LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A4CK4

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-BP071024.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP021156.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.181	29	33	42	rVB	33913	55880	0.86%	0.065%
2	3.470	78	82	89	rBV	69411	103242	1.58%	0.120%
3	3.699	116	121	130	rVB	76518	105223	1.61%	0.123%
4	3.905	151	156	166	rVB	76891	122184	1.87%	0.143%
5	4.799	303	308	318	rVB	57421	88747	1.36%	0.104%
6	5.681	451	458	465	rBV3	24762	47813	0.73%	0.056%
7	6.752	634	640	644	rBV6	19114	38255	0.59%	0.045%
8	6.822	648	652	655	rVV	45228	76403	1.17%	0.089%
9	6.869	655	660	676	rVV	425878	811231	12.45%	0.946%
10	6.999	676	682	693	rVV	372684	626846	9.62%	0.731%
11	7.199	710	716	722	rBV	527845	845266	12.97%	0.986%
12	7.246	722	724	727	rVV	44903	62976	0.97%	0.073%
13	7.281	727	730	739	rVB4	31925	73118	1.12%	0.085%
14	7.646	785	792	806	rVV	769578	1265196	19.41%	1.476%
15	8.381	910	917	929	rBV	545157	983824	15.10%	1.148%
16	8.805	982	989	1003	rBV	420144	847302	13.00%	0.988%
17	9.516	1103	1110	1124	rBV	330753	674758	10.35%	0.787%
18	9.793	1144	1157	1158	rBV5	26149	54376	0.83%	0.063%
19	10.075	1198	1205	1220	rBV	433352	969457	14.87%	1.131%
20	10.369	1250	1255	1258	rVV	26762	42939	0.66%	0.050%
21	10.422	1258	1264	1269	rVV	1000990	1577001	24.20%	1.839%
22	10.469	1269	1272	1277	rVB	89399	133746	2.05%	0.156%
23	10.593	1288	1293	1306	rVB2	49765	126822	1.95%	0.148%
24	11.057	1365	1372	1374	rBV4	18639	41909	0.64%	0.049%
25	11.169	1386	1391	1400	rVV2	123761	257282	3.95%	0.300%
26	11.816	1495	1501	1506	rBV2	31149	49083	0.75%	0.057%
27	12.081	1541	1546	1553	rBV	67032	107775	1.65%	0.126%
28	12.210	1562	1568	1574	rBV3	57176	128489	1.97%	0.150%
29	12.304	1579	1584	1592	rVB	52347	109019	1.67%	0.127%
30	13.081	1712	1716	1718	rBV	31377	40026	0.61%	0.047%
31	13.451	1772	1779	1786	rBV4	38807	96187	1.48%	0.112%
32	13.599	1798	1804	1809	rBV3	55302	106998	1.64%	0.125%
33	13.657	1809	1814	1815	rVV	45560	56859	0.87%	0.066%
34	13.693	1815	1820	1827	rVV	801726	1285338	19.72%	1.499%
35	13.751	1827	1830	1834	rVB2	41504	55538	0.85%	0.065%
36	13.840	1841	1845	1850	rBV2	24725	42493	0.65%	0.050%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

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-BP071024.MA.M
Title : SVOA CALIBRATION

37	13.981	1856	1869	1886	rBV2	1087940	2059433	31.60%	2.402%
38	14.163	1896	1900	1906	rVB2	43291	66754	1.02%	0.078%
39	14.287	1909	1921	1928	rBV	1232440	1963020	30.12%	2.290%
40	14.351	1928	1932	1941	rVB	139266	245371	3.76%	0.286%
41	14.510	1952	1959	1966	rBV	312408	662089	10.16%	0.772%
42	14.598	1968	1974	1986	rBV3	170418	497323	7.63%	0.580%
43	14.704	1987	1992	2001	rVV3	108612	258622	3.97%	0.302%
44	14.775	2001	2004	2009	rVB	36026	50566	0.78%	0.059%
45	14.916	2023	2028	2033	rBV5	37528	68647	1.05%	0.080%
46	14.975	2034	2038	2043	rVB3	27368	43650	0.67%	0.051%
47	15.093	2050	2058	2062	rBV4	37732	93537	1.44%	0.109%
48	15.140	2062	2066	2070	rVB4	41724	64571	0.99%	0.075%
49	15.304	2088	2094	2099	rBV	1434980	1976822	30.33%	2.306%
50	15.357	2099	2103	2108	rVB	123852	187558	2.88%	0.219%
51	15.481	2117	2124	2129	rBV2	304376	487480	7.48%	0.569%
52	15.569	2137	2139	2144	rVB4	23624	40088	0.62%	0.047%
53	15.657	2151	2154	2160	rVB	64268	104872	1.61%	0.122%
54	15.816	2177	2181	2188	rVB	56461	113257	1.74%	0.132%
55	15.887	2189	2193	2201	rVB3	28530	69041	1.06%	0.081%
56	16.045	2212	2220	2231	rVB8	85232	294544	4.52%	0.344%
57	16.393	2272	2279	2284	rBV5	61166	141232	2.17%	0.165%
58	16.440	2284	2287	2292	rVB3	58322	76151	1.17%	0.089%
59	16.493	2292	2296	2300	rBV5	66015	135724	2.08%	0.158%
60	16.663	2322	2325	2332	rVB5	47795	85358	1.31%	0.100%
61	16.763	2337	2342	2349	rVV2	138906	246410	3.78%	0.287%
62	16.834	2349	2354	2360	rVV3	113674	235409	3.61%	0.275%
63	16.916	2364	2368	2375	rVB	144795	235782	3.62%	0.275%
64	17.010	2380	2384	2391	rBV3	106282	183800	2.82%	0.214%
65	17.104	2391	2400	2404	rBV	1555222	2493850	38.26%	2.909%
66	17.145	2404	2407	2412	rVV	2454646	3365807	51.64%	3.926%
67	17.204	2412	2417	2421	rVV2	1255298	2075326	31.84%	2.421%
68	17.245	2421	2424	2429	rVB	492019	668280	10.25%	0.779%
69	17.540	2466	2474	2480	rBV2	247339	578432	8.88%	0.675%
70	17.728	2499	2506	2511	rVB3	203281	445769	6.84%	0.520%
71	17.845	2522	2526	2527	rBV3	95336	124964	1.92%	0.146%
72	18.004	2549	2553	2556	rBV	391362	535304	8.21%	0.624%
73	18.045	2556	2560	2567	rVV2	1240814	2193309	33.65%	2.558%
74	18.110	2568	2571	2574	rVV	148370	157566	2.42%	0.184%
75	18.145	2575	2577	2583	rVB	144428	208815	3.20%	0.244%
76	18.210	2583	2588	2593	rBV2	1042547	1848028	28.36%	2.156%

LSC Area Percent Report

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Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

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-BP071024.MA.M
Title : SVOA CALIBRATION

77	18.263	2594	2597	2602	rVV2	577498	789535	12.11%	0.921%
78	18.316	2602	2606	2615	rVV6	261557	522911	8.02%	0.610%
79	18.504	2634	2638	2641	rBV2	338330	526598	8.08%	0.614%
80	18.534	2641	2643	2650	rVB	328614	525635	8.07%	0.613%
81	18.598	2650	2654	2660	rBV	388769	631779	9.69%	0.737%
82	18.969	2713	2717	2722	rBV2	365836	792545	12.16%	0.924%
83	19.022	2723	2726	2730	rVV4	416630	716656	11.00%	0.836%
84	19.069	2730	2734	2737	rVV2	866752	1247937	19.15%	1.456%
85	19.104	2737	2740	2748	rVB4	789865	1346890	20.67%	1.571%
86	19.251	2758	2765	2770	rBV	3939560	6517467	100.00%	7.602%
87	19.386	2784	2788	2796	rVB3	772177	1368198	20.99%	1.596%
88	19.598	2818	2824	2826	rBV2	2094397	3589699	55.08%	4.187%
89	19.628	2826	2829	2838	rVB	2951222	4422242	67.85%	5.158%
90	19.851	2865	2867	2871	rVB	368327	360969	5.54%	0.421%
91	20.681	3005	3008	3012	rVB	667968	751295	11.53%	0.876%
92	21.216	3095	3099	3106	rVB2	2034270	3362115	51.59%	3.922%
93	21.551	3150	3156	3162	rBV3	2223192	6151043	94.38%	7.175%
94	21.610	3163	3166	3174	rVB	2231383	3649739	56.00%	4.257%
95	21.763	3190	3192	3196	rVB	827065	862014	13.23%	1.005%
96	22.398	3296	3300	3303	rBV2	1023517	1725223	26.47%	2.012%
97	23.833	3541	3544	3552	rVB	1075343	2473523	37.95%	2.885%
98	23.951	3560	3564	3572	rVB	1724819	4079662	62.60%	4.759%
99	24.045	3578	3580	3588	rVB	968774	1526131	23.42%	1.780%
100	30.298	4641	4643	4646	rBV3	261531	300426	4.61%	0.350%

Sum of corrected areas: 85732394

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
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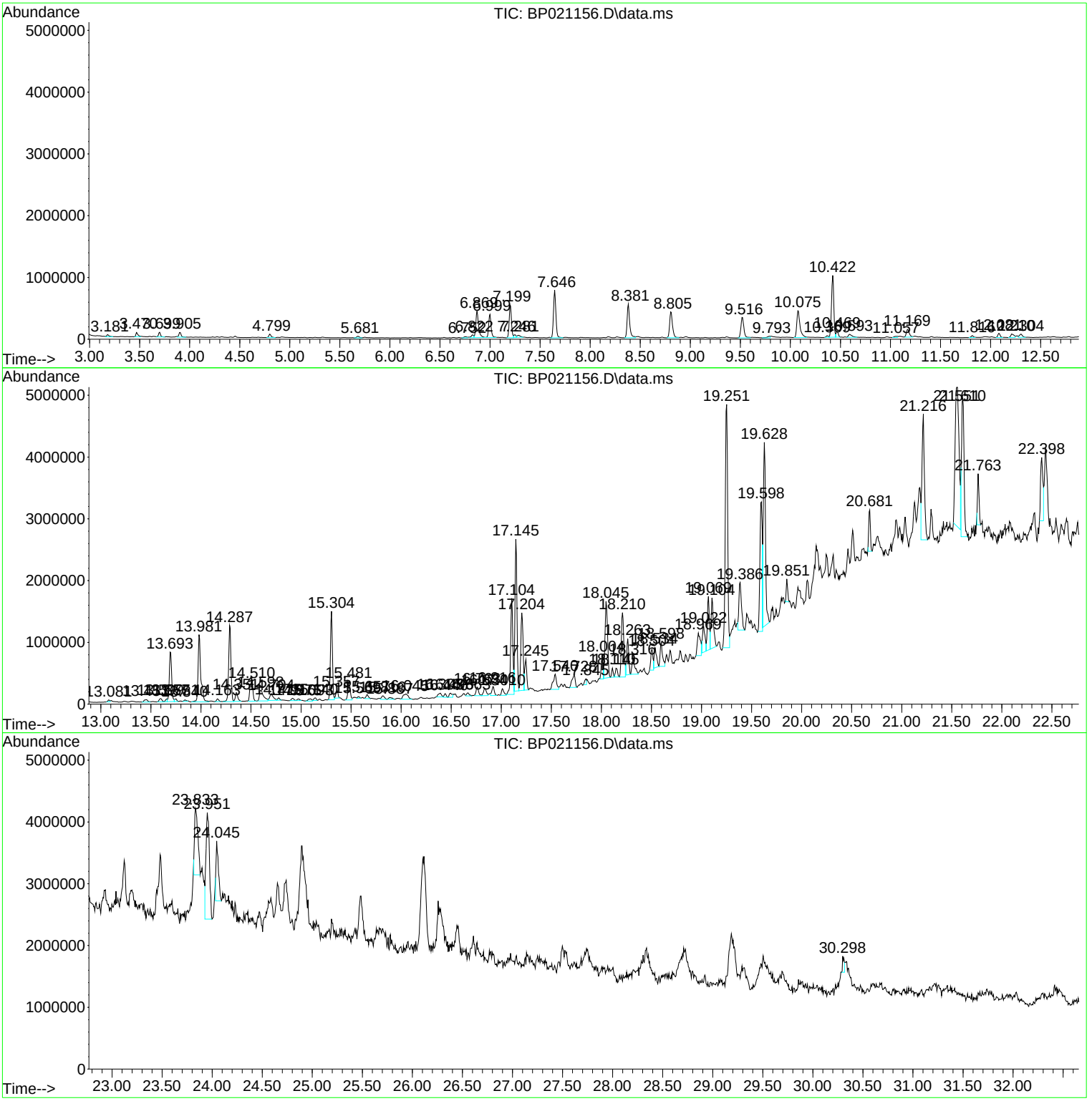
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.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\BP072524\
Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

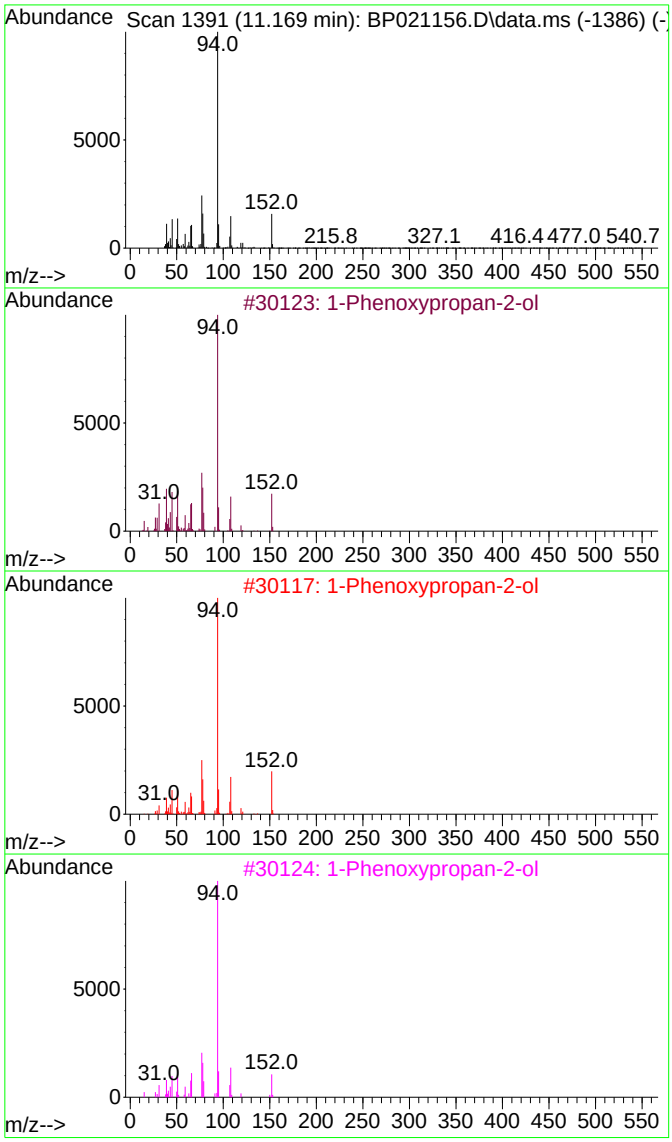
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

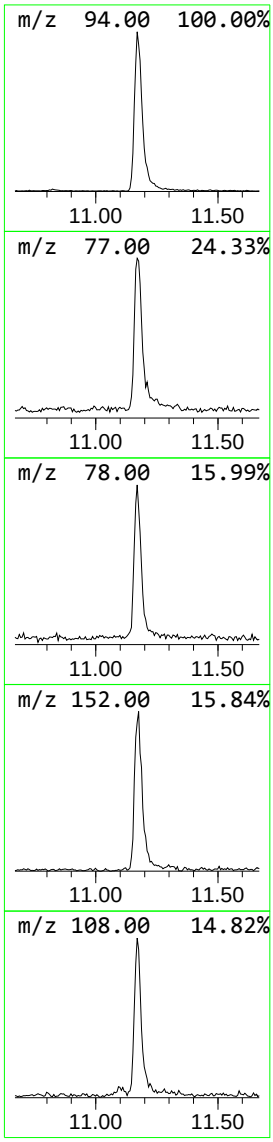
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 1-Phenoxypropan-2-ol Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.169	3.26 ng/ul	257282	Naphthalene-d8	10.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
3	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
4	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
5	1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

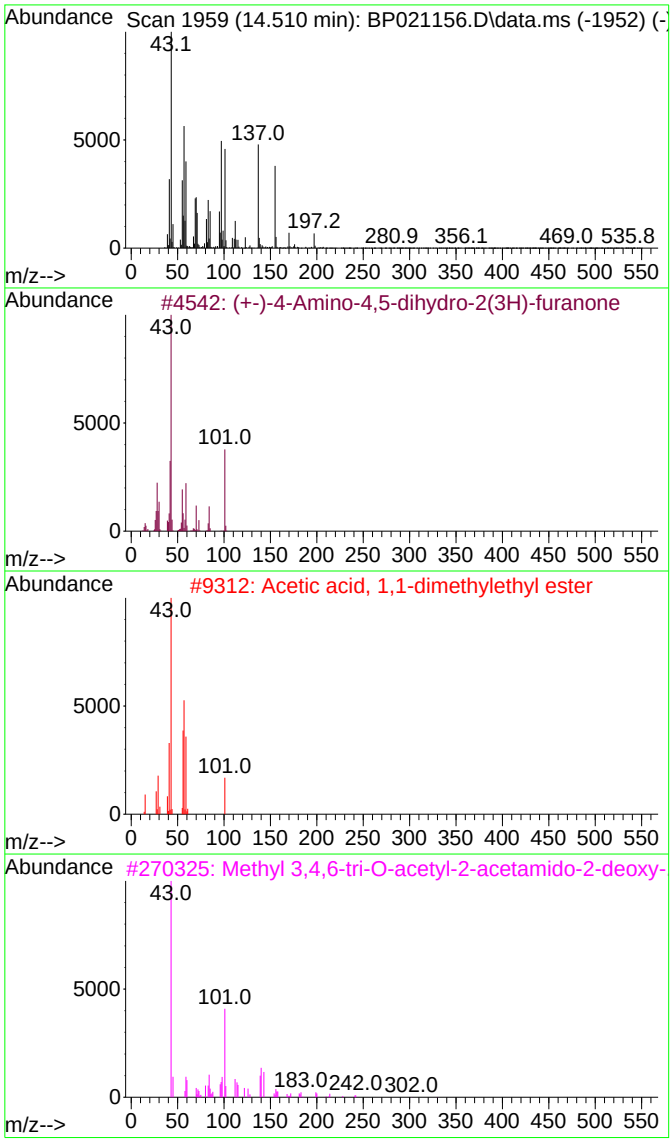
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

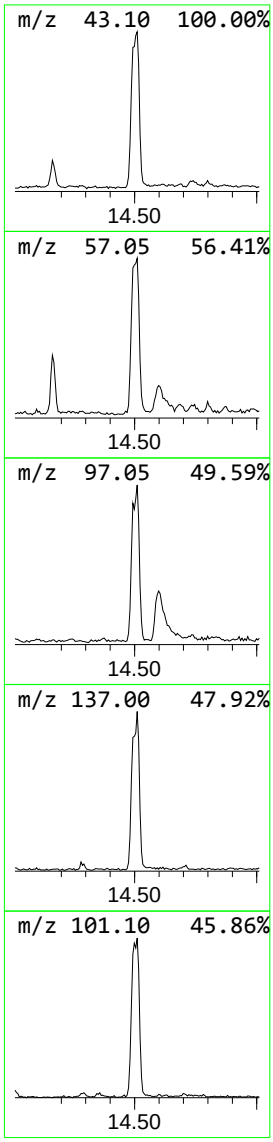
****

Peak Number 2 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.510	6.75 ng/ul	662089	Acenaphthene-d10	14.287

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+-)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	25
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	12
3	Methyl 3,4,6-tri-O-acetyl-2-acet...	361	C15H23NO9	002595-39-3	10
4	trans-5-Isopropyl-6,7-epoxy-8-hy...	228	C13H24O3	058002-07-6	10
5	2-t-Butyl-5-propyl-[1,3]dioxolan...	186	C10H18O3	157733-17-0	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

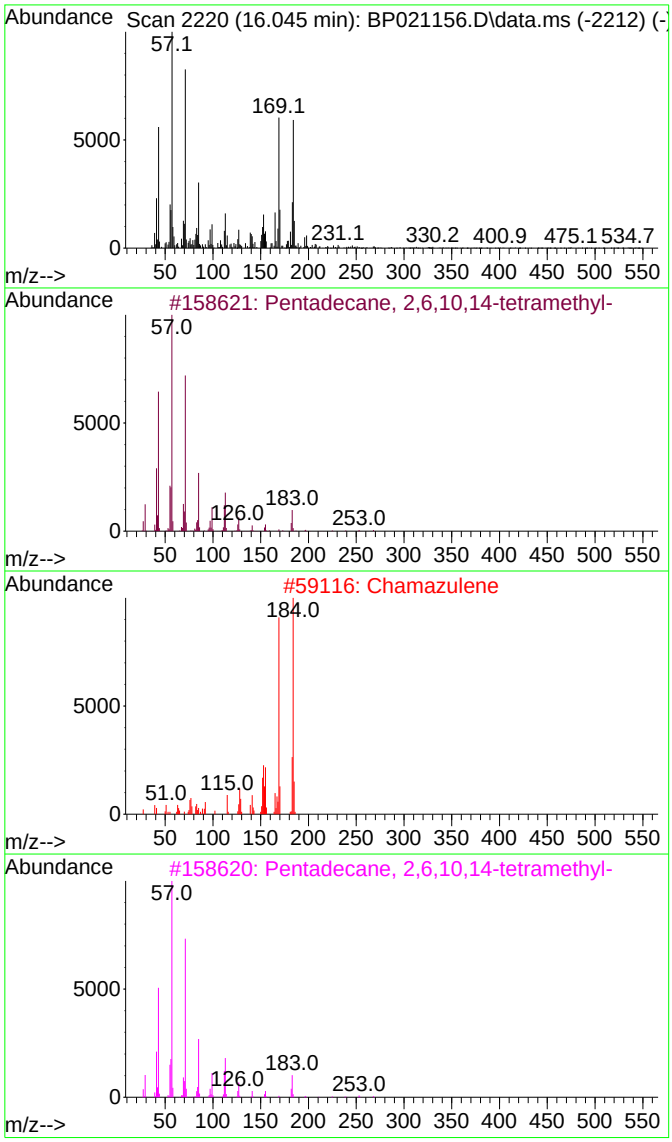
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

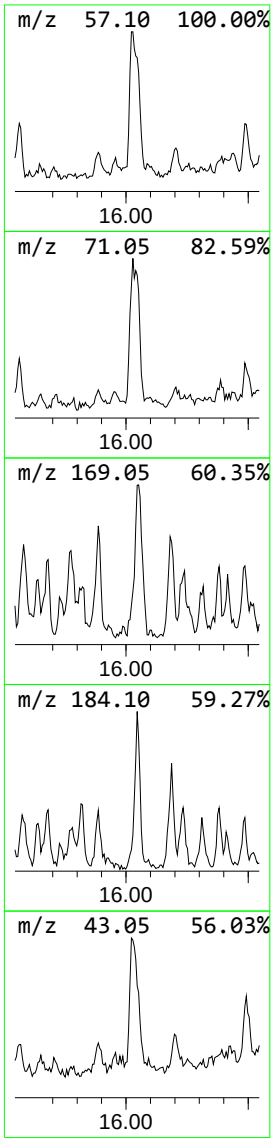
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.046	2.36 ng/ul	294544	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	83
2	Chamazulene	184	C14H16	000529-05-5	64
3	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	60
4	Hexadecane, 2,6,10-trimethyl-	268	C19H40	055000-52-7	45
5	Chamazulene	184	C14H16	000529-05-5	45





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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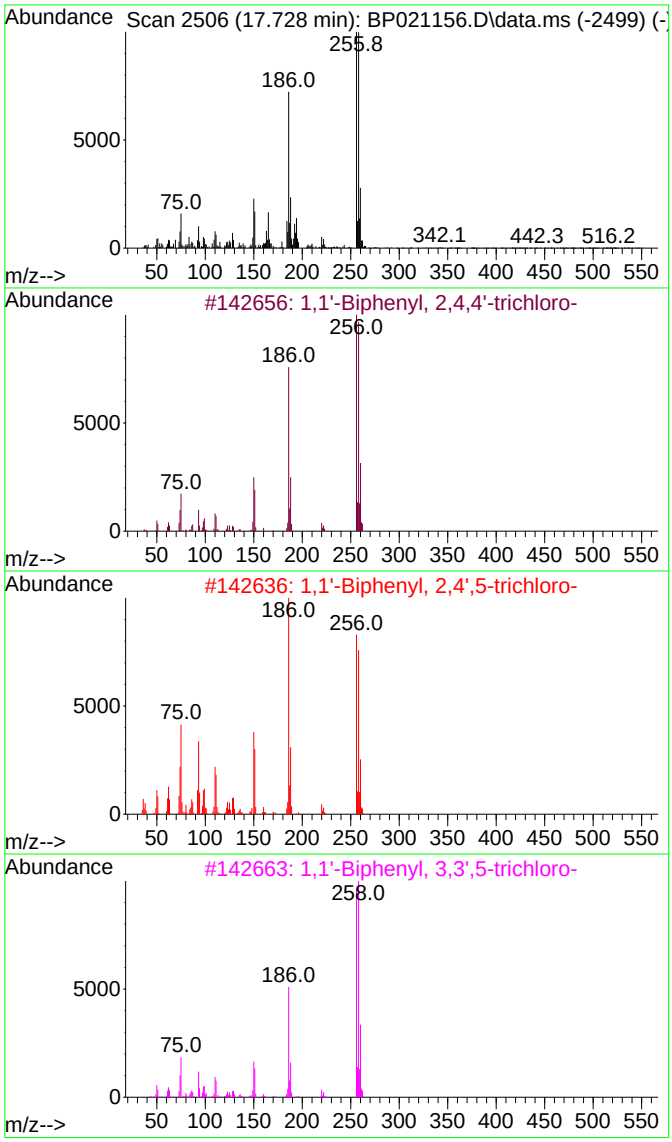
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

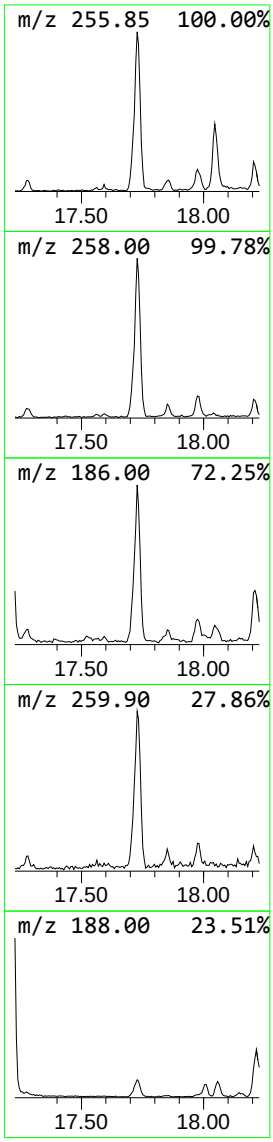
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1,1'-Biphenyl, 2,4,4'-trich... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.728	3.57 ng/ul	445769	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99
2	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	99
3	1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99
4	1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99
5	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

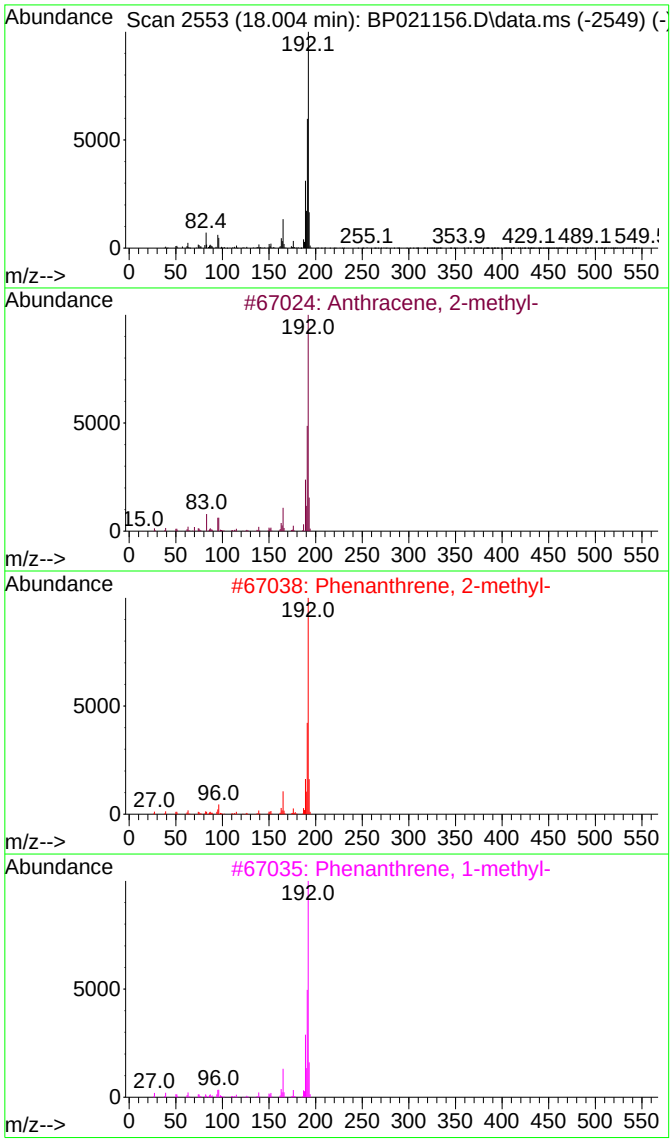
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

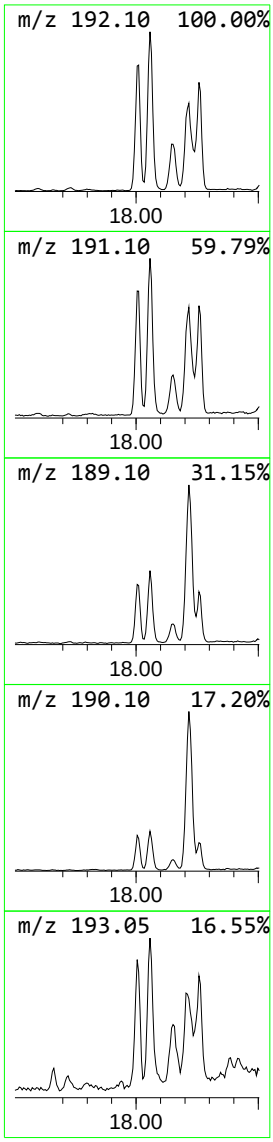
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Anthracene, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.004	4.29 ng/ul	535304	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	94





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Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

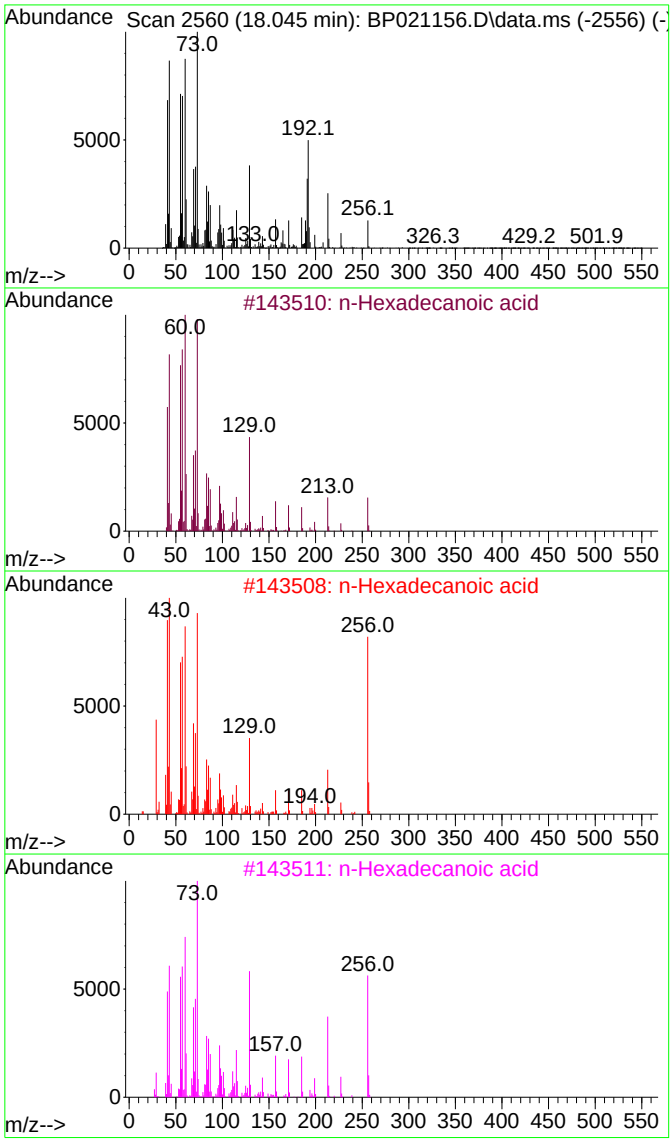
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

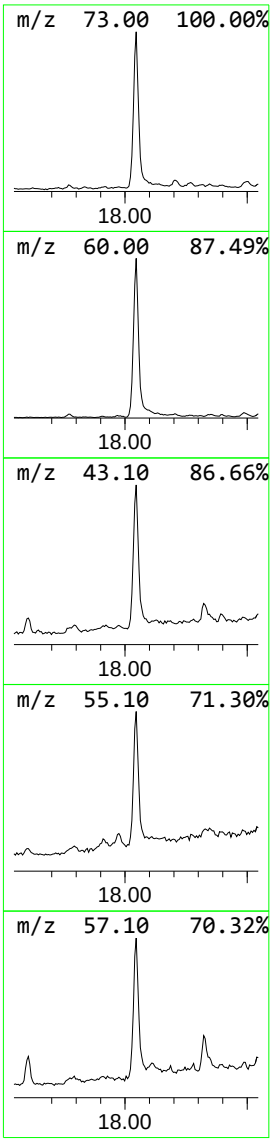
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.045	17.59 ng/ul	2193310	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5	Tridecanoic acid	214	C13H26O2	000638-53-9	96





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Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

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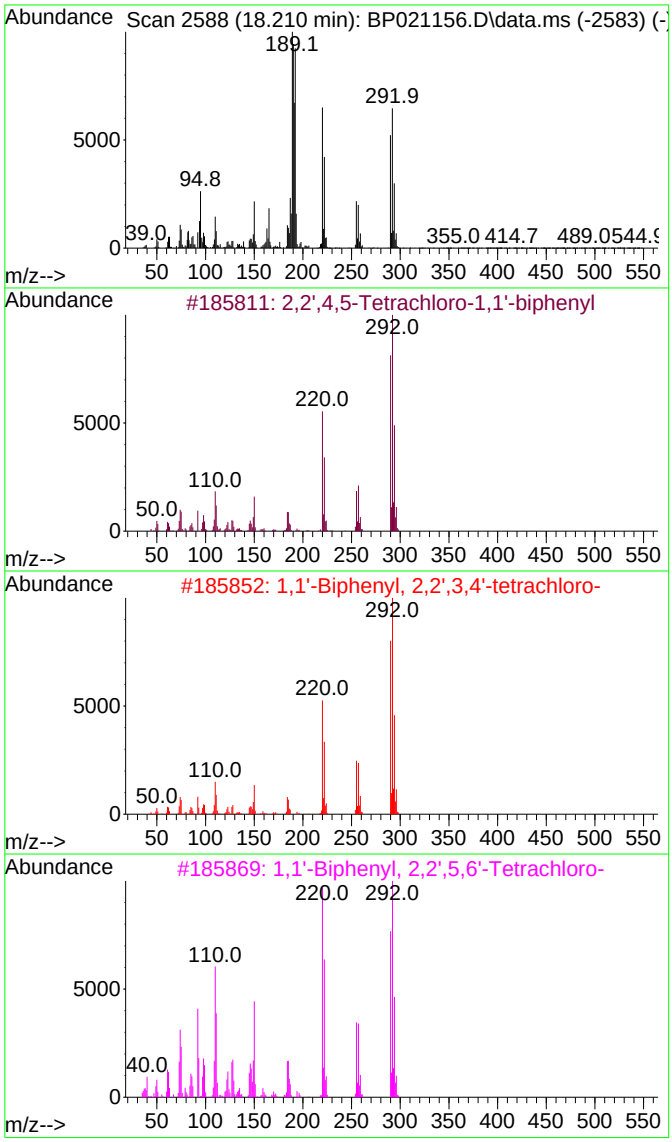
Quant Title : SVOA CALIBRATION

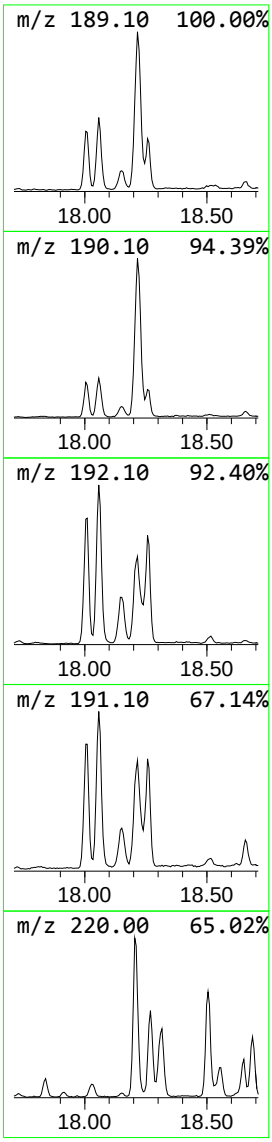
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 2,2',4,5-Tetrachloro-1,1'-b... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.210	14.82 ng/ul	1848030	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	95
2	1,1'-Biphenyl, 2,2',3,4'-tetrach...	290	C12H6Cl4	036559-22-5	95
3	1,1'-Biphenyl, 2,2',5,6'-Tetrach...	290	C12H6Cl4	041464-41-9	95
4	1,1'-Biphenyl, 2,2',3,5'-tetrach...	290	C12H6Cl4	041464-39-5	95
5	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	95





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Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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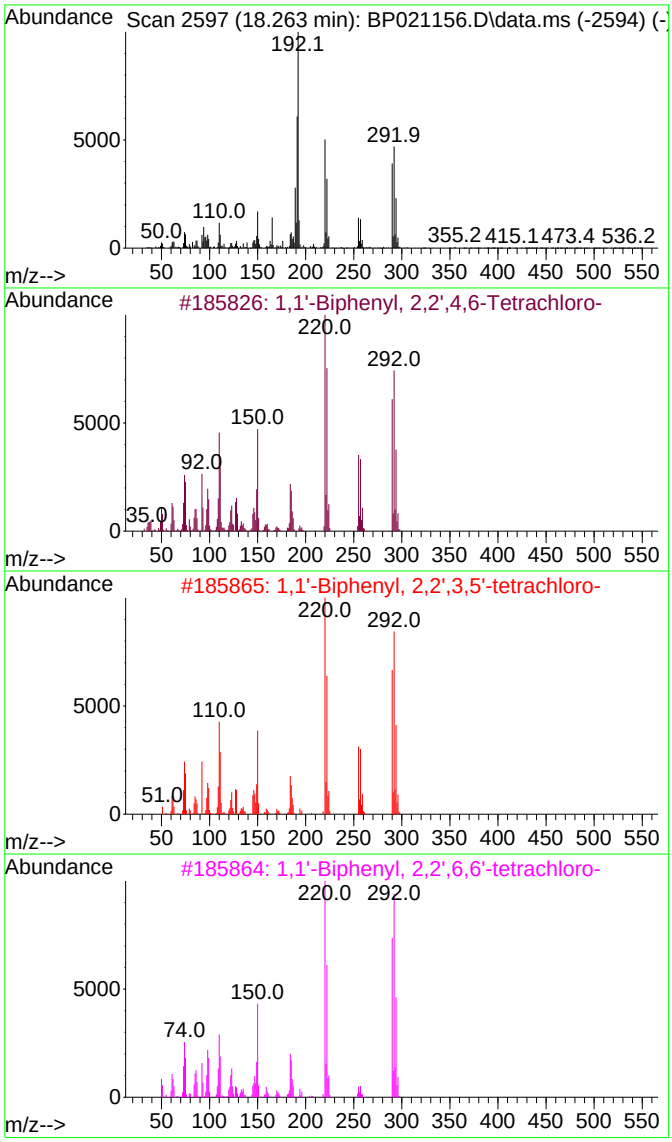
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

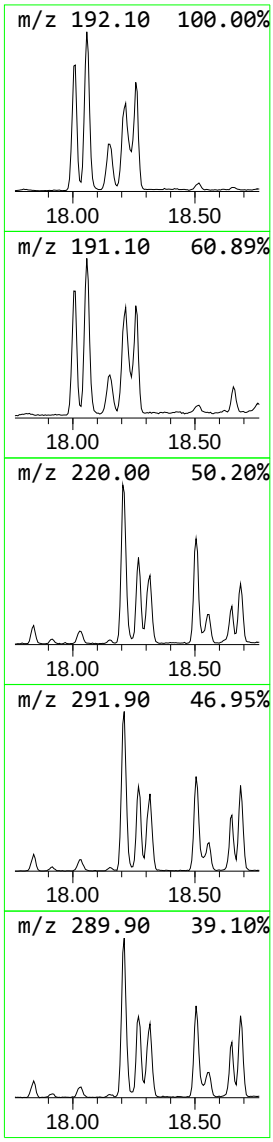
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1,1'-Biphenyl, 2,2',4,6-Tet... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.263	6.33 ng/ul	789535	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,6-Tetrachl...	290	C12H6Cl4	062796-65-0	97
2	1,1'-Biphenyl, 2,2',3,5'-tetrach...	290	C12H6Cl4	041464-39-5	97
3	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	96
4	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	94
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	78





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Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

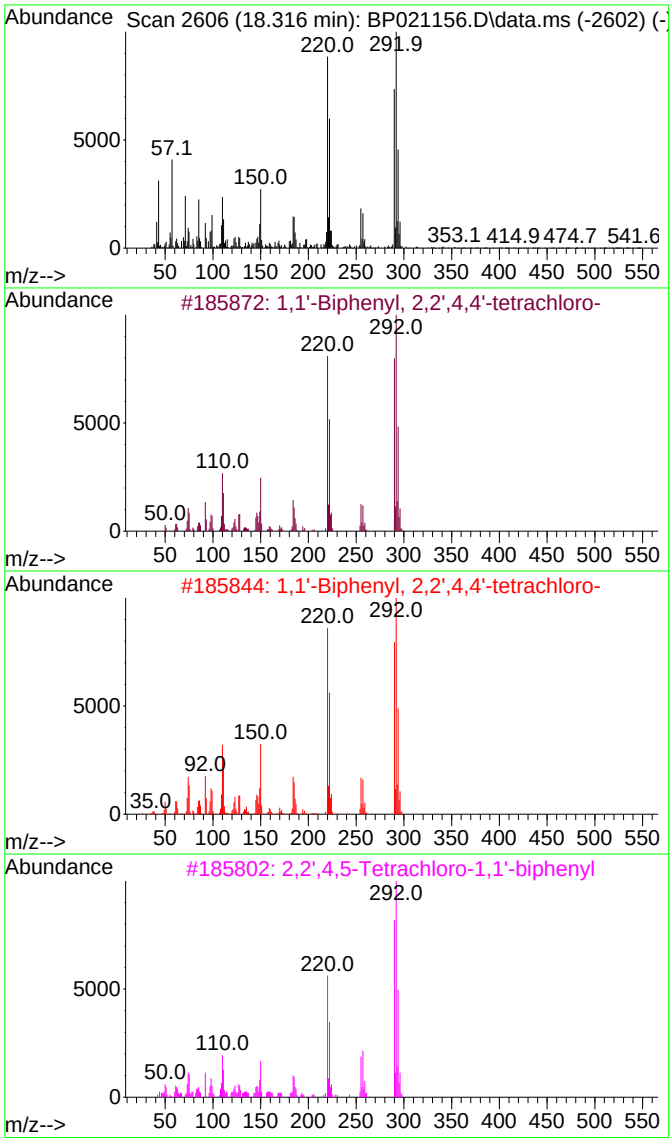
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

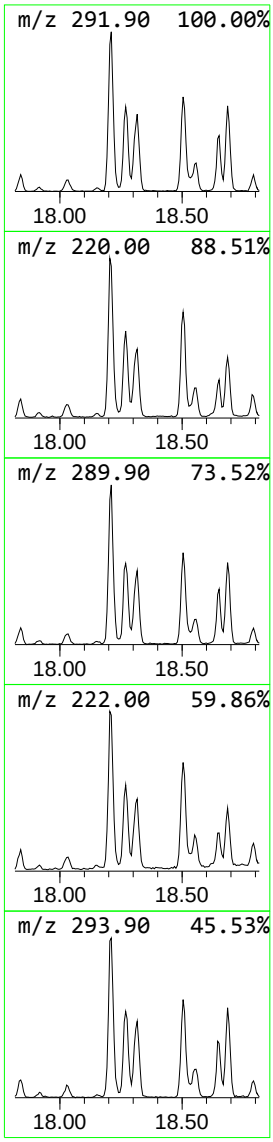
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.316	4.19 ng/ul	522911	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
2	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
3	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
4	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
5	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99





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Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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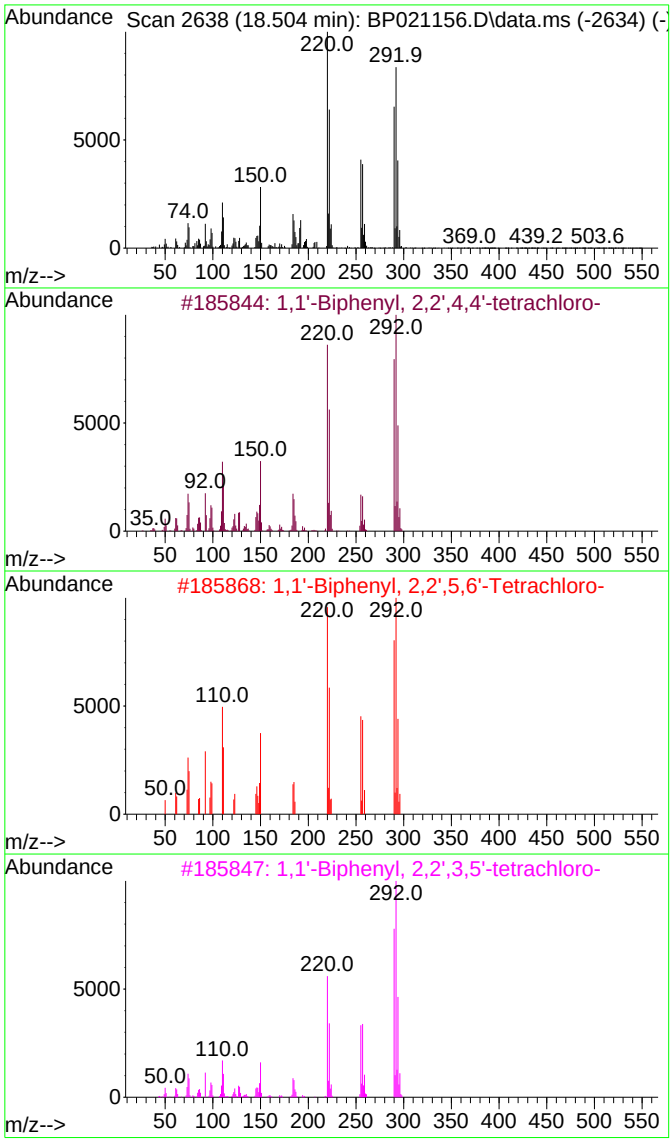
Quant Title : SVOA CALIBRATION

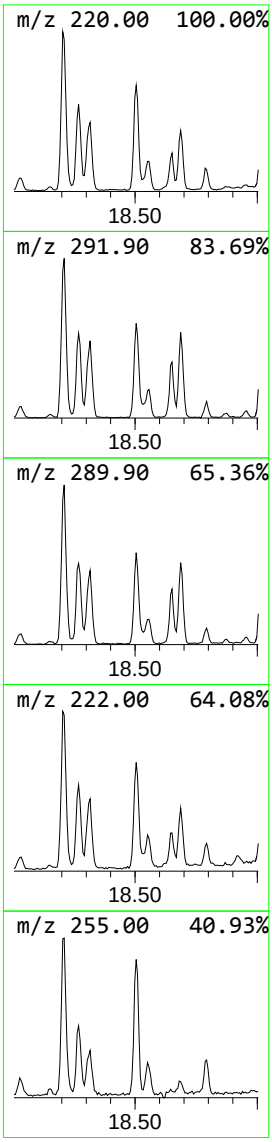
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1,1'-Biphenyl, 2,2',5,6'-Te... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.504	4.22 ng/ul	526598	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
2	1,1'-Biphenyl, 2,2',5,6'-Tetrach...	290	C12H6Cl4	041464-41-9	99
3	1,1'-Biphenyl, 2,2',3,5'-tetrach...	290	C12H6Cl4	041464-39-5	99
4	2,2',3,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-45-7	99
5	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99





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Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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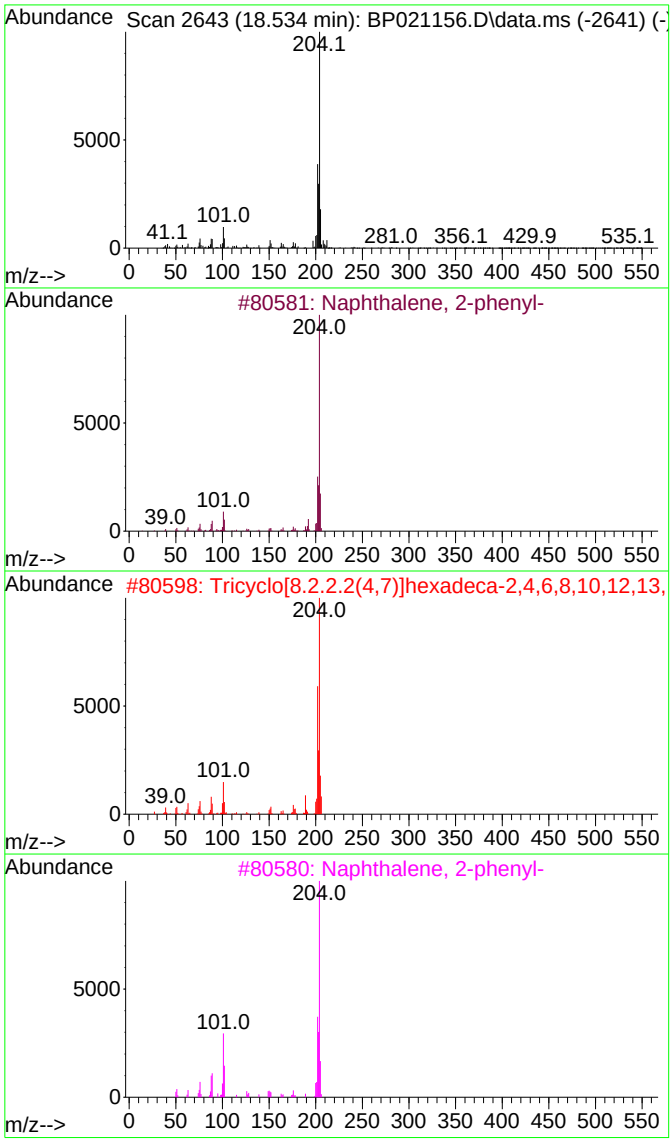
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

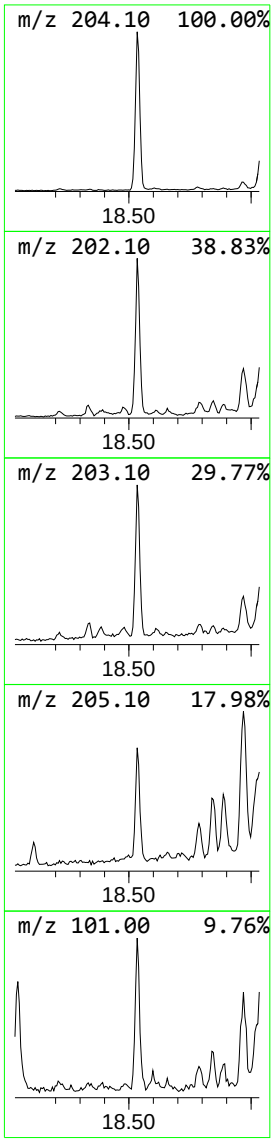
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Naphthalene, 2-phenyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.534	4.22 ng/ul	525635	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	91
2	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	70
3	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	70
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	70
5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	70





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Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

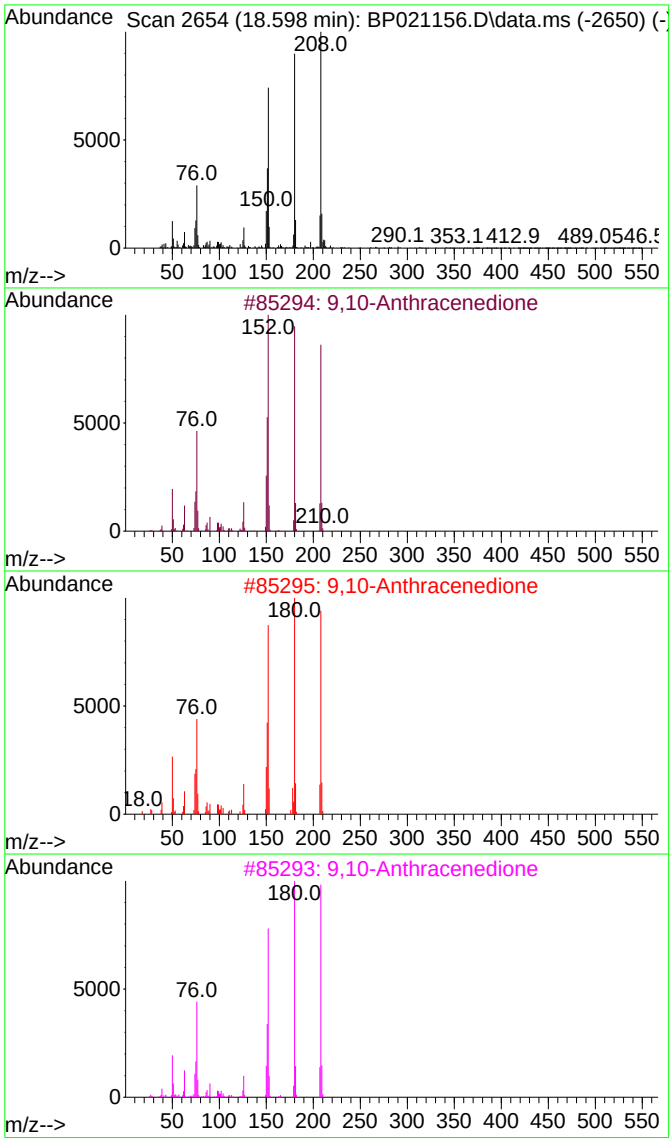
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

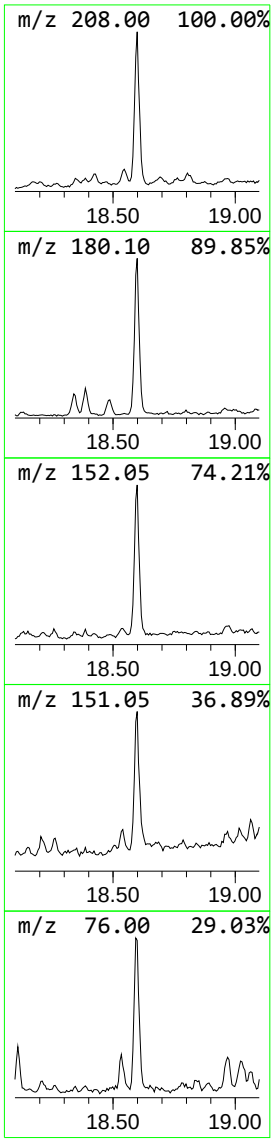
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 9,10-Anthracenedione Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.598	5.07 ng/ul	631779	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
2	9,10-Anthracenedione	208	C14H8O2	000084-65-1	97
3	9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
4	9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
5	1,4-Anthracenedione	208	C14H8O2	000635-12-1	70





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Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

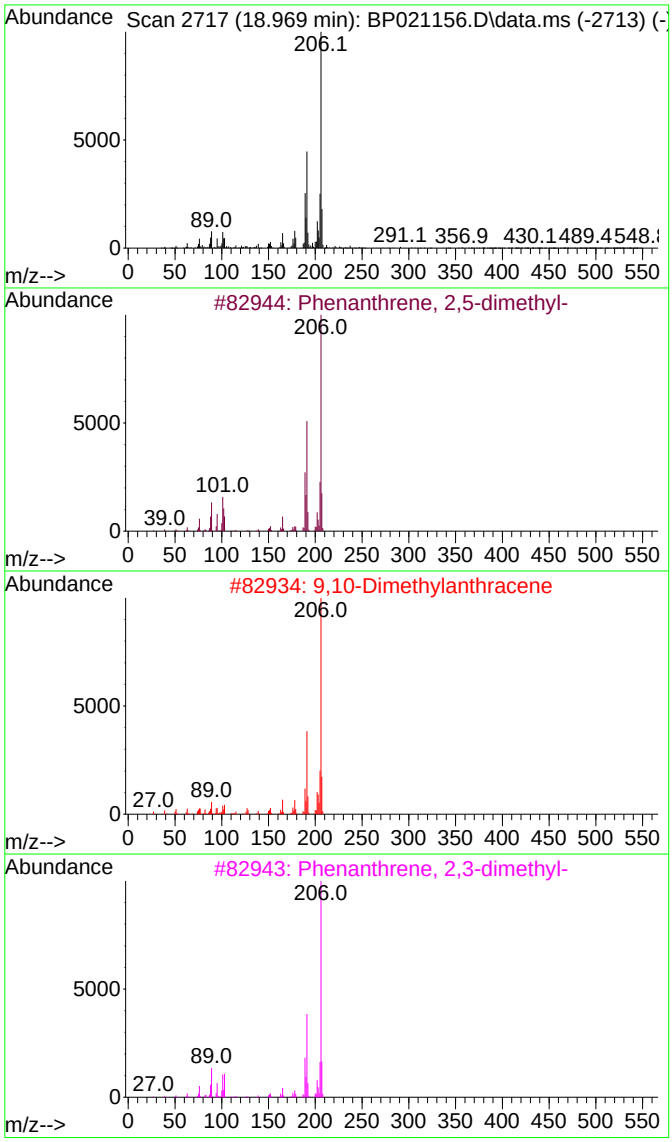
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

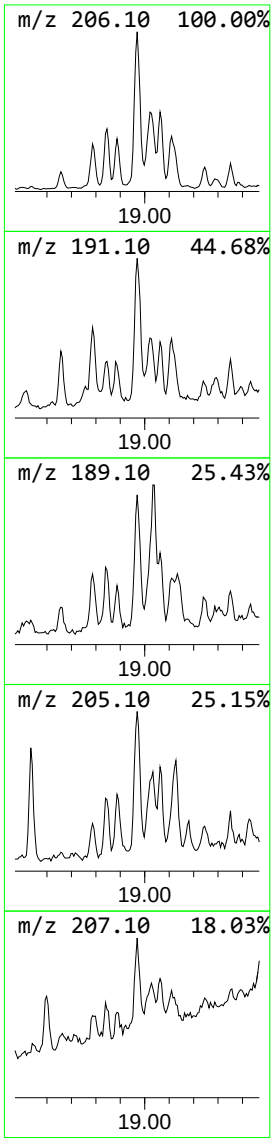
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Phenanthrene, 2,5-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.969	6.36 ng/ul	792545	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	97
2	9,10-Dimethylanthracene	206	C16H14	000781-43-1	96
3	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	91
4	di-p-Tolylacetylene	206	C16H14	002789-88-0	91
5	9,10-Dimethylanthracene	206	C16H14	000781-43-1	91





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Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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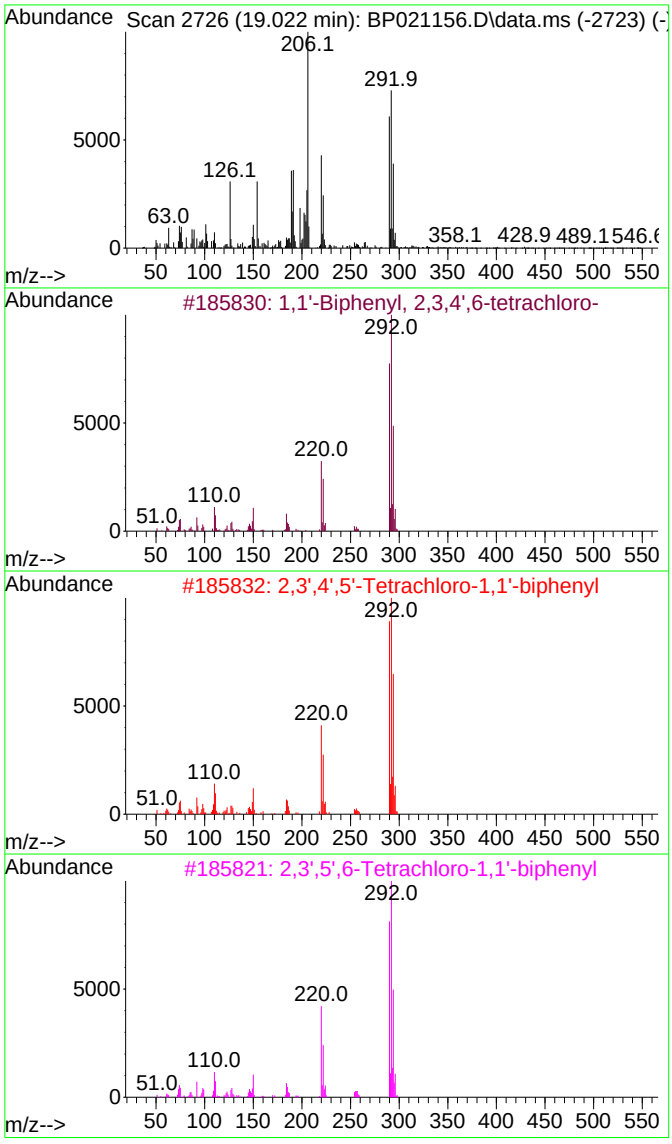
Quant Title : SVOA CALIBRATION

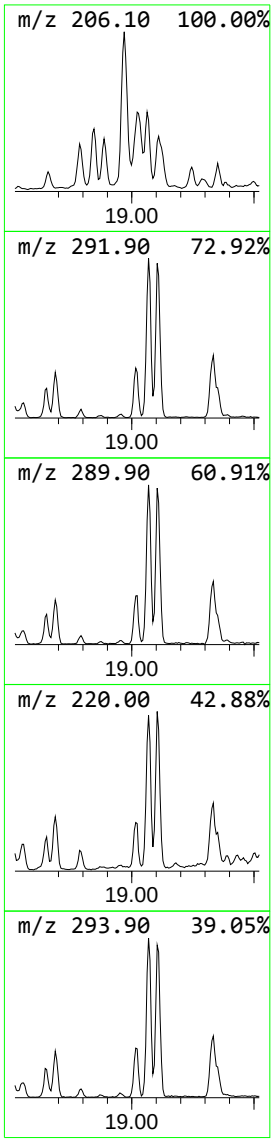
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 1,1'-Biphenyl, 2,3,4',6-tet... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.022	5.75 ng/ul	716656	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,4',6-tetrachl...	290	C12H6Cl4	052663-58-8	93
2	2,3',4',5'-Tetrachloro-1,1'-biph...	290	C12H6Cl4	070362-48-0	93
3	2,3',5',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074338-23-1	93
4	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	92
5	3,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-49-1	92





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Sample : P3263-16
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ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

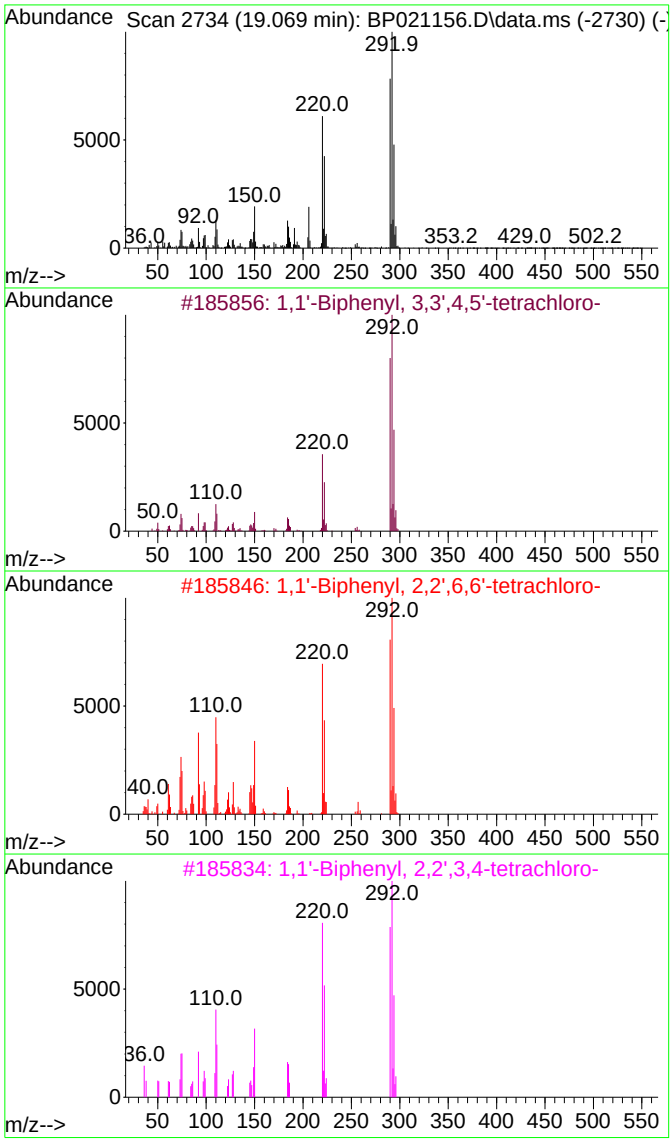
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

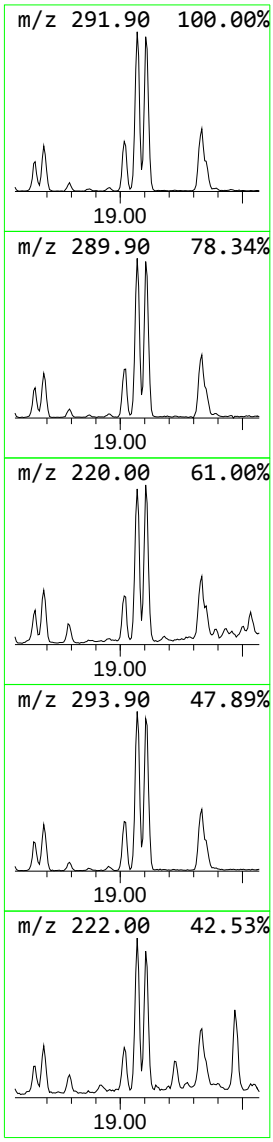
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 1,1'-Biphenyl, 3,3',4,5'-te... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.069	10.01 ng/ul	1247940	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 3,3',4,5'-tetrach...	290	C12H6Cl4	041464-48-6	99
2	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
3	1,1'-Biphenyl, 2,2',3,4-tetrachl...	290	C12H6Cl4	052663-59-9	99
4	1,1'-Biphenyl, 2,3',4',5-tetrach...	290	C12H6Cl4	032598-11-1	98
5	1,1'-Biphenyl, 2,3',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	98





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Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

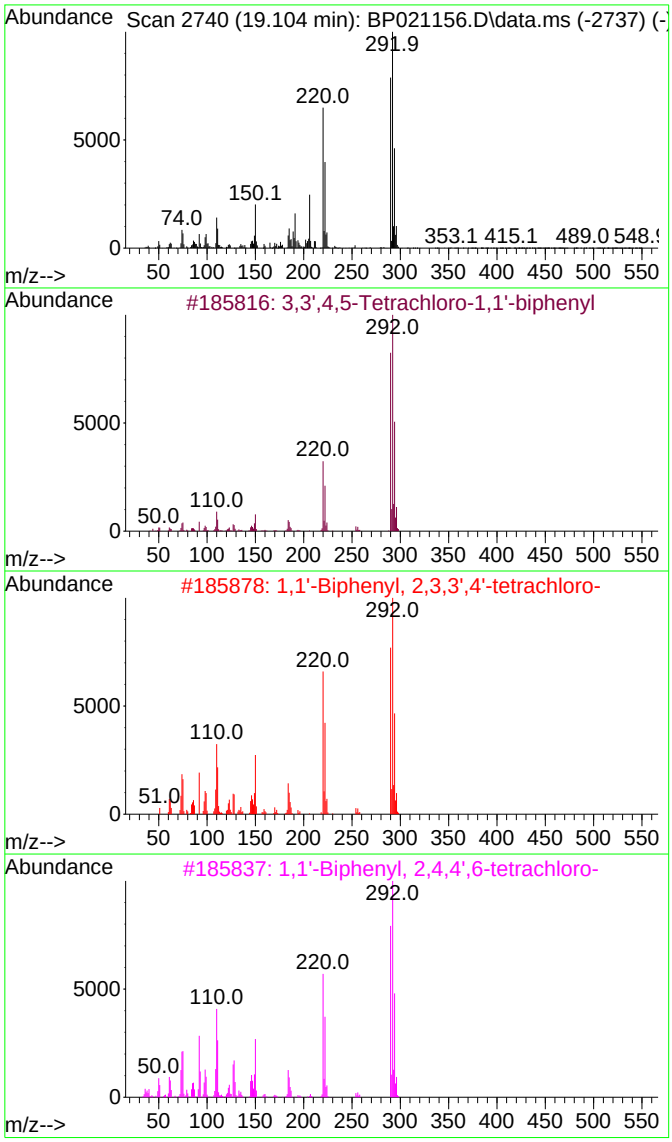
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

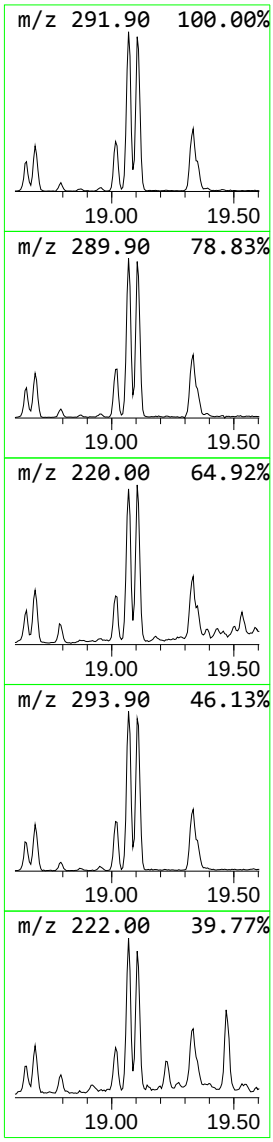
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 3,3',4,5-Tetrachloro-1,1'-b... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.104	10.80 ng/ul	1346890	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-49-1	96
2	1,1'-Biphenyl, 2,3,3',4'-tetrach...	290	C12H6Cl4	041464-43-1	95
3	1,1'-Biphenyl, 2,4,4',6-tetrachl...	290	C12H6Cl4	032598-12-2	95
4	1,1'-Biphenyl, 3,3',4,5'-tetrach...	290	C12H6Cl4	041464-48-6	94
5	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	94





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Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

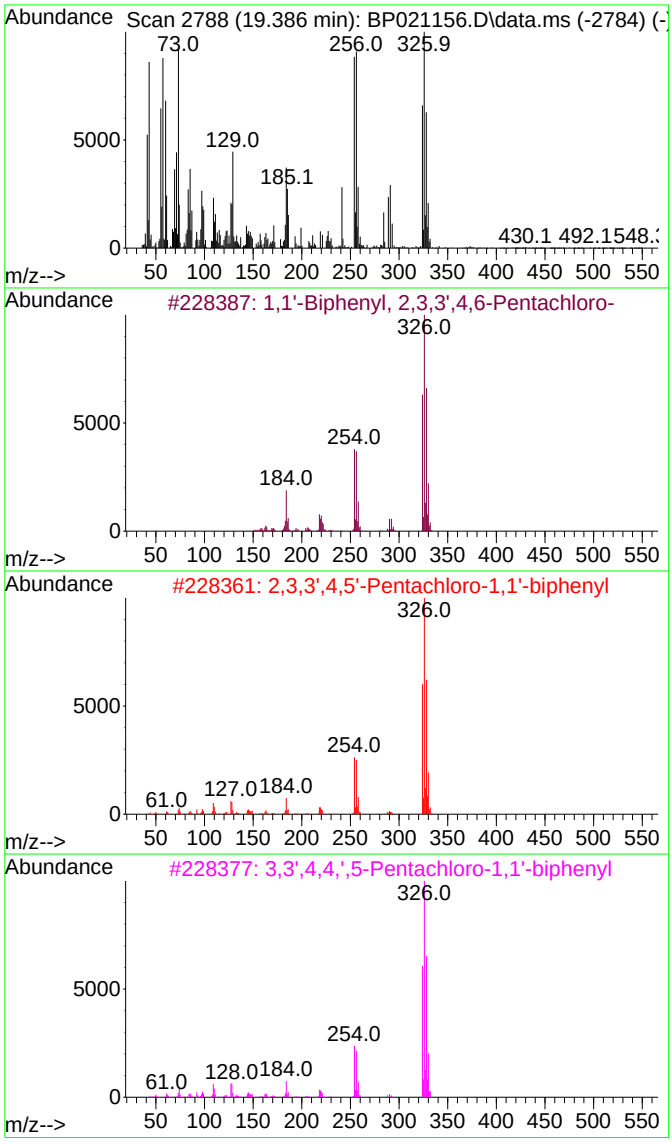
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

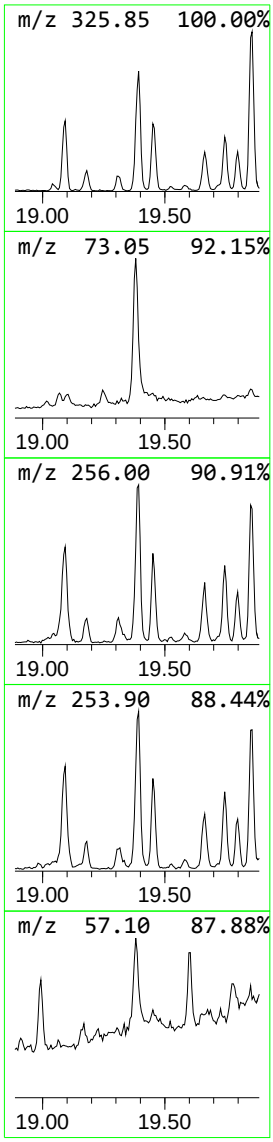
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 1,1'-Biphenyl, 2,3,3',4,6-P... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.387	4.45 ng/ul	1368200	Chrysene-d12	21.557

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3',4,6-Pentac...	324	C12H5Cl5	074472-35-8	99
2	2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	99
3	3,3',4,4',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	99
4	1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	99
5	1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	99





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Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

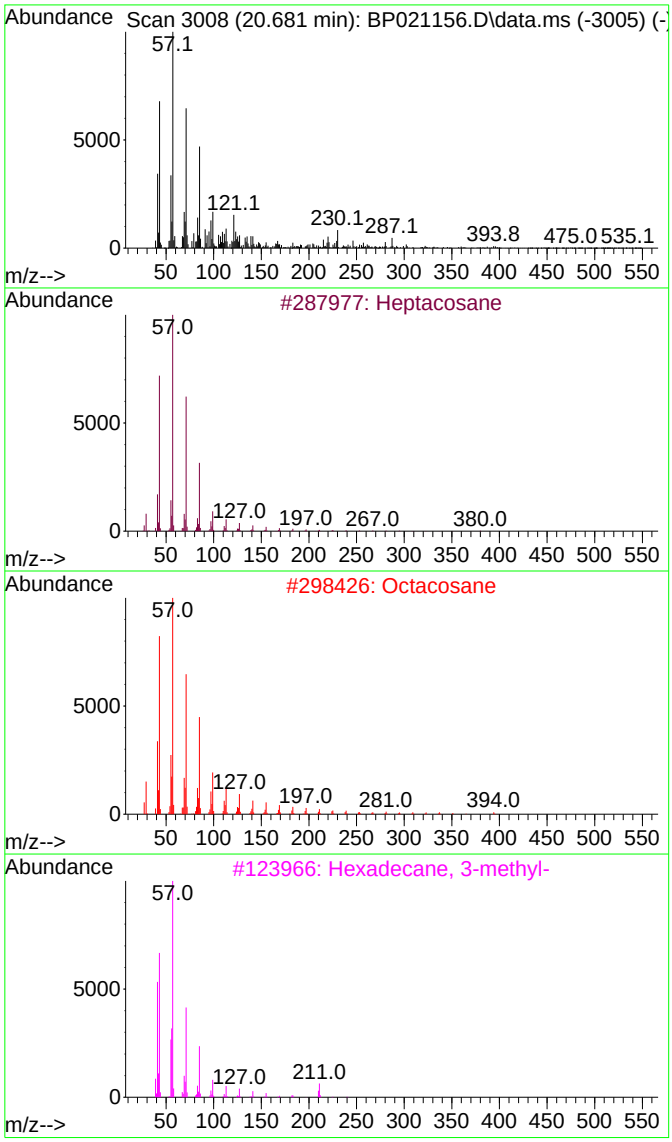
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

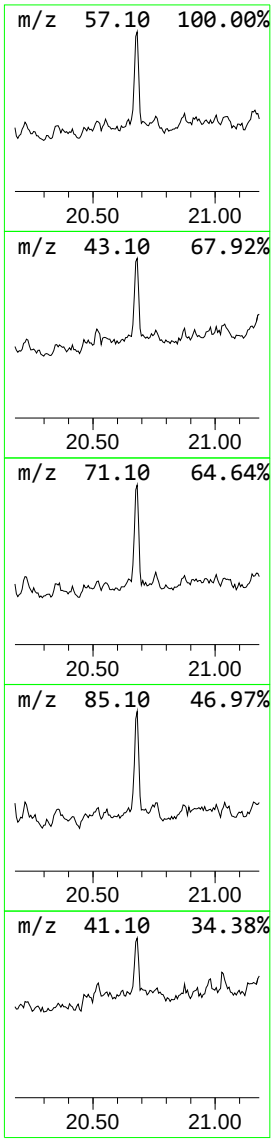
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.680	2.44 ng/ul	751295	Chrysene-d12	21.557

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	86
2	Octacosane	394	C28H58	000630-02-4	83
3	Hexadecane, 3-methyl-	240	C17H36	006418-43-5	76
4	3-Methylheptacosane	394	C28H58	014167-66-9	68
5	Nonacosane, 2-methyl-	422	C30H62	001560-75-4	68





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Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

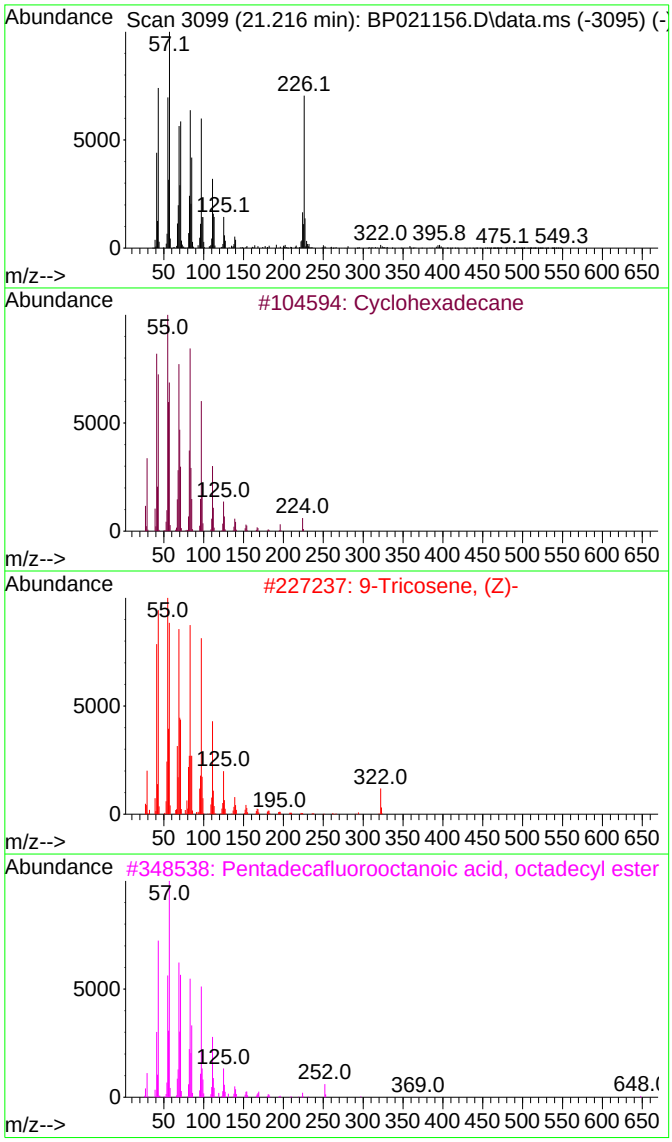
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

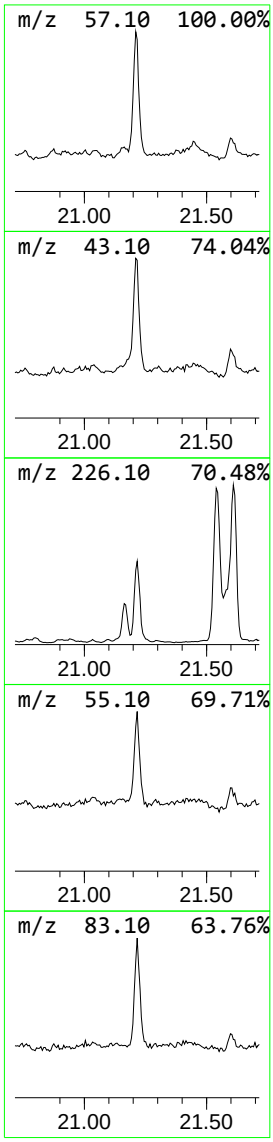
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 (DEL) Alkane: Cyclic21.216 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.216	10.93 ng/ul	3362120	Chrysene-d12	21.557

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	97
2	9-Tricosene, (Z)-	322	C23H46	027519-02-4	89
3	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	89
4	9-Tricosene, (Z)-	322	C23H46	027519-02-4	87
5	Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
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Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

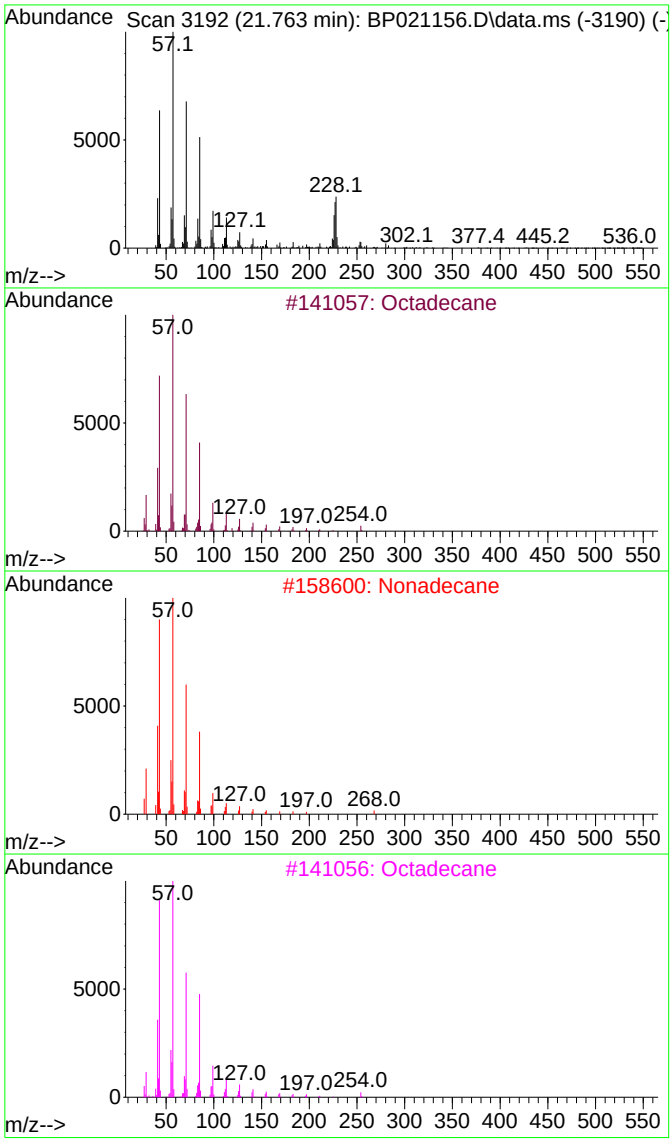
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

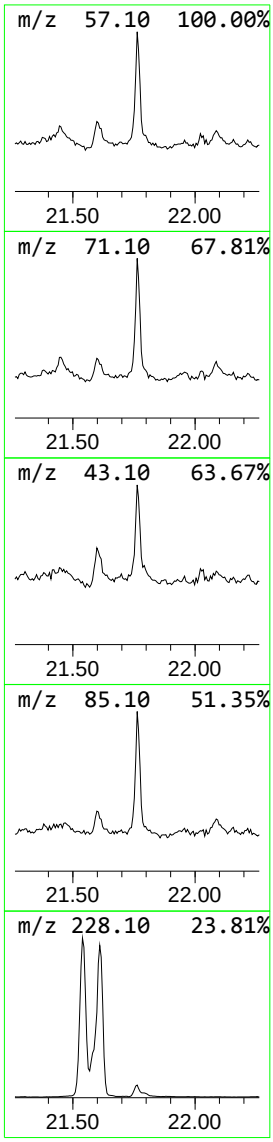
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.763	2.80 ng/ul	862014	Chrysene-d12	21.557

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	98
2	Nonadecane	268	C19H40	000629-92-5	91
3	Octadecane	254	C18H38	000593-45-3	64
4	Tetracosane	338	C24H50	000646-31-1	64
5	Tetracosane, 1-iodo-	464	C24H49I	1000406-32-0	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

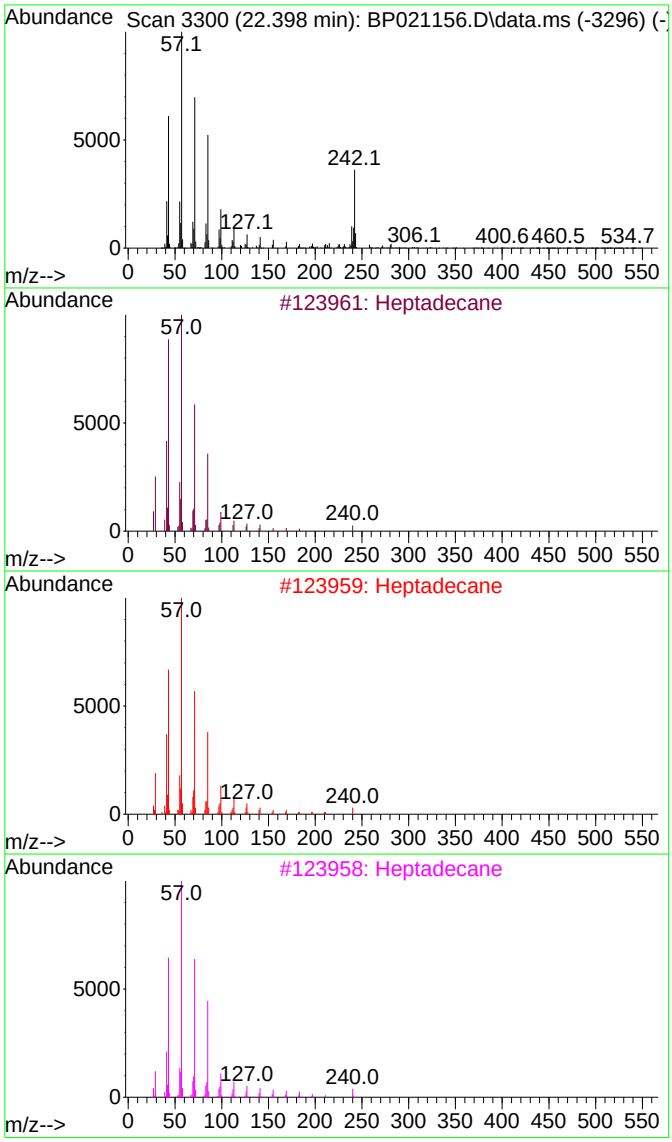
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

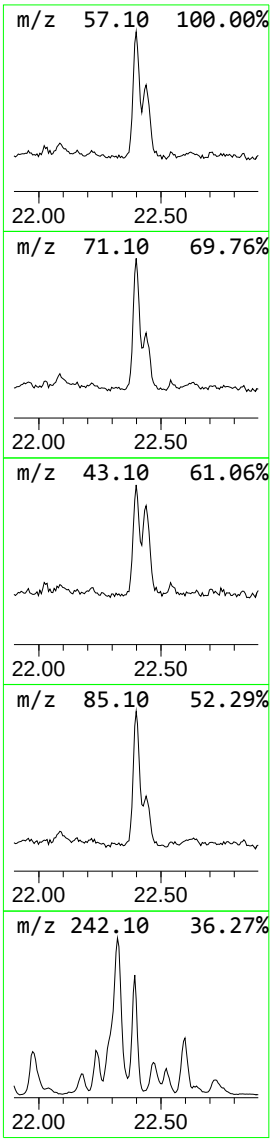
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.398	5.61 ng/ul	1725220	Chrysene-d12	21.557

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	93
2	Heptadecane	240	C17H36	000629-78-7	92
3	Heptadecane	240	C17H36	000629-78-7	91
4	Hexadecane	226	C16H34	000544-76-3	87
5	Heptadecane	240	C17H36	000629-78-7	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

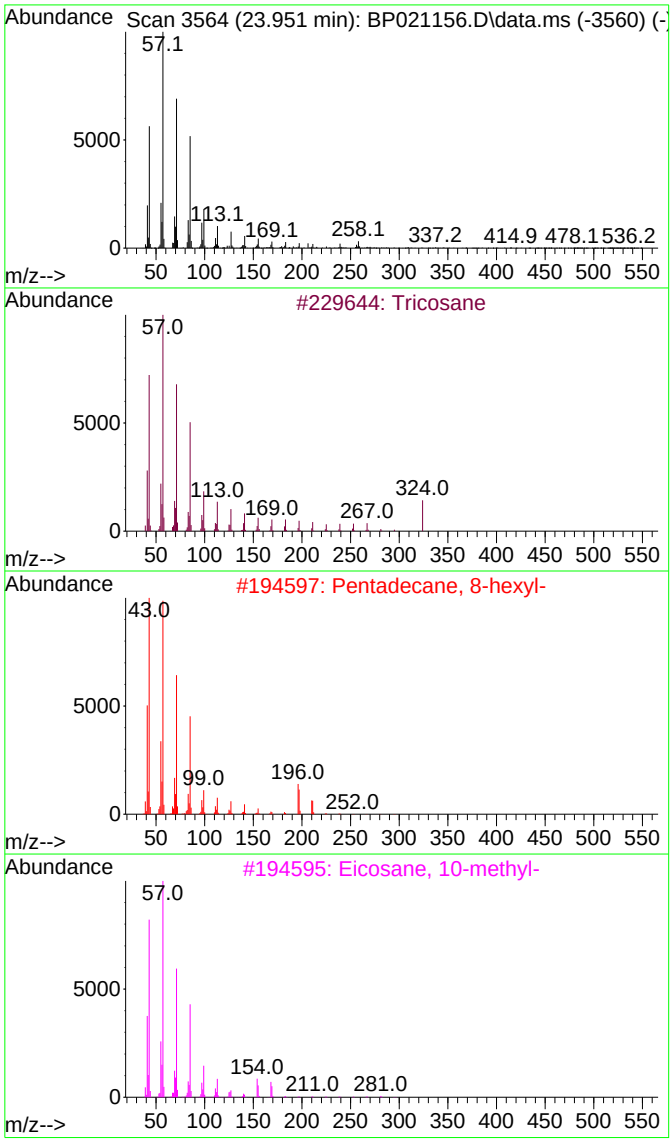
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

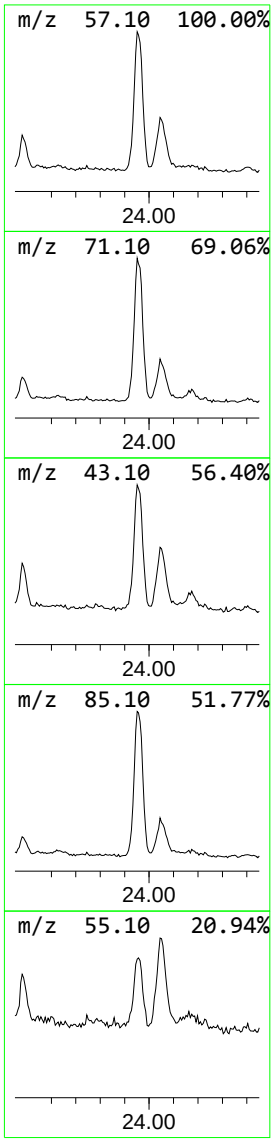
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.951	53.46 ng/ul	4079660	Perylene-d12	24.898

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricosane	324	C23H48	000638-67-5	95
2	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
3	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4	Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
5	Triacontane	422	C30H62	000638-68-6	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

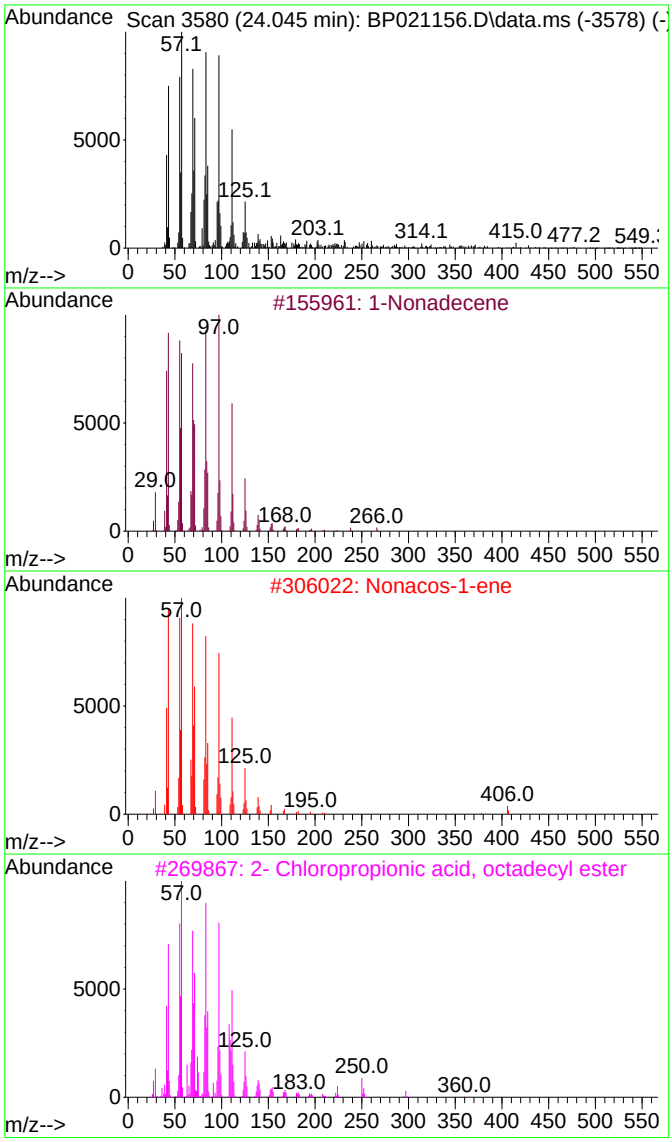
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

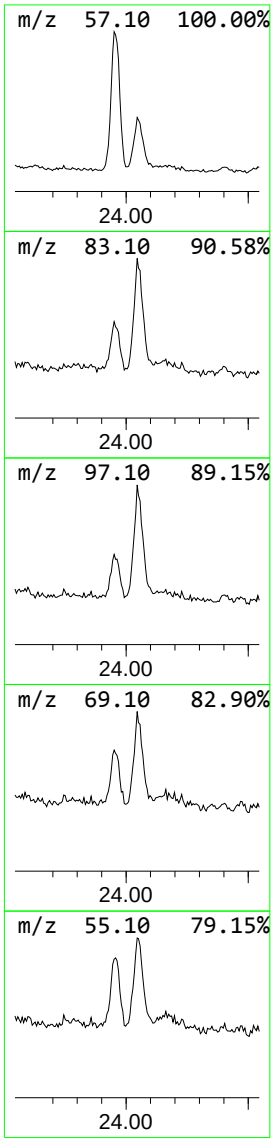
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.045	20.00 ng/ul	1526130	Perylene-d12	24.898

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene	266	C19H38	018435-45-5	98
2	Nonacos-1-ene	406	C29H58	018835-35-3	94
3	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5	1-Triacontanol	438	C30H62O	000593-50-0	93





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Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

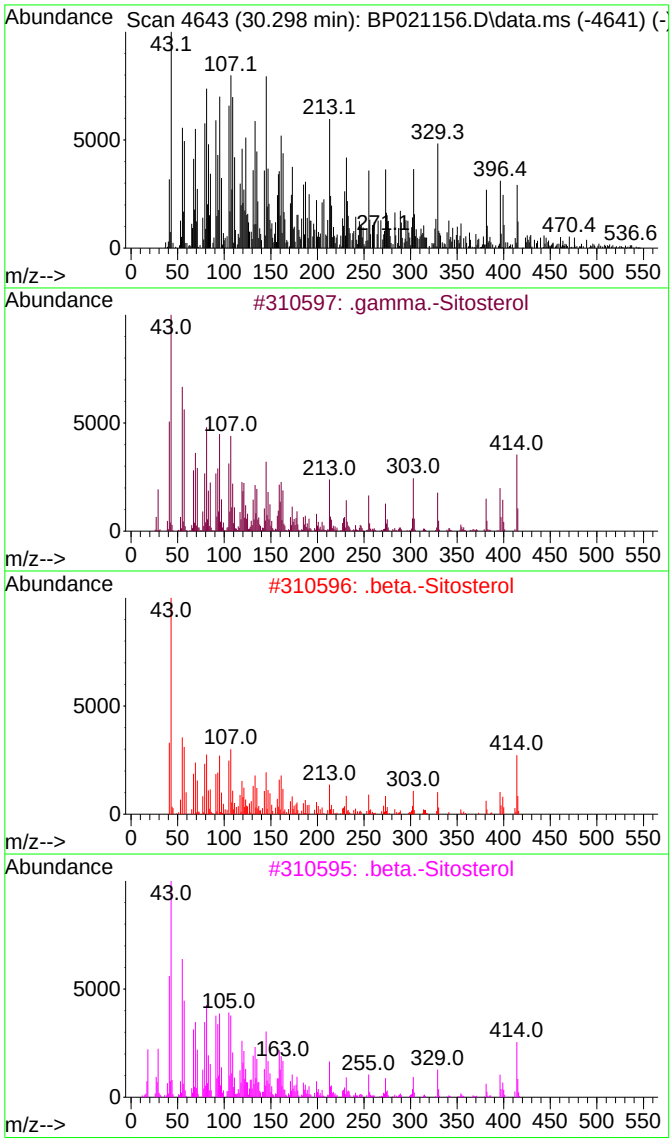
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

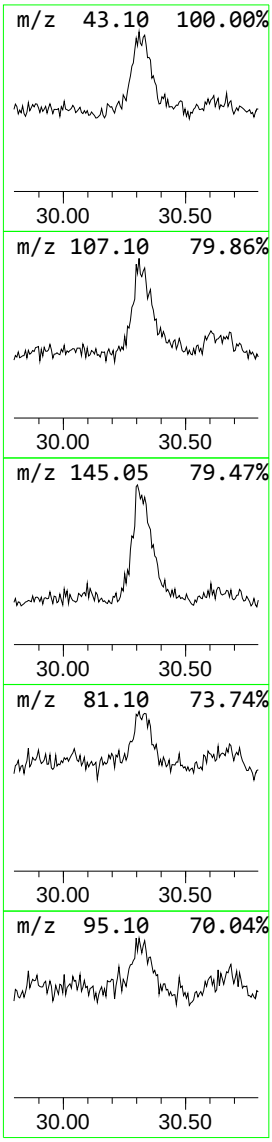
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 .gamma.-Sitosterol Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.298	3.94 ng/ul	300426	Perylene-d12	24.898

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.gamma.-Sitosterol	414	C29H50O	000083-47-6	91
2	.beta.-Sitosterol	414	C29H50O	000083-46-5	60
3	.beta.-Sitosterol	414	C29H50O	000083-46-5	50
4	Acetamide, 2-[[5-(2-methyl-3-fur...	329	C16H15N3O3S	1000350-00-2	38
5	N-(3-Chloro-4-fluoro-phenyl)-2-(...	329	C14H10Cl2FNOS	1000295-12-7	38





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021156.D
Acq On : 25 Jul 2024 21:00
Operator : MA/JU
Sample : P3263-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CK4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
1-Phenoxypropan...	11.169	3.3	ng/ul	257282	2	10.422	1577000	20.0
unknown-01	14.510	6.8	ng/ul	662089	3	14.287	1963020	20.0
(DEL) Alkane: S...	16.046	2.4	ng/ul	294544	4	17.104	2493850	20.0
1,1'-Biphenyl, ...	17.728	3.6	ng/ul	445769	4	17.104	2493850	20.0
Anthracene, 2-m...	18.004	4.3	ng/ul	535304	4	17.104	2493850	20.0
n-Hexadecanoic ...	18.045	17.6	ng/ul	2193310	4	17.104	2493850	20.0
2,2',4,5-Tetrac...	18.210	14.8	ng/ul	1848030	4	17.104	2493850	20.0
1,1'-Biphenyl, ...	18.263	6.3	ng/ul	789535	4	17.104	2493850	20.0
1,1'-Biphenyl, ...	18.316	4.2	ng/ul	522911	4	17.104	2493850	20.0
1,1'-Biphenyl, ...	18.504	4.2	ng/ul	526598	4	17.104	2493850	20.0
Naphthalene, 2-...	18.534	4.2	ng/ul	525635	4	17.104	2493850	20.0
9,10-Anthracene...	18.598	5.1	ng/ul	631779	4	17.104	2493850	20.0
Phenanthrene, 2...	18.969	6.4	ng/ul	792545	4	17.104	2493850	20.0
1,1'-Biphenyl, ...	19.022	5.8	ng/ul	716656	4	17.104	2493850	20.0
1,1'-Biphenyl, ...	19.069	10.0	ng/ul	1247940	4	17.104	2493850	20.0
3,3',4,5-Tetrac...	19.104	10.8	ng/ul	1346890	4	17.104	2493850	20.0
1,1'-Biphenyl, ...	19.387	4.5	ng/ul	1368200	5	21.557	6151040	20.0
(DEL) Alkane: S...	20.680	2.4	ng/ul	751295	5	21.557	6151040	20.0
(DEL) Alkane: C...	21.216	10.9	ng/ul	3362120	5	21.557	6151040	20.0
(DEL) Alkane: S...	21.763	2.8	ng/ul	862014	5	21.557	6151040	20.0
(DEL) Alkane: S...	22.398	5.6	ng/ul	1725220	5	21.557	6151040	20.0
(DEL) Alkane: S...	23.951	53.5	ng/ul	4079660	6	24.898	1526130	20.0
(DEL) Alkane: S...	24.045	20.0	ng/ul	1526130	6	24.898	1526130	20.0
.gamma.-Sitosterol	30.298	3.9	ng/ul	300426	6	24.898	1526130	20.0