LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022215.D
Acq On : 04 Oct 2024 05:22
Operator : RC/JU
Sample : P4017-17DL 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DDCD3DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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

Signal : TIC: BP022215.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.704	622	632	637	rBV6	16061	35151	0.26%	0.106%
2	6.763	637	642	645	rVV2	17960	26415	0.20%	0.080%
3	6.804	645	649	654	rVV	21557	33650	0.25%	0.102%
4	6.869	654	660	667	rVV5	28560	60573	0.45%	0.183%
5	6.934	667	671	677	rVB	16700	25592	0.19%	0.077%
6	7.093	693	698	702	rVV3	16541	25824	0.19%	0.078%
7	7.328	733	738	746	rBV2	98776	172735	1.28%	0.523%
8	7.610	777	786	792	rBV9	9458	28422	0.21%	0.086%
9	7.693	792	800	807	rVV	453290	763910	5.66%	2.312%
10	7.757	807	811	817	rVV2	54479	96141	0.71%	0.291%
11	7.810	817	820	828	rVB3	16987	29638	0.22%	0.090%
12	7.981	846	849	857	rVB6	12653	22696	0.17%	0.069%
13	8.063	857	863	868	rVB	13935	25789	0.19%	0.078%
14	8.122	868	873	877	rBV2	26324	43517	0.32%	0.132%
15	8.216	886	889	896	rVB3	23010	45759	0.34%	0.138%
16	8.275	896	899	903	rVB	20186	25400	0.19%	0.077%
17	8.345	903	911	915	rBV3	36782	90957	0.67%	0.275%
18	8.451	922	929	932	rBV	32781	49184	0.36%	0.149%
19	8.487	932	935	944	rVB2	19594	43147	0.32%	0.131%
20	8.634	954	960	963	rBV2	13927	23274	0.17%	0.070%
21	8.675	963	967	976	rVB3	37999	73178	0.54%	0.221%
22	8.787	976	986	992	rBV3	34097	77545	0.57%	0.235%
23	8.904	998	1006	1011	rVB	127287	196603	1.46%	0.595%
24	9.022	1017	1026	1032	rBV10	15711	41436	0.31%	0.125%
25	9.363	1079	1084	1092	rVB4	10141	24767	0.18%	0.075%
26	9.463	1092	1101	1107	rVB3	25825	52513	0.39%	0.159%
27	9.551	1107	1116	1121	rBV6	15822	36651	0.27%	0.111%
28	9.610	1121	1126	1131	rVV	29771	49156	0.36%	0.149%
29	9.675	1131	1137	1141	rVV3	9701	21735	0.16%	0.066%
30	9.928	1176	1180	1187	rVV3	15335	28999	0.21%	0.088%
31	10.145	1212	1217	1228	rVB2	66754	111601	0.83%	0.338%
32	10.445	1261	1268	1277	rBV2	1012879	1868909	13.84%	5.656%
33	10.569	1283	1289	1296	rBV4	152334	261179	1.93%	0.790%
34	10.739	1312	1318	1321	rBV3	14212	25177	0.19%	0.076%
35	10.828	1327	1333	1342	rVB	208976	343184	2.54%	1.039%
36	11.181	1385	1393	1402	rVB5	23379	51964	0.38%	0.157%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022215.D
Acq On : 04 Oct 2024 05:22
Operator : RC/JU
Sample : P4017-17DL 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DDCD3DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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

37	11.339	1415	1420	1429	rBV4	56075	135436	1.00%	0.410%
38	11.692	1472	1480	1482	rBV3	39117	71865	0.53%	0.217%
39	11.869	1502	1510	1513	rBV2	53667	90726	0.67%	0.275%
40	11.910	1513	1517	1522	rVB	52373	69950	0.52%	0.212%
41	12.110	1547	1551	1558	rVB2	14021	26081	0.19%	0.079%
42	12.622	1632	1638	1645	rVB2	12836	30471	0.23%	0.092%
43	12.869	1676	1680	1687	rVB	54146	89554	0.66%	0.271%
44	13.122	1716	1723	1729	rBV3	47884	87012	0.64%	0.263%
45	13.333	1755	1759	1769	rVB9	12155	33338	0.25%	0.101%
46	13.486	1776	1785	1792	rBV6	32987	78420	0.58%	0.237%
47	13.622	1798	1808	1812	rBV7	30074	86403	0.64%	0.261%
48	13.792	1829	1837	1842	rBV9	25135	59069	0.44%	0.179%
49	13.857	1842	1848	1852	rVV8	18758	33858	0.25%	0.102%
50	13.933	1857	1861	1863	rBV2	39140	49584	0.37%	0.150%
51	14.210	1904	1908	1911	rBV	63558	73226	0.54%	0.222%
52	14.251	1911	1915	1917	rVV4	32296	41943	0.31%	0.127%
53	14.298	1917	1923	1933	rVV	942091	1608527	11.91%	4.868%
54	14.386	1933	1938	1943	rVB2	82284	136264	1.01%	0.412%
55	14.527	1957	1962	1964	rBV4	37104	60911	0.45%	0.184%
56	14.710	1990	1993	1999	rVB7	18094	28840	0.21%	0.087%
57	14.786	2001	2006	2009	rBV4	15857	21882	0.16%	0.066%
58	14.839	2010	2015	2016	rBV4	22227	27707	0.21%	0.084%
59	14.933	2021	2031	2046	rVB	427763	750330	5.55%	2.271%
60	15.098	2054	2059	2065	rBV7	20040	53072	0.39%	0.161%
61	15.174	2065	2072	2077	rVB3	85260	182349	1.35%	0.552%
62	15.304	2090	2094	2098	rBV2	53510	67581	0.50%	0.205%
63	15.586	2136	2142	2148	rVB3	40134	80741	0.60%	0.244%
64	15.686	2154	2159	2165	rBV6	27315	48920	0.36%	0.148%
65	15.739	2165	2168	2173	rVV6	12931	22168	0.16%	0.067%
66	15.798	2175	2178	2180	rVV4	22230	33123	0.25%	0.100%
67	15.833	2182	2184	2189	rVV2	35252	51690	0.38%	0.156%
68	15.886	2189	2193	2197	rVB3	35575	46862	0.35%	0.142%
69	16.051	2216	2221	2225	rBV2	111786	185631	1.37%	0.562%
70	16.221	2247	2250	2255	rVB6	31300	48416	0.36%	0.147%
71	16.274	2255	2259	2265	rVB5	54342	82588	0.61%	0.250%
72	16.398	2277	2280	2284	rVB5	15106	22638	0.17%	0.069%
73	16.521	2299	2301	2305	rVB3	24132	24930	0.18%	0.075%
74	16.610	2312	2316	2321	rBV2	151941	201404	1.49%	0.609%
75	16.751	2332	2340	2350	rBV	8034338	13507606	100.00%	40.876%
76	16.833	2351	2354	2356	rVV2	63537	88691	0.66%	0.268%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022215.D
Acq On : 04 Oct 2024 05:22
Operator : RC/JU
Sample : P4017-17DL 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DDCD3DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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

77	16.863	2356	2359	2363	rVV	136566	180725	1.34%	0.547%
78	16.916	2365	2368	2373	rVB4	68904	103912	0.77%	0.314%
79	17.010	2379	2384	2390	rVB2	68426	109624	0.81%	0.332%
80	17.098	2392	2399	2405	rBV	1371403	2062041	15.27%	6.240%
81	17.198	2413	2416	2422	rVB	52277	83094	0.62%	0.251%
82	17.333	2436	2439	2448	rVB4	44388	78701	0.58%	0.238%
83	17.480	2461	2464	2469	rBV5	24194	42331	0.31%	0.128%
84	17.621	2484	2488	2493	rBV	119939	159158	1.18%	0.482%
85	17.733	2505	2507	2512	rVB5	47064	59441	0.44%	0.180%
86	17.845	2523	2526	2534	rBV7	24887	52963	0.39%	0.160%
87	17.916	2536	2538	2543	rVB5	23278	31112	0.23%	0.094%
88	18.027	2554	2557	2565	rVB4	53846	98842	0.73%	0.299%
89	18.345	2606	2611	2619	rVB	118785	179126	1.33%	0.542%
90	18.563	2645	2648	2653	rBV4	33011	50515	0.37%	0.153%
91	18.998	2719	2722	2729	rVB	116728	140763	1.04%	0.426%
92	19.363	2781	2784	2786	rBV3	17517	22333	0.17%	0.068%
93	19.580	2817	2821	2823	rBV2	52355	64229	0.48%	0.194%
94	19.615	2823	2827	2831	rVB	72624	91416	0.68%	0.277%
95	20.168	2917	2921	2926	rBV	100389	111272	0.82%	0.337%
96	20.686	3007	3009	3013	rVB2	61045	66166	0.49%	0.200%
97	21.204	3093	3097	3104	rVB	326142	401194	2.97%	1.214%
98	21.521	3145	3151	3160	rBV2	1474605	2527111	18.71%	7.647%
99	22.398	3296	3300	3309	rVB3	115941	234968	1.74%	0.711%
100	24.797	3700	3708	3719	rVB	978225	2722491	20.16%	8.239%

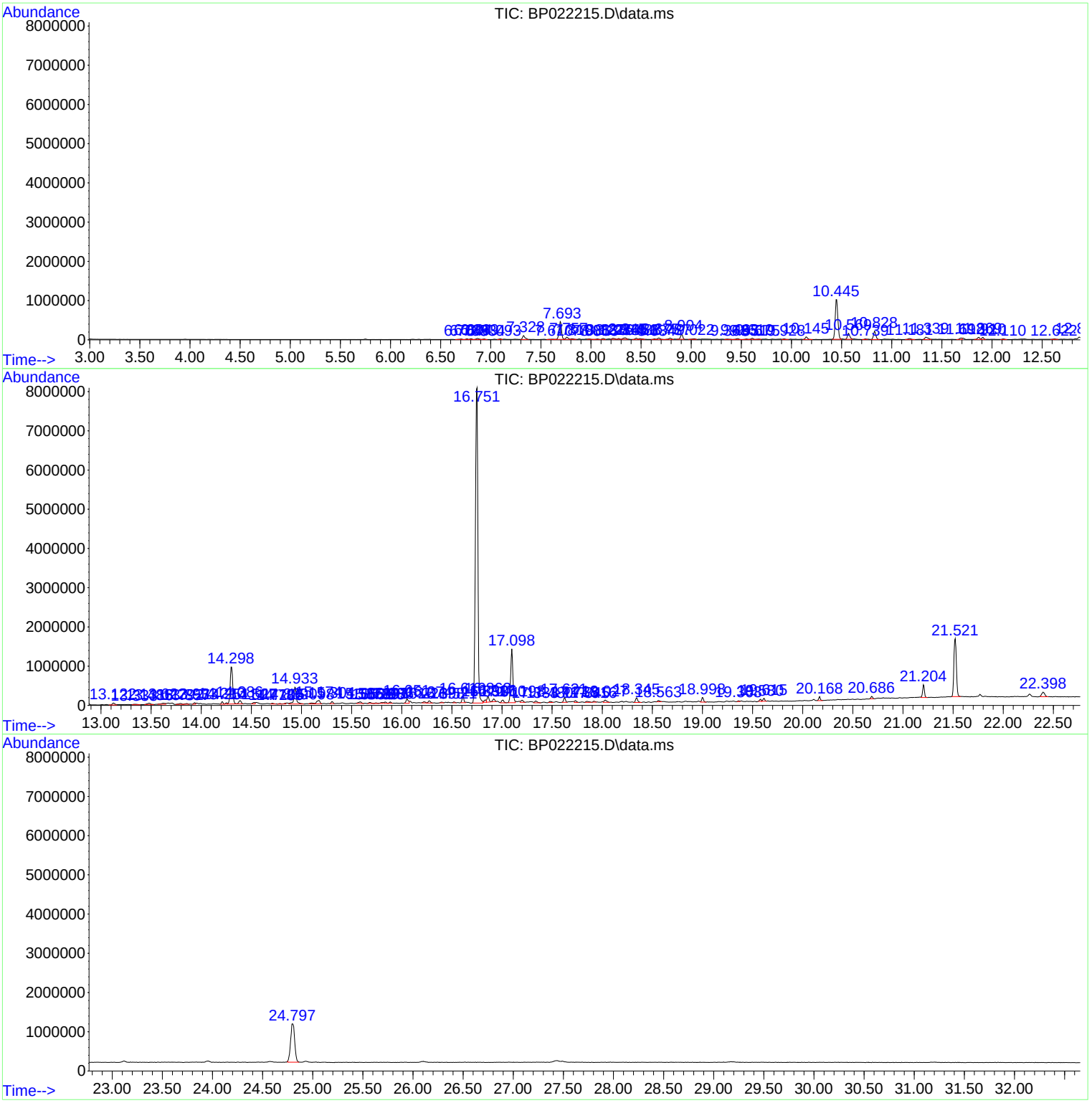
Sum of corrected areas: 33045706

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022215.D
Acq On : 04 Oct 2024 05:22
Operator : RC/JU
Sample : P4017-17DL 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCD3DL

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

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022215.D
Acq On : 04 Oct 2024 05:22
Operator : RC/JU
Sample : P4017-17DL 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCD3DL

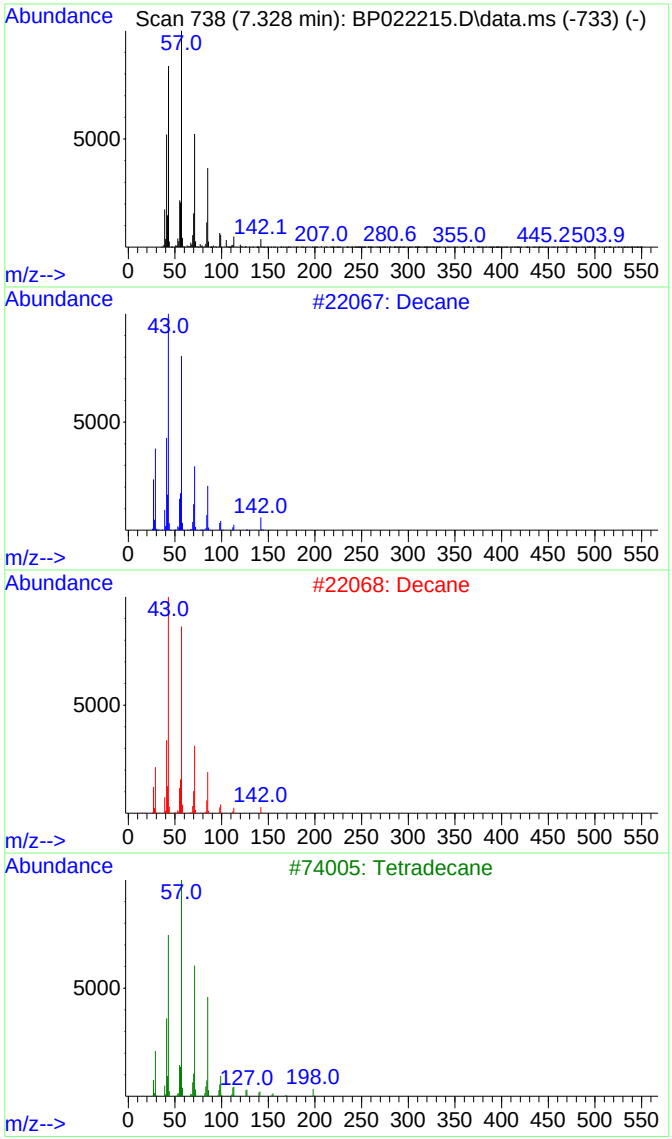
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

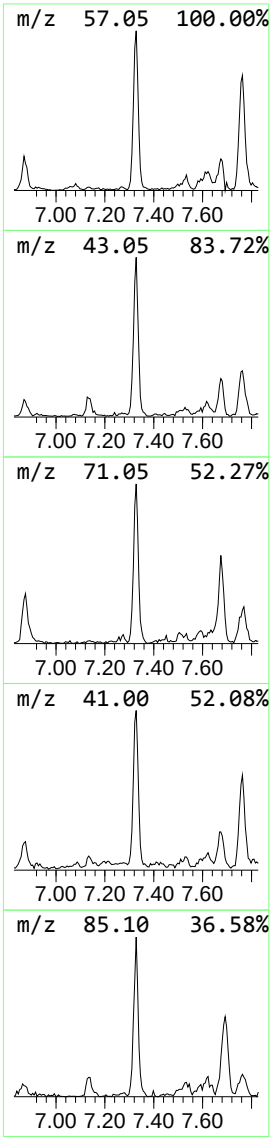
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.328	4.52 ng/ul	172735	1,4-Dichlorobenzene-d4	7.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	94
2	Decane	142	C10H22	000124-18-5	91
3	Tetradecane	198	C14H30	000629-59-4	86
4	Undecane	156	C11H24	001120-21-4	83
5	Tridecane	184	C13H28	000629-50-5	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022215.D
Acq On : 04 Oct 2024 05:22
Operator : RC/JU
Sample : P4017-17DL 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCD3DL

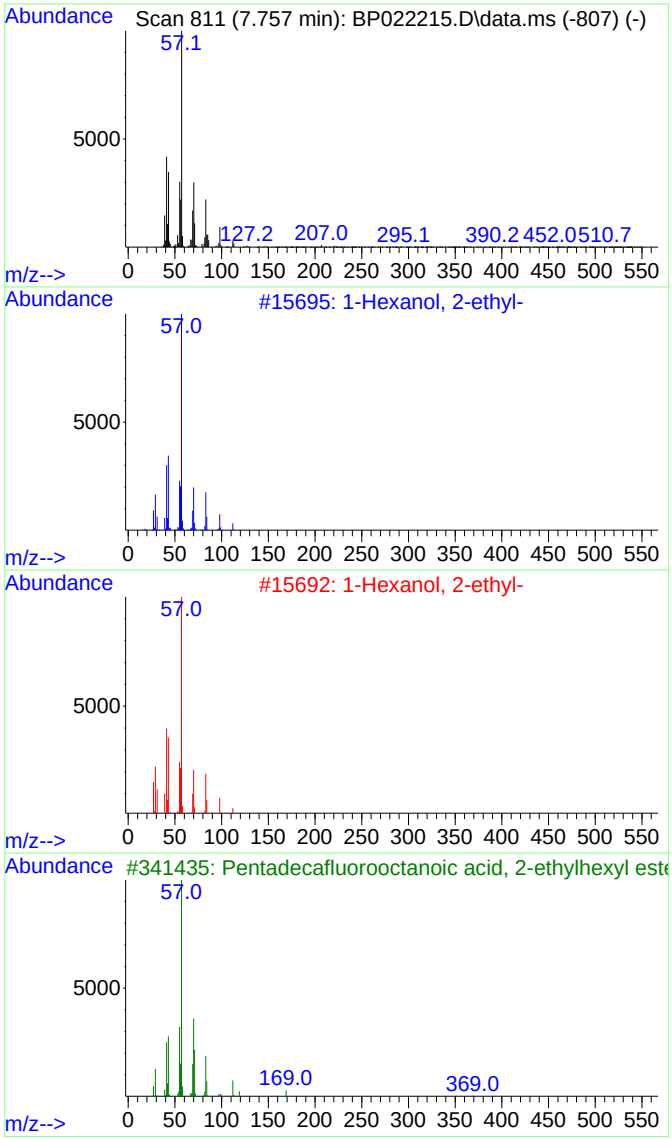
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

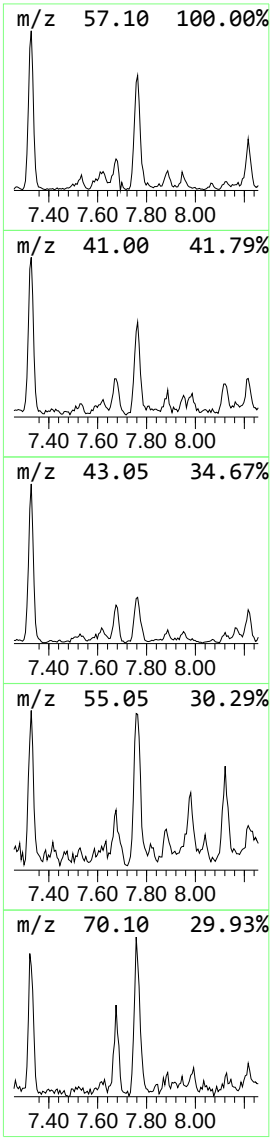
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 1-Hexanol, 2-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.757	2.52 ng/ul	96141	1,4-Dichlorobenzene-d4	7.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64
2	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	59
3	Pentadecafluorooctanoic acid, 2-...	526	C16H17F15O2	1000406-80-9	53
4	2-Propyl-1-pentanol	130	C8H18O	058175-57-8	53
5	Carbonic acid, octyl vinyl ester	200	C11H20O3	1000383-25-5	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022215.D
Acq On : 04 Oct 2024 05:22
Operator : RC/JU
Sample : P4017-17DL 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCD3DL

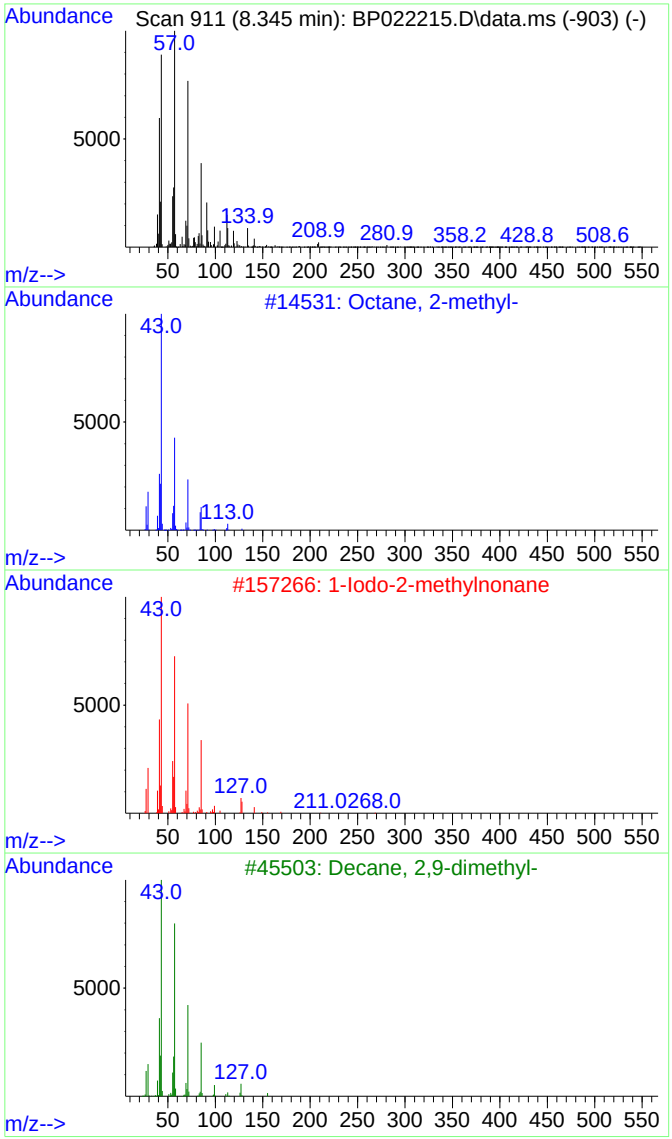
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

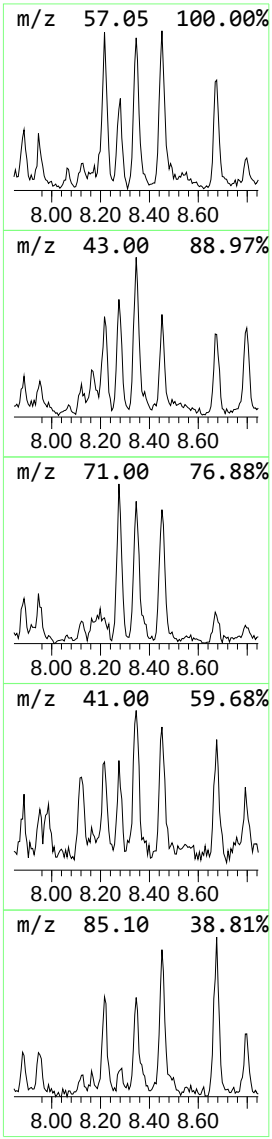
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.345	2.38 ng/ul	90957	1,4-Dichlorobenzene-d4	7.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2-methyl-	128	C9H20	003221-61-2	87
2	1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	87
3	Decane, 2,9-dimethyl-	170	C12H26	001002-17-1	78
4	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	72
5	Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	68





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022215.D
Acq On : 04 Oct 2024 05:22
Operator : RC/JU
Sample : P4017-17DL 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCD3DL

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

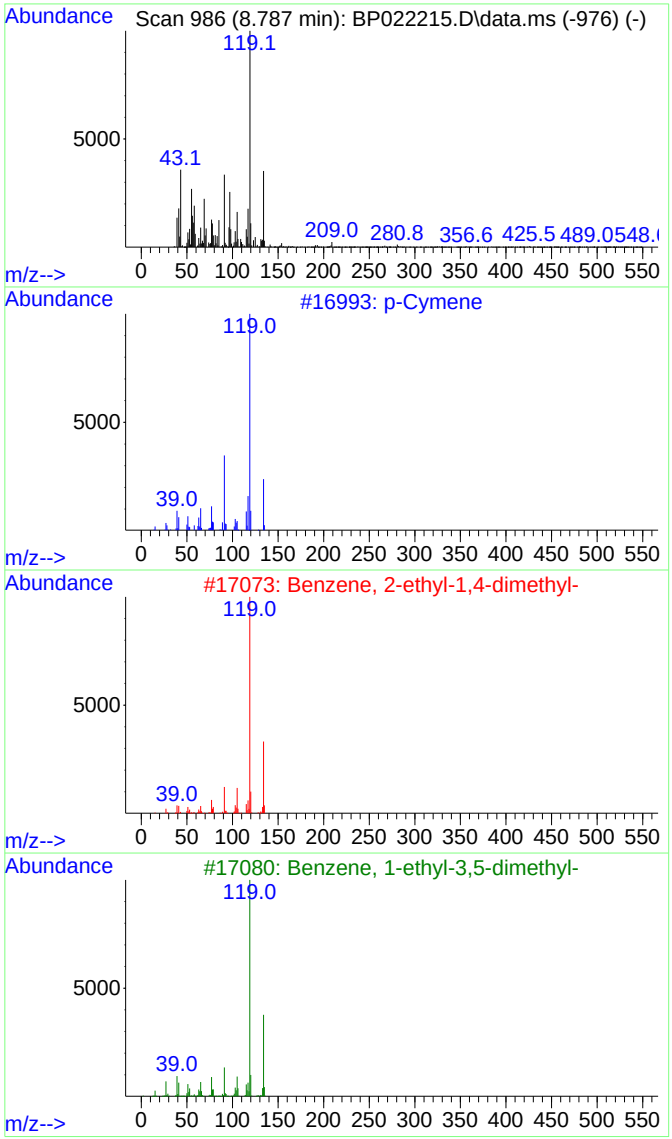
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

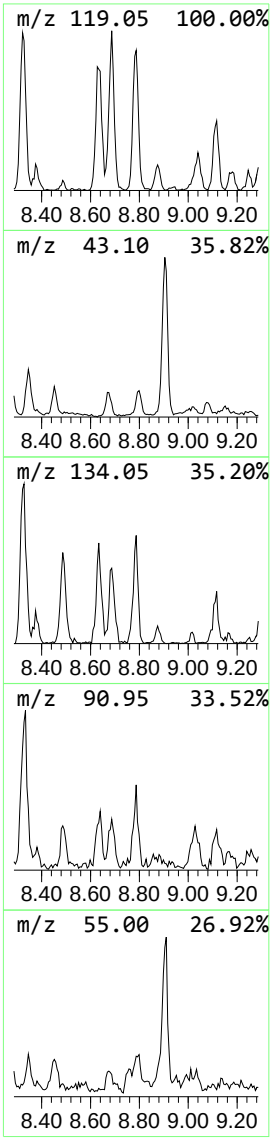
****

Peak Number 4 p-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.787	2.03 ng/ul	77545	1,4-Dichlorobenzene-d4	7.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	64
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	60
3	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	60
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	58
5	p-Cymene	134	C10H14	000099-87-6	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022215.D
Acq On : 04 Oct 2024 05:22
Operator : RC/JU
Sample : P4017-17DL 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCD3DL

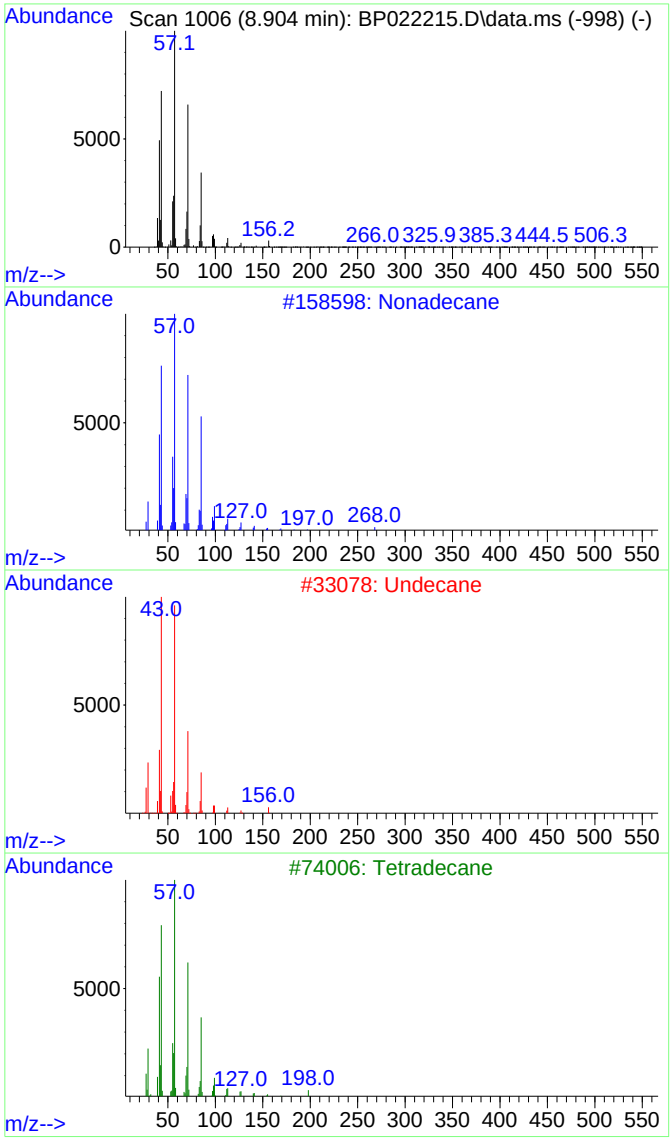
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

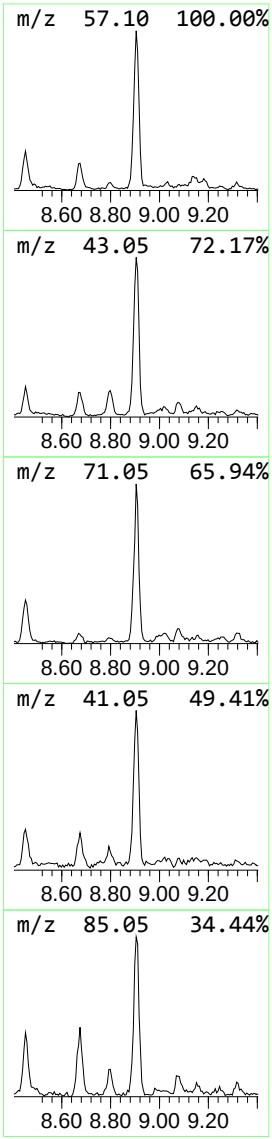
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.904	5.15 ng/ul	196603	1,4-Dichlorobenzene-d4	7.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	90
2	Undecane	156	C11H24	001120-21-4	80
3	Tetradecane	198	C14H30	000629-59-4	78
4	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	72
5	Nonadecane	268	C19H40	000629-92-5	72





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Data File : BP022215.D
Acq On : 04 Oct 2024 05:22
Operator : RC/JU
Sample : P4017-17DL 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCD3DL

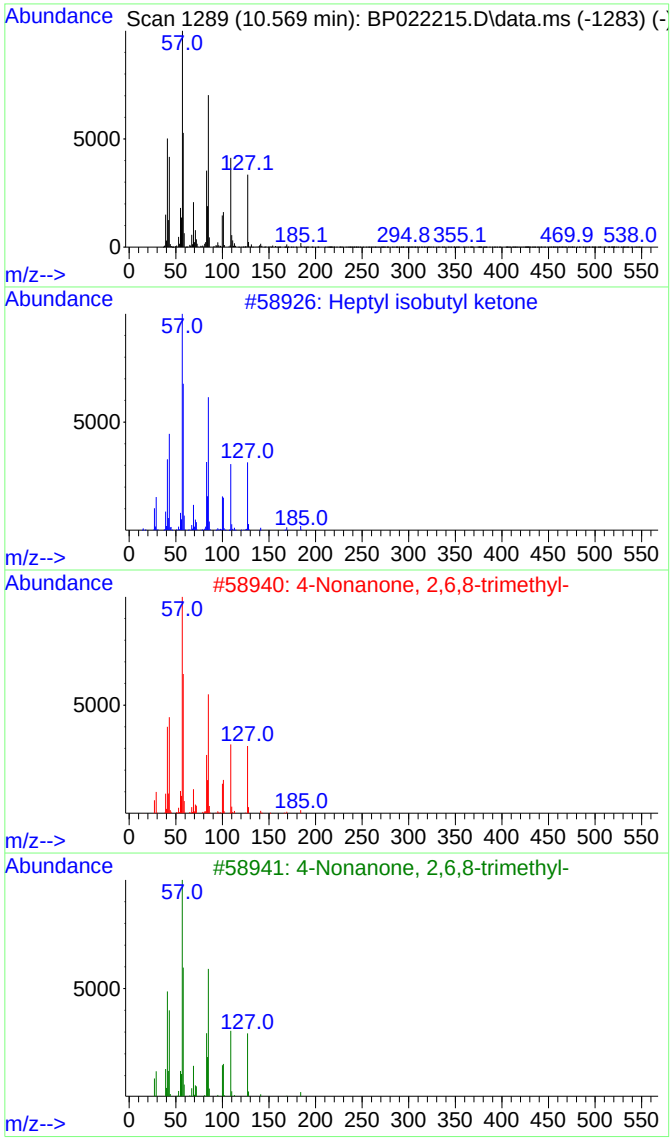
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

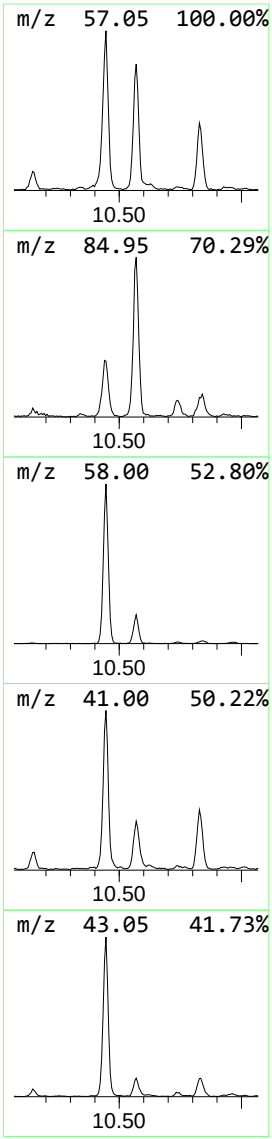
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Heptyl isobutyl ketone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.569	2.79 ng/ul	261179	Naphthalene-d8	10.457

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptyl isobutyl ketone	184	C12H24O	019594-40-2	95
2	4-Nonanone, 2,6,8-trimethyl-	184	C12H24O	000123-18-2	93
3	4-Nonanone, 2,6,8-trimethyl-	184	C12H24O	000123-18-2	93
4	5-Nonanone	142	C9H18O	000502-56-7	38
5	4-Octanone, 2-methyl-	142	C9H18O	007492-38-8	38





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Data File : BP022215.D
Acq On : 04 Oct 2024 05:22
Operator : RC/JU
Sample : P4017-17DL 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCD3DL

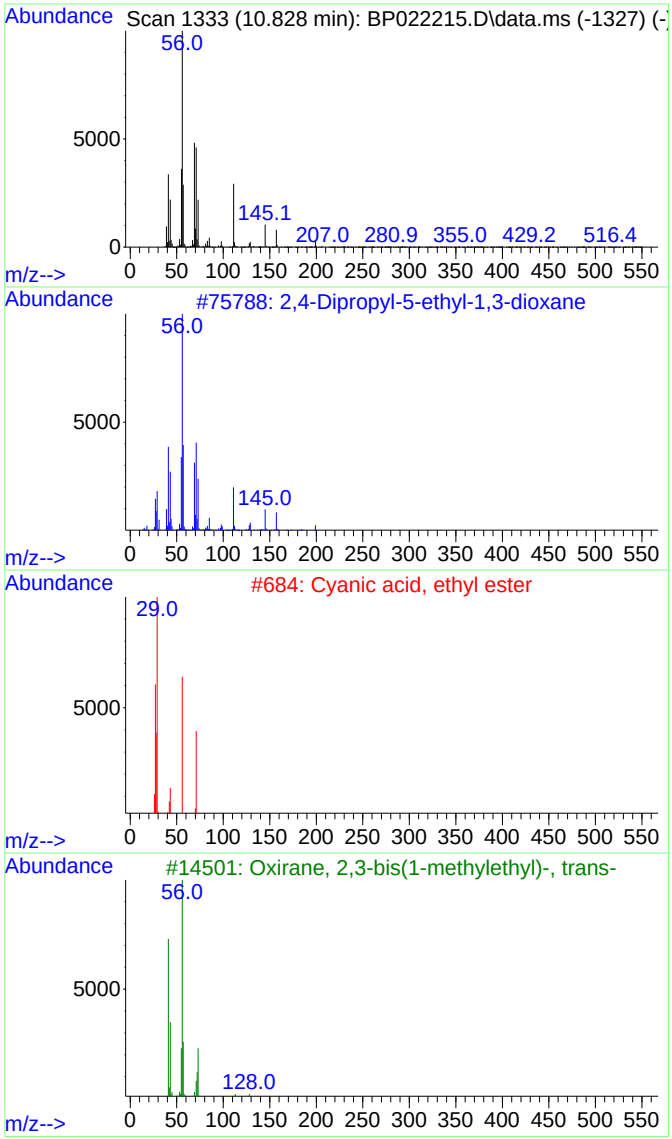
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

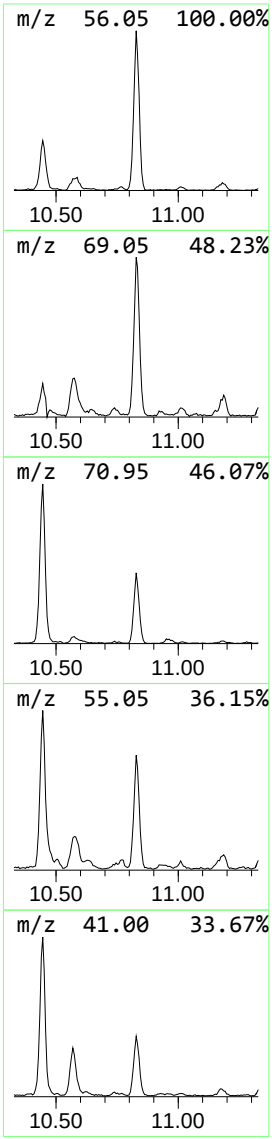
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 2,4-Dipropyl-5-ethyl-1,3-di... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.828	3.67 ng/ul	343184	Naphthalene-d8	10.457

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	72
2	Cyanic acid, ethyl ester	71	C3H5NO	000627-48-5	49
3	Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	43
4	Cyclobutane, 1,2,3,4-tetramethyl-	112	C8H16	069531-57-3	27
5	1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	27





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Data File : BP022215.D
Acq On : 04 Oct 2024 05:22
Operator : RC/JU
Sample : P4017-17DL 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCD3DL

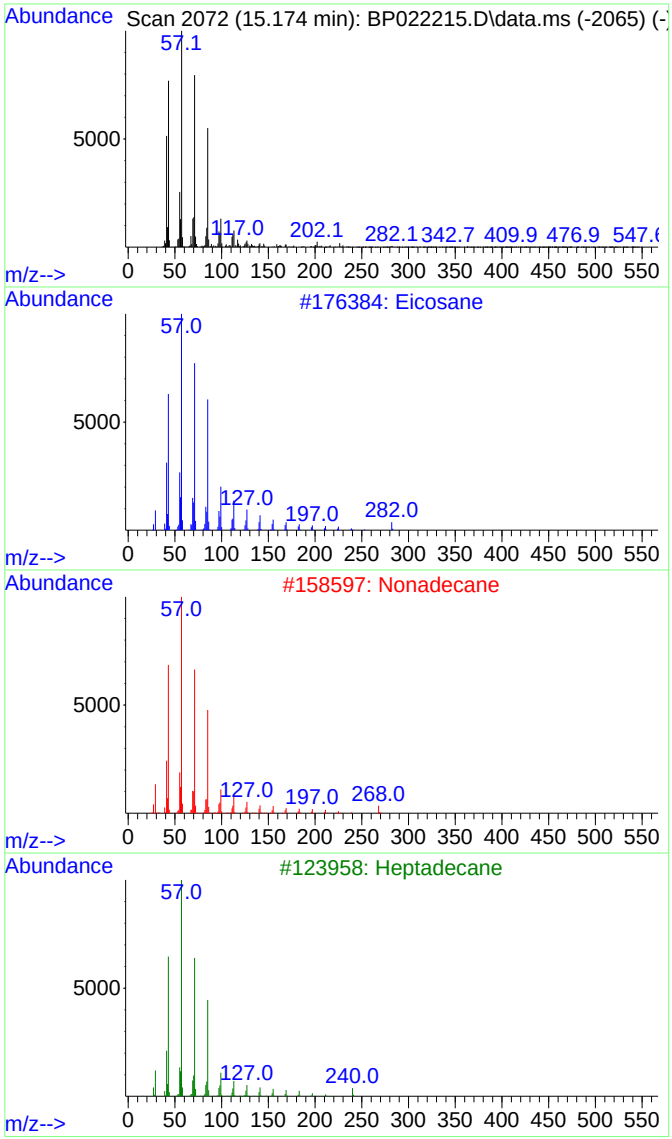
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

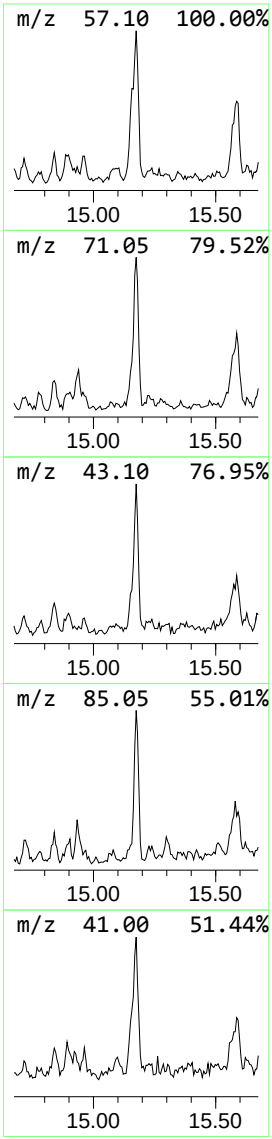
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.175	2.27 ng/ul	182349	Acenaphthene-d10	14.304

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	94
2	Nonadecane	268	C19H40	000629-92-5	90
3	Heptadecane	240	C17H36	000629-78-7	90
4	Eicosane	282	C20H42	000112-95-8	90
5	Tetradecane	198	C14H30	000629-59-4	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022215.D
Acq On : 04 Oct 2024 05:22
Operator : RC/JU
Sample : P4017-17DL 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCD3DL

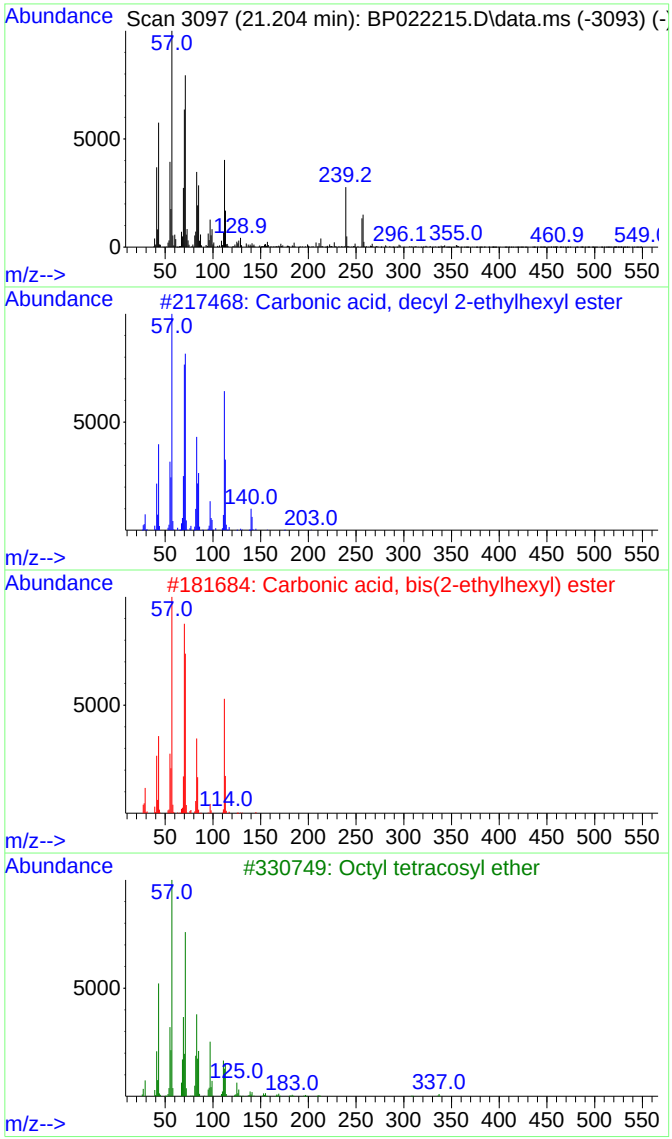
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

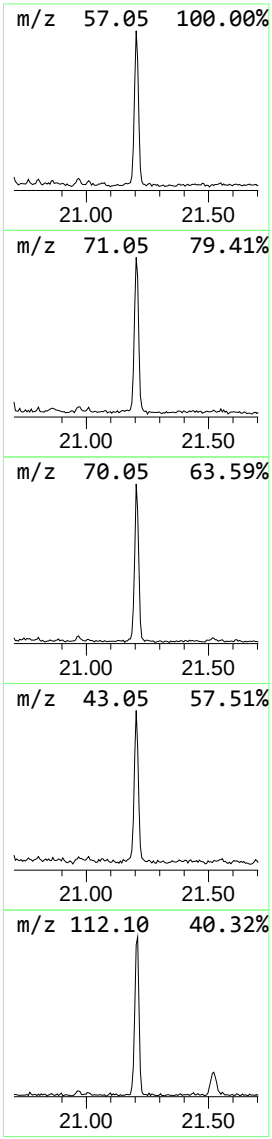
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Carbonic acid, decyl 2-ethy... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.204	3.18 ng/ul	401194	Chrysene-d12	21.521

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, decyl 2-ethylhexy...	314	C19H38O3	1000383-13-7	64
2	Carbonic acid, bis(2-ethylhexyl)...	286	C17H34O3	014858-73-2	58
3	Octyl tetracosyl ether	467	C32H66O	1000406-39-0	46
4	Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	43
5	2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	38





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022215.D
Acq On : 04 Oct 2024 05:22
Operator : RC/JU
Sample : P4017-17DL 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCD3DL

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

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: S...	7.328	4.5	ng/ul	172735	1	7.693	763910	20.0
1-Hexanol, 2-et...	7.757	2.5	ng/ul	96141	1	7.693	763910	20.0
(DEL) Alkane: S...	8.345	2.4	ng/ul	90957	1	7.693	763910	20.0
p-Cymene	8.787	2.0	ng/ul	77545	1	7.693	763910	20.0
(DEL) Alkane: S...	8.904	5.2	ng/ul	196603	1	7.693	763910	20.0
Heptyl isobutyl...	10.569	2.8	ng/ul	261179	2	10.457	1868910	20.0
2,4-Dipropyl-5-...	10.828	3.7	ng/ul	343184	2	10.457	1868910	20.0
(DEL) Alkane: S...	15.175	2.3	ng/ul	182349	3	14.304	1608530	20.0
Carbonic acid, ...	21.204	3.2	ng/ul	401194	5	21.521	2527110	20.0