

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081524\
 Data File : BP021529.D
 Acq On : 16 Aug 2024 01:51
 Operator : RC/JU
 Sample : P3598-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JC-03-081324-A

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % 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\8270E-BP081324.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP021529.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.028	3	7	27	rVB	3177258	5590403	42.02%	5.083%
2	3.552	91	96	105	rVB	197356	303634	2.28%	0.276%
3	4.940	327	332	342	rBV	316189	489386	3.68%	0.445%
4	5.405	404	411	420	rBV	5552998	8490581	63.81%	7.720%
5	5.481	420	424	436	rVB2	82253	220630	1.66%	0.201%
6	5.822	474	482	493	rBV3	149982	435302	3.27%	0.396%
7	6.969	670	677	686	rBV	5345646	8900835	66.90%	8.093%
8	7.399	744	750	755	rBV	97017	151411	1.14%	0.138%
9	7.469	755	762	773	rVB4	130738	310821	2.34%	0.283%
10	7.811	813	820	827	rBV	2036152	3280751	24.66%	2.983%
11	8.187	878	884	890	rVB	77463	134837	1.01%	0.123%
12	8.346	905	911	921	rBV	125091	228626	1.72%	0.208%
13	8.952	998	1014	1023	rBV	3292246	5789513	43.51%	5.264%
14	9.046	1025	1030	1041	rVB	95598	207456	1.56%	0.189%
15	9.510	1098	1109	1116	rBV6	123775	264949	1.99%	0.241%
16	9.793	1152	1157	1165	rBV7	68945	157084	1.18%	0.143%
17	10.593	1285	1293	1299	rBV	2676620	5049407	37.95%	4.591%
18	10.646	1299	1302	1309	rVB	576703	882476	6.63%	0.802%
19	10.775	1320	1324	1329	rVB	148383	224776	1.69%	0.204%
20	11.122	1376	1383	1387	rBV4	82375	207911	1.56%	0.189%
21	11.328	1408	1418	1424	rBV2	391584	801687	6.03%	0.729%
22	11.452	1433	1439	1444	rBV4	92425	175968	1.32%	0.160%
23	11.528	1447	1452	1459	rVB7	88783	217621	1.64%	0.198%
24	11.640	1465	1471	1476	rBV	357652	658375	4.95%	0.599%
25	12.034	1533	1538	1546	rBV2	237562	481047	3.62%	0.437%
26	12.257	1568	1576	1583	rVB3	332908	726410	5.46%	0.661%
27	12.469	1608	1612	1617	rBV2	162074	257411	1.93%	0.234%
28	12.734	1654	1657	1662	rVB	128239	171313	1.29%	0.156%
29	12.951	1688	1694	1696	rBV4	86499	175361	1.32%	0.159%
30	12.999	1697	1702	1706	rVV	498079	772112	5.80%	0.702%
31	13.063	1706	1713	1719	rVB	8562716	12350966	92.83%	11.230%
32	13.275	1744	1749	1758	rVB4	256348	492889	3.70%	0.448%
33	13.375	1763	1766	1771	rBV5	100444	160725	1.21%	0.146%
34	13.604	1801	1805	1807	rBV2	123184	186362	1.40%	0.169%
35	13.763	1828	1832	1836	rBV	182345	284216	2.14%	0.258%
36	13.816	1837	1841	1845	rVB3	186573	263929	1.98%	0.240%

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Stop Thrs : 0

Filtering: 5

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Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M

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37	13.875	1848	1851	1856	rVV4	120704	162125	1.22%	0.147%
38	13.951	1859	1864	1868	rVV2	479578	783179	5.89%	0.712%
39	13.998	1868	1872	1876	rVB3	163513	257117	1.93%	0.234%
40	14.157	1896	1899	1905	rVB4	145489	223761	1.68%	0.203%
41	14.298	1918	1923	1929	rVB6	122521	227606	1.71%	0.207%
42	14.363	1930	1934	1939	rBV3	287783	425800	3.20%	0.387%
43	14.451	1941	1949	1954	rVV	4000542	6312808	47.45%	5.740%
44	14.510	1956	1959	1966	rVB2	218131	368726	2.77%	0.335%
45	14.663	1980	1985	1986	rBV2	206985	290876	2.19%	0.264%
46	14.934	2027	2031	2038	rVB3	131624	217993	1.64%	0.198%
47	14.998	2038	2042	2048	rVV2	260198	374247	2.81%	0.340%
48	15.069	2051	2054	2060	rVV3	143015	276345	2.08%	0.251%
49	15.128	2062	2064	2069	rVB	129210	143752	1.08%	0.131%
50	15.263	2079	2087	2089	rBV6	96509	248787	1.87%	0.226%
51	15.340	2096	2100	2103	rBV	205141	254063	1.91%	0.231%
52	15.504	2125	2128	2134	rVB	138537	204643	1.54%	0.186%
53	15.751	2163	2170	2175	rVB2	280541	541298	4.07%	0.492%
54	15.963	2200	2206	2214	rBV	6564285	8796164	66.11%	7.998%
55	16.210	2244	2248	2250	rBV3	195049	299379	2.25%	0.272%
56	16.240	2250	2253	2262	rVB	401574	766138	5.76%	0.697%
57	17.034	2385	2388	2391	rBV	141960	162423	1.22%	0.148%
58	17.257	2420	2426	2430	rBV	4043871	6268257	47.11%	5.700%
59	17.298	2430	2433	2439	rVB	460408	593572	4.46%	0.540%
60	18.204	2583	2587	2594	rVB2	1083300	1532005	11.51%	1.393%
61	19.386	2784	2788	2791	rBV	484834	739212	5.56%	0.672%
62	19.533	2810	2813	2817	rBV4	594345	963100	7.24%	0.876%
63	19.975	2883	2888	2893	rVB	10860845	13305303	100.00%	12.098%
64	21.380	3125	3127	3132	rBV3	700449	1188365	8.93%	1.081%
65	21.704	3178	3182	3188	rVB2	2959094	4554155	34.23%	4.141%

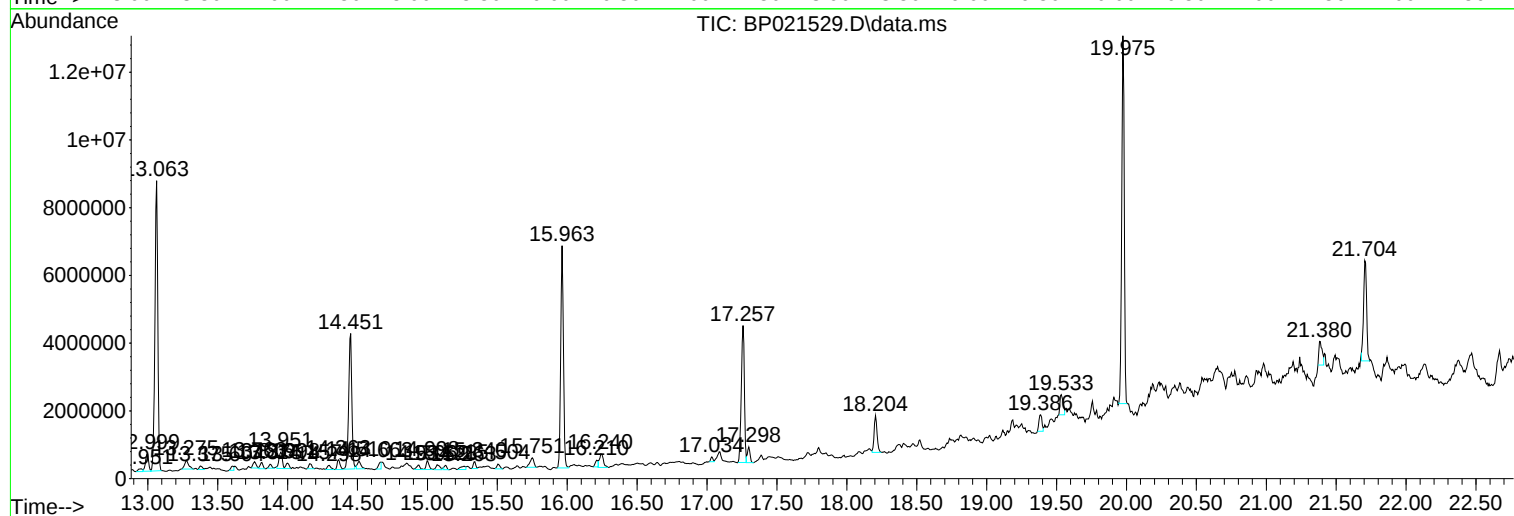
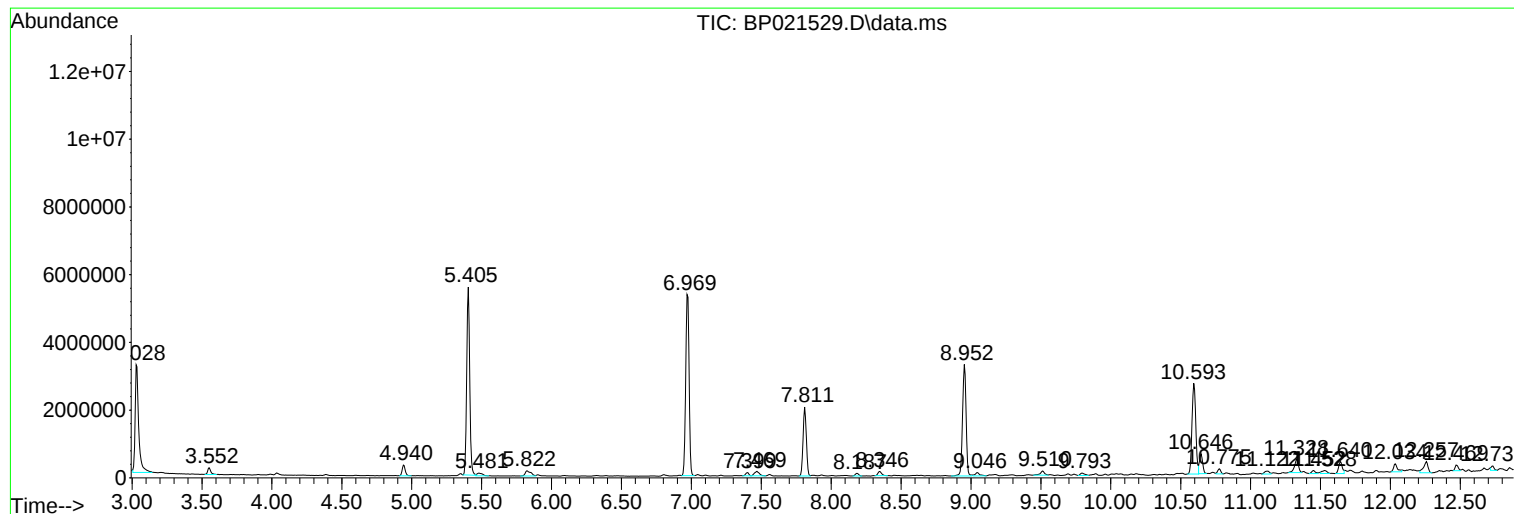
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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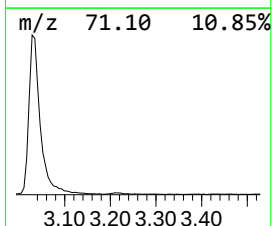
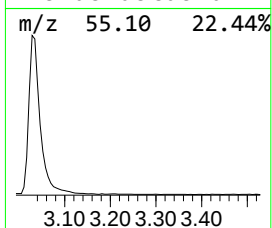
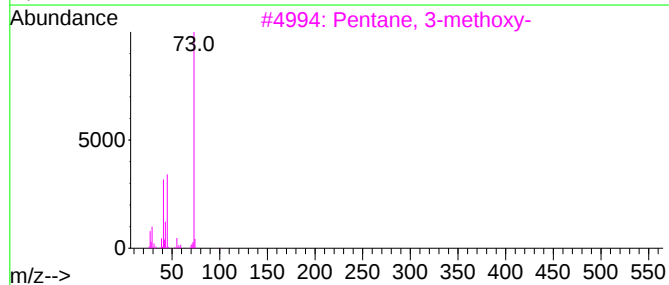
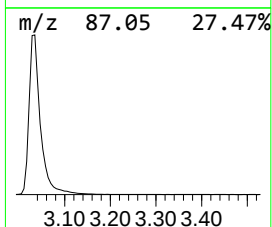
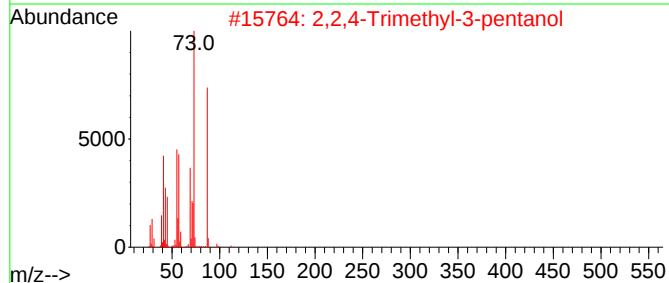
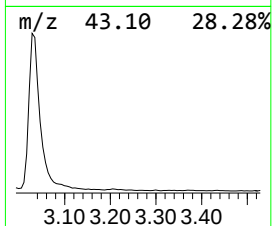
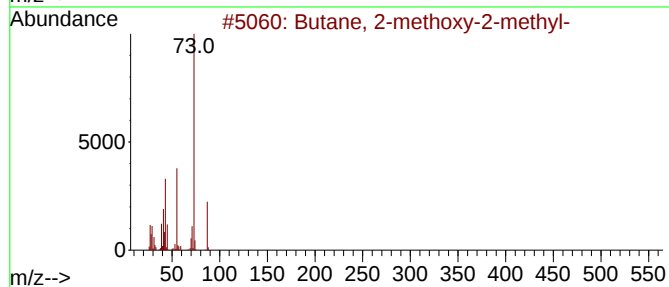
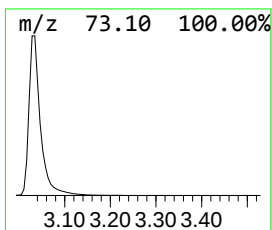
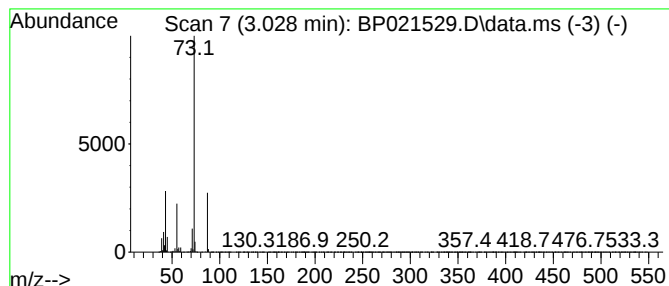
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.028	34.08 ng	5590400	1,4-Dichlorobenzene-d4	7.811

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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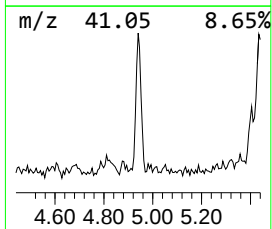
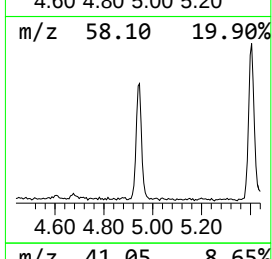
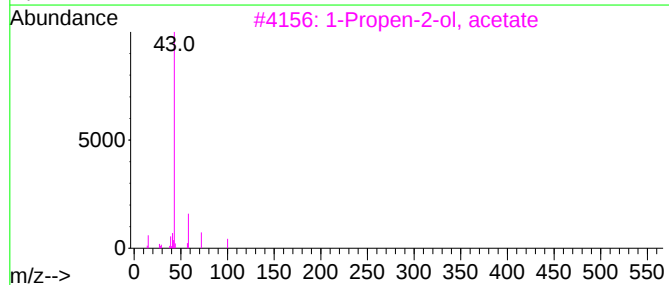
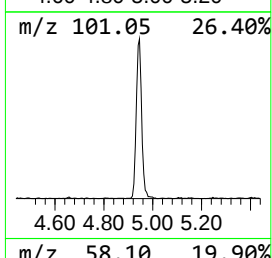
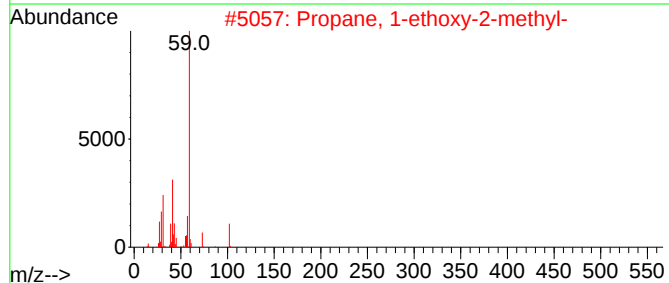
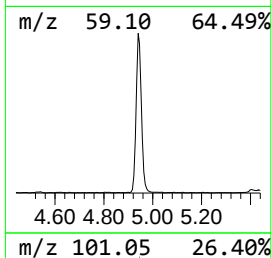
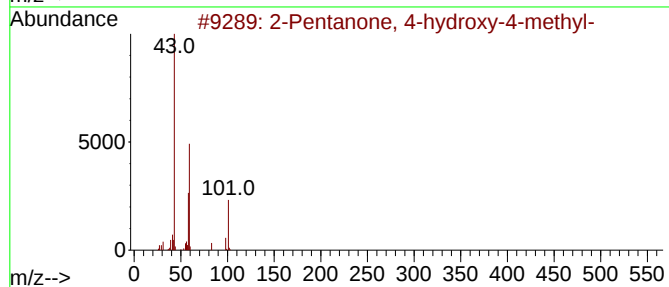
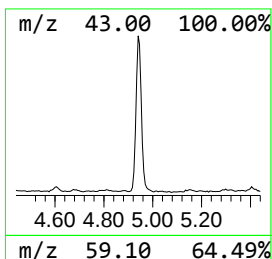
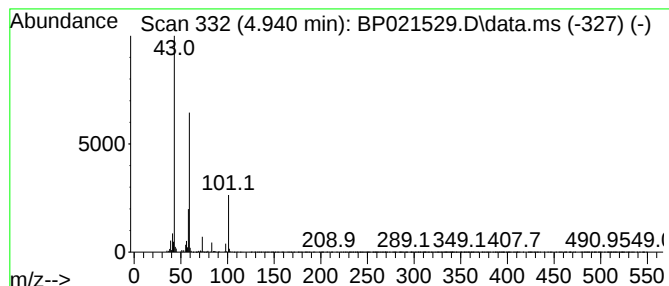
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.940	2.98 ng	489386	1,4-Dichlorobenzene-d4	7.811

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	30
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	10
5			1-Ethylpropyl acetate	130	C7H14O2	000620-11-1	9



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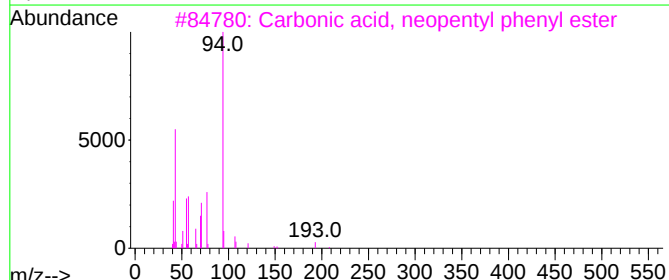
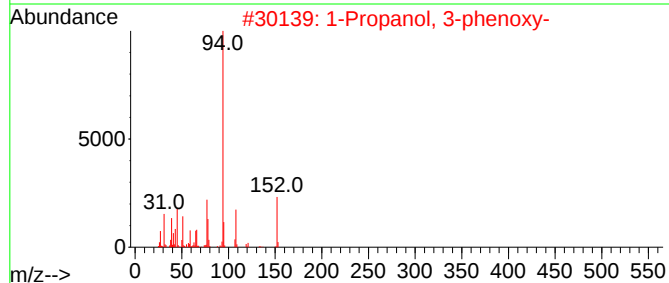
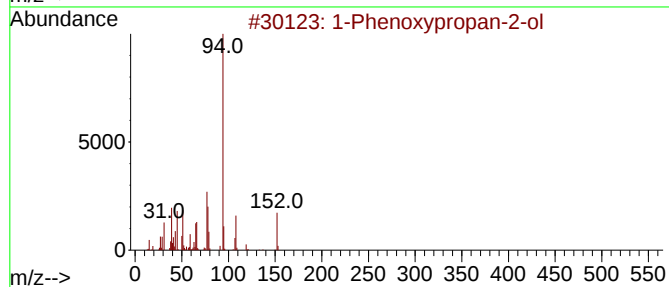
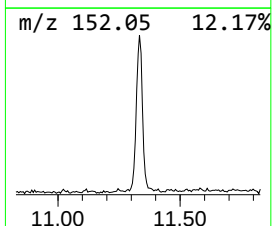
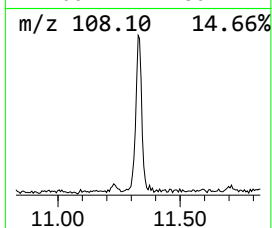
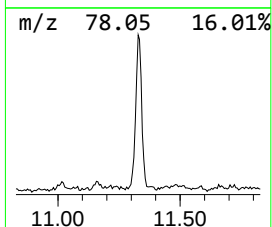
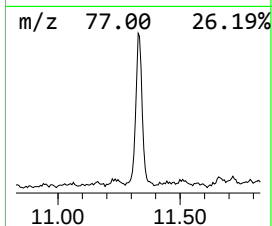
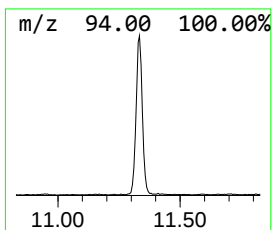
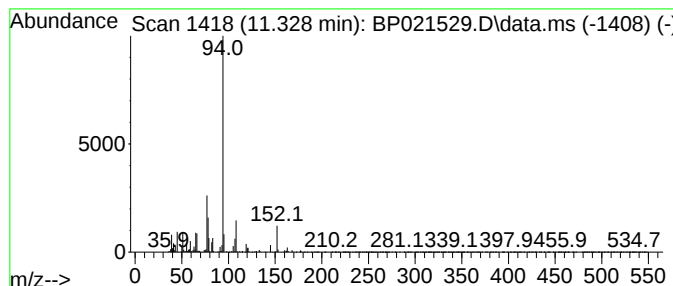
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.328	3.18 ng	801687	Naphthalene-d8	10.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
2			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	83
3			Carbonic acid, neopentyl phenyl ...	208	C12H16O3	013183-19-2	59
4			1-Propanol, 2-phenoxy-	152	C9H12O2	004169-04-4	58
5			Benzene, (2-methylpropoxy)-	150	C10H14O	001126-75-6	50



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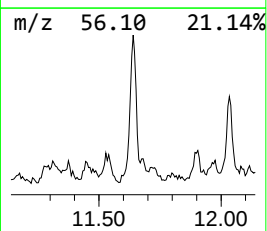
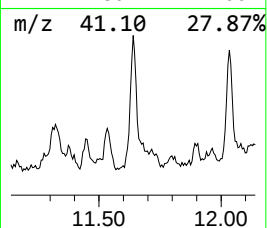
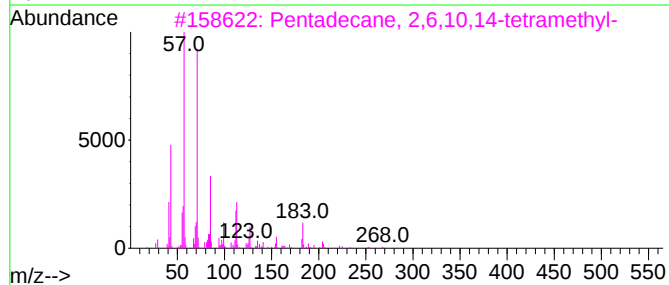
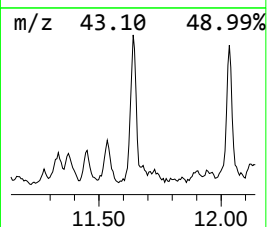
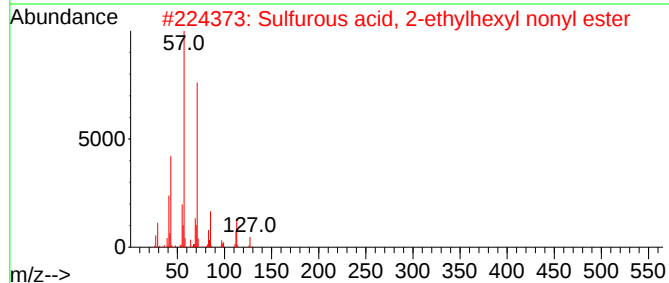
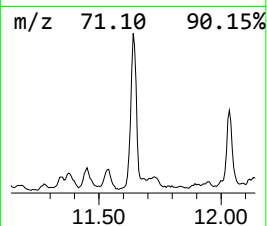
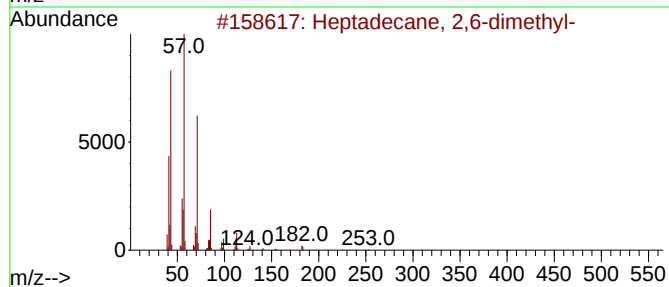
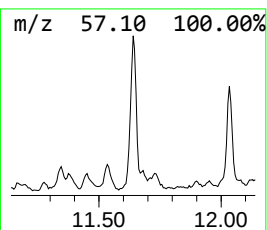
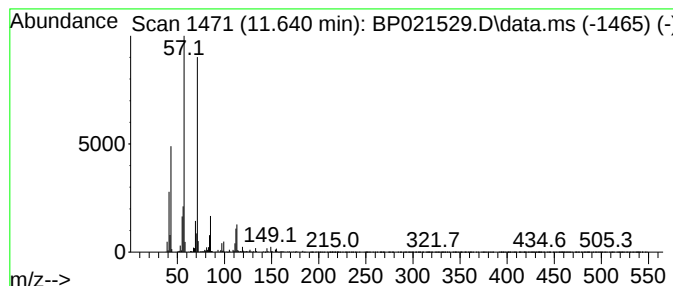
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TIC Library : C:\Database\NIST20.L
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Peak Number 5 Heptadecane, 2,6-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.640	2.61 ng	658375	Naphthalene-d8	10.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	83
2			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	80
3			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	80
4			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	72
5			Nonane, 4-methyl-5-propyl-	184	C13H28	062185-55-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081524\
Data File : BP021529.D
Acq On : 16 Aug 2024 01:51
Operator : RC/JU
Sample : P3598-01
Misc :
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Instrument :
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ClientSampleId :
JC-03-081324-A

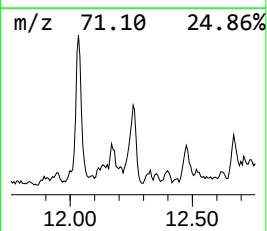
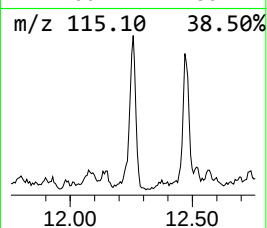
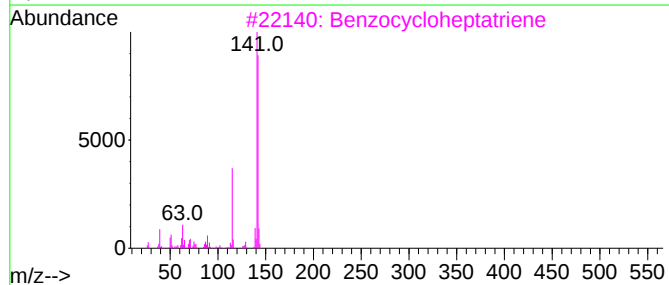
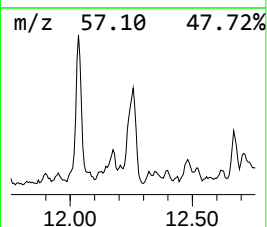
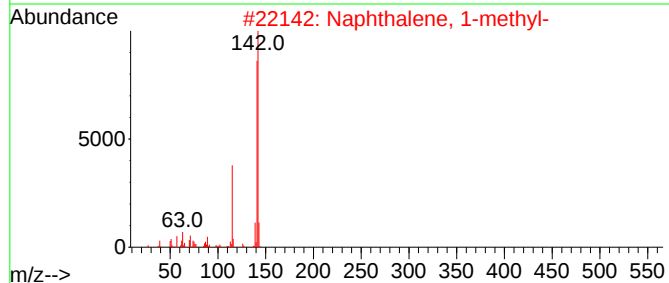
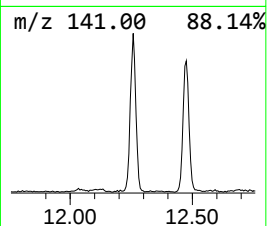
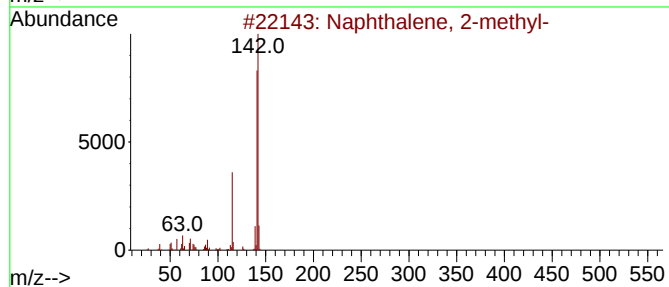
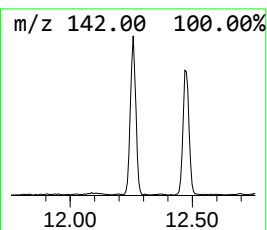
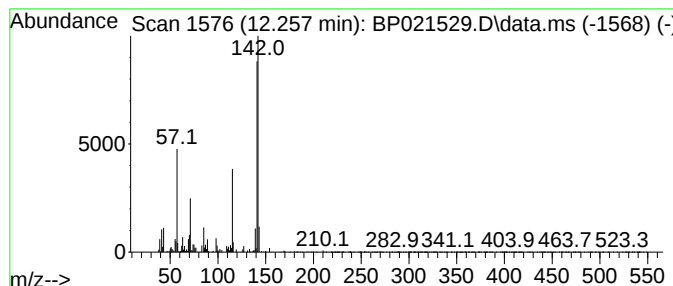
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Naphthalene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.257	2.88 ng	726410	Naphthalene-d8	10.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			Benzocycloheptatriene	142	C11H10	000264-09-5	87
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	87
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081524\
Data File : BP021529.D
Acq On : 16 Aug 2024 01:51
Operator : RC/JU
Sample : P3598-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
JC-03-081324-A

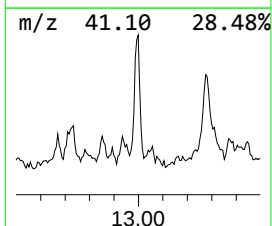
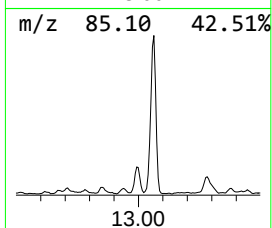
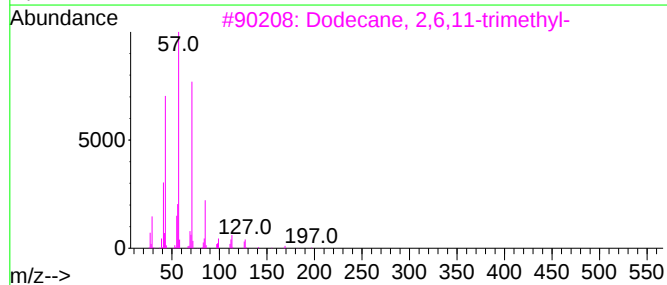
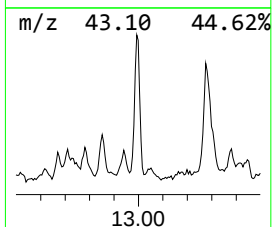
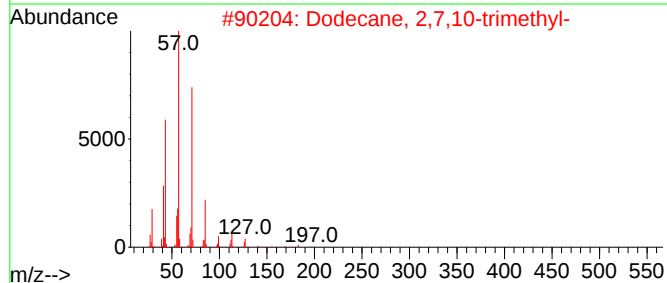
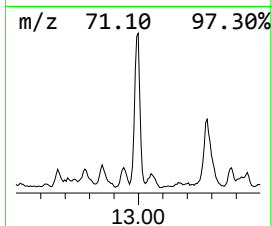
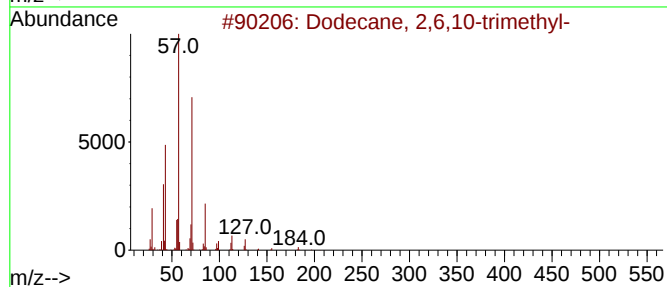
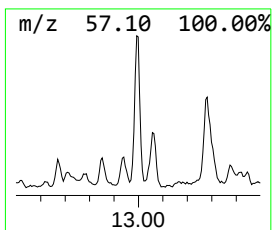
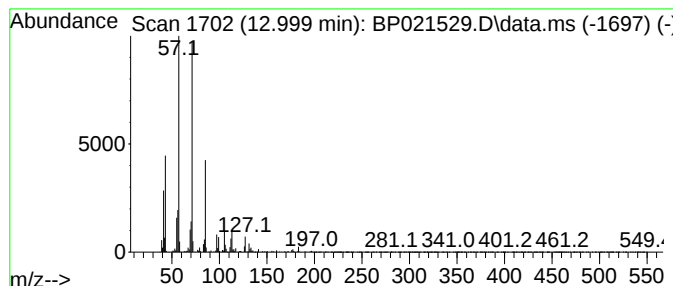
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Dodecane, 2,6,10-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.999	2.45 ng	772112	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	72
2			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72
3			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	64
4			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	62
5			Heptadecane	240	C17H36	000629-78-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081524\
Data File : BP021529.D
Acq On : 16 Aug 2024 01:51
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Sample : P3598-01
Misc :
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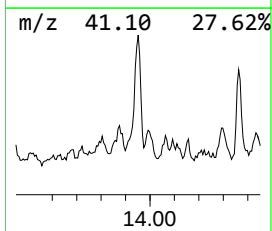
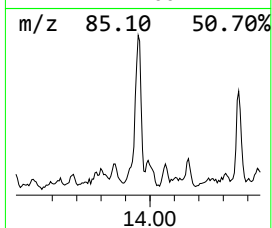
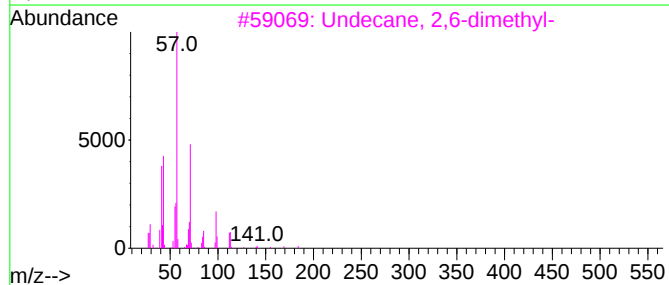
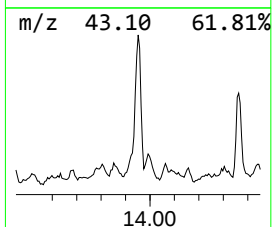
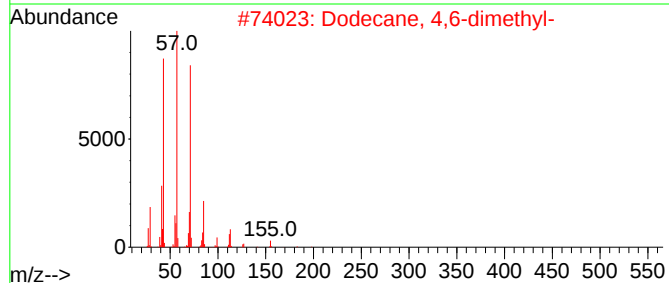
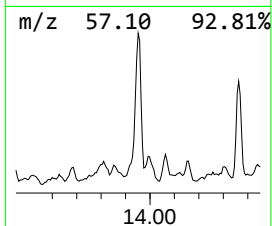
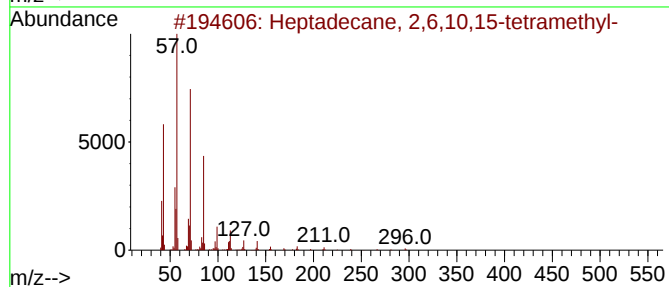
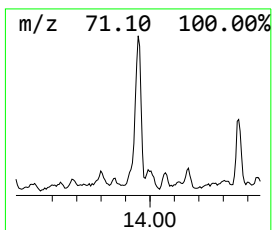
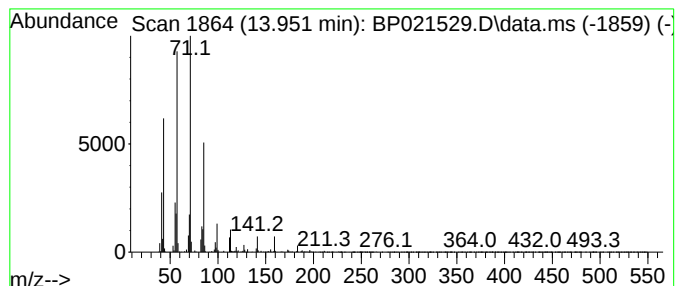
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptadecane, 2,6,10,15-tetr... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.951	2.48 ng	783179	Acenaphthene-d10		14.451	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	86
2	Dodecane, 4,6-dimethyl-		198	C14H30	061141-72-8	64
3	Undecane, 2,6-dimethyl-		184	C13H28	017301-23-4	59
4	Tridecane, 5-propyl-		226	C16H34	055045-11-9	58
5	Sulfurous acid, 2-ethylhexyl pen...		264	C13H28O3S	1000309-18-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081524\
Data File : BP021529.D
Acq On : 16 Aug 2024 01:51
Operator : RC/JU
Sample : P3598-01
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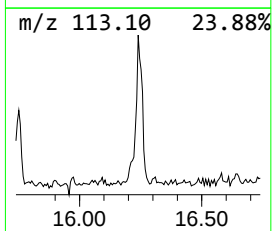
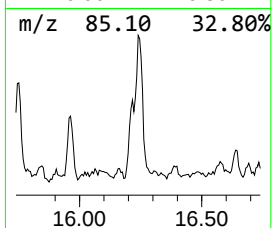
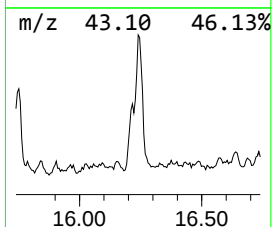
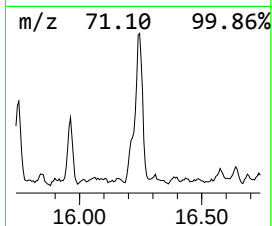
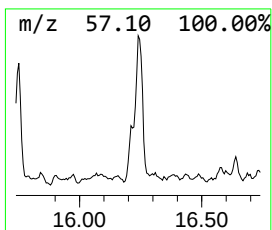
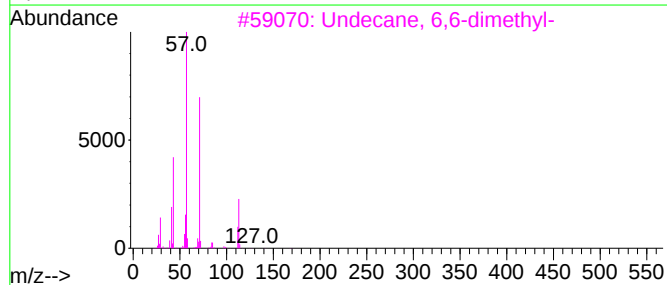
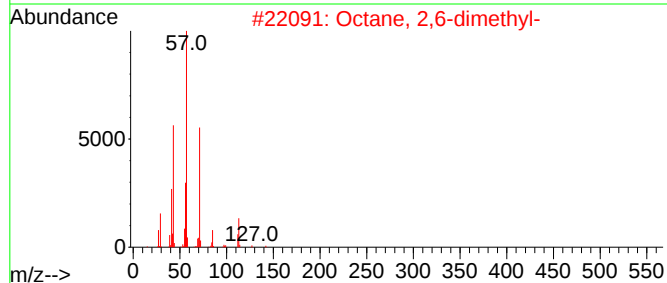
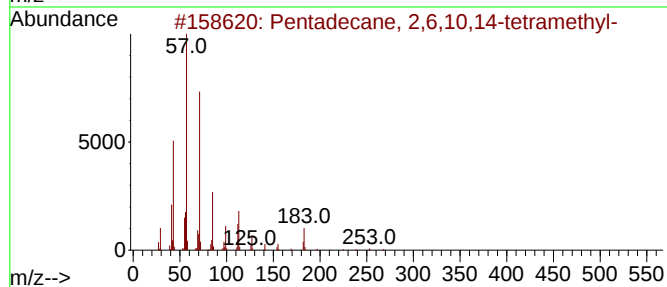
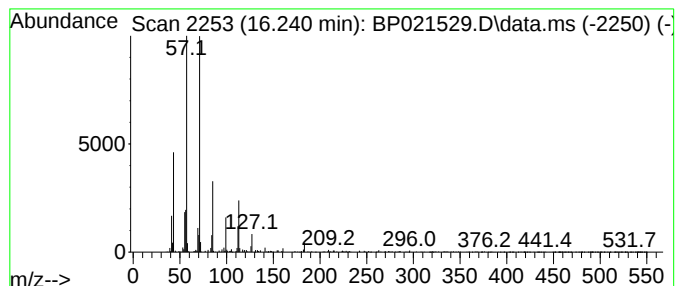
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Pentadecane, 2,6,10,14-tetr... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.240	2.44 ng	766138	Phenanthrene-d10	17.257	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	90
2	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
3	Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	64
4	Heptadecane, 7-methyl-	254	C18H38	020959-33-5	64
5	Sulfurous acid, di(2-ethylhexyl)-	306	C16H34O3S	1000309-19-1	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081524\
Data File : BP021529.D
Acq On : 16 Aug 2024 01:51
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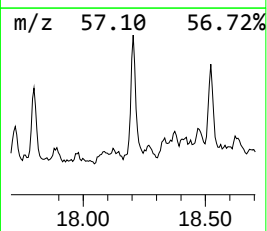
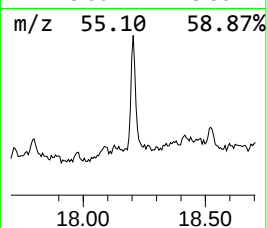
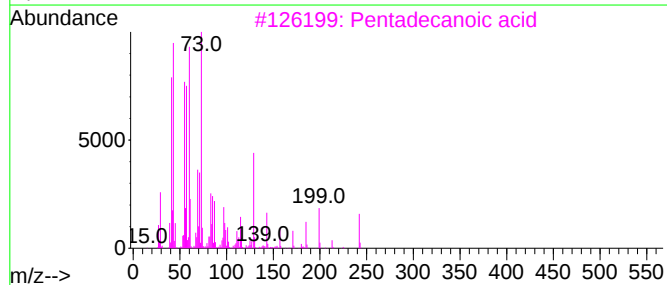
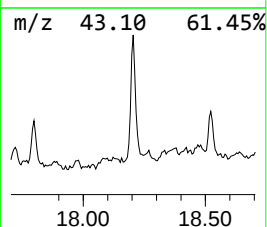
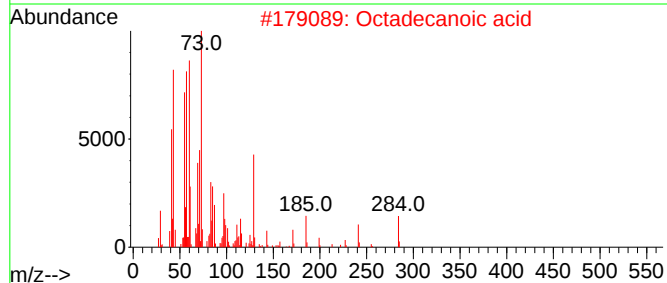
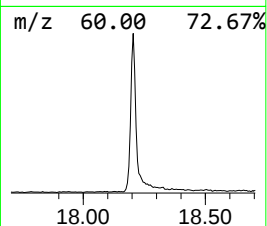
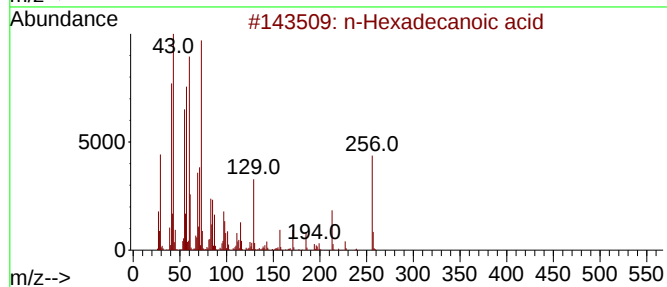
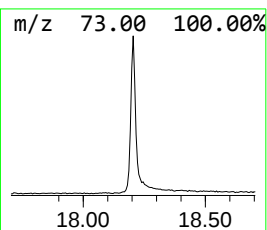
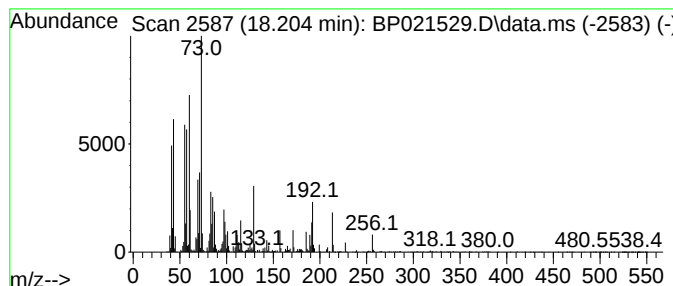
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.204	4.89 ng	1532010	Phenanthrene-d10	17.257

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Octadecanoic acid	284	C18H36O2	000057-11-4	70
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	70
4		Tridecanoic acid	214	C13H26O2	000638-53-9	53
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081524\
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ClientSampleId :
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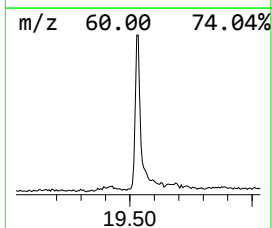
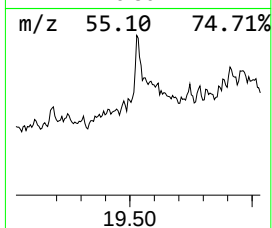
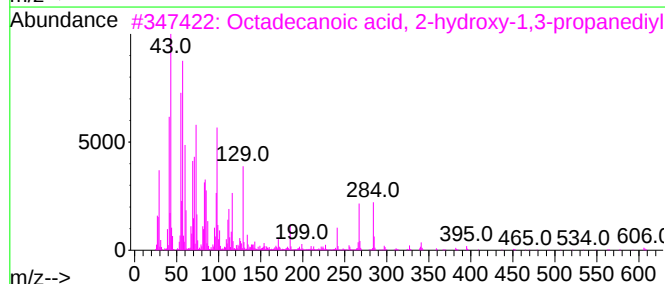
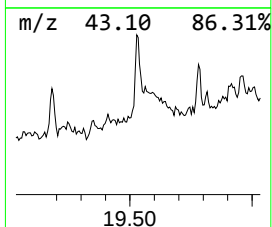
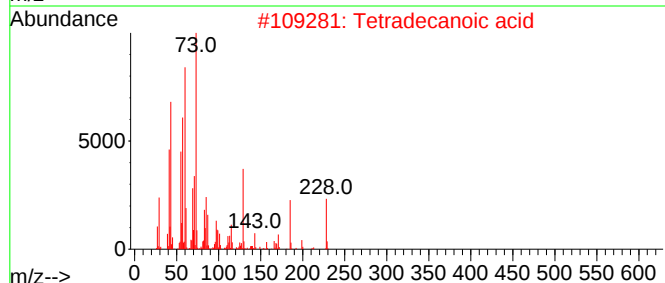
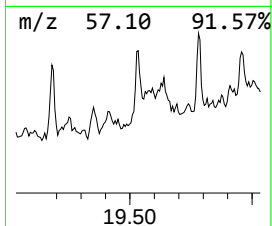
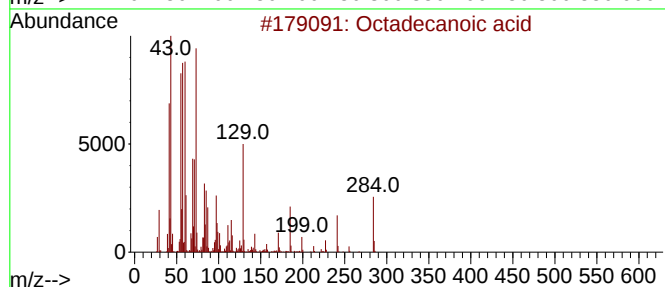
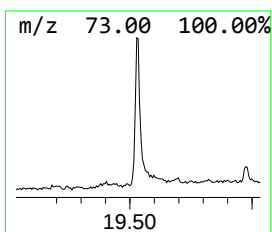
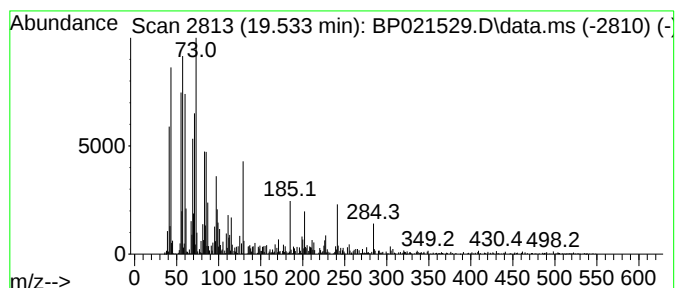
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.534	4.23 ng	963100	Chrysene-d12	21.704

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	94
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	64
3			Octadecanoic acid, 2-hydroxy-1,3...	625	C39H76O5	000504-40-5	64
4			Dodecanoic acid	200	C12H24O2	000143-07-7	43
5			8-Methylnonanoic acid	172	C10H20O2	005963-14-4	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081524\
Data File : BP021529.D
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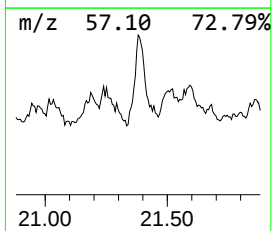
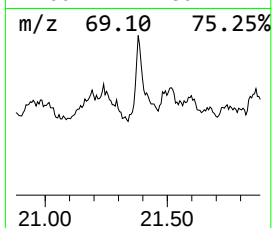
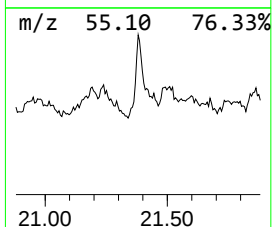
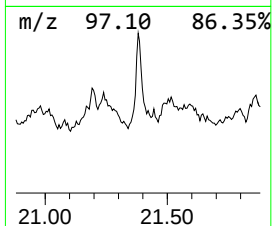
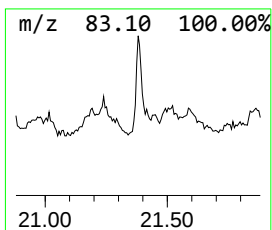
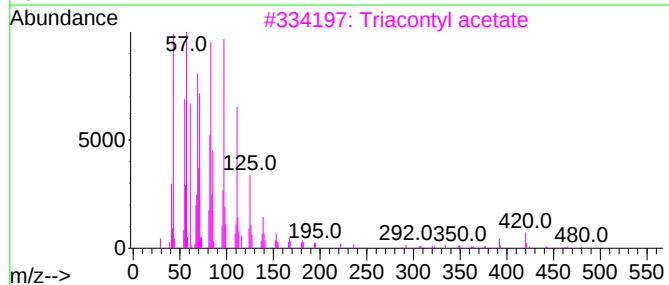
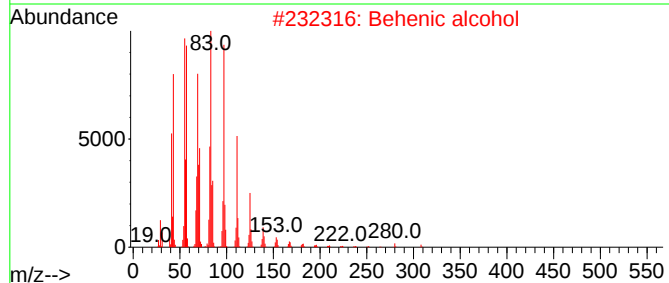
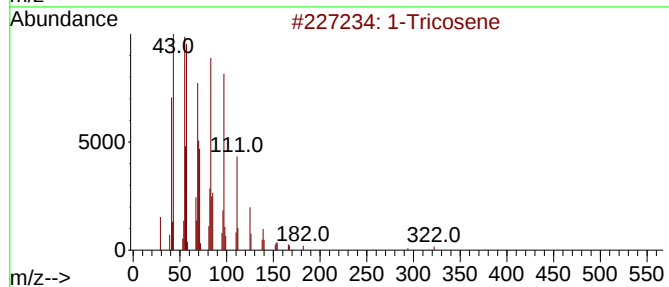
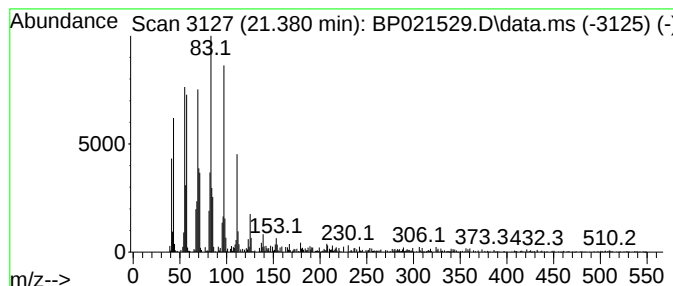
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Tricosene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.380	5.22 ng	1188370	Chrysene-d12	21.704

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tricosene		322	C23H46	018835-32-0	91
2	Behenic alcohol		326	C22H46O	000661-19-8	91
3	Triacontyl acetate		480	C32H64O2	041755-58-2	81
4	Cyclotetracosane		336	C24H48	000297-03-0	76
5	Dichloroacetic acid, heptadecyl ...		366	C19H36Cl2O2	1000282-98-2	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081524\
Data File : BP021529.D
Acq On : 16 Aug 2024 01:51
Operator : RC/JU
Sample : P3598-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JC-03-081324-A

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
Butane, 2-metho...	3.028	34.1	ng	5590400	1	7.811	3280750	20.0
2-Pentanone, 4-...	4.940	3.0	ng	489386	1	7.811	3280750	20.0
1-Phenoxypropan...	11.328	3.2	ng	801687	2	10.593	5049410	20.0
Heptadecane, 2,...	11.640	2.6	ng	658375	2	10.593	5049410	20.0
Naphthalene, 2-...	12.257	2.9	ng	726410	2	10.593	5049410	20.0
Dodecane, 2,6,1...	12.999	2.5	ng	772112	3	14.451	6312810	20.0
Heptadecane, 2,...	13.951	2.5	ng	783179	3	14.451	6312810	20.0
Pentadecane, 2,...	16.240	2.4	ng	766138	4	17.257	6268260	20.0
n-Hexadecanoic ...	18.204	4.9	ng	1532010	4	17.257	6268260	20.0
Octadecanoic acid	19.534	4.2	ng	963100	5	21.704	4554160	20.0
1-Tricosene	21.380	5.2	ng	1188370	5	21.704	4554160	20.0