

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120919\
 Data File : BP001285.D
 Acq On : 09 Dec 2019 22:44
 Operator : JU
 Sample : K6188-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SOIL-COMP-1

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 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\8270-BP112719.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.628	591	602	619	rBV	398817	684747	18.46%	2.319%
2	6.222	696	703	707	rBV	2043606	2785203	75.09%	9.433%
3	7.763	957	965	976	rBV	2074777	3171214	85.50%	10.741%
4	7.952	989	997	1001	rBV	2323797	2950143	79.54%	9.992%
5	8.305	1052	1057	1069	rVB	409940	483159	13.03%	1.636%
6	8.552	1093	1099	1103	rBV	1922188	2255646	60.81%	7.640%
7	9.193	1200	1208	1212	rBV	1461755	1900625	51.24%	6.437%
8	10.340	1397	1403	1410	rBV	501940	576486	15.54%	1.953%
9	12.122	1696	1706	1710	rBV	2758419	3293897	88.81%	11.156%
10	12.787	1814	1819	1826	rBV	142959	162856	4.39%	0.552%
11	13.204	1883	1890	1897	rBV	528879	682987	18.41%	2.313%
12	14.498	2102	2110	2129	rBV	1902997	2509587	67.66%	8.500%
13	15.598	2288	2297	2302	rBV	644223	718869	19.38%	2.435%
14	16.492	2444	2449	2462	rBV2	122500	189962	5.12%	0.643%
15	17.886	2680	2686	2690	rBV	3445251	3709129	100.00%	12.562%
16	19.116	2890	2895	2911	rBV2	249597	596603	16.08%	2.021%
17	19.239	2911	2916	2920	rBV	647209	681884	18.38%	2.309%
18	19.533	2962	2966	2969	rBV	32335	37326	1.01%	0.126%
19	19.922	3025	3032	3035	rBV	63384	84047	2.27%	0.285%
20	20.298	3092	3096	3099	rBV	78677	87638	2.36%	0.297%
21	20.380	3106	3110	3113	rVB	31771	37408	1.01%	0.127%
22	20.692	3158	3163	3167	rBV	119479	144088	3.88%	0.488%
23	21.010	3211	3217	3224	rVB	480863	668120	18.01%	2.263%
24	21.127	3232	3237	3243	rVB	82361	121152	3.27%	0.410%
25	21.622	3316	3321	3327	rBV	94292	159164	4.29%	0.539%
26	21.698	3328	3334	3341	rBV4	63988	200583	5.41%	0.679%
27	22.198	3413	3419	3423	rBV2	40828	79263	2.14%	0.268%
28	22.892	3527	3537	3546	rBV3	206857	497830	13.42%	1.686%
29	23.380	3616	3620	3633	rVB3	16273	55858	1.51%	0.189%

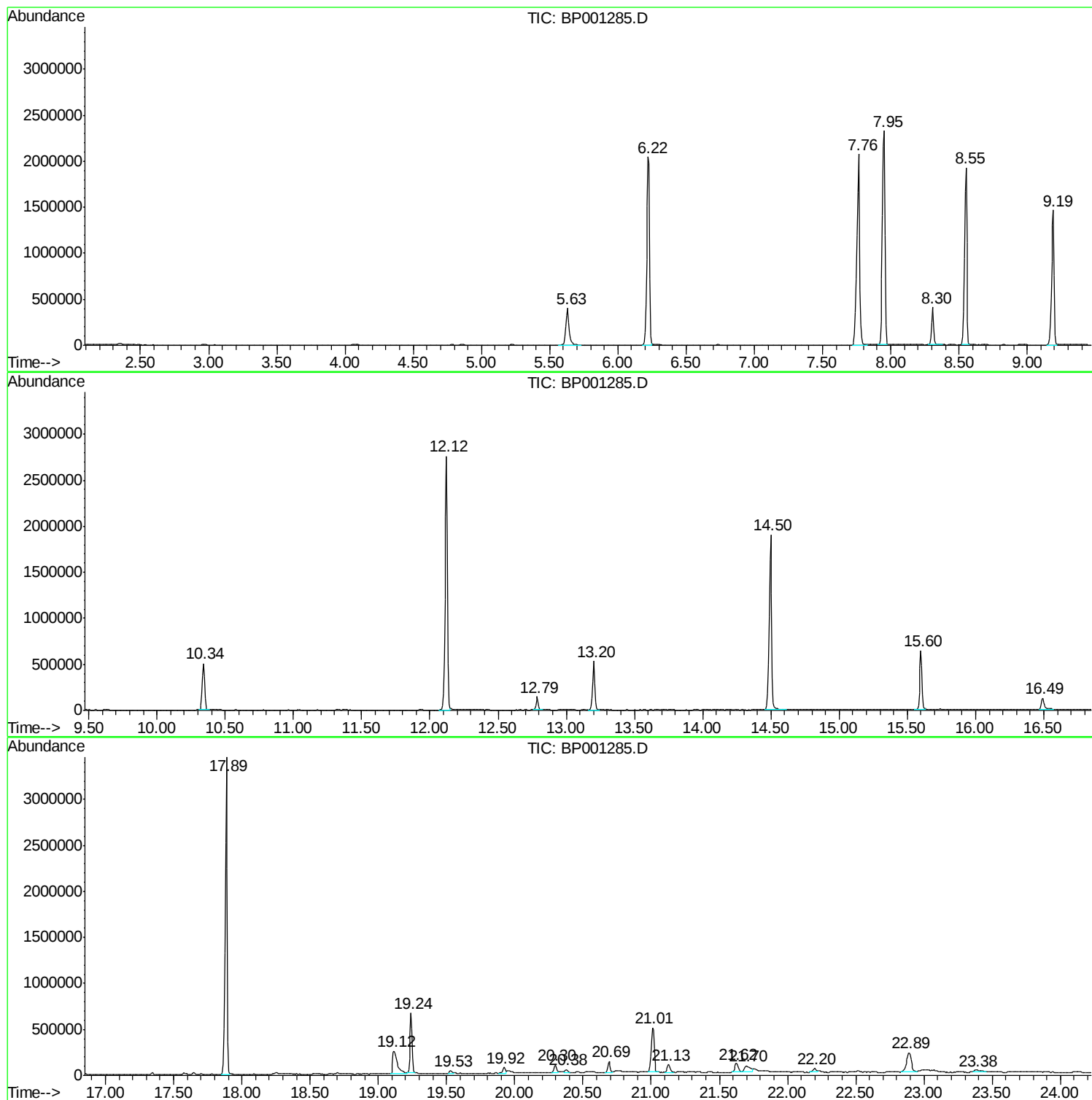
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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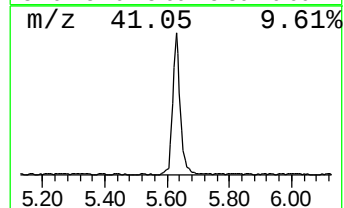
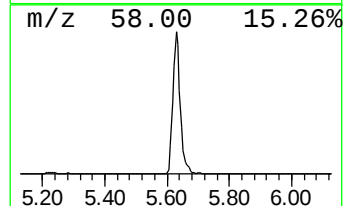
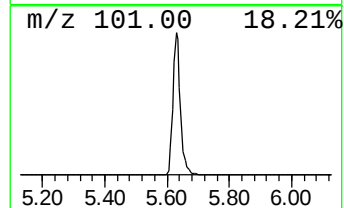
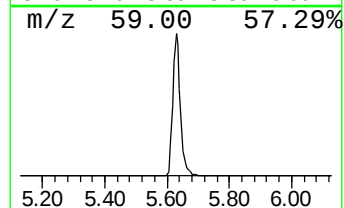
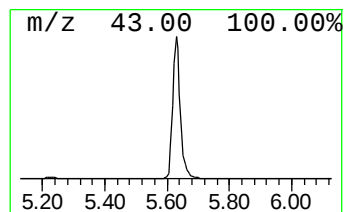
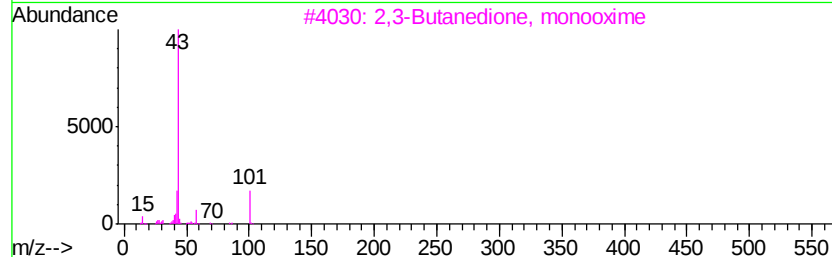
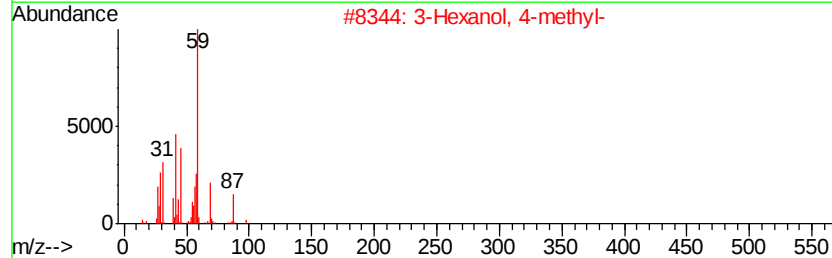
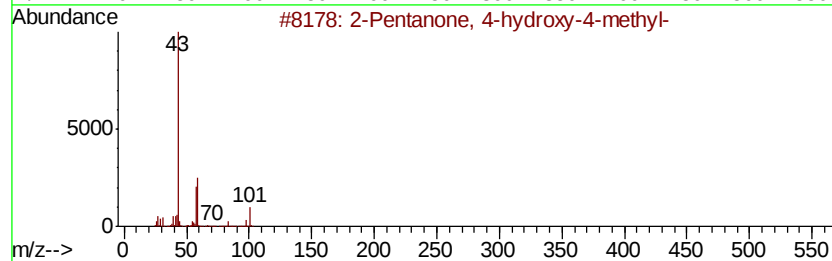
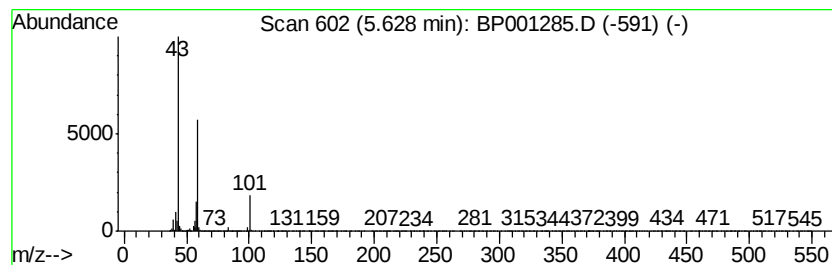
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.63	28.34 ng	684747	1,4-Dichlorobenzene-d4	8.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	9



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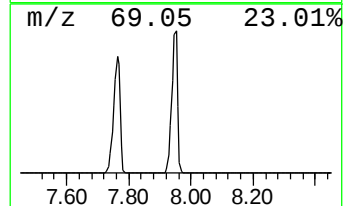
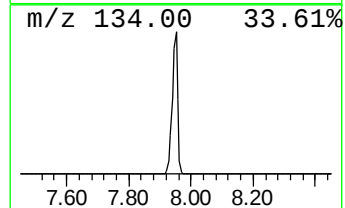
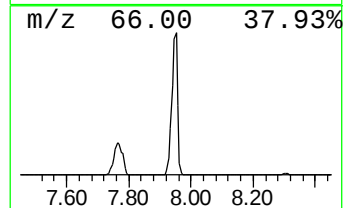
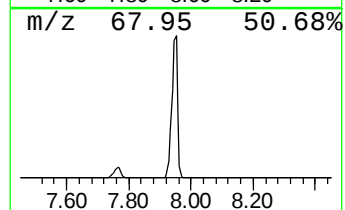
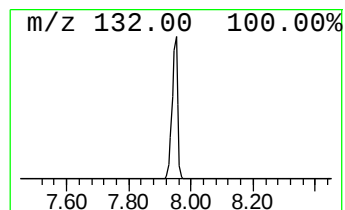
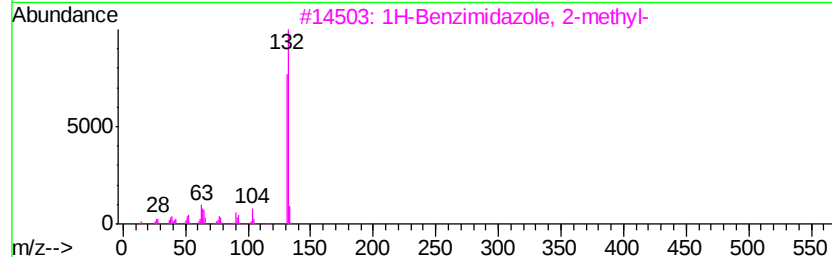
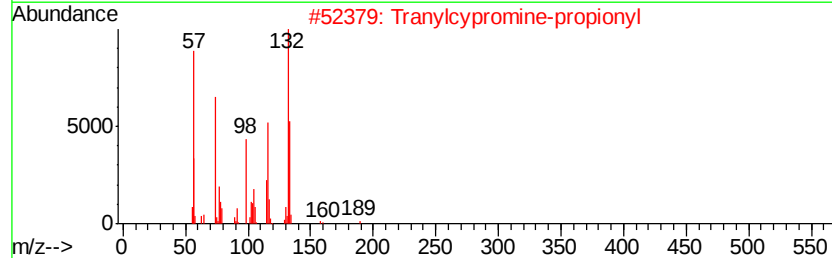
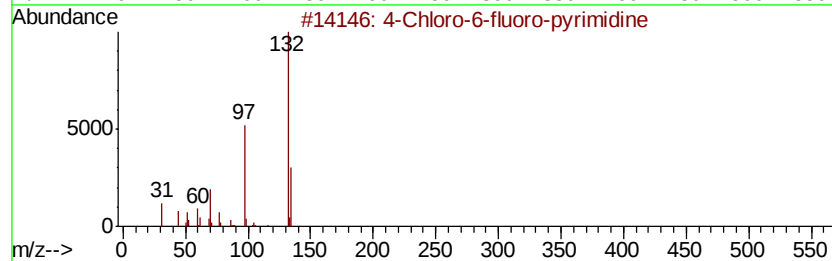
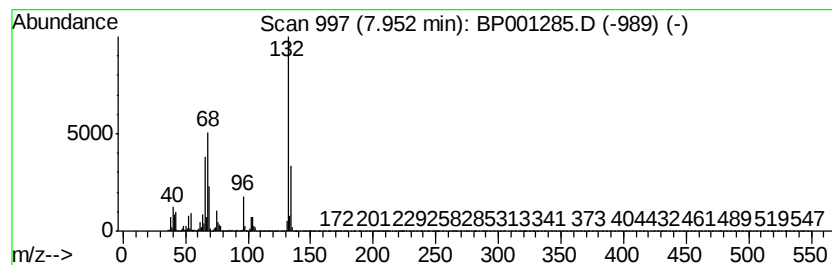
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Peak Number 2 unknown7.95 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.95	122.12 ng	2950140	1,4-Dichlorobenzene-d4	8.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		Tranlycypromine-propionyl	189	C12H15NO	1000123-86-3	22
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	14



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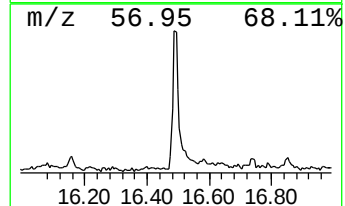
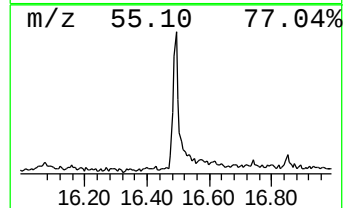
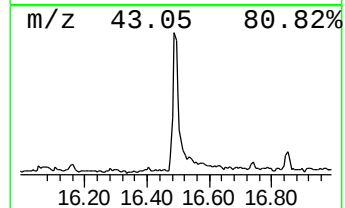
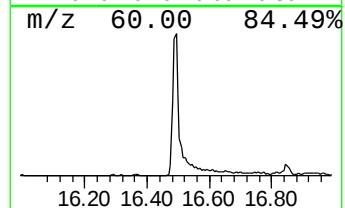
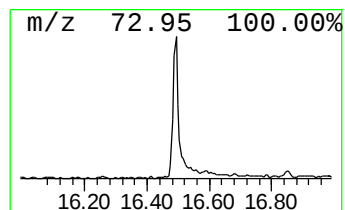
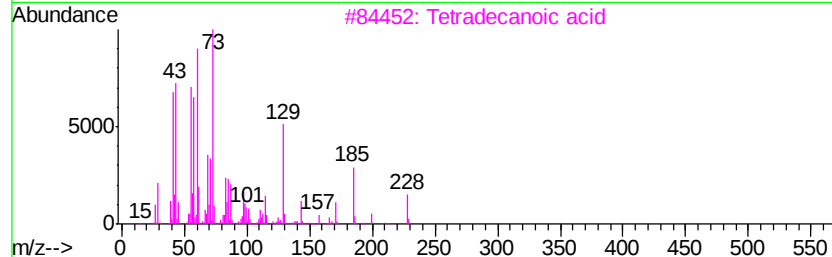
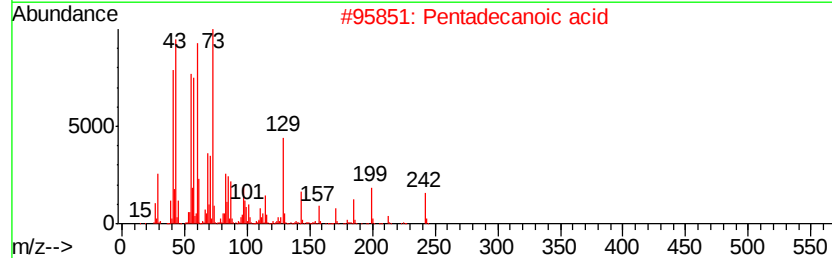
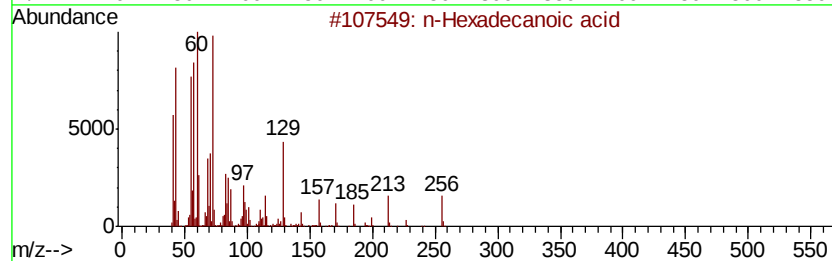
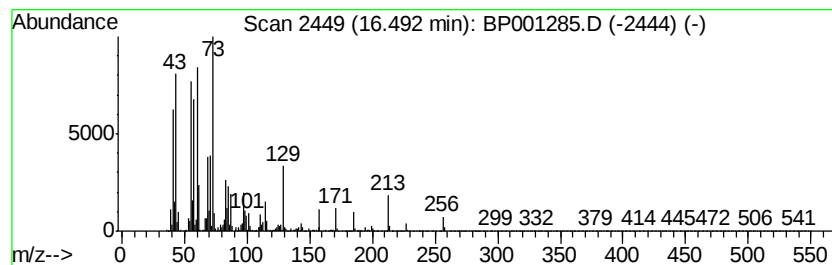
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.49	5.29 ng	189962	Phenanthrene-d10	15.60

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	87
4		Tridecanoic acid	214	C13H26O2	000638-53-9	80
5		11-Bromoundecanoic acid	264	C11H21BrO2	002834-05-1	38



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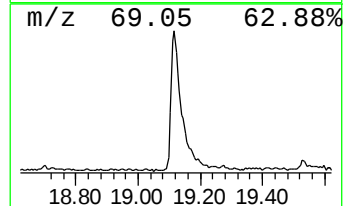
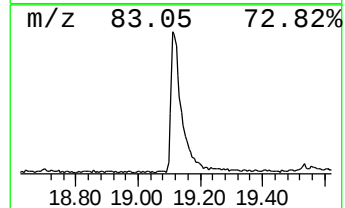
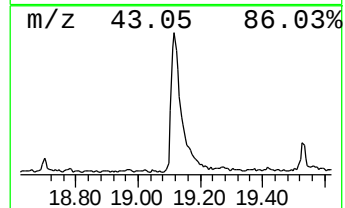
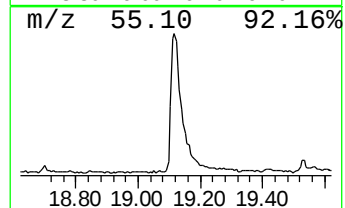
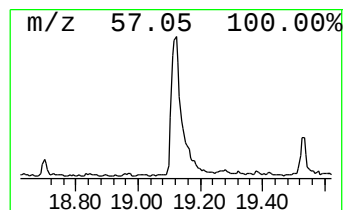
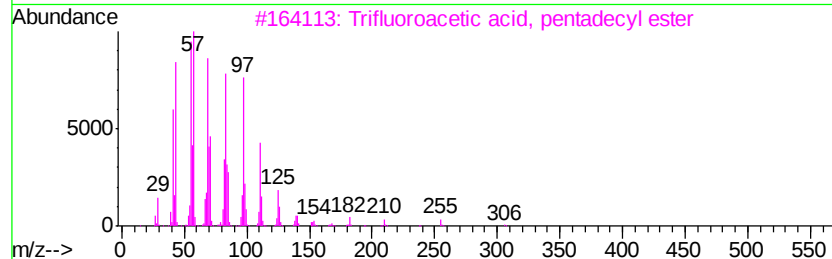
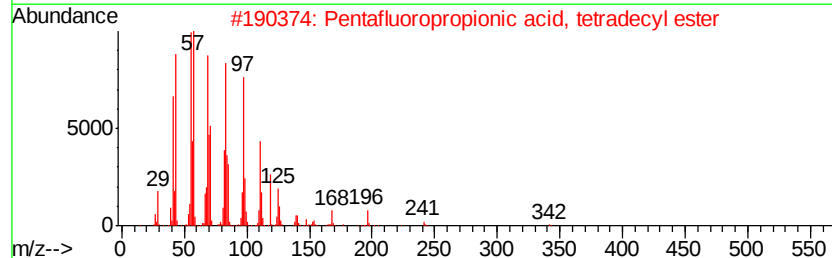
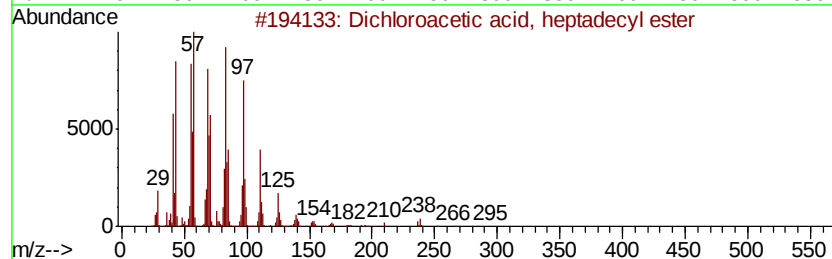
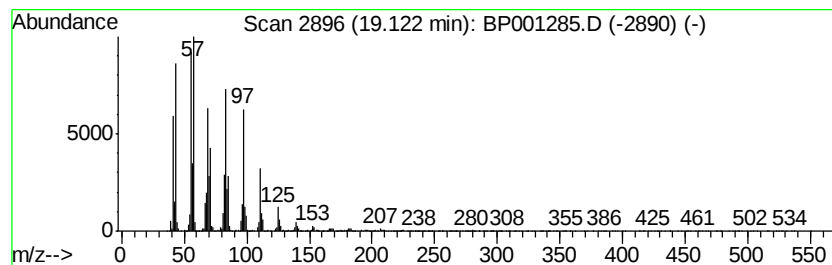
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Peak Number 5 Dichloroacetic acid, heptad... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.12	17.50 ng	596603	Chrysene-d12	19.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
2		Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	94
3		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	94
4		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	94
5		Carbonic acid, hexadecyl 2,2,2-t...	416	C19H35Cl3O3	1000314-56-2	91



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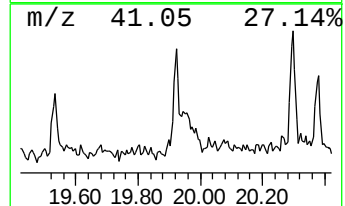
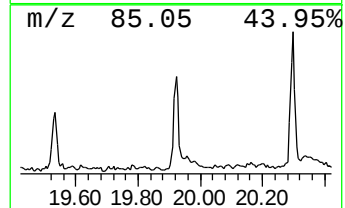
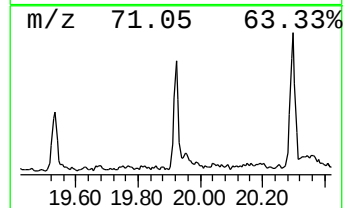
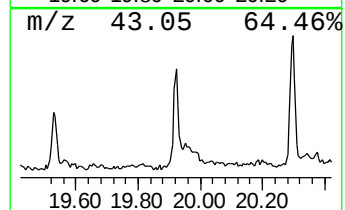
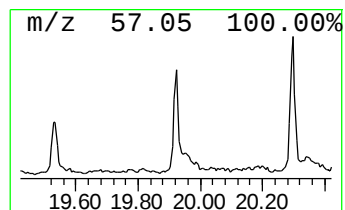
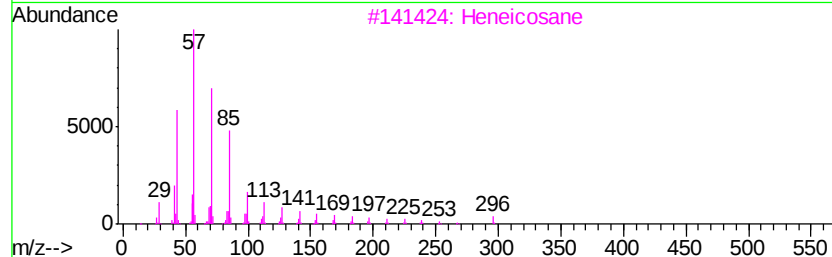
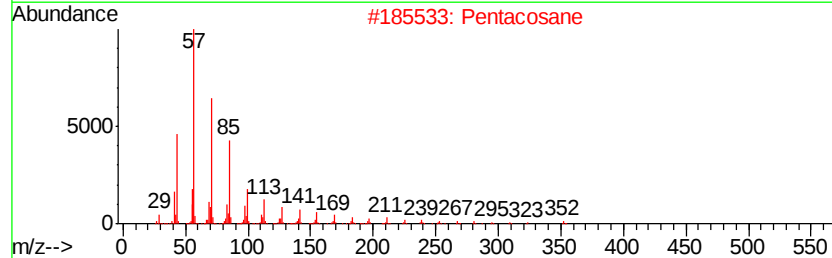
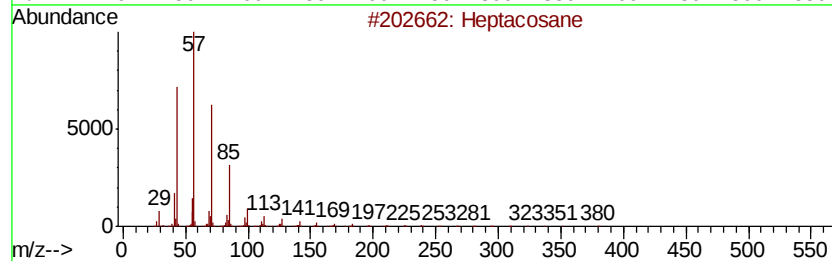
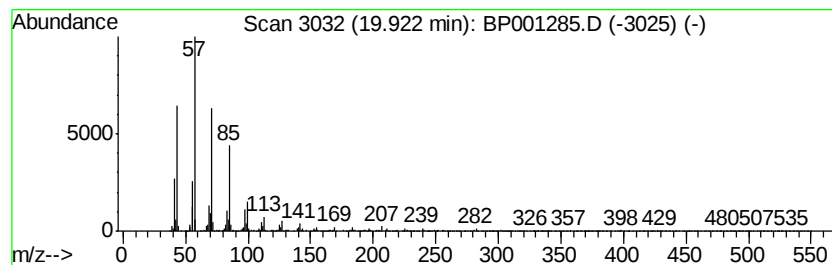
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Peak Number 6 Heptacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.92	2.47 ng	84047	Chrysene-d12	19.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	98
2		Pentacosane	352	C25H52	000629-99-2	97
3		Heneicosane	296	C21H44	000629-94-7	95
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
5		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	95



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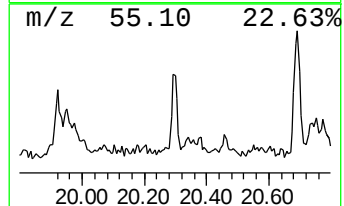
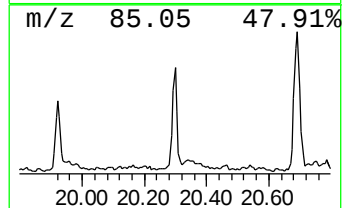
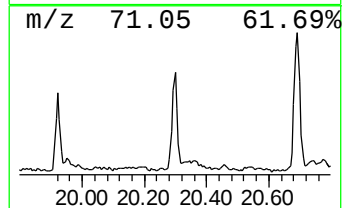
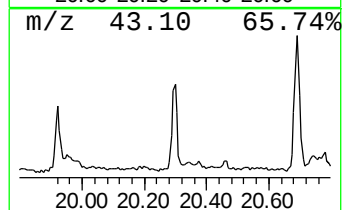
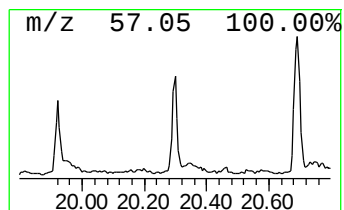
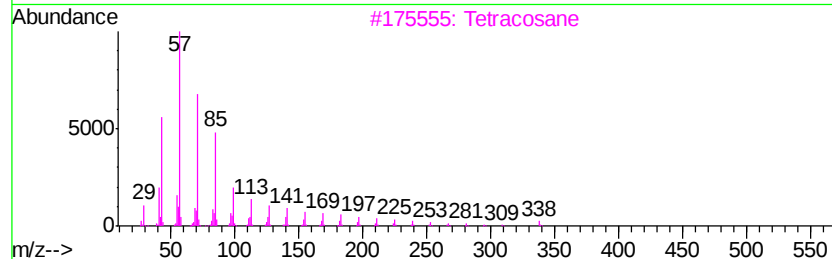
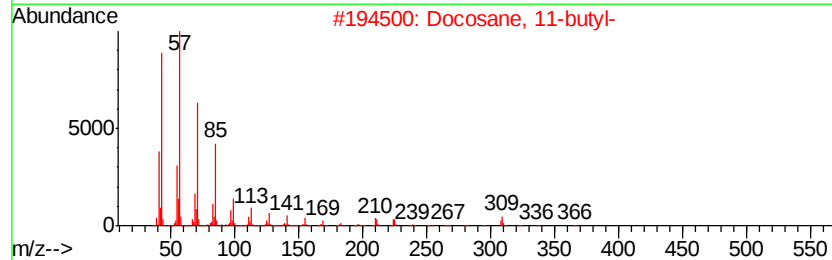
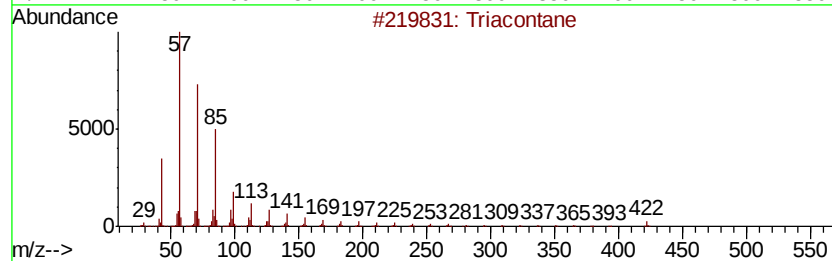
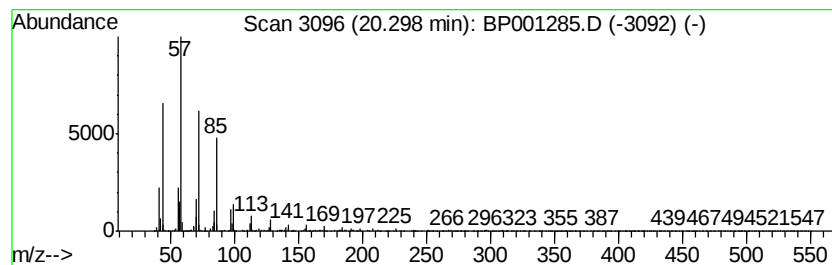
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 7 Triacontane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.30	2.62 ng	87638	Perylene-d12	21.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triacontane	422	C30H62	000638-68-6	94
2		Docosane, 11-butyl-	366	C26H54	013475-76-8	93
3		Tetracosane	338	C24H50	000646-31-1	91
4		Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120919\
Data File : BP001285.D
Acq On : 09 Dec 2019 22:44
Operator : JU
Sample : K6188-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

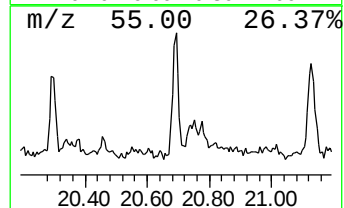
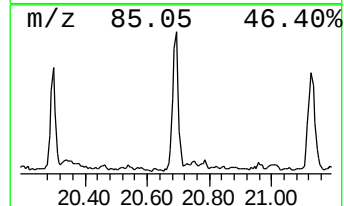
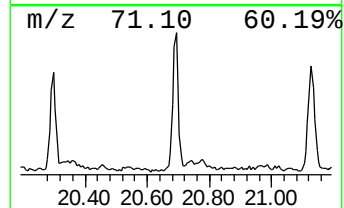
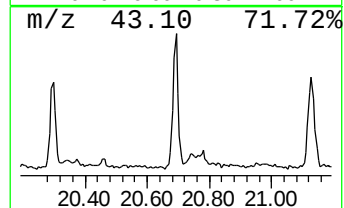
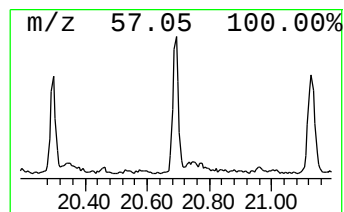
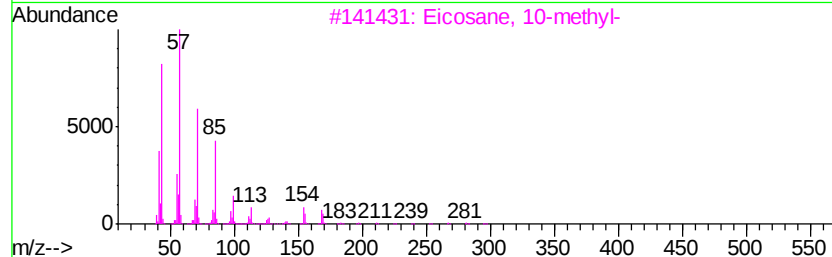
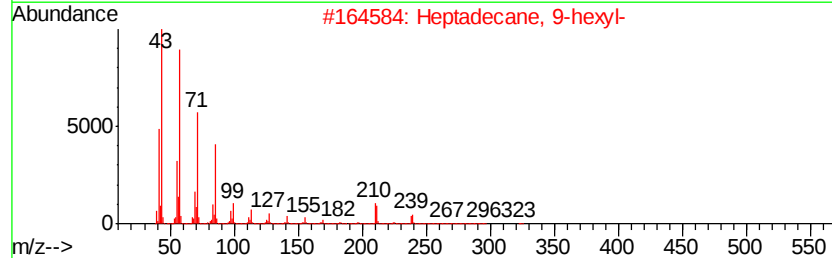
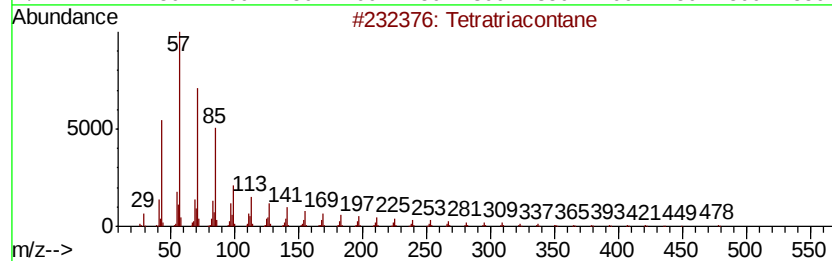
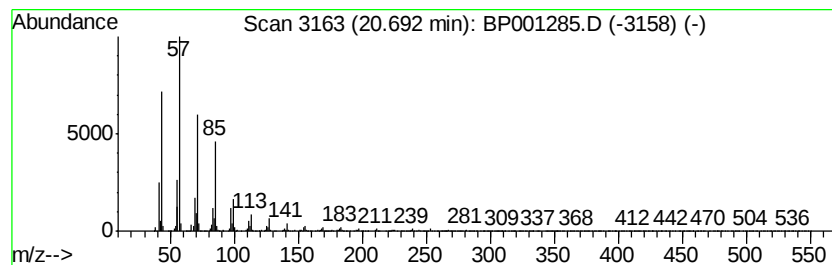
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Tetratriacontane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.69	4.31 ng	144088	Perylene-d12	21.01
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tetratriacontane	479 C34H70	014167-59-0 95
2		Heptadecane, 9-hexyl-	324 C23H48	055124-79-3 94
3		Eicosane, 10-methyl-	296 C21H44	054833-23-7 94
4		2-Bromo dodecane	248 C12H25Br	013187-99-0 93
5		Pentadecane, 8-hexyl-	296 C21H44	013475-75-7 93



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120919\
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Sample : K6188-01
Misc :
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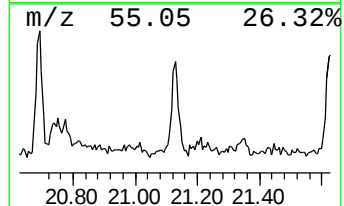
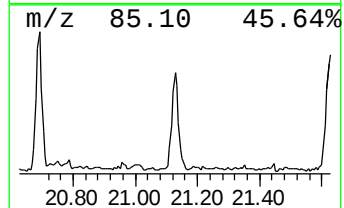
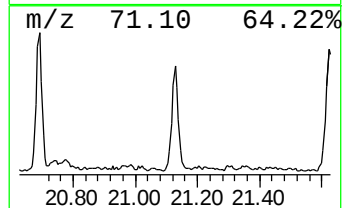
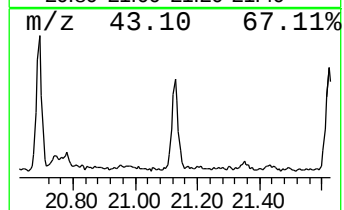
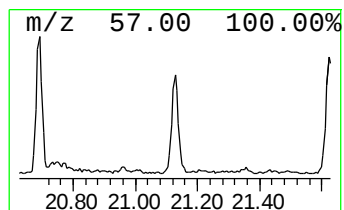
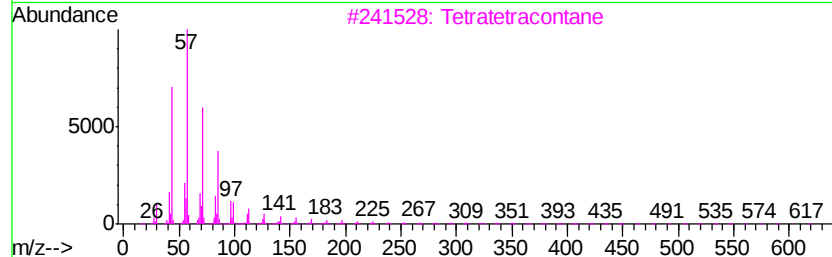
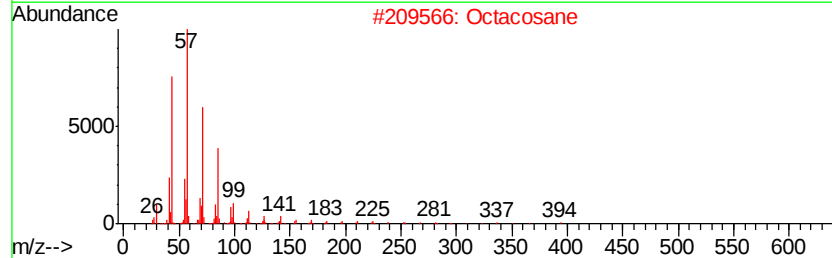
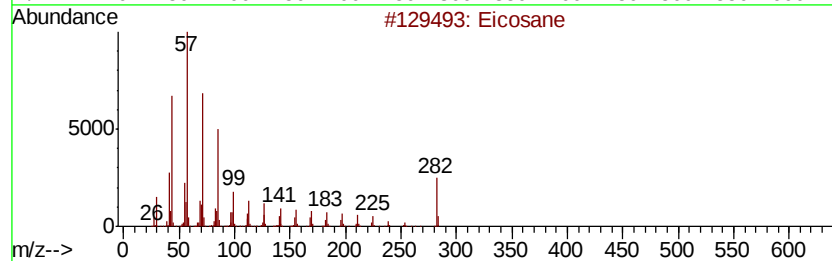
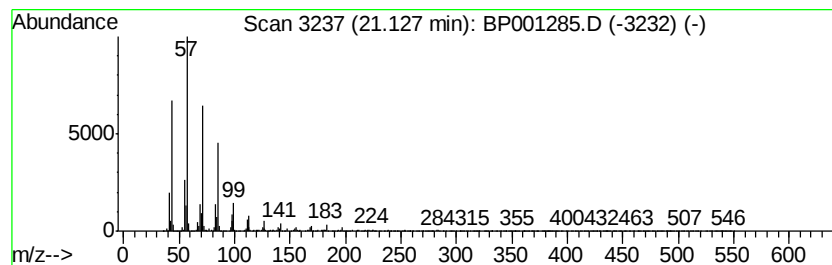
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.13	3.63 ng	121152	Perylene-d12		21.01	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	93
2	Octacosane		394	C28H58	000630-02-4	91
3	Tetratetracontane		619	C44H90	007098-22-8	91
4	Heptacosane		380	C27H56	000593-49-7	91
5	Tetracosane		338	C24H50	000646-31-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120919\
Data File : BP001285.D
Acq On : 09 Dec 2019 22:44
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Misc :
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Instrument :
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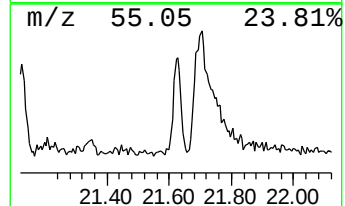
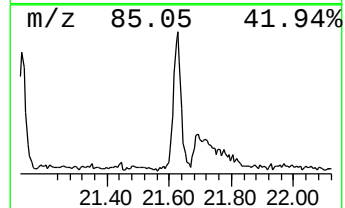
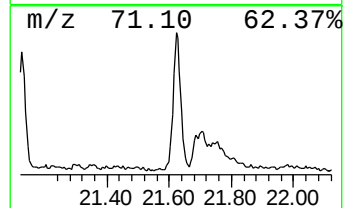
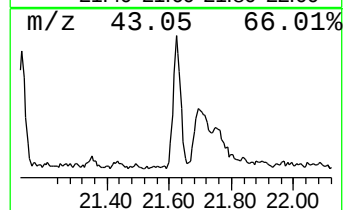
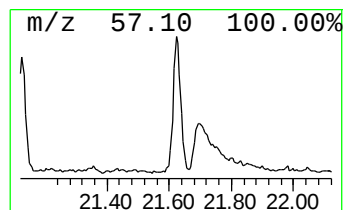
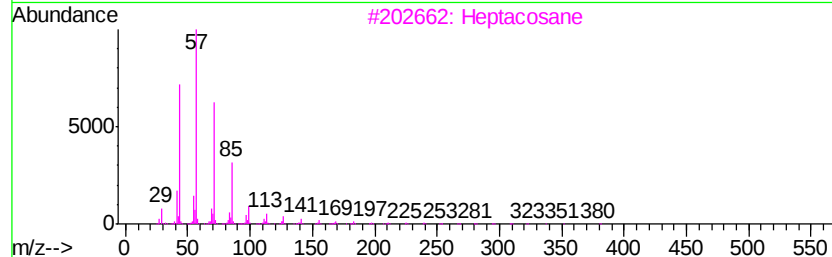
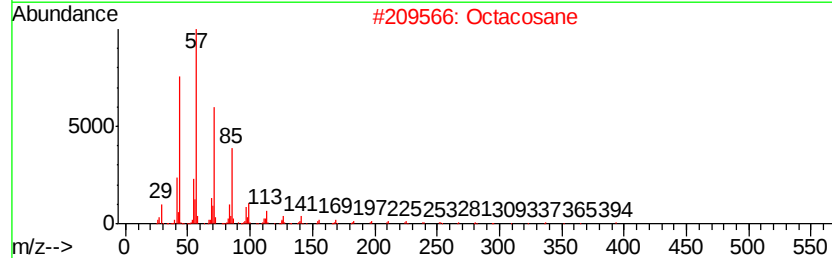
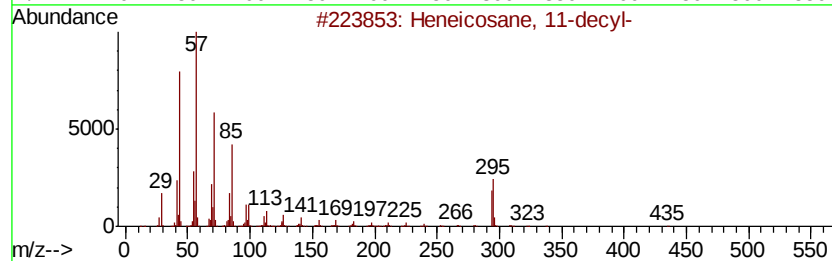
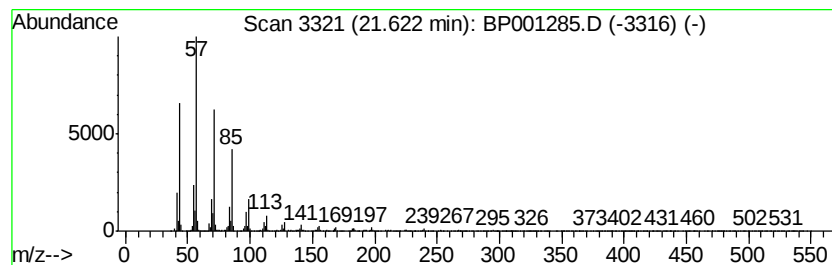
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Heneicosane, 11-decyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.62	4.76 ng	159164	Perylene-d12	21.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane, 11-decyl-	437	C31H64	055320-06-4	92
2			Octacosane	394	C28H58	000630-02-4	91
3			Heptacosane	380	C27H56	000593-49-7	91
4			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91
5			Pentacosane	352	C25H52	000629-99-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120919\
Data File : BP001285.D
Acq On : 09 Dec 2019 22:44
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Instrument :
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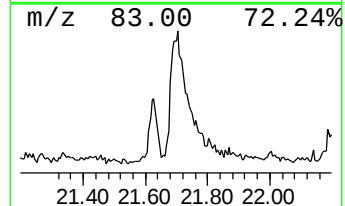
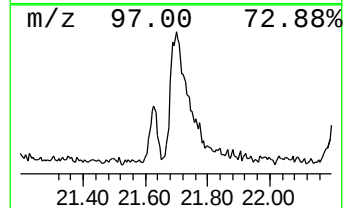
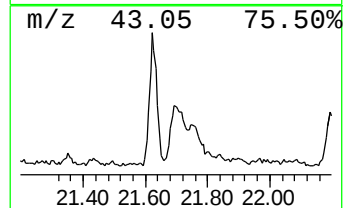
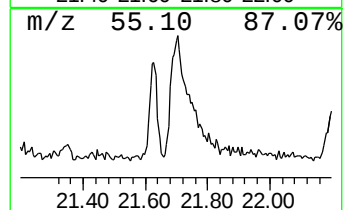
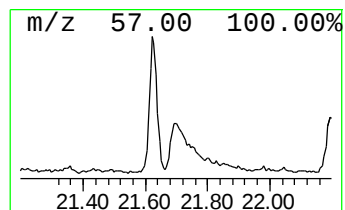
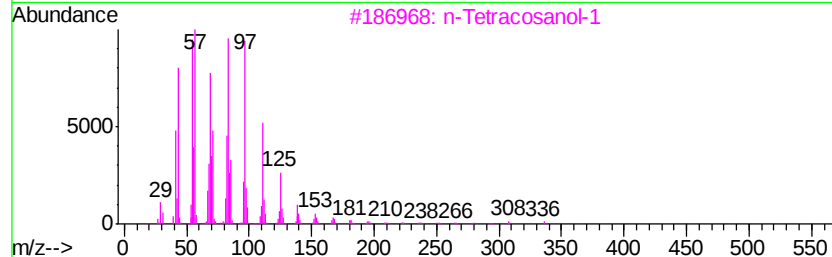
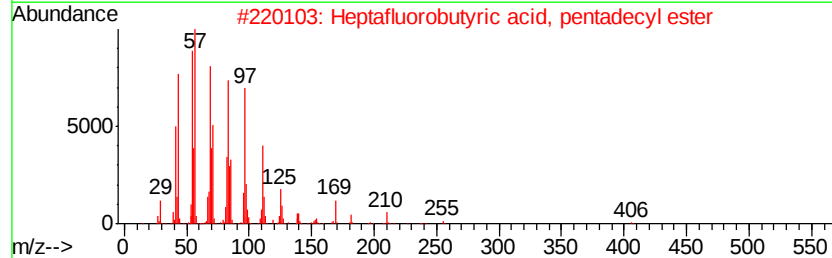
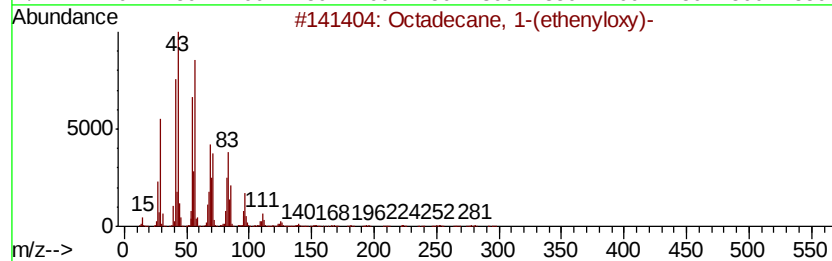
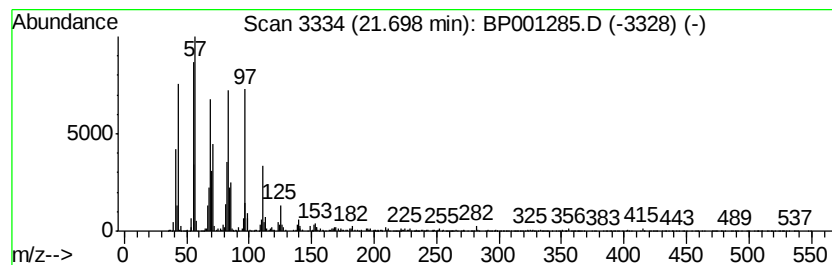
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Octadecane, 1-(ethenyloxy)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.70	6.00 ng	200583	Perylene-d12	21.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane, 1-(ethenyloxy)-	296	C20H40O	000930-02-9	93
2			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	90
3			n-Tetracosanol-1	354	C24H50O	000506-51-4	90
4			E-7-Octadecene	252	C18H36	1000130-92-0	89
5			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120919\
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Acq On : 09 Dec 2019 22:44
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Sample : K6188-01
Misc :
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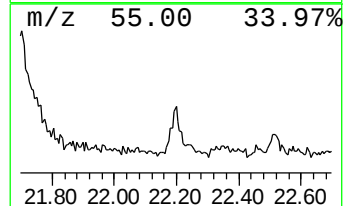
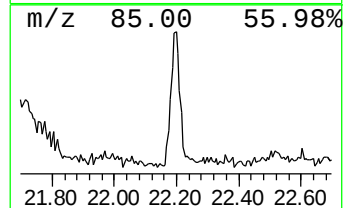
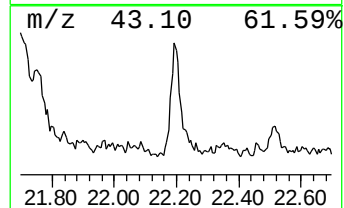
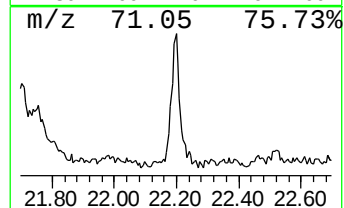
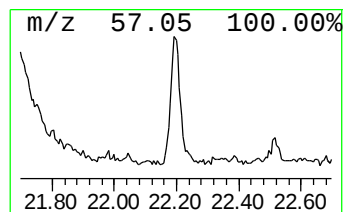
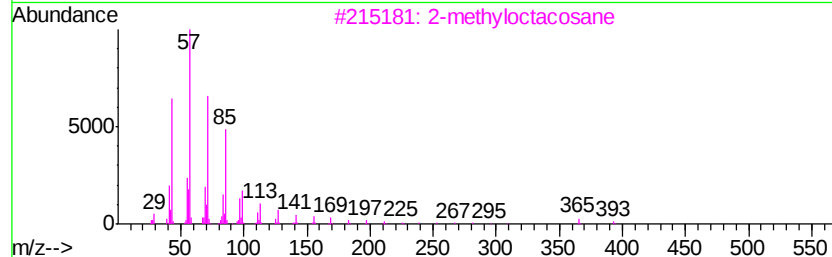
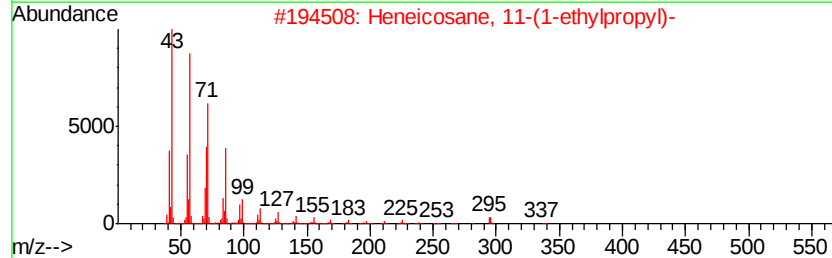
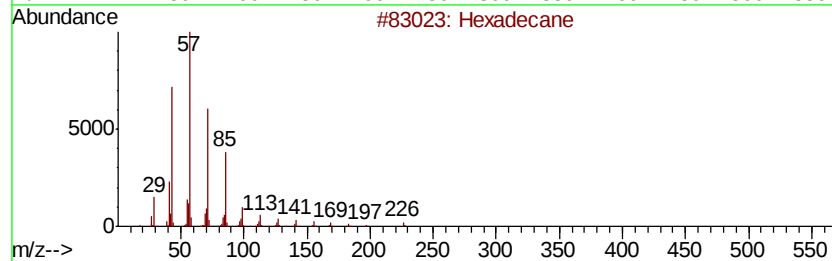
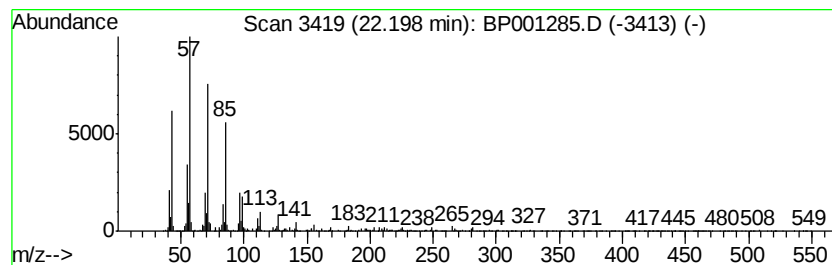
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Hexadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.20	2.37 ng	79263	Perylene-d12	21.01
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226 C16H34	000544-76-3	87
2	Heneicosane, 11-(1-ethylpropyl)-	366 C26H54	055282-11-6	87
3	2-methyloctacosane	408 C29H60	1000376-72-8	87
4	Octacosane	394 C28H58	000630-02-4	83
5	Eicosane, 7-hexyl-	366 C26H54	055333-99-8	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120919\
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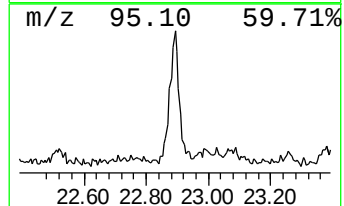
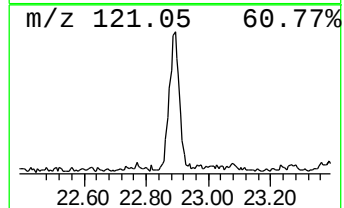
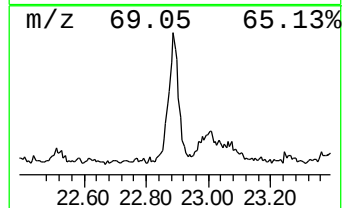
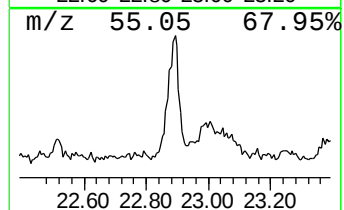
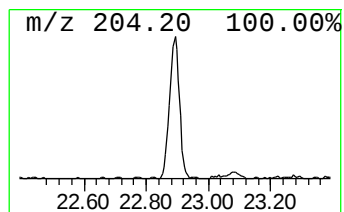
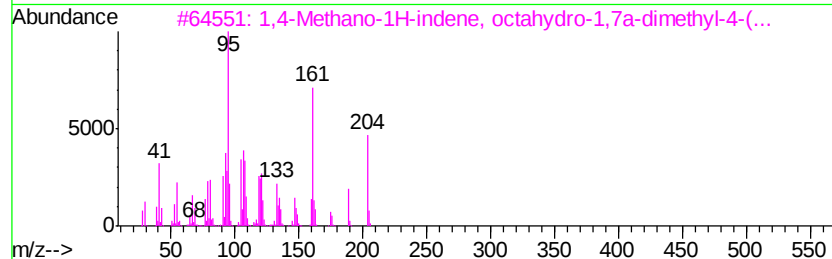
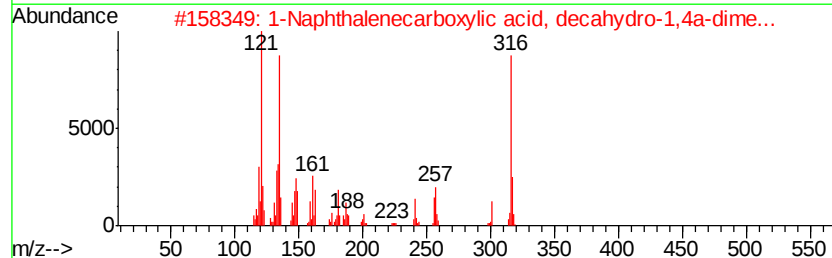
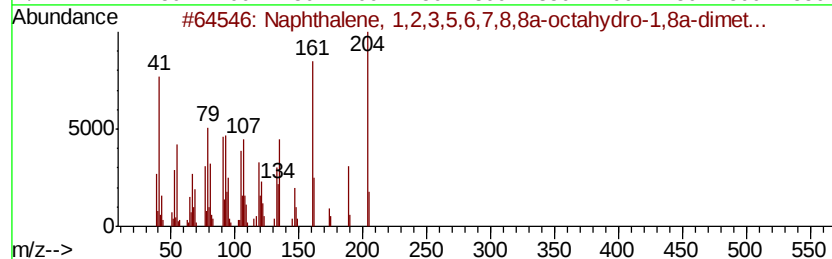
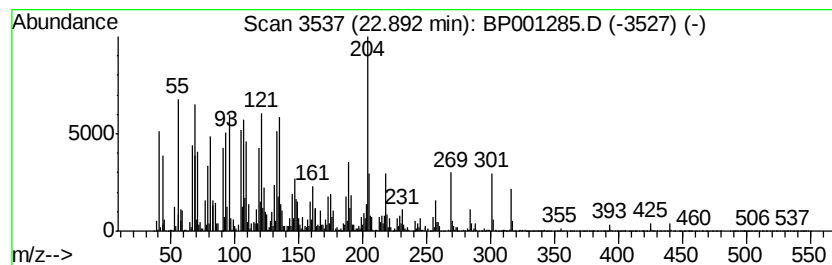
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Naphthalene, 1,2,3,5,6,7,8,... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.89	14.90 ng	497830	Perylene-d12	21.01	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	004630-07-3	58
2	1-Naphthalenecarboxylic acid, de...	316	C21H32O2	001235-39-8	55
3	1,4-Methano-1H-indene, octahydro...	204	C15H24	087064-18-4	49
4	Aromandendrene	204	C15H24	000489-39-4	46
5	1H-Cycloprop[e]azulene, 1a,2,3,5...	204	C15H24	021747-46-6	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP120919\
 Data File : BP001285.D
 Acq On : 09 Dec 2019 22:44
 Operator : JU
 Sample : K6188-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SOIL-COMP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.63	28.3	ng	684747	1	8.30	483159	20.0
unknown7.95	7.95	122.1	ng	2950140	1	8.30	483159	20.0
n-Hexadecanoic acid	16.49	5.3	ng	189962	4	15.60	718869	20.0
Dichloroacetic ac...	19.12	17.5	ng	596603	5	19.24	681884	20.0
Heptacosane	19.92	2.5	ng	84047	5	19.24	681884	20.0
Triacontane	20.30	2.6	ng	87638	6	21.01	668120	20.0
Tetratriacontane	20.69	4.3	ng	144088	6	21.01	668120	20.0
Eicosane	21.13	3.6	ng	121152	6	21.01	668120	20.0
Heneicosane, 11-d...	21.62	4.8	ng	159164	6	21.01	668120	20.0
Octadecane, 1-(et...	21.70	6.0	ng	200583	6	21.01	668120	20.0
Hexadecane	22.20	2.4	ng	79263	6	21.01	668120	20.0
Naphthalene, 1,2,...	22.89	14.9	ng	497830	6	21.01	668120	20.0