

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
 Data File : BP008822.D
 Acq On : 17 Jan 2022 14:14
 Operator : CG/JU
 Sample : N1143-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 WEST-BOILER-ROOM-NORTH-SIDEWALL

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-BP011422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP008822.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.058	7	12	26	rVB	2496699	3195285	3.75%	0.236%
2	4.899	320	325	330	rBV	531711	989540	1.16%	0.073%
3	4.958	330	335	340	rVV2	1550330	2846459	3.34%	0.211%
4	5.064	348	353	359	rBV2	1226468	2270035	2.66%	0.168%
5	5.222	375	380	384	rBV	2067003	3593350	4.22%	0.266%
6	5.311	384	395	404	rVB3	8286247	22368420	26.25%	1.655%
7	5.446	409	418	430	rBV2	11145469	26278864	30.84%	1.944%
8	5.552	430	436	446	rVB3	2339243	5371439	6.30%	0.397%
9	5.664	446	455	462	rBV2	6383036	14860526	17.44%	1.099%
10	5.740	462	468	472	rBV5	6837390	15942411	18.71%	1.179%
11	5.828	477	483	488	rBV	11339604	20214570	23.72%	1.495%
12	5.887	488	493	501	rVB2	9810370	21780667	25.56%	1.611%
13	5.964	501	506	515	rBV3	3347688	6828787	8.01%	0.505%
14	6.058	515	522	529	rBV2	4240235	10507696	12.33%	0.777%
15	6.152	529	538	551	rBV8	18845549	85213506	100.00%	6.304%
16	6.281	553	560	568	rVB2	14308467	45879747	53.84%	3.394%
17	6.375	568	576	582	rBV4	16448139	46945935	55.09%	3.473%
18	6.463	582	591	595	rBV	17331655	51048491	59.91%	3.776%
19	6.534	595	603	607	rBV4	9661122	31636616	37.13%	2.340%
20	6.646	616	622	627	rVB2	12648459	25740050	30.21%	1.904%
21	6.705	627	632	637	rVB3	10281633	18037448	21.17%	1.334%
22	6.787	637	646	653	rBV6	14789166	52725193	61.87%	3.900%
23	6.905	653	666	674	rVB6	19986435	80506696	94.48%	5.955%
24	6.993	674	681	684	rBV3	15697143	34405681	40.38%	2.545%
25	7.034	684	688	699	rVB7	11861314	38420965	45.09%	2.842%
26	7.222	713	720	724	rBV4	10463344	25244412	29.62%	1.867%
27	7.275	725	729	732	rVB2	7695336	12069163	14.16%	0.893%
28	7.316	733	736	740	rBV3	4328820	7509011	8.81%	0.555%
29	7.393	741	749	753	rBV	9515380	26455537	31.05%	1.957%
30	7.581	771	781	784	rBV5	10931171	36629024	42.98%	2.710%
31	7.663	791	795	798	rBV	7416476	10249937	12.03%	0.758%
32	7.710	798	803	805	rBV4	8141607	15562804	18.26%	1.151%
33	7.799	812	818	819	rBV3	8776244	15057656	17.67%	1.114%
34	7.993	842	851	855	rBV3	12344397	38058717	44.66%	2.815%
35	8.105	864	870	875	rVB3	8393201	18527199	21.74%	1.371%
36	8.175	877	882	883	rBV2	9379632	12412075	14.57%	0.918%

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37	8.257	892	896	903	rVB3	8500437	17330072	20.34%	1.282%
38	8.340	905	910	914	rBV3	8312526	15415747	18.09%	1.140%
39	8.487	932	935	939	rVB	6442253	9003387	10.57%	0.666%
40	8.528	939	942	946	rVB2	4146001	5774537	6.78%	0.427%
41	8.581	946	951	957	rBV4	8049933	14105732	16.55%	1.043%
42	8.728	971	976	991	rVB5	15171375	47873469	56.18%	3.541%
43	8.834	991	994	997	rBV	1556084	1689345	1.98%	0.125%
44	8.887	997	1003	1011	rVB5	4108920	9935180	11.66%	0.735%
45	8.993	1011	1021	1031	rBV5	12657894	40198281	47.17%	2.974%
46	9.087	1032	1037	1044	rVB5	9101685	16956686	19.90%	1.254%
47	9.152	1044	1048	1052	rVB2	5612818	8861467	10.40%	0.656%
48	9.205	1052	1057	1059	rBV2	8321045	14996383	17.60%	1.109%
49	9.246	1061	1064	1069	rVB2	9614021	15593544	18.30%	1.154%
50	9.316	1072	1076	1078	rBV3	2996296	4572618	5.37%	0.338%
51	9.352	1078	1082	1091	rVB4	6388292	17130624	20.10%	1.267%
52	9.446	1092	1098	1105	rBV4	4068669	11631735	13.65%	0.860%
53	9.528	1105	1112	1113	rBV2	5102857	9969356	11.70%	0.737%
54	9.646	1129	1132	1137	rVB	2298045	3028576	3.55%	0.224%
55	9.757	1144	1151	1155	rBV3	5625865	11104606	13.03%	0.821%
56	9.799	1155	1158	1163	rVV	4705345	7364255	8.64%	0.545%
57	9.857	1163	1168	1176	rVV5	4985065	11461819	13.45%	0.848%
58	9.952	1176	1184	1189	rVB2	3225039	8436015	9.90%	0.624%
59	10.069	1199	1204	1209	rBV2	1490000	2771193	3.25%	0.205%
60	10.134	1210	1215	1222	rVB3	2011346	4021670	4.72%	0.298%
61	10.204	1223	1227	1232	rBV2	1034663	2048423	2.40%	0.152%
62	10.257	1233	1236	1240	rVB5	890766	1237340	1.45%	0.092%
63	10.369	1251	1255	1258	rVV	838213	1134792	1.33%	0.084%
64	10.451	1265	1269	1274	rVB4	610301	1067073	1.25%	0.079%
65	10.540	1279	1284	1293	rBV5	1336160	4154203	4.88%	0.307%
66	10.657	1296	1304	1309	rVB2	2494998	5827180	6.84%	0.431%
67	10.710	1309	1313	1315	rBV4	738904	1110995	1.30%	0.082%
68	10.787	1321	1326	1330	rBV4	1315356	2104122	2.47%	0.156%
69	10.834	1330	1334	1339	rVB	3176162	4699240	5.51%	0.348%
70	10.893	1340	1344	1352	rVB4	971948	1766689	2.07%	0.131%
71	10.963	1352	1356	1362	rVB4	1141511	2087627	2.45%	0.154%
72	11.063	1366	1373	1382	rBV7	542187	1569356	1.84%	0.116%
73	11.369	1421	1425	1429	rVV	878929	1478039	1.73%	0.109%
74	11.434	1433	1436	1443	rVB2	589303	1073483	1.26%	0.079%
75	11.510	1443	1449	1455	rVB2	632108	1097807	1.29%	0.081%
76	11.581	1455	1461	1469	rVB8	412117	1225968	1.44%	0.091%

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77	11.698	1475	1481	1485	rBV	1613919	2770002	3.25%	0.205%
78	12.310	1578	1585	1591	rVB4	391533	1103550	1.30%	0.082%
79	13.040	1705	1709	1714	rVB	794115	1103878	1.30%	0.082%
80	13.116	1715	1722	1729	rVB	8625942	12494981	14.66%	0.924%
81	13.987	1862	1870	1874	rVV	1036293	1748958	2.05%	0.129%
82	14.492	1948	1956	1966	rVV	2801773	5086561	5.97%	0.376%
83	15.028	2042	2047	2052	rBV2	764678	1144803	1.34%	0.085%
84	15.504	2122	2128	2133	rVV	10810258	14755337	17.32%	1.092%
85	15.763	2165	2172	2177	rBV	2045717	3546406	4.16%	0.262%
86	15.987	2203	2210	2218	rVB3	6224059	10366402	12.17%	0.767%
87	16.245	2246	2254	2259	rBV	3836717	7576802	8.89%	0.560%
88	17.075	2390	2395	2406	rVB	3205873	5400525	6.34%	0.400%
89	17.251	2420	2425	2430	rBV	2943229	4495763	5.28%	0.333%
90	17.681	2494	2498	2503	rVB	711158	1059110	1.24%	0.078%
91	19.622	2825	2828	2836	rVB	587535	1036784	1.22%	0.077%
92	19.810	2855	2860	2867	rVB	12222406	15777442	18.52%	1.167%
93	19.969	2883	2887	2893	rVB	4104345	4650663	5.46%	0.344%
94	20.216	2926	2929	2934	rVB	1003346	1176100	1.38%	0.087%
95	21.057	3068	3072	3081	rVB	904370	1252711	1.47%	0.093%
96	21.257	3102	3106	3114	rVB	7355757	8621367	10.12%	0.638%
97	21.339	3116	3120	3128	rVB	2761744	3715201	4.36%	0.275%
98	21.463	3136	3141	3146	rBV	5027744	7836665	9.20%	0.580%
99	23.768	3527	3533	3546	rVB	1692033	3529182	4.14%	0.261%
100	25.280	3785	3790	3801	rVB2	1362181	3275333	3.84%	0.242%

Sum of corrected areas: 1351817072

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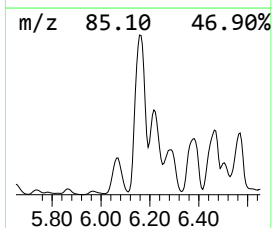
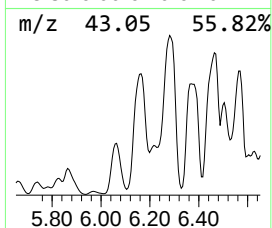
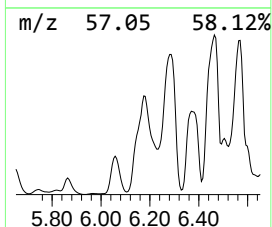
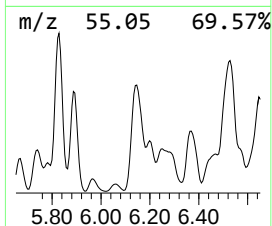
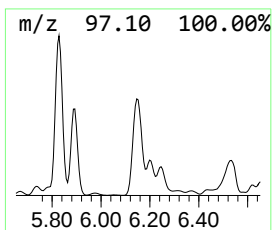
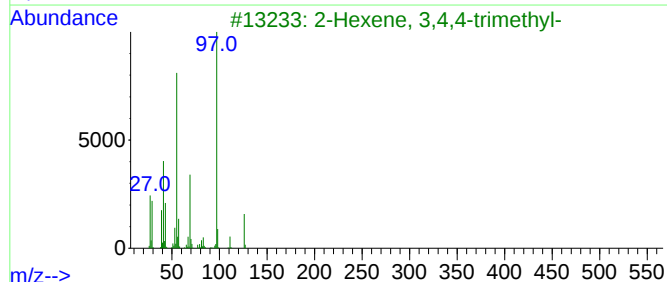
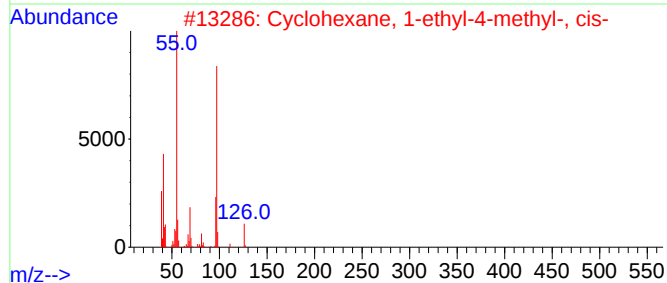
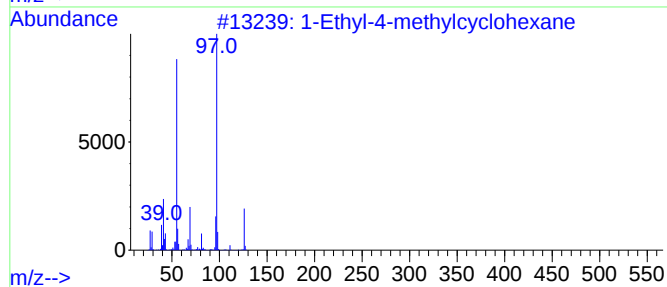
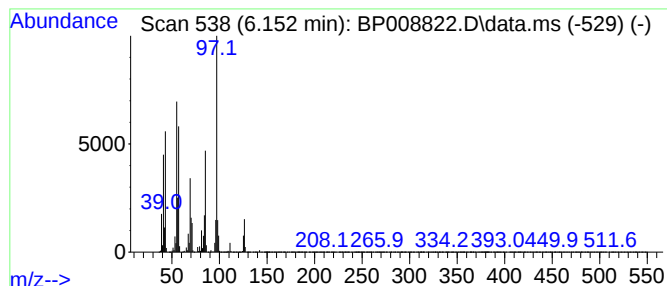
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.152 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.152	113.18 ng	85213500	1,4-Dichlorobenzene-d4	7.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	46
2			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	46
3			2-Hexene, 3,4,4-trimethyl-	126	C9H18	053941-19-8	46
4			Thiophene, 2-propyl-	126	C7H10S	001551-27-5	46
5			Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	46



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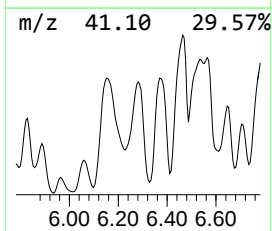
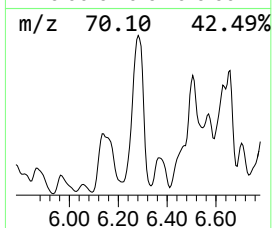
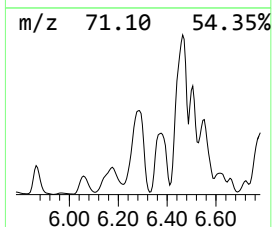
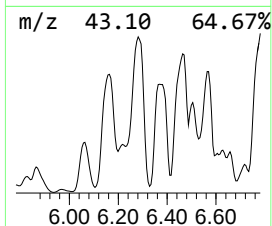
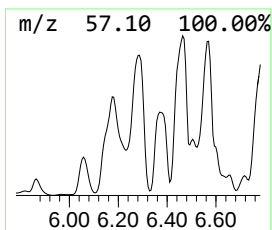
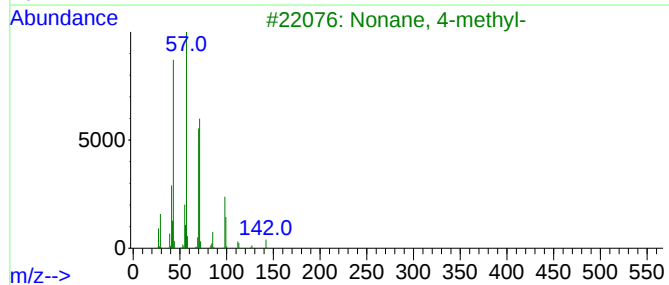
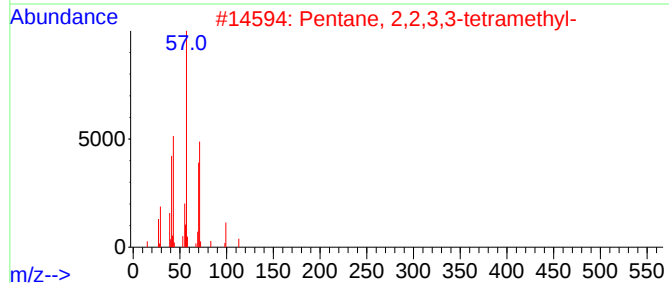
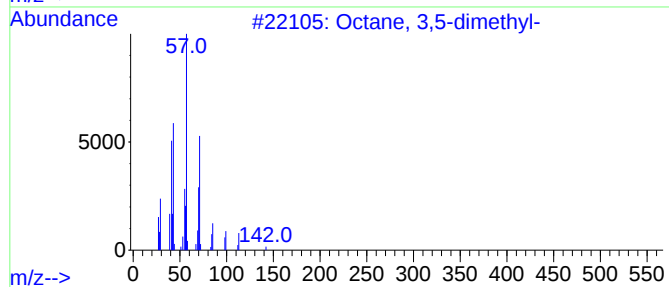
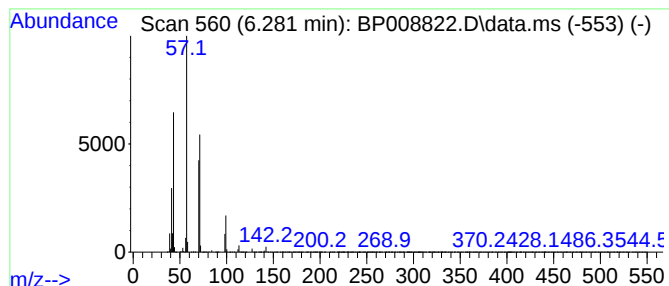
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Octane, 3,5-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.281	60.94 ng	45879700	1,4-Dichlorobenzene-d4	7.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	78
2			Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	72
3			Nonane, 4-methyl-	142	C10H22	017301-94-9	72
4			Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	59
5			Hexane, 3-methyl-	100	C7H16	000589-34-4	53



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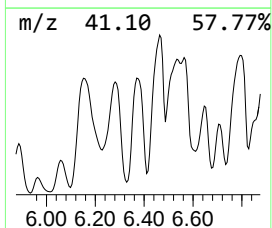
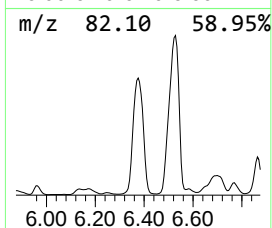
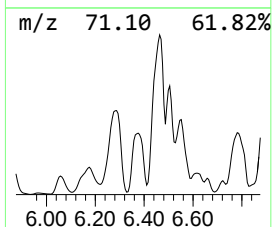
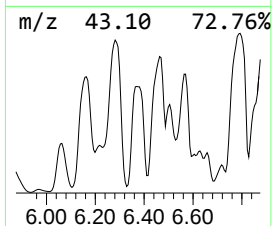
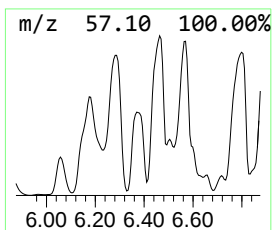
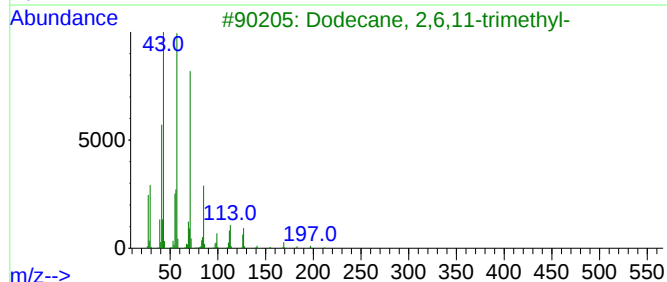
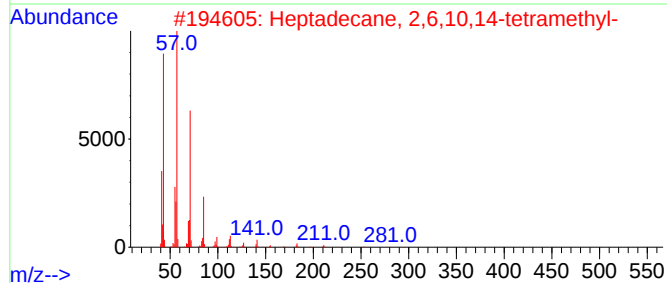
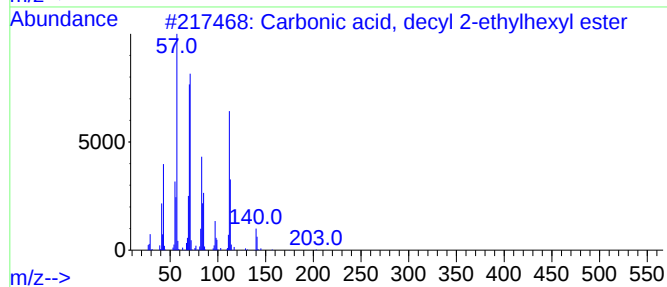
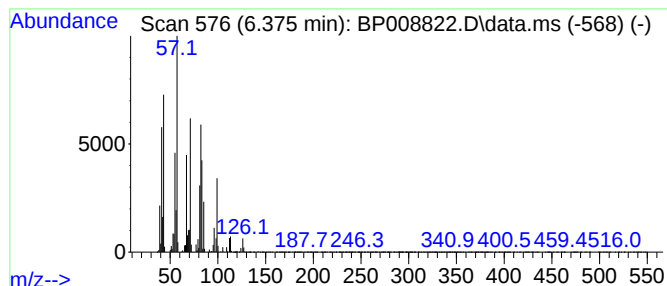
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.375 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.375	62.35 ng	46945900	1,4-Dichlorobenzene-d4	7.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, decyl 2-ethylhexy...	314	C19H38O3	1000383-13-7	47
2			Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	27
3			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	27
4			Octane, 2,7-dimethyl-	142	C10H22	001072-16-8	27
5			1,5,7-Octatrien-3-ol, 3,7-dimethyl-	152	C10H16O	029957-43-5	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

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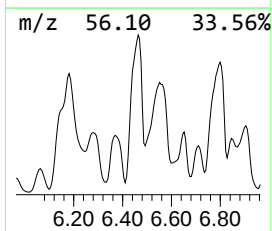
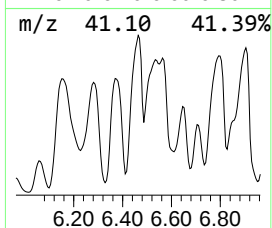
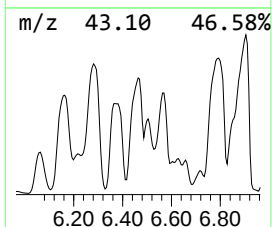
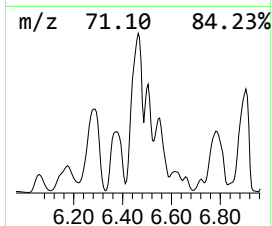
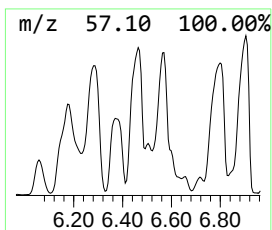
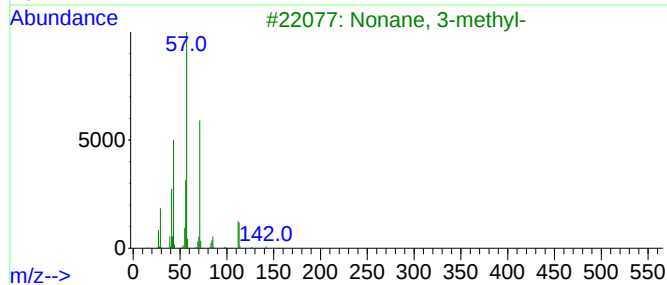
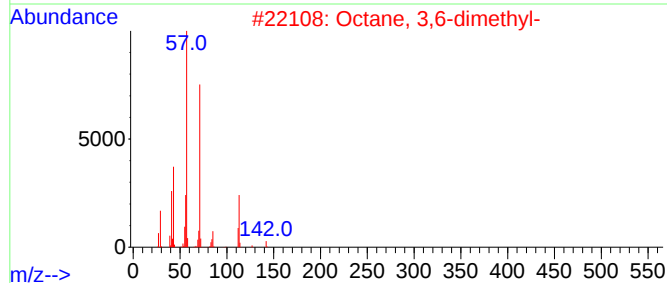
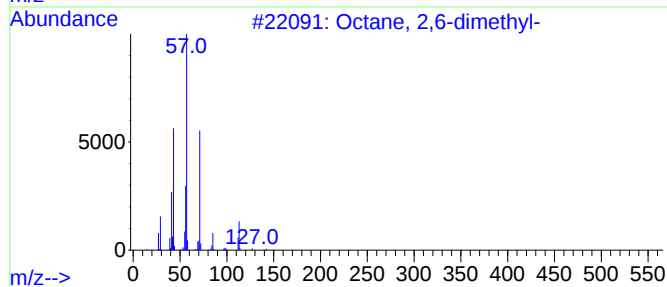
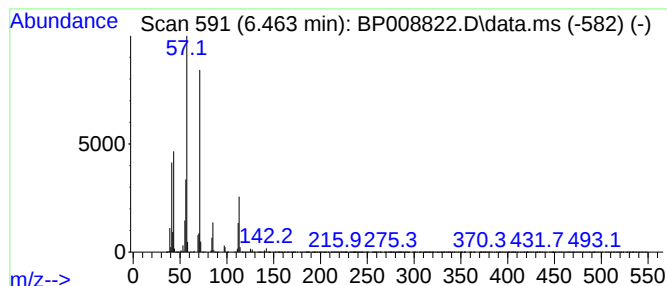
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Octane, 2,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.463	67.80 ng	51048500	1,4-Dichlorobenzene-d4	7.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	95
2			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	91
3			Nonane, 3-methyl-	142	C10H22	005911-04-6	87
4			Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	83
5			Sulfurous acid, di(2-ethylhexyl)...	306	C16H34O3S	1000309-19-1	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

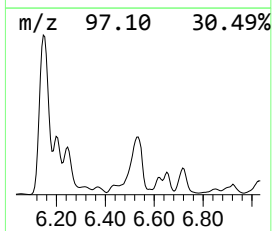
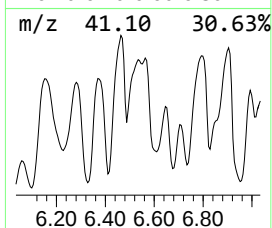
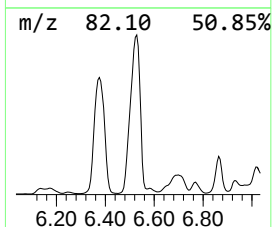
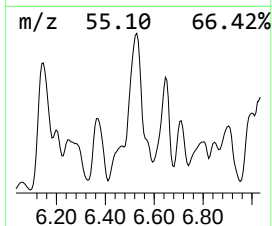
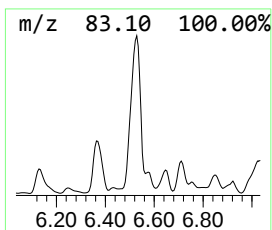
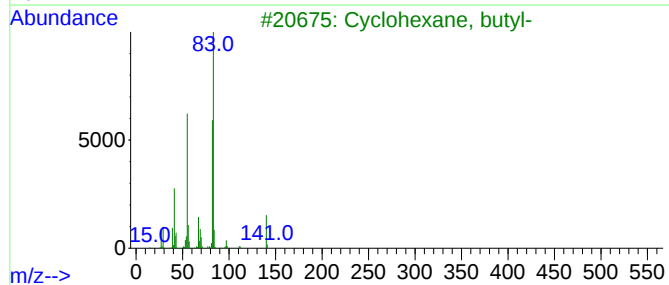
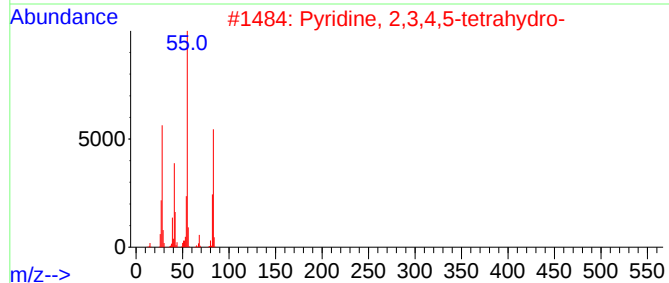
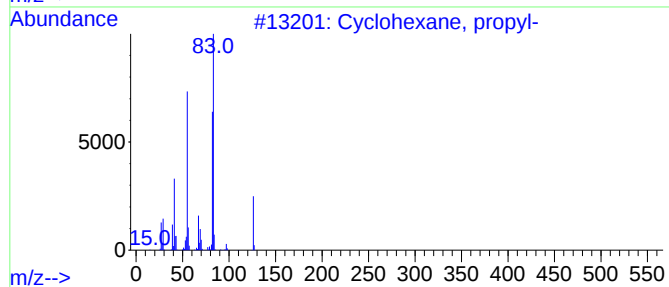
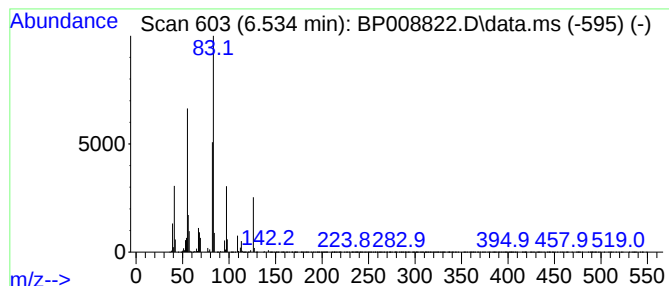
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Cyclohexane, propyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.534	42.02 ng	31636600	1,4-Dichlorobenzene-d4	7.875	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, propyl-	126	C9H18	001678-92-8	68
2	Pyridine, 2,3,4,5-tetrahydro-	83	C5H9N	000505-18-0	53
3	Cyclohexane, butyl-	140	C10H20	001678-93-9	50
4	Ethanone, 1-(1-methylcyclopentyl)-	126	C8H14O	013388-93-7	50
5	Cyclohexane, ethyl-	112	C8H16	001678-91-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
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Instrument :
BNA_P
ClientSampleId :
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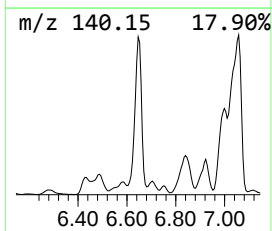
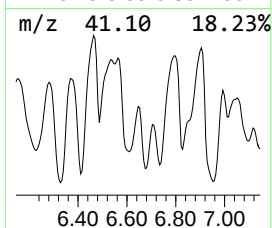
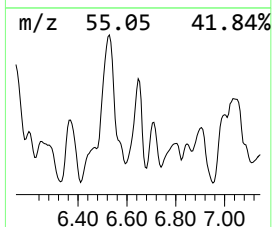
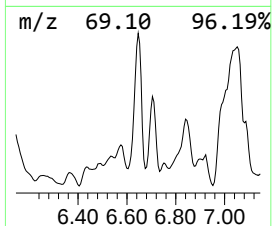
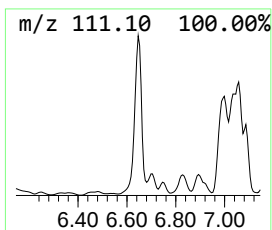
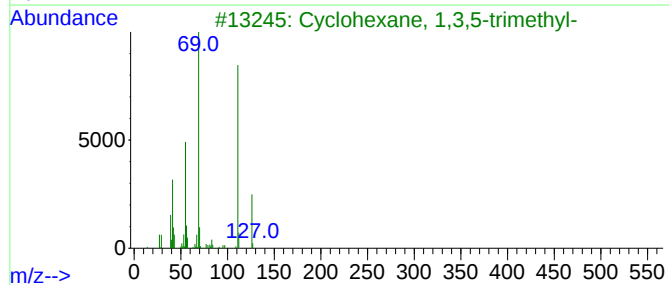
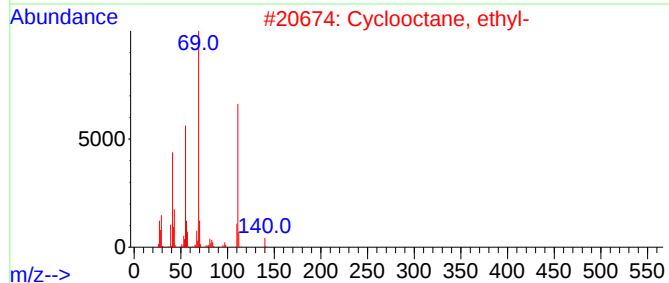
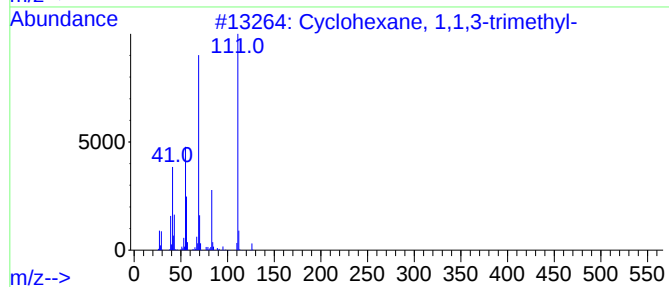
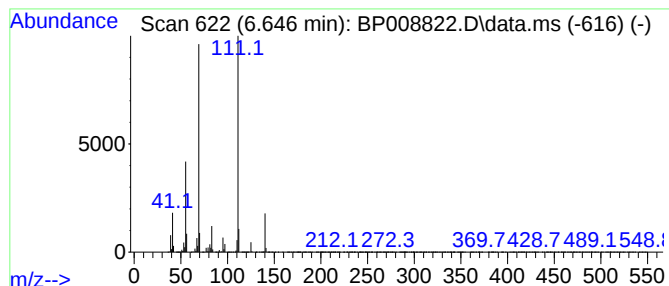
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.646	34.19 ng	25740100	1,4-Dichlorobenzene-d4	7.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	72
2			Cyclooctane, ethyl-	140	C10H20	013152-02-8	72
3			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	72
4			Cyclohexane, 1-ethyl-2,3-dimethyl-	140	C10H20	007058-05-1	64
5			Cyclopentane, 1,1,3-trimethyl-3-...	166	C12H22	074421-09-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

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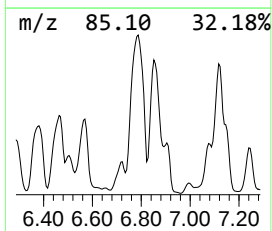
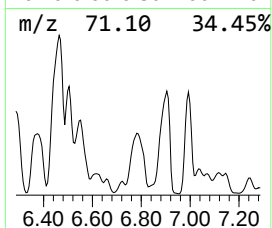
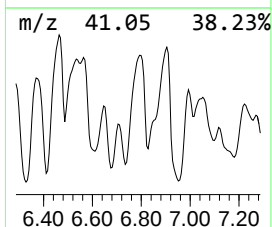
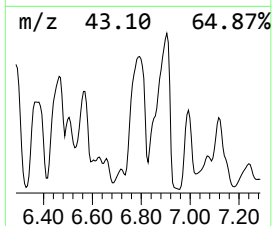
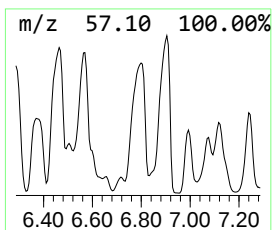
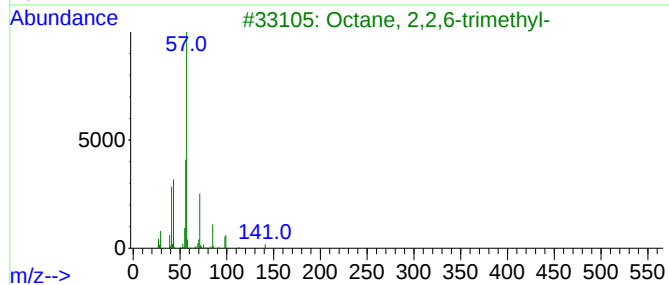
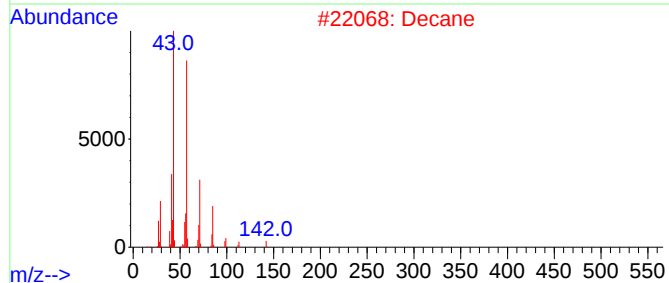
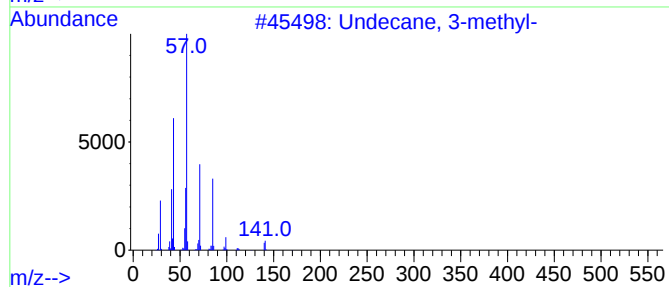
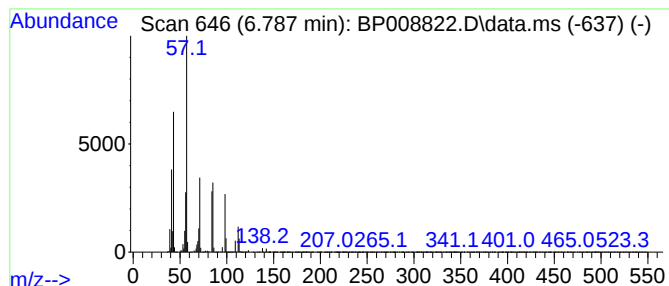
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown6.787 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.787	70.03 ng	52725200	1,4-Dichlorobenzene-d4	7.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 3-methyl-	170	C12H26	001002-43-3	47
2			Decane	142	C10H22	000124-18-5	46
3			Octane, 2,2,6-trimethyl-	156	C11H24	062016-28-8	43
4			Heptane, 3-methyl-	114	C8H18	000589-81-1	38
5			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

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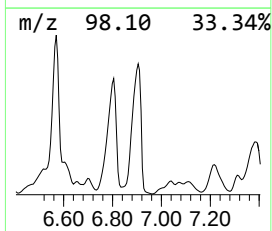
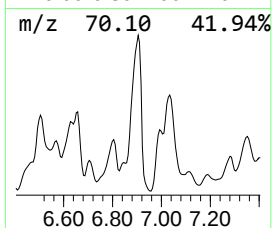
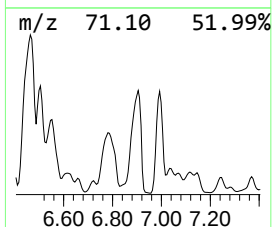
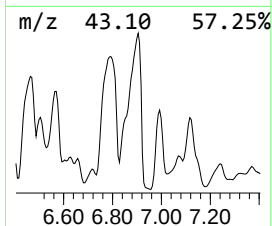
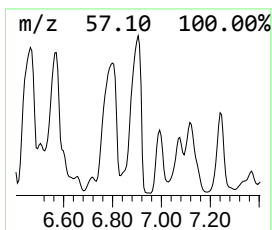
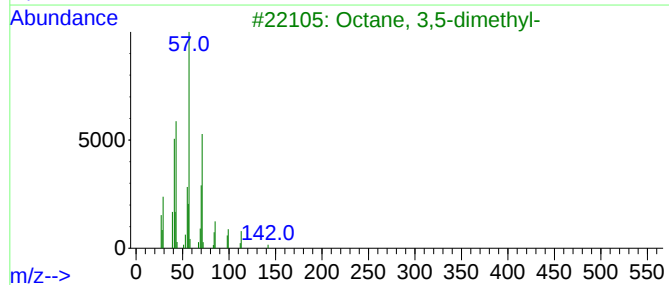
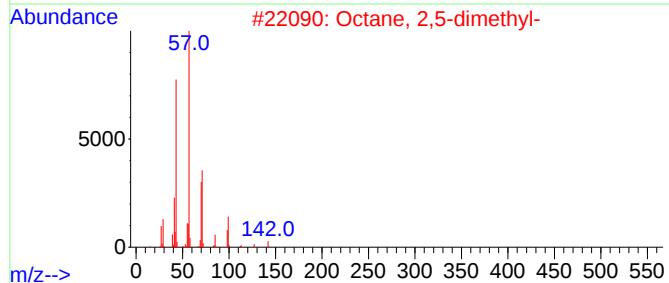
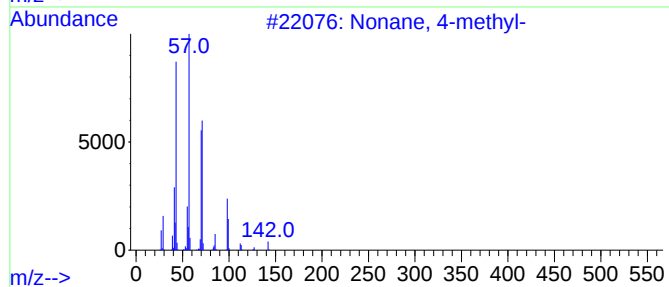
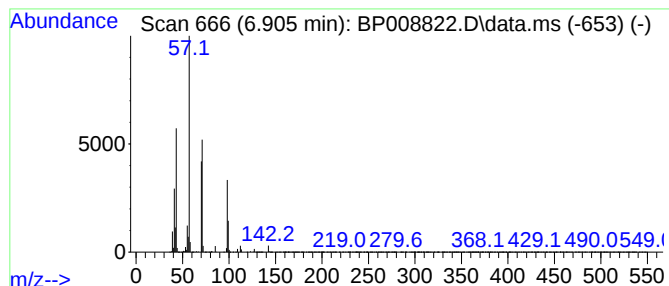
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Nonane, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.905	106.93 ng	80506700	1,4-Dichlorobenzene-d4	7.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 4-methyl-	142	C10H22	017301-94-9	90
2			Octane, 2,5-dimethyl-	142	C10H22	015869-89-3	80
3			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	64
4			Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	50
5			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
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Instrument :
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WEST-BOILER-ROOM-NORTH-SIDEWALL

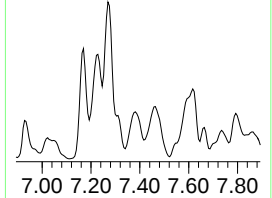
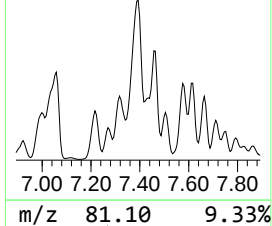
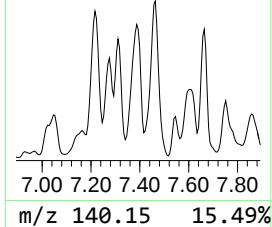
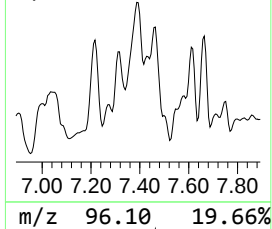
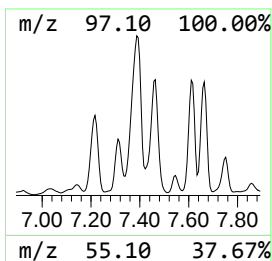
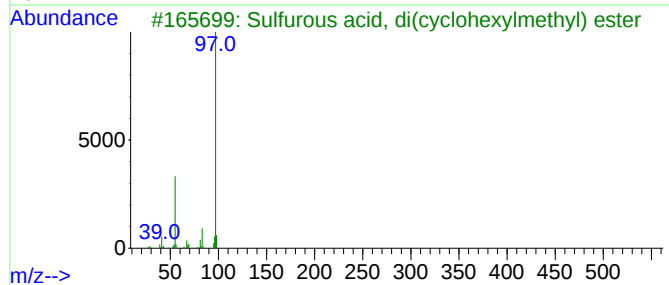
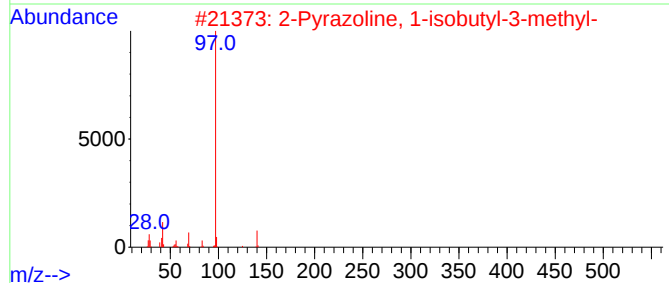
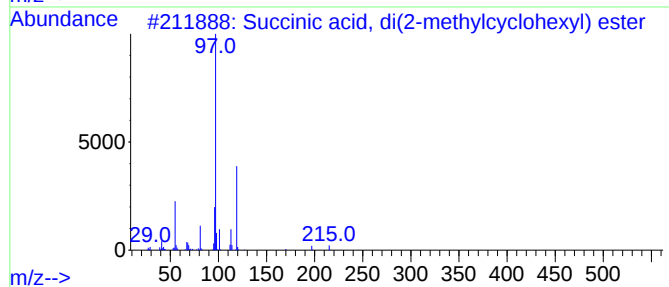
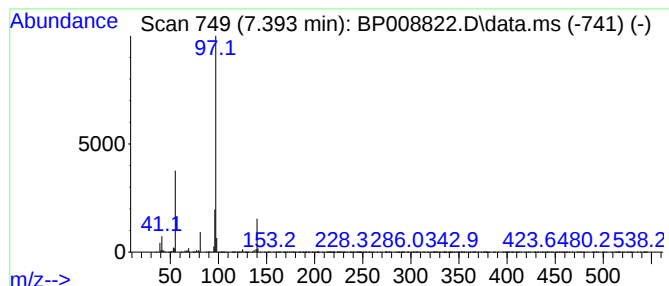
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Succinic acid, di(2-methylc... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.393	35.14 ng	26455500	1,4-Dichlorobenzene-d4	7.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Succinic acid, di(2-methylcyclohexyl) ester	310	C18H30O4	1000329-83-0	56
2			2-Pyrazoline, 1-isobutyl-3-methyl-	140	C8H16N2	026964-53-4	53
3			Sulfurous acid, di(cyclohexylmethyl) ester	274	C14H26O3S	1010309-22-7	53
4			m-Menthane, (1S,3R)-(+)-	140	C10H20	013837-66-6	43
5			Sulfurous acid, cyclohexylmethyl ester	276	C14H28O3S	1000309-21-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
WEST-BOILER-ROOM-NORTH-SIDEWALL

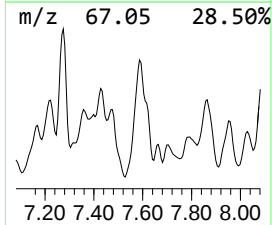
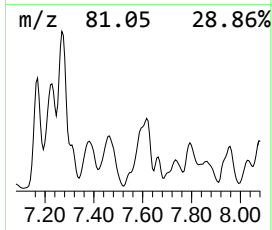
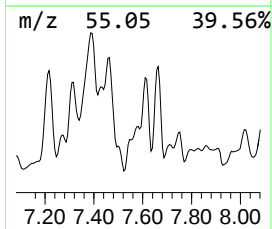
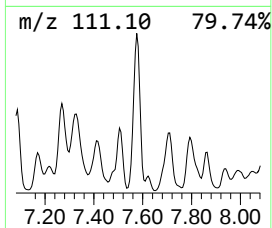
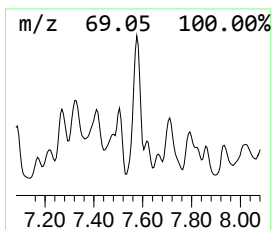
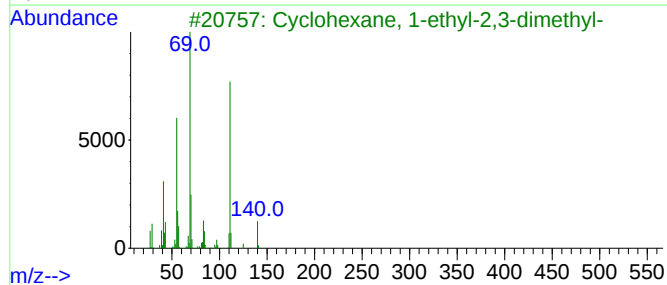
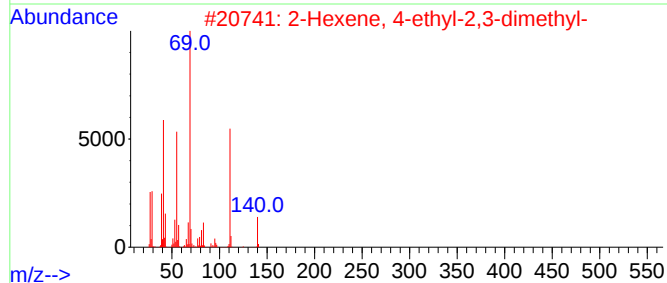
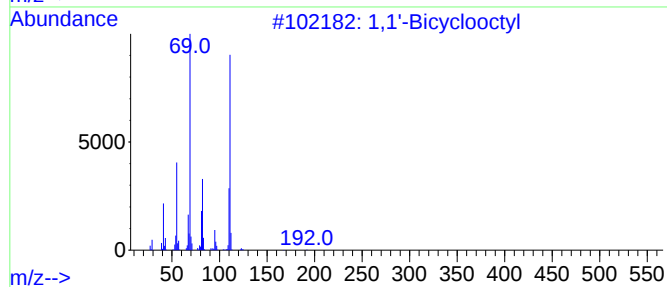
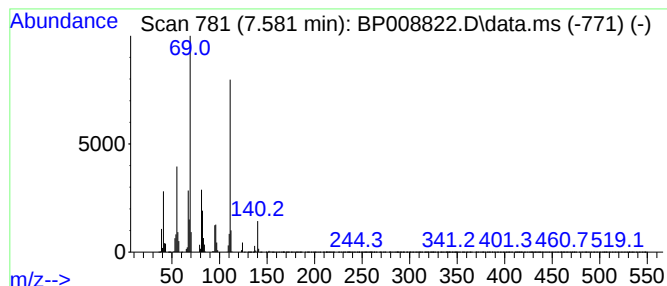
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1,1'-Bicyclooctyl Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.581	48.65 ng	36629000	1,4-Dichlorobenzene-d4	7.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Bicyclooctyl	222	C16H30	006708-17-4	59		
2	2-Hexene, 4-ethyl-2,3-dimethyl-	140	C10H20	1000149-19-7	59		
3	Cyclohexane, 1-ethyl-2,3-dimethyl-	140	C10H20	007058-05-1	58		
4	Trans-1,4-diethylcyclohexane	140	C10H20	013990-93-7	58		
5	2-(Butyliden-2-one)tetrahydrofuran	140	C8H12O2	343270-50-8	53		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
WEST-BOILER-ROOM-NORTH-SIDEWALL

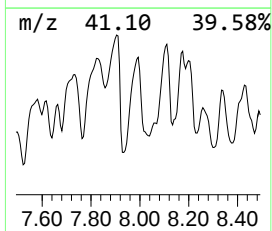
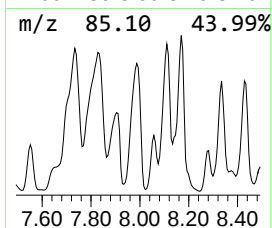
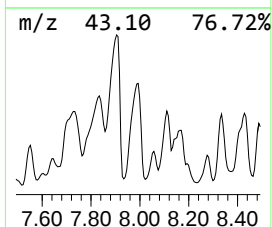
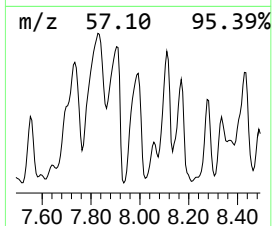
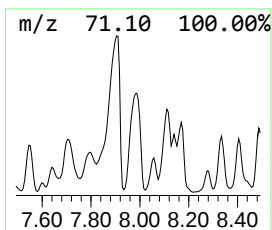
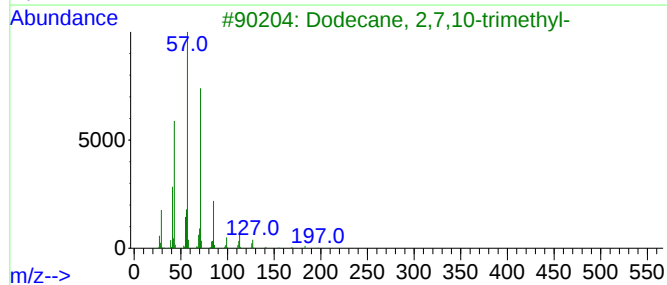
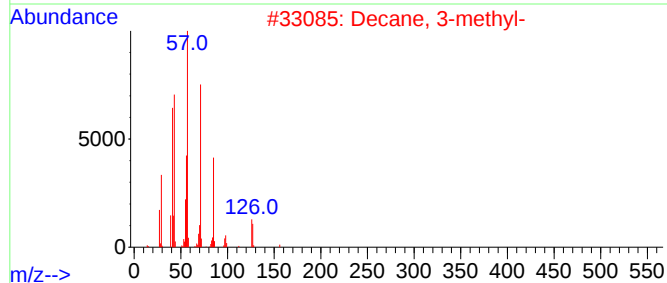
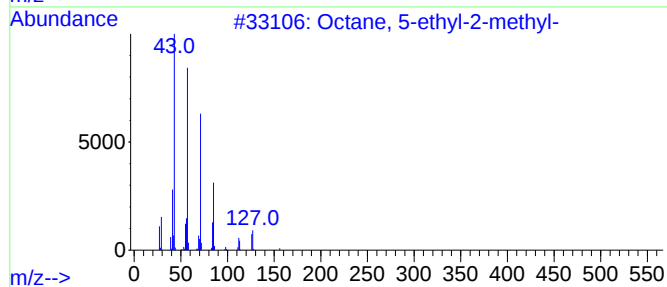
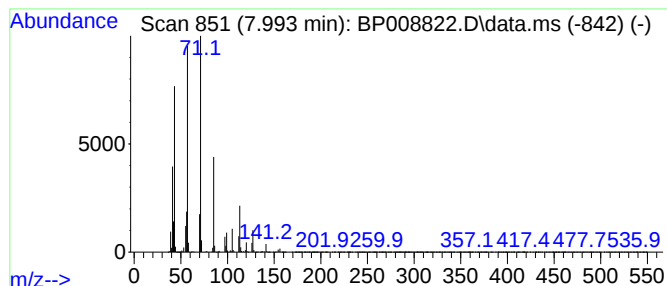
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octane, 5-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.993	50.55 ng	38058700	1,4-Dichlorobenzene-d4	7.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 5-ethyl-2-methyl-	156	C11H24	062016-18-6	74
2			Decane, 3-methyl-	156	C11H24	013151-34-3	68
3			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	59
4			Undecane, 5-methyl-	170	C12H26	001632-70-8	53
5			Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
WEST-BOILER-ROOM-NORTH-SIDEWALL

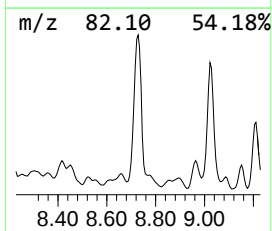
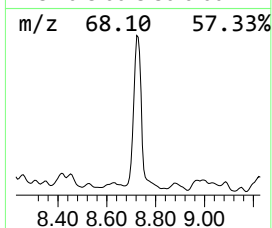
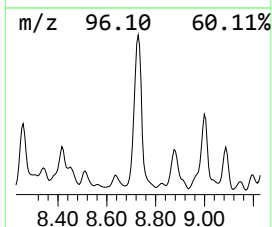
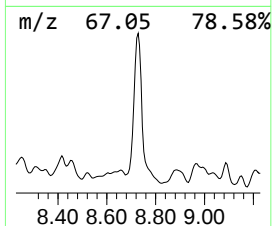
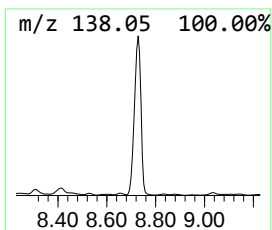
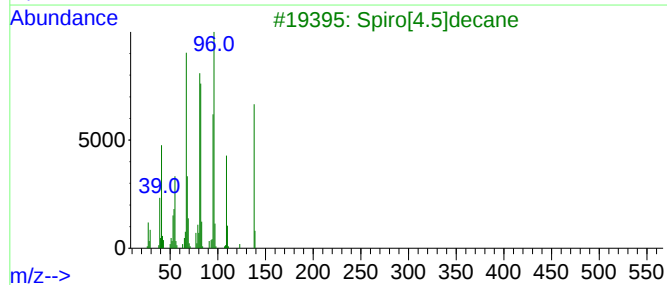
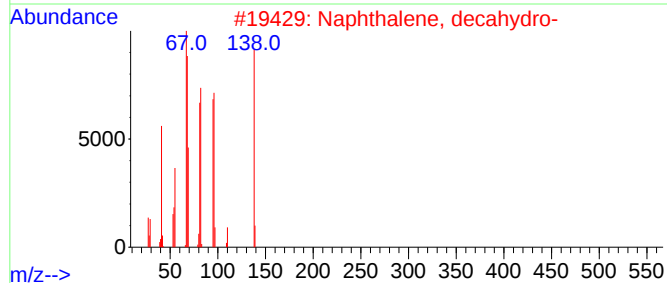
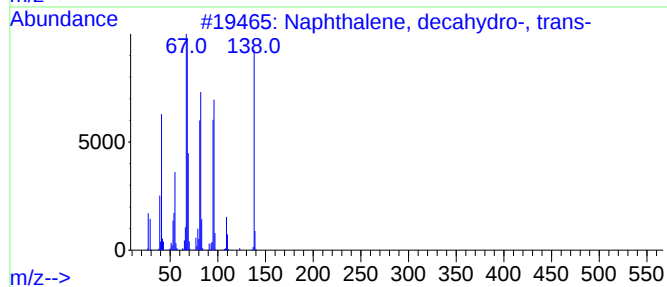
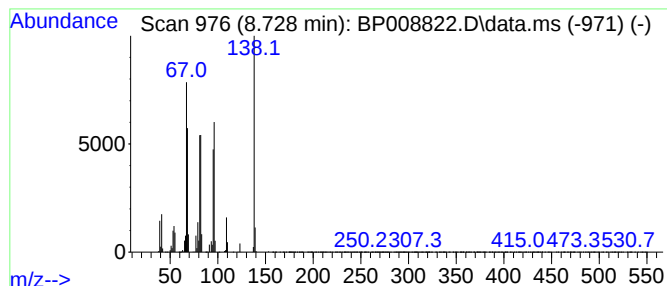
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Naphthalene, decahydro-, tr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.728	63.59 ng	47873500	1,4-Dichlorobenzene-d4	7.875

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	94
2		Naphthalene, decahydro-	138	C10H18	000091-17-8	91
3		Spiro[4.5]decane	138	C10H18	000176-63-6	83
4		Bicyclo[3.1.0]hexan-2-one, 5-(1-...	138	C9H14O	000513-20-2	64
5		Cyclopentene, 1-(3-methylbutyl)-	138	C10H18	037689-15-9	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
WEST-BOILER-ROOM-NORTH-SIDEWALL

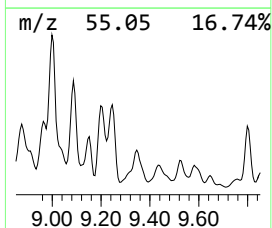
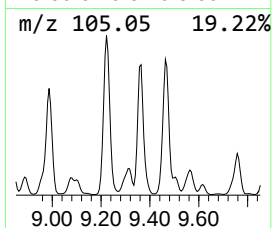
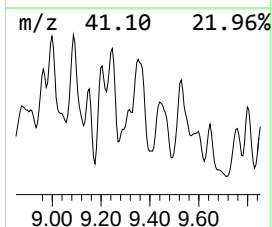
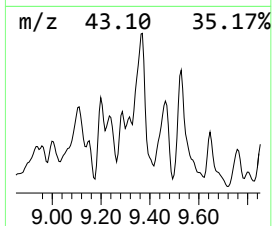
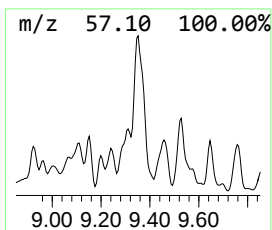
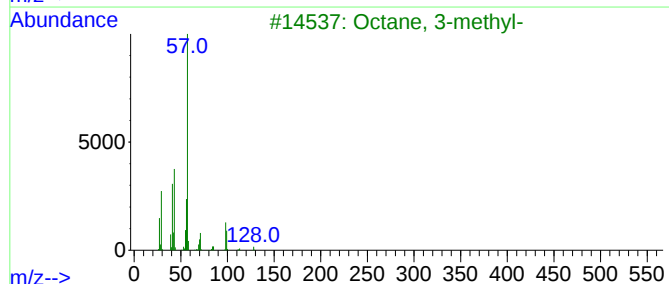
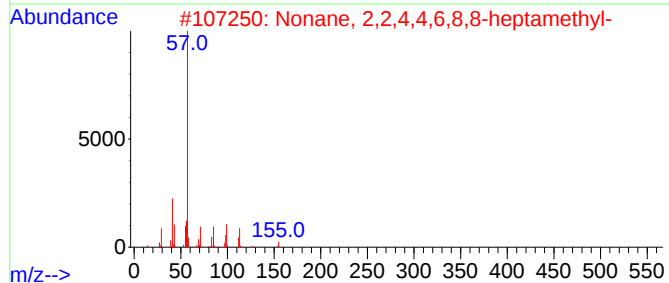
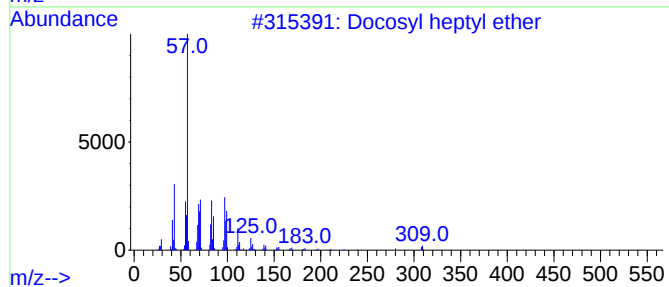
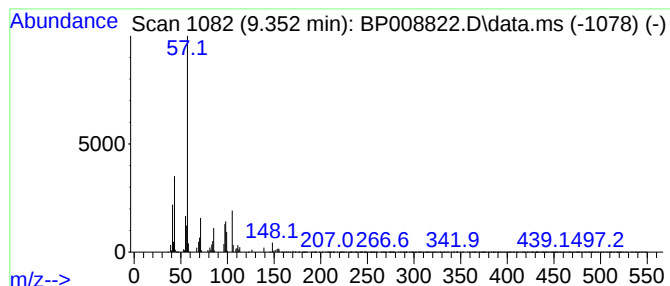
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Docosyl heptyl ether Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.352	58.80 ng	17130600	Naphthalene-d8	10.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosyl heptyl ether	424	C29H60O	1000406-39-7	50
2			Nonane, 2,2,4,4,6,8,8-heptamethyl-	226	C16H34	004390-04-9	43
3			Octane, 3-methyl-	128	C9H20	002216-33-3	38
4			Heptane, 2,3,5-trimethyl-	142	C10H22	020278-85-7	38
5			Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
WEST-BOILER-ROOM-NORTH-SIDEWALL

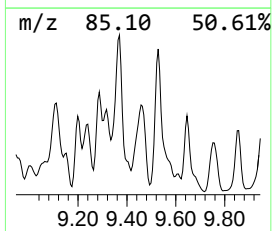
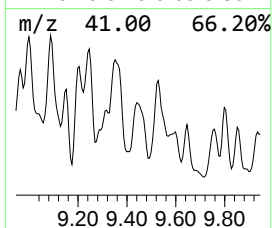
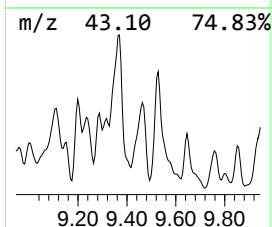
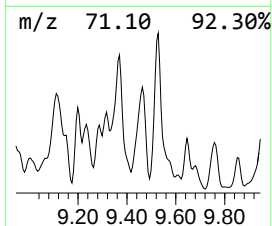
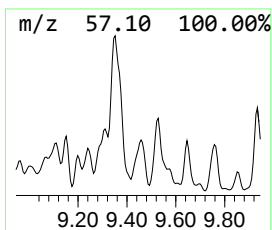
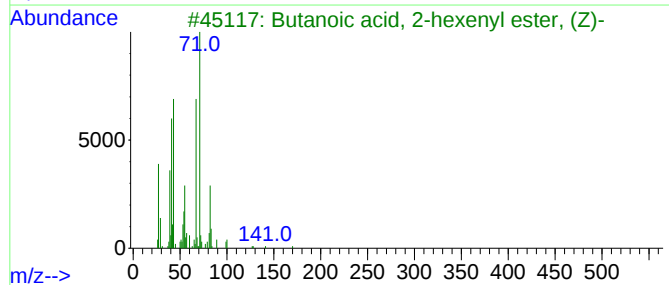
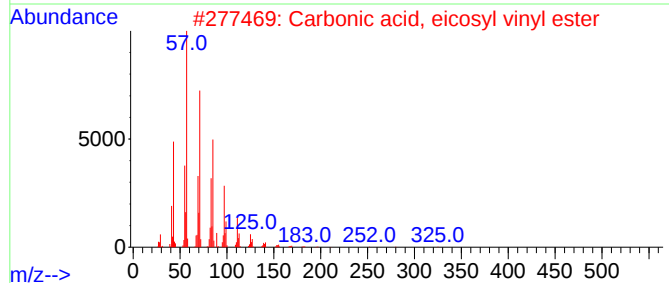
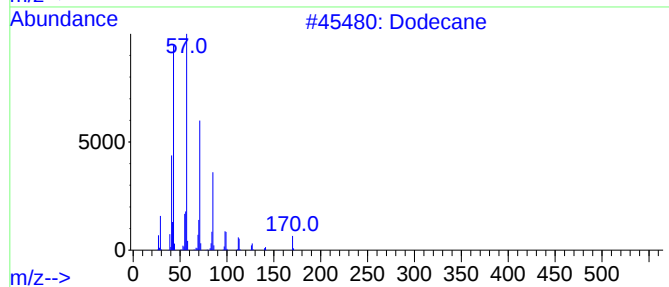
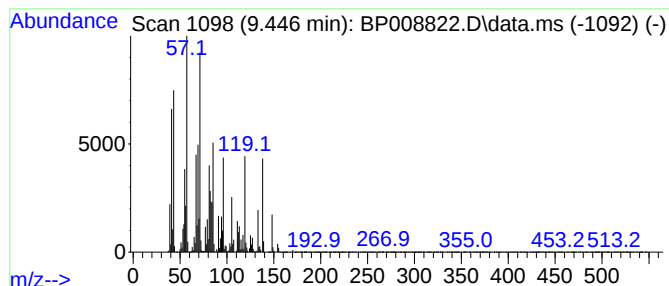
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown9.446 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.446	39.92 ng	11631700	Naphthalene-d8	10.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	38
2			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	30
3			Butanoic acid, 2-hexenyl ester, ...	170	C10H18O2	056922-77-1	30
4			1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	30
5			Tridecane	184	C13H28	000629-50-5	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
WEST-BOILER-ROOM-NORTH-SIDEWALL

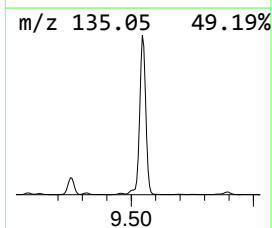
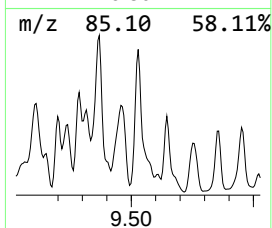
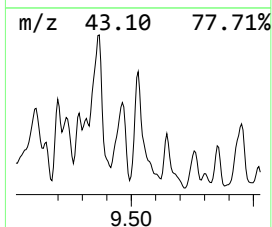
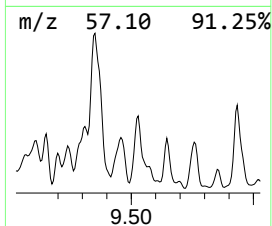
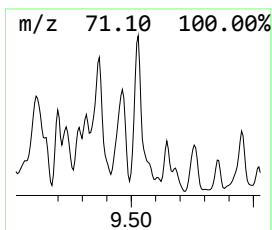
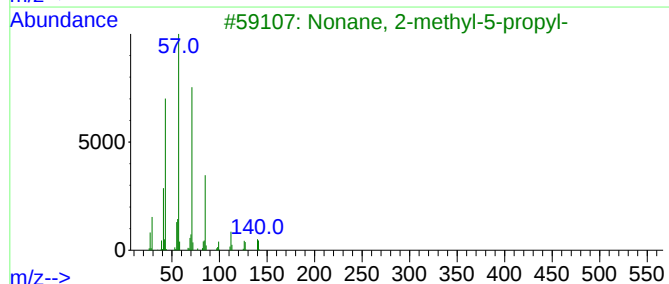
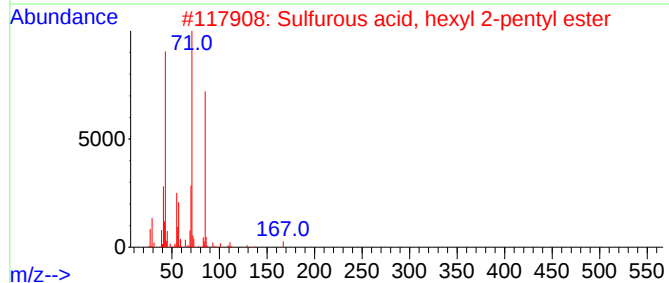
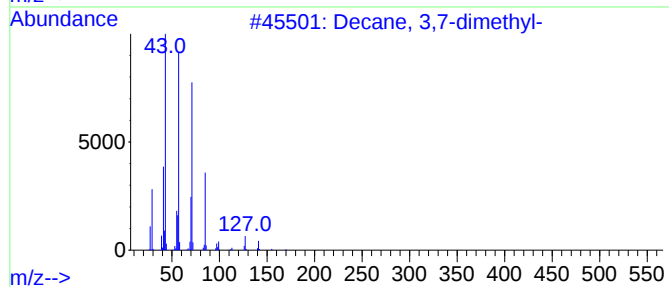
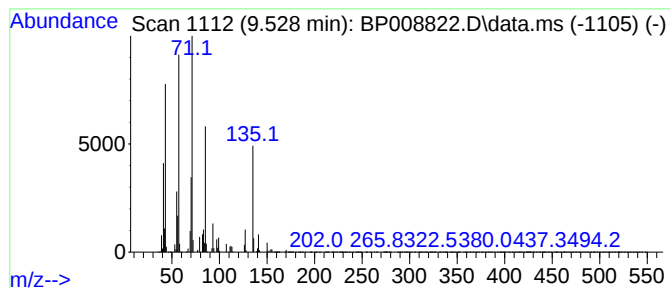
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Decane, 3,7-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.528	34.22 ng	9969360	Naphthalene-d8	10.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3,7-dimethyl-	170	C12H26	017312-54-8	89
2			Sulfurous acid, hexyl 2-pentyl e...	236	C11H24O3S	1000309-15-6	50
3			Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	50
4			Sulfurous acid, decyl pentyl ester	292	C15H32O3S	1000309-14-3	47
5			Sulfurous acid, decyl 2-pentyl e...	292	C15H32O3S	1000309-15-9	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
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Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
WEST-BOILER-ROOM-NORTH-SIDEWALL

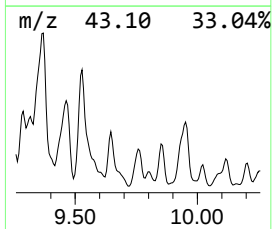
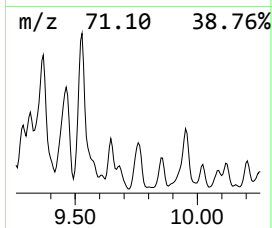
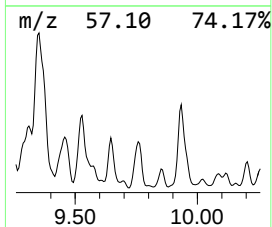
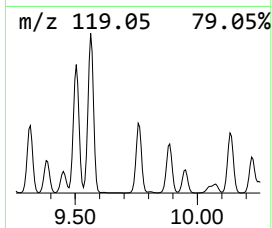
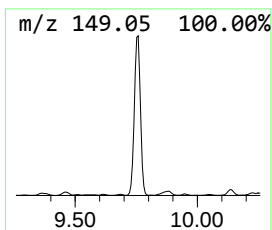
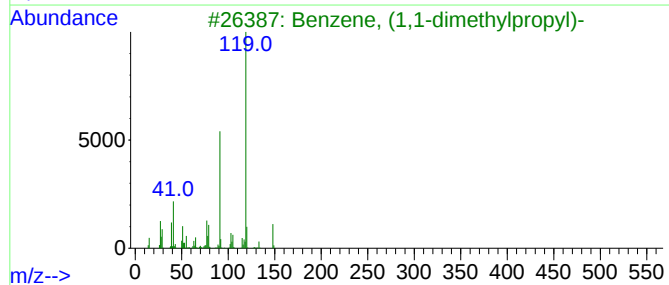
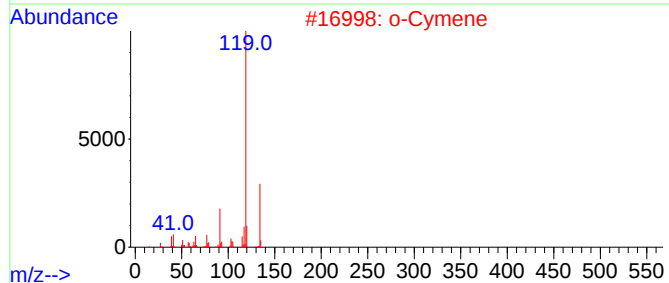
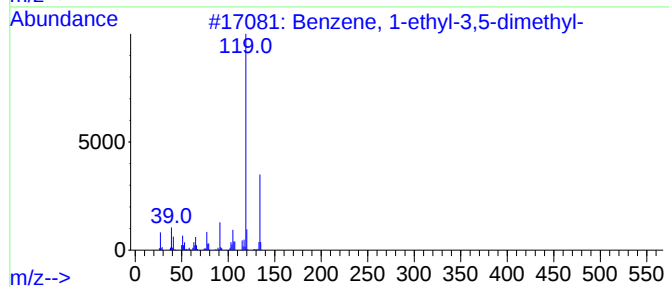
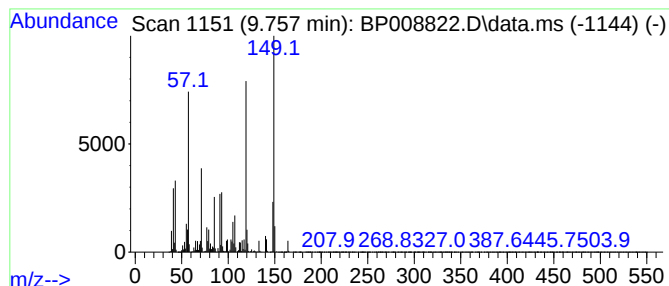
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown9.757 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.757	38.11 ng	11104600	Naphthalene-d8	10.663

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	35
2		o-Cymene	134	C10H14	000527-84-4	35
3		Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	25
4		m-Tolyl isothiocyanate	149	C8H7NS	000621-30-7	22
5		3,5-Dimethylphenylacetic acid	164	C10H12O2	042288-46-0	20



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
WEST-BOILER-ROOM-NORTH-SIDEWALL

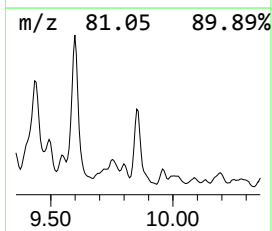
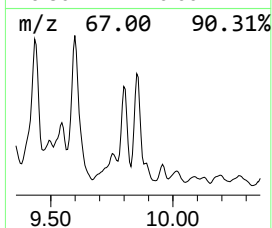
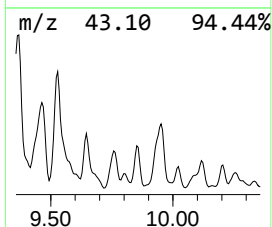
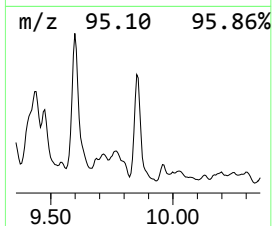
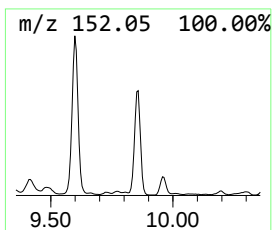
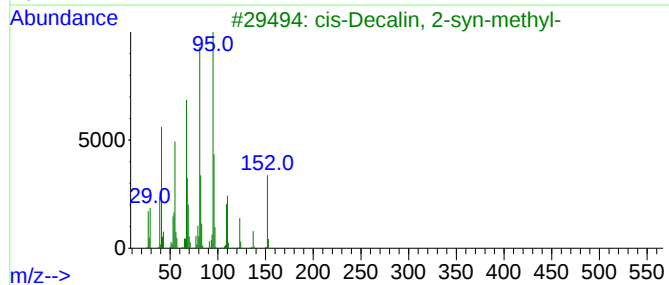
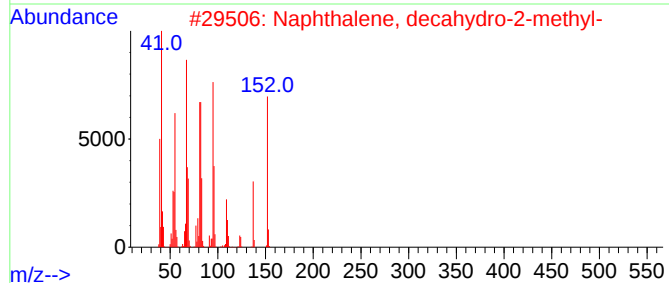
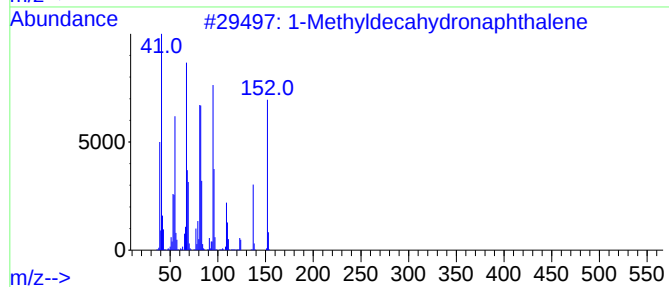
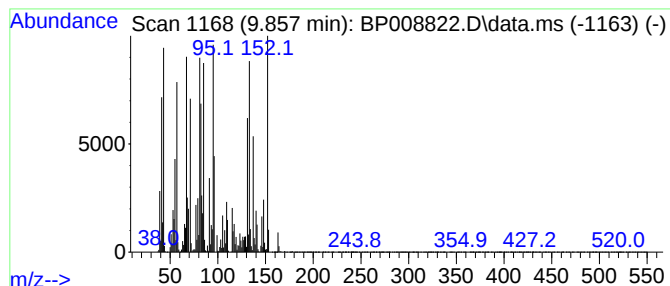
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 1-Methyldecahydronaphthalene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.857	39.34 ng	11461800	Naphthalene-d8	10.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Methyldecahydronaphthalene	152	C11H20	002958-75-0	83
2			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	83
3			cis-Decalin, 2-syn-methyl-	152	C11H20	1000155-85-6	49
4			Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	45
5			Bicyclo[4.1.0]heptan-3-one, 4,7,...	152	C10H16O	004176-01-6	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
WEST-BOILER-ROOM-NORTH-SIDEWALL

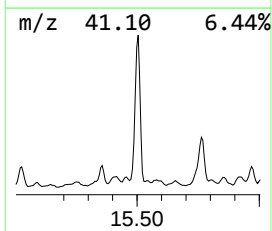
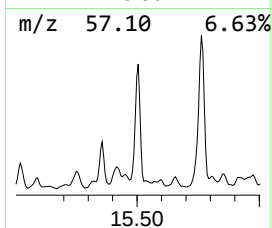
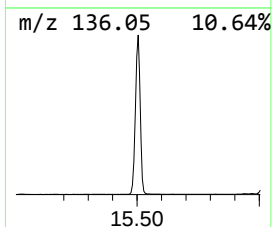
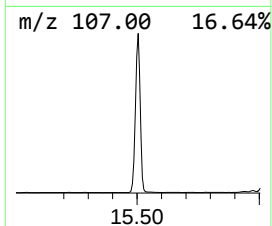
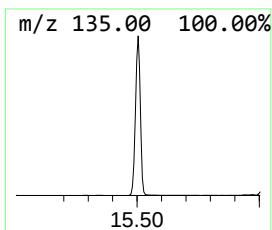
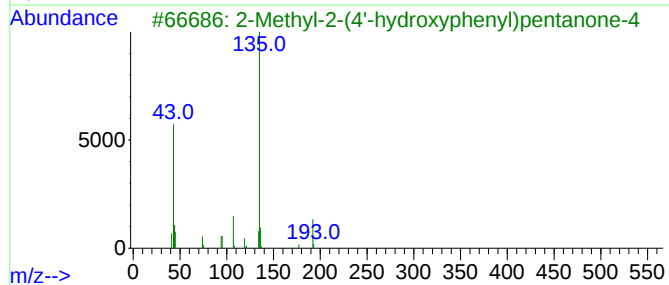
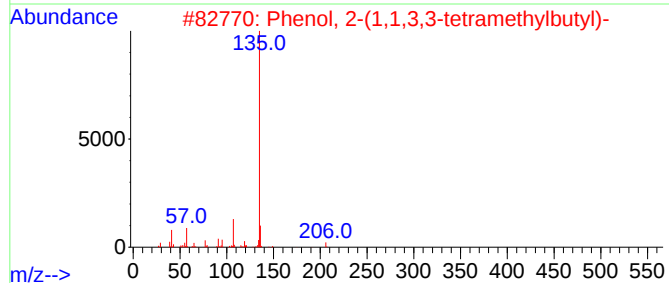
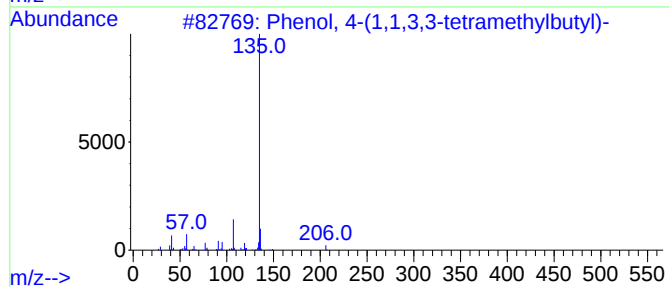
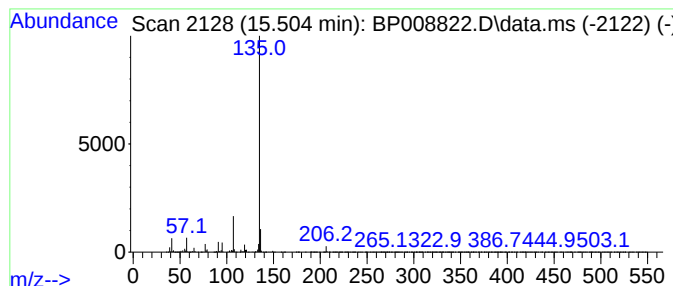
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.504	58.02 ng	14755300	Acenaphthene-d10	14.493

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	96
2		Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	95
3		2-Methyl-2-(4'-hydroxyphenyl)pen...	192	C12H16O2	070205-18-4	78
4		4-tert-Butylphenyl acetate	192	C12H16O2	003056-64-2	78
5		Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

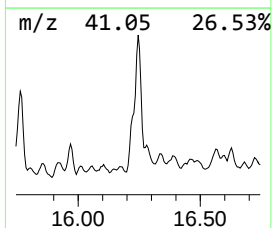
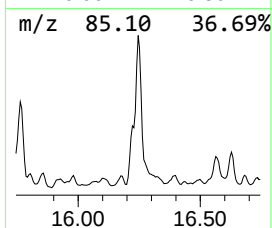
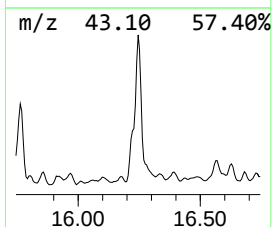
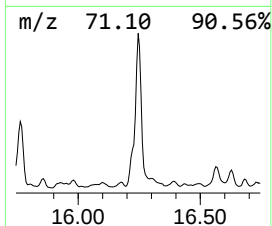
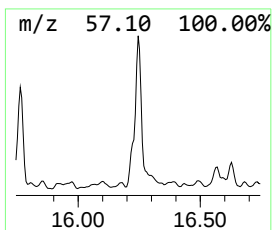
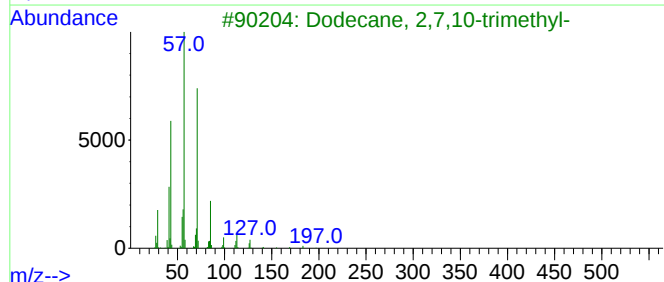
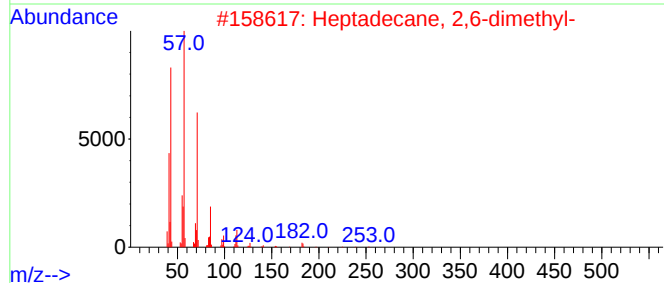
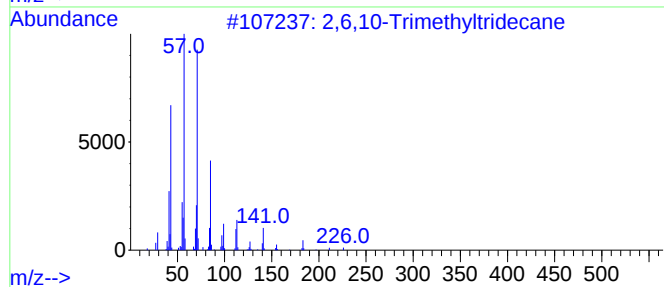
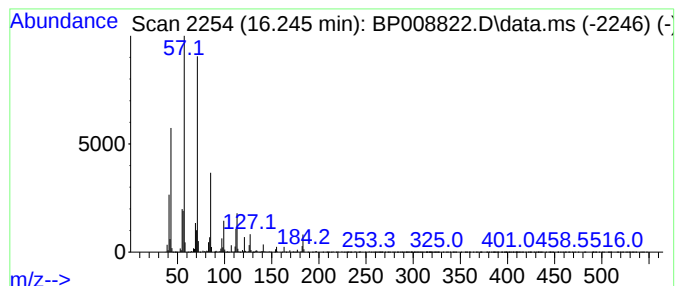
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BNA_P
ClientSampleId :
WEST-BOILER-ROOM-NORTH-SIDEWALL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 2,6,10-Trimethyltridecane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD			R.T.
16.245	33.71 ng	7576800	Phenanthrene-d10			17.251
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2,6,10-Trimethyltridecane	226	C16H34	003891-99-4	90
2		Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	81
3		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	80
4		Pentadecane, 7-methyl-	226	C16H34	006165-40-8	80
5		Undecane, 6-ethyl-	184	C13H28	017312-60-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
Data File : BP008822.D
Acq On : 17 Jan 2022 14:14
Operator : CG/JU
Sample : N1143-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
WEST-BOILER-ROOM-NORTH-SIDEWALL

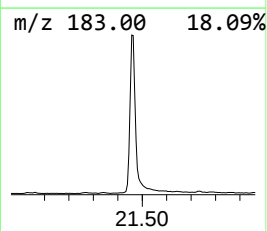
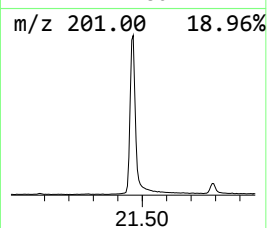
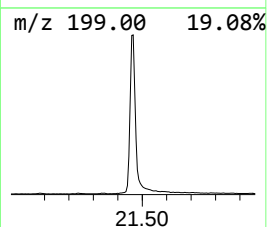
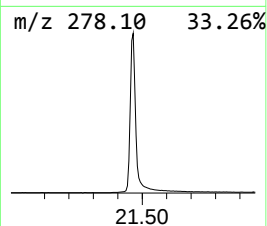
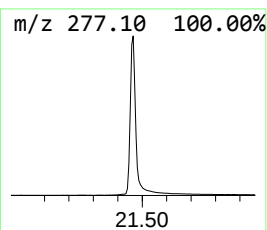
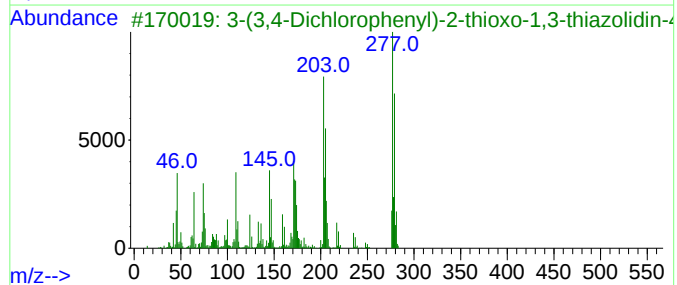
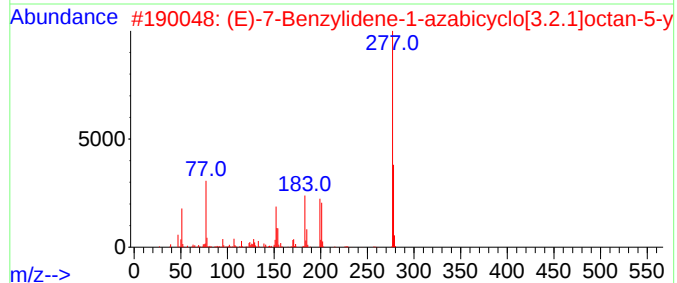
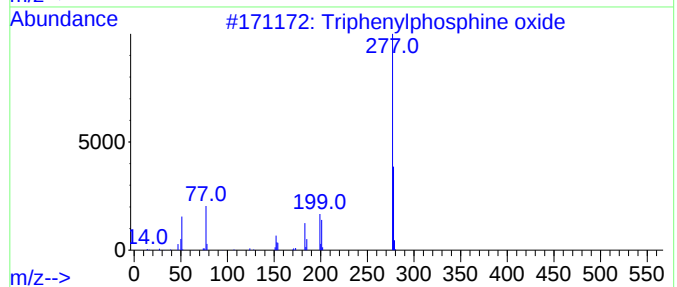
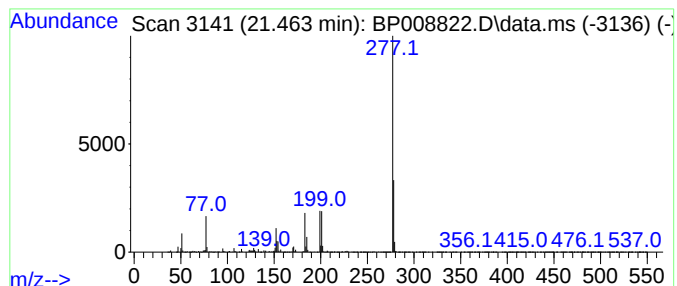
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Triphenylphosphine oxide Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.463	42.19 ng	7836670	Chrysene-d12	21.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	94
2			(E)-7-Benzylidene-1-azabicyclo[3...]	293	C15H19NO3S	1000483-83-4	91
3			3-(3,4-Dichlorophenyl)-2-thioxo-...	277	C9H5Cl2NOS2	017046-34-3	37
4			5-Methyl-8-phenylfuro[3',2':5,6]...	277	C16H11N3O2	1000460-07-9	28
5			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011722\
 Data File : BP008822.D
 Acq On : 17 Jan 2022 14:14
 Operator : CG/JU
 Sample : N1143-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 WEST-BOILER-ROOM-NORTH-SIDEWALL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.152	6.152	113.2	ng	85213500	1	7.875	15057700	20.0
Octane, 3,5-dim...	6.281	60.9	ng	45879700	1	7.875	15057700	20.0
unknown6.375	6.375	62.4	ng	46945900	1	7.875	15057700	20.0
Octane, 2,6-dim...	6.463	67.8	ng	51048500	1	7.875	15057700	20.0
Cyclohexane, pr...	6.534	42.0	ng	31636600	1	7.875	15057700	20.0
Cyclohexane, 1,...	6.646	34.2	ng	25740100	1	7.875	15057700	20.0
unknown6.787	6.787	70.0	ng	52725200	1	7.875	15057700	20.0
Nonane, 4-methyl-	6.905	106.9	ng	80506700	1	7.875	15057700	20.0
Succinic acid, ...	7.393	35.1	ng	26455500	1	7.875	15057700	20.0
1,1'-Bicyclooctyl	7.581	48.6	ng	36629000	1	7.875	15057700	20.0
Octane, 5-ethyl...	7.993	50.5	ng	38058700	1	7.875	15057700	20.0
Naphthalene, de...	8.728	63.6	ng	47873500	1	7.875	15057700	20.0
Docosyl heptyl ...	9.352	58.8	ng	17130600	2	10.663	5827180	20.0
unknown9.446	9.446	39.9	ng	11631700	2	10.663	5827180	20.0
Decane, 3,7-dim...	9.528	34.2	ng	9969360	2	10.663	5827180	20.0
unknown9.757	9.757	38.1	ng	11104600	2	10.663	5827180	20.0
1-Methyldecahyd...	9.857	39.3	ng	11461800	2	10.663	5827180	20.0
Phenol, 4-(1,1,...	15.504	58.0	ng	14755300	3	14.493	5086560	20.0
2,6,10-Trimethy...	16.245	33.7	ng	7576800	4	17.251	4495760	20.0
Triphenylphosph...	21.463	42.2	ng	7836670	5	21.339	3715200	20.0