LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBTE0DL

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-BP051322.M
Title : SVOA CALIBRATION

Signal : TIC: BP010545.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.528	409	415	423	rBV	25874	55462	0.18%	0.075%
2	5.899	469	478	485	rVB3	22407	47550	0.16%	0.064%
3	7.216	695	702	708	rBV	40489	67170	0.22%	0.090%
4	7.422	729	737	746	rBV3	29803	58149	0.19%	0.078%
5	7.534	750	756	763	rVV	54108	89916	0.29%	0.121%
6	7.605	763	768	776	rVB	35031	61535	0.20%	0.083%
7	7.881	807	815	825	rBV	453404	769030	2.52%	1.033%
8	7.999	830	835	843	rVB3	23437	40259	0.13%	0.054%
9	8.258	872	879	887	rBV	192387	313203	1.03%	0.421%
10	8.422	895	907	918	rBV	834038	1425472	4.67%	1.915%
11	8.599	932	937	947	rBV4	20095	39702	0.13%	0.053%
12	9.052	1008	1014	1022	rVV2	43288	105415	0.35%	0.142%
13	9.246	1040	1047	1054	rBV	38583	71176	0.23%	0.096%
14	9.363	1054	1067	1073	rBV	93509	212253	0.69%	0.285%
15	9.428	1073	1078	1085	rVB2	27904	49499	0.16%	0.067%
16	9.769	1130	1136	1142	rVV	31117	56182	0.18%	0.075%
17	9.934	1159	1164	1169	rBV2	23040	40951	0.13%	0.055%
18	10.093	1180	1191	1201	rBV5	64616	196180	0.64%	0.264%
19	10.193	1201	1208	1219	rVV3	43624	142387	0.47%	0.191%
20	10.316	1221	1229	1237	rVV2	50816	112641	0.37%	0.151%
21	10.687	1284	1292	1294	rVV	604341	1056723	3.46%	1.420%
22	10.746	1294	1302	1313	rVV	16758062	30540029	100.00%	41.030%
23	10.852	1313	1320	1331	rVB	873486	1532688	5.02%	2.059%
24	11.793	1474	1480	1485	rBV	29315	50658	0.17%	0.068%
25	12.234	1549	1555	1562	rBV	57158	92003	0.30%	0.124%
26	12.340	1565	1573	1588	rBV	2463644	3955032	12.95%	5.314%
27	12.457	1588	1593	1599	rVB	63582	113144	0.37%	0.152%
28	12.557	1599	1610	1623	rBV	1382157	2279106	7.46%	3.062%
29	12.987	1677	1683	1691	rVB3	35449	68824	0.23%	0.092%
30	13.346	1737	1744	1754	rBV	511492	801174	2.62%	1.076%
31	13.546	1772	1778	1790	rVB2	83718	197573	0.65%	0.265%
32	13.834	1821	1827	1833	rBV2	139112	251499	0.82%	0.338%
33	13.893	1833	1837	1839	rVV	84454	130173	0.43%	0.175%
34	13.922	1839	1842	1851	rVB	139737	213855	0.70%	0.287%
35	14.075	1862	1868	1871	rBV	51935	78767	0.26%	0.106%
36	14.204	1884	1890	1893	rBV	132496	217441	0.71%	0.292%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

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-BP051322.M
Title : SVOA CALIBRATION

37	14.234	1893	1895	1903	rVB	115188	159879	0.52%	0.215%
38	14.516	1932	1943	1948	rBV	1125987	1752981	5.74%	2.355%
39	14.575	1948	1953	1960	rVV	2274189	3396048	11.12%	4.563%
40	14.645	1960	1965	1976	rVB	653921	1031393	3.38%	1.386%
41	14.828	1986	1996	2000	rVB2	31478	57103	0.19%	0.077%
42	14.916	2004	2011	2021	rBV2	1108920	2089419	6.84%	2.807%
43	15.063	2031	2036	2042	rBV2	62439	100763	0.33%	0.135%
44	15.151	2042	2051	2054	rVV10	20973	49260	0.16%	0.066%
45	15.504	2106	2111	2116	rBV	191150	286539	0.94%	0.385%
46	15.563	2116	2121	2130	rVB2	1003116	1439835	4.71%	1.934%
47	15.651	2130	2136	2139	rBV5	44162	90665	0.30%	0.122%
48	15.687	2139	2142	2145	rVV2	50426	66975	0.22%	0.090%
49	15.728	2145	2149	2154	rVB3	39252	58108	0.19%	0.078%
50	15.857	2168	2171	2180	rVB2	59345	95642	0.31%	0.128%
51	16.004	2192	2196	2209	rVB3	75889	150902	0.49%	0.203%
52	16.240	2230	2236	2245	rVB2	85612	152133	0.50%	0.204%
53	16.416	2262	2266	2270	rBV	23174	36207	0.12%	0.049%
54	16.528	2281	2285	2287	rBV2	20582	30928	0.10%	0.042%
55	16.569	2290	2292	2298	rVV3	27344	44033	0.14%	0.059%
56	16.922	2347	2352	2361	rVB	123968	197418	0.65%	0.265%
57	17.087	2375	2380	2386	rVV	96837	147645	0.48%	0.198%
58	17.163	2389	2393	2396	rVV	23590	31258	0.10%	0.042%
59	17.222	2399	2403	2405	rVV2	30138	39709	0.13%	0.053%
60	17.263	2405	2410	2414	rVV	1461319	2208636	7.23%	2.967%
61	17.310	2414	2418	2423	rVV	1032070	1571969	5.15%	2.112%
62	17.363	2423	2427	2431	rVV2	219518	354535	1.16%	0.476%
63	17.398	2431	2433	2438	rVV2	78795	119559	0.39%	0.161%
64	17.669	2470	2479	2491	rBV	2790589	3980291	13.03%	5.347%
65	17.769	2491	2496	2502	rVB5	62069	112434	0.37%	0.151%
66	17.828	2502	2506	2511	rBV	44247	65753	0.22%	0.088%
67	17.922	2518	2522	2526	rVB2	28332	44497	0.15%	0.060%
68	17.975	2526	2531	2535	rBV2	27768	42808	0.14%	0.058%
69	18.098	2547	2552	2556	rBV	60294	92478	0.30%	0.124%
70	18.175	2556	2565	2570	rBV	248743	380824	1.25%	0.512%
71	18.251	2576	2578	2585	rVB	31617	36863	0.12%	0.050%
72	18.322	2585	2590	2594	rBV	70500	106860	0.35%	0.144%
73	18.422	2599	2607	2611	rBV3	75652	212632	0.70%	0.286%
74	18.463	2611	2614	2618	rVV	99369	136884	0.45%	0.184%
75	18.498	2618	2620	2625	rVB2	29127	38016	0.12%	0.051%
76	18.557	2625	2630	2636	rBV2	108936	186200	0.61%	0.250%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

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-BP051322.M
Title : SVOA CALIBRATION

77	18.616	2636	2640	2644	rVV2	45817	63817	0.21%	0.086%
78	19.104	2719	2723	2725	rBV5	22489	33531	0.11%	0.045%
79	19.269	2747	2751	2758	rVB	100599	137944	0.45%	0.185%
80	19.598	2802	2807	2811	rBV	317584	470031	1.54%	0.631%
81	19.992	2870	2874	2880	rBV	79185	118815	0.39%	0.160%
82	20.057	2881	2885	2893	rVV2	43268	82821	0.27%	0.111%
83	21.345	3099	3104	3115	rBV	2044359	2737510	8.96%	3.678%
84	22.463	3291	3294	3303	rVB	103449	134081	0.44%	0.180%
85	23.022	3385	3389	3394	rVB	142358	209376	0.69%	0.281%
86	23.610	3484	3489	3493	rBV	140710	274333	0.90%	0.369%
87	23.657	3493	3497	3503	rVB	214729	366561	1.20%	0.492%
88	23.763	3509	3515	3529	rVB2	1077833	2272818	7.44%	3.053%
89	24.386	3616	3621	3630	rVB	192293	408239	1.34%	0.548%
90	25.245	3762	3767	3777	rVB2	168802	394444	1.29%	0.530%

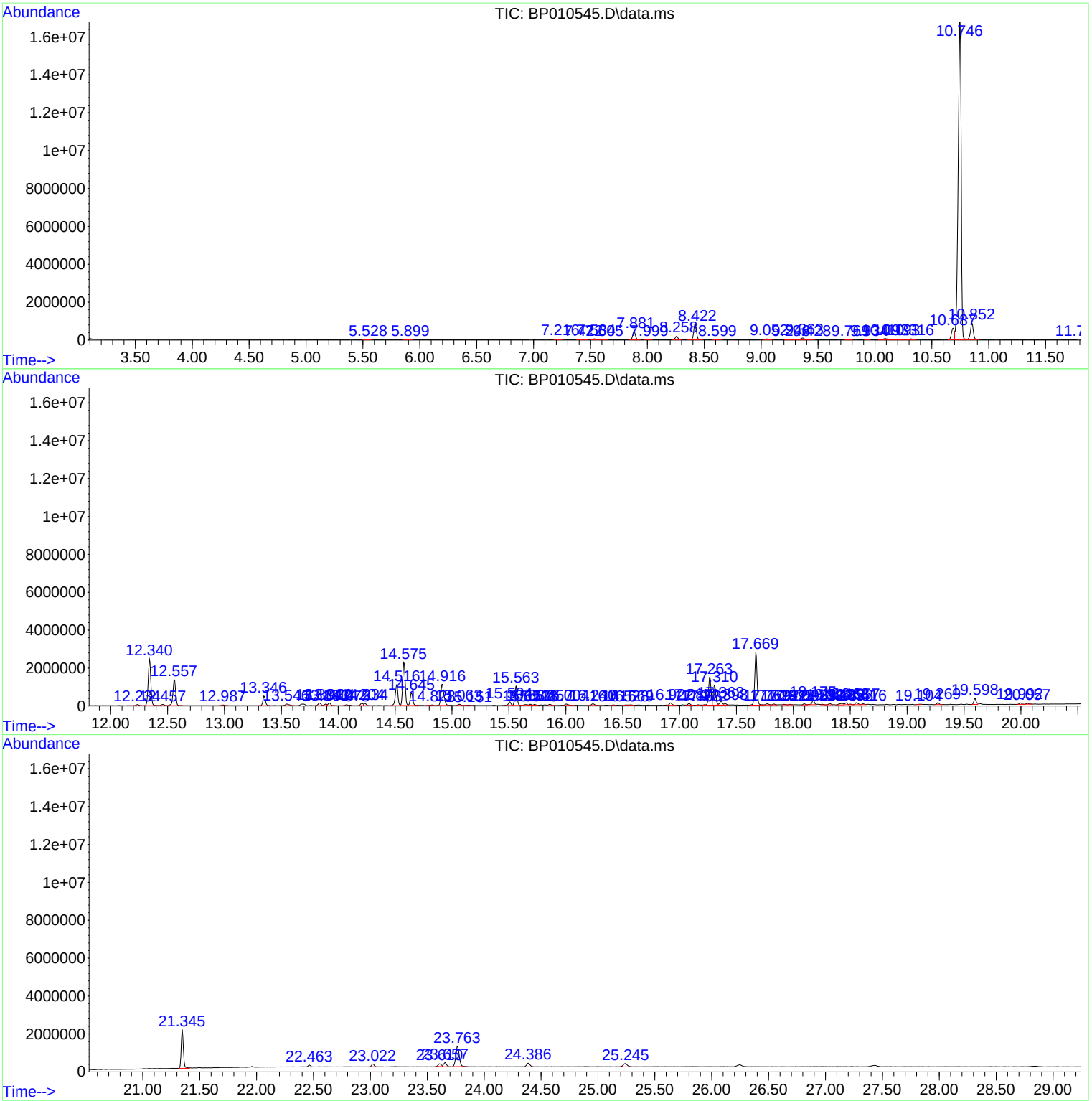
Sum of corrected areas: 74433224

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

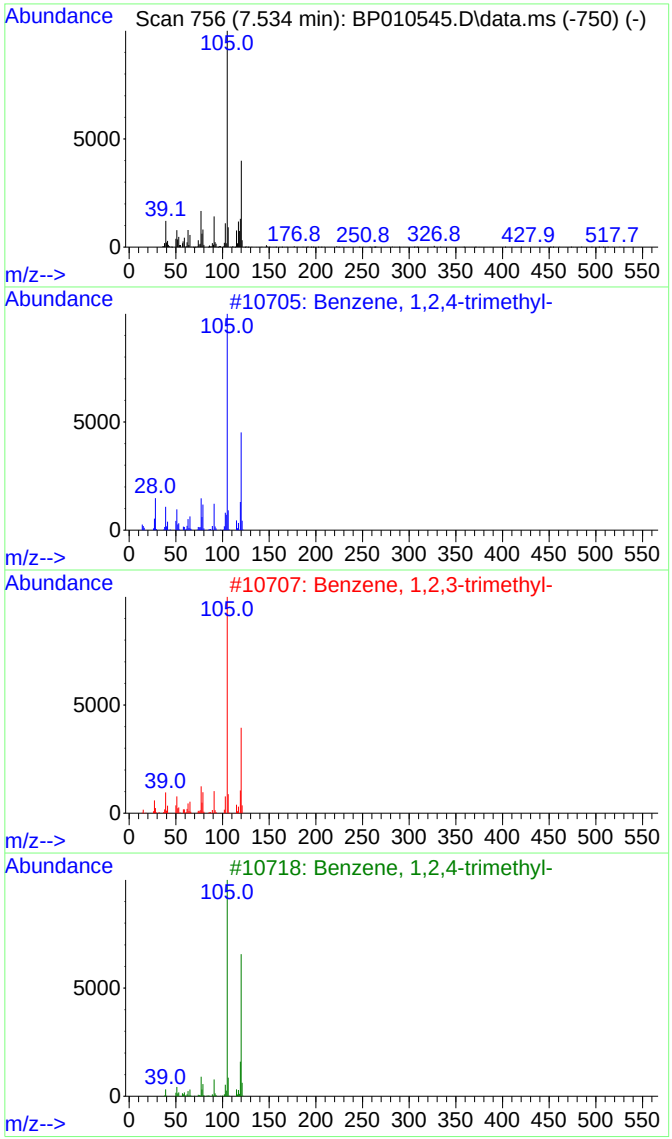
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

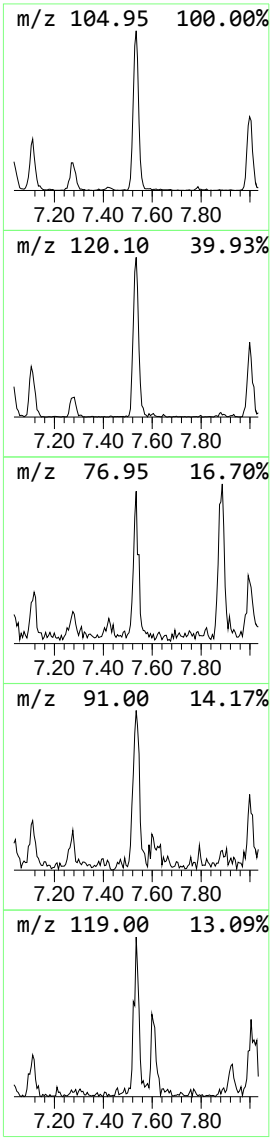
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzene, 1,2,4-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.534	2.34 ng/ul	89916	1,4-Dichlorobenzene-d4	7.881

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	87
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	87
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	87
4	Mesitylene	120	C9H12	000108-67-8	87
5	Mesitylene	120	C9H12	000108-67-8	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

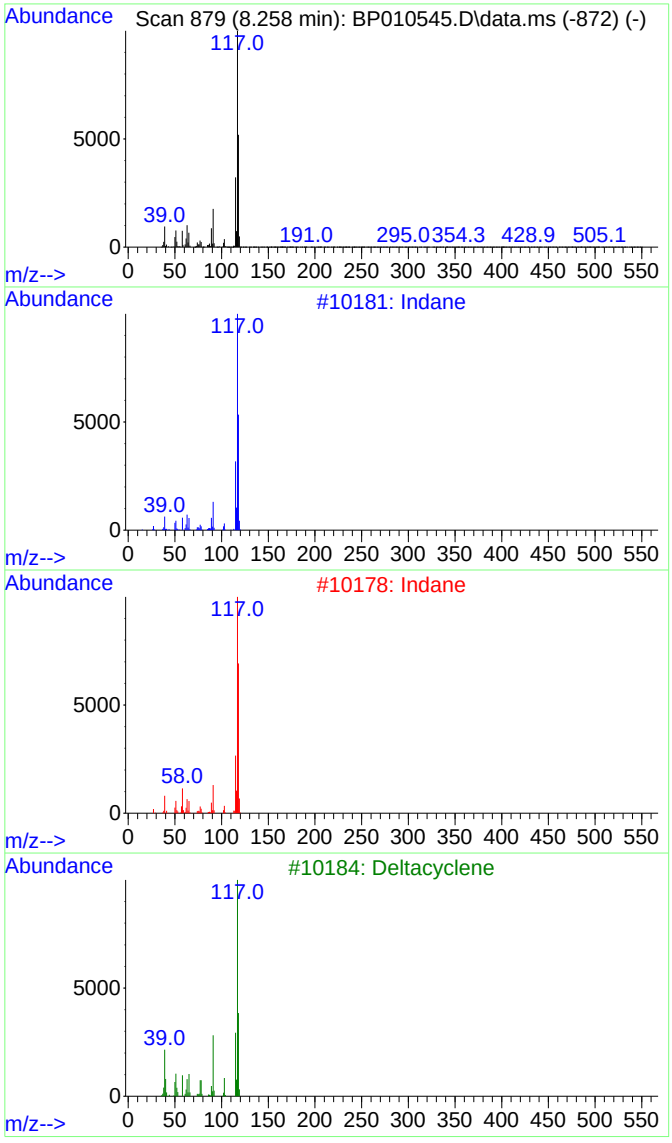
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

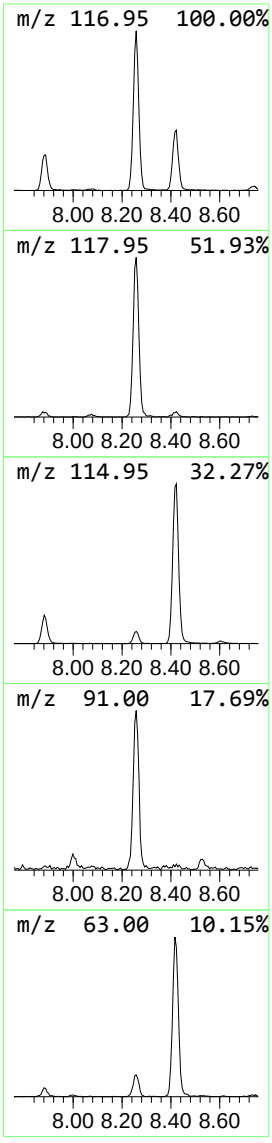
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.258	8.15 ng/ul	313203	1,4-Dichlorobenzene-d4	7.881

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Indane	118	C9H10	000496-11-7	81
3	Deltacyclene	118	C9H10	007785-10-6	68
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5	Indane	118	C9H10	000496-11-7	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

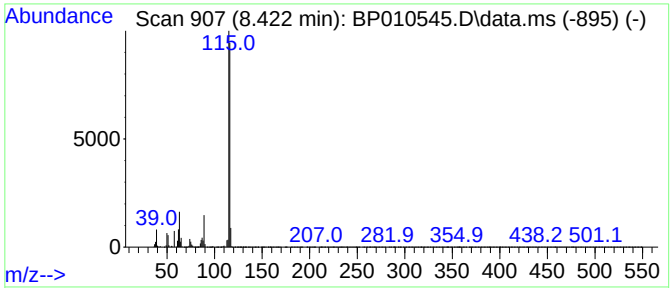
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

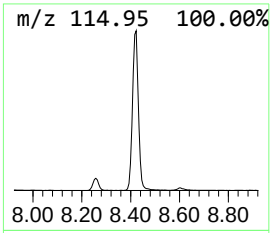
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

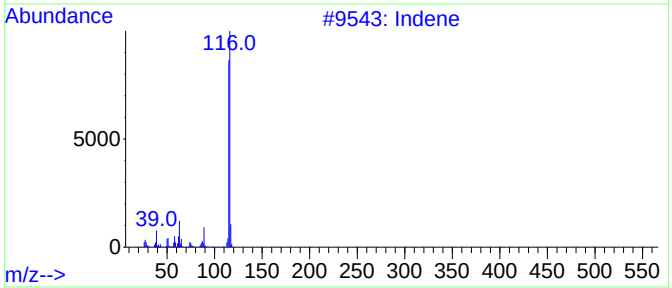
****
Peak Number 3 Indene Concentration Rank 1

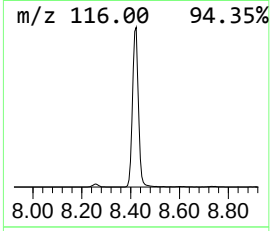
R.T.	EstConc	Area	Relative to ISTD	R.T.
8.422	37.07 ng/ul	1425470	1,4-Dichlorobenzene-d4	7.881

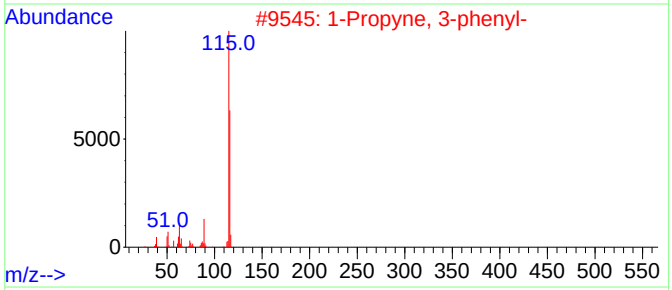
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	96
2	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	95
3	Indene	116	C9H8	000095-13-6	95
4	Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
5	2-Methylphenylacetylene	116	C9H8	000766-47-2	91

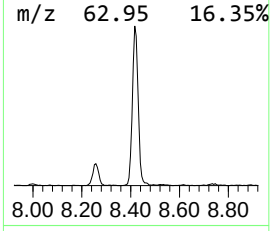


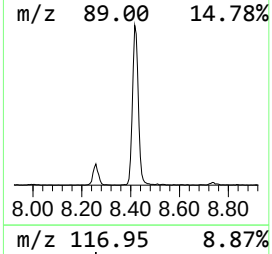


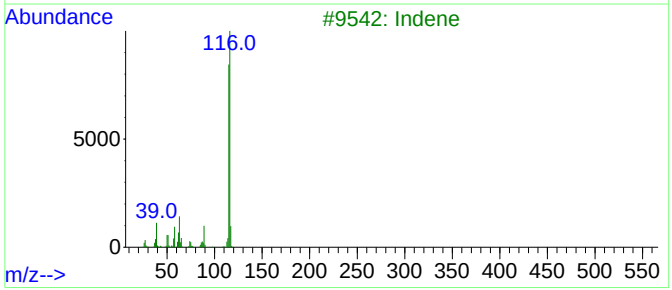


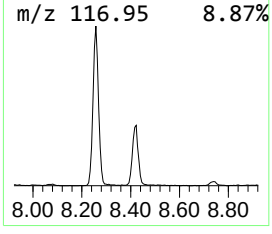












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

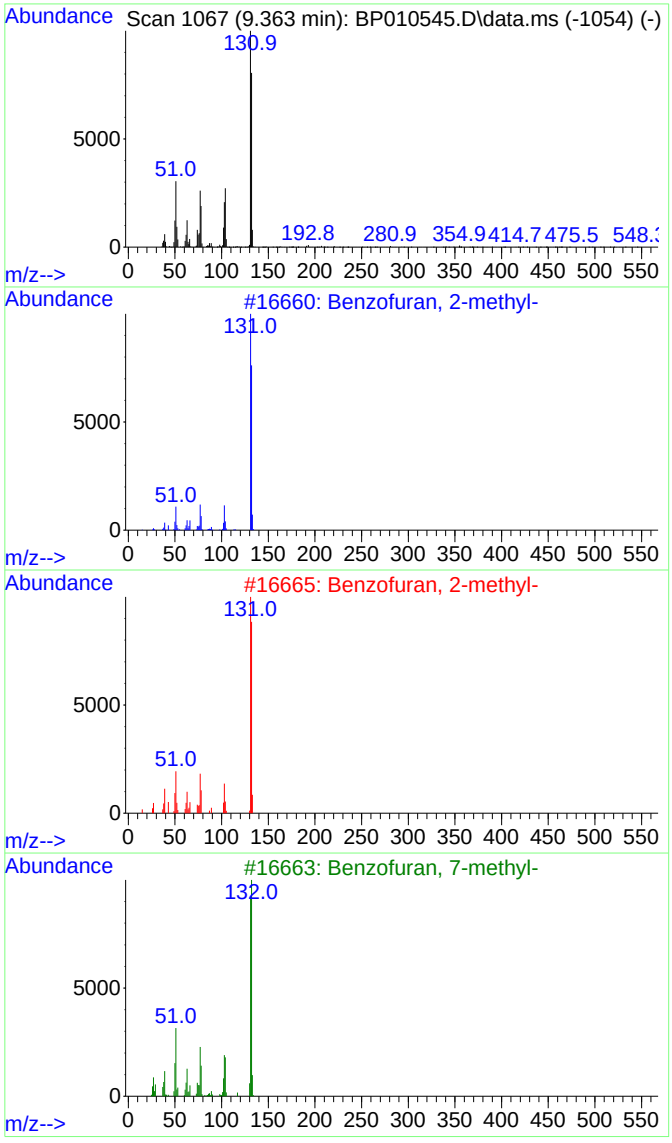
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

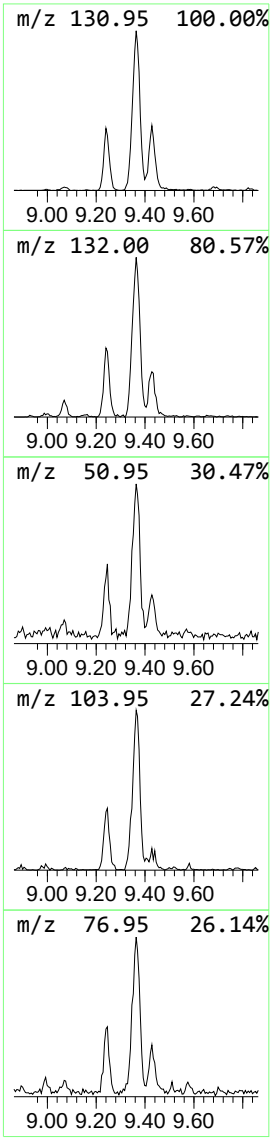
****

Peak Number 4 Benzofuran, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.363	4.02 ng/ul	212253	Naphthalene-d8	10.687

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	95
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
3	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	93
4	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91
5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

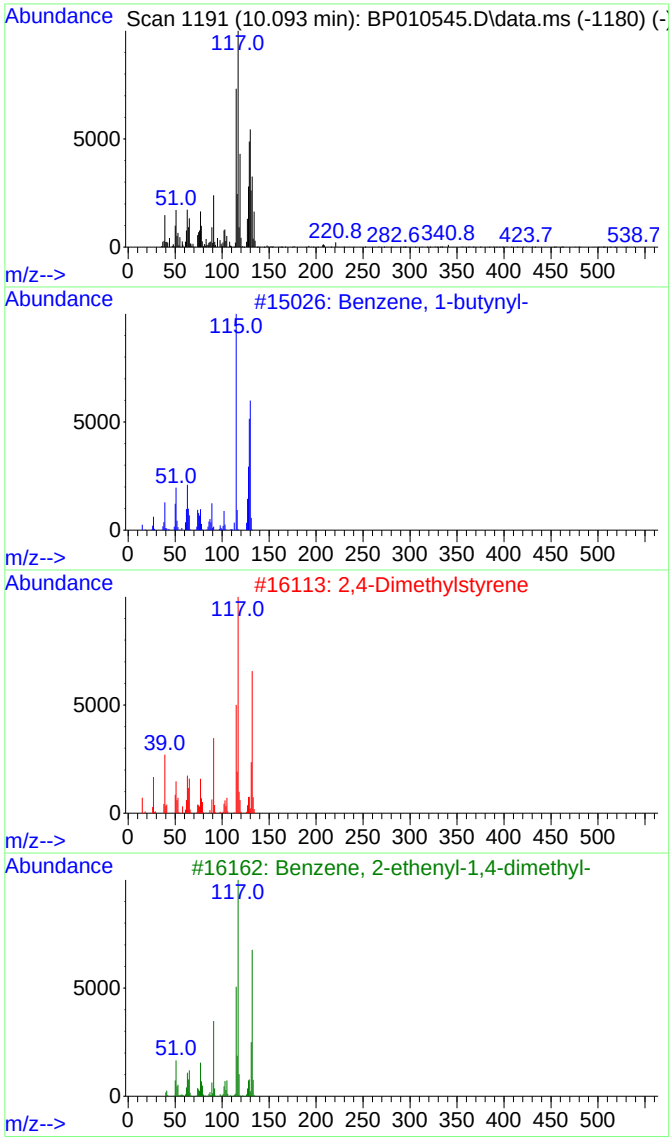
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

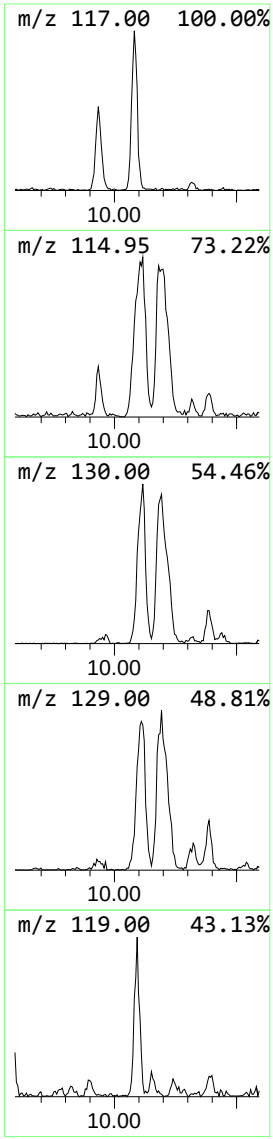
****

Peak Number 5 Benzene, 1-butynyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.093	3.71 ng/ul	196180	Naphthalene-d8	10.687

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-butynyl-	130	C10H10	000622-76-4	60
2	2,4-Dimethylstyrene	132	C10H12	002234-20-0	55
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	55
4	Benzene, 1-butynyl-	130	C10H10	000622-76-4	52
5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	46





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

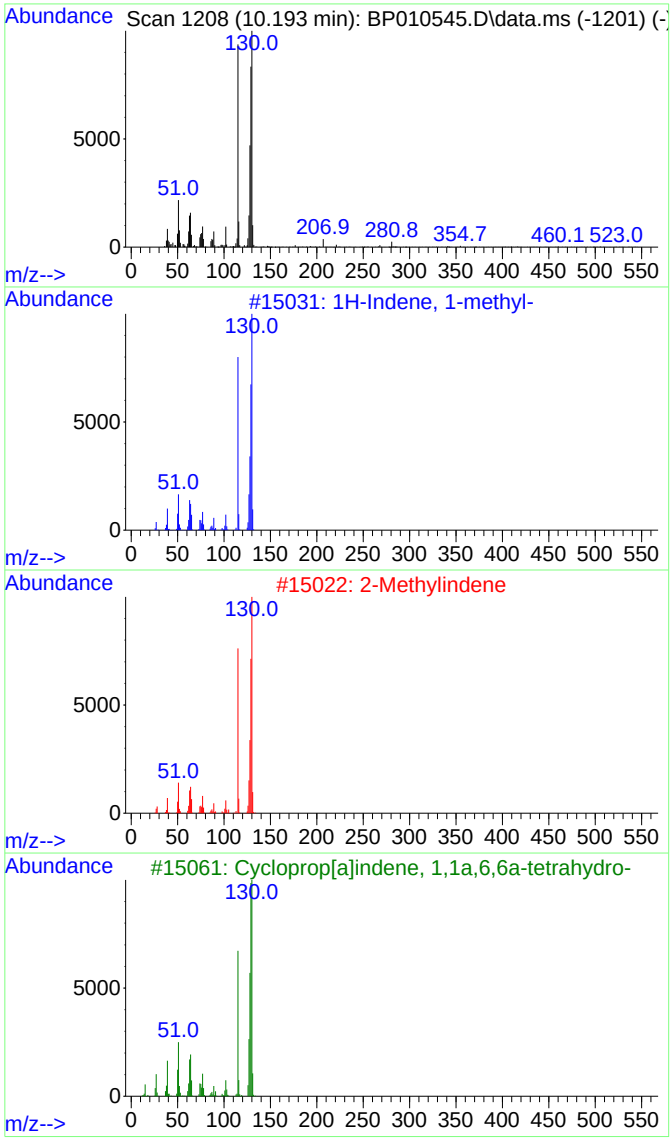
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

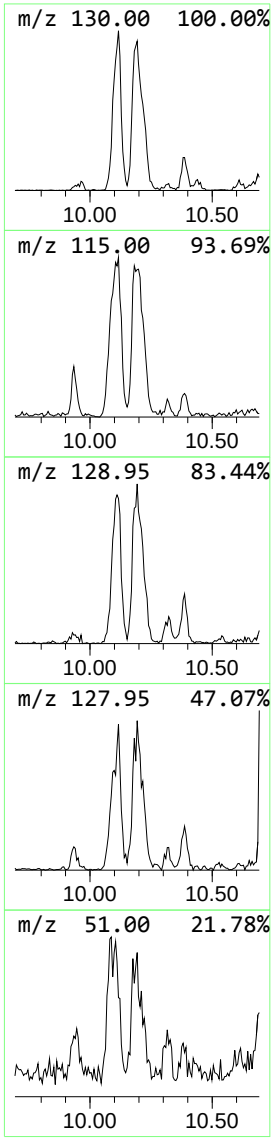
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 1H-Indene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.193	2.69 ng/ul	142387	Naphthalene-d8	10.687

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	95
2	2-Methylindene	130	C10H10	002177-47-1	95
3	Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	95
4	Benzene, 1-butynyl-	130	C10H10	000622-76-4	94
5	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	94





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Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

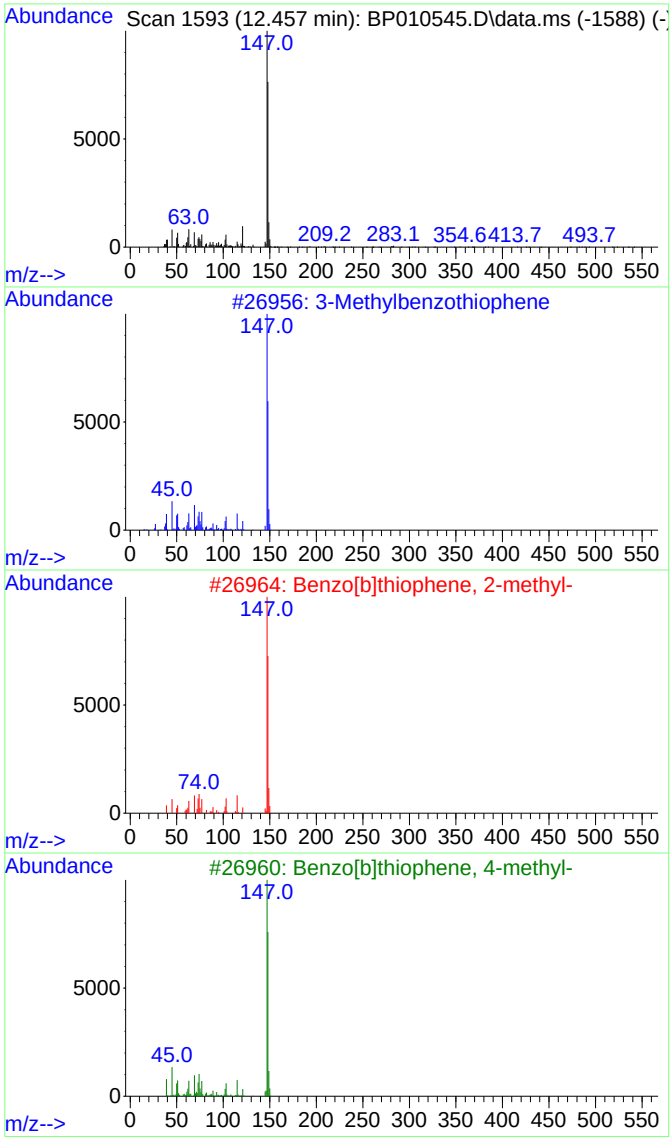
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

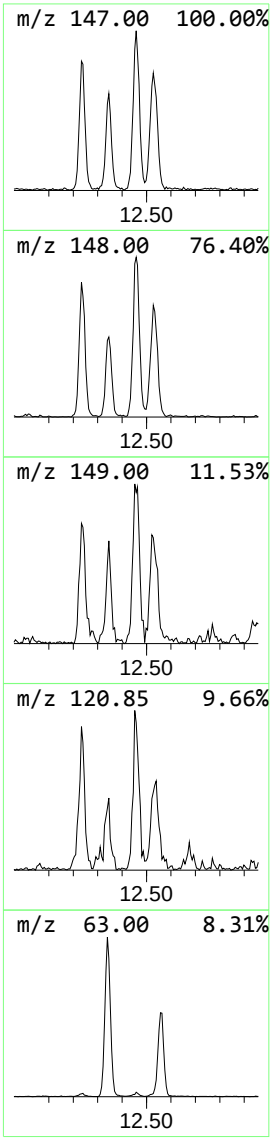
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 3-Methylbenzothiophene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.457	2.14 ng/ul	113144	Naphthalene-d8	10.687

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methylbenzothiophene	148	C9H8S	001455-18-1	91
2	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	90
3	Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	87
4	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	87
5	3-Methylbenzothiophene	148	C9H8S	001455-18-1	87





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Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

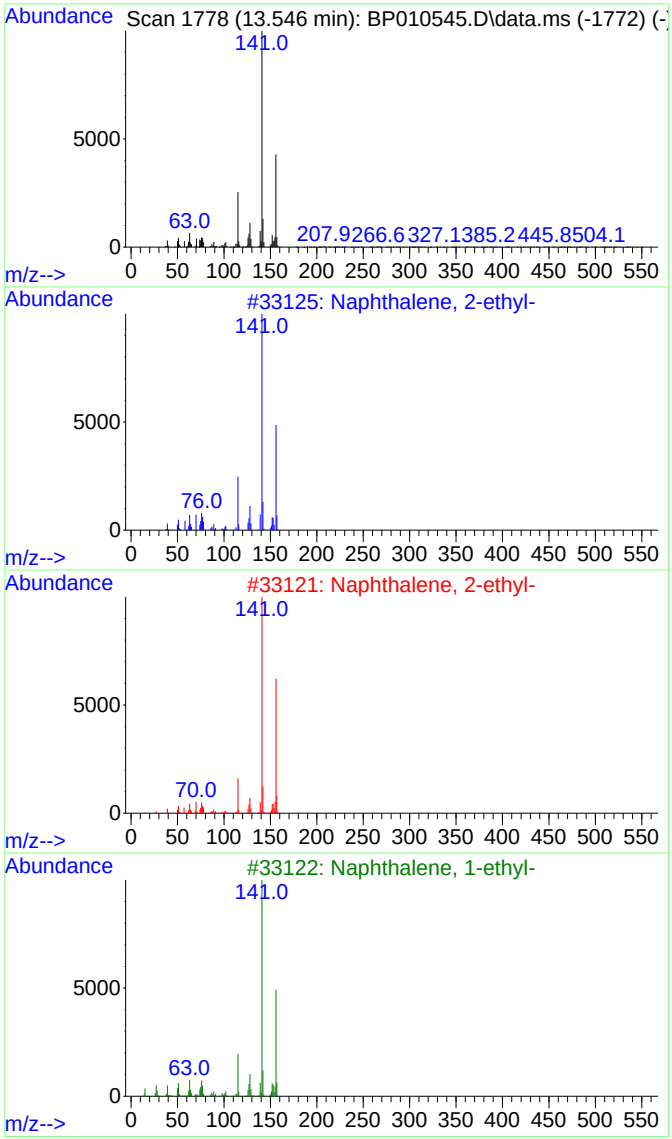
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

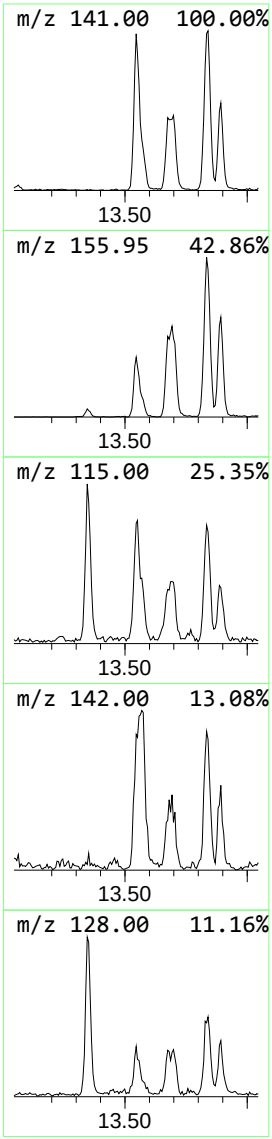
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Naphthalene, 2-ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.546	2.25 ng/ul	197573	Acenaphthene-d10	14.516

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	95
2	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	95
3	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	94
4	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	91
5	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	90





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Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

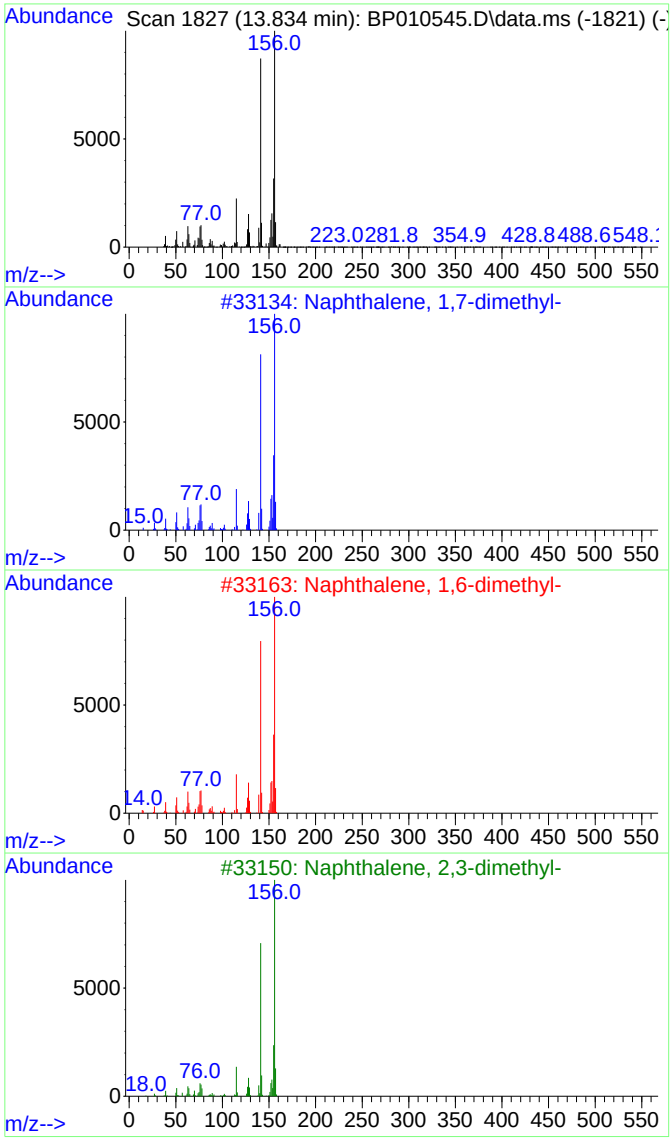
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

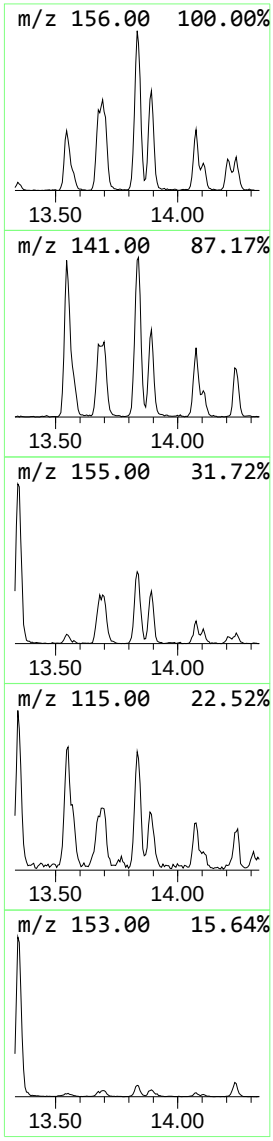
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Naphthalene, 1,7-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.834	2.87 ng/ul	251499	Acenaphthene-d10	14.516

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
2	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97





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Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

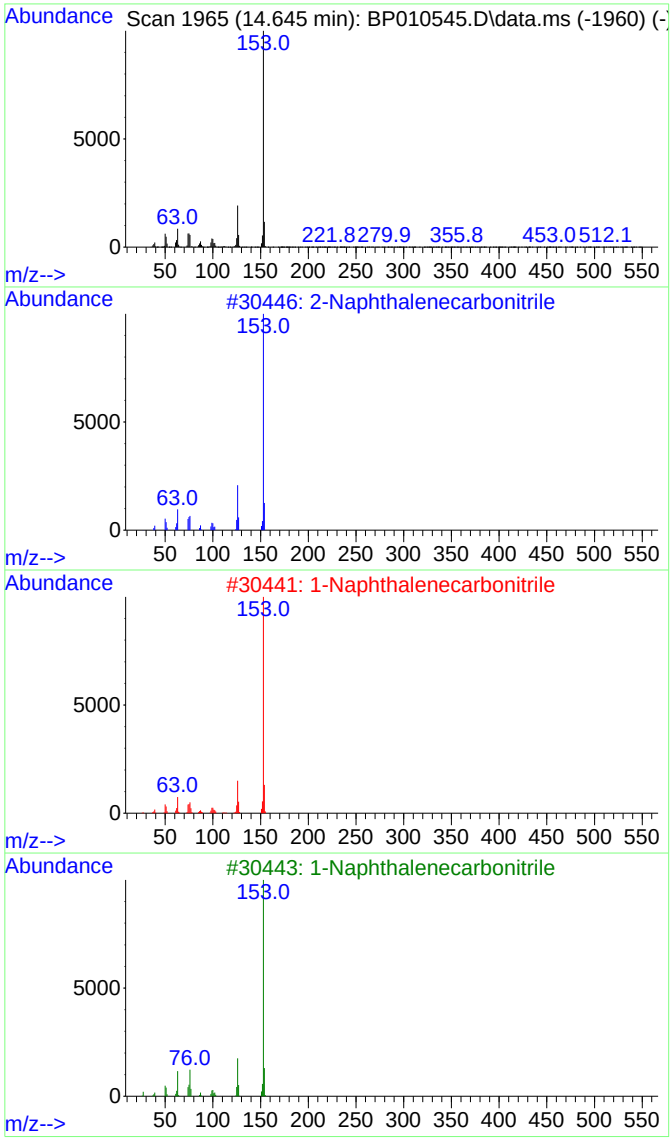
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

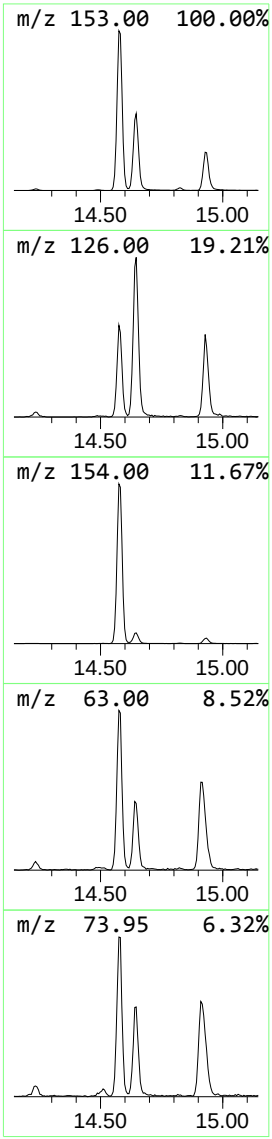
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 2-Naphthalenecarbonitrile Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.646	11.77 ng/ul	1031390	Acenaphthene-d10	14.516

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	97
2	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	95
3	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	94
4	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	91
5	2-Naphthyl isocyanide	153	C11H7N	010124-78-4	91





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Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

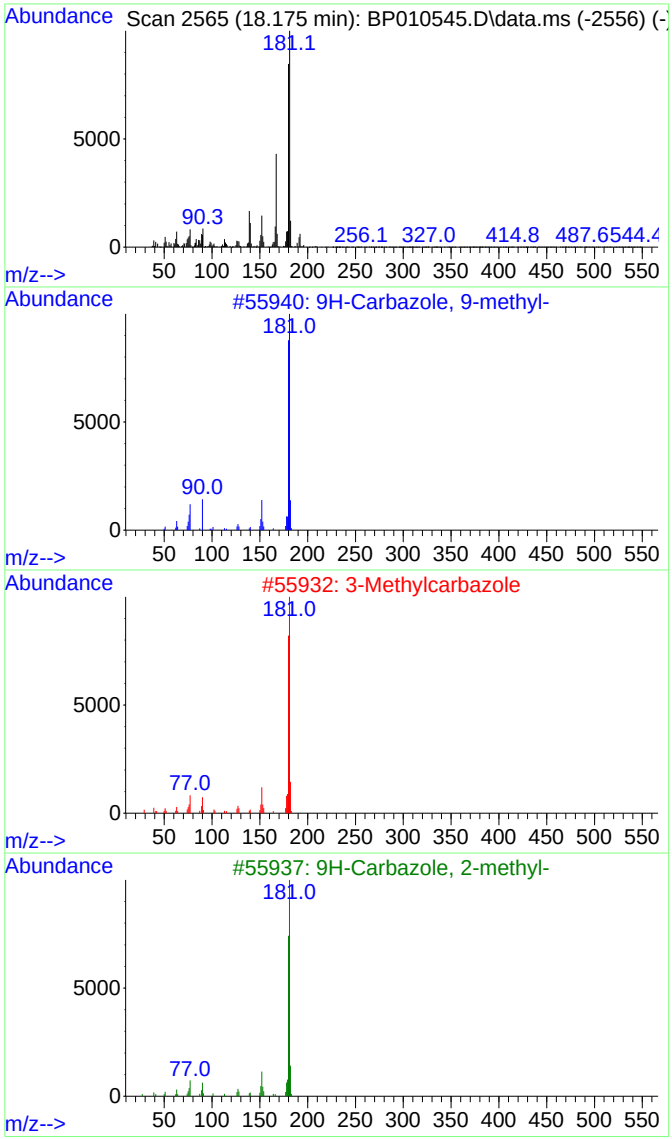
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

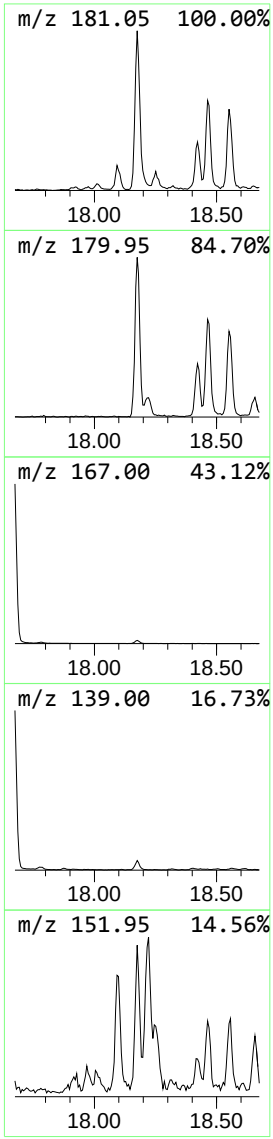
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 9H-Carbazole, 9-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.175	3.45 ng/ul	380824	Phenanthrene-d10	17.263

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Carbazole, 9-methyl-	181	C13H11N	001484-12-4	94
2	3-Methylcarbazole	181	C13H11N	004630-20-0	93
3	9H-Carbazole, 2-methyl-	181	C13H11N	003652-91-3	92
4	1-Methylcarbazole	181	C13H11N	006510-65-2	87
5	2-Fluorenamine	181	C13H11N	000153-78-6	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

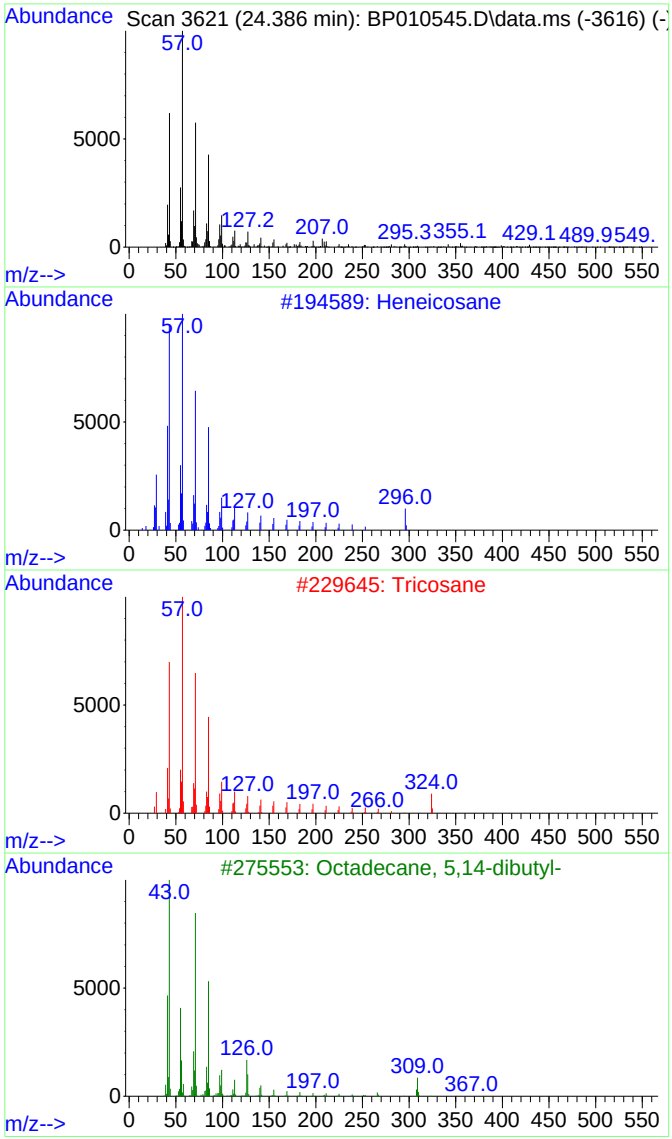
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

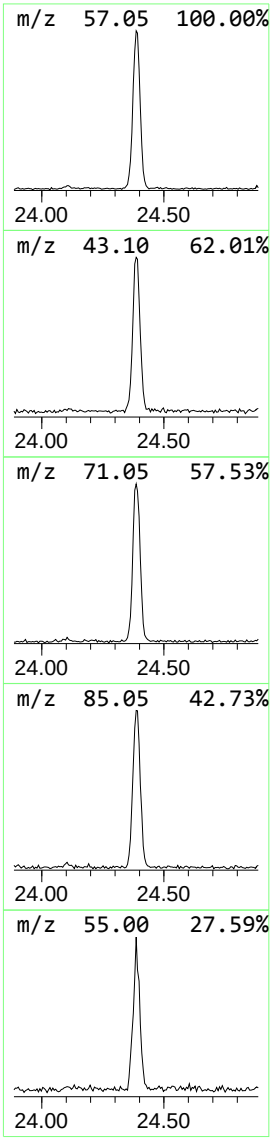
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.386	3.59 ng/ul	408239	Perylene-d12	23.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	95
2	Tricosane	324	C23H48	000638-67-5	95
3	Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	93
4	Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
5	Docosane, 9-butyl-	366	C26H54	055282-14-9	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

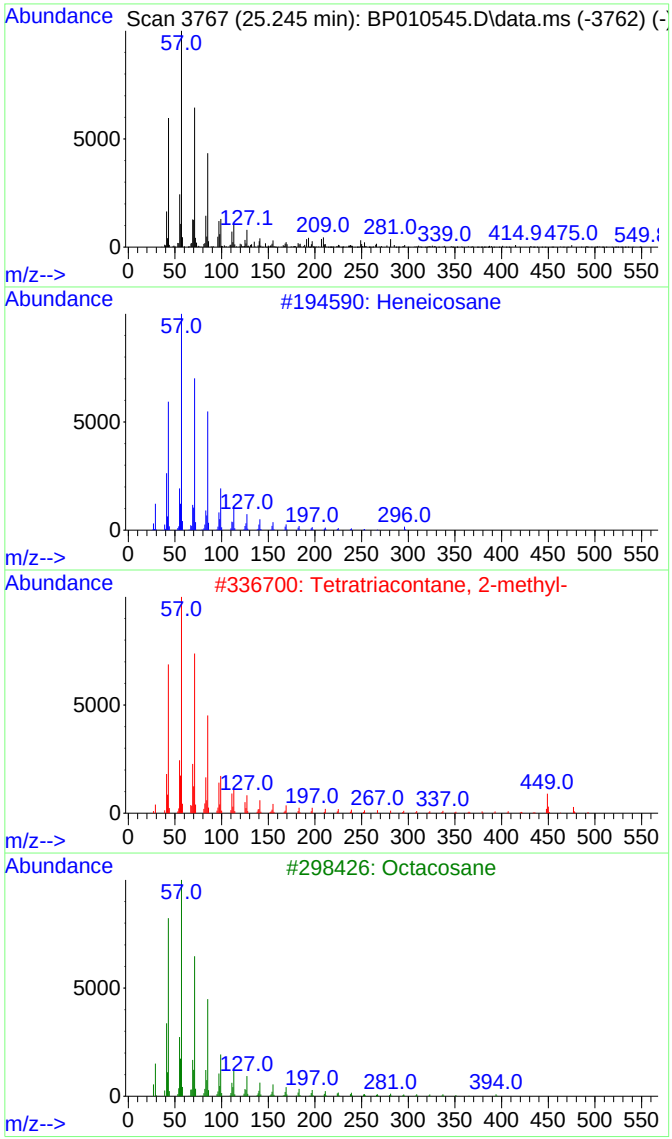
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

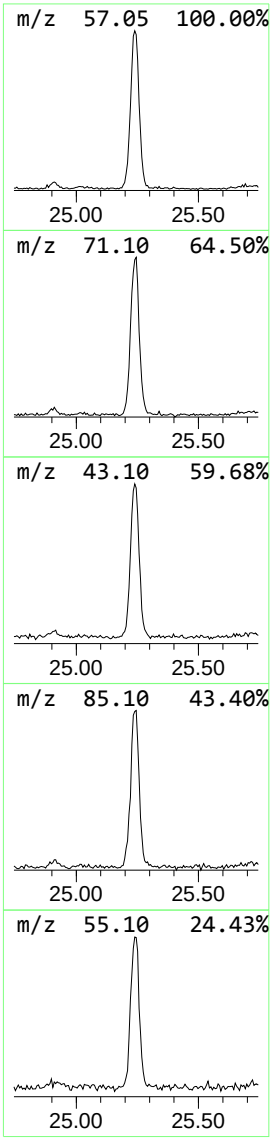
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.245	3.47 ng/ul	394444	Perylene-d12	23.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	96
2	Tetratriacontane, 2-methyl-	493	C35H72	014167-65-8	93
3	Octacosane	394	C28H58	000630-02-4	93
4	Eicosane, 9-octyl-	394	C28H58	013475-77-9	93
5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010545.D
Acq On : 25 May 2022 20:24
Operator : CG/JU
Sample : N2918-03DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE0DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1,2,4-...	7.534	2.3	ng/ul	89916	1	7.881	769030	20.0
Indane	8.258	8.2	ng/ul	313203	1	7.881	769030	20.0
Indene	8.422	37.1	ng/ul	1425470	1	7.881	769030	20.0
Benzofuran, 2-m...	9.363	4.0	ng/ul	212253	2	10.687	1056720	20.0
Benzene, 1-buty...	10.093	3.7	ng/ul	196180	2	10.687	1056720	20.0
1H-Indene, 1-me...	10.193	2.7	ng/ul	142387	2	10.687	1056720	20.0
3-Methylbenzoth...	12.457	2.1	ng/ul	113144	2	10.687	1056720	20.0
Naphthalene, 2-...	13.546	2.3	ng/ul	197573	3	14.516	1752980	20.0
Naphthalene, 1,...	13.834	2.9	ng/ul	251499	3	14.516	1752980	20.0
2-Naphthaleneca...	14.646	11.8	ng/ul	1031390	3	14.516	1752980	20.0
9H-Carbazole, 9...	18.175	3.5	ng/ul	380824	4	17.263	2208640	20.0
(DEL) Alkane: S...	24.386	3.6	ng/ul	408239	6	23.763	2272820	20.0
(DEL) Alkane: S...	25.245	3.5	ng/ul	394444	6	23.763	2272820	20.0