

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082024\
 Data File : BP021580.D
 Acq On : 20 Aug 2024 20:17
 Operator : RC/JU
 Sample : P3600-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 VCPS-B3-081324-10-18

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: BP021580.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	21	rVB	5046665	6302304	37.82%	5.195%
2	3.546	91	95	109	rVB	289582	391700	2.35%	0.323%
3	4.940	326	332	337	rBV	368749	514337	3.09%	0.424%
4	5.399	403	410	419	rBV	6155476	9290214	55.75%	7.658%
5	6.969	669	677	686	rBV	6114654	9844398	59.08%	8.115%
6	7.799	812	818	825	rBV	1323700	2162869	12.98%	1.783%
7	8.946	997	1013	1023	rBV	3767948	6386265	38.32%	5.264%
8	9.040	1025	1029	1039	rVB3	90369	191617	1.15%	0.158%
9	9.510	1105	1109	1121	rVB4	116048	223801	1.34%	0.184%
10	10.593	1284	1293	1299	rBV	1700612	3166972	19.01%	2.611%
11	10.651	1299	1303	1312	rVB2	144683	258488	1.55%	0.213%
12	11.328	1405	1418	1425	rBV	365933	757522	4.55%	0.624%
13	11.640	1465	1471	1480	rBV2	88708	185110	1.11%	0.153%
14	12.993	1696	1701	1706	rVB	134729	177079	1.06%	0.146%
15	13.063	1706	1713	1726	rBV	7568086	12454917	74.74%	10.267%
16	13.275	1744	1749	1757	rVB5	93615	182468	1.10%	0.150%
17	13.945	1854	1863	1868	rBV2	165187	281683	1.69%	0.232%
18	14.451	1942	1949	1956	rBV	2789056	4346721	26.08%	3.583%
19	14.669	1981	1986	1987	rBV	166207	233818	1.40%	0.193%
20	14.687	1987	1989	1999	rVB	211668	270204	1.62%	0.223%
21	15.010	2038	2044	2049	rBV2	163040	361866	2.17%	0.298%
22	15.516	2126	2130	2134	rBV	124518	176361	1.06%	0.145%
23	15.751	2166	2170	2178	rVB	110195	184845	1.11%	0.152%
24	15.963	2199	2206	2215	rBV	7196271	10882179	65.30%	8.971%
25	16.239	2250	2253	2259	rVB	175900	248540	1.49%	0.205%
26	17.092	2394	2398	2407	rVB	374079	639972	3.84%	0.528%
27	17.251	2419	2425	2430	rBV	3019885	5319887	31.92%	4.385%
28	17.304	2430	2434	2439	rVB	1038232	1585233	9.51%	1.307%
29	17.392	2445	2449	2455	rVB	276891	395031	2.37%	0.326%
30	17.722	2501	2505	2510	rBV	287106	455955	2.74%	0.376%
31	18.204	2583	2587	2592	rBV2	1154611	1636507	9.82%	1.349%
32	18.469	2630	2632	2637	rVB3	259817	310982	1.87%	0.256%
33	18.998	2719	2722	2729	rBV	874649	1298672	7.79%	1.071%
34	19.116	2739	2742	2747	rBV2	610227	857643	5.15%	0.707%
35	19.386	2785	2788	2794	rVB	1693525	2428600	14.57%	2.002%
36	19.457	2797	2800	2803	rBV	657462	820982	4.93%	0.677%

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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

37	19.639	2829	2831	2834	rBV2	621696	703777	4.22%	0.580%
38	19.675	2834	2837	2844	rVB	1077762	1446113	8.68%	1.192%
39	19.757	2848	2851	2859	rVB2	1466553	2586262	15.52%	2.132%
40	19.980	2882	2889	2895	rBV	10104817	16663711	100.00%	13.737%
41	20.204	2924	2927	2931	rVB	2761914	3384457	20.31%	2.790%
42	21.716	3181	3184	3189	rVB2	3047293	4853524	29.13%	4.001%
43	23.386	3465	3468	3473	rBV	580620	934267	5.61%	0.770%
44	25.168	3763	3771	3780	rVB2	1833243	5510359	33.07%	4.542%

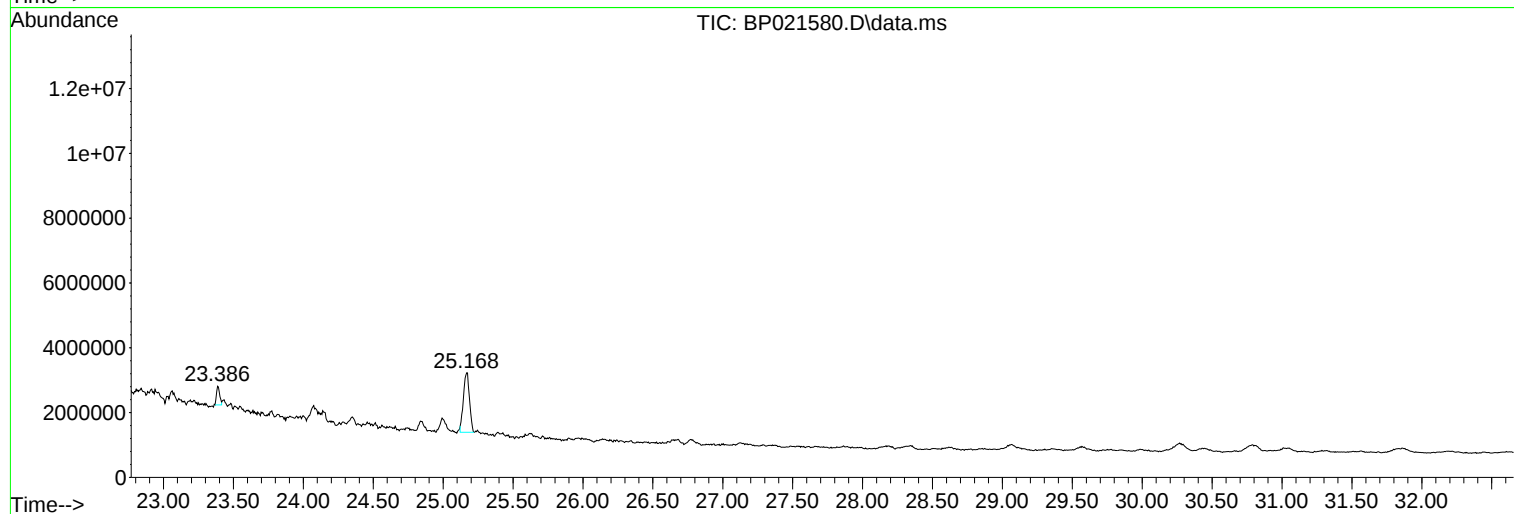
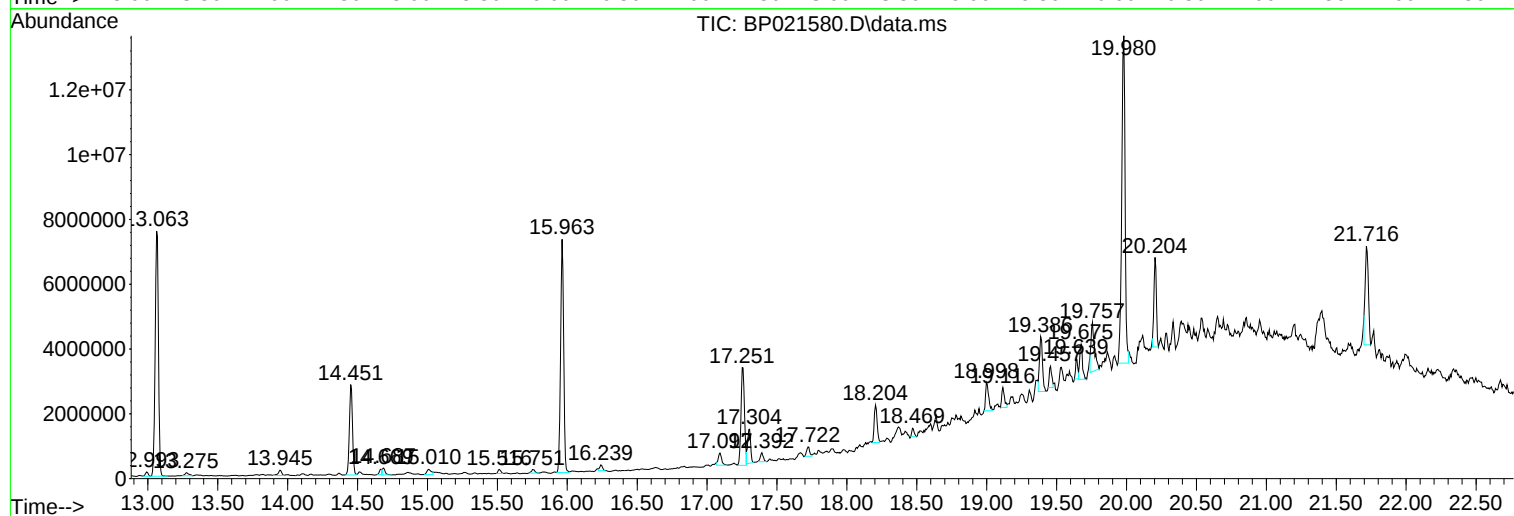
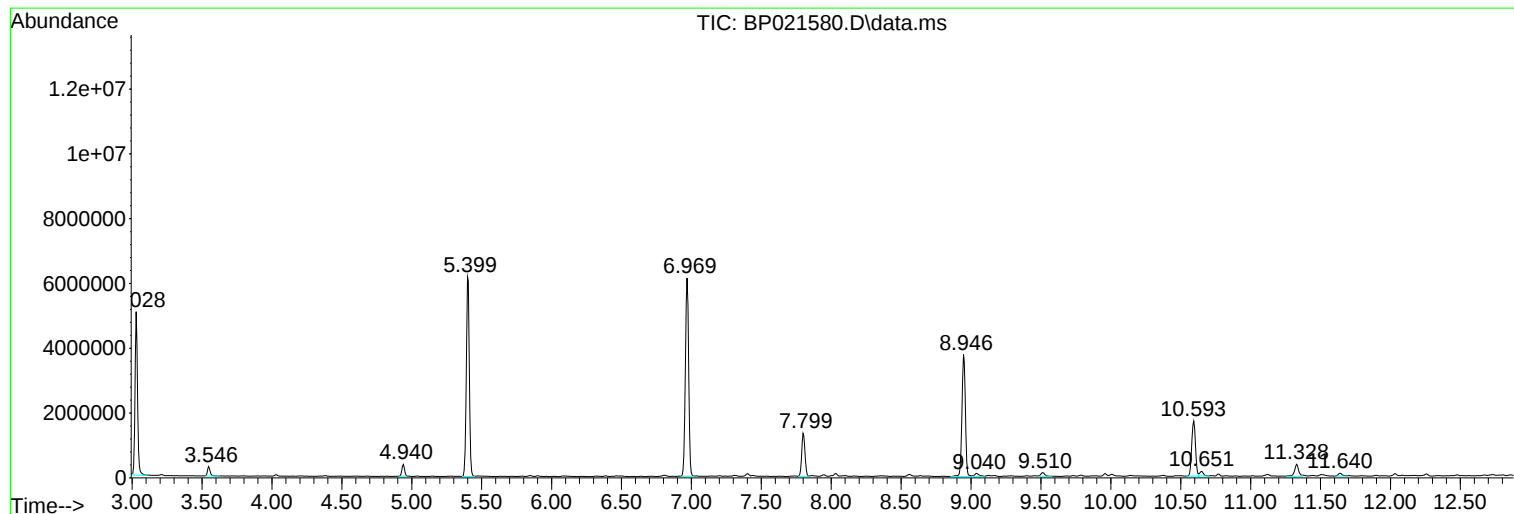
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Sample : P3600-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
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VCPS-B3-081324-10-18

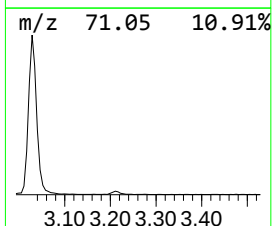
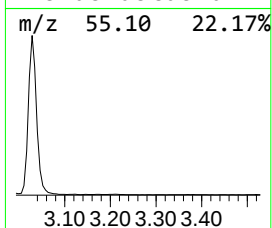
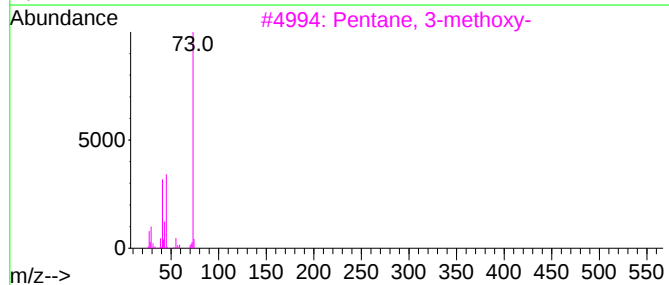
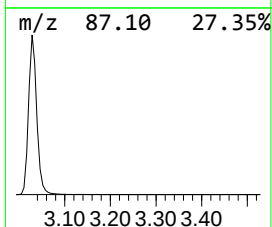
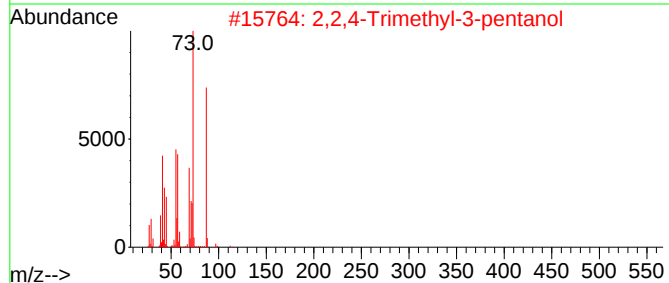
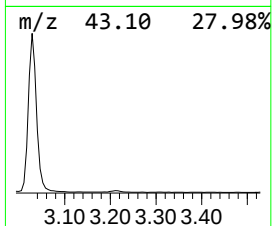
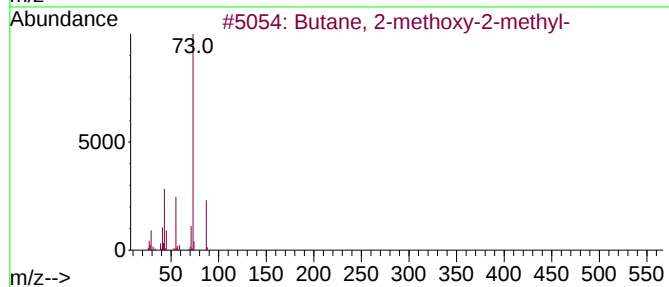
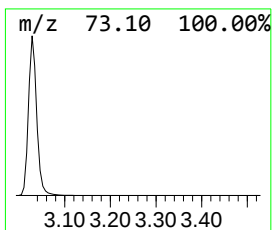
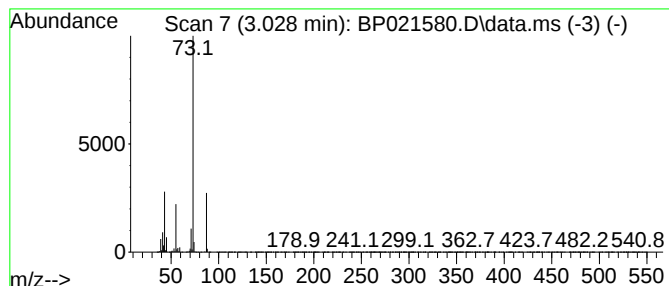
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.028	58.28 ng	6302300	1,4-Dichlorobenzene-d4	7.799

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	86
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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Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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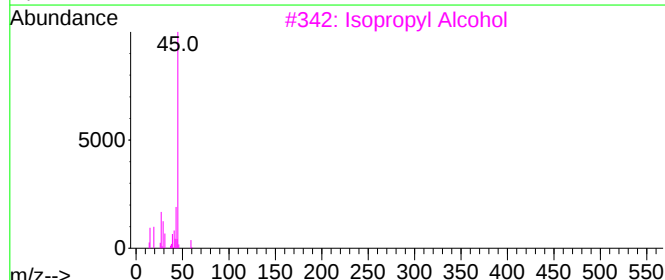
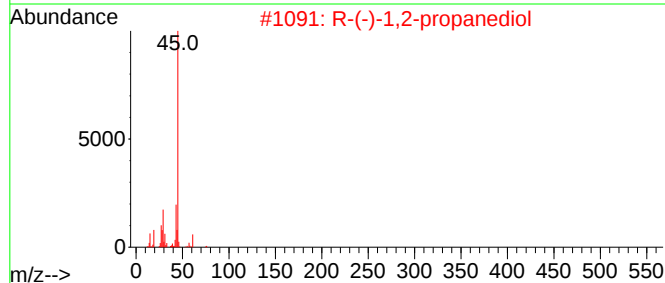
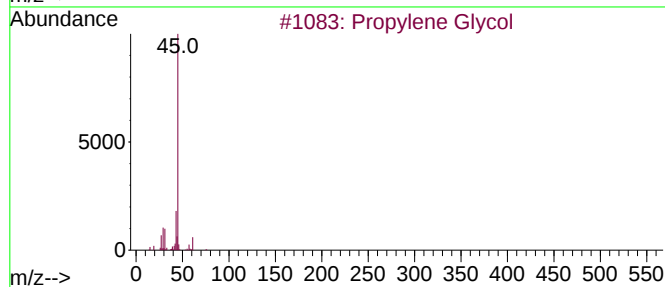
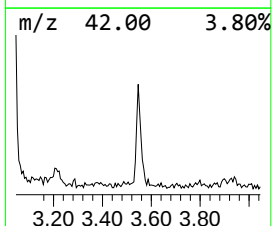
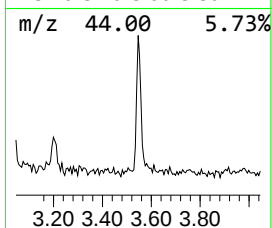
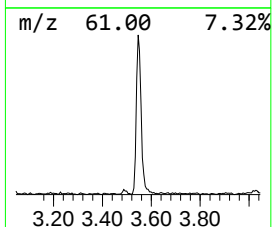
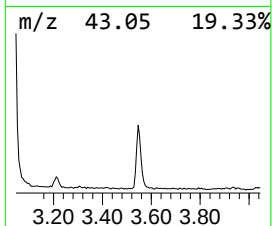
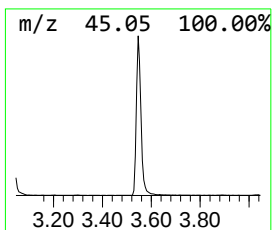
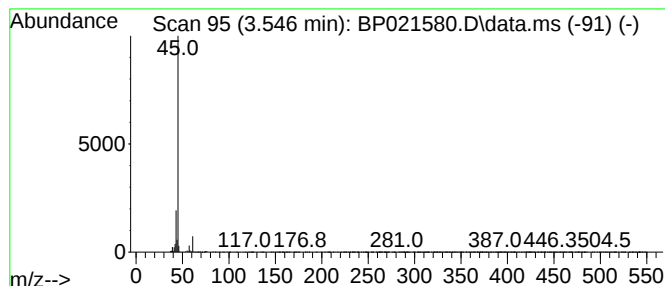
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Propylene Glycol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.546	3.62 ng	391700	1,4-Dichlorobenzene-d4	7.799

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	86
2			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	83
3			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
4			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
5			2-Heptanol	116	C7H16O	000543-49-7	9



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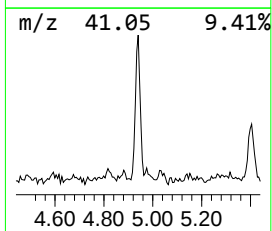
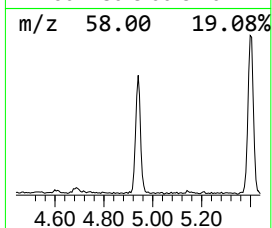
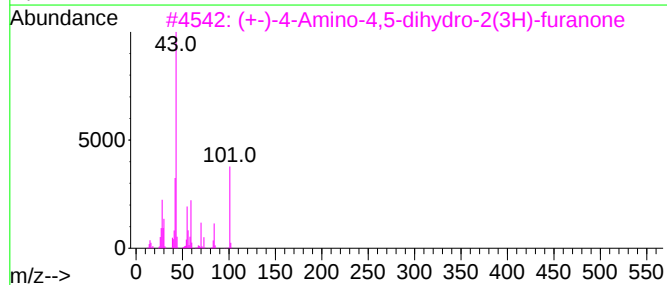
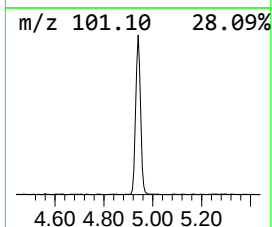
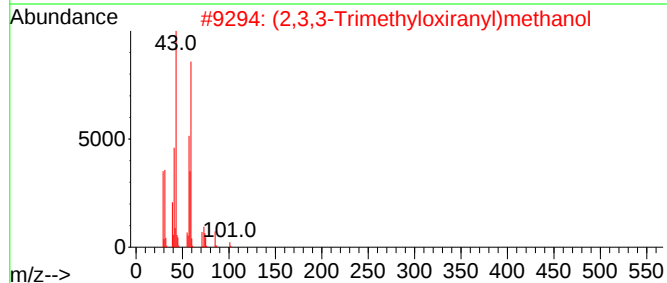
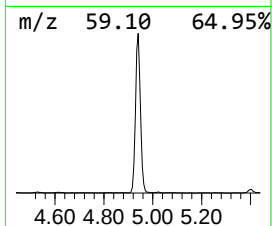
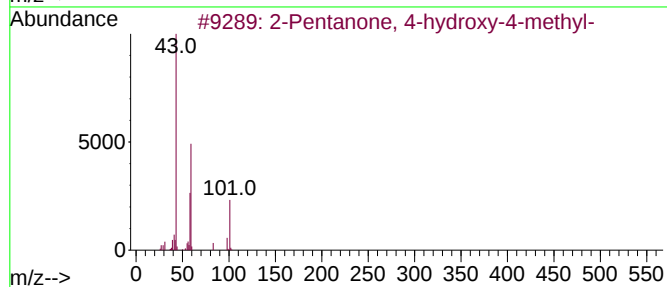
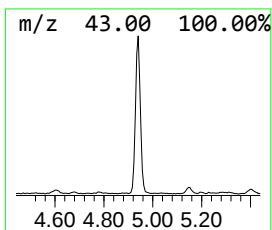
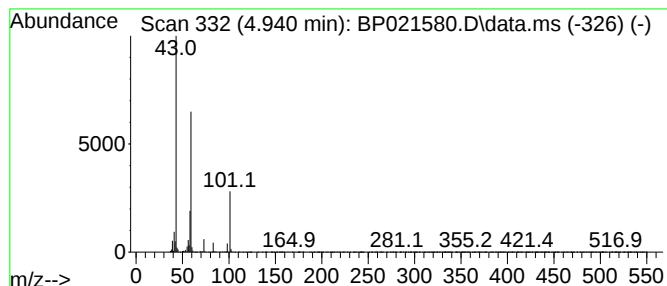
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.940	4.76 ng	514337	1,4-Dichlorobenzene-d4	7.799

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			(2,3,3-Trimethyloxiranyl)methanol	116	C6H12O2	1000306-64-7	33
3			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	32
4			2-Butanol, 1-methoxy-	104	C5H12O2	053778-73-7	23
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	16



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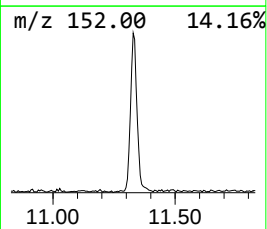
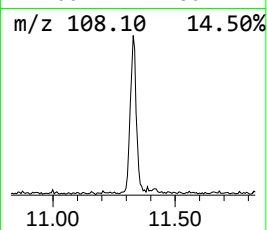
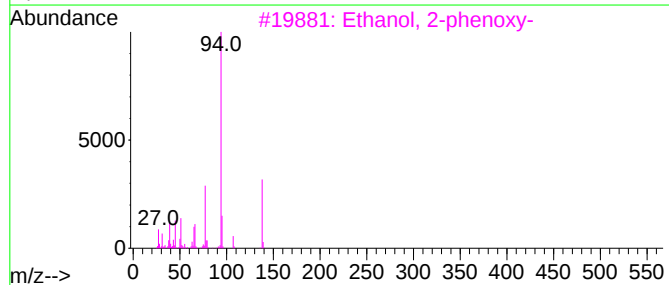
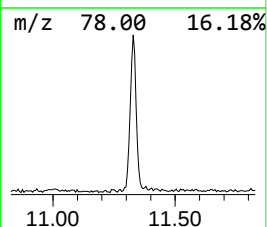
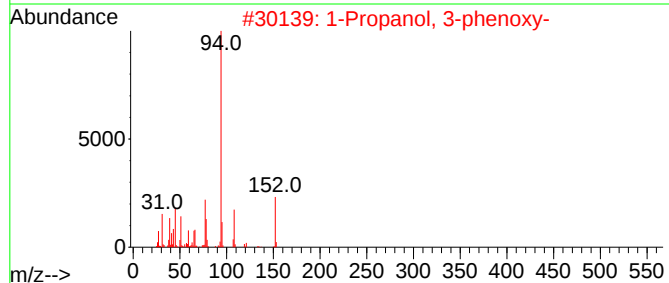
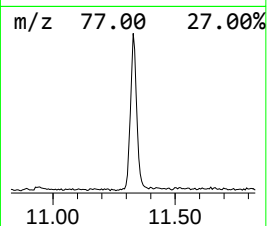
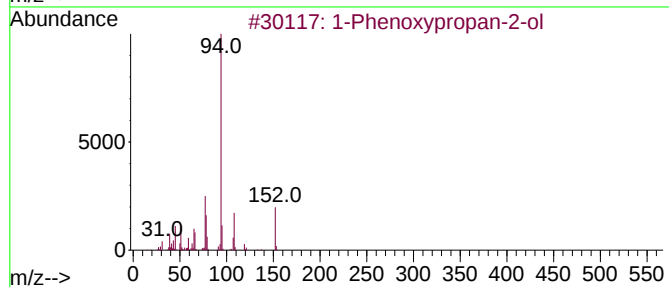
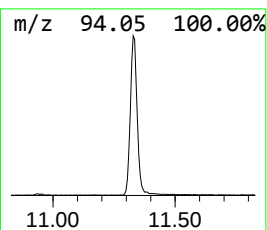
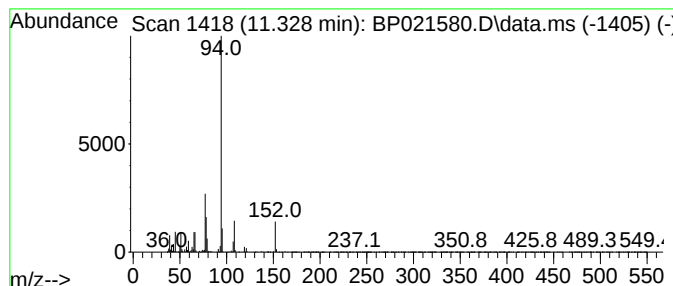
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.328	4.78 ng	757522	Naphthalene-d8	10.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
2			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	83
3			Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64
4			1-Propanol, 2-phenoxy-	152	C9H12O2	004169-04-4	59
5			Carbonic acid, neopentyl phenyl ...	208	C12H16O3	013183-19-2	59



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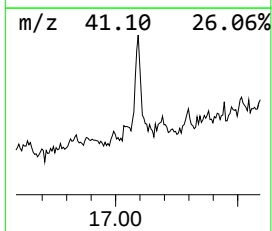
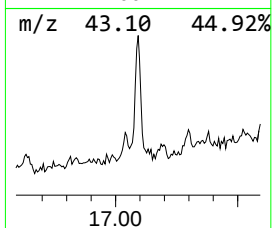
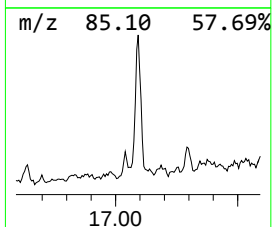
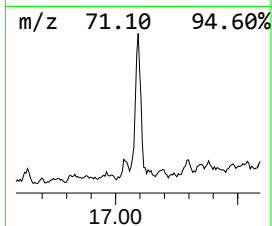
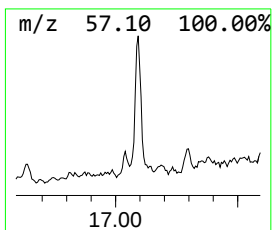
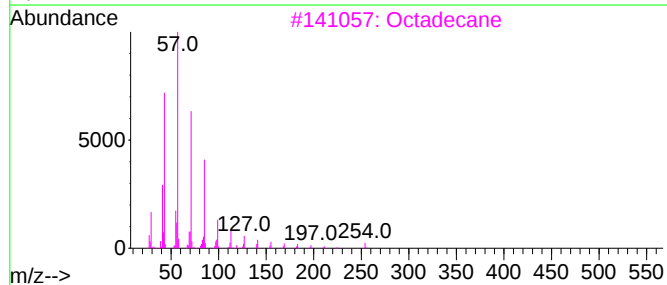
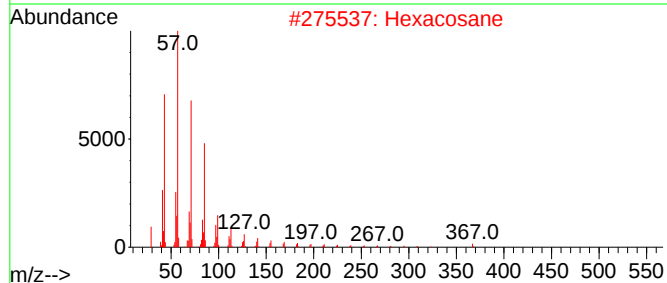
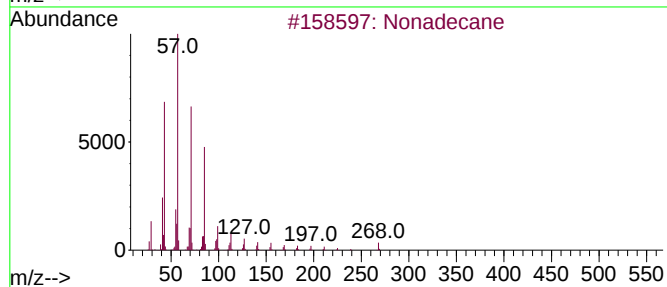
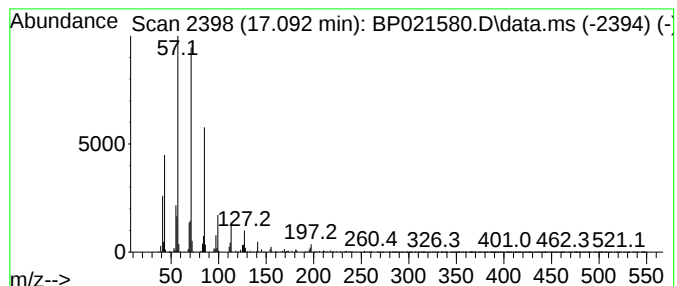
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Nonadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.092	2.41 ng	639972	Phenanthrene-d10	17.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	91
2			Hexacosane	366	C26H54	000630-01-3	90
3			Octadecane	254	C18H38	000593-45-3	90
4			Heneicosane	296	C21H44	000629-94-7	90
5			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082024\
Data File : BP021580.D
Acq On : 20 Aug 2024 20:17
Operator : RC/JU
Sample : P3600-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
VCPS-B3-081324-10-18

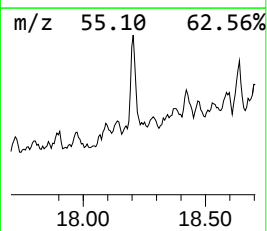
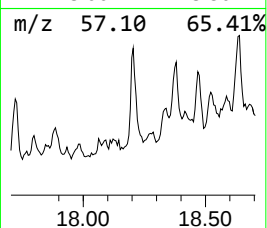
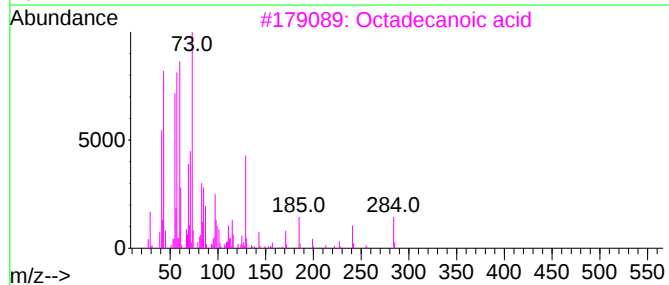
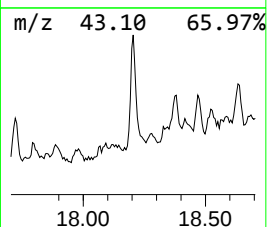
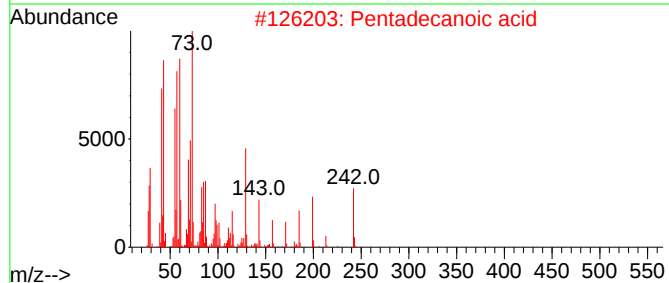
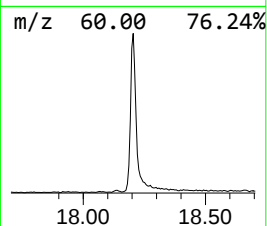
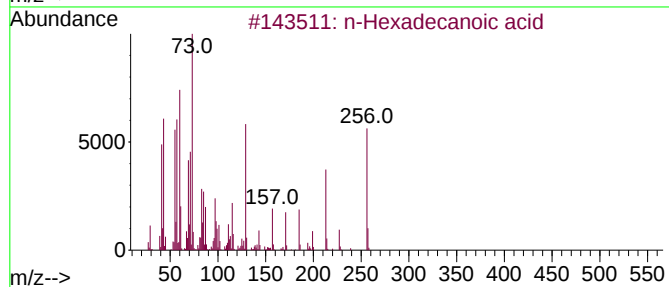
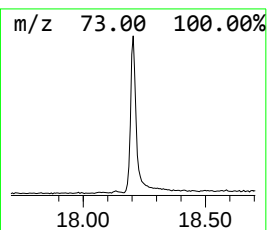
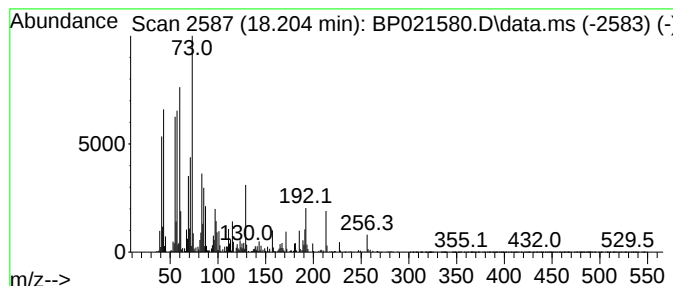
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.204	6.15 ng	1636510	Phenanthrene-d10	17.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	74
3			Octadecanoic acid	284	C18H36O2	000057-11-4	64
4			Dodecanoic acid	200	C12H24O2	000143-07-7	50
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082024\
Data File : BP021580.D
Acq On : 20 Aug 2024 20:17
Operator : RC/JU
Sample : P3600-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

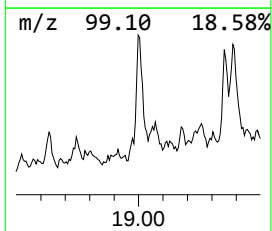
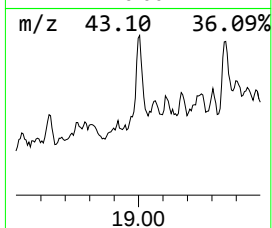
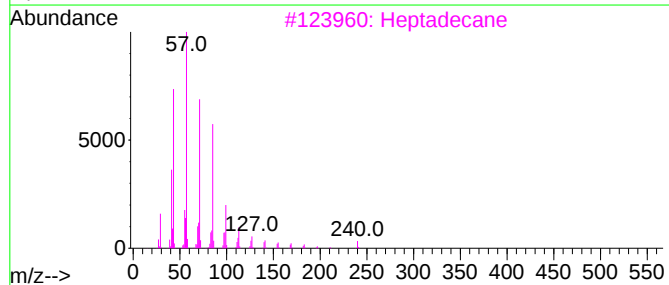
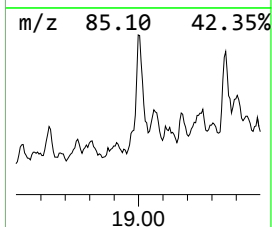
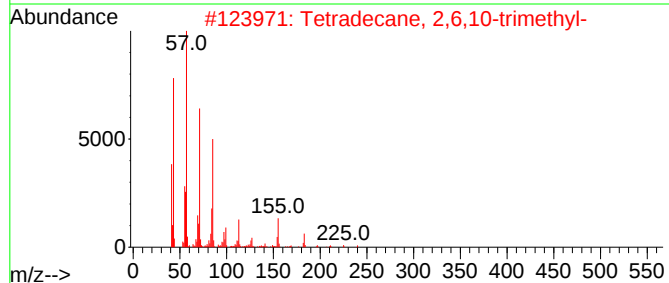
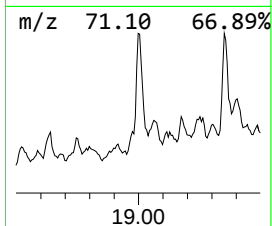
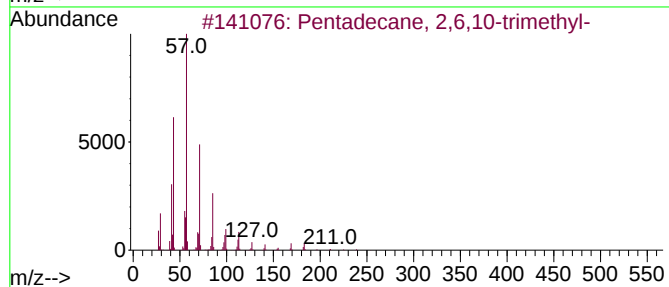
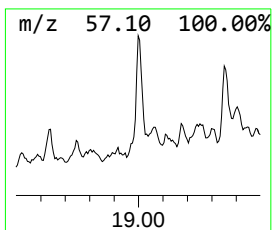
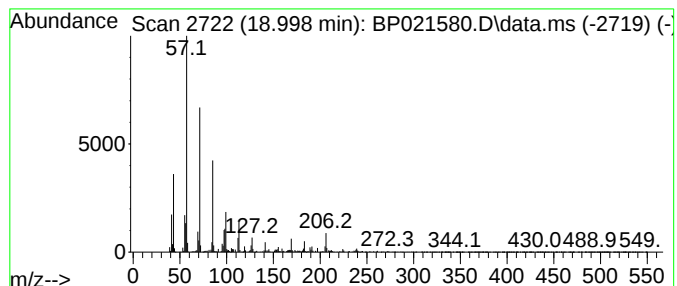
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ClientSampleId :
VCPS-B3-081324-10-18

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 7 Pentadecane, 2,6,10-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.998	4.88 ng	1298670	Phenanthrene-d10		17.251	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10-trimethyl-		254	C18H38	003892-00-0	90
2	Tetradecane, 2,6,10-trimethyl-		240	C17H36	014905-56-7	87
3	Heptadecane		240	C17H36	000629-78-7	87
4	Octadecane		254	C18H38	000593-45-3	86
5	Hexacosane		366	C26H54	000630-01-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082024\
Data File : BP021580.D
Acq On : 20 Aug 2024 20:17
Operator : RC/JU
Sample : P3600-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
VCPS-B3-081324-10-18

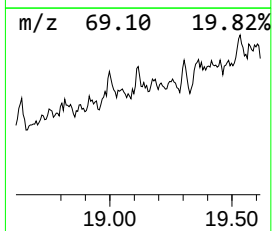
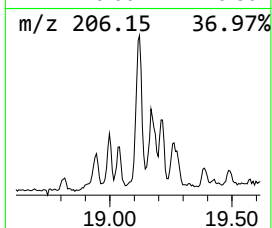
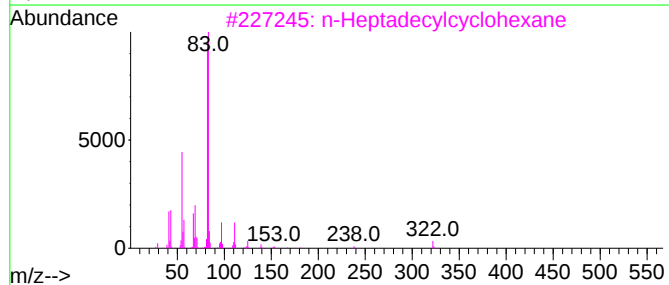
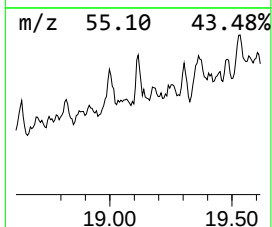
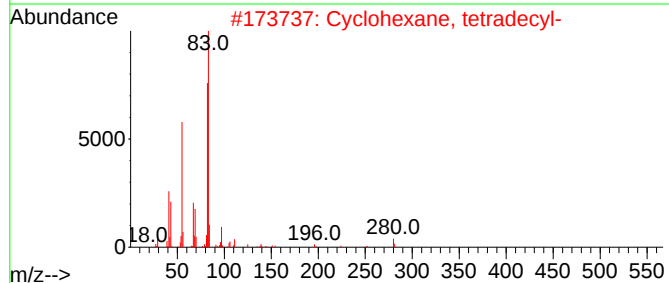
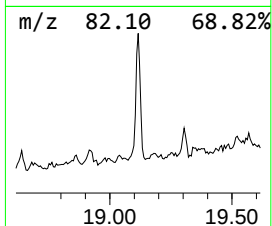
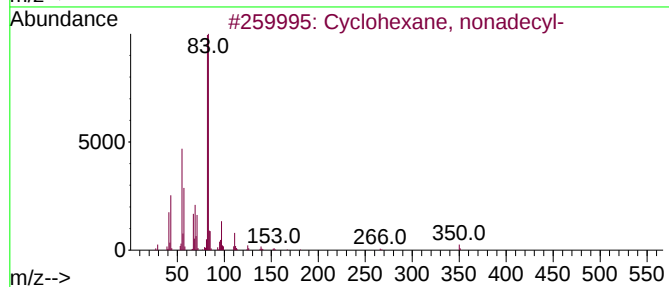
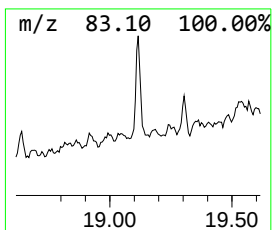
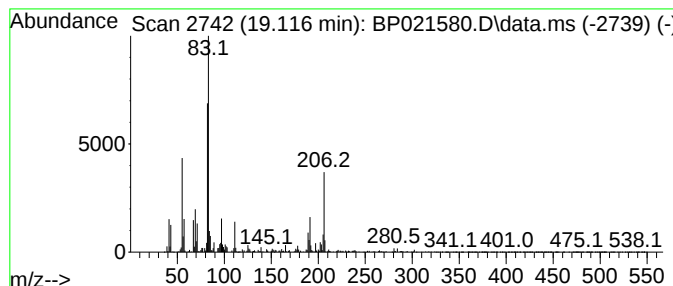
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Cyclohexane, nonadecyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.116	3.22 ng	857643	Phenanthrene-d10	17.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, nonadecyl-	350	C25H50	022349-03-7	76
2			Cyclohexane, tetradecyl-	280	C20H40	001795-18-2	70
3			n-Heptadecylcyclohexane	322	C23H46	019781-73-8	64
4			Cyclohexane, undecyl-	238	C17H34	054105-66-7	62
5			Heptylcyclohexane	182	C13H26	005617-41-4	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082024\
Data File : BP021580.D
Acq On : 20 Aug 2024 20:17
Operator : RC/JU
Sample : P3600-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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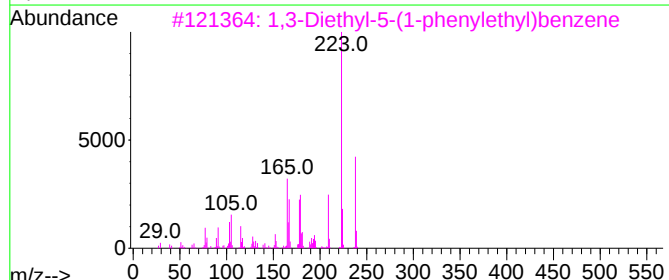
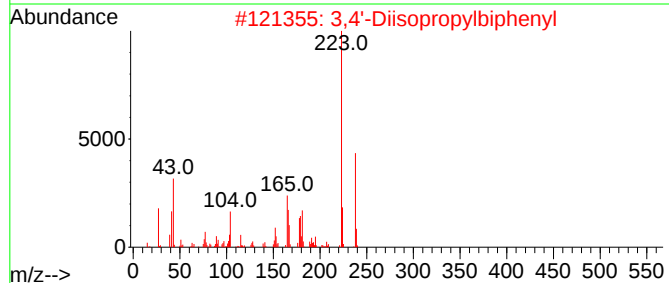
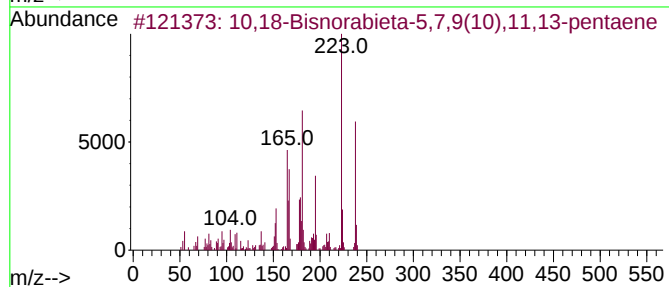
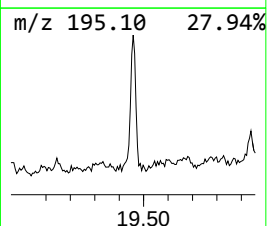
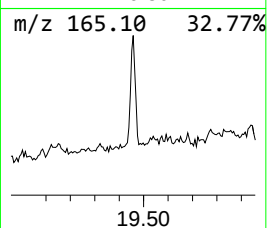
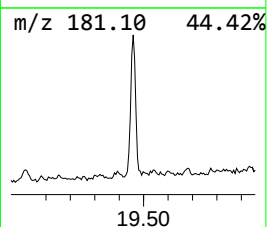
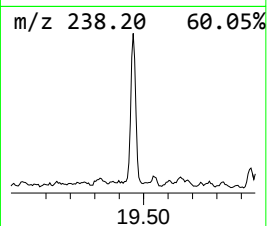
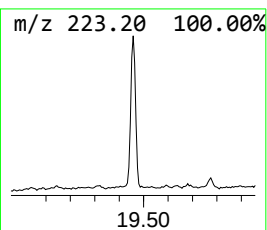
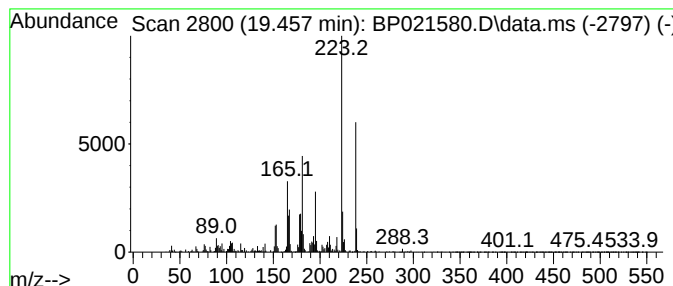
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.457	3.09 ng	820982	Phenanthrene-d10	17.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99		
2	3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	86		
3	1,3-Diethyl-5-(1-phenylethyl)ben...	238	C18H22	1000482-32-6	62		
4	4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	53		
5	4,4'-Diacetyl biphenyl	238	C16H14O2	000787-69-9	49		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082024\
Data File : BP021580.D
Acq On : 20 Aug 2024 20:17
Operator : RC/JU
Sample : P3600-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
VCPS-B3-081324-10-18

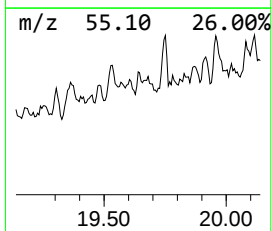
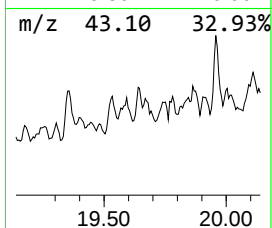
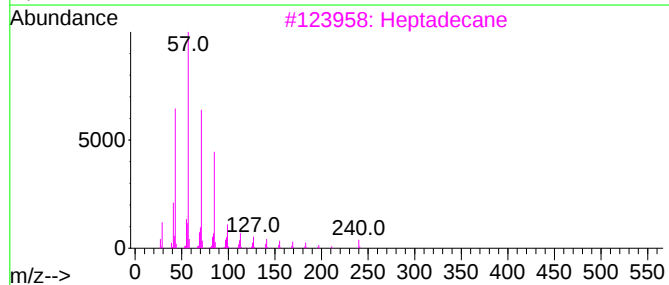
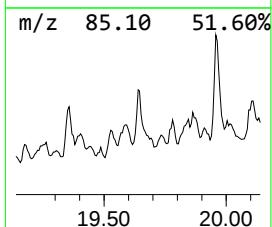
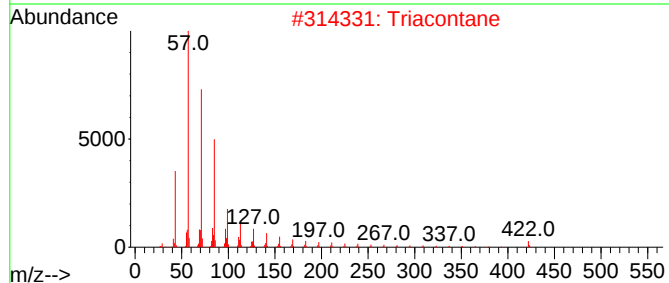
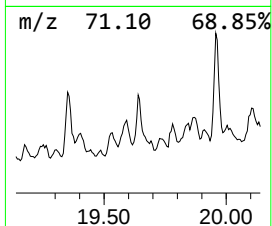
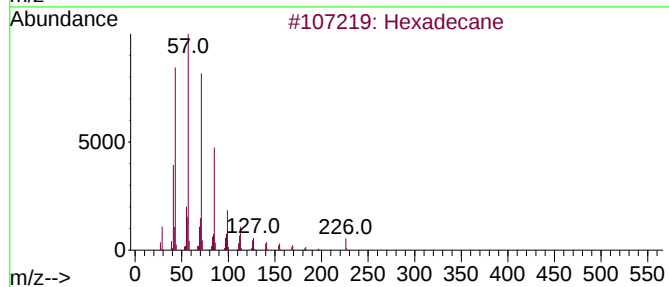
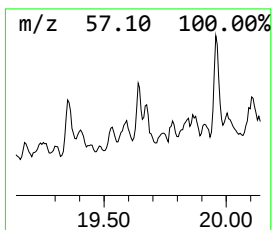
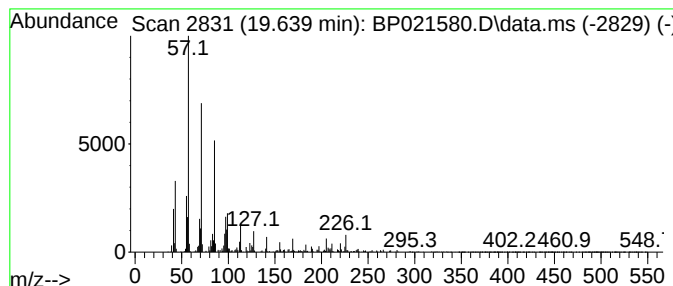
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.639	2.90 ng	703777	Chrysene-d12	21.721

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	93
2		Triacontane	422	C30H62	000638-68-6	87
3		Heptadecane	240	C17H36	000629-78-7	87
4		Octadecane	254	C18H38	000593-45-3	87
5		Docosane	310	C22H46	000629-97-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082024\
Data File : BP021580.D
Acq On : 20 Aug 2024 20:17
Operator : RC/JU
Sample : P3600-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
VCPS-B3-081324-10-18

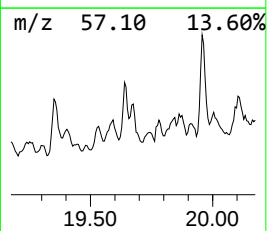
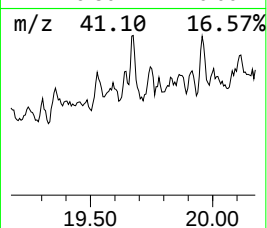
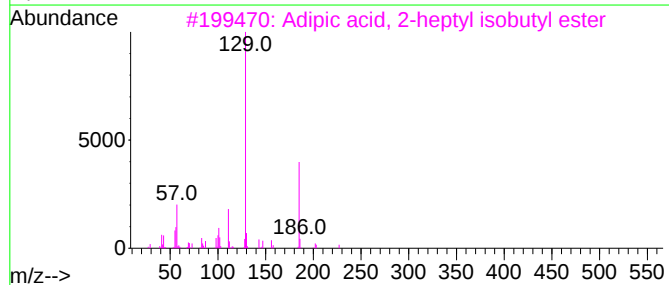
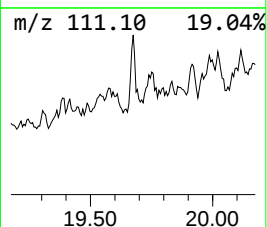
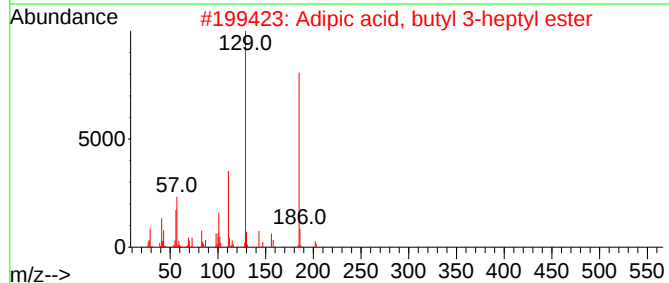
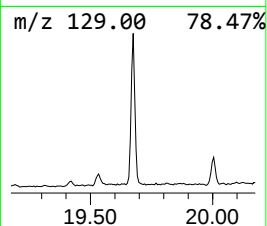
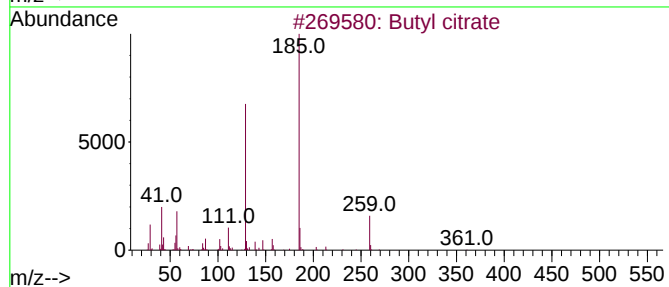
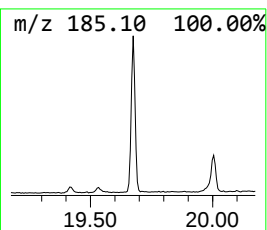
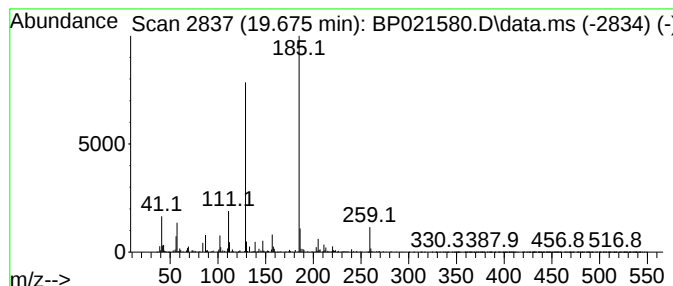
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Butyl citrate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.674	5.96 ng	1446110	Chrysene-d12	21.721

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butyl citrate	360	C18H32O7	000077-94-1	58
2			Adipic acid, butyl 3-heptyl ester	300	C17H32O4	1010353-64-9	56
3			Adipic acid, 2-heptyl isobutyl e...	300	C17H32O4	1000353-64-3	56
4			Adipic acid, cycloheptyl isobuty...	298	C17H30O4	1000324-71-5	50
5			Adipic acid, butyl 4-methylpent...	286	C16H30O4	1000353-50-2	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082024\
Data File : BP021580.D
Acq On : 20 Aug 2024 20:17
Operator : RC/JU
Sample : P3600-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
VCPS-B3-081324-10-18

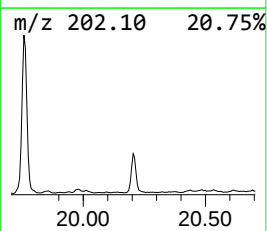
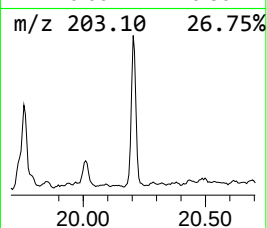
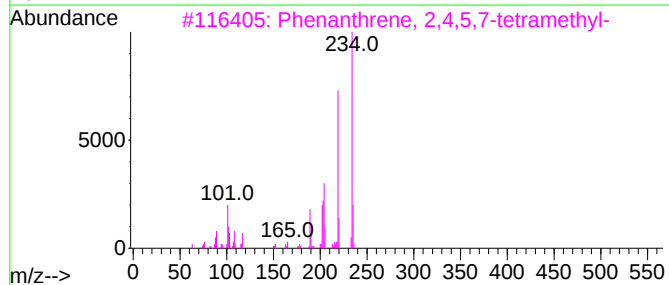
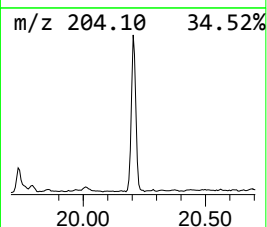
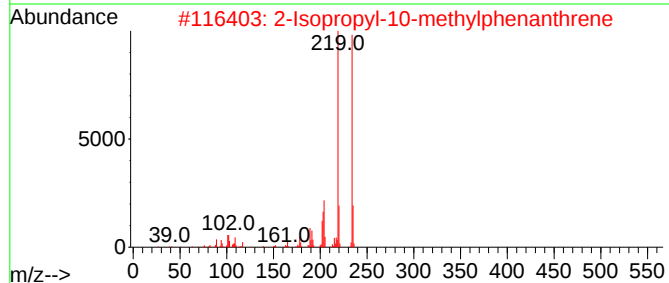
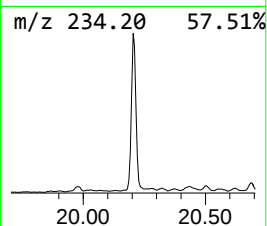
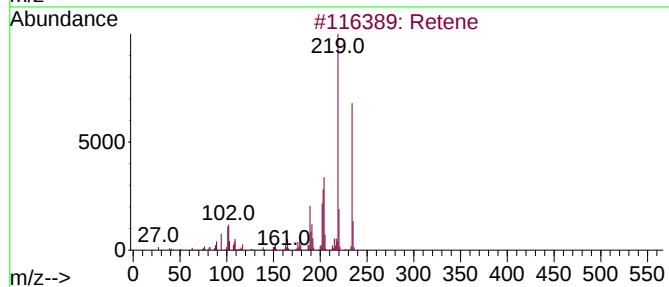
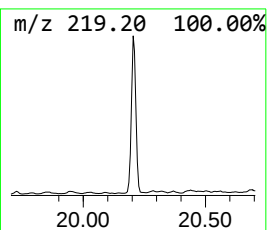
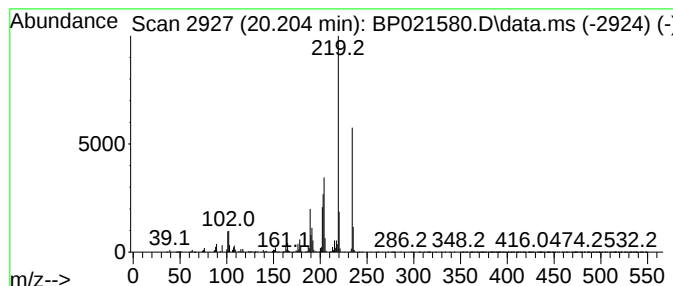
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Retene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.204	13.95 ng	3384460	Chrysene-d12	21.721

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Retene	234	C18H18	000483-65-8	99
2			2-Isopropyl-10-methylphenanthrene	234	C18H18	066552-97-4	95
3			Phenanthrene, 2,4,5,7-tetramethyl-	234	C18H18	007396-38-5	76
4			Ethanone, 1-(4,6-dihydroxy-2,3,5...	234	C13H14O4	021987-07-5	53
5			2-[[1-(4-Methylphenyl)ethylidene...	234	C16H14N2	1000326-46-0	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082024\
Data File : BP021580.D
Acq On : 20 Aug 2024 20:17
Operator : RC/JU
Sample : P3600-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
VCPS-B3-081324-10-18

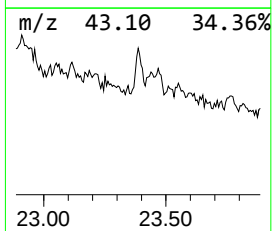
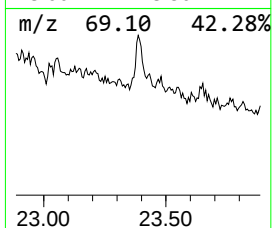
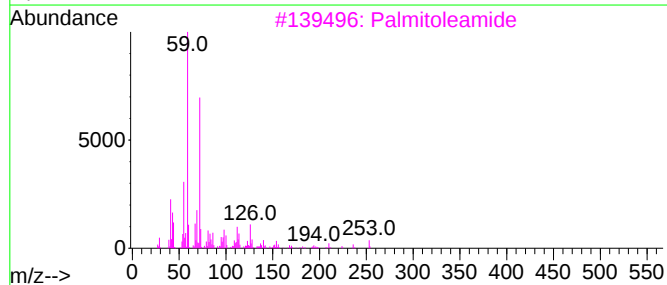
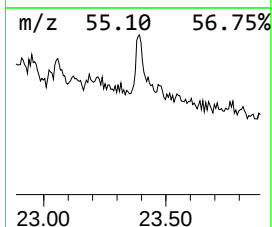
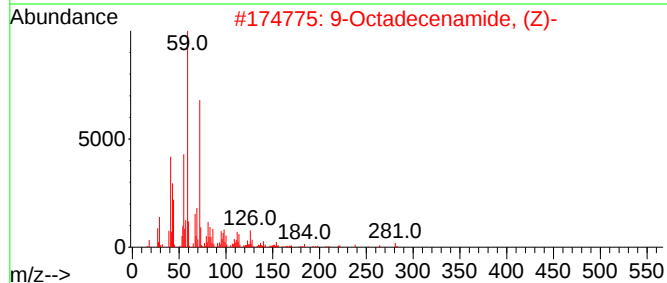
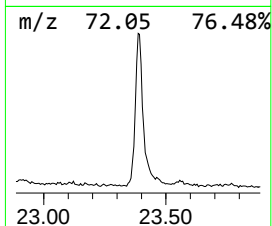
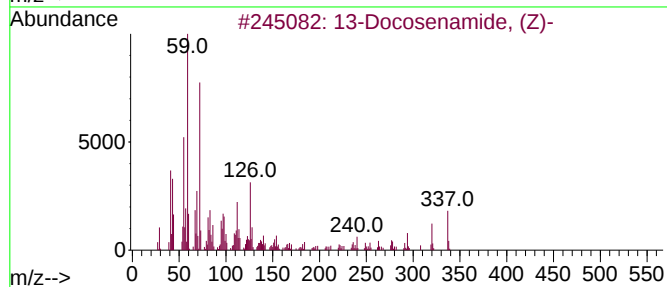
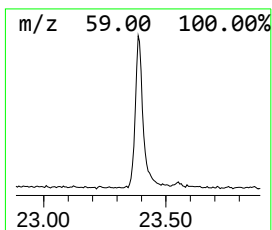
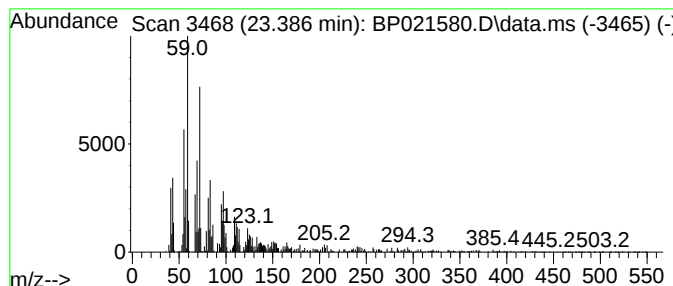
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 13-Docosenamide, (Z)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.386	3.85 ng	934267	Chrysene-d12	21.721

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	74
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	74
3			Palmitoleamide	253	C16H31NO	106010-22-4	59
4			Guanidine, N-[3-[(2-bromophenyl)...	268	C10H13BrN4	1000349-85-6	43
5			Dodecanoylhydrazide, N2-(1-methy...	296	C18H36N2O	1010262-91-8	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082024\
Data File : BP021580.D
Acq On : 20 Aug 2024 20:17
Operator : RC/JU
Sample : P3600-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
VCPS-B3-081324-10-18

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	58.3	ng	6302300	1	7.799	2162870	20.0
Propylene Glycol	3.546	3.6	ng	391700	1	7.799	2162870	20.0
2-Pentanone, 4-...	4.940	4.8	ng	514337	1	7.799	2162870	20.0
1-Phenoxypropan...	11.328	4.8	ng	757522	2	10.593	3166970	20.0
Nonadecane	17.092	2.4	ng	639972	4	17.251	5319890	20.0
n-Hexadecanoic ...	18.204	6.2	ng	1636510	4	17.251	5319890	20.0
Pentadecane, 2,...	18.998	4.9	ng	1298670	4	17.251	5319890	20.0
Cyclohexane, no...	19.116	3.2	ng	857643	4	17.251	5319890	20.0
10,18-Bisnorabi...	19.457	3.1	ng	820982	4	17.251	5319890	20.0
Hexadecane	19.639	2.9	ng	703777	5	21.721	4853520	20.0
Butyl citrate	19.674	6.0	ng	1446110	5	21.721	4853520	20.0
Retene	20.204	13.9	ng	3384460	5	21.721	4853520	20.0
13-Docosenamide...	23.386	3.9	ng	934267	5	21.721	4853520	20.0