

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
 Data File : BP008524.D
 Acq On : 23 Dec 2021 14:53
 Operator : CG/JU
 Sample : M5121-04ME
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGL53ME

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-BP122321.M
 Title : SVOA CALIBRATION

Signal : TIC: BP008524.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.023	3	6	14	rVB	1456778	2100681	8.23%	1.097%
2	3.093	14	18	38	rVB	2411409	3326934	13.03%	1.737%
3	3.352	58	62	68	rVB	84708	120013	0.47%	0.063%
4	3.758	128	131	146	rVB	307625	454574	1.78%	0.237%
5	4.099	186	189	197	rVB3	30209	45908	0.18%	0.024%
6	7.058	683	692	703	rBV	1361003	2231061	8.74%	1.165%
7	7.234	714	722	731	rBV	1152629	1863111	7.30%	0.973%
8	7.434	748	756	768	rBV	1438583	2290055	8.97%	1.196%
9	7.905	829	836	844	rVB	1663547	2675713	10.48%	1.397%
10	8.605	948	955	966	rBV	1756384	2838479	11.12%	1.482%
11	8.722	966	975	984	rVB3	28504	68559	0.27%	0.036%
12	9.057	1020	1032	1041	rBV	1373760	2198261	8.61%	1.148%
13	9.157	1045	1049	1057	rVB2	29118	55744	0.22%	0.029%
14	9.534	1106	1113	1118	rBV	54104	91003	0.36%	0.048%
15	9.787	1149	1156	1163	rBV	813810	1383434	5.42%	0.722%
16	10.322	1239	1247	1258	rBV	1800989	2935854	11.50%	1.533%
17	10.569	1284	1289	1296	rBV2	41275	74050	0.29%	0.039%
18	10.710	1305	1313	1318	rBV	2183067	3953632	15.48%	2.064%
19	10.757	1318	1321	1327	rVB	287026	447466	1.75%	0.234%
20	10.834	1327	1334	1343	rBV	1345964	2307165	9.03%	1.205%
21	11.757	1483	1491	1496	rBV4	32422	65853	0.26%	0.034%
22	12.151	1551	1558	1563	rBV	75501	120298	0.47%	0.063%
23	12.293	1574	1582	1588	rBV2	44182	74436	0.29%	0.039%
24	12.369	1589	1595	1602	rVB2	40550	76992	0.30%	0.040%
25	12.963	1691	1696	1700	rBV3	34091	62561	0.24%	0.033%
26	13.046	1706	1710	1715	rBV7	26420	44656	0.17%	0.023%
27	13.104	1715	1720	1724	rVV	65754	91671	0.36%	0.048%
28	13.340	1755	1760	1763	rBV3	30483	48446	0.19%	0.025%
29	13.387	1763	1768	1777	rVB2	257314	366565	1.44%	0.191%
30	13.863	1845	1849	1853	rBV6	35394	53787	0.21%	0.028%
31	13.945	1854	1863	1872	rVV	3207504	4585412	17.96%	2.394%
32	14.045	1876	1880	1884	rVV2	124896	174929	0.69%	0.091%
33	14.087	1884	1887	1895	rVB3	48998	84475	0.33%	0.044%
34	14.228	1902	1911	1918	rVV	4076505	6022719	23.59%	3.145%
35	14.381	1934	1937	1945	rVB3	92324	151615	0.59%	0.079%
36	14.457	1945	1950	1954	rBV	332889	457014	1.79%	0.239%

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Integrator: RTE

Smoothing : OFF

Filtering: 5

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
 Title : SVOA CALIBRATION

37	14.534	1956	1963	1971	rVV	3545493	5423915	21.24%	2.832%
38	14.710	1987	1993	2002	rBV	1068947	1546672	6.06%	0.808%
39	14.945	2030	2033	2039	rVB4	80863	145393	0.57%	0.076%
40	15.075	2052	2055	2063	rVB3	101616	148044	0.58%	0.077%
41	15.145	2063	2067	2070	rBV4	58095	100706	0.39%	0.053%
42	15.345	2096	2101	2107	rBV5	150884	285560	1.12%	0.149%
43	15.404	2107	2111	2116	rBV	343769	464792	1.82%	0.243%
44	15.528	2125	2132	2139	rVB	4969825	7350334	28.78%	3.838%
45	15.592	2140	2143	2145	rBV2	44553	61238	0.24%	0.032%
46	15.634	2145	2150	2155	rVB	657897	951159	3.72%	0.497%
47	15.710	2160	2163	2166	rBV2	99833	134343	0.53%	0.070%
48	15.745	2167	2169	2174	rVB	105804	112417	0.44%	0.059%
49	15.816	2177	2181	2185	rBV2	221385	354114	1.39%	0.185%
50	15.898	2191	2195	2202	rVB5	116614	214752	0.84%	0.112%
51	15.969	2203	2207	2211	rBV2	176493	285660	1.12%	0.149%
52	16.181	2241	2243	2246	rVB	95278	88477	0.35%	0.046%
53	16.269	2254	2258	2260	rBV2	432482	573357	2.25%	0.299%
54	16.298	2260	2263	2267	rVB	455848	611091	2.39%	0.319%
55	16.387	2274	2278	2283	rVB3	148593	241477	0.95%	0.126%
56	16.539	2300	2304	2307	rBV	228401	291258	1.14%	0.152%
57	16.581	2308	2311	2315	rVV	137208	174281	0.68%	0.091%
58	17.069	2390	2394	2398	rBV	382149	460867	1.80%	0.241%
59	17.122	2400	2403	2413	rVB	287615	433591	1.70%	0.226%
60	17.286	2425	2431	2441	rBV	3940839	5944144	23.28%	3.104%
61	17.381	2441	2447	2454	rVB	5098734	7507006	29.40%	3.920%
62	17.804	2516	2519	2524	rVB	343903	396881	1.55%	0.207%
63	18.475	2630	2633	2637	rVB	322021	354139	1.39%	0.185%
64	18.710	2669	2673	2679	rBV	350472	453313	1.78%	0.237%
65	18.822	2689	2692	2697	rVB	415093	479140	1.88%	0.250%
66	19.080	2733	2736	2741	rBV	324826	411195	1.61%	0.215%
67	19.233	2759	2762	2766	rBV2	312696	431007	1.69%	0.225%
68	19.610	2821	2826	2836	rBV	5526791	8612572	33.73%	4.497%
69	20.163	2916	2920	2924	rVB	623682	697790	2.73%	0.364%
70	20.410	2959	2962	2965	rVB	265858	278411	1.09%	0.145%
71	20.522	2977	2981	2987	rVB	1440585	1599676	6.26%	0.835%
72	20.645	2998	3002	3005	rBV	937364	997161	3.90%	0.521%
73	20.869	3034	3040	3046	rVB2	1375402	2103316	8.24%	1.098%
74	20.933	3048	3051	3056	rVB	308117	373424	1.46%	0.195%
75	20.998	3058	3062	3067	rVB	911110	1012516	3.97%	0.529%
76	21.098	3075	3079	3084	rVB	1317993	1598268	6.26%	0.835%

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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 >
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 Title : SVOA CALIBRATION

77	21.292	3107	3112	3120	rVV	21087024	25535866	100.00%	13.333%
78	21.369	3120	3125	3132	rVB	3842096	5421988	21.23%	2.831%
79	21.516	3147	3150	3151	rBV	492875	503322	1.97%	0.263%
80	21.545	3151	3155	3158	rVB	1357827	1624845	6.36%	0.848%
81	21.698	3178	3181	3185	rBV	428037	552053	2.16%	0.288%
82	21.745	3185	3189	3194	rVB	684548	889604	3.48%	0.465%
83	21.839	3201	3205	3209	rVB	1468148	1922280	7.53%	1.004%
84	21.892	3209	3214	3221	rVB3	1409397	2190944	8.58%	1.144%
85	22.010	3230	3234	3239	rBV	1407712	1956981	7.66%	1.022%
86	22.239	3267	3273	3284	rVB	4039215	6981810	27.34%	3.646%
87	22.527	3317	3322	3333	rVB	1469863	2526156	9.89%	1.319%
88	22.763	3358	3362	3367	rVB	765570	1096060	4.29%	0.572%
89	22.857	3373	3378	3387	rVB	1968841	3650690	14.30%	1.906%
90	22.957	3390	3395	3406	rVB	1145080	2361861	9.25%	1.233%
91	23.098	3414	3419	3424	rBV	1128489	1879808	7.36%	0.982%
92	23.357	3457	3463	3465	rBV	3117237	5504025	21.55%	2.874%
93	23.645	3505	3512	3520	rVV2	3680510	7399234	28.98%	3.863%
94	23.745	3524	3529	3533	rVV2	1196785	2495177	9.77%	1.303%
95	23.804	3533	3539	3554	rVB2	2776916	6922511	27.11%	3.615%
96	24.121	3588	3593	3597	rBV	649138	1273950	4.99%	0.665%
97	24.498	3652	3657	3665	rVB	738095	1489813	5.83%	0.778%
98	24.815	3703	3711	3723	rVB	2518502	6218936	24.35%	3.247%
99	25.374	3801	3806	3814	rVB2	462659	1028492	4.03%	0.537%
100	26.786	4039	4046	4061	rVB	771499	2443937	9.57%	1.276%

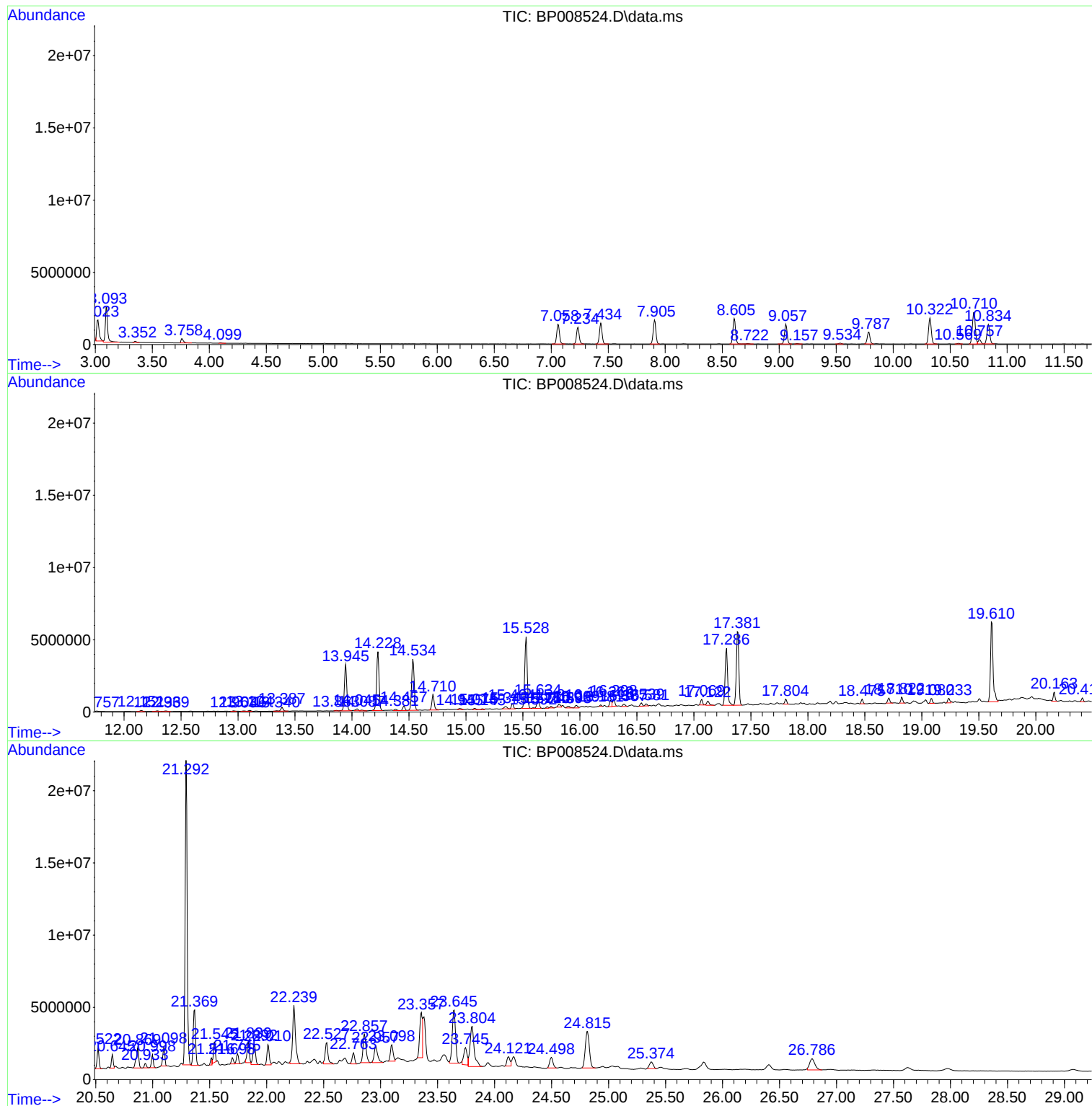
Sum of corrected areas: 191516926

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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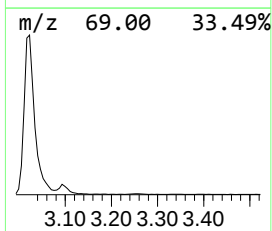
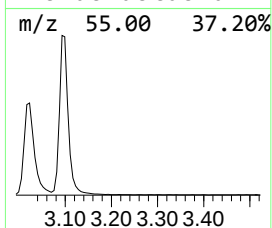
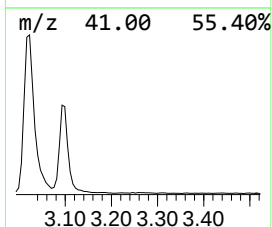
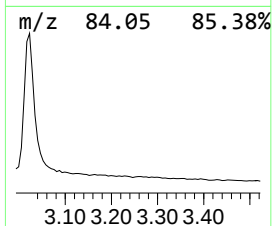
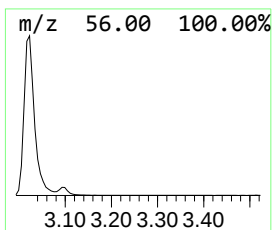
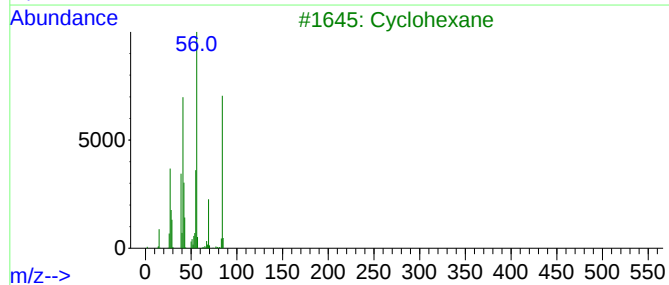
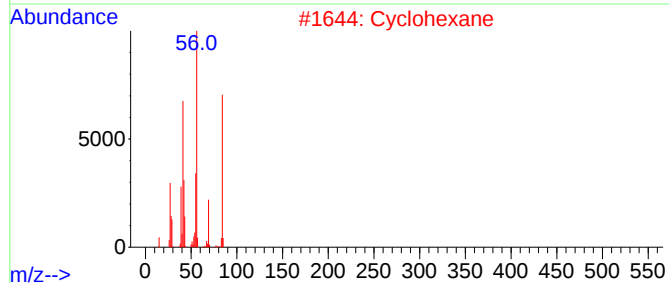
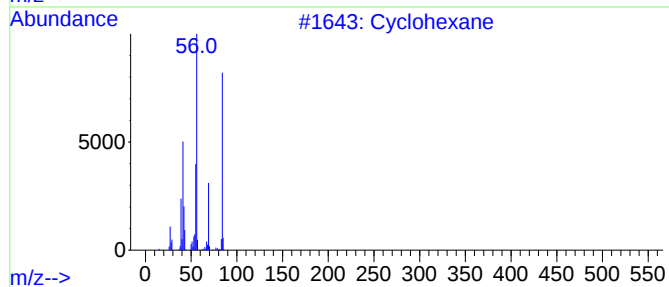
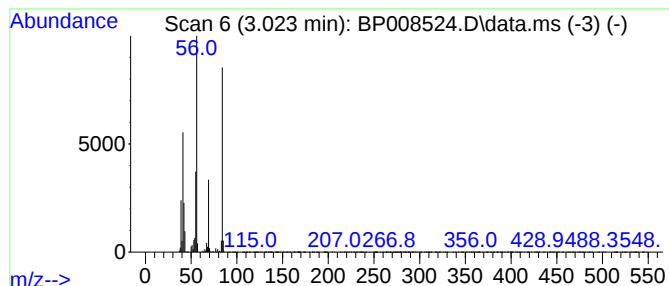
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.023 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.023	15.70 ng/ul	2100680	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane	84	C6H12	000110-82-7	95
2			Cyclohexane	84	C6H12	000110-82-7	91
3			Cyclohexane	84	C6H12	000110-82-7	90
4			Cyclopentane, methyl-	84	C6H12	000096-37-7	83
5			Cyclohexane	84	C6H12	000110-82-7	80



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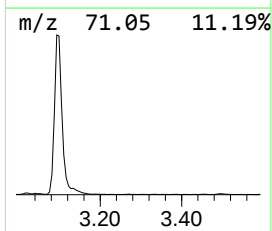
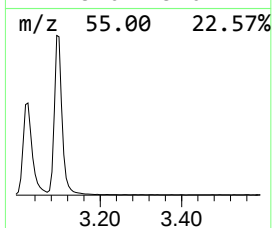
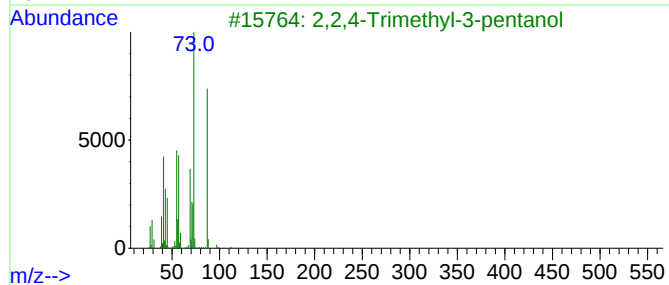
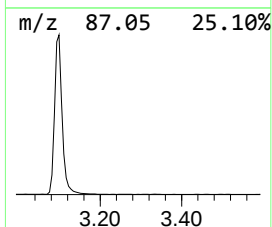
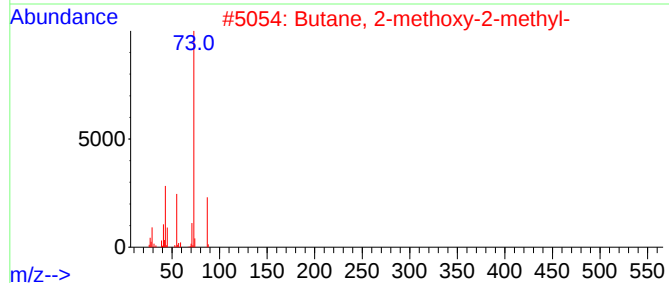
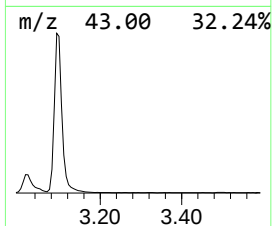
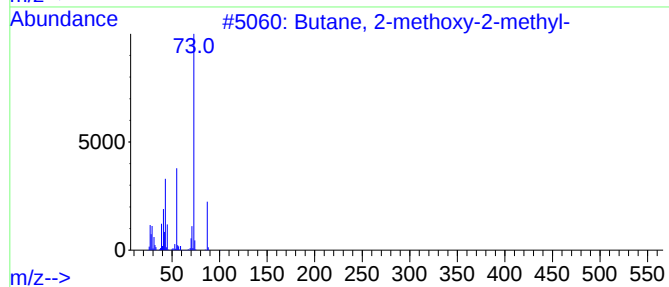
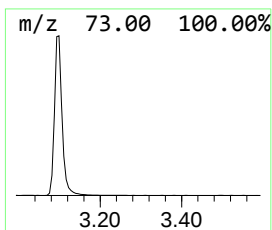
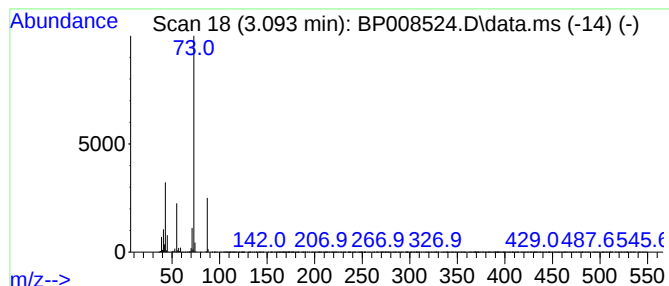
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.093	24.87 ng/ul	3326930	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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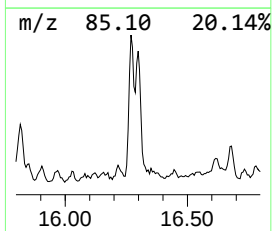
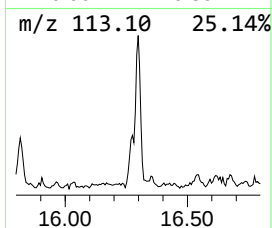
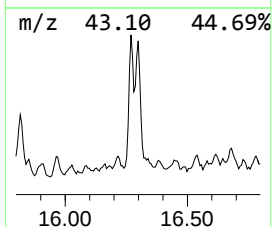
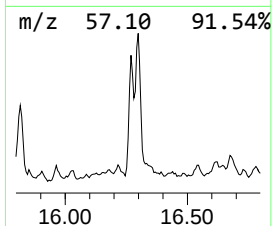
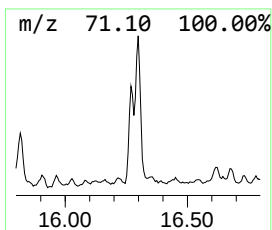
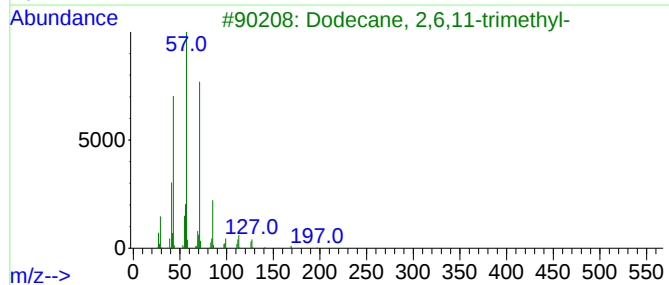
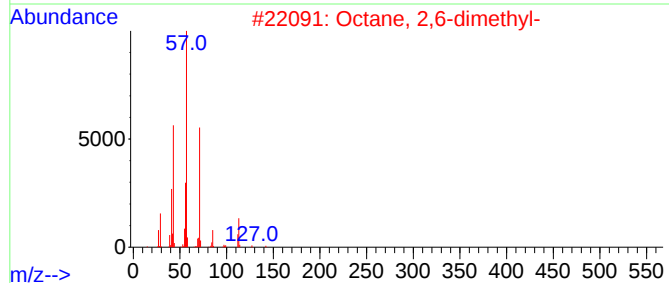
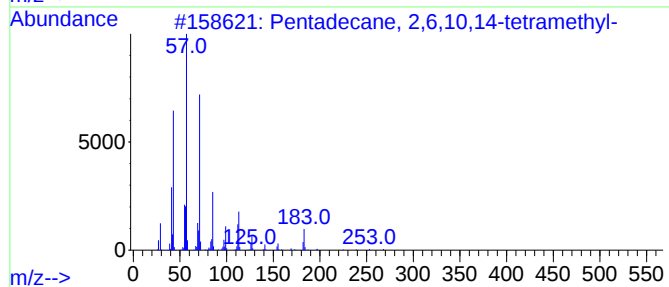
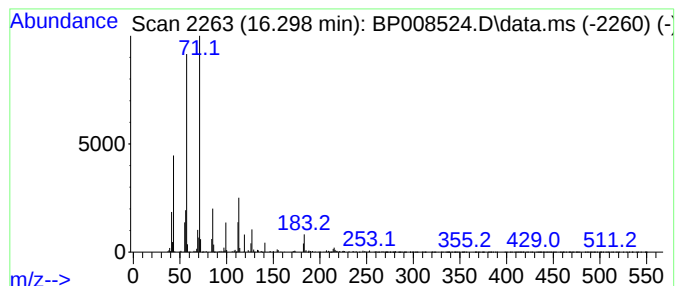
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.298	2.06 ng/ul	611091	Phenanthrene-d10	17.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	86
2			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	80
3			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	64
4			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
5			Dodecane	170	C12H26	000112-40-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
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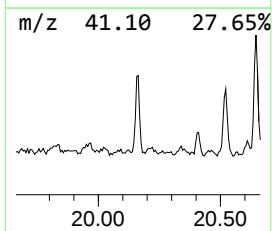
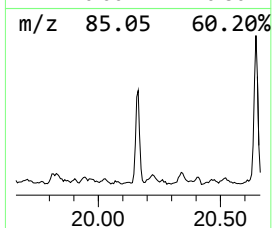
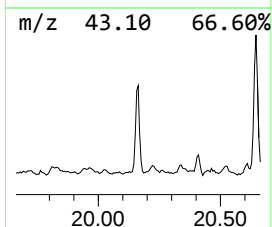
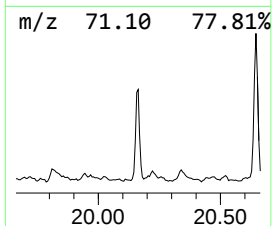
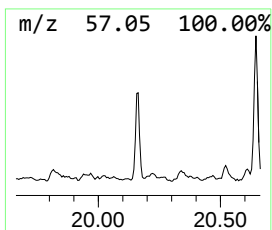
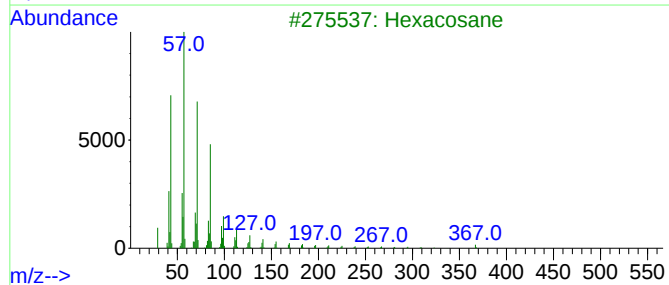
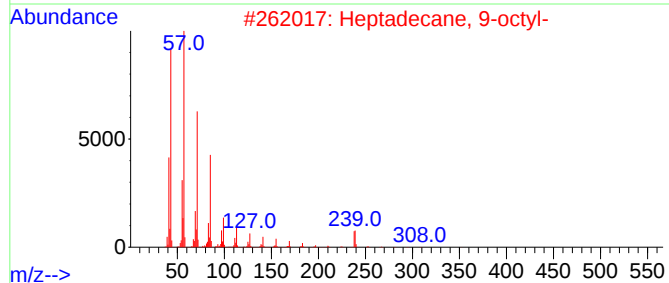
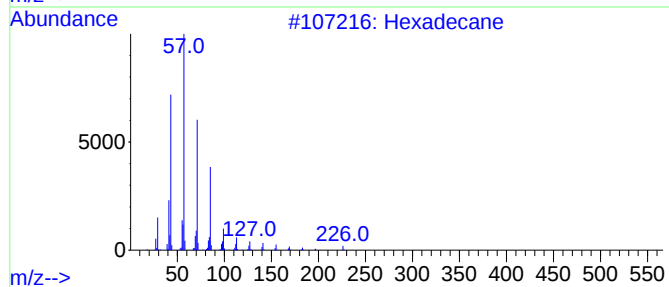
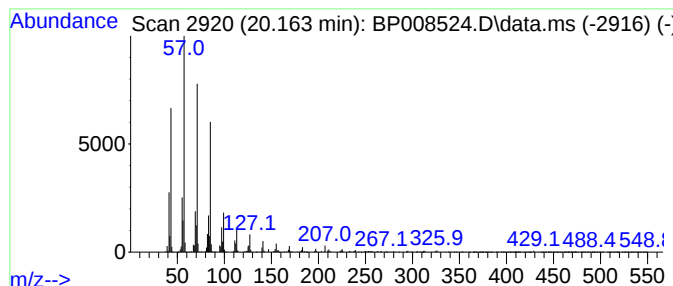
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.163	2.57 ng/ul	697790	Chrysene-d12		21.363	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	95
2	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	95
3	Hexacosane		366	C26H54	000630-01-3	91
4	Octacosane, 2-methyl-		408	C29H60	001560-98-1	91
5	Eicosane		282	C20H42	000112-95-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
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Operator : CG/JU
Sample : M5121-04ME
Misc :
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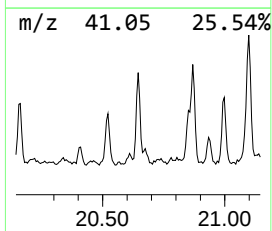
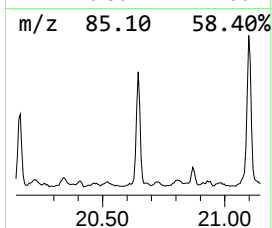
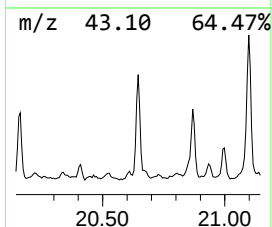
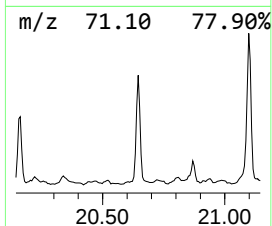
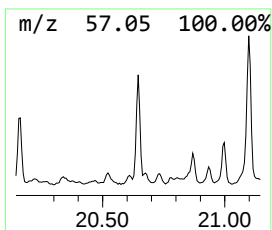
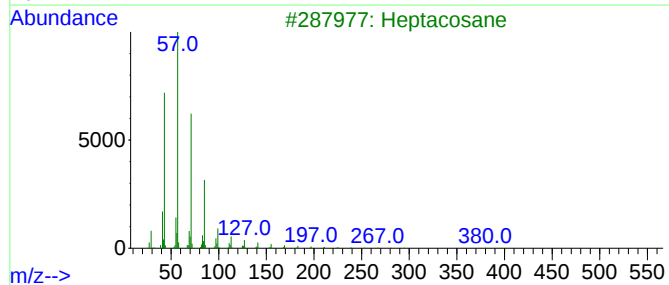
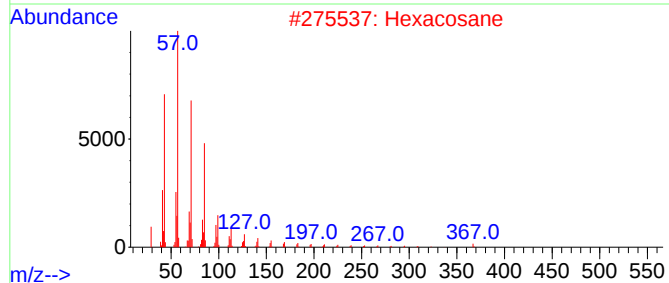
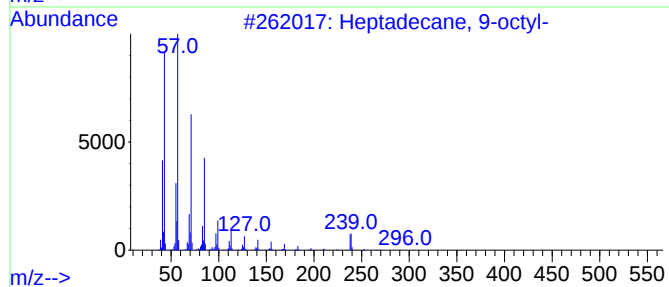
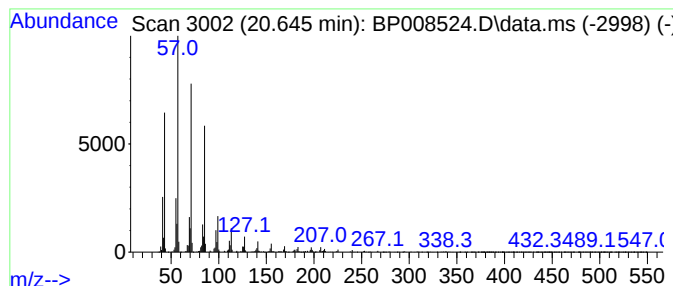
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.645	3.68 ng/ul	997161	Chrysene-d12	21.363	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	96
2	Hexacosane	366	C26H54	000630-01-3	94
3	Heptacosane	380	C27H56	000593-49-7	91
4	Pentacosane	352	C25H52	000629-99-2	91
5	Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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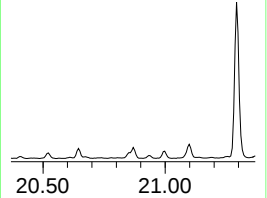
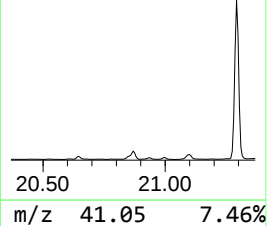
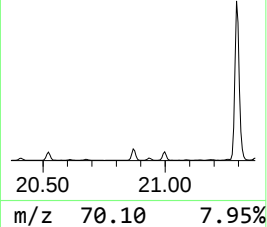
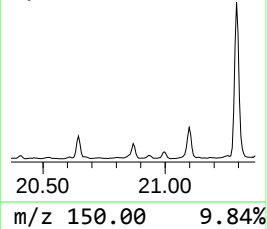
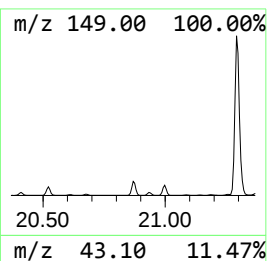
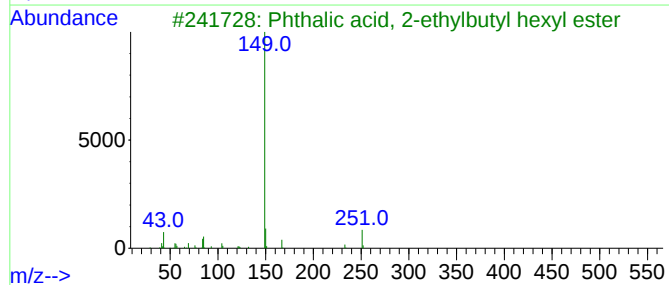
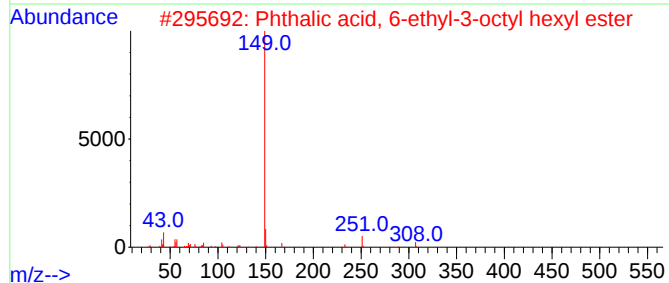
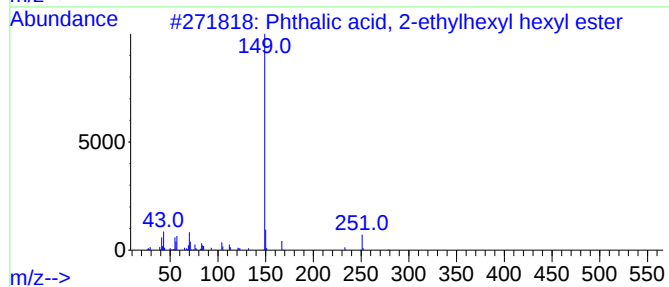
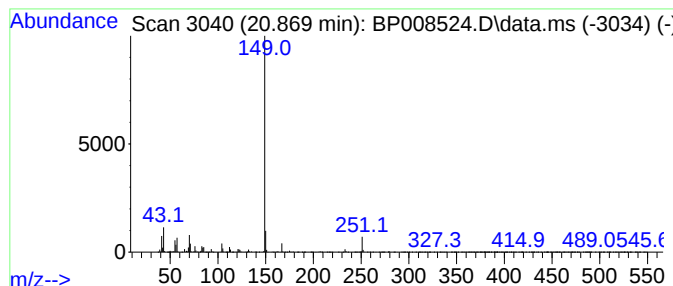
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phthalic acid, 2-ethylhexyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.869	7.76 ng/ul	2103320	Chrysene-d12	21.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, 2-ethylhexyl hexy...	362	C22H34O4	1000309-02-5	91
2			Phthalic acid, 6-ethyl-3-octyl h...	390	C24H38O4	1000315-17-6	91
3			Phthalic acid, 2-ethylbutyl hexy...	334	C20H30O4	1000356-89-1	90
4			Phthalic acid, isohexyl 2-methyl...	320	C19H28O4	1000322-81-3	86
5			1,2-Benzenedicarboxylic acid, de...	390	C24H38O4	025724-58-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
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Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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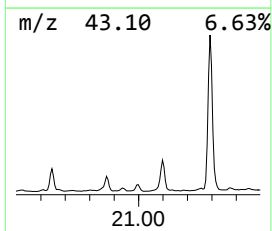
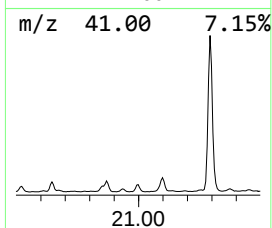
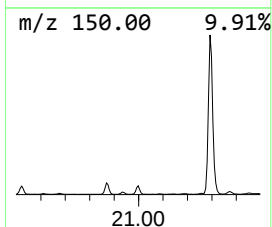
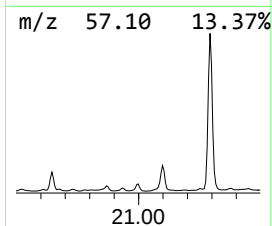
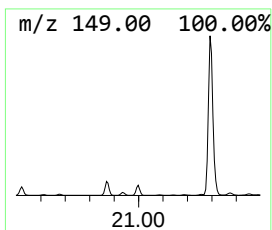
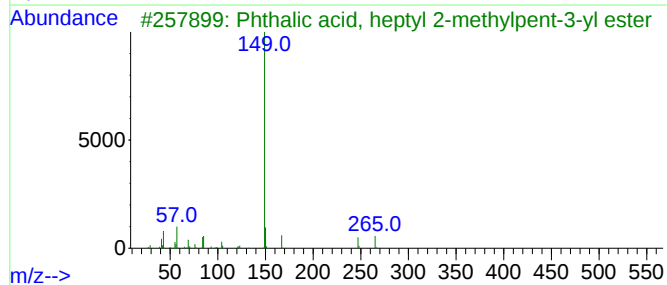
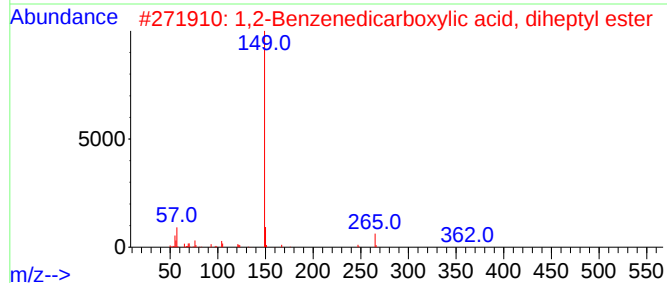
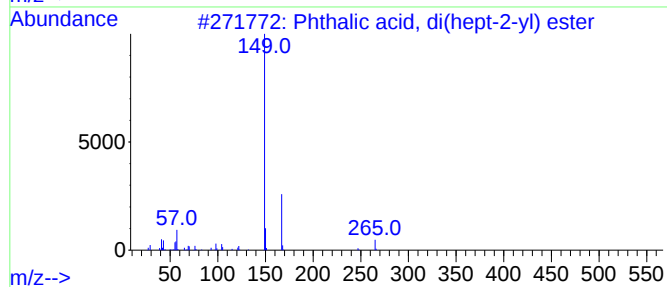
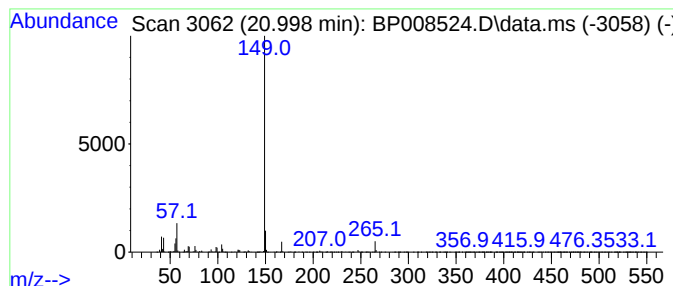
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Phthalic acid, di(hept-2-yl... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.998	3.73 ng/ul	1012520	Chrysene-d12	21.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, di(hept-2-yl) ester	362	C22H34O4	1000356-98-3	87
2			1,2-Benzenedicarboxylic acid, di...	362	C22H34O4	003648-21-3	87
3			Phthalic acid, heptyl 2-methylpe...	348	C21H32O4	1000356-91-2	86
4			1,2-Benzenedicarboxylic acid, di...	362	C22H34O4	003648-21-3	86
5			Phthalic acid, 2-ethylbutyl hept...	348	C21H32O4	1000356-89-2	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
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Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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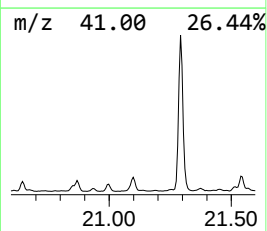
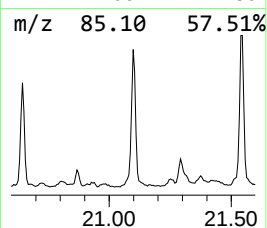
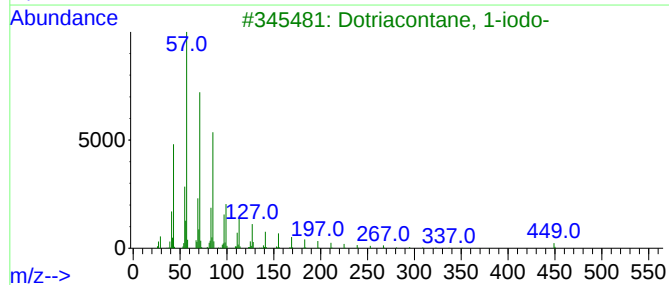
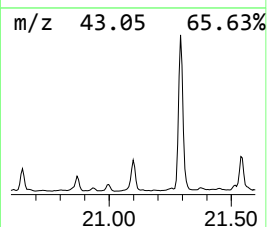
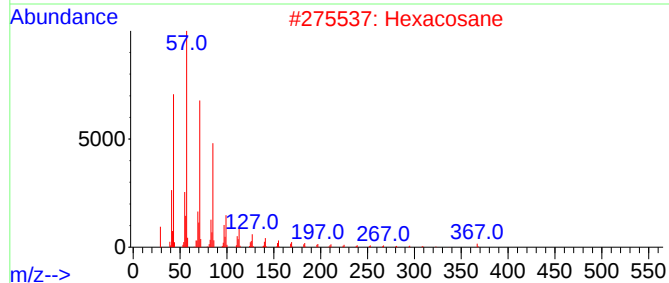
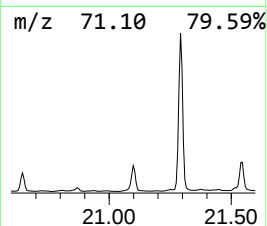
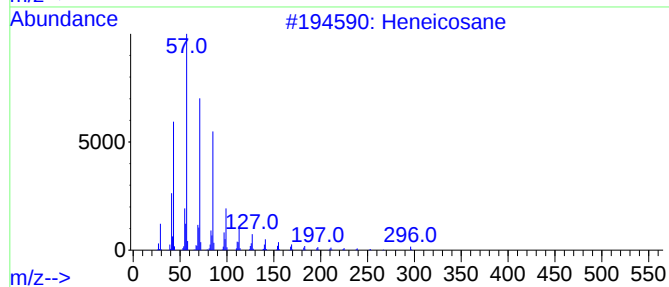
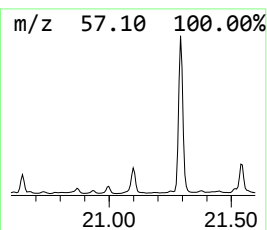
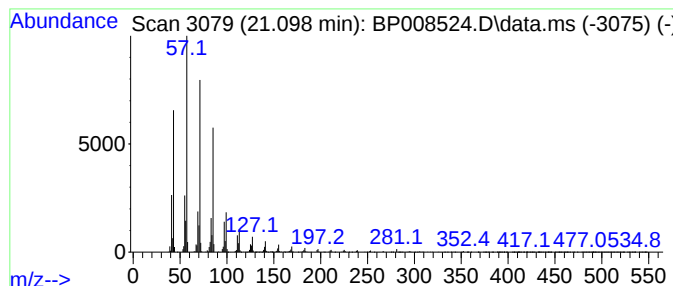
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.098	5.90 ng/ul	1598270	Chrysene-d12	21.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Hexacosane	366	C26H54	000630-01-3	91
3			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
4			Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
5			Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
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Sample : M5121-04ME
Misc :
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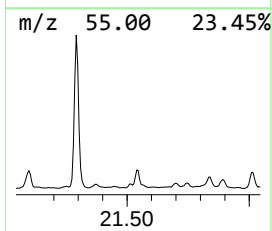
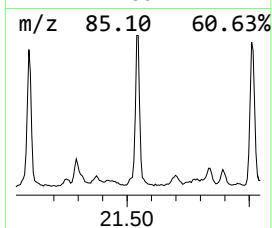
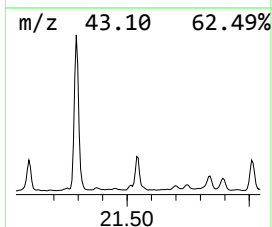
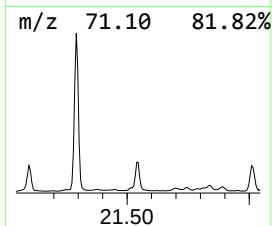
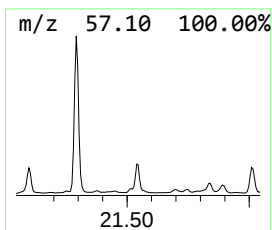
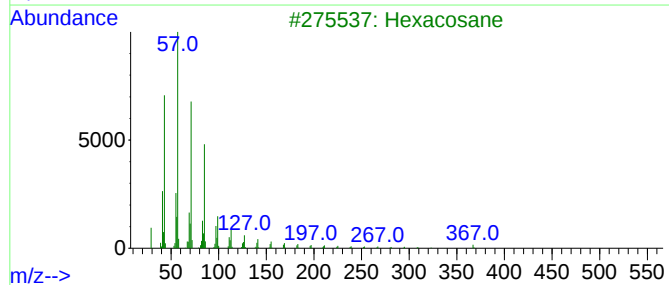
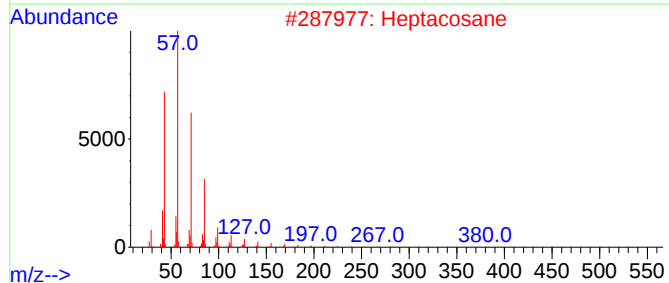
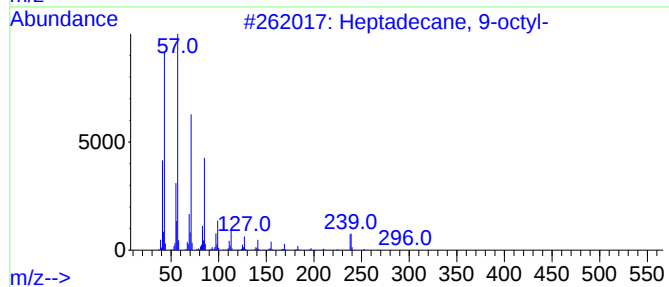
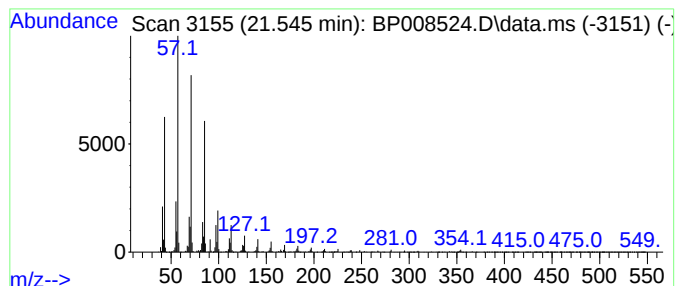
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.545	5.99 ng/ul	1624850	Chrysene-d12	21.363	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2	Heptacosane	380	C27H56	000593-49-7	91
3	Hexacosane	366	C26H54	000630-01-3	91
4	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5	Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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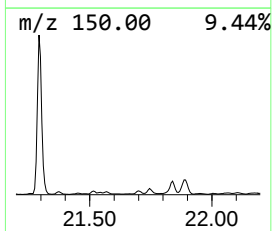
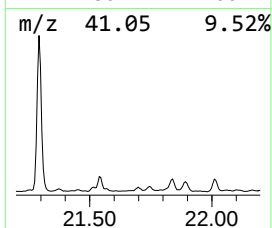
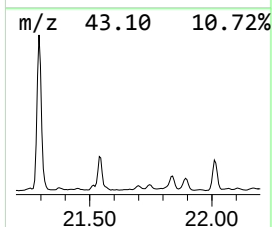
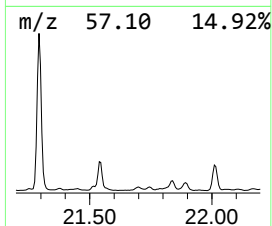
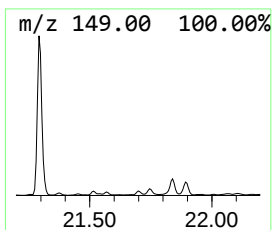
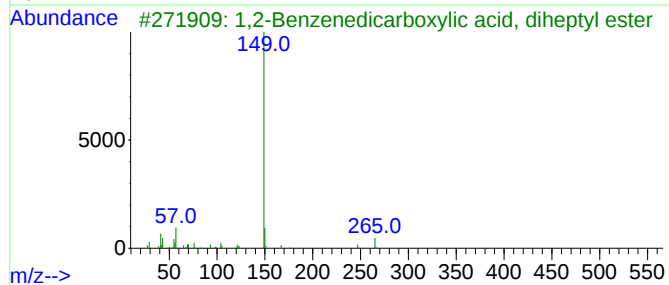
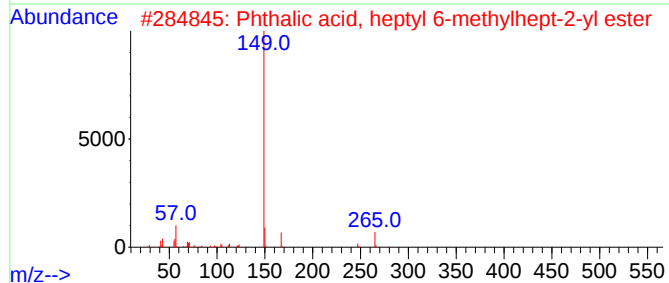
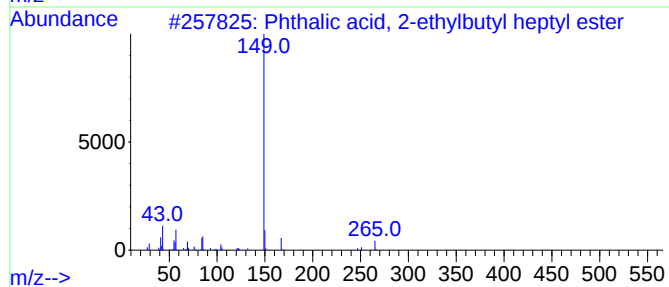
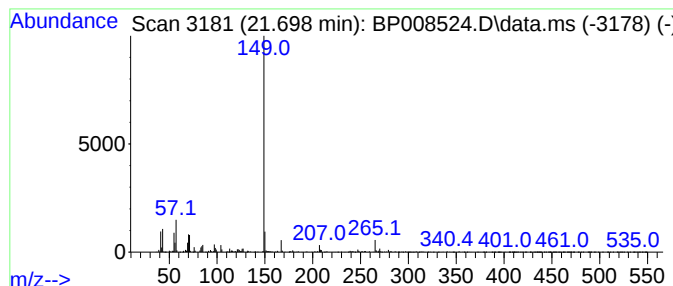
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Phthalic acid, 2-ethylbutyl... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.698	2.04 ng/ul	552053	Chrysene-d12	21.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, 2-ethylbutyl hept...	348	C21H32O4	1000356-89-2	86
2			Phthalic acid, heptyl 6-methylhe...	376	C23H36O4	1000377-96-1	86
3			1,2-Benzenedicarboxylic acid, di...	362	C22H34O4	003648-21-3	80
4			Phthalic acid, 6-methylhept-2-yl...	460	C29H48O4	1010377-96-7	80
5			Phthalic acid, heptyl 2-methylbu...	334	C20H30O4	1010322-81-5	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGL53ME

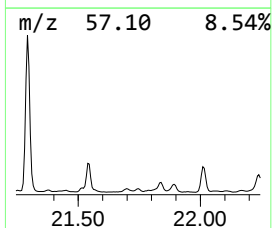
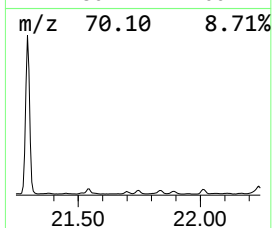
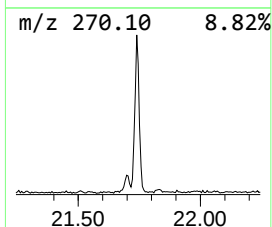
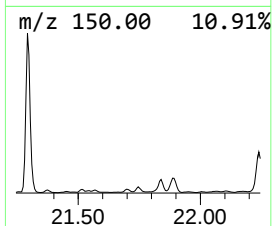
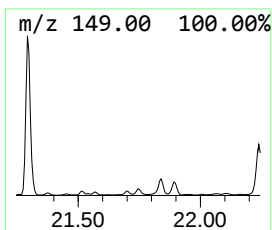
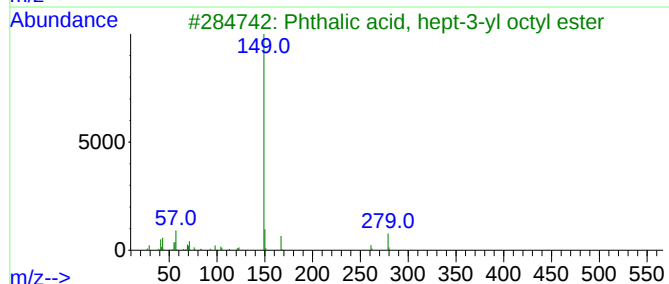
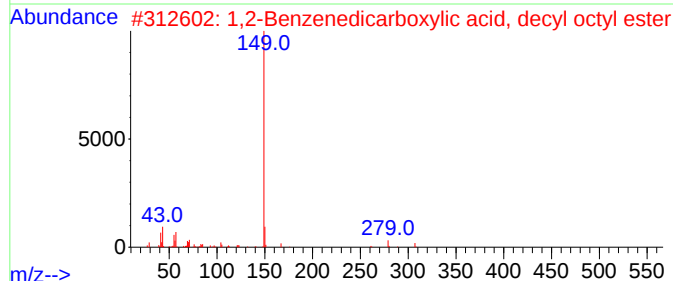
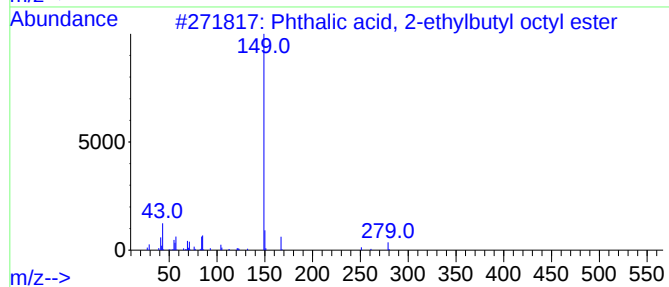
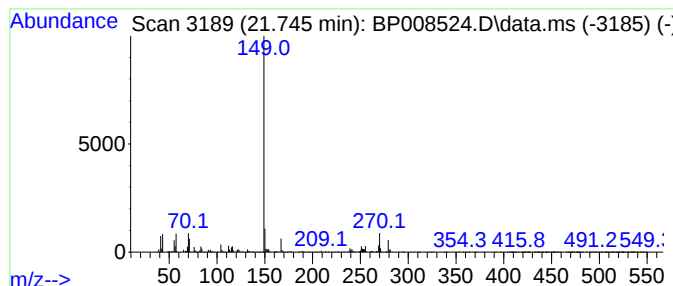
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Phthalic acid, 2-ethylbutyl... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.745	3.28 ng/ul	889604	Chrysene-d12	21.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, 2-ethylbutyl octyl...	362	C22H34O4	1000356-89-3	86
2			1,2-Benzenedicarboxylic acid, de...	418	C26H42O4	000119-07-3	80
3			Phthalic acid, hept-3-yl octyl e...	376	C23H36O4	1000356-99-2	80
4			Phthalic acid, 2-methylpent-3-yl...	362	C22H34O4	1000356-91-3	80
5			Phthalic acid, 2,4-dimethylpent...	376	C23H36O4	1000356-84-7	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGL53ME

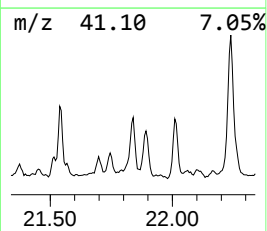
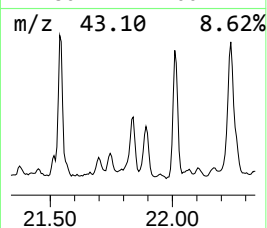
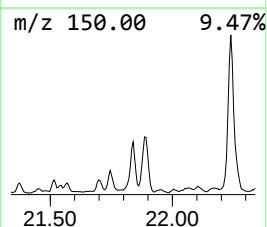
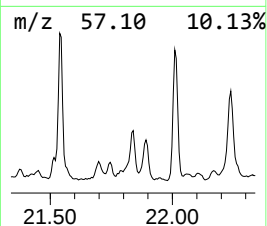
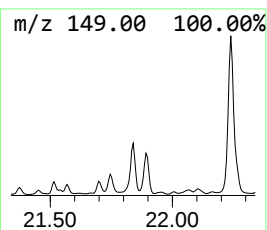
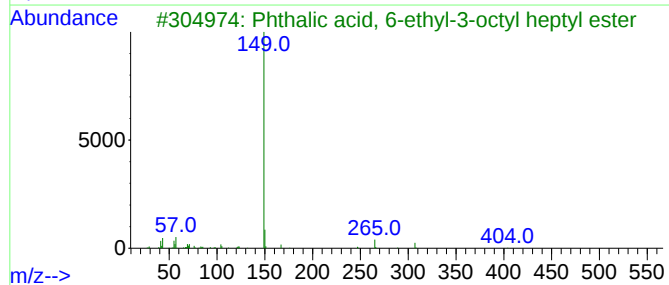
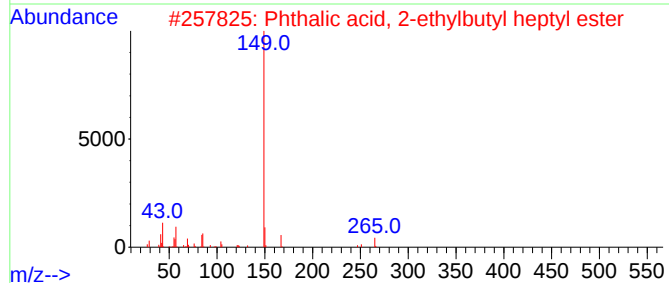
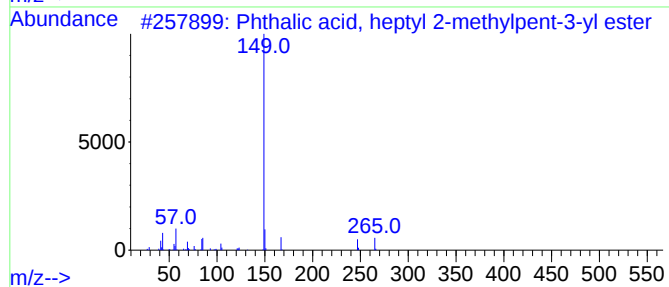
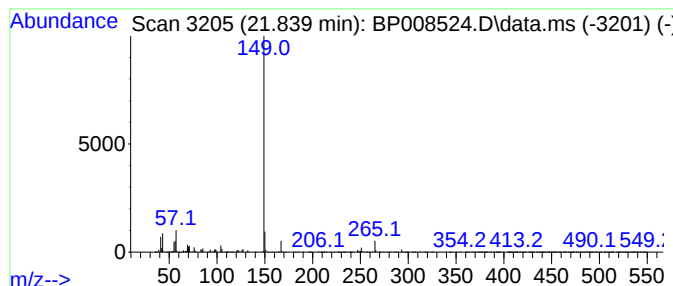
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Phthalic acid, heptyl 2-met... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.839	7.09 ng/ul	1922280	Chrysene-d12	21.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, heptyl 2-methylpe...	348	C21H32O4	1000356-91-2	86
2			Phthalic acid, 2-ethylbutyl hept...	348	C21H32O4	1000356-89-2	86
3			Phthalic acid, 6-ethyl-3-octyl h...	404	C25H40O4	1000315-17-7	86
4			Phthalic acid, heptyl 2-methylbu...	334	C20H30O4	1010322-81-5	86
5			Phthalic acid, heptyl hept-4-yl ...	362	C22H34O4	1000356-78-8	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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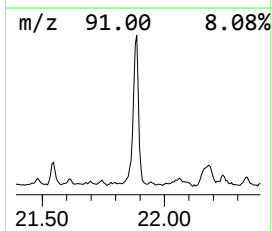
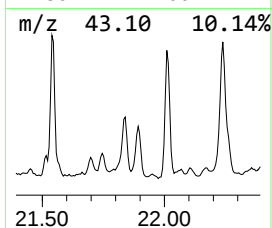
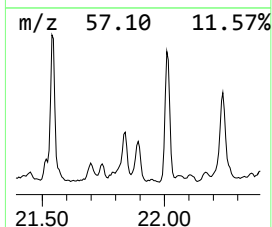
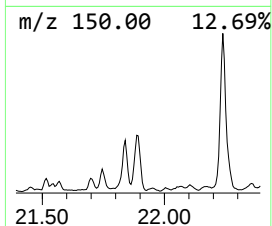
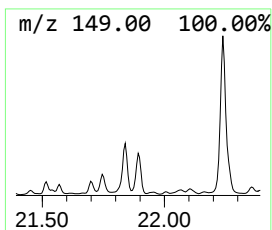
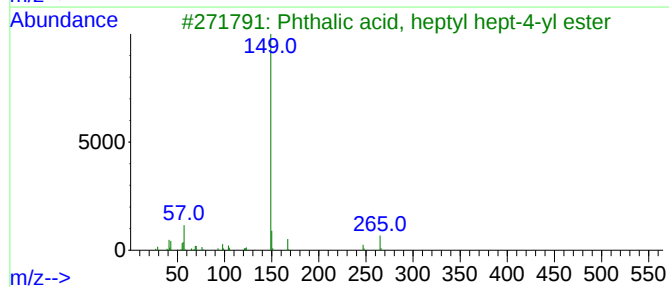
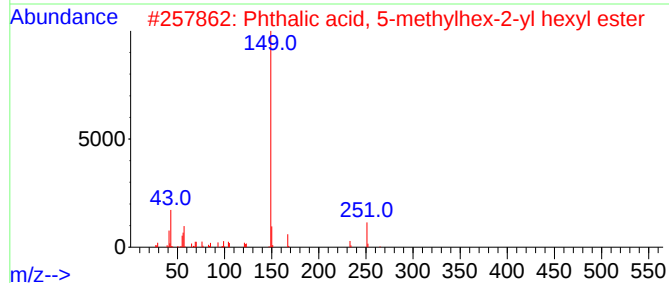
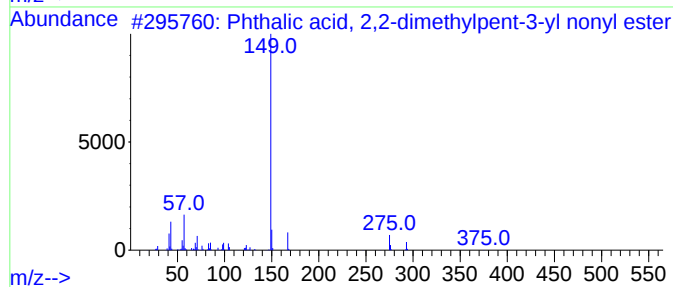
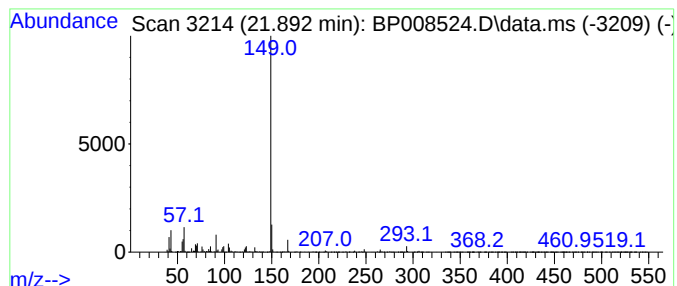
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Phthalic acid, 2,2-dimethyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.892	8.08 ng/ul	2190940	Chrysene-d12	21.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, 2,2-dimethylpent-...	390	C24H38O4	1000415-53-2	86
2			Phthalic acid, 5-methylhex-2-yl ...	348	C21H32O4	1000371-08-8	86
3			Phthalic acid, heptyl hept-4-yl ...	362	C22H34O4	1000356-78-8	86
4			Phthalic acid, dec-2-yl heptyl e...	404	C25H40O4	1000356-94-3	86
5			Phthalic acid, hexyl 6-methylhep...	362	C22H34O4	1000377-96-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
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Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

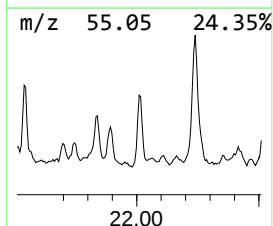
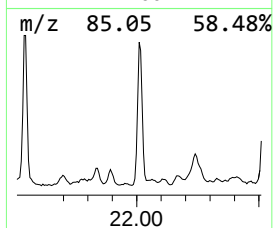
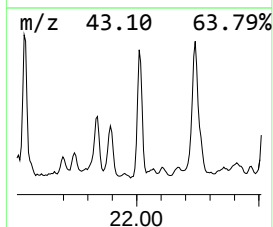
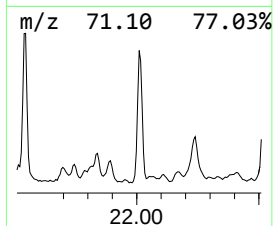
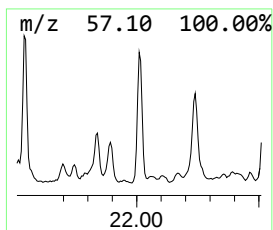
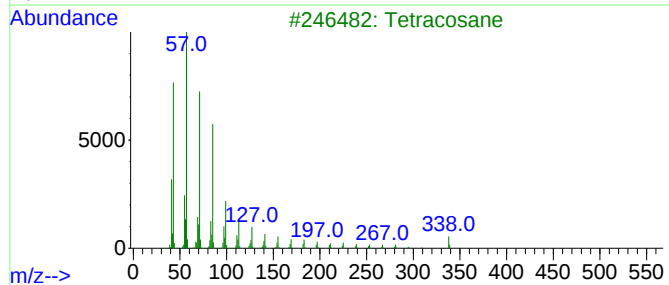
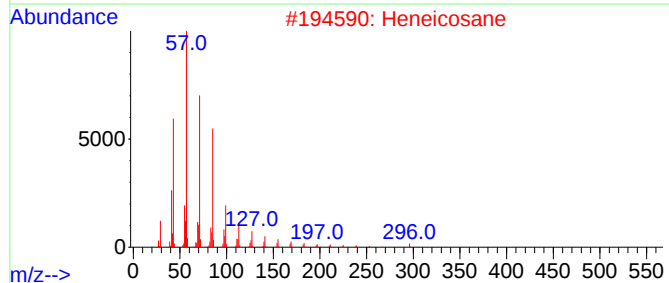
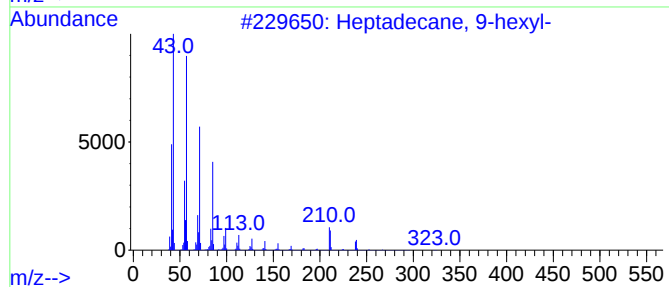
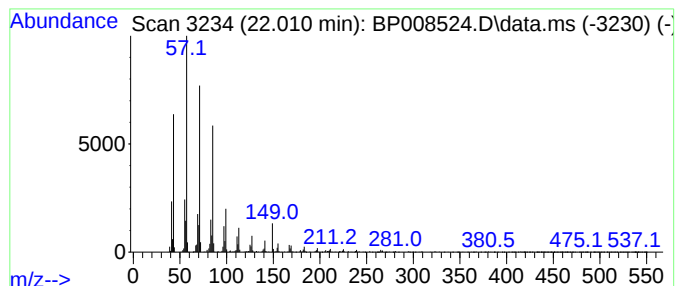
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.010	7.22 ng/ul	1956980	Chrysene-d12	21.363	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	91
2	Heneicosane	296	C21H44	000629-94-7	90
3	Tetracosane	338	C24H50	000646-31-1	90
4	Octacosane, 2-methyl-	408	C29H60	001560-98-1	90
5	Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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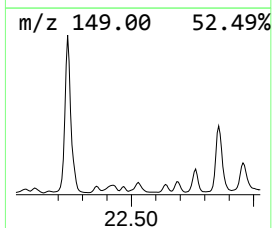
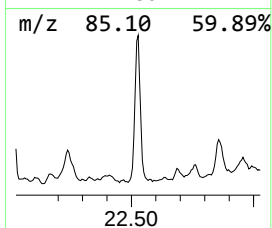
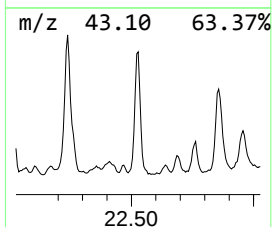
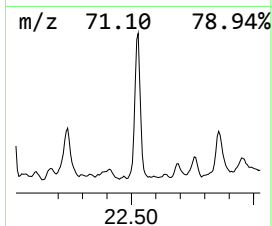
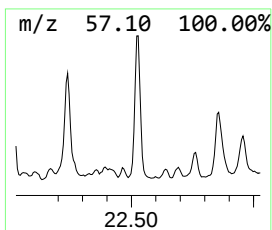
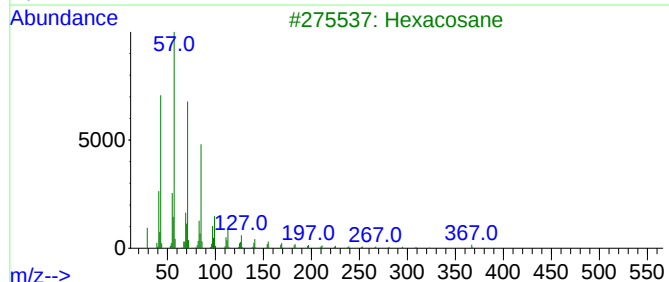
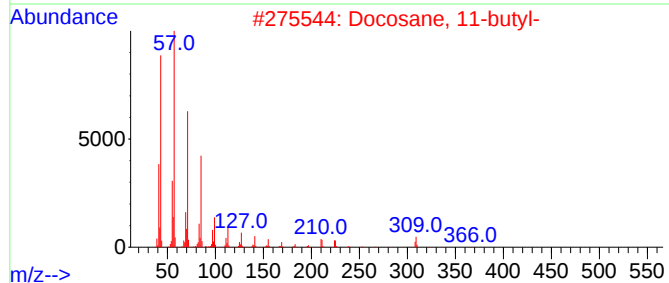
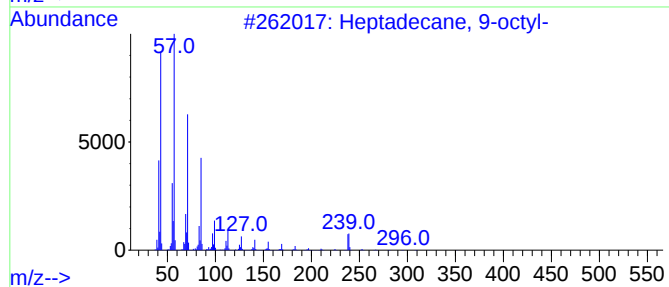
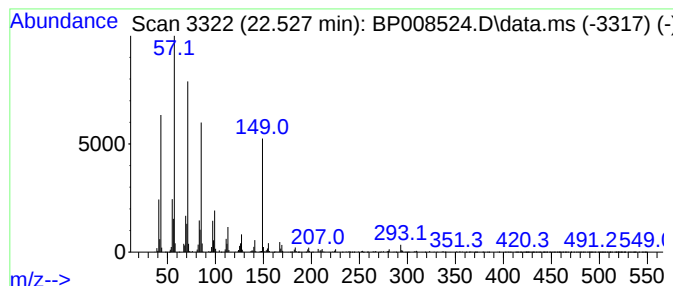
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.527	9.32 ng/ul	2526160	Chrysene-d12	21.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
2			Docosane, 11-butyl-	366	C26H54	013475-76-8	93
3			Hexacosane	366	C26H54	000630-01-3	93
4			Tetracosane	338	C24H50	000646-31-1	89
5			Eicosane, 9-octyl-	394	C28H58	013475-77-9	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGL53ME

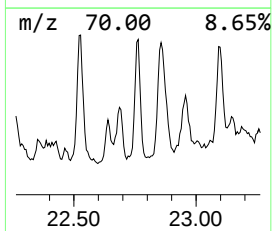
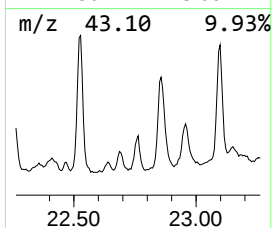
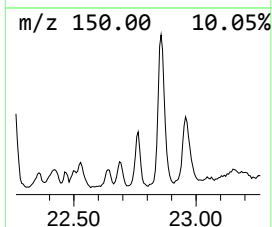
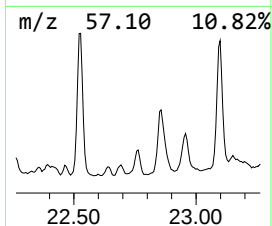
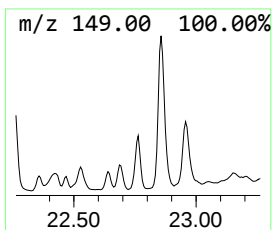
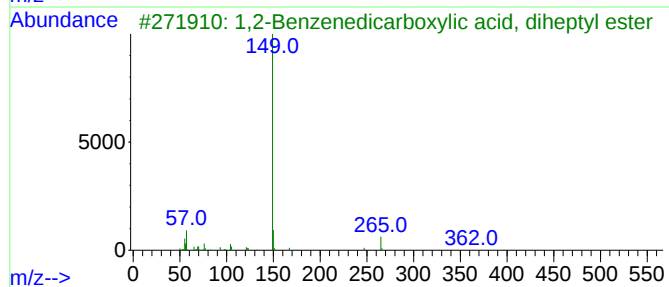
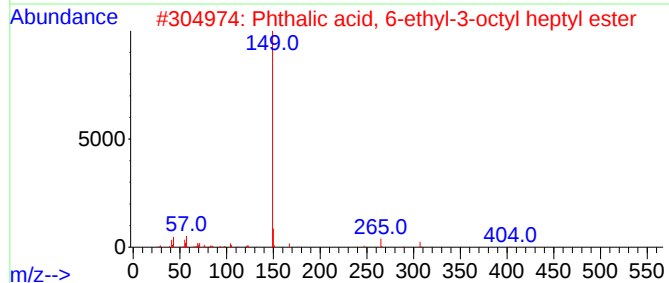
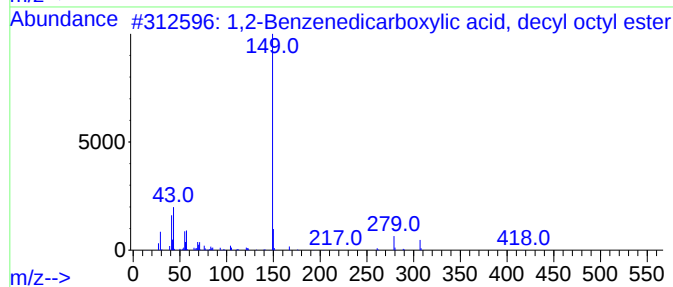
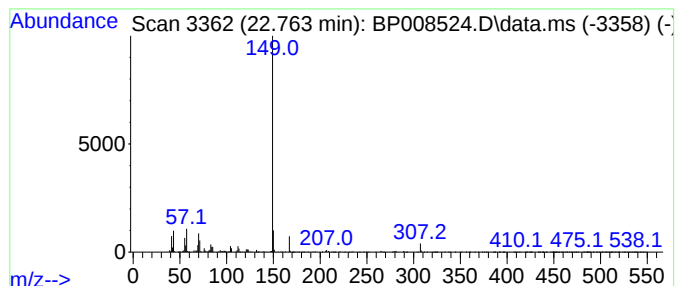
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 1,2-Benzenedicarboxylic aci... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.763	3.17 ng/ul	1096060	Perylene-d12	23.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Benzenedicarboxylic acid, de...	418	C26H42O4	000119-07-3	86
2			Phthalic acid, 6-ethyl-3-octyl h...	404	C25H40O4	1000315-17-7	86
3			1,2-Benzenedicarboxylic acid, di...	362	C22H34O4	003648-21-3	81
4			Phthalic acid, 2-ethylhexyl hexy...	362	C22H34O4	1000309-02-5	80
5			Phthalic acid, hept-3-yl octyl e...	376	C23H36O4	1000356-99-2	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
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Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BGL53ME

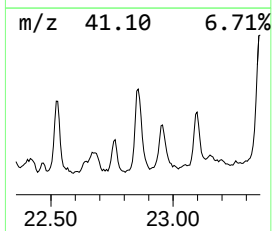
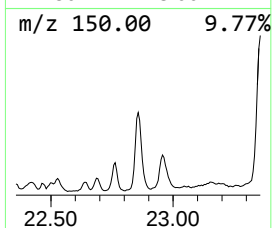
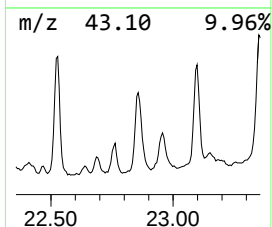
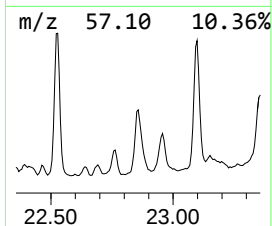
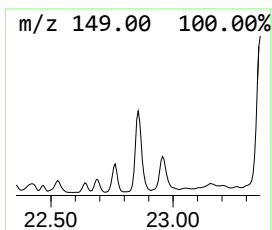
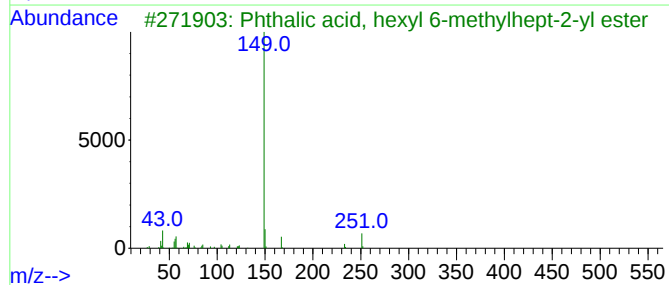
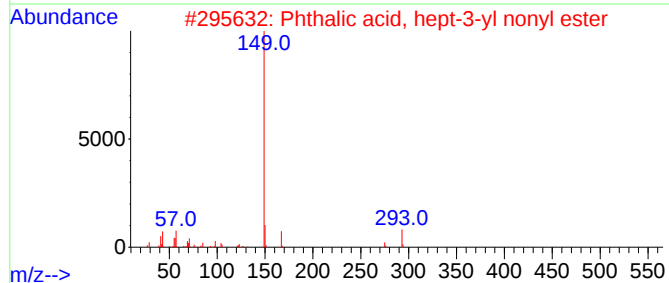
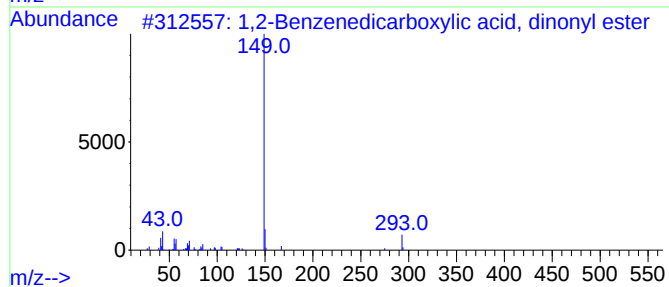
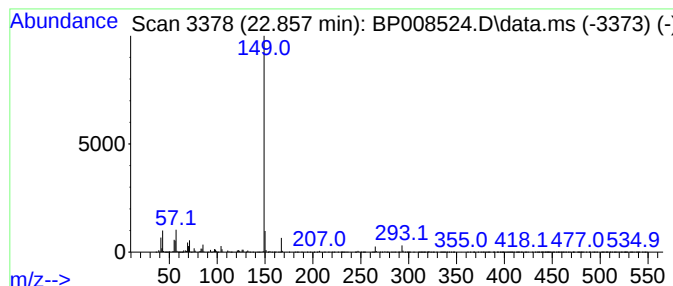
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 1,2-Benzenedicarboxylic aci... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.857	10.55 ng/ul	3650690	Perylene-d12	23.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	87
2			Phthalic acid, hept-3-yl nonyl e...	390	C24H38O4	1000356-99-3	86
3			Phthalic acid, hexyl 6-methylhep...	362	C22H34O4	1000377-96-0	86
4			Phthalic acid, hept-2-yl undecyl...	418	C26H42O4	1000356-97-8	80
5			Phthalic acid, heptyl 4-octyl ester	376	C23H36O4	1000314-84-1	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGL53ME

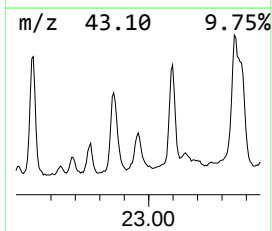
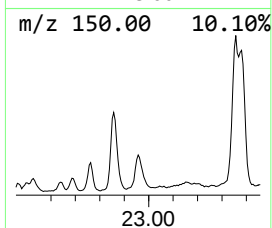
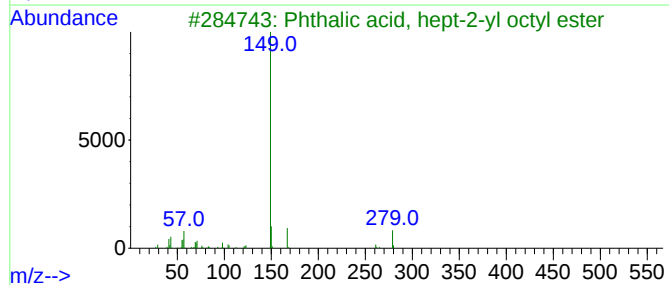
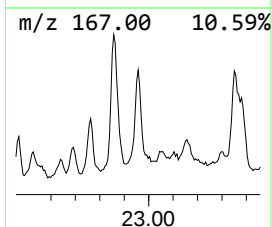
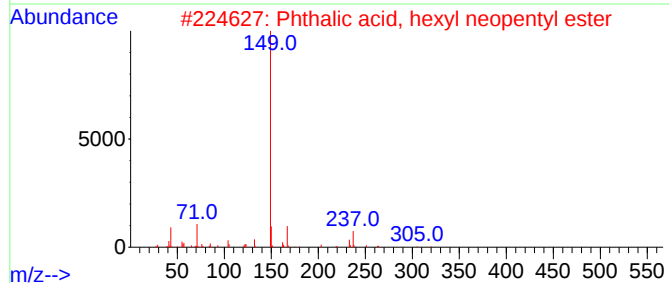
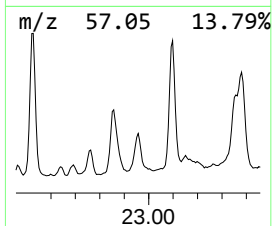
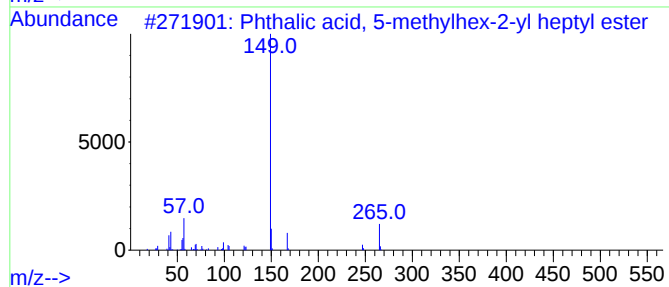
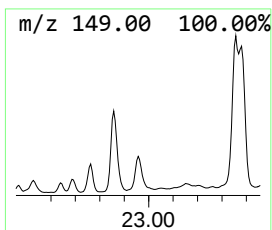
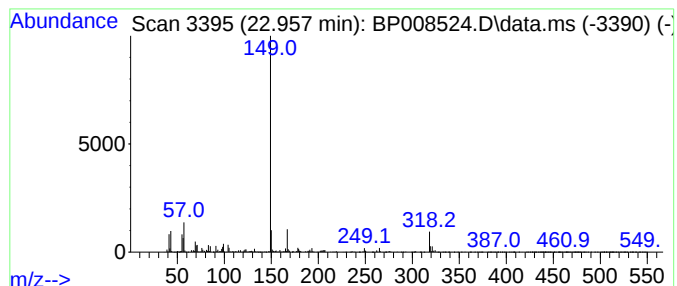
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Phthalic acid, 5-methylhex-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.957	6.82 ng/ul	2361860	Perylene-d12	23.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, 5-methylhex-2-yl ...	362	C22H34O4	1000371-08-7	80
2			Phthalic acid, hexyl neopentyl e...	320	C19H28O4	1000315-29-9	72
3			Phthalic acid, hept-2-yl octyl e...	376	C23H36O4	1000356-97-5	72
4			Phthalic acid, isopropyl octyl e...	320	C19H28O4	1000315-00-2	72
5			Di-sec-butyl phthalate	278	C16H22O4	1000373-65-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGL53ME

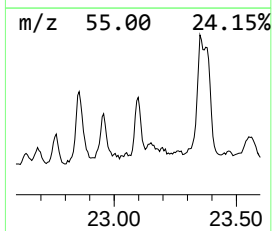
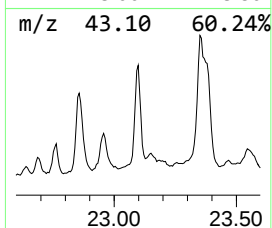
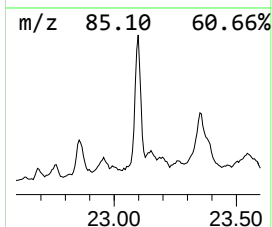
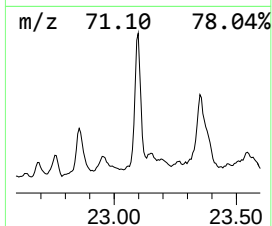
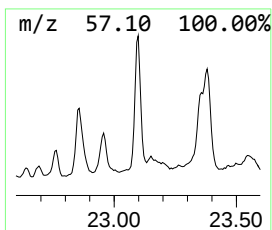
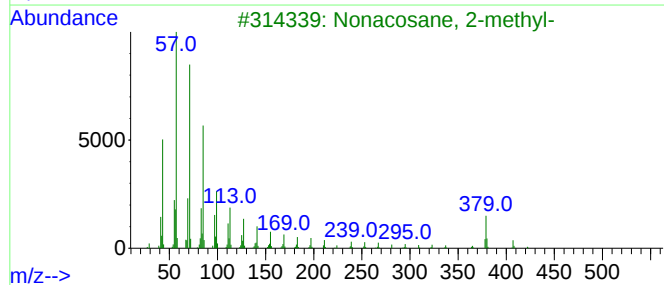
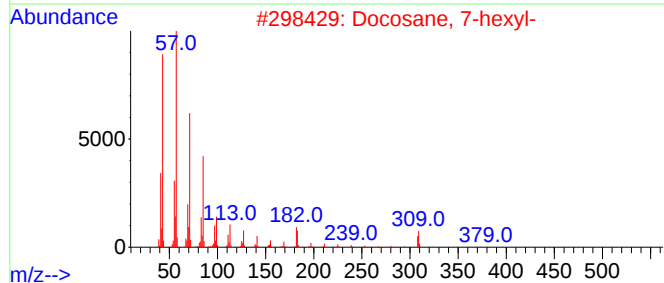
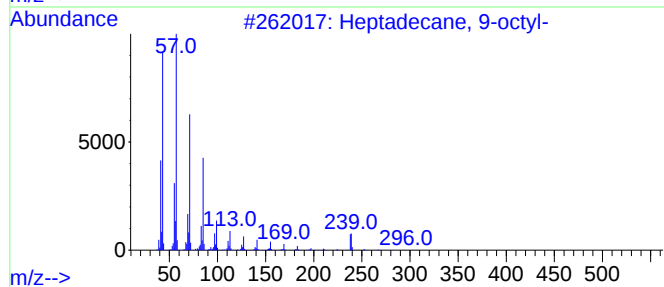
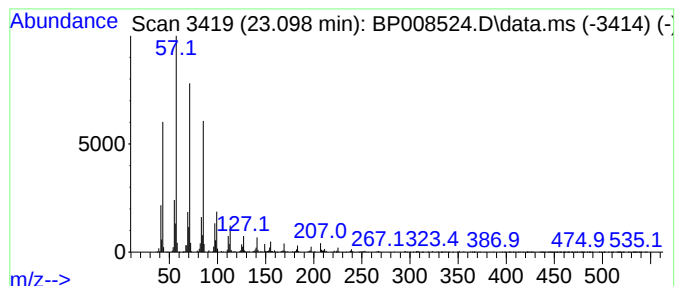
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.098	5.43 ng/ul	1879810	Perylene-d12	23.798

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2		Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
3		Nonacosane, 2-methyl-	422	C30H62	001560-75-4	93
4		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5		Octadecane, 2-methyl-	268	C19H40	001560-88-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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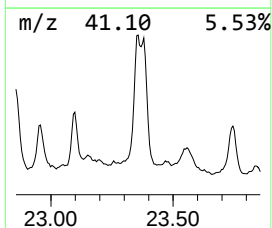
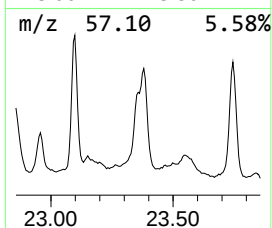
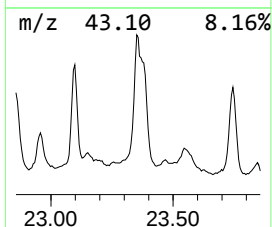
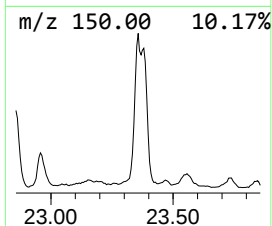
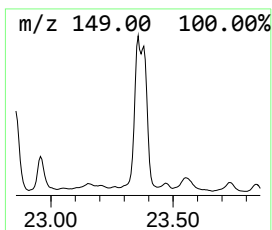
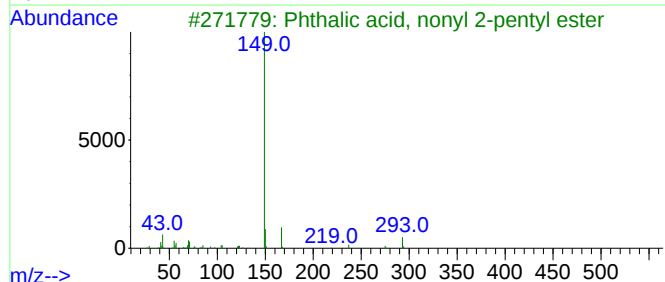
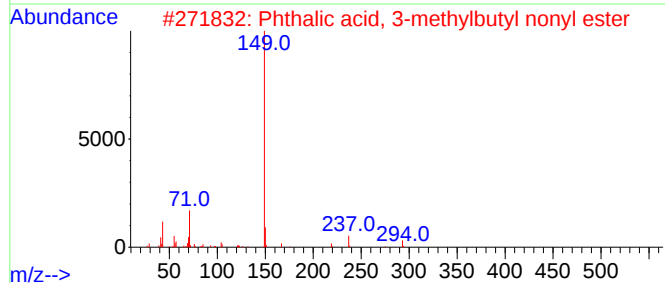
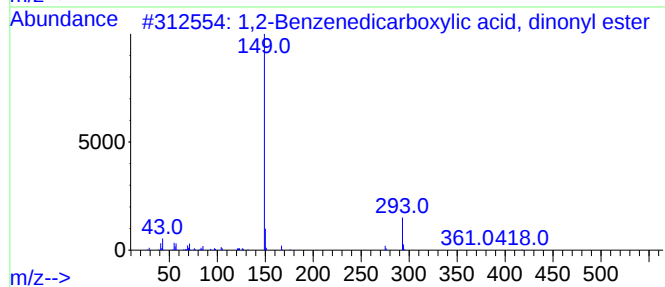
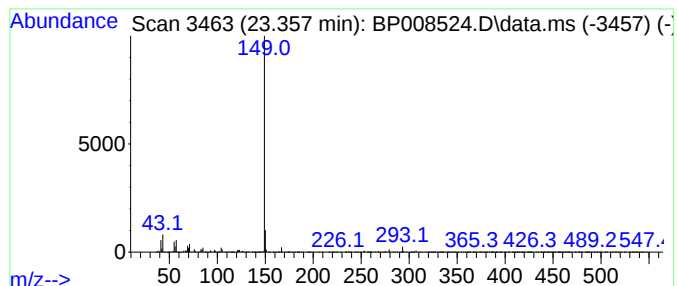
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Phthalic acid, 3-methylbuty... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.357	15.90 ng/ul	5504030	Perylene-d12	23.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	80
2			Phthalic acid, 3-methylbutyl non...	362	C22H34O4	1000356-81-0	80
3			Phthalic acid, nonyl 2-pentyl ester	362	C22H34O4	1000315-48-1	80
4			1,2-Benzenedicarboxylic acid, de...	418	C26H42O4	000119-07-3	72
5			Di-n-octyl phthalate	390	C24H38O4	000117-84-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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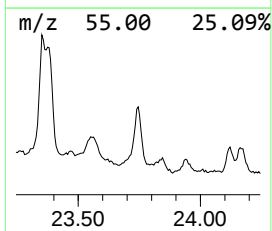
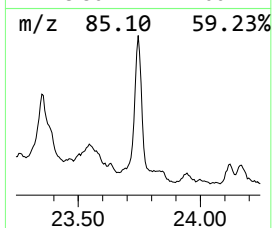
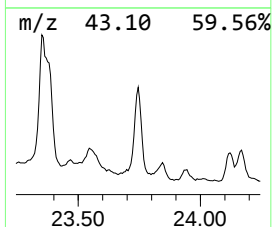
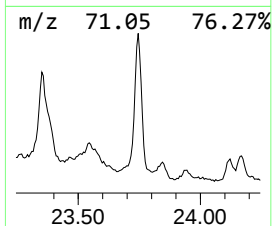
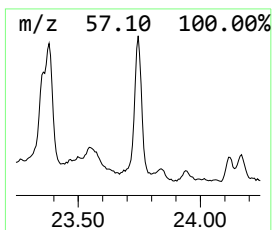
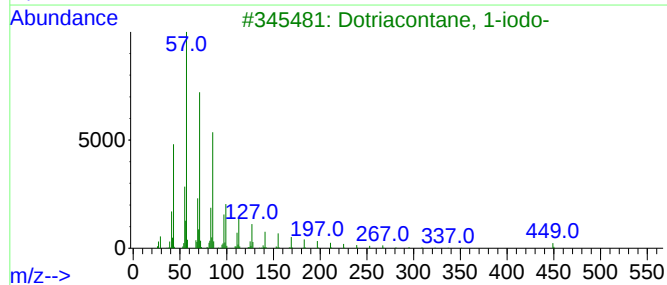
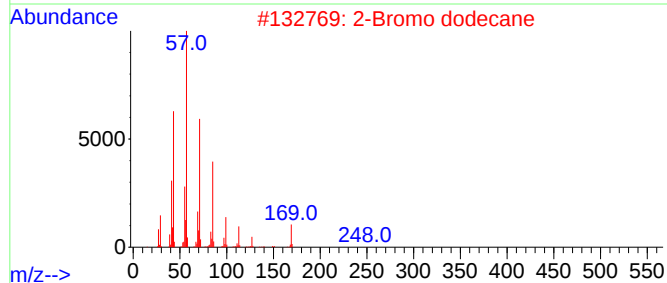
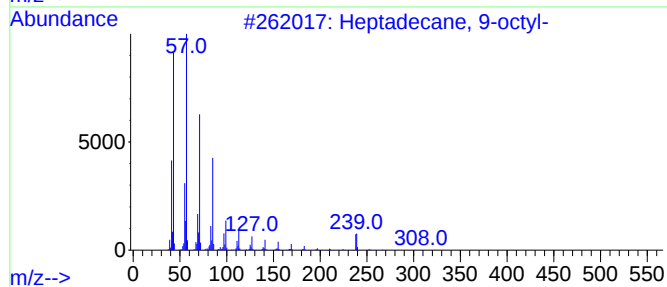
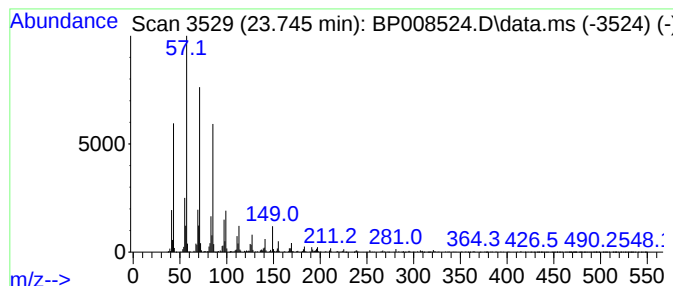
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.745	7.21 ng/ul	2495180	Perylene-d12	23.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
2			2-Bromo dodecane	248	C12H25Br	013187-99-0	91
3			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	90
4			Hexacosane	366	C26H54	000630-01-3	90
5			1,3-Propanediol, ethyl tetracosy...	440	C29H60O2	1000406-35-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGL53ME

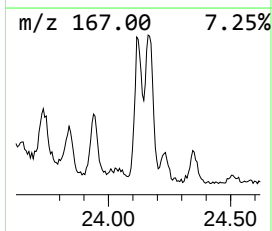
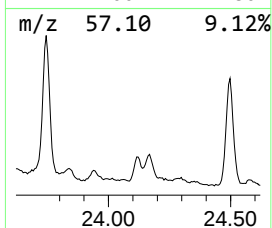
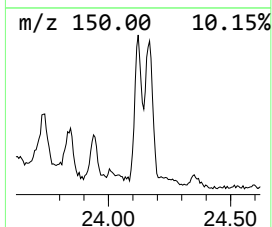
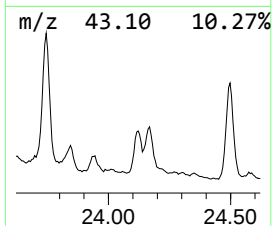
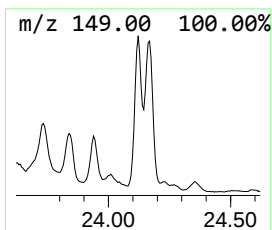
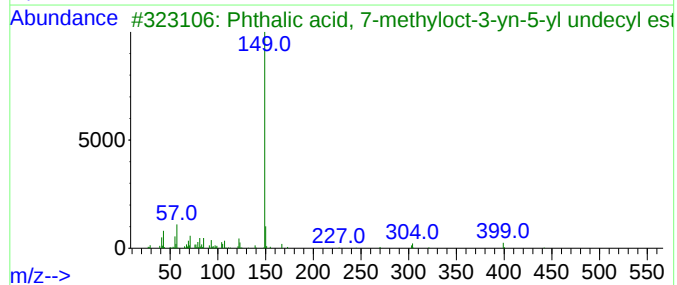
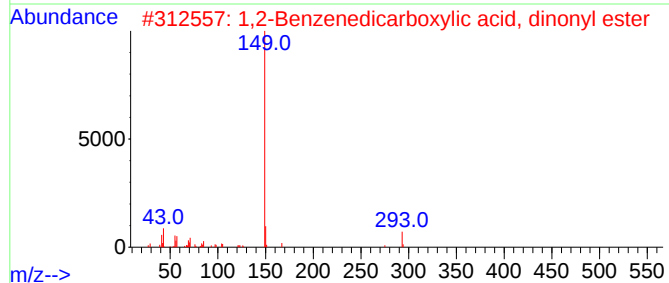
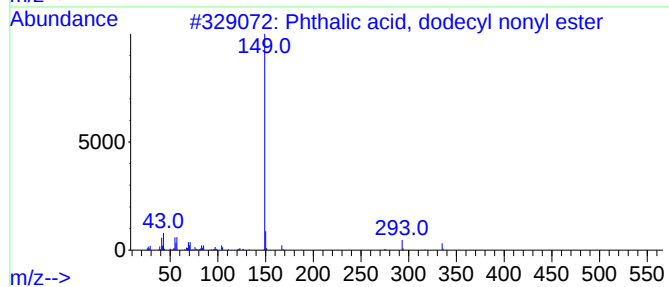
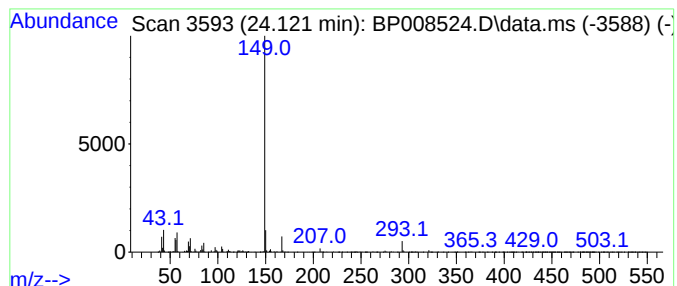
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Phthalic acid, dodecyl nonyl... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.121	3.68 ng/ul	1273950	Perylene-d12	23.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, dodecyl nonyl ester	460	C29H48O4	1000308-92-6	86
2			1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	86
3			Phthalic acid, 7-methyloct-3-yn-...	442	C28H42O4	1000315-43-3	80
4			Phthalic acid, hexyl 6-methylhep...	362	C22H34O4	1000377-96-0	80
5			Phthalic acid, isobutyl pentadec...	432	C27H44O4	1000309-07-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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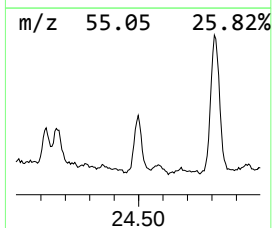
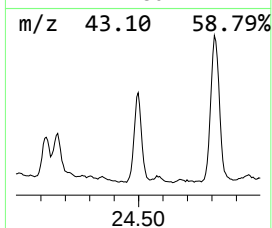
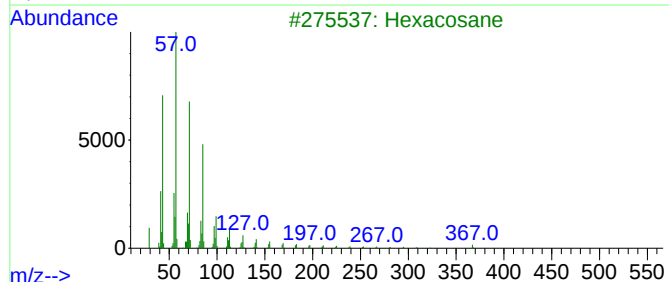
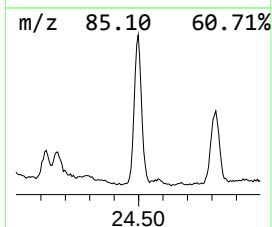
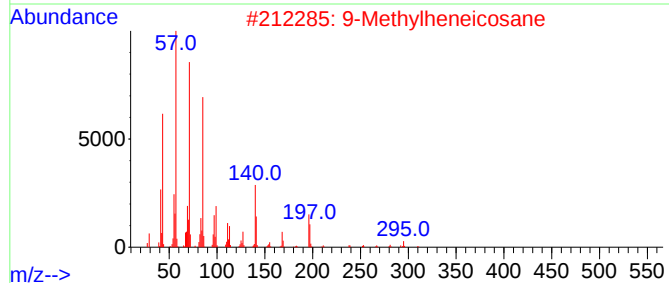
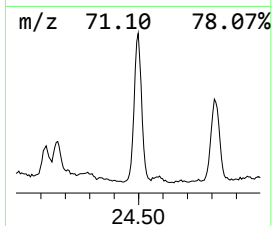
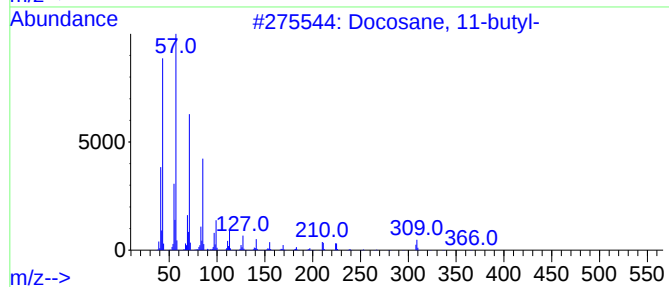
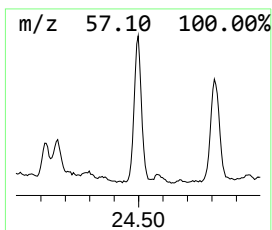
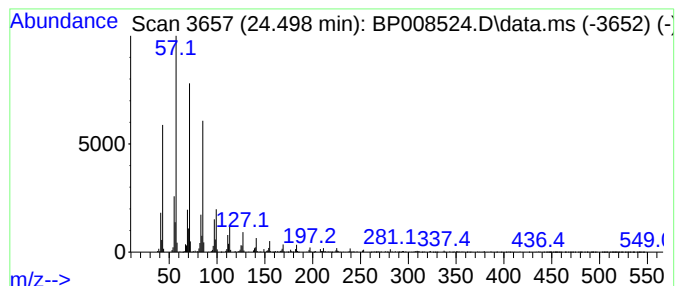
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.498	4.30 ng/ul	1489810	Perylene-d12	23.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosane, 11-butyl-	366	C26H54	013475-76-8	95
2			9-Methylheneicosane	310	C22H46	070475-51-3	94
3			Hexacosane	366	C26H54	000630-01-3	91
4			Heptacosane	380	C27H56	000593-49-7	91
5			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91



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Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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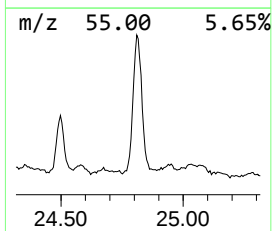
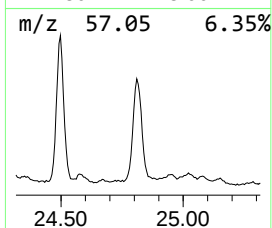
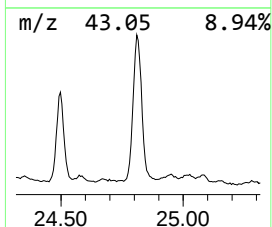
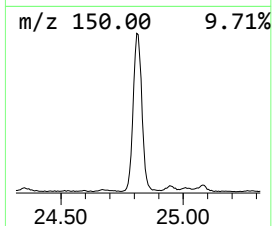
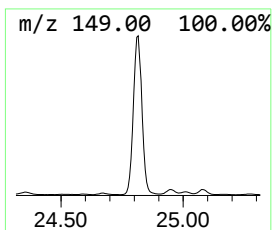
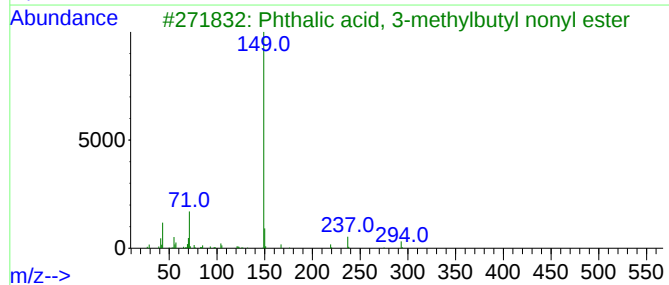
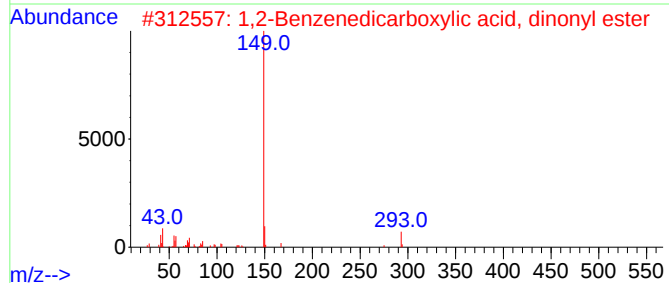
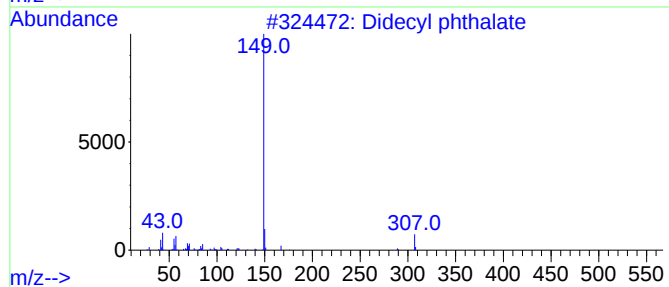
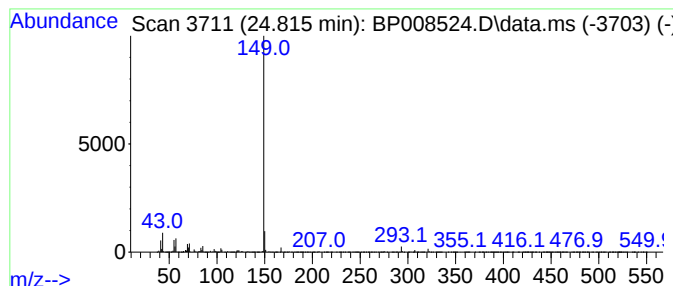
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 Didecyl phthalate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.816	17.97 ng/ul	6218940	Perylene-d12	23.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Didecyl phthalate	446	C28H46O4	000084-77-5	80
2			1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	80
3			Phthalic acid, 3-methylbutyl non...	362	C22H34O4	1000356-81-0	78
4			1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	78
5			Phthalic acid, cyclohexyl pentyl...	318	C19H26O4	1000315-33-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGL53ME

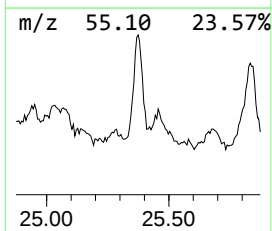
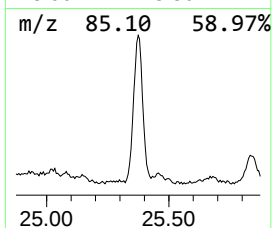
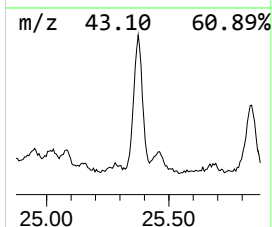
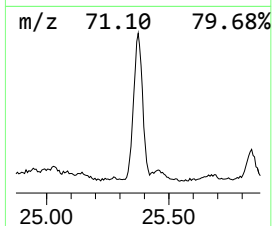
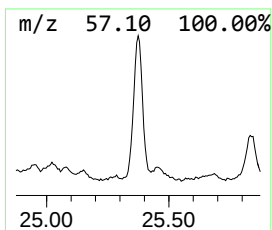
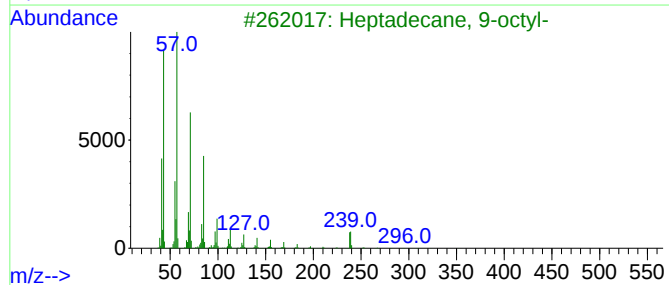
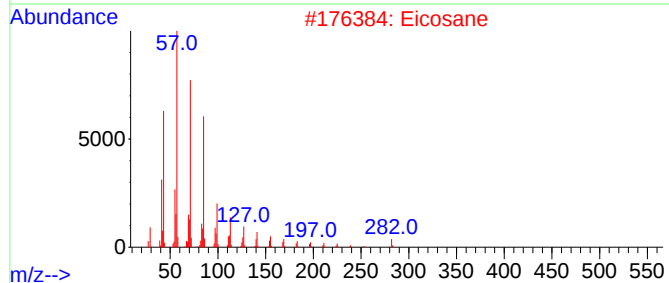
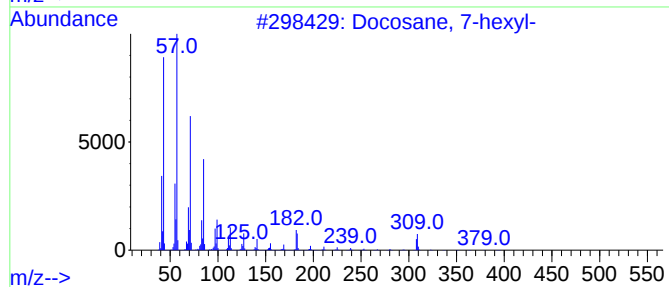
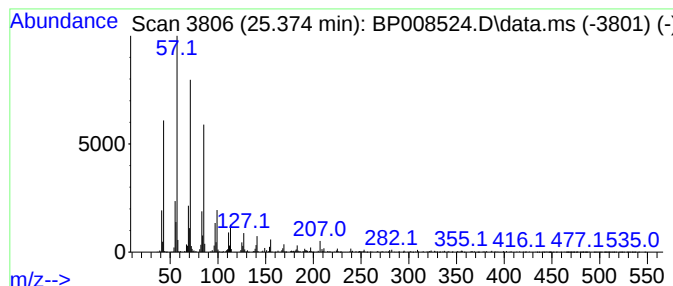
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.374	2.97 ng/ul	1028490	Perylene-d12	23.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosane, 7-hexyl-	394	C28H58	055373-86-9	95
2			Eicosane	282	C20H42	000112-95-8	94
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
4			Hentriacontane, 3-methyl-	451	C32H66	004981-99-1	93
5			Eicosane, 10-methyl-	296	C21H44	054833-23-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008524.D
Acq On : 23 Dec 2021 14:53
Operator : CG/JU
Sample : M5121-04ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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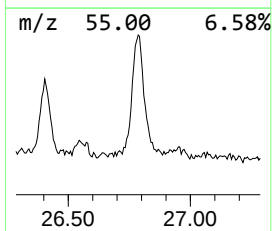
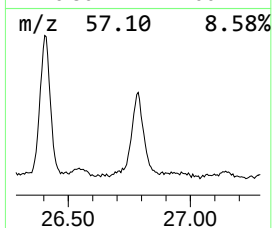
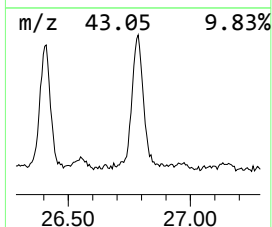
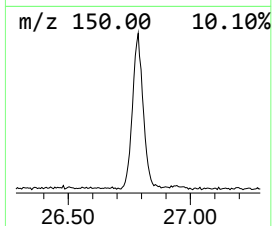
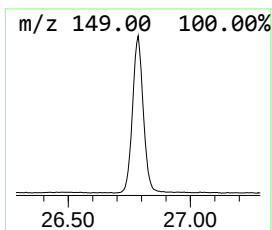
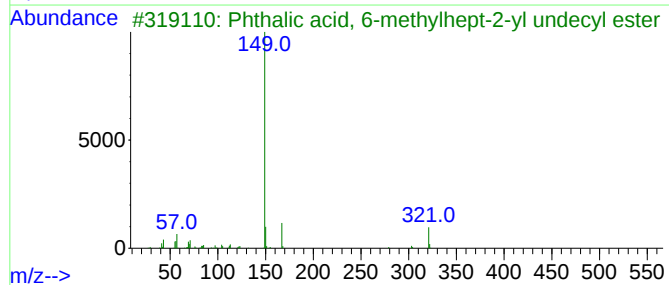
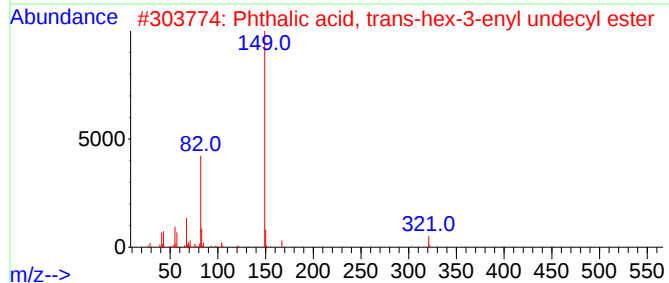
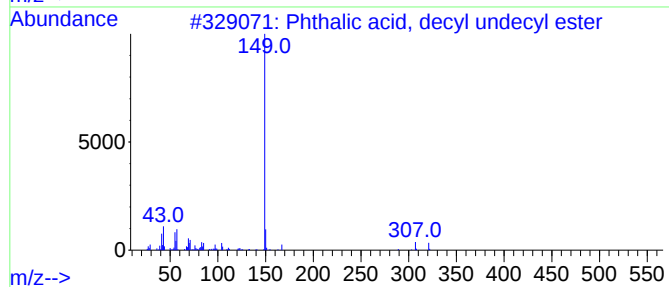
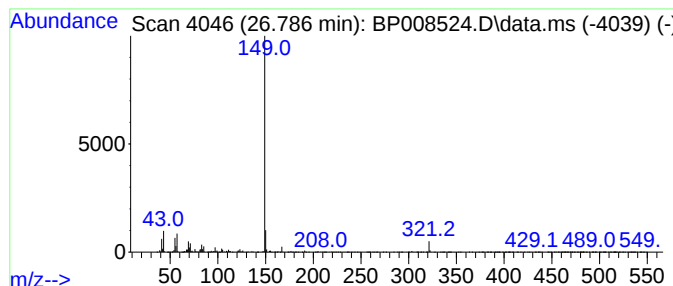
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 Phthalic acid, decyl undecy... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.786	7.06 ng/ul	2443940	Perylene-d12	23.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, decyl undecyl ester	460	C29H48O4	1000308-91-9	90
2			Phthalic acid, trans-hex-3-enyl ...	402	C25H38O4	1000360-48-9	80
3			Phthalic acid, 6-methylhept-2-yl...	432	C27H44O4	1000377-96-5	80
4			Phthalic acid, hex-3-yl undecyl ...	404	C25H40O4	1000356-96-2	80
5			Didecyl phthalate	446	C28H46O4	000084-77-5	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
 Data File : BP008524.D
 Acq On : 23 Dec 2021 14:53
 Operator : CG/JU
 Sample : M5121-04ME
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGL53ME

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: C...	3.023	15.7	ng/ul	2100680	1	7.905	2675710	20.0
Butane, 2-metho...	3.093	24.9	ng/ul	3326930	1	7.905	2675710	20.0
(DEL) Alkane: S...	16.298	2.1	ng/ul	611091	4	17.287	5944140	20.0
(DEL) Alkane: S...	20.163	2.6	ng/ul	697790	5	21.363	5421990	20.0
(DEL) Alkane: S...	20.645	3.7	ng/ul	997161	5	21.363	5421990	20.0
Phthalic acid, ...	20.869	7.8	ng/ul	2103320	5	21.363	5421990	20.0
Phthalic acid, ...	20.998	3.7	ng/ul	1012520	5	21.363	5421990	20.0
(DEL) Alkane: S...	21.098	5.9	ng/ul	1598270	5	21.363	5421990	20.0
(DEL) Alkane: S...	21.545	6.0	ng/ul	1624850	5	21.363	5421990	20.0
Phthalic acid, ...	21.698	2.0	ng/ul	552053	5	21.363	5421990	20.0
Phthalic acid, ...	21.745	3.3	ng/ul	889604	5	21.363	5421990	20.0
Phthalic acid, ...	21.839	7.1	ng/ul	1922280	5	21.363	5421990	20.0
Phthalic acid, ...	21.892	8.1	ng/ul	2190940	5	21.363	5421990	20.0
(DEL) Alkane: S...	22.010	7.2	ng/ul	1956980	5	21.363	5421990	20.0
(DEL) Alkane: S...	22.527	9.3	ng/ul	2526160	5	21.363	5421990	20.0
1,2-Benzenedica...	22.763	3.2	ng/ul	1096060	6	23.798	6922510	20.0
1,2-Benzenedica...	22.857	10.6	ng/ul	3650690	6	23.798	6922510	20.0
Phthalic acid, ...	22.957	6.8	ng/ul	2361860	6	23.798	6922510	20.0
(DEL) Alkane: S...	23.098	5.4	ng/ul	1879810	6	23.798	6922510	20.0
Phthalic acid, ...	23.357	15.9	ng/ul	5504030	6	23.798	6922510	20.0
(DEL) Alkane: S...	23.745	7.2	ng/ul	2495180	6	23.798	6922510	20.0
Phthalic acid, ...	24.121	3.7	ng/ul	1273950	6	23.798	6922510	20.0
(DEL) Alkane: S...	24.498	4.3	ng/ul	1489810	6	23.798	6922510	20.0
Didecyl phthalate	24.816	18.0	ng/ul	6218940	6	23.798	6922510	20.0
(DEL) Alkane: S...	25.374	3.0	ng/ul	1028490	6	23.798	6922510	20.0
Phthalic acid, ...	26.786	7.1	ng/ul	2443940	6	23.798	6922510	20.0