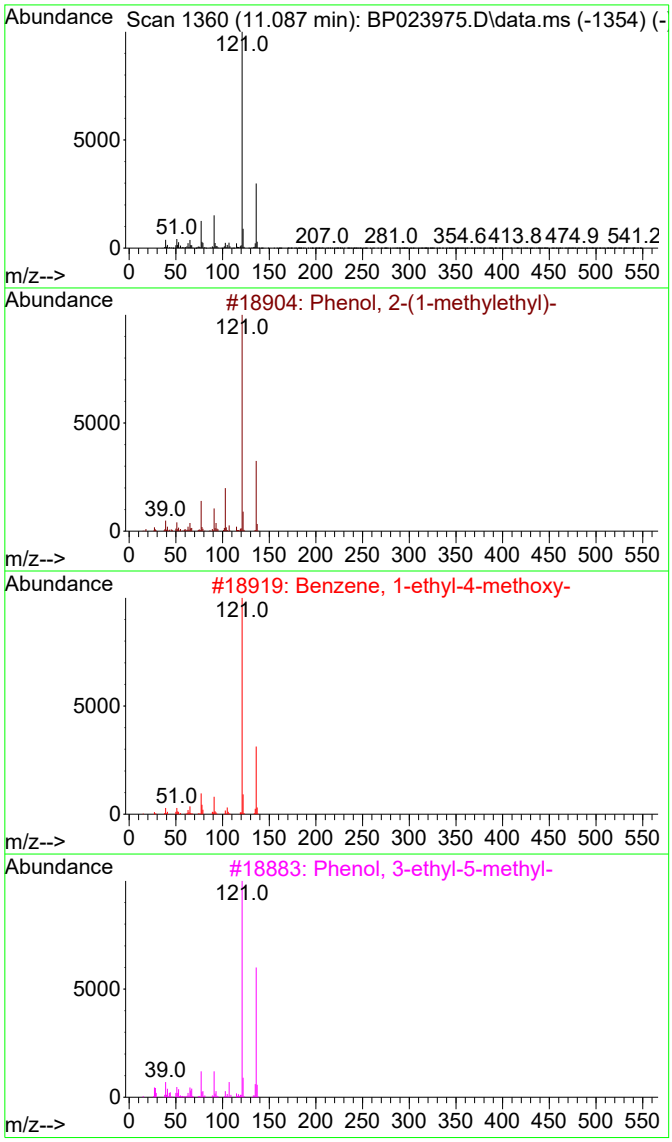
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

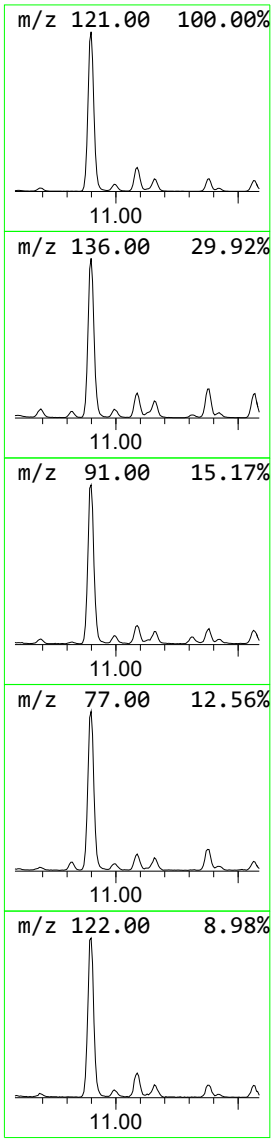
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.087	2.47 ng/ul	443060	Naphthalene-d8	10.546

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	91
2	Benzene, 1-ethyl-4-methoxy-	136	C9H12O	001515-95-3	91
3	Phenol, 3-ethyl-5-methyl-	136	C9H12O	000698-71-5	90
4	Phenol, 4-ethyl-3-methyl-	136	C9H12O	001123-94-0	90
5	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023975.D
Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

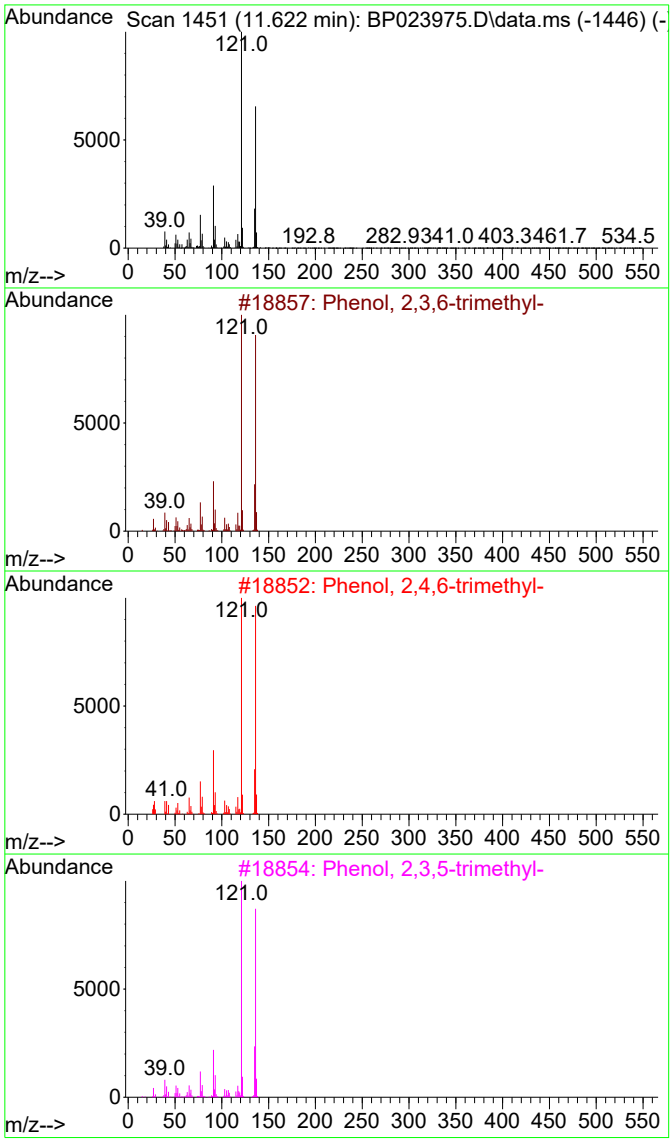
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

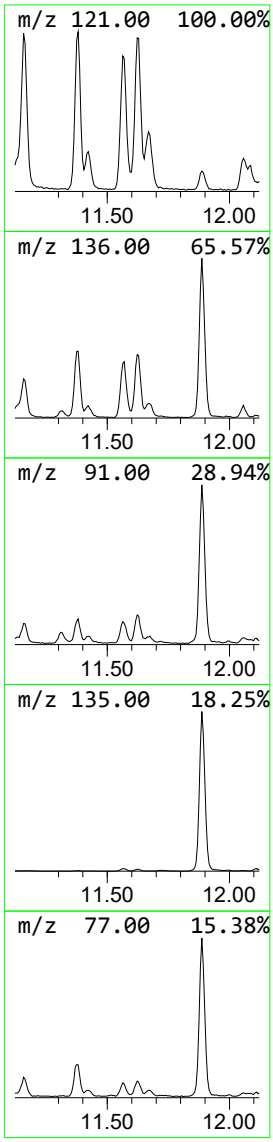
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Phenol, 2,3,6-trimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.622	2.04 ng/ul	366706	Naphthalene-d8	10.546

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





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Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

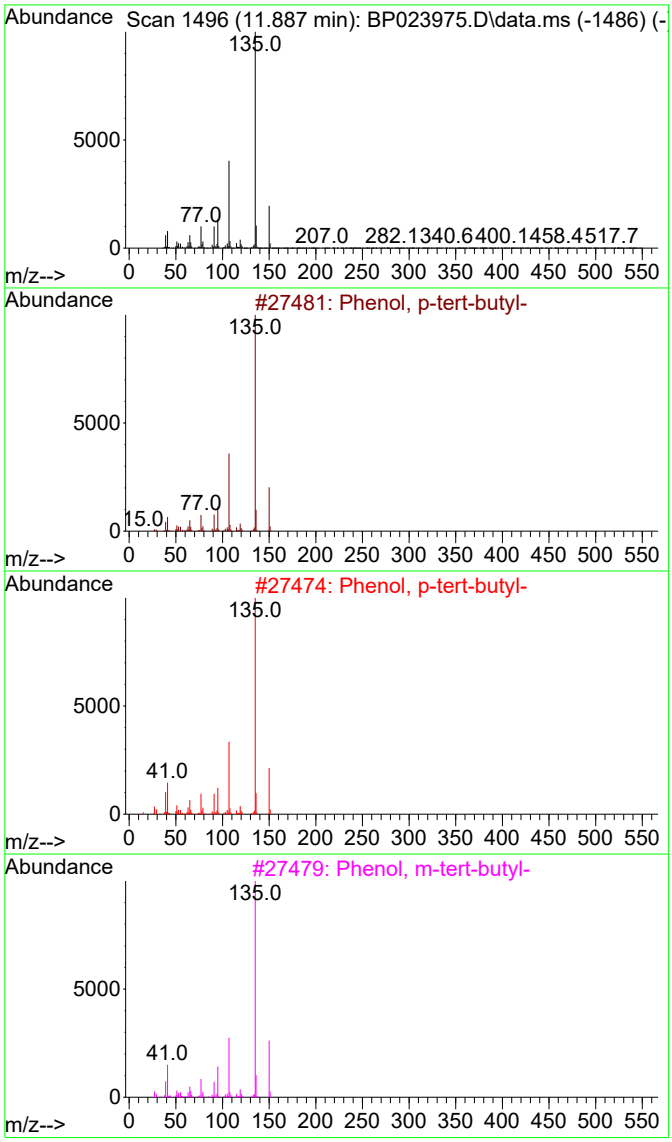
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

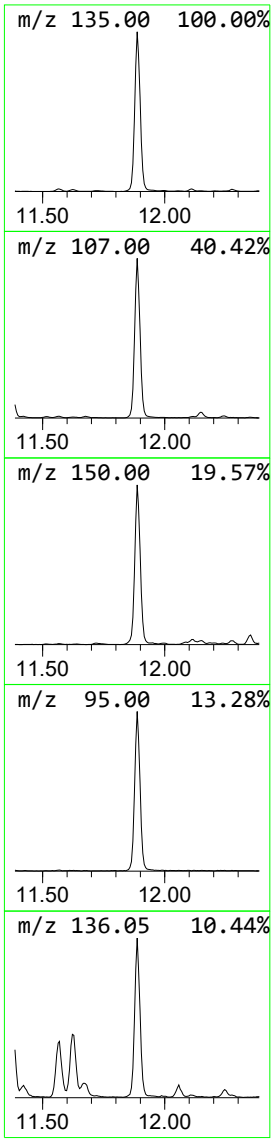
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Phenol, p-tert-butyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.887	22.63 ng/ul	4061670	Naphthalene-d8	10.546

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, m-tert-butyl-	150	C10H14O	000585-34-2	87





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Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

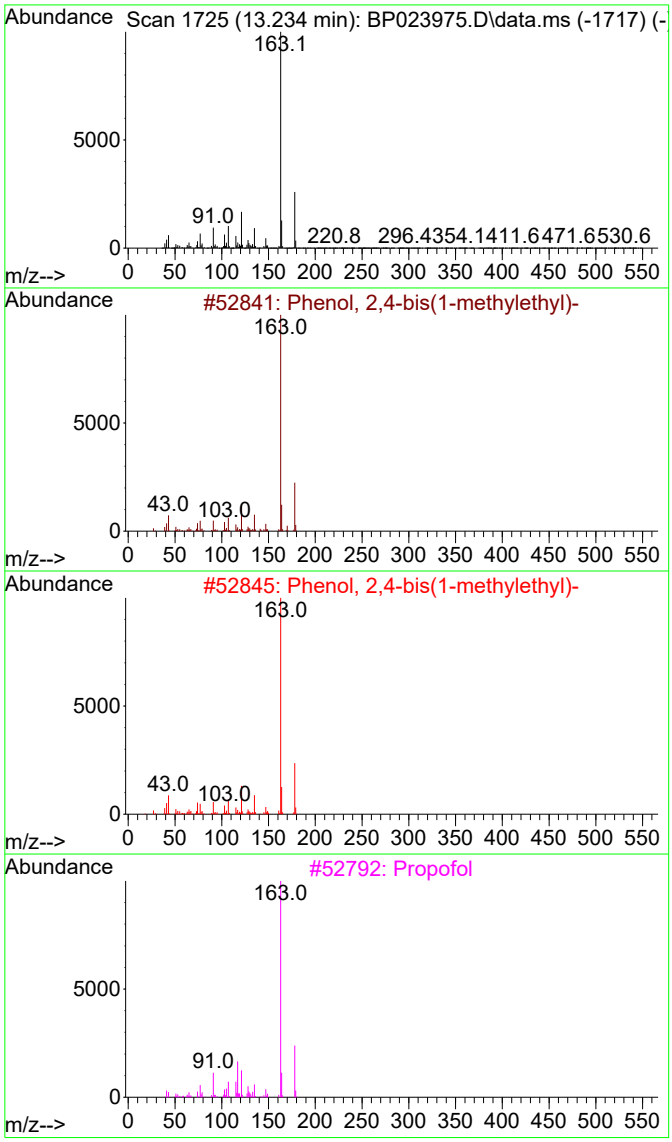
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

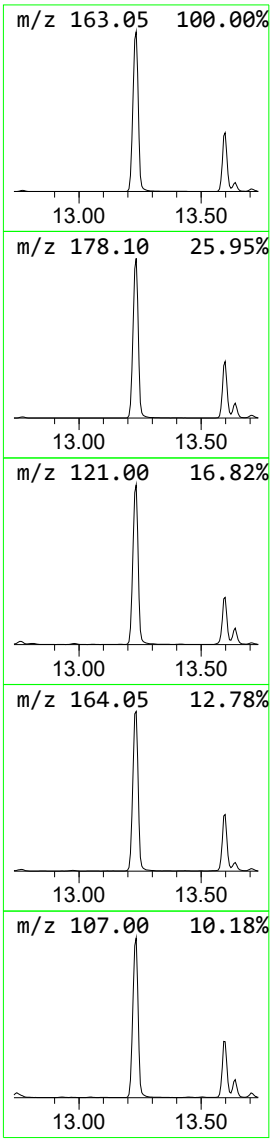
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.234	124.88 ng/ul	29499400	Acenaphthene-d10	14.387

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	94
3	Propofol	178	C12H18O	002078-54-8	94
4	Propofol	178	C12H18O	002078-54-8	94
5	Propofol	178	C12H18O	002078-54-8	94





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Data File : BP023975.D
Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

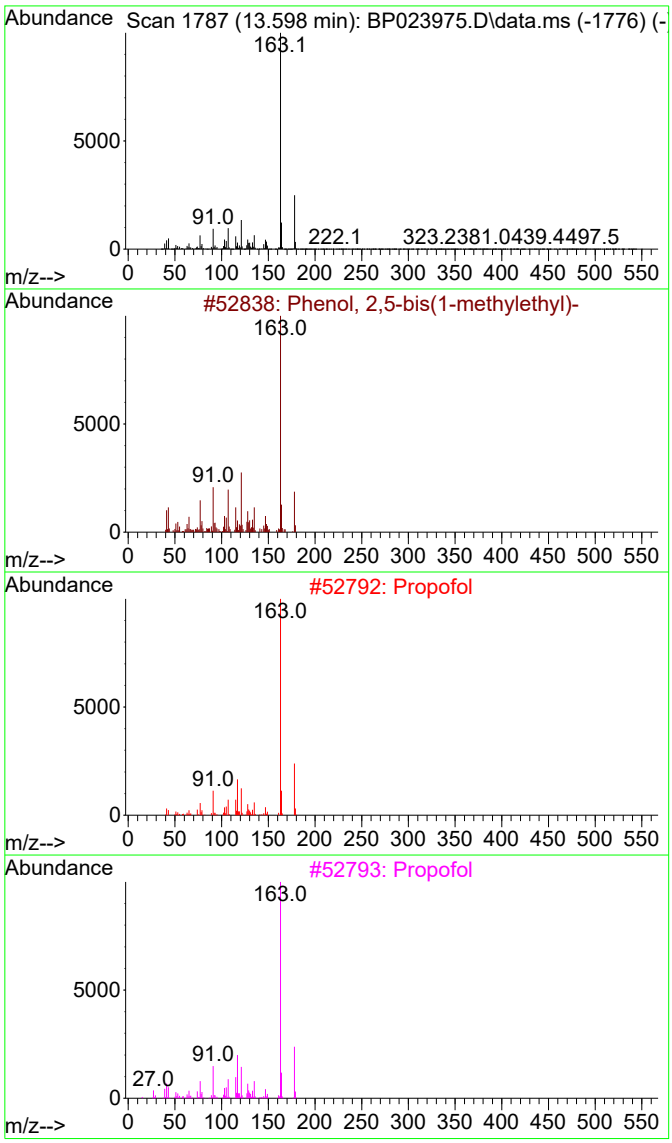
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

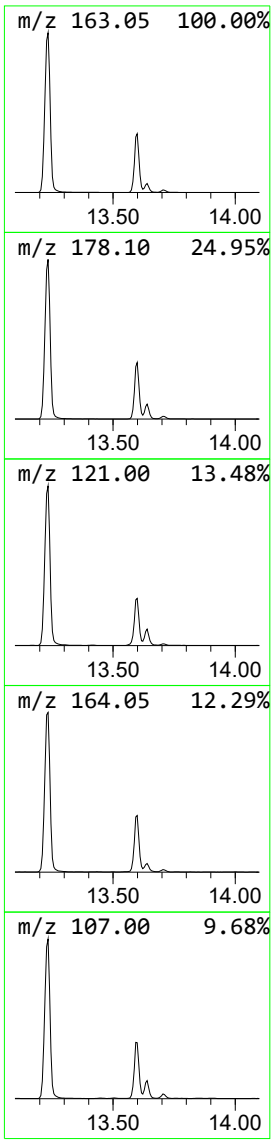
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Phenol, 2,5-bis(1-methyleth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.599	40.82 ng/ul	9641870	Acenaphthene-d10	14.387

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





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023975.D
Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

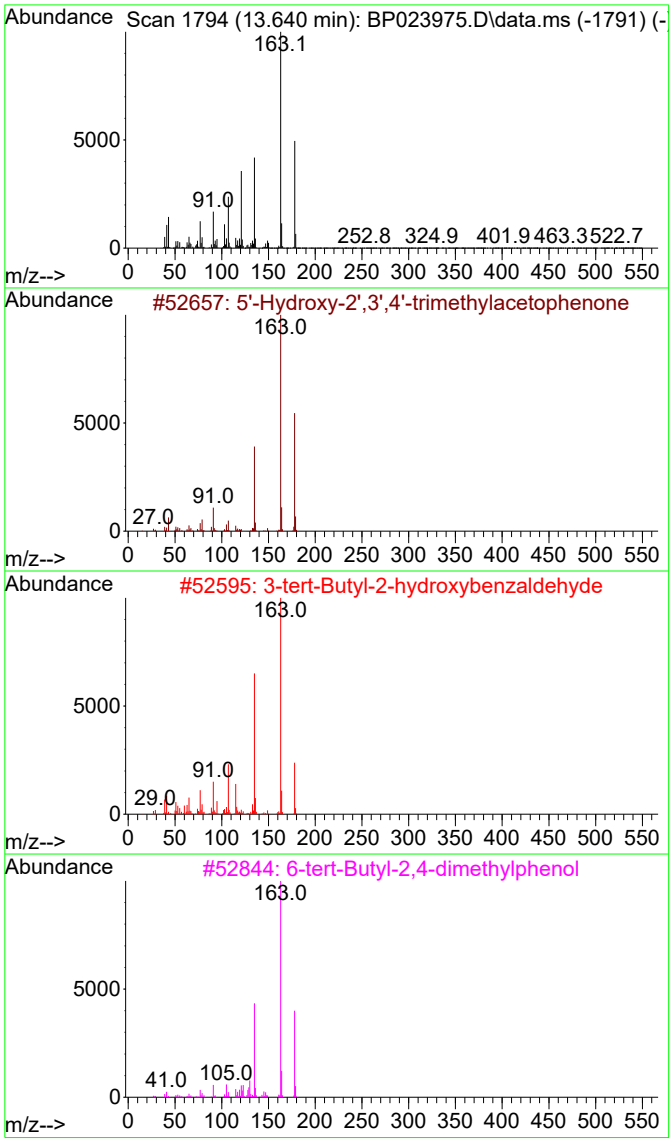
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

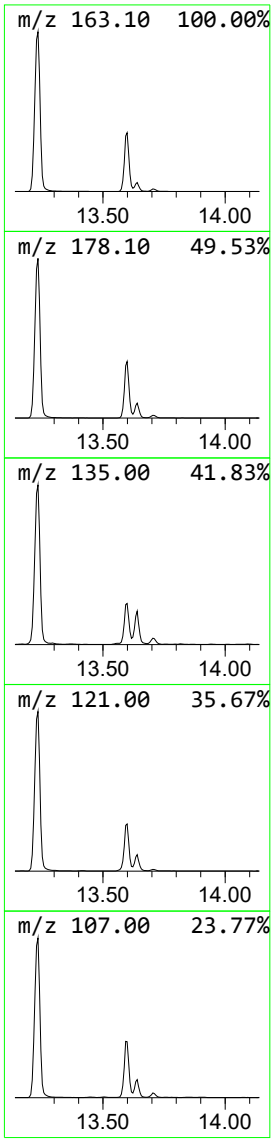
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 5'-Hydroxy-2',3',4'-trimeth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.640	8.37 ng/ul	1976540	Acenaphthene-d10	14.387

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5'-Hydroxy-2',3',4'-trimethylace...	178	C11H14O2	013667-28-2	87
2	3-tert-Butyl-2-hydroxybenzaldehyde	178	C11H14O2	024623-65-2	86
3	6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	81
4	Propofol	178	C12H18O	002078-54-8	76
5	Propofol	178	C12H18O	002078-54-8	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023975.D
Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

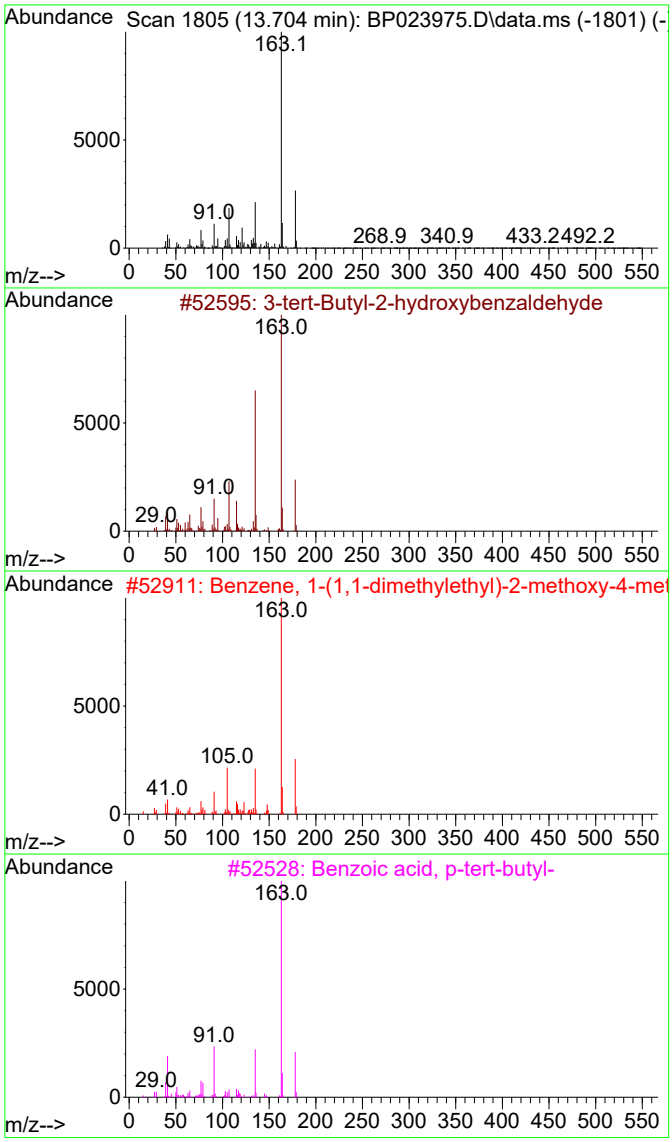
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

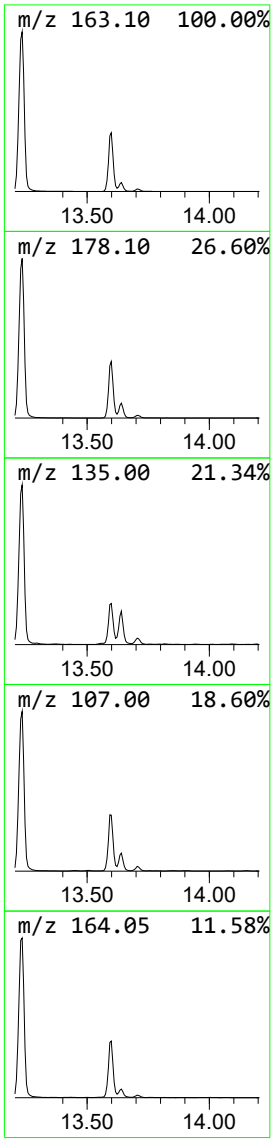
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 3-tert-Butyl-2-hydroxybenza... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.704	2.37 ng/ul	560339	Acenaphthene-d10	14.387

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-tert-Butyl-2-hydroxybenzaldehyde	178	C11H14O2	024623-65-2	87
2	Benzene, 1-(1,1-dimethylethyl)-2...	178	C12H18O	000088-40-4	87
3	Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	81
4	Propofol	178	C12H18O	002078-54-8	74
5	Propofol	178	C12H18O	002078-54-8	72





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023975.D
Acq On : 06 Feb 2025 18:01
Operator : RC/JU
Sample : Q1202-17DL 20X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29B0DL

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	6.3	ng/ul	1132130	2	10.546	3590270	20.0
Phenol, 3,5-dim...	10.034	5.8	ng/ul	1031860	2	10.546	3590270	20.0
Phenol, 3,4-dim...	10.416	2.2	ng/ul	402421	2	10.546	3590270	20.0
Phenol, 2-(1-me...	10.469	66.7	ng/ul	11981600	2	10.546	3590270	20.0
p-Cumenol	10.899	19.3	ng/ul	3469260	2	10.546	3590270	20.0
Benzene, 1-ethy...	11.087	2.5	ng/ul	443060	2	10.546	3590270	20.0
Phenol, 2,3,6-t...	11.622	2.0	ng/ul	366706	2	10.546	3590270	20.0
Phenol, p-tert-...	11.887	22.6	ng/ul	4061670	2	10.546	3590270	20.0
Phenol, 2,4-bis...	13.234	124.9	ng/ul	29499400	3	14.387	4724480	20.0
Phenol, 2,5-bis...	13.599	40.8	ng/ul	9641870	3	14.387	4724480	20.0
5'-Hydroxy-2',3...	13.640	8.4	ng/ul	1976540	3	14.387	4724480	20.0
3-tert-Butyl-2-...	13.704	2.4	ng/ul	560339	3	14.387	4724480	20.0